Energy Laboratories Inc

ANALYTICAL RUN Summary

28-Feb-22

Run ID VOA5975C.I_220104A

Run Start Date: 1/4/2022
Analyst: Melissa Chavez
Ical:
Column ID:
Comments:

Instrument ID	Description
Bal #22	Balance

Std ID	Std Name	Std Amount	Std Units	Samp Amount	Samp Units	SampType	Expiration Date
VOCF3473	Calibration Surrogates		ul	42	ml	CAL	3/14/2022
VOCF3517	Internal Standard / Surrogates (INT/SURR)	8.4	ul	42	ml	MBLK, ICV (	12/31/2022
VOCF3529B	2nd Source MtBE	1.05	ul	42	ml	ICV	1/29/2022
VOCF3546A	Liquids		ul	42	ml	CAL	1/13/2022
VOCF3549	2nd Source Ketones	1.05	ul	42	ml	ICV	1/15/2022
VOCF3550	Ketones		ul	42	ml	CAL	1/16/2022
VOCF3558B	2nd Source Liquids	1.05	ul	42	ml	ICV	2/27/2022
VOCF3559A	MtBE		ul	42	ml	CAL	1/27/2022
VOCF3562A	Gases		ul	42	ml	CAL	1/10/2022
VOCF3563	Internals	8.4	ul	42	ml	CAL	7/3/2022
VOCF3566A	2nd Source Gases	1.05	ul	42	ml	ICV	1/11/2022

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970784	04JAN08_D_TU	VOC-8260-BFB	TUNE	)A5975C\VG010	1/4/2022 2:38:00	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
173, % of mass 174	A	%	0	0		100	0	0	0	0	0	0%	0	1.99	0%	
174, % of mass 95	A	%	95.2	95.2		100	0	0	0	0	0	95%	50	99.99	0%	
175, % of mass 174	A	%	6.6	6.6		100	0	0	0	0	0	7%	5	9	0%	
176, % of mass 174	A	%	95.7	95.7		100	0	0	0	0	0	96%	95	101	0%	
177, % of mass 176	A	%	6.7	6.7		100	0	0	0	0	0	7%	5	9	0%	
50, % of mass 95	A	%	21.2	21.2		100	0	0	0	0	0	21%	15	40	0%	
75, % of mass 95	A	%	51	51		100	0	0	0	0	0	51%	30	60	0%	
95, Base Peak	A	%	100	100		100	0	0	0	0	0	100%	0	100	0%	
96, % of mass 95	A	%	5.4	5.4		100	0	0	0	0	0	5%	5	9	0%	

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Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970898	MBLK010422_	VOC-8260-W-Q	MBLK	)A5975C\VG010	1/4/2022 3:05:37	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	0.5	500	0%	0	0	0%	
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	0.5	500	0%	0	0	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	0.5	500	0%	0	0	0%	
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	0.5	500	0%	0	0	0%	
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	0.5	500	0%	0	0	0%	
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	0.5	500	0%	0	0	0%	
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	0.5	500	0%	0	0	0%	
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	0.5	500	0%	0	0	0%	
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	0.5	500	0%	0	0	0%	
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	0.5	500	0%	0	0	0%	
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	0.5	500	0%	0	0	0%	
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	0.5	500	0%	0	0	0%	
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	0.5	500	0%	0	0	0%	
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	0.5	500	0%	0	0	0%	
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	0.5	500	0%	0	0	0%	
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	0.5	500	0%	0	0	0%	
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	0.5	500	0%	0	0	0%	
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	0.5	500	0%	0	0	0%	
Benzene	A	ug/L	0.12327	0		0	0	0	0.0914	0.5	500	0%	0	0	0%	
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	0.5	500	0%	0	0	0%	
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	0.5	500	0%	0	0	0%	
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	0.5	500	0%	0	0	0%	
Bromoform	A	ug/L	0	0		0	0	0	0.119	0.5	500	0%	0	0	0%	
Bromomethane	A	ug/L	0	0		0	0	0	0.253	0.5	500	0%	0	0	0%	
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	0.5	500	0%	0	0	0%	
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	0.5	500	0%	0	0	0%	
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	0.5	500	0%	0	0	0%	
Chloroethane	A	ug/L	0	0		0	0	0	0.169	0.5	500	0%	0	0	0%	
Chloroform	A	ug/L	0	0		0	0	0	0.0789	0.5	500	0%	0	0	0%	
Chloromethane	A	ug/L	0	0		0	0	0	0.162	0.5	500	0%	0	0	0%	
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	0.5	500	0%	0	0	0%	
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	0.5	500	0%	0	0	0%	
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	0.5	500	0%	0	0	0%	
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	0.5	500	0%	0	0	0%	
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	0.5	500	0%	0	0	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970898	MBLK010422_	VOC-8260-W-Q	MBLK	)A5975C\VG010	1/4/2022 3:05:37	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	0.5	1000	0%	0	0	0%	
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	10	5000	0%	0	0	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	0.5	500	0%	0	0	0%	
Methylene chloride	A	ug/L	1.44235	0		0	0	0	0.338	0.5	500	0%	0	0	0%	
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	0.5	500	0%	0	0	0%	
Styrene	A	ug/L	0	0		0	0	0	0.067	0.5	500	0%	0	0	0%	
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	0.5	500	0%	0	0	0%	
Toluene	A	ug/L	0	0		0	0	0	0.0679	0.5	500	0%	0	0	0%	
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	0.5	500	0%	0	0	0%	
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	0.5	500	0%	0	0	0%	
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	0.5	500	0%	0	0	0%	
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	0.5	500	0%	0	0	0%	
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	0.5	500	0%	0	0	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	0.5	1500	0%	0	0	0%	
1,2-Dichloroethane-d4	S	ug/L	279.39635	11.175854		10	0	0	0.229	0.5	500	112%	70	130	0%	
Dibromofluoromethane	S	ug/L	278.46353	11.1385412		10	0	0	0.129	0.5	500	111%	77	126	0%	
p-Bromofluorobenzene	S	ug/L	267.28149	10.6912596		10	0	0	0.149	0.5	500	107%	76	127	0%	
Toluene-d8	S	ug/L	265.34358	10.6137432		10	0	0	0.23	0.5	500	106%	79	122	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970899	ICAL010422_1	VOC-8260-W-Q	CAL1	)A5975C\VG010	1/4/2022 3:33:04	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	2.73073	0.1092292		0.1	0	0	0.0746	0.5	500	109%	50	150	0%	
1,2-Dichloroethane	A	ug/L	2.90899	0.1163596		0.1	0	0	0.116	0.5	500	116%	50	150	0%	
1,3-Dichlorobenzene	A	ug/L	2.6327	0.105308		0.1	0	0	0.0803	0.5	500	105%	50	150	0%	
1,4-Dichlorobenzene	A	ug/L	2.76134	0.1104536		0.1	0	0	0.0858	0.5	500	110%	50	150	0%	
Benzene	A	ug/L	2.73933	0.1095732		0.1	0	0	0.0914	0.5	500	110%	50	150	0%	
Chloroform	A	ug/L	2.89464	0.1157856		0.1	0	0	0.0789	0.5	500	116%	50	150	0%	
Ethylbenzene	A	ug/L	2.53666	0.1014664		0.1	0	0	0.0836	0.5	500	101%	50	150	0%	
m+p-Xylenes	A	ug/L	5.07121	0.2028484		0.2	0	0	0.15	0.5	1000	101%	50	150	0%	
Styrene	A	ug/L	2.16254	0.0865016		0.1	0	0	0.067	0.5	500	87%	50	150	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970899	ICAL010422_1	VOC-8260-W-Q	CAL1	)A5975C\VG010	1/4/2022 3:33:04	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Tetrachloroethene	A	ug/L	2.67723	0.1070892		0.1	0	0	0.0671	0.5	500	107%	50	150	0%	
Toluene	A	ug/L	2.6145	0.10458		0.1	0	0	0.0679	0.5	500	105%	50	150	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970901	ICAL010422_2	VOC-8260-W-Q	CAL2	)A5975C\VG010	1/4/2022 4:00:35	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	12.82253	0.5129012		0.5	0	0	0.101	0.5	500	103%	50	150	0%	
1,1,1-Trichloroethane	A	ug/L	12.18907	0.4875628		0.5	0	0	0.131	0.5	500	98%	50	150	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	12.84375	0.51375		0.5	0	0	0.0872	0.5	500	103%	50	150	0%	
1,1,2-Trichloroethane	A	ug/L	13.23404	0.5293616		0.5	0	0	0.108	0.5	500	106%	50	150	0%	
1,1-Dichloroethane	A	ug/L	12.06522	0.4826088		0.5	0	0	0.135	0.5	500	97%	50	150	0%	
1,1-Dichloroethene	A	ug/L	11.90807	0.4763228		0.5	0	0	0.141	0.5	500	95%	50	150	0%	
1,1-Dichloropropene	A	ug/L	11.33971	0.4535884		0.5	0	0	0.083	0.5	500	91%	50	150	0%	
1,2,3-Trichloropropane	A	ug/L	13.70838	0.5483352		0.5	0	0	0.235	0.5	500	110%	50	150	0%	
1,2-Dibromoethane	A	ug/L	12.86397	0.5145588		0.5	0	0	0.0916	0.5	500	103%	50	150	0%	
1,2-Dichlorobenzene	A	ug/L	12.14234	0.4856936		0.5	0	0	0.0746	0.5	500	97%	70	130	0%	
1,2-Dichloroethane	A	ug/L	12.39059	0.4956236		0.5	0	0	0.116	0.5	500	99%	70	130	0%	
1,2-Dichloropropane	A	ug/L	12.0602	0.482408		0.5	0	0	0.0847	0.5	500	96%	50	150	0%	
1,3-Dichlorobenzene	A	ug/L	11.84726	0.4738904		0.5	0	0	0.0803	0.5	500	95%	70	130	0%	
1,3-Dichloropropane	A	ug/L	11.85262	0.4741048		0.5	0	0	0.0791	0.5	500	95%	50	150	0%	
1,4-Dichlorobenzene	A	ug/L	11.96618	0.4786472		0.5	0	0	0.0858	0.5	500	96%	70	130	0%	
2,2-Dichloropropane	A	ug/L	12.48201	0.4992804		0.5	0	0	0.186	0.5	500	100%	50	150	0%	
2-Chlorotoluene	A	ug/L	11.19768	0.4479072		0.5	0	0	0.0876	0.5	500	90%	50	150	0%	
4-Chlorotoluene	A	ug/L	11.22327	0.4489308		0.5	0	0	0.0728	0.5	500	90%	50	150	0%	
Benzene	A	ug/L	12.18007	0.4872028		0.5	0	0	0.0914	0.5	500	97%	70	130	0%	
Bromobenzene	A	ug/L	12.331	0.49324		0.5	0	0	0.0831	0.5	500	99%	50	150	0%	
Bromochloromethane	A	ug/L	12.9568	0.518272		0.5	0	0	0.141	0.5	500	104%	50	150	0%	
Bromodichloromethane	A	ug/L	12.60141	0.5040564		0.5	0	0	0.12	0.5	500	101%	50	150	0%	
Bromoform	A	ug/L	11.78598	0.4714392		0.5	0	0	0.119	0.5	500	94%	50	150	0%	
Bromomethane	A	ug/L	12.04638	0.4818552		0.5	0	0	0.253	0.5	500	96%	50	150	0%	
Carbon tetrachloride	A	ug/L	12.2545	0.49018		0.5	0	0	0.143	0.5	500	98%	50	150	0%	
Chlorobenzene	A	ug/L	12.52043	0.5008172		0.5	0	0	0.0914	0.5	500	100%	50	150	0%	
Chlorodibromomethane	A	ug/L	12.83929	0.5135716		0.5	0	0	0.0841	0.5	500	103%	50	150	0%	
Chloroethane	A	ug/L	14.86697	0.5946788		0.5	0	0	0.169	0.5	500	119%	50	150	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970901	ICAL010422_2	VOC-8260-W-Q	CAL2	)A5975C\VG010	1/4/2022 4:00:35	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Chloroform	A	ug/L	13.06683	0.5226732		0.5	0	0	0.0789	0.5	500	105%	70	130	0%	
Chloromethane	A	ug/L	13.86612	0.5546448		0.5	0	0	0.162	0.5	500	111%	50	150	0%	
cis-1,2-Dichloroethene	A	ug/L	12.56593	0.5026372		0.5	0	0	0.108	0.5	500	101%	50	150	0%	
cis-1,3-Dichloropropene	A	ug/L	12.07376	0.4829504		0.5	0	0	0.073	0.5	500	97%	50	150	0%	
Dibromomethane	A	ug/L	14.06189	0.5624756		0.5	0	0	0.147	0.5	500	112%	50	150	0%	
Dichlorodifluoromethane	A	ug/L	12.06625	0.48265		0.5	0	0	0.175	0.5	500	97%	50	150	0%	
Ethylbenzene	A	ug/L	11.04112	0.4416448		0.5	0	0	0.0836	0.5	500	88%	70	130	0%	
m+p-Xylenes	A	ug/L	22.14096	0.8856384		1	0	0	0.15	0.5	1000	89%	70	130	0%	
Methyl ethyl ketone	A	ug/L	122.052	4.88208		5	0	0	1.77	10	5000	98%	50	150	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	12.32545	0.493018		0.5	0	0	0.101	0.5	500	99%	50	150	0%	
Methylene chloride	A	ug/L	15.62358	0.6249432		0.5	0	0	0.338	0.5	500	125%	50	150	0%	
o-Xylene	A	ug/L	10.66119	0.4264476		0.5	0	0	0.0604	0.5	500	85%	50	150	0%	
Styrene	A	ug/L	11.49684	0.4598736		0.5	0	0	0.067	0.5	500	92%	70	130	0%	
Tetrachloroethene	A	ug/L	11.73024	0.4692096		0.5	0	0	0.0671	0.5	500	94%	70	130	0%	
Toluene	A	ug/L	11.28985	0.451594		0.5	0	0	0.0679	0.5	500	90%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	12.50224	0.5000896		0.5	0	0	0.125	0.5	500	100%	50	150	0%	
trans-1,3-Dichloropropene	A	ug/L	11.75888	0.4703552		0.5	0	0	0.0846	0.5	500	94%	50	150	0%	
Trichloroethene	A	ug/L	11.67527	0.4670108		0.5	0	0	0.0993	0.5	500	93%	50	150	0%	
Trichlorofluoromethane	A	ug/L	11.36372	0.4545488		0.5	0	0	0.134	0.5	500	91%	50	150	0%	
Vinyl chloride	A	ug/L	12.54456	0.5017824		0.5	0	0	0.153	0.5	500	100%	50	150	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	50	150	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	50	150	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	50	150	0%	
Xylenes, Total	M	ug/L	32.80215	1.312086		1.5	0	0	0.0604	0.5	1500	87%	50	150	0%	
1,2-Dichloroethane-d4	S	ug/L	12.66005	0.506402		0.5	0	0	0.229	0.5	500	101%	50	150	0%	
Dibromofluoromethane	S	ug/L	12.59997	0.5039988		0.5	0	0	0.129	0.5	500	101%	50	150	0%	
p-Bromofluorobenzene	S	ug/L	11.33932	0.4535728		0.5	0	0	0.149	0.5	500	91%	50	150	0%	
Toluene-d8	S	ug/L	11.30891	0.4523564		0.5	0	0	0.23	0.5	500	90%	50	150	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970902	ICAL010422_3	VOC-8260-W-Q	CAL3	)A5975C\VG010	1/4/2022 4:28:05	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970902	ICAL010422_3	VOC-8260-W-Q	CAL3	)A5975C\VG010	1/4/2022 4:28:05	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	24.15093	0.9660372		1	0	0	0.101	0.5	500	97%	70	130	0%	
1,1,1-Trichloroethane	A	ug/L	25.18087	1.0072348		1	0	0	0.131	0.5	500	101%	70	130	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	27.78828	1.1115312		1	0	0	0.0872	0.5	500	111%	70	130	0%	
1,1,2-Trichloroethane	A	ug/L	25.84	1.0336		1	0	0	0.108	0.5	500	103%	70	130	0%	
1,1-Dichloroethane	A	ug/L	25.68346	1.0273384		1	0	0	0.135	0.5	500	103%	70	130	0%	
1,1-Dichloroethene	A	ug/L	25.88489	1.0353956		1	0	0	0.141	0.5	500	104%	70	130	0%	
1,1-Dichloropropene	A	ug/L	24.36174	0.9744696		1	0	0	0.083	0.5	500	97%	70	130	0%	
1,2,3-Trichloropropane	A	ug/L	26.71444	1.0685776		1	0	0	0.235	0.5	500	107%	70	130	0%	
1,2-Dibromoethane	A	ug/L	24.36006	0.9744024		1	0	0	0.0916	0.5	500	97%	70	130	0%	
1,2-Dichlorobenzene	A	ug/L	24.94023	0.9976092		1	0	0	0.0746	0.5	500	100%	70	130	0%	
1,2-Dichloroethane	A	ug/L	23.46155	0.938462		1	0	0	0.116	0.5	500	94%	70	130	0%	
1,2-Dichloropropane	A	ug/L	25.11474	1.0045896		1	0	0	0.0847	0.5	500	100%	70	130	0%	
1,3-Dichlorobenzene	A	ug/L	25.77252	1.0309008		1	0	0	0.0803	0.5	500	103%	70	130	0%	
1,3-Dichloropropane	A	ug/L	24.38386	0.9753544		1	0	0	0.0791	0.5	500	98%	70	130	0%	
1,4-Dichlorobenzene	A	ug/L	25.32843	1.0131372		1	0	0	0.0858	0.5	500	101%	70	130	0%	
2,2-Dichloropropane	A	ug/L	26.26917	1.0507668		1	0	0	0.186	0.5	500	105%	70	130	0%	
2-Chlorotoluene	A	ug/L	25.05504	1.0022016		1	0	0	0.0876	0.5	500	100%	70	130	0%	
4-Chlorotoluene	A	ug/L	24.39357	0.9757428		1	0	0	0.0728	0.5	500	98%	70	130	0%	
Benzene	A	ug/L	23.79187	0.9516748		1	0	0	0.0914	0.5	500	95%	70	130	0%	
Bromobenzene	A	ug/L	24.76128	0.9904512		1	0	0	0.0831	0.5	500	99%	70	130	0%	
Bromochloromethane	A	ug/L	25.4383	1.017532		1	0	0	0.141	0.5	500	102%	70	130	0%	
Bromodichloromethane	A	ug/L	24.39404	0.9757616		1	0	0	0.12	0.5	500	98%	70	130	0%	
Bromoform	A	ug/L	25.92121	1.0368484		1	0	0	0.119	0.5	500	104%	70	130	0%	
Bromomethane	A	ug/L	25.77927	1.0311708		1	0	0	0.253	0.5	500	103%	70	130	0%	
Carbon tetrachloride	A	ug/L	24.77733	0.9910932		1	0	0	0.143	0.5	500	99%	70	130	0%	
Chlorobenzene	A	ug/L	24.70152	0.9880608		1	0	0	0.0914	0.5	500	99%	70	130	0%	
Chlorodibromomethane	A	ug/L	24.3492	0.973968		1	0	0	0.0841	0.5	500	97%	70	130	0%	
Chloroethane	A	ug/L	26.12501	1.0450004		1	0	0	0.169	0.5	500	105%	70	130	0%	
Chloroform	A	ug/L	24.17337	0.9669348		1	0	0	0.0789	0.5	500	97%	70	130	0%	
Chloromethane	A	ug/L	26.34224	1.0536896		1	0	0	0.162	0.5	500	105%	70	130	0%	
cis-1,2-Dichloroethene	A	ug/L	24.5653	0.982612		1	0	0	0.108	0.5	500	98%	70	130	0%	
cis-1,3-Dichloropropene	A	ug/L	23.25283	0.9301132		1	0	0	0.073	0.5	500	93%	70	130	0%	
Dibromomethane	A	ug/L	23.84392	0.9537568		1	0	0	0.147	0.5	500	95%	70	130	0%	
Dichlorodifluoromethane	A	ug/L	25.67929	1.0271716		1	0	0	0.175	0.5	500	103%	70	130	0%	
Ethylbenzene	A	ug/L	23.74212	0.9496848		1	0	0	0.0836	0.5	500	95%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970902	ICAL010422_3	VOC-8260-W-Q	CAL3	)A5975C\VG010	1/4/2022 4:28:05	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	45.78355	1.831342		2	0	0	0.15	0.5	1000	92%	70	130	0%	
Methyl ethyl ketone	A	ug/L	235.05043	9.4020172		10	0	0	1.77	10	5000	94%	70	130	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	23.04184	0.9216736		1	0	0	0.101	0.5	500	92%	70	130	0%	
Methylene chloride	A	ug/L	26.30581	1.0522324		1	0	0	0.338	0.5	500	105%	70	130	0%	
o-Xylene	A	ug/L	23.64197	0.9456788		1	0	0	0.0604	0.5	500	95%	70	130	0%	
Styrene	A	ug/L	23.41194	0.9364776		1	0	0	0.067	0.5	500	94%	70	130	0%	
Tetrachloroethene	A	ug/L	25.39483	1.0157932		1	0	0	0.0671	0.5	500	102%	70	130	0%	
Toluene	A	ug/L	23.63186	0.9452744		1	0	0	0.0679	0.5	500	95%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	25.46407	1.0185628		1	0	0	0.125	0.5	500	102%	70	130	0%	
trans-1,3-Dichloropropene	A	ug/L	23.78943	0.9515772		1	0	0	0.0846	0.5	500	95%	70	130	0%	
Trichloroethene	A	ug/L	24.14841	0.9659364		1	0	0	0.0993	0.5	500	97%	70	130	0%	
Trichlorofluoromethane	A	ug/L	26.65307	1.0661228		1	0	0	0.134	0.5	500	107%	70	130	0%	
Vinyl chloride	A	ug/L	25.64884	1.0259536		1	0	0	0.153	0.5	500	103%	70	130	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Xylenes, Total	M	ug/L	69.42552	2.7770208		3	0	0	0.0604	0.5	1500	93%	70	130	0%	
1,2-Dichloroethane-d4	S	ug/L	25.72803	1.0291212		1	0	0	0.229	0.5	500	103%	70	130	0%	
Dibromofluoromethane	S	ug/L	25.62188	1.0248752		1	0	0	0.129	0.5	500	102%	70	130	0%	
p-Bromofluorobenzene	S	ug/L	25.28989	1.0115956		1	0	0	0.149	0.5	500	101%	70	130	0%	
Toluene-d8	S	ug/L	23.3046	0.932184		1	0	0	0.23	0.5	500	93%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970904	ICAL010422_4	VOC-8260-W-Q	CAL4	)A5975C\VG010	1/4/2022 4:55:32	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	47.50287	1.9001148		2	0	0	0.101	0.5	500	95%	70	130	0%	
1,1,1-Trichloroethane	A	ug/L	48.26875	1.93075		2	0	0	0.131	0.5	500	97%	70	130	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	48.61239	1.9444956		2	0	0	0.0872	0.5	500	97%	70	130	0%	
1,1,2-Trichloroethane	A	ug/L	48.47589	1.9390356		2	0	0	0.108	0.5	500	97%	70	130	0%	
1,1-Dichloroethane	A	ug/L	49.18279	1.9673116		2	0	0	0.135	0.5	500	98%	70	130	0%	
1,1-Dichloroethene	A	ug/L	48.80561	1.9522244		2	0	0	0.141	0.5	500	98%	70	130	0%	
1,1-Dichloropropene	A	ug/L	47.76266	1.9105064		2	0	0	0.083	0.5	500	96%	70	130	0%	
1,2,3-Trichloropropane	A	ug/L	49.19244	1.9676976		2	0	0	0.235	0.5	500	98%	70	130	0%	
1,2-Dibromoethane	A	ug/L	49.38886	1.9755544		2	0	0	0.0916	0.5	500	99%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970904	ICAL010422_4	VOC-8260-W-Q	CAL4	)A5975C\VG010	1/4/2022 4:55:32	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	48.54976	1.9419904		2	0	0	0.0746	0.5	500	97%	70	130	0%	
1,2-Dichloroethane	A	ug/L	48.98798	1.9595192		2	0	0	0.116	0.5	500	98%	70	130	0%	
1,2-Dichloropropane	A	ug/L	47.52725	1.90109		2	0	0	0.0847	0.5	500	95%	70	130	0%	
1,3-Dichlorobenzene	A	ug/L	47.38535	1.895414		2	0	0	0.0803	0.5	500	95%	70	130	0%	
1,3-Dichloropropane	A	ug/L	48.8841	1.955364		2	0	0	0.0791	0.5	500	98%	70	130	0%	
1,4-Dichlorobenzene	A	ug/L	48.01064	1.9204256		2	0	0	0.0858	0.5	500	96%	70	130	0%	
2,2-Dichloropropane	A	ug/L	50.38039	2.0152156		2	0	0	0.186	0.5	500	101%	70	130	0%	
2-Chlorotoluene	A	ug/L	47.44663	1.8978652		2	0	0	0.0876	0.5	500	95%	70	130	0%	
4-Chlorotoluene	A	ug/L	48.3865	1.93546		2	0	0	0.0728	0.5	500	97%	70	130	0%	
Benzene	A	ug/L	48.00539	1.9202156		2	0	0	0.0914	0.5	500	96%	70	130	0%	
Bromobenzene	A	ug/L	47.5759	1.903036		2	0	0	0.0831	0.5	500	95%	70	130	0%	
Bromochloromethane	A	ug/L	51.62325	2.06493		2	0	0	0.141	0.5	500	103%	70	130	0%	
Bromodichloromethane	A	ug/L	47.2409	1.889636		2	0	0	0.12	0.5	500	94%	70	130	0%	
Bromoform	A	ug/L	50.51704	2.0206816		2	0	0	0.119	0.5	500	101%	70	130	0%	
Bromomethane	A	ug/L	47.59212	1.9036848		2	0	0	0.253	0.5	500	95%	70	130	0%	
Carbon tetrachloride	A	ug/L	47.75203	1.9100812		2	0	0	0.143	0.5	500	96%	70	130	0%	
Chlorobenzene	A	ug/L	47.39593	1.8958372		2	0	0	0.0914	0.5	500	95%	70	130	0%	
Chlorodibromomethane	A	ug/L	46.24113	1.8496452		2	0	0	0.0841	0.5	500	92%	70	130	0%	
Chloroethane	A	ug/L	46.22429	1.8489716		2	0	0	0.169	0.5	500	92%	70	130	0%	
Chloroform	A	ug/L	48.20314	1.9281256		2	0	0	0.0789	0.5	500	96%	70	130	0%	
Chloromethane	A	ug/L	49.79828	1.9919312		2	0	0	0.162	0.5	500	100%	70	130	0%	
cis-1,2-Dichloroethene	A	ug/L	48.41535	1.936614		2	0	0	0.108	0.5	500	97%	70	130	0%	
cis-1,3-Dichloropropene	A	ug/L	46.52826	1.8611304		2	0	0	0.073	0.5	500	93%	70	130	0%	
Dibromomethane	A	ug/L	47.4844	1.899376		2	0	0	0.147	0.5	500	95%	70	130	0%	
Dichlorodifluoromethane	A	ug/L	49.48348	1.9793392		2	0	0	0.175	0.5	500	99%	70	130	0%	
Ethylbenzene	A	ug/L	46.80795	1.872318		2	0	0	0.0836	0.5	500	94%	70	130	0%	
m+p-Xylenes	A	ug/L	92.53468	3.7013872		4	0	0	0.15	0.5	1000	93%	70	130	0%	
Methyl ethyl ketone	A	ug/L	479.42958	19.1771832		20	0	0	1.77	10	5000	96%	70	130	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	47.53006	1.9012024		2	0	0	0.101	0.5	500	95%	70	130	0%	
Methylene chloride	A	ug/L	50.44212	2.0176848		2	0	0	0.338	0.5	500	101%	70	130	0%	
o-Xylene	A	ug/L	47.5086	1.900344		2	0	0	0.0604	0.5	500	95%	70	130	0%	
Styrene	A	ug/L	46.70518	1.8682072		2	0	0	0.067	0.5	500	93%	70	130	0%	
Tetrachloroethene	A	ug/L	46.29317	1.8517268		2	0	0	0.0671	0.5	500	93%	70	130	0%	
Toluene	A	ug/L	47.01163	1.8804652		2	0	0	0.0679	0.5	500	94%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	49.51777	1.9807108		2	0	0	0.125	0.5	500	99%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970904	ICAL010422_4	VOC-8260-W-Q	CAL4	)A5975C\VG010	1/4/2022 4:55:32	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
trans-1,3-Dichloropropene	A	ug/L	47.0378	1.881512		2	0	0	0.0846	0.5	500	94%	70	130	0%	
Trichloroethene	A	ug/L	47.11894	1.8847576		2	0	0	0.0993	0.5	500	94%	70	130	0%	
Trichlorofluoromethane	A	ug/L	49.31283	1.9725132		2	0	0	0.134	0.5	500	99%	70	130	0%	
Vinyl chloride	A	ug/L	48.95796	1.9583184		2	0	0	0.153	0.5	500	98%	70	130	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Xylenes, Total	M	ug/L	140.04328	5.6017312		6	0	0	0.0604	0.5	1500	93%	70	130	0%	
1,2-Dichloroethane-d4	S	ug/L	48.12519	1.9250076		2	0	0	0.229	0.5	500	96%	70	130	0%	
Dibromofluoromethane	S	ug/L	48.16607	1.9266428		2	0	0	0.129	0.5	500	96%	70	130	0%	
p-Bromofluorobenzene	S	ug/L	46.6647	1.866588		2	0	0	0.149	0.5	500	93%	70	130	0%	
Toluene-d8	S	ug/L	47.14406	1.8857624		2	0	0	0.23	0.5	500	94%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970905	ICAL010422_5	VOC-8260-W-Q	CAL5	)A5975C\VG010	1/4/2022 5:50:25	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	119.0492	4.761968		5	0	0	0.101	0.5	500	95%	70	130	0%	
1,1,1-Trichloroethane	A	ug/L	118.57641	4.7430564		5	0	0	0.131	0.5	500	95%	70	130	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	115.61793	4.6247172		5	0	0	0.0872	0.5	500	92%	70	130	0%	
1,1,2-Trichloroethane	A	ug/L	117.41297	4.6965188		5	0	0	0.108	0.5	500	94%	70	130	0%	
1,1-Dichloroethane	A	ug/L	118.11254	4.7245016		5	0	0	0.135	0.5	500	94%	70	130	0%	
1,1-Dichloroethene	A	ug/L	119.87977	4.7951908		5	0	0	0.141	0.5	500	96%	70	130	0%	
1,1-Dichloropropene	A	ug/L	119.80016	4.7920064		5	0	0	0.083	0.5	500	96%	70	130	0%	
1,2,3-Trichloropropane	A	ug/L	112.62609	4.5050436		5	0	0	0.235	0.5	500	90%	70	130	0%	
1,2-Dibromoethane	A	ug/L	119.23942	4.7695768		5	0	0	0.0916	0.5	500	95%	70	130	0%	
1,2-Dichlorobenzene	A	ug/L	115.43227	4.6172908		5	0	0	0.0746	0.5	500	92%	70	130	0%	
1,2-Dichloroethane	A	ug/L	118.21434	4.7285736		5	0	0	0.116	0.5	500	95%	70	130	0%	
1,2-Dichloropropane	A	ug/L	121.98902	4.8795608		5	0	0	0.0847	0.5	500	98%	70	130	0%	
1,3-Dichlorobenzene	A	ug/L	117.4899	4.699596		5	0	0	0.0803	0.5	500	94%	70	130	0%	
1,3-Dichloropropane	A	ug/L	123.01316	4.9205264		5	0	0	0.0791	0.5	500	98%	70	130	0%	
1,4-Dichlorobenzene	A	ug/L	118.7699	4.750796		5	0	0	0.0858	0.5	500	95%	70	130	0%	
2,2-Dichloropropane	A	ug/L	118.32027	4.7328108		5	0	0	0.186	0.5	500	95%	70	130	0%	
2-Chlorotoluene	A	ug/L	120.26748	4.8106992		5	0	0	0.0876	0.5	500	96%	70	130	0%	
4-Chlorotoluene	A	ug/L	121.05908	4.8423632		5	0	0	0.0728	0.5	500	97%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970905	ICAL010422_5	VOC-8260-W-Q	CAL5	)A5975C\VG010	1/4/2022 5:50:25	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Benzene	A	ug/L	116.95526	4.6782104		5	0	0	0.0914	0.5	500	94%	70	130	0%	
Bromobenzene	A	ug/L	119.48008	4.7792032		5	0	0	0.0831	0.5	500	96%	70	130	0%	
Bromochloromethane	A	ug/L	118.06829	4.7227316		5	0	0	0.141	0.5	500	94%	70	130	0%	
Bromodichloromethane	A	ug/L	121.97488	4.8789952		5	0	0	0.12	0.5	500	98%	70	130	0%	
Bromoform	A	ug/L	115.7218	4.628872		5	0	0	0.119	0.5	500	93%	70	130	0%	
Bromomethane	A	ug/L	123.65037	4.9460148		5	0	0	0.253	0.5	500	99%	70	130	0%	
Carbon tetrachloride	A	ug/L	119.4667	4.778668		5	0	0	0.143	0.5	500	96%	70	130	0%	
Chlorobenzene	A	ug/L	120.69031	4.8276124		5	0	0	0.0914	0.5	500	97%	70	130	0%	
Chlorodibromomethane	A	ug/L	120.74537	4.8298148		5	0	0	0.0841	0.5	500	97%	70	130	0%	
Chloroethane	A	ug/L	122.40855	4.896342		5	0	0	0.169	0.5	500	98%	70	130	0%	
Chloroform	A	ug/L	114.59119	4.5836476		5	0	0	0.0789	0.5	500	92%	70	130	0%	
Chloromethane	A	ug/L	122.61785	4.904714		5	0	0	0.162	0.5	500	98%	70	130	0%	
cis-1,2-Dichloroethene	A	ug/L	116.61895	4.664758		5	0	0	0.108	0.5	500	93%	70	130	0%	
cis-1,3-Dichloropropene	A	ug/L	120.71159	4.8284636		5	0	0	0.073	0.5	500	97%	70	130	0%	
Dibromomethane	A	ug/L	118.24252	4.7297008		5	0	0	0.147	0.5	500	95%	70	130	0%	
Dichlorodifluoromethane	A	ug/L	127.81927	5.1127708		5	0	0	0.175	0.5	500	102%	70	130	0%	
Ethylbenzene	A	ug/L	122.52434	4.9009736		5	0	0	0.0836	0.5	500	98%	70	130	0%	
m+p-Xylenes	A	ug/L	250.25869	10.0103476		10	0	0	0.15	0.5	1000	100%	70	130	0%	
Methyl ethyl ketone	A	ug/L	1159.30194	46.3720776		50	0	0	1.77	10	5000	93%	70	130	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	127.13745	5.085498		5	0	0	0.101	0.5	500	102%	70	130	0%	
Methylene chloride	A	ug/L	110.6249	4.424996		5	0	0	0.338	0.5	500	88%	70	130	0%	
o-Xylene	A	ug/L	123.23778	4.9295112		5	0	0	0.0604	0.5	500	99%	70	130	0%	
Styrene	A	ug/L	127.19102	5.0876408		5	0	0	0.067	0.5	500	102%	70	130	0%	
Tetrachloroethene	A	ug/L	119.90031	4.7960124		5	0	0	0.0671	0.5	500	96%	70	130	0%	
Toluene	A	ug/L	122.65711	4.9062844		5	0	0	0.0679	0.5	500	98%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	118.65107	4.7460428		5	0	0	0.125	0.5	500	95%	70	130	0%	
trans-1,3-Dichloropropene	A	ug/L	121.49288	4.8597152		5	0	0	0.0846	0.5	500	97%	70	130	0%	
Trichloroethene	A	ug/L	123.46463	4.9385852		5	0	0	0.0993	0.5	500	99%	70	130	0%	
Trichlorofluoromethane	A	ug/L	129.06871	5.1627484		5	0	0	0.134	0.5	500	103%	70	130	0%	
Vinyl chloride	A	ug/L	125.88087	5.0352348		5	0	0	0.153	0.5	500	101%	70	130	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Xylenes, Total	M	ug/L	373.49647	14.9398588		15	0	0	0.0604	0.5	1500	100%	70	130	0%	
1,2-Dichloroethane-d4	S	ug/L	116.64203	4.6656812		5	0	0	0.229	0.5	500	93%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970905	ICAL010422_5	VOC-8260-W-Q	CAL5	)A5975C\VG010	1/4/2022 5:50:25	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Dibromofluoromethane	S	ug/L	115.11464	4.6045856		5	0	0	0.129	0.5	500	92%	70	130	0%	
p-Bromofluorobenzene	S	ug/L	117.93503	4.7174012		5	0	0	0.149	0.5	500	94%	70	130	0%	
Toluene-d8	S	ug/L	121.27495	4.850998		5	0	0	0.23	0.5	500	97%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970906	ICAL010422_6	VOC-8260-W-Q	CAL6	)A5975C\VG010	1/4/2022 6:45:10	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	254.82737	10.1930948		10	0	0	0.101	0.5	500	102%	70	130	0%	
1,1,1-Trichloroethane	A	ug/L	258.72281	10.3489124		10	0	0	0.131	0.5	500	103%	70	130	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	250.15769	10.0063076		10	0	0	0.0872	0.5	500	100%	70	130	0%	
1,1,2-Trichloroethane	A	ug/L	248.28816	9.9315264		10	0	0	0.108	0.5	500	99%	70	130	0%	
1,1-Dichloroethane	A	ug/L	258.43252	10.3373008		10	0	0	0.135	0.5	500	103%	70	130	0%	
1,1-Dichloroethene	A	ug/L	258.09028	10.3236112		10	0	0	0.141	0.5	500	103%	70	130	0%	
1,1-Dichloropropene	A	ug/L	264.6638	10.586552		10	0	0	0.083	0.5	500	106%	70	130	0%	
1,2,3-Trichloropropane	A	ug/L	249.26347	9.9705388		10	0	0	0.235	0.5	500	100%	70	130	0%	
1,2-Dibromoethane	A	ug/L	257.88869	10.3155476		10	0	0	0.0916	0.5	500	103%	70	130	0%	
1,2-Dichlorobenzene	A	ug/L	257.65242	10.3060968		10	0	0	0.0746	0.5	500	103%	70	130	0%	
1,2-Dichloroethane	A	ug/L	251.96754	10.0787016		10	0	0	0.116	0.5	500	101%	70	130	0%	
1,2-Dichloropropane	A	ug/L	254.71606	10.1886424		10	0	0	0.0847	0.5	500	102%	70	130	0%	
1,3-Dichlorobenzene	A	ug/L	258.62971	10.3451884		10	0	0	0.0803	0.5	500	103%	70	130	0%	
1,3-Dichloropropane	A	ug/L	263.47539	10.5390156		10	0	0	0.0791	0.5	500	105%	70	130	0%	
1,4-Dichlorobenzene	A	ug/L	254.91697	10.1966788		10	0	0	0.0858	0.5	500	102%	70	130	0%	
2,2-Dichloropropane	A	ug/L	253.03965	10.121586		10	0	0	0.186	0.5	500	101%	70	130	0%	
2-Chlorotoluene	A	ug/L	267.26165	10.690466		10	0	0	0.0876	0.5	500	107%	70	130	0%	
4-Chlorotoluene	A	ug/L	267.44092	10.6976368		10	0	0	0.0728	0.5	500	107%	70	130	0%	
Benzene	A	ug/L	257.54165	10.301666		10	0	0	0.0914	0.5	500	103%	70	130	0%	
Bromobenzene	A	ug/L	263.29438	10.5317752		10	0	0	0.0831	0.5	500	105%	70	130	0%	
Bromochloromethane	A	ug/L	247.05862	9.8823448		10	0	0	0.141	0.5	500	99%	70	130	0%	
Bromodichloromethane	A	ug/L	257.22856	10.2891424		10	0	0	0.12	0.5	500	103%	70	130	0%	
Bromoform	A	ug/L	257.5099	10.300396		10	0	0	0.119	0.5	500	103%	70	130	0%	
Bromomethane	A	ug/L	251.76065	10.070426		10	0	0	0.253	0.5	500	101%	70	130	0%	
Carbon tetrachloride	A	ug/L	260.87744	10.4350976		10	0	0	0.143	0.5	500	104%	70	130	0%	
Chlorobenzene	A	ug/L	258.25445	10.330178		10	0	0	0.0914	0.5	500	103%	70	130	0%	
Chlorodibromomethane	A	ug/L	258.35353	10.3341412		10	0	0	0.0841	0.5	500	103%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970906	ICAL010422_6	VOC-8260-W-Q	CAL6	)A5975C\VG010	1/4/2022 6:45:10	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Chloroethane	A	ug/L	231.74321	9.2697284		10	0	0	0.169	0.5	500	93%	70	130	0%	
Chloroform	A	ug/L	248.08043	9.9232172		10	0	0	0.0789	0.5	500	99%	70	130	0%	
Chloromethane	A	ug/L	240.2183	9.608732		10	0	0	0.162	0.5	500	96%	70	130	0%	
cis-1,2-Dichloroethene	A	ug/L	261.87064	10.4748256		10	0	0	0.108	0.5	500	105%	70	130	0%	
cis-1,3-Dichloropropene	A	ug/L	265.28626	10.6114504		10	0	0	0.073	0.5	500	106%	70	130	0%	
Dibromomethane	A	ug/L	252.27336	10.0909344		10	0	0	0.147	0.5	500	101%	70	130	0%	
Dichlorodifluoromethane	A	ug/L	252.15586	10.0862344		10	0	0	0.175	0.5	500	101%	70	130	0%	
Ethylbenzene	A	ug/L	266.81931	10.6727724		10	0	0	0.0836	0.5	500	107%	70	130	0%	
m+p-Xylenes	A	ug/L	543.42617	21.7370468		20	0	0	0.15	0.5	1000	109%	70	130	0%	
Methyl ethyl ketone	A	ug/L	2688.24739	107.529896		100	0	0	1.77	10	5000	108%	70	130	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	258.95351	10.3581404		10	0	0	0.101	0.5	500	104%	70	130	0%	
Methylene chloride	A	ug/L	235.46573	9.4186292		10	0	0	0.338	0.5	500	94%	70	130	0%	
o-Xylene	A	ug/L	270.46357	10.8185428		10	0	0	0.0604	0.5	500	108%	70	130	0%	
Styrene	A	ug/L	278.0455	11.12182		10	0	0	0.067	0.5	500	111%	70	130	0%	
Tetrachloroethene	A	ug/L	259.74185	10.389674		10	0	0	0.0671	0.5	500	104%	70	130	0%	
Toluene	A	ug/L	263.13299	10.5253196		10	0	0	0.0679	0.5	500	105%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	254.6608	10.186432		10	0	0	0.125	0.5	500	102%	70	130	0%	
trans-1,3-Dichloropropene	A	ug/L	263.80268	10.5521072		10	0	0	0.0846	0.5	500	106%	70	130	0%	
Trichloroethene	A	ug/L	262.29307	10.4917228		10	0	0	0.0993	0.5	500	105%	70	130	0%	
Trichlorofluoromethane	A	ug/L	259.05024	10.3620096		10	0	0	0.134	0.5	500	104%	70	130	0%	
Vinyl chloride	A	ug/L	248.65325	9.94613		10	0	0	0.153	0.5	500	99%	70	130	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Xylenes, Total	M	ug/L	813.88974	32.5555896		30	0	0	0.0604	0.5	1500	109%	70	130	0%	
1,2-Dichloroethane-d4	S	ug/L	258.23239	10.3292956		10	0	0	0.229	0.5	500	103%	70	130	0%	
Dibromofluoromethane	S	ug/L	259.02233	10.3608932		10	0	0	0.129	0.5	500	104%	70	130	0%	
p-Bromofluorobenzene	S	ug/L	267.31855	10.692742		10	0	0	0.149	0.5	500	107%	70	130	0%	
Toluene-d8	S	ug/L	270.0265	10.80106		10	0	0	0.23	0.5	500	108%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970907	ICAL010422_7	VOC-8260-W-Q	CAL7	)A5975C\VG010	1/4/2022 7:39:45	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970907	ICAL010422_7	VOC-8260-W-Q	CAL7	)A5975C\VG010	1/4/2022 7:39:45	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	392.18595	15.687438		15	0	0	0.101	0.5	500	105%	70	130	0%	
1,1,1-Trichloroethane	A	ug/L	386.6625	15.4665		15	0	0	0.131	0.5	500	103%	70	130	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	367.42759	14.6971036		15	0	0	0.0872	0.5	500	98%	70	130	0%	
1,1,2-Trichloroethane	A	ug/L	373.25341	14.9301364		15	0	0	0.108	0.5	500	100%	70	130	0%	
1,1-Dichloroethane	A	ug/L	380.44366	15.2177464		15	0	0	0.135	0.5	500	101%	70	130	0%	
1,1-Dichloroethene	A	ug/L	380.37253	15.2149012		15	0	0	0.141	0.5	500	101%	70	130	0%	
1,1-Dichloropropene	A	ug/L	397.13223	15.8852892		15	0	0	0.083	0.5	500	106%	70	130	0%	
1,2,3-Trichloropropane	A	ug/L	363.67316	14.5469264		15	0	0	0.235	0.5	500	97%	70	130	0%	
1,2-Dibromoethane	A	ug/L	377.7698	15.110792		15	0	0	0.0916	0.5	500	101%	70	130	0%	
1,2-Dichlorobenzene	A	ug/L	375.32826	15.0131304		15	0	0	0.0746	0.5	500	100%	70	130	0%	
1,2-Dichloroethane	A	ug/L	366.9787	14.679148		15	0	0	0.116	0.5	500	98%	70	130	0%	
1,2-Dichloropropane	A	ug/L	388.85021	15.5540084		15	0	0	0.0847	0.5	500	104%	70	130	0%	
1,3-Dichlorobenzene	A	ug/L	383.62247	15.3448988		15	0	0	0.0803	0.5	500	102%	70	130	0%	
1,3-Dichloropropane	A	ug/L	389.34421	15.5737684		15	0	0	0.0791	0.5	500	104%	70	130	0%	
1,4-Dichlorobenzene	A	ug/L	371.39689	14.8558756		15	0	0	0.0858	0.5	500	99%	70	130	0%	
2,2-Dichloropropane	A	ug/L	369.84356	14.7937424		15	0	0	0.186	0.5	500	99%	70	130	0%	
2-Chlorotoluene	A	ug/L	391.82688	15.6730752		15	0	0	0.0876	0.5	500	104%	70	130	0%	
4-Chlorotoluene	A	ug/L	396.27563	15.8510252		15	0	0	0.0728	0.5	500	106%	70	130	0%	
Benzene	A	ug/L	385.85261	15.4341044		15	0	0	0.0914	0.5	500	103%	70	130	0%	
Bromobenzene	A	ug/L	386.44198	15.4576792		15	0	0	0.0831	0.5	500	103%	70	130	0%	
Bromochloromethane	A	ug/L	371.80037	14.8720148		15	0	0	0.141	0.5	500	99%	70	130	0%	
Bromodichloromethane	A	ug/L	386.19404	15.4477616		15	0	0	0.12	0.5	500	103%	70	130	0%	
Bromoform	A	ug/L	378.22002	15.1288008		15	0	0	0.119	0.5	500	101%	70	130	0%	
Bromomethane	A	ug/L	385.12594	15.4050376		15	0	0	0.253	0.5	500	103%	70	130	0%	
Carbon tetrachloride	A	ug/L	386.9014	15.476056		15	0	0	0.143	0.5	500	103%	70	130	0%	
Chlorobenzene	A	ug/L	386.94547	15.4778188		15	0	0	0.0914	0.5	500	103%	70	130	0%	
Chlorodibromomethane	A	ug/L	387.68121	15.5072484		15	0	0	0.0841	0.5	500	103%	70	130	0%	
Chloroethane	A	ug/L	364.45728	14.5782912		15	0	0	0.169	0.5	500	97%	70	130	0%	
Chloroform	A	ug/L	366.93889	14.6775556		15	0	0	0.0789	0.5	500	98%	70	130	0%	
Chloromethane	A	ug/L	352.08363	14.0833452		15	0	0	0.162	0.5	500	94%	70	130	0%	
cis-1,2-Dichloroethene	A	ug/L	386.72365	15.468946		15	0	0	0.108	0.5	500	103%	70	130	0%	
cis-1,3-Dichloropropene	A	ug/L	400.79296	16.0317184		15	0	0	0.073	0.5	500	107%	70	130	0%	
Dibromomethane	A	ug/L	380.65469	15.2261876		15	0	0	0.147	0.5	500	102%	70	130	0%	
Dichlorodifluoromethane	A	ug/L	373.94485	14.957794		15	0	0	0.175	0.5	500	100%	70	130	0%	
Ethylbenzene	A	ug/L	404.7587	16.190348		15	0	0	0.0836	0.5	500	108%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970907	ICAL010422_7	VOC-8260-W-Q	CAL7	)A5975C\VG010	1/4/2022 7:39:45	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	812.85557	32.5142228		30	0	0	0.15	0.5	1000	108%	70	130	0%	
Methyl ethyl ketone	A	ug/L	3961.341	158.45364		150	0	0	1.77	10	5000	106%	70	130	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	391.17667	15.6470668		15	0	0	0.101	0.5	500	104%	70	130	0%	
Methylene chloride	A	ug/L	348.06663	13.9226652		15	0	0	0.338	0.5	500	93%	70	130	0%	
o-Xylene	A	ug/L	408.20432	16.3281728		15	0	0	0.0604	0.5	500	109%	70	130	0%	
Styrene	A	ug/L	413.75947	16.5503788		15	0	0	0.067	0.5	500	110%	70	130	0%	
Tetrachloroethene	A	ug/L	382.87963	15.3151852		15	0	0	0.0671	0.5	500	102%	70	130	0%	
Toluene	A	ug/L	397.01061	15.8804244		15	0	0	0.0679	0.5	500	106%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	376.13673	15.0454692		15	0	0	0.125	0.5	500	100%	70	130	0%	
trans-1,3-Dichloropropene	A	ug/L	402.10977	16.0843908		15	0	0	0.0846	0.5	500	107%	70	130	0%	
Trichloroethene	A	ug/L	394.48959	15.7795836		15	0	0	0.0993	0.5	500	105%	70	130	0%	
Trichlorofluoromethane	A	ug/L	371.42899	14.8571596		15	0	0	0.134	0.5	500	99%	70	130	0%	
Vinyl chloride	A	ug/L	372.35639	14.8942556		15	0	0	0.153	0.5	500	99%	70	130	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Xylenes, Total	M	ug/L	1221.05989	48.8423956		45	0	0	0.0604	0.5	1500	109%	70	130	0%	
1,2-Dichloroethane-d4	S	ug/L	378.33349	15.1333396		15	0	0	0.229	0.5	500	101%	70	130	0%	
Dibromofluoromethane	S	ug/L	384.7503	15.390012		15	0	0	0.129	0.5	500	103%	70	130	0%	
p-Bromofluorobenzene	S	ug/L	394.65655	15.786262		15	0	0	0.149	0.5	500	105%	70	130	0%	
Toluene-d8	S	ug/L	405.5583	16.222332		15	0	0	0.23	0.5	500	108%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970908	ICAL010422_8	VOC-8260-W-Q	CAL8	)A5975C\VG010	1/4/2022 8:34:31	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	520.28551	20.8114204		20	0	0	0.101	0.5	500	104%	70	130	0%	
1,1,1-Trichloroethane	A	ug/L	518.83124	20.7532496		20	0	0	0.131	0.5	500	104%	70	130	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	491.66999	19.6667996		20	0	0	0.0872	0.5	500	98%	70	130	0%	
1,1,2-Trichloroethane	A	ug/L	505.18031	20.2072124		20	0	0	0.108	0.5	500	101%	70	130	0%	
1,1-Dichloroethane	A	ug/L	515.3207	20.612828		20	0	0	0.135	0.5	500	103%	70	130	0%	
1,1-Dichloroethene	A	ug/L	515.06031	20.6024124		20	0	0	0.141	0.5	500	103%	70	130	0%	
1,1-Dichloropropene	A	ug/L	543.51208	21.7404832		20	0	0	0.083	0.5	500	109%	70	130	0%	
1,2,3-Trichloropropane	A	ug/L	491.52294	19.6609176		20	0	0	0.235	0.5	500	98%	70	130	0%	
1,2-Dibromoethane	A	ug/L	507.9234	20.316936		20	0	0	0.0916	0.5	500	102%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970908	ICAL010422_8	VOC-8260-W-Q	CAL8	)A5975C\VG010	1/4/2022 8:34:31	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	506.38707	20.2554828		20	0	0	0.0746	0.5	500	101%	70	130	0%	
1,2-Dichloroethane	A	ug/L	497.36991	19.8947964		20	0	0	0.116	0.5	500	99%	70	130	0%	
1,2-Dichloropropane	A	ug/L	524.16945	20.966778		20	0	0	0.0847	0.5	500	105%	70	130	0%	
1,3-Dichlorobenzene	A	ug/L	511.55042	20.4620168		20	0	0	0.0803	0.5	500	102%	70	130	0%	
1,3-Dichloropropane	A	ug/L	511.24793	20.4499172		20	0	0	0.0791	0.5	500	102%	70	130	0%	
1,4-Dichlorobenzene	A	ug/L	502.30007	20.0920028		20	0	0	0.0858	0.5	500	100%	70	130	0%	
2,2-Dichloropropane	A	ug/L	499.04726	19.9618904		20	0	0	0.186	0.5	500	100%	70	130	0%	
2-Chlorotoluene	A	ug/L	538.49638	21.5398552		20	0	0	0.0876	0.5	500	108%	70	130	0%	
4-Chlorotoluene	A	ug/L	531.84706	21.2738824		20	0	0	0.0728	0.5	500	106%	70	130	0%	
Benzene	A	ug/L	511.66576	20.4666304		20	0	0	0.0914	0.5	500	102%	70	130	0%	
Bromobenzene	A	ug/L	516.0104	20.640416		20	0	0	0.0831	0.5	500	103%	70	130	0%	
Bromochloromethane	A	ug/L	494.60544	19.7842176		20	0	0	0.141	0.5	500	99%	70	130	0%	
Bromodichloromethane	A	ug/L	518.37176	20.7348704		20	0	0	0.12	0.5	500	104%	70	130	0%	
Bromoform	A	ug/L	522.76605	20.910642		20	0	0	0.119	0.5	500	105%	70	130	0%	
Bromomethane	A	ug/L	515.01414	20.6005656		20	0	0	0.253	0.5	500	103%	70	130	0%	
Carbon tetrachloride	A	ug/L	521.26297	20.8505188		20	0	0	0.143	0.5	500	104%	70	130	0%	
Chlorobenzene	A	ug/L	515.99575	20.63983		20	0	0	0.0914	0.5	500	103%	70	130	0%	
Chlorodibromomethane	A	ug/L	520.43607	20.8174428		20	0	0	0.0841	0.5	500	104%	70	130	0%	
Chloroethane	A	ug/L	481.51432	19.2605728		20	0	0	0.169	0.5	500	96%	70	130	0%	
Chloroform	A	ug/L	489.12212	19.5648848		20	0	0	0.0789	0.5	500	98%	70	130	0%	
Chloromethane	A	ug/L	480.17469	19.2069876		20	0	0	0.162	0.5	500	96%	70	130	0%	
cis-1,2-Dichloroethene	A	ug/L	516.05445	20.642178		20	0	0	0.108	0.5	500	103%	70	130	0%	
cis-1,3-Dichloropropene	A	ug/L	538.90085	21.556034		20	0	0	0.073	0.5	500	108%	70	130	0%	
Dibromomethane	A	ug/L	500.74556	20.0298224		20	0	0	0.147	0.5	500	100%	70	130	0%	
Dichlorodifluoromethane	A	ug/L	494.74738	19.7898952		20	0	0	0.175	0.5	500	99%	70	130	0%	
Ethylbenzene	A	ug/L	544.68805	21.787522		20	0	0	0.0836	0.5	500	109%	70	130	0%	
m+p-Xylenes	A	ug/L	1087.40818	43.4963272		40	0	0	0.15	0.5	1000	109%	70	130	0%	
Methyl ethyl ketone	A	ug/L	5327.12526	213.085010		200	0	0	1.77	10	5000	107%	70	130	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	522.81865	20.912746		20	0	0	0.101	0.5	500	105%	70	130	0%	
Methylene chloride	A	ug/L	466.99932	18.6799728		20	0	0	0.338	0.5	500	93%	70	130	0%	
o-Xylene	A	ug/L	547.47638	21.8990552		20	0	0	0.0604	0.5	500	109%	70	130	0%	
Styrene	A	ug/L	555.79455	22.231782		20	0	0	0.067	0.5	500	111%	70	130	0%	
Tetrachloroethene	A	ug/L	514.92548	20.5970192		20	0	0	0.0671	0.5	500	103%	70	130	0%	
Toluene	A	ug/L	536.51007	21.4604028		20	0	0	0.0679	0.5	500	107%	70	130	0%	
trans-1,2-Dichloroethene	A	ug/L	510.00974	20.4003896		20	0	0	0.125	0.5	500	102%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970908	ICAL010422_8	VOC-8260-W-Q	CAL8	)A5975C\VG010	1/4/2022 8:34:31	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
trans-1,3-Dichloropropene	A	ug/L	533.75507	21.3502028		20	0	0	0.0846	0.5	500	107%	70	130	0%	
Trichloroethene	A	ug/L	534.40073	21.3760292		20	0	0	0.0993	0.5	500	107%	70	130	0%	
Trichlorofluoromethane	A	ug/L	489.6475	19.5859		20	0	0	0.134	0.5	500	98%	70	130	0%	
Vinyl chloride	A	ug/L	498.3563	19.934252		20	0	0	0.153	0.5	500	100%	70	130	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	70	130	0%	
Xylenes, Total	M	ug/L	1634.88456	65.3953824		60	0	0	0.0604	0.5	1500	109%	70	130	0%	
1,2-Dichloroethane-d4	S	ug/L	510.30803	20.4123212		20	0	0	0.229	0.5	500	102%	70	130	0%	
Dibromofluoromethane	S	ug/L	510.39915	20.415966		20	0	0	0.129	0.5	500	102%	70	130	0%	
p-Bromofluorobenzene	S	ug/L	541.3964	21.655856		20	0	0	0.149	0.5	500	108%	70	130	0%	
Toluene-d8	S	ug/L	544.21357	21.7685428		20	0	0	0.23	0.5	500	109%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970909	ICV010422	VOC-8260-W-Q	ICV	)A5975C\VG010	1/4/2022 9:29:14	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	126.66575	5.06663		5	0	0	0.101	0.5	500	101%	80	120	0%	
1,1,1-Trichloroethane	A	ug/L	128.25238	5.1300952		5	0	0	0.131	0.5	500	103%	80	120	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	127.47217	5.0988868		5	0	0	0.0872	0.5	500	102%	80	120	0%	
1,1,2-Trichloroethane	A	ug/L	123.03611	4.9214444		5	0	0	0.108	0.5	500	98%	80	120	0%	
1,1-Dichloroethane	A	ug/L	135.803	5.43212		5	0	0	0.135	0.5	500	109%	80	120	0%	
1,1-Dichloroethene	A	ug/L	134.45663	5.3782652		5	0	0	0.141	0.5	500	108%	80	120	0%	
1,1-Dichloropropene	A	ug/L	124.18526	4.9674104		5	0	0	0.083	0.5	500	99%	80	120	0%	
1,2,3-Trichloropropane	A	ug/L	122.95232	4.9180928		5	0	0	0.235	0.5	500	98%	80	120	0%	
1,2-Dibromoethane	A	ug/L	124.27642	4.9710568		5	0	0	0.0916	0.5	500	99%	80	120	0%	
1,2-Dichlorobenzene	A	ug/L	128.71039	5.1484156		5	0	0	0.0746	0.5	500	103%	80	120	0%	
1,2-Dichloroethane	A	ug/L	120.79914	4.8319656		5	0	0	0.116	0.5	500	97%	80	120	0%	
1,2-Dichloropropane	A	ug/L	125.66265	5.026506		5	0	0	0.0847	0.5	500	101%	80	120	0%	
1,3-Dichlorobenzene	A	ug/L	135.11854	5.4047416		5	0	0	0.0803	0.5	500	108%	80	120	0%	
1,3-Dichloropropane	A	ug/L	121.84417	4.8737668		5	0	0	0.0791	0.5	500	97%	80	120	0%	
1,4-Dichlorobenzene	A	ug/L	129.88123	5.1952492		5	0	0	0.0858	0.5	500	104%	80	120	0%	
2,2-Dichloropropane	A	ug/L	131.40305	5.256122		5	0	0	0.186	0.5	500	105%	80	120	0%	
2-Chlorotoluene	A	ug/L	131.29475	5.25179		5	0	0	0.0876	0.5	500	105%	80	120	0%	
4-Chlorotoluene	A	ug/L	137.07902	5.4831608		5	0	0	0.0728	0.5	500	110%	80	120	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970909	ICV010422	VOC-8260-W-Q	ICV	)A5975C\VG010	1/4/2022 9:29:14	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Benzene	A	ug/L	131.31393	5.2525572		5	0	0	0.0914	0.5	500	105%	80	120	0%	
Bromobenzene	A	ug/L	131.67879	5.2671516		5	0	0	0.0831	0.5	500	105%	80	120	0%	
Bromochloromethane	A	ug/L	123.60094	4.9440376		5	0	0	0.141	0.5	500	99%	80	120	0%	
Bromodichloromethane	A	ug/L	128.87588	5.1550352		5	0	0	0.12	0.5	500	103%	80	120	0%	
Bromoform	A	ug/L	129.9644	5.198576		5	0	0	0.119	0.5	500	104%	80	120	0%	
Bromomethane	A	ug/L	116.91567	4.6766268		5	0	0	0.253	0.5	500	94%	80	120	0%	
Carbon tetrachloride	A	ug/L	128.79275	5.15171		5	0	0	0.143	0.5	500	103%	80	120	0%	
Chlorobenzene	A	ug/L	131.63517	5.2654068		5	0	0	0.0914	0.5	500	105%	80	120	0%	
Chlorodibromomethane	A	ug/L	125.11031	5.0044124		5	0	0	0.0841	0.5	500	100%	80	120	0%	
Chloroethane	A	ug/L	115.59324	4.6237296		5	0	0	0.169	0.5	500	92%	80	120	0%	
Chloroform	A	ug/L	120.42358	4.8169432		5	0	0	0.0789	0.5	500	96%	80	120	0%	
Chloromethane	A	ug/L	108.77392	4.3509568		5	0	0	0.162	0.5	500	87%	80	120	0%	
cis-1,2-Dichloroethene	A	ug/L	130.12309	5.2049236		5	0	0	0.108	0.5	500	104%	80	120	0%	
cis-1,3-Dichloropropene	A	ug/L	121.55615	4.862246		5	0	0	0.073	0.5	500	97%	80	120	0%	
Dibromomethane	A	ug/L	125.30472	5.0121888		5	0	0	0.147	0.5	500	100%	80	120	0%	
Dichlorodifluoromethane	A	ug/L	111.37489	4.4549956		5	0	0	0.175	0.5	500	89%	80	120	0%	
Ethylbenzene	A	ug/L	131.91134	5.2764536		5	0	0	0.0836	0.5	500	106%	80	120	0%	
m+p-Xylenes	A	ug/L	262.75886	10.5103544		10	0	0	0.15	0.5	1000	105%	80	120	0%	
Methyl ethyl ketone	A	ug/L	1198.44392	47.9377568		50	0	0	1.77	10	5000	96%	80	120	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	134.72237	5.3888948		5	0	0	0.101	0.5	500	108%	80	120	0%	
Methylene chloride	A	ug/L	121.52968	4.8611872		5	0	0	0.338	0.5	500	97%	80	120	0%	
o-Xylene	A	ug/L	132.22141	5.2888564		5	0	0	0.0604	0.5	500	106%	80	120	0%	
Styrene	A	ug/L	137.49736	5.4998944		5	0	0	0.067	0.5	500	110%	80	120	0%	
Tetrachloroethene	A	ug/L	126.01413	5.0405652		5	0	0	0.0671	0.5	500	101%	80	120	0%	
Toluene	A	ug/L	132.0244	5.280976		5	0	0	0.0679	0.5	500	106%	80	120	0%	
trans-1,2-Dichloroethene	A	ug/L	134.70283	5.3881132		5	0	0	0.125	0.5	500	108%	80	120	0%	
trans-1,3-Dichloropropene	A	ug/L	129.02156	5.1608624		5	0	0	0.0846	0.5	500	103%	80	120	0%	
Trichloroethene	A	ug/L	131.10958	5.2443832		5	0	0	0.0993	0.5	500	105%	80	120	0%	
Trichlorofluoromethane	A	ug/L	121.7847	4.871388		5	0	0	0.134	0.5	500	97%	80	120	0%	
Vinyl chloride	A	ug/L	120.15175	4.80607		5	0	0	0.153	0.5	500	96%	80	120	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	394.98027	15.7992108		15	0	0	0.0604	0.5	1500	105%	80	120	0%	
1,2-Dichloroethane-d4	S	ug/L	280.28858	11.2115432		10	0	0	0.229	0.5	500	112%	80	120	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14970909	ICV010422	VOC-8260-W-Q	ICV	)A5975C\VG010	1/4/2022 9:29:14	1	R372940		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Dibromofluoromethane	S	ug/L	271.19937	10.8479748		10	0	0	0.129	0.5	500	108%	80	120	0%	
p-Bromofluorobenzene	S	ug/L	269.89759	10.7959036		10	0	0	0.149	0.5	500	108%	80	120	0%	
Toluene-d8	S	ug/L	276.91062	11.0764248		10	0	0	0.23	0.5	500	111%	80	120	0%	

```
Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN01.D
Sample Name    : PRIMER
Operator       : MSC
Date injected  :  4 Jan 2022   9:44 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 1
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN02.D
Sample Name    : BFB010422_
Operator       : MSC
Date injected  :  4 Jan 2022  10:11 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 2
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN03.D
Sample Name    : CCV010422_
Operator       : MSC
Date injected  :  4 Jan 2022  10:56 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 3
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN04.D
Sample Name    : PRIMER
Misc. Info.    : Replaced purge trap
Operator       : MSC
Date injected  :  4 Jan 2022  12:17 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 4
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN05.D
Sample Name    : BFB010422_
```

Operator   : MSC
Date injected : 4 Jan 2022 12:44 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.840
End Time   : 16.498
Vial Number   : 5

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN06.D
Sample Name   : CCV010422_
Operator   : MSC
Date injected : 4 Jan 2022   1:24 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.840
End Time   : 16.498
Vial Number   : 6

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN07.D
Sample Name   : BLK
Operator   : MSC
Date injected : 4 Jan 2022   2:09 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.839
End Time   : 16.498
Vial Number   : 7

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN08.D
Sample Name   : BFB010422_
Operator   : MSC
Date injected : 4 Jan 2022   2:38 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.840
End Time   : 16.498
Vial Number   : 8

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN09.D
Sample Name   : MBLK010422_
Operator   : MSC
Date injected : 4 Jan 2022   3:05 pm
Instrument   : VOA5975C

```
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 9
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN10.D
Sample Name    : ICAL010422_1
Operator       : MSC
Date injected  :  4 Jan 2022   3:33 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 10
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN11.D
Sample Name    : ICAL010422_2
Operator       : MSC
Date injected  :  4 Jan 2022   4:00 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 11
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN12.D
Sample Name    : ICAL010422_3
Operator       : MSC
Date injected  :  4 Jan 2022   4:28 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 12
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN13.D
Sample Name    : ICAL010422_4
Operator       : MSC
Date injected  :  4 Jan 2022   4:55 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
```

```
End Time       : 16.498
Vial Number    : 13
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN14.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  4 Jan 2022   5:22 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 14
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN15.D
Sample Name    : ICAL010422_5
Operator       : MSC
Date injected  :  4 Jan 2022   5:50 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 15
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN16.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  4 Jan 2022   6:17 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 16
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN17.D
Sample Name    : ICAL010422_6
Operator       : MSC
Date injected  :  4 Jan 2022   6:45 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 17
---------------------------------------------------------------
```

```
Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN18.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  4 Jan 2022   7:12 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 18
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN19.D
Sample Name    : ICAL010422_7
Operator       : MSC
Date injected  :  4 Jan 2022   7:39 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 19
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN20.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  4 Jan 2022   8:07 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.839
End Time       : 16.498
Vial Number    : 20
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN21.D
Sample Name    : ICAL010422_8
Operator       : MSC
Date injected  :  4 Jan 2022   8:34 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 21
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN22.D
Sample Name    : BLK
```

Operator   : MSC
Date injected : 4 Jan 2022   9:01 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.839
End Time   : 16.498
Vial Number   : 22

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN23.D
Sample Name   : ICV010422
Operator   : MSC
Date injected : 4 Jan 2022   9:29 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.840
End Time   : 16.498
Vial Number   : 23

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN24.D
Sample Name   : BLK
Operator   : MSC
Date injected : 4 Jan 2022   9:56 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.839
End Time   : 16.498
Vial Number   : 24

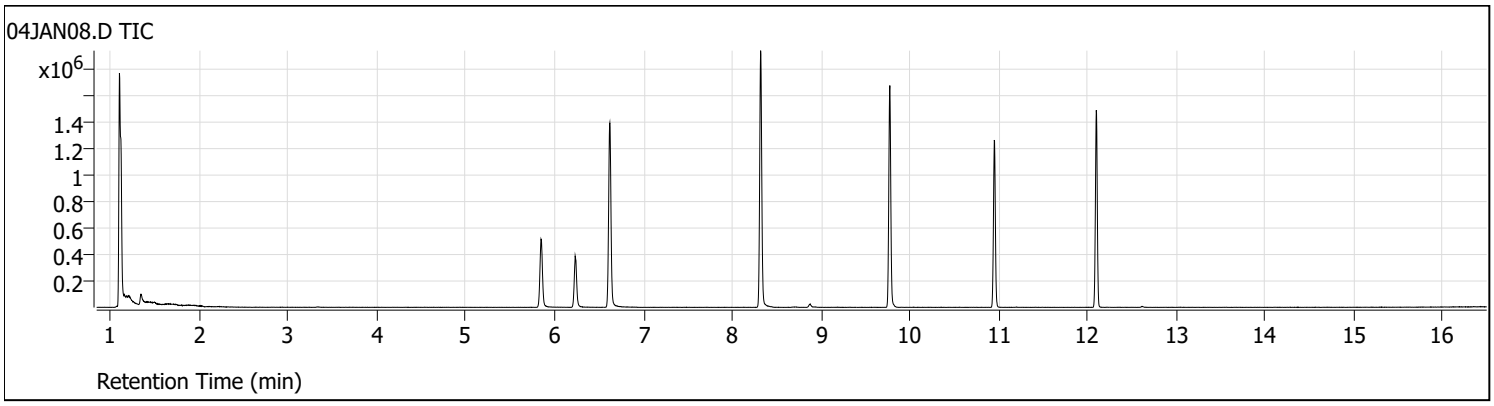
Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN25.D
Sample Name   : MDL010422_Q1_1
Operator   : MSC
Date injected : 4 Jan 2022 10:23 pm
Instrument   : VOA5975C
Method used   : 5975CACQF
No of spectra : 5616
Start Time   : 0.840
End Time   : 16.498
Vial Number   : 25

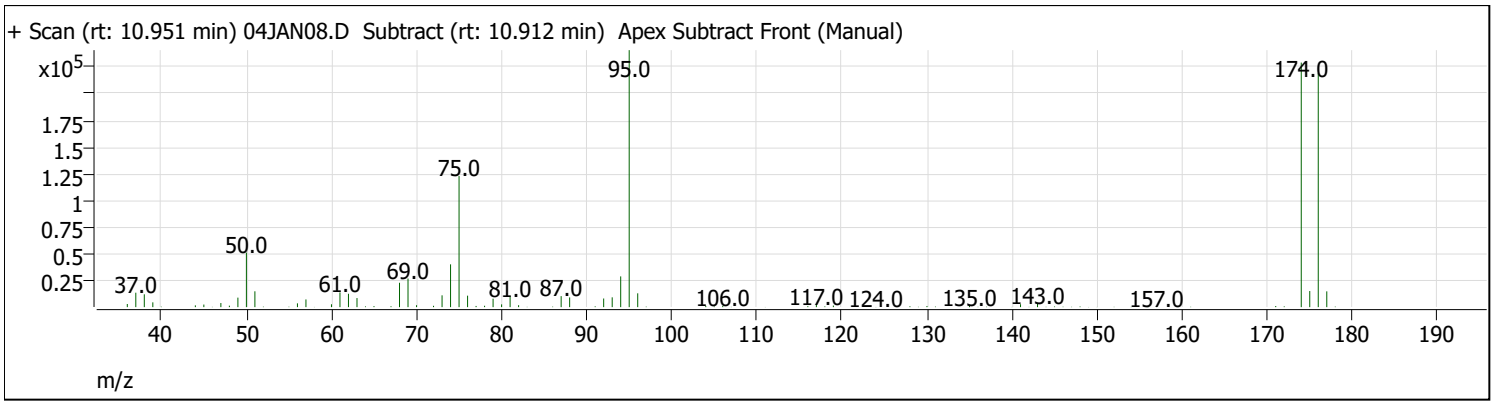
Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN26.D
Sample Name   : LOD010422_Q1_HalfCAL2
Operator   : MSC
Date injected : 4 Jan 2022 10:51 pm
Instrument   : VOA5975C

```
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.839
End Time       : 16.498
Vial Number    : 26
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN27.D
Sample Name    : MDL010422_Q1_2xCAL1
Operator       : MSC
Date injected  :  4 Jan 2022  11:18 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 27
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010422\04JAN28.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  4 Jan 2022  11:45 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 28
---------------------------------------------------------------
```

Tune Evaluation Report

Data Path:	D:\Org\Data\VOA5975C\VG010422\04JAN08.D
Acq on:	1/4/2022 2:38:09 PM
Operator:	MSC
Sample:	BFB010422_
Inst Name:	VOA5975C
ALS Vial:	8
Method:	\\MASSHUNTER\Org\Data\Methods\BFBapex.m





Target Mass	Rel. To Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Pass/Fail
50	95	15	40	21.2	51080	Pass
75	95	30	60	51.0	122824	Pass
95	95	100	100	100.0	240768	Pass
96	95	5	9	5.4	12961	Pass
173	174	0	2	0.0	0	Pass
174	95	50	100	95.2	229120	Pass
175	174	5	9	6.6	15102	Pass
176	174	95	101	95.7	219264	Pass
177	176	5	9	6.7	14796	Pass

Quantitative Analysis Results Summary Report

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:48 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Sequence Table

Data File	sample Name	Sample Type	Vial Position	Inj Vol	Level	Acq Method File
04JAN09.D	MBLK010422_	Method Blank	9	0		5975CACQF.M
04JAN10.D	ICAL010422_1	Cal	10	0	1	5975CACQF.M
04JAN11.D	ICAL010422_2	Cal	11	0	2	5975CACQF.M
04JAN12.D	ICAL010422_3	Cal	12	0	3	5975CACQF.M
04JAN13.D	ICAL010422_4	Cal	13	0	4	5975CACQF.M
04JAN15.D	ICAL010422_5	Cal	15	0	5	5975CACQF.M
04JAN17.D	ICAL010422_6	Cal	17	0	6	5975CACQF.M
04JAN19.D	ICAL010422_7	Cal	19	0	7	5975CACQF.M
04JAN21.D	ICAL010422_8	Cal	21	0	8	5975CACQF.M
04JAN23.D	ICV010422	QC	23	0	QC	5975CACQF.M

Quantitation Results

Compound: Dichlorodifluoromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	1.241	4353	770895	0.0056	4.3090	2.5000	172.4
04JAN11.D	Calibration	Fluorobenzene	1.244	12087	764419	0.0158	12.0663	12.5000	96.5
04JAN12.D	Calibration	Fluorobenzene	1.244	26627	791270	0.0337	25.6793	25.0000	102.7
04JAN13.D	Calibration	Fluorobenzene	1.241	50457	778120	0.0648	49.4835	50.0000	99.0
04JAN15.D	Calibration	Fluorobenzene	1.241	137933	823488	0.1675	127.8193	125.0000	102.3
04JAN17.D	Calibration	Fluorobenzene	1.241	276334	836278	0.3304	252.1559	250.0000	100.9
04JAN19.D	Calibration	Fluorobenzene	1.241	412544	841876	0.4900	373.9449	375.0000	99.7
04JAN21.D	Calibration	Fluorobenzene	1.241	545484	841364	0.6483	494.7474	500.0000	98.9
04JAN23.D	QC	Fluorobenzene	1.241	116936	801210	0.1459	111.3749	125.0000	

Compound: Chloromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene	1.333	0	775552	0.0000	ND		
04JAN10.D	Calibration	Fluorobenzene	1.406	7435	770895	0.0096	6.0637	2.5000	242.5
04JAN11.D	Calibration	Fluorobenzene	1.406	16859	764419	0.0221	13.8661	12.5000	110.9
04JAN12.D	Calibration	Fluorobenzene	1.406	33153	791270	0.0419	26.3422	25.0000	105.4
04JAN13.D	Calibration	Fluorobenzene	1.408	61632	778120	0.0792	49.7983	50.0000	99.6
04JAN15.D	Calibration	Fluorobenzene	1.409	160604	823488	0.1950	122.6179	125.0000	98.1
04JAN17.D	Calibration	Fluorobenzene	1.408	319523	836278	0.3821	240.2183	250.0000	96.1
04JAN19.D	Calibration	Fluorobenzene	1.409	471454	841876	0.5600	352.0836	375.0000	93.9
04JAN21.D	Calibration	Fluorobenzene	1.406	642582	841364	0.7637	480.1747	500.0000	96.0

Compound: Chloromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN23.D	QC	Fluorobenzene	1.406	138617	801210	0.1730	108.7739	125.0000	

Compound: Vinyl chloride

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene	1.489	0	775552	0.0000	ND		
04JAN10.D	Calibration	Fluorobenzene	1.495	4274	770895	0.0055	3.8739	2.5000	155.0
04JAN11.D	Calibration	Fluorobenzene	1.498	13724	764419	0.0180	12.5446	12.5000	100.4
04JAN12.D	Calibration	Fluorobenzene	1.495	29046	791270	0.0367	25.6488	25.0000	102.6
04JAN13.D	Calibration	Fluorobenzene	1.495	54521	778120	0.0701	48.9580	50.0000	97.9
04JAN15.D	Calibration	Fluorobenzene	1.495	148358	823488	0.1802	125.8809	125.0000	100.7
04JAN17.D	Calibration	Fluorobenzene	1.498	297604	836278	0.3559	248.6532	250.0000	99.5
04JAN19.D	Calibration	Fluorobenzene	1.498	448643	841876	0.5329	372.3564	375.0000	99.3
04JAN21.D	Calibration	Fluorobenzene	1.495	600092	841364	0.7132	498.3563	500.0000	99.7
04JAN23.D	QC	Fluorobenzene	1.495	137775	801210	0.1720	120.1518	125.0000	

Compound: Bromomethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	1.796	1902	770895	0.0025	3.8547	2.5000	154.2
04JAN11.D	Calibration	Fluorobenzene	1.796	5893	764419	0.0077	12.0464	12.5000	96.4
04JAN12.D	Calibration	Fluorobenzene	1.796	13054	791270	0.0165	25.7793	25.0000	103.1
04JAN13.D	Calibration	Fluorobenzene	1.799	23699	778120	0.0305	47.5921	50.0000	95.2
04JAN15.D	Calibration	Fluorobenzene	1.799	65163	823488	0.0791	123.6504	125.0000	98.9
04JAN17.D	Calibration	Fluorobenzene	1.799	134737	836278	0.1611	251.7606	250.0000	100.7
04JAN19.D	Calibration	Fluorobenzene	1.796	207491	841876	0.2465	385.1259	375.0000	102.7
04JAN21.D	Calibration	Fluorobenzene	1.793	277301	841364	0.3296	515.0141	500.0000	103.0
04JAN23.D	QC	Fluorobenzene	1.796	59947	801210	0.0748	116.9157	125.0000	

Compound: Chloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	1.899	2178	770895	0.0028	3.9871	2.5000	159.5
04JAN11.D	Calibration	Fluorobenzene	1.897	8052	764419	0.0105	14.8670	12.5000	118.9
04JAN12.D	Calibration	Fluorobenzene	1.897	14646	791270	0.0185	26.1250	25.0000	104.5
04JAN13.D	Calibration	Fluorobenzene	1.897	25484	778120	0.0328	46.2243	50.0000	92.4
04JAN15.D	Calibration	Fluorobenzene	1.894	71420	823488	0.0867	122.4086	125.0000	97.9
04JAN17.D	Calibration	Fluorobenzene	1.894	137312	836278	0.1642	231.7432	250.0000	92.7
04JAN19.D	Calibration	Fluorobenzene	1.897	217393	841876	0.2582	364.4573	375.0000	97.2
04JAN21.D	Calibration	Fluorobenzene	1.894	287041	841364	0.3412	481.5143	500.0000	96.3
04JAN23.D	QC	Fluorobenzene	1.897	65619	801210	0.0819	115.5932	125.0000	

Compound: Trichlorofluoromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	2.153	5030	770895	0.0065	3.6731	2.5000	146.9
04JAN11.D	Calibration	Fluorobenzene	2.142	15431	764419	0.0202	11.3637	12.5000	90.9
04JAN12.D	Calibration	Fluorobenzene	2.142	37464	791270	0.0473	26.6531	25.0000	106.6
04JAN13.D	Calibration	Fluorobenzene	2.145	68163	778120	0.0876	49.3128	50.0000	98.6
04JAN15.D	Calibration	Fluorobenzene	2.142	188808	823488	0.2293	129.0687	125.0000	103.3
04JAN17.D	Calibration	Fluorobenzene	2.145	384837	836278	0.4602	259.0502	250.0000	103.6
04JAN19.D	Calibration	Fluorobenzene	2.145	555477	841876	0.6598	371.4290	375.0000	99.0
04JAN21.D	Calibration	Fluorobenzene	2.145	731829	841364	0.8698	489.6475	500.0000	97.9
04JAN23.D	QC	Fluorobenzene	2.145	173333	801210	0.2163	121.7847	125.0000	

Compound: 1,1-Dichloroethene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	2.700	2084	770895	0.0027	2.6839	2.5000	107.4
04JAN11.D	Calibration	Fluorobenzene	2.700	9169	764419	0.0120	11.9081	12.5000	95.3
04JAN12.D	Calibration	Fluorobenzene	2.700	20631	791270	0.0261	25.8849	25.0000	103.5
04JAN13.D	Calibration	Fluorobenzene	2.702	38253	778120	0.0492	48.8056	50.0000	97.6
04JAN15.D	Calibration	Fluorobenzene	2.697	99438	823488	0.1208	119.8798	125.0000	95.9
04JAN17.D	Calibration	Fluorobenzene	2.702	217406	836278	0.2600	258.0903	250.0000	103.2
04JAN19.D	Calibration	Fluorobenzene	2.700	322557	841876	0.3831	380.3725	375.0000	101.4
04JAN21.D	Calibration	Fluorobenzene	2.700	436507	841364	0.5188	515.0603	500.0000	103.0
04JAN23.D	QC	Fluorobenzene	2.702	108512	801210	0.1354	134.4566	125.0000	

Compound: Methylene chloride

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene	3.335	1661	775552	0.0021	1.4424		
04JAN10.D	Calibration	Fluorobenzene	3.324	4095	770895	0.0053	3.5774	2.5000	143.1
04JAN11.D	Calibration	Fluorobenzene	3.338	17734	764419	0.0232	15.6236	12.5000	125.0
04JAN12.D	Calibration	Fluorobenzene	3.333	30908	791270	0.0391	26.3058	25.0000	105.2
04JAN13.D	Calibration	Fluorobenzene	3.335	58282	778120	0.0749	50.4421	50.0000	100.9
04JAN15.D	Calibration	Fluorobenzene	3.336	135271	823488	0.1643	110.6249	125.0000	88.5
04JAN17.D	Calibration	Fluorobenzene	3.333	292397	836278	0.3496	235.4657	250.0000	94.2
04JAN19.D	Calibration	Fluorobenzene	3.330	435116	841876	0.5168	348.0666	375.0000	92.8
04JAN21.D	Calibration	Fluorobenzene	3.330	583438	841364	0.6934	466.9993	500.0000	93.4
04JAN23.D	QC	Fluorobenzene	3.330	144585	801210	0.1805	121.5297	125.0000	

Compound: trans-1,2-Dichloroethene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	3.723	2146	770895	0.0028	2.7090	2.5000	108.4
04JAN11.D	Calibration	Fluorobenzene	3.720	9821	764419	0.0128	12.5022	12.5000	100.0

Compound: trans-1,2-Dichloroethene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN12.D	Calibration	Fluorobenzene	3.712	20706	791270	0.0262	25.4641	25.0000	101.9
04JAN13.D	Calibration	Fluorobenzene	3.717	39596	778120	0.0509	49.5178	50.0000	99.0
04JAN15.D	Calibration	Fluorobenzene	3.718	100409	823488	0.1219	118.6511	125.0000	94.9
04JAN17.D	Calibration	Fluorobenzene	3.715	218855	836278	0.2617	254.6608	250.0000	101.9
04JAN19.D	Calibration	Fluorobenzene	3.715	325415	841876	0.3865	376.1367	375.0000	100.3
04JAN21.D	Calibration	Fluorobenzene	3.718	440967	841364	0.5241	510.0097	500.0000	102.0
04JAN23.D	QC	Fluorobenzene	3.715	110909	801210	0.1384	134.7028	125.0000	

Compound: Methyl tert-butyl ether (MTBE)

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	3.759	2717	770895	0.0035	2.6532	2.5000	106.1
04JAN11.D	Calibration	Fluorobenzene	3.762	12515	764419	0.0164	12.3255	12.5000	98.6
04JAN12.D	Calibration	Fluorobenzene	3.754	24218	791270	0.0306	23.0418	25.0000	92.2
04JAN13.D	Calibration	Fluorobenzene	3.757	49126	778120	0.0631	47.5301	50.0000	95.1
04JAN15.D	Calibration	Fluorobenzene	3.754	139068	823488	0.1689	127.1375	125.0000	101.7
04JAN17.D	Calibration	Fluorobenzene	3.751	287653	836278	0.3440	258.9535	250.0000	103.6
04JAN19.D	Calibration	Fluorobenzene	3.751	437439	841876	0.5196	391.1767	375.0000	104.3
04JAN21.D	Calibration	Fluorobenzene	3.754	584294	841364	0.6945	522.8187	500.0000	104.6
04JAN23.D	QC	Fluorobenzene	3.754	143378	801210	0.1790	134.7224	125.0000	

Compound: 1,1-Dichloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	4.376	3892	770895	0.0050	2.6393	2.5000	105.6
04JAN11.D	Calibration	Fluorobenzene	4.378	17642	764419	0.0231	12.0652	12.5000	96.5
04JAN12.D	Calibration	Fluorobenzene	4.379	38874	791270	0.0491	25.6835	25.0000	102.7
04JAN13.D	Calibration	Fluorobenzene	4.381	73205	778120	0.0941	49.1828	50.0000	98.4
04JAN15.D	Calibration	Fluorobenzene	4.378	186052	823488	0.2259	118.1125	125.0000	94.5
04JAN17.D	Calibration	Fluorobenzene	4.384	413408	836278	0.4943	258.4325	250.0000	103.4
04JAN19.D	Calibration	Fluorobenzene	4.381	612660	841876	0.7277	380.4437	375.0000	101.5
04JAN21.D	Calibration	Fluorobenzene	4.378	829359	841364	0.9857	515.3207	500.0000	103.1
04JAN23.D	QC	Fluorobenzene	4.376	208131	801210	0.2598	135.8030	125.0000	

Compound: 2,2-Dichloropropane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	5.196	2930	770895	0.0038	2.6520	2.5000	106.1
04JAN11.D	Calibration	Fluorobenzene	5.196	13676	764419	0.0179	12.4820	12.5000	99.9
04JAN12.D	Calibration	Fluorobenzene	5.190	29793	791270	0.0377	26.2692	25.0000	105.1
04JAN13.D	Calibration	Fluorobenzene	5.193	56189	778120	0.0722	50.3804	50.0000	100.8
04JAN15.D	Calibration	Fluorobenzene	5.196	139656	823488	0.1696	118.3203	125.0000	94.7

Compound: 2,2-Dichloropropane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN17.D	Calibration	Fluorobenzene	5.190	303307	836278	0.3627	253.0397	250.0000	101.2
04JAN19.D	Calibration	Fluorobenzene	5.190	446282	841876	0.5301	369.8436	375.0000	98.6
04JAN21.D	Calibration	Fluorobenzene	5.190	601823	841364	0.7153	499.0473	500.0000	99.8
04JAN23.D	QC	Fluorobenzene	5.190	150902	801210	0.1883	131.4031	125.0000	

Compound: cis-1,2-Dichloroethene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	5.212	2376	770895	0.0031	2.9581	2.5000	118.3
04JAN11.D	Calibration	Fluorobenzene	5.221	10008	764419	0.0131	12.5659	12.5000	100.5
04JAN12.D	Calibration	Fluorobenzene	5.212	20252	791270	0.0256	24.5653	25.0000	98.3
04JAN13.D	Calibration	Fluorobenzene	5.209	39251	778120	0.0504	48.4154	50.0000	96.8
04JAN15.D	Calibration	Fluorobenzene	5.215	100057	823488	0.1215	116.6190	125.0000	93.3
04JAN17.D	Calibration	Fluorobenzene	5.215	228170	836278	0.2728	261.8706	250.0000	104.7
04JAN19.D	Calibration	Fluorobenzene	5.212	339211	841876	0.4029	386.7236	375.0000	103.1
04JAN21.D	Calibration	Fluorobenzene	5.212	452377	841364	0.5377	516.0544	500.0000	103.2
04JAN23.D	QC	Fluorobenzene	5.209	108623	801210	0.1356	130.1231	125.0000	

Compound: Methyl ethyl ketone

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	5.302	3035	770895	0.0039	27.8967	25.0000	111.6
04JAN11.D	Calibration	Fluorobenzene	5.288	13167	764419	0.0172	122.0520	125.0000	97.6
04JAN12.D	Calibration	Fluorobenzene	5.282	26248	791270	0.0332	235.0504	250.0000	94.0
04JAN13.D	Calibration	Fluorobenzene	5.285	52648	778120	0.0677	479.4296	500.0000	95.9
04JAN15.D	Calibration	Fluorobenzene	5.282	134730	823488	0.1636	1159.3019	1250.0000	92.7
04JAN17.D	Calibration	Fluorobenzene	5.279	317271	836278	0.3794	2688.2474	2500.0000	107.5
04JAN19.D	Calibration	Fluorobenzene	5.279	470653	841876	0.5591	3961.3410	3750.0000	105.6
04JAN21.D	Calibration	Fluorobenzene	5.279	632539	841364	0.7518	5327.1253	5000.0000	106.5
04JAN23.D	QC	Fluorobenzene	5.282	135511	801210	0.1691	1198.4439	1250.0000	

Compound: Bromochloromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	5.522	807	770895	0.0010	2.4260	2.5000	97.0
04JAN11.D	Calibration	Fluorobenzene	5.516	4275	764419	0.0056	12.9568	12.5000	103.7
04JAN12.D	Calibration	Fluorobenzene	5.522	8688	791270	0.0110	25.4383	25.0000	101.8
04JAN13.D	Calibration	Fluorobenzene	5.516	17338	778120	0.0223	51.6233	50.0000	103.2
04JAN15.D	Calibration	Fluorobenzene	5.519	41966	823488	0.0510	118.0683	125.0000	94.5
04JAN17.D	Calibration	Fluorobenzene	5.519	89178	836278	0.1066	247.0586	250.0000	98.8
04JAN19.D	Calibration	Fluorobenzene	5.516	135103	841876	0.1605	371.8004	375.0000	99.1
04JAN21.D	Calibration	Fluorobenzene	5.519	179618	841364	0.2135	494.6054	500.0000	98.9

Compound: Bromochloromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN23.D	QC	Fluorobenzene	5.513	42744	801210	0.0533	123.6009	125.0000	

Compound: Chloroform

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	5.659	4248	770895	0.0055	2.8946	2.5000	115.8
04JAN11.D	Calibration	Fluorobenzene	5.656	19015	764419	0.0249	13.0668	12.5000	104.5
04JAN12.D	Calibration	Fluorobenzene	5.653	36413	791270	0.0460	24.1734	25.0000	96.7
04JAN13.D	Calibration	Fluorobenzene	5.650	71403	778120	0.0918	48.2031	50.0000	96.4
04JAN15.D	Calibration	Fluorobenzene	5.653	179640	823488	0.2181	114.5912	125.0000	91.7
04JAN17.D	Calibration	Fluorobenzene	5.653	394946	836278	0.4723	248.0804	250.0000	99.2
04JAN19.D	Calibration	Fluorobenzene	5.650	588080	841876	0.6985	366.9389	375.0000	97.9
04JAN21.D	Calibration	Fluorobenzene	5.653	783422	841364	0.9311	489.1221	500.0000	97.8
04JAN23.D	QC	Fluorobenzene	5.647	183676	801210	0.2292	120.4236	125.0000	

Compound: 1,1,1-Trichloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	5.831	3510	770895	0.0046	2.5521	2.5000	102.1
04JAN11.D	Calibration	Fluorobenzene	5.837	16623	764419	0.0217	12.1891	12.5000	97.5
04JAN12.D	Calibration	Fluorobenzene	5.826	35547	791270	0.0449	25.1809	25.0000	100.7
04JAN13.D	Calibration	Fluorobenzene	5.834	67007	778120	0.0861	48.2688	50.0000	96.5
04JAN15.D	Calibration	Fluorobenzene	5.834	174206	823488	0.2115	118.5764	125.0000	94.9
04JAN17.D	Calibration	Fluorobenzene	5.831	386005	836278	0.4616	258.7228	250.0000	103.5
04JAN19.D	Calibration	Fluorobenzene	5.831	580748	841876	0.6898	386.6625	375.0000	103.1
04JAN21.D	Calibration	Fluorobenzene	5.834	778785	841364	0.9256	518.8312	500.0000	103.8
04JAN23.D	QC	Fluorobenzene	5.831	183324	801210	0.2288	128.2524	125.0000	

Compound: Dibromofluoromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene	5.848	203459	775552	0.2623	278.4635		
04JAN10.D	Calibration	Fluorobenzene	5.851	2508	770895	0.0033	3.4533	2.5000	138.1
04JAN11.D	Calibration	Fluorobenzene	5.845	9074	764419	0.0119	12.6000	12.5000	100.8
04JAN12.D	Calibration	Fluorobenzene	5.845	19100	791270	0.0241	25.6219	25.0000	102.5
04JAN13.D	Calibration	Fluorobenzene	5.848	35309	778120	0.0454	48.1661	50.0000	96.3
04JAN15.D	Calibration	Fluorobenzene	5.845	89307	823488	0.1084	115.1146	125.0000	92.1
04JAN17.D	Calibration	Fluorobenzene	5.845	204073	836278	0.2440	259.0223	250.0000	103.6
04JAN19.D	Calibration	Fluorobenzene	5.848	305158	841876	0.3625	384.7503	375.0000	102.6
04JAN21.D	Calibration	Fluorobenzene	5.845	404568	841364	0.4808	510.3991	500.0000	102.1
04JAN23.D	QC	Fluorobenzene	5.848	204707	801210	0.2555	271.1994	250.0000	

Compound: Carbon tetrachloride

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	6.029	4342	770895	0.0056	3.2043	2.5000	128.2
04JAN11.D	Calibration	Fluorobenzene	6.021	16466	764419	0.0215	12.2545	12.5000	98.0
04JAN12.D	Calibration	Fluorobenzene	6.024	34462	791270	0.0436	24.7773	25.0000	99.1
04JAN13.D	Calibration	Fluorobenzene	6.026	65313	778120	0.0839	47.7520	50.0000	95.5
04JAN15.D	Calibration	Fluorobenzene	6.024	172928	823488	0.2100	119.4667	125.0000	95.6
04JAN17.D	Calibration	Fluorobenzene	6.026	383485	836278	0.4586	260.8774	250.0000	104.4
04JAN19.D	Calibration	Fluorobenzene	6.024	572545	841876	0.6801	386.9014	375.0000	103.2
04JAN21.D	Calibration	Fluorobenzene	6.024	770907	841364	0.9163	521.2630	500.0000	104.3
04JAN23.D	QC	Fluorobenzene	6.027	181384	801210	0.2264	128.7928	125.0000	

Compound: 1,1-Dichloropropene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	6.041	2830	770895	0.0037	2.4201	2.5000	96.8
04JAN11.D	Calibration	Fluorobenzene	6.038	13149	764419	0.0172	11.3397	12.5000	90.7
04JAN12.D	Calibration	Fluorobenzene	6.038	29241	791270	0.0370	24.3617	25.0000	97.4
04JAN13.D	Calibration	Fluorobenzene	6.035	56376	778120	0.0725	47.7627	50.0000	95.5
04JAN15.D	Calibration	Fluorobenzene	6.038	149649	823488	0.1817	119.8002	125.0000	95.8
04JAN17.D	Calibration	Fluorobenzene	6.038	335741	836278	0.4015	264.6638	250.0000	105.9
04JAN19.D	Calibration	Fluorobenzene	6.040	507157	841876	0.6024	397.1322	375.0000	105.9
04JAN21.D	Calibration	Fluorobenzene	6.038	693669	841364	0.8245	543.5121	500.0000	108.7
04JAN23.D	QC	Fluorobenzene	6.038	150930	801210	0.1884	124.1853	125.0000	

Compound: 1,2-Dichloroethane-d4

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene	6.233	88174	775552	0.1137	279.3964		
04JAN10.D	Calibration	Fluorobenzene	6.233	923	770895	0.0012	2.9438	2.5000	117.8
04JAN11.D	Calibration	Fluorobenzene	6.227	3938	764419	0.0052	12.6600	12.5000	101.3
04JAN12.D	Calibration	Fluorobenzene	6.236	8284	791270	0.0105	25.7280	25.0000	102.9
04JAN13.D	Calibration	Fluorobenzene	6.233	15238	778120	0.0196	48.1252	50.0000	96.3
04JAN15.D	Calibration	Fluorobenzene	6.233	39086	823488	0.0475	116.6420	125.0000	93.3
04JAN17.D	Calibration	Fluorobenzene	6.236	87876	836278	0.1051	258.2324	250.0000	103.3
04JAN19.D	Calibration	Fluorobenzene	6.233	129608	841876	0.1540	378.3335	375.0000	100.9
04JAN21.D	Calibration	Fluorobenzene	6.233	174713	841364	0.2077	510.3080	500.0000	102.1
04JAN23.D	QC	Fluorobenzene	6.230	91382	801210	0.1141	280.2886	250.0000	

Compound: Benzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene	6.266	381	775552	0.0005	0.1233		
04JAN10.D	Calibration	Fluorobenzene	6.278	8408	770895	0.0109	2.7393	2.5000	109.6
04JAN11.D	Calibration	Fluorobenzene	6.278	37071	764419	0.0485	12.1801	12.5000	97.4

Compound: Benzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN12.D	Calibration	Fluorobenzene	6.278	74956	791270	0.0947	23.7919	25.0000	95.2
04JAN13.D	Calibration	Fluorobenzene	6.277	148727	778120	0.1911	48.0054	50.0000	96.0
04JAN15.D	Calibration	Fluorobenzene	6.278	383469	823488	0.4657	116.9553	125.0000	93.6
04JAN17.D	Calibration	Fluorobenzene	6.280	857534	836278	1.0254	257.5416	250.0000	103.0
04JAN19.D	Calibration	Fluorobenzene	6.278	1293370	841876	1.5363	385.8526	375.0000	102.9
04JAN21.D	Calibration	Fluorobenzene	6.280	1714050	841364	2.0372	511.6658	500.0000	102.3
04JAN23.D	QC	Fluorobenzene	6.280	418900	801210	0.5228	131.3139	125.0000	

Compound: 1,2-Dichloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Fluorobenzene			775552		ND		
04JAN10.D	Calibration	Fluorobenzene	6.322	2415	770895	0.0031	2.9090	2.5000	116.4
04JAN11.D	Calibration	Fluorobenzene	6.322	10202	764419	0.0133	12.3906	12.5000	99.1
04JAN12.D	Calibration	Fluorobenzene	6.322	19996	791270	0.0253	23.4616	25.0000	93.8
04JAN13.D	Calibration	Fluorobenzene	6.325	41058	778120	0.0528	48.9880	50.0000	98.0
04JAN15.D	Calibration	Fluorobenzene	6.322	104855	823488	0.1273	118.2143	125.0000	94.6
04JAN17.D	Calibration	Fluorobenzene	6.322	226964	836278	0.2714	251.9675	250.0000	100.8
04JAN19.D	Calibration	Fluorobenzene	6.322	332775	841876	0.3953	366.9787	375.0000	97.9
04JAN21.D	Calibration	Fluorobenzene	6.322	450739	841364	0.5357	497.3699	500.0000	99.5
04JAN23.D	QC	Fluorobenzene	6.325	104249	801210	0.1301	120.7991	125.0000	

Compound: Trichloroethene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	7.033	2372	296081	0.0080	2.6564	2.5000	106.3
04JAN11.D	Calibration	Chlorobenzene-d5	7.025	10442	296554	0.0352	11.6753	12.5000	93.4
04JAN12.D	Calibration	Chlorobenzene-d5	7.028	21946	301338	0.0728	24.1484	25.0000	96.6
04JAN13.D	Calibration	Chlorobenzene-d5	7.030	42682	300356	0.1421	47.1189	50.0000	94.2
04JAN15.D	Calibration	Chlorobenzene-d5	7.030	114123	306491	0.3724	123.4646	125.0000	98.8
04JAN17.D	Calibration	Chlorobenzene-d5	7.030	250285	316399	0.7910	262.2931	250.0000	104.9
04JAN19.D	Calibration	Chlorobenzene-d5	7.028	374370	314668	1.1897	394.4896	375.0000	105.2
04JAN21.D	Calibration	Chlorobenzene-d5	7.028	505400	313585	1.6117	534.4007	500.0000	106.9
04JAN23.D	QC	Chlorobenzene-d5	7.025	121734	307868	0.3954	131.1096	125.0000	

Compound: 1,2-Dichloropropane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	7.273	2148	296081	0.0073	2.7347	2.5000	109.4
04JAN11.D	Calibration	Chlorobenzene-d5	7.270	9488	296554	0.0320	12.0602	12.5000	96.5
04JAN12.D	Calibration	Chlorobenzene-d5	7.270	20077	301338	0.0666	25.1147	25.0000	100.5
04JAN13.D	Calibration	Chlorobenzene-d5	7.273	37870	300356	0.1261	47.5273	50.0000	95.1
04JAN15.D	Calibration	Chlorobenzene-d5	7.270	99187	306491	0.3236	121.9890	125.0000	97.6

Compound: 1,2-Dichloropropane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN17.D	Calibration	Chlorobenzene-d5	7.270	213800	316399	0.6757	254.7161	250.0000	101.9
04JAN19.D	Calibration	Chlorobenzene-d5	7.270	324602	314668	1.0316	388.8502	375.0000	103.7
04JAN21.D	Calibration	Chlorobenzene-d5	7.270	436057	313585	1.3906	524.1695	500.0000	104.8
04JAN23.D	QC	Chlorobenzene-d5	7.270	102633	307868	0.3334	125.6626	125.0000	

Compound: Dibromomethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	7.396	902	296081	0.0030	2.7162	2.5000	108.6
04JAN11.D	Calibration	Chlorobenzene-d5	7.399	4675	296554	0.0158	14.0619	12.5000	112.5
04JAN12.D	Calibration	Chlorobenzene-d5	7.393	8055	301338	0.0267	23.8439	25.0000	95.4
04JAN13.D	Calibration	Chlorobenzene-d5	7.396	15989	300356	0.0532	47.4844	50.0000	95.0
04JAN15.D	Calibration	Chlorobenzene-d5	7.399	40628	306491	0.1326	118.2425	125.0000	94.6
04JAN17.D	Calibration	Chlorobenzene-d5	7.396	89483	316399	0.2828	252.2734	250.0000	100.9
04JAN19.D	Calibration	Chlorobenzene-d5	7.396	134282	314668	0.4267	380.6547	375.0000	101.5
04JAN21.D	Calibration	Chlorobenzene-d5	7.396	176038	313585	0.5614	500.7456	500.0000	100.1
04JAN23.D	QC	Chlorobenzene-d5	7.393	43248	307868	0.1405	125.3047	125.0000	

Compound: Bromodichloromethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	7.597	2536	296081	0.0086	2.7684	2.5000	110.7
04JAN11.D	Calibration	Chlorobenzene-d5	7.585	11562	296554	0.0390	12.6014	12.5000	100.8
04JAN12.D	Calibration	Chlorobenzene-d5	7.583	22743	301338	0.0755	24.3940	25.0000	97.6
04JAN13.D	Calibration	Chlorobenzene-d5	7.585	43900	300356	0.1462	47.2409	50.0000	94.5
04JAN15.D	Calibration	Chlorobenzene-d5	7.585	115664	306491	0.3774	121.9749	125.0000	97.6
04JAN17.D	Calibration	Chlorobenzene-d5	7.582	251805	316399	0.7958	257.2286	250.0000	102.9
04JAN19.D	Calibration	Chlorobenzene-d5	7.585	375983	314668	1.1949	386.1940	375.0000	103.0
04JAN21.D	Calibration	Chlorobenzene-d5	7.585	502929	313585	1.6038	518.3718	500.0000	103.7
04JAN23.D	QC	Chlorobenzene-d5	7.585	122757	307868	0.3987	128.8759	125.0000	

Compound: cis-1,3-Dichloropropene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	8.054	2583	296081	0.0087	2.4939	2.5000	99.8
04JAN11.D	Calibration	Chlorobenzene-d5	8.062	12525	296554	0.0422	12.0738	12.5000	96.6
04JAN12.D	Calibration	Chlorobenzene-d5	8.057	24511	301338	0.0813	23.2528	25.0000	93.0
04JAN13.D	Calibration	Chlorobenzene-d5	8.057	48886	300356	0.1628	46.5283	50.0000	93.1
04JAN15.D	Calibration	Chlorobenzene-d5	8.059	129419	306491	0.4223	120.7116	125.0000	96.6
04JAN17.D	Calibration	Chlorobenzene-d5	8.057	293617	316399	0.9280	265.2863	250.0000	106.1
04JAN19.D	Calibration	Chlorobenzene-d5	8.057	441168	314668	1.4020	400.7930	375.0000	106.9
04JAN21.D	Calibration	Chlorobenzene-d5	8.059	591147	313585	1.8851	538.9008	500.0000	107.8

Compound: cis-1,3-Dichloropropene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN23.D	QC	Chlorobenzene-d5	8.054	130910	307868	0.4252	121.5561	125.0000	

Compound: Toluene-d8

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5	8.319	770154	301196	2.5570	265.3436		
04JAN10.D	Calibration	Chlorobenzene-d5	8.322	7777	296081	0.0263	2.7257	2.5000	109.0
04JAN11.D	Calibration	Chlorobenzene-d5	8.322	32318	296554	0.1090	11.3089	12.5000	90.5
04JAN12.D	Calibration	Chlorobenzene-d5	8.319	67673	301338	0.2246	23.3046	25.0000	93.2
04JAN13.D	Calibration	Chlorobenzene-d5	8.319	136453	300356	0.4543	47.1441	50.0000	94.3
04JAN15.D	Calibration	Chlorobenzene-d5	8.319	358186	306491	1.1687	121.2749	125.0000	97.0
04JAN17.D	Calibration	Chlorobenzene-d5	8.319	823306	316399	2.6021	270.0265	250.0000	108.0
04JAN19.D	Calibration	Chlorobenzene-d5	8.322	1229775	314668	3.9082	405.5583	375.0000	108.1
04JAN21.D	Calibration	Chlorobenzene-d5	8.319	1644540	313585	5.2443	544.2136	500.0000	108.8
04JAN23.D	QC	Chlorobenzene-d5	8.322	821531	307868	2.6685	276.9106	250.0000	

Compound: Toluene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	8.380	5039	296081	0.0170	2.6145	2.5000	104.6
04JAN11.D	Calibration	Chlorobenzene-d5	8.388	21794	296554	0.0735	11.2899	12.5000	90.3
04JAN12.D	Calibration	Chlorobenzene-d5	8.389	46355	301338	0.1538	23.6319	25.0000	94.5
04JAN13.D	Calibration	Chlorobenzene-d5	8.388	91915	300356	0.3060	47.0116	50.0000	94.0
04JAN15.D	Calibration	Chlorobenzene-d5	8.389	244712	306491	0.7984	122.6571	125.0000	98.1
04JAN17.D	Calibration	Chlorobenzene-d5	8.386	541945	316399	1.7129	263.1330	250.0000	105.3
04JAN19.D	Calibration	Chlorobenzene-d5	8.388	813204	314668	2.5843	397.0106	375.0000	105.9
04JAN21.D	Calibration	Chlorobenzene-d5	8.389	1095161	313585	3.4924	536.5101	500.0000	107.3
04JAN23.D	QC	Chlorobenzene-d5	8.386	264584	307868	0.8594	132.0244	125.0000	

Compound: trans-1,3-Dichloropropene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	8.634	1470	296081	0.0050	1.9942	2.5000	79.8
04JAN11.D	Calibration	Chlorobenzene-d5	8.645	8683	296554	0.0293	11.7589	12.5000	94.1
04JAN12.D	Calibration	Chlorobenzene-d5	8.634	17850	301338	0.0592	23.7894	25.0000	95.2
04JAN13.D	Calibration	Chlorobenzene-d5	8.639	35179	300356	0.1171	47.0378	50.0000	94.1
04JAN15.D	Calibration	Chlorobenzene-d5	8.637	92719	306491	0.3025	121.4929	125.0000	97.2
04JAN17.D	Calibration	Chlorobenzene-d5	8.639	207833	316399	0.6569	263.8027	250.0000	105.5
04JAN19.D	Calibration	Chlorobenzene-d5	8.639	315063	314668	1.0013	402.1098	375.0000	107.2
04JAN21.D	Calibration	Chlorobenzene-d5	8.637	416771	313585	1.3291	533.7551	500.0000	106.8
04JAN23.D	QC	Chlorobenzene-d5	8.637	98907	307868	0.3213	129.0216	125.0000	

Compound: 1,1,2-Trichloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	8.810	960	296081	0.0032	2.5012	2.5000	100.0
04JAN11.D	Calibration	Chlorobenzene-d5	8.824	5090	296554	0.0172	13.2340	12.5000	105.9
04JAN12.D	Calibration	Chlorobenzene-d5	8.815	10099	301338	0.0335	25.8400	25.0000	103.4
04JAN13.D	Calibration	Chlorobenzene-d5	8.815	18884	300356	0.0629	48.4759	50.0000	97.0
04JAN15.D	Calibration	Chlorobenzene-d5	8.818	46673	306491	0.1523	117.4130	125.0000	93.9
04JAN17.D	Calibration	Chlorobenzene-d5	8.815	101888	316399	0.3220	248.2882	250.0000	99.3
04JAN19.D	Calibration	Chlorobenzene-d5	8.818	152331	314668	0.4841	373.2534	375.0000	99.5
04JAN21.D	Calibration	Chlorobenzene-d5	8.815	205463	313585	0.6552	505.1803	500.0000	101.0
04JAN23.D	QC	Chlorobenzene-d5	8.815	49128	307868	0.1596	123.0361	125.0000	

Compound: Tetrachloroethene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	8.932	2105	296081	0.0071	2.6772	2.5000	107.1
04JAN11.D	Calibration	Chlorobenzene-d5	8.935	9238	296554	0.0312	11.7302	12.5000	93.8
04JAN12.D	Calibration	Chlorobenzene-d5	8.935	20322	301338	0.0674	25.3948	25.0000	101.6
04JAN13.D	Calibration	Chlorobenzene-d5	8.935	36925	300356	0.1229	46.2932	50.0000	92.6
04JAN15.D	Calibration	Chlorobenzene-d5	8.935	97590	306491	0.3184	119.9003	125.0000	95.9
04JAN17.D	Calibration	Chlorobenzene-d5	8.938	218245	316399	0.6898	259.7419	250.0000	103.9
04JAN19.D	Calibration	Chlorobenzene-d5	8.938	319950	314668	1.0168	382.8796	375.0000	102.1
04JAN21.D	Calibration	Chlorobenzene-d5	8.938	428812	313585	1.3675	514.9255	500.0000	103.0
04JAN23.D	QC	Chlorobenzene-d5	8.935	103027	307868	0.3346	126.0141	125.0000	

Compound: 1,3-Dichloropropane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	8.977	2257	296081	0.0076	2.9881	2.5000	119.5
04JAN11.D	Calibration	Chlorobenzene-d5	8.985	8967	296554	0.0302	11.8526	12.5000	94.8
04JAN12.D	Calibration	Chlorobenzene-d5	8.983	18745	301338	0.0622	24.3839	25.0000	97.5
04JAN13.D	Calibration	Chlorobenzene-d5	8.980	37457	300356	0.1247	48.8841	50.0000	97.8
04JAN15.D	Calibration	Chlorobenzene-d5	8.980	96183	306491	0.3138	123.0132	125.0000	98.4
04JAN17.D	Calibration	Chlorobenzene-d5	8.980	212669	316399	0.6722	263.4754	250.0000	105.4
04JAN19.D	Calibration	Chlorobenzene-d5	8.980	312547	314668	0.9933	389.3442	375.0000	103.8
04JAN21.D	Calibration	Chlorobenzene-d5	8.980	408993	313585	1.3042	511.2479	500.0000	102.2
04JAN23.D	QC	Chlorobenzene-d5	8.980	95697	307868	0.3108	121.8442	125.0000	

Compound: Chlorodibromomethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	9.203	1468	296081	0.0050	2.4461	2.5000	97.8
04JAN11.D	Calibration	Chlorobenzene-d5	9.206	7718	296554	0.0260	12.8393	12.5000	102.7

Compound: Chlorodibromomethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN12.D	Calibration	Chlorobenzene-d5	9.197	14873	301338	0.0494	24.3492	25.0000	97.4
04JAN13.D	Calibration	Chlorobenzene-d5	9.203	28153	300356	0.0937	46.2411	50.0000	92.5
04JAN15.D	Calibration	Chlorobenzene-d5	9.206	75015	306491	0.2448	120.7454	125.0000	96.6
04JAN17.D	Calibration	Chlorobenzene-d5	9.203	165695	316399	0.5237	258.3535	250.0000	103.3
04JAN19.D	Calibration	Chlorobenzene-d5	9.203	247279	314668	0.7858	387.6812	375.0000	103.4
04JAN21.D	Calibration	Chlorobenzene-d5	9.203	330813	313585	1.0549	520.4361	500.0000	104.1
04JAN23.D	QC	Chlorobenzene-d5	9.203	78076	307868	0.2536	125.1103	125.0000	

Compound: 1,2-Dibromoethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	9.300	1299	296081	0.0044	3.0943	2.5000	123.8
04JAN11.D	Calibration	Chlorobenzene-d5	9.300	5410	296554	0.0182	12.8640	12.5000	102.9
04JAN12.D	Calibration	Chlorobenzene-d5	9.309	10410	301338	0.0345	24.3601	25.0000	97.4
04JAN13.D	Calibration	Chlorobenzene-d5	9.303	21037	300356	0.0700	49.3889	50.0000	98.8
04JAN15.D	Calibration	Chlorobenzene-d5	9.306	51827	306491	0.1691	119.2394	125.0000	95.4
04JAN17.D	Calibration	Chlorobenzene-d5	9.306	115714	316399	0.3657	257.8887	250.0000	103.2
04JAN19.D	Calibration	Chlorobenzene-d5	9.306	168577	314668	0.5357	377.7698	375.0000	100.7
04JAN21.D	Calibration	Chlorobenzene-d5	9.303	225877	313585	0.7203	507.9234	500.0000	101.6
04JAN23.D	QC	Chlorobenzene-d5	9.306	54259	307868	0.1762	124.2764	125.0000	

Compound: Chlorobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	9.805	5771	296081	0.0195	2.7350	2.5000	109.4
04JAN11.D	Calibration	Chlorobenzene-d5	9.802	26461	296554	0.0892	12.5204	12.5000	100.2
04JAN12.D	Calibration	Chlorobenzene-d5	9.802	53047	301338	0.1760	24.7015	25.0000	98.8
04JAN13.D	Calibration	Chlorobenzene-d5	9.802	101452	300356	0.3378	47.3959	50.0000	94.8
04JAN15.D	Calibration	Chlorobenzene-d5	9.802	263617	306491	0.8601	120.6903	125.0000	96.6
04JAN17.D	Calibration	Chlorobenzene-d5	9.802	582326	316399	1.8405	258.2544	250.0000	103.3
04JAN19.D	Calibration	Chlorobenzene-d5	9.802	867732	314668	2.7576	386.9455	375.0000	103.2
04JAN21.D	Calibration	Chlorobenzene-d5	9.802	1153147	313585	3.6773	515.9957	500.0000	103.2
04JAN23.D	QC	Chlorobenzene-d5	9.802	288815	307868	0.9381	131.6352	125.0000	

Compound: 1,1,1,2-Tetrachloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	9.889	1893	296081	0.0064	2.5659	2.5000	102.6
04JAN11.D	Calibration	Chlorobenzene-d5	9.889	9473	296554	0.0319	12.8225	12.5000	102.6
04JAN12.D	Calibration	Chlorobenzene-d5	9.889	18130	301338	0.0602	24.1509	25.0000	96.6
04JAN13.D	Calibration	Chlorobenzene-d5	9.889	35544	300356	0.1183	47.5029	50.0000	95.0
04JAN15.D	Calibration	Chlorobenzene-d5	9.889	90898	306491	0.2966	119.0492	125.0000	95.2

Compound: 1,1,1,2-Tetrachloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN17.D	Calibration	Chlorobenzene-d5	9.891	200859	316399	0.6348	254.8274	250.0000	101.9
04JAN19.D	Calibration	Chlorobenzene-d5	9.892	307436	314668	0.9770	392.1859	375.0000	104.6
04JAN21.D	Calibration	Chlorobenzene-d5	9.892	406450	313585	1.2961	520.2855	500.0000	104.1
04JAN23.D	QC	Chlorobenzene-d5	9.892	97148	307868	0.3156	126.6657	125.0000	

Compound: Ethylbenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	9.920	9283	296081	0.0314	2.5367	2.5000	101.5
04JAN11.D	Calibration	Chlorobenzene-d5	9.917	40470	296554	0.1365	11.0411	12.5000	88.3
04JAN12.D	Calibration	Chlorobenzene-d5	9.917	88428	301338	0.2935	23.7421	25.0000	95.0
04JAN13.D	Calibration	Chlorobenzene-d5	9.917	173769	300356	0.5785	46.8079	50.0000	93.6
04JAN15.D	Calibration	Chlorobenzene-d5	9.920	464148	306491	1.5144	122.5243	125.0000	98.0
04JAN17.D	Calibration	Chlorobenzene-d5	9.919	1043443	316399	3.2979	266.8193	250.0000	106.7
04JAN19.D	Calibration	Chlorobenzene-d5	9.919	1574219	314668	5.0028	404.7587	375.0000	107.9
04JAN21.D	Calibration	Chlorobenzene-d5	9.919	2111152	313585	6.7323	544.6881	500.0000	108.9
04JAN23.D	QC	Chlorobenzene-d5	9.917	501953	307868	1.6304	131.9113	125.0000	

Compound: m+p-Xylenes

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	10.045	7212	296081	0.0244	5.0712	5.0000	101.4
04JAN11.D	Calibration	Chlorobenzene-d5	10.037	31538	296554	0.1063	22.1410	25.0000	88.6
04JAN12.D	Calibration	Chlorobenzene-d5	10.039	66267	301338	0.2199	45.7836	50.0000	91.6
04JAN13.D	Calibration	Chlorobenzene-d5	10.039	133498	300356	0.4445	92.5347	100.0000	92.5
04JAN15.D	Calibration	Chlorobenzene-d5	10.039	368418	306491	1.2021	250.2587	250.0000	100.1
04JAN17.D	Calibration	Chlorobenzene-d5	10.039	825866	316399	2.6102	543.4262	500.0000	108.7
04JAN19.D	Calibration	Chlorobenzene-d5	10.039	1228570	314668	3.9043	812.8556	750.0000	108.4
04JAN21.D	Calibration	Chlorobenzene-d5	10.039	1637879	313585	5.2231	1087.4082	1000.0000	108.7
04JAN23.D	QC	Chlorobenzene-d5	10.039	388558	307868	1.2621	262.7589	250.0000	

Compound: o-Xylene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	10.430	3330	296081	0.0112	2.6303	2.5000	105.2
04JAN11.D	Calibration	Chlorobenzene-d5	10.430	13519	296554	0.0456	10.6612	12.5000	85.3
04JAN12.D	Calibration	Chlorobenzene-d5	10.427	30463	301338	0.1011	23.6420	25.0000	94.6
04JAN13.D	Calibration	Chlorobenzene-d5	10.430	61016	300356	0.2031	47.5086	50.0000	95.0
04JAN15.D	Calibration	Chlorobenzene-d5	10.430	161509	306491	0.5270	123.2378	125.0000	98.6
04JAN17.D	Calibration	Chlorobenzene-d5	10.430	365914	316399	1.1565	270.4636	250.0000	108.2
04JAN19.D	Calibration	Chlorobenzene-d5	10.433	549244	314668	1.7455	408.2043	375.0000	108.9
04JAN21.D	Calibration	Chlorobenzene-d5	10.430	734101	313585	2.3410	547.4764	500.0000	109.5

Compound: o-Xylene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN23.D	QC	Chlorobenzene-d5	10.430	174061	307868	0.5654	132.2214	125.0000	

Compound: Styrene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	Chlorobenzene-d5			301196		ND		
04JAN10.D	Calibration	Chlorobenzene-d5	10.444	4408	296081	0.0149	2.1625	2.5000	86.5
04JAN11.D	Calibration	Chlorobenzene-d5	10.449	23472	296554	0.0791	11.4968	12.5000	92.0
04JAN12.D	Calibration	Chlorobenzene-d5	10.447	48569	301338	0.1612	23.4119	25.0000	93.6
04JAN13.D	Calibration	Chlorobenzene-d5	10.444	96576	300356	0.3215	46.7052	50.0000	93.4
04JAN15.D	Calibration	Chlorobenzene-d5	10.447	268375	306491	0.8756	127.1910	125.0000	101.8
04JAN17.D	Calibration	Chlorobenzene-d5	10.446	605646	316399	1.9142	278.0455	250.0000	111.2
04JAN19.D	Calibration	Chlorobenzene-d5	10.447	896331	314668	2.8485	413.7595	375.0000	110.3
04JAN21.D	Calibration	Chlorobenzene-d5	10.449	1199879	313585	3.8263	555.7946	500.0000	111.2
04JAN23.D	QC	Chlorobenzene-d5	10.449	291425	307868	0.9466	137.4974	125.0000	

Compound: Bromoform

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	10.625	708	227879	0.0031	2.4287	2.5000	97.1
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	10.625	3652	242142	0.0151	11.7860	12.5000	94.3
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	10.625	7972	240335	0.0332	25.9212	25.0000	103.7
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	10.625	16073	248636	0.0646	50.5170	50.0000	101.0
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	10.628	39165	264477	0.1481	115.7218	125.0000	92.6
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	10.628	87836	266553	0.3295	257.5099	250.0000	103.0
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	10.628	129038	266611	0.4840	378.2200	375.0000	100.9
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	10.625	175918	262971	0.6690	522.7660	500.0000	104.6
04JAN23.D	QC	1,4-Dichlorobenzene-d4	10.628	42560	255907	0.1663	129.9644	125.0000	

Compound: p-Bromofluorobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4	10.951	226743	231562	0.9792	267.2815		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	10.951	2719	227879	0.0119	3.2569	2.5000	130.3
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	10.948	10059	242142	0.0415	11.3393	12.5000	90.7
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	10.951	22267	240335	0.0926	25.2899	25.0000	101.2
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	10.951	42506	248636	0.1710	46.6647	50.0000	93.3
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	10.954	114269	264477	0.4321	117.9350	125.0000	94.3
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	10.951	261042	266553	0.9793	267.3186	250.0000	106.9
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	10.951	385474	266611	1.4458	394.6566	375.0000	105.2
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	10.949	521580	262971	1.9834	541.3964	500.0000	108.3
04JAN23.D	QC	1,4-Dichlorobenzene-d4	10.951	253034	255907	0.9888	269.8976	250.0000	

Compound: Bromobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	11.088	2024	227879	0.0089	2.7439	2.5000	109.8
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	11.096	9663	242142	0.0399	12.3310	12.5000	98.6
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	11.094	19259	240335	0.0801	24.7613	25.0000	99.0
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	11.093	38282	248636	0.1540	47.5759	50.0000	95.2
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	11.094	102265	264477	0.3867	119.4801	125.0000	95.6
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	11.093	227127	266553	0.8521	263.2944	250.0000	105.3
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	11.093	333431	266611	1.2506	386.4420	375.0000	103.1
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	11.094	439147	262971	1.6699	516.0104	500.0000	103.2
04JAN23.D	QC	1,4-Dichlorobenzene-d4	11.093	109054	255907	0.4261	131.6788	125.0000	

Compound: 1,1,2,2-Tetrachloroethane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	11.113	1142	227879	0.0050	2.6916	2.5000	107.7
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	11.116	5793	242142	0.0239	12.8437	12.5000	102.7
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	11.110	12440	240335	0.0518	27.7883	25.0000	111.2
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	11.105	22514	248636	0.0906	48.6124	50.0000	97.2
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	11.116	56958	264477	0.2154	115.6179	125.0000	92.5
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	11.113	124205	266553	0.4660	250.1577	250.0000	100.1
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	11.110	182470	266611	0.6844	367.4276	375.0000	98.0
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	11.113	240837	262971	0.9158	491.6700	500.0000	98.3
04JAN23.D	QC	1,4-Dichlorobenzene-d4	11.113	60763	255907	0.2374	127.4722	125.0000	

Compound: 1,2,3-Trichloropropane

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4			227879		ND	2.5000	
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	11.144	1654	242142	0.0068	13.7084	12.5000	109.7
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	11.149	3200	240335	0.0133	26.7144	25.0000	106.9
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	11.146	6096	248636	0.0245	49.1924	50.0000	98.4
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	11.147	14846	264477	0.0561	112.6261	125.0000	90.1
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	11.152	33115	266553	0.1242	249.2635	250.0000	99.7
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	11.149	48325	266611	0.1813	363.6732	375.0000	97.0
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	11.146	64422	262971	0.2450	491.5229	500.0000	98.3
04JAN23.D	QC	1,4-Dichlorobenzene-d4	11.146	15682	255907	0.0613	122.9523	125.0000	

Compound: 2-Chlorotoluene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	11.292	1844	227879	0.0081	2.5124	2.5000	100.5
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	11.289	8731	242142	0.0361	11.1977	12.5000	89.6

Compound: 2-Chlorotoluene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	11.286	19390	240335	0.0807	25.0550	25.0000	100.2
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	11.289	37987	248636	0.1528	47.4466	50.0000	94.9
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	11.292	102424	264477	0.3873	120.2675	125.0000	96.2
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	11.291	229396	266553	0.8606	267.2616	250.0000	106.9
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	11.291	336386	266611	1.2617	391.8269	375.0000	104.5
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	11.292	455991	262971	1.7340	538.4964	500.0000	107.7
04JAN23.D	QC	1,4-Dichlorobenzene-d4	11.291	108192	255907	0.4228	131.2948	125.0000	

Compound: 4-Chlorotoluene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	11.400	5419	227879	0.0238	2.2650	2.5000	90.6
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	11.400	28532	242142	0.1178	11.2233	12.5000	89.8
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	11.400	61551	240335	0.2561	24.3936	25.0000	97.6
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	11.400	126308	248636	0.5080	48.3865	50.0000	96.8
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	11.400	336146	264477	1.2710	121.0591	125.0000	96.8
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	11.400	748435	266553	2.8078	267.4409	250.0000	107.0
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	11.400	1109221	266611	4.1604	396.2756	375.0000	105.7
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	11.400	1468376	262971	5.5838	531.8471	500.0000	106.4
04JAN23.D	QC	1,4-Dichlorobenzene-d4	11.400	368295	255907	1.4392	137.0790	125.0000	

Compound: 1,3-Dichlorobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	12.033	3541	227879	0.0155	2.6327	2.5000	105.3
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	12.036	16932	242142	0.0699	11.8473	12.5000	94.8
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	12.033	36559	240335	0.1521	25.7725	25.0000	103.1
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	12.033	69539	248636	0.2797	47.3853	50.0000	94.8
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	12.031	183404	264477	0.6935	117.4899	125.0000	94.0
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	12.033	406895	266553	1.5265	258.6297	250.0000	103.5
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	12.033	603674	266611	2.2643	383.6225	375.0000	102.3
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	12.033	793993	262971	3.0193	511.5504	500.0000	102.3
04JAN23.D	QC	1,4-Dichlorobenzene-d4	12.033	204088	255907	0.7975	135.1185	125.0000	

Compound: 1,4-Dichlorobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	12.125	3787	227879	0.0166	2.7613	2.5000	110.5
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	12.122	17438	242142	0.0720	11.9662	12.5000	95.7
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	12.125	36635	240335	0.1524	25.3284	25.0000	101.3
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	12.125	71841	248636	0.2889	48.0106	50.0000	96.0
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	12.125	189045	264477	0.7148	118.7699	125.0000	95.0

Compound: 1,4-Dichlorobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	12.122	408934	266553	1.5342	254.9170	250.0000	102.0
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	12.125	595919	266611	2.2352	371.3969	375.0000	99.0
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	12.125	794954	262971	3.0230	502.3001	500.0000	100.5
04JAN23.D	QC	1,4-Dichlorobenzene-d4	12.122	200032	255907	0.7817	129.8812	125.0000	

Compound: 1,2-Dichlorobenzene

Data File	Sample Type	ISTD	RT	Resp	ISTD Resp	Resp Ratio	Final Conc	Exp. Conc	Accuracy
04JAN09.D	Blank	1,4-Dichlorobenzene-d4			231562		ND		
04JAN10.D	Calibration	1,4-Dichlorobenzene-d4	12.499	3104	227879	0.0136	2.7307	2.5000	109.2
04JAN11.D	Calibration	1,4-Dichlorobenzene-d4	12.493	14666	242142	0.0606	12.1423	12.5000	97.1
04JAN12.D	Calibration	1,4-Dichlorobenzene-d4	12.488	29899	240335	0.1244	24.9402	25.0000	99.8
04JAN13.D	Calibration	1,4-Dichlorobenzene-d4	12.491	60213	248636	0.2422	48.5498	50.0000	97.1
04JAN15.D	Calibration	1,4-Dichlorobenzene-d4	12.493	152284	264477	0.5758	115.4323	125.0000	92.3
04JAN17.D	Calibration	1,4-Dichlorobenzene-d4	12.493	342576	266553	1.2852	257.6524	250.0000	103.1
04JAN19.D	Calibration	1,4-Dichlorobenzene-d4	12.493	499147	266611	1.8722	375.3283	375.0000	100.1
04JAN21.D	Calibration	1,4-Dichlorobenzene-d4	12.493	664247	262971	2.5259	506.3871	500.0000	101.3
04JAN23.D	QC	1,4-Dichlorobenzene-d4	12.493	164299	255907	0.6420	128.7104	125.0000	

Initial Calibration Report - VOA5975C

Method Path	\\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL
Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m
Batch Name	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin
Last Calib Update	1/9/2022 8:59:52 PM

Level Name	Calibration Files	Acq. Date-Time	Level Last Update Time
1	D:\Org\Data\VOA5975C\VG010422\04JAN10.D	1/4/2022 3:33:04 PM	1/5/2022 11:05:51 AM
2	D:\Org\Data\VOA5975C\VG010422\04JAN11.D	1/4/2022 4:00:35 PM	1/5/2022 11:05:51 AM
3	D:\Org\Data\VOA5975C\VG010422\04JAN12.D	1/4/2022 4:28:05 PM	1/5/2022 11:05:51 AM
4	D:\Org\Data\VOA5975C\VG010422\04JAN13.D	1/4/2022 4:55:32 PM	1/5/2022 11:05:51 AM
5	D:\Org\Data\VOA5975C\VG010422\04JAN15.D	1/4/2022 5:50:25 PM	1/5/2022 11:05:51 AM
6	D:\Org\Data\VOA5975C\VG010422\04JAN17.D	1/4/2022 6:45:10 PM	1/5/2022 11:05:51 AM
7	D:\Org\Data\VOA5975C\VG010422\04JAN19.D	1/4/2022 7:39:45 PM	1/5/2022 11:05:51 AM
8	D:\Org\Data\VOA5975C\VG010422\04JAN21.D	1/4/2022 8:34:31 PM	1/5/2022 11:05:51 AM

Compound	Curve Fit	1	2	3	4	5	6	7	8	Avg RF	%RSD
I Fluorobenzene						ISTD					
T Dichlorodifluoromethane	Avg RF		0.3162	0.3365	0.3242	0.3350	0.3304	0.3267	0.3242	0.3276	2.141
T Chloromethane	Avg RF		0.4411	0.4190	0.3960	0.3901	0.3821	0.3733	0.3819	0.3976	6.061
T Vinyl chloride	Avg RF		0.3591	0.3671	0.3503	0.3603	0.3559	0.3553	0.3566	0.3578	1.449
T Bromomethane	Avg RF		0.1542	0.1650	0.1523	0.1583	0.1611	0.1643	0.1648	0.1600	3.267
T Chloroethane	Avg RF		0.2107	0.1851	0.1638	0.1735	0.1642	0.1721	0.1706	0.1771	9.265
T Trichlorofluoromethane	Avg RF		0.4037	0.4735	0.4380	0.4586	0.4602	0.4399	0.4349	0.4441	5.118
T 1,1-Dichloroethene	Avg RF		0.2399	0.2607	0.2458	0.2415	0.2600	0.2554	0.2594	0.2518	3.629
T Methylene chloride	Avg RF		0.4640	0.3906	0.3745	0.3285	0.3496	0.3446	0.3467	0.3712	12.340
T trans-1,2-Dichloroethene	Avg RF		0.2570	0.2617	0.2544	0.2439	0.2617	0.2577	0.2621	0.2569	2.508
T Methyl tert-butyl ether (MTBE)	Avg RF		0.3274	0.3061	0.3157	0.3378	0.3440	0.3464	0.3472	0.3321	4.880
T 1,1-Dichloroethane	Avg RF		0.4616	0.4913	0.4704	0.4519	0.4943	0.4852	0.4929	0.4782	3.547
T 2,2-Dichloropropane	Avg RF		0.3578	0.3765	0.3611	0.3392	0.3627	0.3534	0.3576	0.3583	3.118
T cis-1,2-Dichloroethene	Avg RF		0.2618	0.2559	0.2522	0.2430	0.2728	0.2686	0.2688	0.2605	4.101
T Methyl ethyl ketone	Avg RF		0.0344	0.0332	0.0338	0.0327	0.0379	0.0373	0.0376	0.0353 #	6.353
T Bromochloromethane	Avg RF		0.1118	0.1098	0.1114	0.1019	0.1066	0.1070	0.1067	0.1079	3.188
T Chloroform	Avg RF	0.5510	0.4975	0.4602	0.4588	0.4363	0.4723	0.4657	0.4656	0.4759	7.299
T 1,1,1-Trichloroethane	Avg RF		0.4349	0.4492	0.4306	0.4231	0.4616	0.4599	0.4628	0.4460	3.677
S Dibromofluoromethane	Avg RF		0.2374	0.2414	0.2269	0.2169	0.2440	0.2416	0.2404	0.2355	4.222
T Carbon tetrachloride	Avg RF		0.4308	0.4355	0.4197	0.4200	0.4586	0.4534	0.4581	0.4394	3.906
T 1,1-Dichloropropene	Avg RF		0.3440	0.3695	0.3623	0.3635	0.4015	0.4016	0.4122	0.3792	6.770
S 1,2-Dichloroethane-d4	Avg RF		0.1030	0.1047	0.0979	0.0949	0.1051	0.1026	0.1038	0.1017	3.759
T Benzene	Avg RF	1.0907	0.9699	0.9473	0.9557	0.9313	1.0254	1.0242	1.0186	0.9954	5.369
T 1,2-Dichloroethane	Avg RF	0.3133	0.2669	0.2527	0.2638	0.2547	0.2714	0.2635	0.2679	0.2693	7.024
I Chlorobenzene-d5						ISTD					
T Trichloroethene	Avg RF		0.7042	0.7283	0.7105	0.7447	0.7910	0.7932	0.8058	0.7540	5.603
T 1,2-Dichloropropane	Avg RF		0.6399	0.6663	0.6304	0.6472	0.6757	0.6877	0.6953	0.6632	3.729
T Dibromomethane	Avg RF		0.3153	0.2673	0.2662	0.2651	0.2828	0.2845	0.2807	0.2803	6.261
T Bromodichloromethane	Avg RF		0.7798	0.7547	0.7308	0.7548	0.7958	0.7966	0.8019	0.7735	3.503

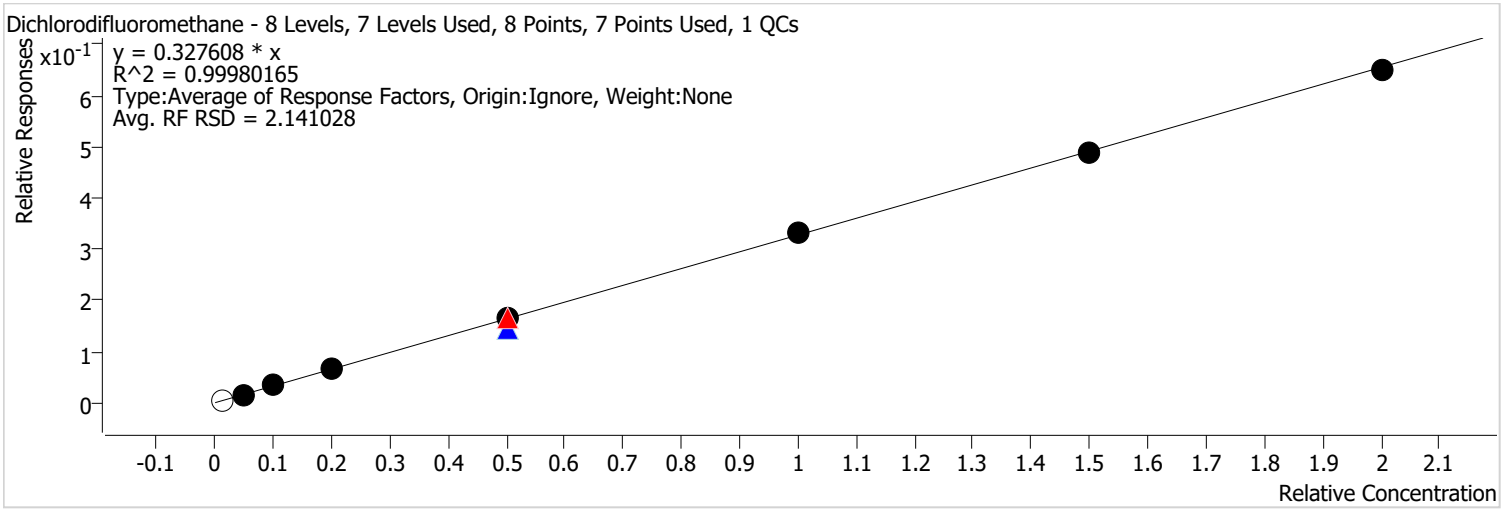
Compound	Curve Fit	1	2	3	4	5	6	7	8	Avg RF	%RSD
T cis-1,3-Dichloropropene	Avg RF		0.8447	0.8134	0.8138	0.8445	0.9280	0.9347	0.9426	0.8745	6.654
S Toluene-d8	Avg RF		2.1796	2.2458	2.2715	2.3373	2.6021	2.6054	2.6222	2.4091	8.032
T Toluene	Avg RF	1.7019	1.4698	1.5383	1.5301	1.5969	1.7129	1.7229	1.7462	1.6274	6.538
T trans-1,3-Dichloropropene	Avg RF		0.5856	0.5924	0.5856	0.6050	0.6569	0.6675	0.6645	0.6225	6.190
T 1,1,2-Trichloroethane	Avg RF		0.3433	0.3351	0.3144	0.3046	0.3220	0.3227	0.3276	0.3242	3.951
T Tetrachloroethene	Avg RF	0.7110	0.6230	0.6744	0.6147	0.6368	0.6898	0.6779	0.6837	0.6639	5.221
T 1,3-Dichloropropane	Avg RF		0.6047	0.6221	0.6235	0.6276	0.6722	0.6622	0.6521	0.6378	3.855
T Chlorodibromomethane	Avg RF		0.5205	0.4936	0.4687	0.4895	0.5237	0.5239	0.5275	0.5068	4.501
T 1,2-Dibromoethane	Avg RF		0.3649	0.3455	0.3502	0.3382	0.3657	0.3572	0.3602	0.3545	2.909
T Chlorobenzene	Avg RF		1.7846	1.7604	1.6889	1.7202	1.8405	1.8384	1.8387	1.7817	3.458
T 1,1,1,2-Tetrachloroethane	Avg RF		0.6389	0.6016	0.5917	0.5932	0.6348	0.6513	0.6481	0.6228	4.223
T Ethylbenzene	Avg RF	3.1353	2.7294	2.9345	2.8927	3.0288	3.2979	3.3352	3.3662	3.0900	7.526
T m+p-Xylenes	Avg RF	1.2179	1.0635	1.0995	1.1112	1.2021	1.3051	1.3014	1.3058	1.2008	8.296
T o-Xylene	Avg RF		0.9117	1.0109	1.0157	1.0539	1.1565	1.1636	1.1705	1.0690	9.204
T Styrene	Avg RF	1.4888	1.5830	1.6118	1.6077	1.7513	1.9142	1.8990	1.9132	1.7211	9.933
I 1,4-Dichlorobenzene-d4				---------- ISTD ----------							
T Bromoform	Avg RF		0.3016	0.3317	0.3232	0.2962	0.3295	0.3227	0.3345	0.3199	4.706
S p-Bromofluorobenzene	Avg RF		0.8308	0.9265	0.8548	0.8641	0.9793	0.9639	0.9917	0.9159	7.165
T Bromobenzene	Avg RF		0.7981	0.8013	0.7698	0.7733	0.8521	0.8338	0.8350	0.8091	3.949
T 1,1,2,2-Tetrachloroethane	Avg RF		0.4785	0.5176	0.4528	0.4307	0.4660	0.4563	0.4579	0.4657	5.814
T 1,2,3-Trichloropropane	Avg RF		0.1366	0.1331	0.1226	0.1123	0.1242	0.1208	0.1225	0.1246	6.496
T 2-Chlorotoluene	Avg RF		0.7211	0.8068	0.7639	0.7745	0.8606	0.8411	0.8670	0.8050	6.783
T 4-Chlorotoluene	Avg RF		2.3566	2.5611	2.5400	2.5420	2.8078	2.7736	2.7919	2.6247	6.481
T 1,3-Dichlorobenzene	Avg RF	1.5539	1.3985	1.5212	1.3984	1.3869	1.5265	1.5095	1.5097	1.4756	4.644
T 1,4-Dichlorobenzene	Avg RF	1.6618	1.4403	1.5243	1.4447	1.4296	1.5342	1.4901	1.5115	1.5046	4.999
T 1,2-Dichlorobenzene	Avg RF	1.3621	1.2114	1.2441	1.2109	1.1516	1.2852	1.2481	1.2630	1.2470	4.949

(RedFont and #) = Outlier Flag; (I) = Internal Standard; (T) = Target; (S) = Surrogate; (M) = Matrix Spike

Calibration Report

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:39 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Dichlorodifluoromethane %RSE = 2.1



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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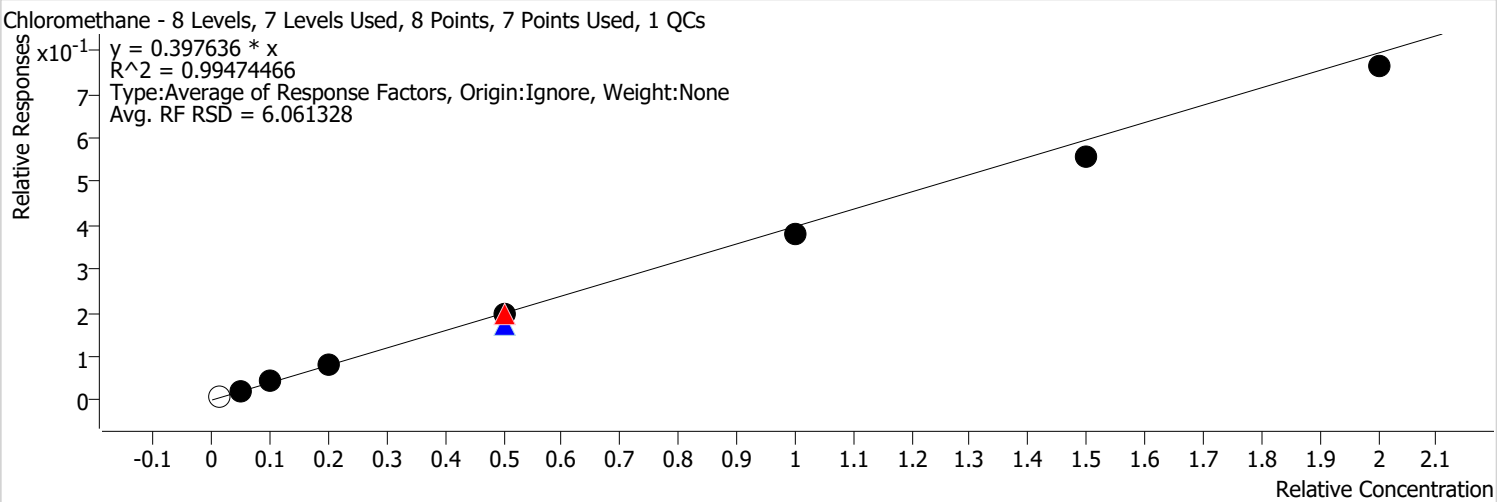
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	12087	12.5000	0.3162	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	26627	25.0000	0.3365	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	50457	50.0000	0.3242	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	116936	125.0000	0.2919	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	137933	125.0000	0.3350	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	137933	125.0000	0.3350	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	276334	250.0000	0.3304	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	412544	375.0000	0.3267	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	545484	500.0000	0.3242	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:42 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Chloromethane %RSE = 6.1



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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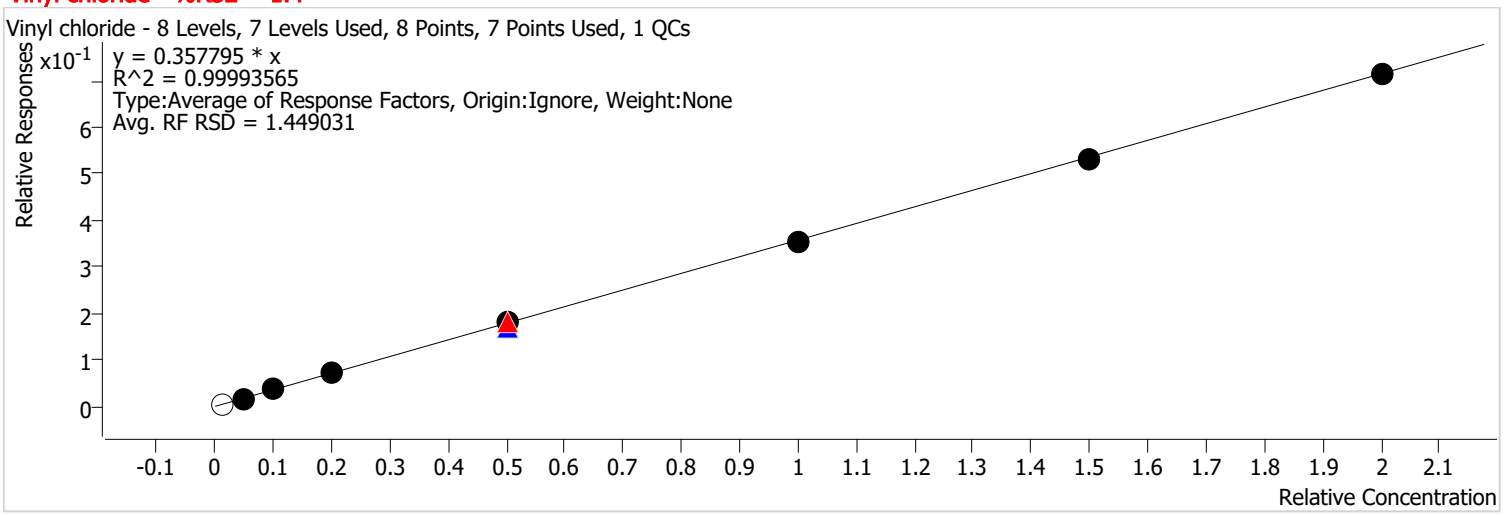
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	16859	12.5000	0.4411	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	33153	25.0000	0.4190	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	61632	50.0000	0.3960	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	138617	125.0000	0.3460	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	160604	125.0000	0.3901	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	160604	125.0000	0.3901	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	319523	250.0000	0.3821	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	471454	375.0000	0.3733	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	642582	500.0000	0.3819	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:42 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Vinyl chloride %RSE = 1.4



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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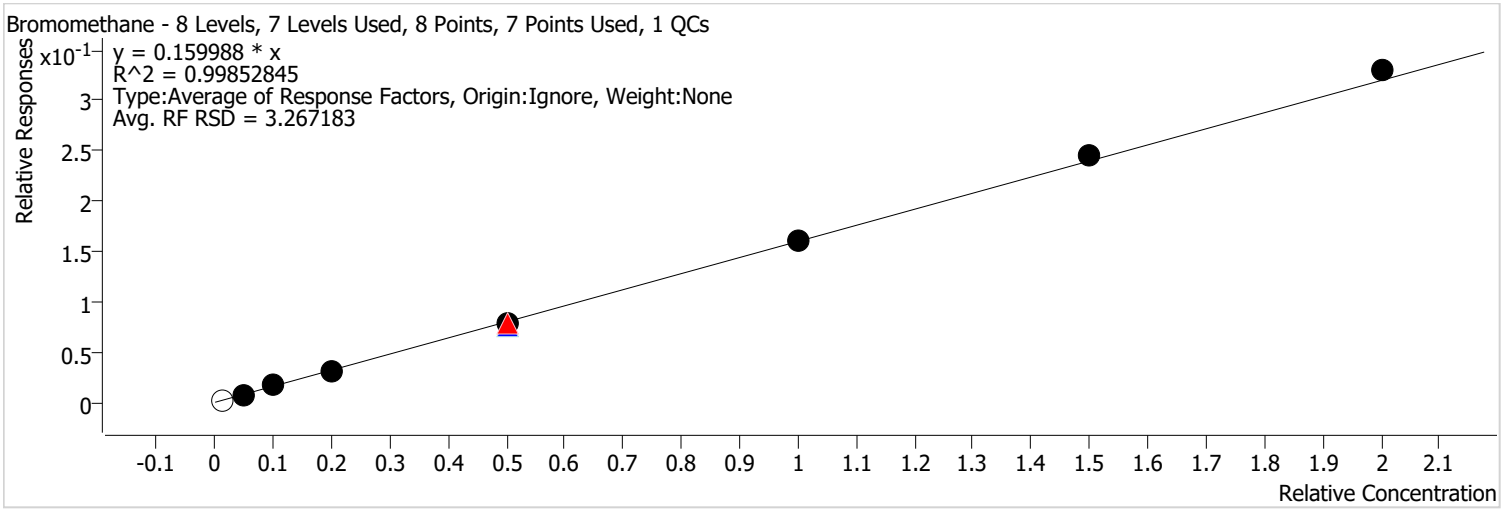
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	29046	25.0000	0.3671	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	54521	50.0000	0.3503	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	137775	125.0000	0.3439	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	148358	125.0000	0.3603	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	148358	125.0000	0.3603	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	297604	250.0000	0.3559	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	448643	375.0000	0.3553	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	600092	500.0000	0.3566	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:42 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Bromomethane %RSE = 3.3



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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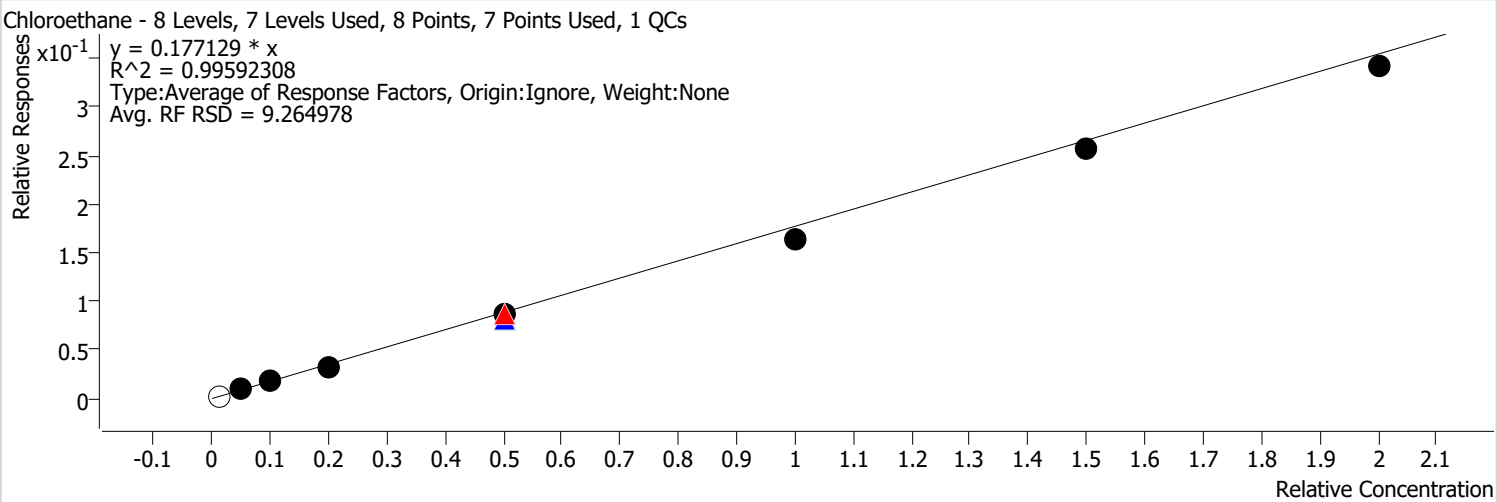
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	5893	12.5000	0.1542	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	13054	25.0000	0.1650	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	23699	50.0000	0.1523	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	59947	125.0000	0.1496	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	65163	125.0000	0.1583	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	65163	125.0000	0.1583	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	134737	250.0000	0.1611	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	207491	375.0000	0.1643	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	277301	500.0000	0.1648	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:42 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Chloroethane %RSE = 9.3



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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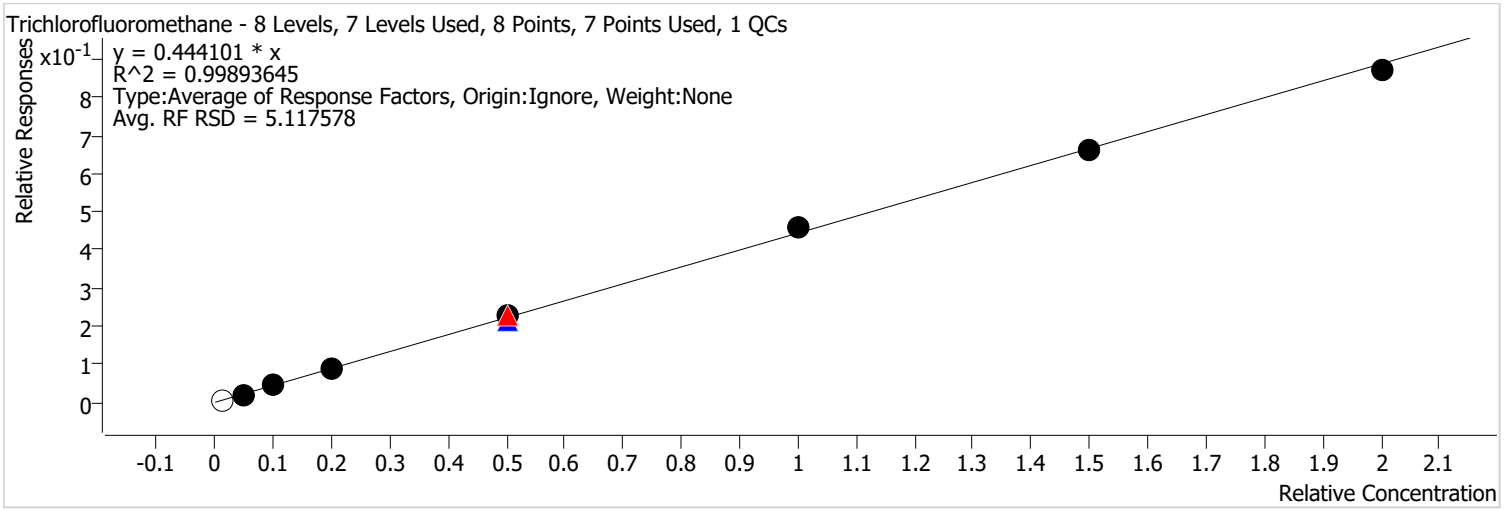
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	14646	25.0000	0.1851	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	25484	50.0000	0.1638	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	65619	125.0000	0.1638	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	71420	125.0000	0.1735	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	71420	125.0000	0.1735	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	137312	250.0000	0.1642	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	217393	375.0000	0.1721	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	287041	500.0000	0.1706	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Trichlorofluoromethane %RSE = 5.1



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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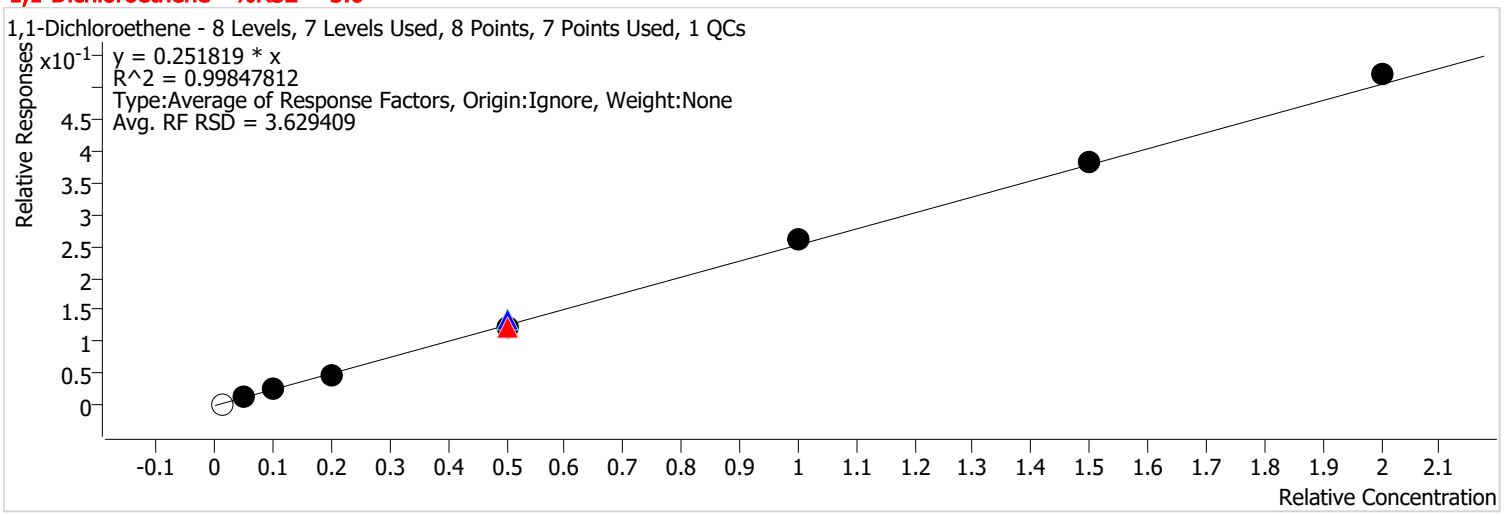
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	37464	25.0000	0.4735	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	68163	50.0000	0.4380	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	173333	125.0000	0.4327	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	188808	125.0000	0.4586	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	188808	125.0000	0.4586	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	384837	250.0000	0.4602	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	555477	375.0000	0.4399	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	731829	500.0000	0.4349	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,1-Dichloroethene %RSE = 3.6



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	20631	25.0000	0.2607	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	38253	50.0000	0.2458	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	108512	125.0000	0.2709	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	99438	125.0000	0.2415	
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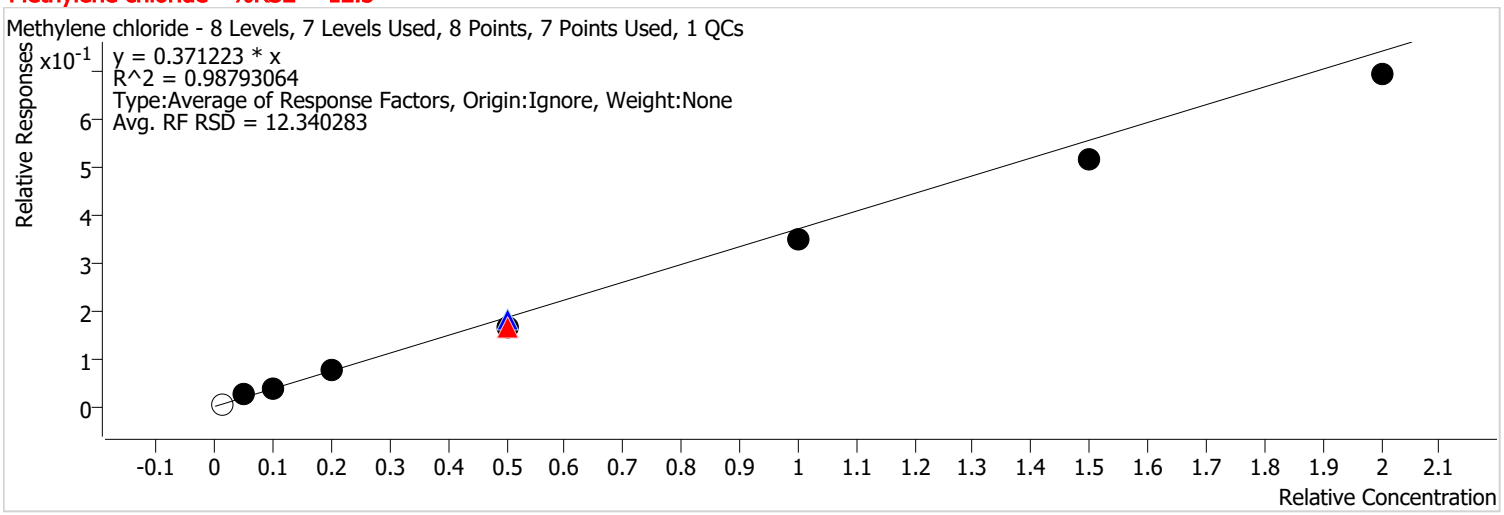
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	217406	250.0000	0.2600	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	322557	375.0000	0.2554	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	436507	500.0000	0.2594	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Methylene chloride %RSE = 12.3



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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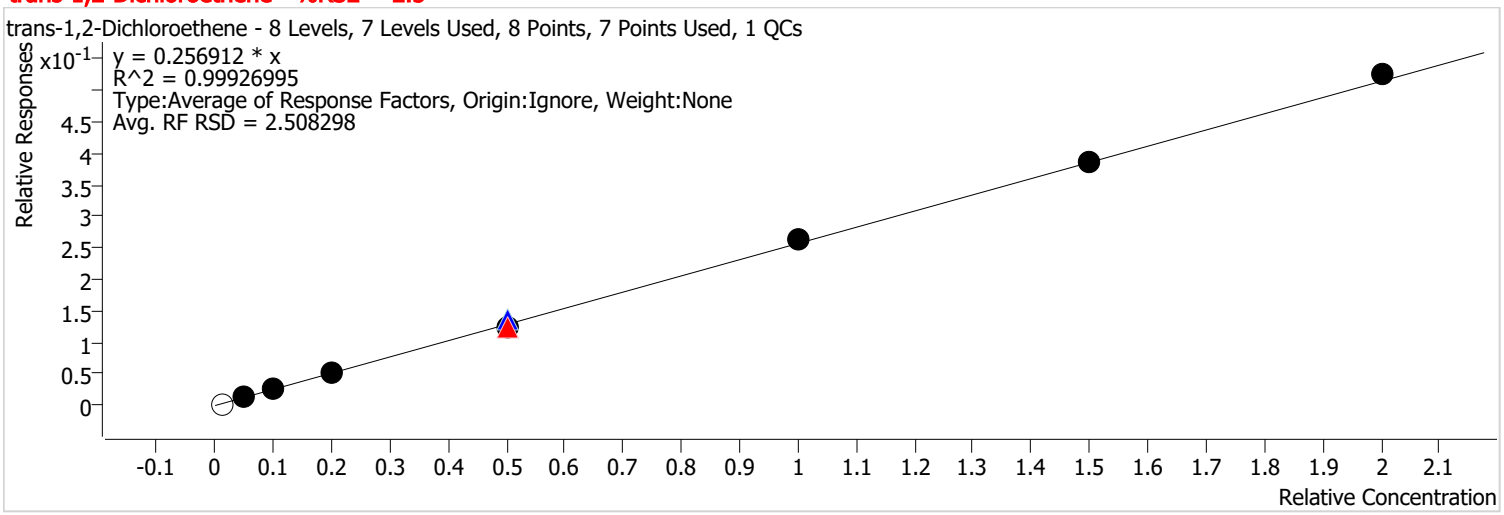
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	17734	12.5000	0.4640	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	30908	25.0000	0.3906	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	58282	50.0000	0.3745	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	144585	125.0000	0.3609	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	135271	125.0000	0.3285	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	135271	125.0000	0.3285	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	292397	250.0000	0.3496	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	435116	375.0000	0.3446	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	583438	500.0000	0.3467	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

trans-1,2-Dichloroethene %RSE = 2.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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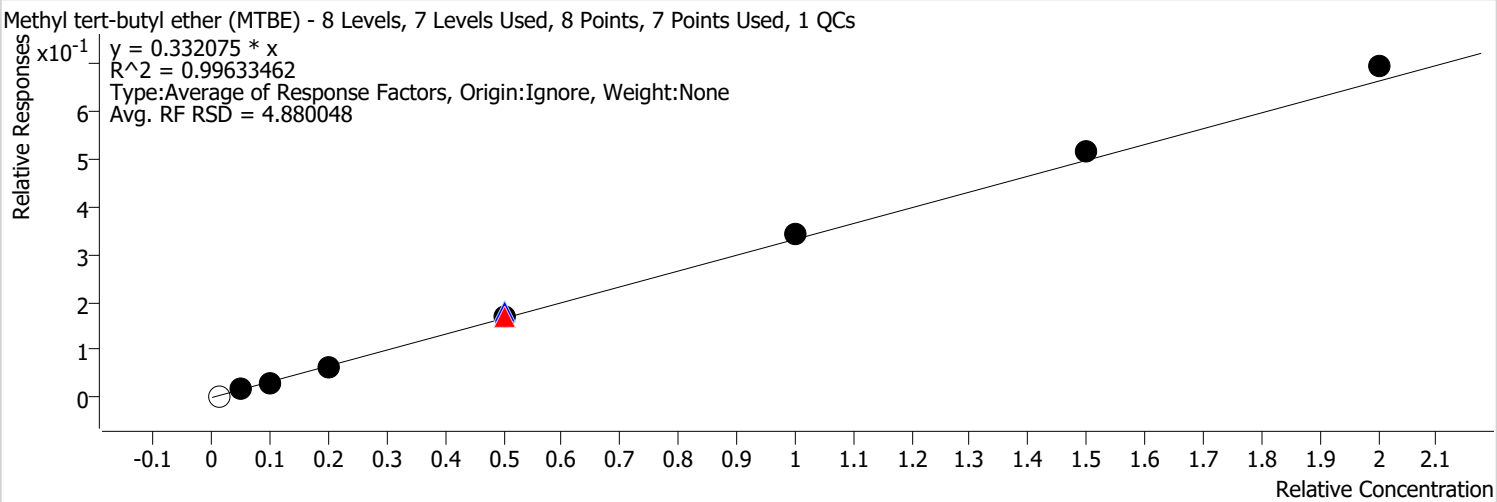
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	9821	12.5000	0.2570	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	20706	25.0000	0.2617	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	39596	50.0000	0.2544	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	110909	125.0000	0.2769	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	100409	125.0000	0.2439	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	100409	125.0000	0.2439	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	218855	250.0000	0.2617	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	325415	375.0000	0.2577	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	440967	500.0000	0.2621	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Methyl tert-butyl ether (MTBE) %RSE = 4.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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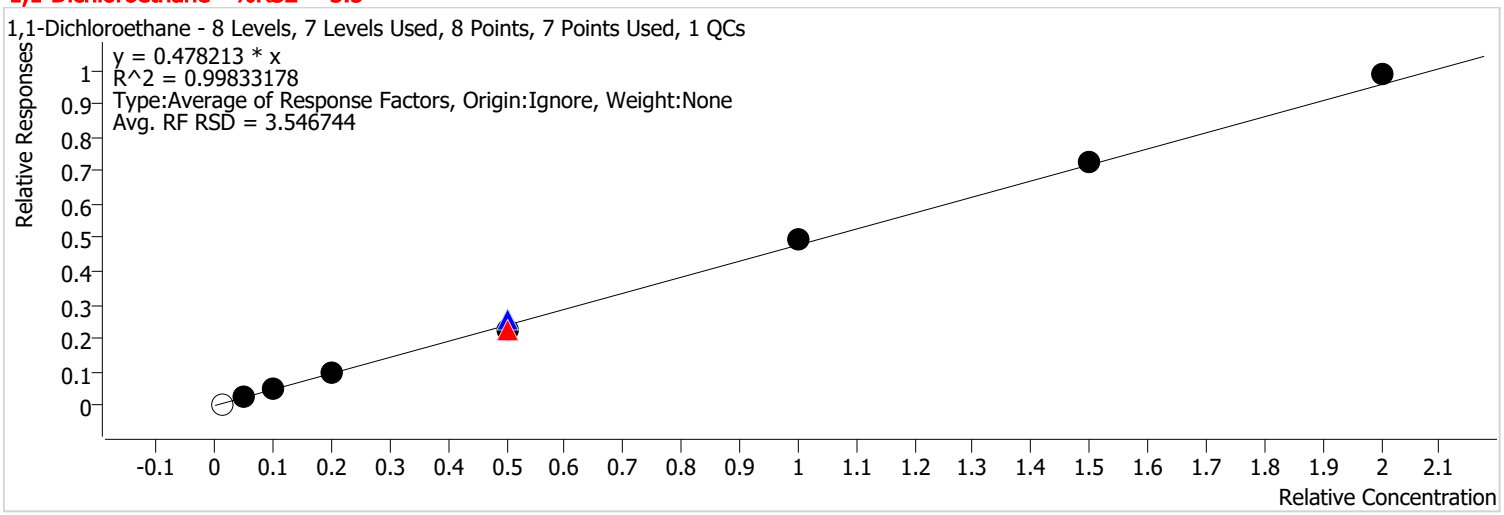
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	12515	12.5000	0.3274	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	24218	25.0000	0.3061	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	49126	50.0000	0.3157	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	143378	125.0000	0.3579	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	139068	125.0000	0.3378	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	139068	125.0000	0.3378	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	287653	250.0000	0.3440	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	437439	375.0000	0.3464	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	584294	500.0000	0.3472	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,1-Dichloroethane %RSE = 3.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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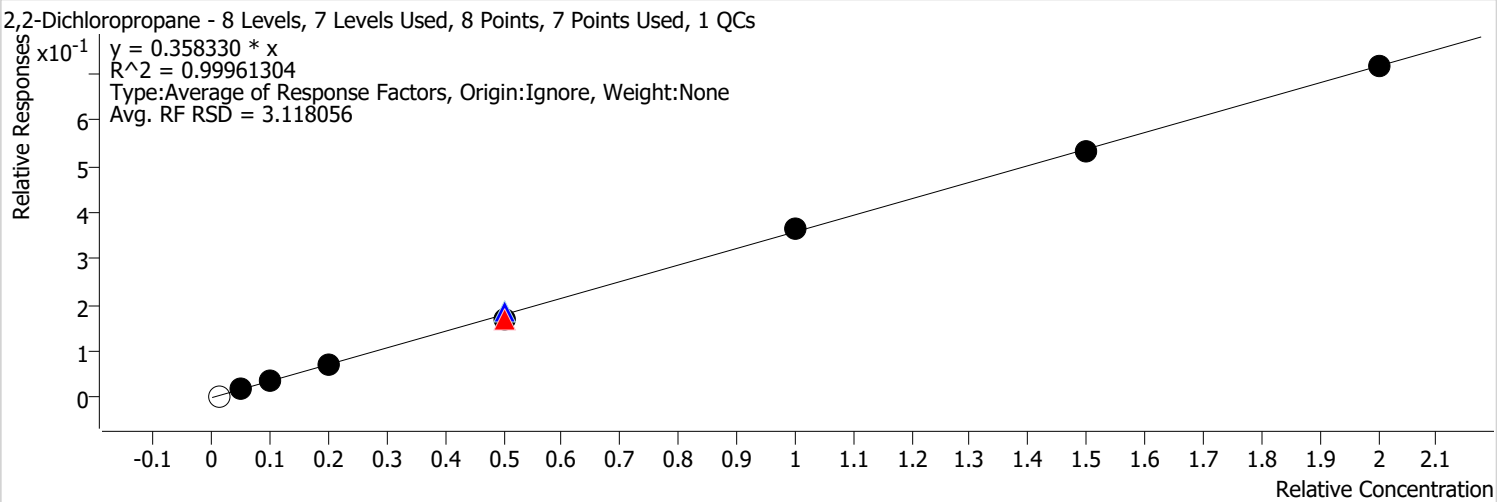
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	17642	12.5000	0.4616	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	38874	25.0000	0.4913	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	73205	50.0000	0.4704	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	208131	125.0000	0.5195	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	186052	125.0000	0.4519	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	186052	125.0000	0.4519	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	413408	250.0000	0.4943	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	612660	375.0000	0.4852	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	829359	500.0000	0.4929	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

2,2-Dichloropropane %RSE = 3.1



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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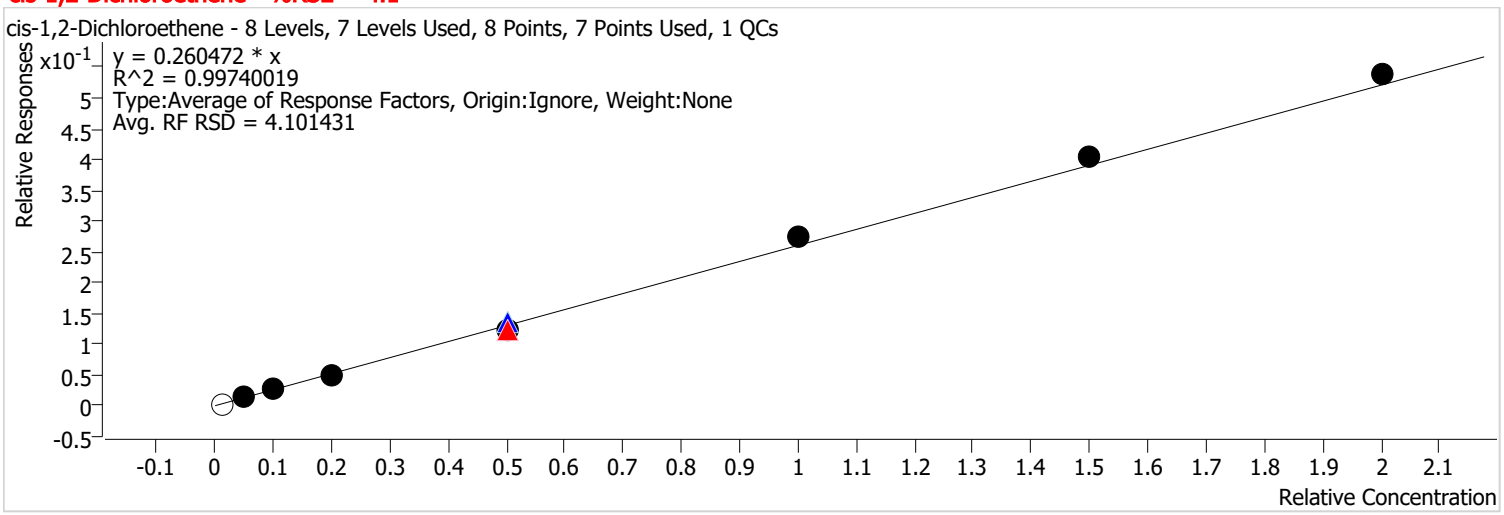
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	13676	12.5000	0.3578	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	29793	25.0000	0.3765	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	56189	50.0000	0.3611	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	150902	125.0000	0.3767	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	139656	125.0000	0.3392	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	139656	125.0000	0.3392	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	303307	250.0000	0.3627	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	446282	375.0000	0.3534	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	601823	500.0000	0.3576	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

cis-1,2-Dichloroethene %RSE = 4.1



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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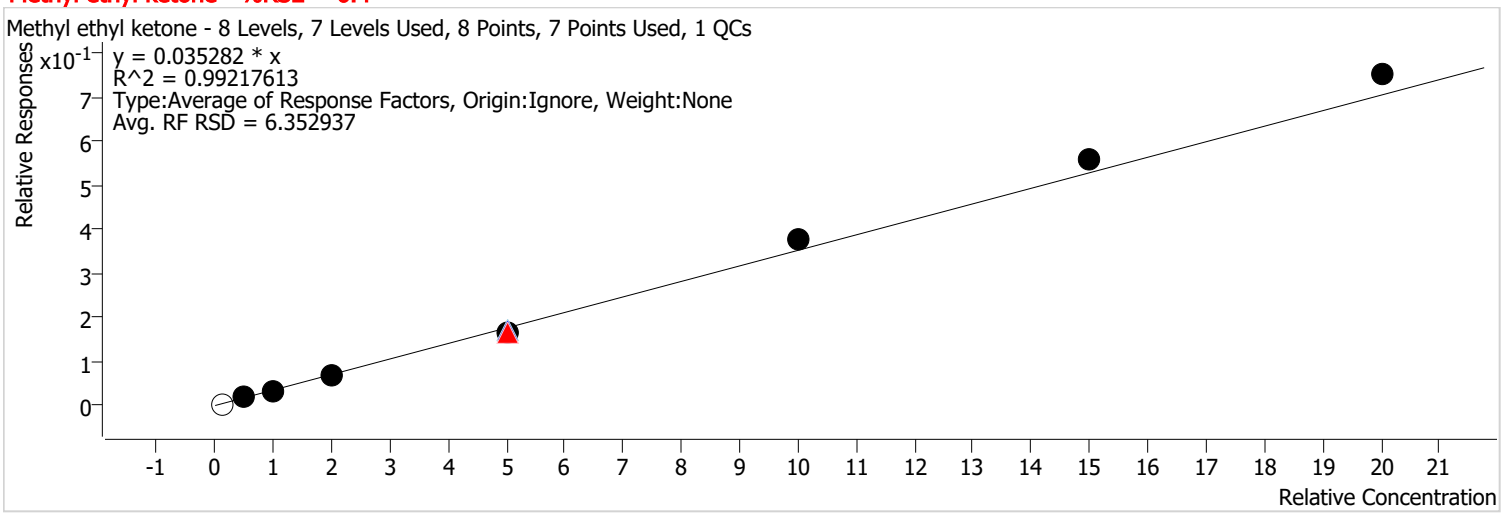
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	10008	12.5000	0.2618	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	20252	25.0000	0.2559	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	39251	50.0000	0.2522	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	108623	125.0000	0.2711	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	100057	125.0000	0.2430	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	100057	125.0000	0.2430	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	228170	250.0000	0.2728	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	339211	375.0000	0.2686	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	452377	500.0000	0.2688	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Methyl ethyl ketone %RSE = 6.4



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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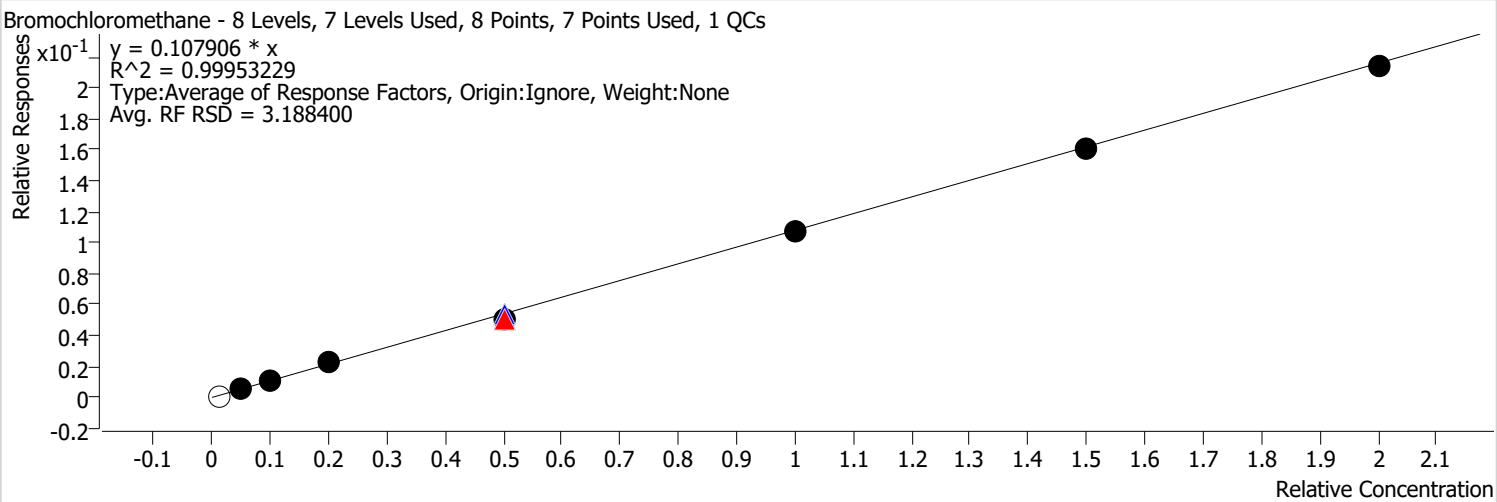
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	13167	125.0000	0.0344	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	26248	250.0000	0.0332	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	52648	500.0000	0.0338	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	135511	1250.0000	0.0338	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	134730	1250.0000	0.0327	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	134730	1250.0000	0.0327	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	317271	2500.0000	0.0379	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	470653	3750.0000	0.0373	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	632539	5000.0000	0.0376	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Bromochloromethane %RSE = 3.2



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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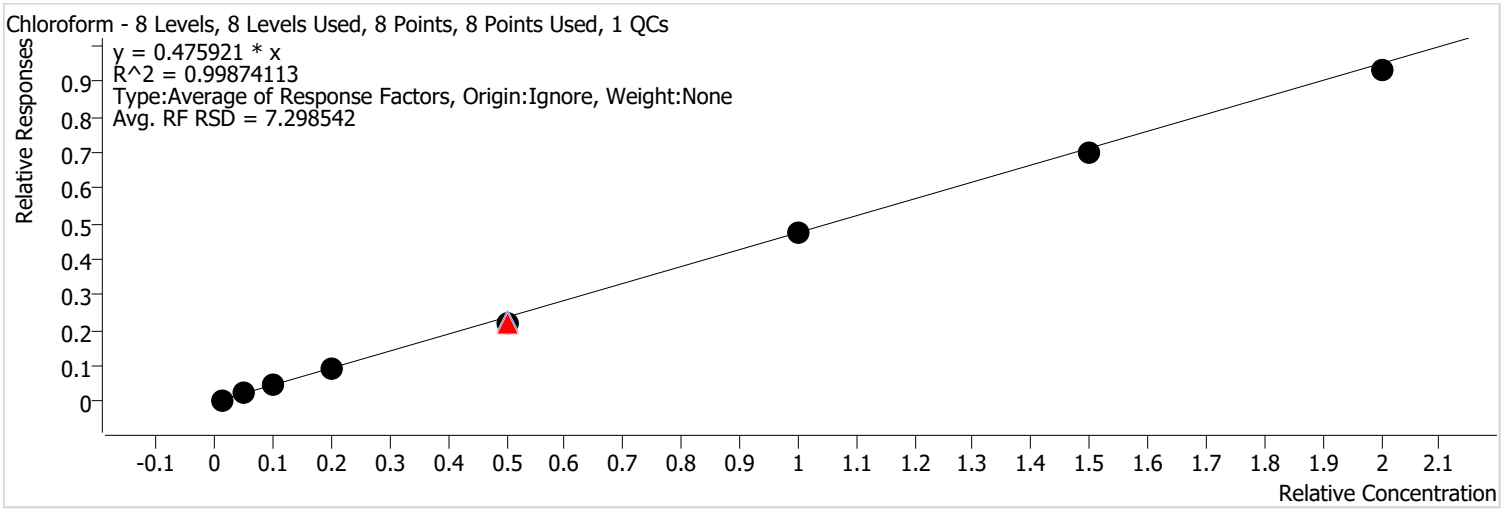
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	4275	12.5000	0.1118	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	8688	25.0000	0.1098	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	17338	50.0000	0.1114	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	42744	125.0000	0.1067	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	41966	125.0000	0.1019	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	41966	125.0000	0.1019	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	89178	250.0000	0.1066	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	135103	375.0000	0.1070	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	179618	500.0000	0.1067	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

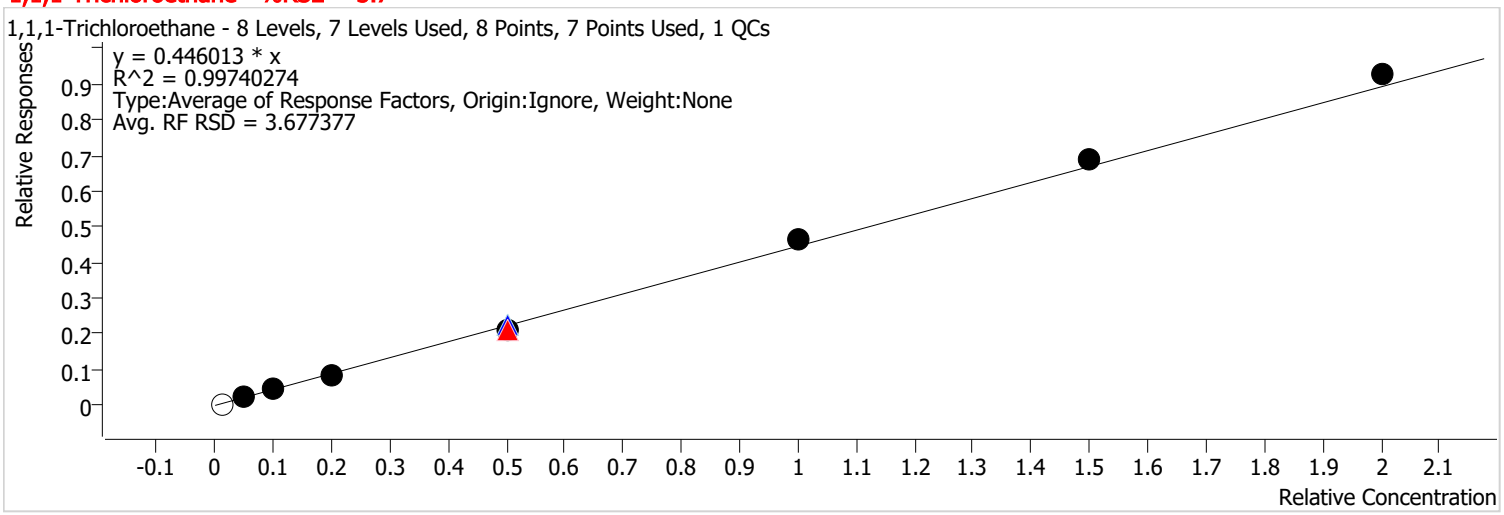
Chloroform %RSE = 7.3



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
D:\Org\Data\VOA5975C\VG010422\04JAN10.D	Calibration	1	x	4248	2.5000	0.5510	
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	19015	12.5000	0.4975	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	36413	25.0000	0.4602	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	71403	50.0000	0.4588	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	183676	125.0000	0.4585	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	179640	125.0000	0.4363	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	179640	125.0000	0.4363	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	394946	250.0000	0.4723	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	588080	375.0000	0.4657	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	783422	500.0000	0.4656	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,1,1-Trichloroethane %RSE = 3.7



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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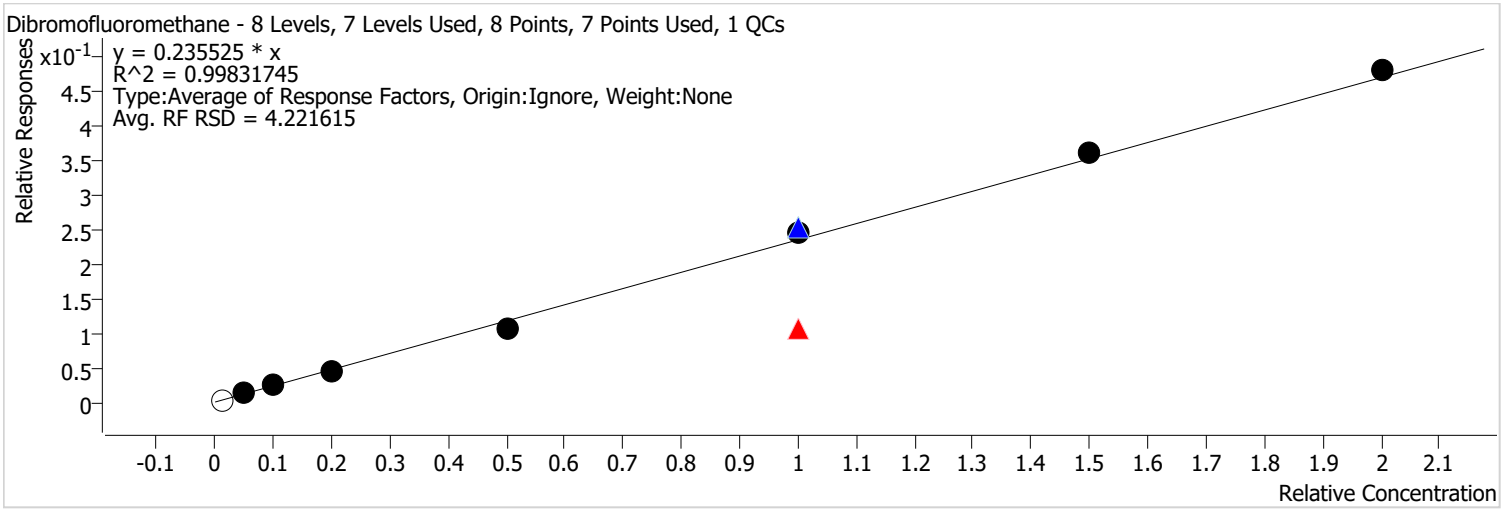
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	16623	12.5000	0.4349	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	35547	25.0000	0.4492	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	67007	50.0000	0.4306	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	183324	125.0000	0.4576	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	174206	125.0000	0.4231	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	174206	125.0000	0.4231	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	386005	250.0000	0.4616	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	580748	375.0000	0.4599	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	778785	500.0000	0.4628	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

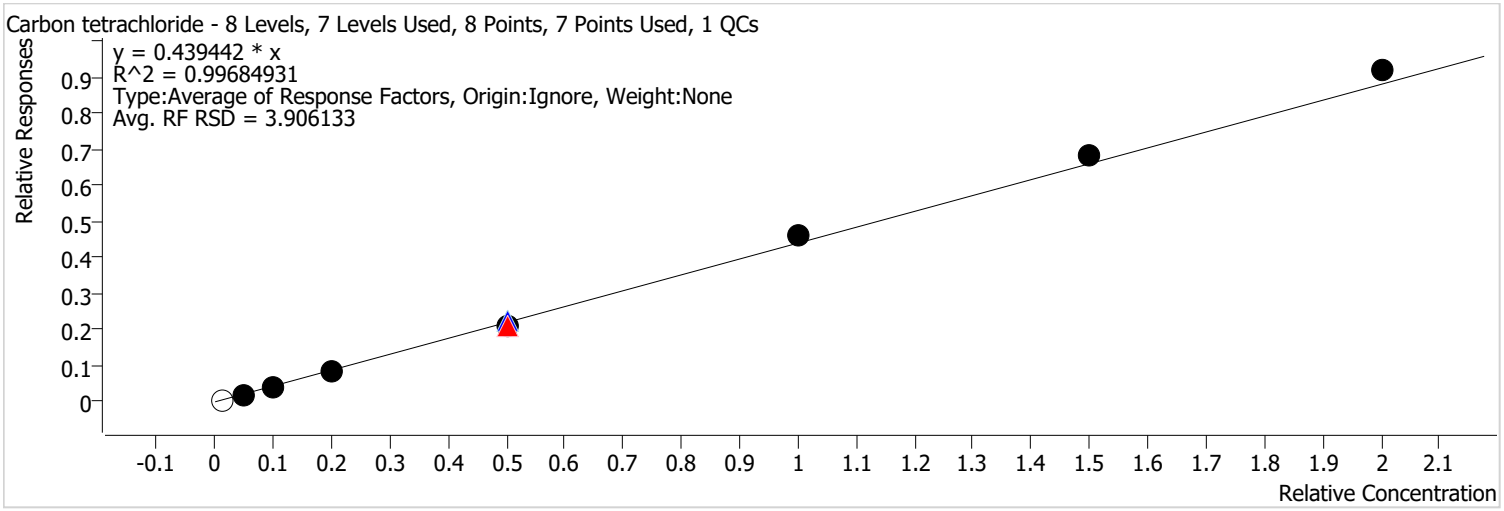
Dibromofluoromethane %RSE =



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
D:\Org\Data\VOA5975C\VG010422\04JAN10.D	Calibration	1		2508	2.5000	0.3253	
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	9074	12.5000	0.2374	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	19100	25.0000	0.2414	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	35309	50.0000	0.2269	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	89307	125.0000	0.2169	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	204707	250.0000	0.2555	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	204073	250.0000	0.2440	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	89307	250.0000	0.1084	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	305158	375.0000	0.2416	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	404568	500.0000	0.2404	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:43 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Carbon tetrachloride %RSE = 3.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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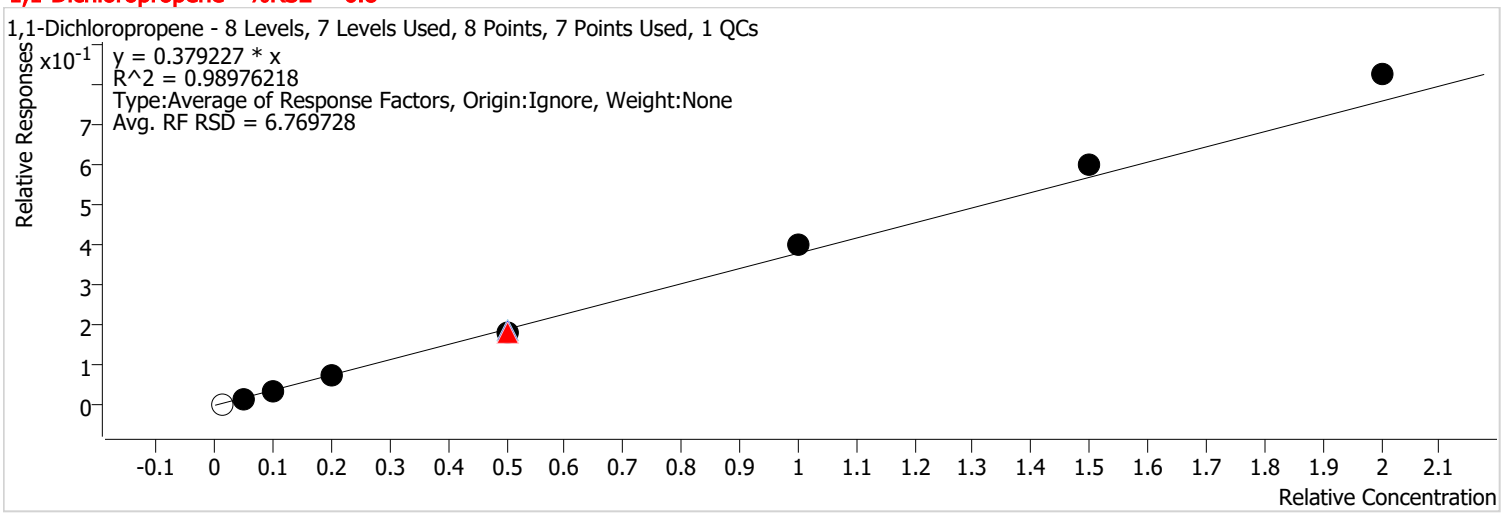
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	16466	12.5000	0.4308	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	34462	25.0000	0.4355	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	65313	50.0000	0.4197	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	181384	125.0000	0.4528	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	172928	125.0000	0.4200	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	172928	125.0000	0.4200	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	383485	250.0000	0.4586	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	572545	375.0000	0.4534	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	770907	500.0000	0.4581	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

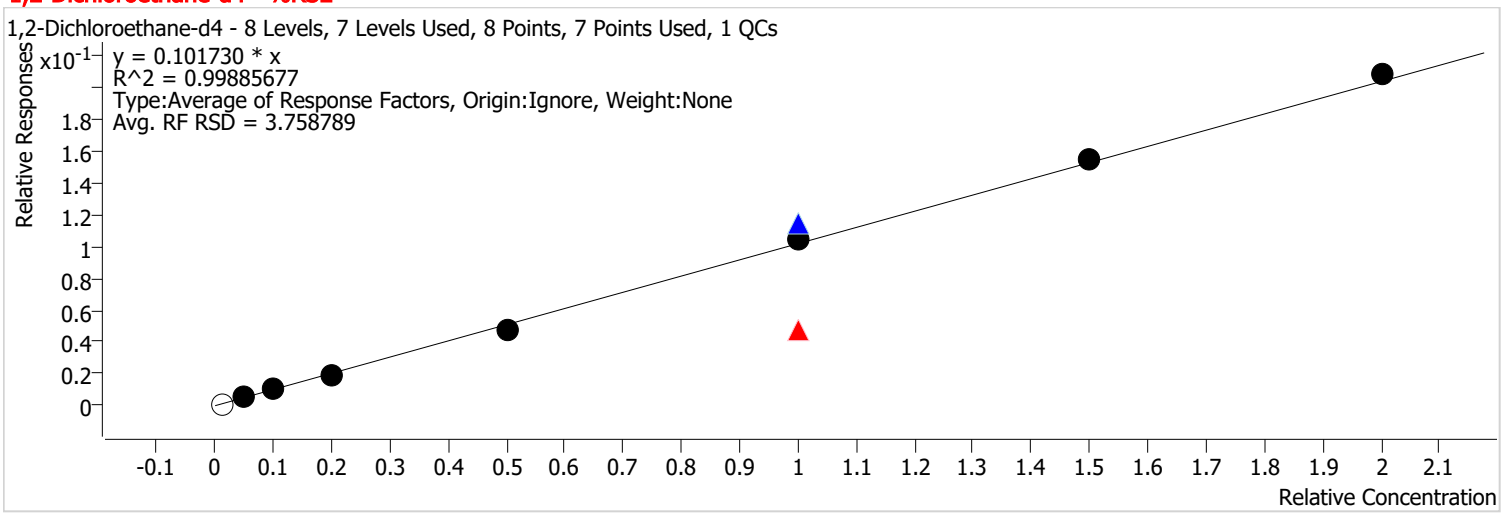
1,1-Dichloropropene %RSE = 6.8



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
D:\Org\Data\VOA5975C\VG010422\04JAN10.D	Calibration	1		2830	2.5000	0.3671	
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	13149	12.5000	0.3440	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	29241	25.0000	0.3695	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	56376	50.0000	0.3623	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	150930	125.0000	0.3768	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	149649	125.0000	0.3635	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	149649	125.0000	0.3635	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	335741	250.0000	0.4015	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	507157	375.0000	0.4016	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	693669	500.0000	0.4122	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,2-Dichloroethane-d4 %RSE =



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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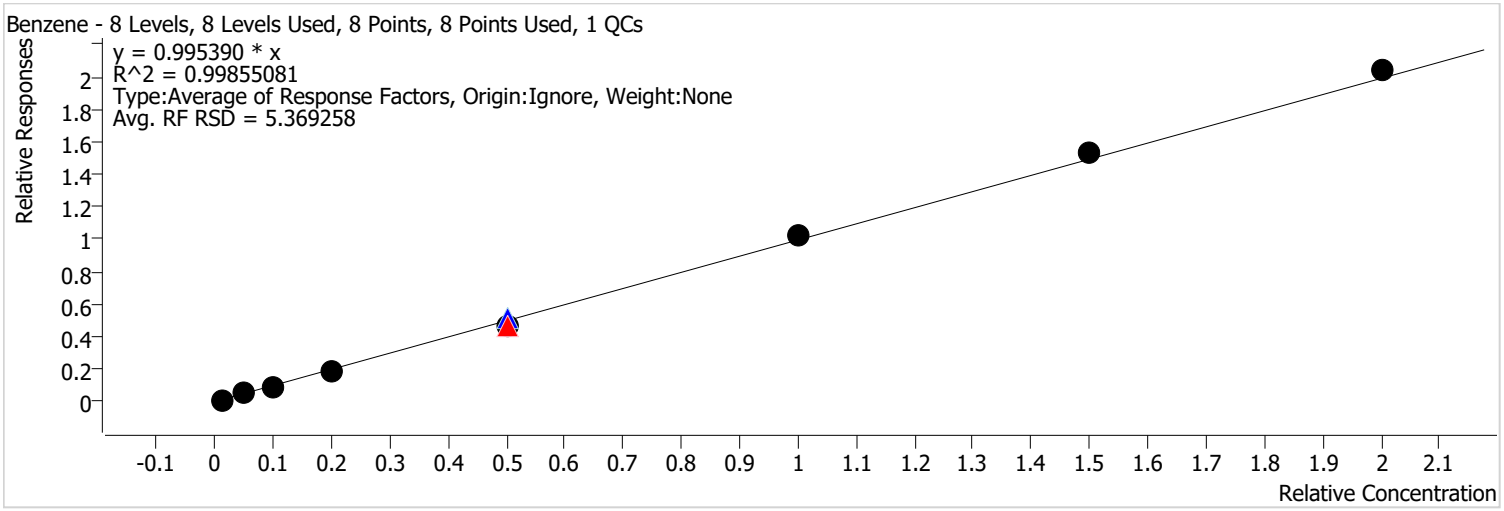
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	3938	12.5000	0.1030	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	8284	25.0000	0.1047	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	15238	50.0000	0.0979	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	39086	125.0000	0.0949	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	91382	250.0000	0.1141	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	87876	250.0000	0.1051	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	39086	250.0000	0.0475	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	129608	375.0000	0.1026	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	174713	500.0000	0.1038	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

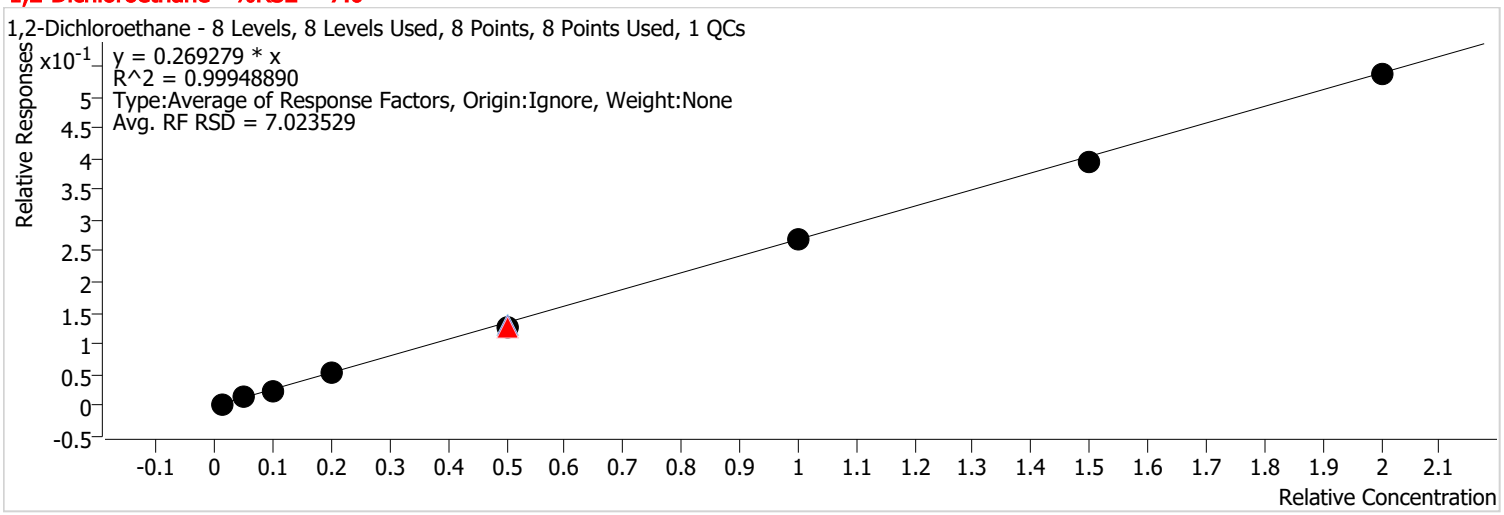
Benzene %RSE = 5.4



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
D:\Org\Data\VOA5975C\VG010422\04JAN10.D	Calibration	1	x	8408	2.5000	1.0907	
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	37071	12.5000	0.9699	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	74956	25.0000	0.9473	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	148727	50.0000	0.9557	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	418900	125.0000	1.0457	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	383469	125.0000	0.9313	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	383469	125.0000	0.9313	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	857534	250.0000	1.0254	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	1293370	375.0000	1.0242	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1714050	500.0000	1.0186	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,2-Dichloroethane %RSE = 7.0



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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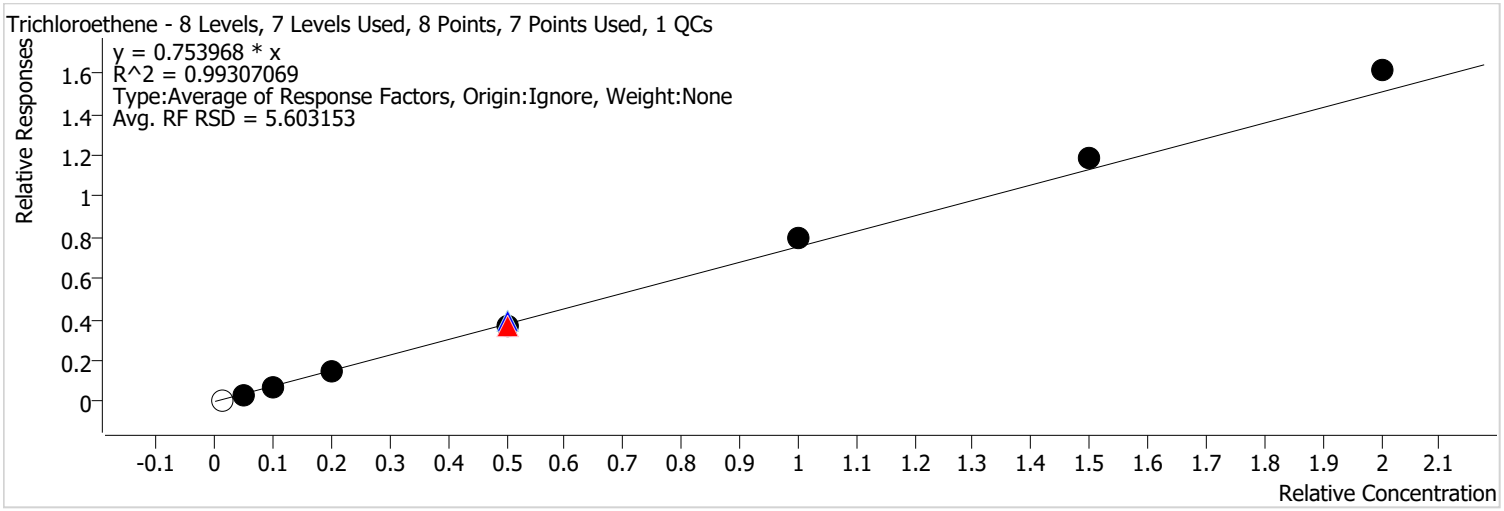
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	10202	12.5000	0.2669	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	19996	25.0000	0.2527	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	41058	50.0000	0.2638	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	104249	125.0000	0.2602	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	104855	125.0000	0.2547	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	104855	125.0000	0.2547	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	226964	250.0000	0.2714	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	332775	375.0000	0.2635	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	450739	500.0000	0.2679	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Trichloroethene %RSE = 5.6



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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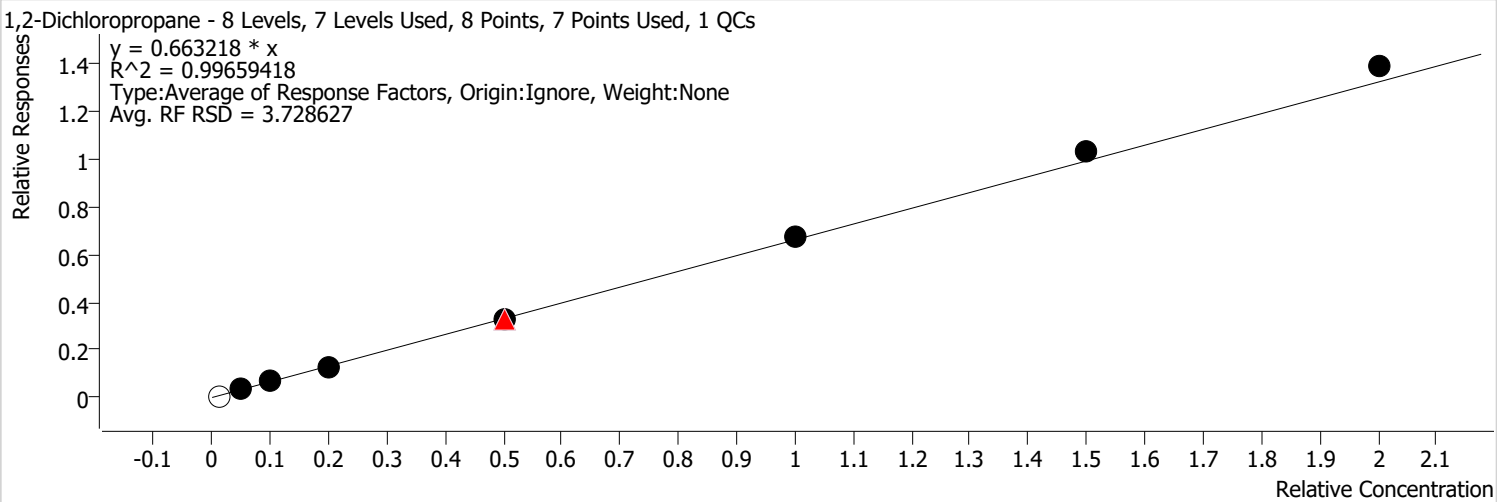
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	10442	12.5000	0.7042	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	21946	25.0000	0.7283	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	42682	50.0000	0.7105	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	121734	125.0000	0.7908	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	114123	125.0000	0.7447	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	114123	125.0000	0.7447	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	250285	250.0000	0.7910	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	374370	375.0000	0.7932	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	505400	500.0000	0.8058	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,2-Dichloropropane %RSE = 3.7



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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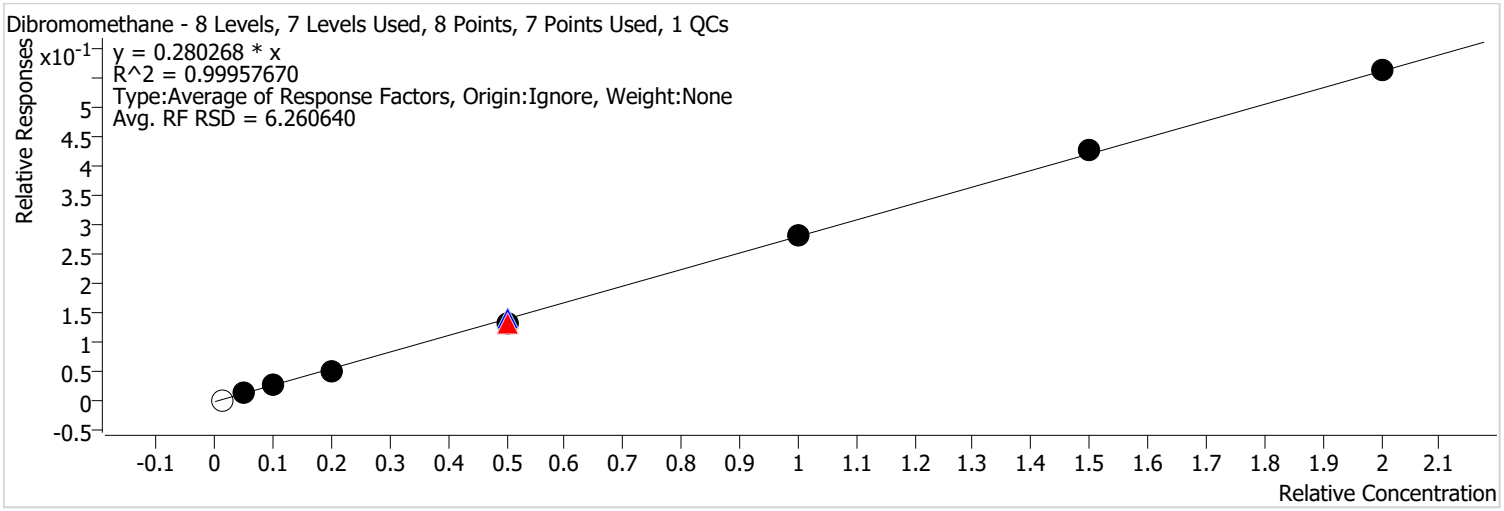
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	9488	12.5000	0.6399	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	20077	25.0000	0.6663	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	37870	50.0000	0.6304	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	102633	125.0000	0.6667	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	99187	125.0000	0.6472	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	99187	125.0000	0.6472	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	213800	250.0000	0.6757	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	324602	375.0000	0.6877	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	436057	500.0000	0.6953	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Dibromomethane %RSE = 6.3



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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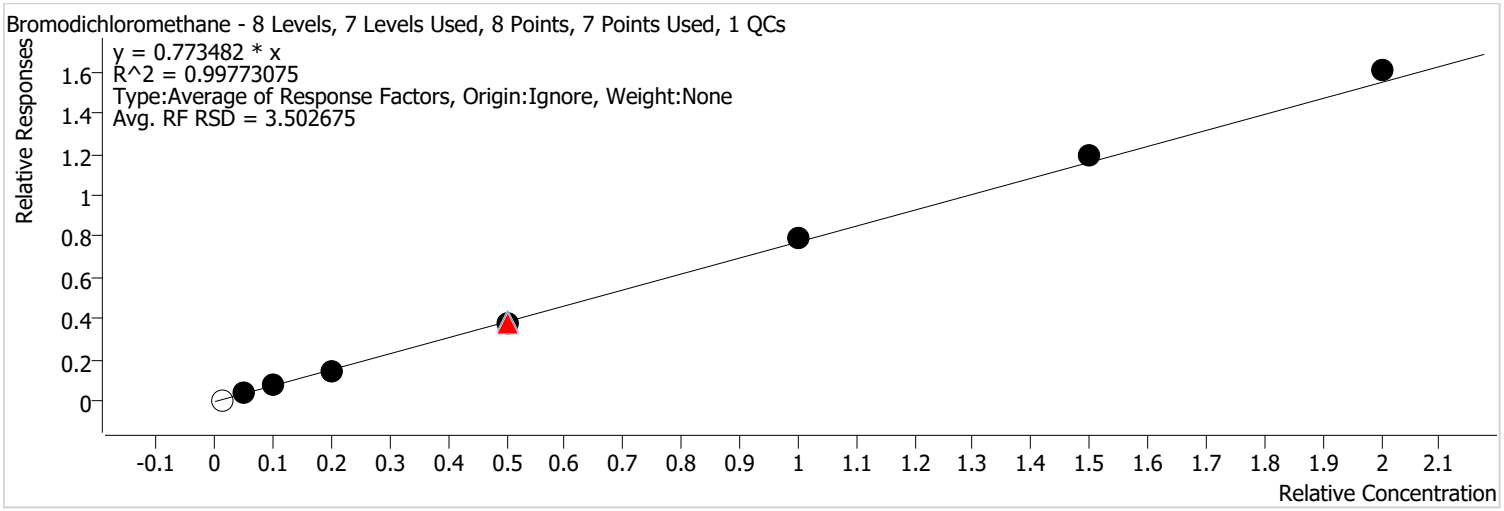
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	4675	12.5000	0.3153	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	8055	25.0000	0.2673	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	15989	50.0000	0.2662	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	43248	125.0000	0.2810	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	40628	125.0000	0.2651	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	40628	125.0000	0.2651	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	89483	250.0000	0.2828	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	134282	375.0000	0.2845	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	176038	500.0000	0.2807	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Bromodichloromethane %RSE = 3.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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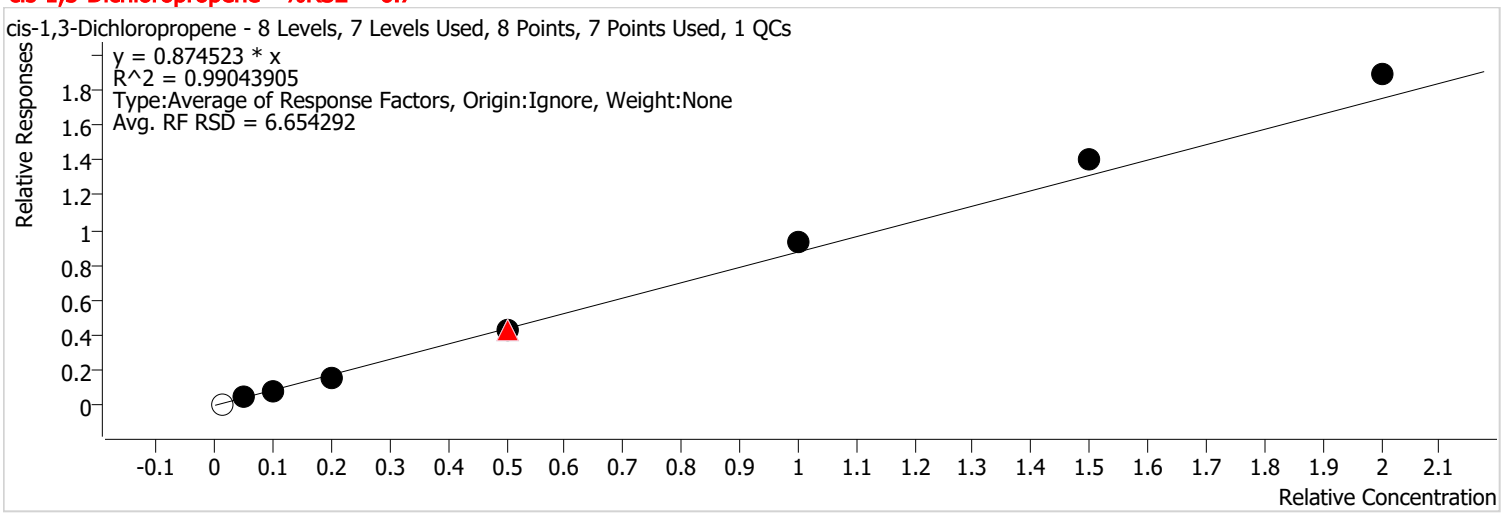
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	11562	12.5000	0.7798	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	22743	25.0000	0.7547	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	43900	50.0000	0.7308	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	122757	125.0000	0.7975	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	115664	125.0000	0.7548	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	115664	125.0000	0.7548	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	251805	250.0000	0.7958	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	375983	375.0000	0.7966	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	502929	500.0000	0.8019	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

cis-1,3-Dichloropropene %RSE = 6.7



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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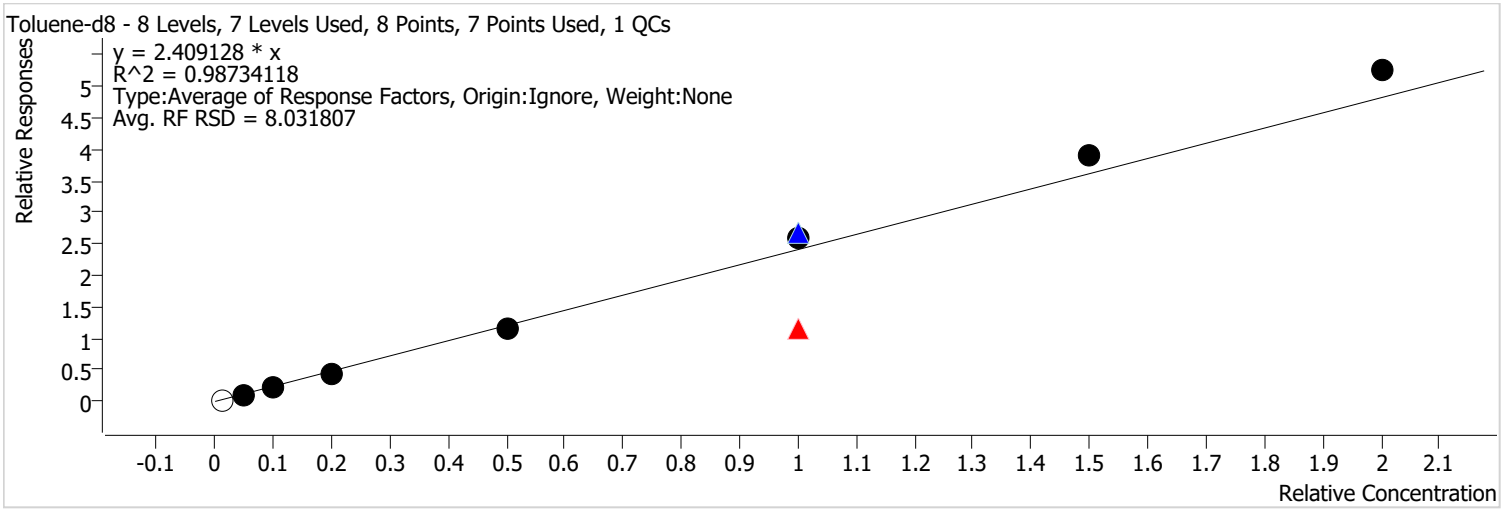
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	12525	12.5000	0.8447	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	24511	25.0000	0.8134	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	48886	50.0000	0.8138	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	130910	125.0000	0.8504	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	129419	125.0000	0.8445	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	129419	125.0000	0.8445	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	293617	250.0000	0.9280	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	441168	375.0000	0.9347	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	591147	500.0000	0.9426	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Toluene-d8 %RSE =



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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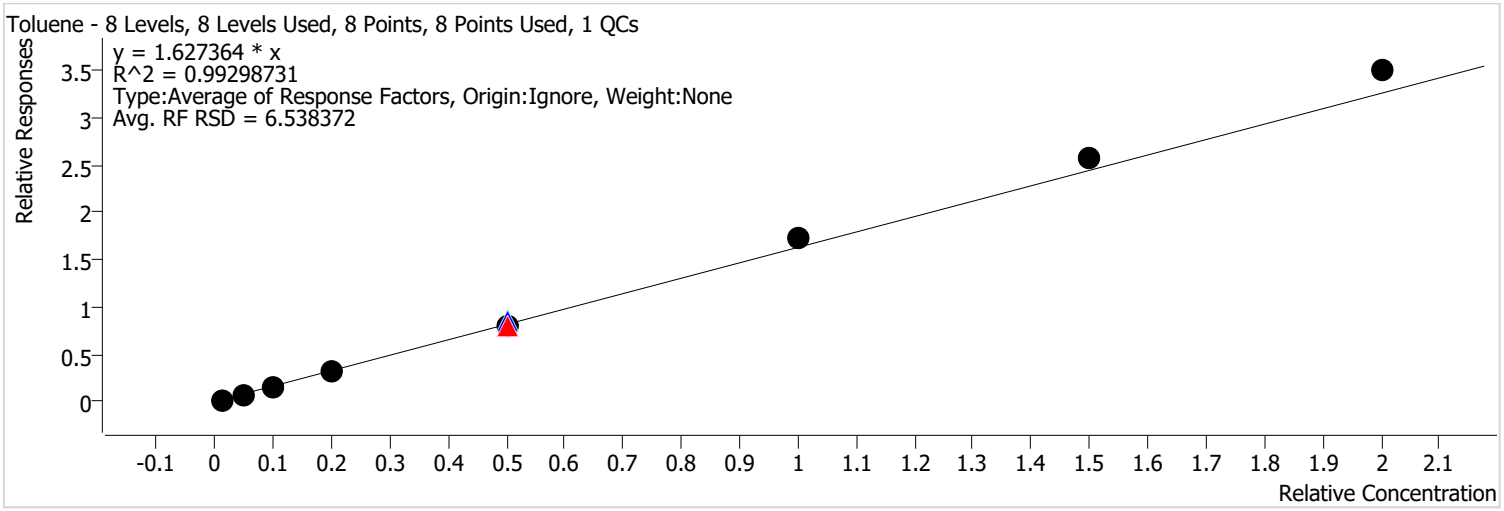
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	32318	12.5000	2.1796	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	67673	25.0000	2.2458	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	136453	50.0000	2.2715	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	358186	125.0000	2.3373	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	821531	250.0000	2.6685	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	823306	250.0000	2.6021	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	358186	250.0000	1.1687	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	1229775	375.0000	2.6054	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1644540	500.0000	2.6222	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Toluene %RSE = 6.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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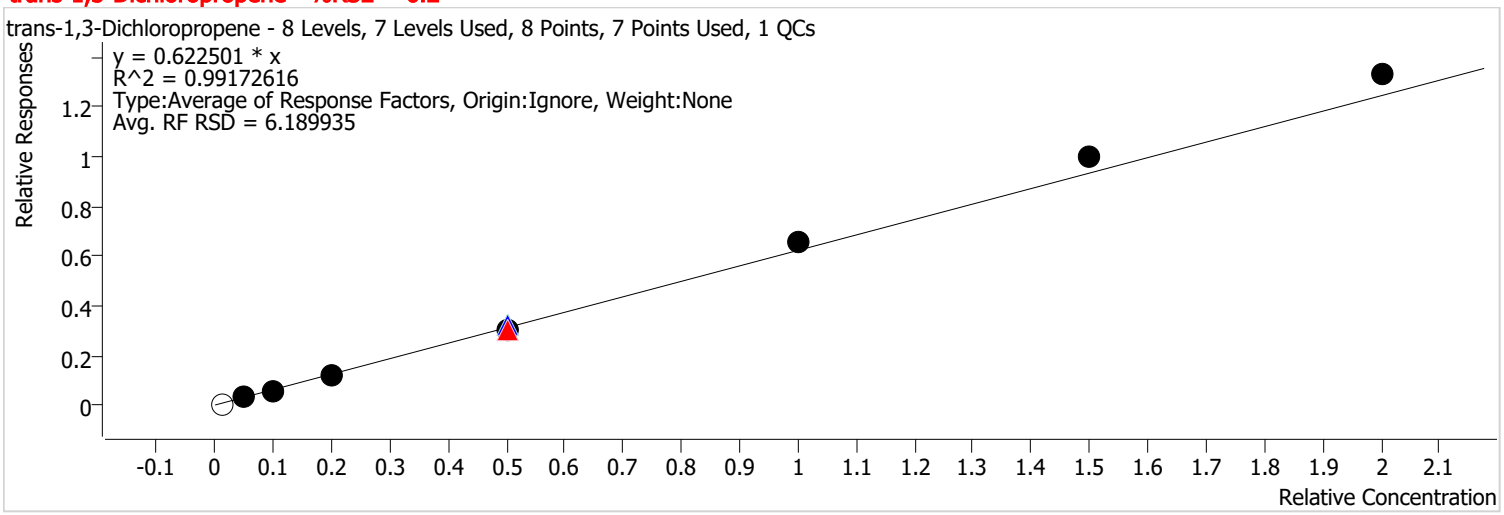
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	21794	12.5000	1.4698	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	46355	25.0000	1.5383	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	91915	50.0000	1.5301	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	264584	125.0000	1.7188	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	244712	125.0000	1.5969	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	244712	125.0000	1.5969	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	541945	250.0000	1.7129	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	813204	375.0000	1.7229	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1095161	500.0000	1.7462	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

trans-1,3-Dichloropropene %RSE = 6.2



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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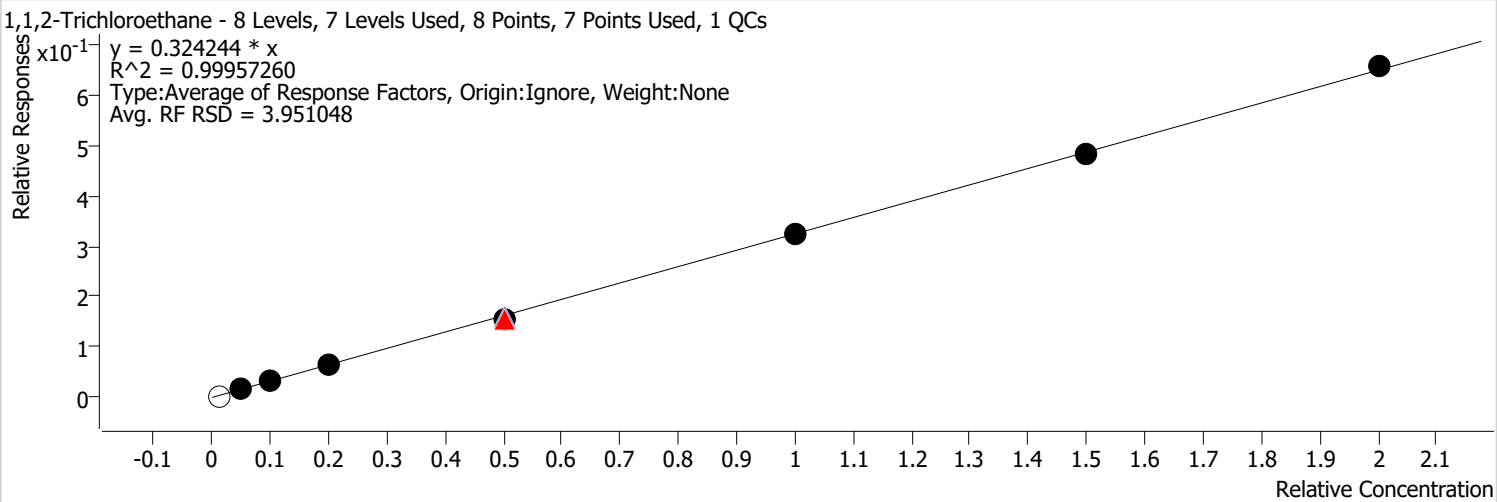
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	8683	12.5000	0.5856	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	17850	25.0000	0.5924	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	35179	50.0000	0.5856	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	98907	125.0000	0.6425	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	92719	125.0000	0.6050	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	92719	125.0000	0.6050	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	207833	250.0000	0.6569	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	315063	375.0000	0.6675	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	416771	500.0000	0.6645	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,1,2-Trichloroethane %RSE = 4.0



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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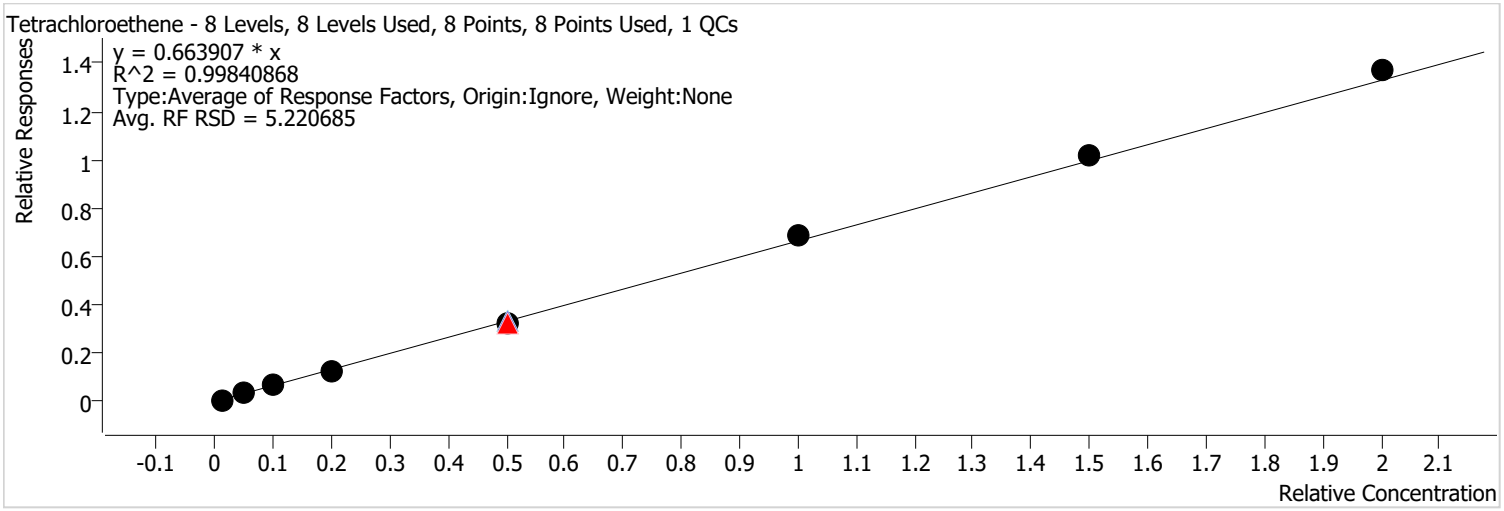
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	5090	12.5000	0.3433	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	10099	25.0000	0.3351	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	18884	50.0000	0.3144	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	49128	125.0000	0.3191	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	46673	125.0000	0.3046	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	46673	125.0000	0.3046	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	101888	250.0000	0.3220	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	152331	375.0000	0.3227	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	205463	500.0000	0.3276	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Tetrachloroethene %RSE = 5.2



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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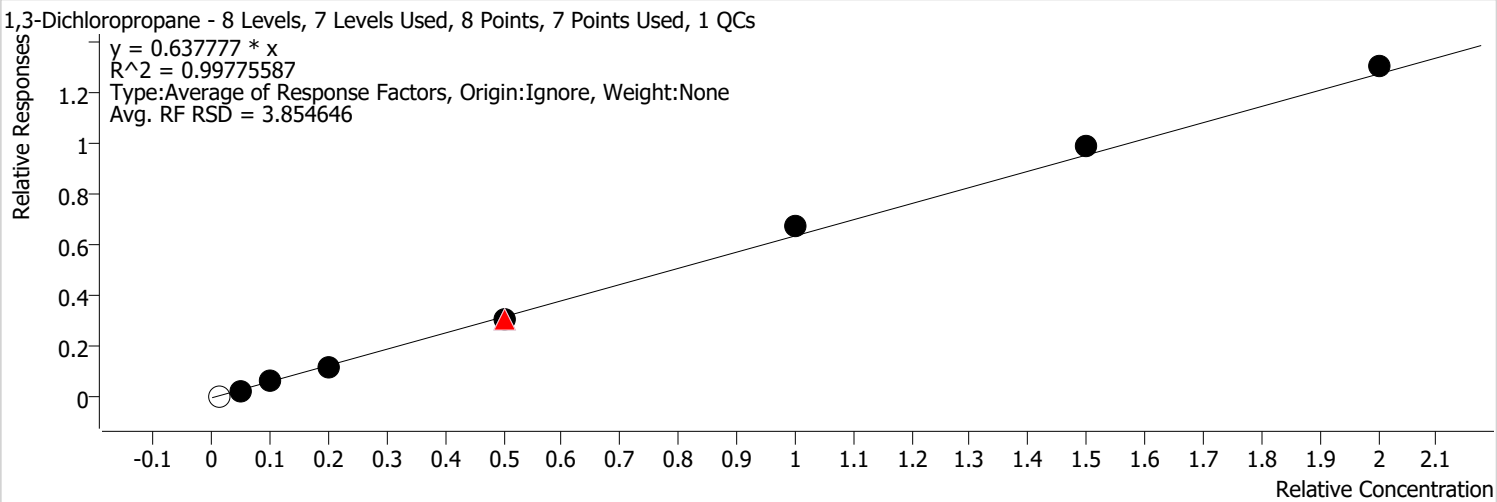
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	9238	12.5000	0.6230	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	20322	25.0000	0.6744	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	36925	50.0000	0.6147	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	103027	125.0000	0.6693	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	97590	125.0000	0.6368	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	97590	125.0000	0.6368	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	218245	250.0000	0.6898	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	319950	375.0000	0.6779	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	428812	500.0000	0.6837	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,3-Dichloropropane %RSE = 3.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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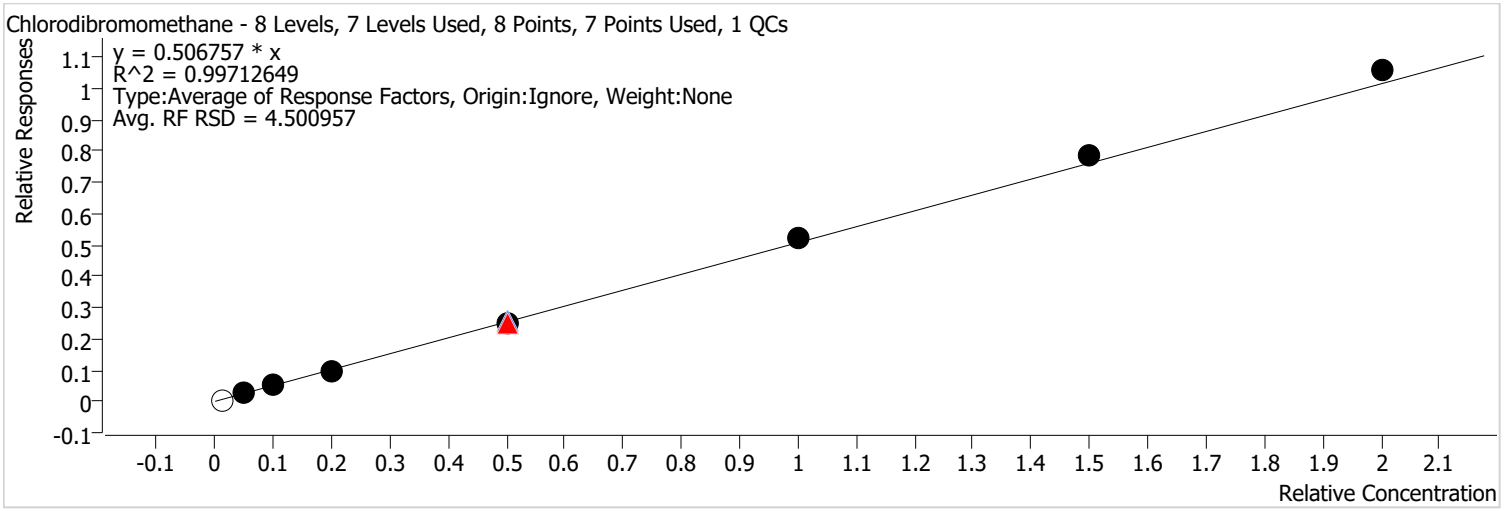
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	8967	12.5000	0.6047	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	18745	25.0000	0.6221	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	37457	50.0000	0.6235	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	95697	125.0000	0.6217	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	96183	125.0000	0.6276	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	96183	125.0000	0.6276	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	212669	250.0000	0.6722	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	312547	375.0000	0.6622	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	408993	500.0000	0.6521	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Chlorodibromomethane %RSE = 4.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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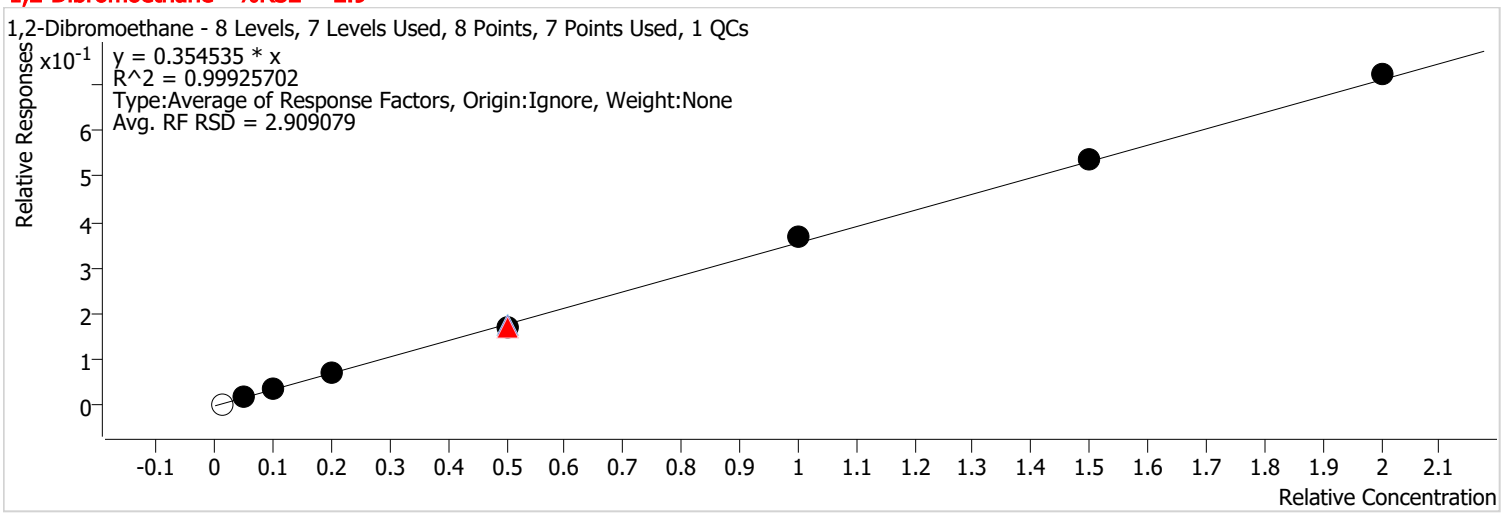
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	7718	12.5000	0.5205	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	14873	25.0000	0.4936	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	28153	50.0000	0.4687	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	78076	125.0000	0.5072	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	75015	125.0000	0.4895	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	75015	125.0000	0.4895	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	165695	250.0000	0.5237	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	247279	375.0000	0.5239	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	330813	500.0000	0.5275	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,2-Dibromoethane %RSE = 2.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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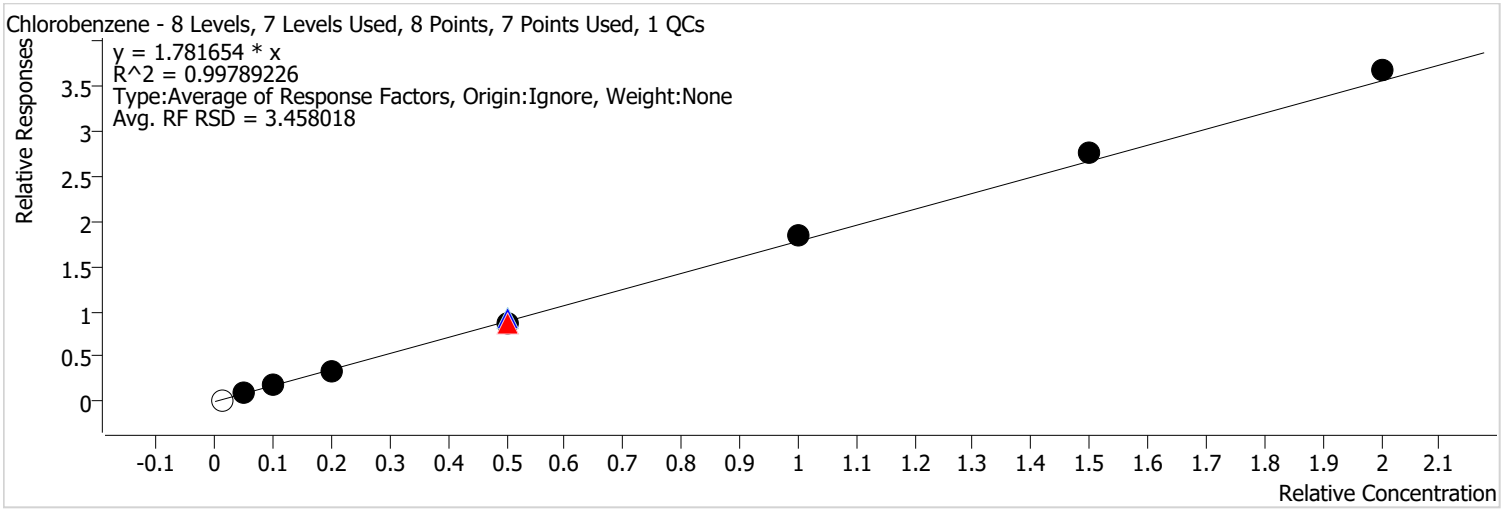
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	5410	12.5000	0.3649	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	10410	25.0000	0.3455	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	21037	50.0000	0.3502	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	54259	125.0000	0.3525	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	51827	125.0000	0.3382	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	51827	125.0000	0.3382	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	115714	250.0000	0.3657	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	168577	375.0000	0.3572	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	225877	500.0000	0.3602	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Chlorobenzene %RSE = 3.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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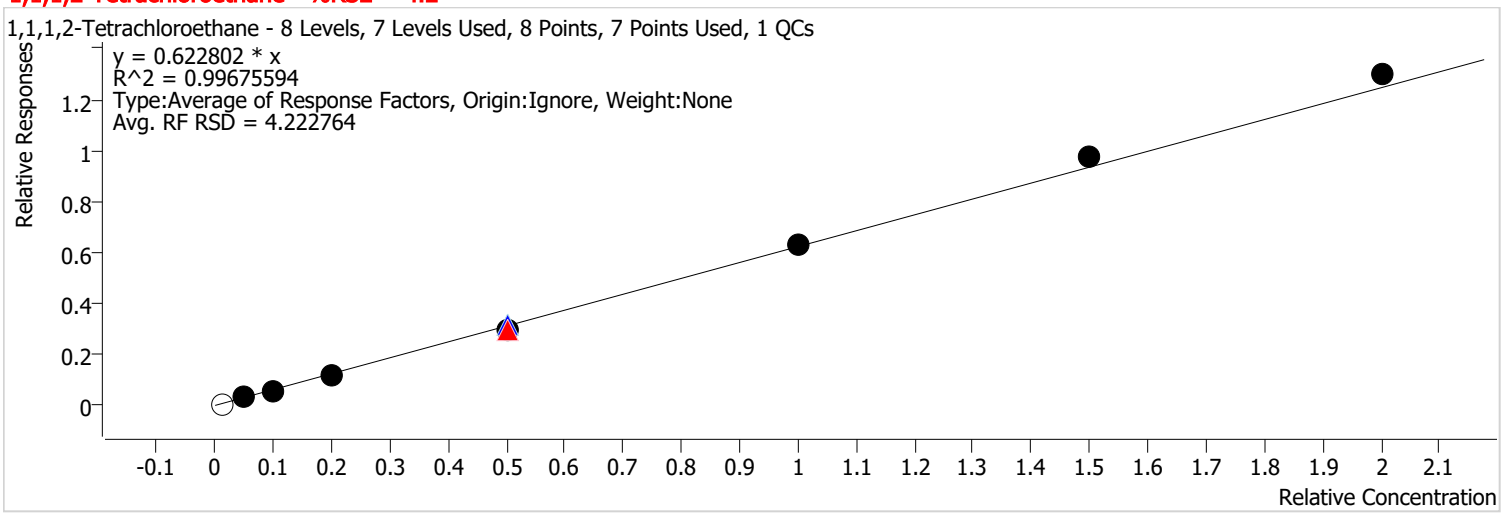
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	26461	12.5000	1.7846	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	53047	25.0000	1.7604	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	101452	50.0000	1.6889	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	288815	125.0000	1.8762	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	263617	125.0000	1.7202	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	263617	125.0000	1.7202	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	582326	250.0000	1.8405	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	867732	375.0000	1.8384	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1153147	500.0000	1.8387	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:44 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,1,1,2-Tetrachloroethane %RSE = 4.2



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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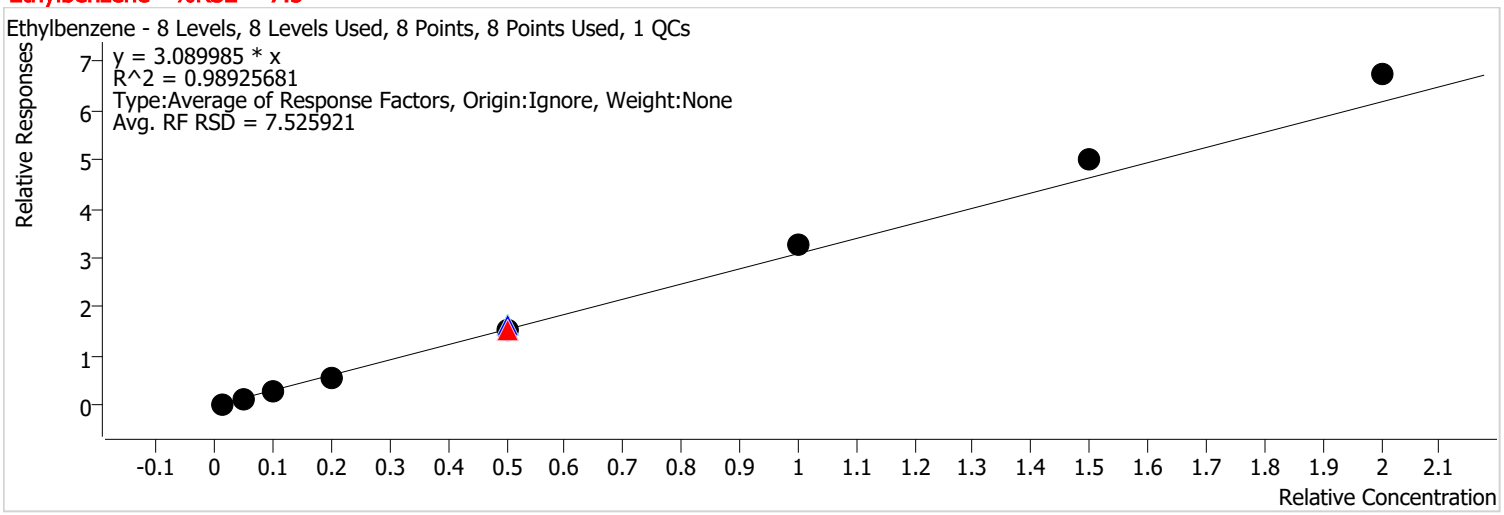
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	9473	12.5000	0.6389	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	18130	25.0000	0.6016	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	35544	50.0000	0.5917	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	97148	125.0000	0.6311	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	90898	125.0000	0.5932	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	90898	125.0000	0.5932	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	200859	250.0000	0.6348	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	307436	375.0000	0.6513	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	406450	500.0000	0.6481	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Ethylbenzene %RSE = 7.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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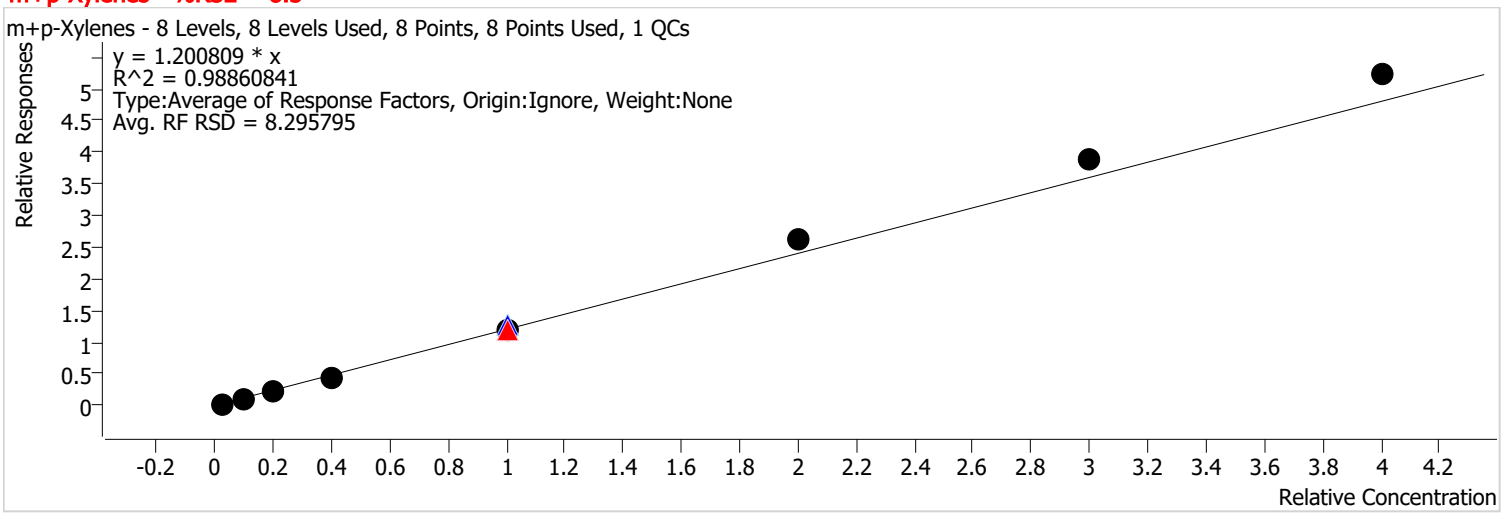
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	40470	12.5000	2.7294	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	88428	25.0000	2.9345	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	173769	50.0000	2.8927	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	501953	125.0000	3.2608	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	464148	125.0000	3.0288	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	464148	125.0000	3.0288	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	1043443	250.0000	3.2979	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	1574219	375.0000	3.3352	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	2111152	500.0000	3.3662	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

m+p-Xylenes %RSE = 8.3



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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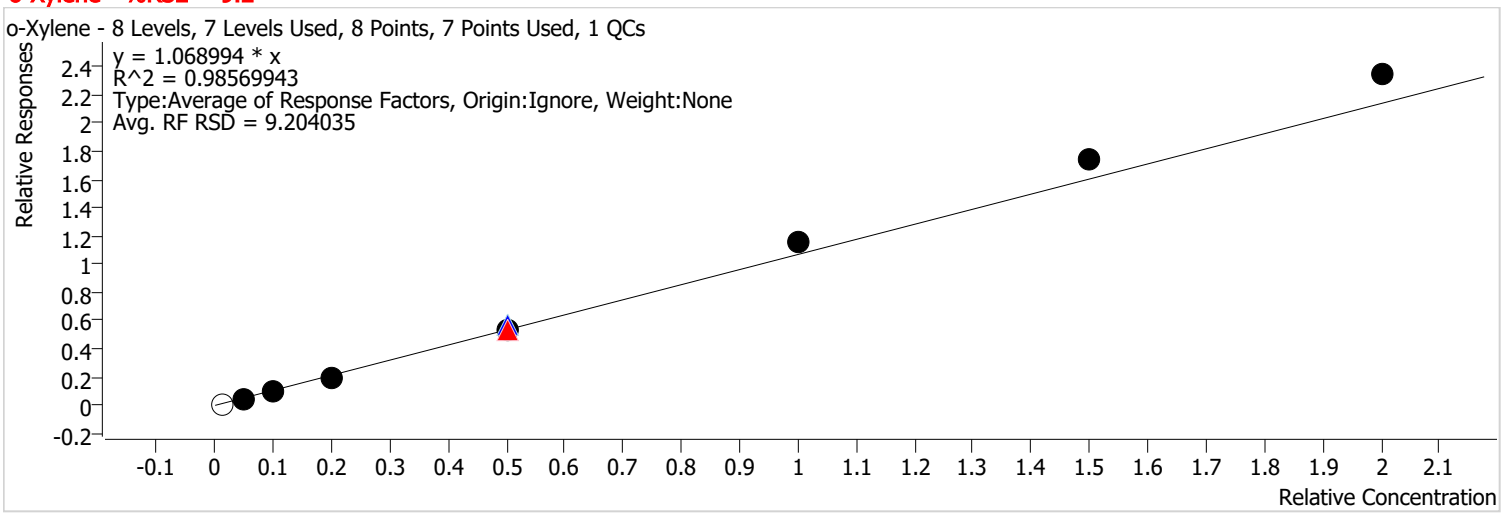
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	31538	25.0000	1.0635	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	66267	50.0000	1.0995	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	133498	100.0000	1.1112	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	388558	250.0000	1.2621	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	368418	250.0000	1.2021	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	368418	250.0000	1.2021	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	825866	500.0000	1.3051	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	1228570	750.0000	1.3014	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1637879	1000.0000	1.3058	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

o-Xylene %RSE = 9.2



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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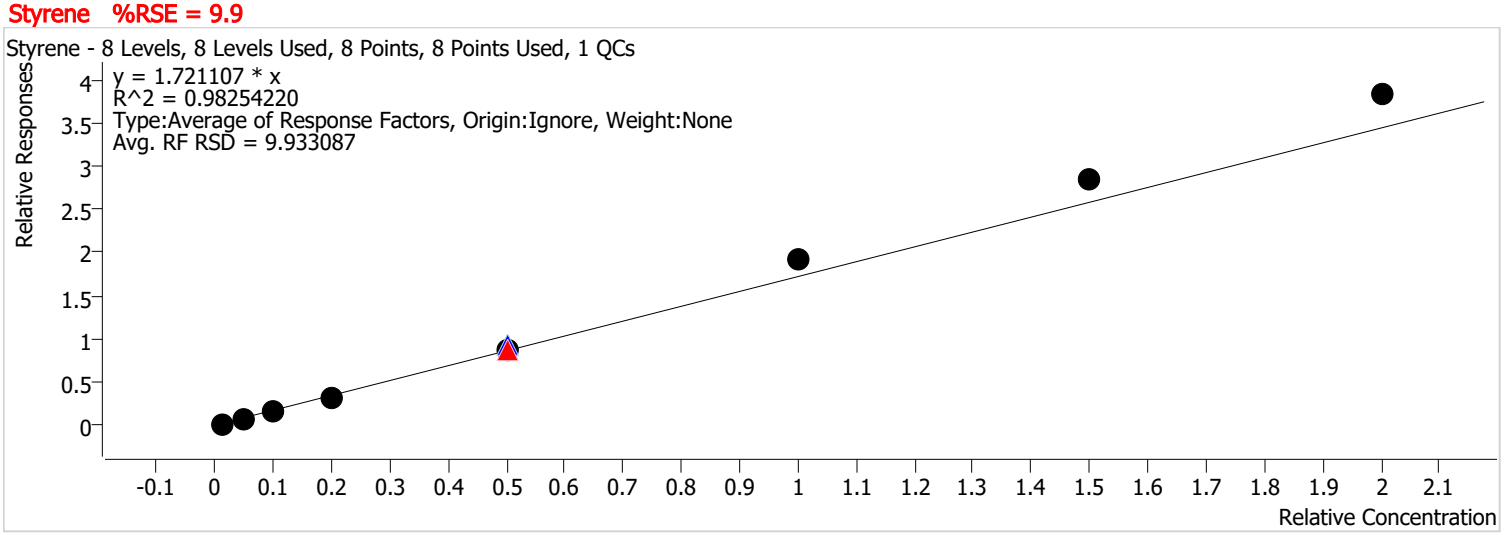
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	13519	12.5000	0.9117	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	30463	25.0000	1.0109	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	61016	50.0000	1.0157	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	174061	125.0000	1.1308	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	161509	125.0000	1.0539	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	161509	125.0000	1.0539	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	365914	250.0000	1.1565	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	549244	375.0000	1.1636	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	734101	500.0000	1.1705	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Styrene %RSE = 9.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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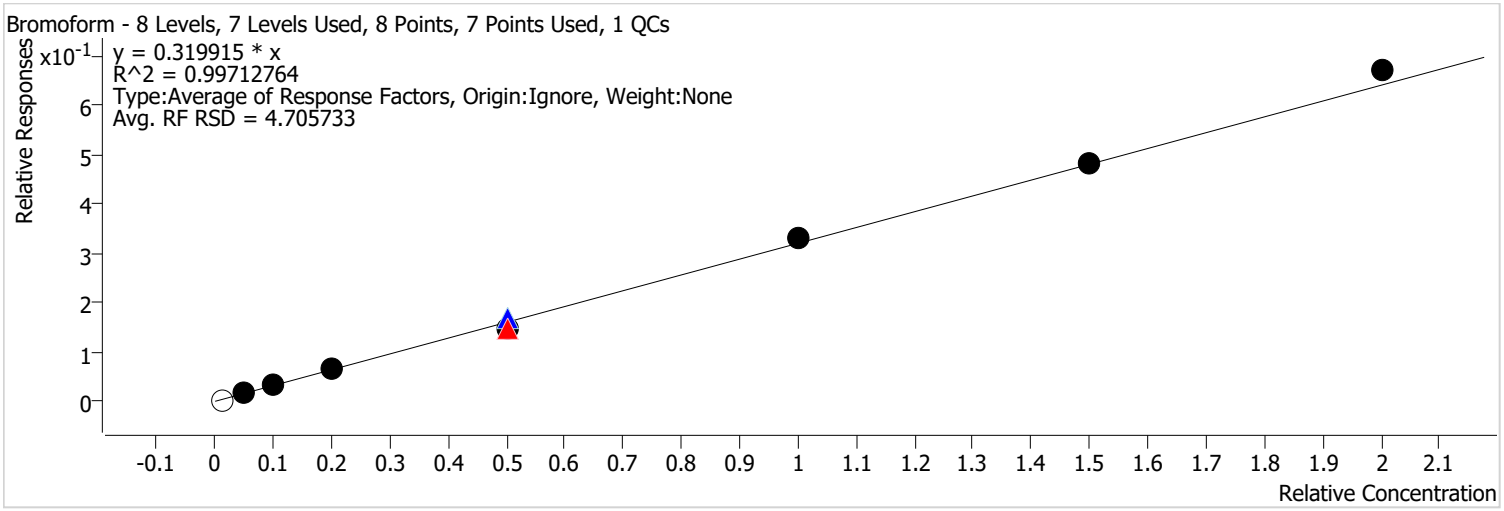
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	23472	12.5000	1.5830	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	48569	25.0000	1.6118	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	96576	50.0000	1.6077	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	291425	125.0000	1.8932	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	268375	125.0000	1.7513	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	268375	125.0000	1.7513	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	605646	250.0000	1.9142	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	896331	375.0000	1.8990	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1199879	500.0000	1.9132	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Bromoform %RSE = 4.7



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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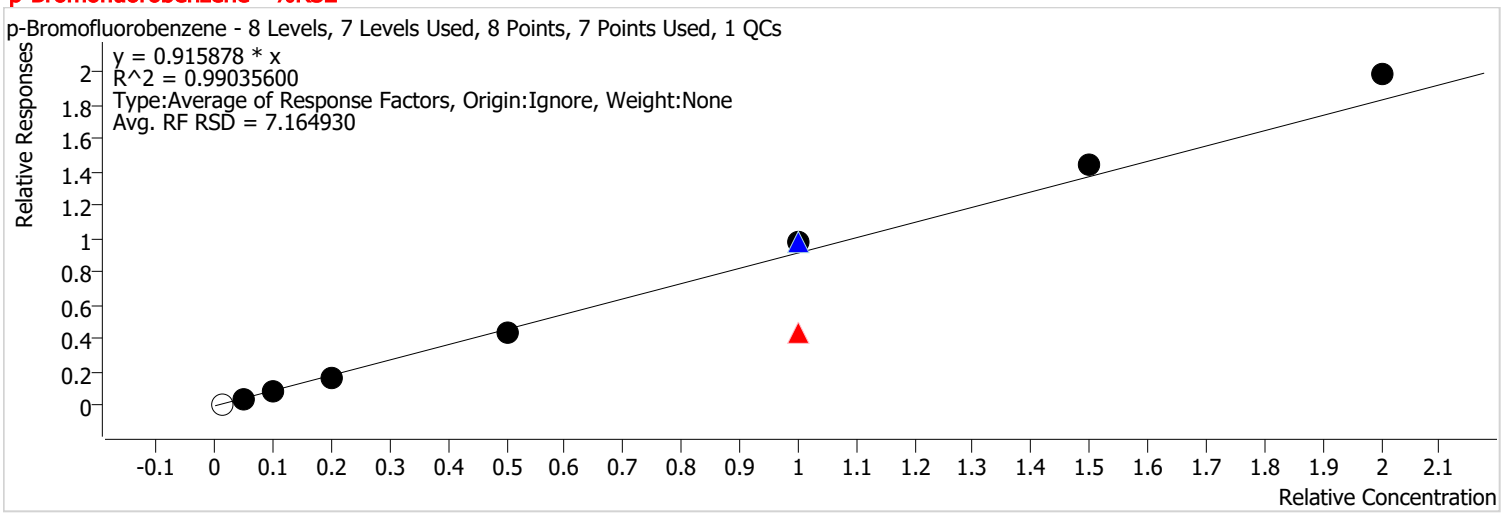
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	3652	12.5000	0.3016	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	7972	25.0000	0.3317	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	16073	50.0000	0.3232	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	42560	125.0000	0.3326	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	39165	125.0000	0.2962	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	39165	125.0000	0.2962	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	87836	250.0000	0.3295	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	129038	375.0000	0.3227	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	175918	500.0000	0.3345	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

p-Bromofluorobenzene %RSE =



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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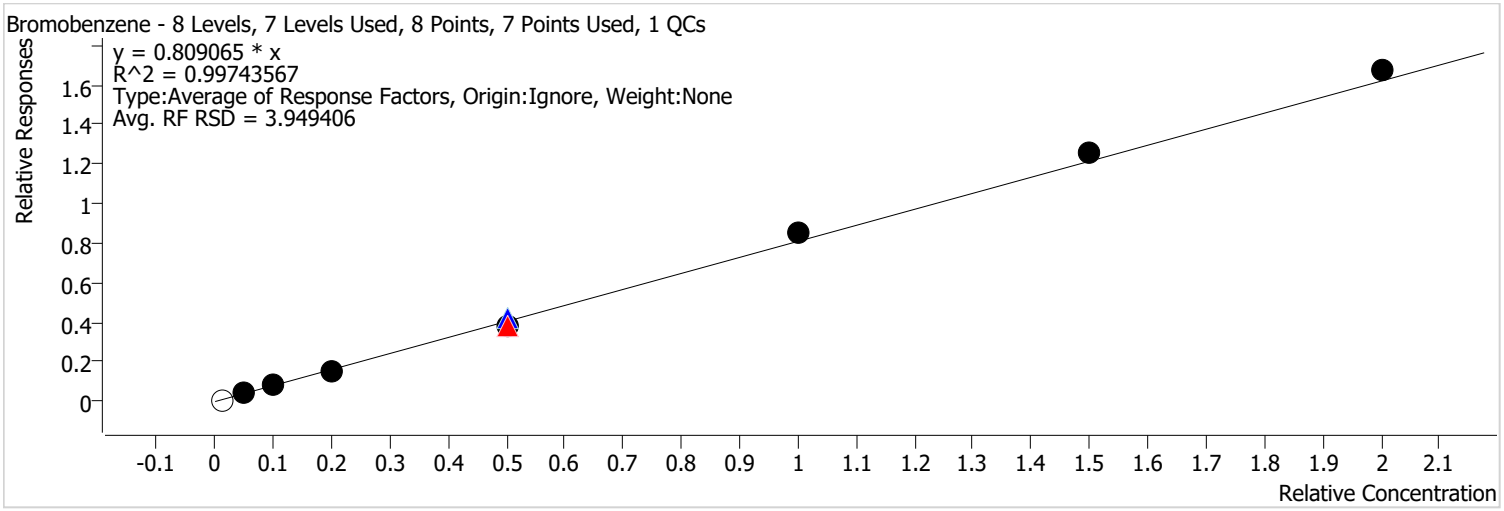
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	10059	12.5000	0.8308	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	22267	25.0000	0.9265	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	42506	50.0000	0.8548	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	114269	125.0000	0.8641	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	253034	250.0000	0.9888	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	261042	250.0000	0.9793	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	114269	250.0000	0.4321	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	385474	375.0000	0.9639	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	521580	500.0000	0.9917	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

Bromobenzene %RSE = 3.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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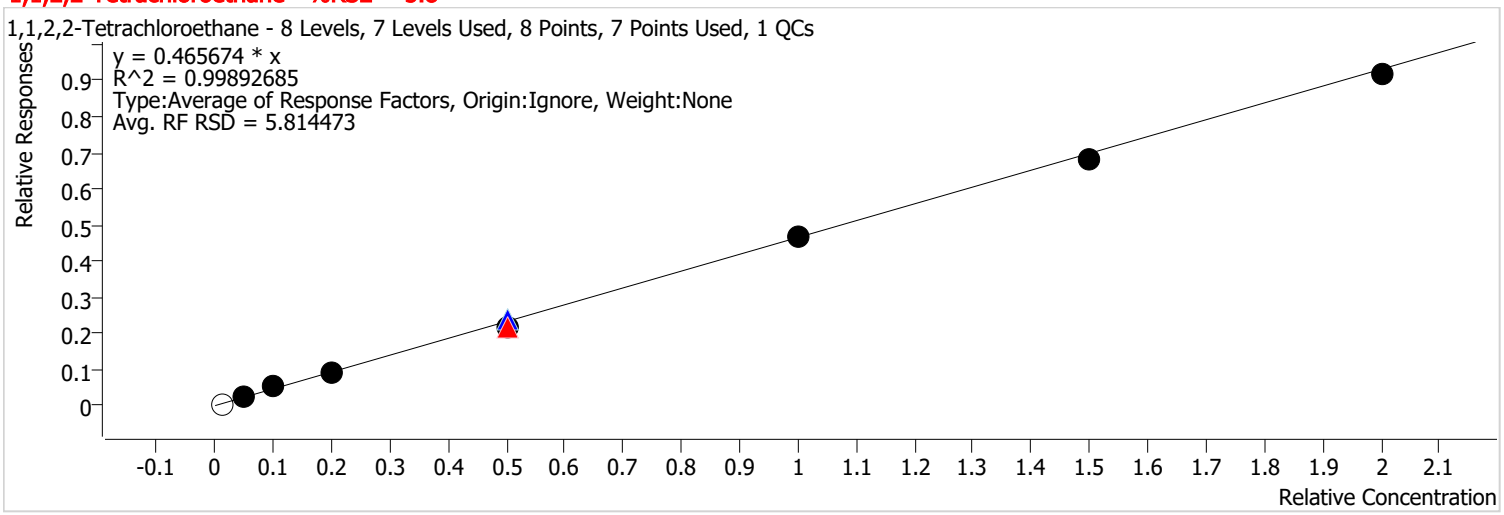
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	9663	12.5000	0.7981	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	19259	25.0000	0.8013	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	38282	50.0000	0.7698	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	109054	125.0000	0.8523	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	102265	125.0000	0.7733	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	102265	125.0000	0.7733	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	227127	250.0000	0.8521	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	333431	375.0000	0.8338	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	439147	500.0000	0.8350	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,1,2,2-Tetrachloroethane %RSE = 5.8



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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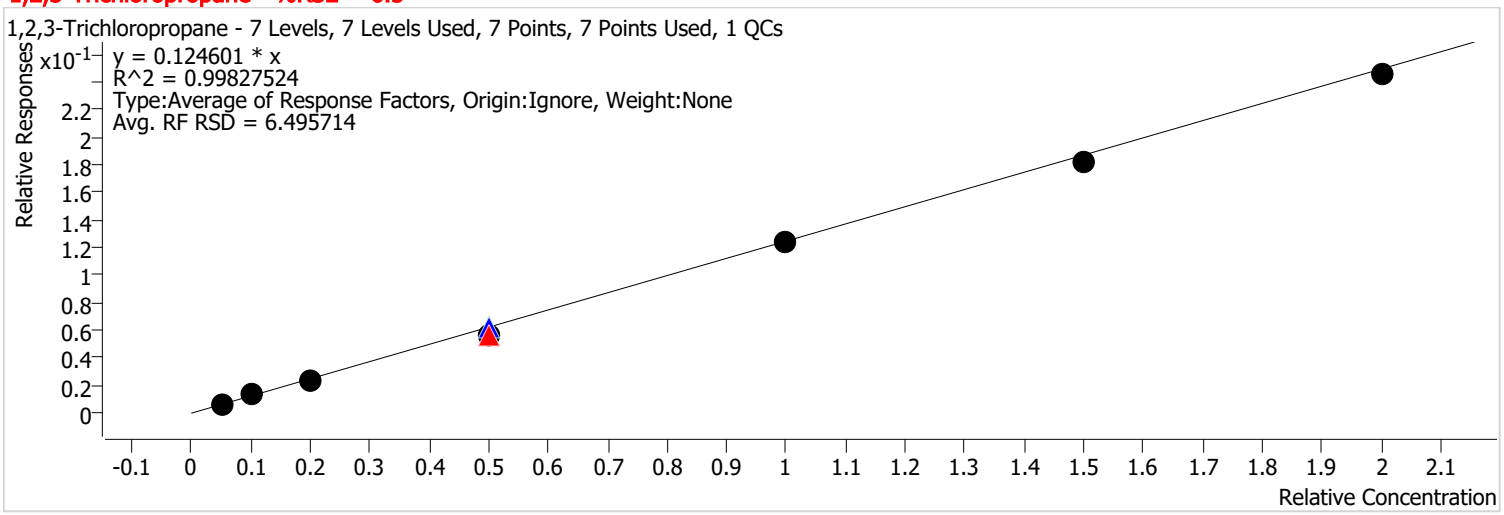
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	5793	12.5000	0.4785	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	12440	25.0000	0.5176	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	22514	50.0000	0.4528	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	60763	125.0000	0.4749	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	56958	125.0000	0.4307	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	56958	125.0000	0.4307	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	124205	250.0000	0.4660	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	182470	375.0000	0.4563	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	240837	500.0000	0.4579	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,2,3-Trichloropropane %RSE = 6.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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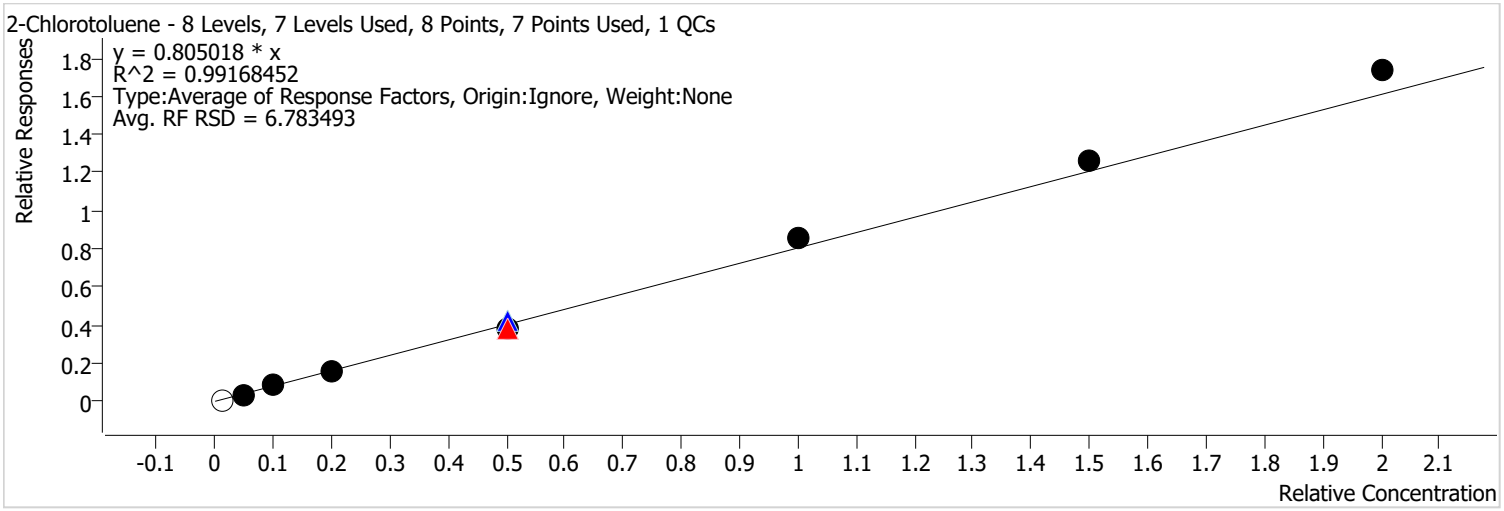
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	3200	25.0000	0.1331	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	6096	50.0000	0.1226	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	15682	125.0000	0.1226	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	14846	125.0000	0.1123	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	14846	125.0000	0.1123	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	33115	250.0000	0.1242	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	48325	375.0000	0.1208	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	64422	500.0000	0.1225	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

2-Chlorotoluene %RSE = 6.8



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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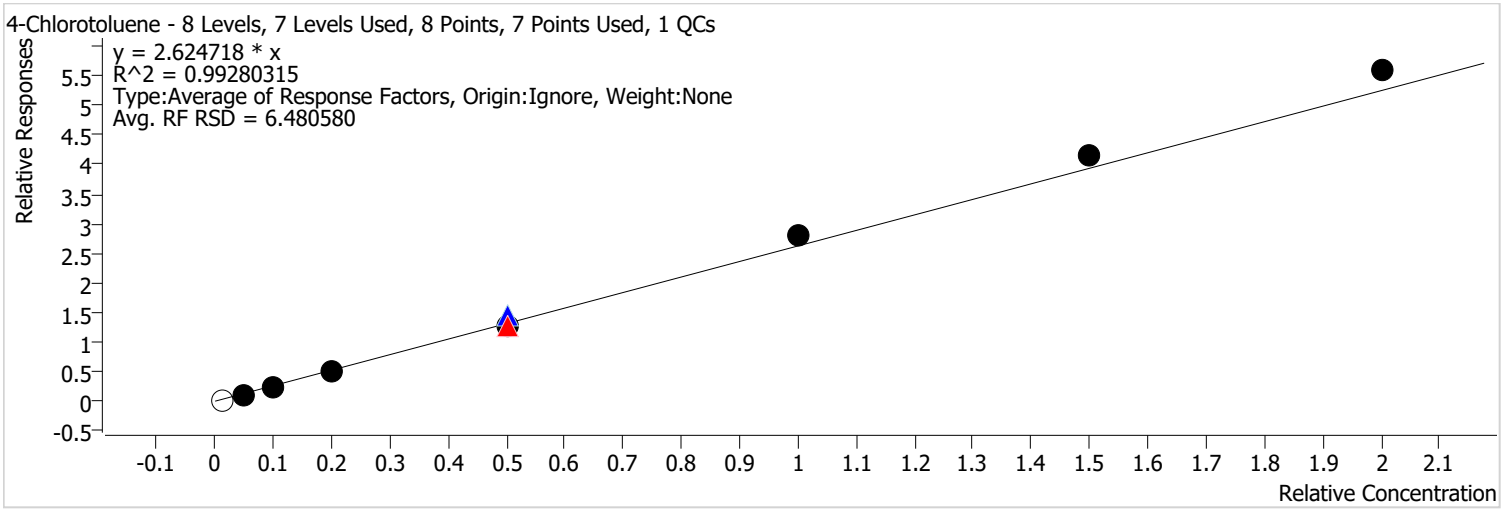
D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	8731	12.5000	0.7211	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	19390	25.0000	0.8068	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	37987	50.0000	0.7639	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	108192	125.0000	0.8456	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	102424	125.0000	0.7745	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	102424	125.0000	0.7745	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	229396	250.0000	0.8606	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	336386	375.0000	0.8411	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	455991	500.0000	0.8670	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

4-Chlorotoluene %RSE = 6.5



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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D:\Org\Data\VOA5975C\VG010422\04JAN11.D	Calibration	2	x	28532	12.5000	2.3566	
D:\Org\Data\VOA5975C\VG010422\04JAN12.D	Calibration	3	x	61551	25.0000	2.5611	
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	126308	50.0000	2.5400	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	368295	125.0000	2.8784	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	336146	125.0000	2.5420	
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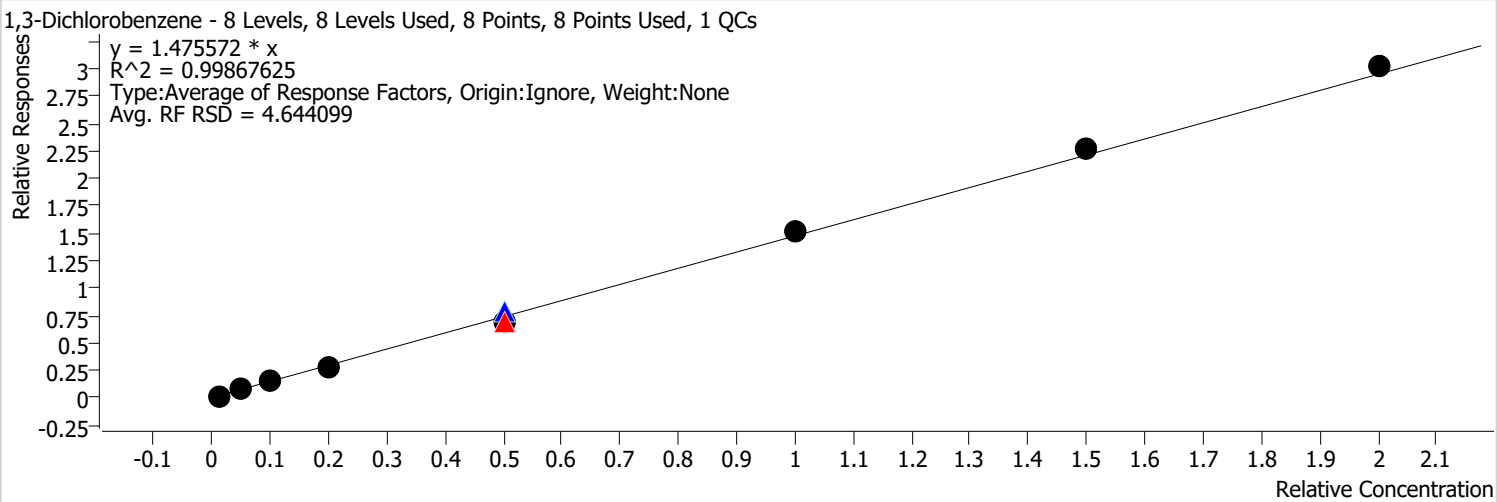
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	748435	250.0000	2.8078	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	1109221	375.0000	2.7736	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	1468376	500.0000	2.7919	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,3-Dichlorobenzene %RSE = 4.6



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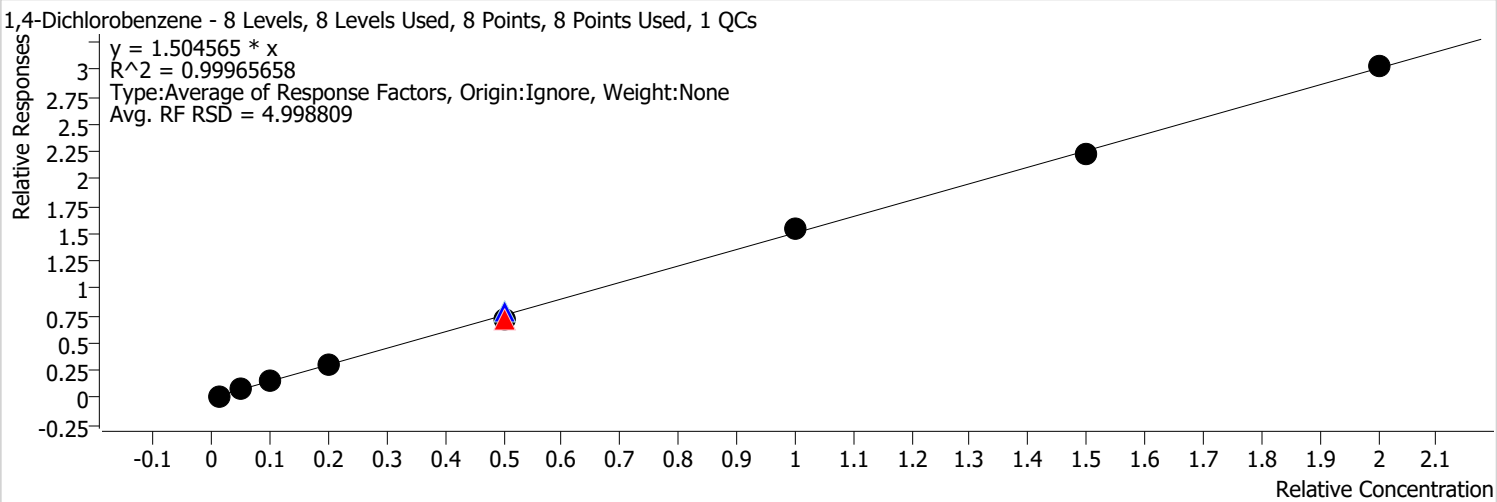
D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	69539	50.0000	1.3984	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	204088	125.0000	1.5950	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	183404	125.0000	1.3869	
D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D	CC	CC	x	183404	125.0000	1.3869	
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	406895	250.0000	1.5265	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	603674	375.0000	1.5095	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	793993	500.0000	1.5097	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,4-Dichlorobenzene %RSE = 5.0



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	71841	50.0000	1.4447	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	200032	125.0000	1.5633	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	189045	125.0000	1.4296	
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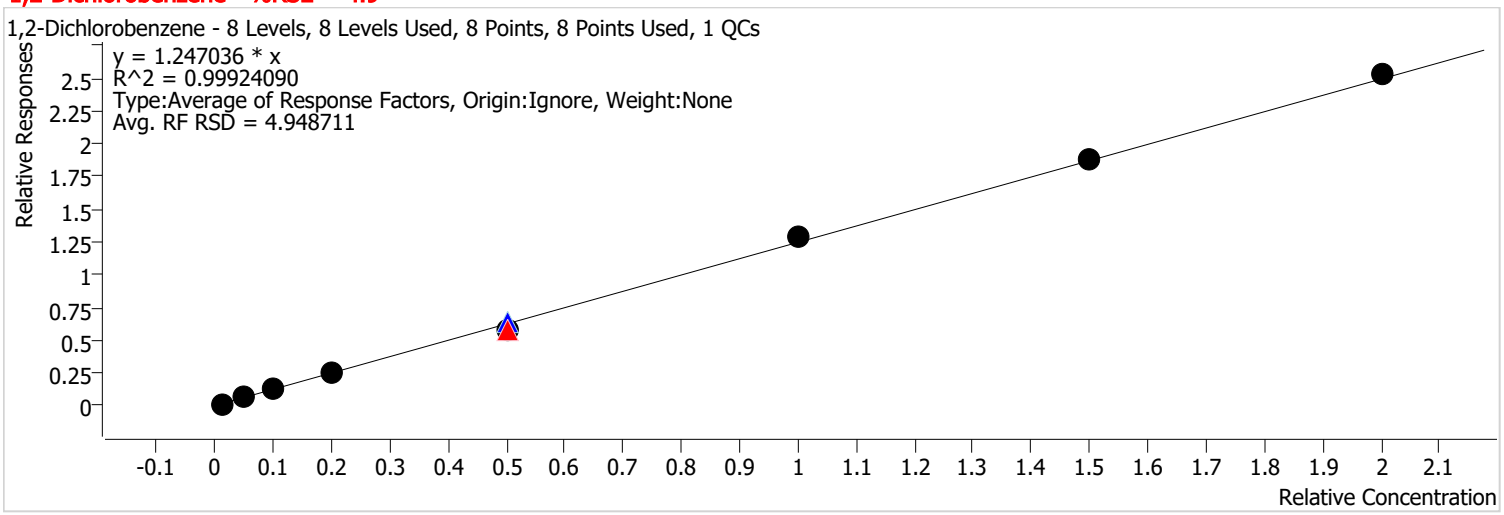
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	408934	250.0000	1.5342	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	595919	375.0000	1.4901	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	794954	500.0000	1.5115	

Batch Path	D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin		
Analysis Time	2/28/2022 1:57 PM	**Analyst Name**	BL2000\mchavez
Report Time	2/28/2022 2:00:45 PM	**Reporter Name**	BL2000\mchavez
Last Calib Update	1/9/2022 8:59 PM	**Batch State**	Processed
Quant Batch Version	10.0	**Quant Report Version**	10.0

1,2-Dichlorobenzene %RSE = 4.9



Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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D:\Org\Data\VOA5975C\VG010422\04JAN13.D	Calibration	4	x	60213	50.0000	1.2109	
D:\Org\Data\VOA5975C\VG010422\04JAN23.D	QC	QC	x	164299	125.0000	1.2841	
D:\Org\Data\VOA5975C\VG010422\04JAN15.D	Calibration	5	x	152284	125.0000	1.1516	
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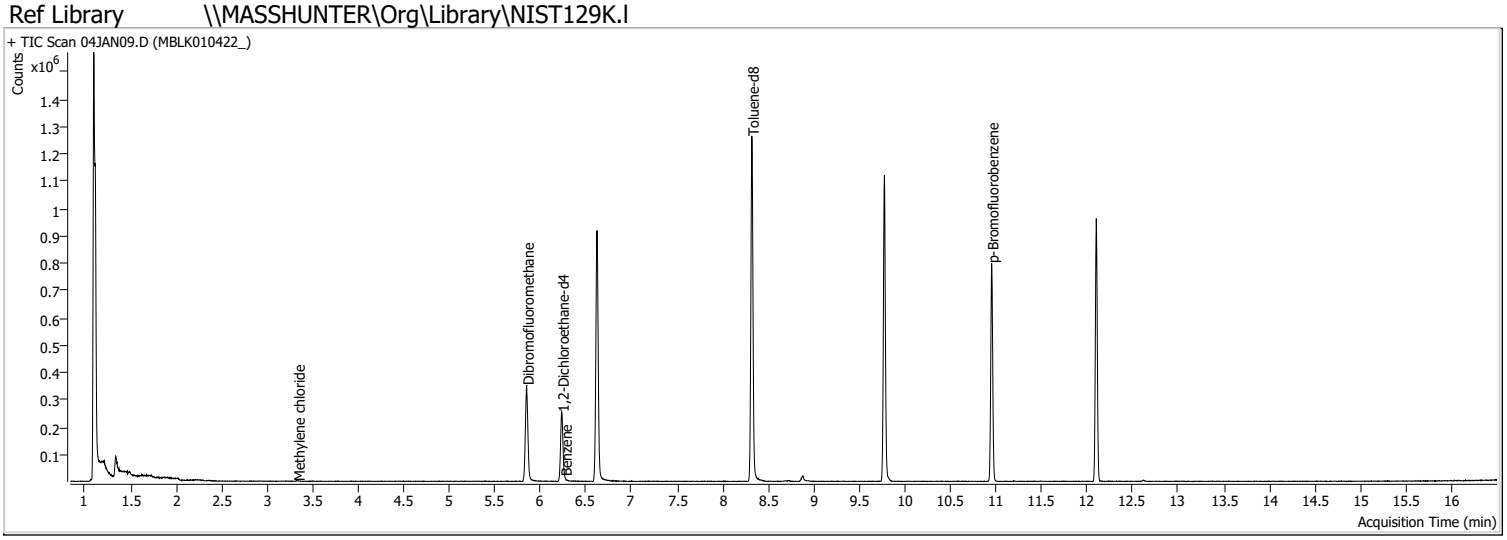
D:\Org\Data\VOA5975C\VG010422\04JAN17.D	Calibration	6	x	342576	250.0000	1.2852	
D:\Org\Data\VOA5975C\VG010422\04JAN19.D	Calibration	7	x	499147	375.0000	1.2481	
D:\Org\Data\VOA5975C\VG010422\04JAN21.D	Calibration	8	x	664247	500.0000	1.2630	

Quantitation Results Report (QT Reviewed)

Agilent | Trusted Answers

Data File	04JAN09.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/4/2022 3:05:37 PM
Sample Name	MBLK010422_	Instrument	VOA5975C
Vial	9	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

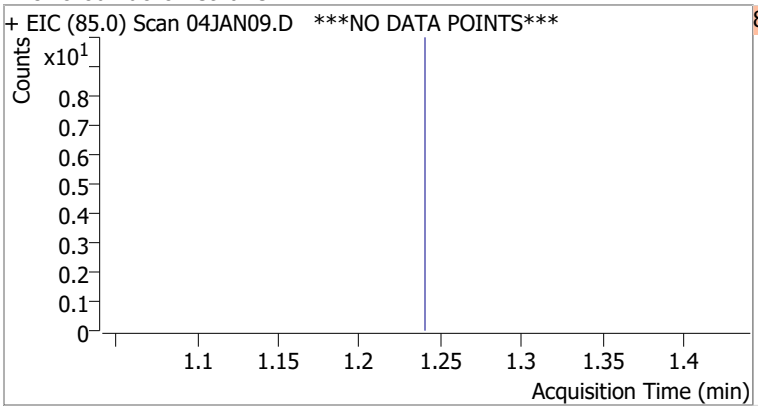


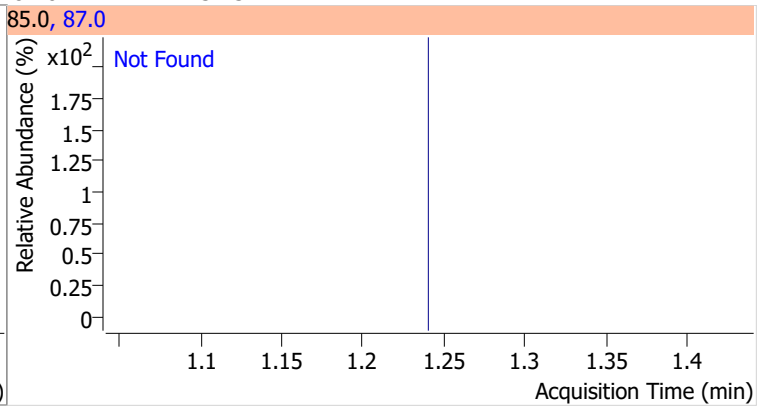
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.620	96.0	775552	250.0000	ng		-0.003
M Chlorobenzene-d5	9.774	82.0	301196	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	231562	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.848	113.0	203459	278.4635	ng		0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 111.39%			
S 1,2-Dichloroethane-d4	6.233	67.0	88174	279.3964	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 111.76%			
S Toluene-d8	8.319	98.0	770154	265.3436	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 106.14%			
S p-Bromofluorobenzene	10.951	95.0	226743	267.2815	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 106.91%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.333	50.0	0		ng	md	1
T Vinyl chloride	1.489	62.0	0		ng	md	1
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.335	49.0	1661	1.4424	ng	m	97
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	0.000		0	N.D.			

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.266	78.0	381	0.1233	ng	m	98
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	0.000		0	N.D.			
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	0.000		0	N.D.			
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	0.000		0	N.D.			
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	0.000		0	N.D.			
T m+p-Xylenes	0.000		0	N.D.			
T o-Xylene	0.000		0	N.D.			
T Styrene	0.000		0	N.D.			
T Bromoform	0.000		0	N.D.			
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

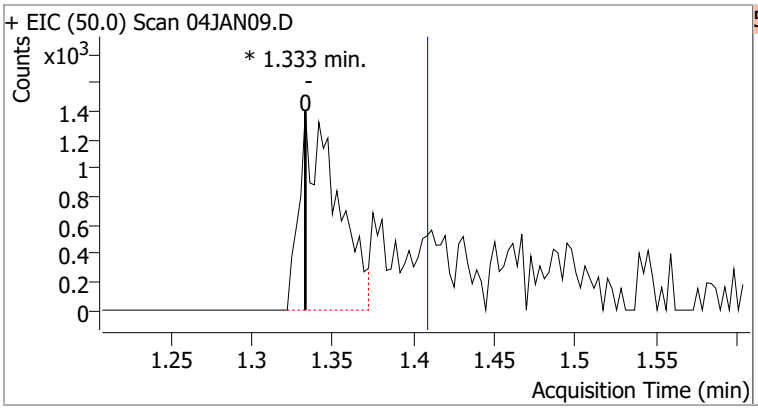
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

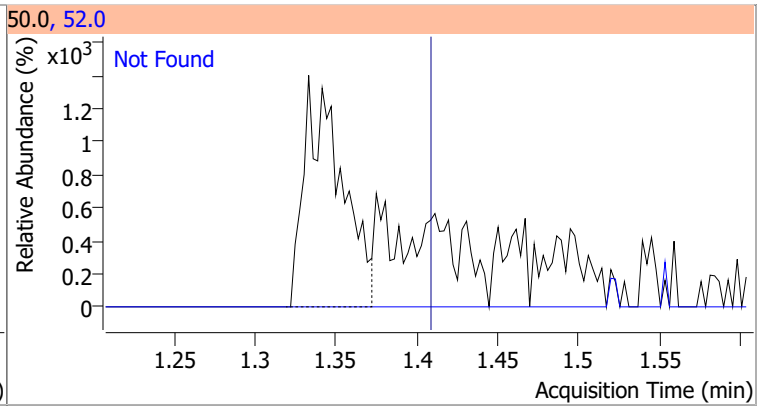
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3



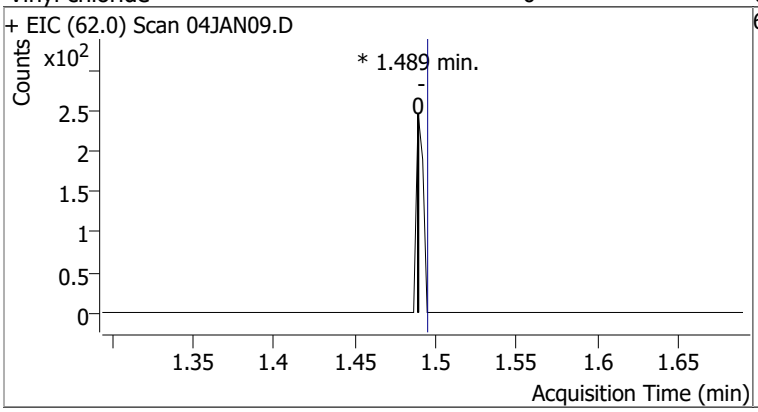


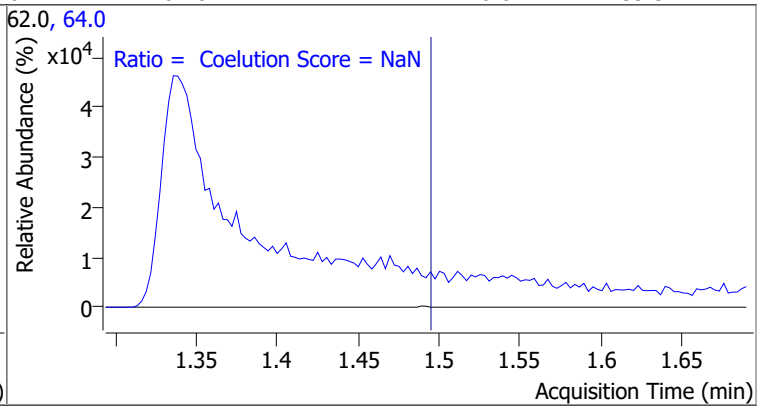
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane		0		0	52.0		2.1	62.1



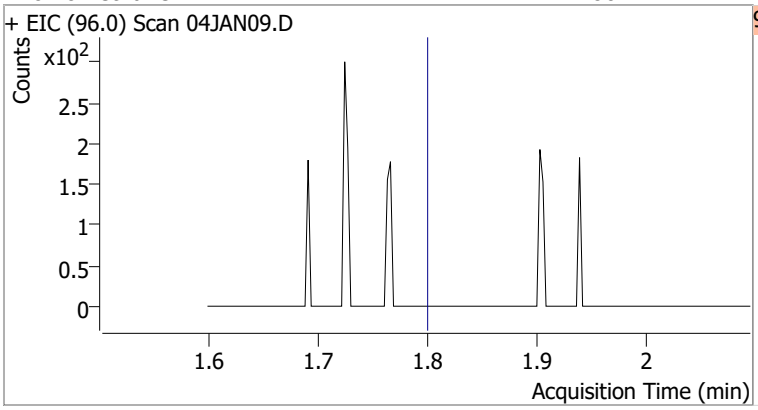


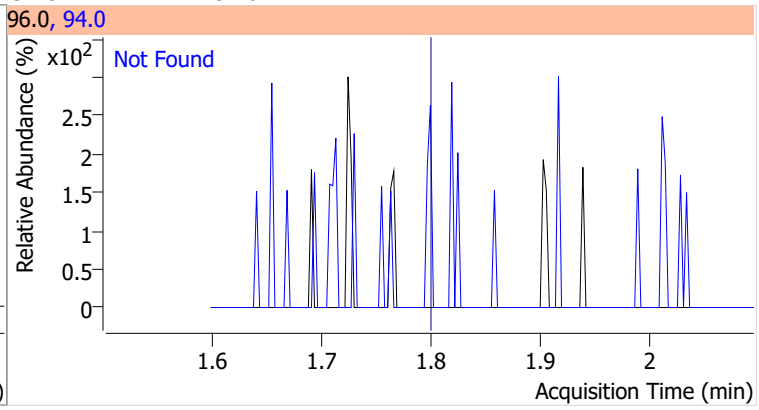
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride		0		0	64.0		0.0	59.9



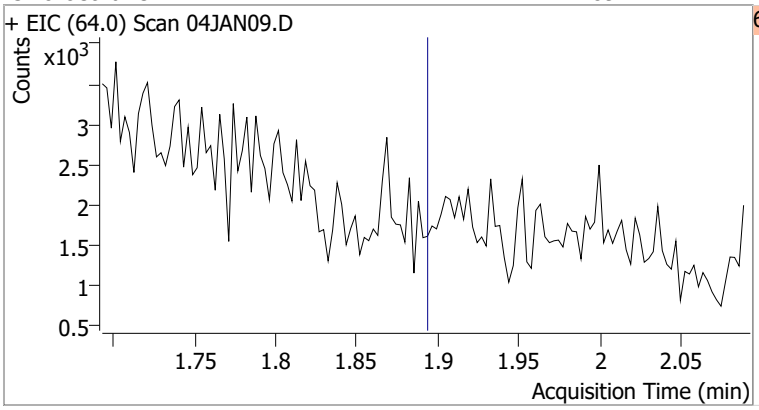


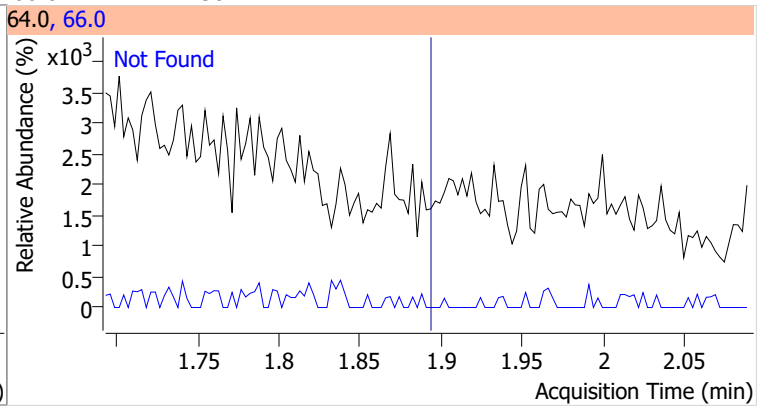
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



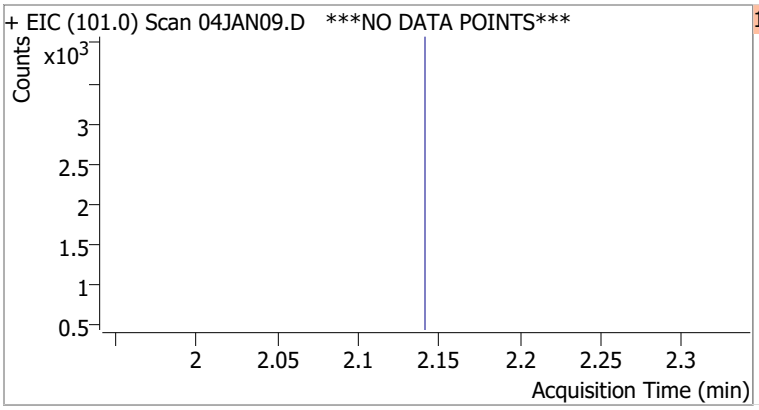


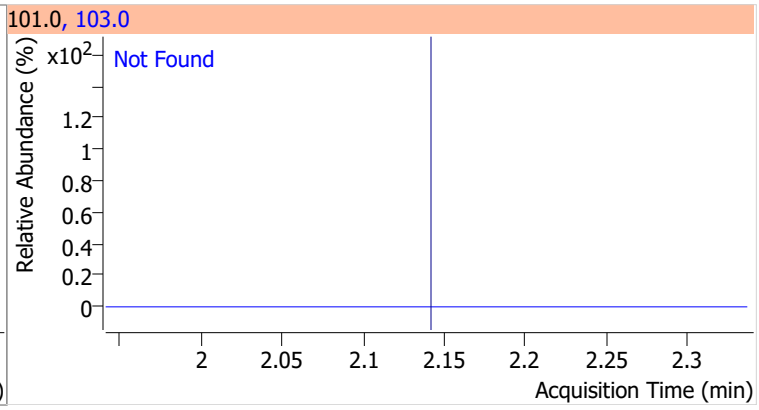
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



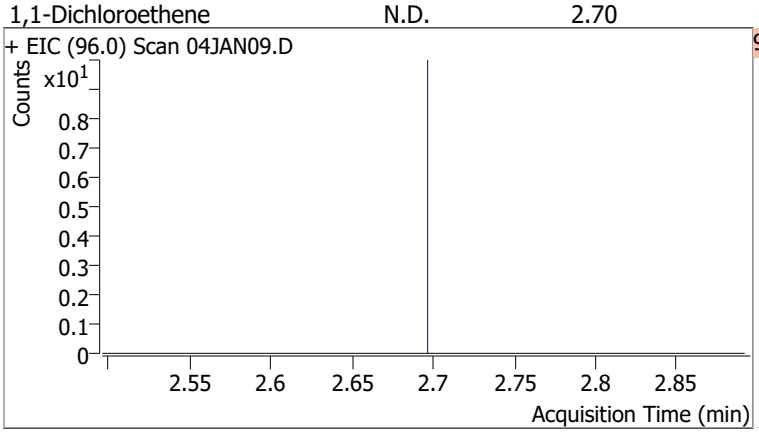


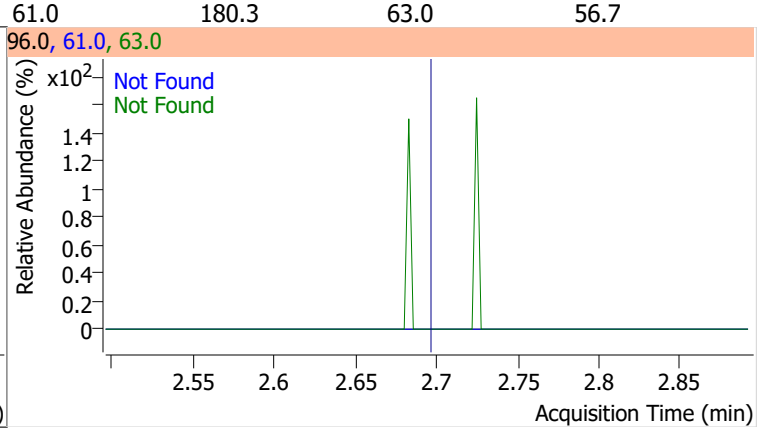
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



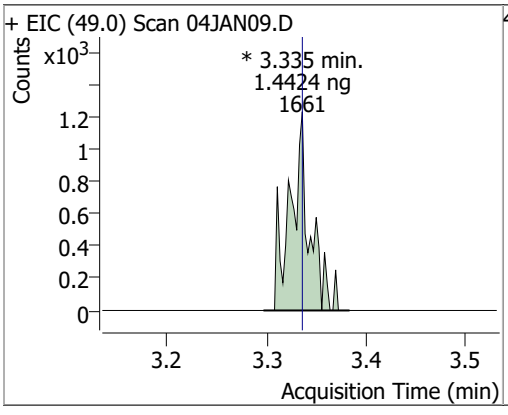


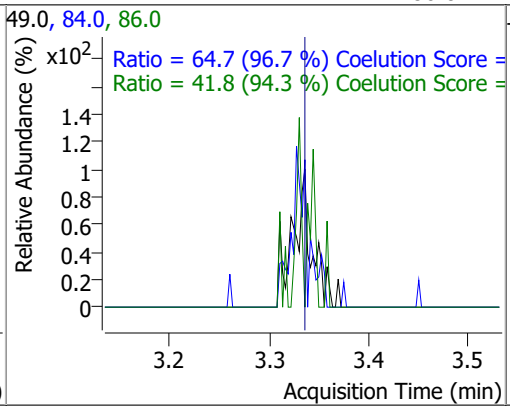
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

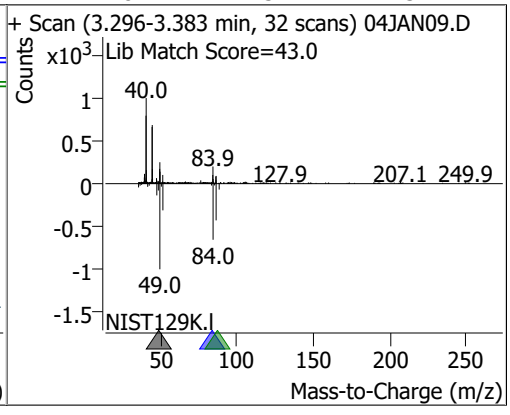




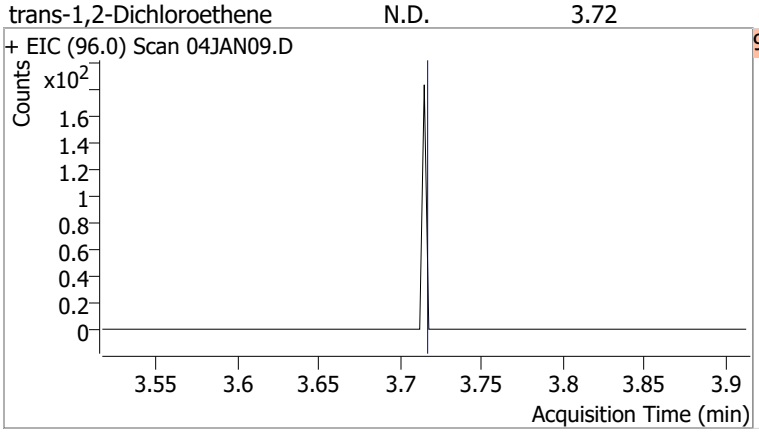
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	1.4424	3.34	0.00	1661 (m)	84.0	64.7	36.9	96.9
					86.0	41.8	14.3	74.3

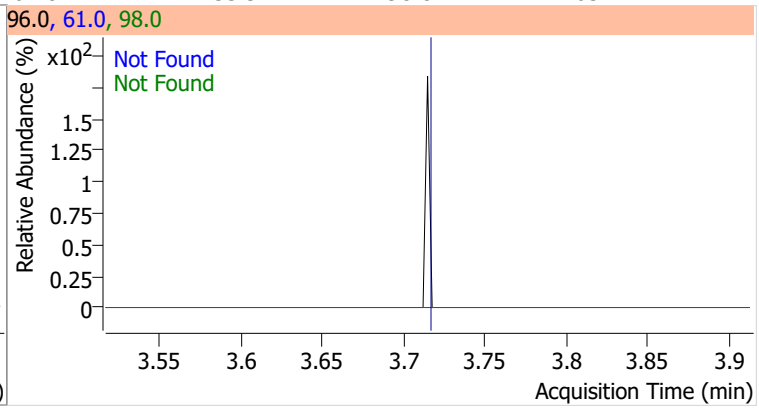




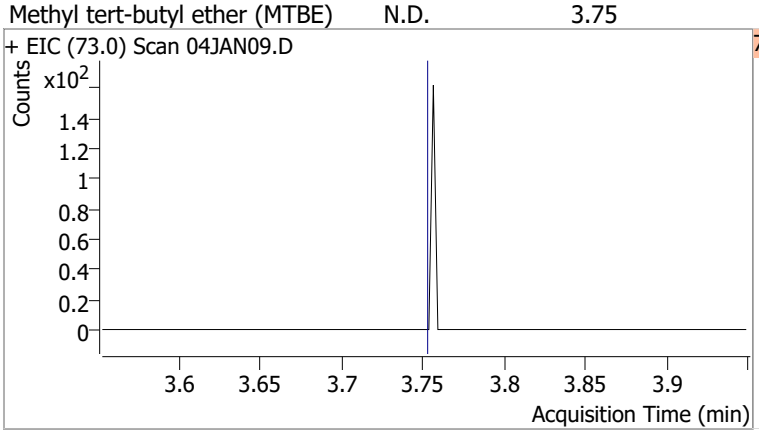


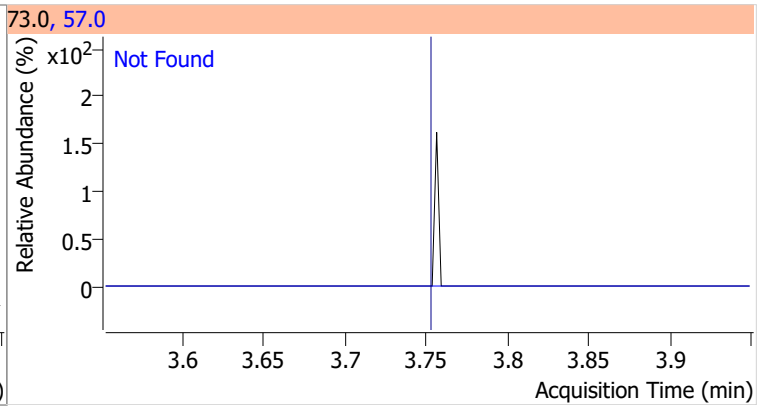
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



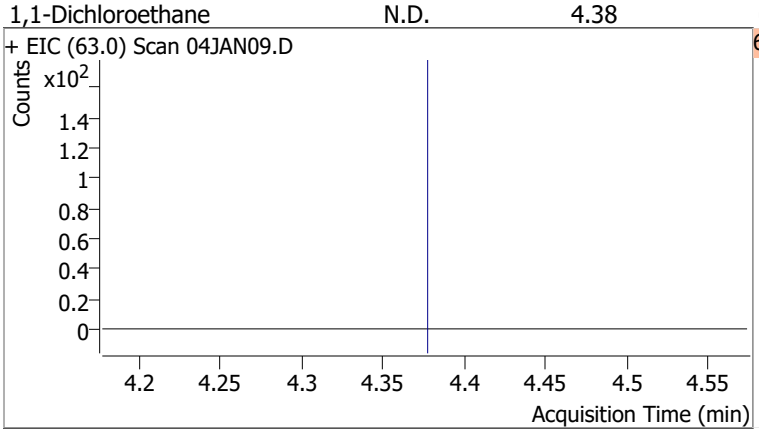


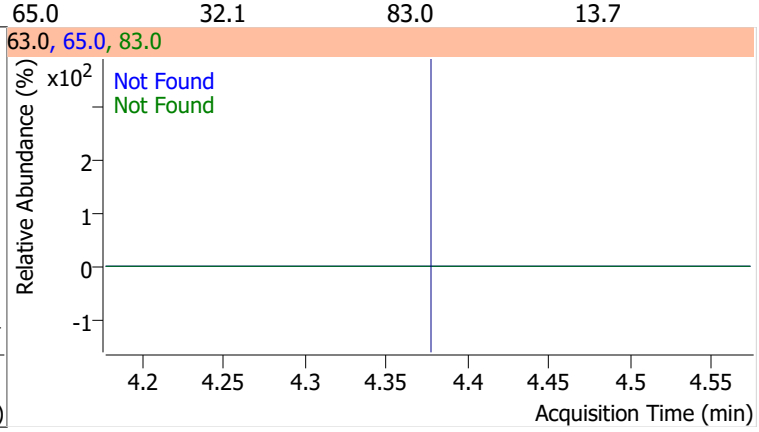
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



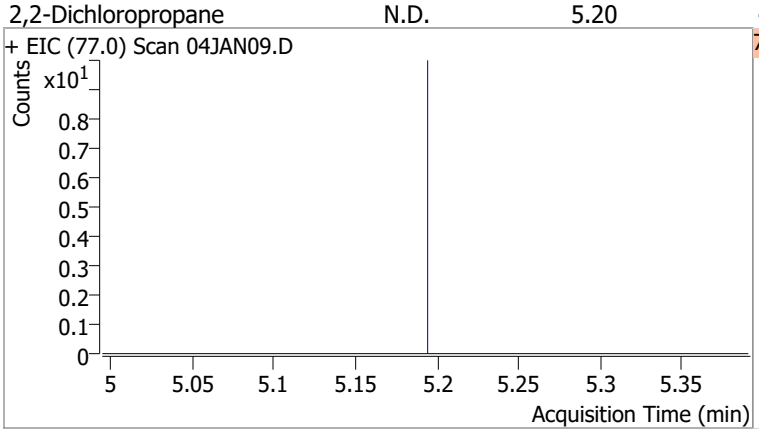


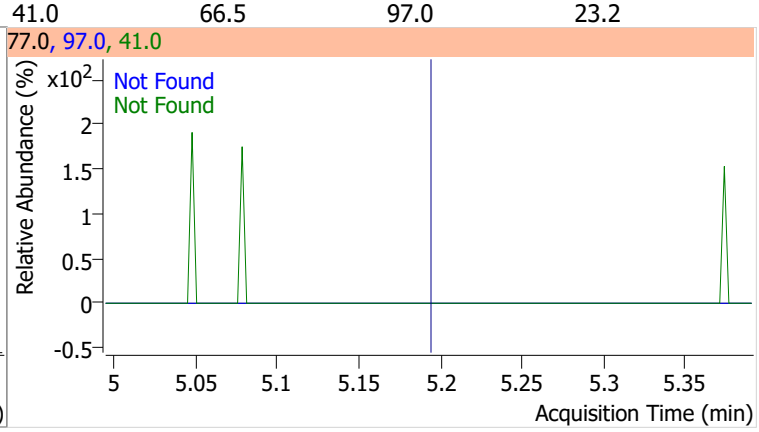
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



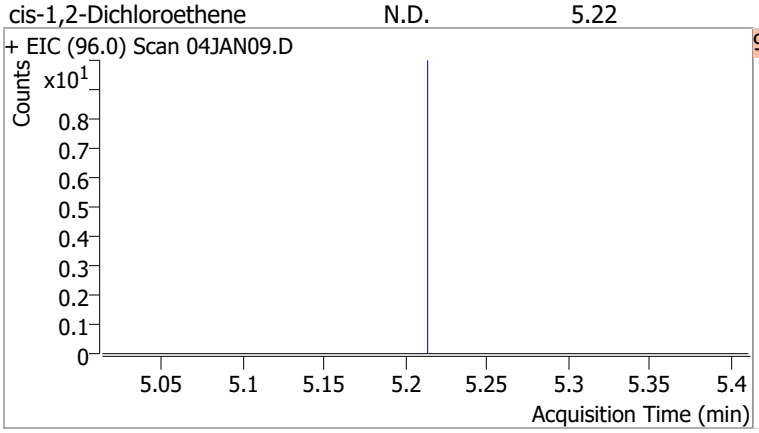


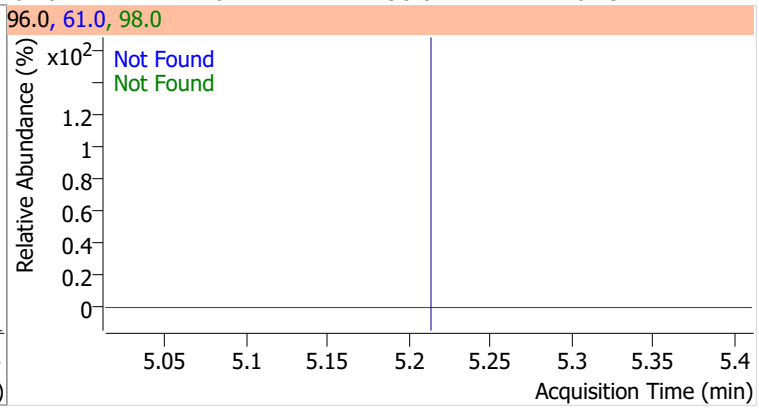
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



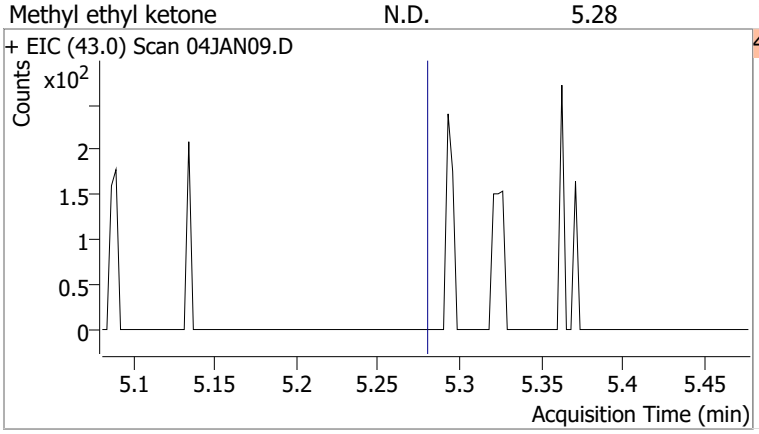


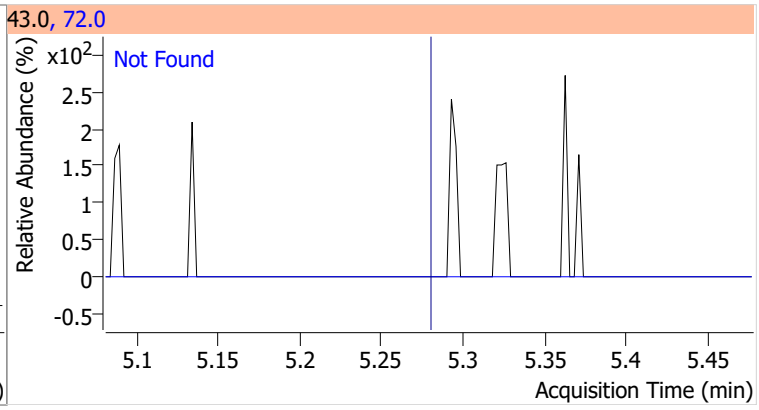
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



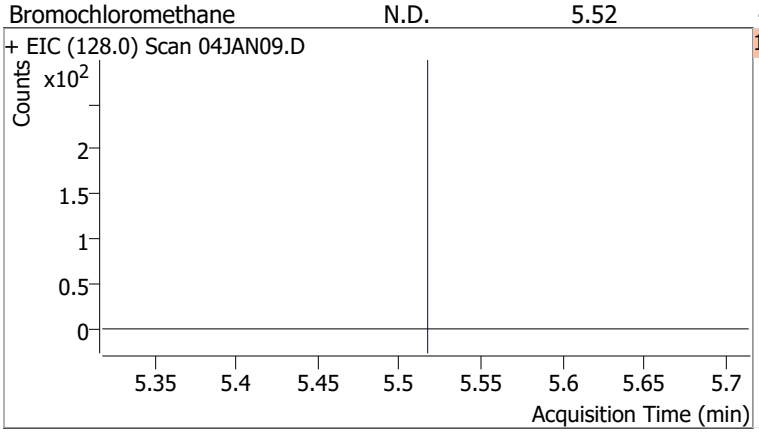


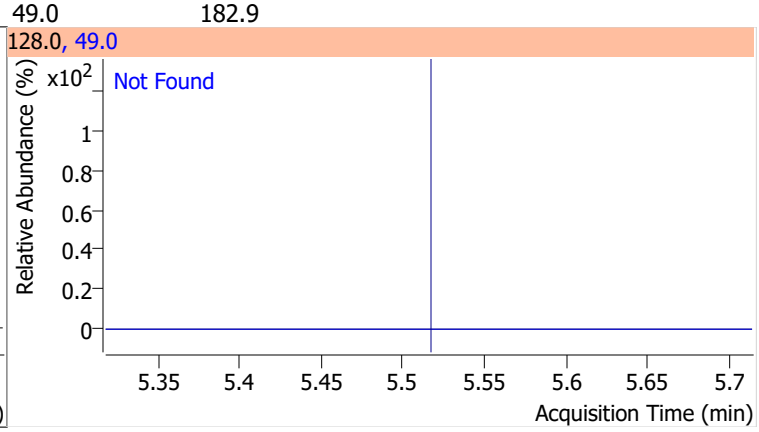
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



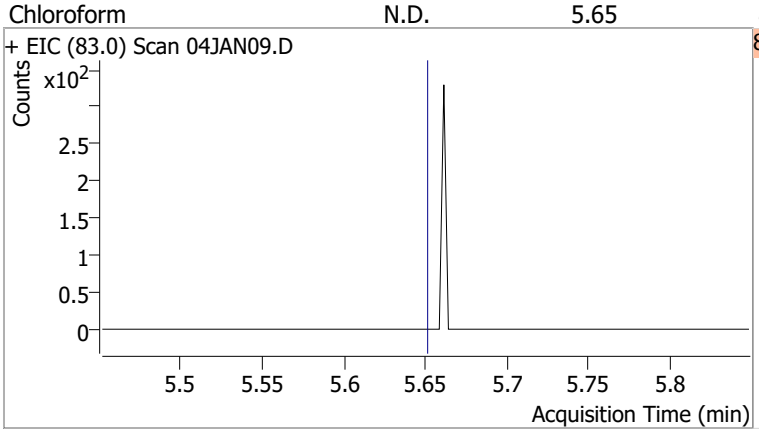


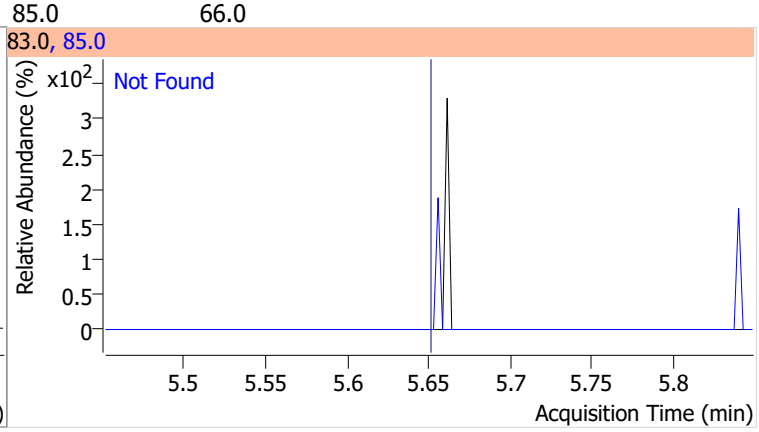
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9



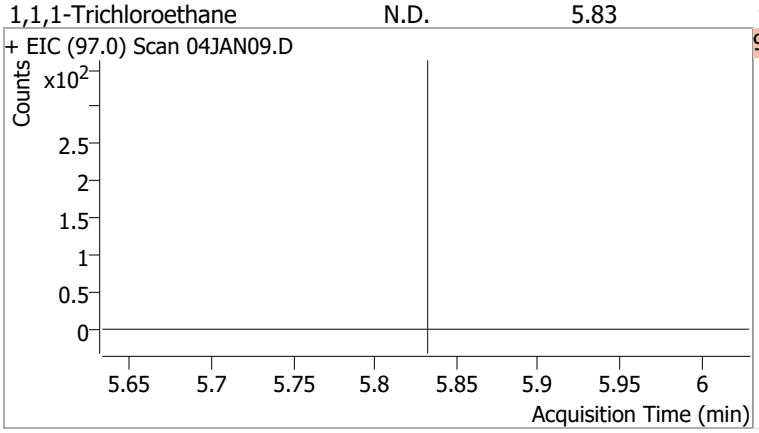


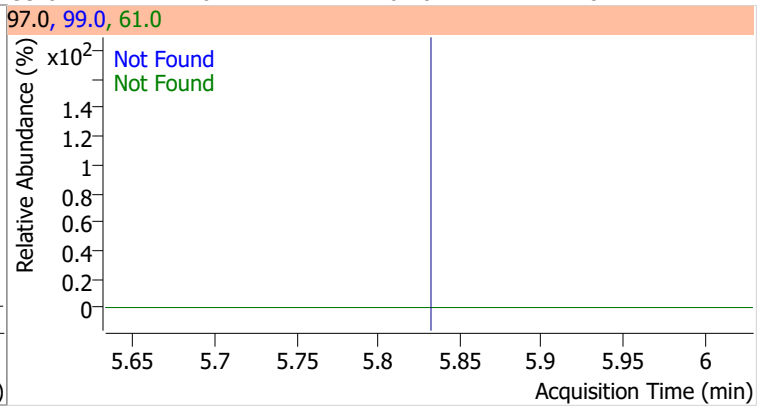
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroform	N.D.	5.65	85.0	66.0



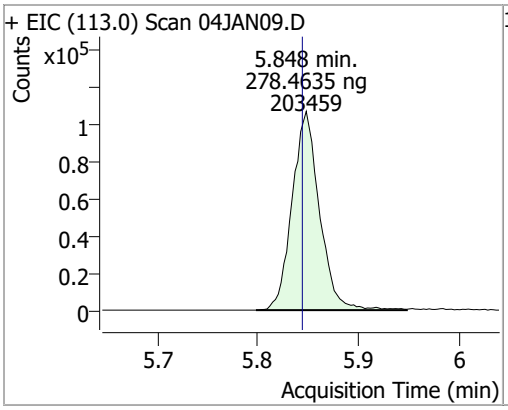


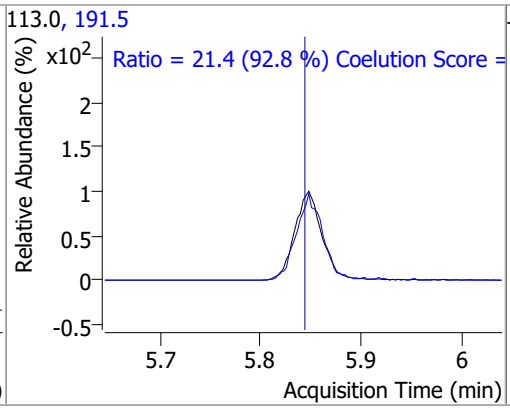
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

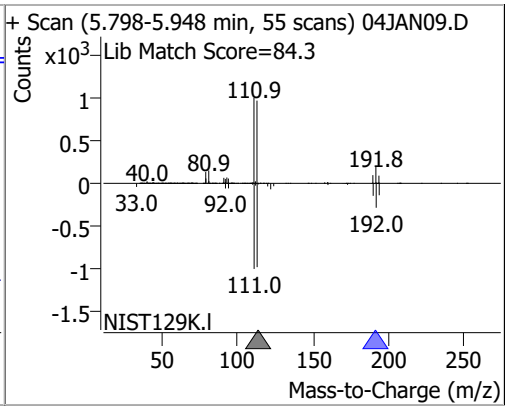




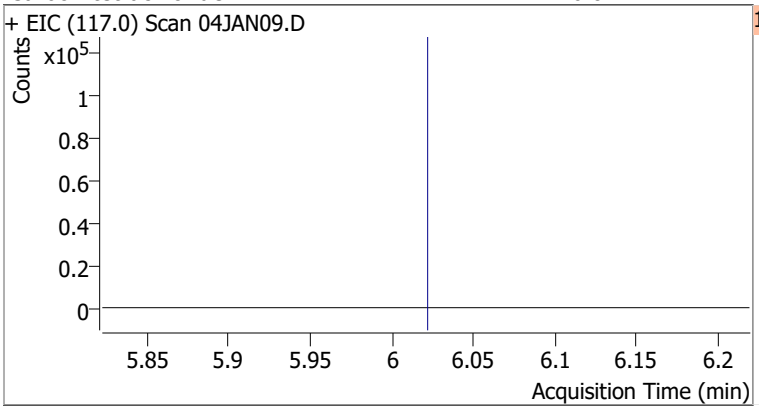
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	278.4635	5.85	0.00	203459	191.5	21.4	0.0	53.1

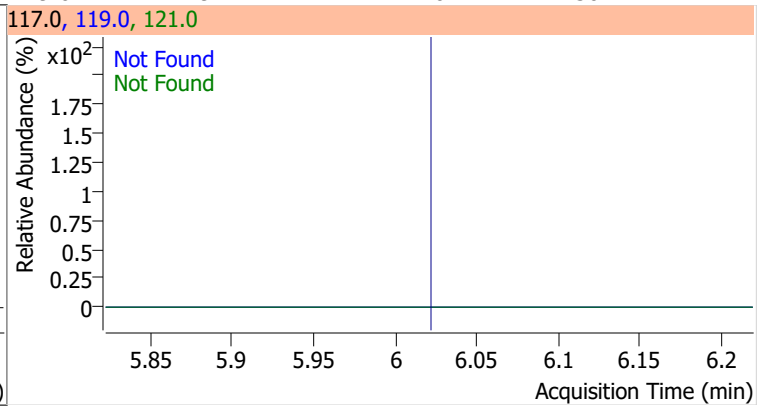




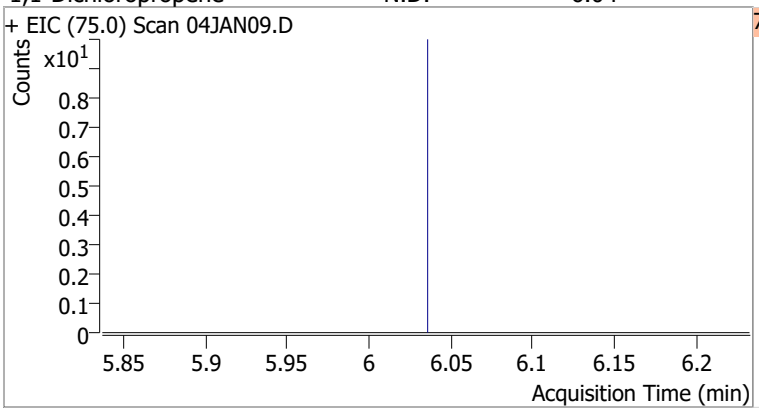


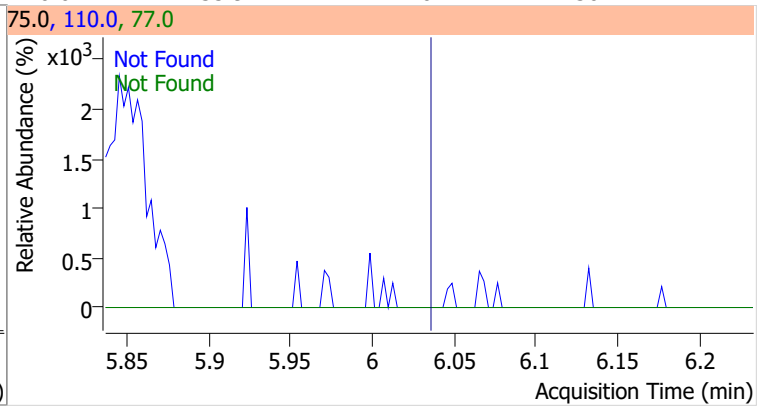
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



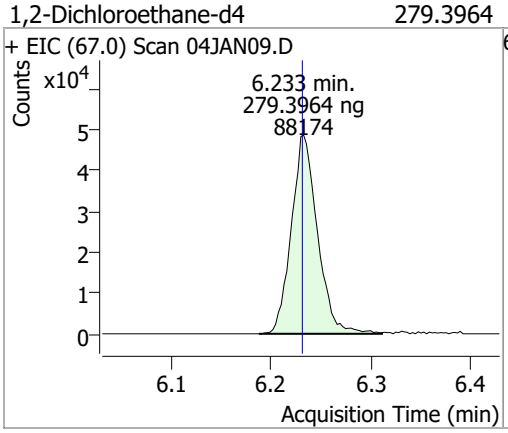


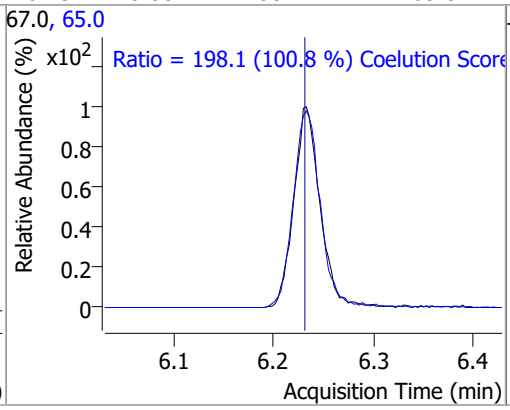
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

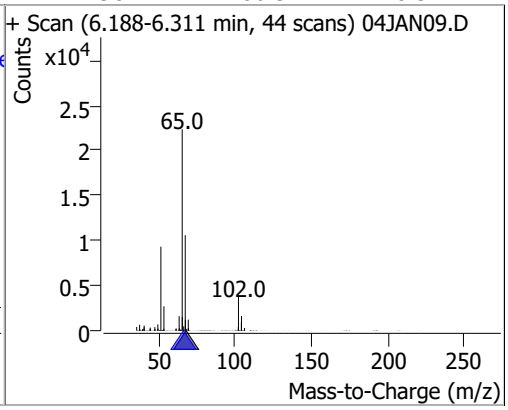




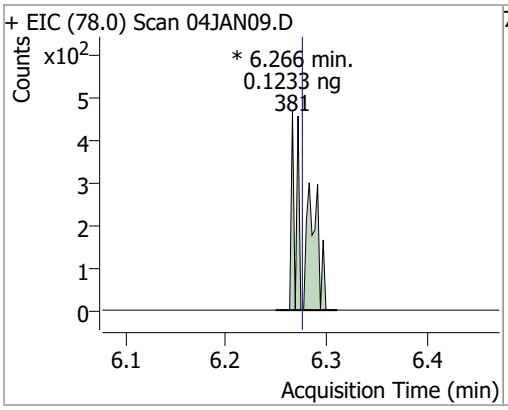
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	279.3964	6.23	0.00	88174	65.0	198.1	166.5	226.5

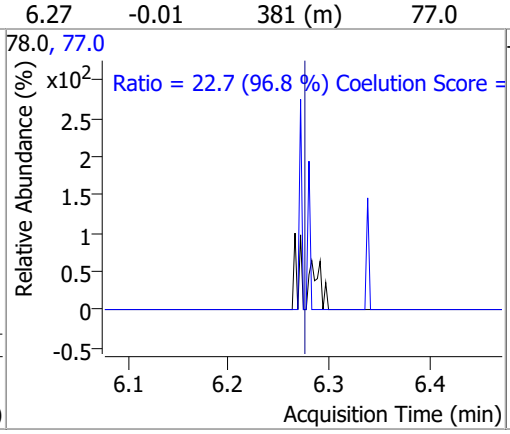


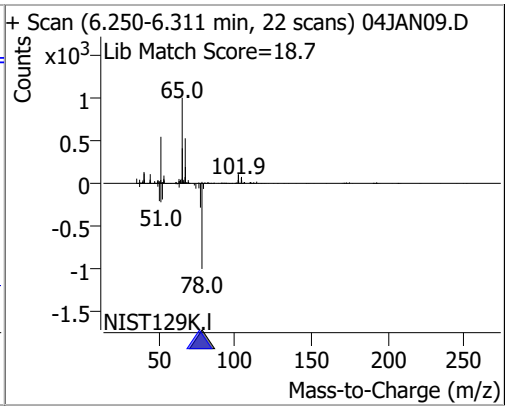




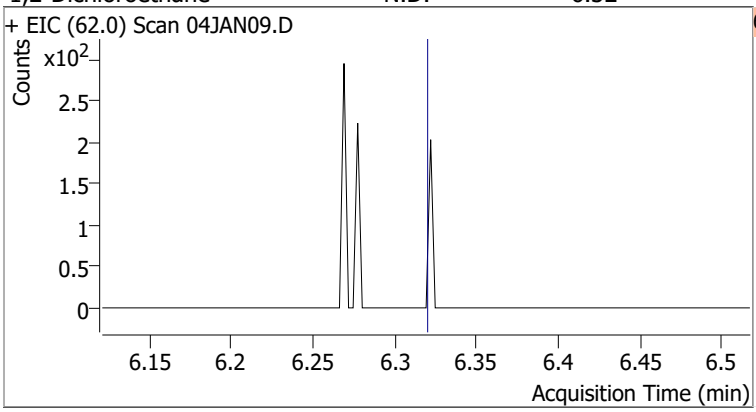
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.1233	6.27	-0.01	381 (m)	77.0	22.7	0.0	53.5

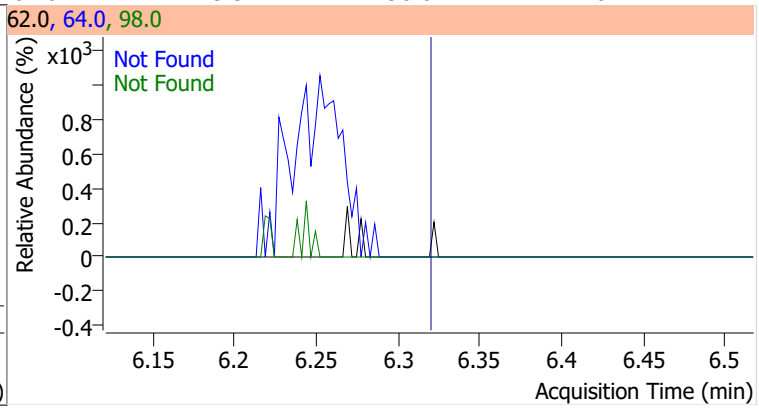




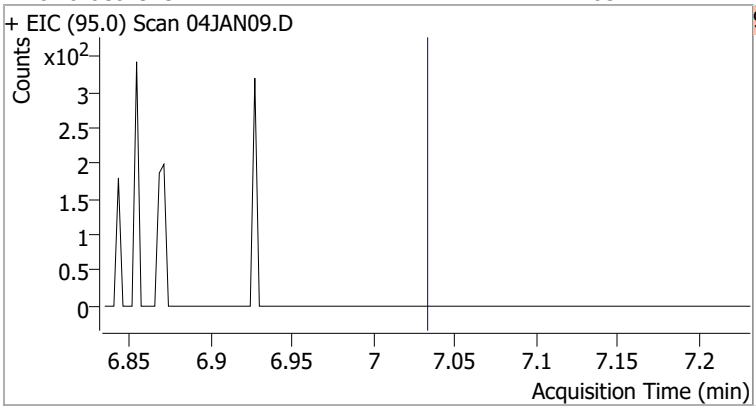


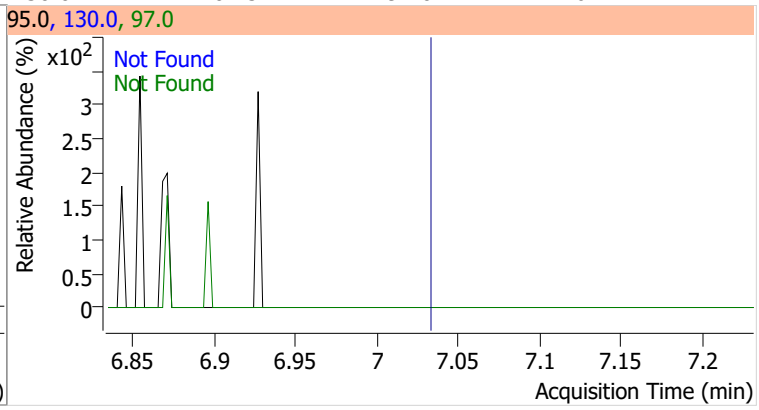
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



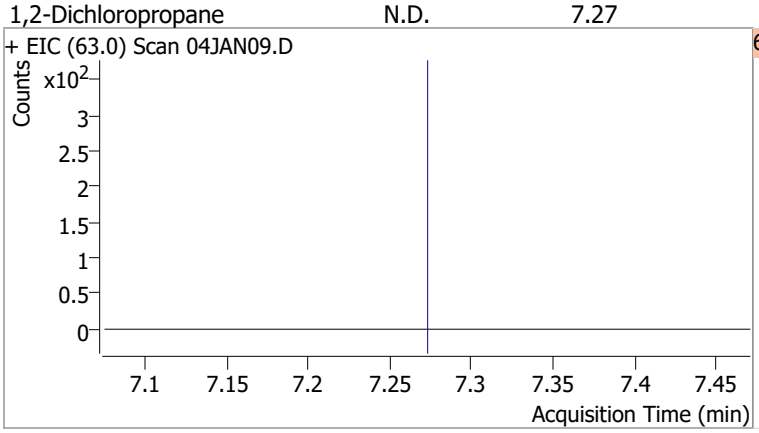


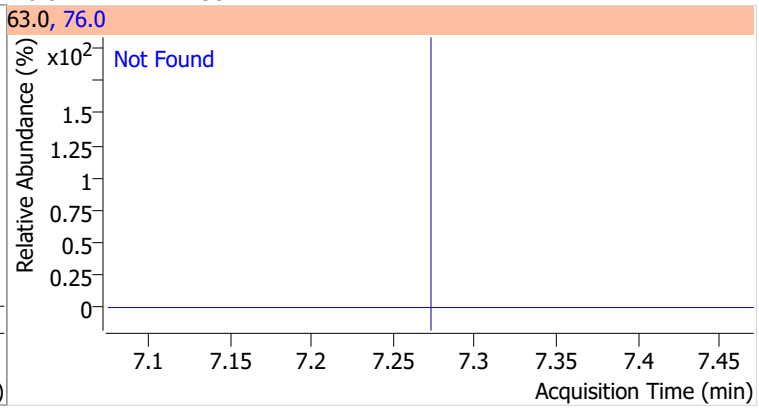
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



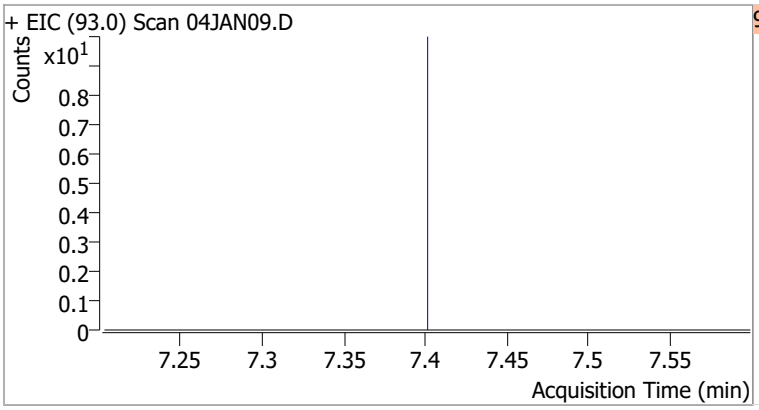


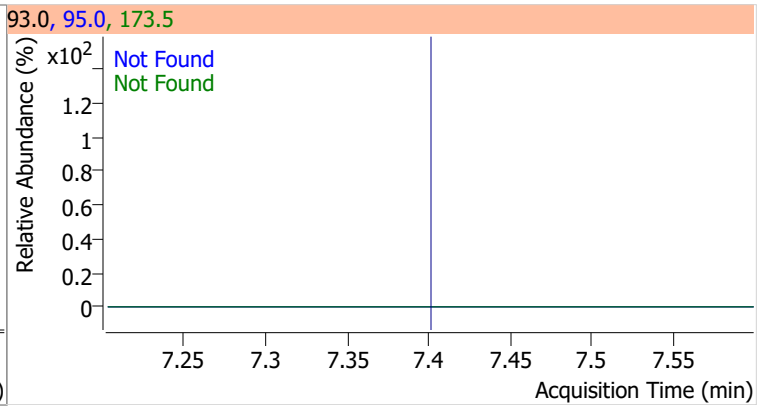
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



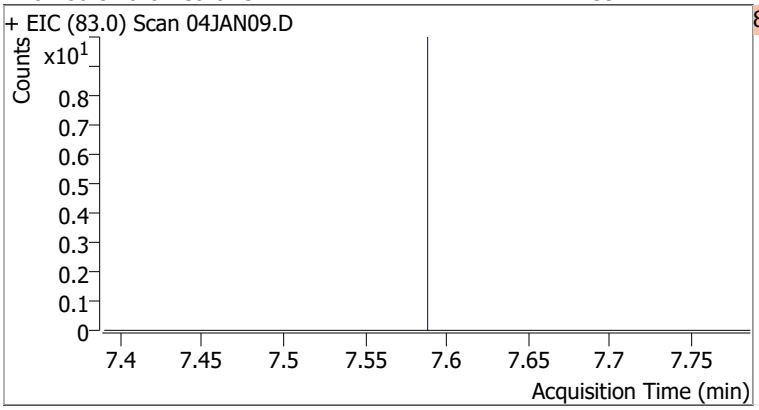


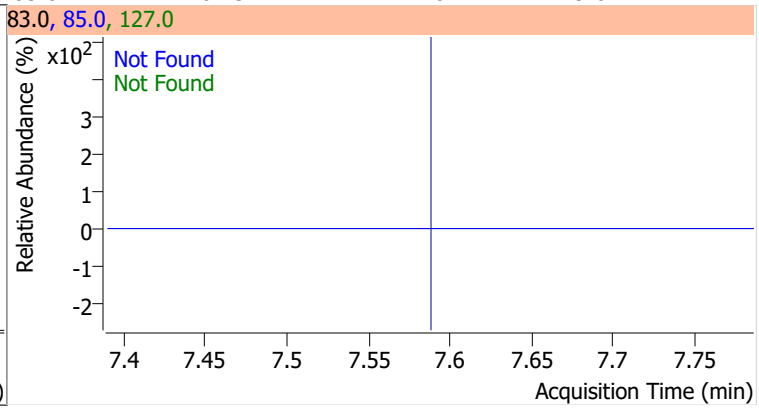
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



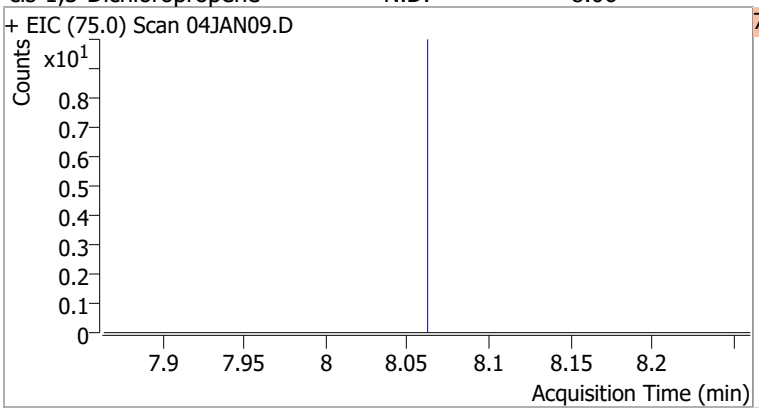


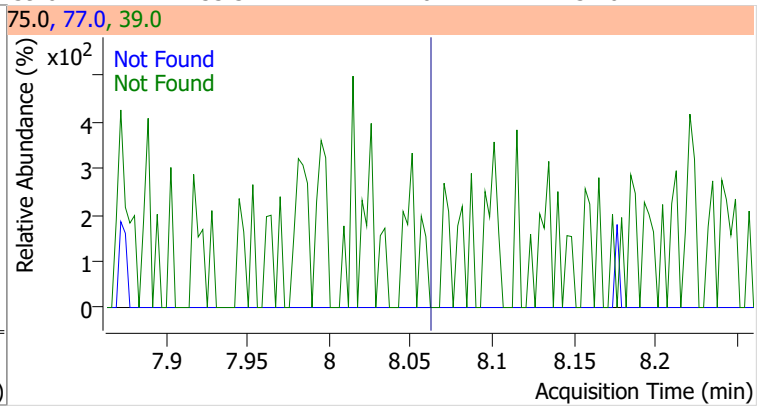
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



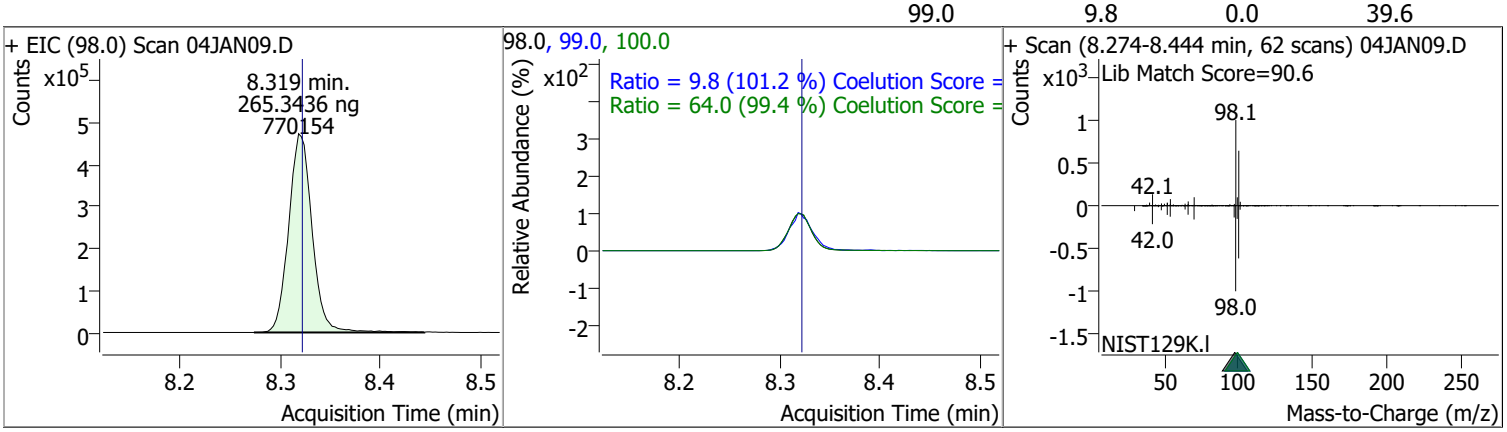


Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

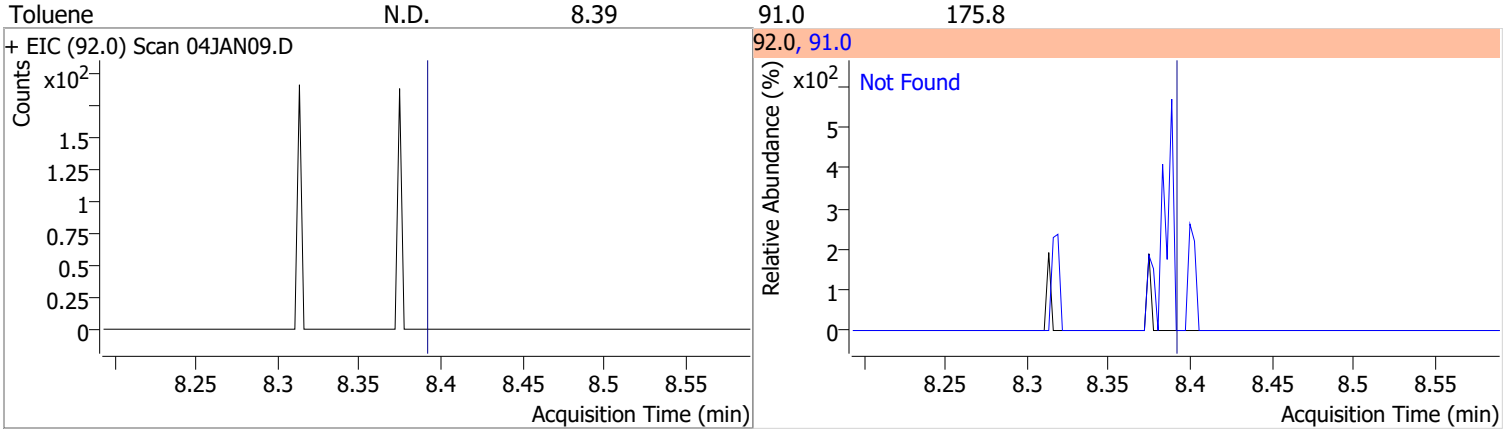




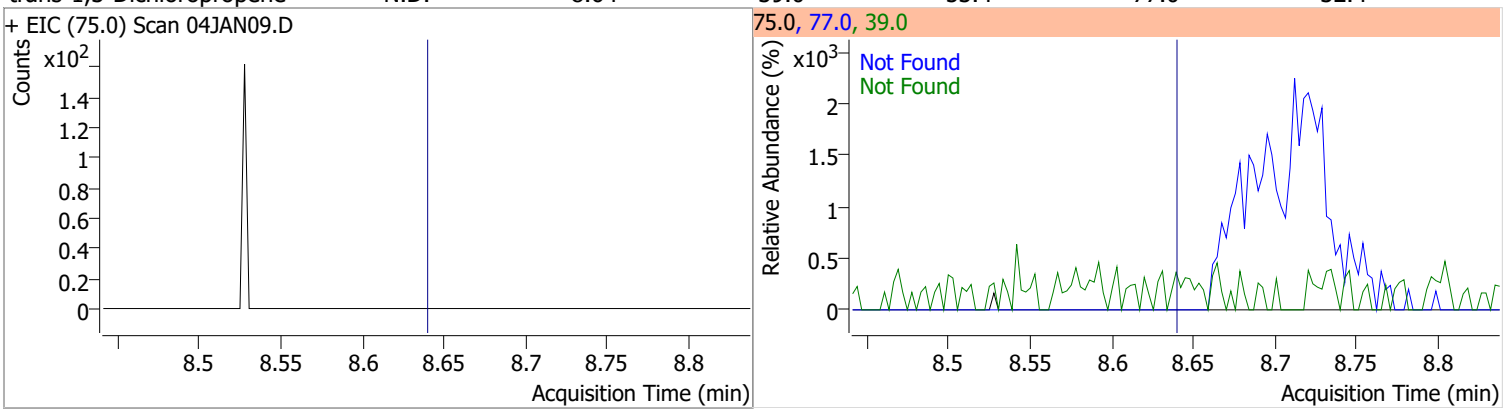
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	265.3436	8.32	0.00	770154	100.0	64.0	34.4	94.4
					99.0	9.8	0.0	39.6



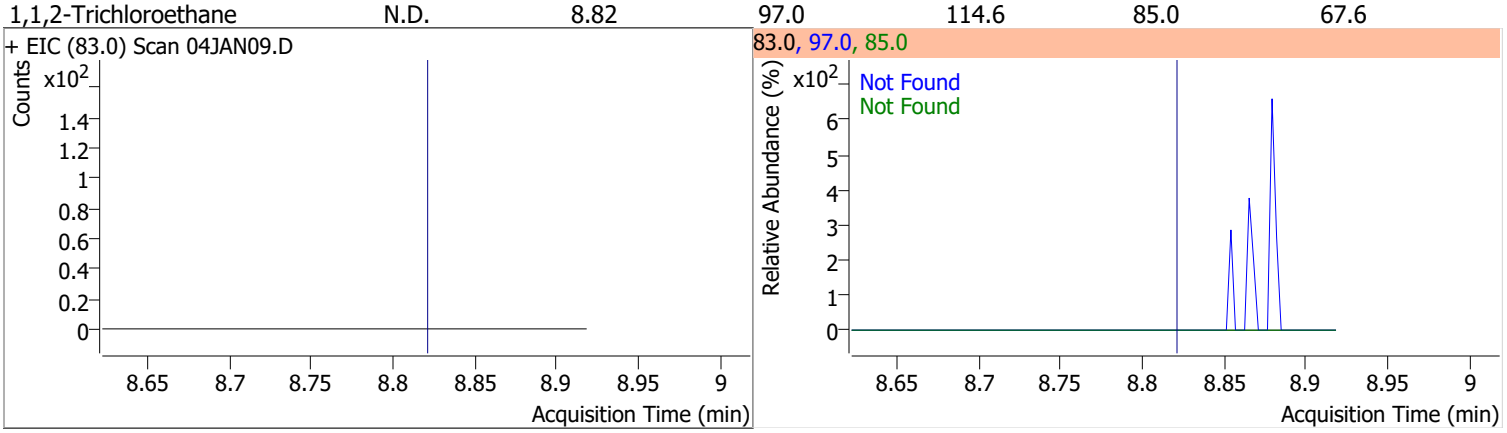
Compound	Conc.	Exp RT	QIon	Exp Ratio
Toluene	N.D.	8.39	91.0	175.8



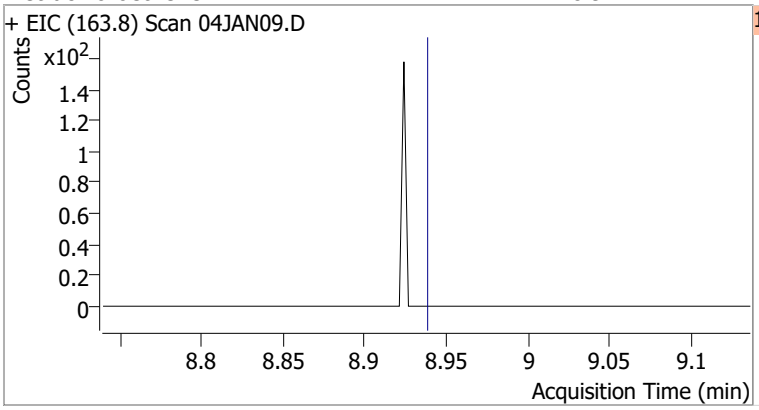
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4

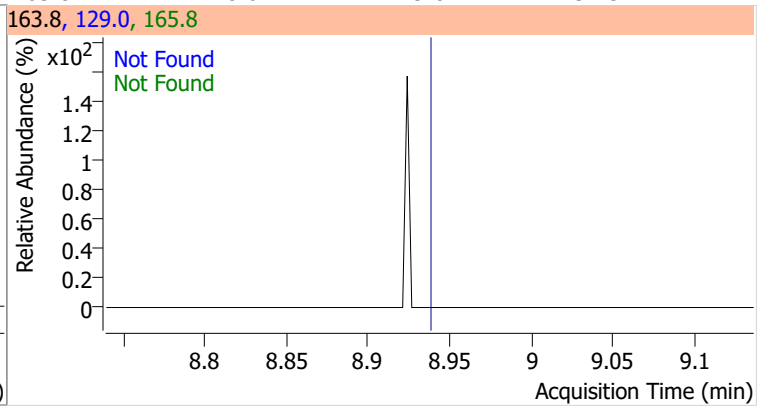


Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6

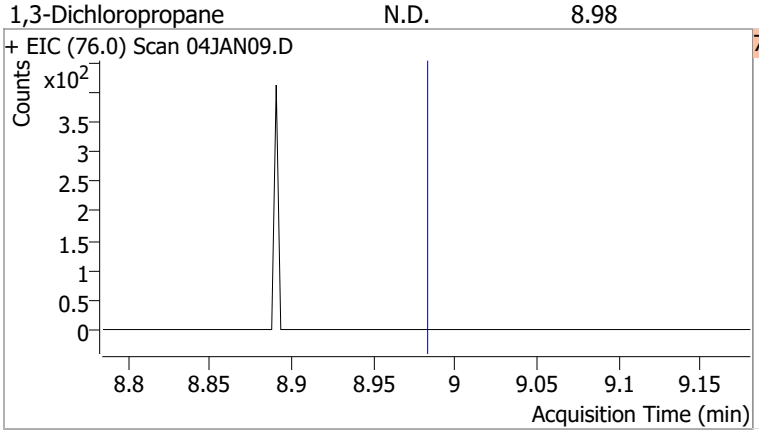


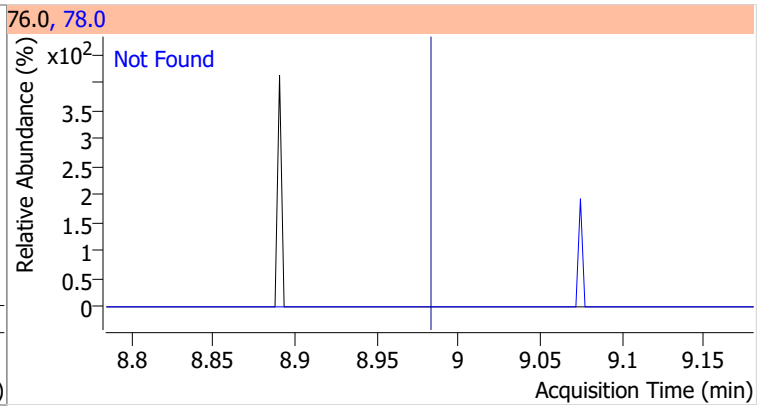
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



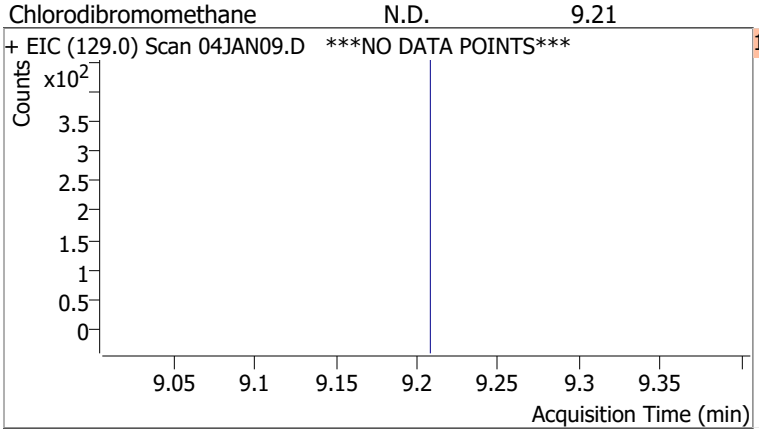


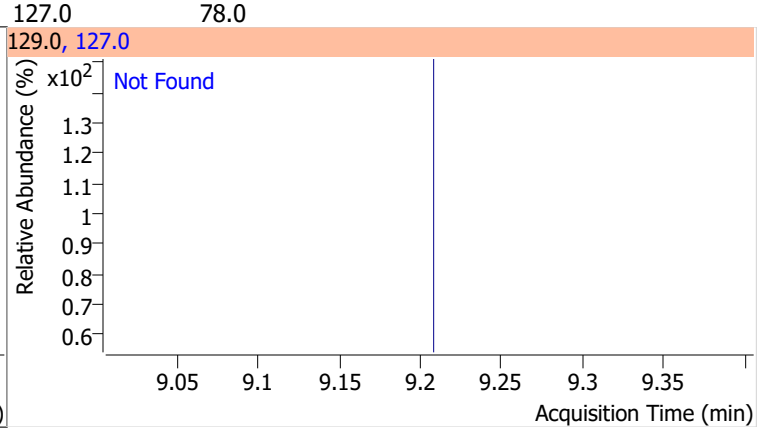
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



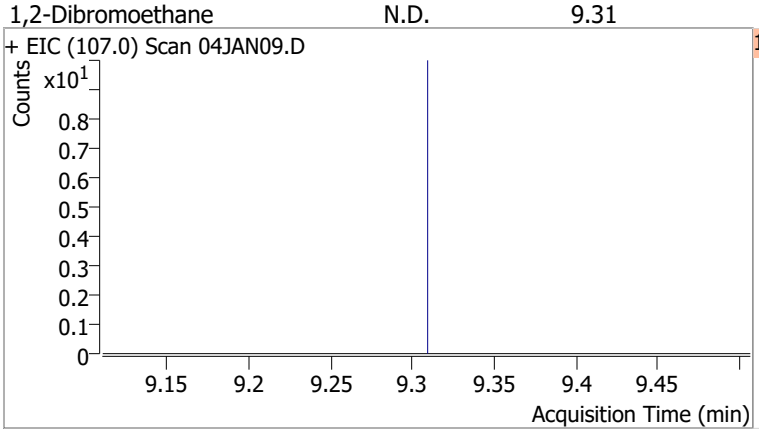


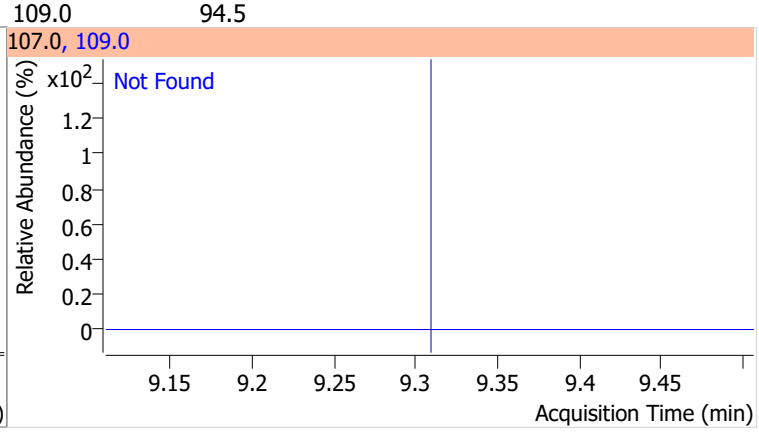
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



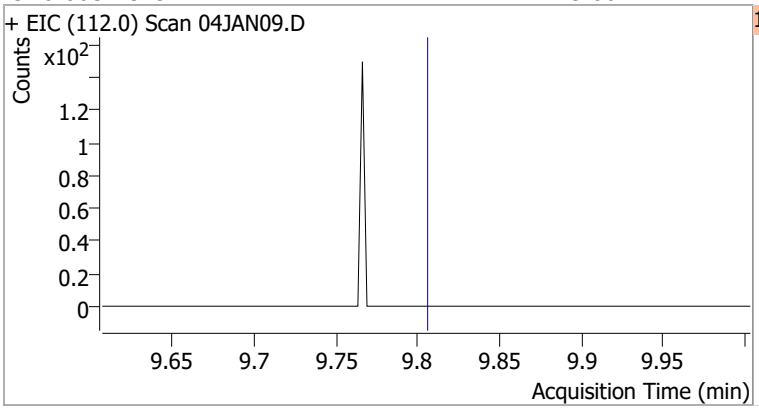


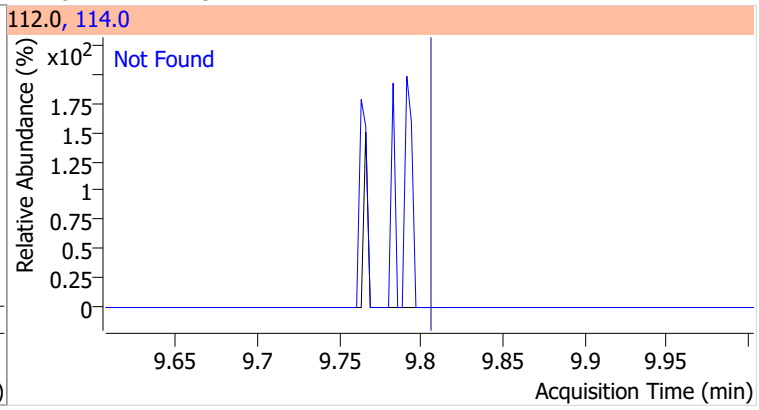
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



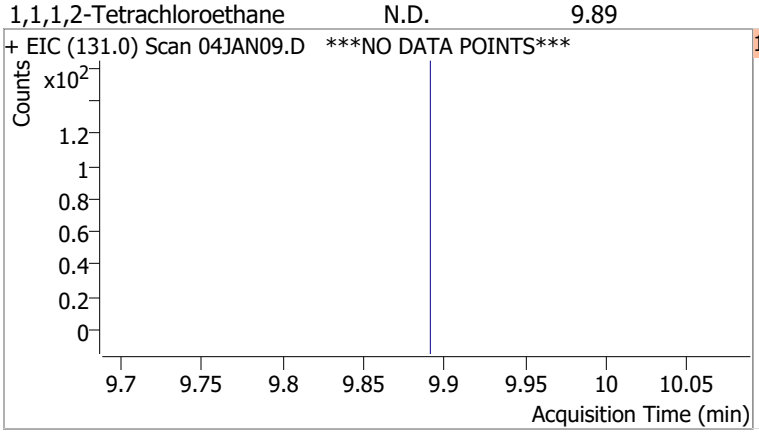


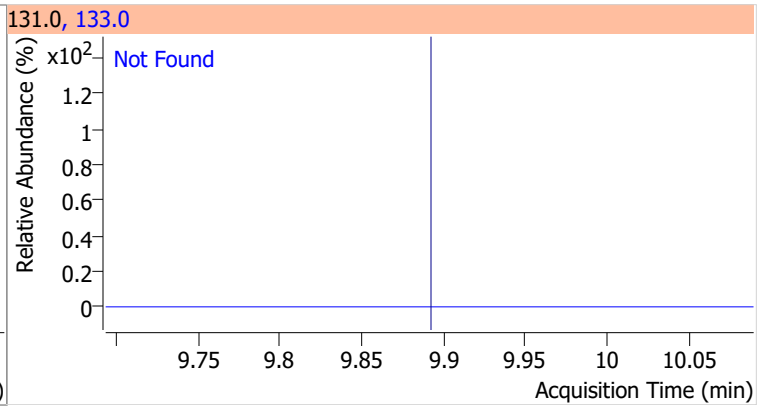
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



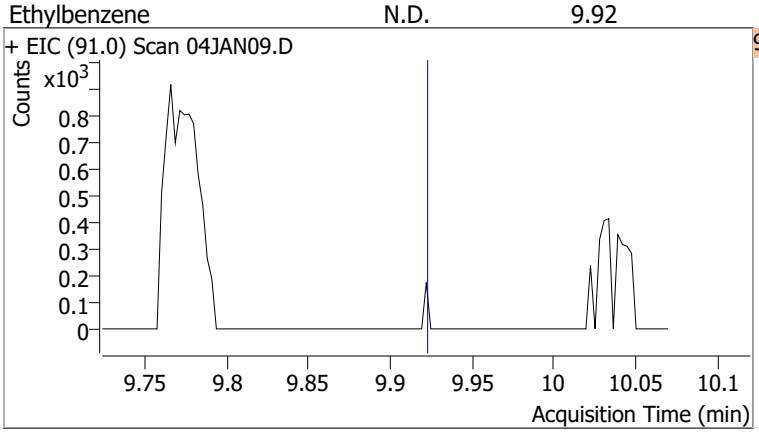


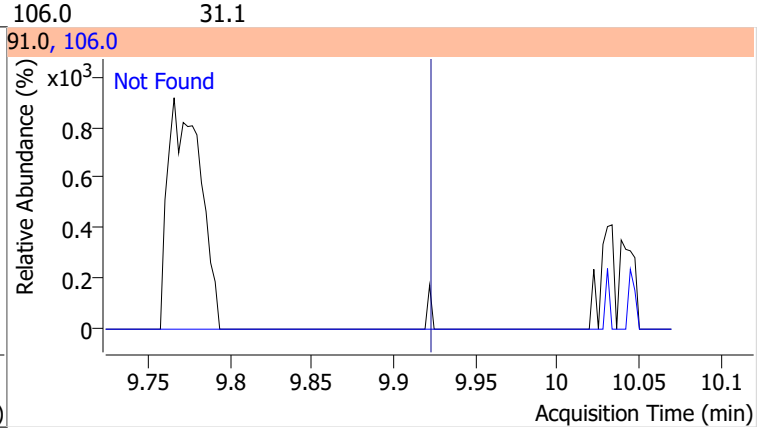
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



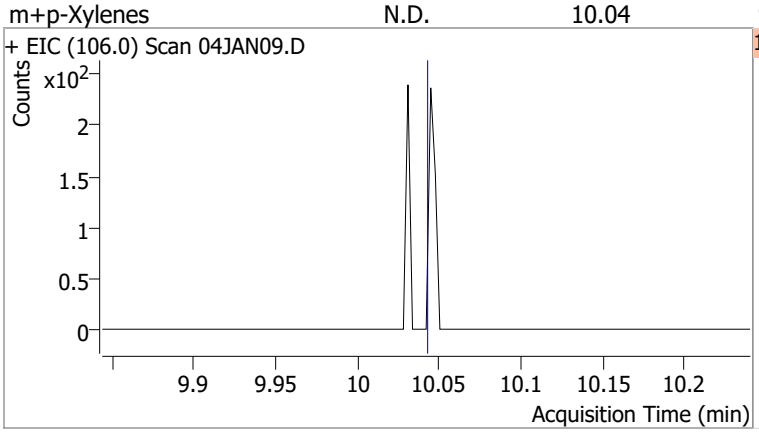


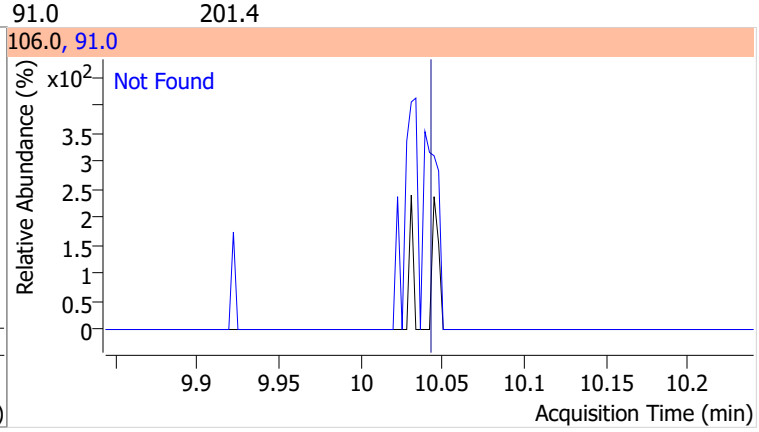
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1



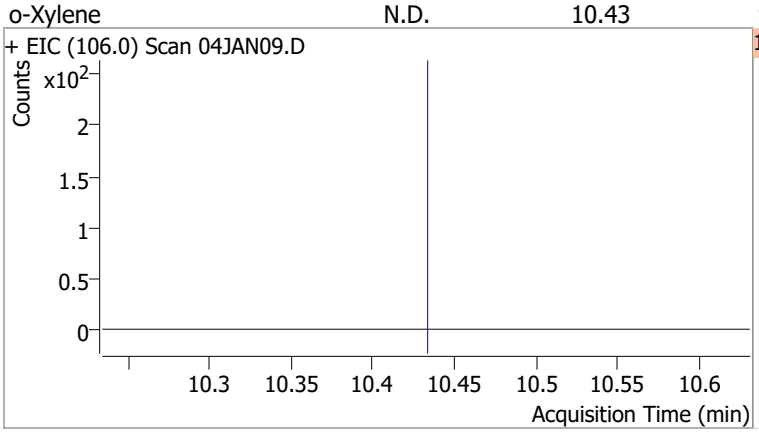


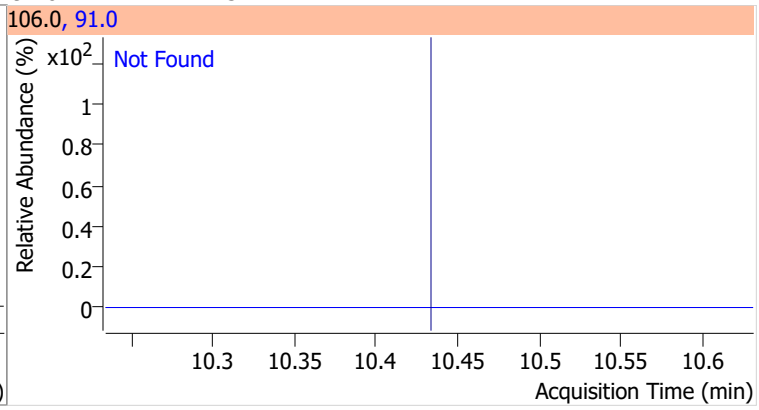
Compound	Conc.	Exp RT	QIon	Exp Ratio
m+p-Xylenes	N.D.	10.04	91.0	201.4



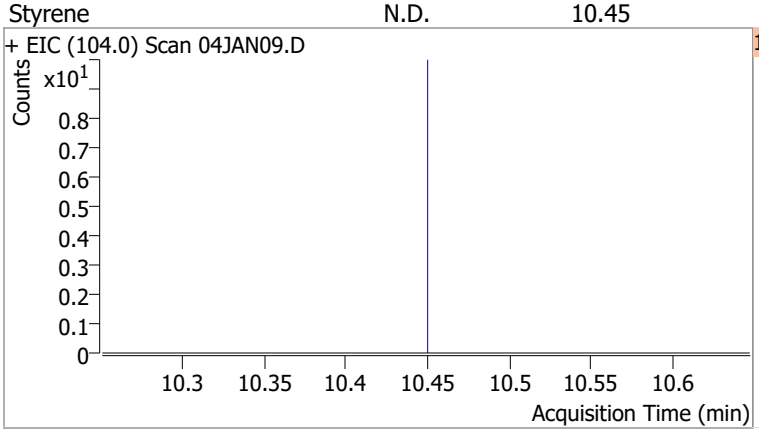


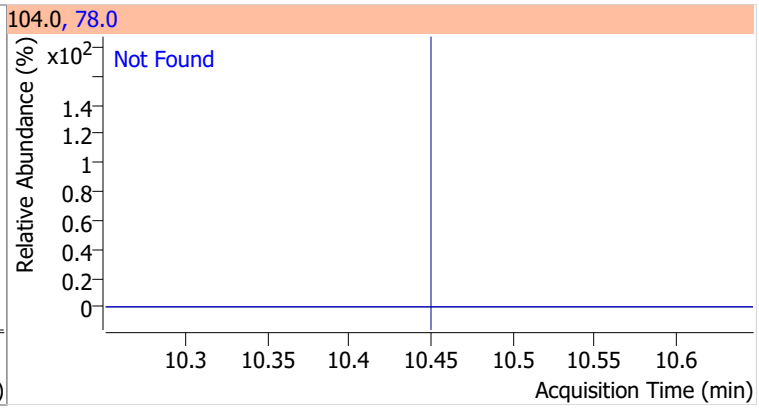
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



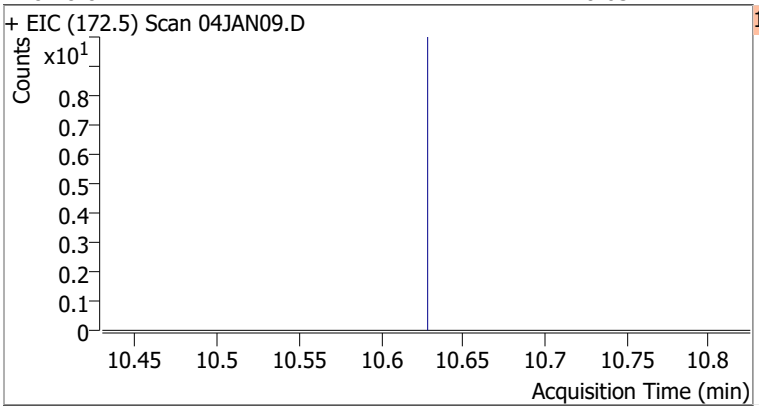


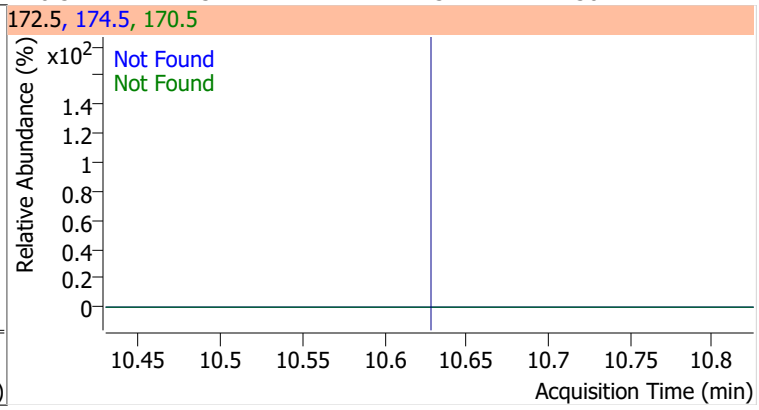
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



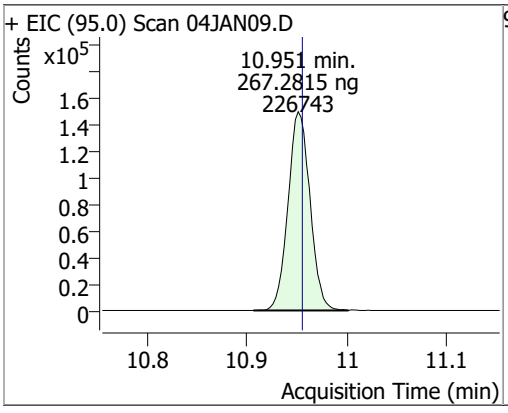


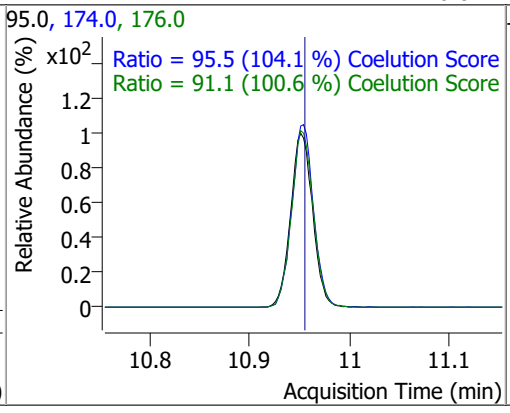
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

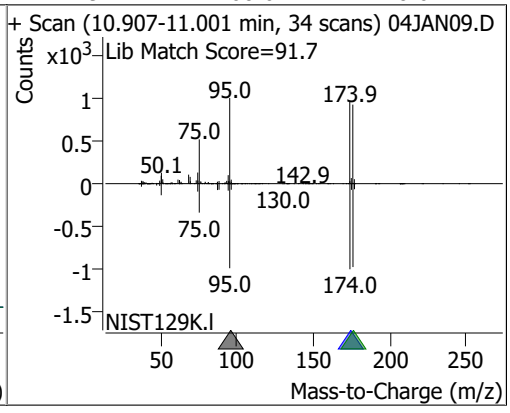




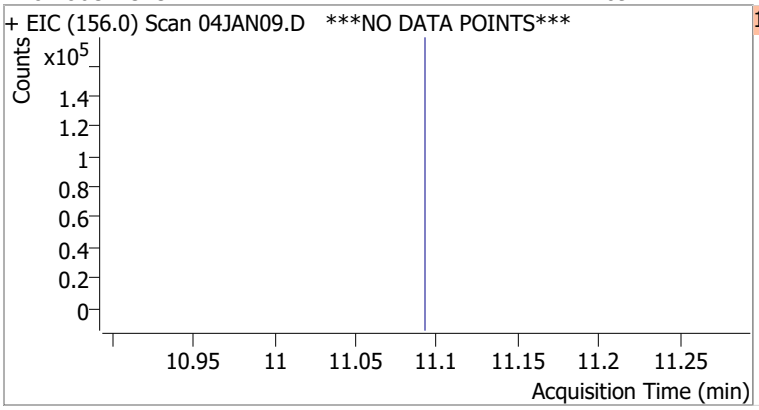
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	267.2815	10.95	0.00	226743	174.0	95.5	61.7	121.7
					176.0	91.1	60.6	120.6

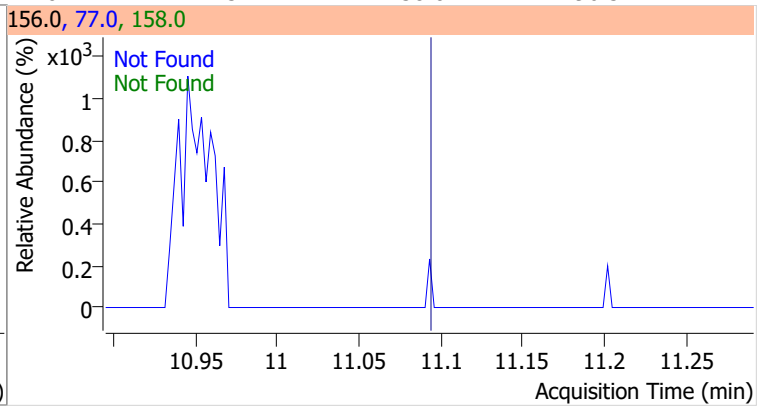




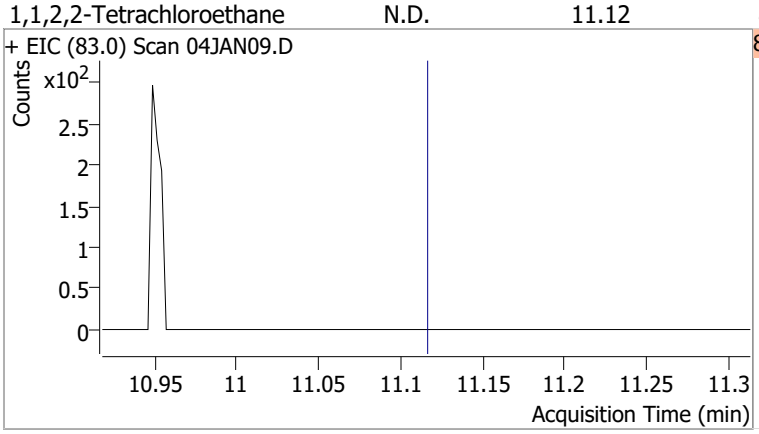


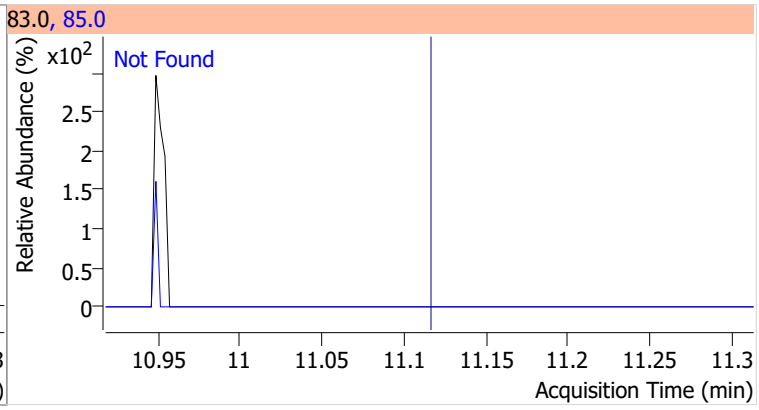
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



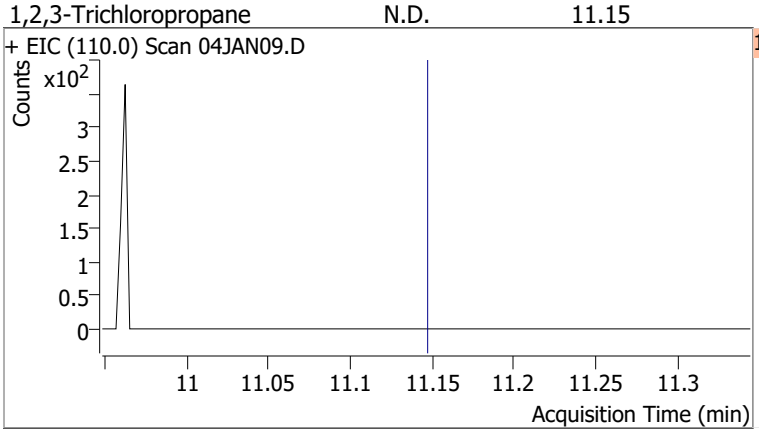


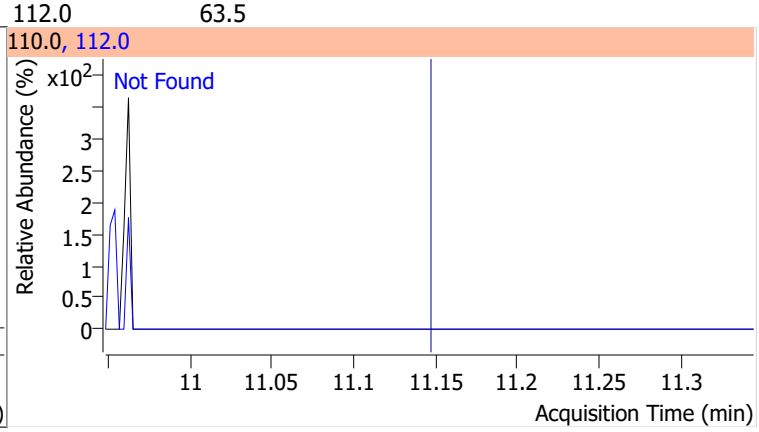
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



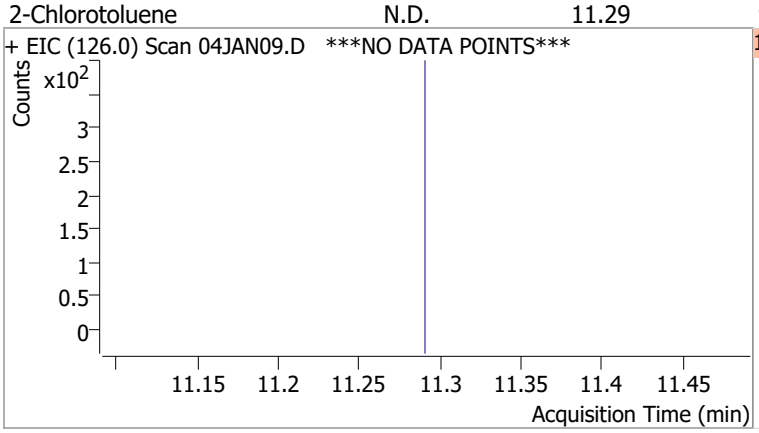


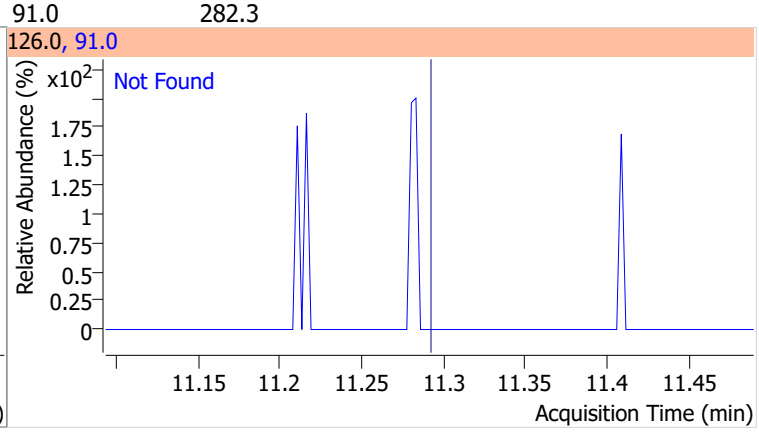
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



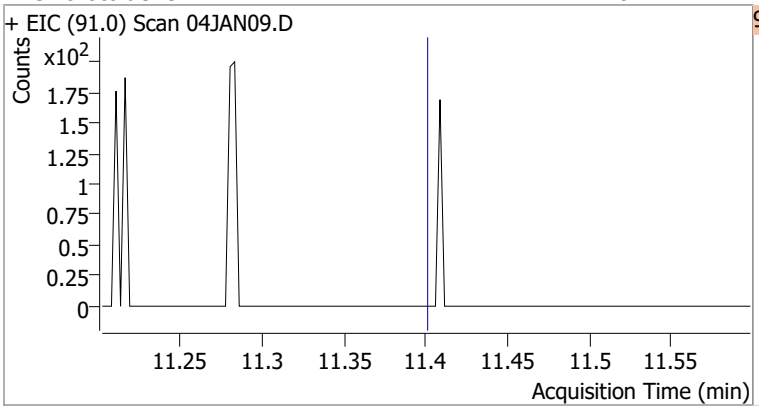


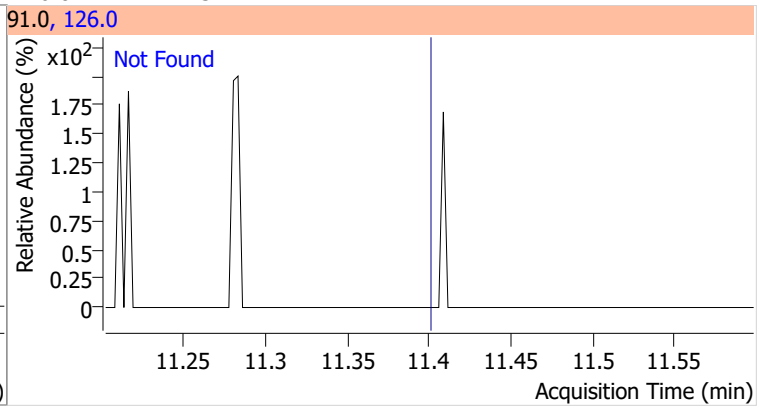
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



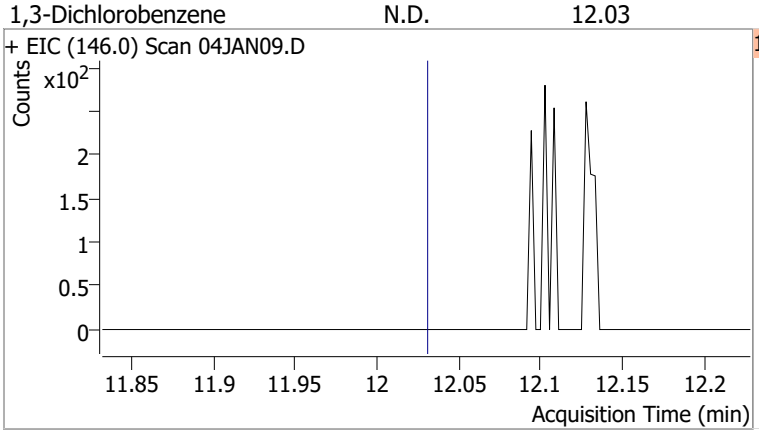


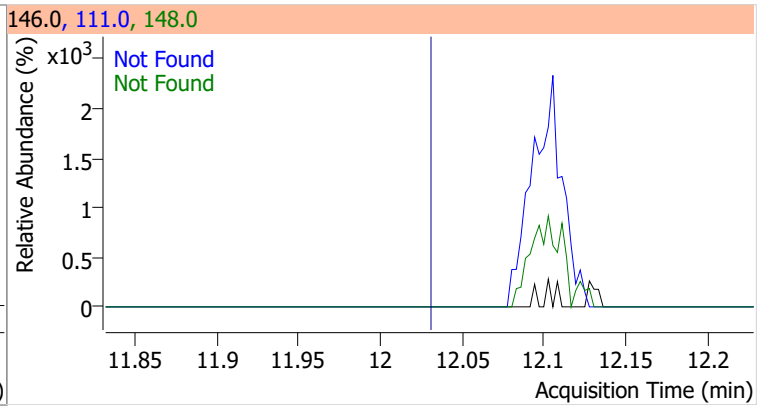
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



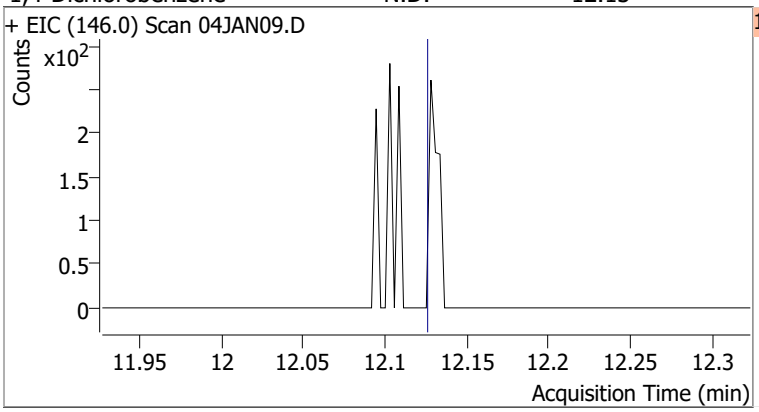


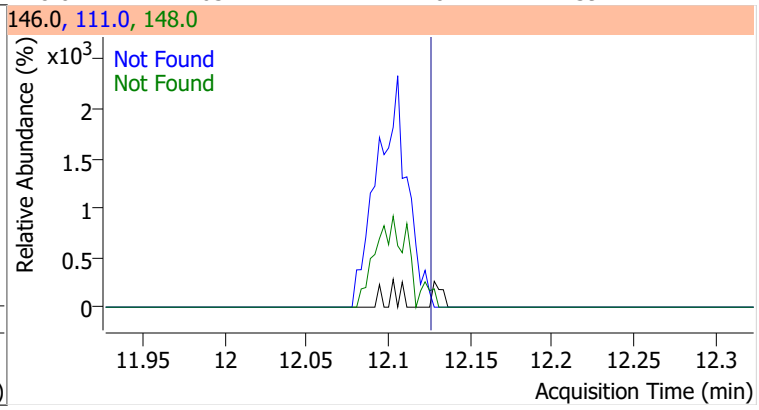
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



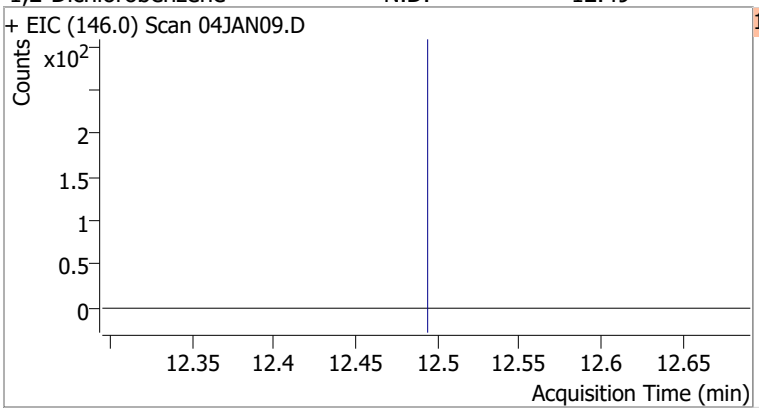


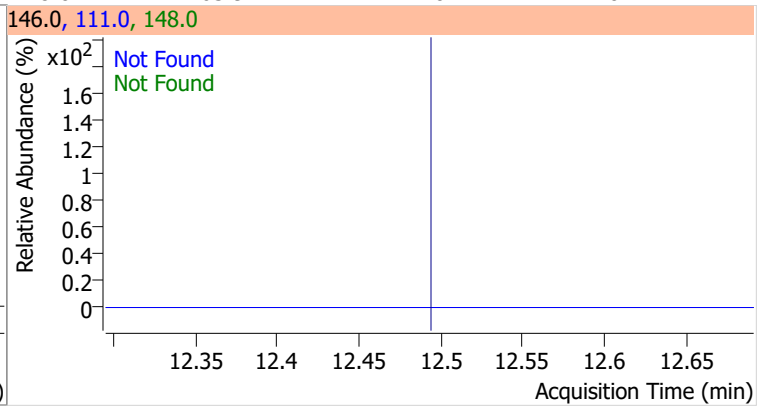
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





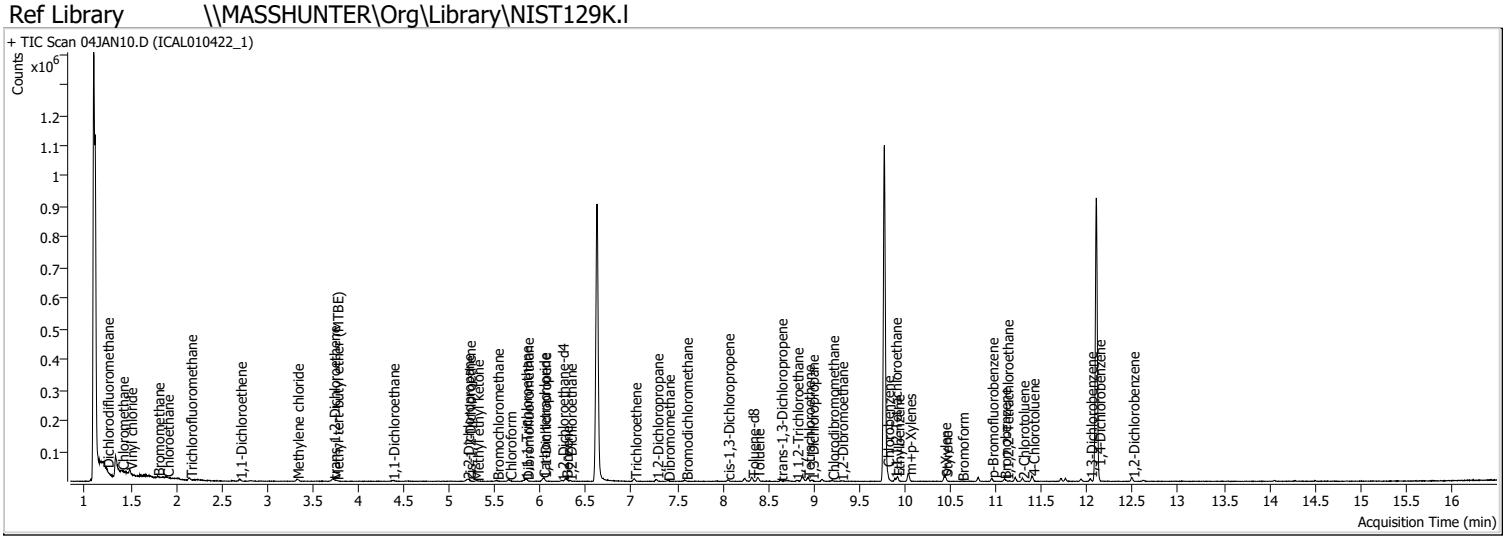
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	04JAN10.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/4/2022 3:33:04 PM
Sample Name	ICAL010422_1	Instrument	VOA5975C
Vial	10	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

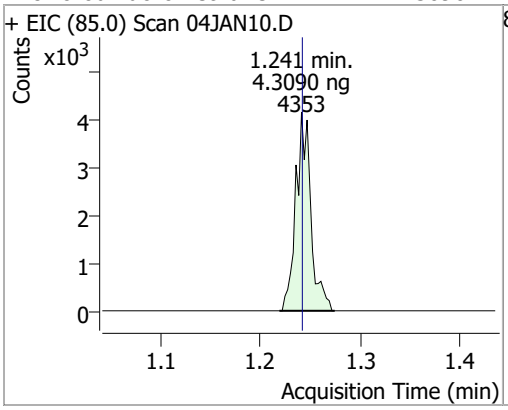


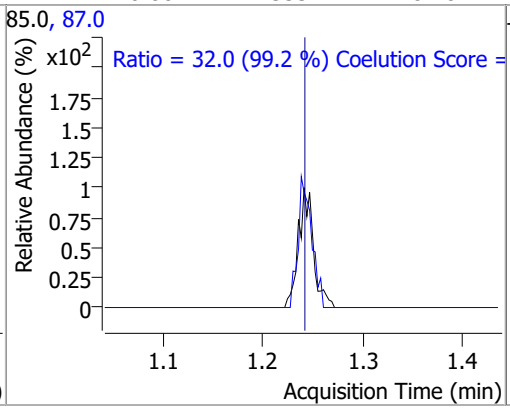
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	770895	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	296081	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	227879	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.851	113.0	2508	3.4533	ng	m	0.006
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 1.38%		*	
S 1,2-Dichloroethane-d4	6.233	67.0	923	2.9438	ng	m	0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 1.18%		*	
S Toluene-d8	8.322	98.0	7777	2.7257	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 1.09%		*	
S p-Bromofluorobenzene	10.951	95.0	2719	3.2569	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 1.30%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.241	85.0	4353	4.3090	ng		99
T Chloromethane	1.406	50.0	7435	6.0637	ng		83
T Vinyl chloride	1.495	62.0	4274	3.8739	ng		94
T Bromomethane	1.796	96.0	1902	3.8547	ng	m	88
T Chloroethane	1.899	64.0	2178	3.9871	ng	m	86
T Trichlorofluoromethane	2.153	101.0	5030	3.6731	ng		91
T 1,1-Dichloroethene	2.700	96.0	2084	2.6839	ng	m	95
T Methylene chloride	3.324	49.0	4095	3.5774	ng		88
T trans-1,2-Dichloroethene	3.723	96.0	2146	2.7090	ng	m	100
T Methyl tert-butyl ether (MTBE)	3.759	73.0	2717	2.6532	ng	m	90
T 1,1-Dichloroethane	4.376	63.0	3892	2.6393	ng		91
T 2,2-Dichloropropane	5.196	77.0	2930	2.6520	ng	m	88
T cis-1,2-Dichloroethene	5.212	96.0	2376	2.9581	ng	m	95
T Methyl ethyl ketone	5.302	43.0	3035	27.8967	ng		85
T Bromochloromethane	5.522	128.0	807	2.4260	ng	m	82
T Chloroform	5.659	83.0	4248	2.8946	ng		97

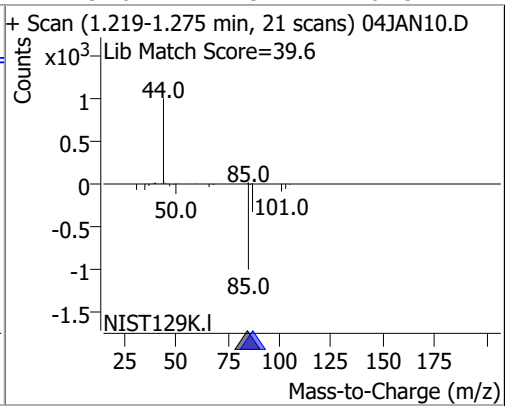
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	5.831	97.0	3510	2.5521	ng		99
T Carbon tetrachloride	6.029	117.0	4342	3.2043	ng		77
T 1,1-Dichloropropene	6.041	75.0	2830	2.4201	ng		91
T Benzene	6.278	78.0	8408	2.7393	ng		98
T 1,2-Dichloroethane	6.322	62.0	2415	2.9090	ng	m	96
T Trichloroethene	7.033	95.0	2372	2.6564	ng	m	93
T 1,2-Dichloropropane	7.273	63.0	2148	2.7347	ng		93
T Dibromomethane	7.396	93.0	902	2.7162	ng	m	88
T Bromodichloromethane	7.597	83.0	2536	2.7684	ng		98
T cis-1,3-Dichloropropene	8.054	75.0	2583	2.4939	ng		94
T Toluene	8.380	92.0	5039	2.6145	ng		93
T trans-1,3-Dichloropropene	8.634	75.0	1470	1.9942	ng	m	83
T 1,1,2-Trichloroethane	8.810	83.0	960	2.5012	ng	m	89
T Tetrachloroethene	8.932	163.8	2105	2.6772	ng	m	95
T 1,3-Dichloropropane	8.977	76.0	2257	2.9881	ng		77
T Chlorodibromomethane	9.203	129.0	1468	2.4461	ng	m	100
T 1,2-Dibromoethane	9.300	107.0	1299	3.0943	ng	m	85
T Chlorobenzene	9.805	112.0	5771	2.7350	ng		100
T 1,1,1,2-Tetrachloroethane	9.889	131.0	1893	2.5659	ng	m	98
T Ethylbenzene	9.920	91.0	9283	2.5367	ng		93
T m+p-Xylenes	10.045	106.0	7212	5.0712	ng		88
T o-Xylene	10.430	106.0	3330	2.6303	ng	#	80
T Styrene	10.444	104.0	4408	2.1625	ng		98
T Bromoform	10.625	172.5	708	2.4287	ng	m	87
T Bromobenzene	11.088	156.0	2024	2.7439	ng	m	94
T 1,1,2,2-Tetrachloroethane	11.113	83.0	1142	2.6916	ng	m	92
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	11.292	126.0	1844	2.5124	ng	m	97
T 4-Chlorotoluene	11.400	91.0	5419	2.2650	ng		96
T 1,3-Dichlorobenzene	12.033	146.0	3541	2.6327	ng		99
T 1,4-Dichlorobenzene	12.125	146.0	3787	2.7613	ng		90
T 1,2-Dichlorobenzene	12.499	146.0	3104	2.7307	ng		96

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

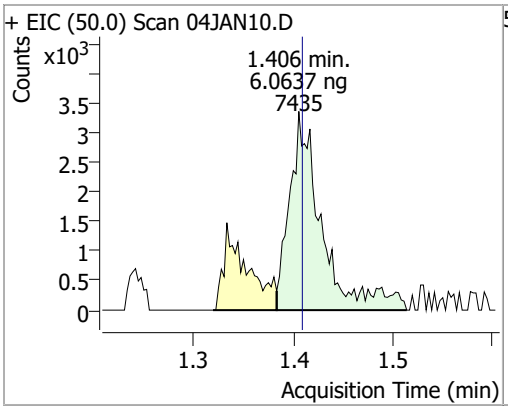
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	4.3090	1.24	0.00	4353	87.0	32.0	2.3	62.3

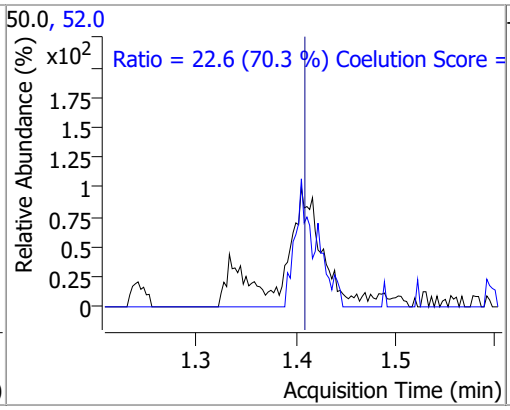


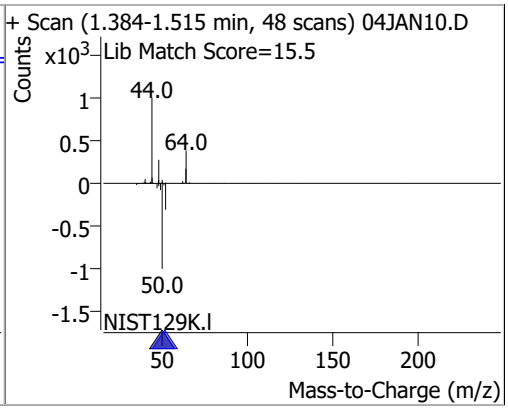




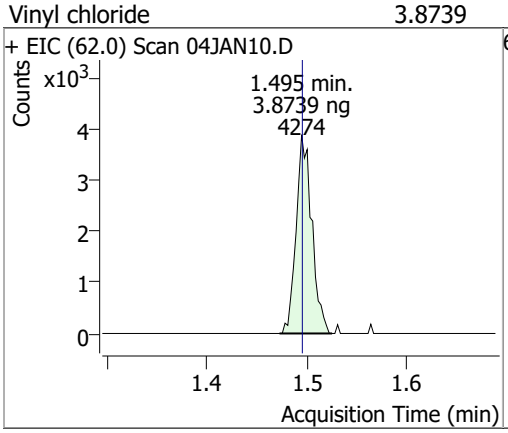
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	6.0637	1.41	0.00	7435	52.0	22.6	2.1	62.1

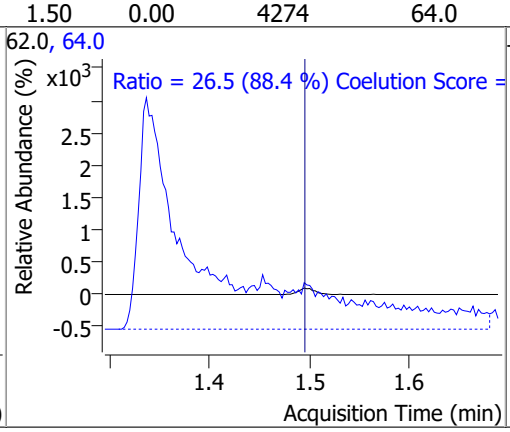


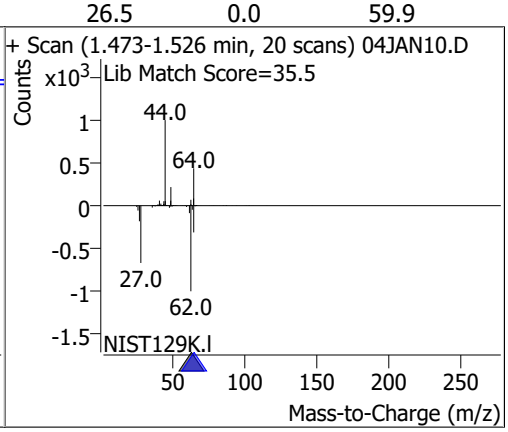




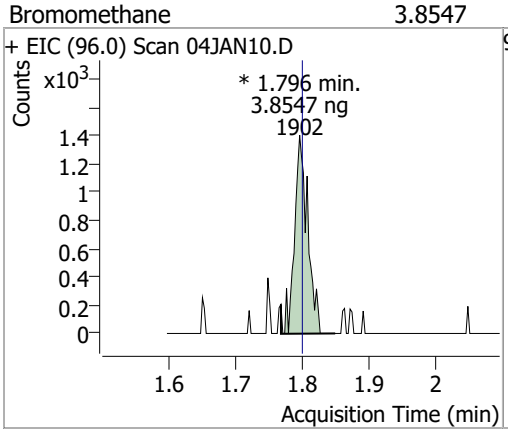
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	3.8739	1.50	0.00	4274	64.0	26.5	0.0	59.9

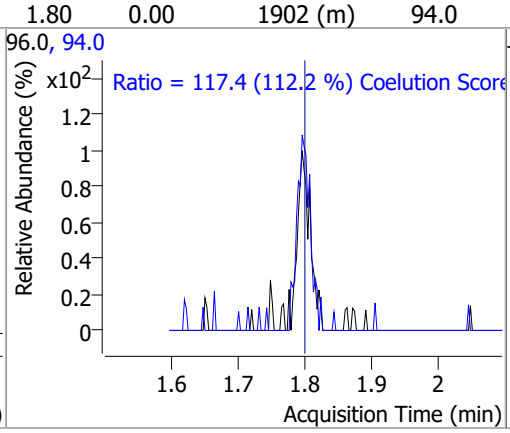


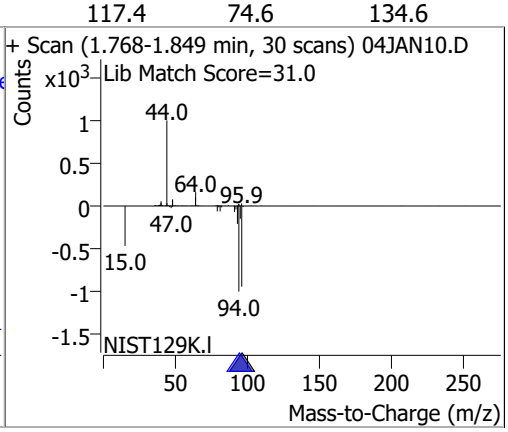




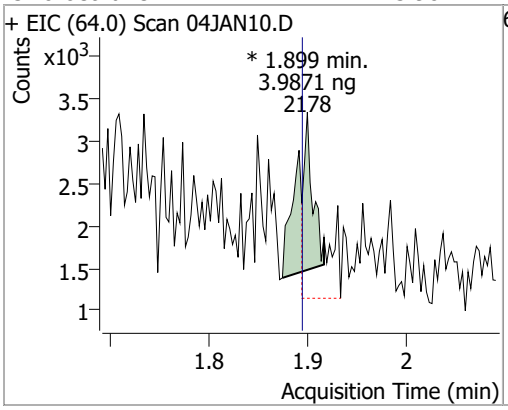
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	3.8547	1.80	0.00	1902 (m)	94.0	117.4	74.6	134.6

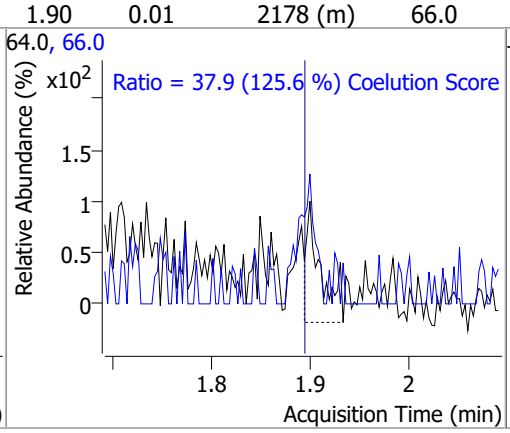


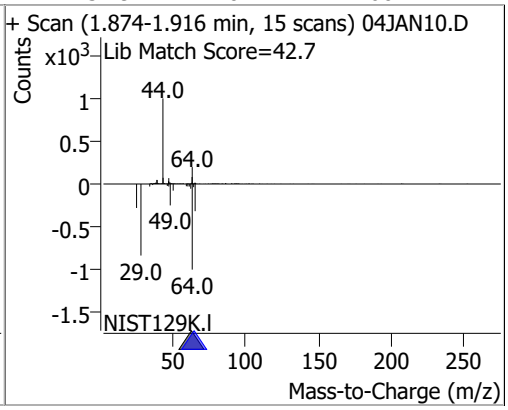




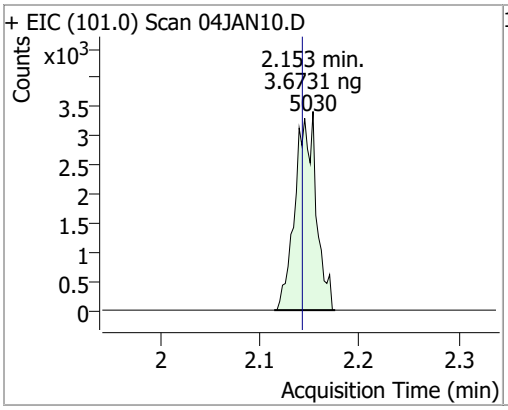
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	3.9871	1.90	0.01	2178 (m)	66.0	37.9	0.1	60.1

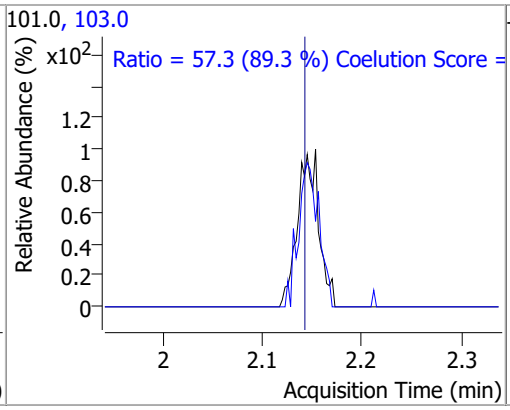


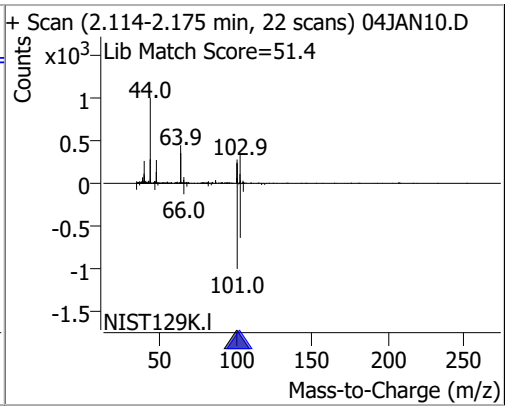




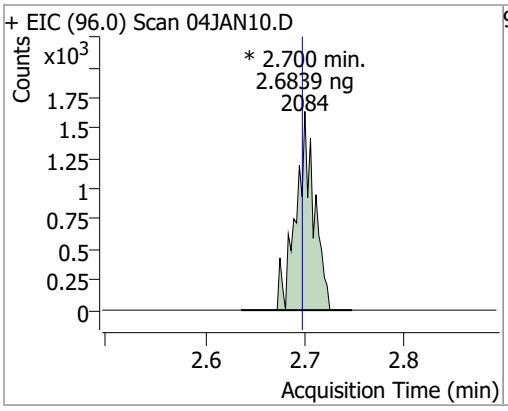
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	3.6731	2.15	0.01	5030	103.0	57.3	34.2	94.2

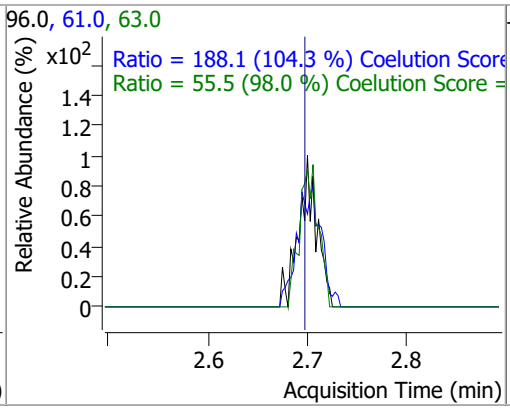


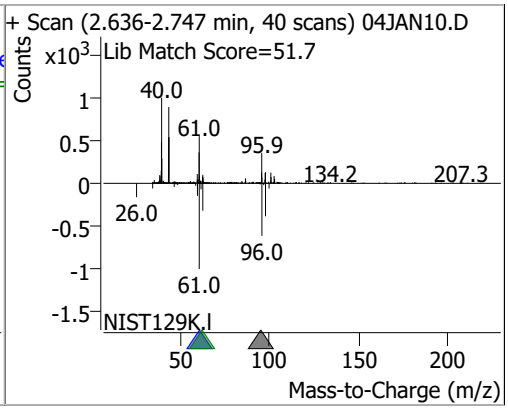




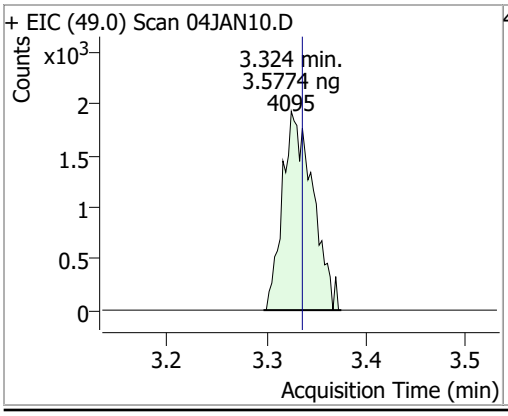
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	2.6839	2.70	0.00	2084 (m)	61.0	188.1	150.3	210.3
					63.0	55.5	26.7	86.7

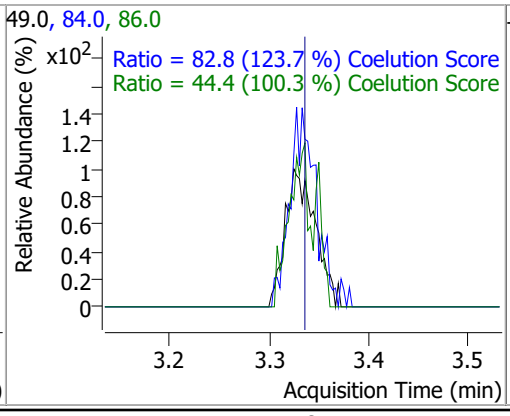


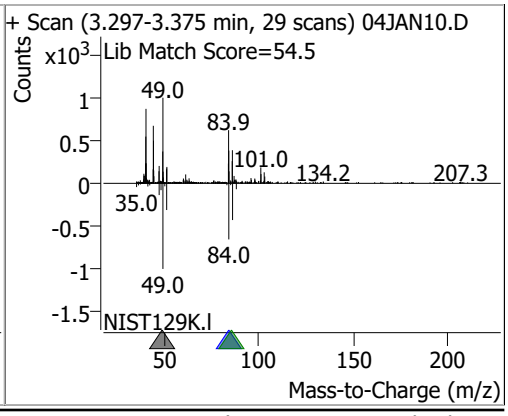




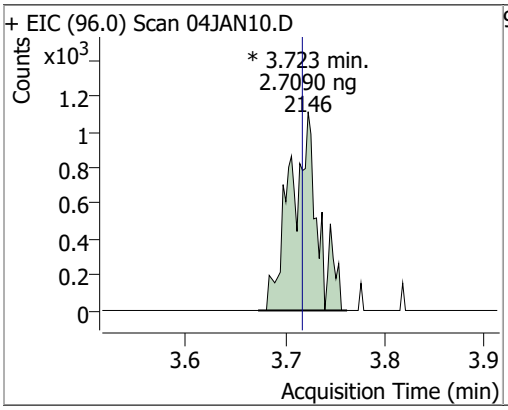
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	3.5774	3.32	-0.01	4095	84.0	82.8	36.9	96.9
					86.0	44.4	14.3	74.3

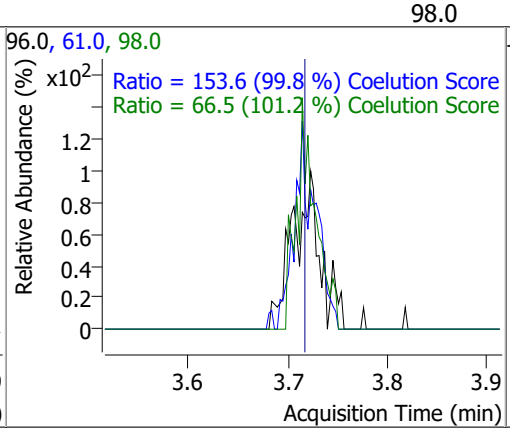


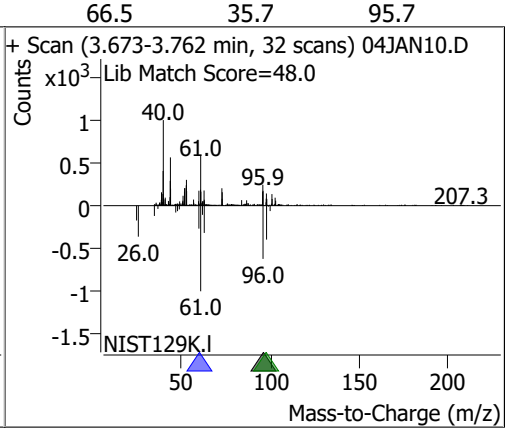




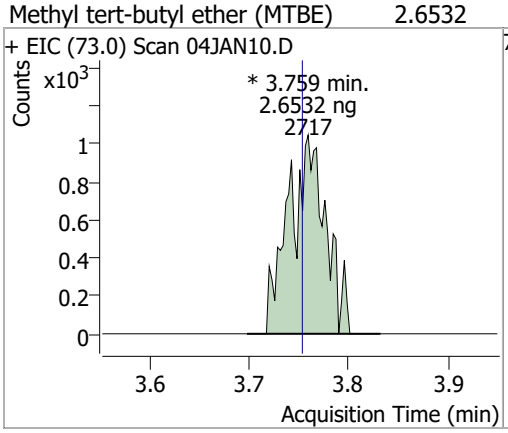
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	2.7090	3.72	0.01	2146 (m)	61.0	153.6	123.9	183.9
					98.0	66.5	35.7	95.7

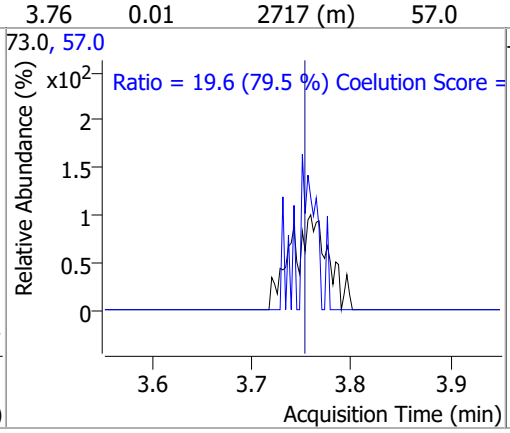


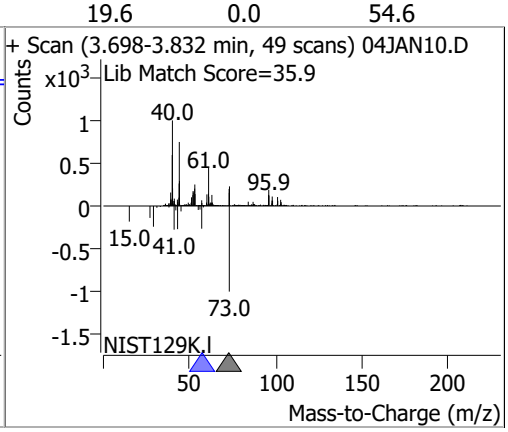




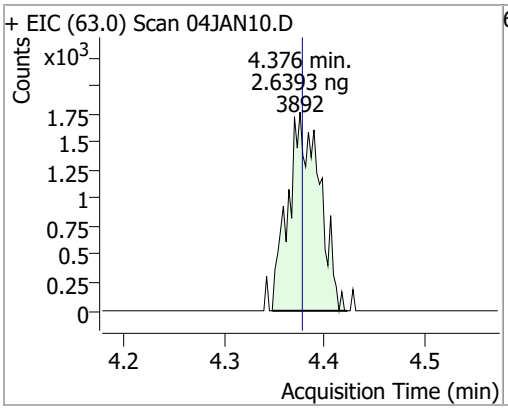
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	2.6532	3.76	0.01	2717 (m)	57.0	19.6	0.0	54.6

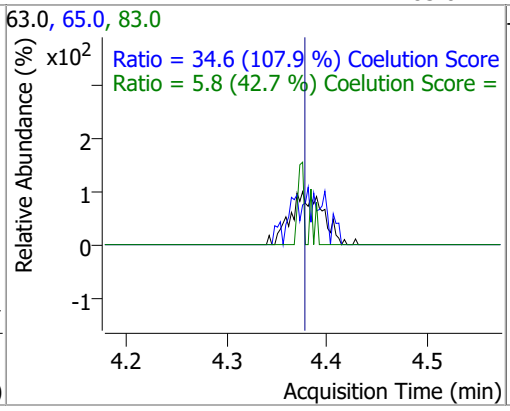


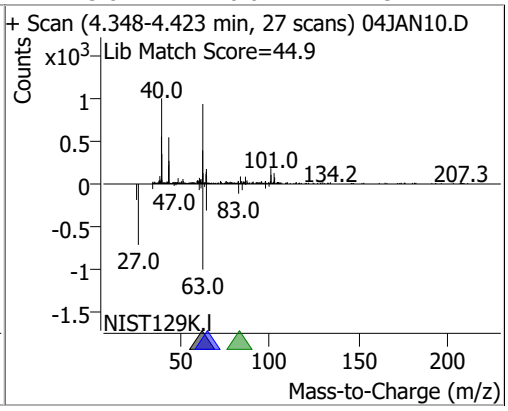




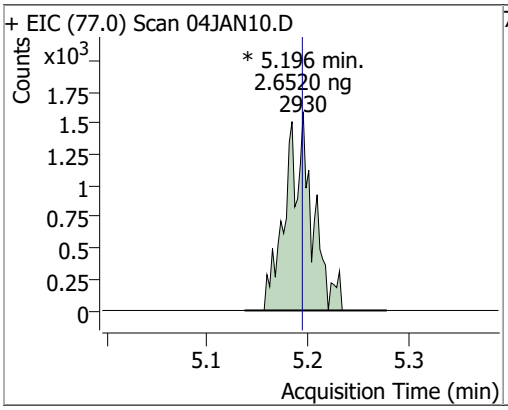
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	2.6393	4.38	0.00	3892	65.0	34.6	2.1	62.1
					83.0	5.8	0.0	43.7

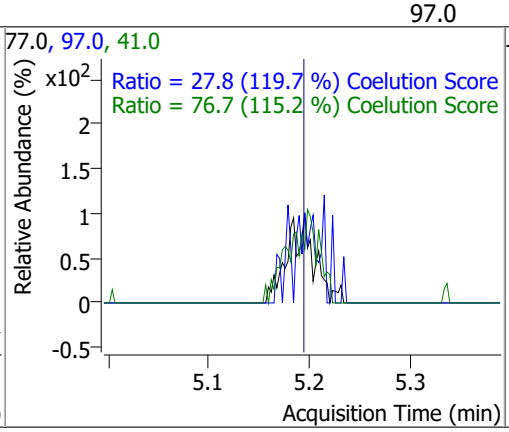


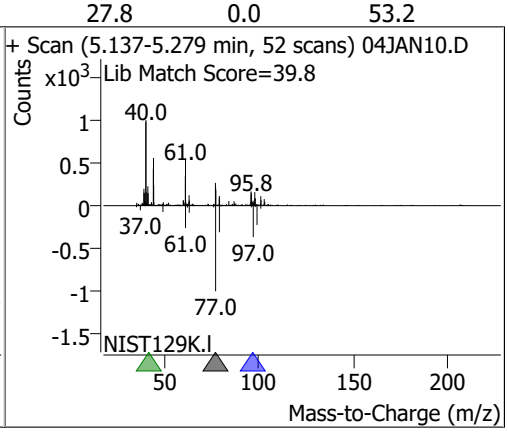




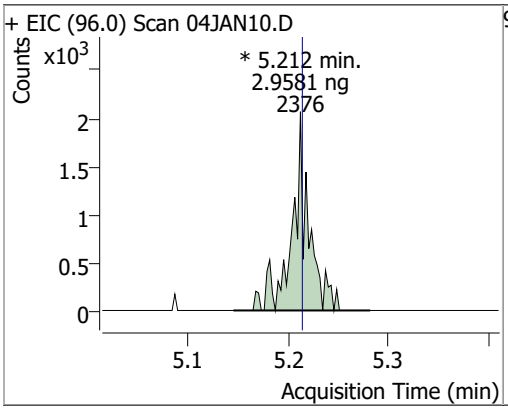
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	2.6520	5.20	0.00	2930 (m)	41.0	76.7	36.5	96.5
					97.0	27.8	0.0	53.2

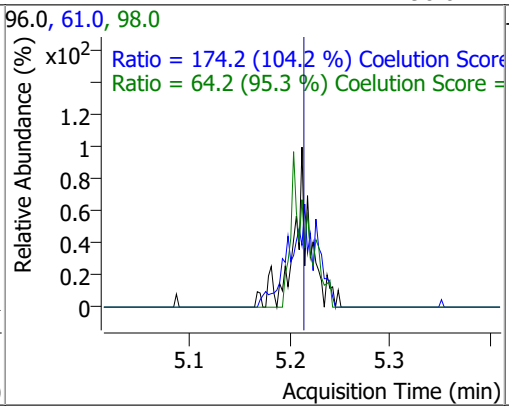


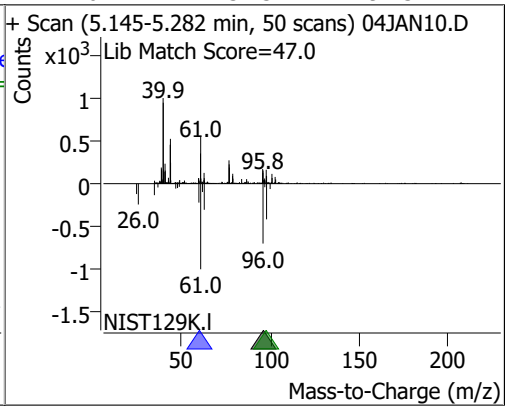




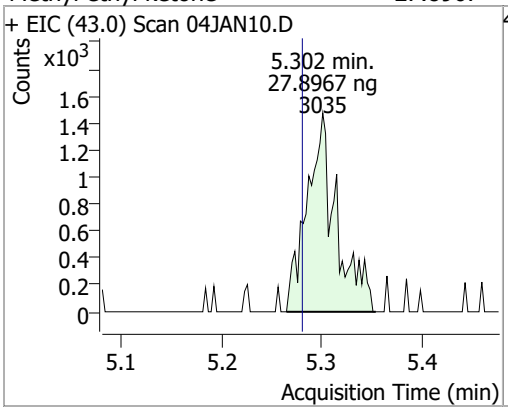
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	2.9581	5.21	0.00	2376 (m)	61.0	174.2	137.2	197.2
					98.0	64.2	37.3	97.3

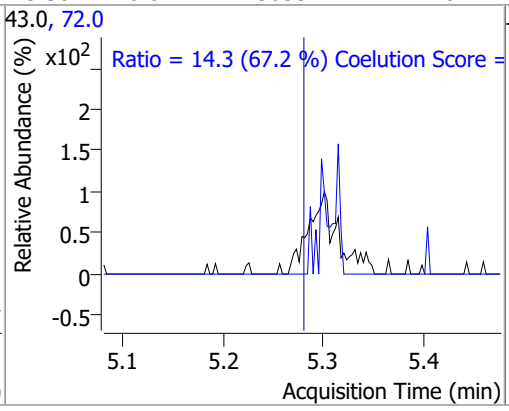


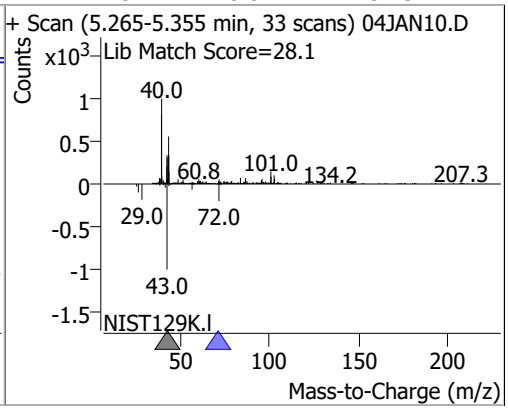




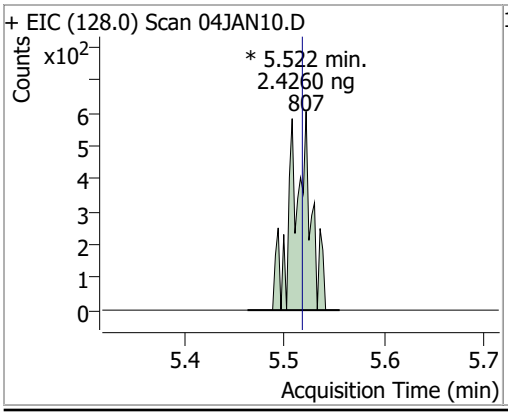
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	27.8967	5.30	0.02	3035	72.0	14.3	0.0	51.3

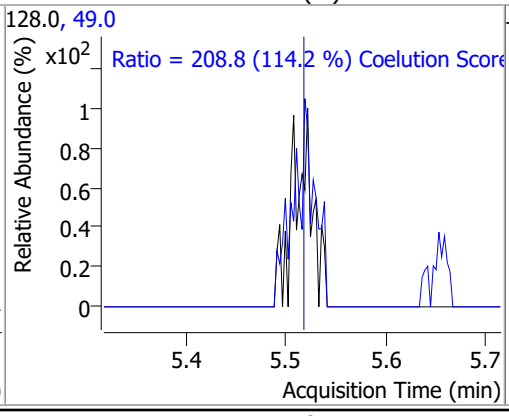


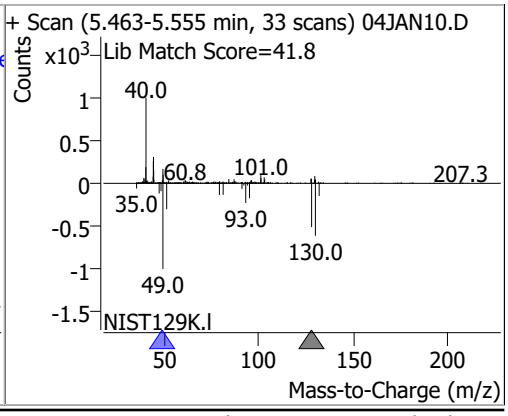




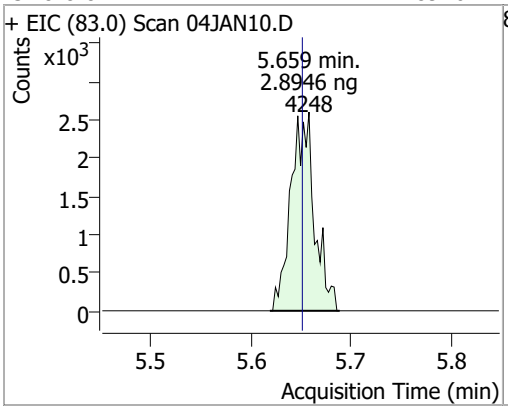
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	2.4260	5.52	0.00	807 (m)	49.0	208.8	152.9	212.9

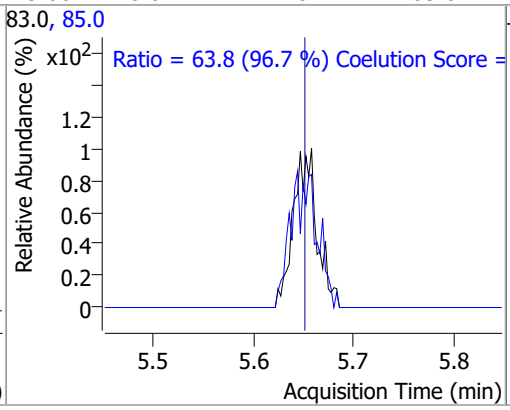


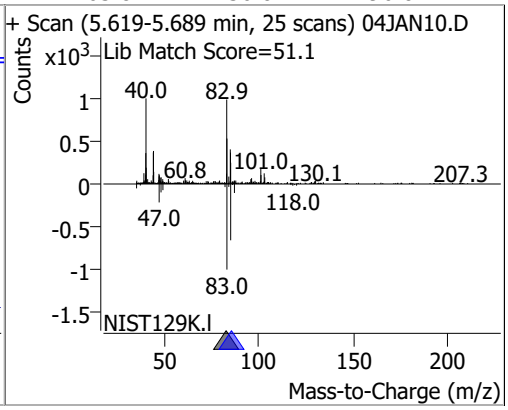




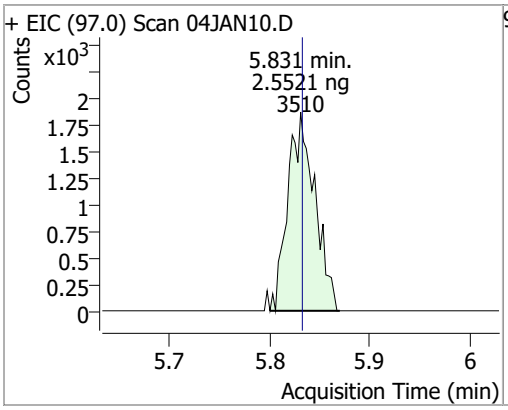
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	2.8946	5.66	0.01	4248	85.0	63.8	36.0	96.0

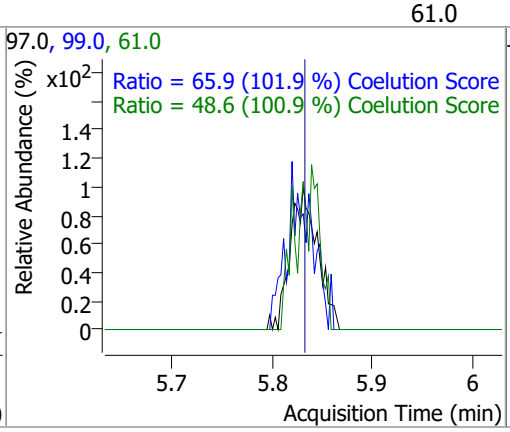


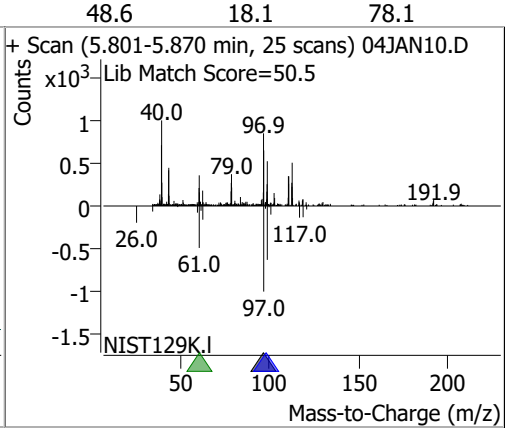




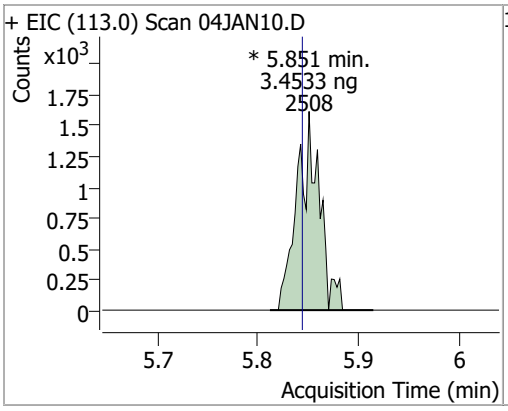
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	2.5521	5.83	0.00	3510	99.0	65.9	34.7	94.7
					61.0	48.6	18.1	78.1

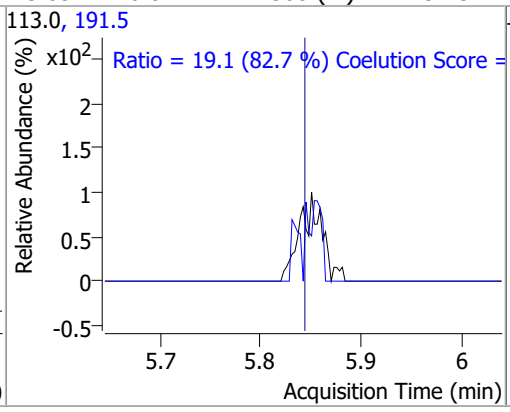


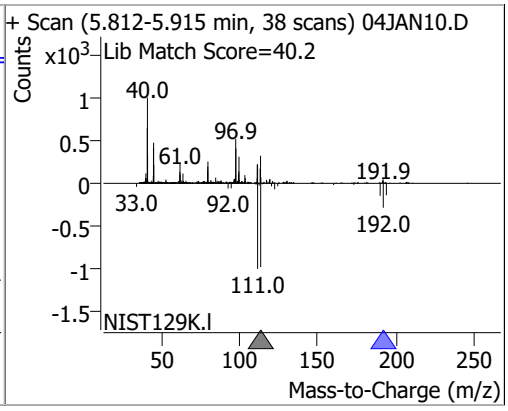




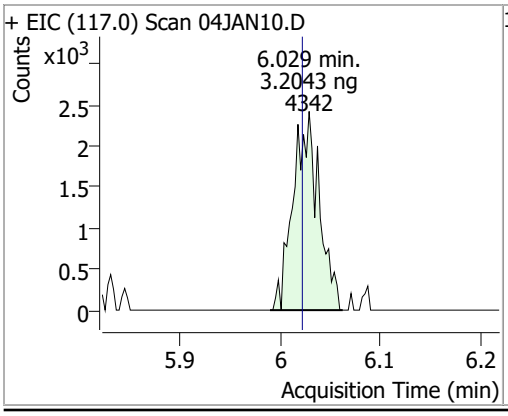
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	3.4533	5.85	0.01	2508 (m)	191.5	19.1	0.0	53.1

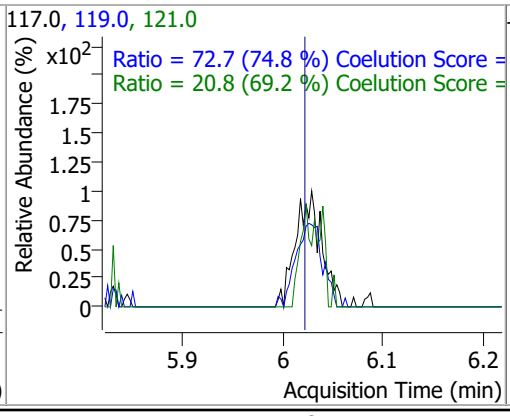


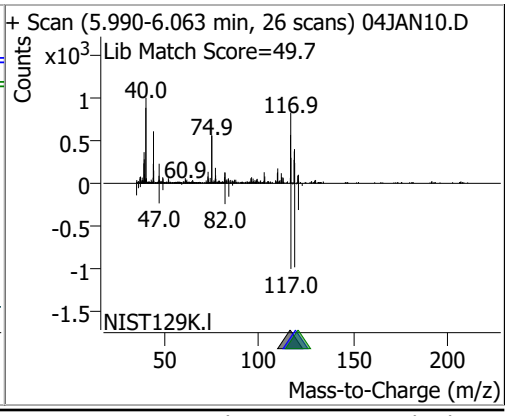




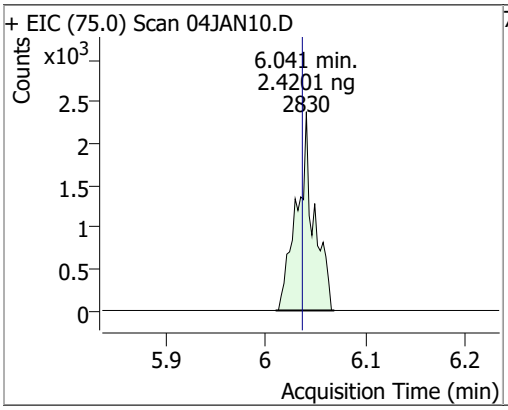
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	3.2043	6.03	0.01	4342	119.0	72.7	67.2	127.2
					121.0	20.8	0.1	60.1

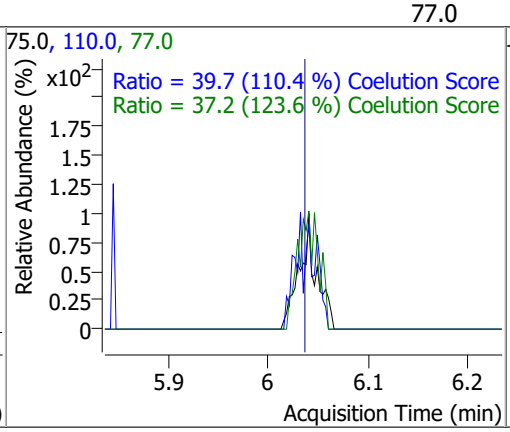


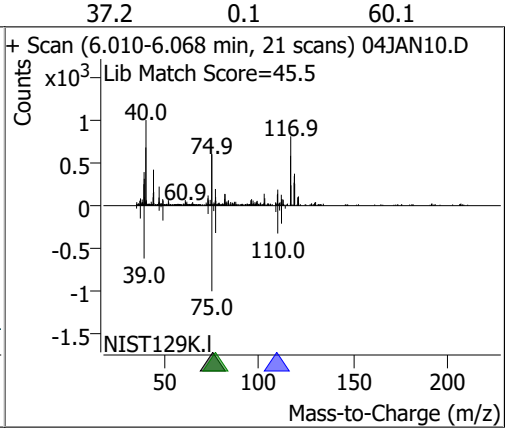




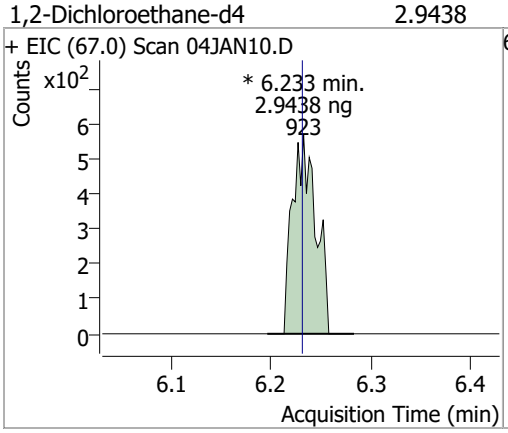
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	2.4201	6.04	0.00	2830	110.0	39.7	5.9	65.9
					77.0	37.2	0.1	60.1

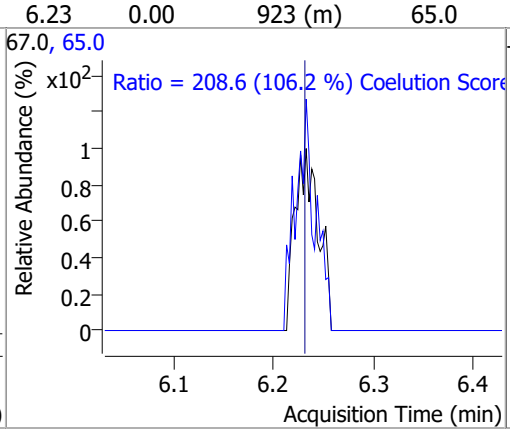


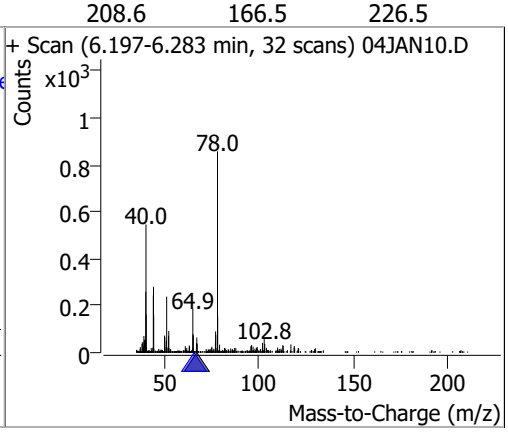




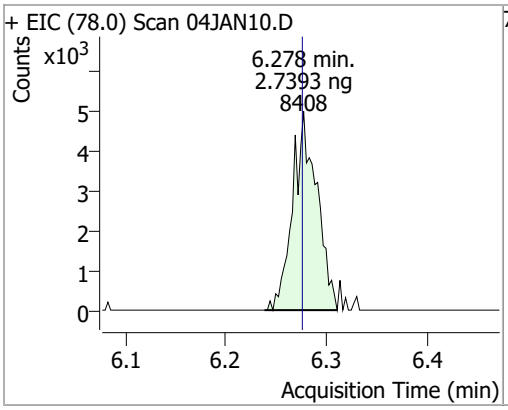
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	2.9438	6.23	0.00	923 (m)	65.0	208.6	166.5	226.5

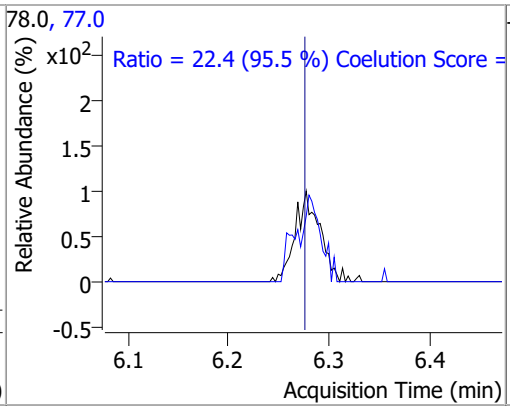


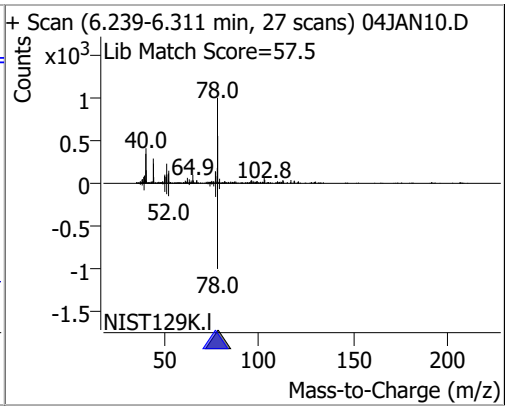




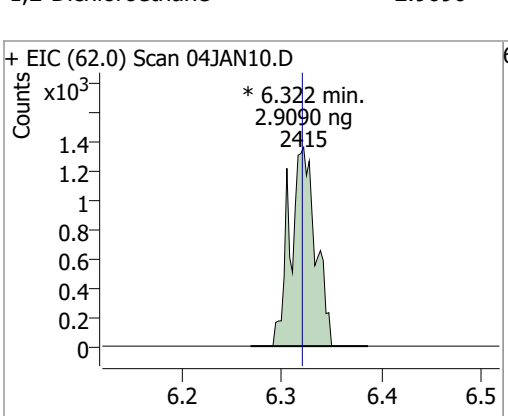
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	2.7393	6.28	0.00	8408	77.0	22.4	0.0	53.5

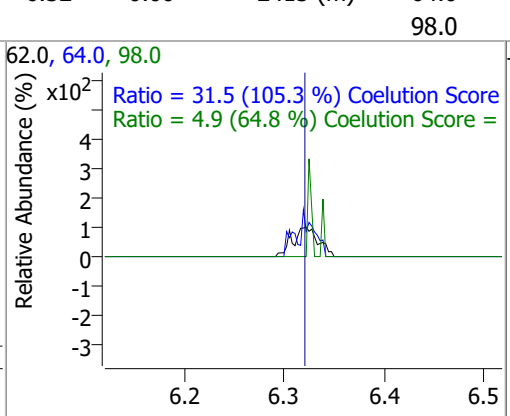


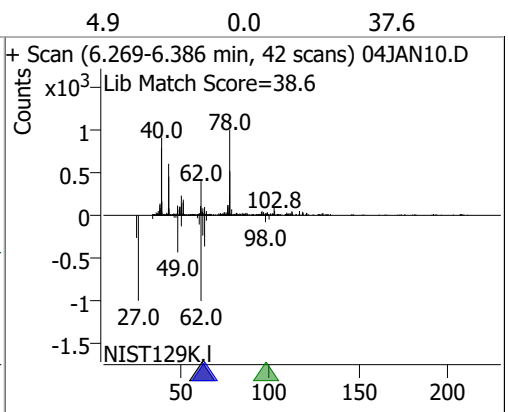




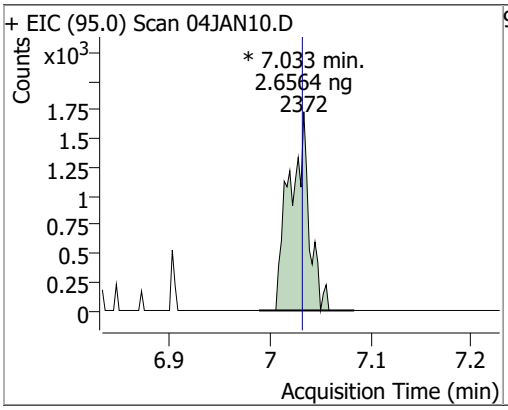
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	2.9090	6.32	0.00	2415 (m)	64.0	31.5	0.0	59.9
					98.0	4.9	0.0	37.6

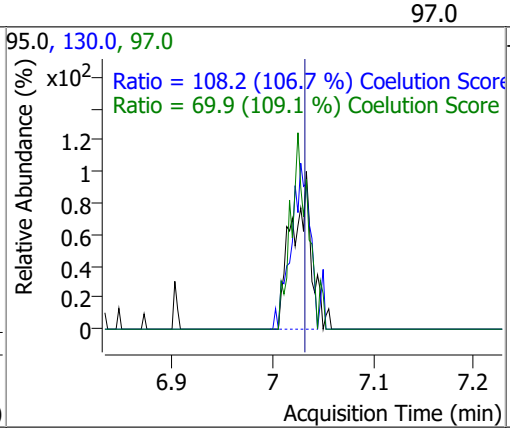


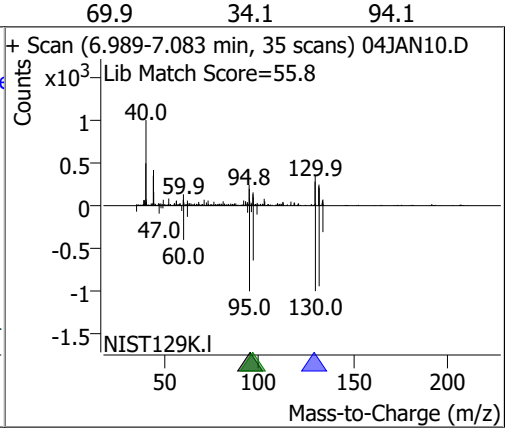




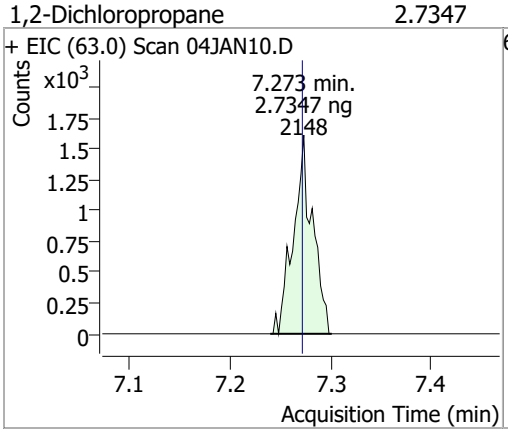
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	2.6564	7.03	0.00	2372 (m)	130.0	108.2	71.5	131.5
					97.0	69.9	34.1	94.1

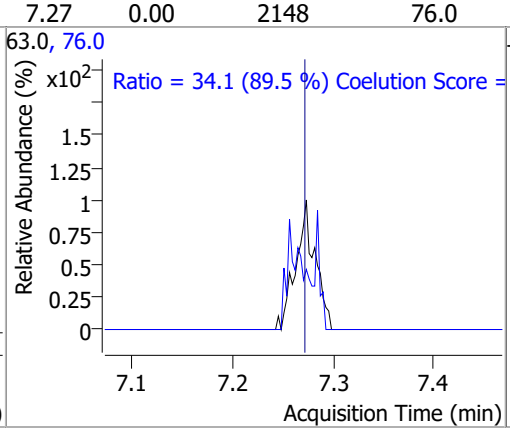


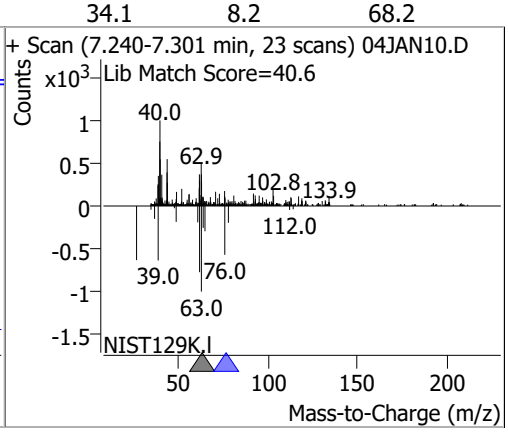




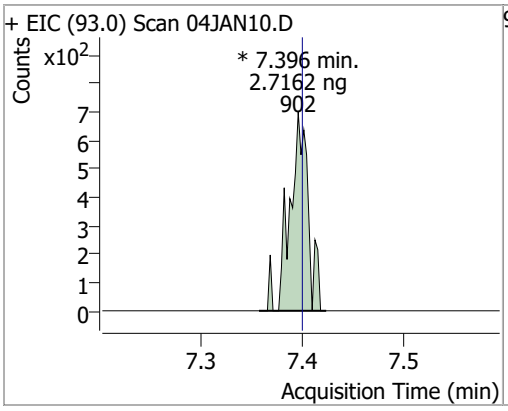
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	2.7347	7.27	0.00	2148	76.0	34.1	8.2	68.2

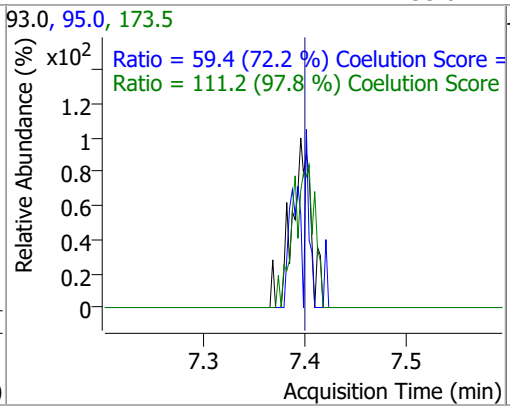


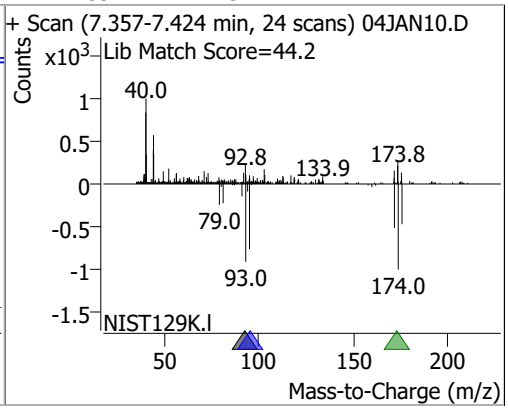




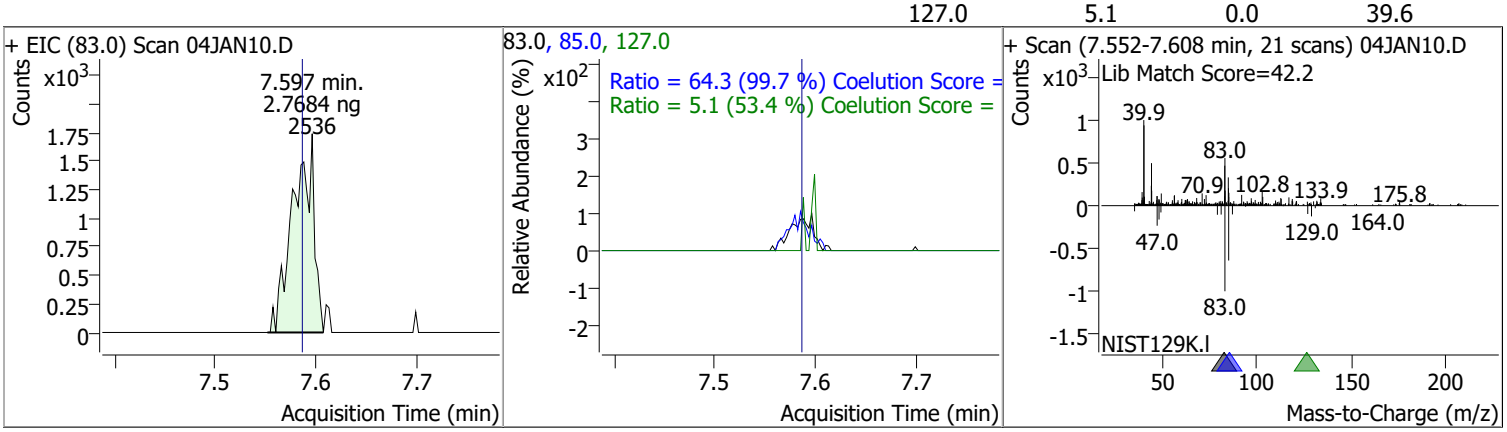
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	2.7162	7.40	0.00	902 (m)	173.5	111.2	83.7	143.7
					95.0	59.4	52.2	112.2



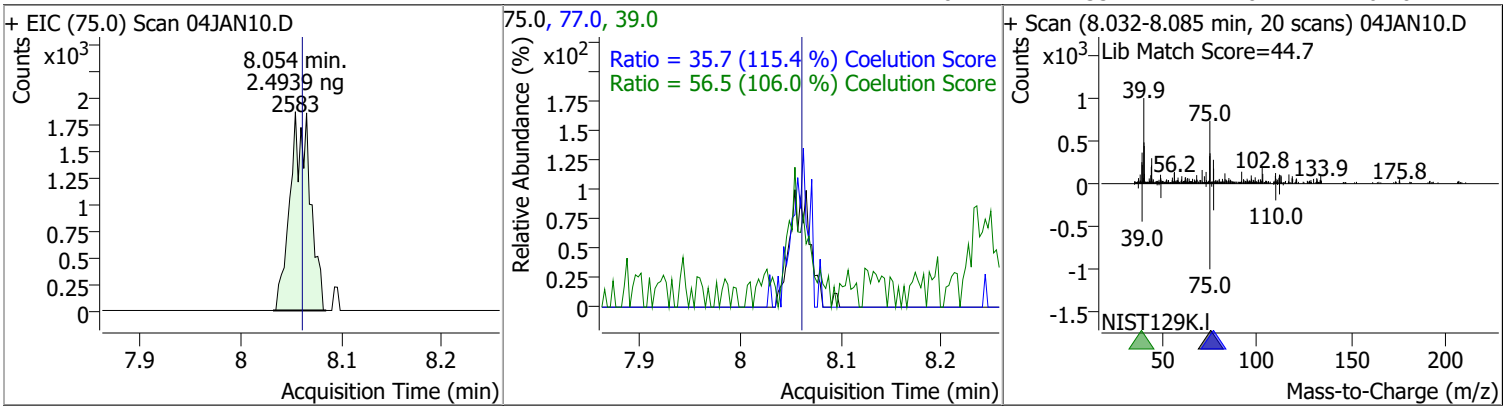




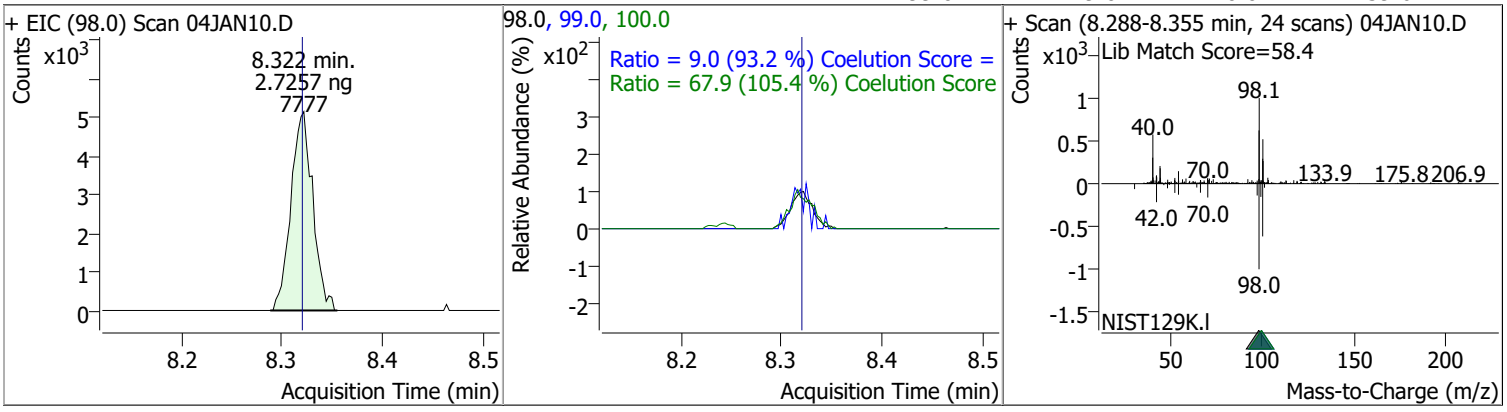
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	2.7684	7.60	0.01	2536	85.0	64.3	34.5	94.5
					127.0	5.1	0.0	39.6



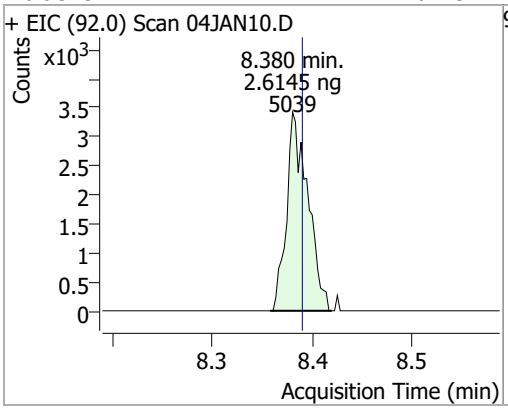
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	2.4939	8.05	-0.01	2583	39.0	56.5	23.3	83.3
					77.0	35.7	1.0	61.0

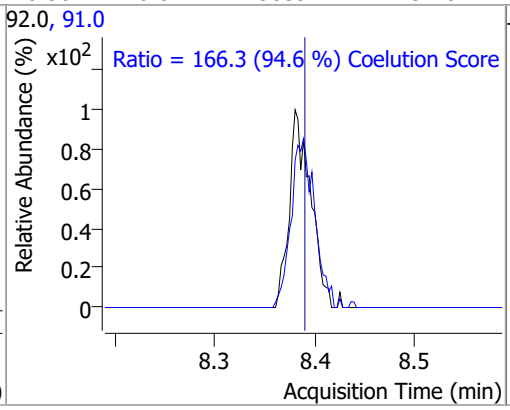


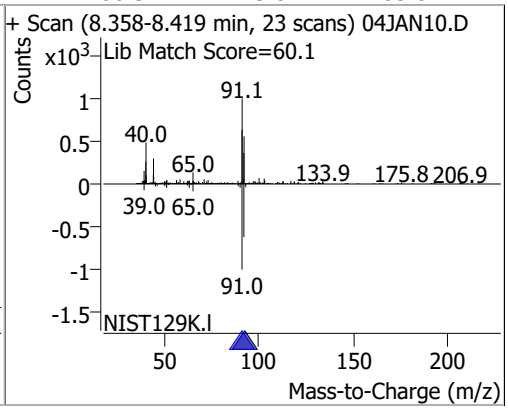
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	2.7257	8.32	0.00	7777	100.0	67.9	34.4	94.4
					99.0	9.0	0.0	39.6



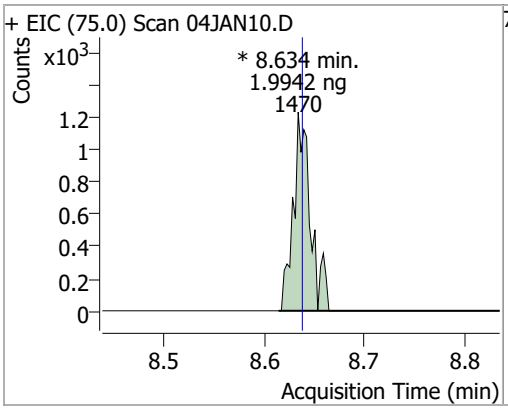
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	2.6145	8.38	-0.01	5039	91.0	166.3	145.8	205.8

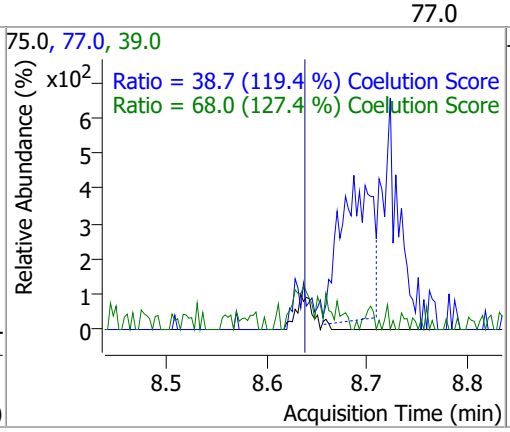


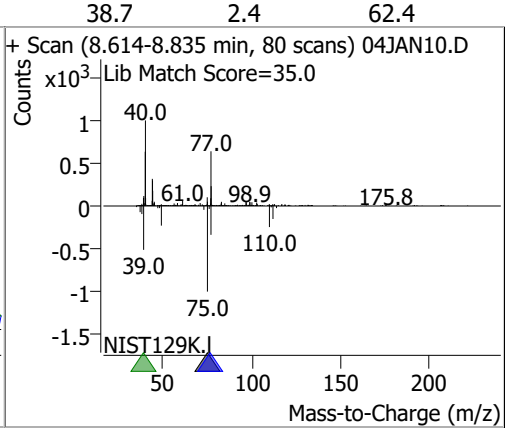




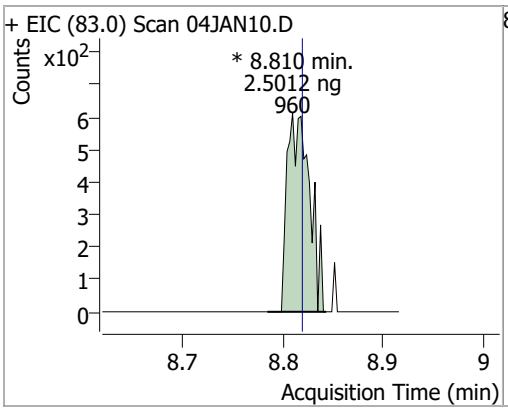
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	1.9942	8.63	0.00	1470 (m)	39.0	68.0	23.4	83.4
					77.0	38.7	2.4	62.4

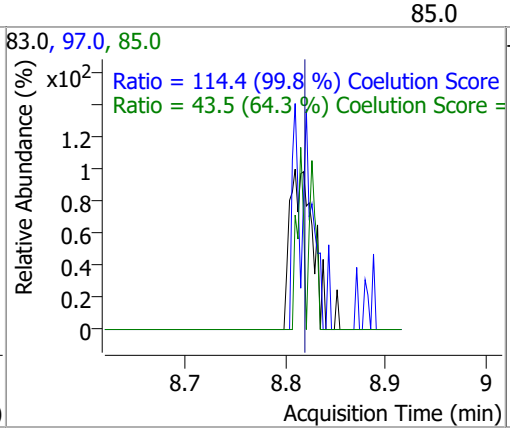


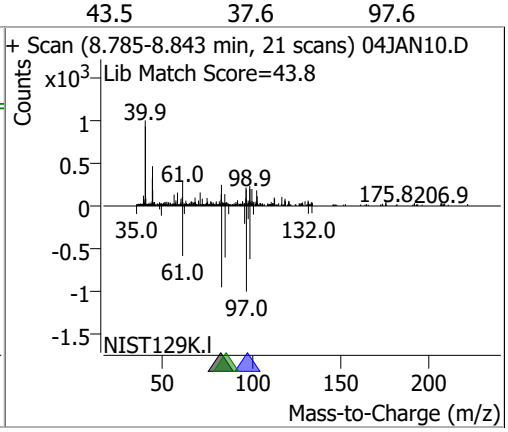




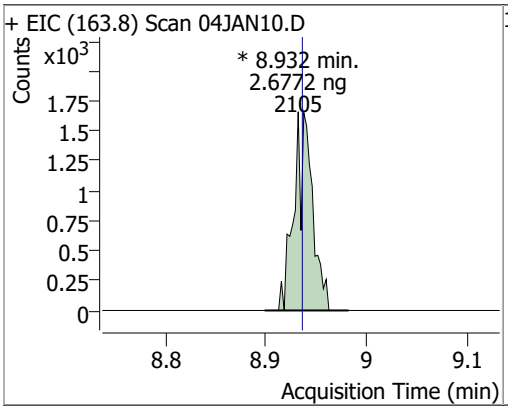
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	2.5012	8.81	-0.01	960 (m)	97.0	114.4	84.6	144.6
					85.0	43.5	37.6	97.6

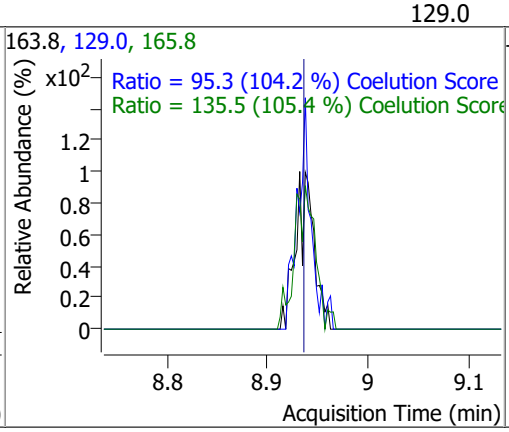


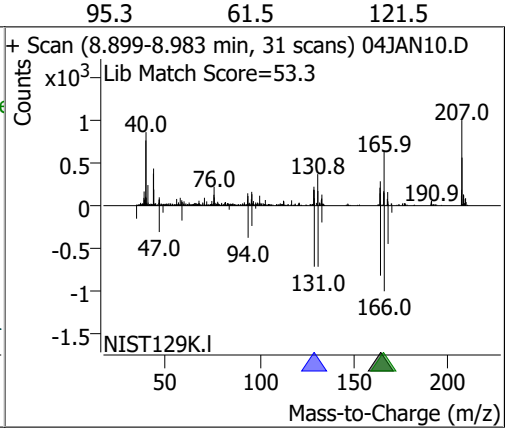




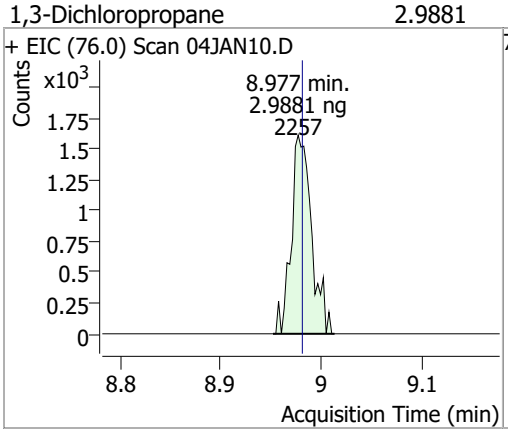
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	2.6772	8.93	0.00	2105 (m)	165.8	135.5	98.6	158.6
					129.0	95.3	61.5	121.5

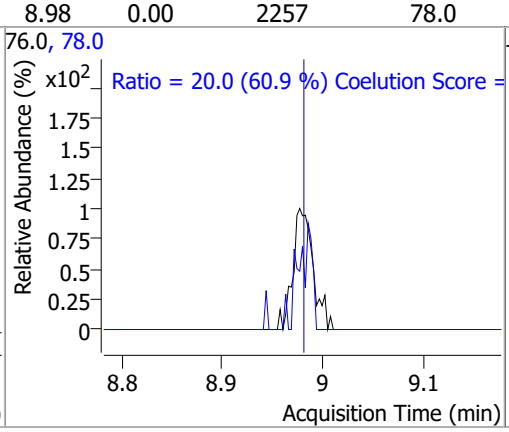


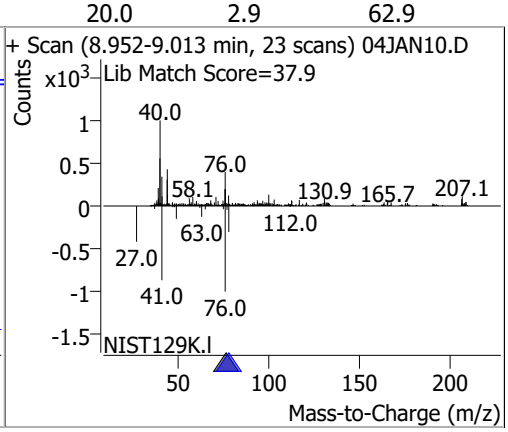




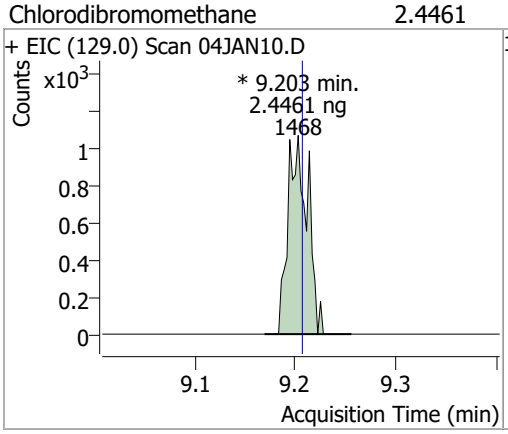
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	2.9881	8.98	0.00	2257	78.0	20.0	2.9	62.9

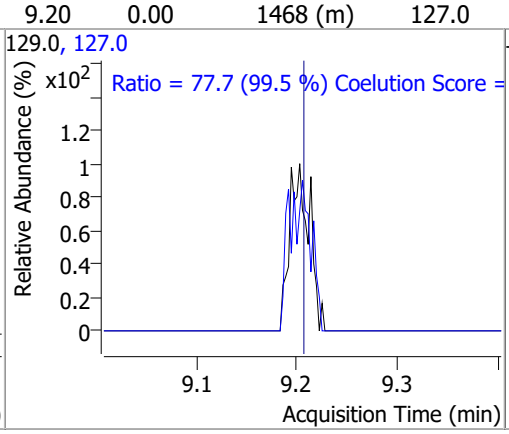


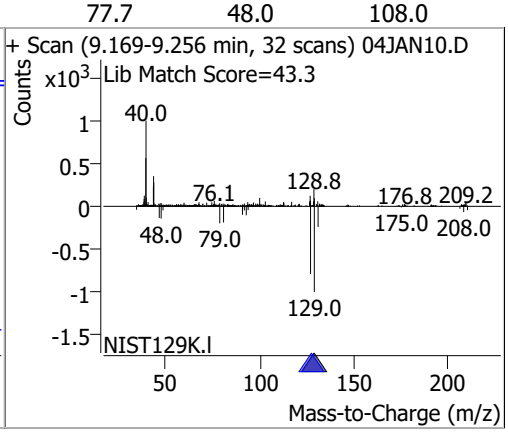




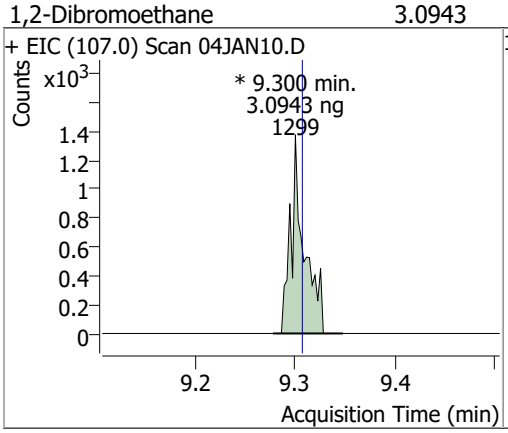
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	2.4461	9.20	0.00	1468 (m)	127.0	77.7	48.0	108.0

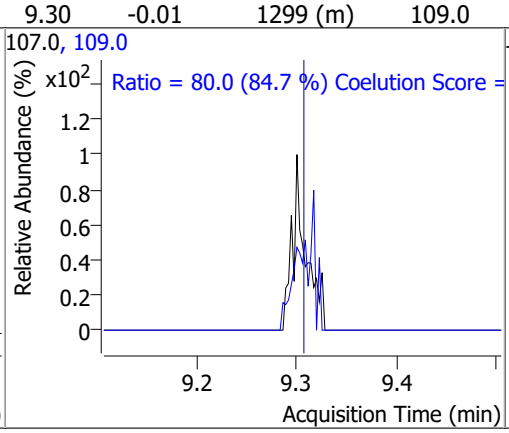


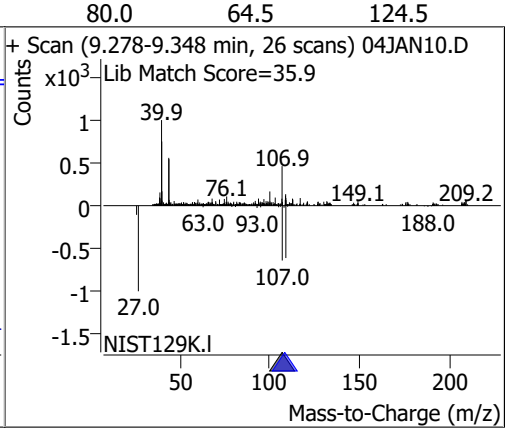




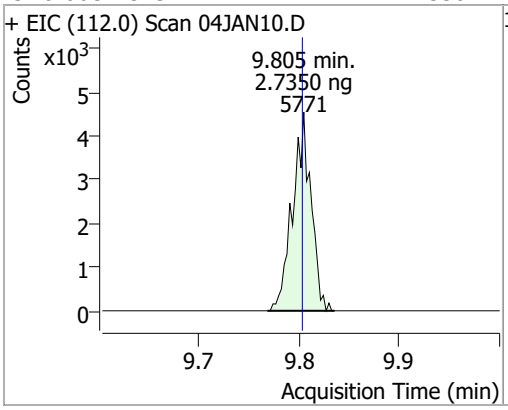
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	3.0943	9.30	-0.01	1299 (m)	109.0	80.0	64.5	124.5

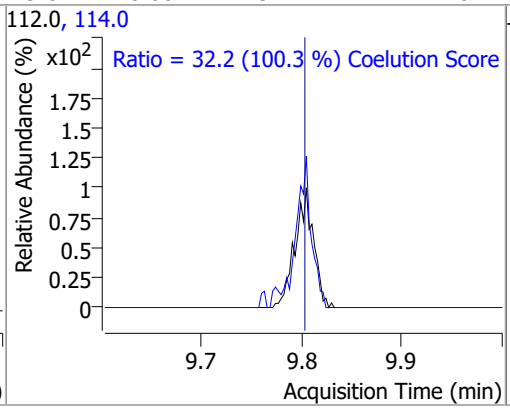


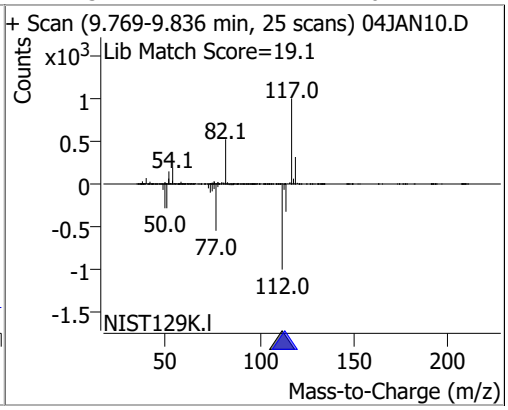




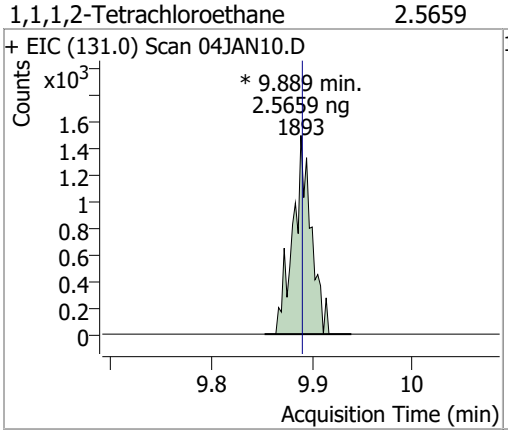
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	2.7350	9.81	0.00	5771	114.0	32.2	2.1	62.1

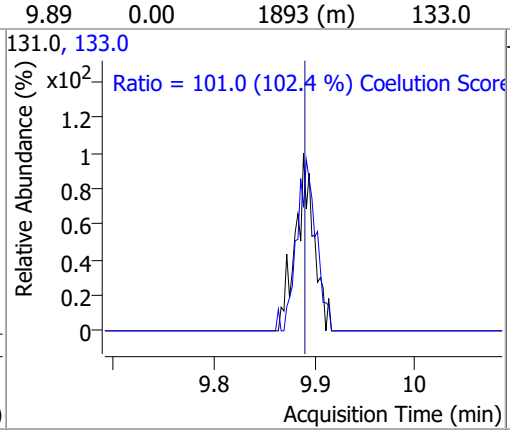


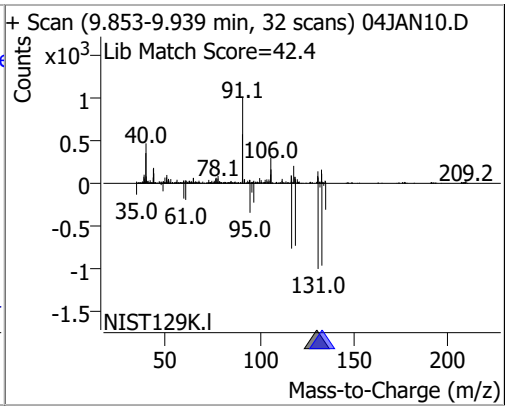




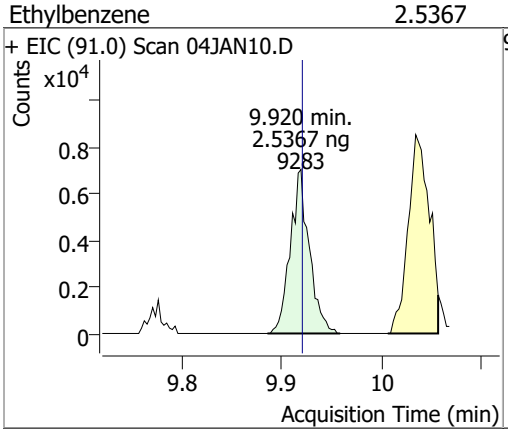
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	2.5659	9.89	0.00	1893 (m)	133.0	101.0	68.6	128.6

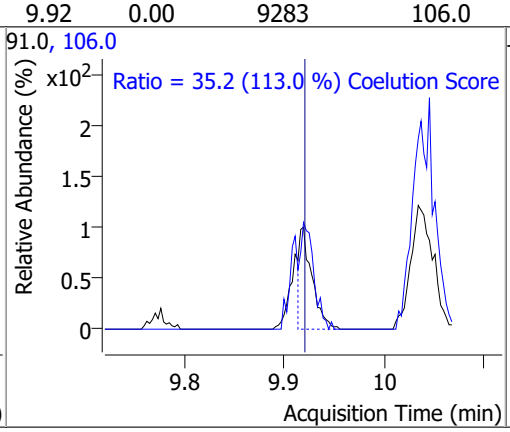


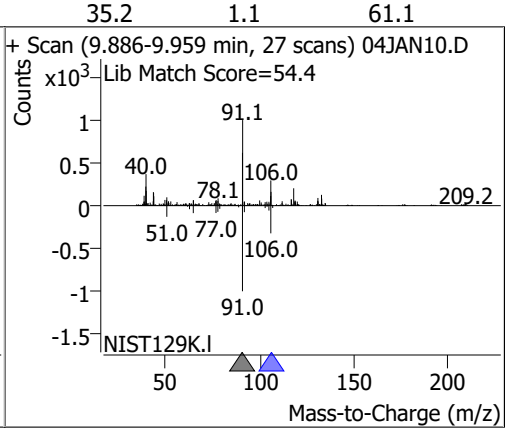




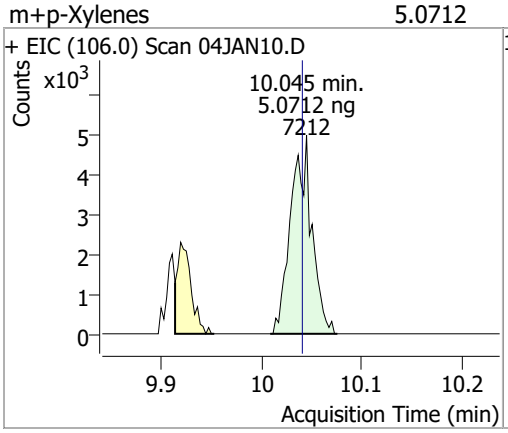
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	2.5367	9.92	0.00	9283	106.0	35.2	1.1	61.1

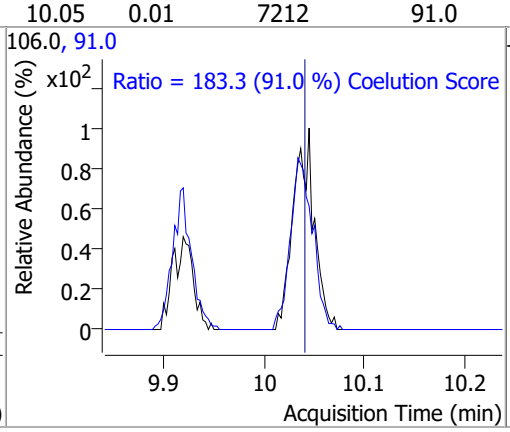


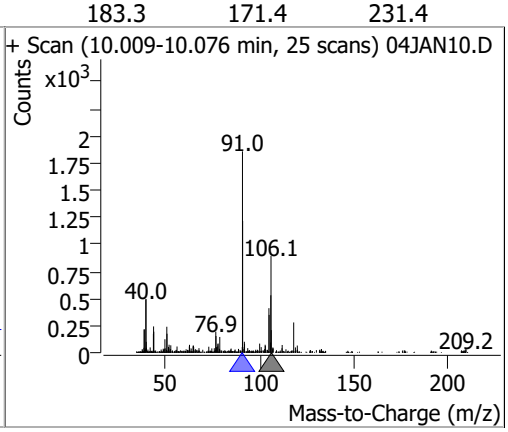




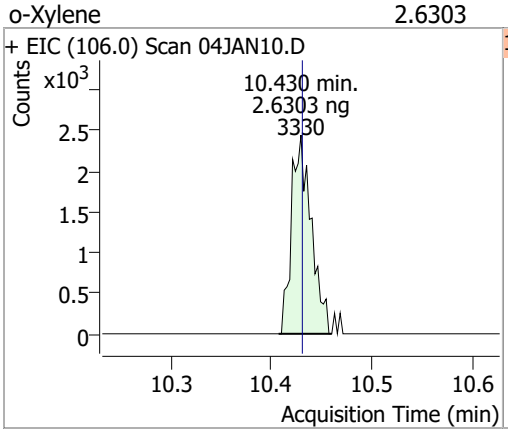
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	5.0712	10.05	0.01	7212	91.0	183.3	171.4	231.4

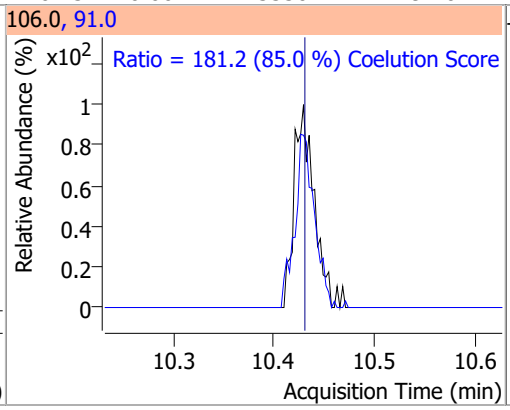


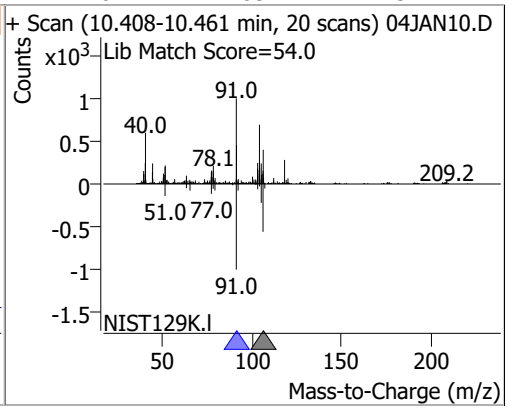




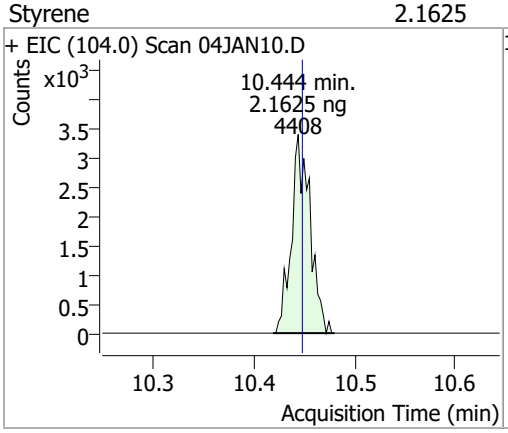
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	2.6303	10.43	0.00	3330	91.0	181.2	183.1	243.1

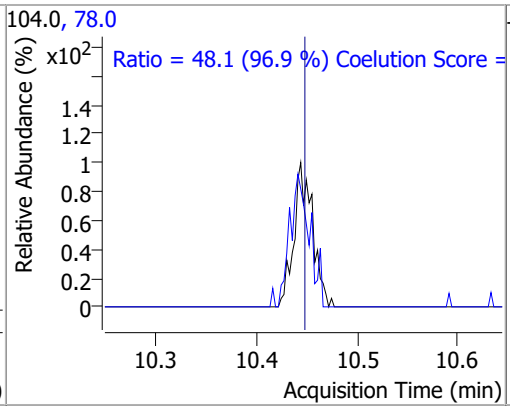


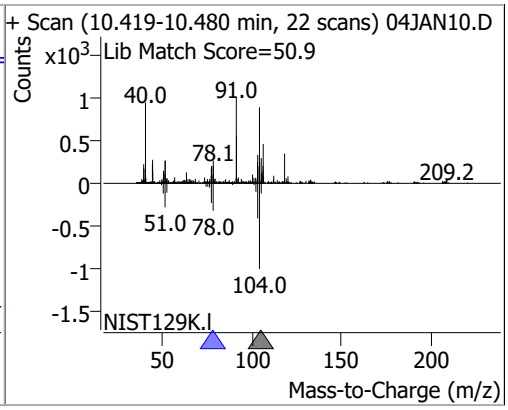




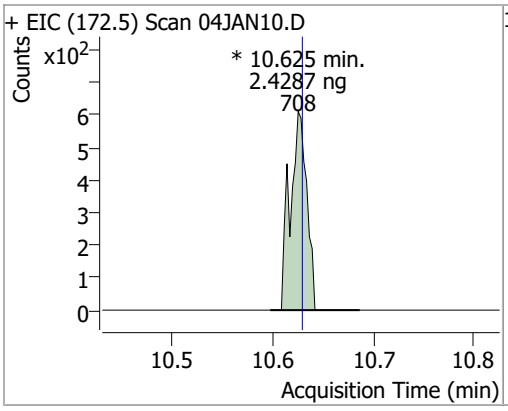
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	2.1625	10.44	0.00	4408	78.0	48.1	19.6	79.6

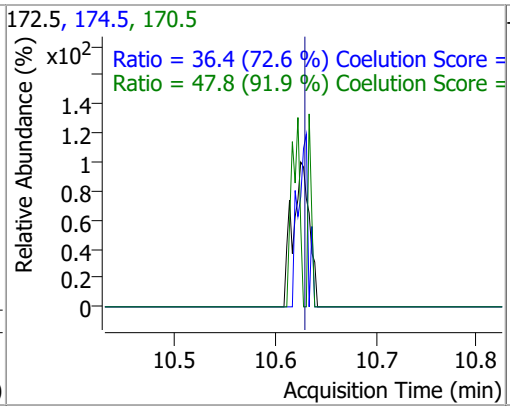


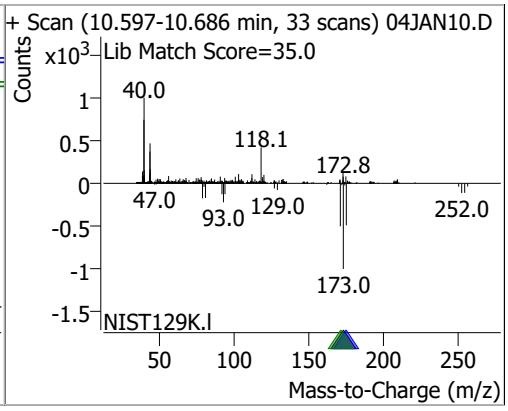




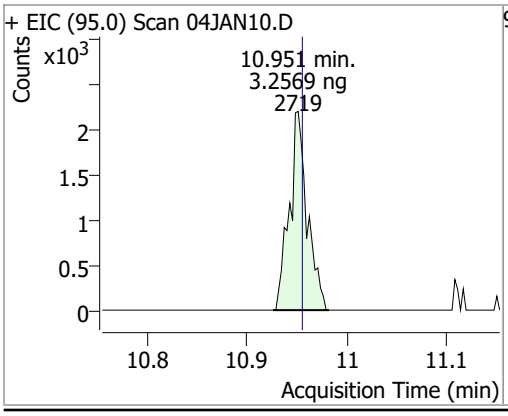
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	2.4287	10.63	0.00	708 (m)	170.5	47.8	22.1	82.1
					174.5	36.4	20.1	80.1

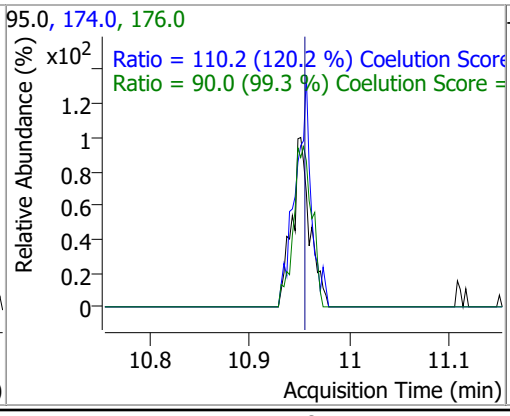


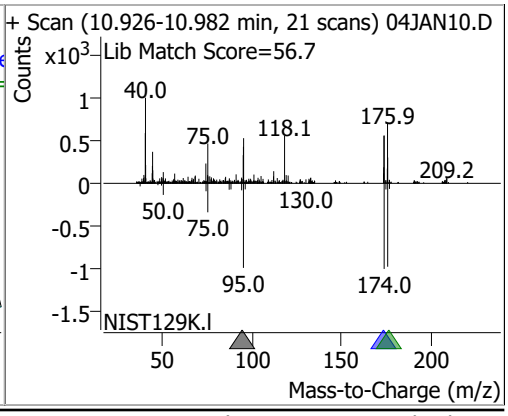




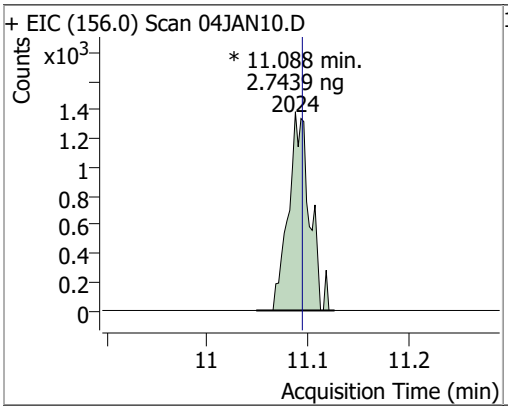
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	3.2569	10.95	0.00	2719	174.0	110.2	61.7	121.7
					176.0	90.0	60.6	120.6

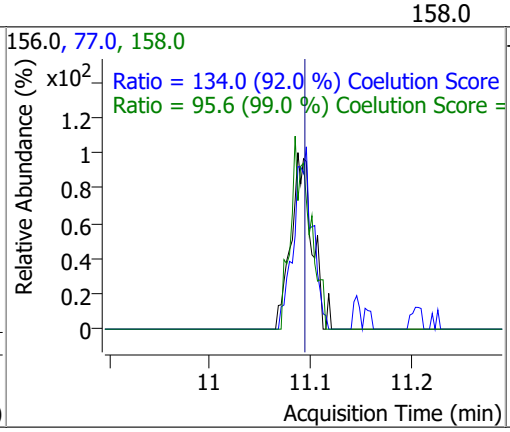


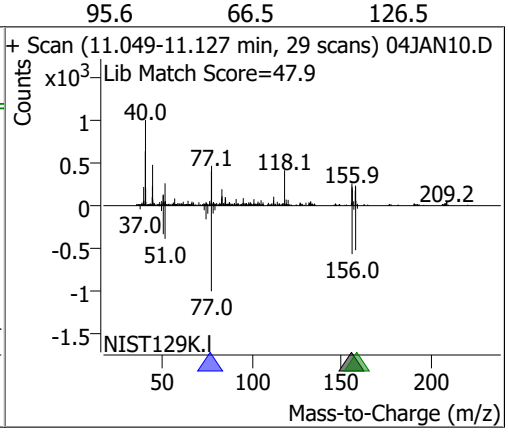




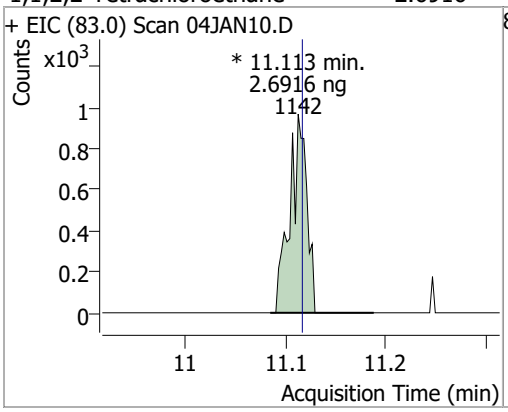
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	2.7439	11.09	-0.01	2024 (m)	77.0	134.0	115.7	175.7
					158.0	95.6	66.5	126.5

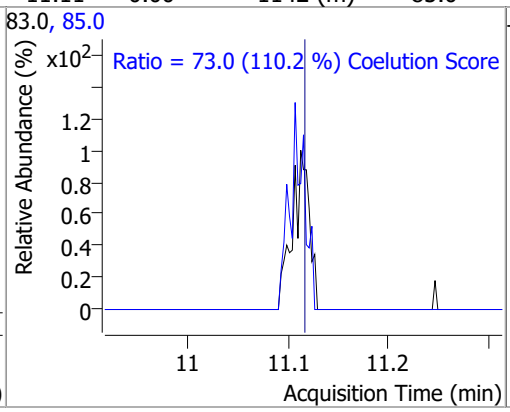


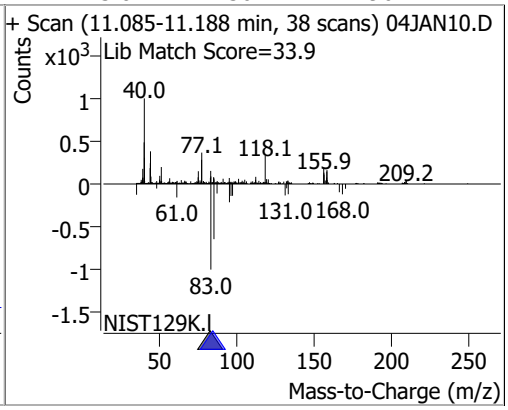




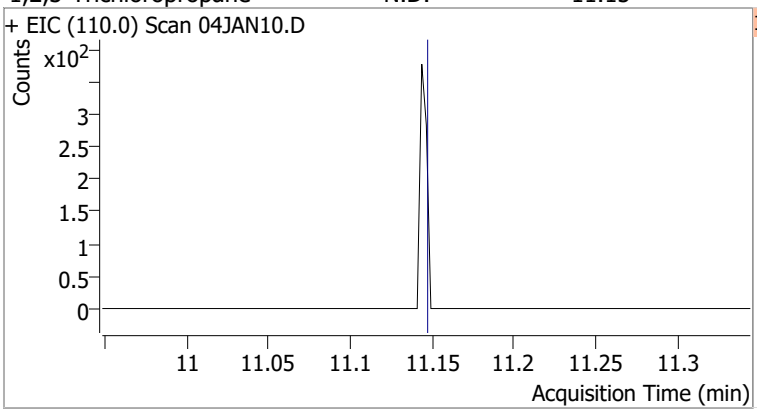
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	2.6916	11.11	0.00	1142 (m)	85.0	73.0	36.2	96.2

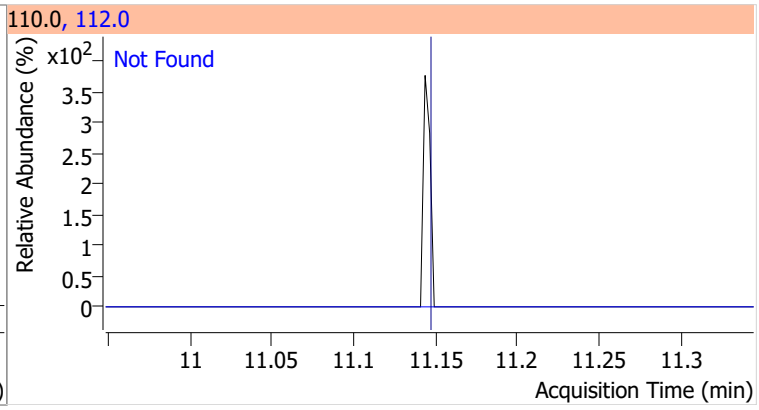




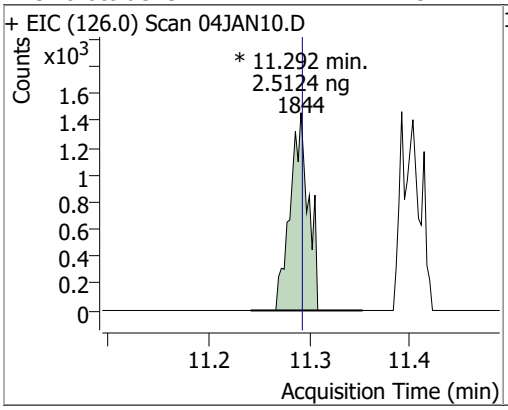


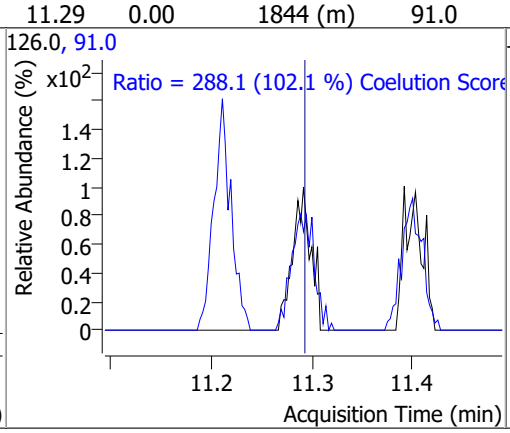
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5

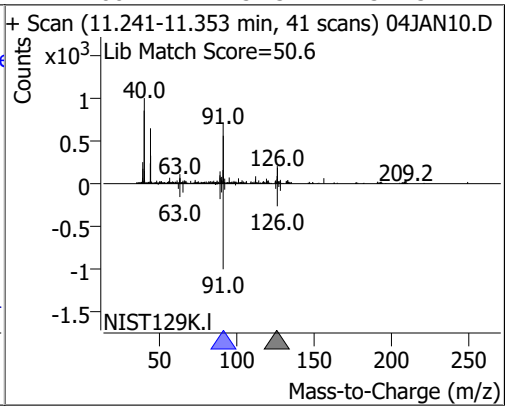




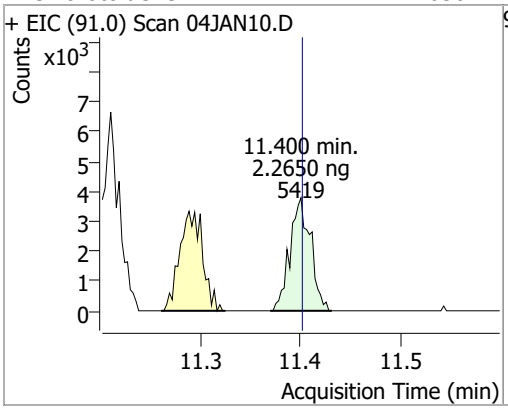
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	2.5124	11.29	0.00	1844 (m)	91.0	288.1	252.3	312.3

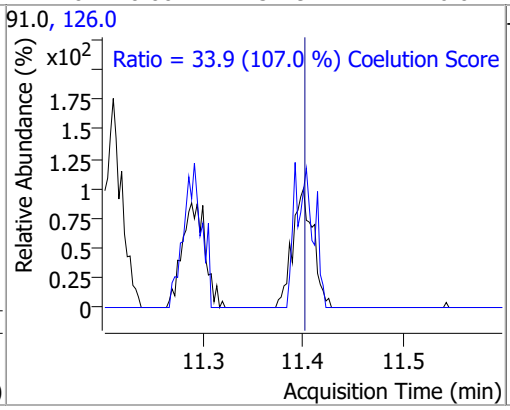


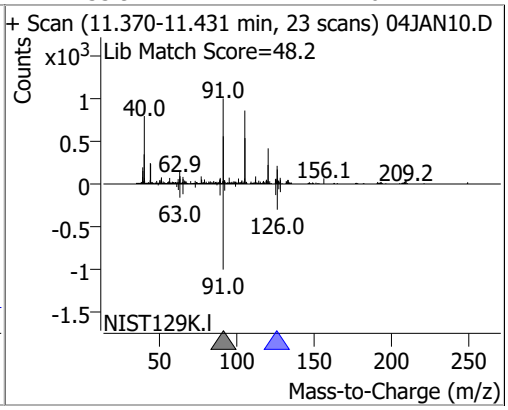




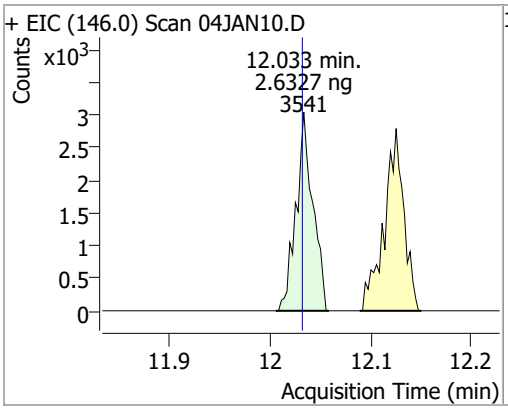
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	2.2650	11.40	0.00	5419	126.0	33.9	1.7	61.7

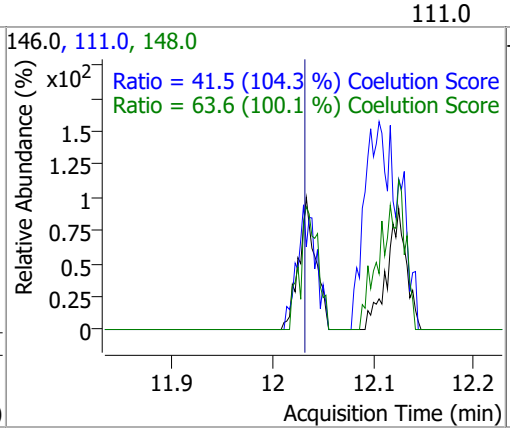


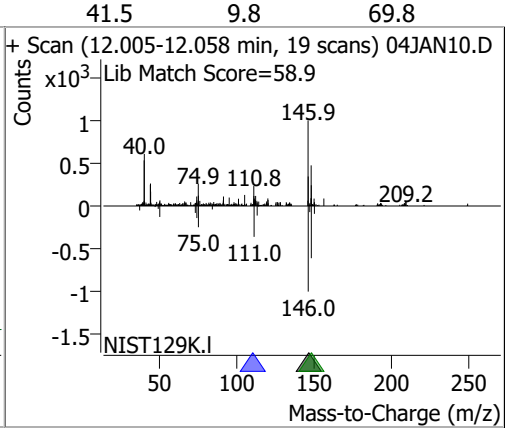




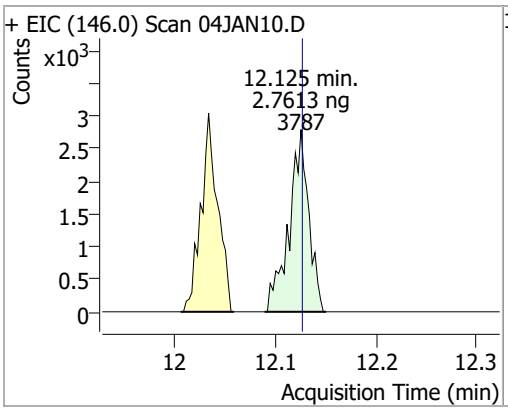
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	2.6327	12.03	0.00	3541	148.0	63.6	33.6	93.6
					111.0	41.5	9.8	69.8

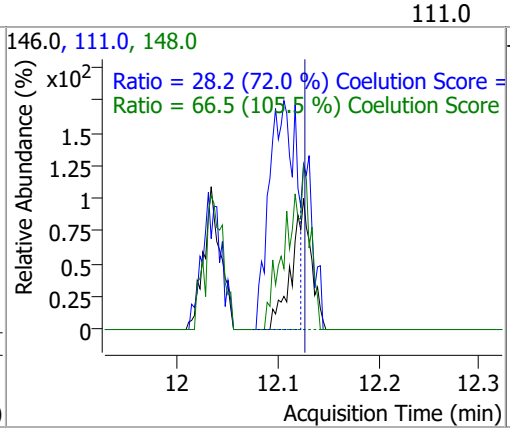


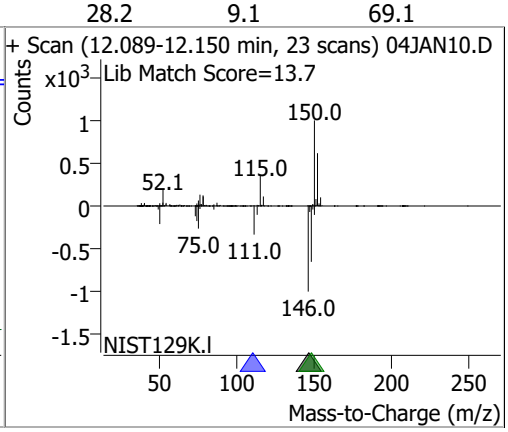




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	2.7613	12.13	0.00	3787	148.0	66.5	33.1	93.1
					111.0	28.2	9.1	69.1

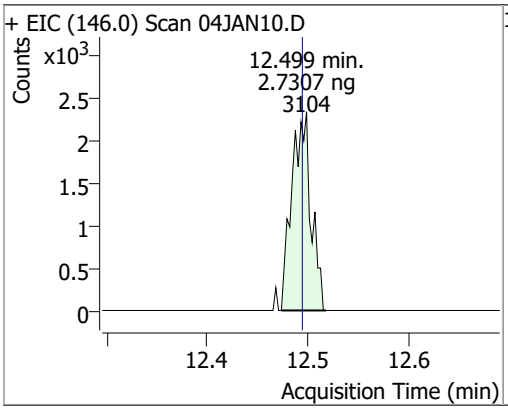


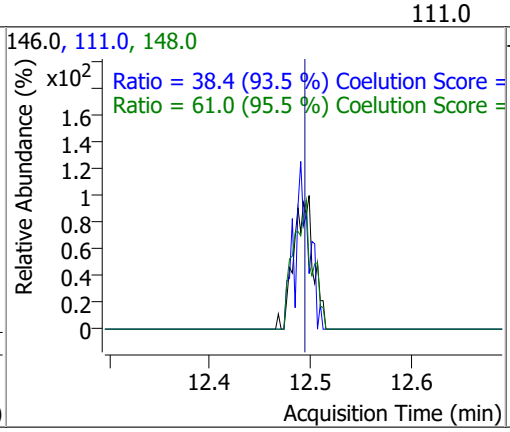


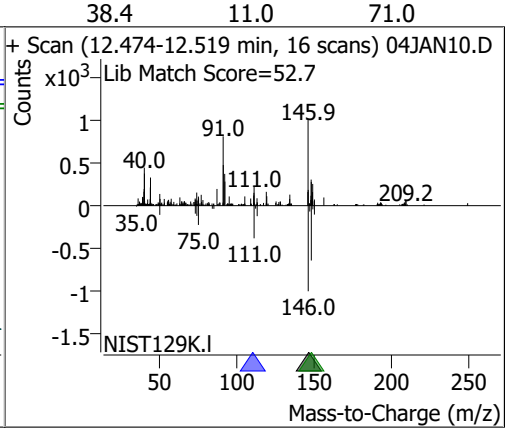


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	2.7307	12.50	0.01	3104	148.0	61.0	33.9	93.9
					111.0	38.4	11.0	71.0



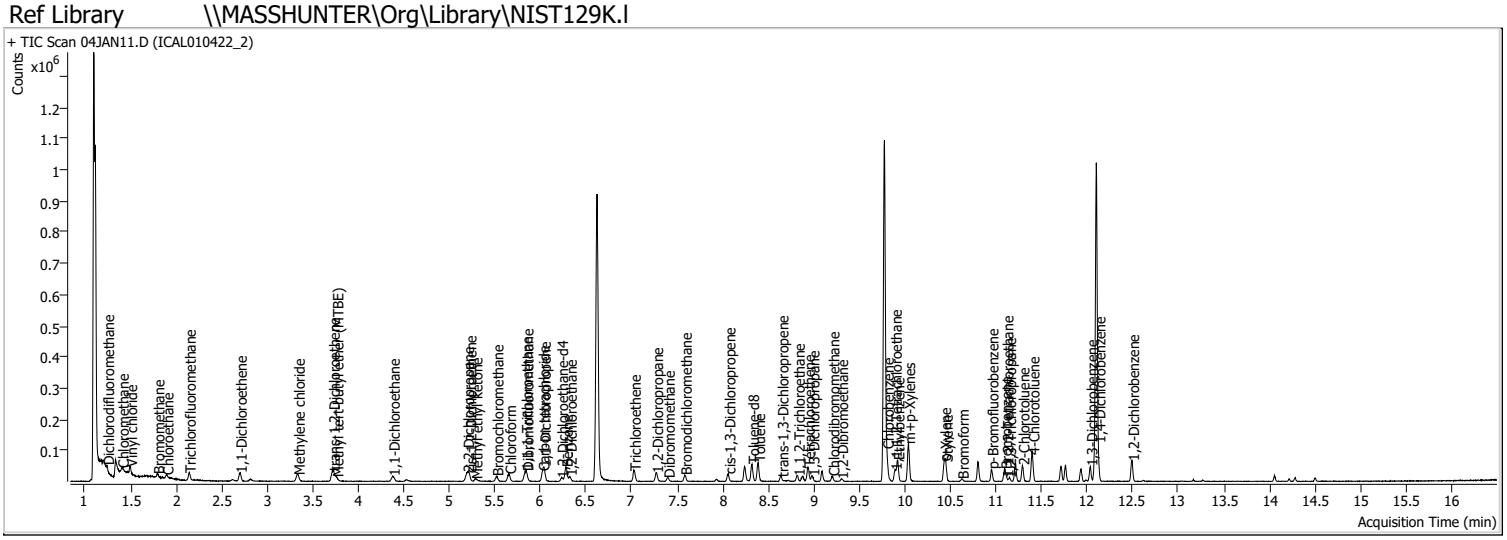




Quantitation Results Report (QT Reviewed)

Agilent | Trusted Answers

Data File	04JAN11.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/4/2022 4:00:35 PM
Sample Name	ICAL010422_2	Instrument	VOA5975C
Vial	11	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

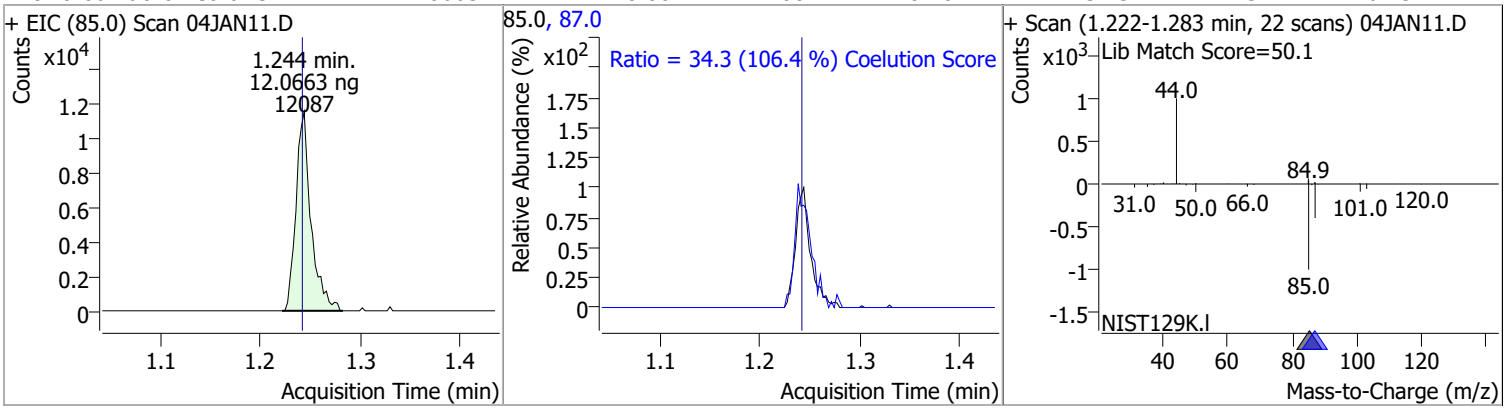


Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	764419	250.0000	ng		-0.003
M Chlorobenzene-d5	9.774	82.0	296554	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.103	152.0	242142	250.0000	ng		0.003
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	9074	12.6000	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 5.04%		*	
S 1,2-Dichloroethane-d4	6.227	67.0	3938	12.6600	ng		-0.006
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 5.06%		*	
S Toluene-d8	8.322	98.0	32318	11.3089	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 4.52%		*	
S p-Bromofluorobenzene	10.948	95.0	10059	11.3393	ng		-0.006
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 4.54%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.244	85.0	12087	12.0663	ng		96
T Chloromethane	1.406	50.0	16859	13.8661	ng		100
T Vinyl chloride	1.498	62.0	13724	12.5446	ng		75
T Bromomethane	1.796	96.0	5893	12.0464	ng		97
T Chloroethane	1.897	64.0	8052	14.8670	ng	m	98
T Trichlorofluoromethane	2.142	101.0	15431	11.3637	ng		95
T 1,1-Dichloroethene	2.700	96.0	9169	11.9081	ng		96
T Methylene chloride	3.338	49.0	17734	15.6236	ng		93
T trans-1,2-Dichloroethene	3.720	96.0	9821	12.5022	ng	m	95
T Methyl tert-butyl ether (MTBE)	3.762	73.0	12515	12.3255	ng	m	99
T 1,1-Dichloroethane	4.378	63.0	17642	12.0652	ng		94
T 2,2-Dichloropropane	5.196	77.0	13676	12.4820	ng		95
T cis-1,2-Dichloroethene	5.221	96.0	10008	12.5659	ng		95
T Methyl ethyl ketone	5.288	43.0	13167	122.0520	ng		95
T Bromochloromethane	5.516	128.0	4275	12.9568	ng		91
T Chloroform	5.656	83.0	19015	13.0668	ng		98

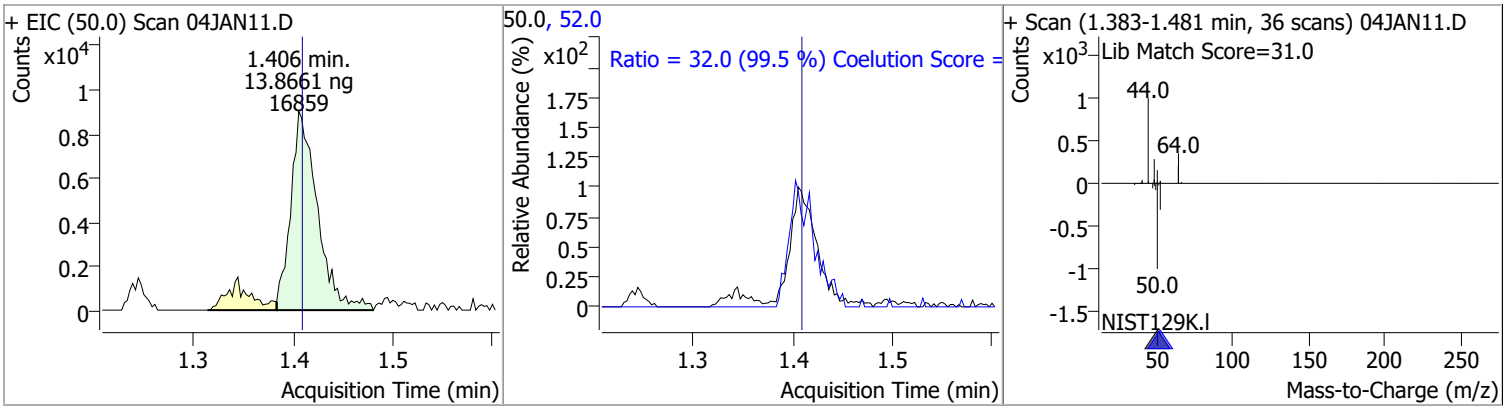
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	5.837	97.0	16623	12.1891	ng		97
T Carbon tetrachloride	6.021	117.0	16466	12.2545	ng		98
T 1,1-Dichloropropene	6.038	75.0	13149	11.3397	ng		94
T Benzene	6.278	78.0	37071	12.1801	ng		99
T 1,2-Dichloroethane	6.322	62.0	10202	12.3906	ng		92
T Trichloroethene	7.025	95.0	10442	11.6753	ng		94
T 1,2-Dichloropropane	7.270	63.0	9488	12.0602	ng		99
T Dibromomethane	7.399	93.0	4675	14.0619	ng		93
T Bromodichloromethane	7.585	83.0	11562	12.6014	ng		97
T cis-1,3-Dichloropropene	8.062	75.0	12525	12.0738	ng		94
T Toluene	8.388	92.0	21794	11.2899	ng		97
T trans-1,3-Dichloropropene	8.645	75.0	8683	11.7589	ng		98
T 1,1,2-Trichloroethane	8.824	83.0	5090	13.2340	ng	m	91
T Tetrachloroethene	8.935	163.8	9238	11.7302	ng		99
T 1,3-Dichloropropane	8.985	76.0	8967	11.8526	ng		97
T Chlorodibromomethane	9.206	129.0	7718	12.8393	ng		97
T 1,2-Dibromoethane	9.300	107.0	5410	12.8640	ng		100
T Chlorobenzene	9.802	112.0	26461	12.5204	ng		99
T 1,1,1,2-Tetrachloroethane	9.889	131.0	9473	12.8225	ng		88
T Ethylbenzene	9.917	91.0	40470	11.0411	ng		99
T m+p-Xylenes	10.037	106.0	31538	22.1410	ng		100
T o-Xylene	10.430	106.0	13519	10.6612	ng		92
T Styrene	10.449	104.0	23472	11.4968	ng		100
T Bromoform	10.625	172.5	3652	11.7860	ng		92
T Bromobenzene	11.096	156.0	9663	12.3310	ng		96
T 1,1,2,2-Tetrachloroethane	11.116	83.0	5793	12.8437	ng		99
T 1,2,3-Trichloropropane	11.144	110.0	1654	13.7084	ng	m	99
T 2-Chlorotoluene	11.289	126.0	8731	11.1977	ng		94
T 4-Chlorotoluene	11.400	91.0	28532	11.2233	ng		100
T 1,3-Dichlorobenzene	12.036	146.0	16932	11.8473	ng		99
T 1,4-Dichlorobenzene	12.122	146.0	17438	11.9662	ng		94
T 1,2-Dichlorobenzene	12.493	146.0	14666	12.1423	ng		98

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

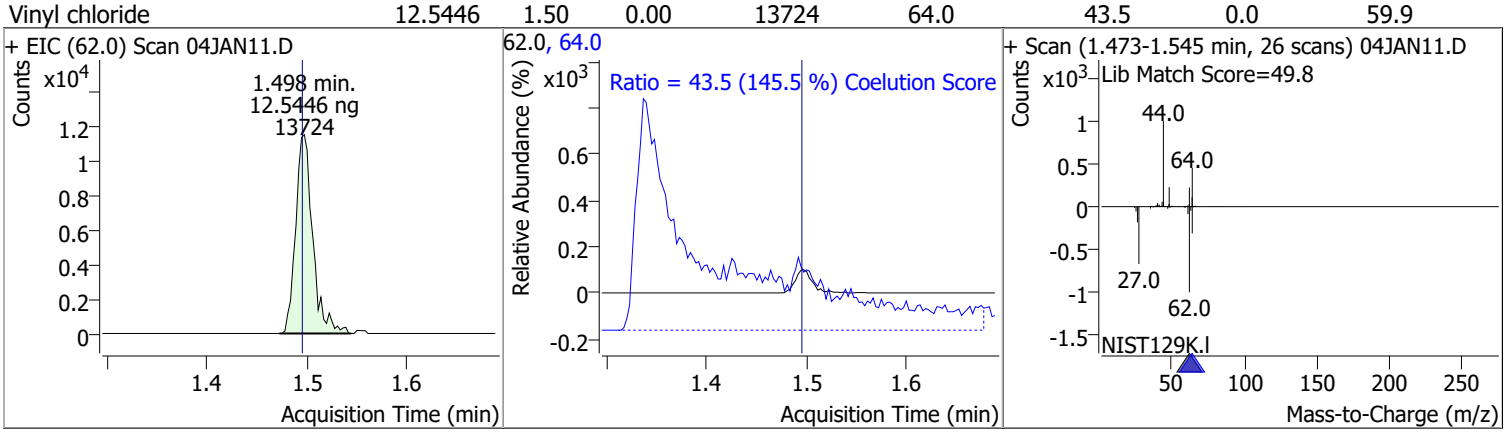
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	12.0663	1.24	0.00	12087	87.0	34.3	2.3	62.3



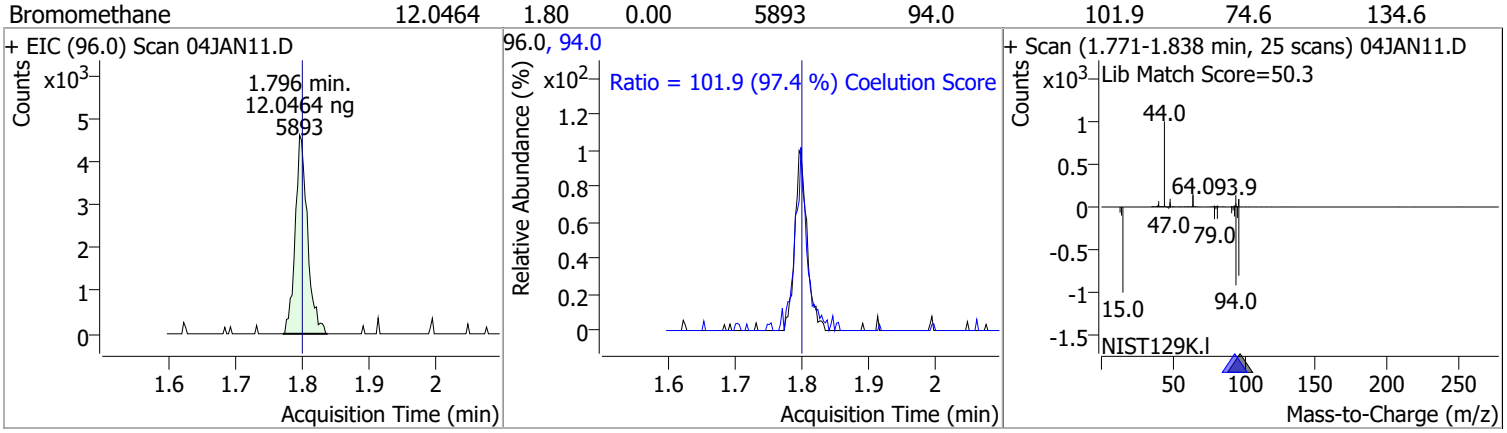
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	13.8661	1.41	0.00	16859	52.0	32.0	2.1	62.1



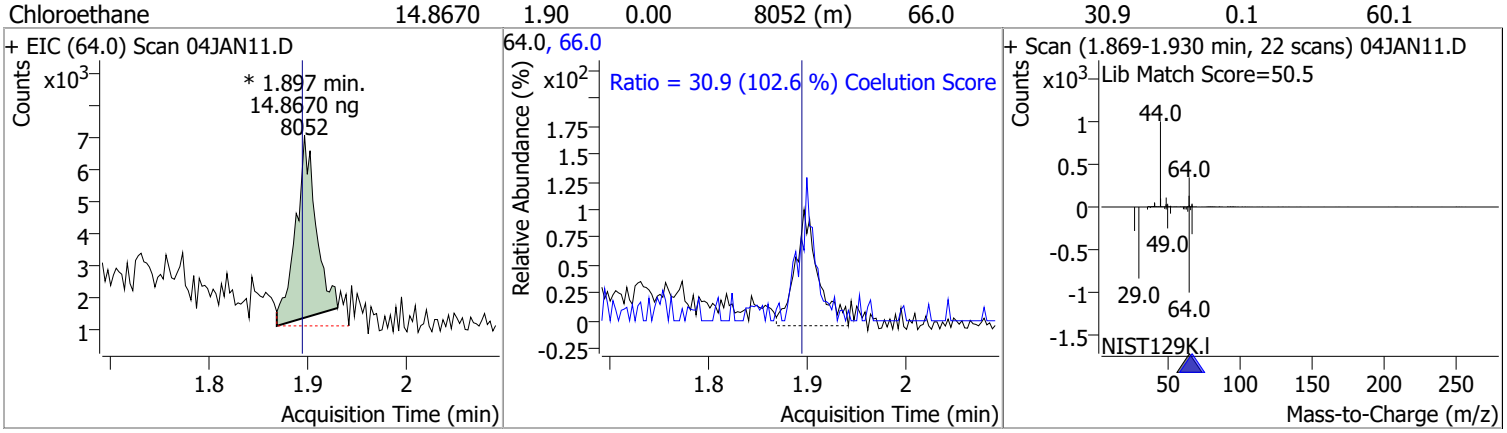
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	12.5446	1.50	0.00	13724	64.0	43.5	0.0	59.9



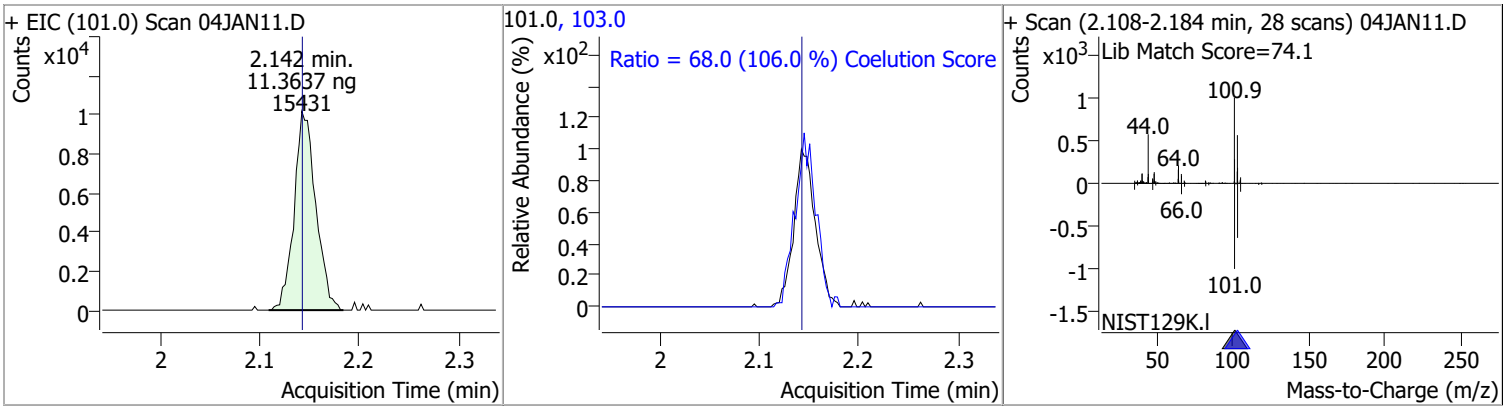
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	12.0464	1.80	0.00	5893	94.0	101.9	74.6	134.6



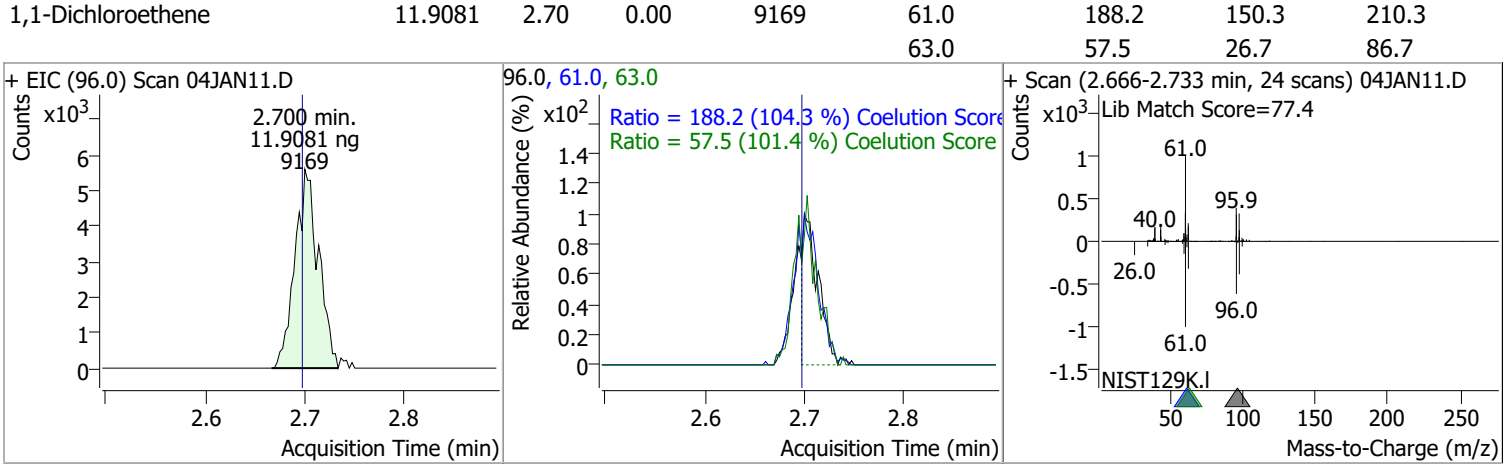
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	14.8670	1.90	0.00	8052 (m)	66.0	30.9	0.1	60.1



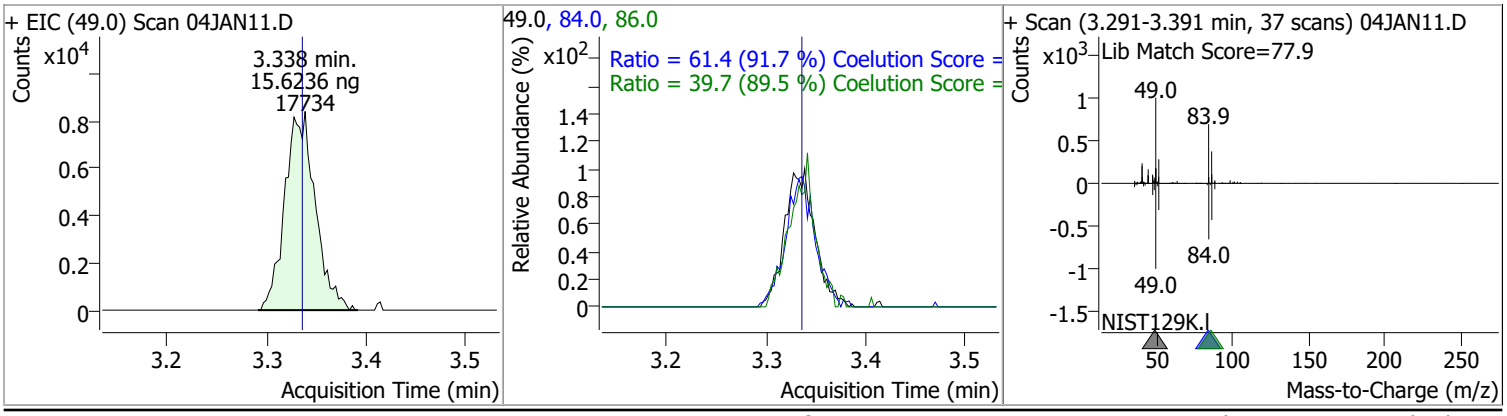
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	11.3637	2.14	0.00	15431	103.0	68.0	34.2	94.2



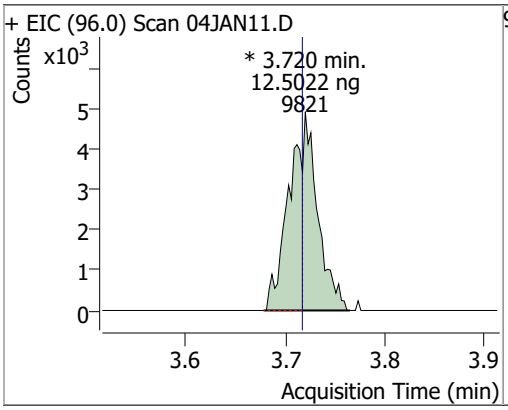
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	11.9081	2.70	0.00	9169	61.0	188.2	150.3	210.3
					63.0	57.5	26.7	86.7

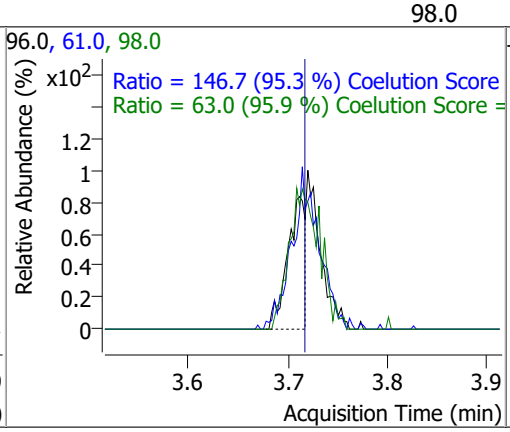


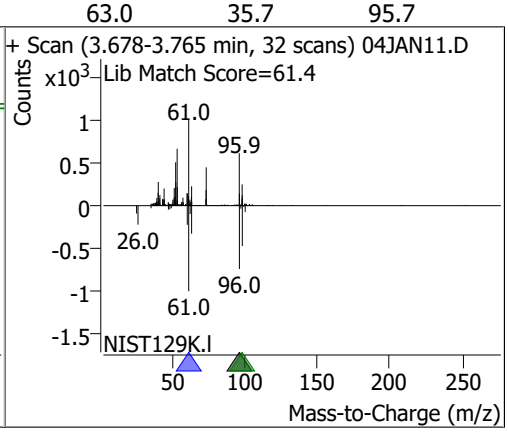
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	15.6236	3.34	0.00	17734	84.0	61.4	36.9	96.9
					86.0	39.7	14.3	74.3



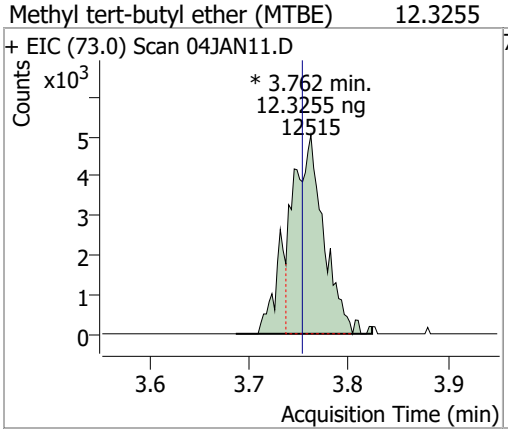
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	12.5022	3.72	0.00	9821 (m)	61.0	146.7	123.9	183.9
					98.0	63.0	35.7	95.7

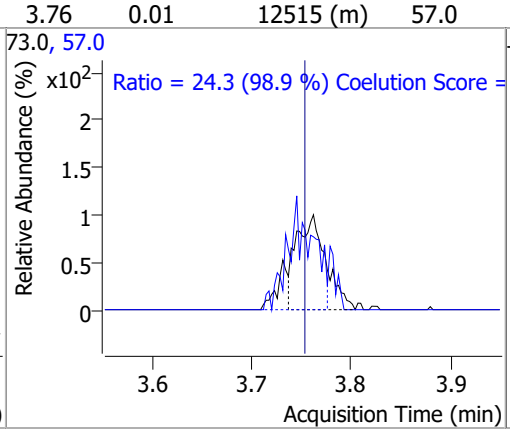


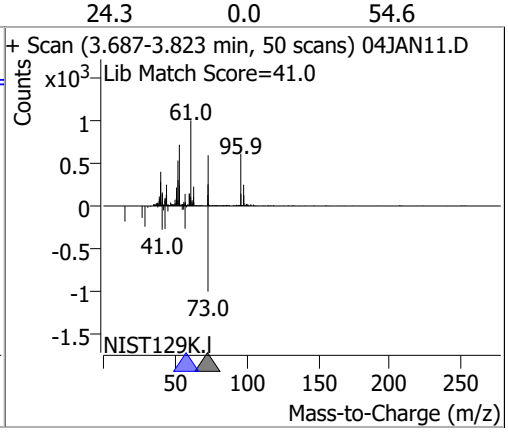




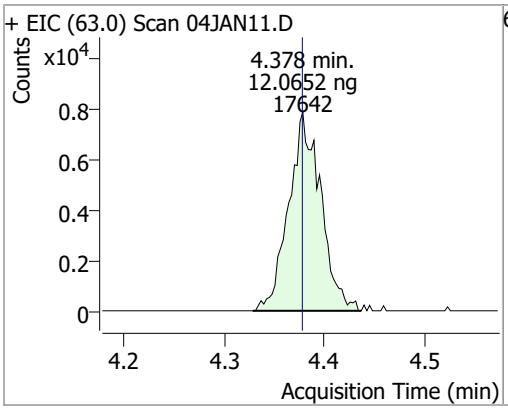
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	12.3255	3.76	0.01	12515 (m)	57.0	24.3	0.0	54.6

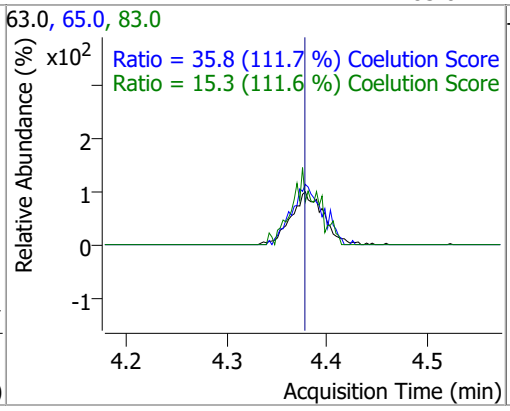


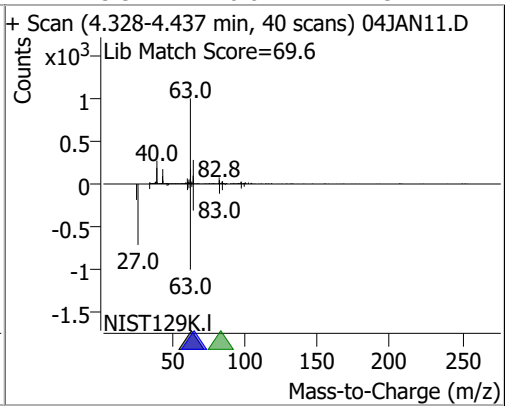




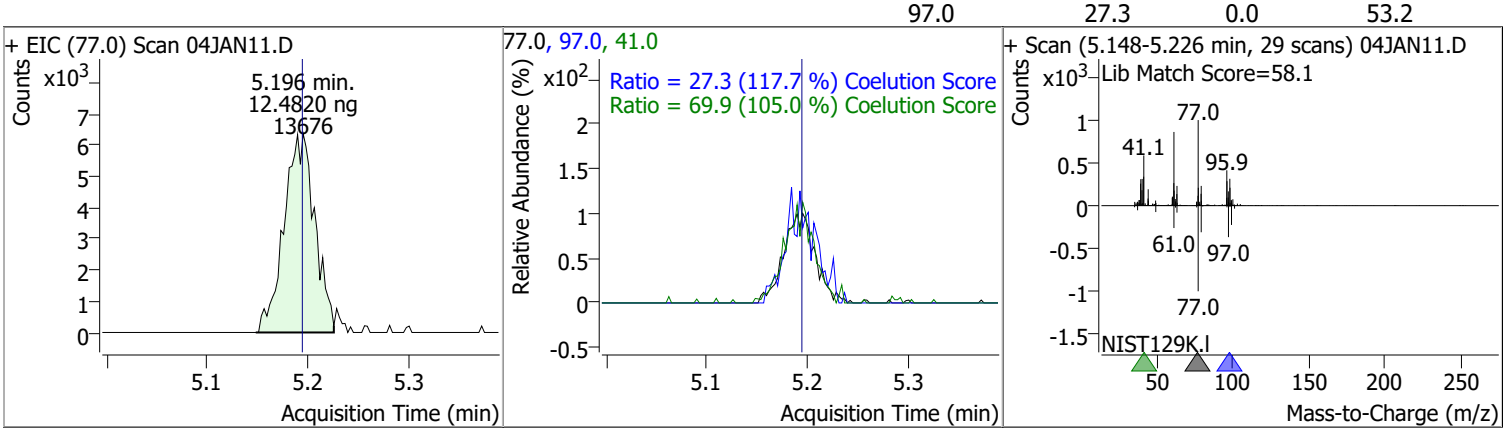
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	12.0652	4.38	0.00	17642	65.0	35.8	2.1	62.1
					83.0	15.3	0.0	43.7



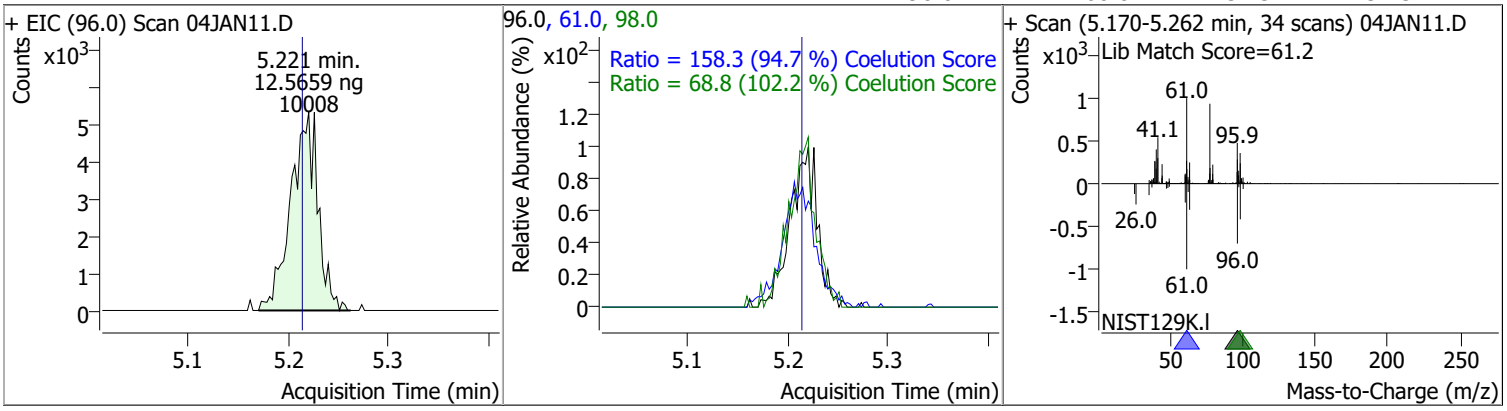




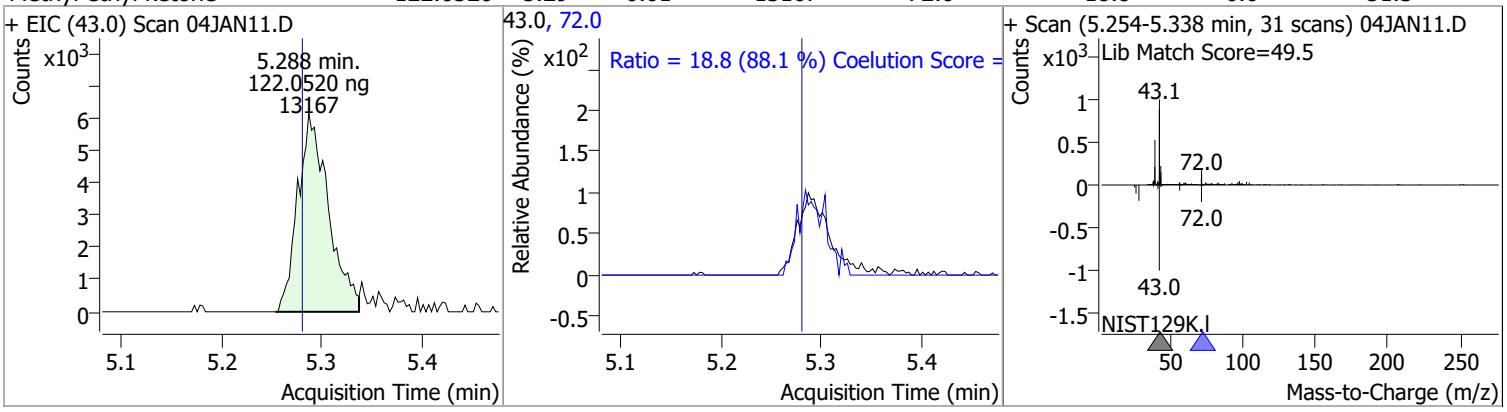
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	12.4820	5.20	0.00	13676	41.0	69.9	36.5	96.5
					97.0	27.3	0.0	53.2



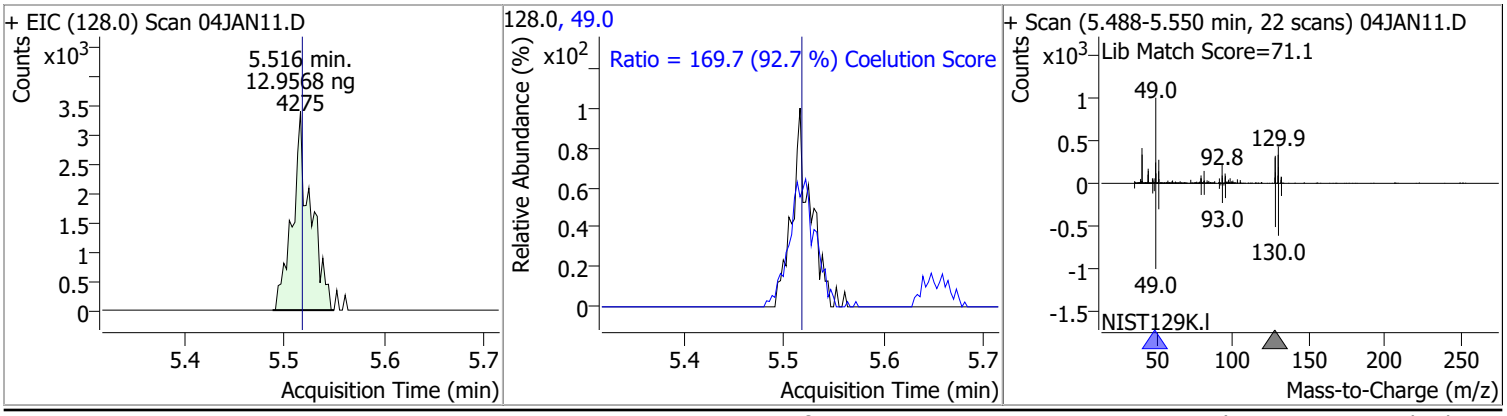
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	12.5659	5.22	0.01	10008	61.0	158.3	137.2	197.2
					98.0	68.8	37.3	97.3



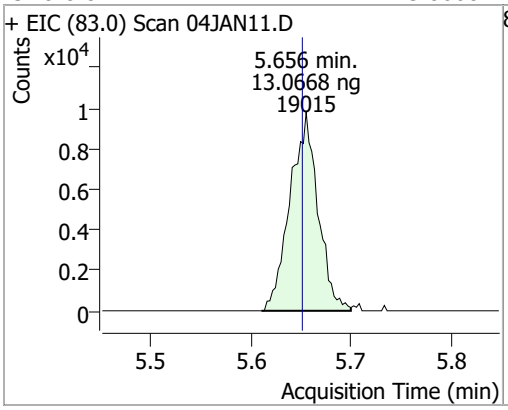
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	122.0520	5.29	0.01	13167	72.0	18.8	0.0	51.3

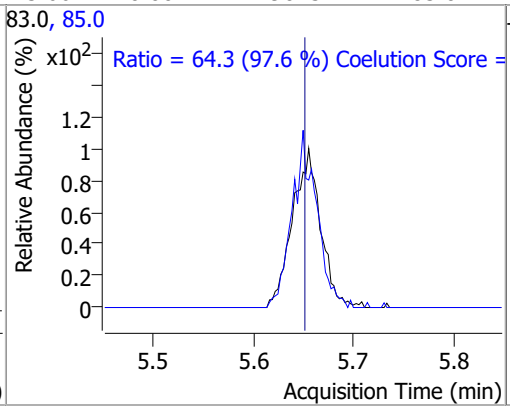


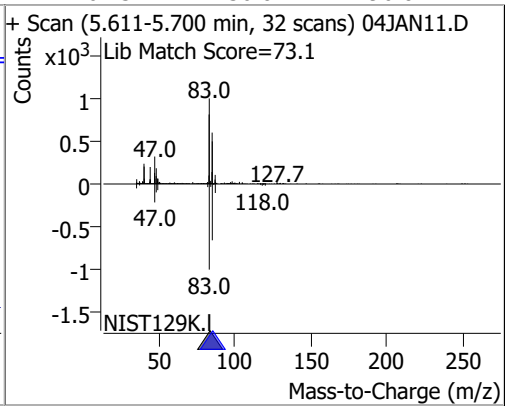
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	12.9568	5.52	0.00	4275	49.0	169.7	152.9	212.9



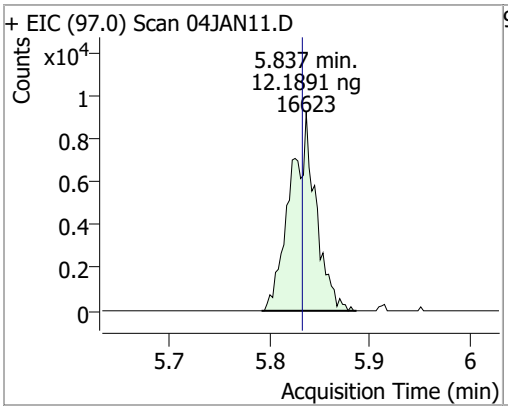
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	13.0668	5.66	0.00	19015	85.0	64.3	36.0	96.0

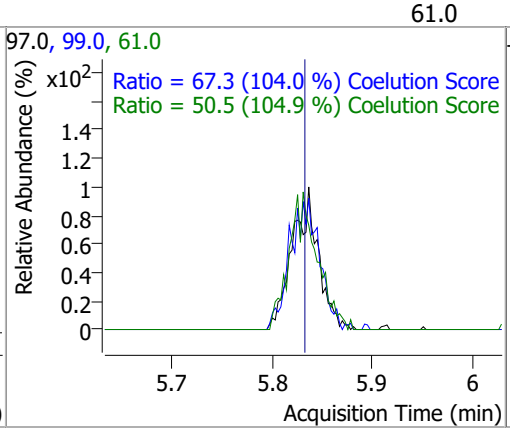


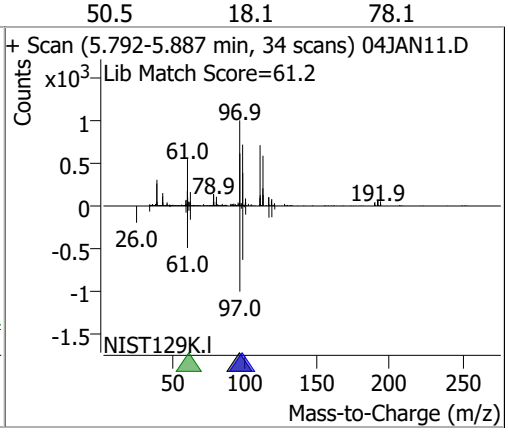




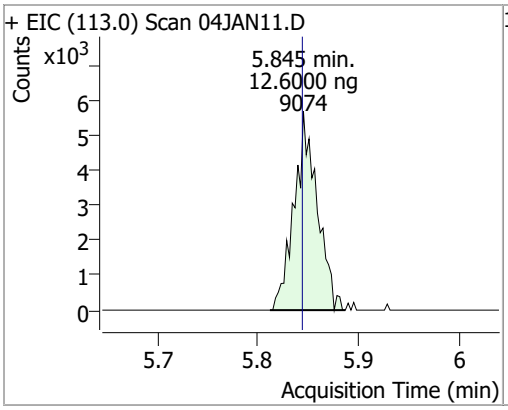
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	12.1891	5.84	0.00	16623	99.0	67.3	34.7	94.7
					61.0	50.5	18.1	78.1

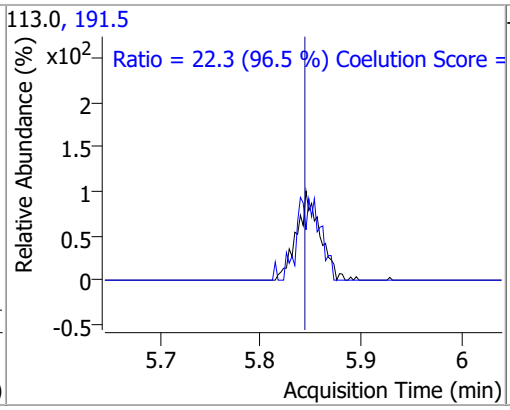


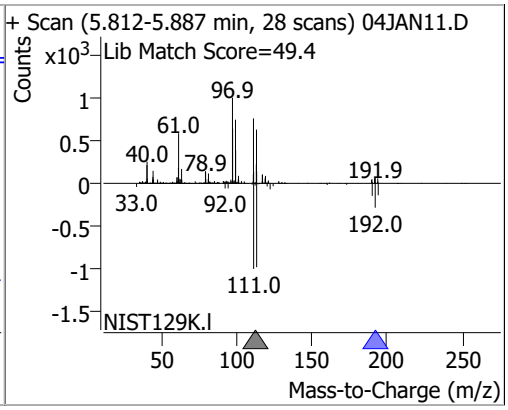




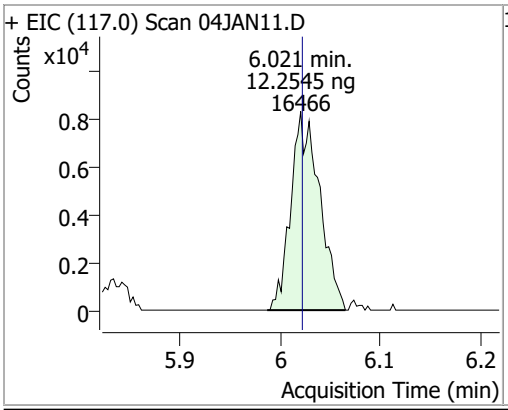
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	12.6000	5.85	0.00	9074	191.5	22.3	0.0	53.1

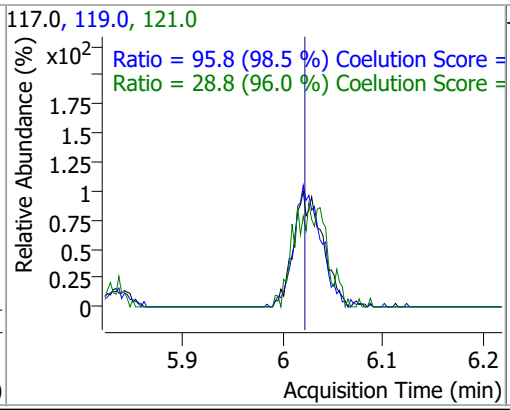


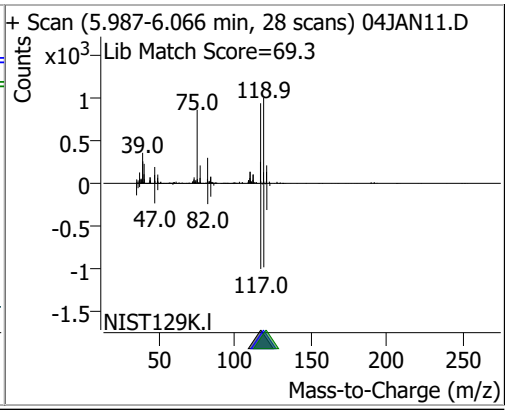




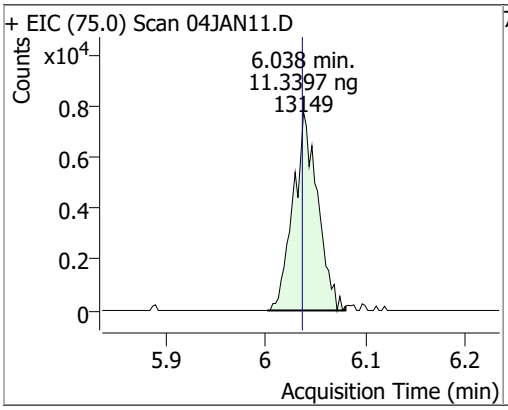
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	12.2545	6.02	0.00	16466	119.0	95.8	67.2	127.2
					121.0	28.8	0.1	60.1

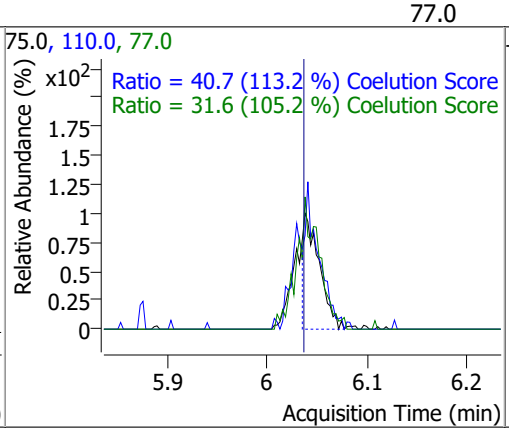


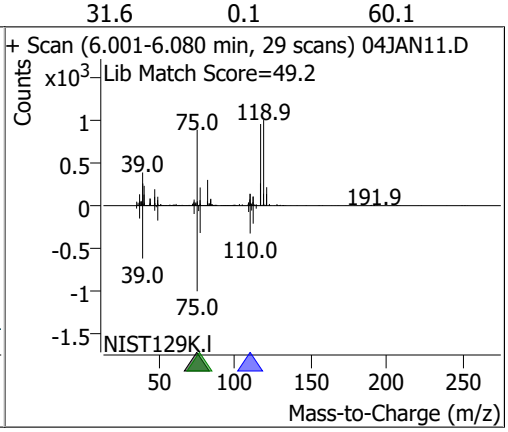




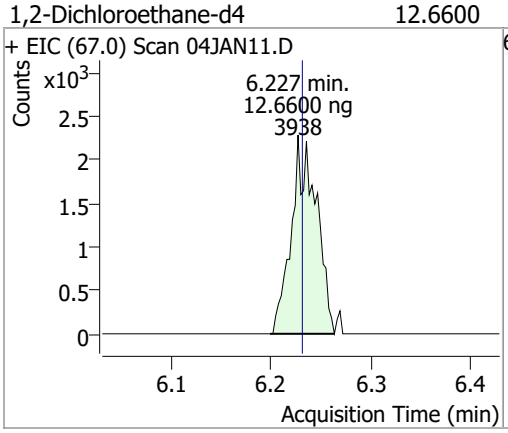
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	11.3397	6.04	0.00	13149	110.0	40.7	5.9	65.9
					77.0	31.6	0.1	60.1

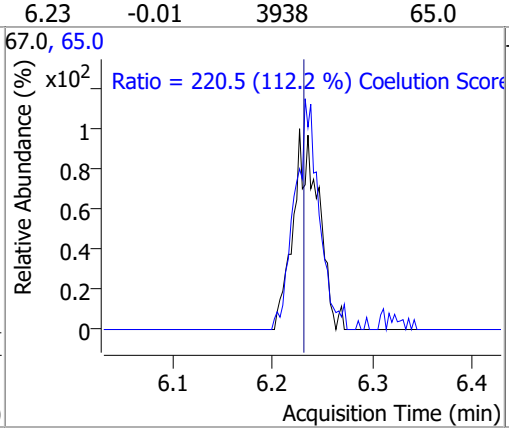


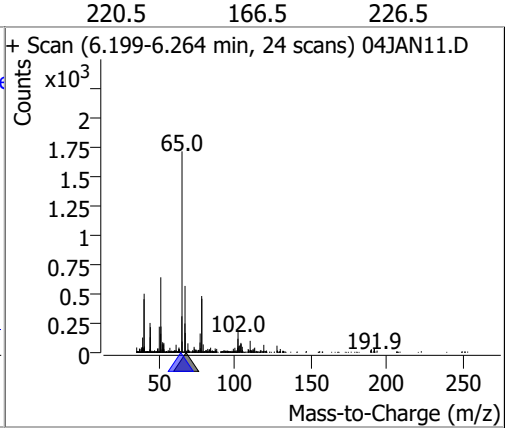




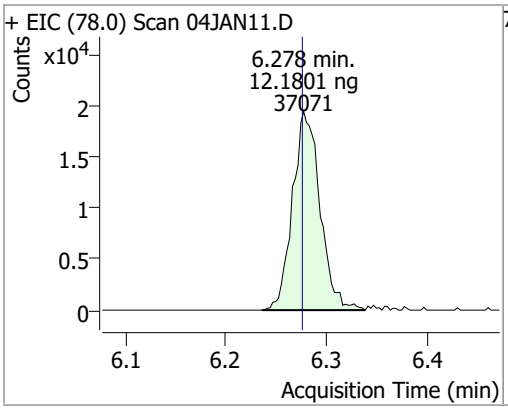
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	12.6600	6.23	-0.01	3938	65.0	220.5	166.5	226.5

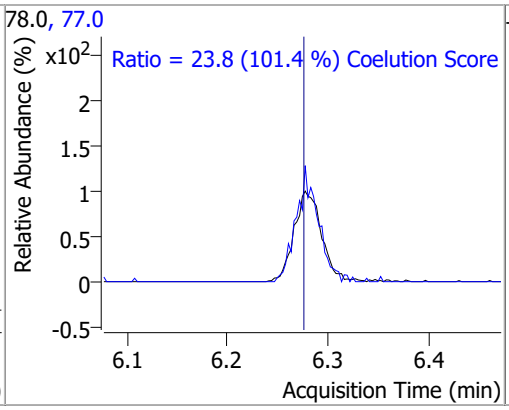


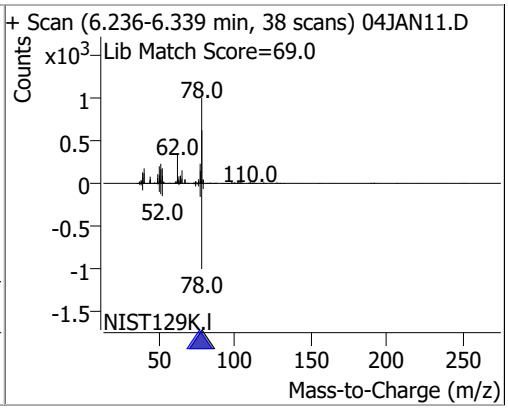




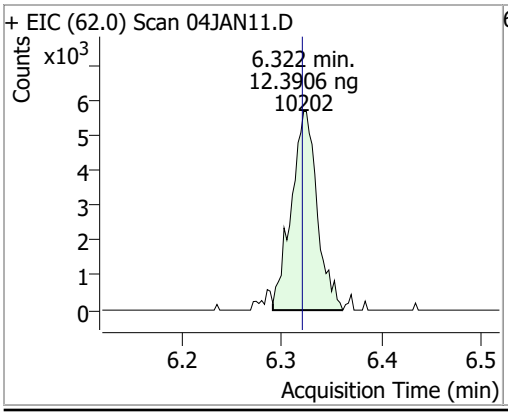
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	12.1801	6.28	0.00	37071	77.0	23.8	0.0	53.5

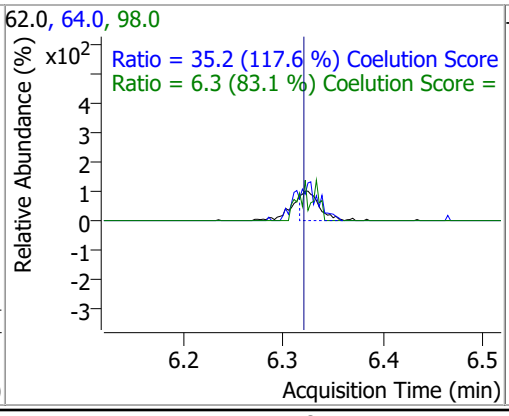


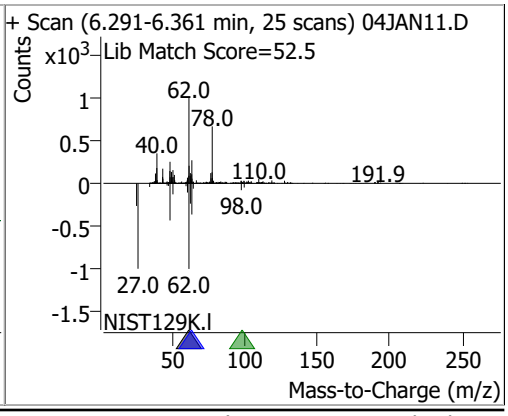




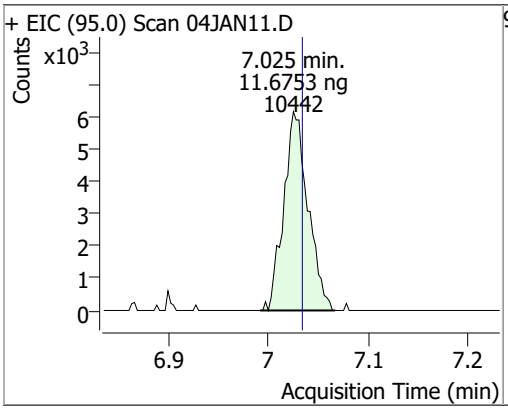
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	12.3906	6.32	0.00	10202	64.0	35.2	0.0	59.9
					98.0	6.3	0.0	37.6

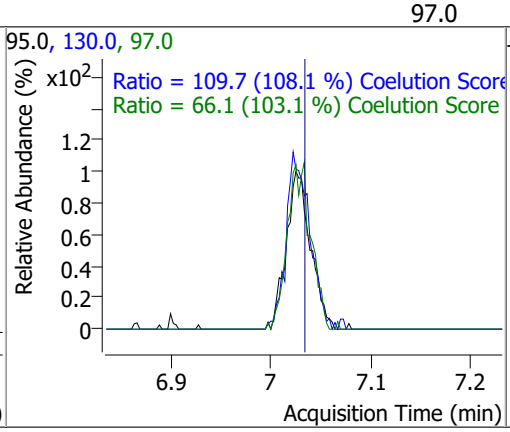


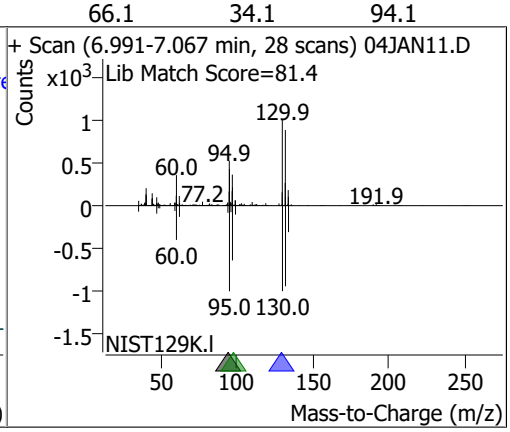




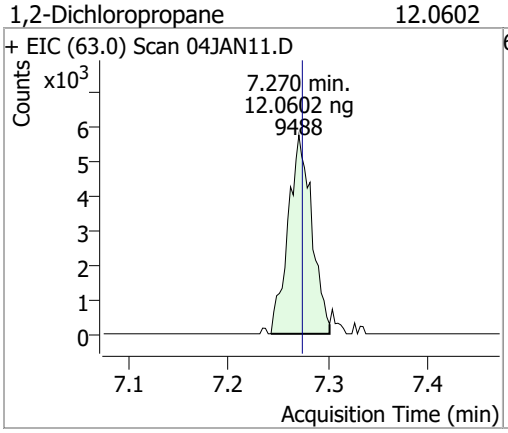
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	11.6753	7.02	-0.01	10442	130.0	109.7	71.5	131.5
					97.0	66.1	34.1	94.1

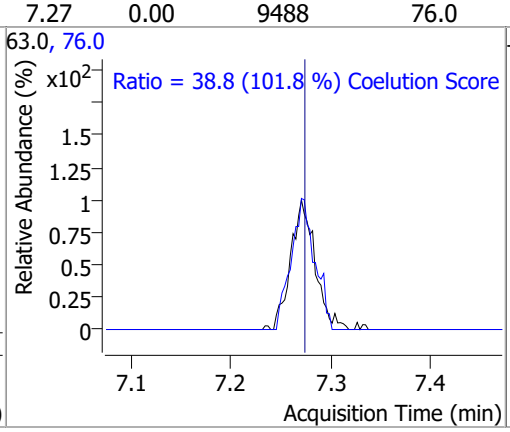


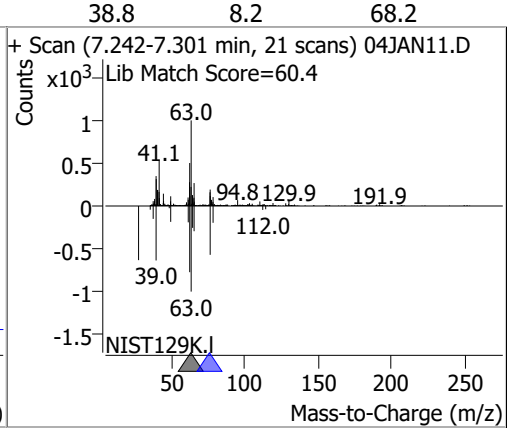




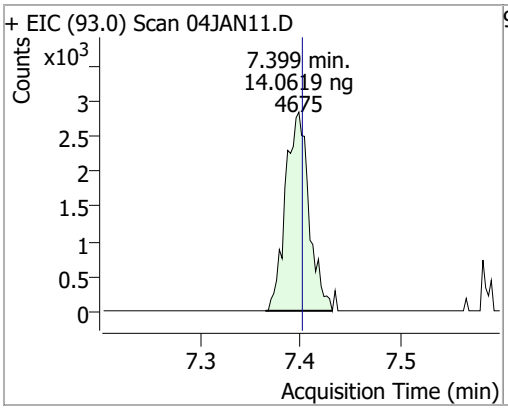
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	12.0602	7.27	0.00	9488	76.0	38.8	8.2	68.2

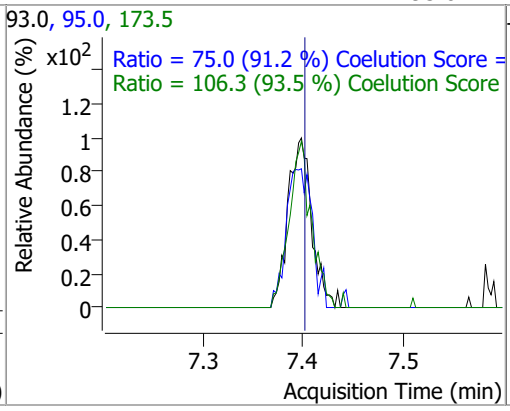


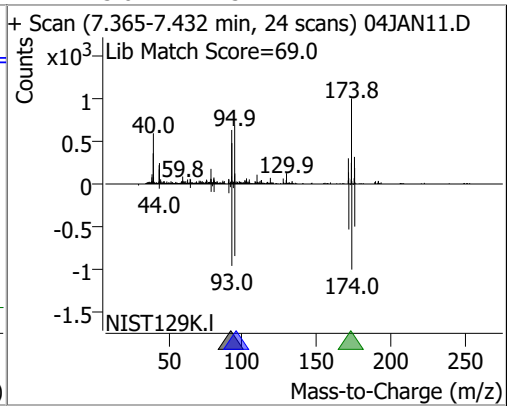




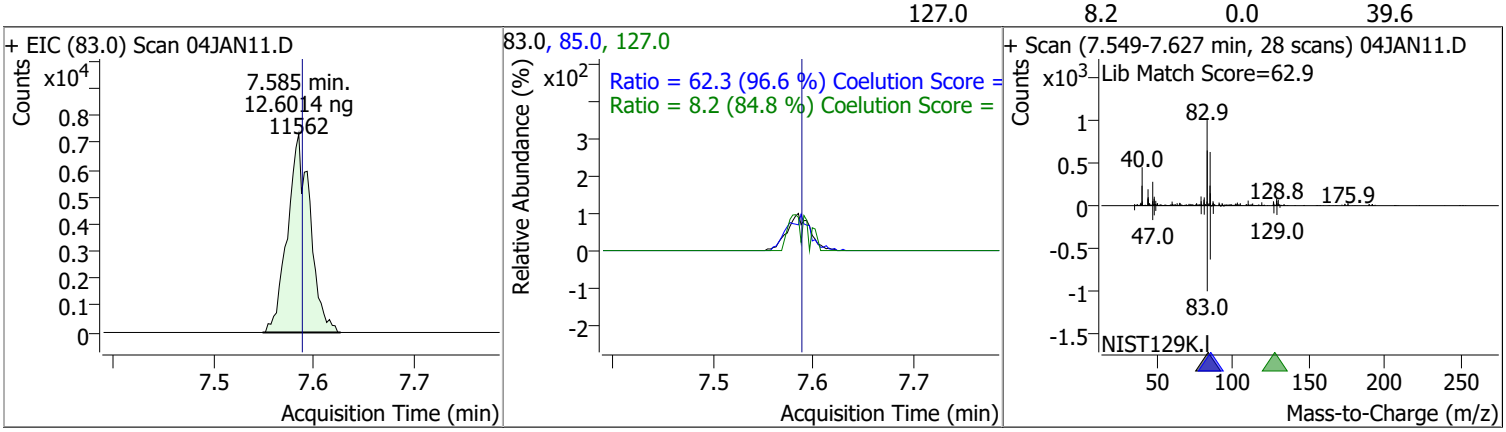
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	14.0619	7.40	0.00	4675	173.5	106.3	83.7	143.7
					95.0	75.0	52.2	112.2



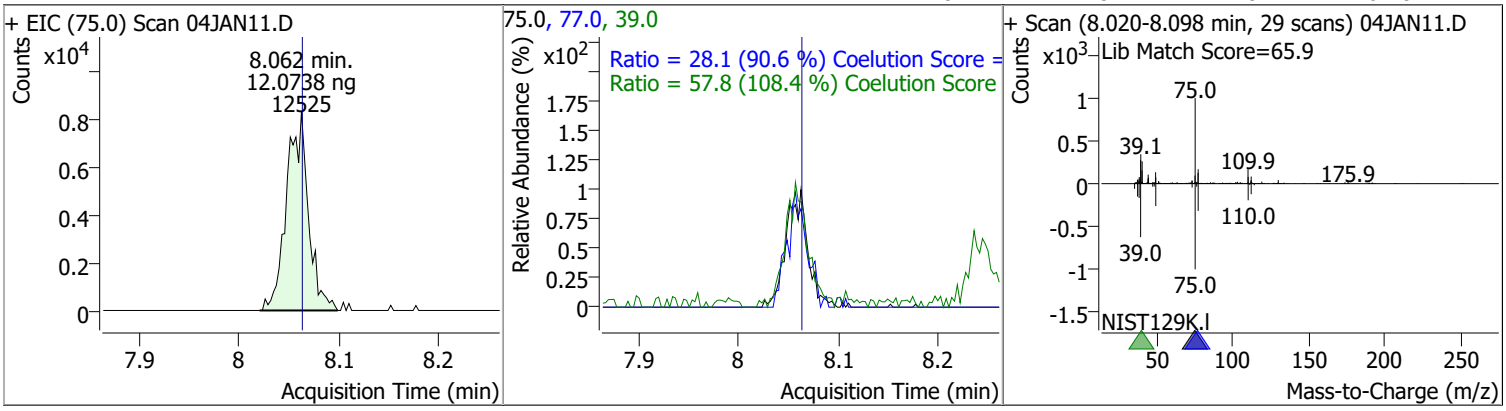




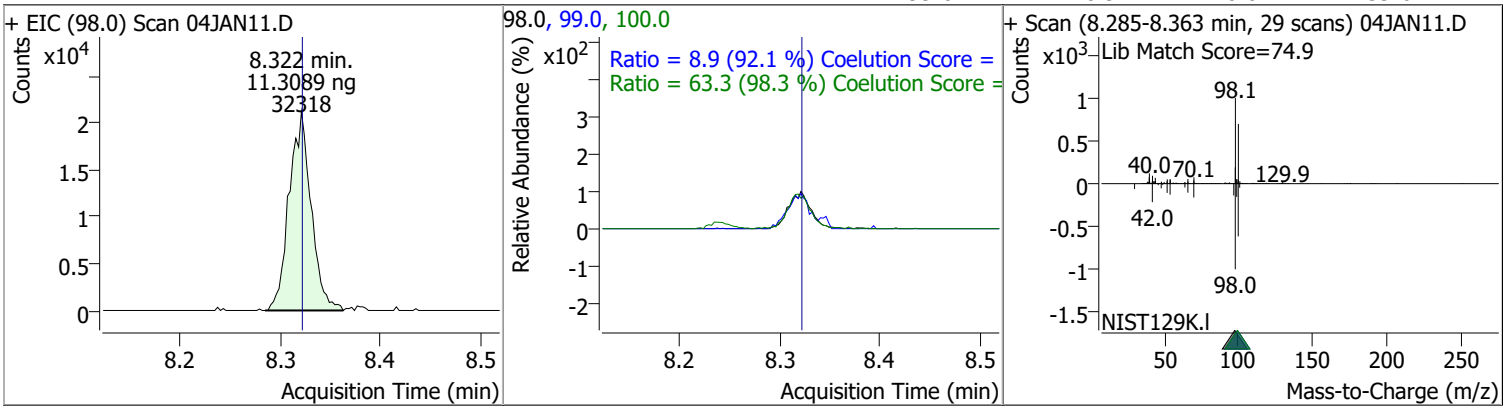
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	12.6014	7.59	0.00	11562	85.0	62.3	34.5	94.5
					127.0	8.2	0.0	39.6



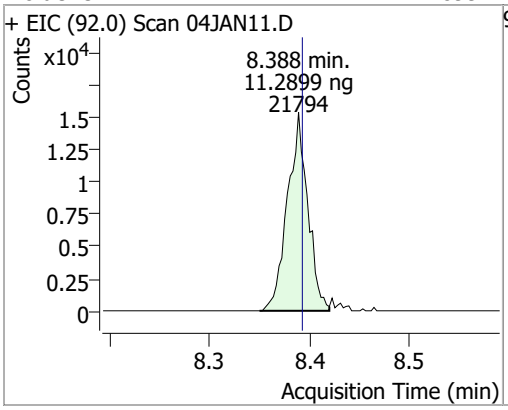
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	12.0738	8.06	0.00	12525	39.0	57.8	23.3	83.3
					77.0	28.1	1.0	61.0

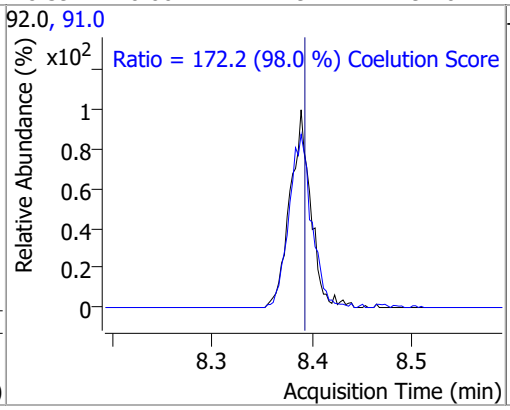


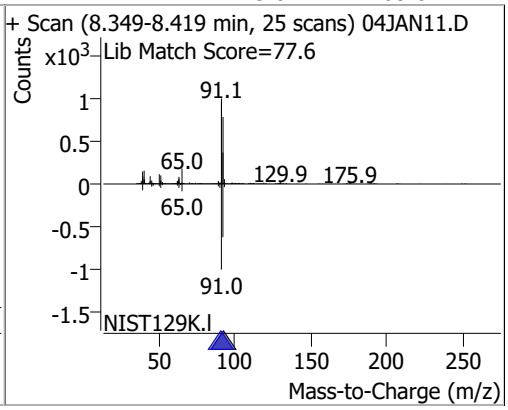
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	11.3089	8.32	0.00	32318	100.0	63.3	34.4	94.4
					99.0	8.9	0.0	39.6



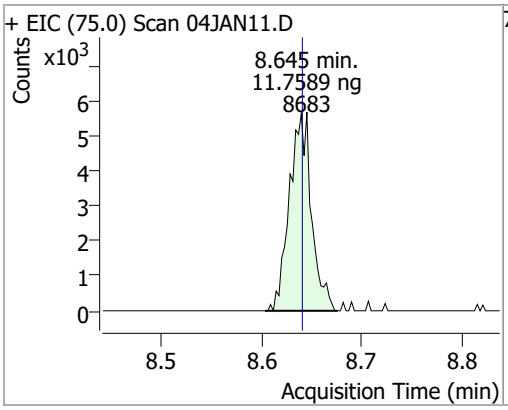
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	11.2899	8.39	0.00	21794	91.0	172.2	145.8	205.8

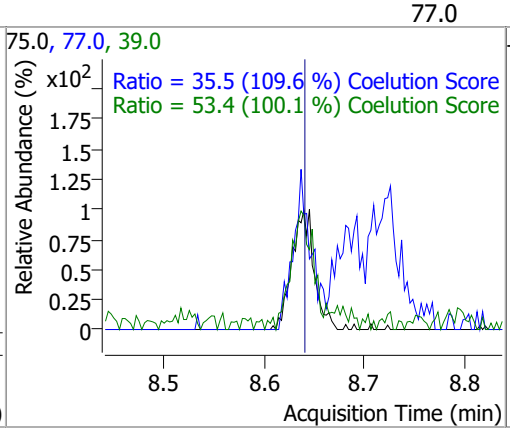


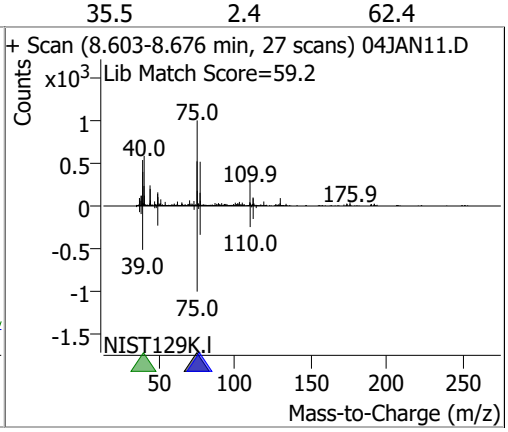




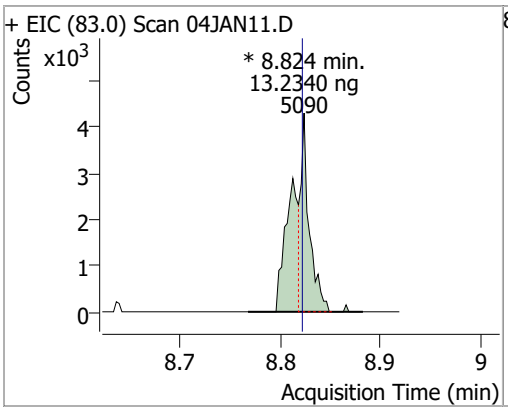
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	11.7589	8.65	0.01	8683	39.0	53.4	23.4	83.4
					77.0	35.5	2.4	62.4

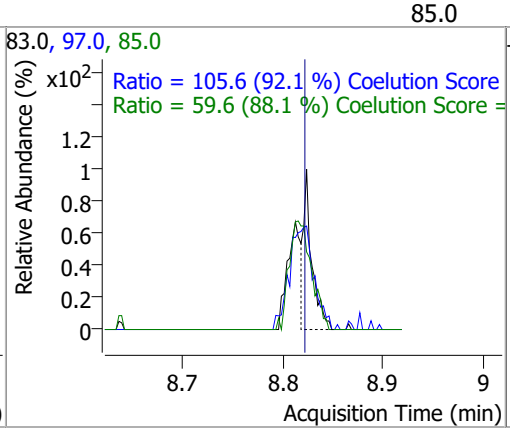


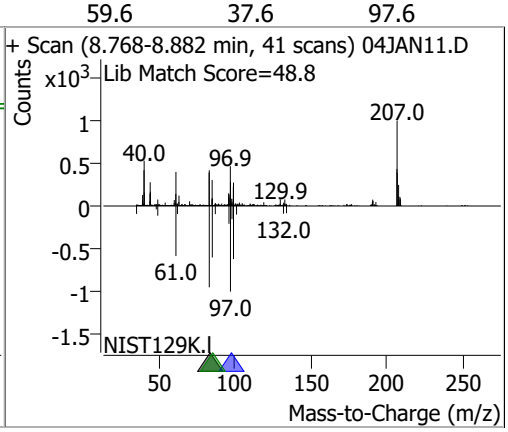




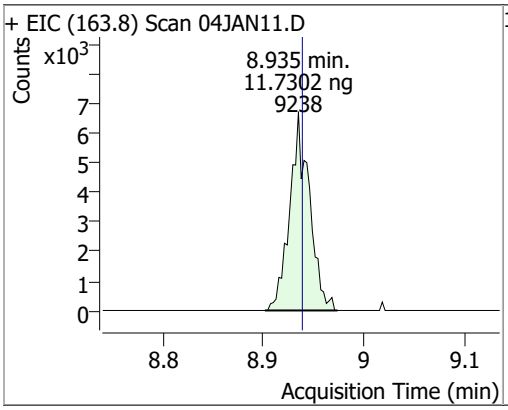
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	13.2340	8.82	0.01	5090 (m)	97.0	105.6	84.6	144.6
					85.0	59.6	37.6	97.6

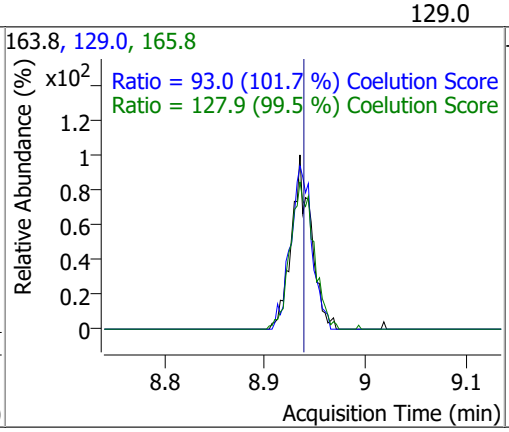


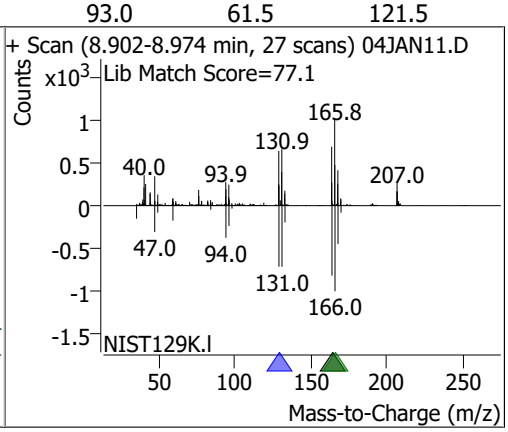




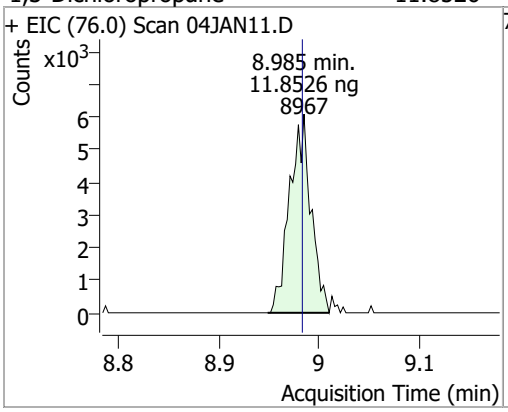
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	11.7302	8.94	0.00	9238	165.8	127.9	98.6	158.6
					129.0	93.0	61.5	121.5

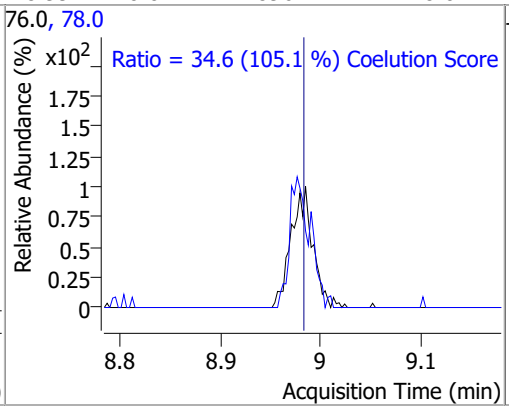


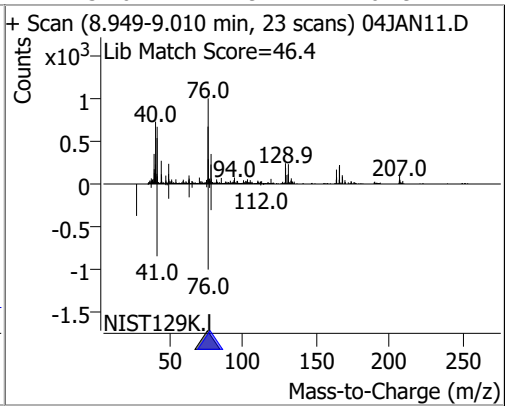




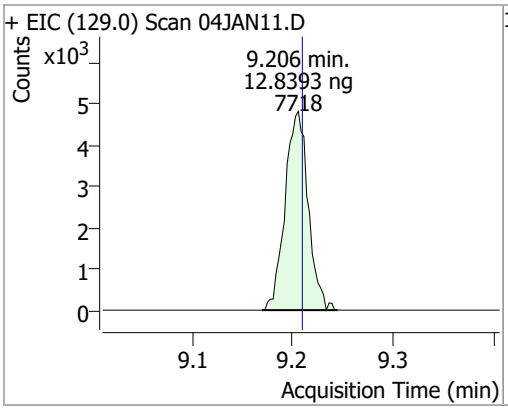
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	11.8526	8.99	0.01	8967	78.0	34.6	2.9	62.9

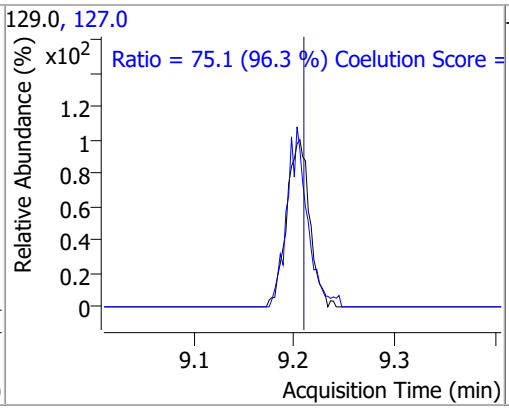


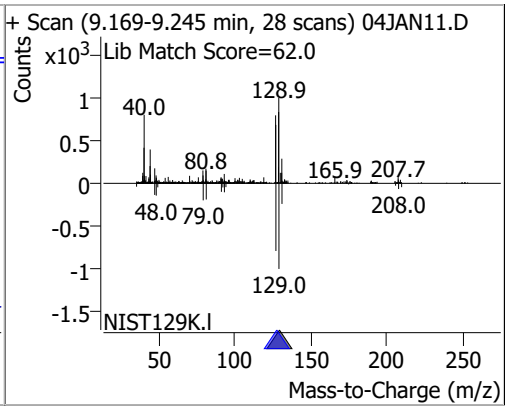




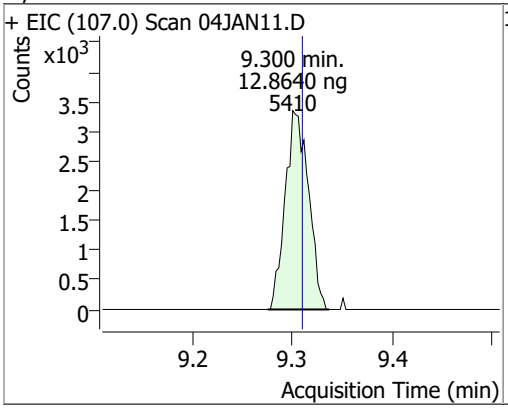
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	12.8393	9.21	0.00	7718	127.0	75.1	48.0	108.0

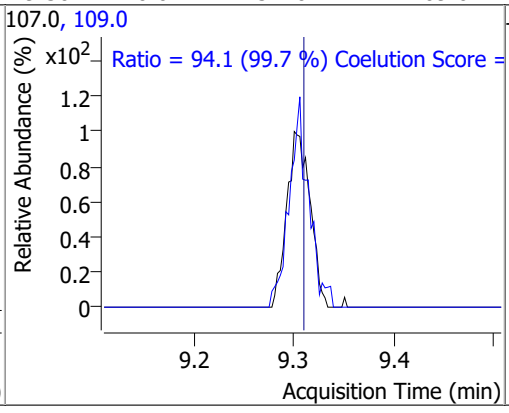


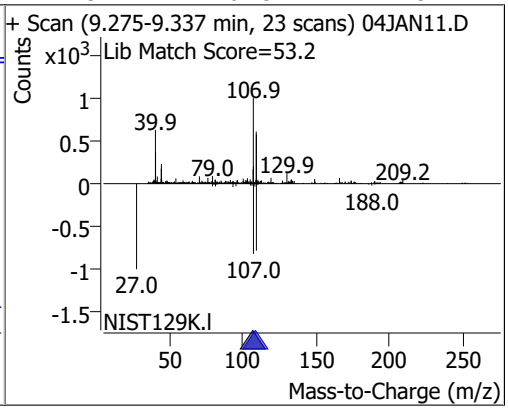




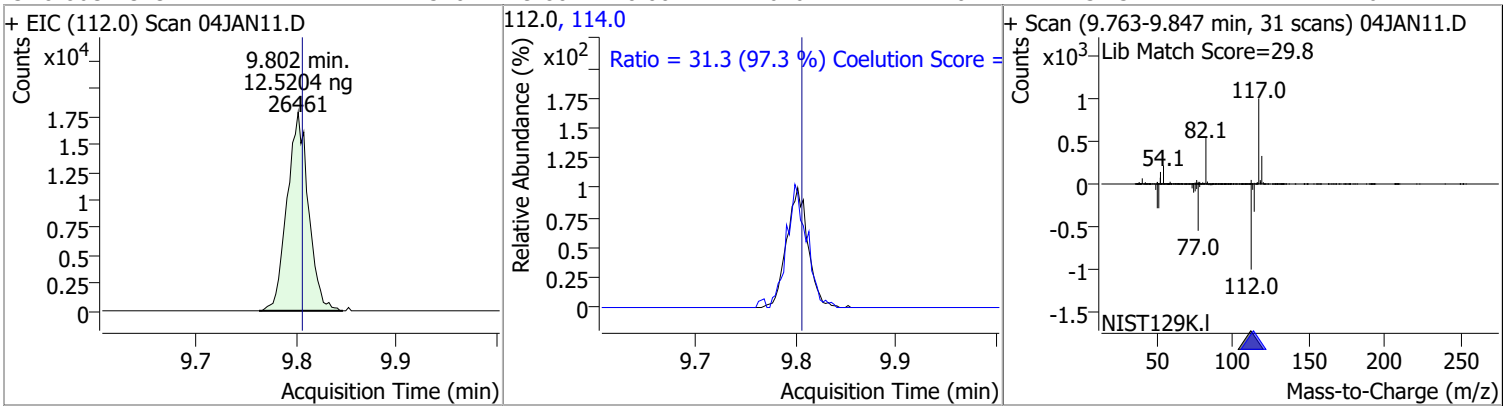
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	12.8640	9.30	-0.01	5410	109.0	94.1	64.5	124.5



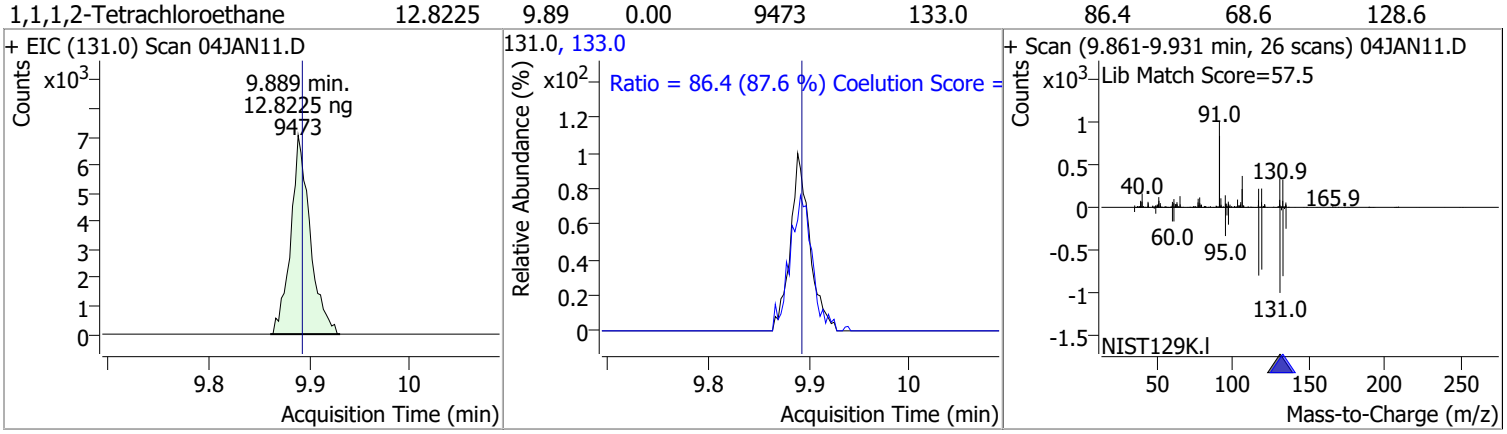




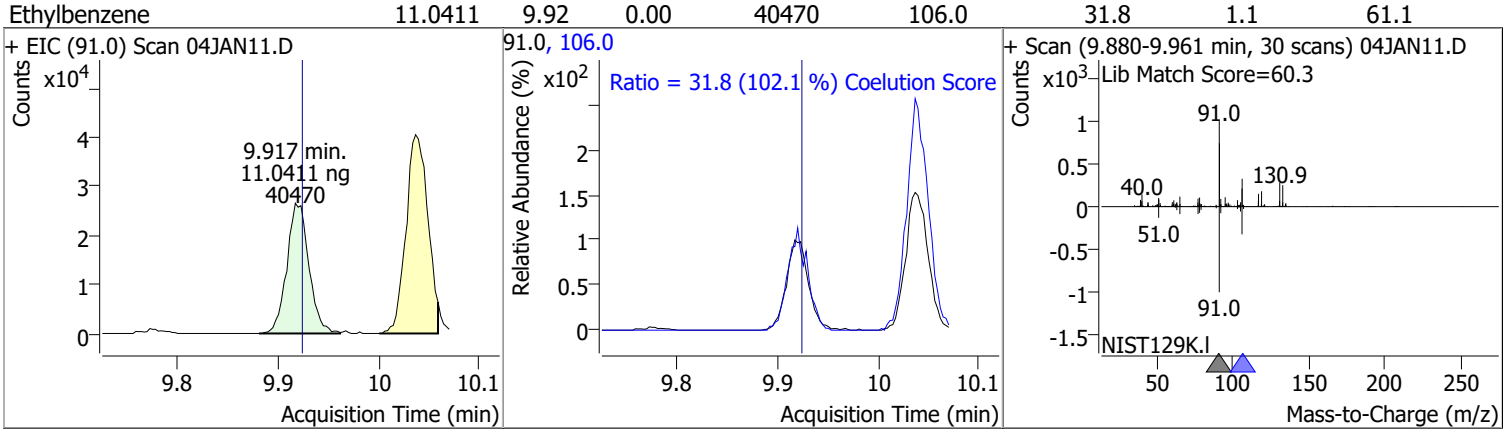
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	12.5204	9.80	0.00	26461	114.0	31.3	2.1	62.1



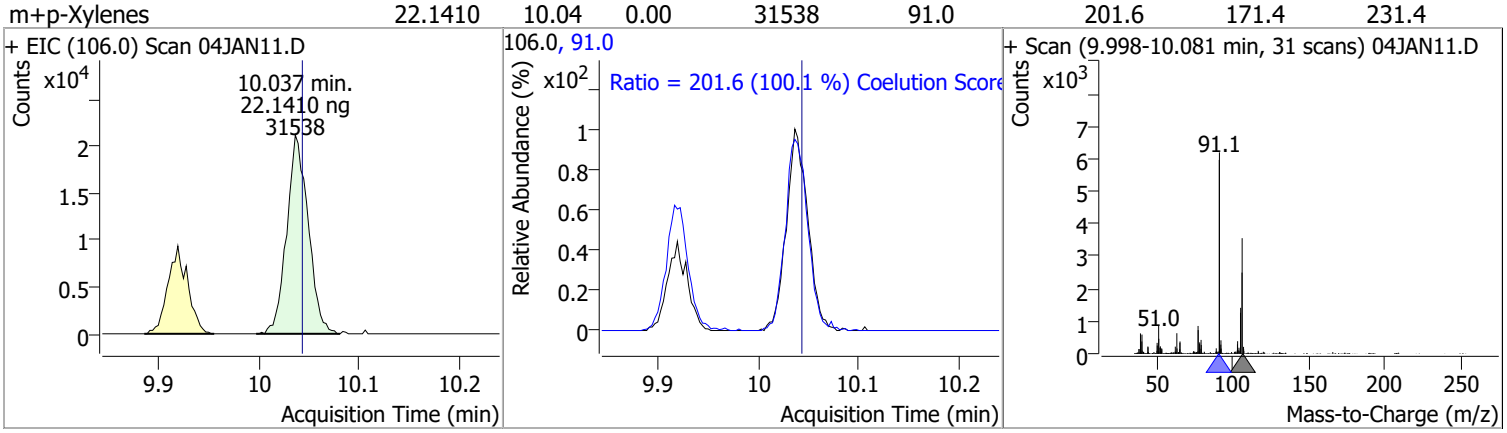
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	12.8225	9.89	0.00	9473	133.0	86.4	68.6	128.6



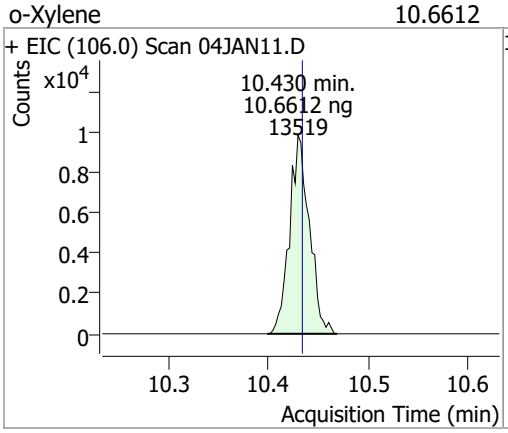
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	11.0411	9.92	0.00	40470	106.0	31.8	1.1	61.1

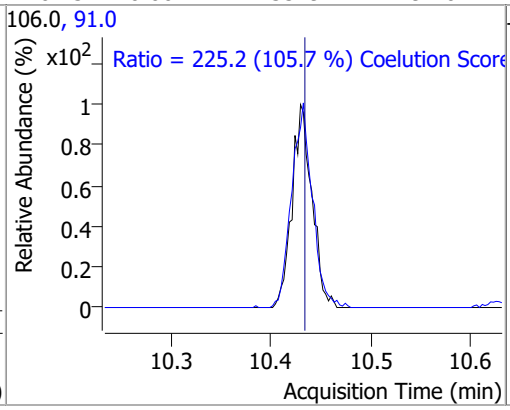


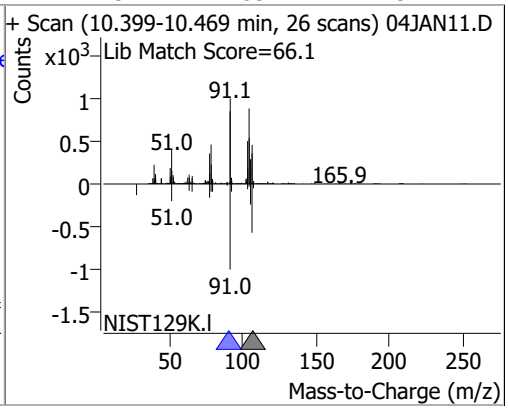
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	22.1410	10.04	0.00	31538	91.0	201.6	171.4	231.4



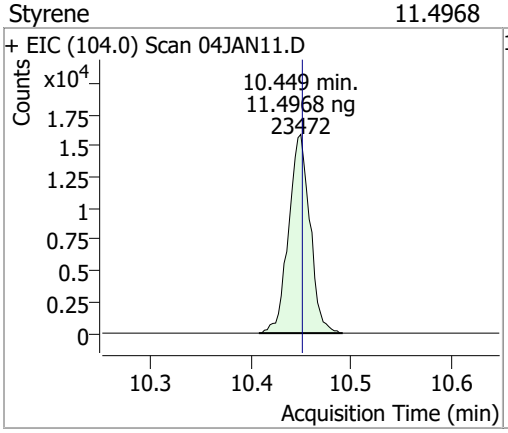
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	10.6612	10.43	0.00	13519	91.0	225.2	183.1	243.1

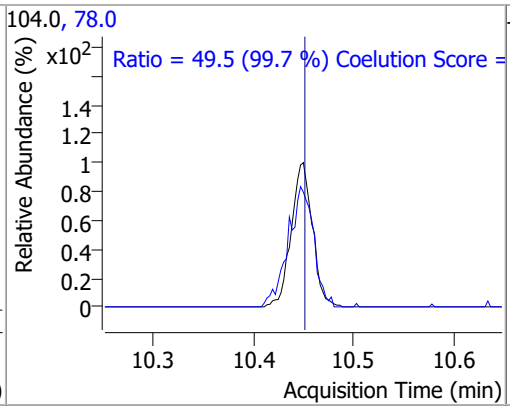


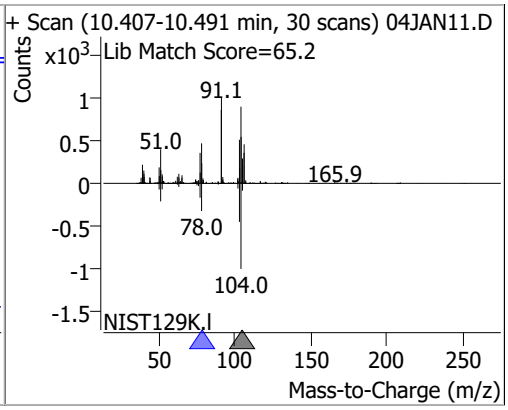




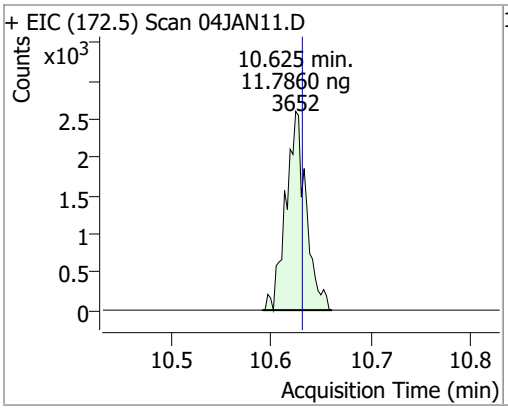
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	11.4968	10.45	0.00	23472	78.0	49.5	19.6	79.6

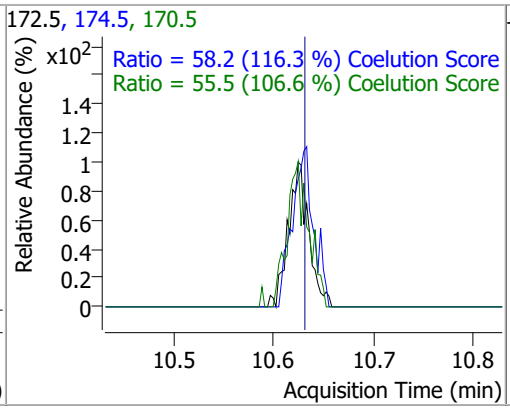


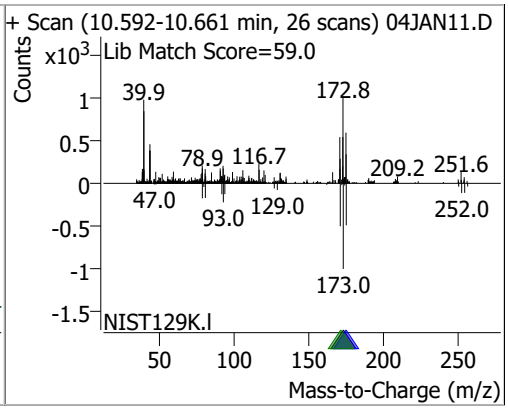




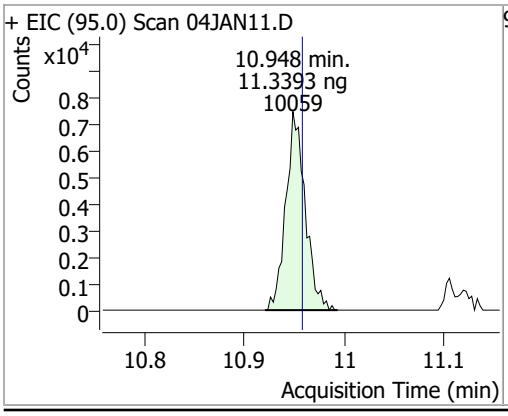
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	11.7860	10.62	0.00	3652	170.5	55.5	22.1	82.1
					174.5	58.2	20.1	80.1

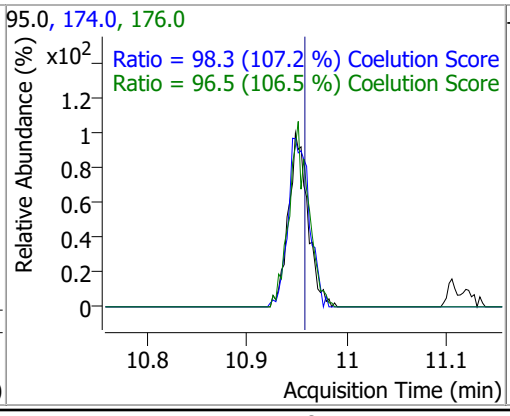


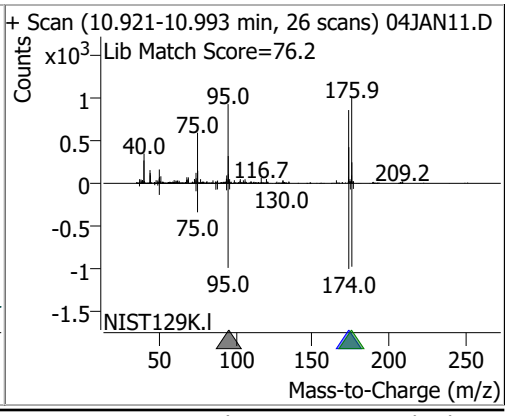




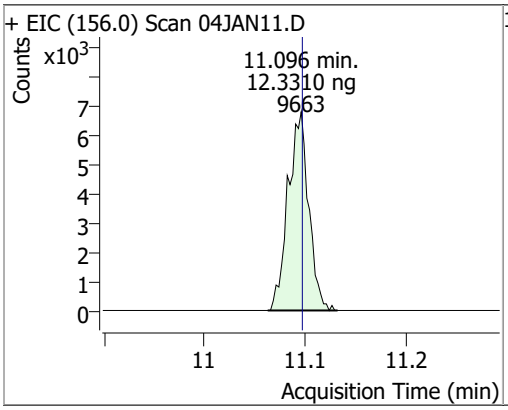
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	11.3393	10.95	-0.01	10059	174.0	98.3	61.7	121.7
					176.0	96.5	60.6	120.6

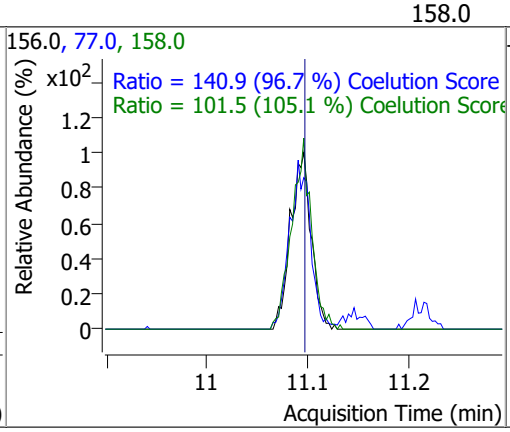


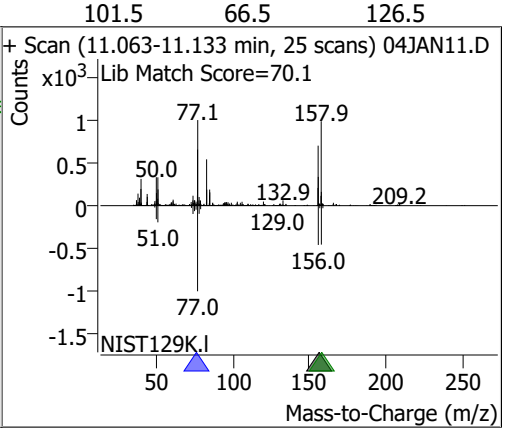




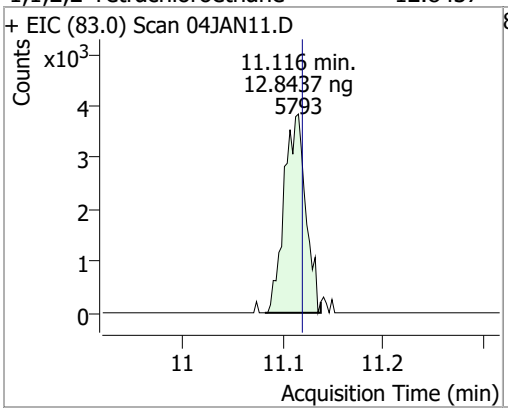
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	12.3310	11.10	0.00	9663	77.0	140.9	115.7	175.7
					158.0	101.5	66.5	126.5

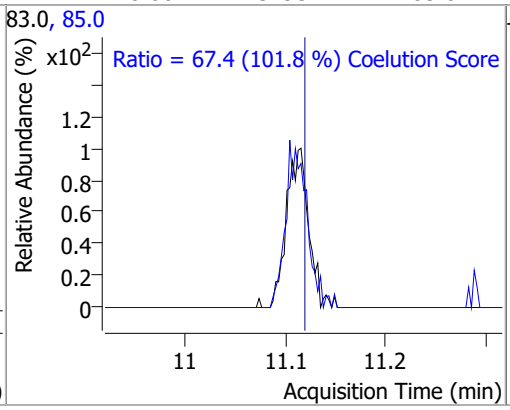


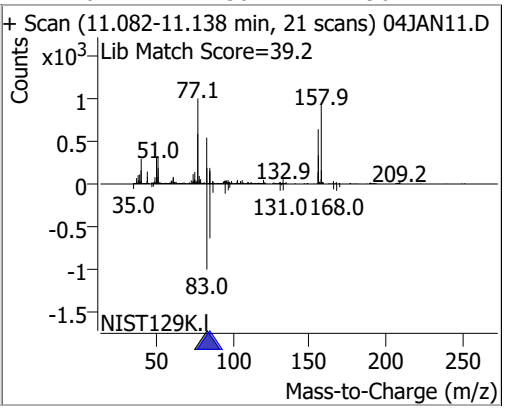




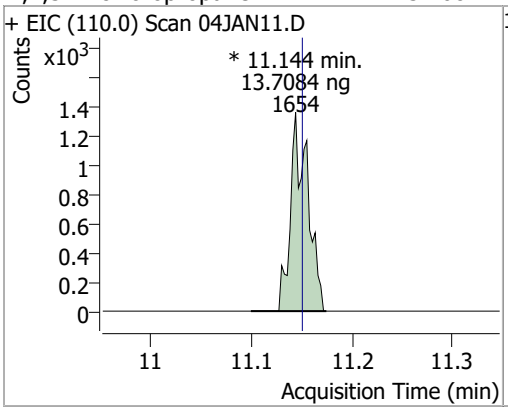
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	12.8437	11.12	0.00	5793	85.0	67.4	36.2	96.2

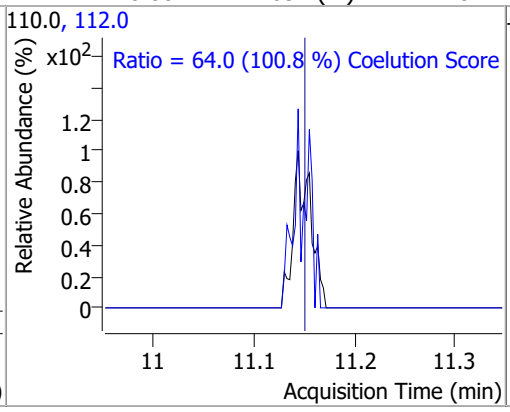


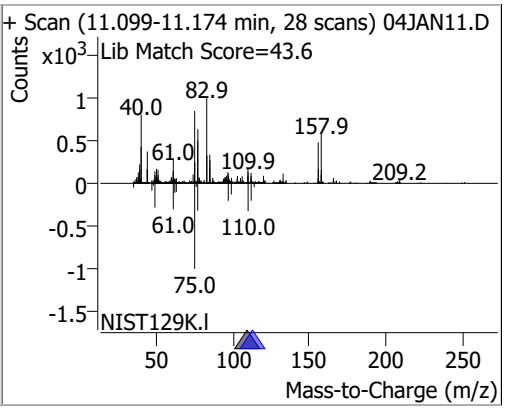




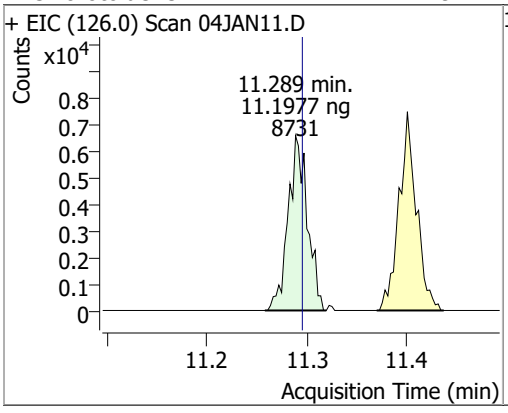
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	13.7084	11.14	0.00	1654 (m)	112.0	64.0	33.5	93.5

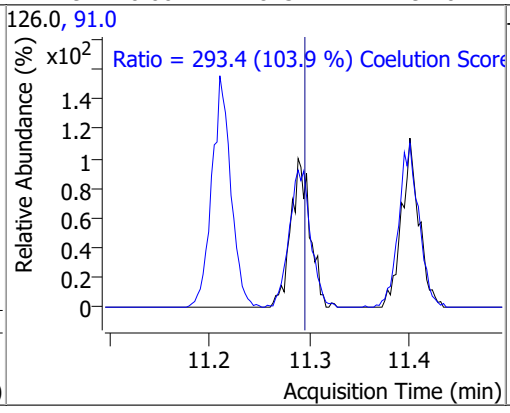


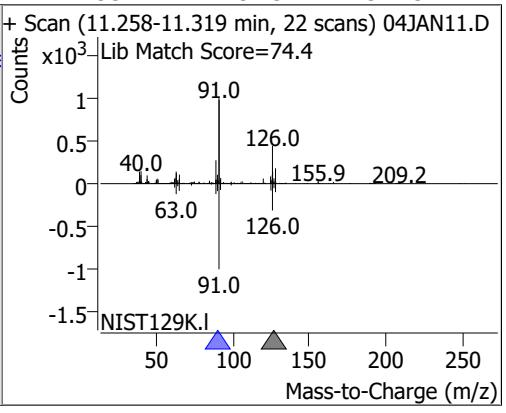




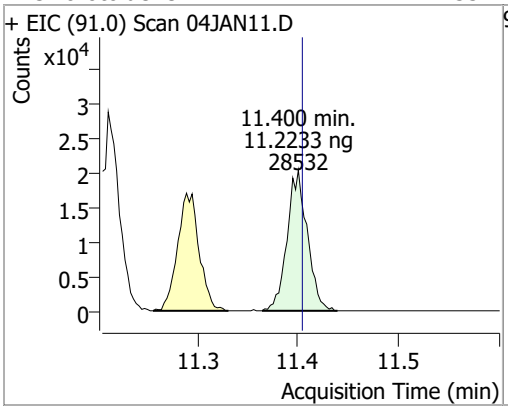
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	11.1977	11.29	0.00	8731	91.0	293.4	252.3	312.3

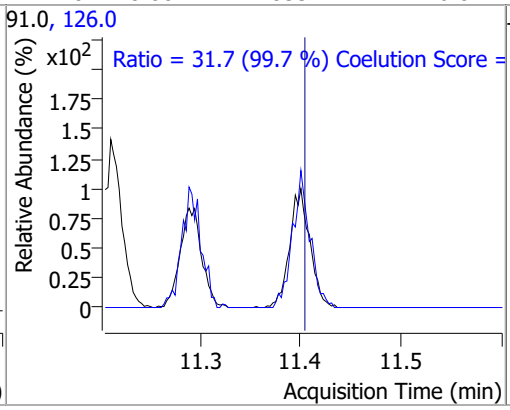


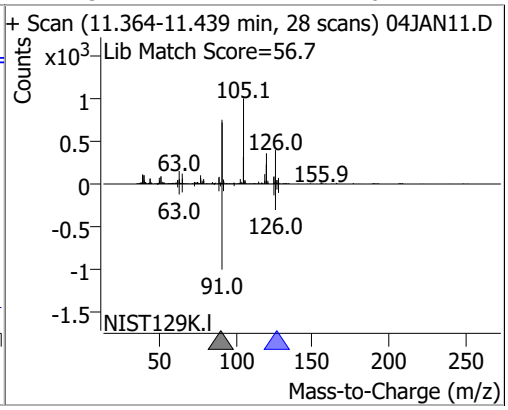




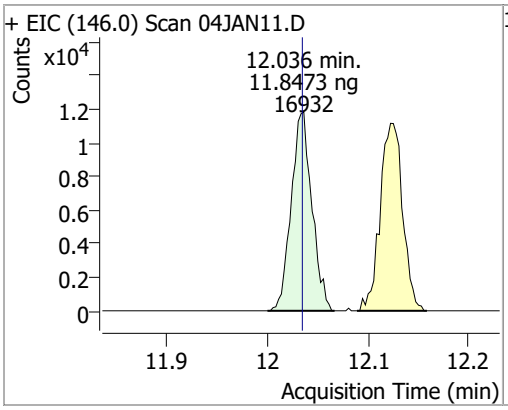
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	11.2233	11.40	0.00	28532	126.0	31.7	1.7	61.7

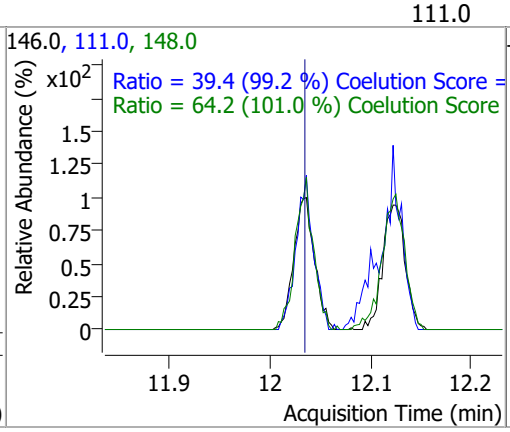


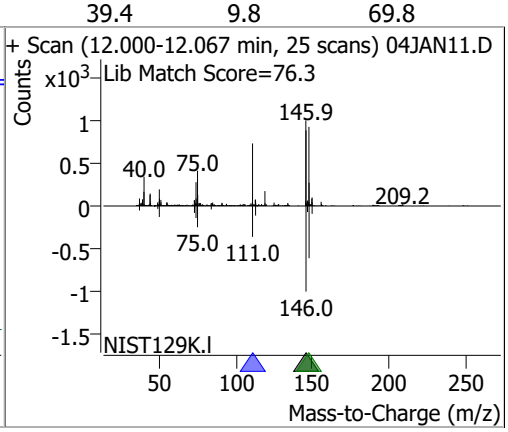




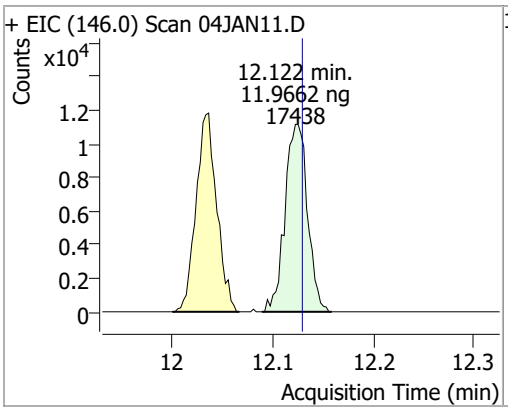
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	11.8473	12.04	0.01	16932	148.0	64.2	33.6	93.6
					111.0	39.4	9.8	69.8

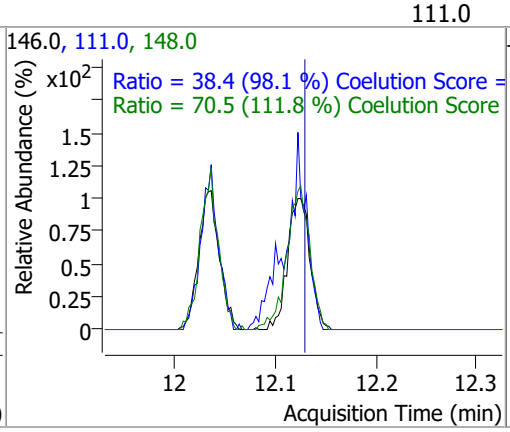


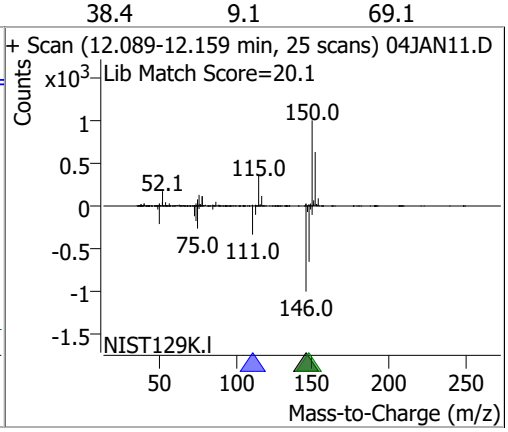




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	11.9662	12.12	0.00	17438	148.0	70.5	33.1	93.1
					111.0	38.4	9.1	69.1

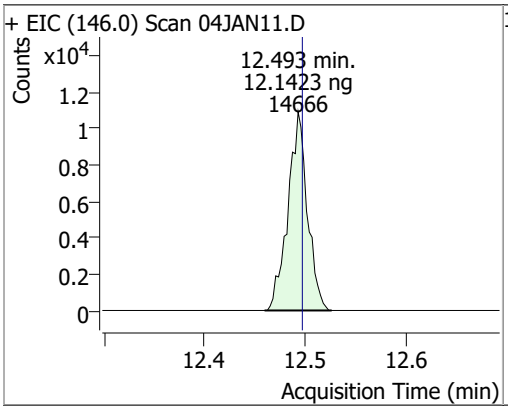


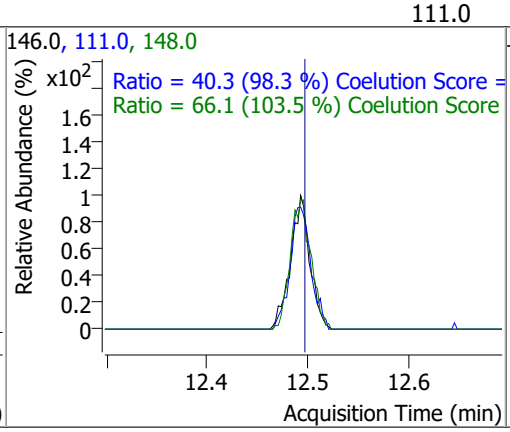


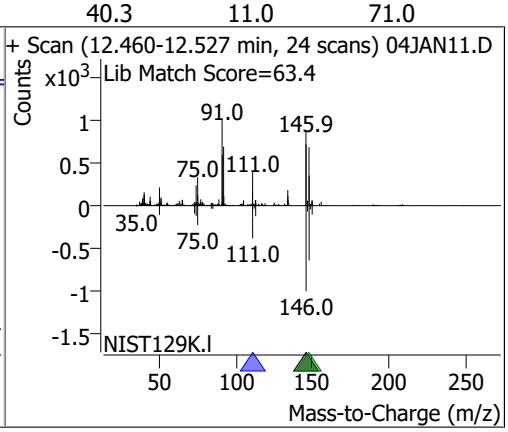


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	12.1423	12.49	0.00	14666	148.0	66.1	33.9	93.9
					111.0	40.3	11.0	71.0



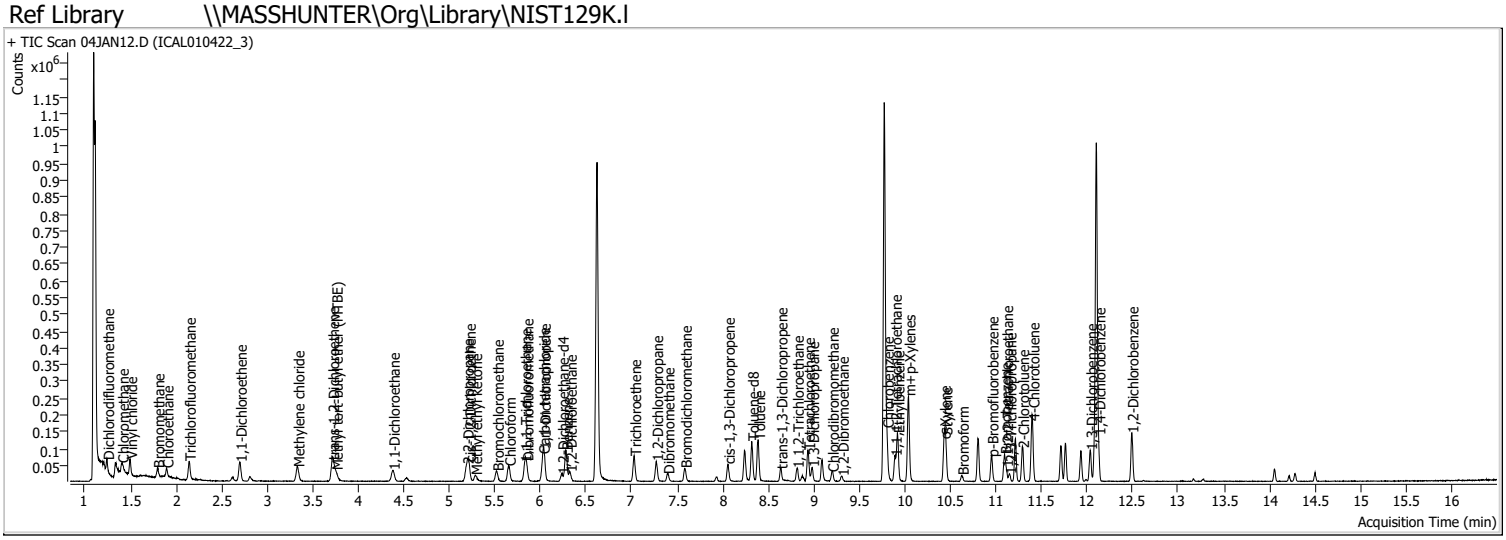




Quantitation Results Report (QT Reviewed)

Agilent | Trusted Answers

Data File	04JAN12.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/4/2022 4:28:05 PM
Sample Name	ICAL010422_3	Instrument	VOA5975C
Vial	12	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

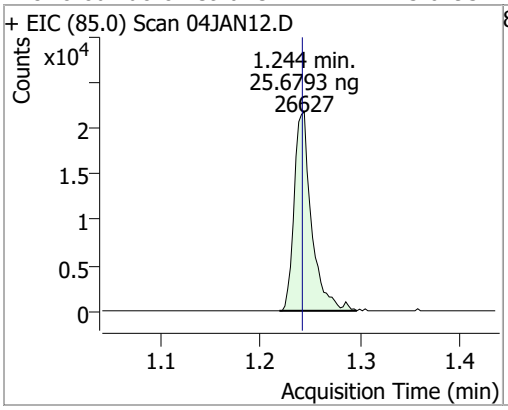


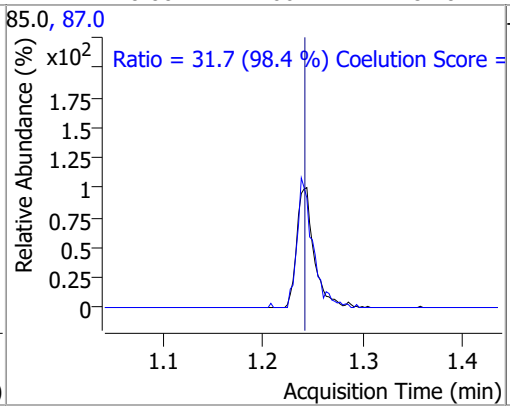
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	791270	250.0000	ng		-0.003
M Chlorobenzene-d5	9.775	82.0	301338	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.103	152.0	240335	250.0000	ng		0.003
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	19100	25.6219	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 10.25%		*	
S 1,2-Dichloroethane-d4	6.236	67.0	8284	25.7280	ng		0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 10.29%		*	
S Toluene-d8	8.319	98.0	67673	23.3046	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 9.32%		*	
S p-Bromofluorobenzene	10.951	95.0	22267	25.2899	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 10.12%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.244	85.0	26627	25.6793	ng		99
T Chloromethane	1.406	50.0	33153	26.3422	ng		95
T Vinyl chloride	1.495	62.0	29046	25.6488	ng		80
T Bromomethane	1.796	96.0	13054	25.7793	ng		99
T Chloroethane	1.897	64.0	14646	26.1250	ng	m	92
T Trichlorofluoromethane	2.142	101.0	37464	26.6531	ng		98
T 1,1-Dichloroethene	2.700	96.0	20631	25.8849	ng		93
T Methylene chloride	3.333	49.0	30908	26.3058	ng		100
T trans-1,2-Dichloroethene	3.712	96.0	20706	25.4641	ng		97
T Methyl tert-butyl ether (MTBE)	3.754	73.0	24218	23.0418	ng		92
T 1,1-Dichloroethane	4.379	63.0	38874	25.6835	ng		97
T 2,2-Dichloropropane	5.190	77.0	29793	26.2692	ng		100
T cis-1,2-Dichloroethene	5.212	96.0	20252	24.5653	ng		98
T Methyl ethyl ketone	5.282	43.0	26248	235.0504	ng		98
T Bromochloromethane	5.522	128.0	8688	25.4383	ng		98
T Chloroform	5.653	83.0	36413	24.1734	ng		97

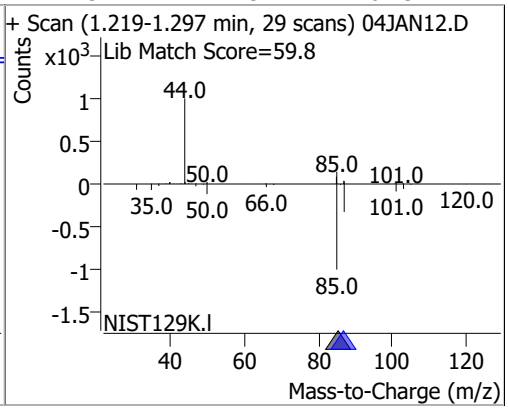
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	5.826	97.0	35547	25.1809	ng		96
T Carbon tetrachloride	6.024	117.0	34462	24.7773	ng		99
T 1,1-Dichloropropene	6.038	75.0	29241	24.3617	ng		96
T Benzene	6.278	78.0	74956	23.7919	ng		97
T 1,2-Dichloroethane	6.322	62.0	19996	23.4616	ng		97
T Trichloroethene	7.028	95.0	21946	24.1484	ng		98
T 1,2-Dichloropropane	7.270	63.0	20077	25.1147	ng		98
T Dibromomethane	7.393	93.0	8055	23.8439	ng		97
T Bromodichloromethane	7.583	83.0	22743	24.3940	ng		98
T cis-1,3-Dichloropropene	8.057	75.0	24511	23.2528	ng		97
T Toluene	8.389	92.0	46355	23.6319	ng		99
T trans-1,3-Dichloropropene	8.634	75.0	17850	23.7894	ng		97
T 1,1,2-Trichloroethane	8.815	83.0	10099	25.8400	ng		95
T Tetrachloroethene	8.935	163.8	20322	25.3948	ng		97
T 1,3-Dichloropropane	8.983	76.0	18745	24.3839	ng		99
T Chlorodibromomethane	9.197	129.0	14873	24.3492	ng		99
T 1,2-Dibromoethane	9.309	107.0	10410	24.3601	ng		95
T Chlorobenzene	9.802	112.0	53047	24.7015	ng		100
T 1,1,1,2-Tetrachloroethane	9.889	131.0	18130	24.1509	ng		100
T Ethylbenzene	9.917	91.0	88428	23.7421	ng		100
T m+p-Xylenes	10.039	106.0	66267	45.7836	ng		98
T o-Xylene	10.427	106.0	30463	23.6420	ng		98
T Styrene	10.447	104.0	48569	23.4119	ng		97
T Bromoform	10.625	172.5	7972	25.9212	ng		96
T Bromobenzene	11.094	156.0	19259	24.7613	ng		94
T 1,1,2,2-Tetrachloroethane	11.110	83.0	12440	27.7883	ng		93
T 1,2,3-Trichloropropane	11.149	110.0	3200	26.7144	ng	m	96
T 2-Chlorotoluene	11.286	126.0	19390	25.0550	ng		94
T 4-Chlorotoluene	11.400	91.0	61551	24.3936	ng		100
T 1,3-Dichlorobenzene	12.033	146.0	36559	25.7725	ng		97
T 1,4-Dichlorobenzene	12.125	146.0	36635	25.3284	ng		92
T 1,2-Dichlorobenzene	12.488	146.0	29899	24.9402	ng		98

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

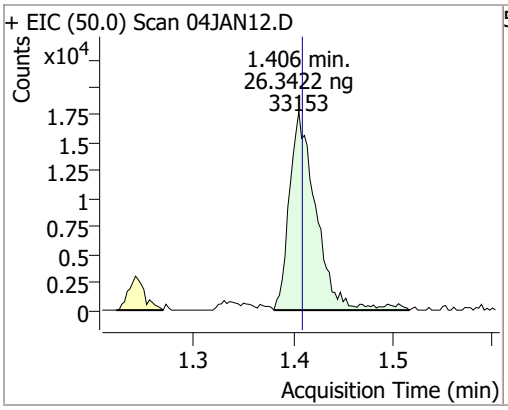
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	25.6793	1.24	0.00	26627	87.0	31.7	2.3	62.3

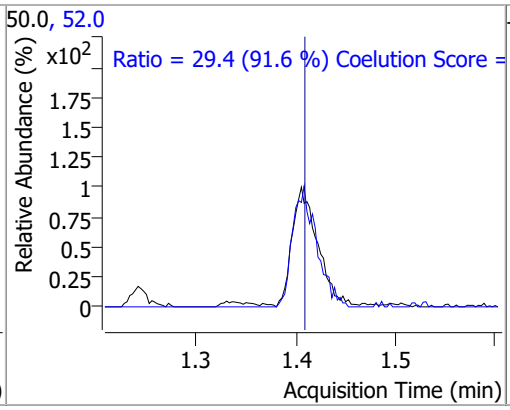


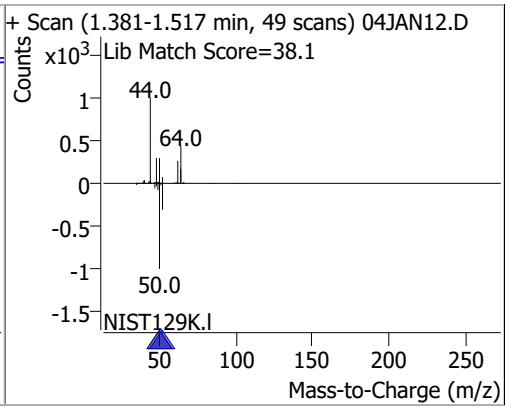




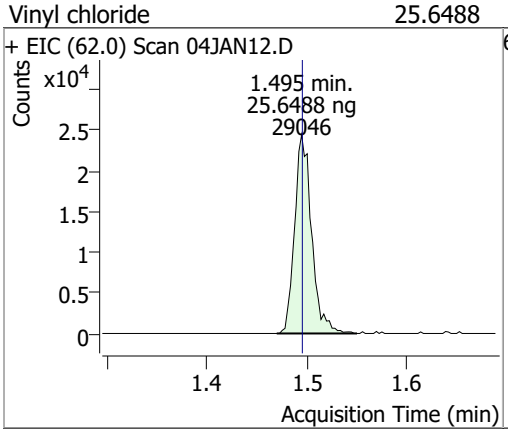
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	26.3422	1.41	0.00	33153	52.0	29.4	2.1	62.1

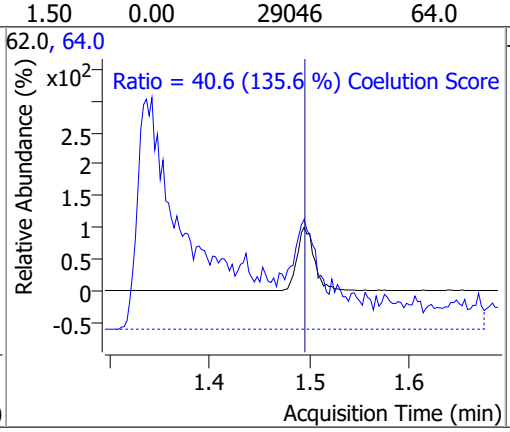


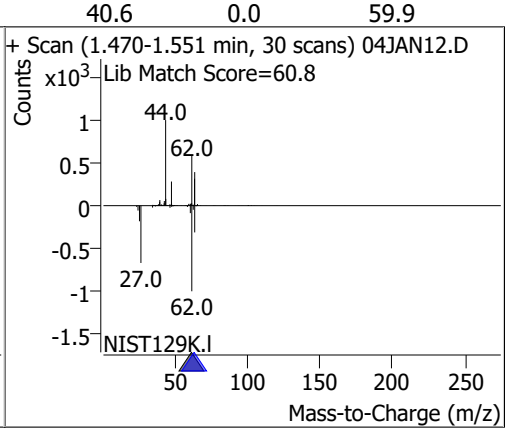




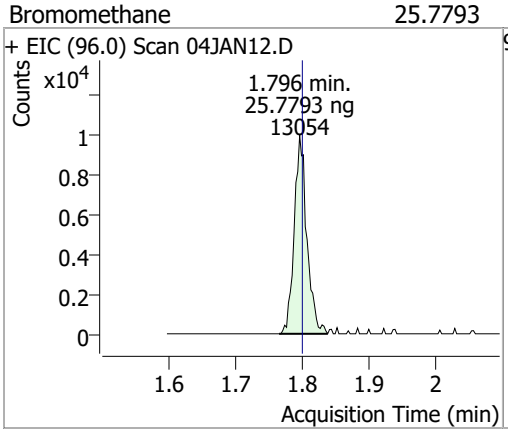
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	25.6488	1.50	0.00	29046	64.0	40.6	0.0	59.9

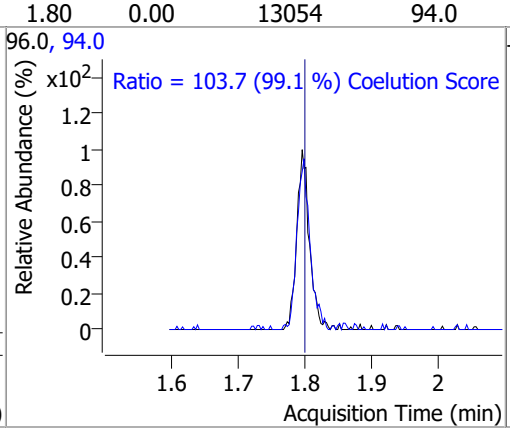


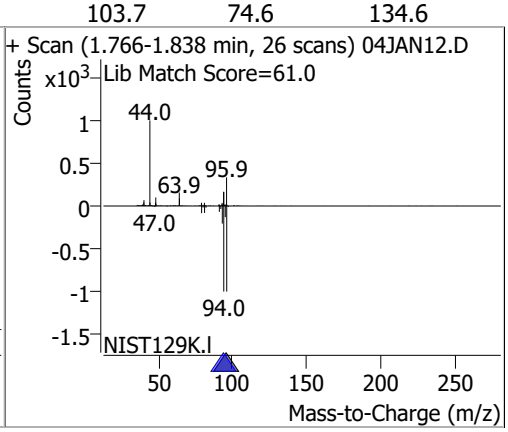




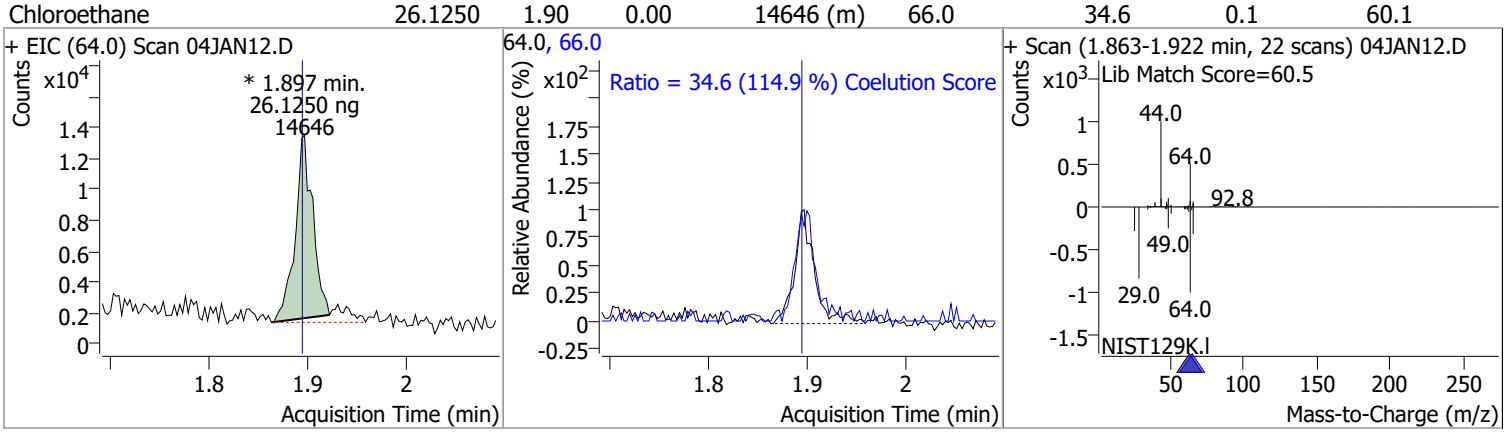
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	25.7793	1.80	0.00	13054	94.0	103.7	74.6	134.6



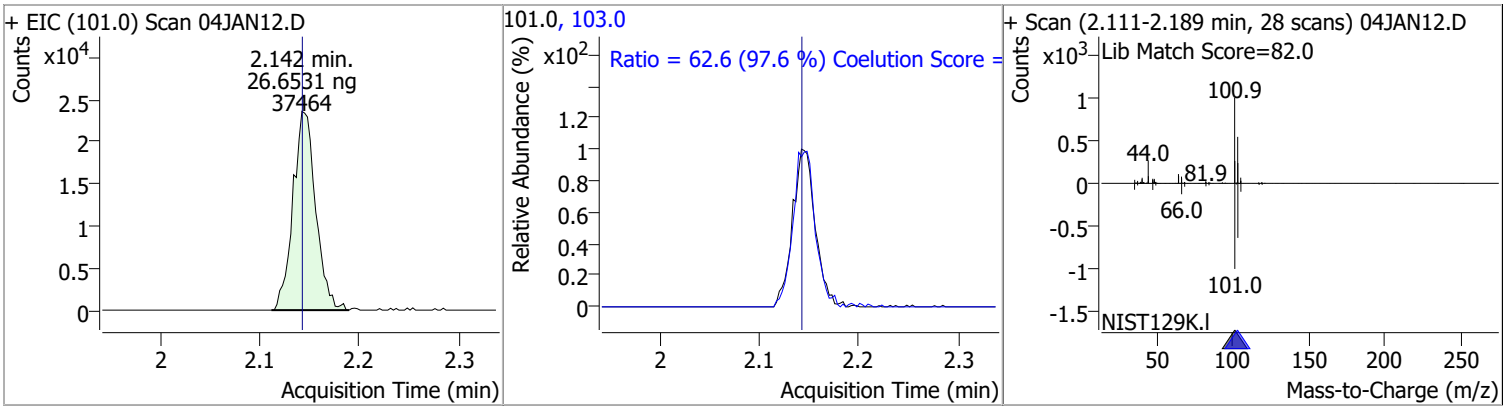




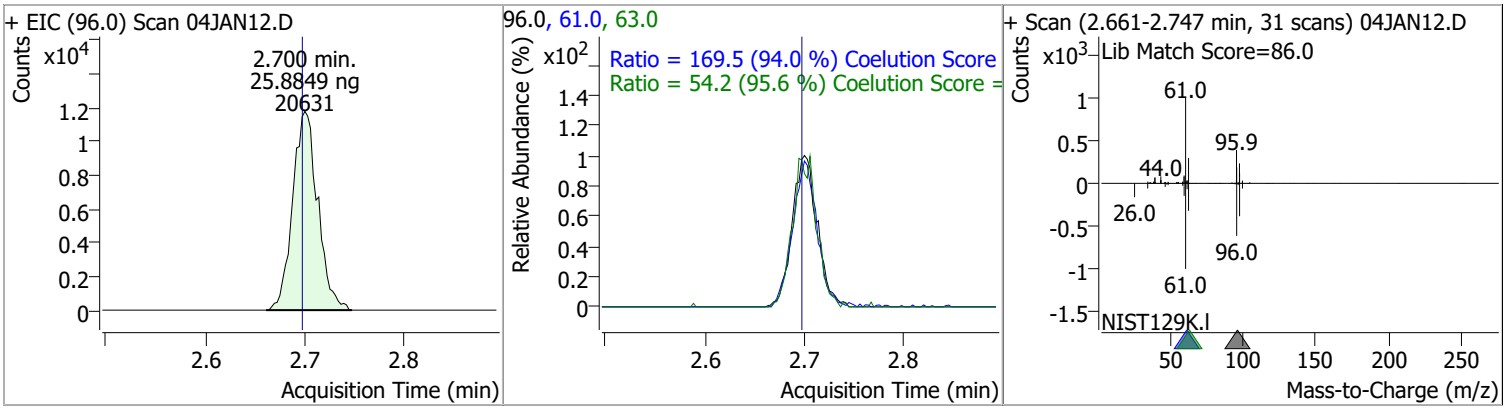
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	26.1250	1.90	0.00	14646 (m)	66.0	34.6	0.1	60.1



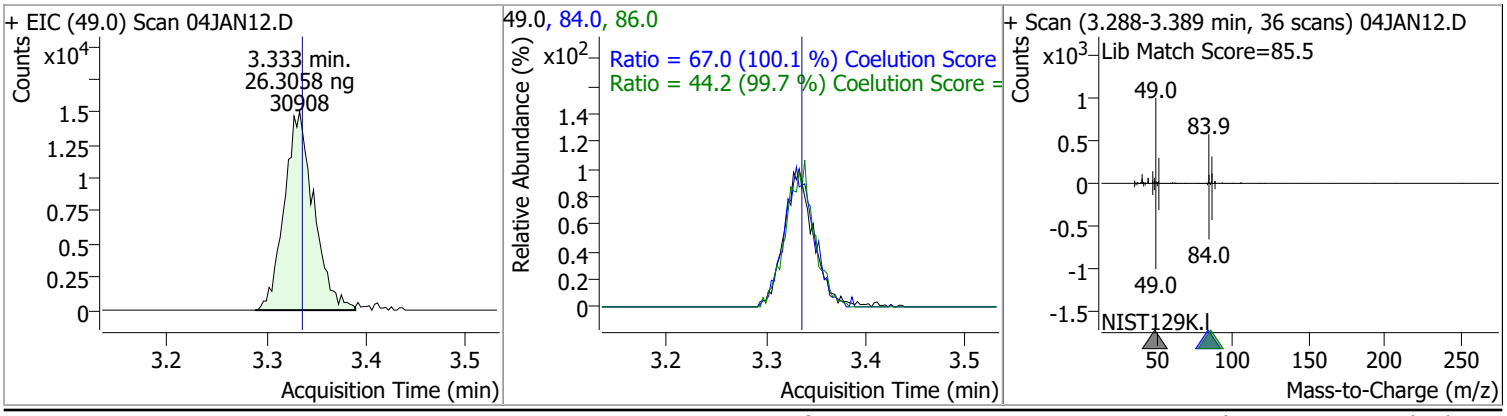
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	26.6531	2.14	0.00	37464	103.0	62.6	34.2	94.2



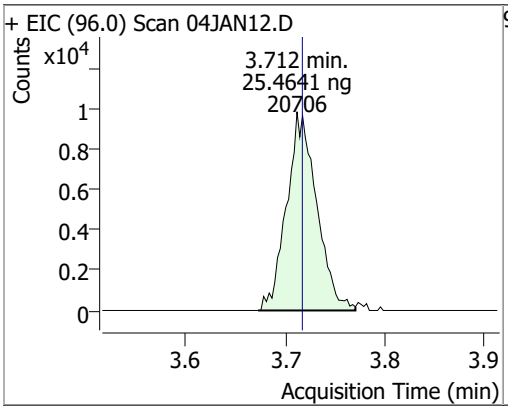
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	25.8849	2.70	0.00	20631	61.0	169.5	150.3	210.3
					63.0	54.2	26.7	86.7

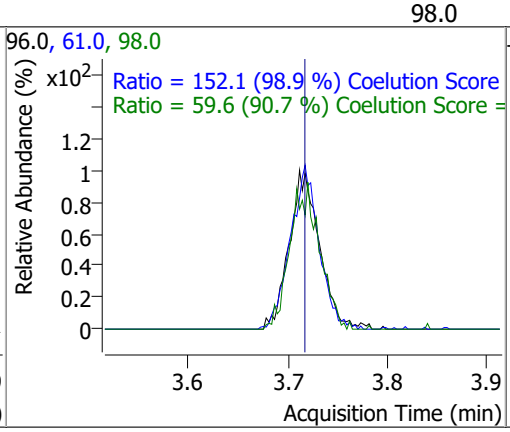


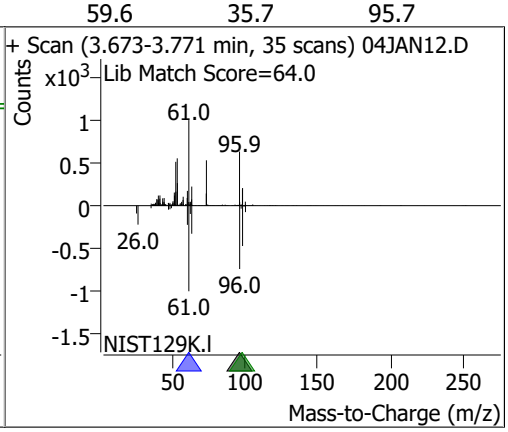
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	26.3058	3.33	0.00	30908	84.0	67.0	36.9	96.9
					86.0	44.2	14.3	74.3



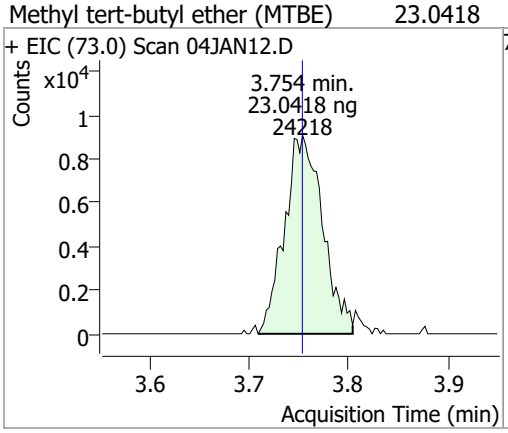
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	25.4641	3.71	-0.01	20706	61.0	152.1	123.9	183.9
					98.0	59.6	35.7	95.7

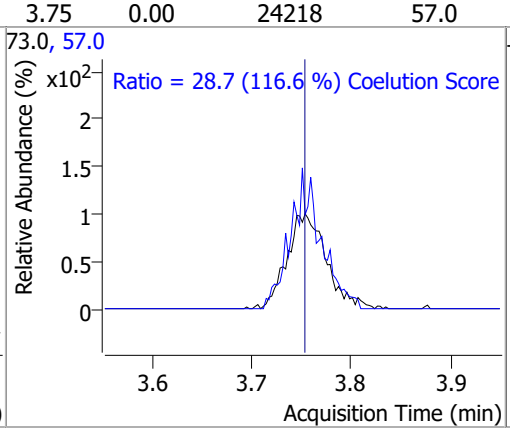


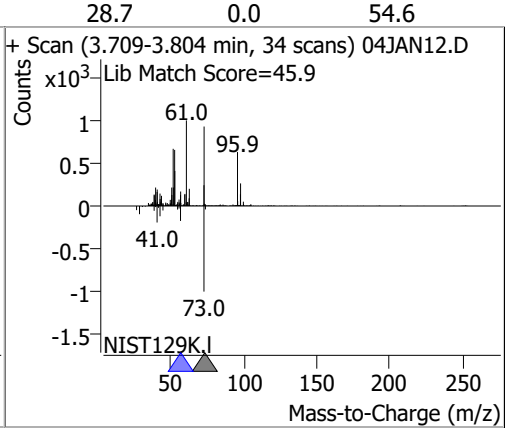




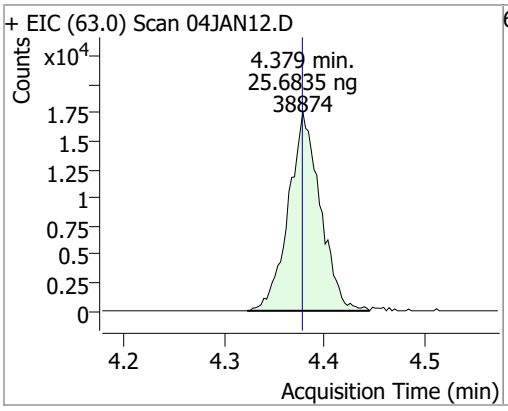
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	23.0418	3.75	0.00	24218	57.0	28.7	0.0	54.6

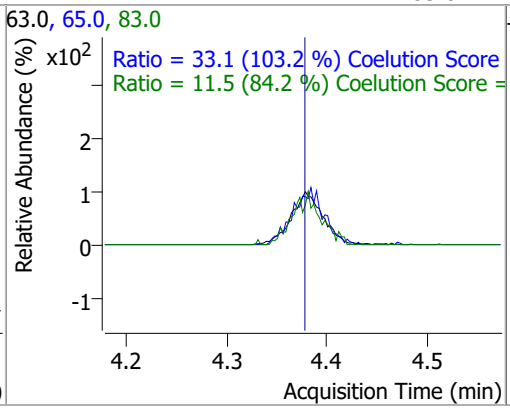


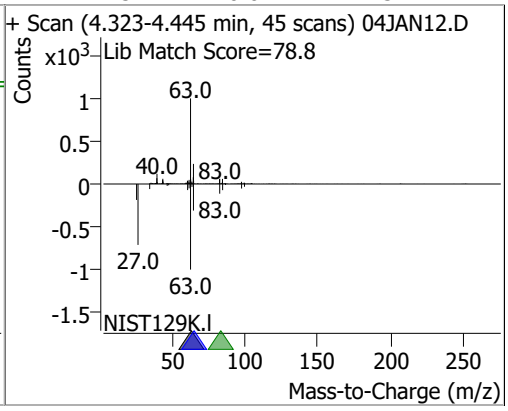




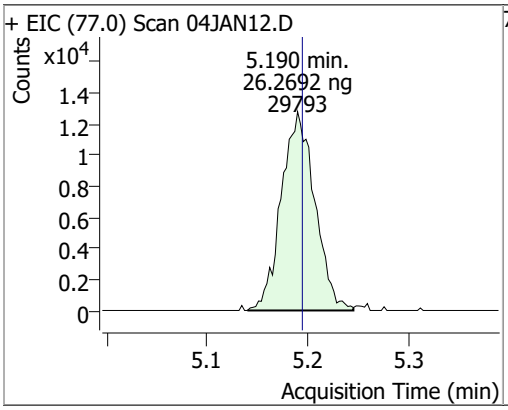
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	25.6835	4.38	0.00	38874	65.0	33.1	2.1	62.1
					83.0	11.5	0.0	43.7

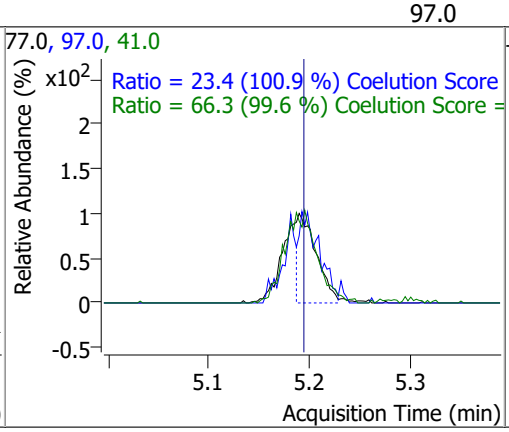


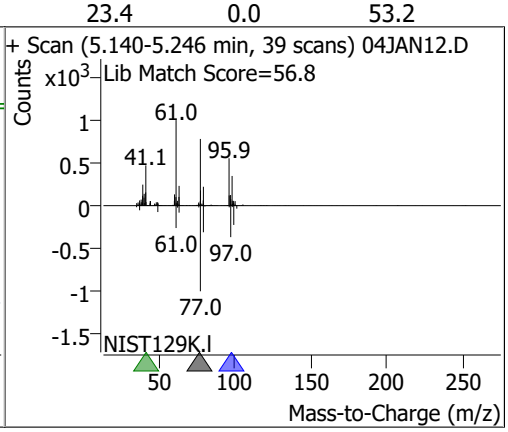




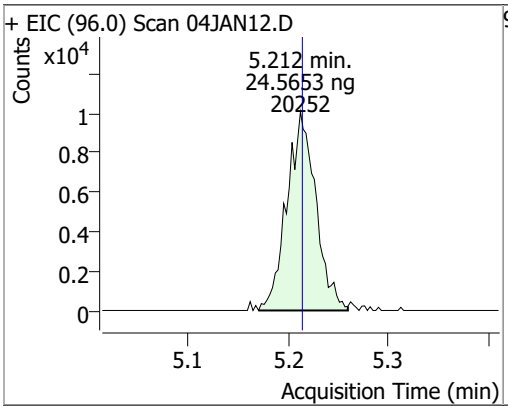
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	26.2692	5.19	-0.01	29793	41.0	66.3	36.5	96.5
					97.0	23.4	0.0	53.2

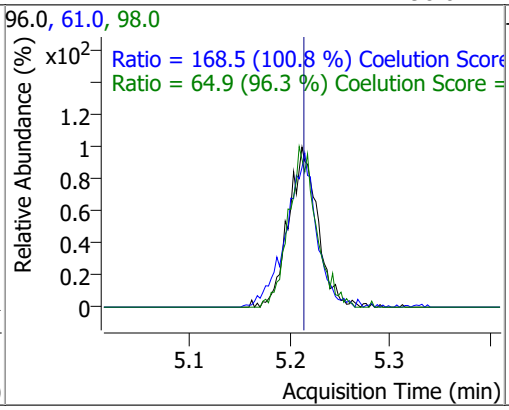


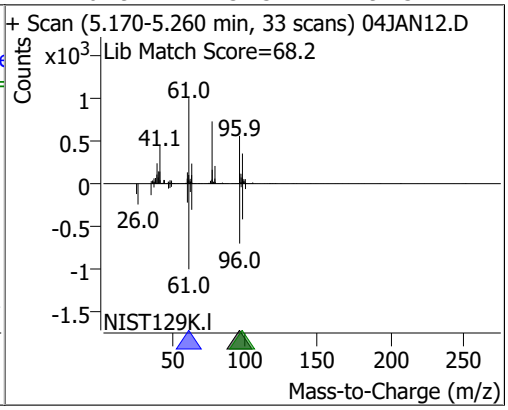




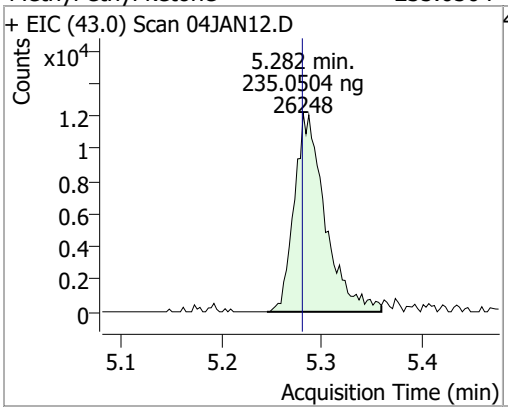
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	24.5653	5.21	0.00	20252	61.0	168.5	137.2	197.2
					98.0	64.9	37.3	97.3

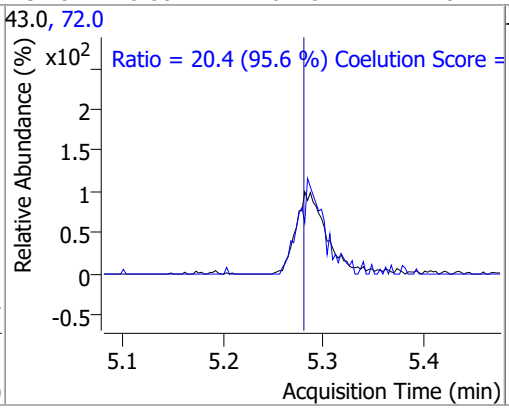


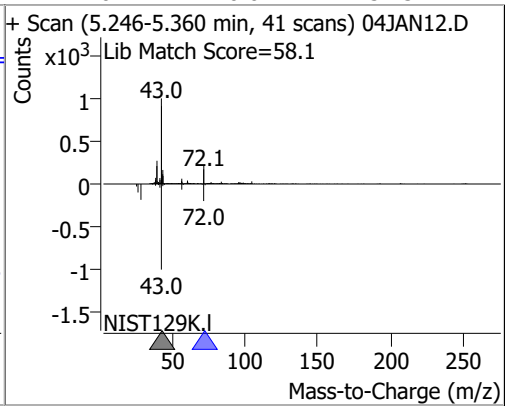




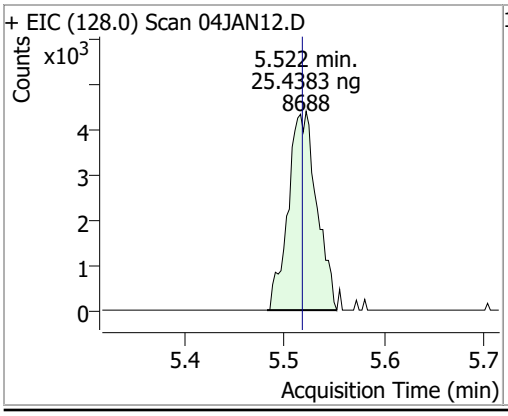
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	235.0504	5.28	0.00	26248	72.0	20.4	0.0	51.3

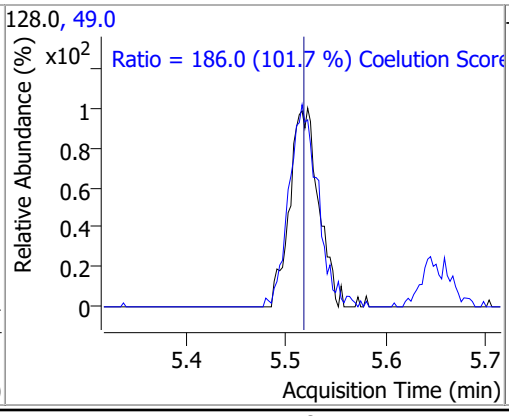


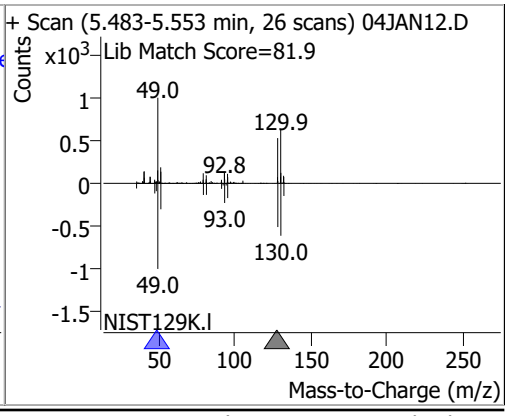




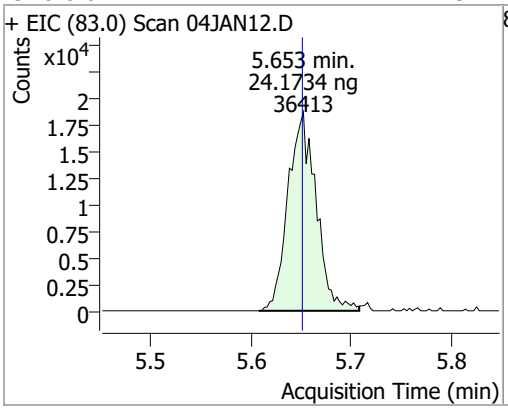
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	25.4383	5.52	0.00	8688	49.0	186.0	152.9	212.9

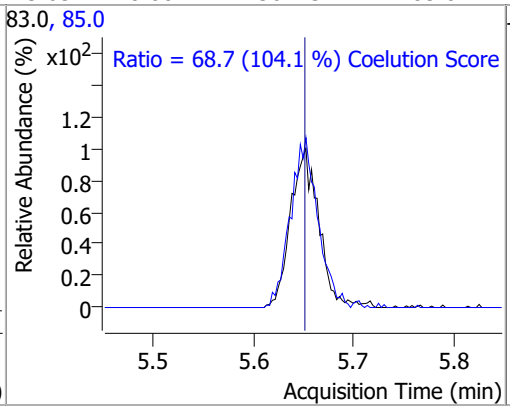


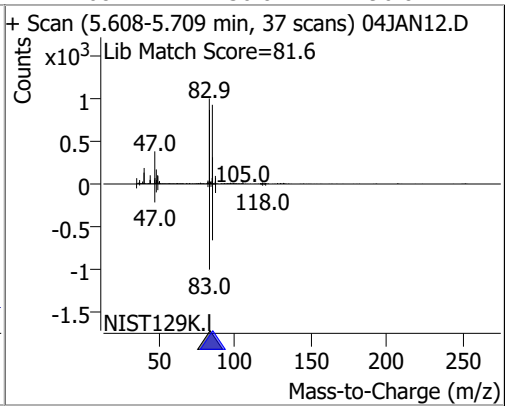




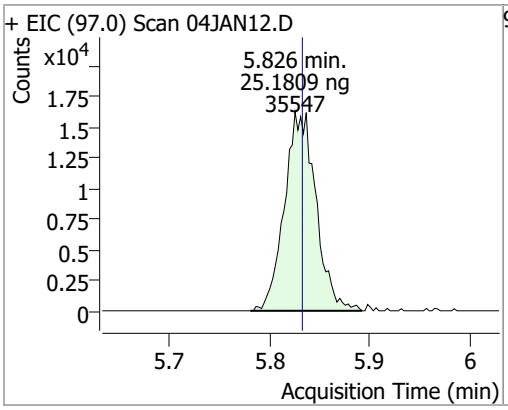
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	24.1734	5.65	0.00	36413	85.0	68.7	36.0	96.0

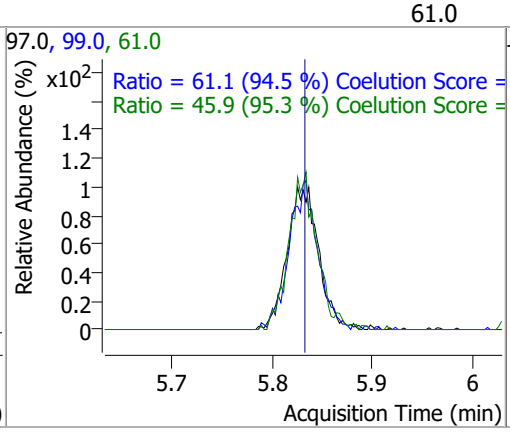


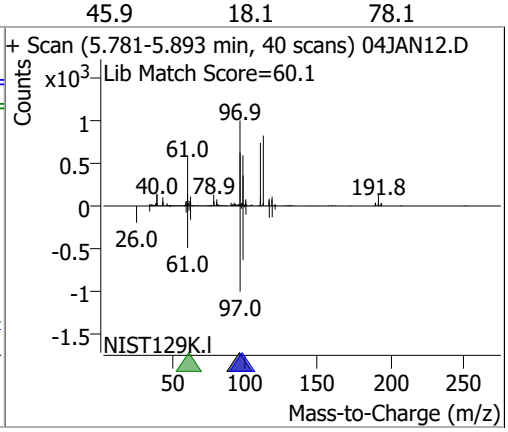




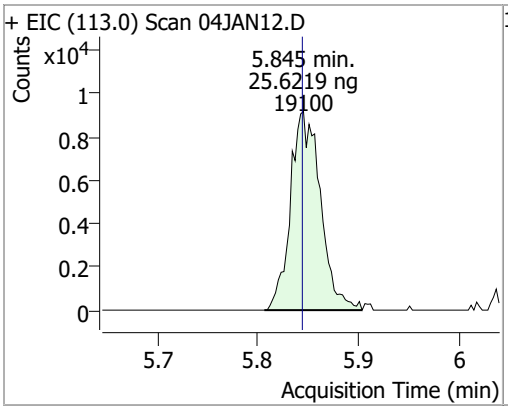
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	25.1809	5.83	-0.01	35547	99.0	61.1	34.7	94.7
					61.0	45.9	18.1	78.1

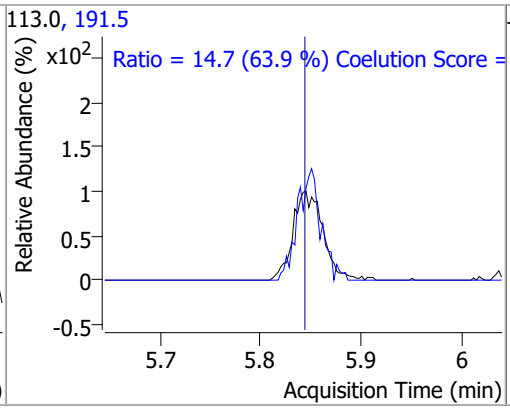


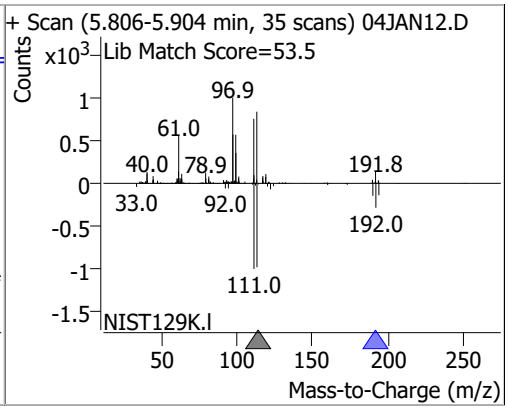




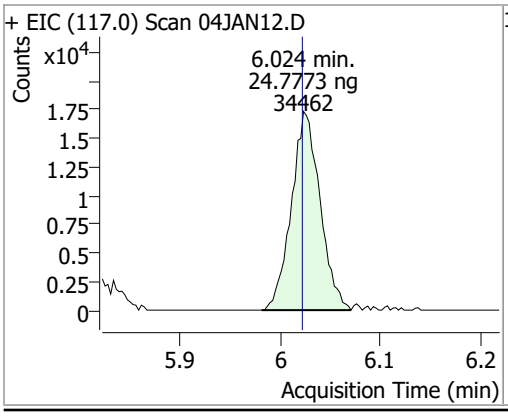
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	25.6219	5.85	0.00	19100	191.5	14.7	0.0	53.1

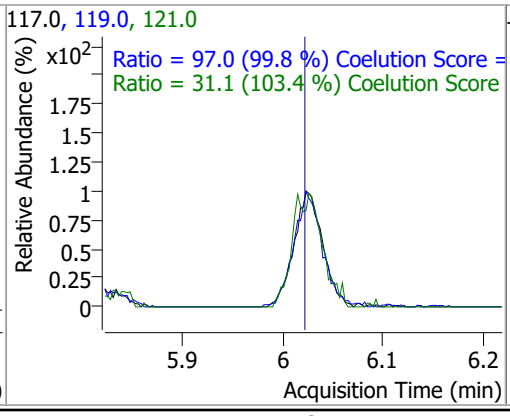


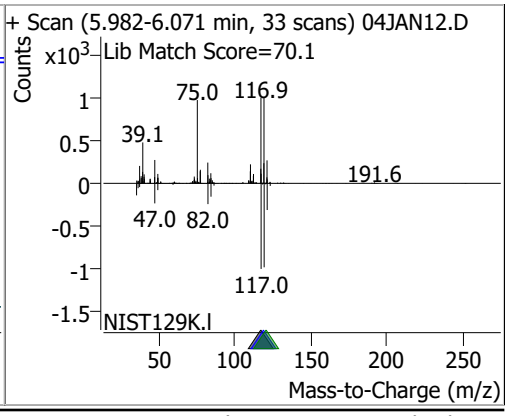




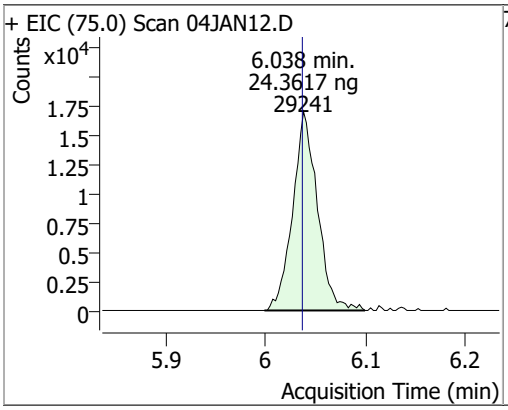
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	24.7773	6.02	0.00	34462	119.0	97.0	67.2	127.2
					121.0	31.1	0.1	60.1

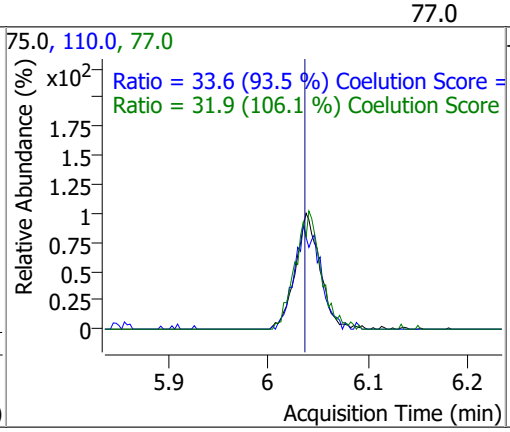


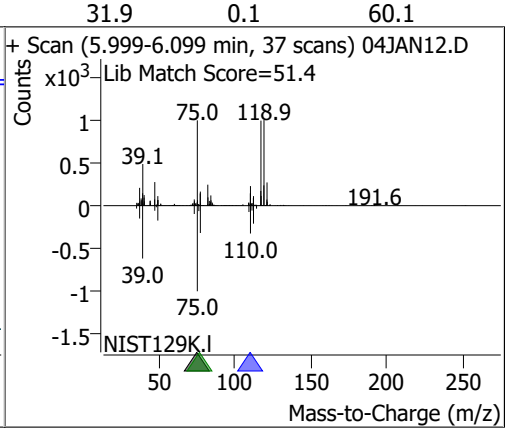




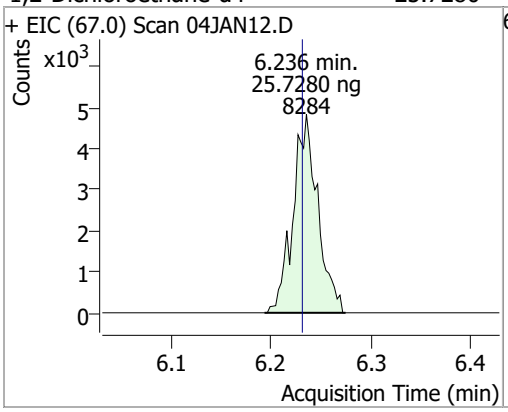
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	24.3617	6.04	0.00	29241	110.0	33.6	5.9	65.9
					77.0	31.9	0.1	60.1

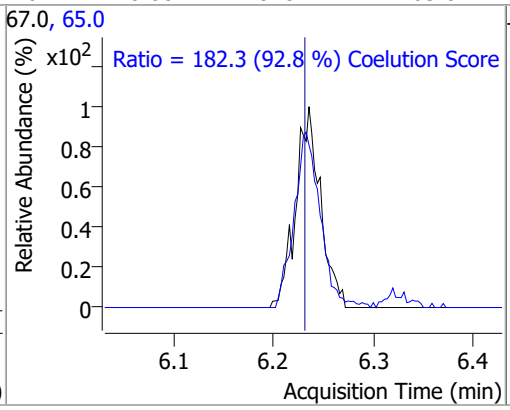


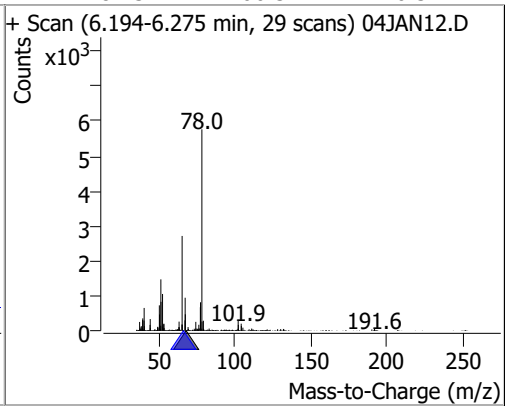




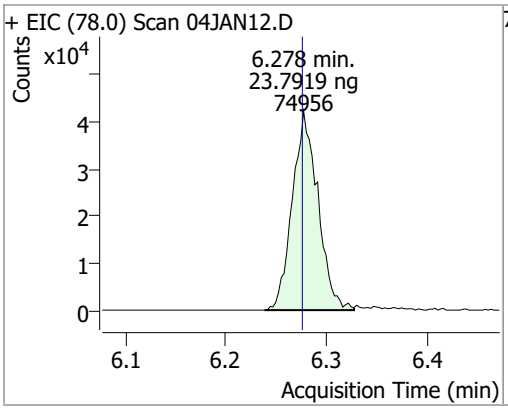
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	25.7280	6.24	0.00	8284	65.0	182.3	166.5	226.5

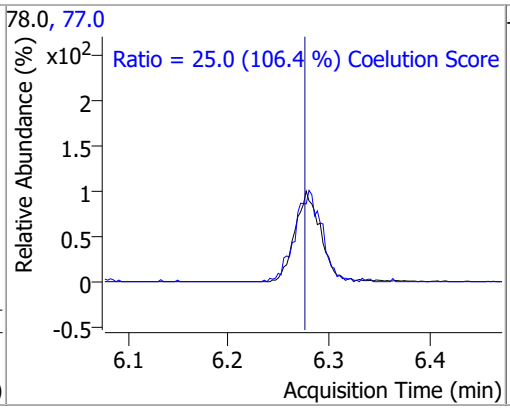


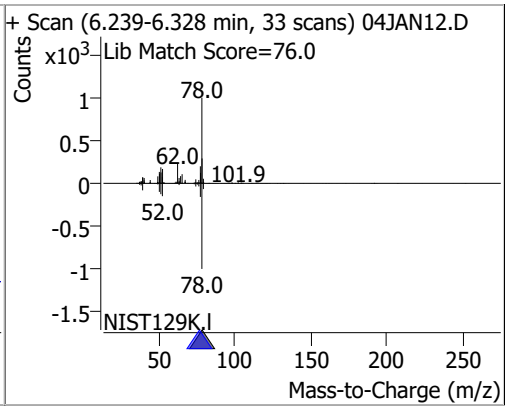




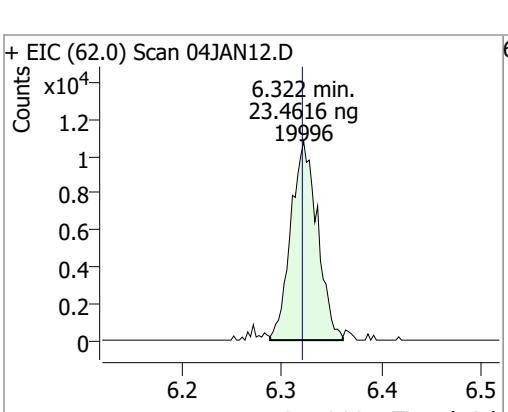
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	23.7919	6.28	0.00	74956	77.0	25.0	0.0	53.5

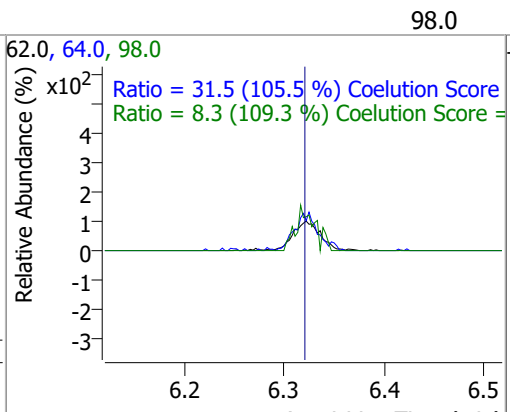


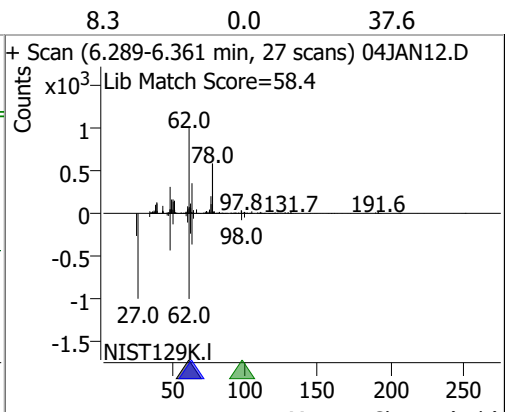




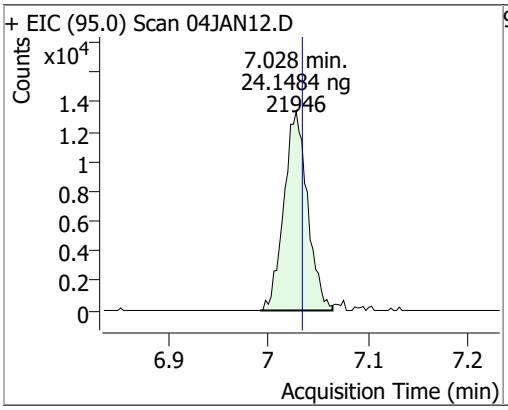
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	23.4616	6.32	0.00	19996	64.0	31.5	0.0	59.9
					98.0	8.3	0.0	37.6

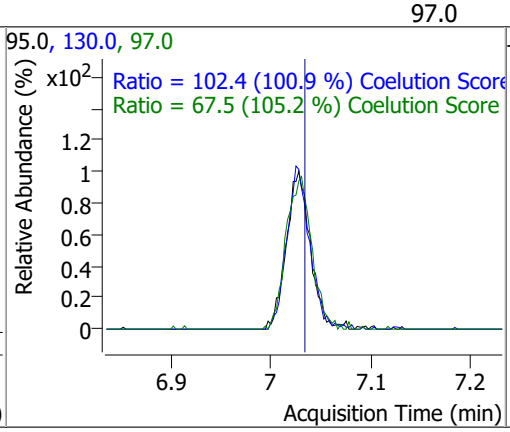


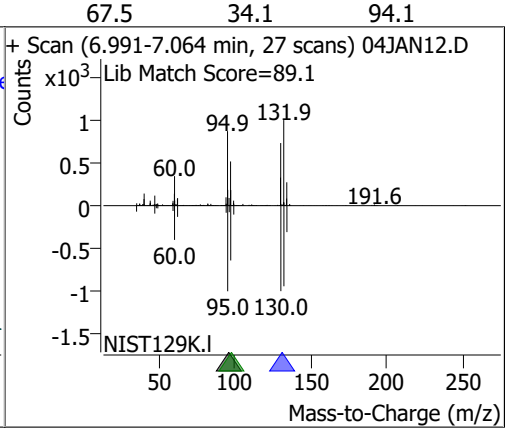




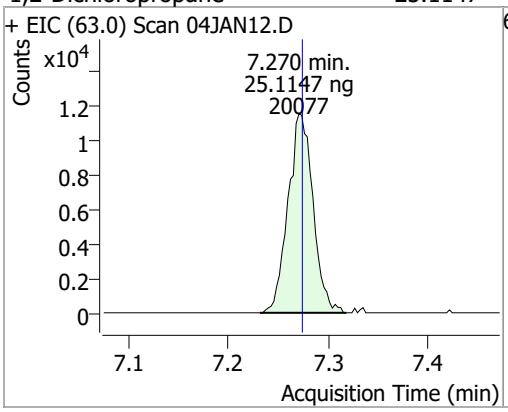
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	24.1484	7.03	0.00	21946	130.0	102.4	71.5	131.5
					97.0	67.5	34.1	94.1

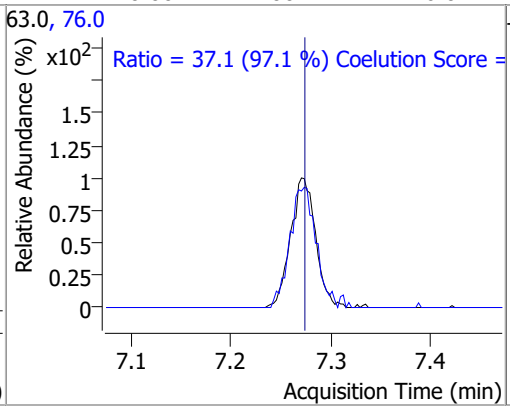


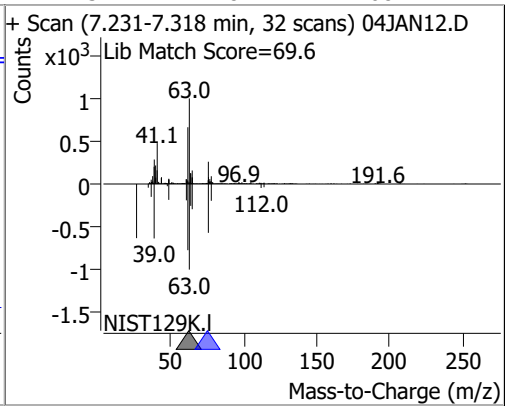




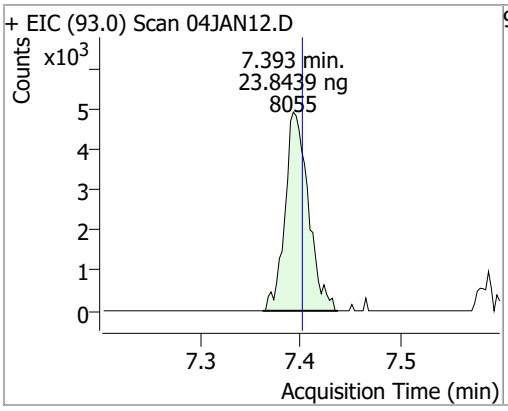
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	25.1147	7.27	0.00	20077	76.0	37.1	8.2	68.2

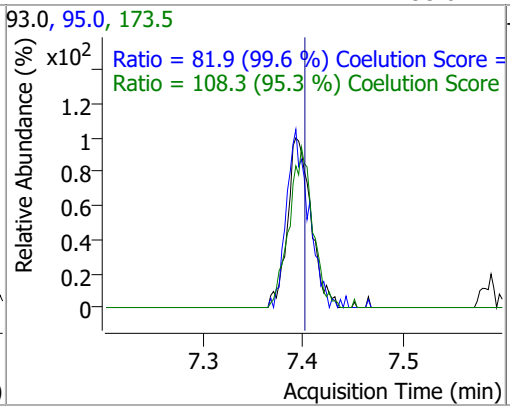


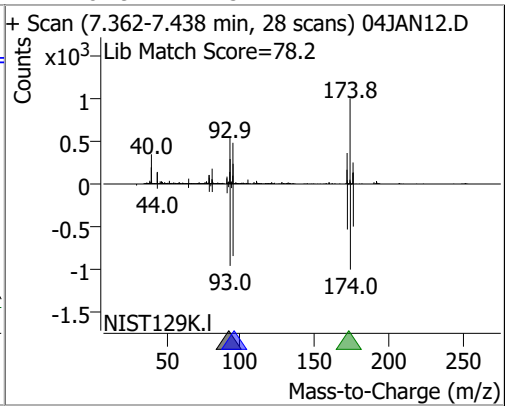




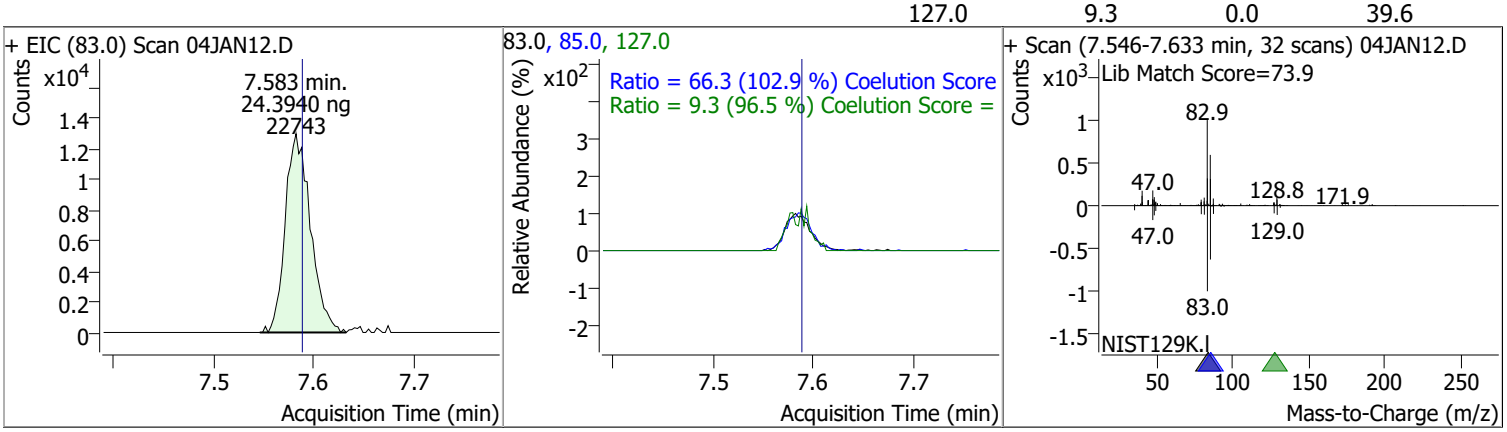
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	23.8439	7.39	-0.01	8055	173.5	108.3	83.7	143.7
					95.0	81.9	52.2	112.2



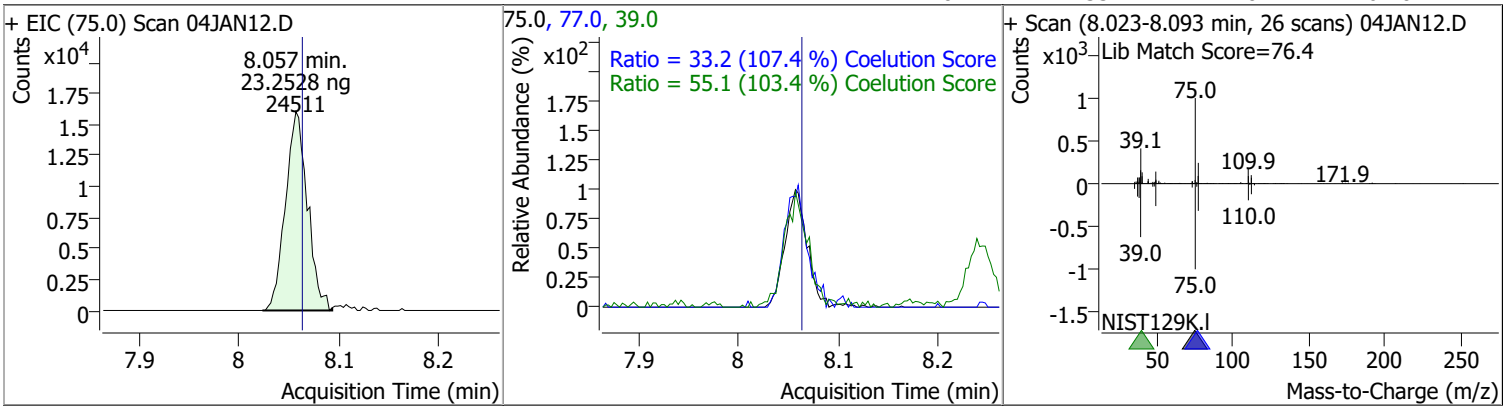




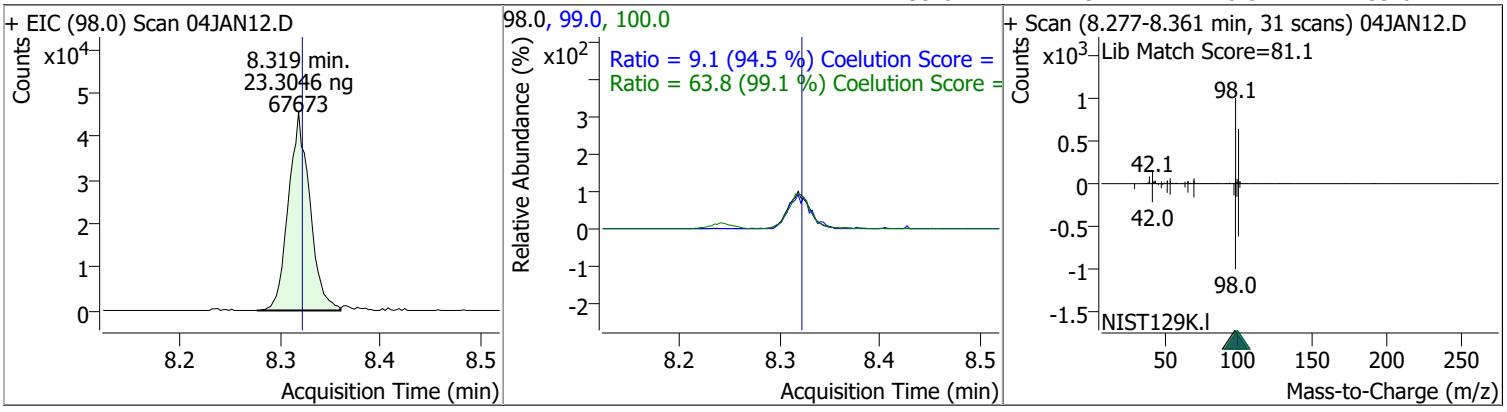
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	24.3940	7.58	0.00	22743	85.0	66.3	34.5	94.5
					127.0	9.3	0.0	39.6



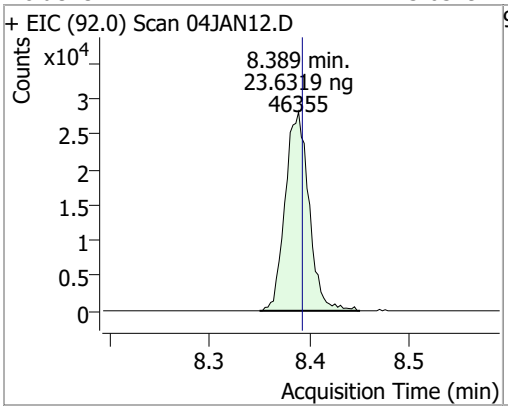
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	23.2528	8.06	0.00	24511	39.0	55.1	23.3	83.3
					77.0	33.2	1.0	61.0

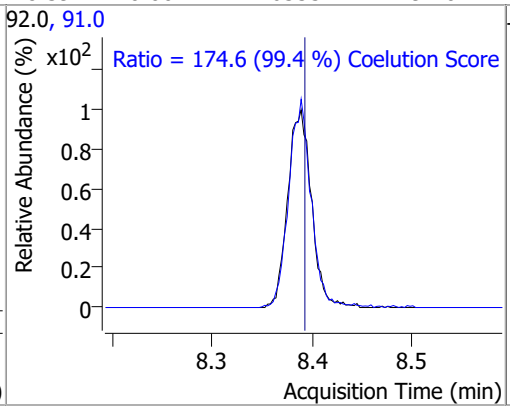


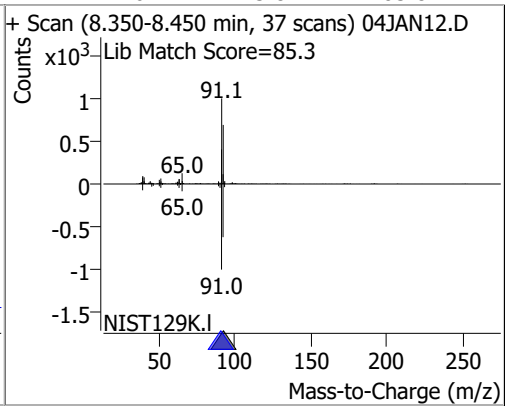
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	23.3046	8.32	0.00	67673	100.0	63.8	34.4	94.4
					99.0	9.1	0.0	39.6



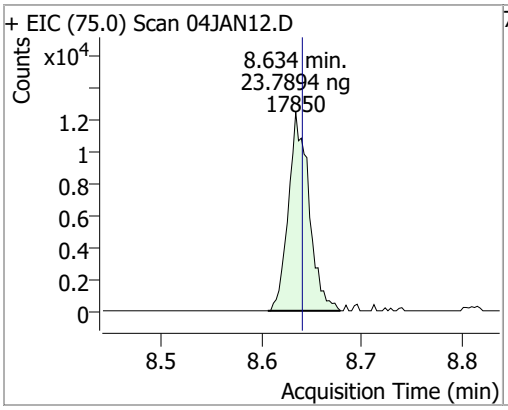
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	23.6319	8.39	0.00	46355	91.0	174.6	145.8	205.8

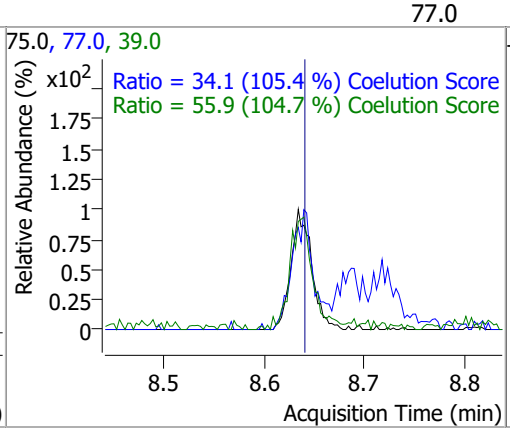


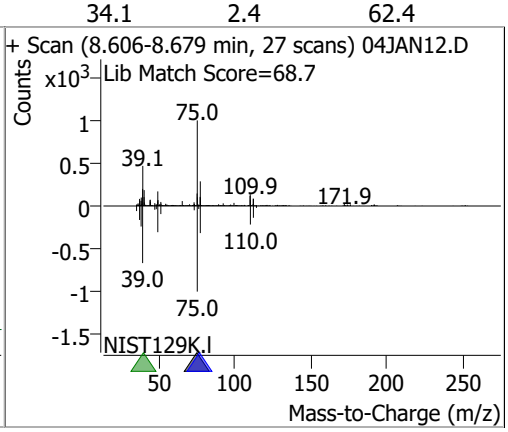




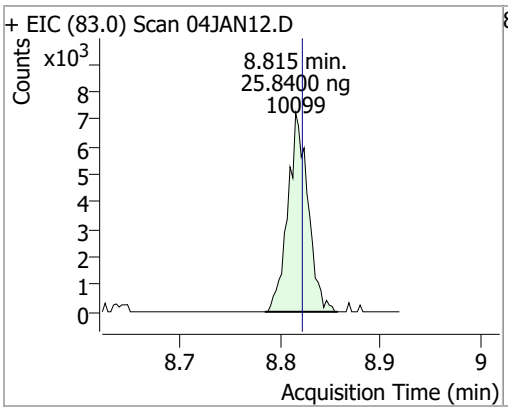
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	23.7894	8.63	0.00	17850	39.0	55.9	23.4	83.4
					77.0	34.1	2.4	62.4

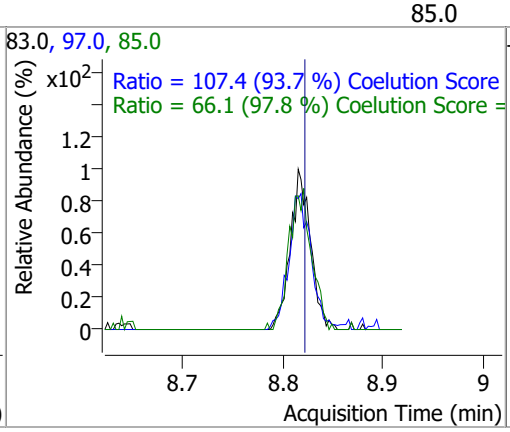


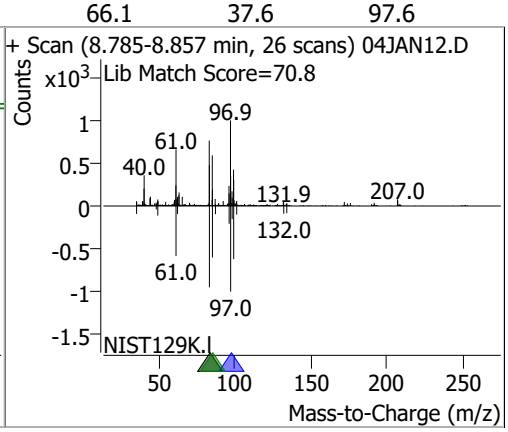




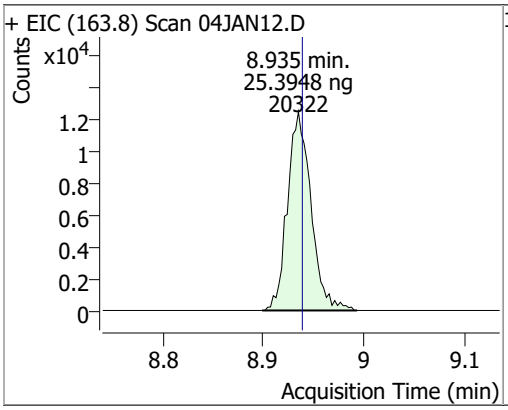
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	25.8400	8.82	0.00	10099	97.0	107.4	84.6	144.6
					85.0	66.1	37.6	97.6

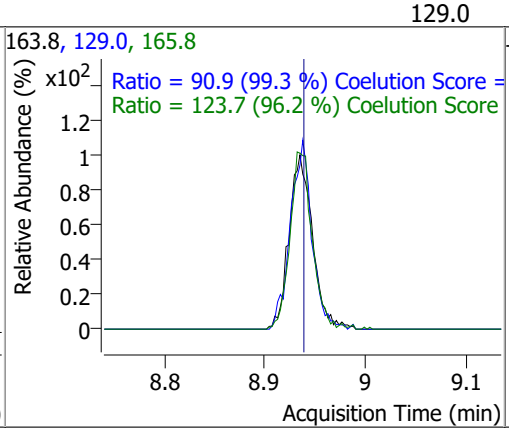


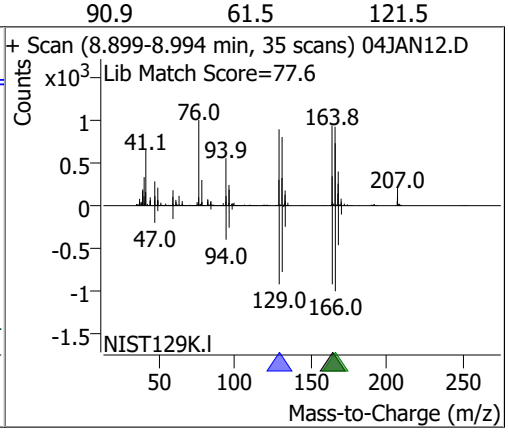




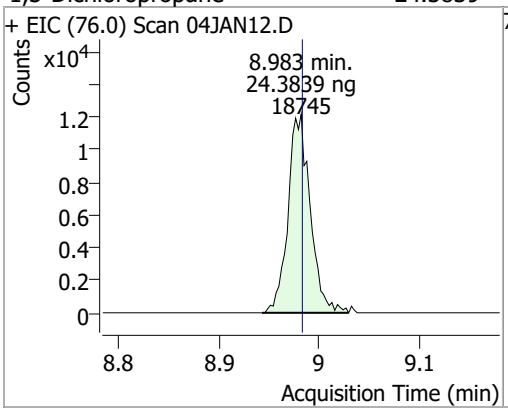
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	25.3948	8.94	0.00	20322	165.8	123.7	98.6	158.6
					129.0	90.9	61.5	121.5

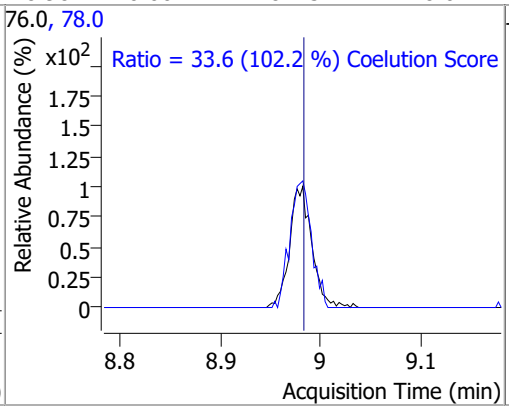


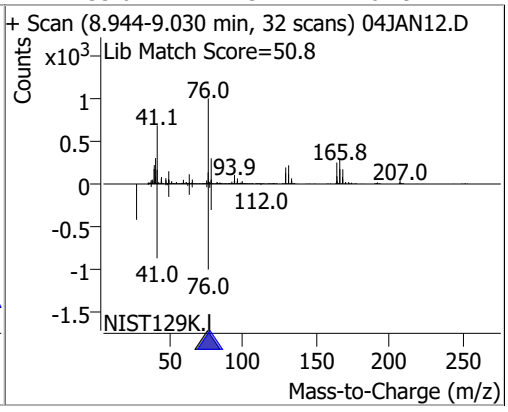




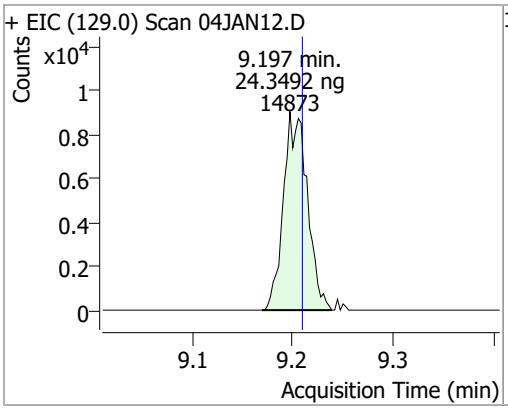
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	24.3839	8.98	0.00	18745	78.0	33.6	2.9	62.9

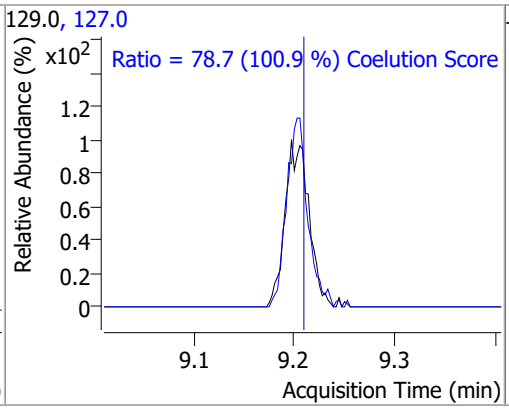


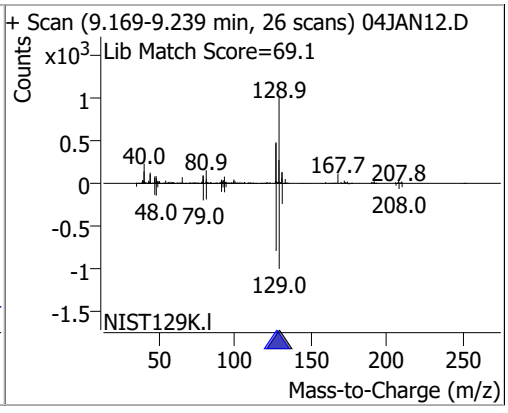




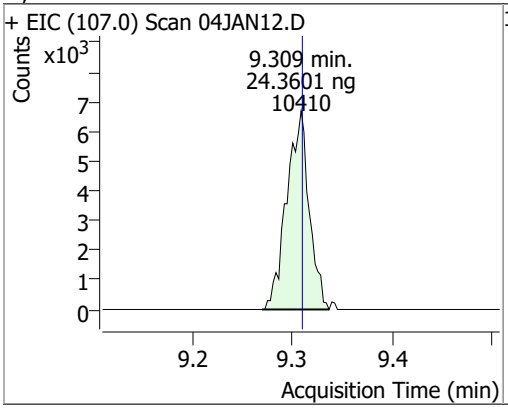
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	24.3492	9.20	-0.01	14873	127.0	78.7	48.0	108.0

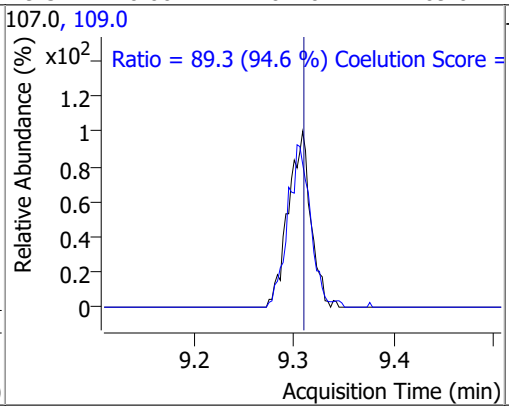


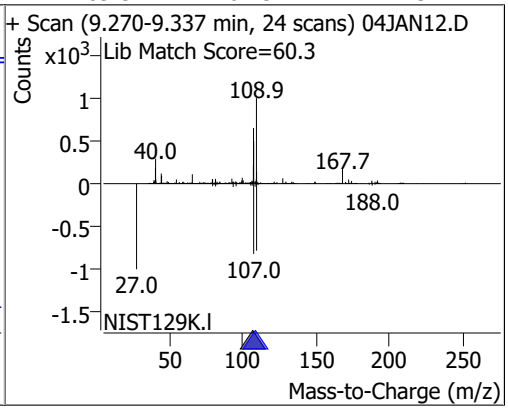




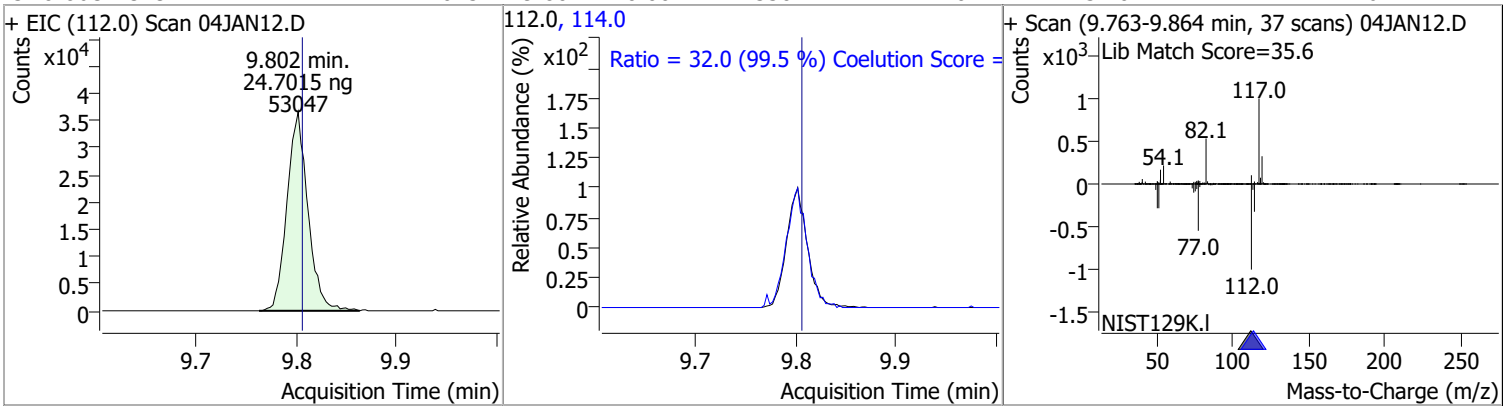
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	24.3601	9.31	0.00	10410	109.0	89.3	64.5	124.5



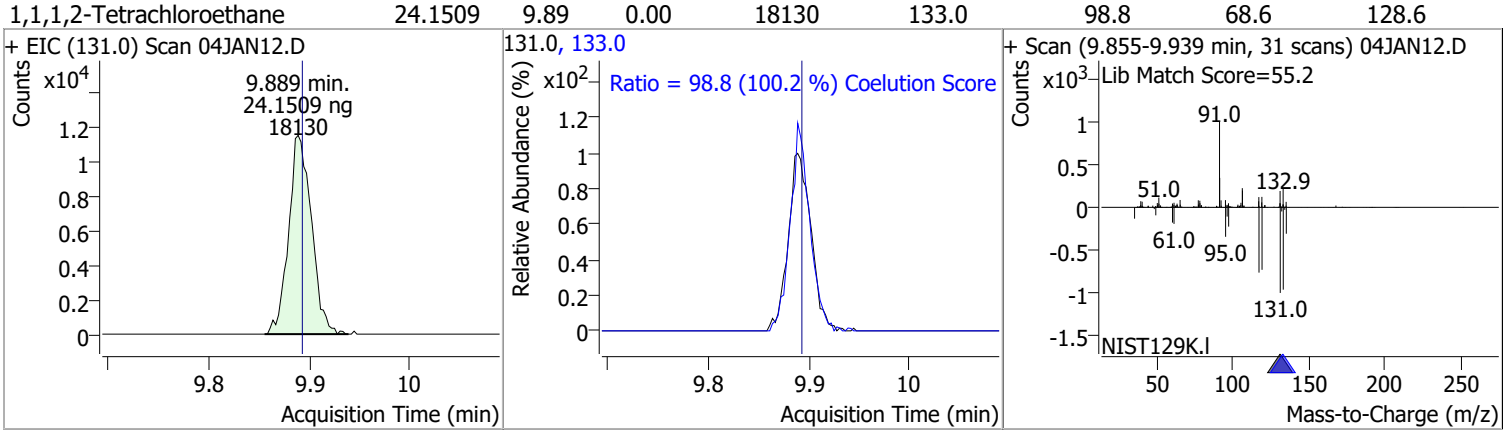




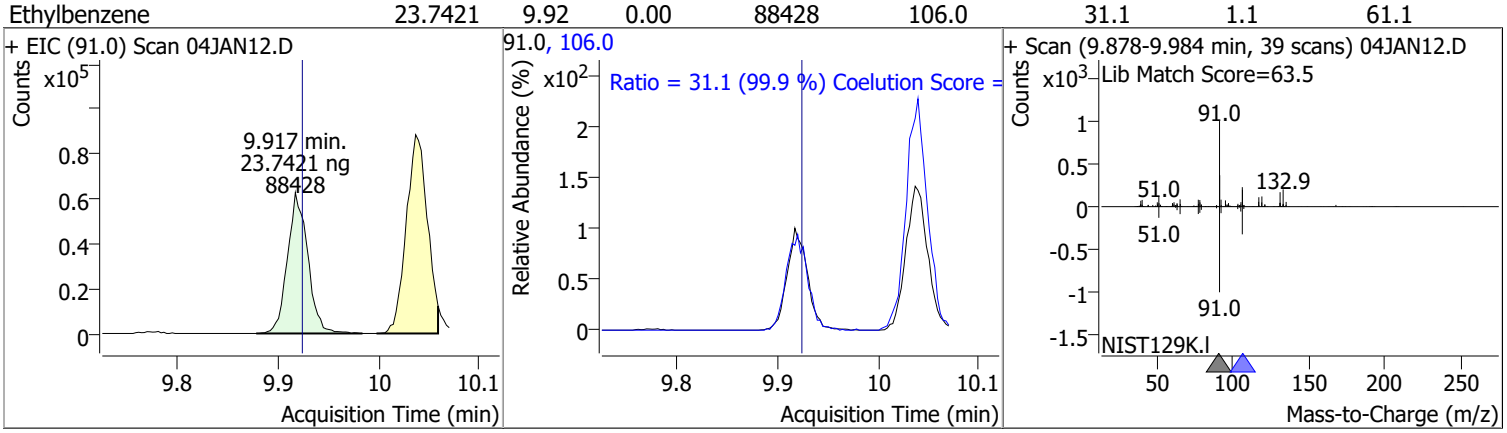
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	24.7015	9.80	0.00	53047	114.0	32.0	2.1	62.1



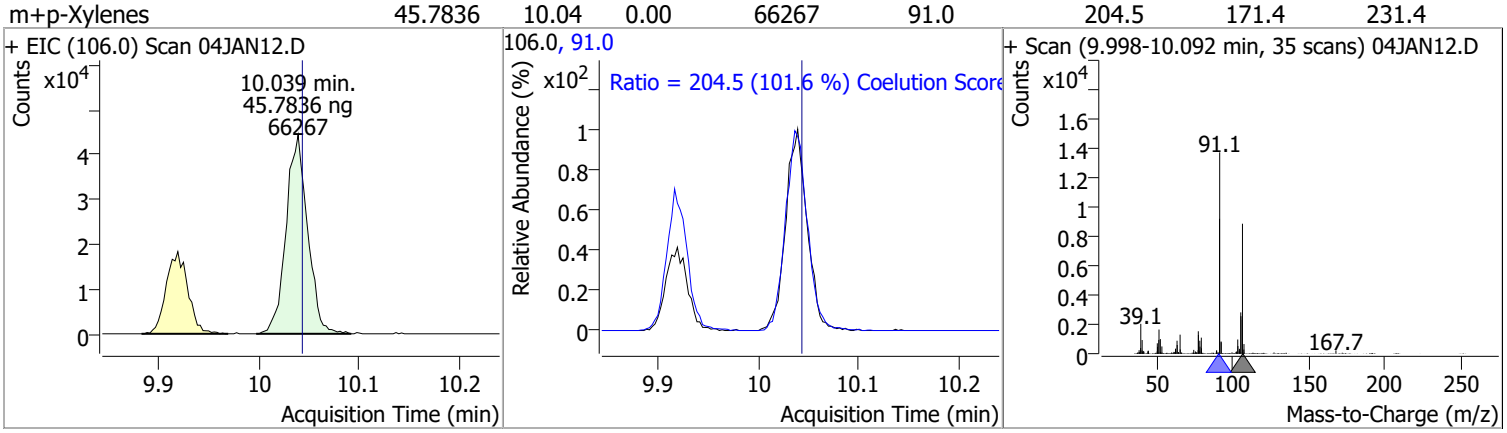
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	24.1509	9.89	0.00	18130	133.0	98.8	68.6	128.6



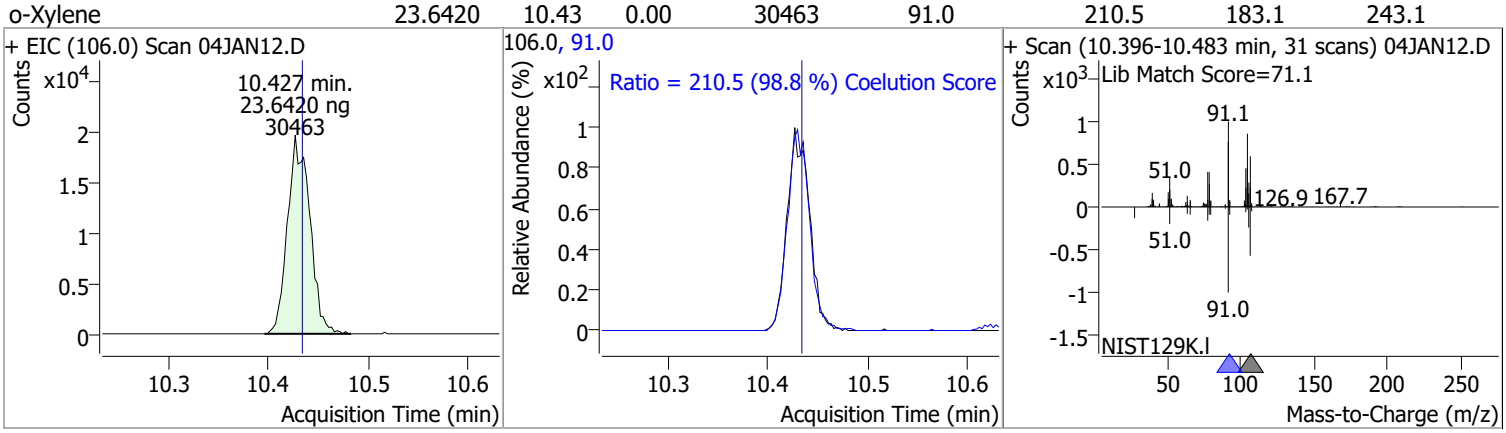
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	23.7421	9.92	0.00	88428	106.0	31.1	1.1	61.1



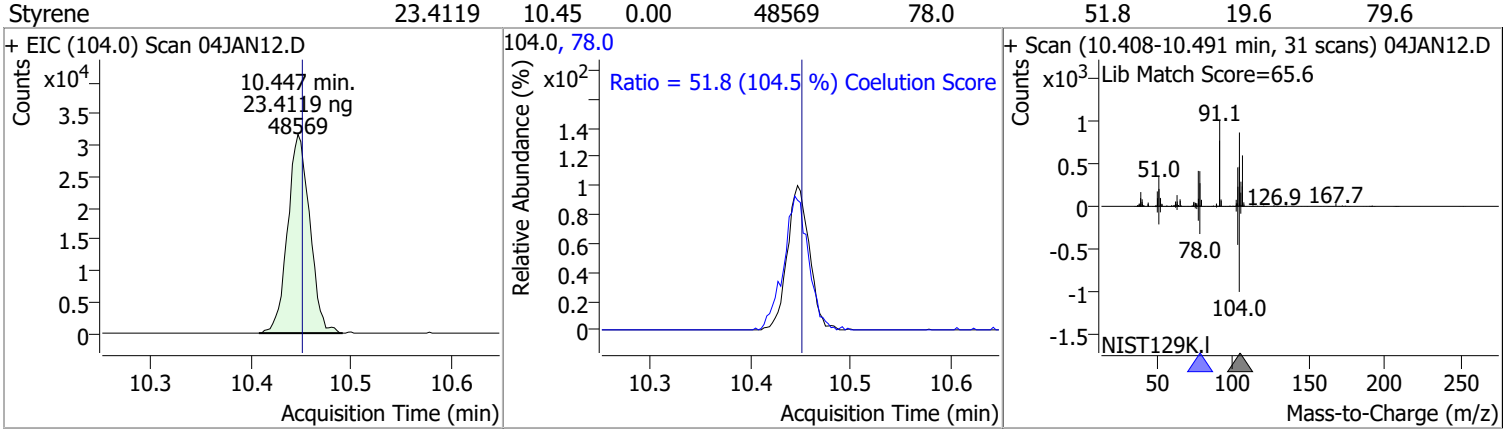
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	45.7836	10.04	0.00	66267	91.0	204.5	171.4	231.4



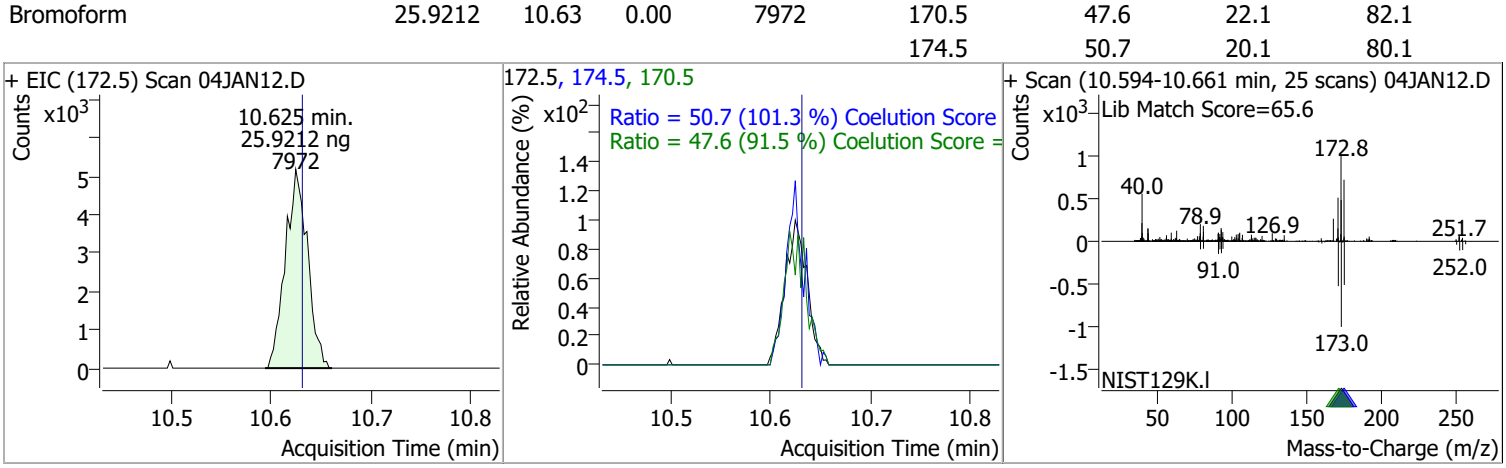
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	23.6420	10.43	0.00	30463	91.0	210.5	183.1	243.1



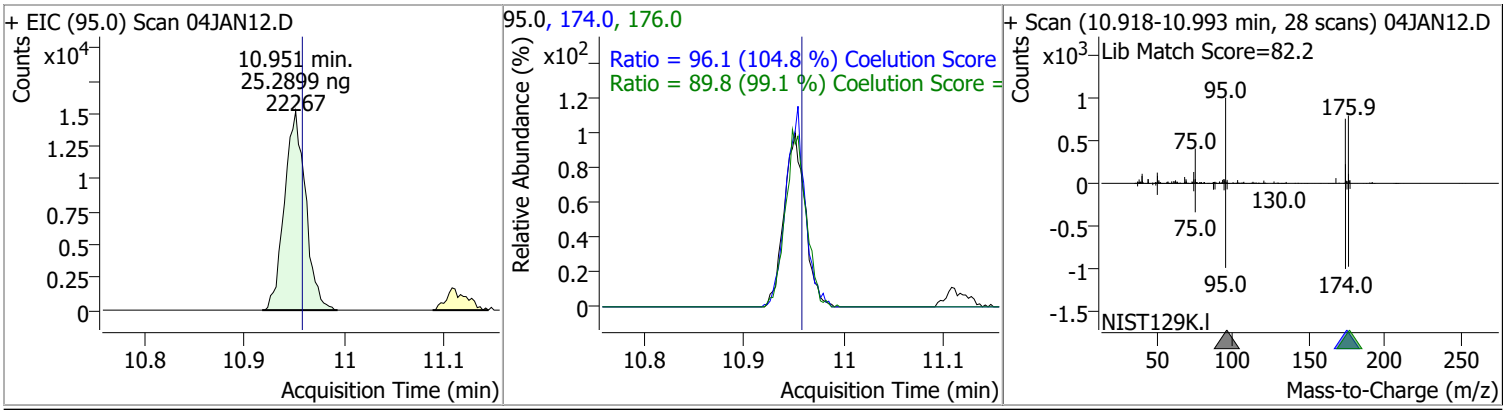
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	23.4119	10.45	0.00	48569	78.0	51.8	19.6	79.6



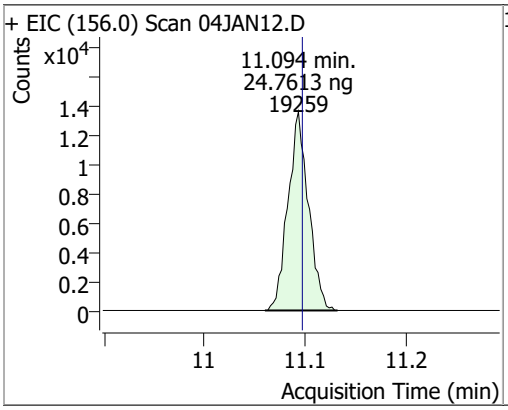
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	25.9212	10.63	0.00	7972	170.5	47.6	22.1	82.1
					174.5	50.7	20.1	80.1

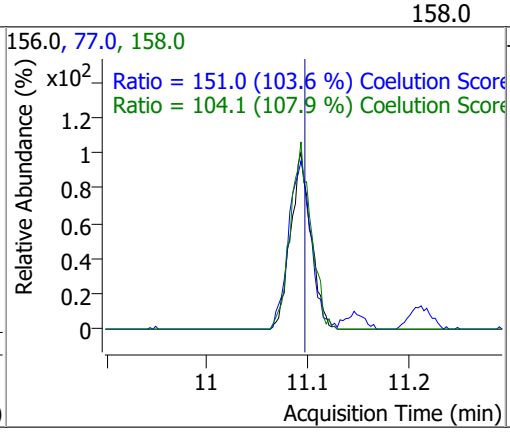


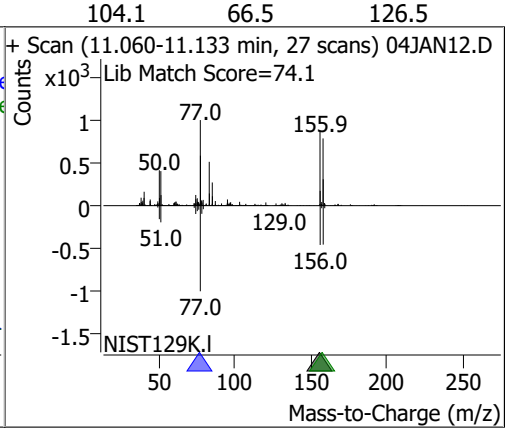
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	25.2899	10.95	0.00	22267	174.0	96.1	61.7	121.7
					176.0	89.8	60.6	120.6



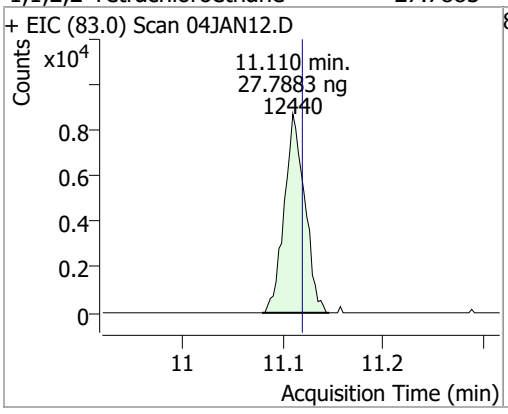
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	24.7613	11.09	0.00	19259	77.0	151.0	115.7	175.7
					158.0	104.1	66.5	126.5

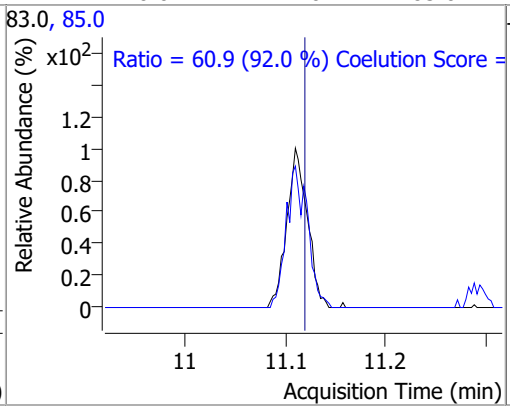


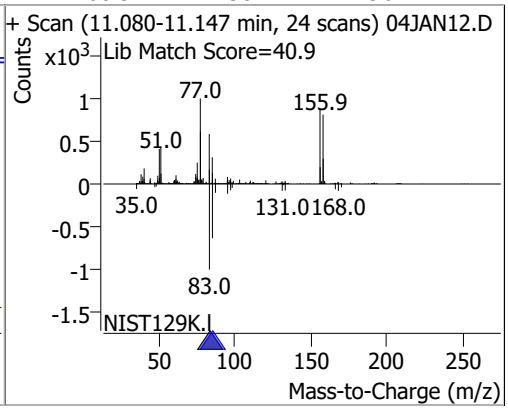




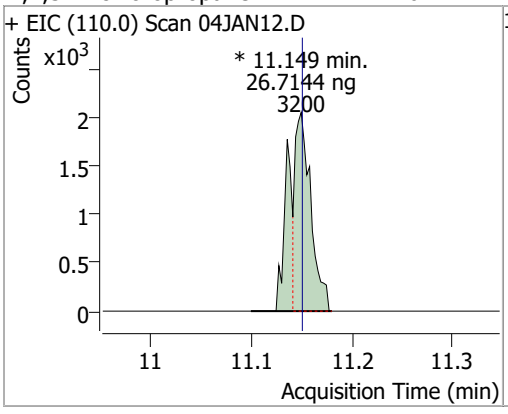
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	27.7883	11.11	-0.01	12440	85.0	60.9	36.2	96.2

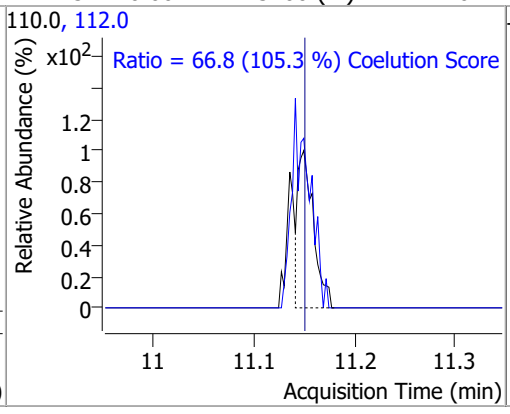


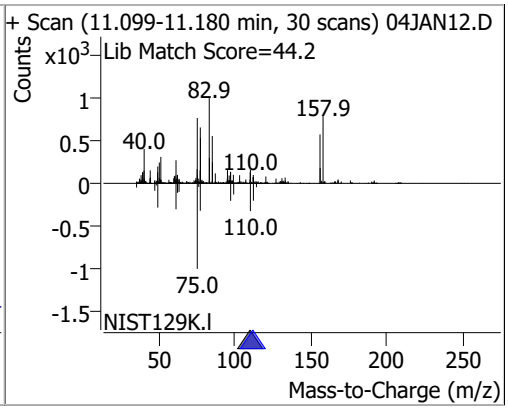




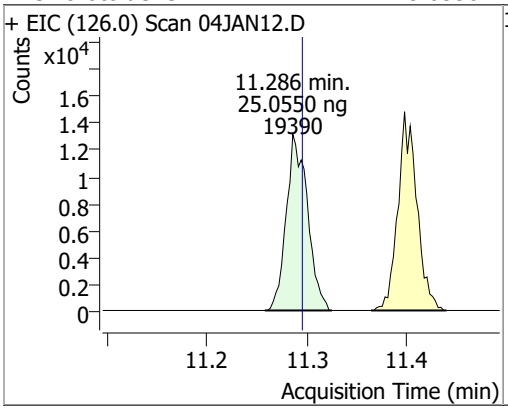
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	26.7144	11.15	0.00	3200 (m)	112.0	66.8	33.5	93.5

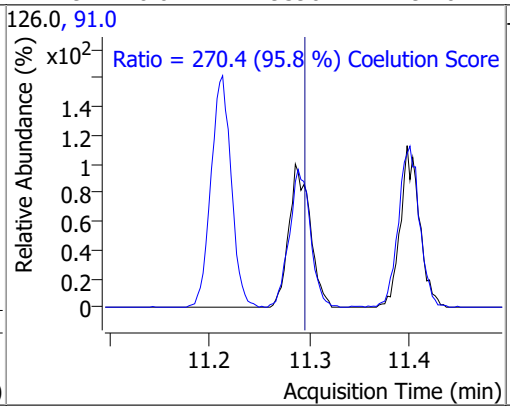


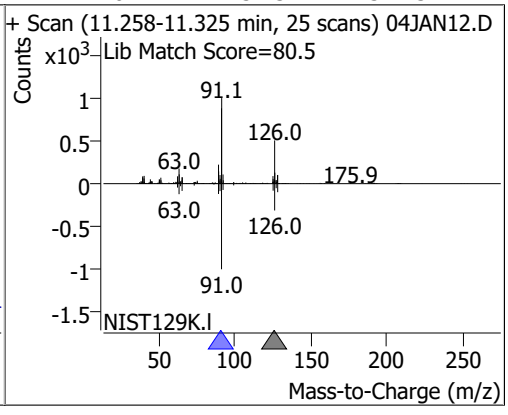




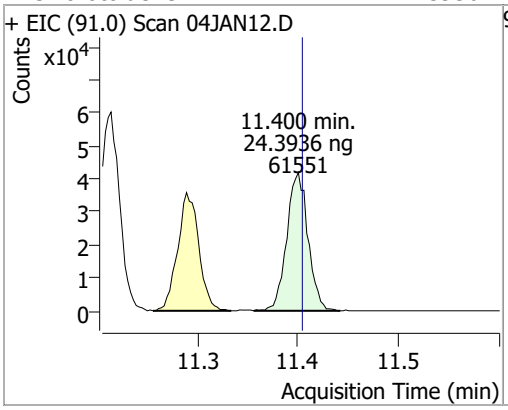
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	25.0550	11.29	-0.01	19390	91.0	270.4	252.3	312.3

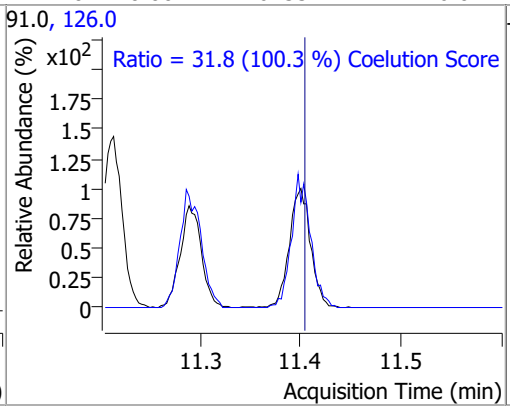


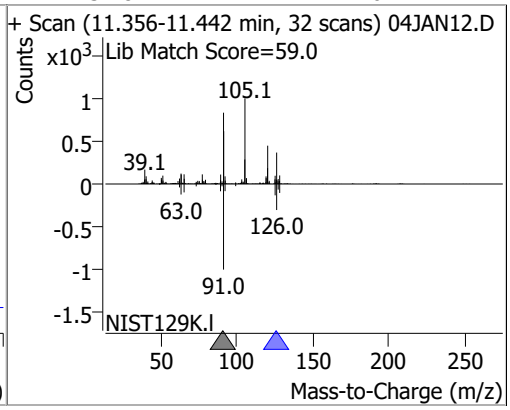




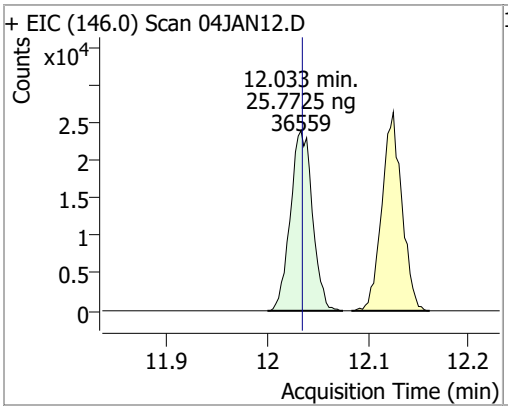
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	24.3936	11.40	0.00	61551	126.0	31.8	1.7	61.7

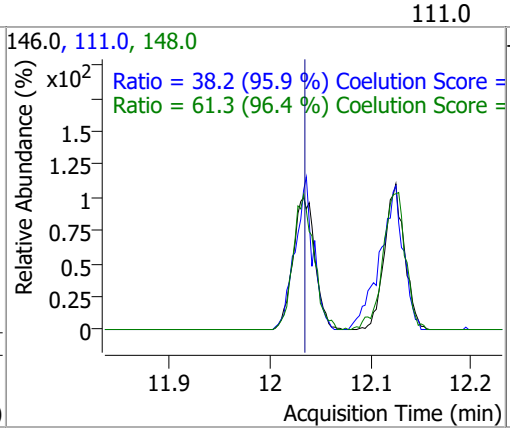


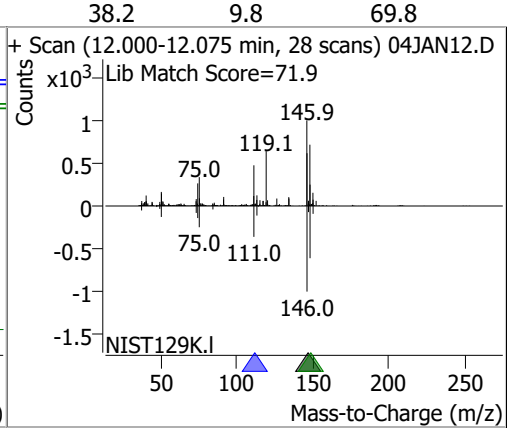




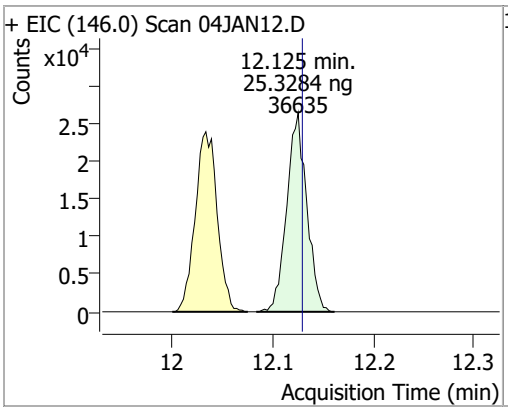
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	25.7725	12.03	0.00	36559	148.0	61.3	33.6	93.6
					111.0	38.2	9.8	69.8

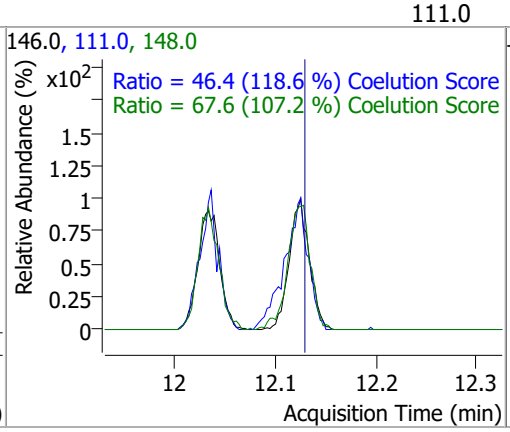


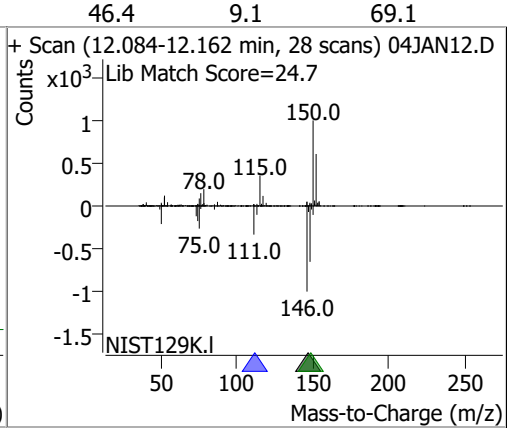




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	25.3284	12.13	0.00	36635	148.0	67.6	33.1	93.1
					111.0	46.4	9.1	69.1

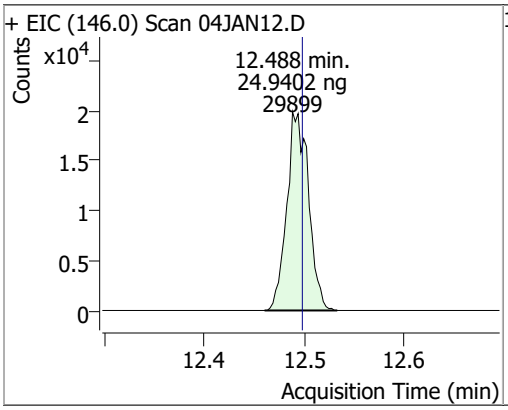


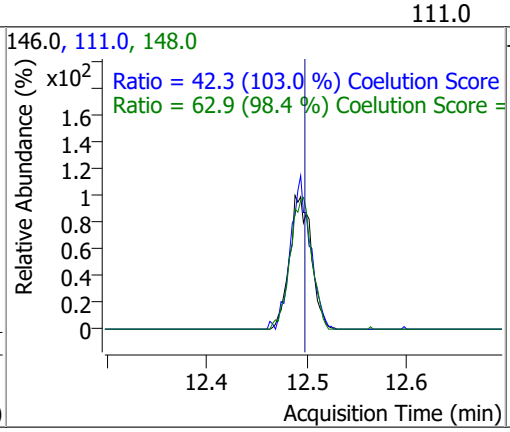


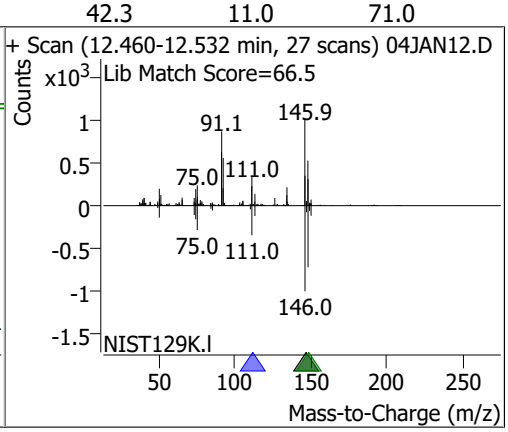


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	24.9402	12.49	-0.01	29899	148.0	62.9	33.9	93.9
					111.0	42.3	11.0	71.0



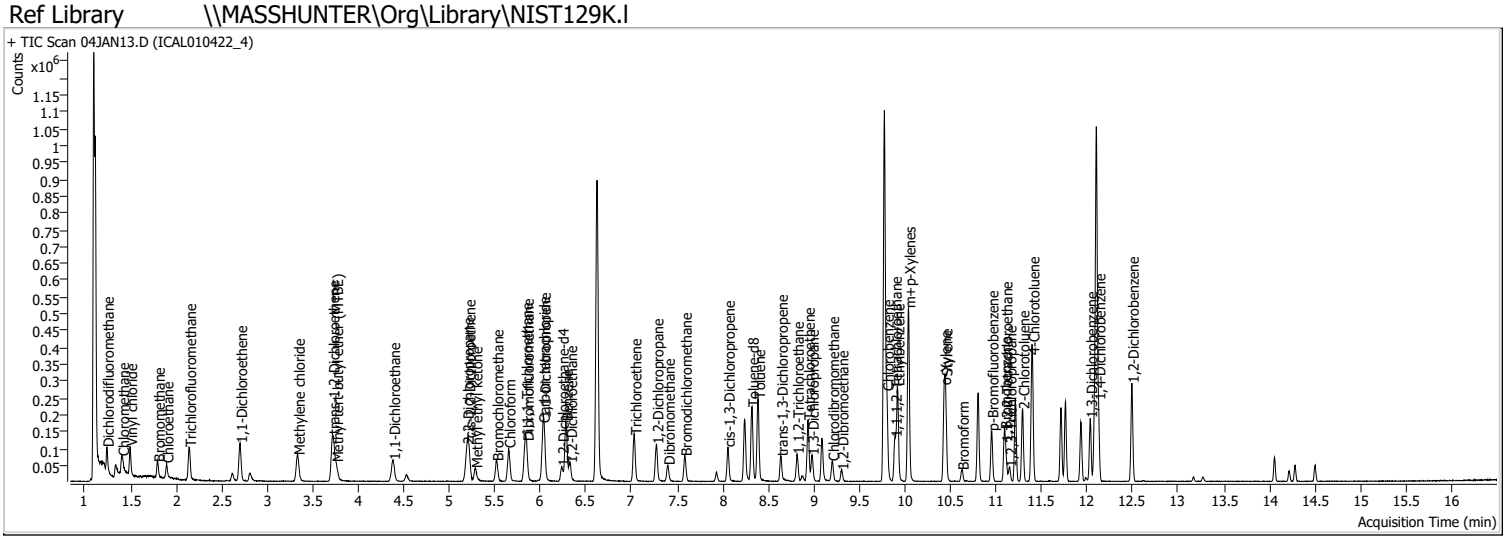




Quantitation Results Report (QT Reviewed)

Agilent | Trusted Answers

Data File	04JAN13.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/4/2022 4:55:32 PM
Sample Name	ICAL010422_4	Instrument	VOA5975C
Vial	13	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

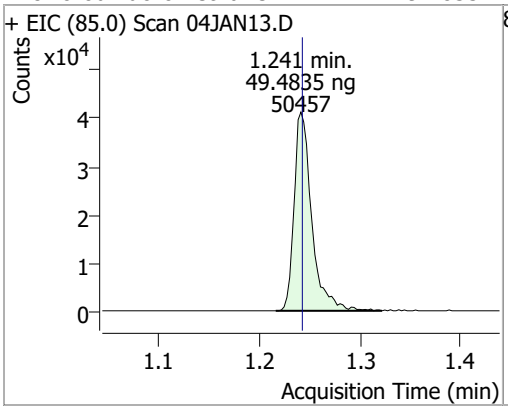


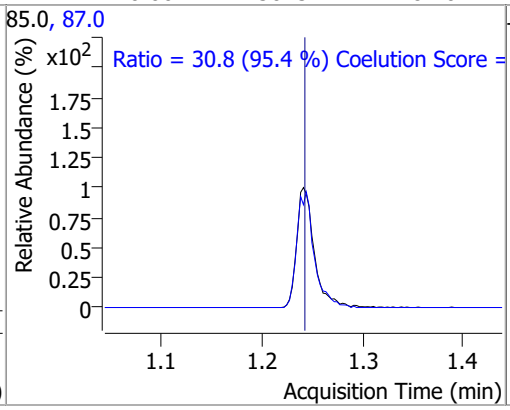
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.623	96.0	778120	250.0000	ng		0.000
M Chlorobenzene-d5	9.774	82.0	300356	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	248636	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.848	113.0	35309	48.1661	ng		0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 19.27%		*	
S 1,2-Dichloroethane-d4	6.233	67.0	15238	48.1252	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 19.25%		*	
S Toluene-d8	8.319	98.0	136453	47.1441	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 18.86%		*	
S p-Bromofluorobenzene	10.951	95.0	42506	46.6647	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 18.67%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.241	85.0	50457	49.4835	ng		97
T Chloromethane	1.408	50.0	61632	49.7983	ng		99
T Vinyl chloride	1.495	62.0	54521	48.9580	ng		96
T Bromomethane	1.799	96.0	23699	47.5921	ng		100
T Chloroethane	1.897	64.0	25484	46.2243	ng		98
T Trichlorofluoromethane	2.145	101.0	68163	49.3128	ng		98
T 1,1-Dichloroethene	2.702	96.0	38253	48.8056	ng		100
T Methylene chloride	3.335	49.0	58282	50.4421	ng		99
T trans-1,2-Dichloroethene	3.717	96.0	39596	49.5178	ng		99
T Methyl tert-butyl ether (MTBE)	3.757	73.0	49126	47.5301	ng		100
T 1,1-Dichloroethane	4.381	63.0	73205	49.1828	ng		100
T 2,2-Dichloropropane	5.193	77.0	56189	50.3804	ng		100
T cis-1,2-Dichloroethene	5.209	96.0	39251	48.4154	ng		99
T Methyl ethyl ketone	5.285	43.0	52648	479.4296	ng		99
T Bromochloromethane	5.516	128.0	17338	51.6233	ng		96
T Chloroform	5.650	83.0	71403	48.2031	ng		99

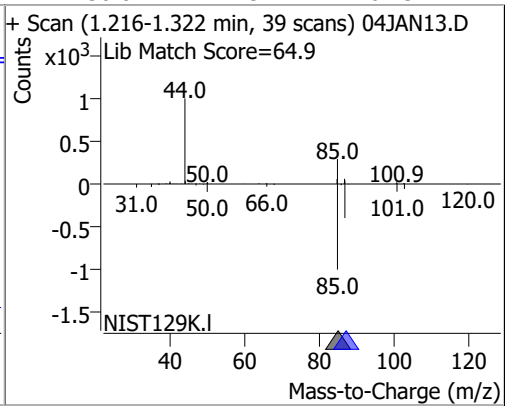
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.834	97.0	67007	48.2688	ng	98
T Carbon tetrachloride	6.026	117.0	65313	47.7520	ng	98
T 1,1-Dichloropropene	6.035	75.0	56376	47.7627	ng	99
T Benzene	6.277	78.0	148727	48.0054	ng	100
T 1,2-Dichloroethane	6.325	62.0	41058	48.9880	ng	97
T Trichloroethene	7.030	95.0	42682	47.1189	ng	98
T 1,2-Dichloropropane	7.273	63.0	37870	47.5273	ng	96
T Dibromomethane	7.396	93.0	15989	47.4844	ng	97
T Bromodichloromethane	7.585	83.0	43900	47.2409	ng	97
T cis-1,3-Dichloropropene	8.057	75.0	48886	46.5283	ng	97
T Toluene	8.388	92.0	91915	47.0116	ng	100
T trans-1,3-Dichloropropene	8.639	75.0	35179	47.0378	ng	100
T 1,1,2-Trichloroethane	8.815	83.0	18884	48.4759	ng	99
T Tetrachloroethene	8.935	163.8	36925	46.2932	ng	97
T 1,3-Dichloropropane	8.980	76.0	37457	48.8841	ng	98
T Chlorodibromomethane	9.203	129.0	28153	46.2411	ng	99
T 1,2-Dibromoethane	9.303	107.0	21037	49.3889	ng	93
T Chlorobenzene	9.802	112.0	101452	47.3959	ng	99
T 1,1,1,2-Tetrachloroethane	9.889	131.0	35544	47.5029	ng	99
T Ethylbenzene	9.917	91.0	173769	46.8079	ng	99
T m+p-Xylenes	10.039	106.0	133498	92.5347	ng	98
T o-Xylene	10.430	106.0	61016	47.5086	ng	98
T Styrene	10.444	104.0	96576	46.7052	ng	100
T Bromoform	10.625	172.5	16073	50.5170	ng	96
T Bromobenzene	11.093	156.0	38282	47.5759	ng	98
T 1,1,2,2-Tetrachloroethane	11.105	83.0	22514	48.6124	ng	99
T 1,2,3-Trichloropropane	11.146	110.0	6096	49.1924	ng	97
T 2-Chlorotoluene	11.289	126.0	37987	47.4466	ng	99
T 4-Chlorotoluene	11.400	91.0	126308	48.3865	ng	100
T 1,3-Dichlorobenzene	12.033	146.0	69539	47.3853	ng	99
T 1,4-Dichlorobenzene	12.125	146.0	71841	48.0106	ng	97
T 1,2-Dichlorobenzene	12.491	146.0	60213	48.5498	ng	98

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

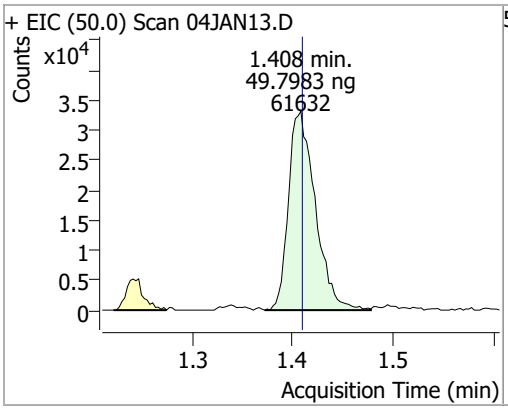
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	49.4835	1.24	0.00	50457	87.0	30.8	2.3	62.3

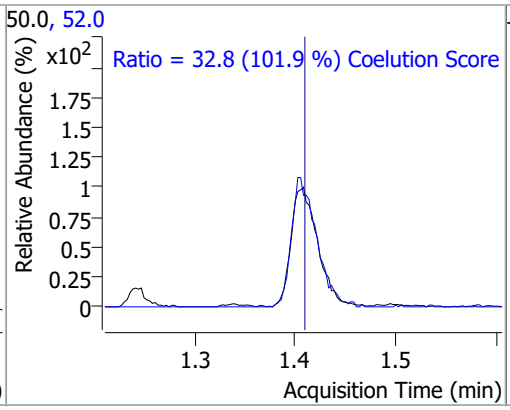


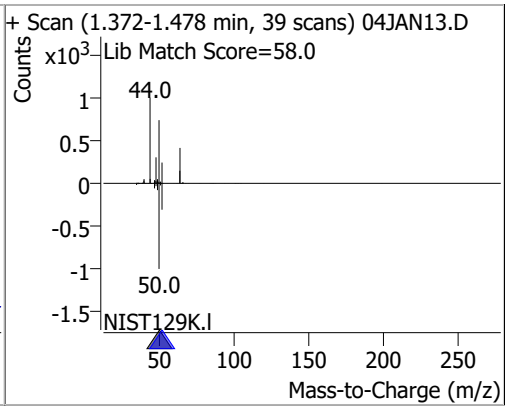




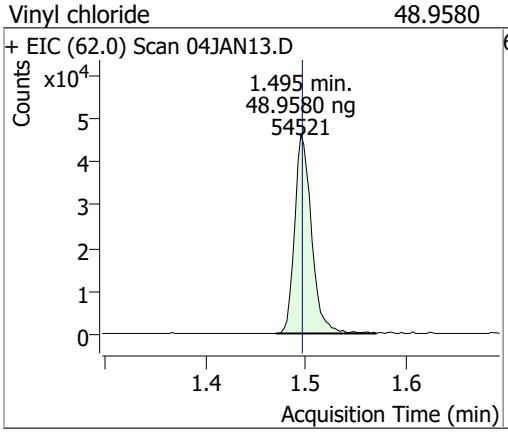
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	49.7983	1.41	0.00	61632	52.0	32.8	2.1	62.1

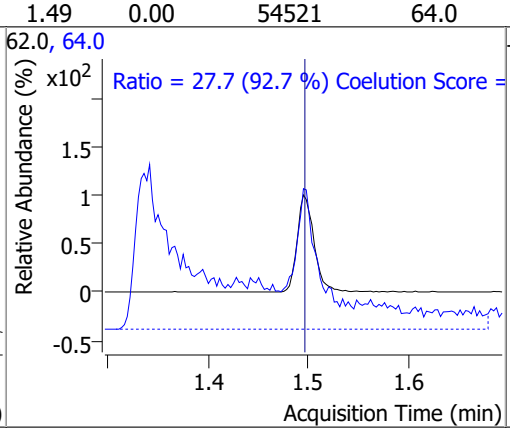


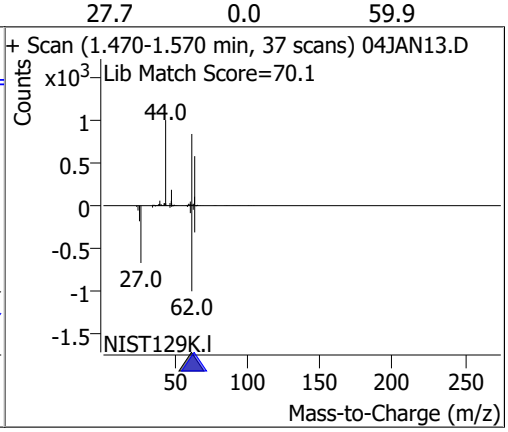




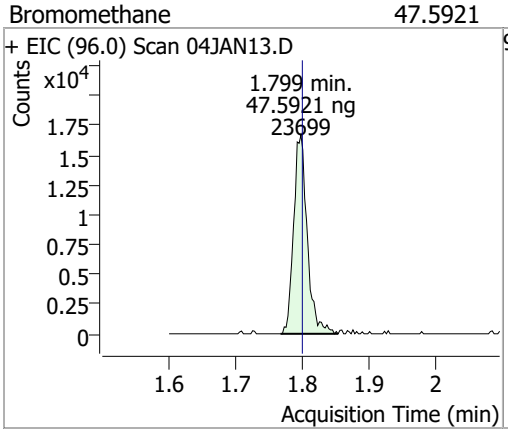
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	48.9580	1.49	0.00	54521	64.0	27.7	0.0	59.9

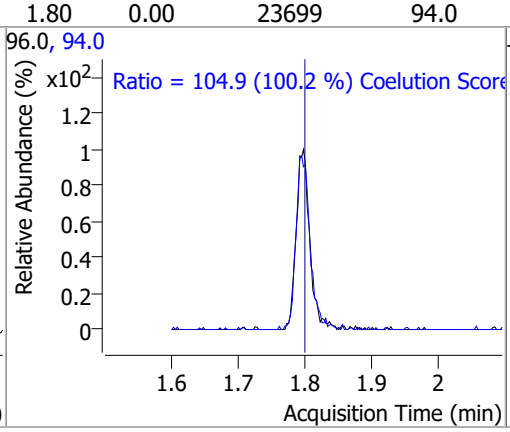


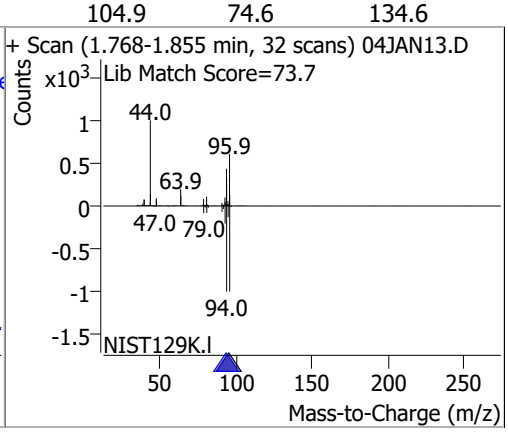




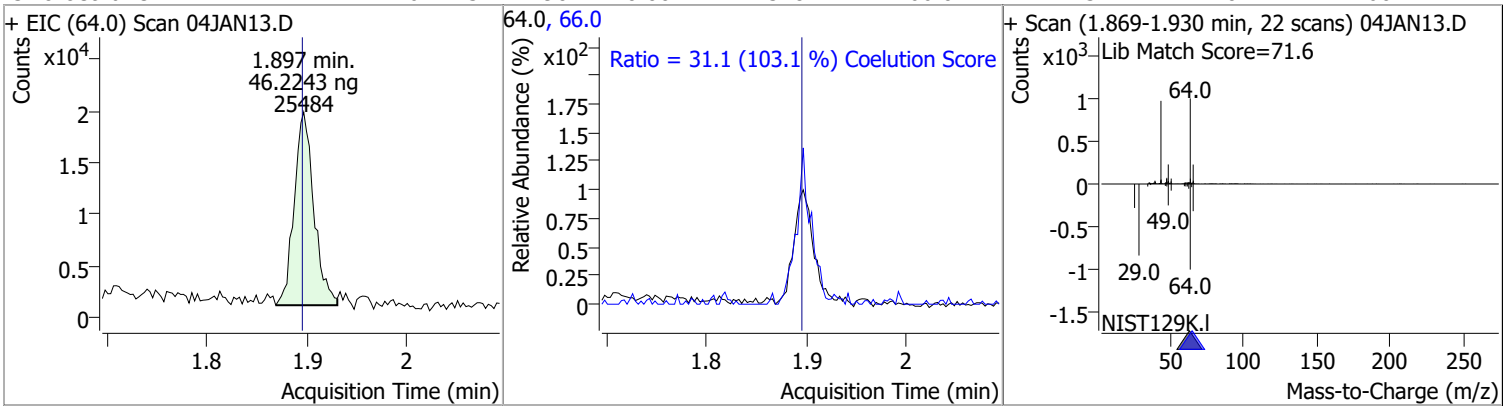
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	47.5921	1.80	0.00	23699	94.0	104.9	74.6	134.6



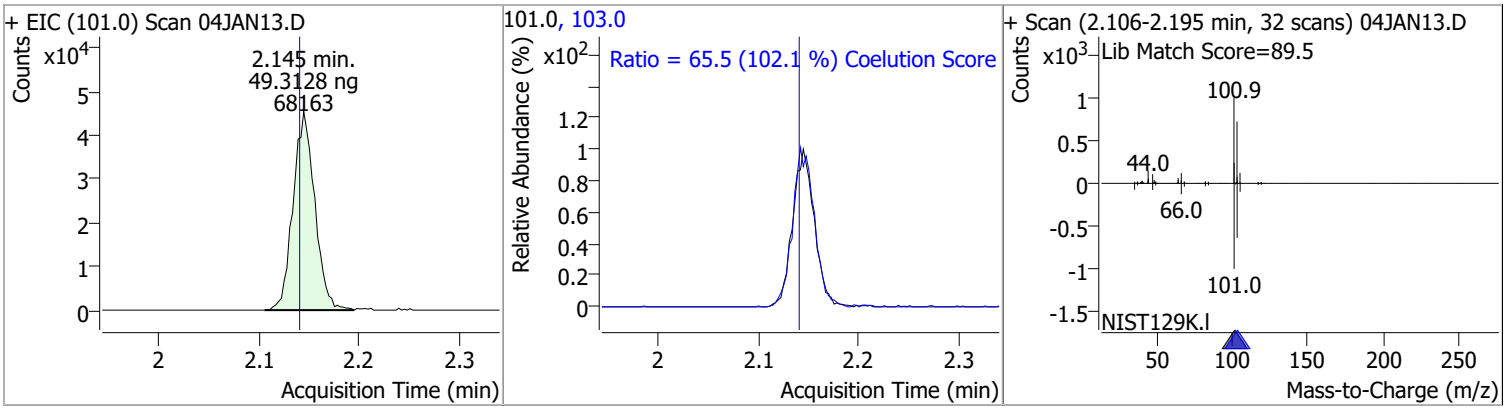




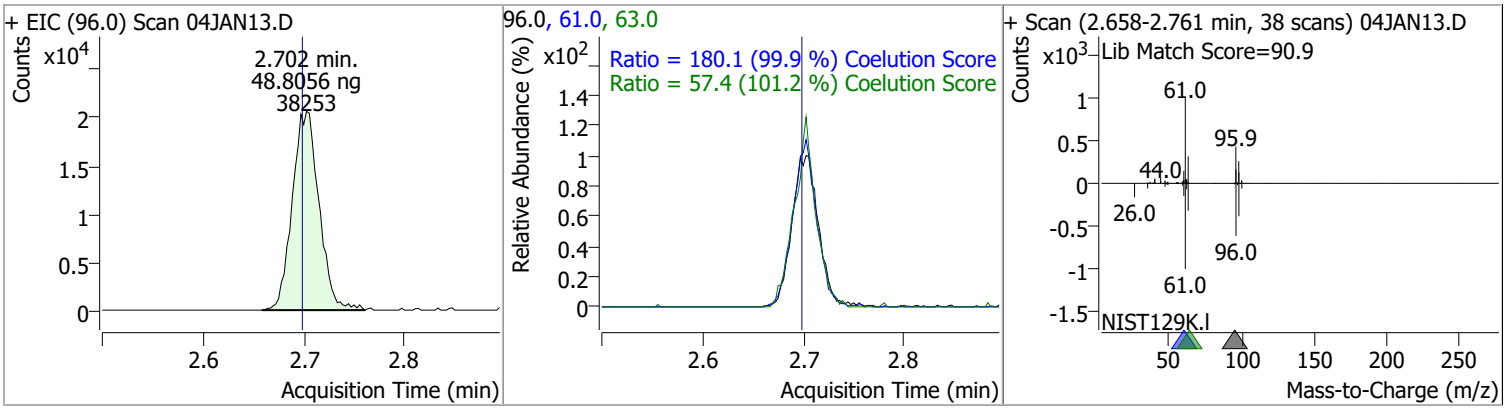
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	46.2243	1.90	0.00	25484	66.0	31.1	0.1	60.1



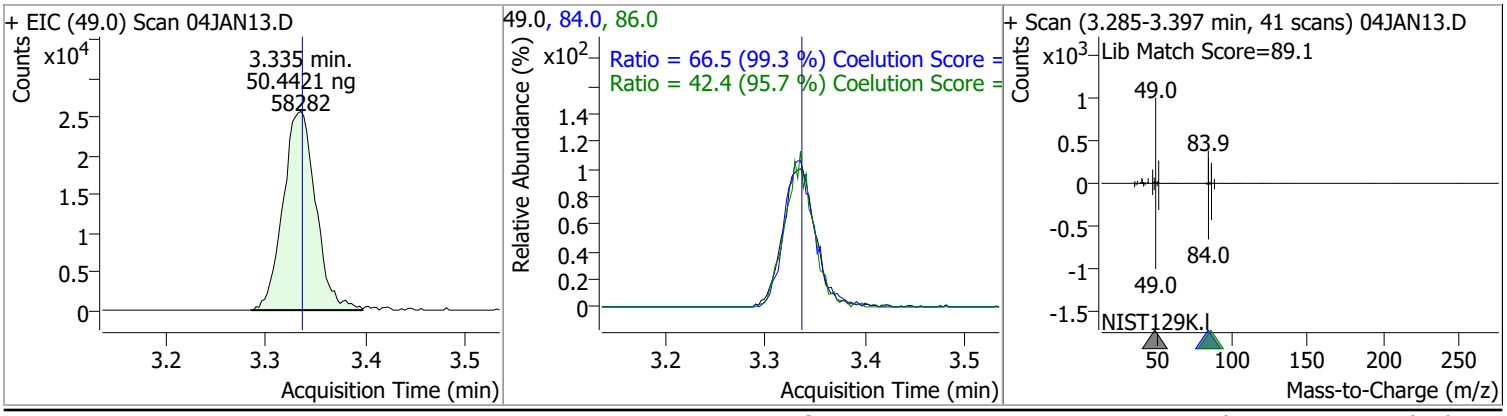
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	49.3128	2.14	0.00	68163	103.0	65.5	34.2	94.2



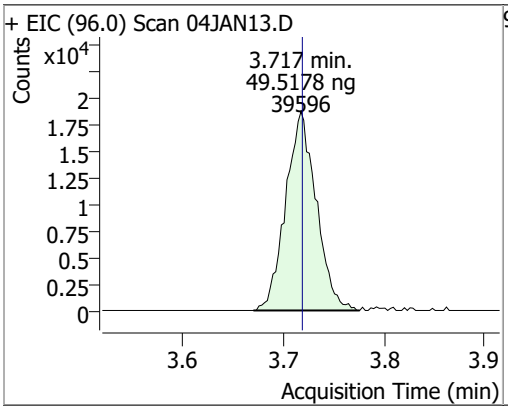
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	48.8056	2.70	0.01	38253	61.0	180.1	150.3	210.3
					63.0	57.4	26.7	86.7

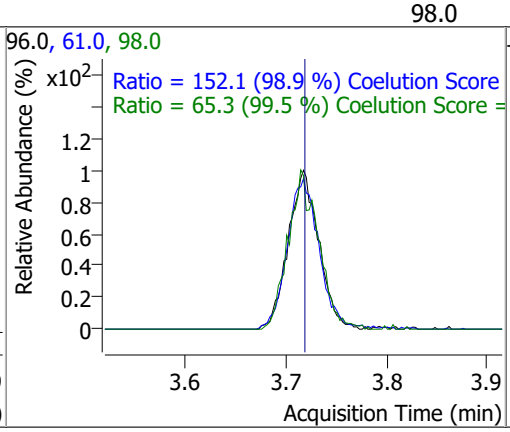


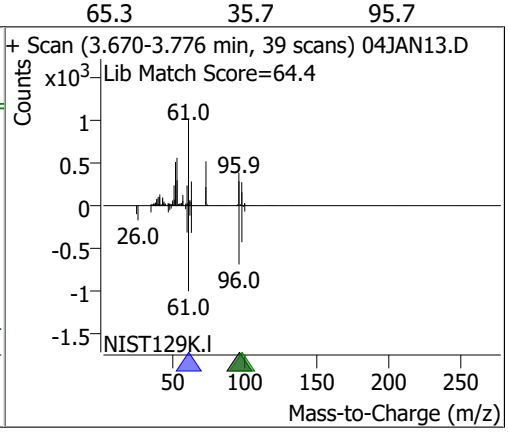
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	50.4421	3.34	0.00	58282	84.0	66.5	36.9	96.9
					86.0	42.4	14.3	74.3



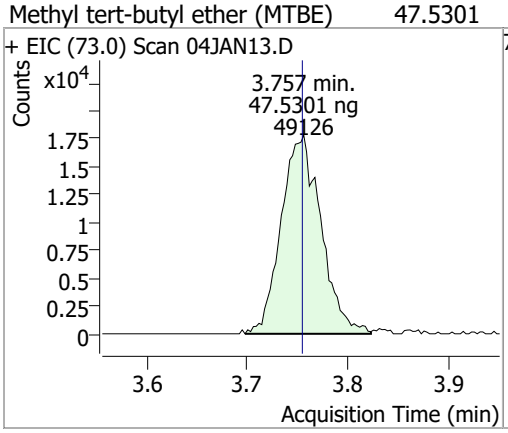
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	49.5178	3.72	0.00	39596	61.0	152.1	123.9	183.9
					98.0	65.3	35.7	95.7

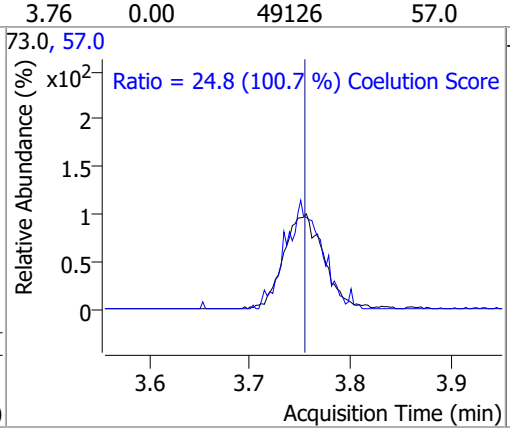


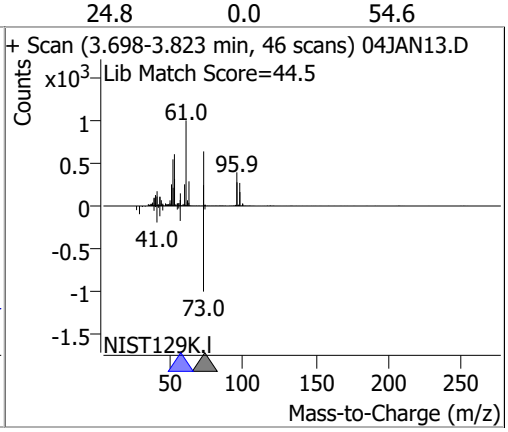




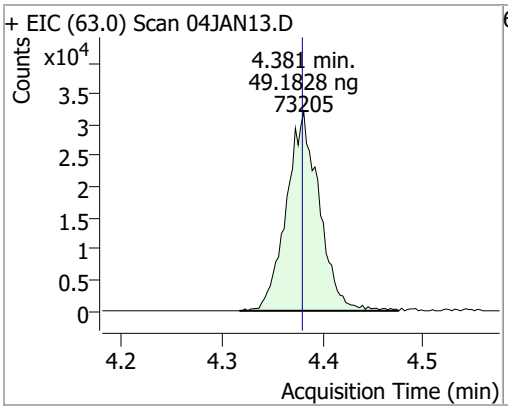
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	47.5301	3.76	0.00	49126	57.0	24.8	0.0	54.6

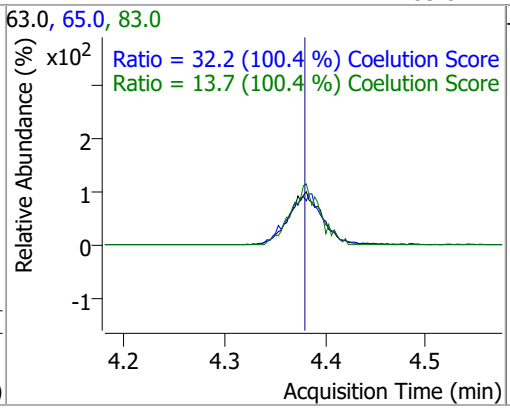


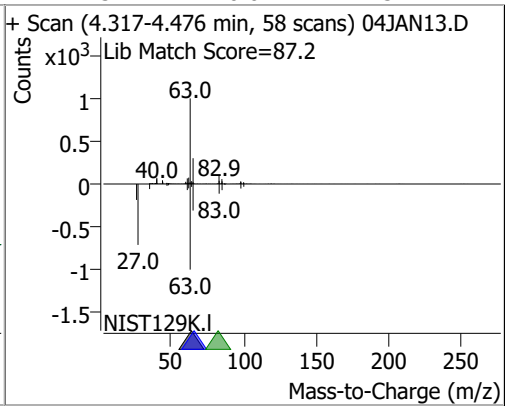




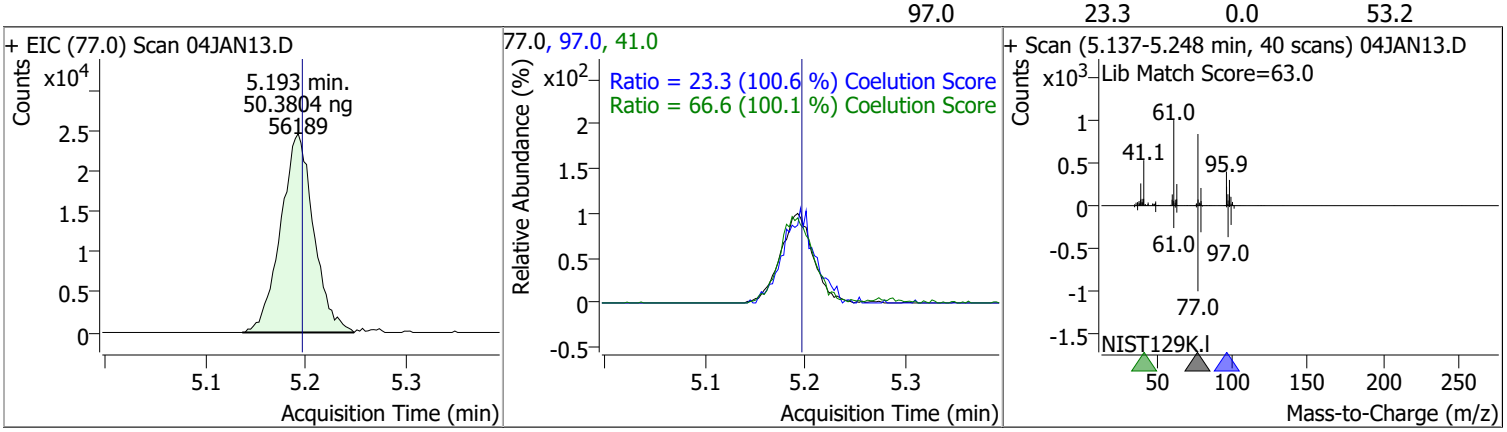
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	49.1828	4.38	0.00	73205	65.0	32.2	2.1	62.1
					83.0	13.7	0.0	43.7



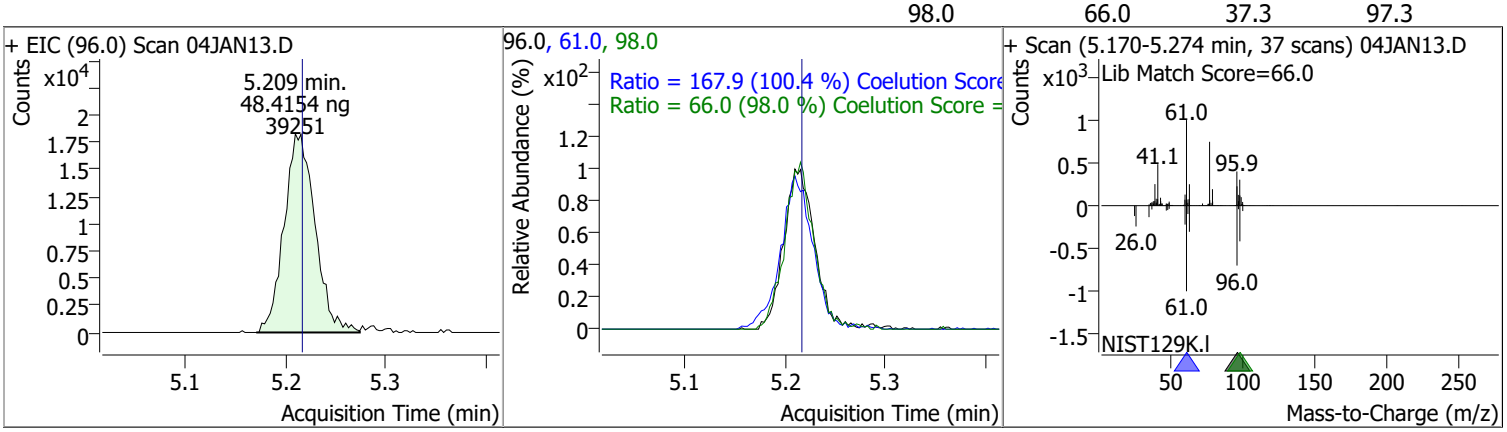




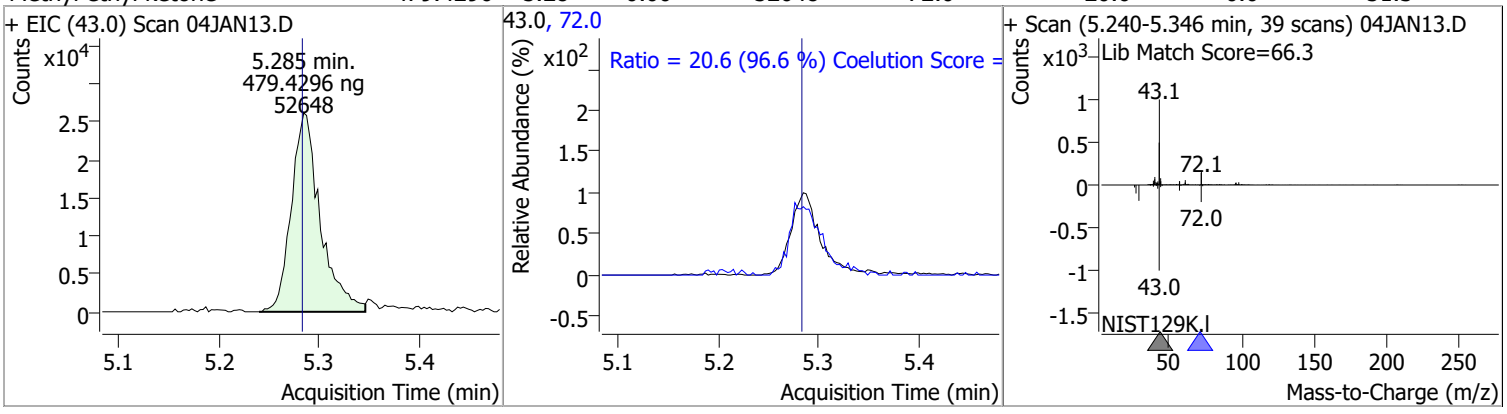
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	50.3804	5.19	0.00	56189	41.0	66.6	36.5	96.5
					97.0	23.3	0.0	53.2



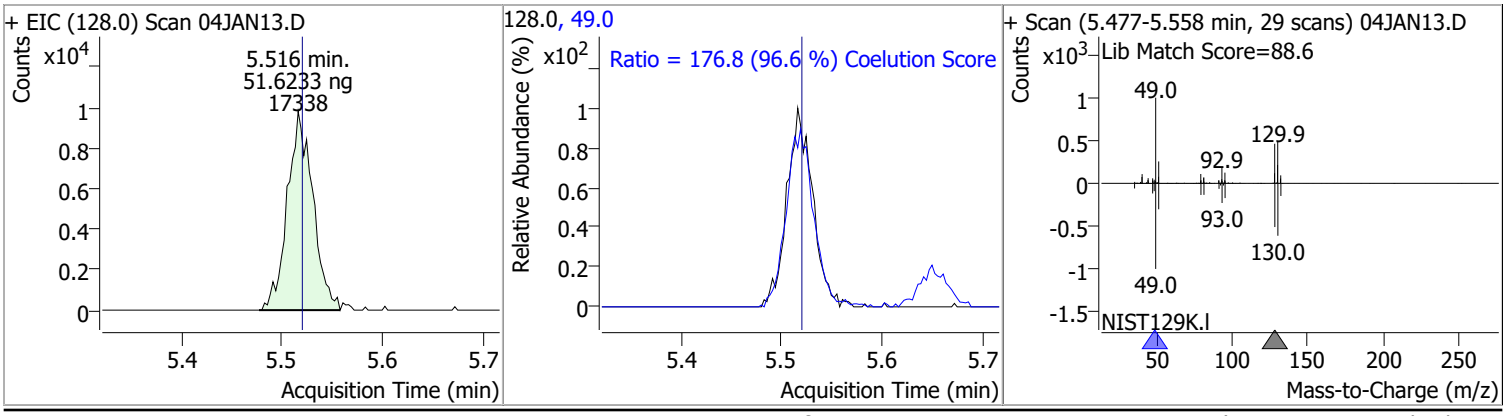
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	48.4154	5.21	-0.01	39251	61.0	167.9	137.2	197.2
					98.0	66.0	37.3	97.3



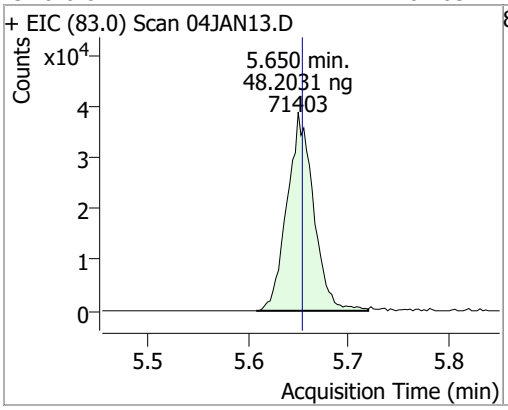
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	479.4296	5.28	0.00	52648	72.0	20.6	0.0	51.3

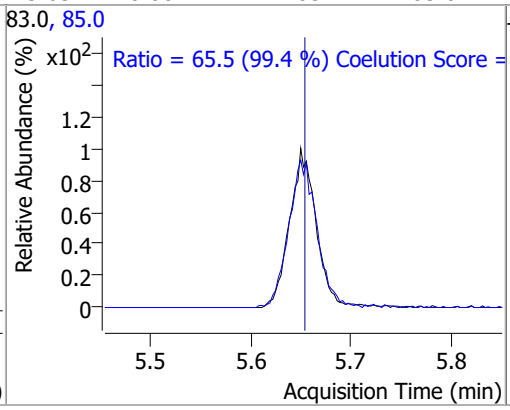


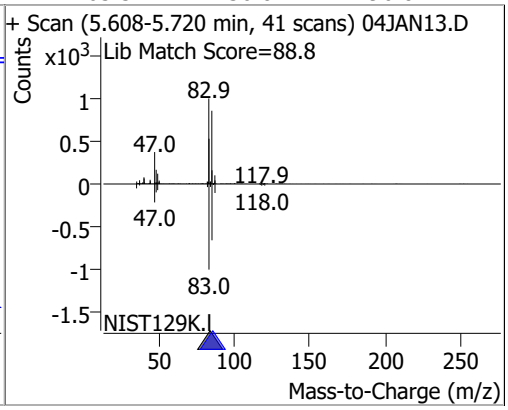
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	51.6233	5.52	0.00	17338	49.0	176.8	152.9	212.9



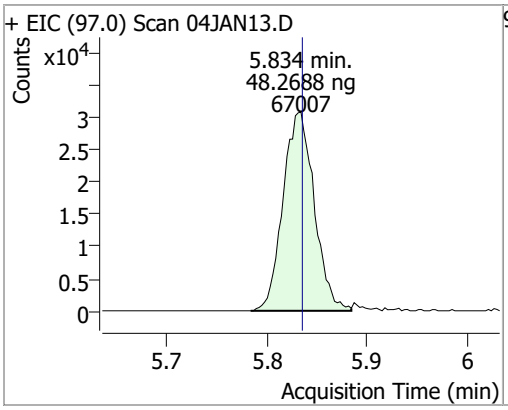
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	48.2031	5.65	0.00	71403	85.0	65.5	36.0	96.0

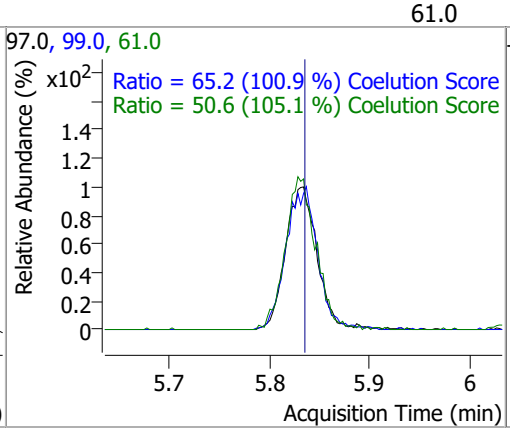


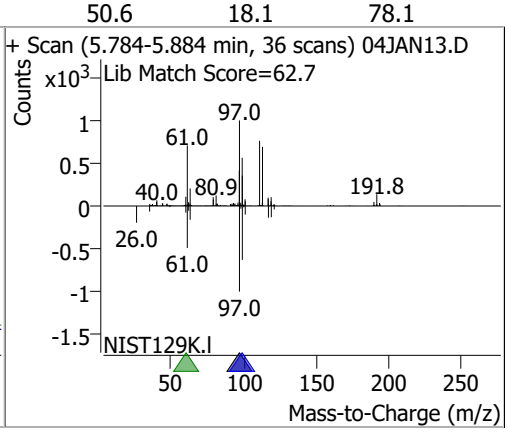




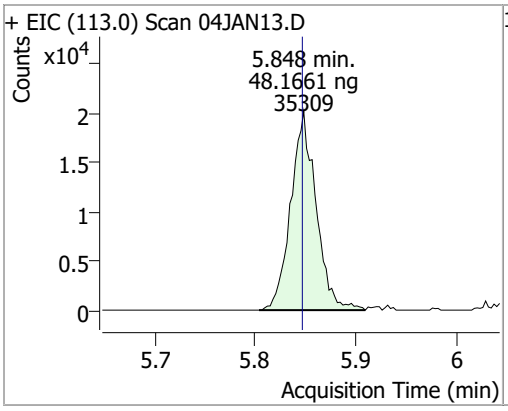
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	48.2688	5.83	0.00	67007	99.0	65.2	34.7	94.7
					61.0	50.6	18.1	78.1

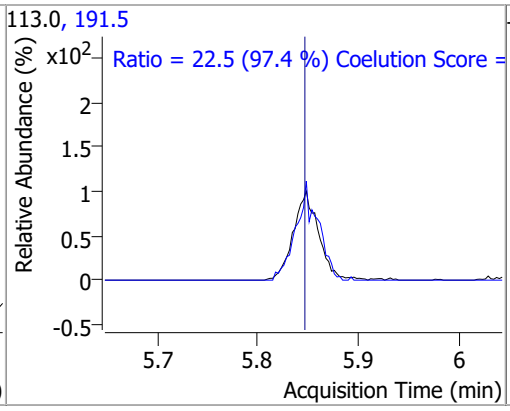


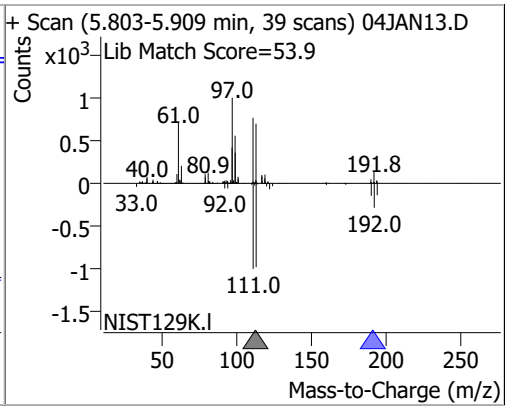




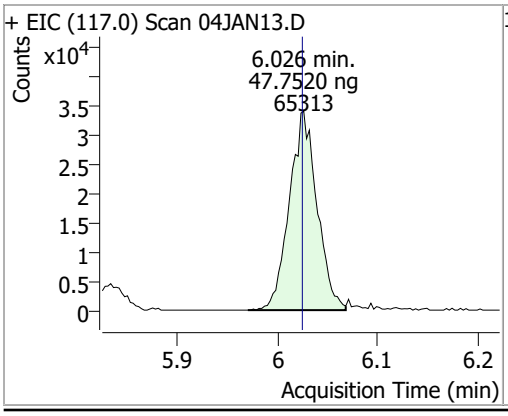
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	48.1661	5.85	0.00	35309	191.5	22.5	0.0	53.1

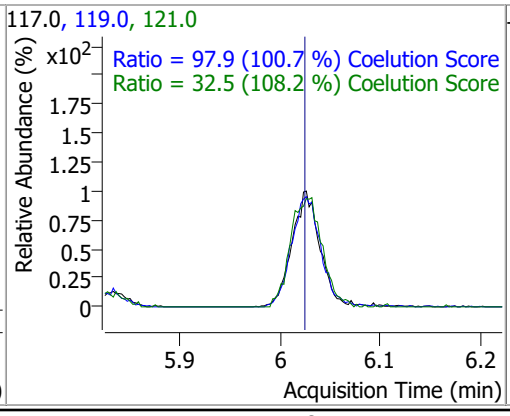


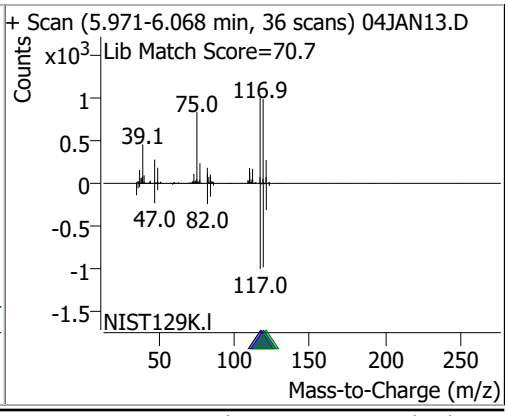




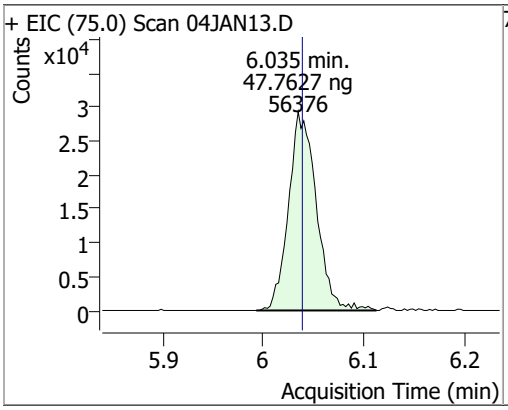
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	47.7520	6.03	0.00	65313	119.0	97.9	67.2	127.2
					121.0	32.5	0.1	60.1

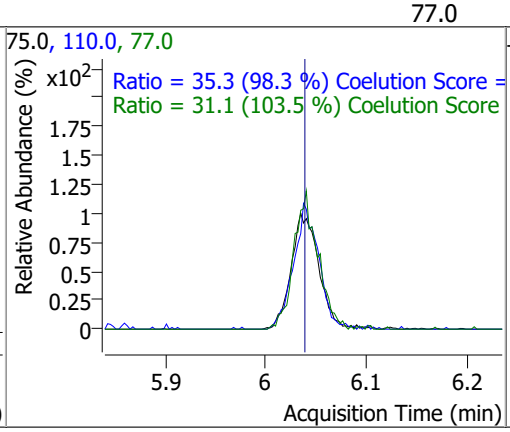


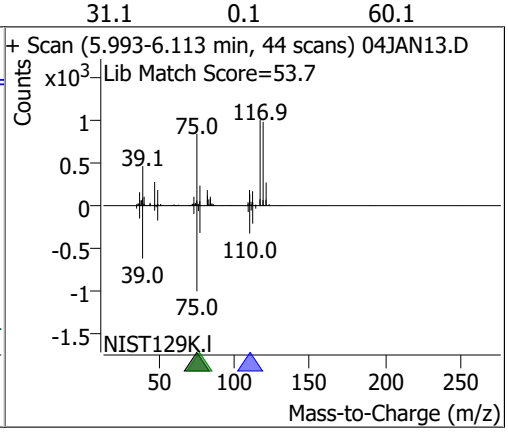




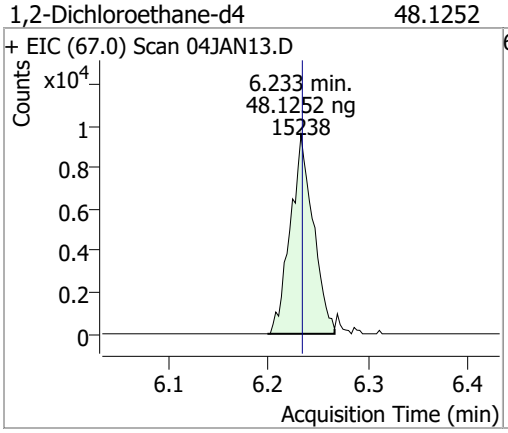
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	47.7627	6.03	0.00	56376	110.0	35.3	5.9	65.9
					77.0	31.1	0.1	60.1

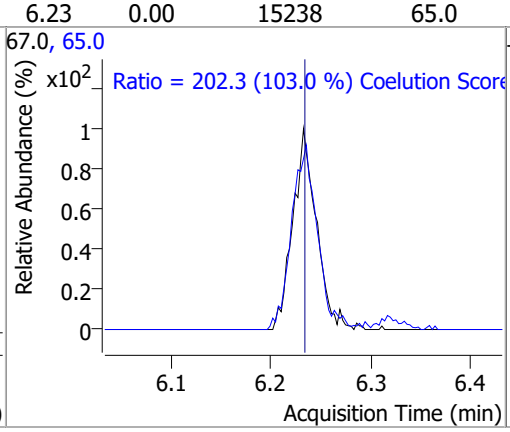


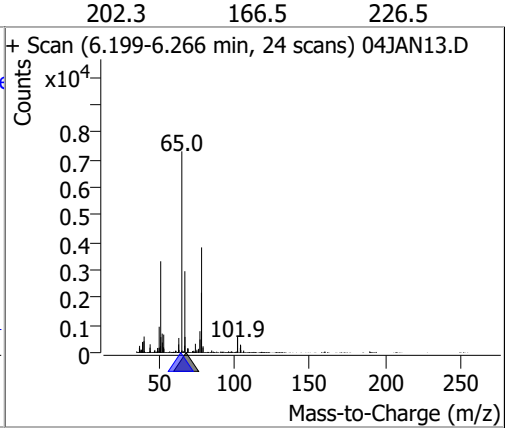




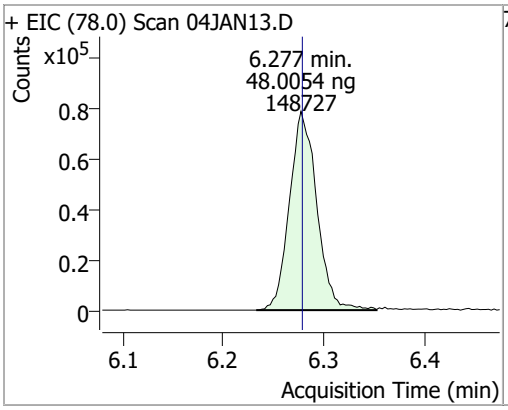
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	48.1252	6.23	0.00	15238	65.0	202.3	166.5	226.5

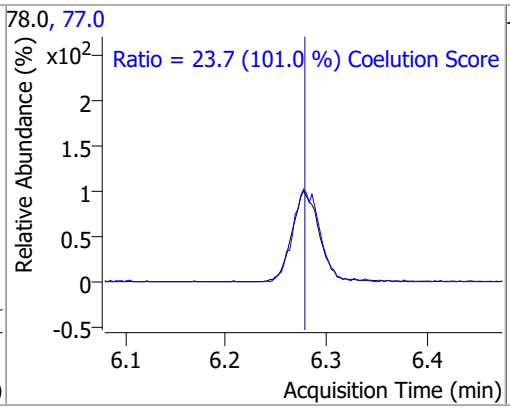


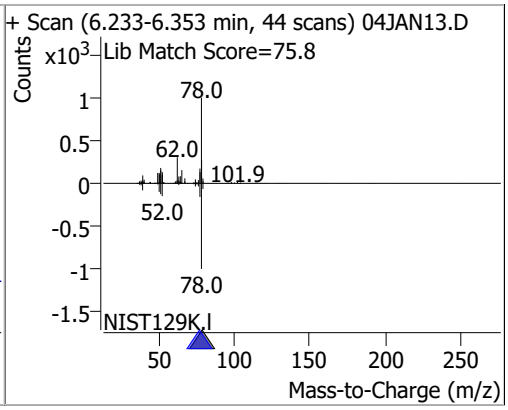




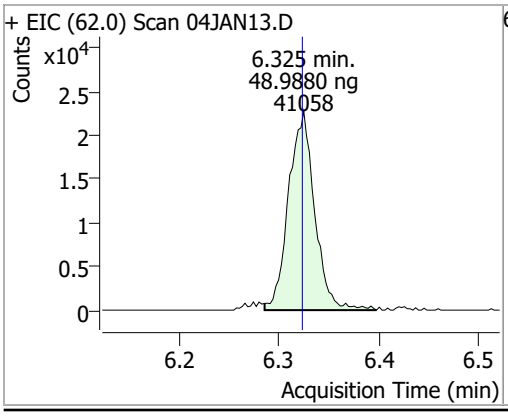
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	48.0054	6.28	0.00	148727	77.0	23.7	0.0	53.5

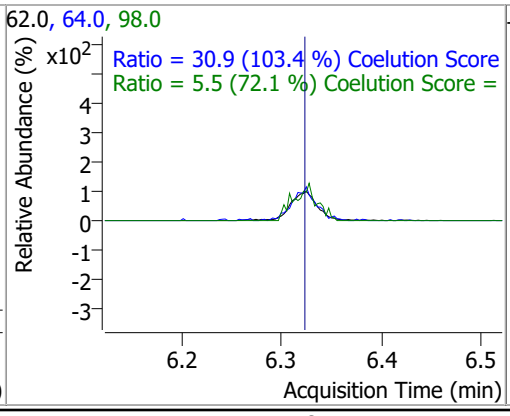


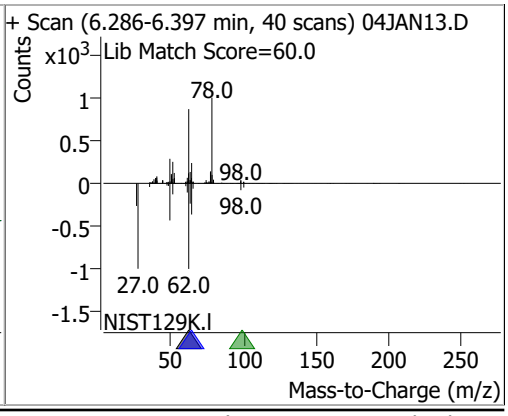




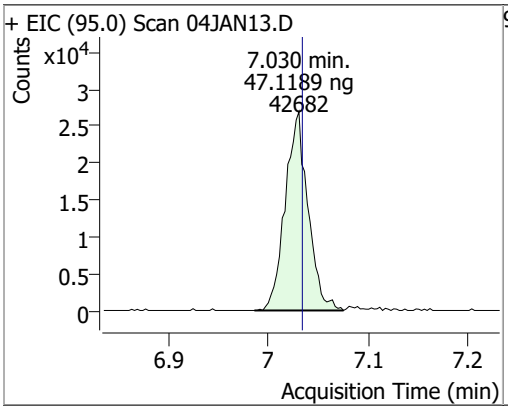
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	48.9880	6.32	0.00	41058	64.0	30.9	0.0	59.9
					98.0	5.5	0.0	37.6

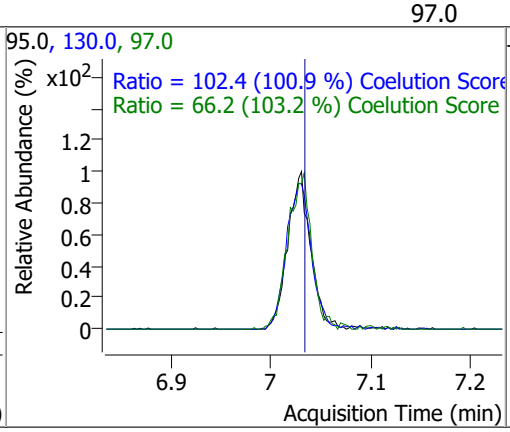


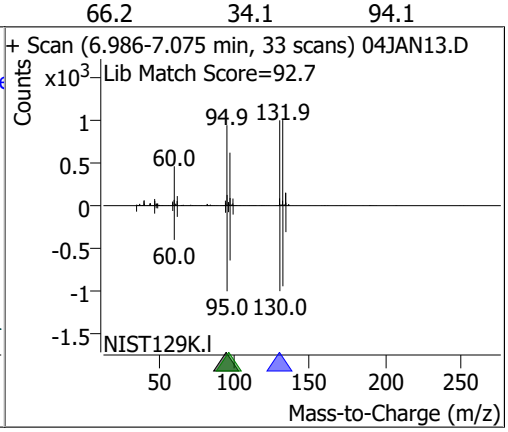




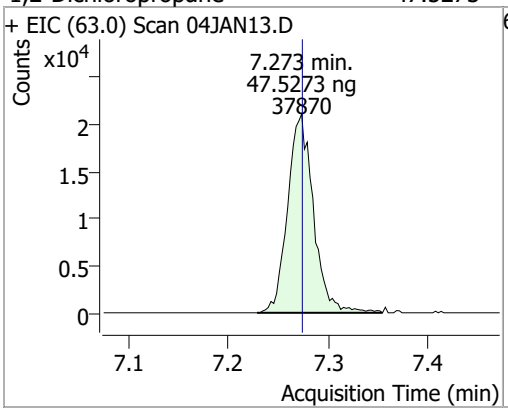
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	47.1189	7.03	0.00	42682	130.0	102.4	71.5	131.5
					97.0	66.2	34.1	94.1

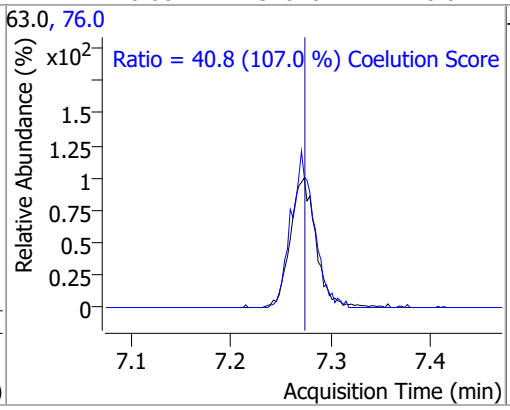


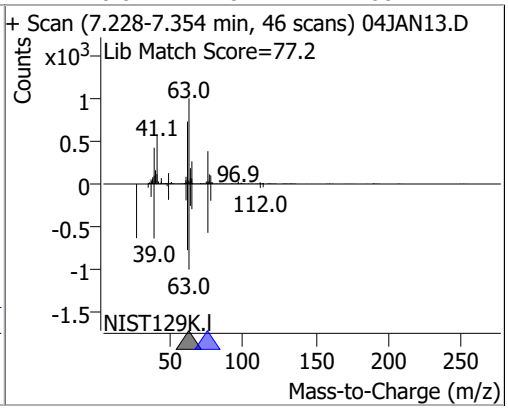




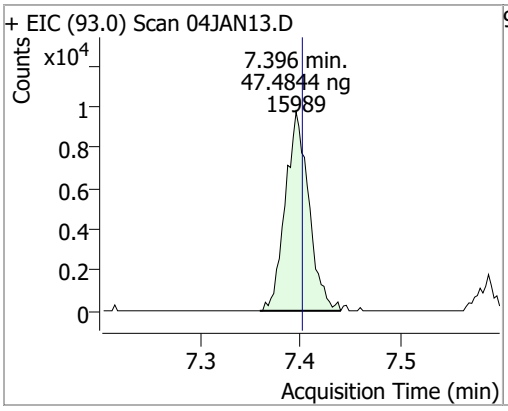
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	47.5273	7.27	0.00	37870	76.0	40.8	8.2	68.2

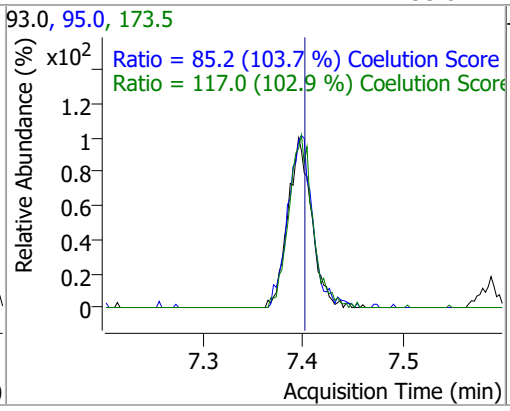


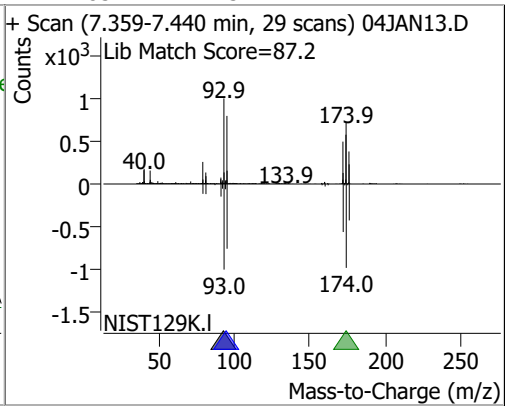




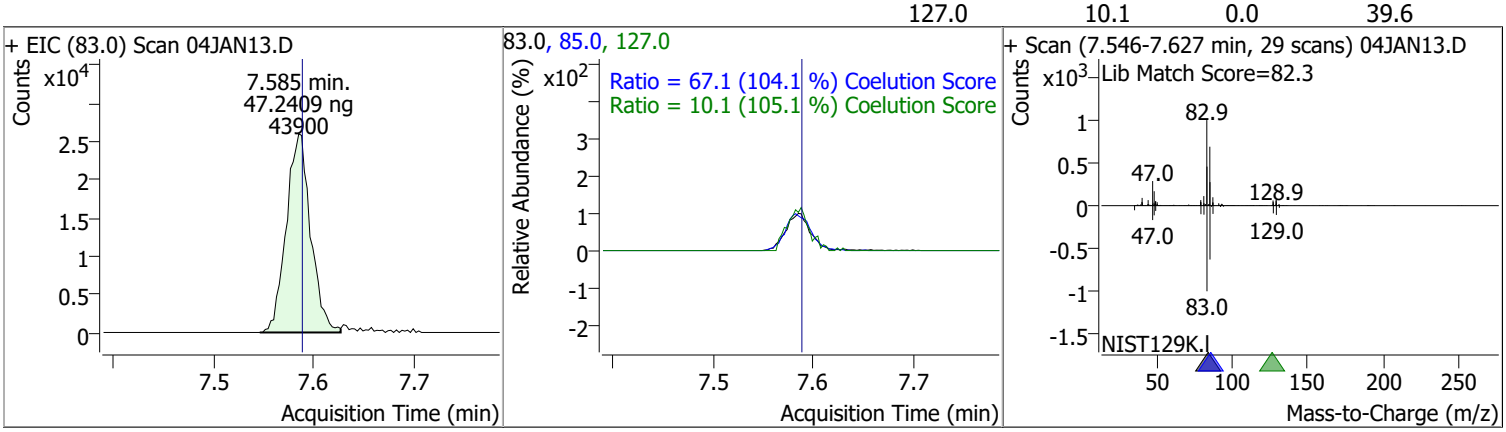
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	47.4844	7.40	0.00	15989	173.5	117.0	83.7	143.7
					95.0	85.2	52.2	112.2



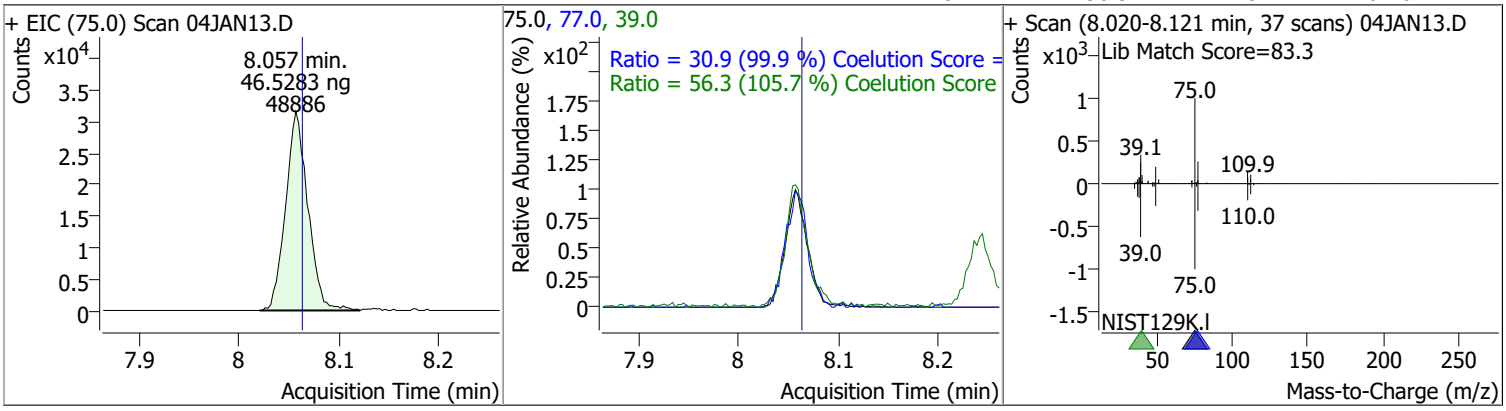




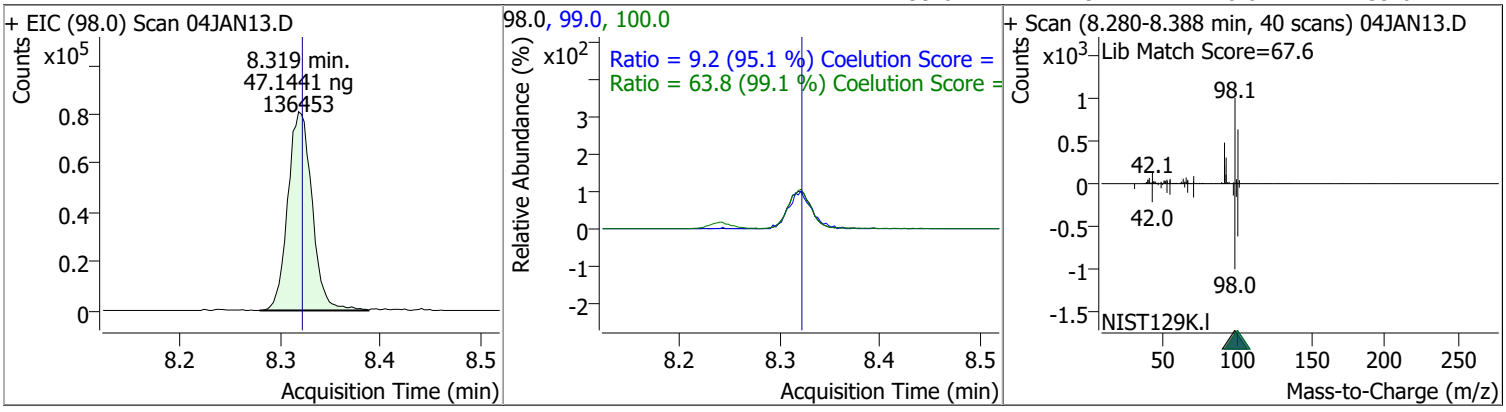
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	47.2409	7.59	0.00	43900	85.0	67.1	34.5	94.5
					127.0	10.1	0.0	39.6



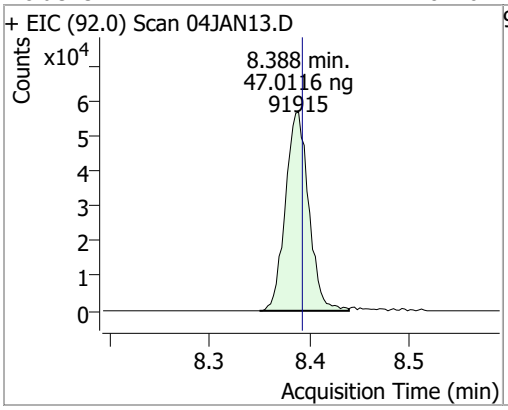
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	46.5283	8.06	0.00	48886	39.0	56.3	23.3	83.3
					77.0	30.9	1.0	61.0

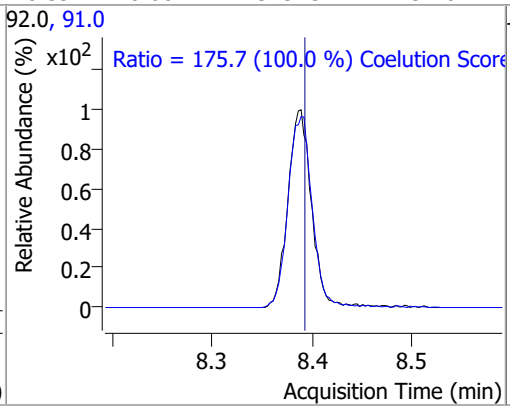


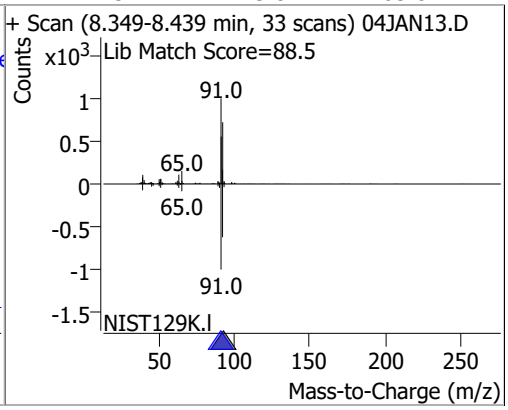
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	47.1441	8.32	0.00	136453	100.0	63.8	34.4	94.4
					99.0	9.2	0.0	39.6



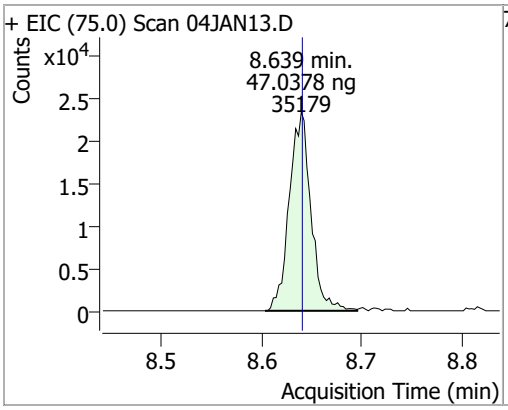
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	47.0116	8.39	0.00	91915	91.0	175.7	145.8	205.8

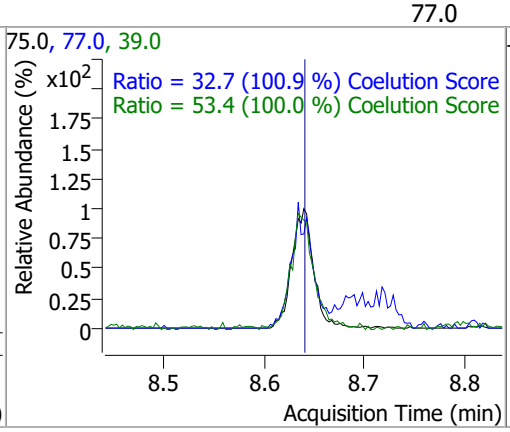


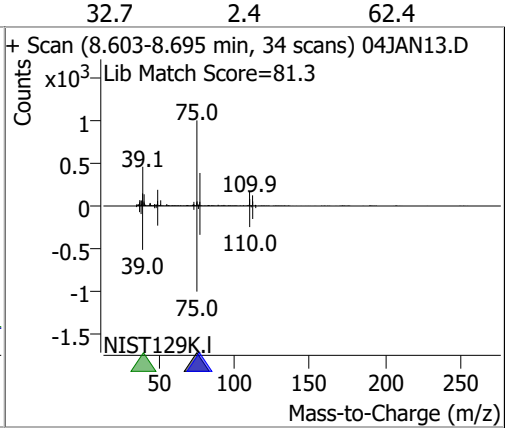




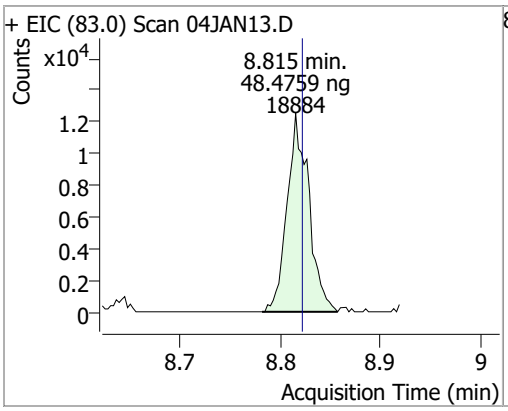
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	47.0378	8.64	0.00	35179	39.0	53.4	23.4	83.4
					77.0	32.7	2.4	62.4

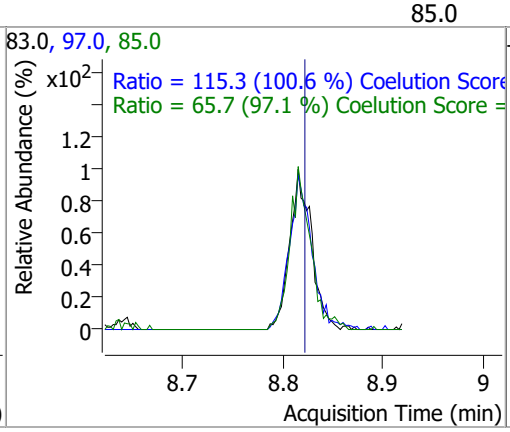


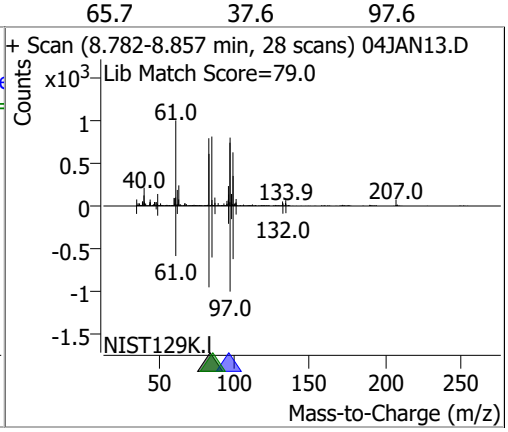




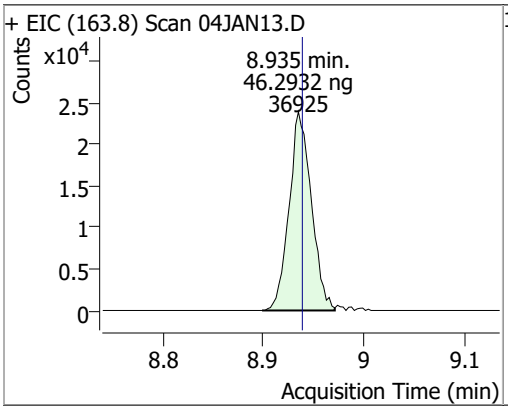
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	48.4759	8.82	0.00	18884	97.0	115.3	84.6	144.6
					85.0	65.7	37.6	97.6

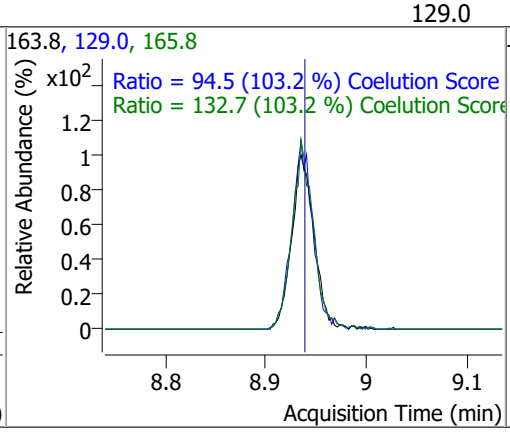


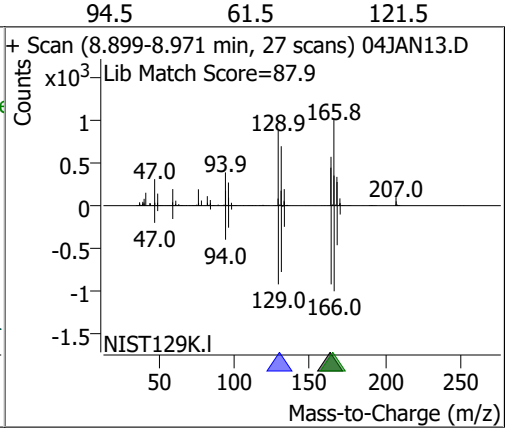




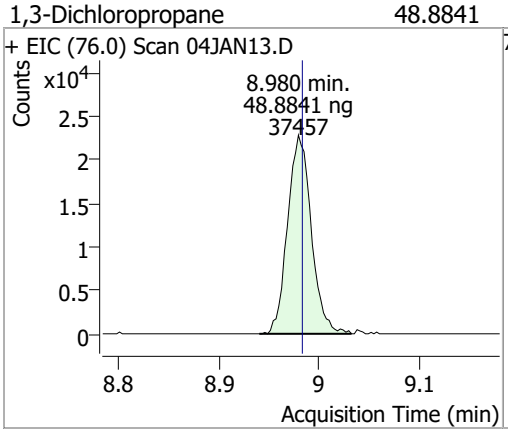
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	46.2932	8.94	0.00	36925	165.8	132.7	98.6	158.6
					129.0	94.5	61.5	121.5

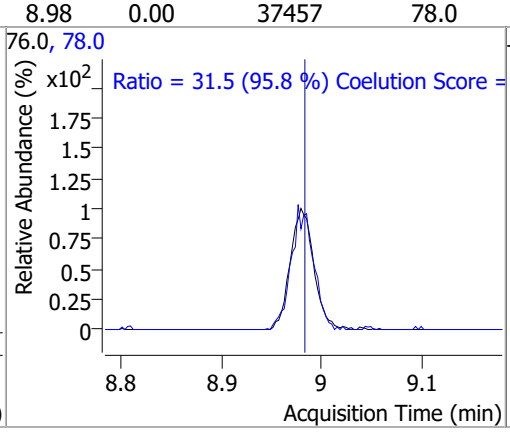


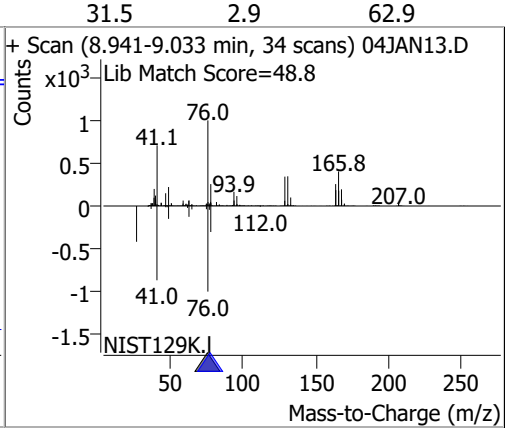




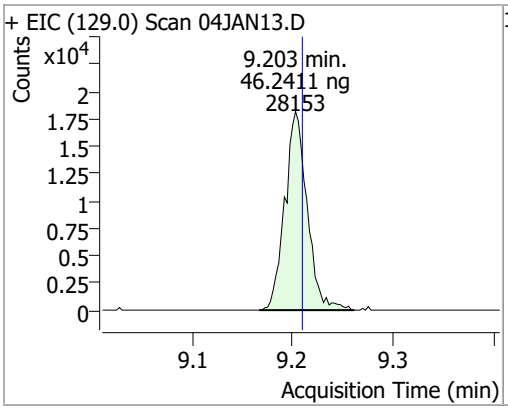
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	48.8841	8.98	0.00	37457	78.0	31.5	2.9	62.9

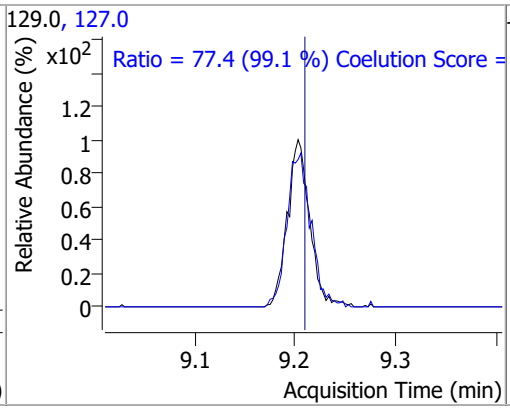


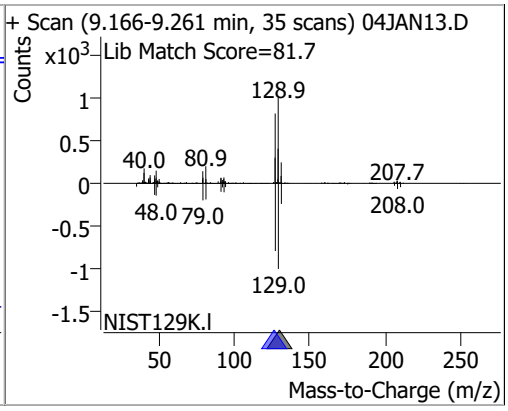




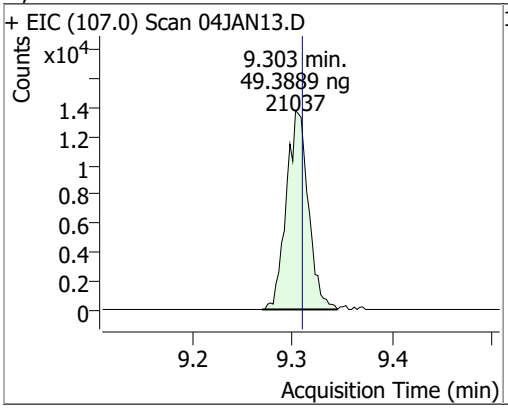
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	46.2411	9.20	0.00	28153	127.0	77.4	48.0	108.0

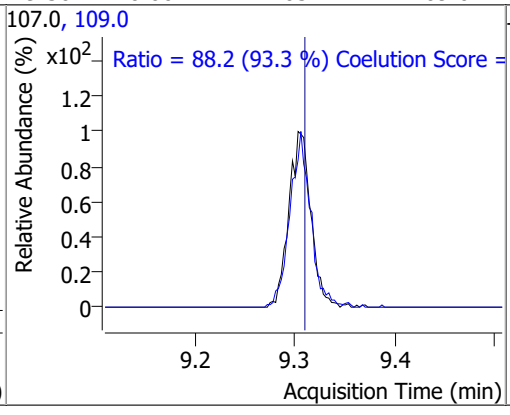


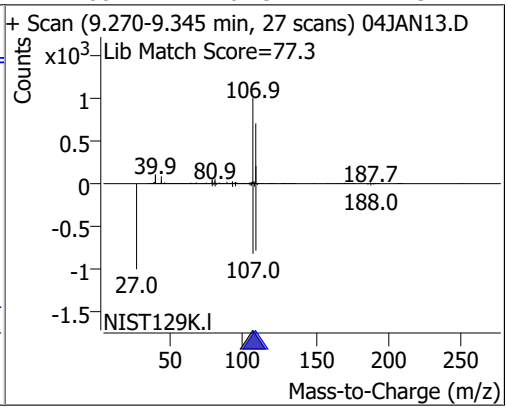




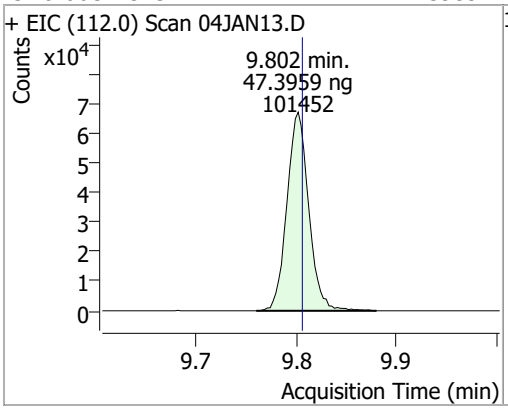
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	49.3889	9.30	0.00	21037	109.0	88.2	64.5	124.5

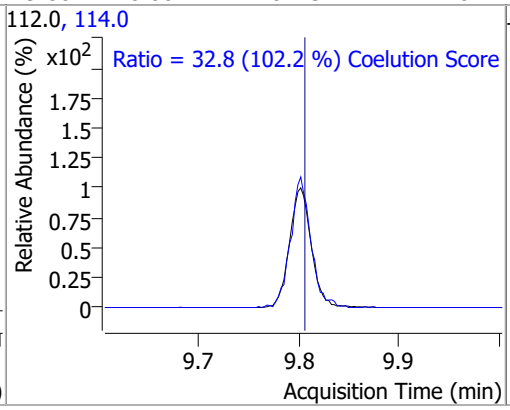


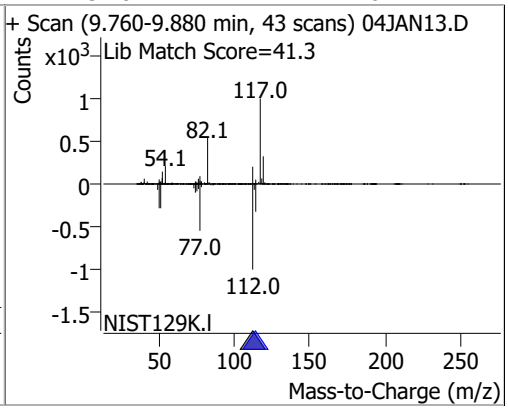




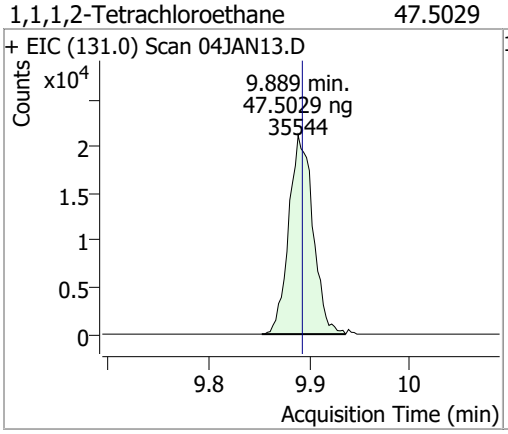
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	47.3959	9.80	0.00	101452	114.0	32.8	2.1	62.1

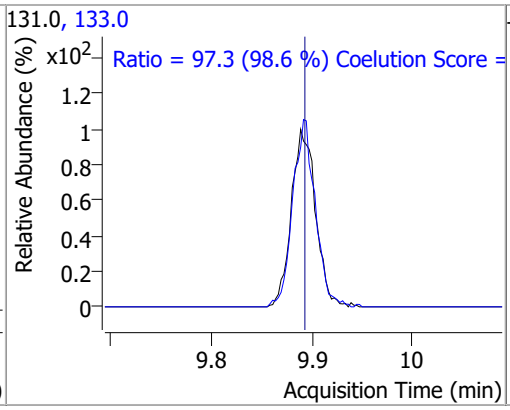


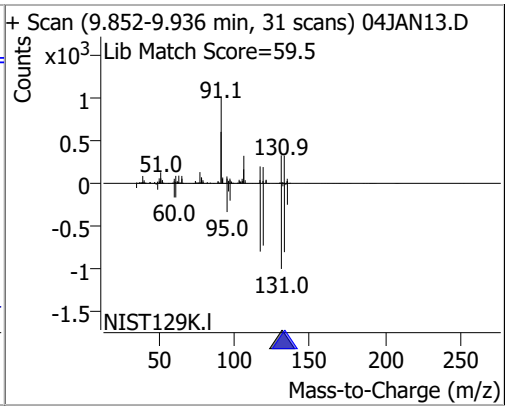




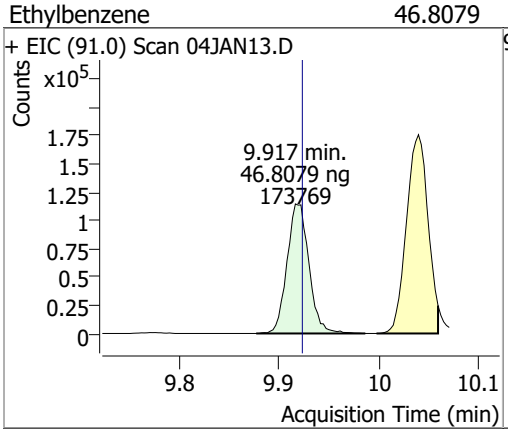
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	47.5029	9.89	0.00	35544	133.0	97.3	68.6	128.6

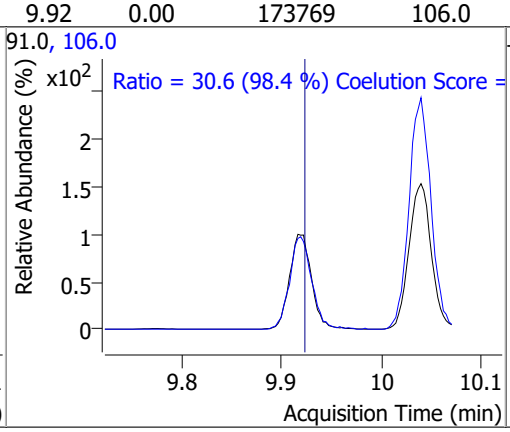


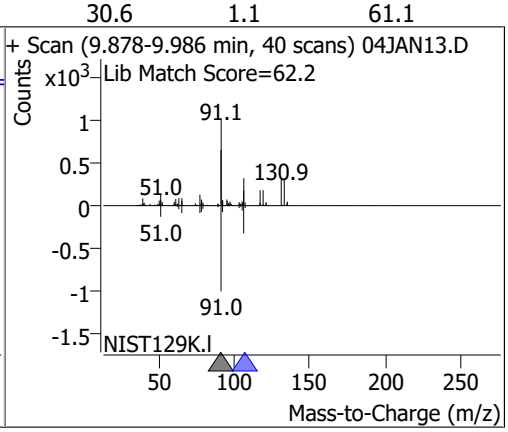




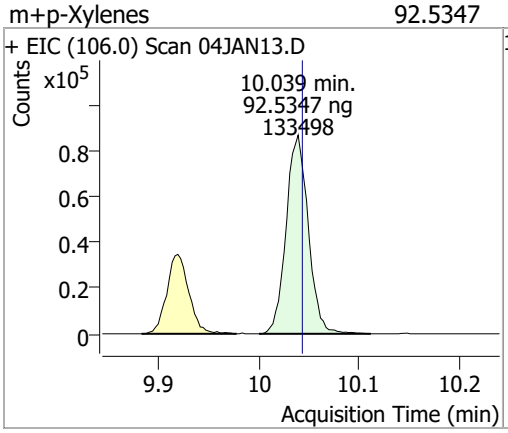
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	46.8079	9.92	0.00	173769	106.0	30.6	1.1	61.1

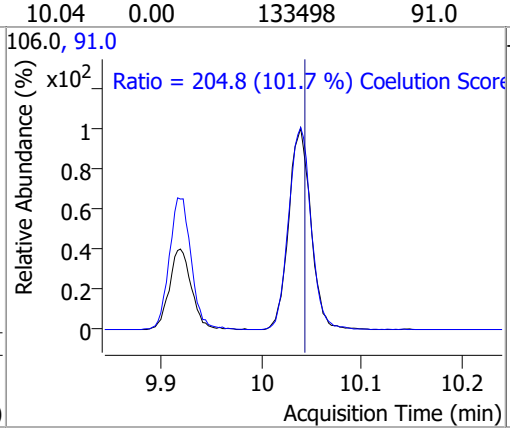


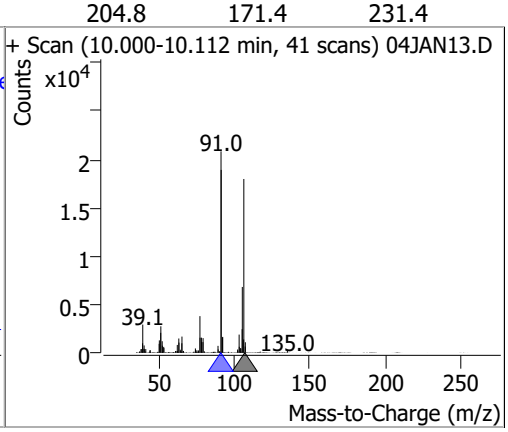




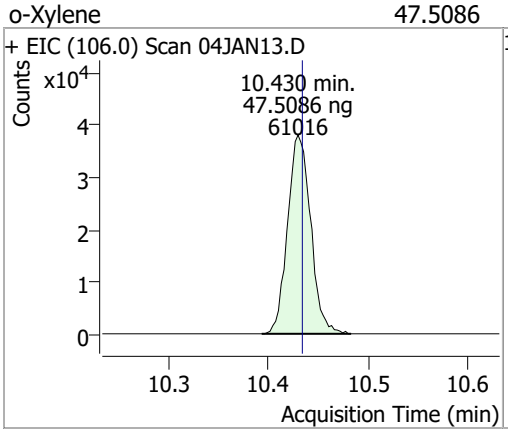
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	92.5347	10.04	0.00	133498	91.0	204.8	171.4	231.4

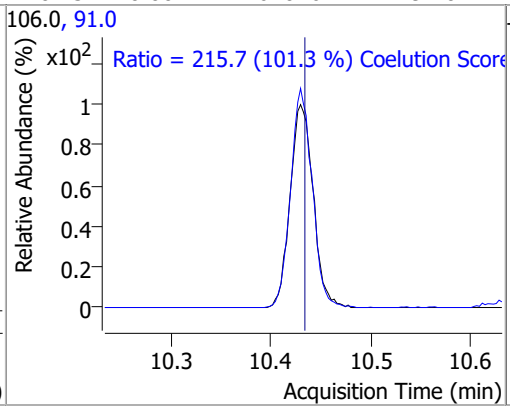


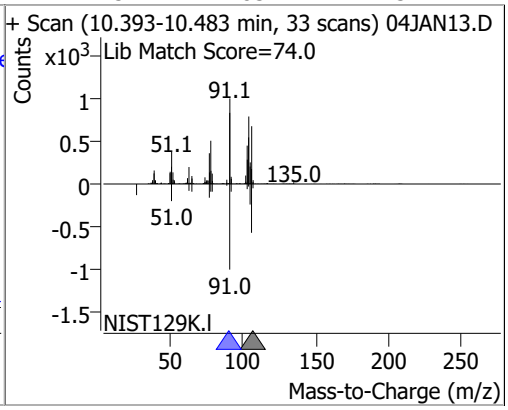




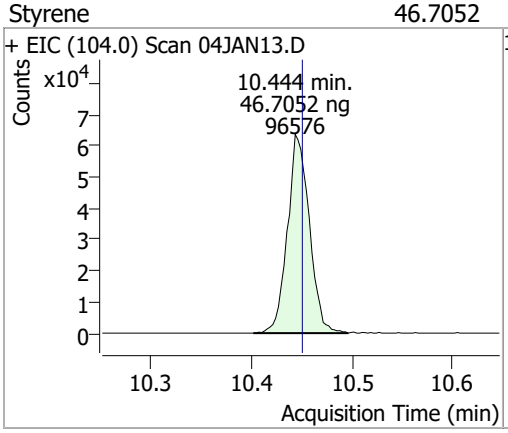
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	47.5086	10.43	0.00	61016	91.0	215.7	183.1	243.1

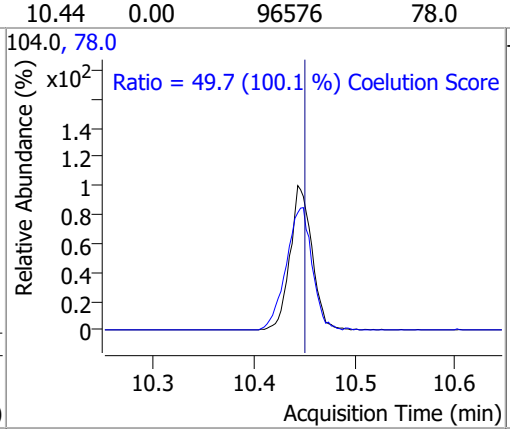


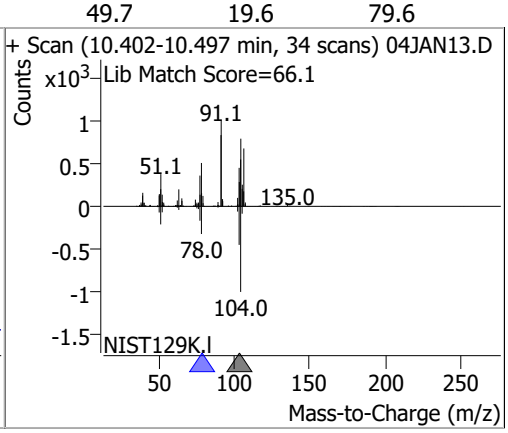




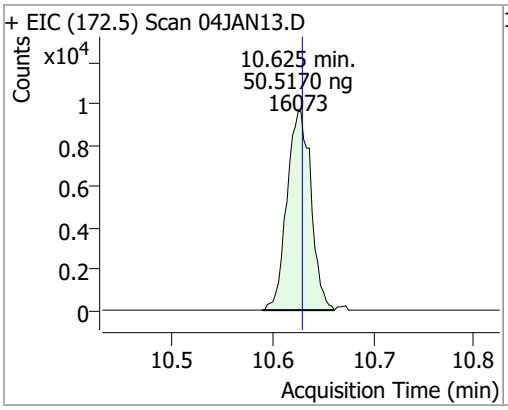
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	46.7052	10.44	0.00	96576	78.0	49.7	19.6	79.6

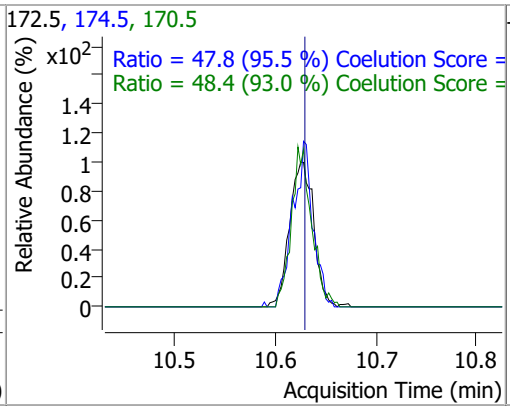


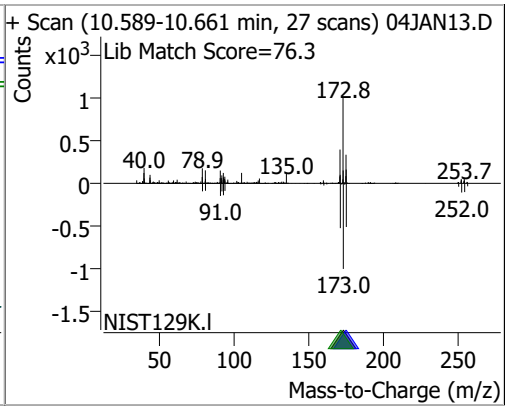




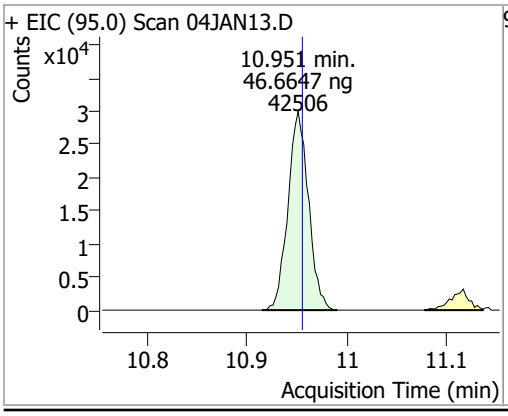
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	50.5170	10.62	0.00	16073	170.5	48.4	22.1	82.1
					174.5	47.8	20.1	80.1

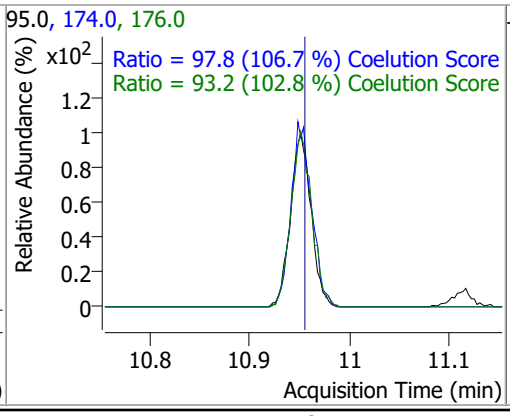


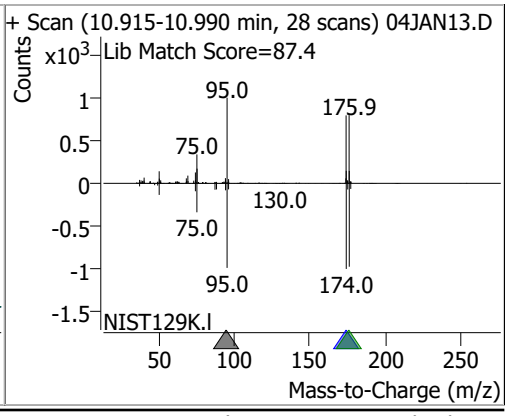




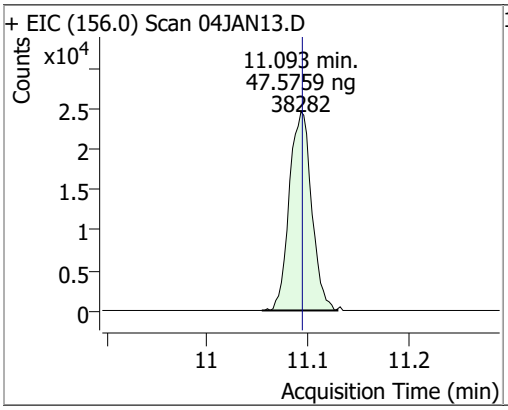
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	46.6647	10.95	0.00	42506	174.0	97.8	61.7	121.7
					176.0	93.2	60.6	120.6

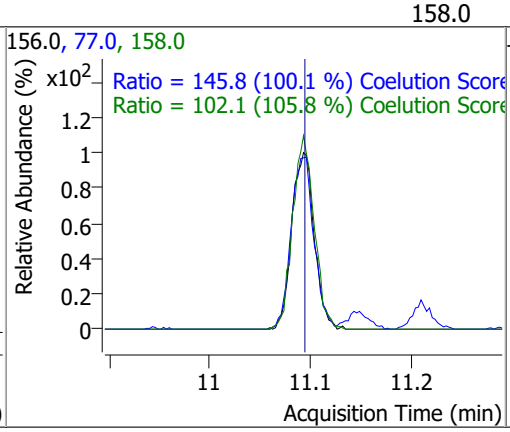


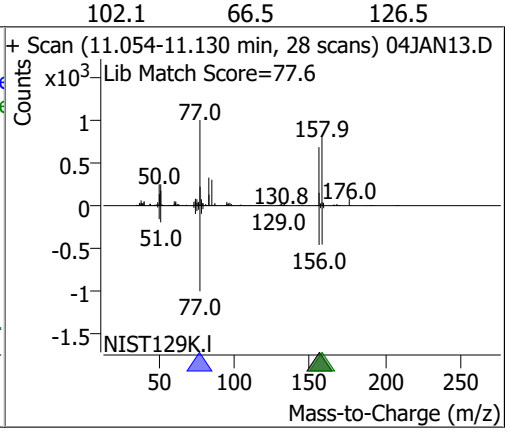




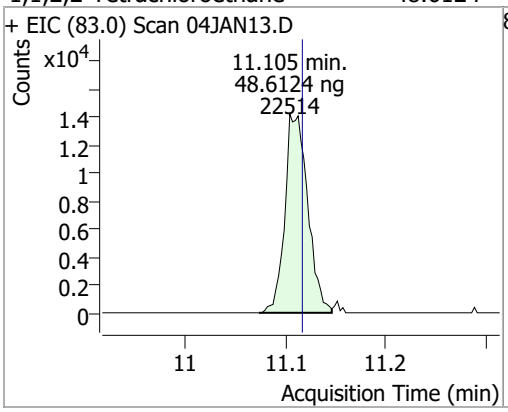
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	47.5759	11.09	0.00	38282	77.0	145.8	115.7	175.7
					158.0	102.1	66.5	126.5

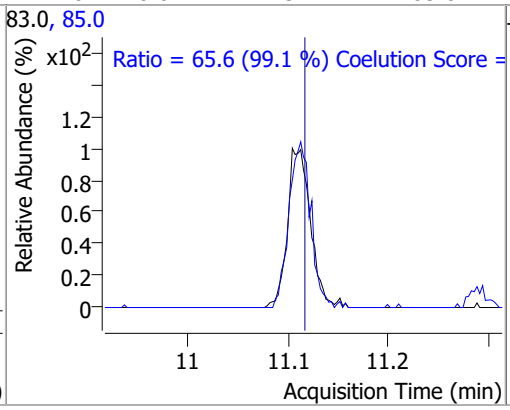


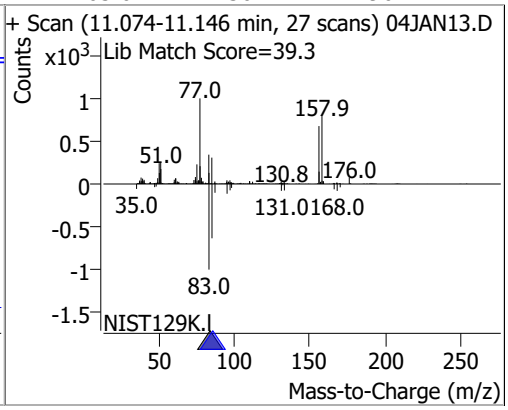




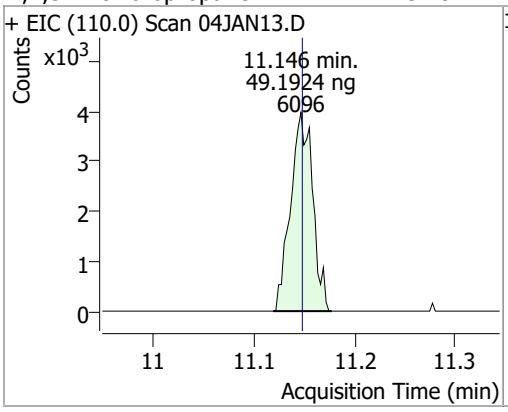
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	48.6124	11.10	-0.01	22514	85.0	65.6	36.2	96.2

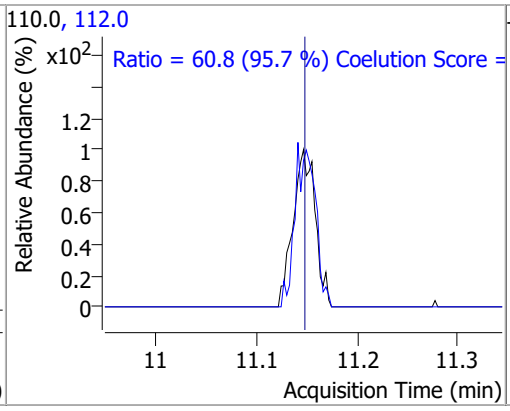


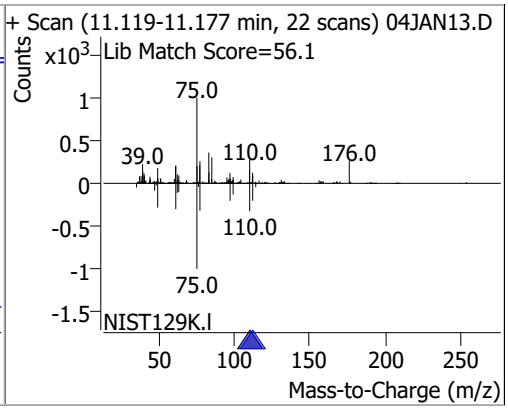




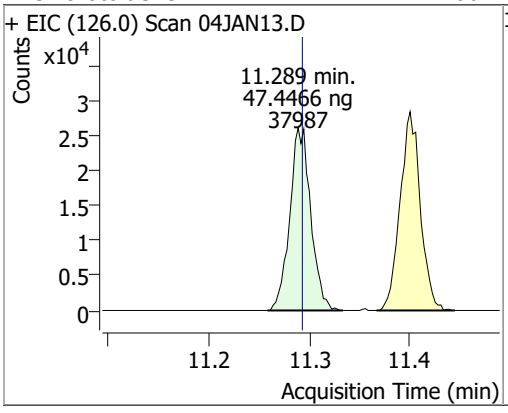
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	49.1924	11.15	0.00	6096	112.0	60.8	33.5	93.5

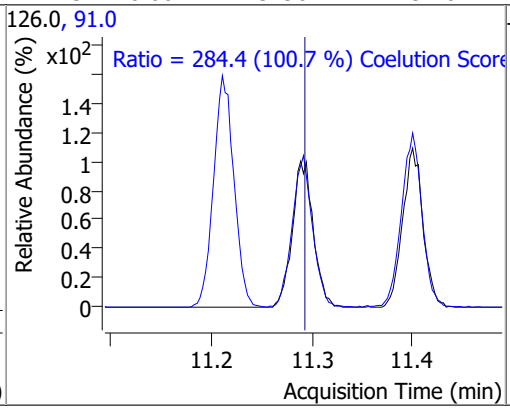


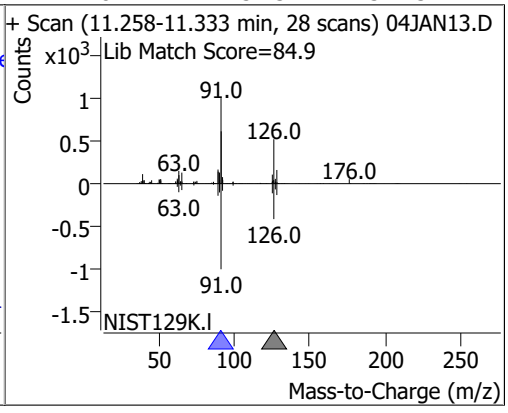




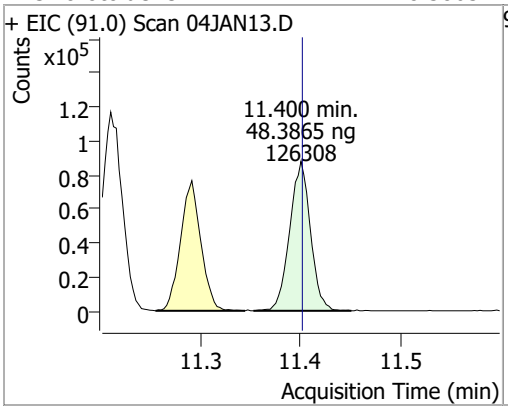
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	47.4466	11.29	0.00	37987	91.0	284.4	252.3	312.3

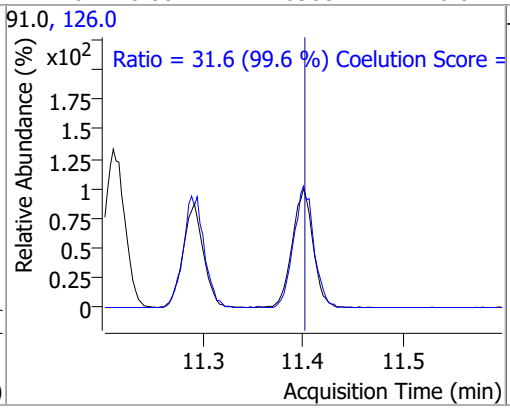


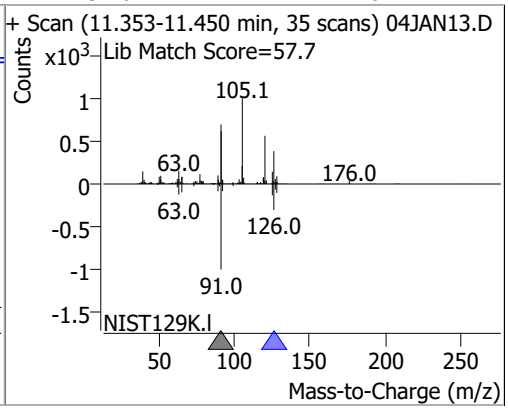




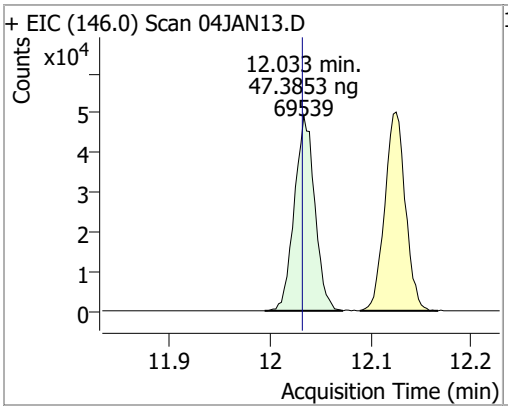
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	48.3865	11.40	0.00	126308	126.0	31.6	1.7	61.7

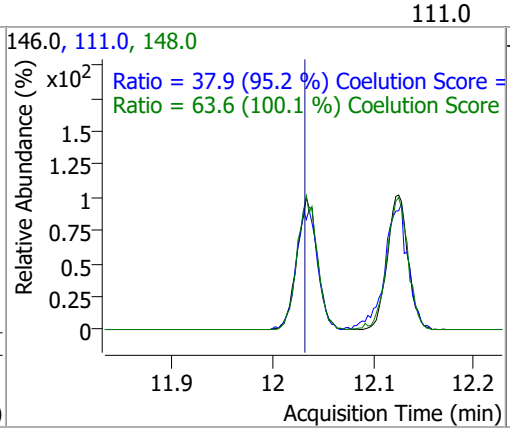


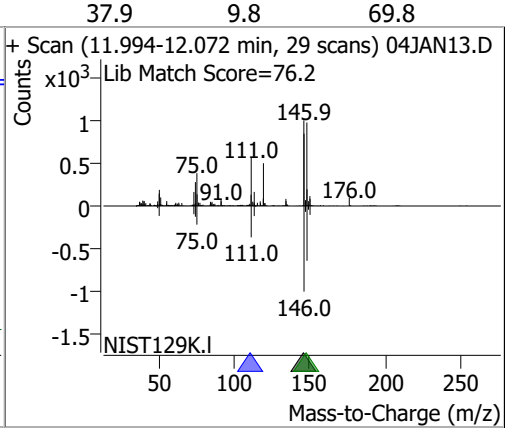




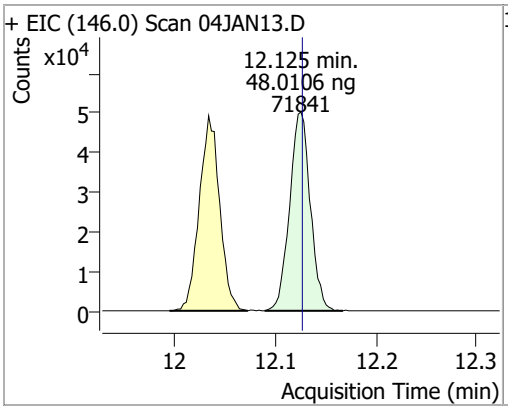
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	47.3853	12.03	0.00	69539	148.0	63.6	33.6	93.6
					111.0	37.9	9.8	69.8

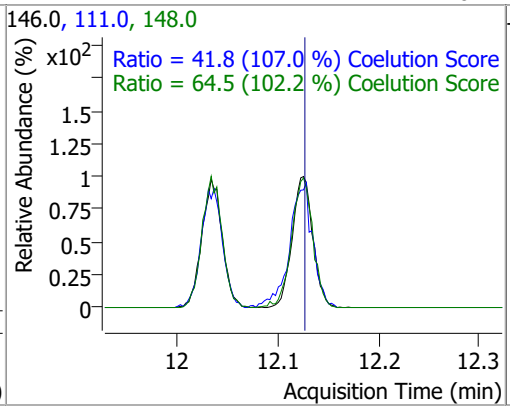


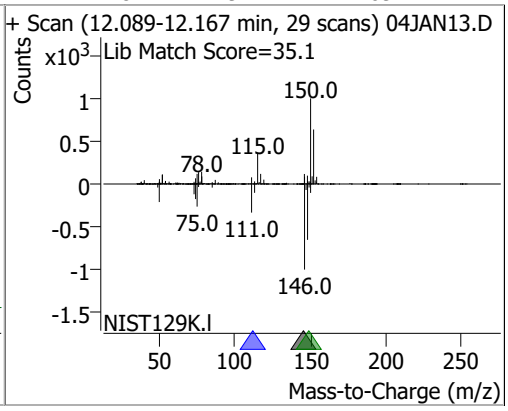




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	48.0106	12.13	0.00	71841	148.0	64.5	33.1	93.1
					111.0	41.8	9.1	69.1

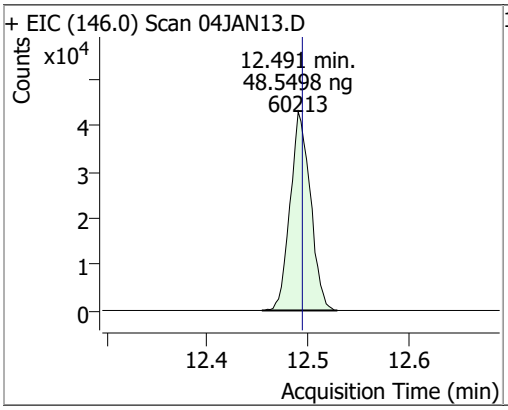


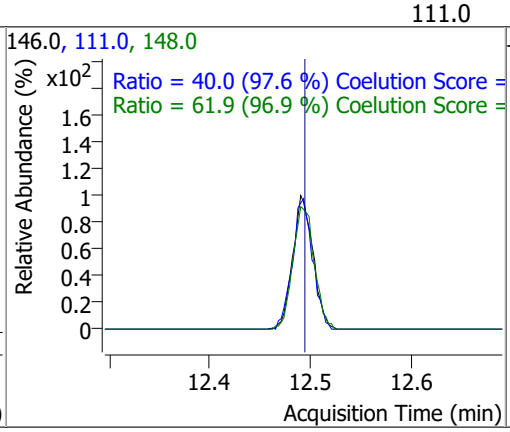


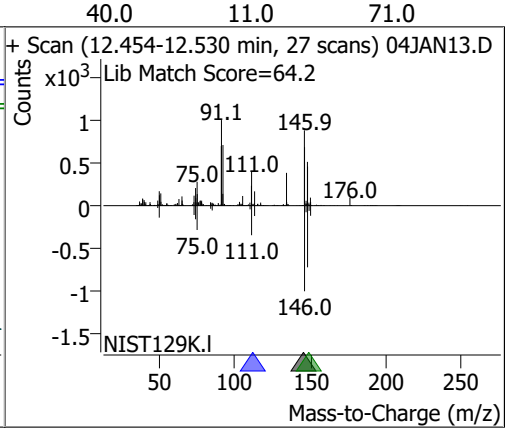


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	48.5498	12.49	0.00	60213	148.0	61.9	33.9	93.9
					111.0	40.0	11.0	71.0







Quantitation Results Report (QT Reviewed)

Data File	04JAN15.D	Operator	MSC
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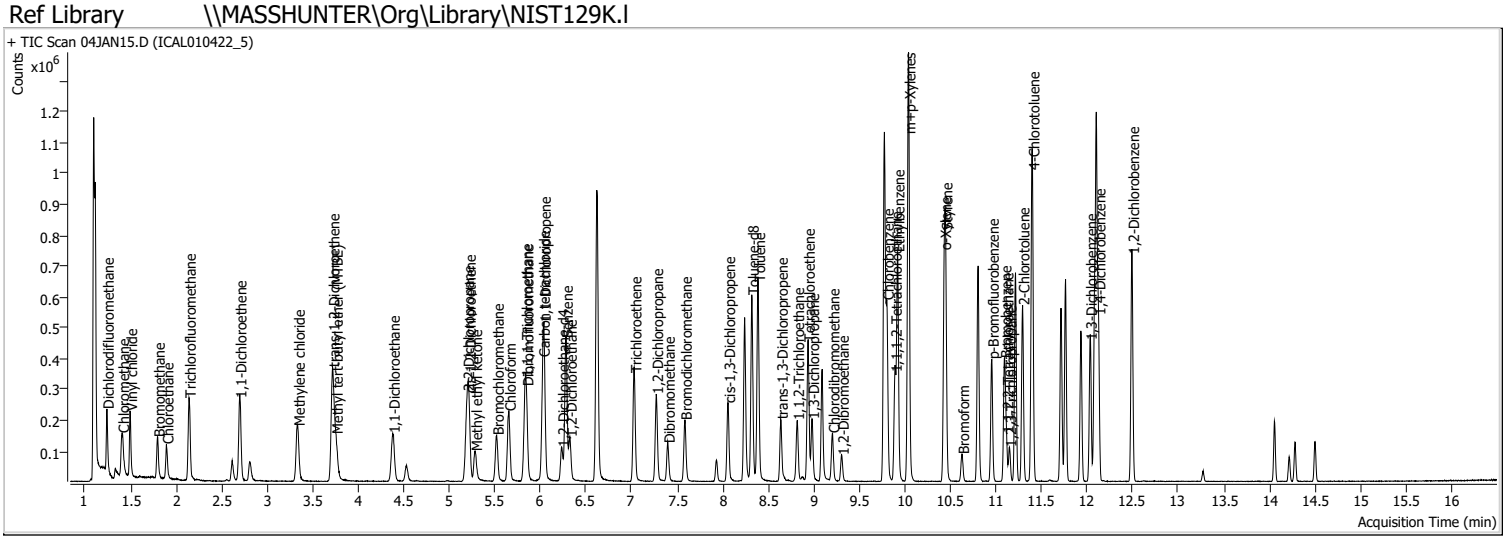
Sample Name	ICAL010422_5	Instrument	VOA5975C
Vial	15	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

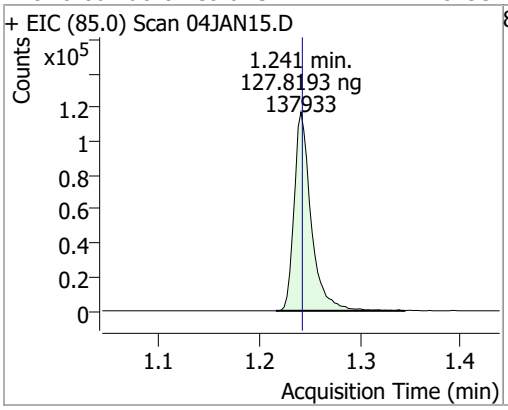


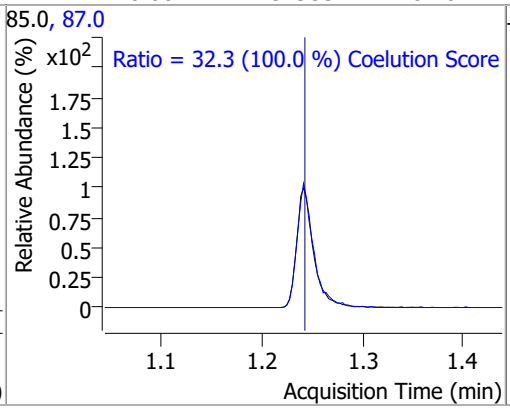
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.623	96.0	823488	250.0000	ng		0.000
M Chlorobenzene-d5	9.772	82.0	306491	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	264477	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	89307	115.1146	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 46.05%		*	
S 1,2-Dichloroethane-d4	6.233	67.0	39086	116.6420	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 46.66%		*	
S Toluene-d8	8.319	98.0	358186	121.2749	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 48.51%		*	
S p-Bromofluorobenzene	10.954	95.0	114269	117.9350	ng		0.000
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 47.17%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.241	85.0	137933	127.8193	ng		100
T Chloromethane	1.409	50.0	160604	122.6179	ng		100
T Vinyl chloride	1.495	62.0	148358	125.8809	ng		100
T Bromomethane	1.799	96.0	65163	123.6504	ng		100
T Chloroethane	1.894	64.0	71420	122.4086	ng		100
T Trichlorofluoromethane	2.142	101.0	188808	129.0687	ng		100
T 1,1-Dichloroethene	2.697	96.0	99438	119.8798	ng		100
T Methylene chloride	3.336	49.0	135271	110.6249	ng		100
T trans-1,2-Dichloroethene	3.718	96.0	100409	118.6511	ng		100
T Methyl tert-butyl ether (MTBE)	3.754	73.0	139068	127.1375	ng		100
T 1,1-Dichloroethane	4.378	63.0	186052	118.1125	ng		100
T 2,2-Dichloropropane	5.196	77.0	139656	118.3203	ng		100
T cis-1,2-Dichloroethene	5.215	96.0	100057	116.6190	ng		100
T Methyl ethyl ketone	5.282	43.0	134730	1159.3019	ng		100
T Bromochloromethane	5.519	128.0	41966	118.0683	ng		100
T Chloroform	5.653	83.0	179640	114.5912	ng		100

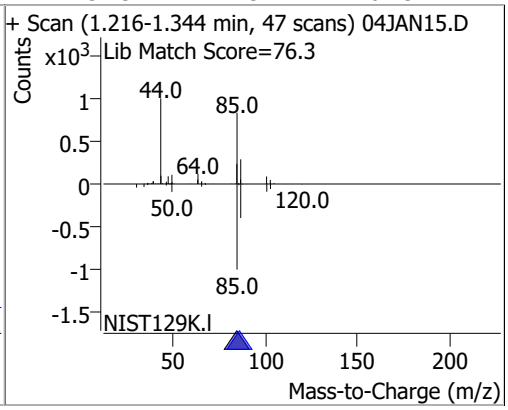
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.834	97.0	174206	118.5764	ng	100
T Carbon tetrachloride	6.024	117.0	172928	119.4667	ng	100
T 1,1-Dichloropropene	6.038	75.0	149649	119.8002	ng	100
T Benzene	6.278	78.0	383469	116.9553	ng	100
T 1,2-Dichloroethane	6.322	62.0	104855	118.2143	ng	100
T Trichloroethene	7.030	95.0	114123	123.4646	ng	100
T 1,2-Dichloropropane	7.270	63.0	99187	121.9890	ng	100
T Dibromomethane	7.399	93.0	40628	118.2425	ng	100
T Bromodichloromethane	7.585	83.0	115664	121.9749	ng	100
T cis-1,3-Dichloropropene	8.059	75.0	129419	120.7116	ng	100
T Toluene	8.389	92.0	244712	122.6571	ng	100
T trans-1,3-Dichloropropene	8.637	75.0	92719	121.4929	ng	100
T 1,1,2-Trichloroethane	8.818	83.0	46673	117.4130	ng	100
T Tetrachloroethene	8.935	163.8	97590	119.9003	ng	100
T 1,3-Dichloropropane	8.980	76.0	96183	123.0132	ng	100
T Chlorodibromomethane	9.206	129.0	75015	120.7454	ng	100
T 1,2-Dibromoethane	9.306	107.0	51827	119.2394	ng	100
T Chlorobenzene	9.802	112.0	263617	120.6903	ng	100
T 1,1,1,2-Tetrachloroethane	9.889	131.0	90898	119.0492	ng	100
T Ethylbenzene	9.920	91.0	464148	122.5243	ng	100
T m+p-Xylenes	10.039	106.0	368418	250.2587	ng	100
T o-Xylene	10.430	106.0	161509	123.2378	ng	100
T Styrene	10.447	104.0	268375	127.1910	ng	100
T Bromoform	10.628	172.5	39165	115.7218	ng	100
T Bromobenzene	11.094	156.0	102265	119.4801	ng	100
T 1,1,2,2-Tetrachloroethane	11.116	83.0	56958	115.6179	ng	100
T 1,2,3-Trichloropropane	11.147	110.0	14846	112.6261	ng	100
T 2-Chlorotoluene	11.292	126.0	102424	120.2675	ng	100
T 4-Chlorotoluene	11.400	91.0	336146	121.0591	ng	100
T 1,3-Dichlorobenzene	12.031	146.0	183404	117.4899	ng	100
T 1,4-Dichlorobenzene	12.125	146.0	189045	118.7699	ng	100
T 1,2-Dichlorobenzene	12.493	146.0	152284	115.4323	ng	100

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

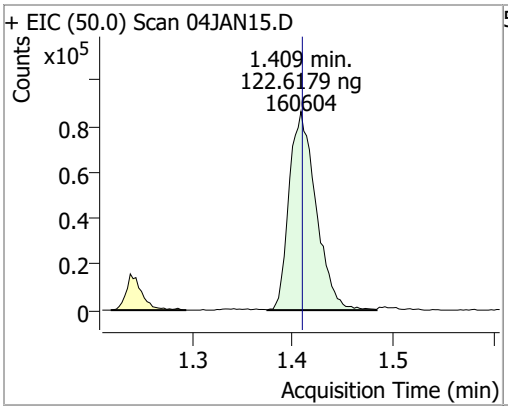
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	127.8193	1.24	0.00	137933	87.0	32.3	2.3	62.3

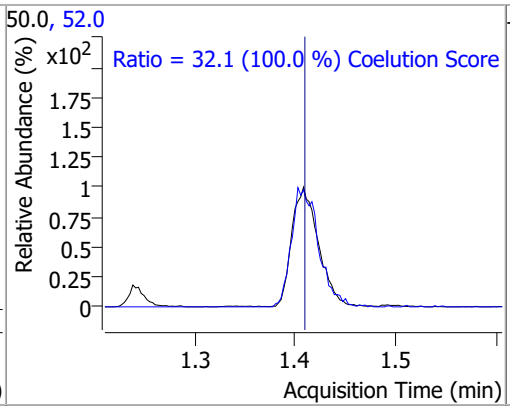


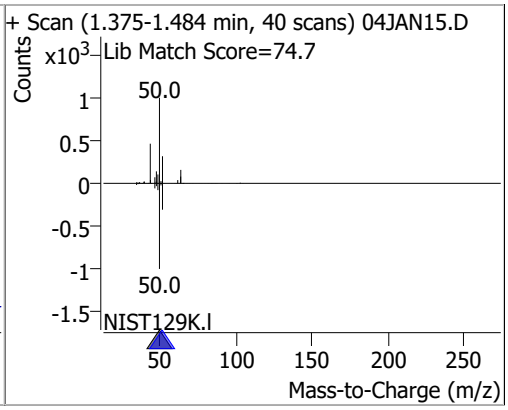




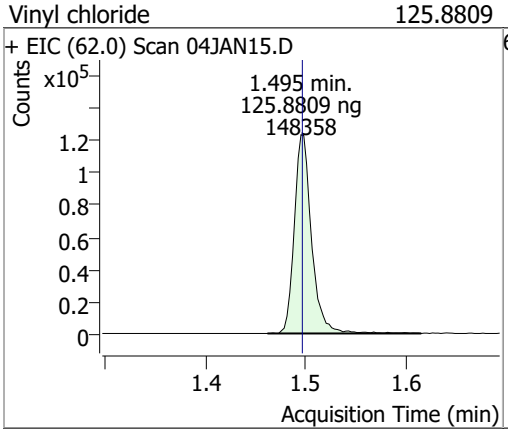
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	122.6179	1.41	0.00	160604	52.0	32.1	2.1	62.1

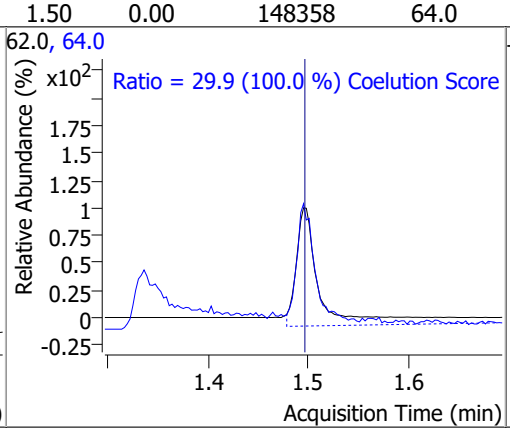


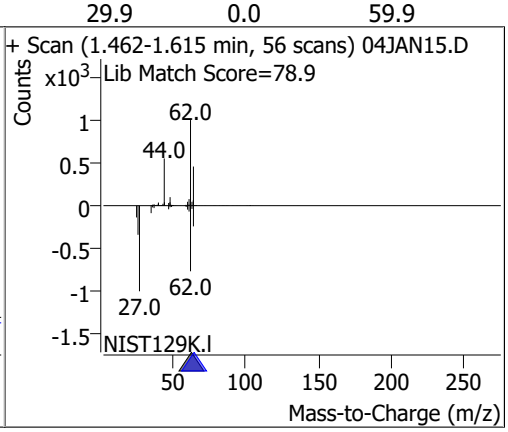




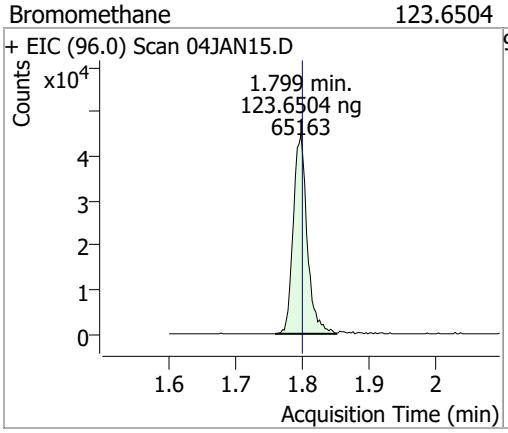
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	125.8809	1.50	0.00	148358	64.0	29.9	0.0	59.9

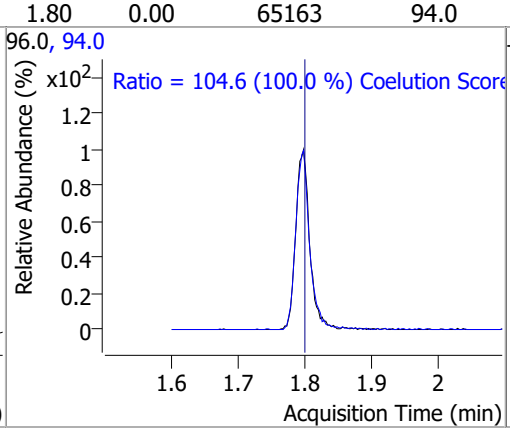


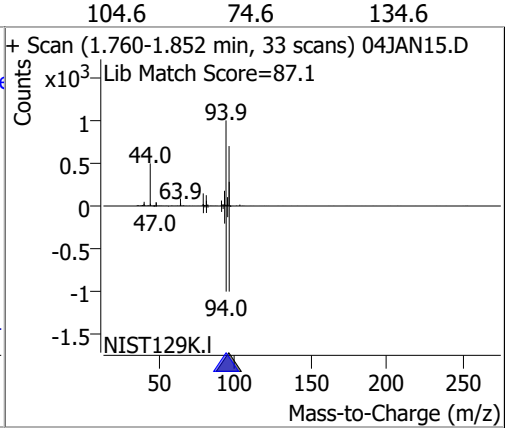




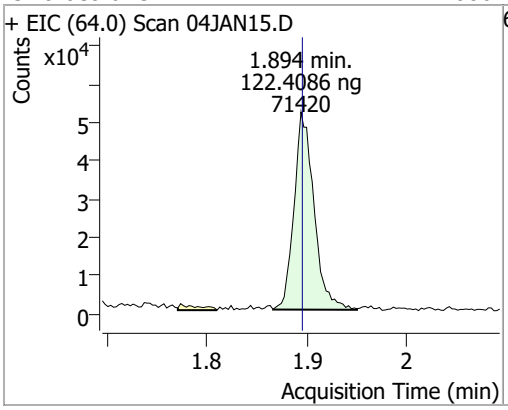
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	123.6504	1.80	0.00	65163	94.0	104.6	74.6	134.6

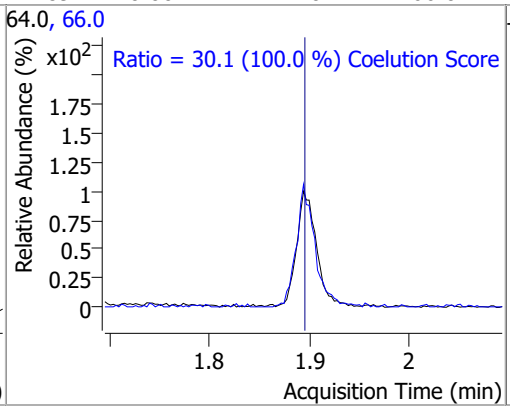


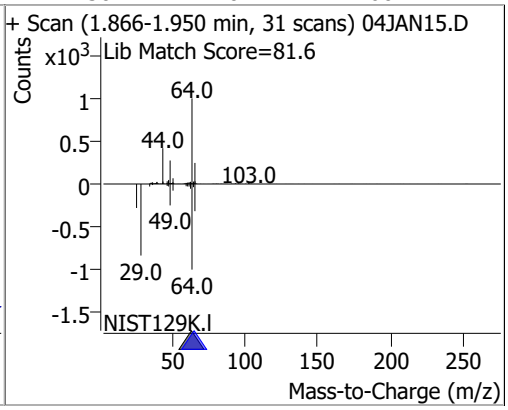




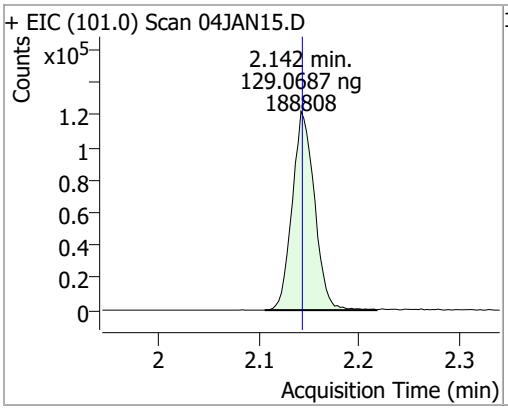
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	122.4086	1.89	0.00	71420	66.0	30.1	0.1	60.1

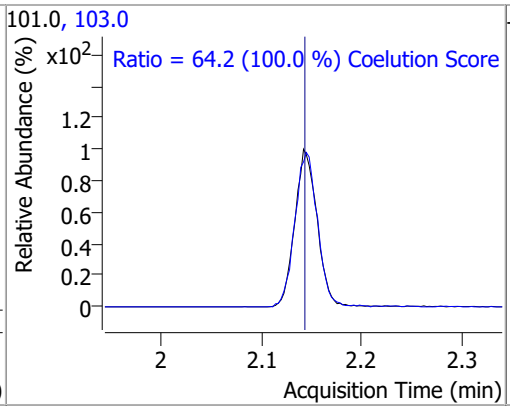


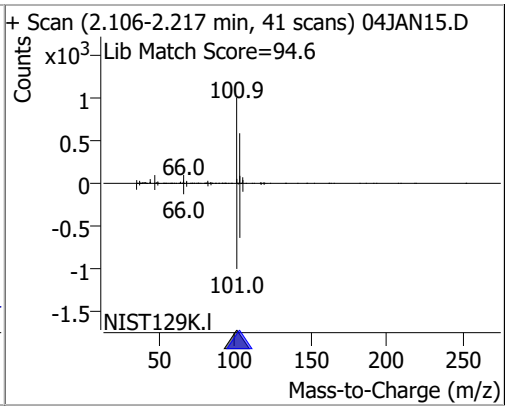




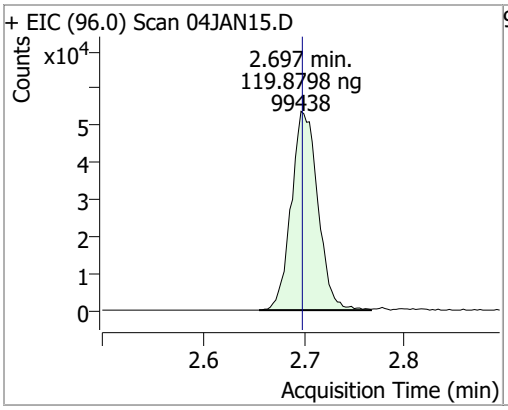
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	129.0687	2.14	0.00	188808	103.0	64.2	34.2	94.2

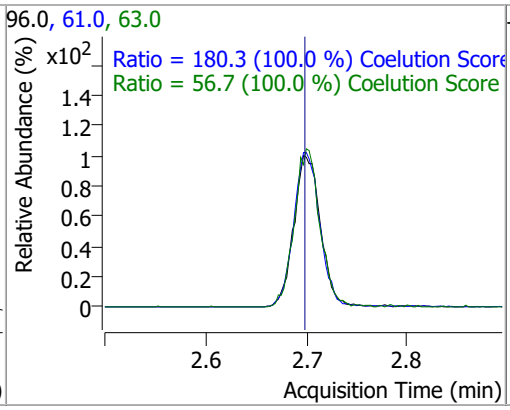


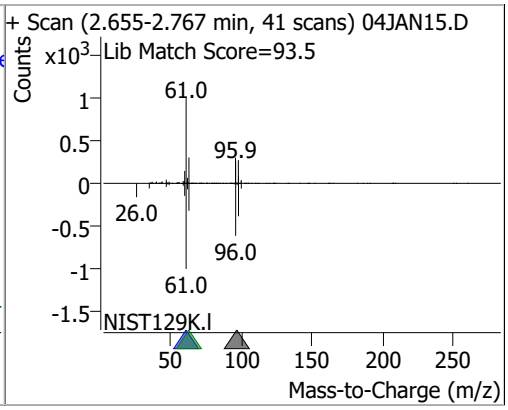




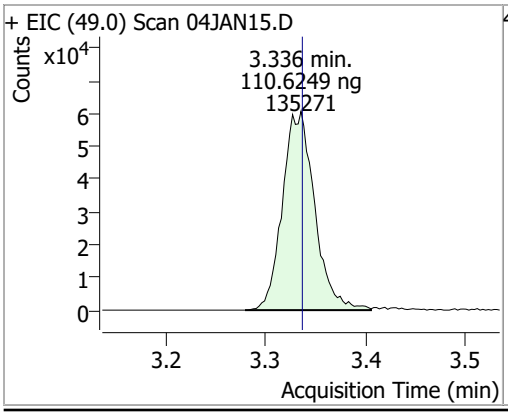
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	119.8798	2.70	0.00	99438	61.0	180.3	150.3	210.3
					63.0	56.7	26.7	86.7

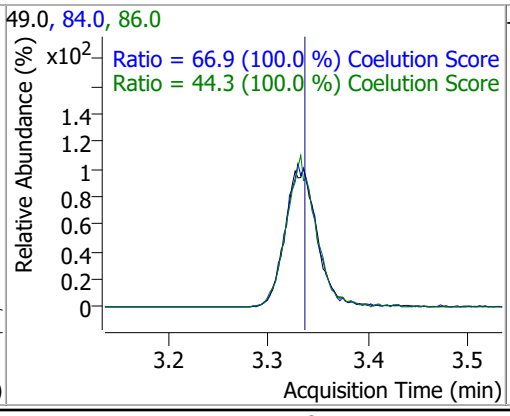


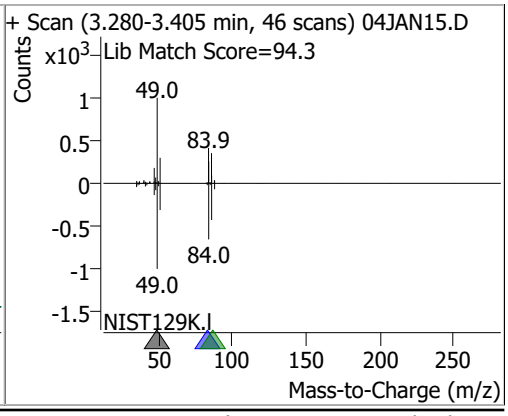




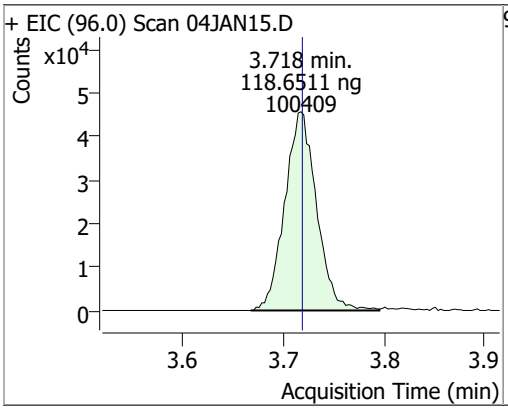
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	110.6249	3.34	0.00	135271	84.0	66.9	36.9	96.9
					86.0	44.3	14.3	74.3

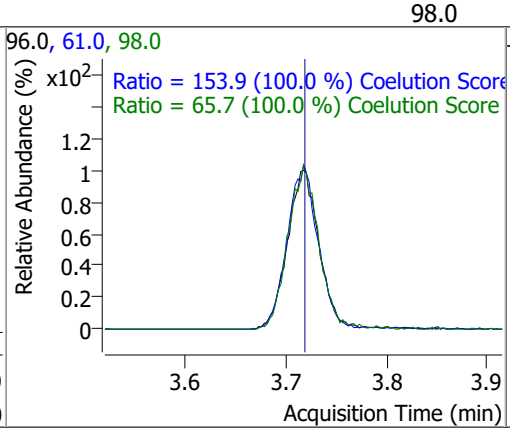


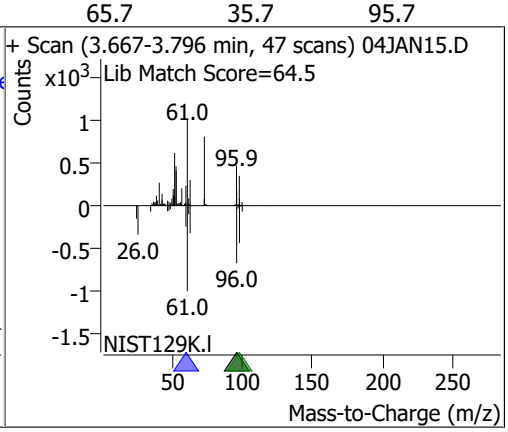




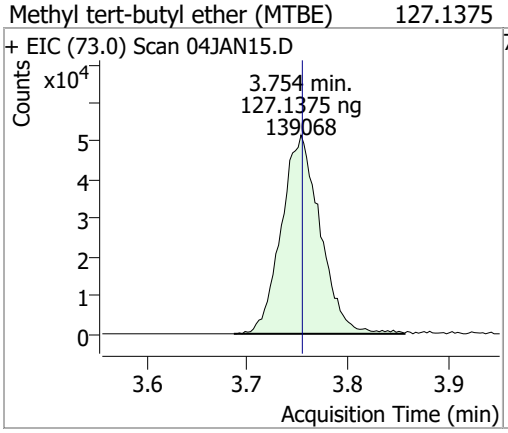
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	118.6511	3.72	0.00	100409	61.0	153.9	123.9	183.9
					98.0	65.7	35.7	95.7

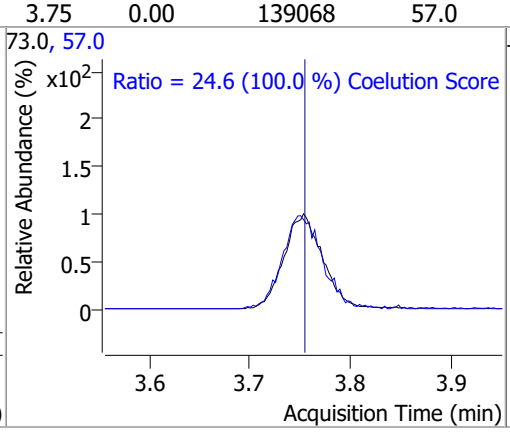


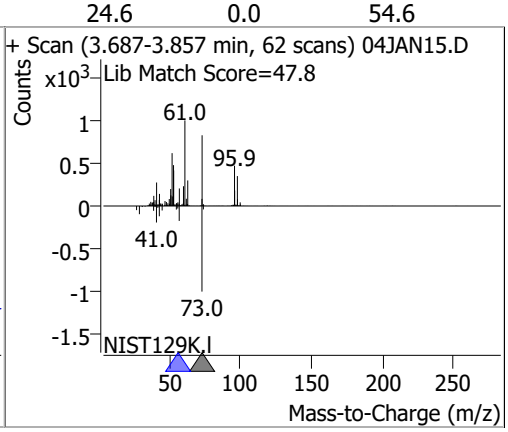




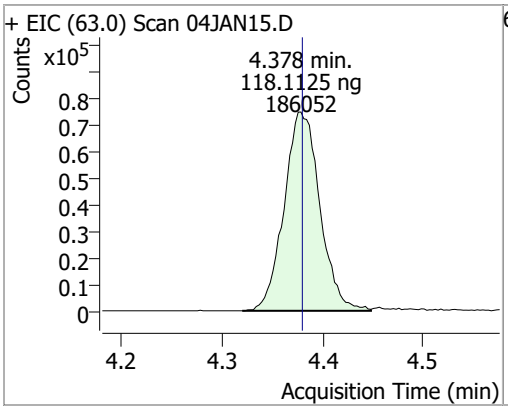
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	127.1375	3.75	0.00	139068	57.0	24.6	0.0	54.6

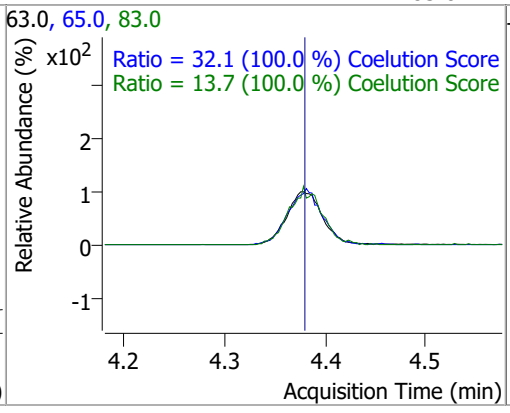


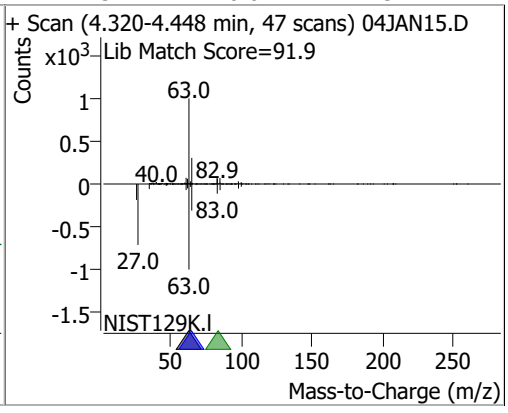




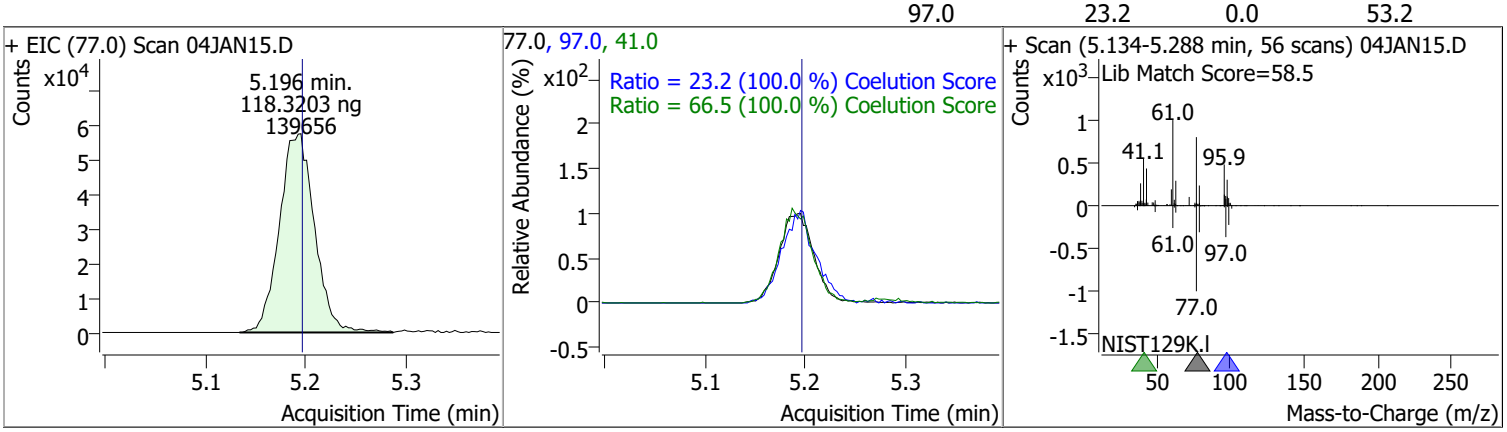
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	118.1125	4.38	0.00	186052	65.0	32.1	2.1	62.1
					83.0	13.7	0.0	43.7



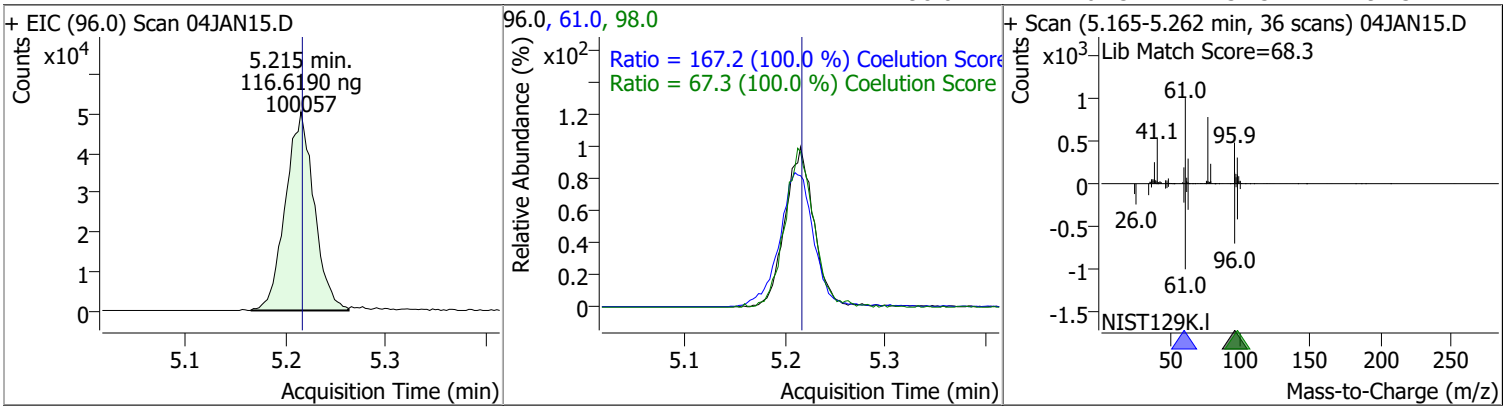




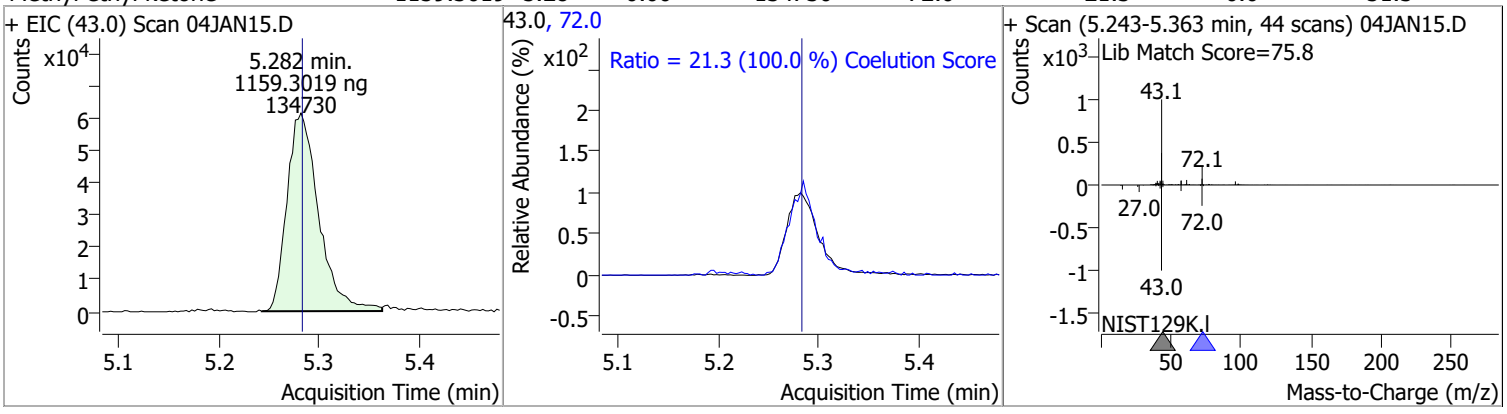
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	118.3203	5.20	0.00	139656	41.0	66.5	36.5	96.5
					97.0	23.2	0.0	53.2



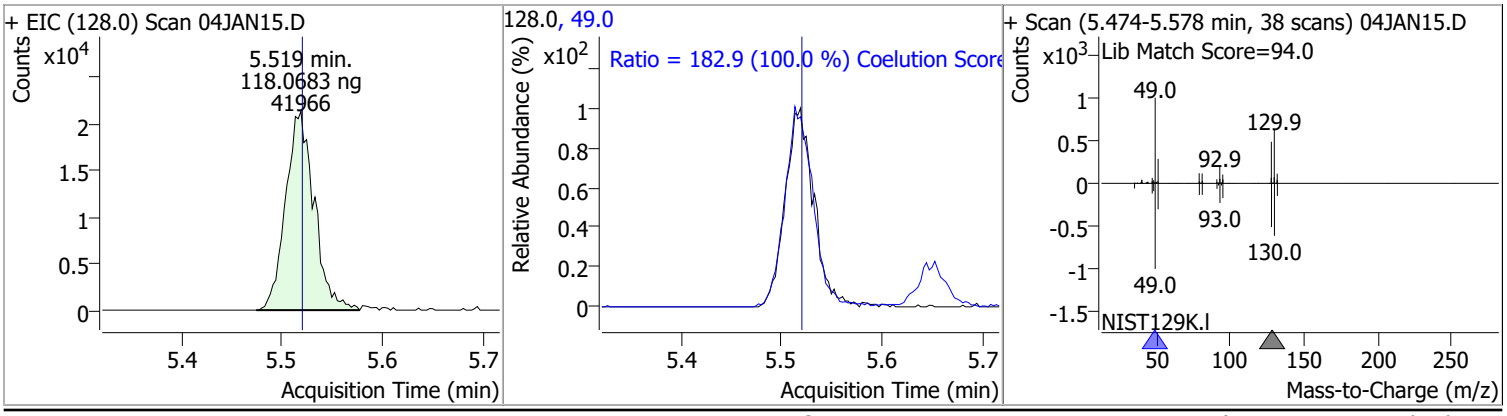
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	116.6190	5.22	0.00	100057	61.0	167.2	137.2	197.2
					98.0	67.3	37.3	97.3



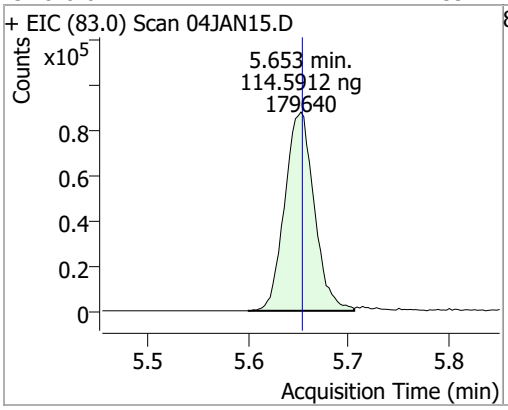
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1159.3019	5.28	0.00	134730	72.0	21.3	0.0	51.3

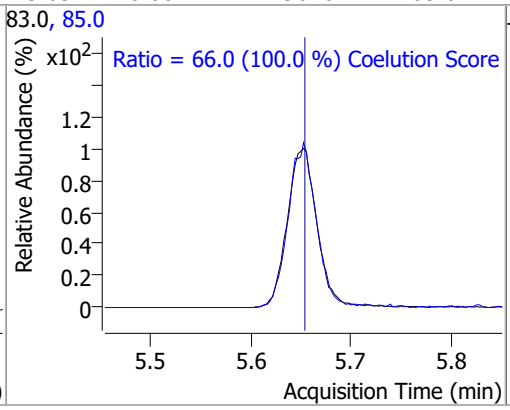


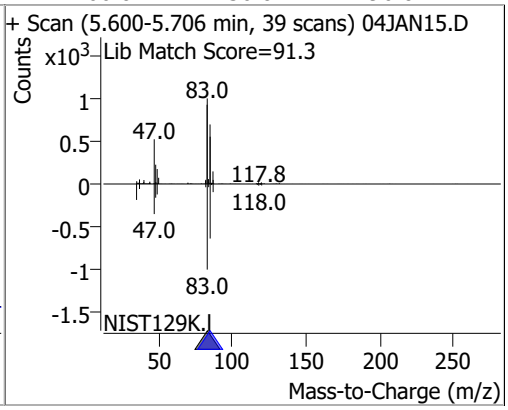
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	118.0683	5.52	0.00	41966	49.0	182.9	152.9	212.9



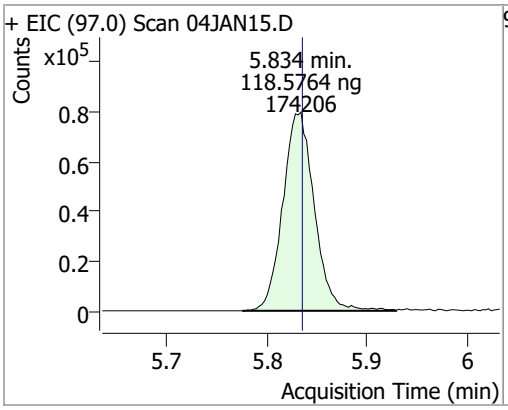
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	114.5912	5.65	0.00	179640	85.0	66.0	36.0	96.0

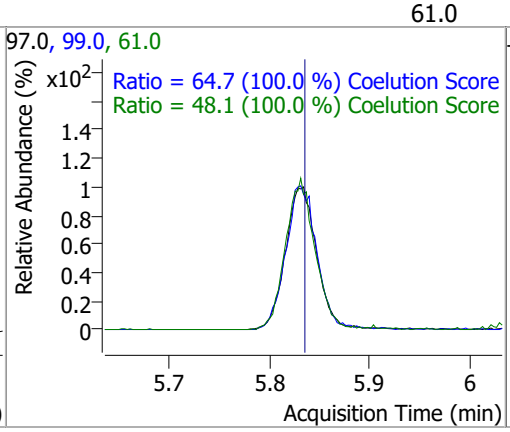


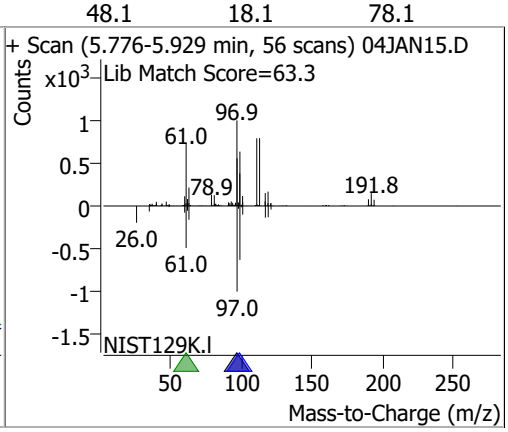




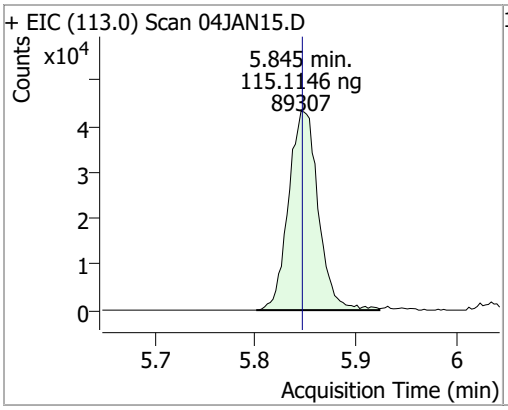
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	118.5764	5.83	0.00	174206	99.0	64.7	34.7	94.7
					61.0	48.1	18.1	78.1

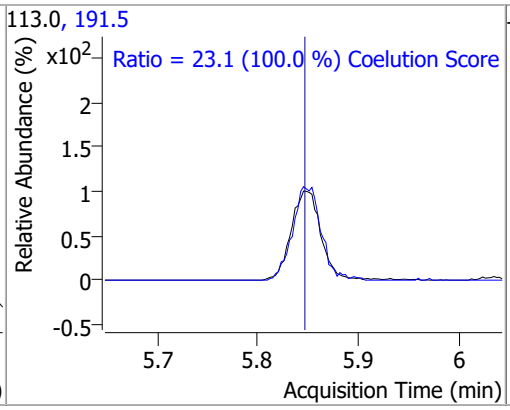


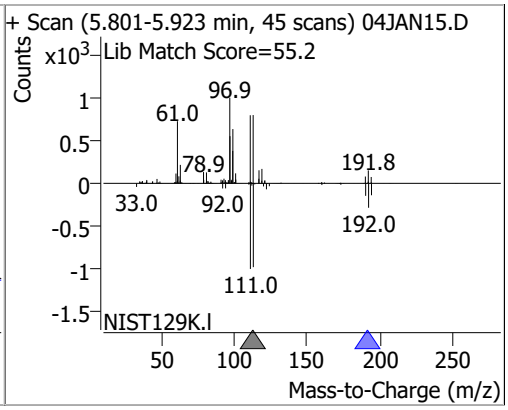




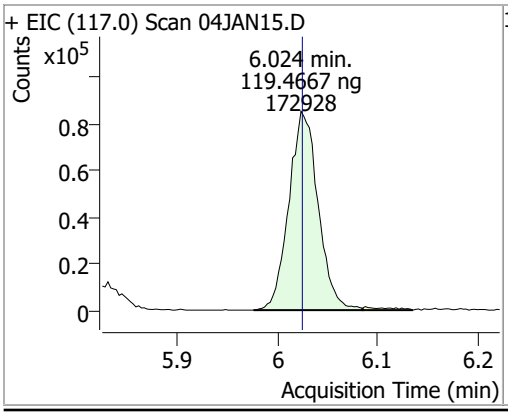
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	115.1146	5.85	0.00	89307	191.5	23.1	0.0	53.1

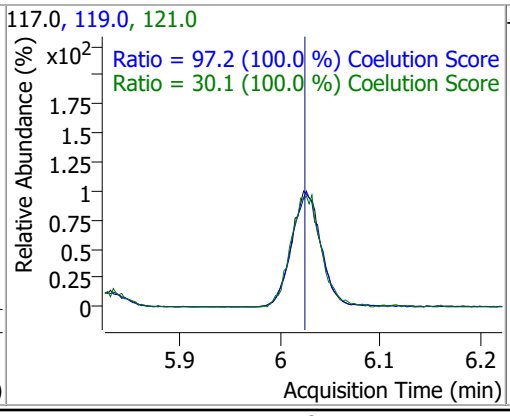


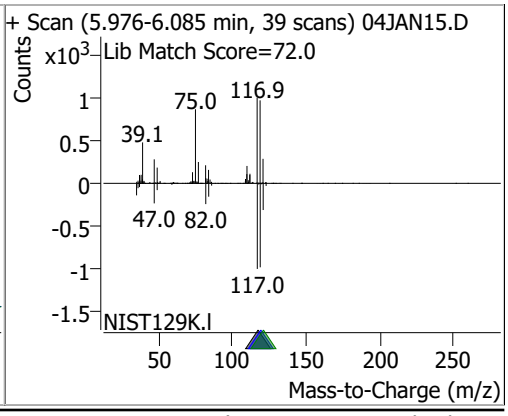




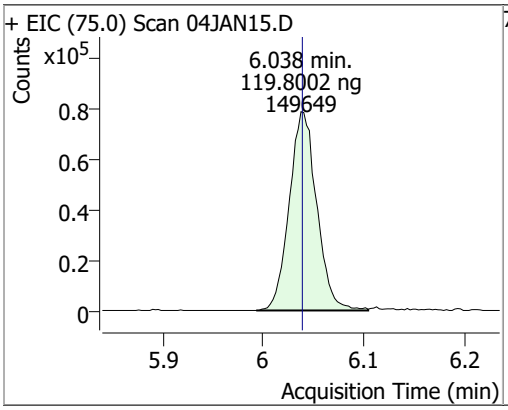
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	119.4667	6.02	0.00	172928	119.0	97.2	67.2	127.2
					121.0	30.1	0.1	60.1

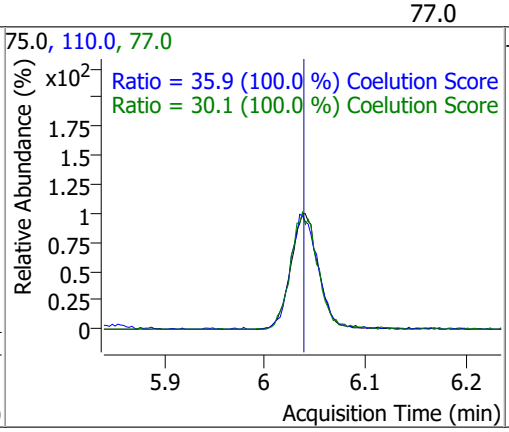


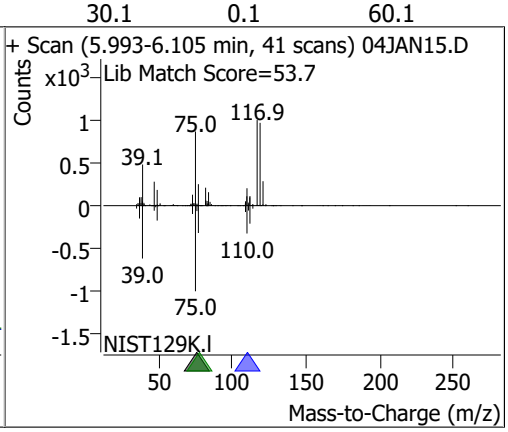




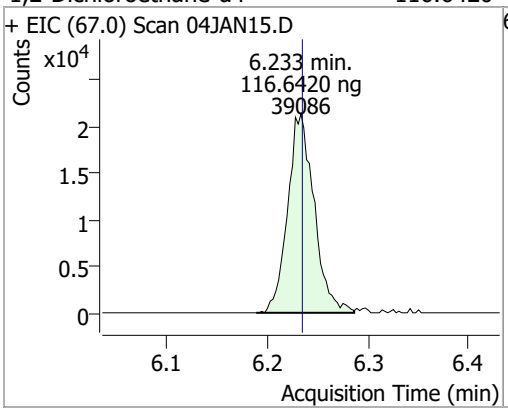
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	119.8002	6.04	0.00	149649	110.0	35.9	5.9	65.9
					77.0	30.1	0.1	60.1

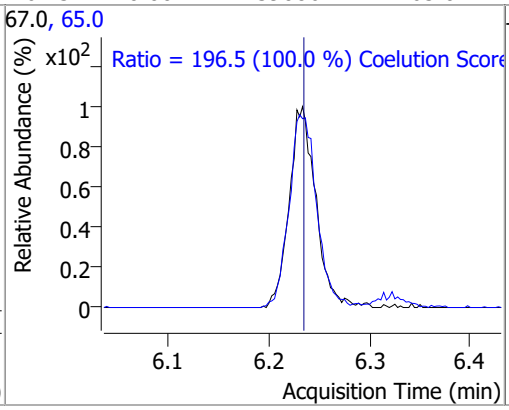


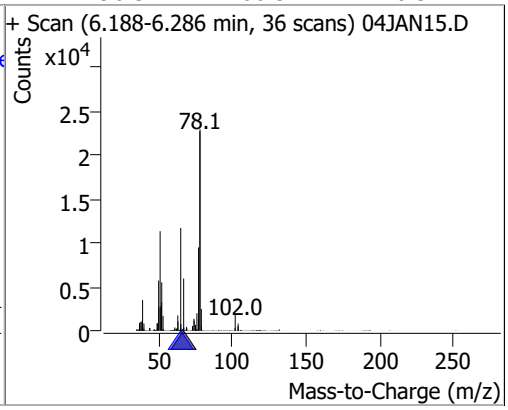




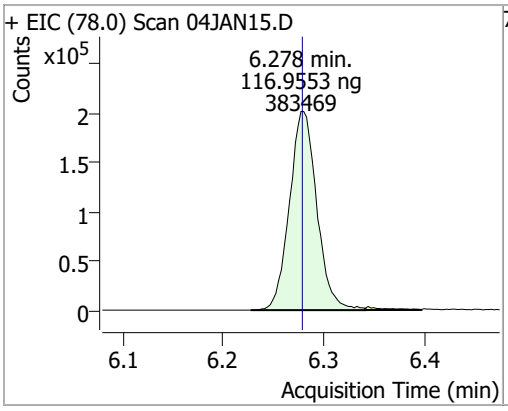
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	116.6420	6.23	0.00	39086	65.0	196.5	166.5	226.5

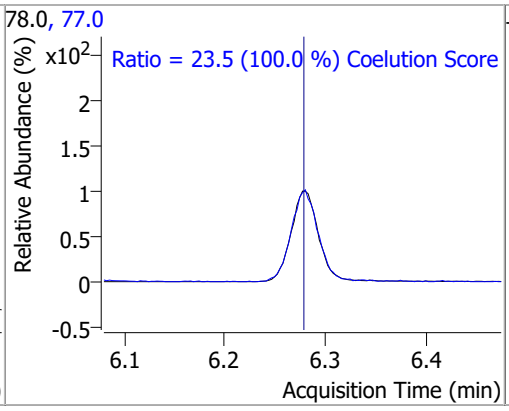


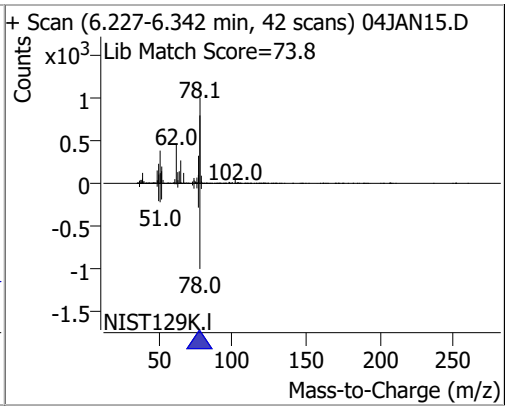




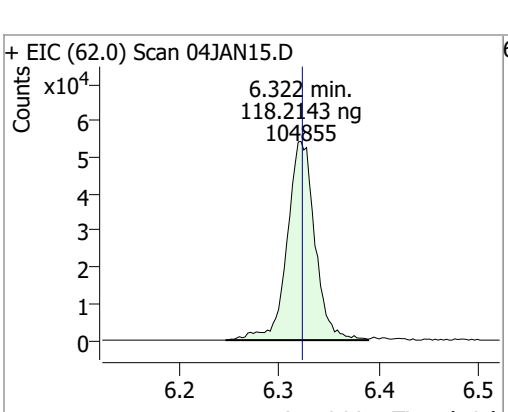
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	116.9553	6.28	0.00	383469	77.0	23.5	0.0	53.5

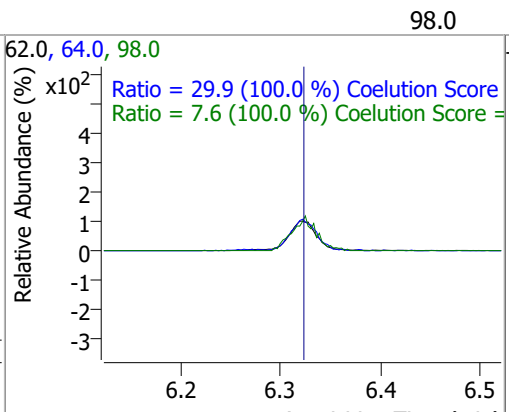


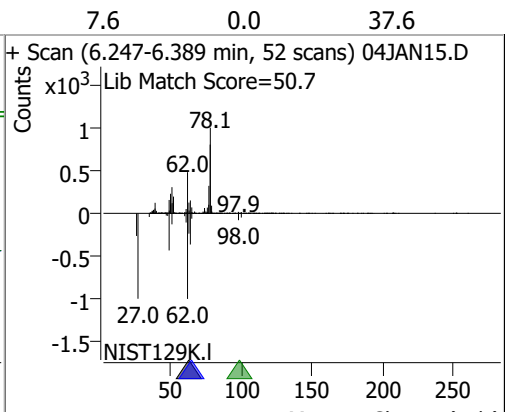




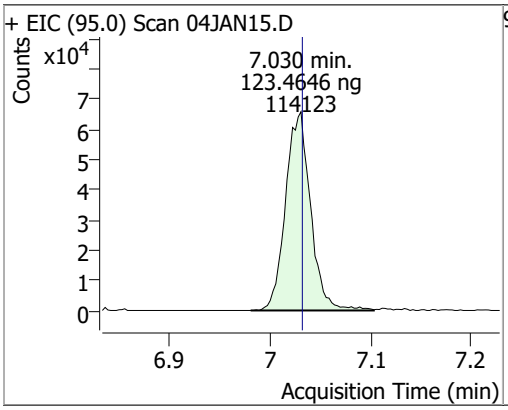
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	118.2143	6.32	0.00	104855	64.0	29.9	0.0	59.9
					98.0	7.6	0.0	37.6

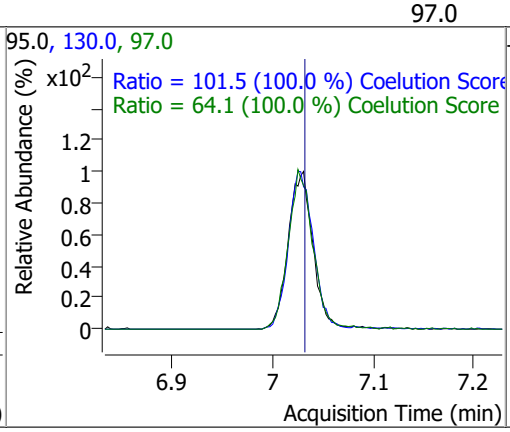


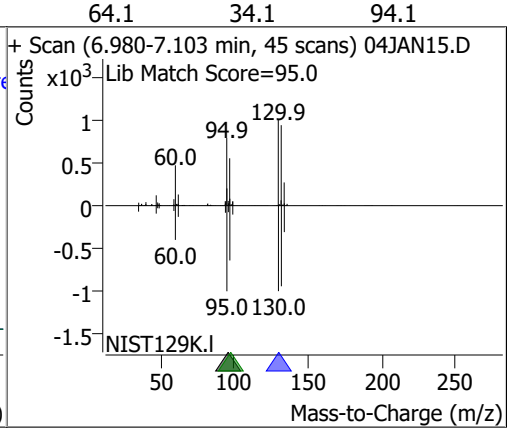




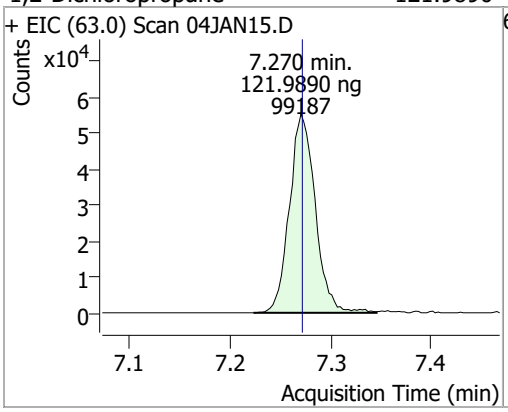
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	123.4646	7.03	0.00	114123	130.0	101.5	71.5	131.5
					97.0	64.1	34.1	94.1

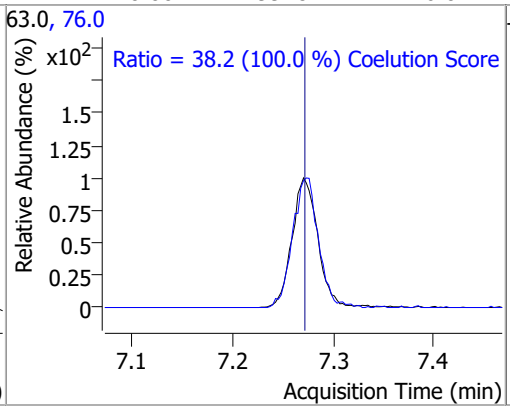


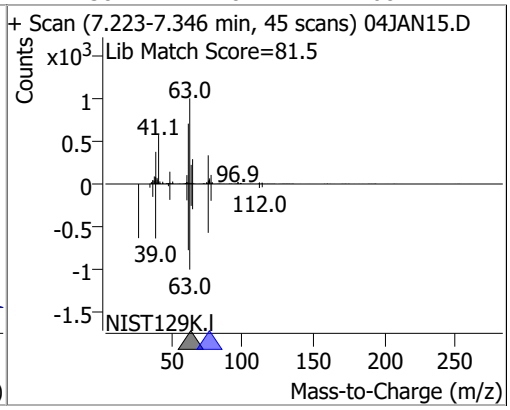




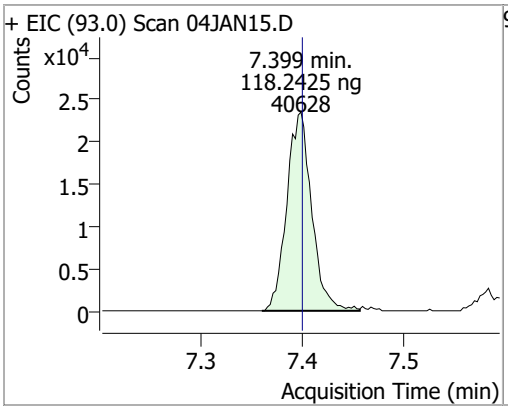
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	121.9890	7.27	0.00	99187	76.0	38.2	8.2	68.2

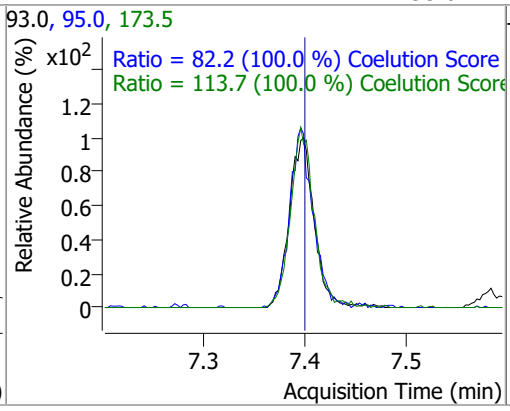


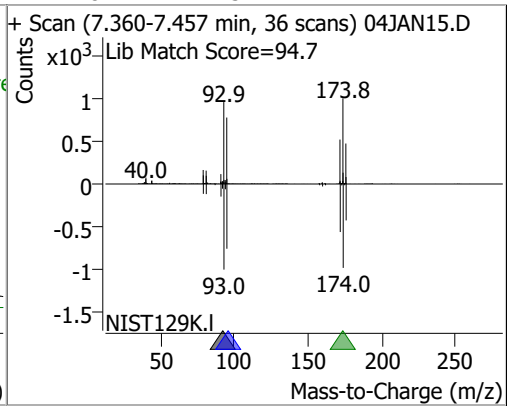




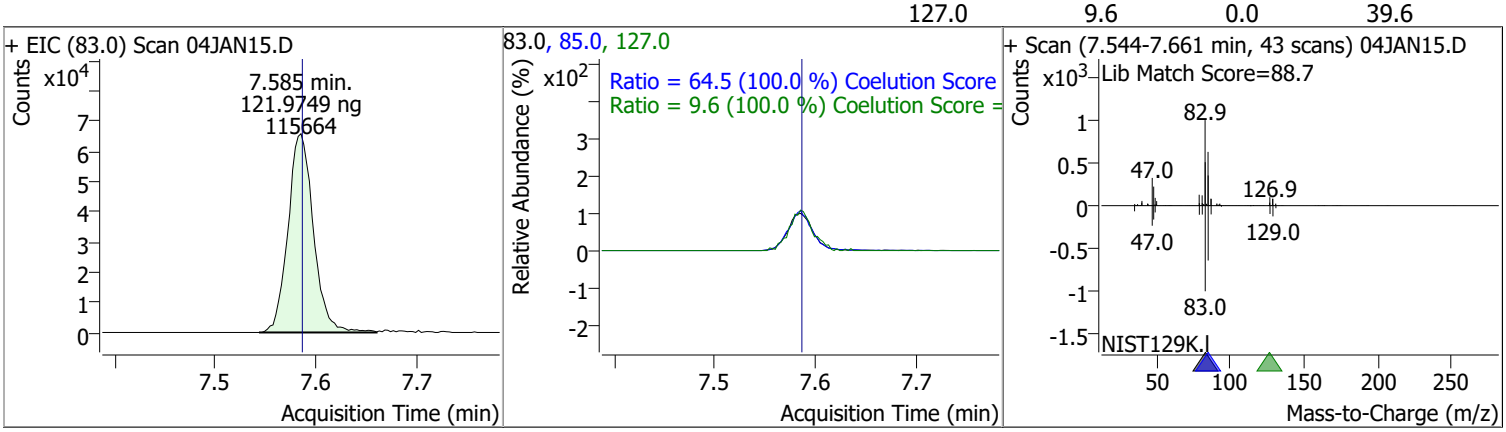
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	118.2425	7.40	0.00	40628	173.5	113.7	83.7	143.7
					95.0	82.2	52.2	112.2



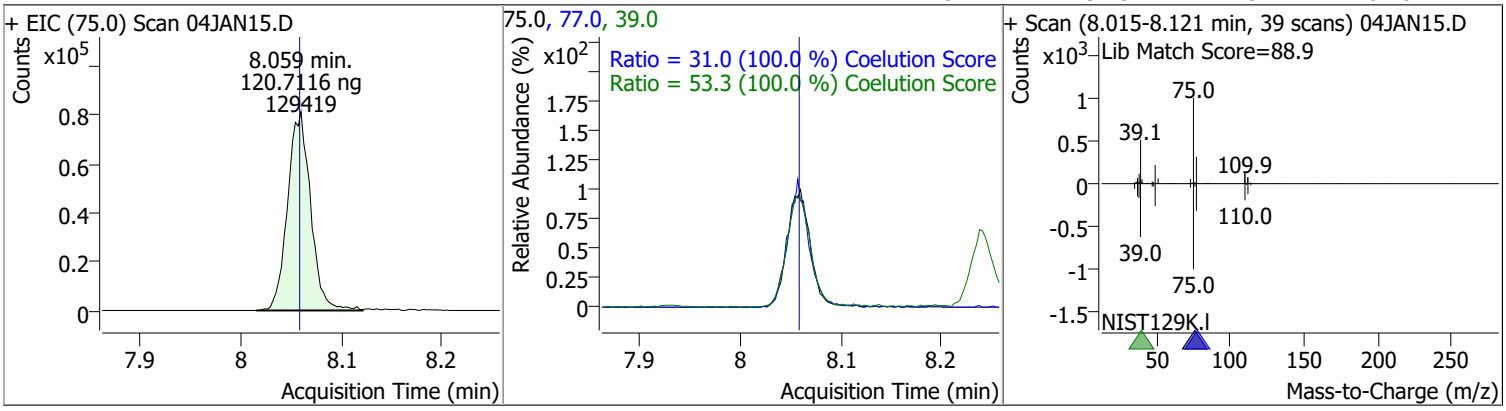




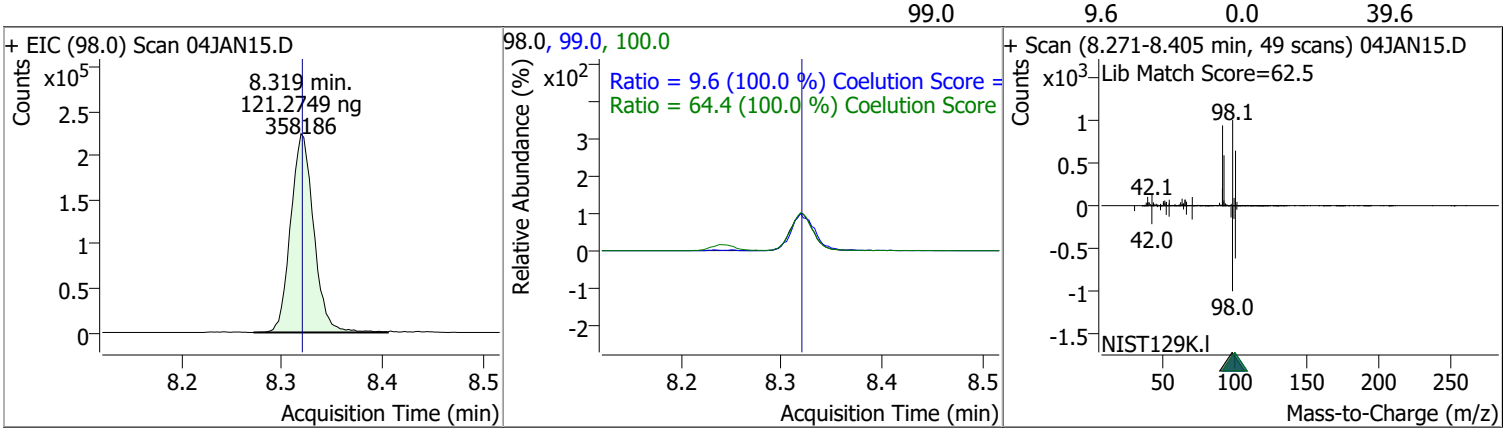
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	121.9749	7.59	0.00	115664	85.0	64.5	34.5	94.5
					127.0	9.6	0.0	39.6



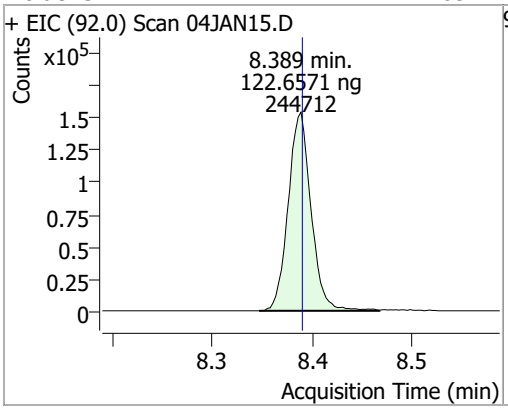
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	120.7116	8.06	0.00	129419	39.0	53.3	23.3	83.3
					77.0	31.0	1.0	61.0

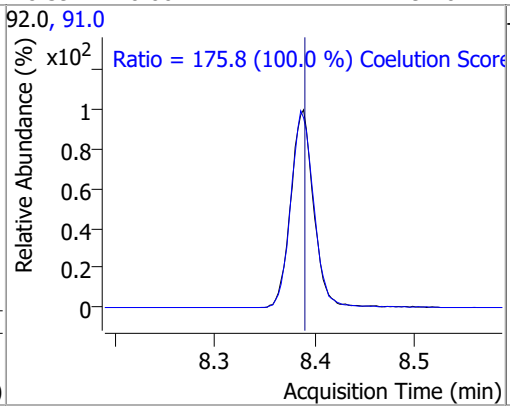


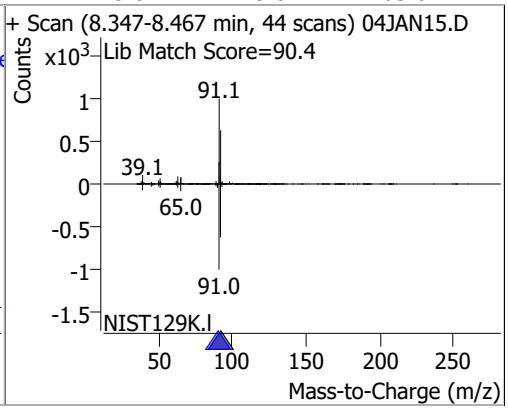
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	121.2749	8.32	0.00	358186	100.0	64.4	34.4	94.4
					99.0	9.6	0.0	39.6



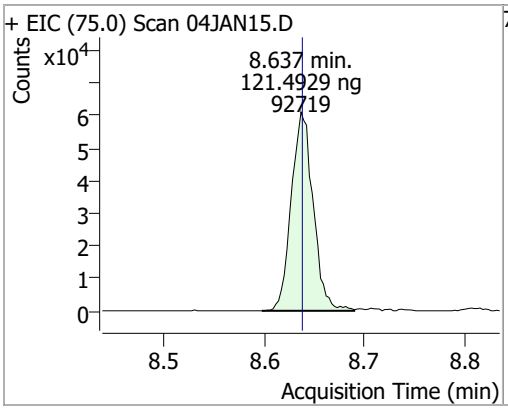
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	122.6571	8.39	0.00	244712	91.0	175.8	145.8	205.8

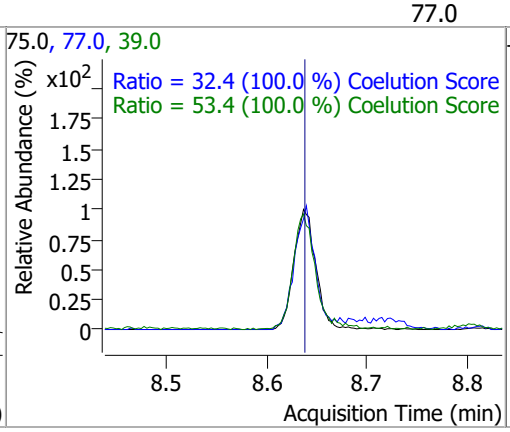


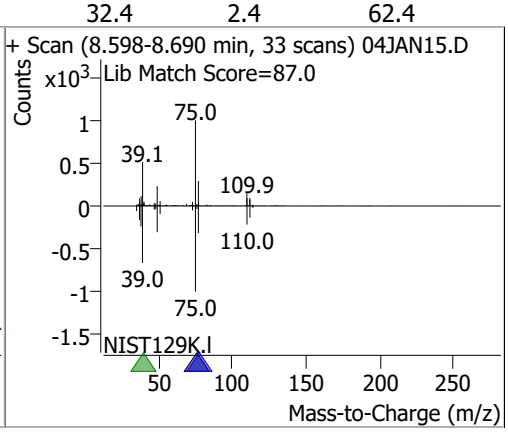




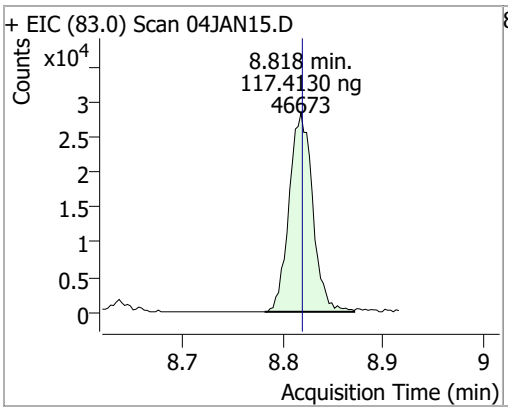
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	121.4929	8.64	0.00	92719	39.0	53.4	23.4	83.4
					77.0	32.4	2.4	62.4

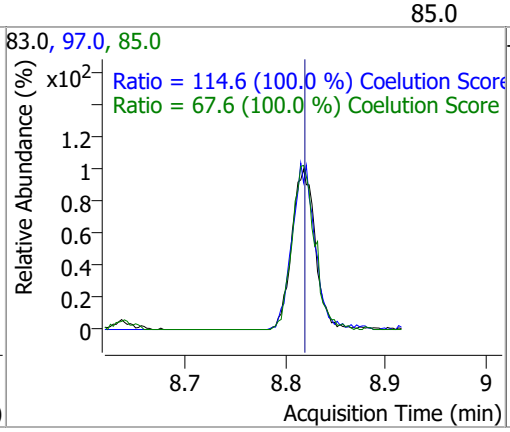


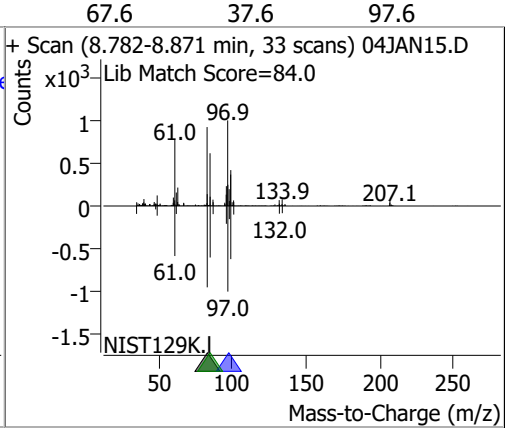




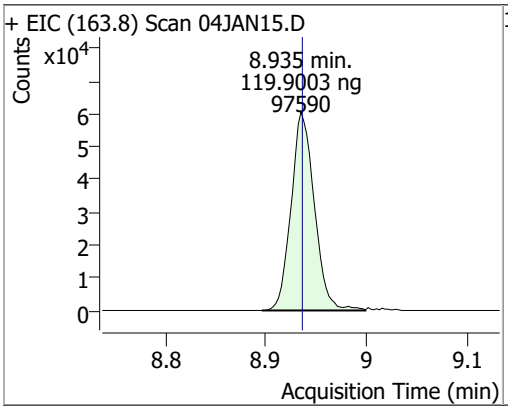
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	117.4130	8.82	0.00	46673	97.0	114.6	84.6	144.6
					85.0	67.6	37.6	97.6

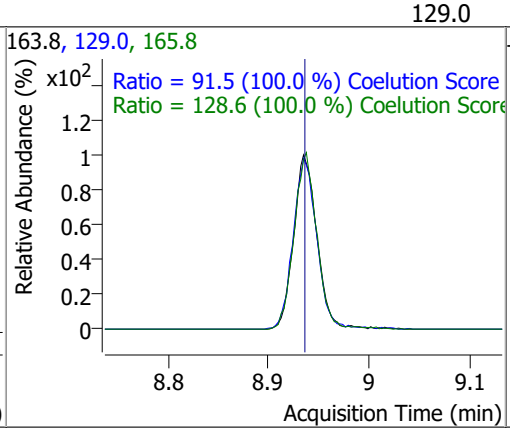


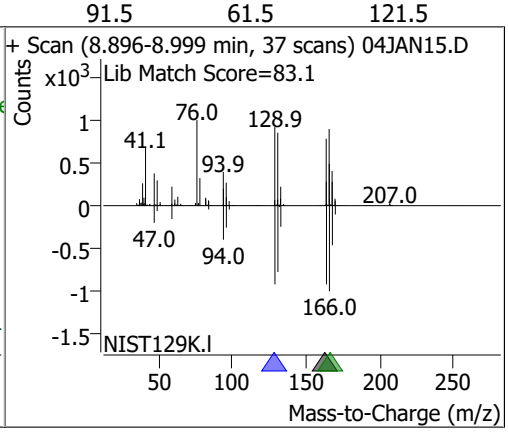




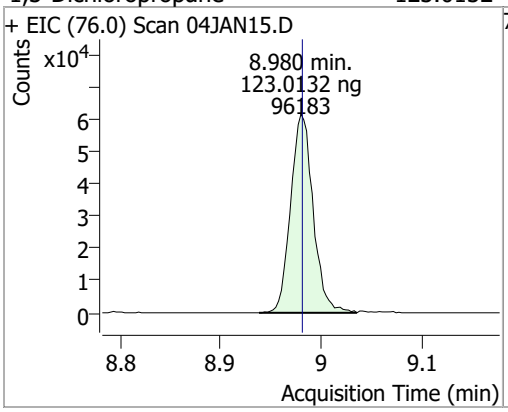
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	119.9003	8.94	0.00	97590	165.8	128.6	98.6	158.6
					129.0	91.5	61.5	121.5

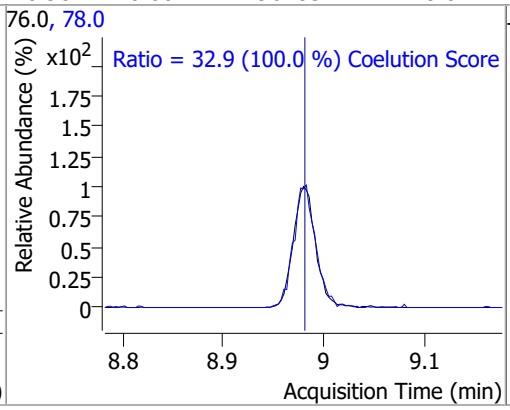


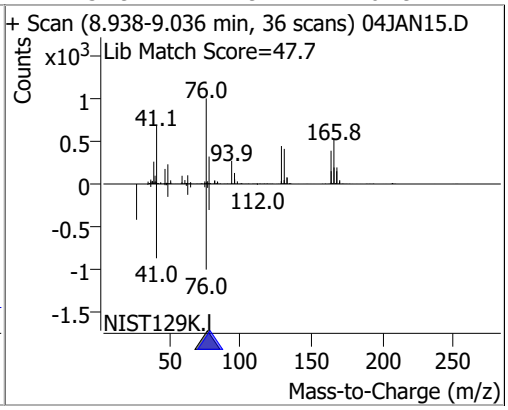




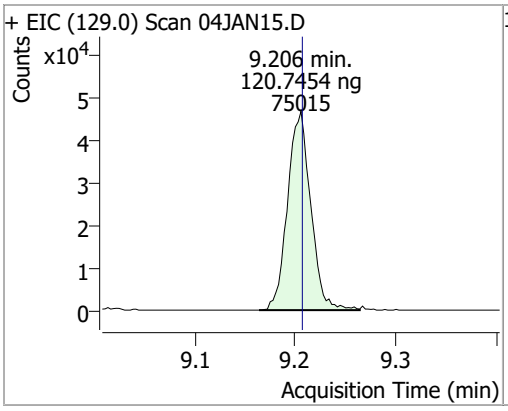
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	123.0132	8.98	0.00	96183	78.0	32.9	2.9	62.9

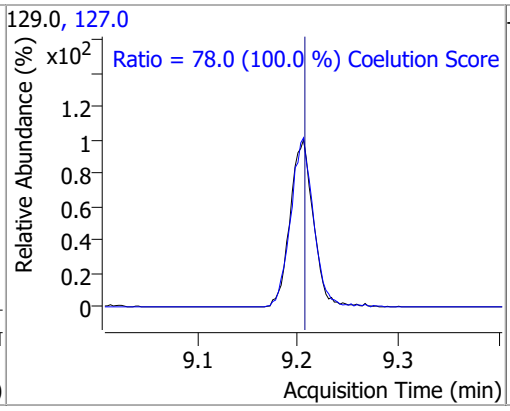


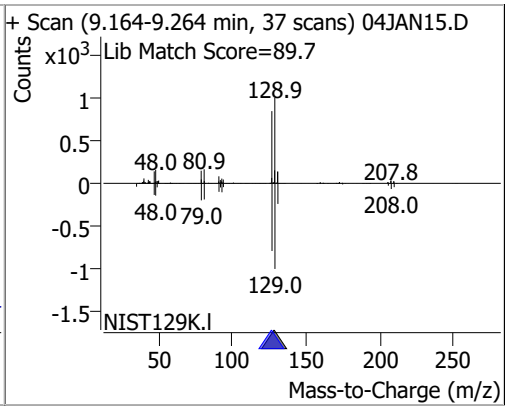




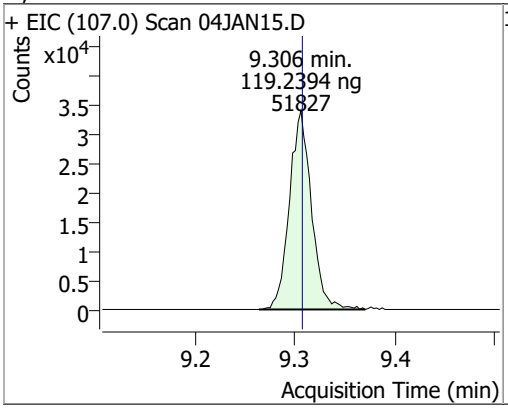
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	120.7454	9.21	0.00	75015	127.0	78.0	48.0	108.0

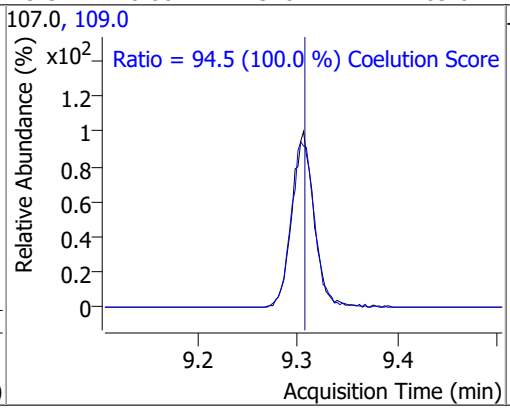


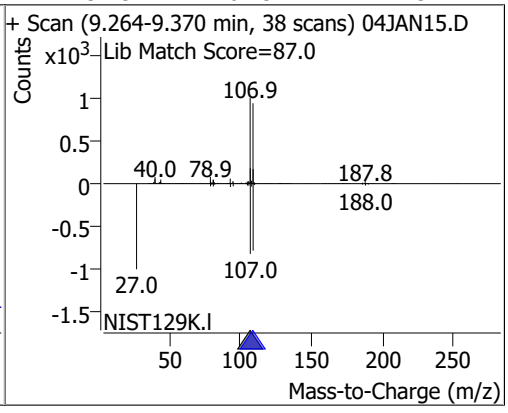




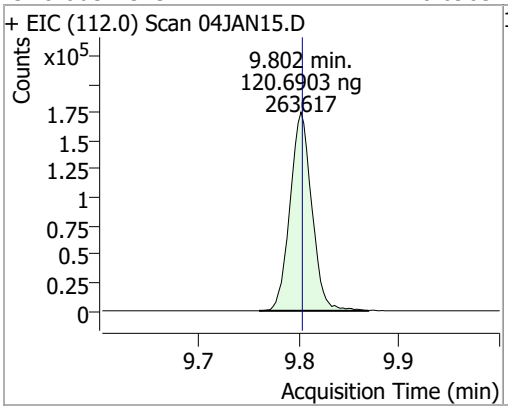
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	119.2394	9.31	0.00	51827	109.0	94.5	64.5	124.5

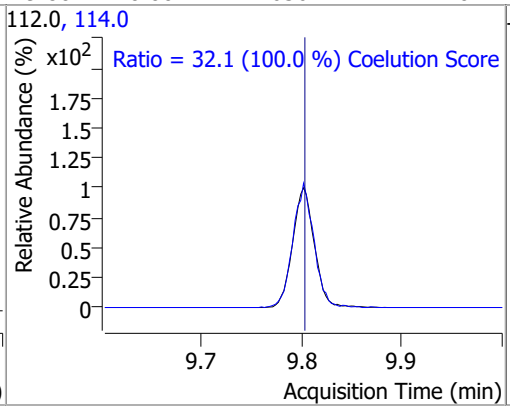


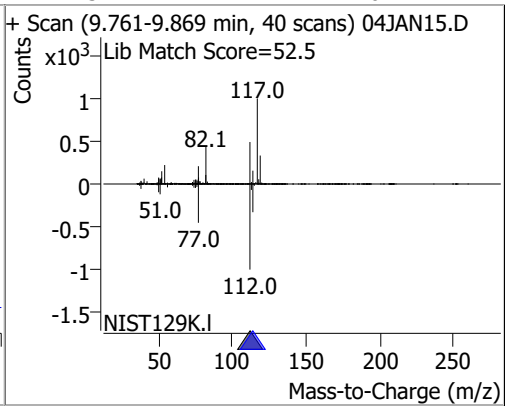




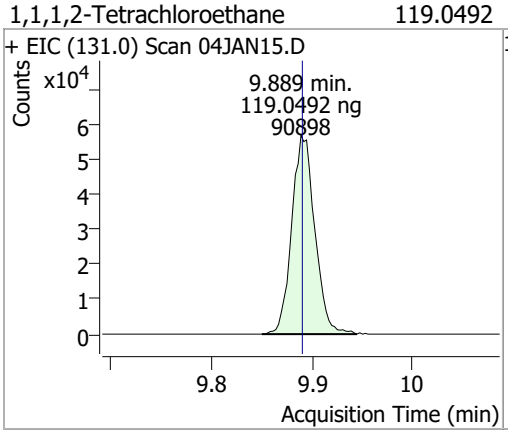
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	120.6903	9.80	0.00	263617	114.0	32.1	2.1	62.1

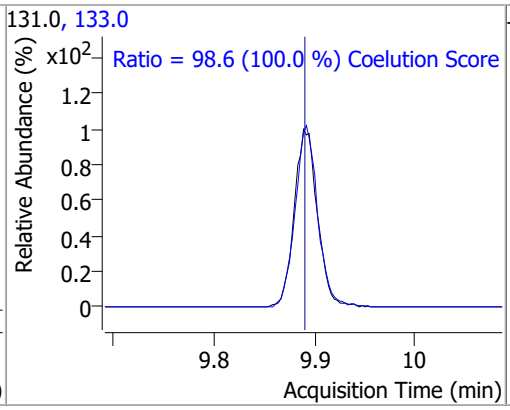


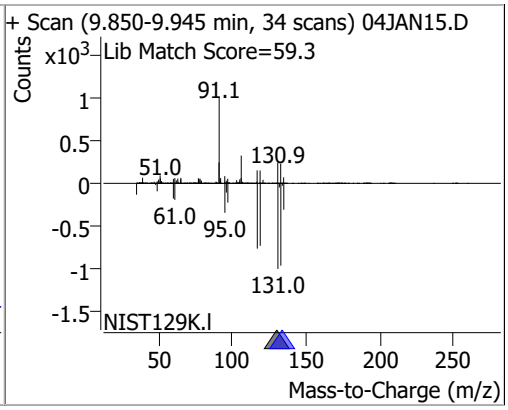




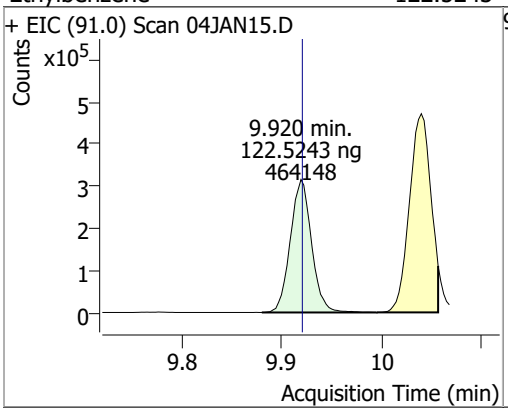
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	119.0492	9.89	0.00	90898	133.0	98.6	68.6	128.6

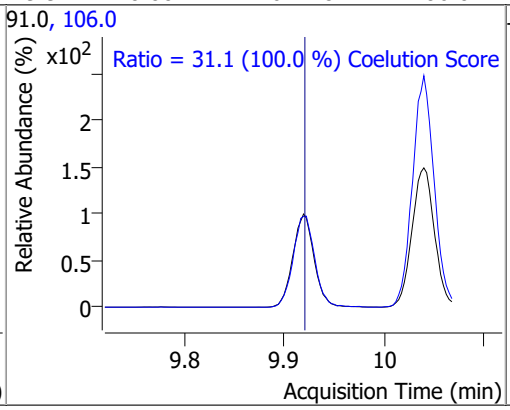


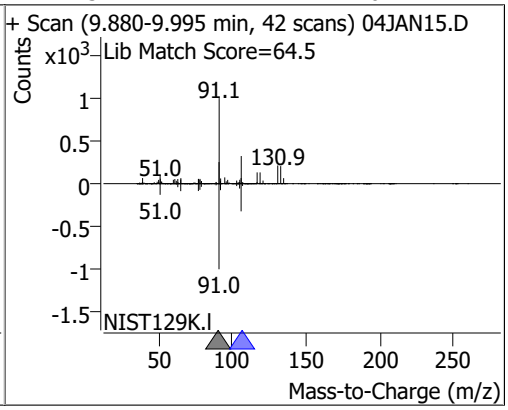




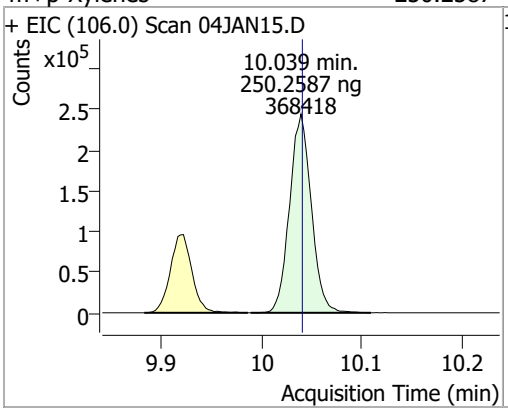
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	122.5243	9.92	0.00	464148	106.0	31.1	1.1	61.1

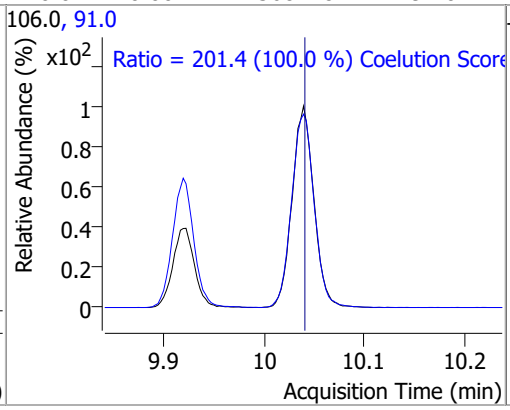


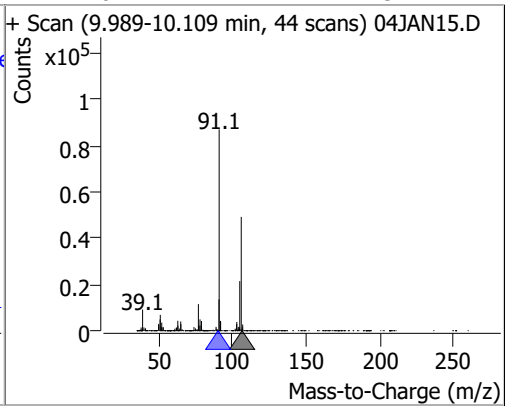




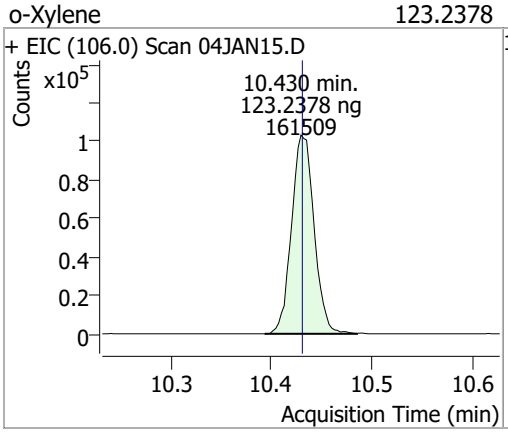
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	250.2587	10.04	0.00	368418	91.0	201.4	171.4	231.4

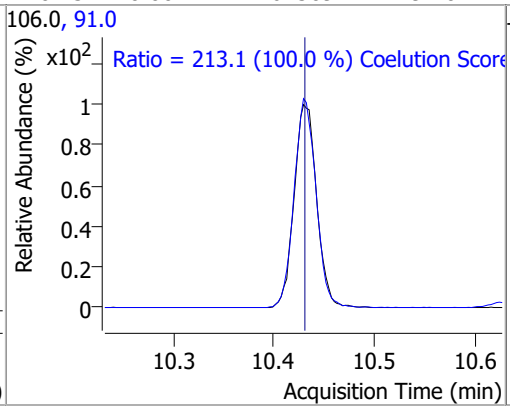


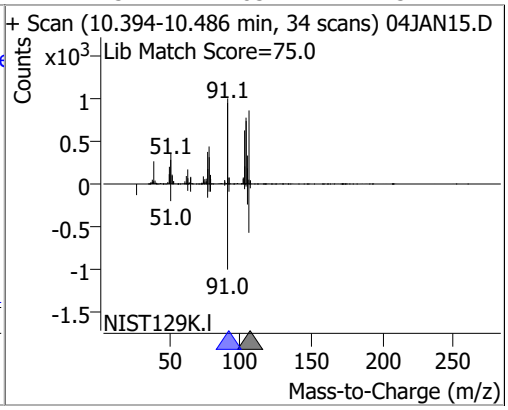




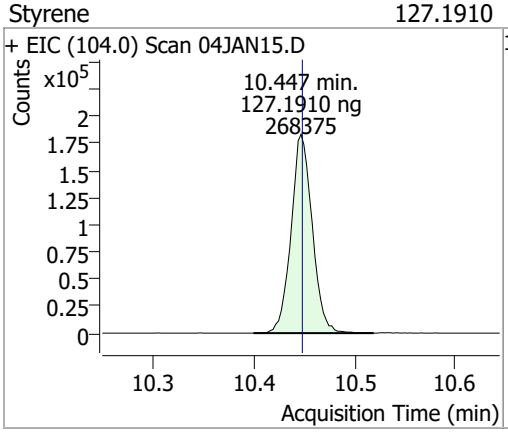
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	123.2378	10.43	0.00	161509	91.0	213.1	183.1	243.1

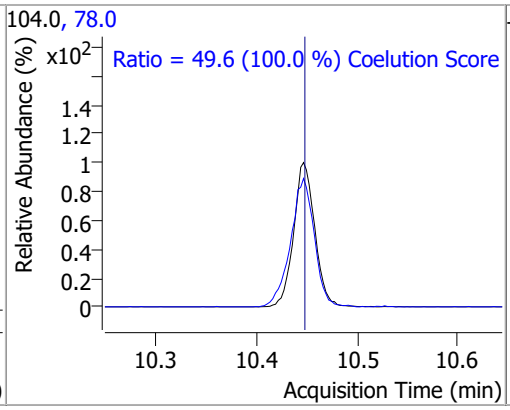


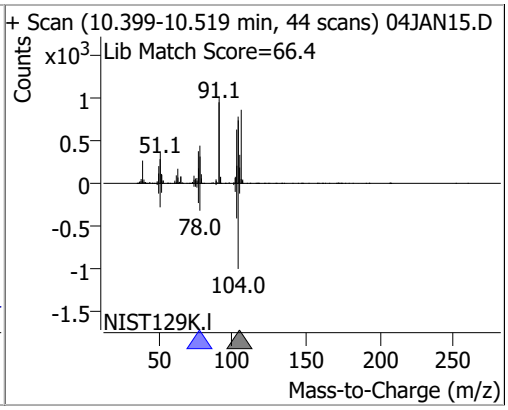




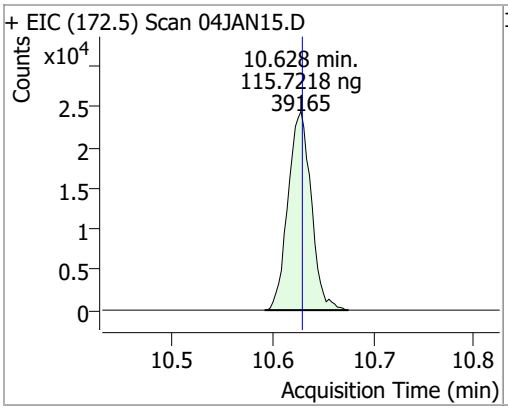
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	127.1910	10.45	0.00	268375	78.0	49.6	19.6	79.6

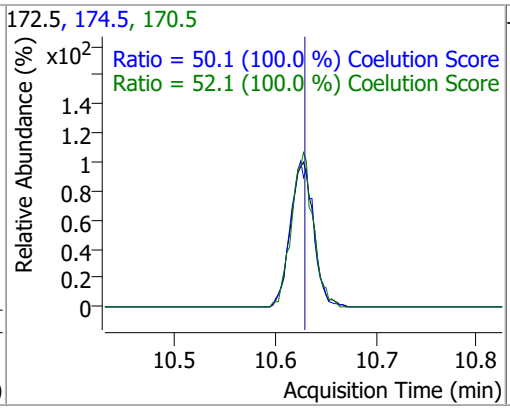


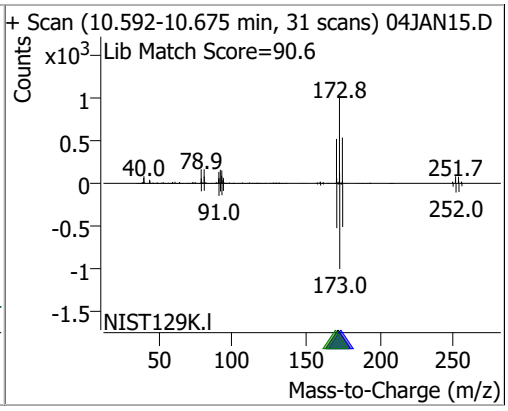




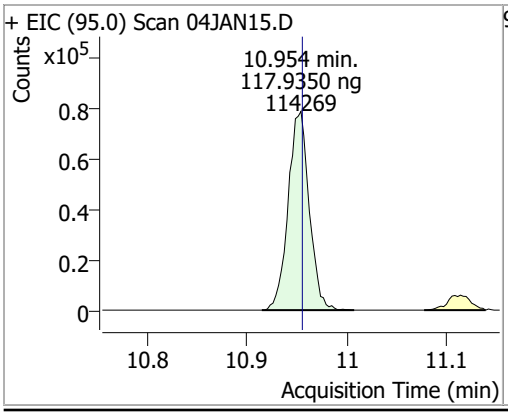
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	115.7218	10.63	0.00	39165	170.5	52.1	22.1	82.1
					174.5	50.1	20.1	80.1

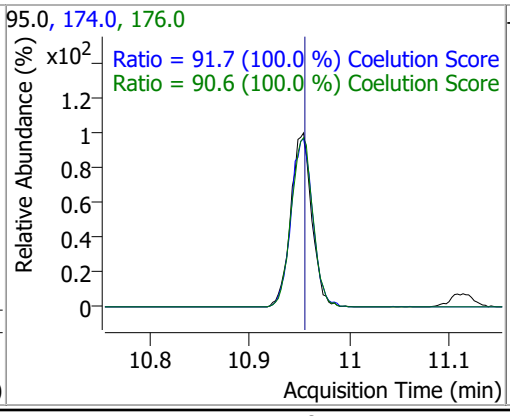


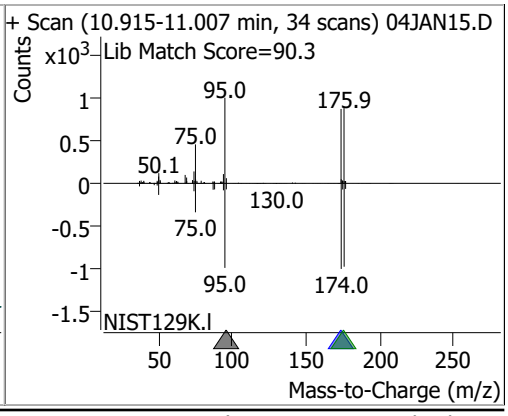




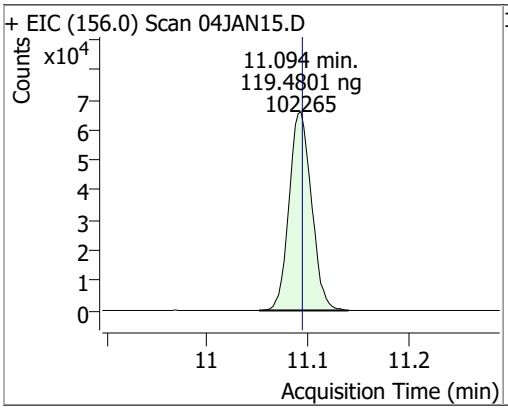
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	117.9350	10.95	0.00	114269	174.0	91.7	61.7	121.7
					176.0	90.6	60.6	120.6

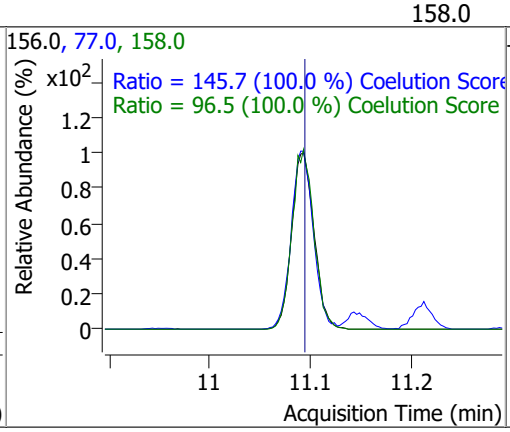


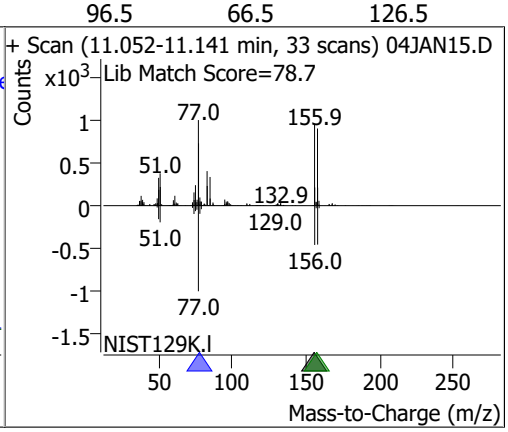




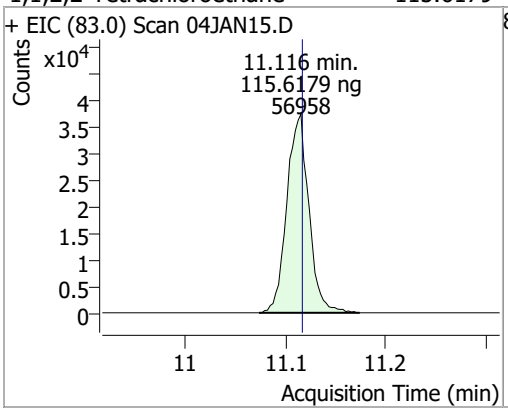
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	119.4801	11.09	0.00	102265	77.0	145.7	115.7	175.7
					158.0	96.5	66.5	126.5

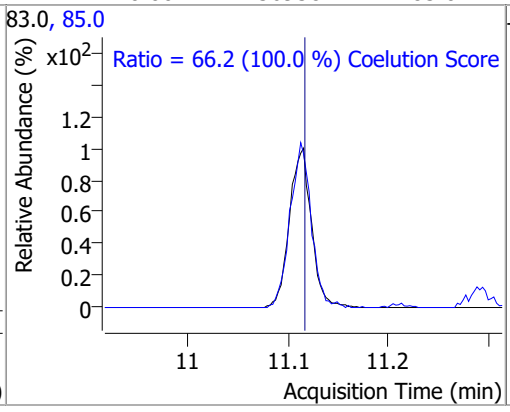


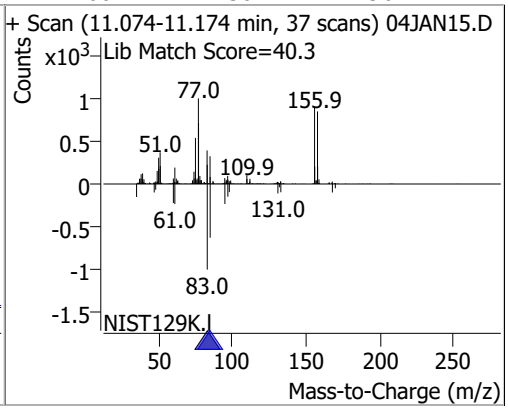




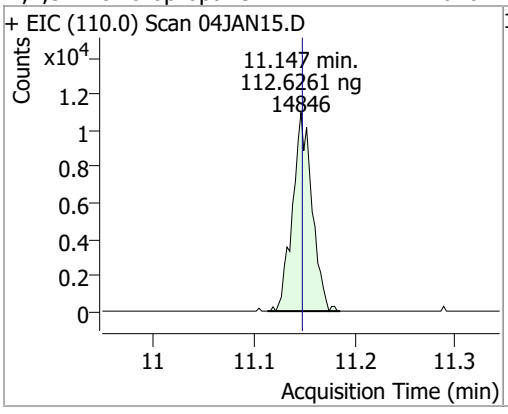
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	115.6179	11.12	0.00	56958	85.0	66.2	36.2	96.2

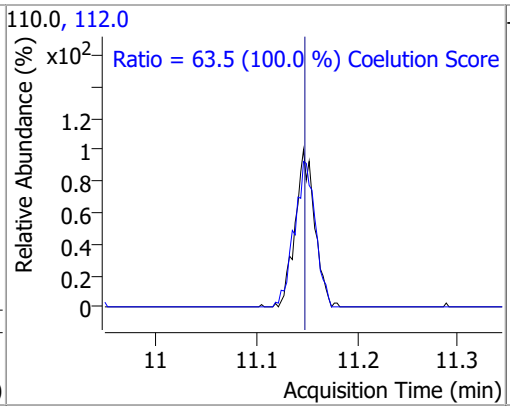


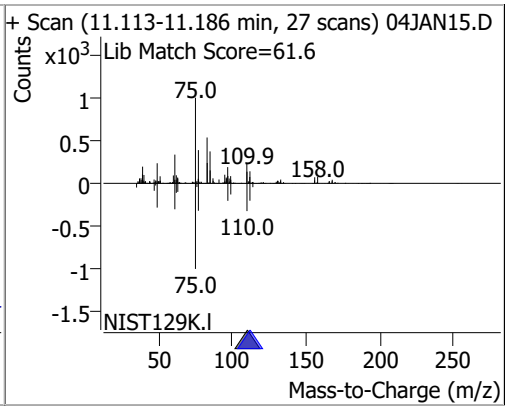




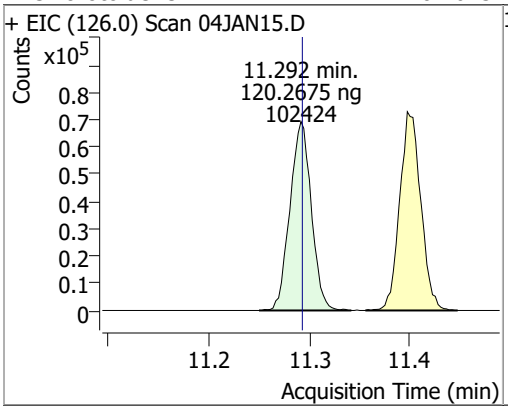
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	112.6261	11.15	0.00	14846	112.0	63.5	33.5	93.5

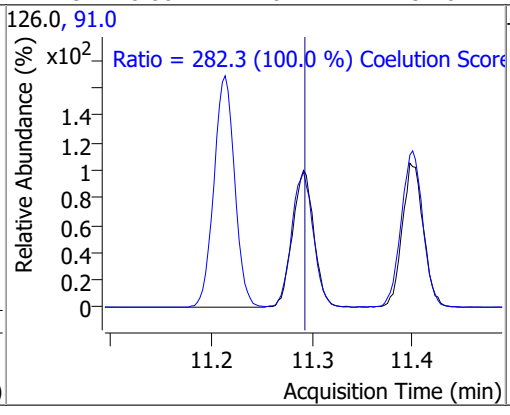


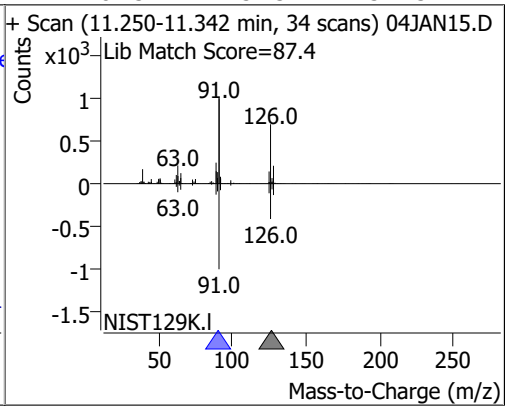




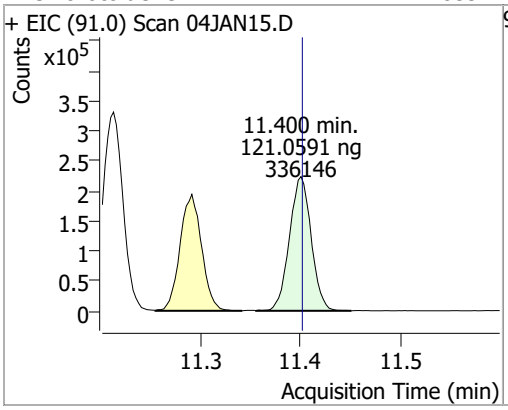
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	120.2675	11.29	0.00	102424	91.0	282.3	252.3	312.3

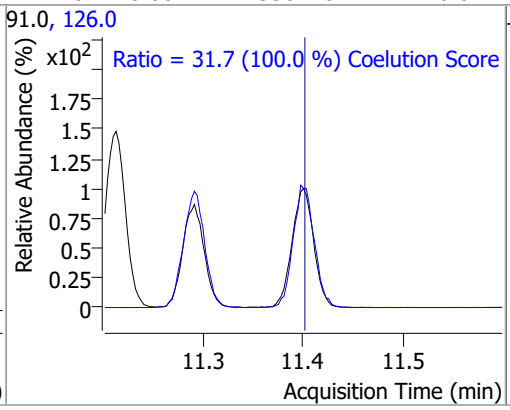


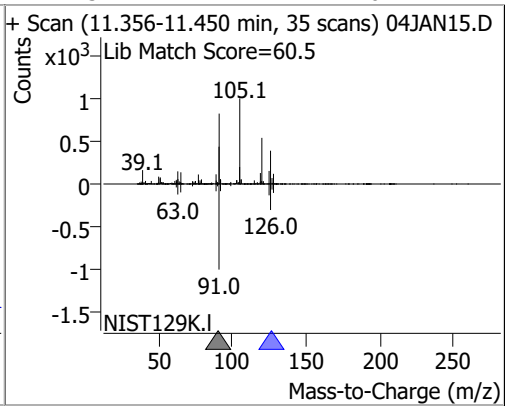




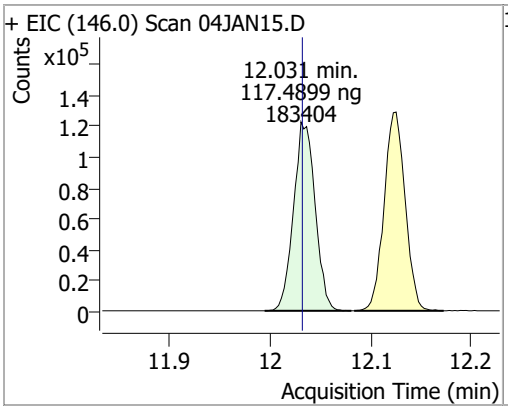
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	121.0591	11.40	0.00	336146	126.0	31.7	1.7	61.7

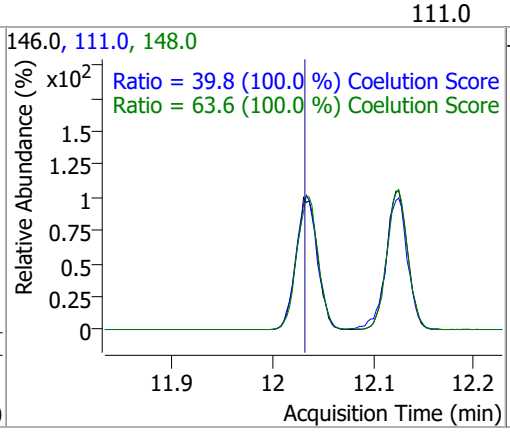


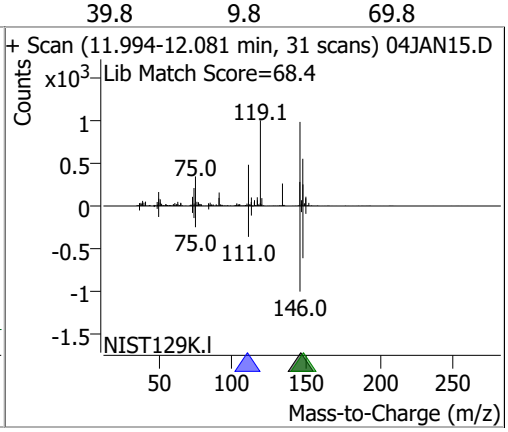




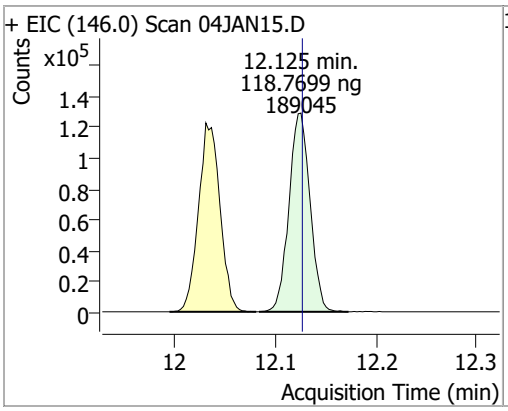
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	117.4899	12.03	0.00	183404	148.0	63.6	33.6	93.6
					111.0	39.8	9.8	69.8

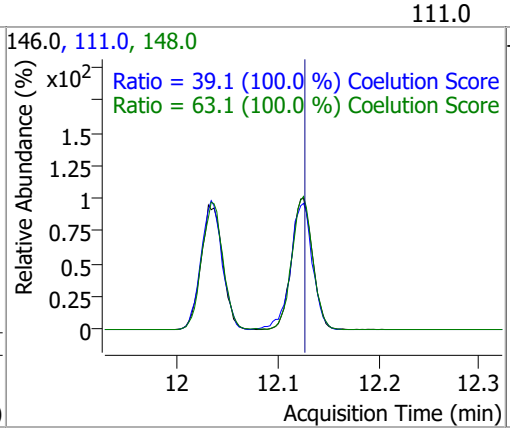


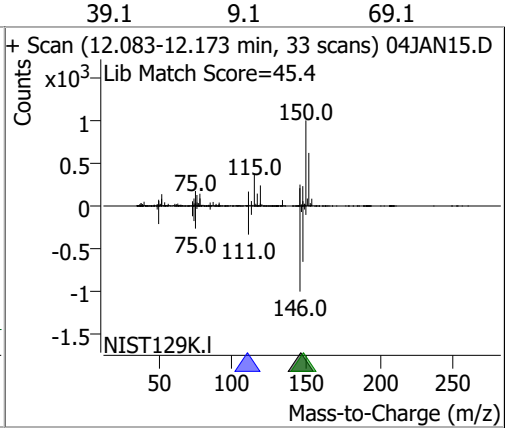




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	118.7699	12.13	0.00	189045	148.0	63.1	33.1	93.1
					111.0	39.1	9.1	69.1

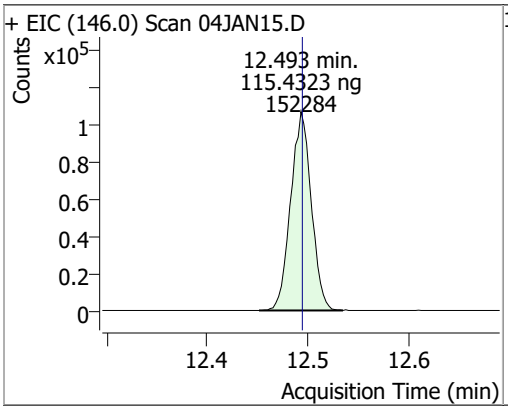


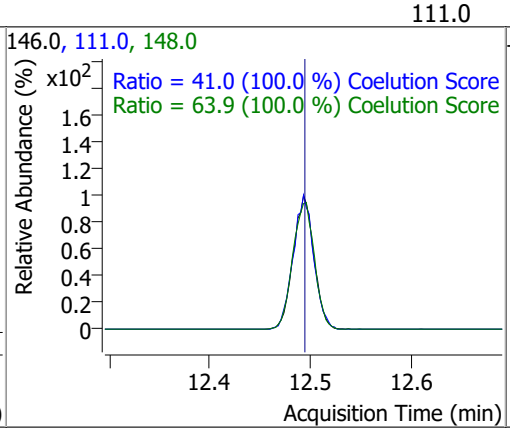


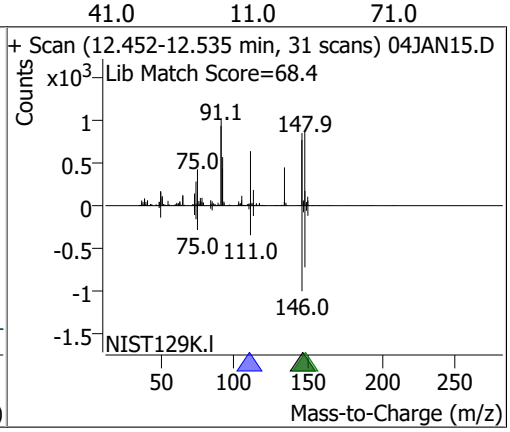


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	115.4323	12.49	0.00	152284	148.0	63.9	33.9	93.9
					111.0	41.0	11.0	71.0







Quantitation Results Report (QT Reviewed)

Agilent | Trusted Answers

Data File	04JAN17.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/4/2022 6:45:10 PM
Sample Name	ICAL010422_6	Instrument	VOA5975C
Vial	17	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
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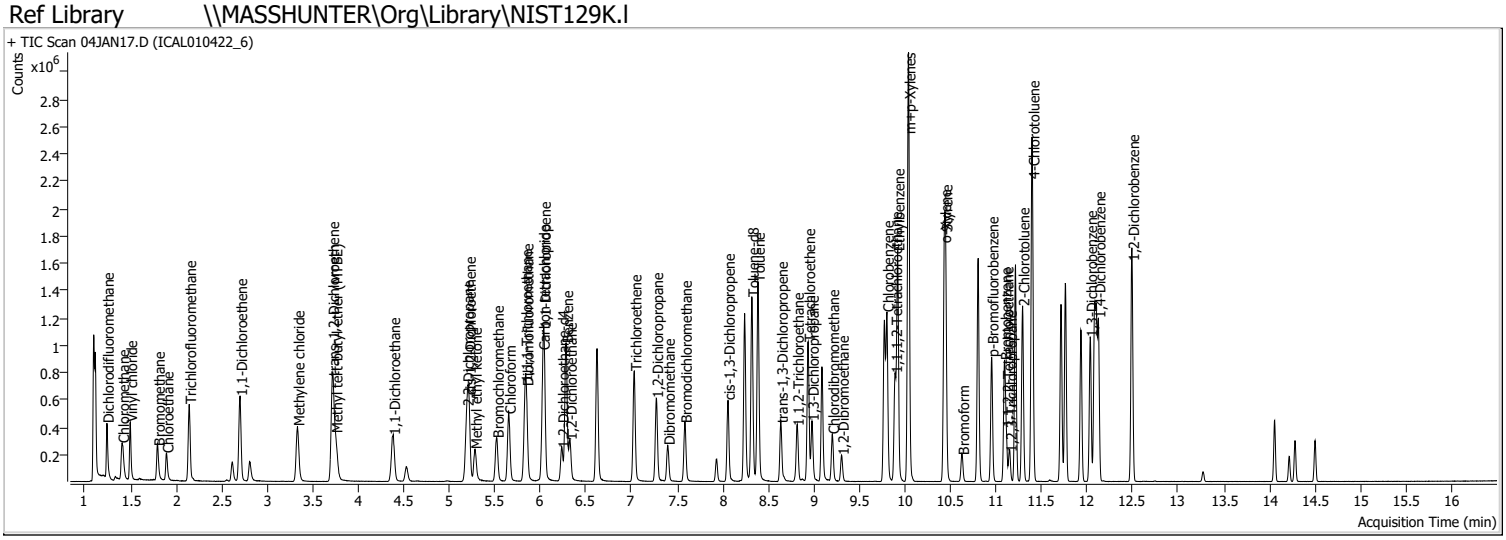
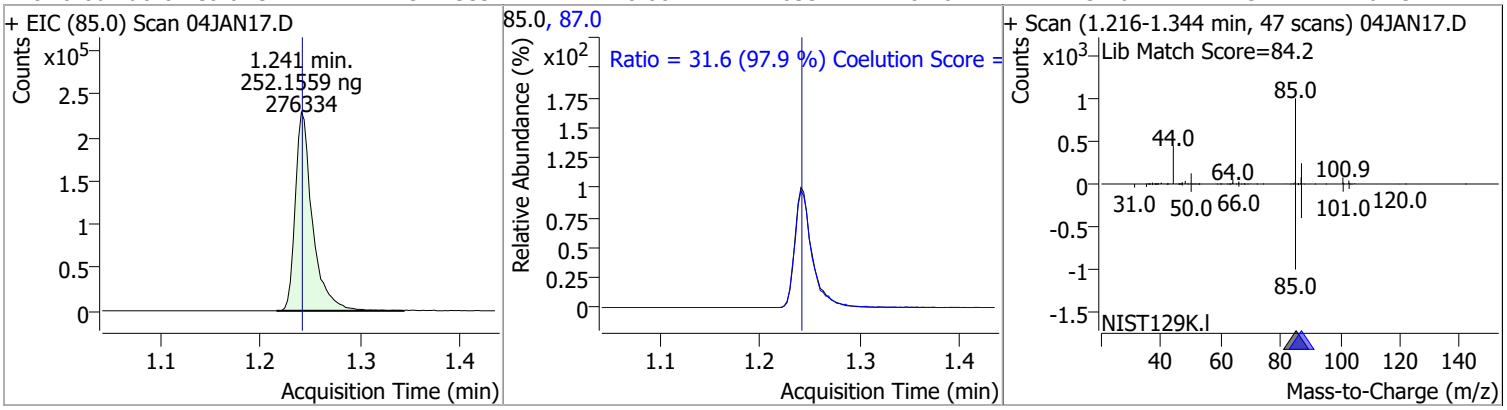


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.620	96.0	836278	250.0000	ng	-0.003
M Chlorobenzene-d5	9.774	82.0	316399	250.0000	ng	0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	266553	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.845	113.0	204073	259.0223	ng	0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 103.61%		
S 1,2-Dichloroethane-d4	6.236	67.0	87876	258.2324	ng	0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 103.29%		
S Toluene-d8	8.319	98.0	823306	270.0265	ng	0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 108.01%		
S p-Bromofluorobenzene	10.951	95.0	261042	267.3186	ng	-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 106.93%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.241	85.0	276334	252.1559	ng	99
T Chloromethane	1.408	50.0	319523	240.2183	ng	99
T Vinyl chloride	1.498	62.0	297604	248.6532	ng	86
T Bromomethane	1.799	96.0	134737	251.7606	ng	97
T Chloroethane	1.894	64.0	137312	231.7432	ng	98
T Trichlorofluoromethane	2.145	101.0	384837	259.0502	ng	98
T 1,1-Dichloroethene	2.702	96.0	217406	258.0903	ng	99
T Methylene chloride	3.333	49.0	292397	235.4657	ng	99
T trans-1,2-Dichloroethene	3.715	96.0	218855	254.6608	ng	99
T Methyl tert-butyl ether (MTBE)	3.751	73.0	287653	258.9535	ng	99
T 1,1-Dichloroethane	4.384	63.0	413408	258.4325	ng	99
T 2,2-Dichloropropane	5.190	77.0	303307	253.0397	ng	99
T cis-1,2-Dichloroethene	5.215	96.0	228170	261.8706	ng	96
T Methyl ethyl ketone	5.279	43.0	317271	2688.2474	ng	99
T Bromochloromethane	5.519	128.0	89178	247.0586	ng	95
T Chloroform	5.653	83.0	394946	248.0804	ng	99

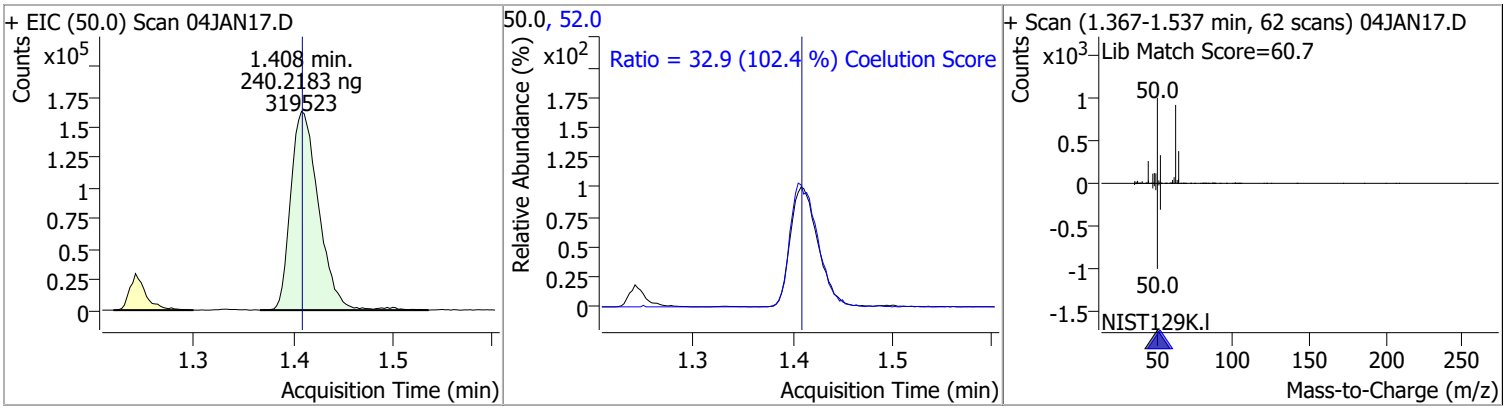
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.831	97.0	386005	258.7228	ng	99
T Carbon tetrachloride	6.026	117.0	383485	260.8774	ng	99
T 1,1-Dichloropropene	6.038	75.0	335741	264.6638	ng	99
T Benzene	6.280	78.0	857534	257.5416	ng	100
T 1,2-Dichloroethane	6.322	62.0	226964	251.9675	ng	99
T Trichloroethene	7.030	95.0	250285	262.2931	ng	100
T 1,2-Dichloropropane	7.270	63.0	213800	254.7161	ng	100
T Dibromomethane	7.396	93.0	89483	252.2734	ng	97
T Bromodichloromethane	7.582	83.0	251805	257.2286	ng	100
T cis-1,3-Dichloropropene	8.057	75.0	293617	265.2863	ng	99
T Toluene	8.386	92.0	541945	263.1330	ng	100
T trans-1,3-Dichloropropene	8.639	75.0	207833	263.8027	ng	98
T 1,1,2-Trichloroethane	8.815	83.0	101888	248.2882	ng	99
T Tetrachloroethene	8.938	163.8	218245	259.7419	ng	98
T 1,3-Dichloropropane	8.980	76.0	212669	263.4754	ng	98
T Chlorodibromomethane	9.203	129.0	165695	258.3535	ng	100
T 1,2-Dibromoethane	9.306	107.0	115714	257.8887	ng	100
T Chlorobenzene	9.802	112.0	582326	258.2544	ng	100
T 1,1,1,2-Tetrachloroethane	9.891	131.0	200859	254.8274	ng	100
T Ethylbenzene	9.919	91.0	1043443	266.8193	ng	100
T m+p-Xylenes	10.039	106.0	825866	543.4262	ng	100
T o-Xylene	10.430	106.0	365914	270.4636	ng	100
T Styrene	10.446	104.0	605646	278.0455	ng	99
T Bromoform	10.628	172.5	87836	257.5099	ng	100
T Bromobenzene	11.093	156.0	227127	263.2944	ng	99
T 1,1,2,2-Tetrachloroethane	11.113	83.0	124205	250.1577	ng	97
T 1,2,3-Trichloropropane	11.152	110.0	33115	249.2635	ng	97
T 2-Chlorotoluene	11.291	126.0	229396	267.2616	ng	99
T 4-Chlorotoluene	11.400	91.0	748435	267.4409	ng	99
T 1,3-Dichlorobenzene	12.033	146.0	406895	258.6297	ng	99
T 1,4-Dichlorobenzene	12.122	146.0	408934	254.9170	ng	98
T 1,2-Dichlorobenzene	12.493	146.0	342576	257.6524	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

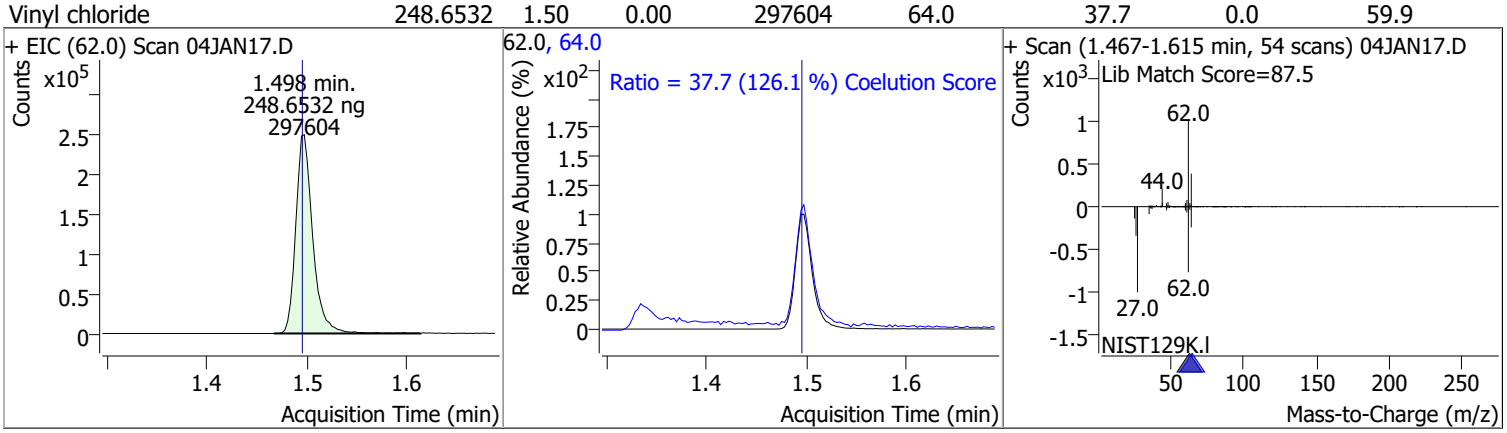
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	252.1559	1.24	0.00	276334	87.0	31.6	2.3	62.3



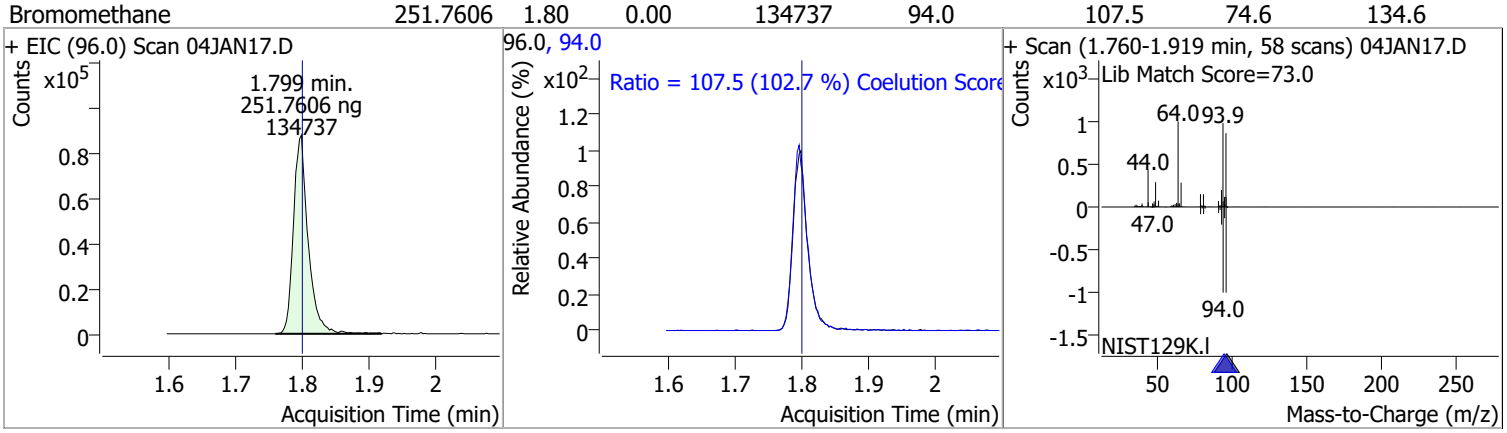
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	240.2183	1.41	0.00	319523	52.0	32.9	2.1	62.1



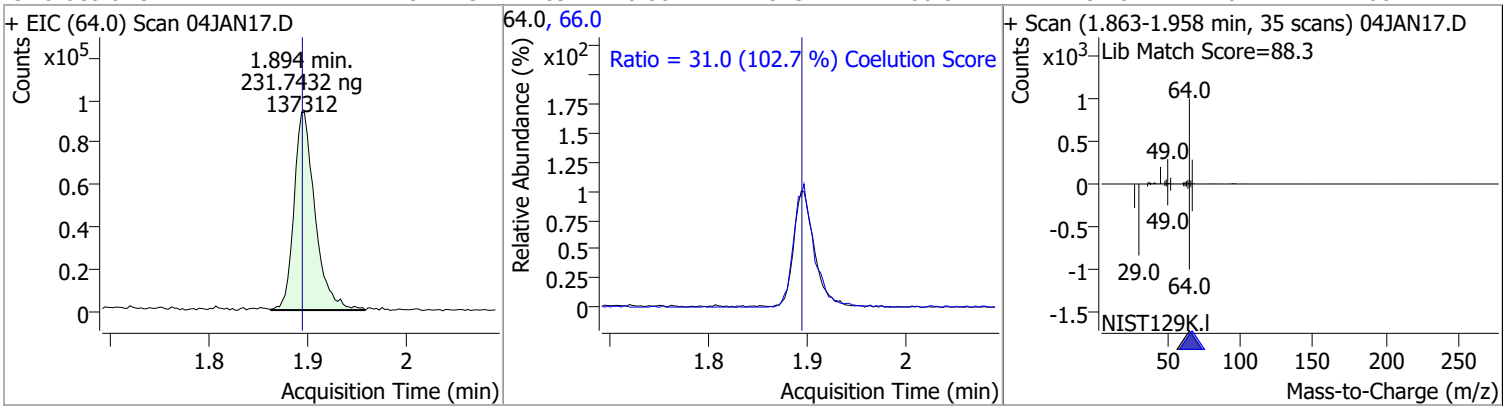
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	248.6532	1.50	0.00	297604	64.0	37.7	0.0	59.9



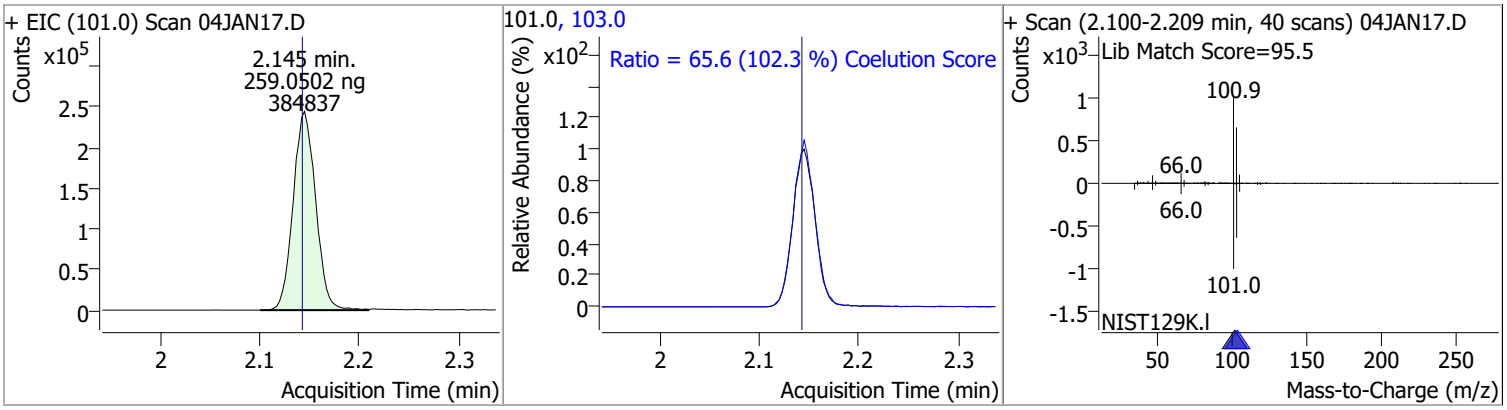
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	251.7606	1.80	0.00	134737	94.0	107.5	74.6	134.6



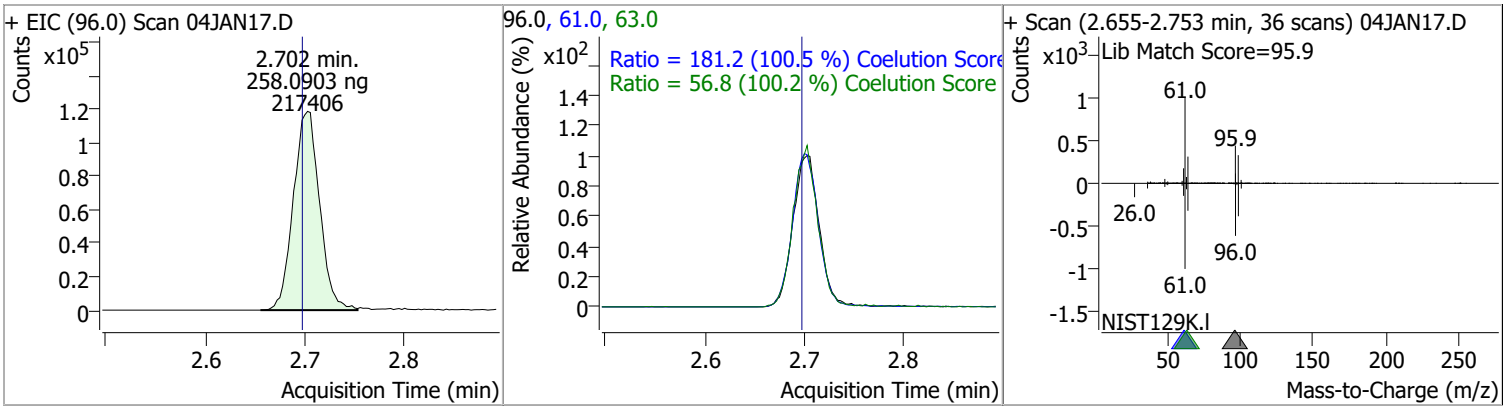
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	231.7432	1.89	0.00	137312	66.0	31.0	0.1	60.1



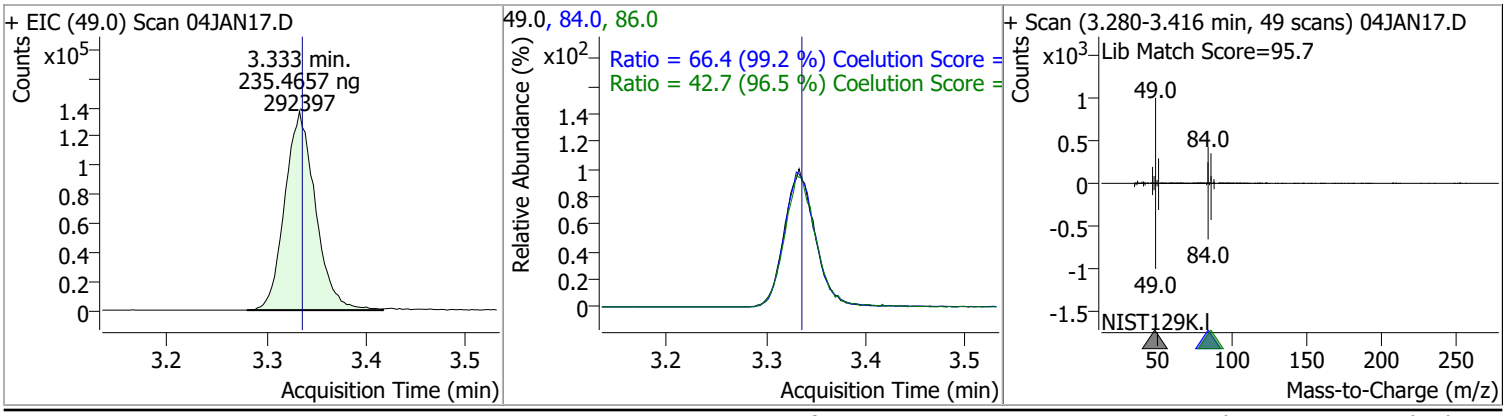
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	259.0502	2.14	0.00	384837	103.0	65.6	34.2	94.2



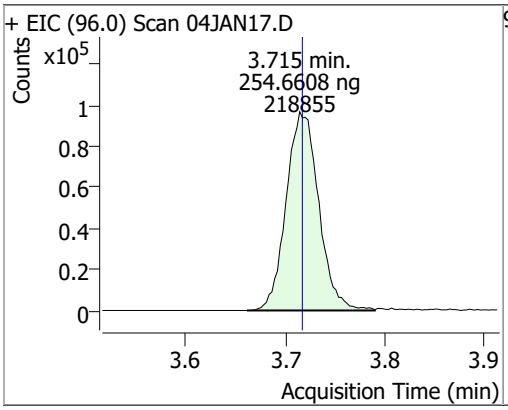
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	258.0903	2.70	0.01	217406	61.0	181.2	150.3	210.3
					63.0	56.8	26.7	86.7

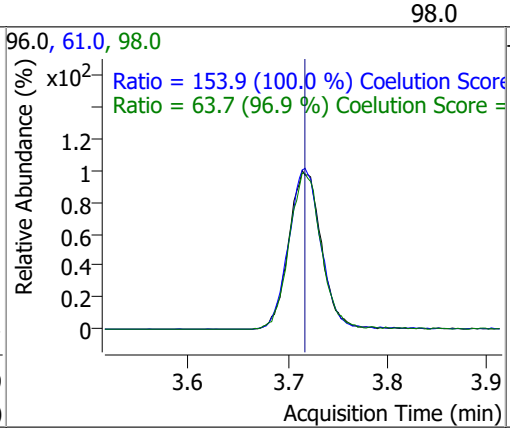


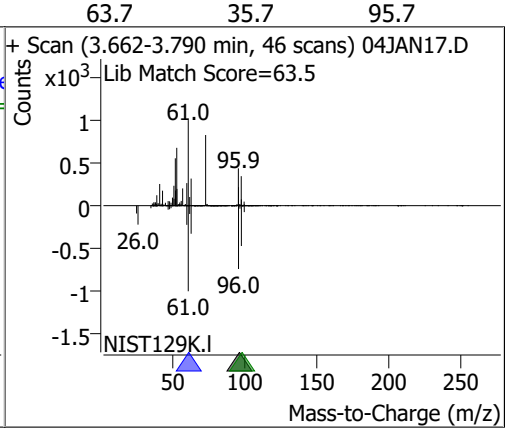
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	235.4657	3.33	0.00	292397	84.0	66.4	36.9	96.9
					86.0	42.7	14.3	74.3



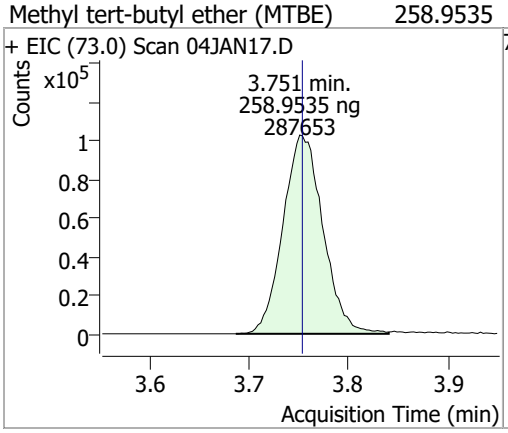
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	254.6608	3.71	0.00	218855	61.0	153.9	123.9	183.9
					98.0	63.7	35.7	95.7

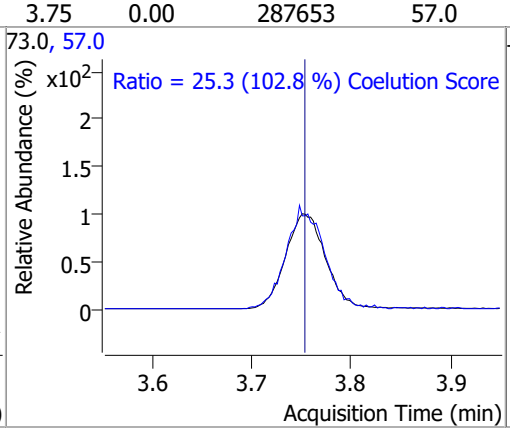


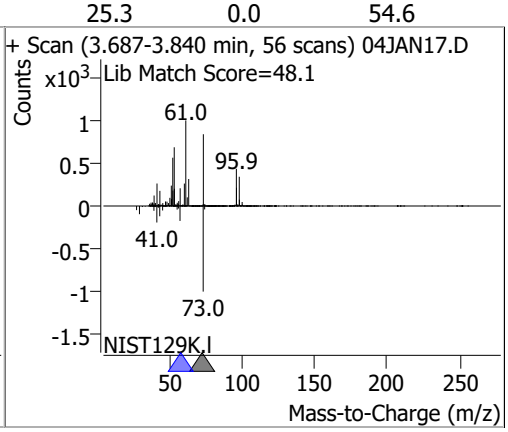




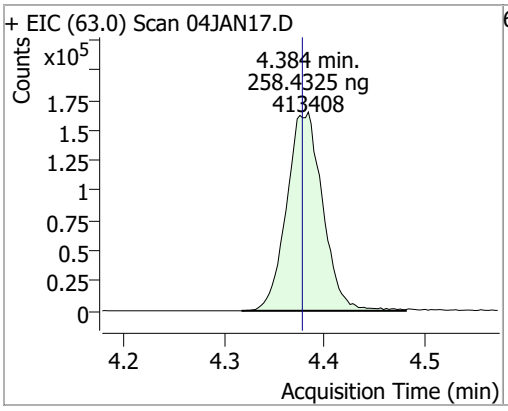
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	258.9535	3.75	0.00	287653	57.0	25.3	0.0	54.6

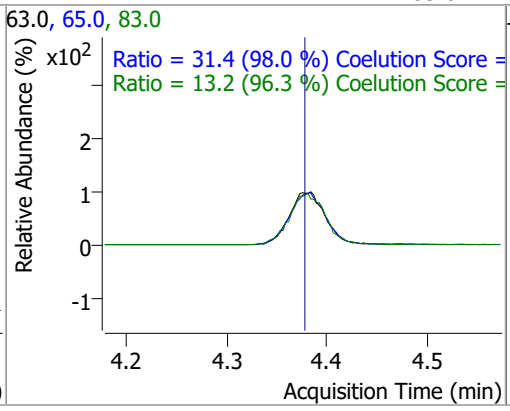


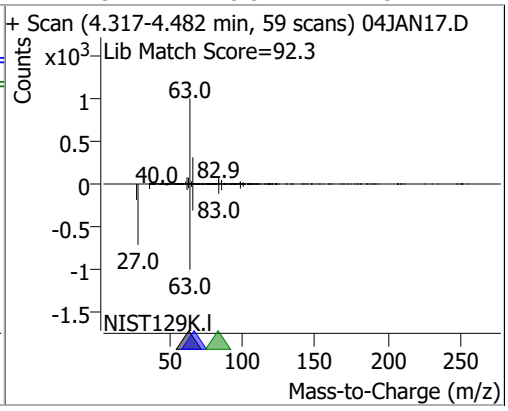




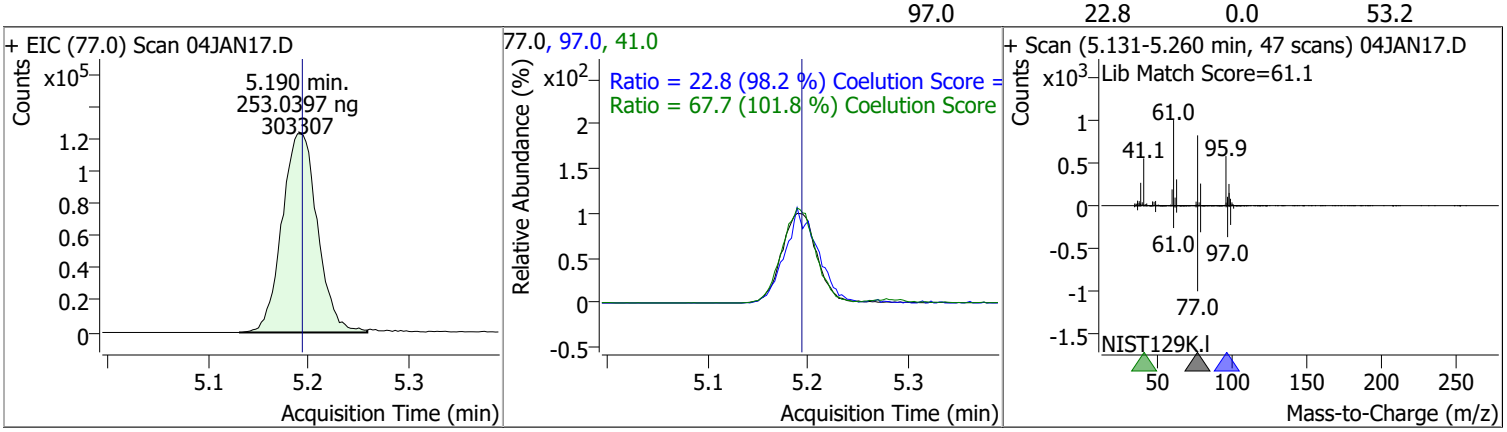
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	258.4325	4.38	0.01	413408	65.0	31.4	2.1	62.1
					83.0	13.2	0.0	43.7



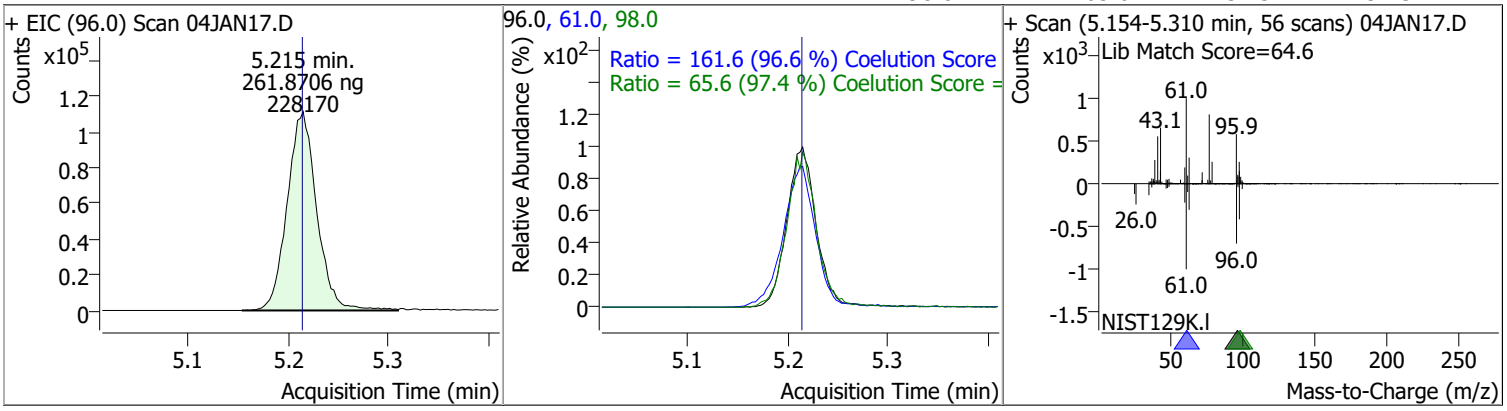




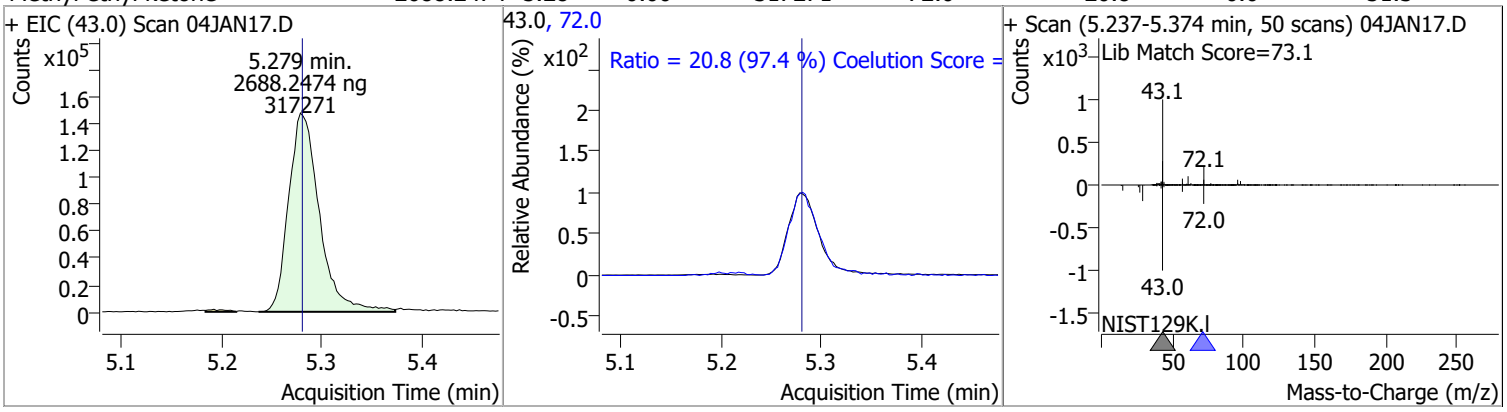
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	253.0397	5.19	-0.01	303307	41.0	67.7	36.5	96.5
					97.0	22.8	0.0	53.2



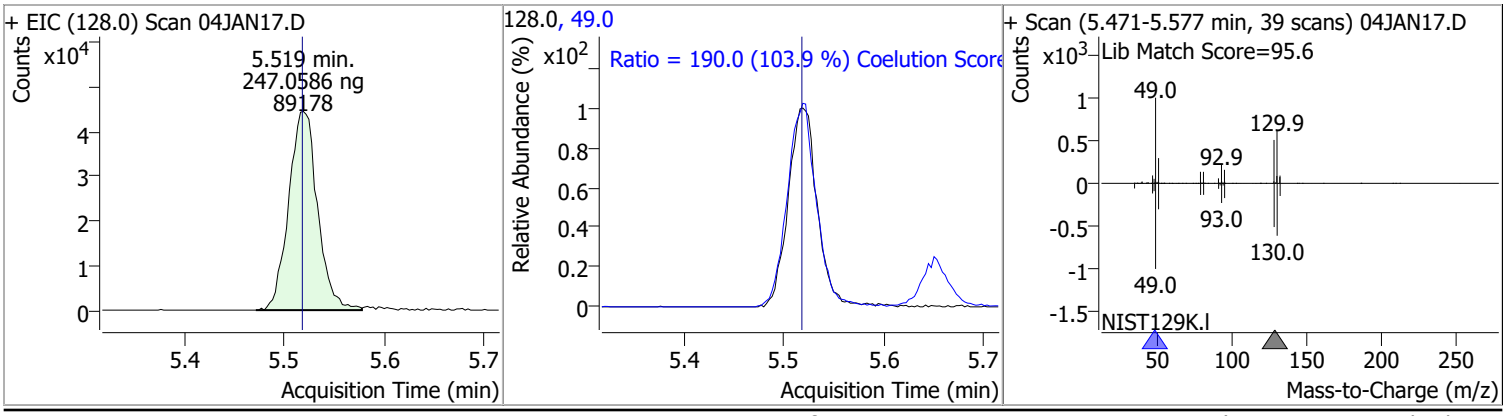
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	261.8706	5.21	0.00	228170	61.0	161.6	137.2	197.2
					98.0	65.6	37.3	97.3



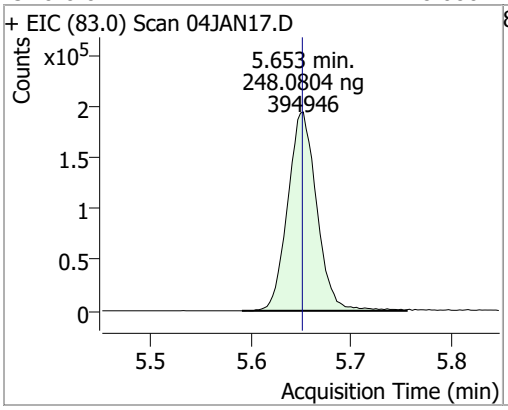
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	2688.2474	5.28	0.00	317271	72.0	20.8	0.0	51.3

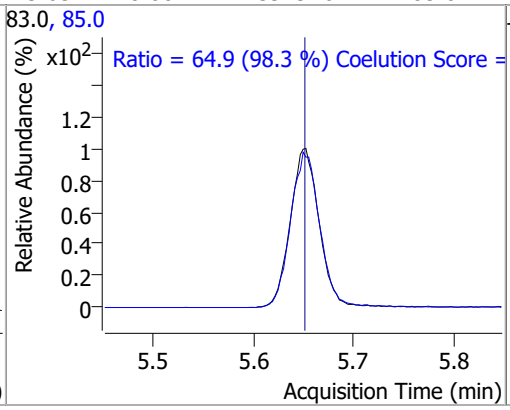


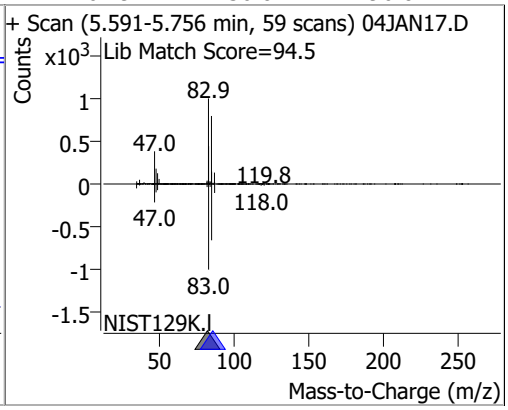
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	247.0586	5.52	0.00	89178	49.0	190.0	152.9	212.9



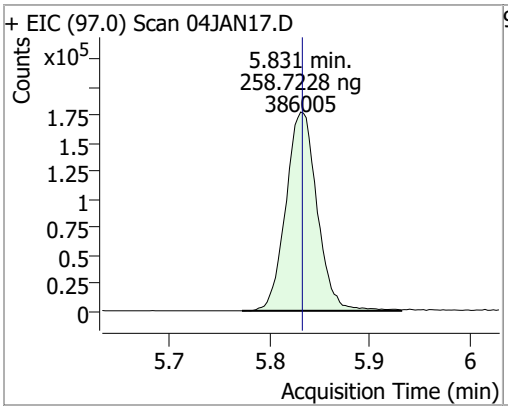
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	248.0804	5.65	0.00	394946	85.0	64.9	36.0	96.0

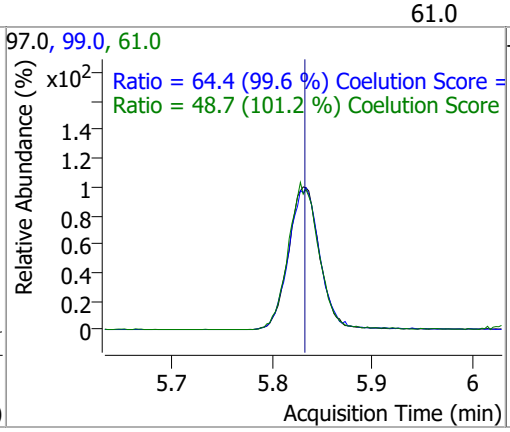


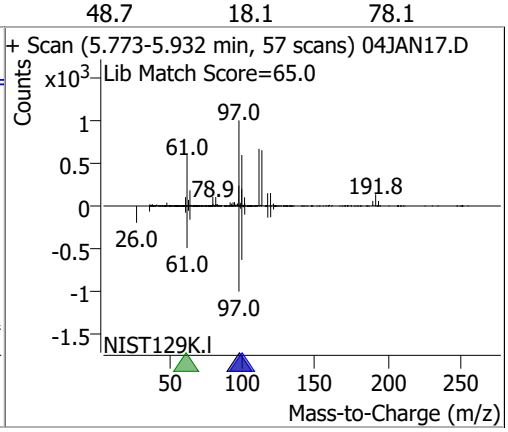




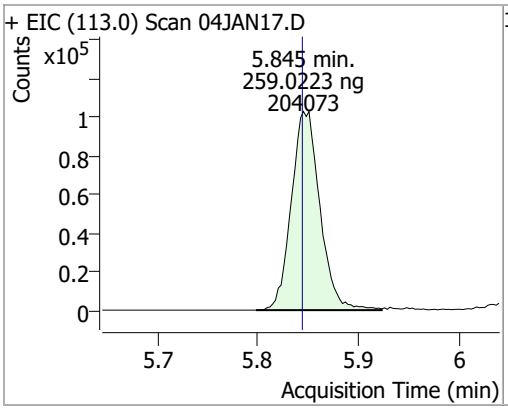
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	258.7228	5.83	0.00	386005	99.0	64.4	34.7	94.7
					61.0	48.7	18.1	78.1

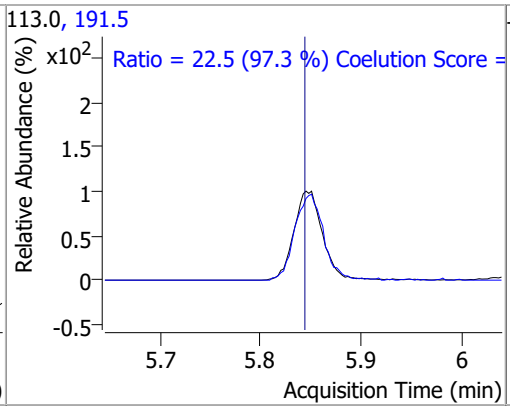


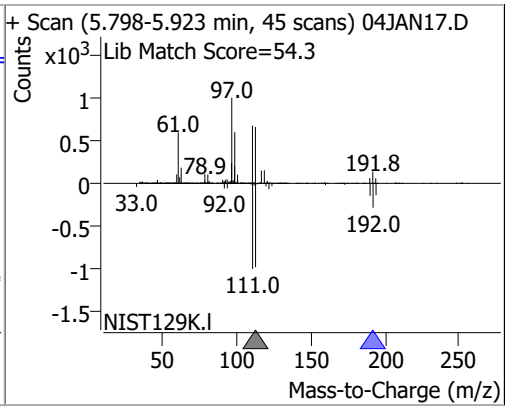




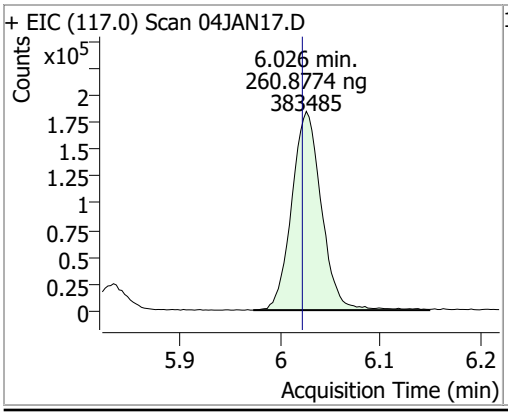
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	259.0223	5.85	0.00	204073	191.5	22.5	0.0	53.1

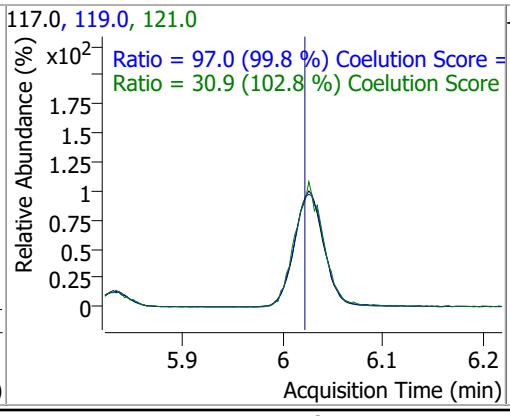


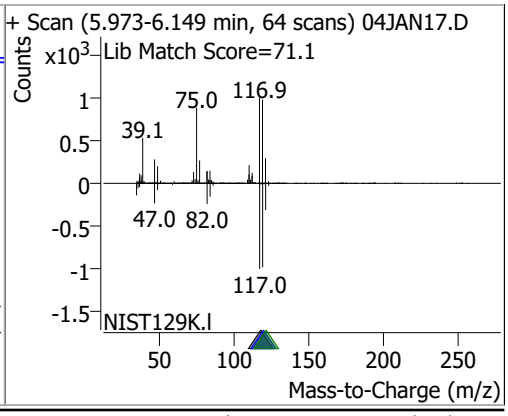




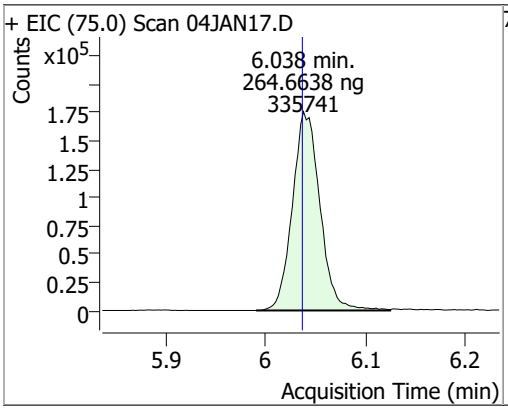
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	260.8774	6.03	0.00	383485	119.0	97.0	67.2	127.2
					121.0	30.9	0.1	60.1

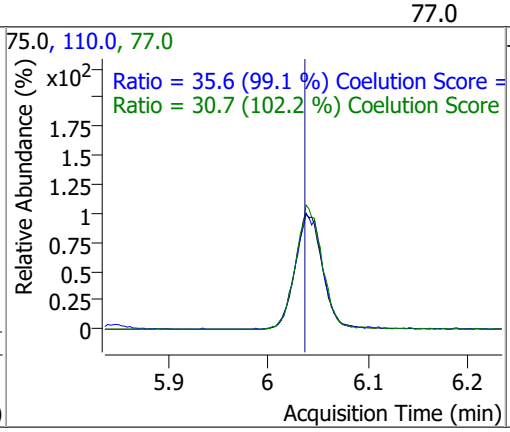


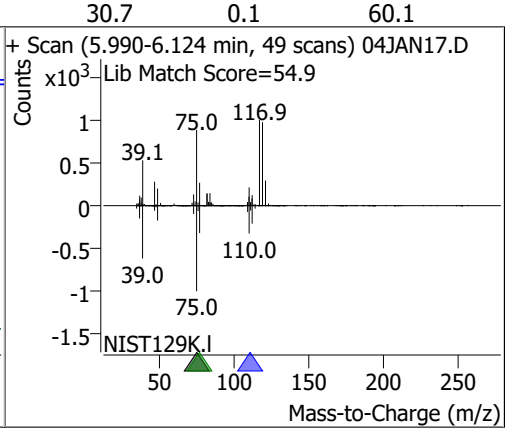




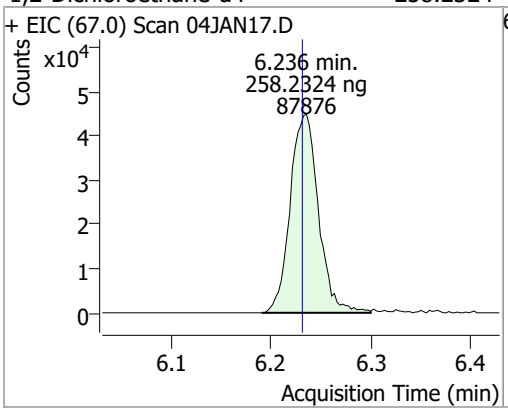
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	264.6638	6.04	0.00	335741	110.0	35.6	5.9	65.9
					77.0	30.7	0.1	60.1

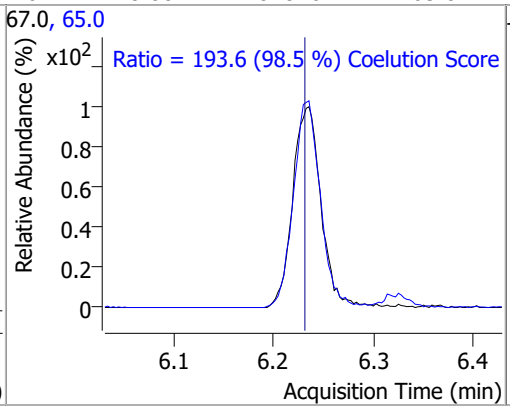


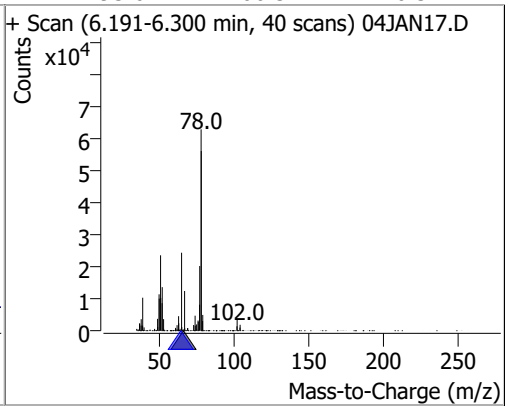




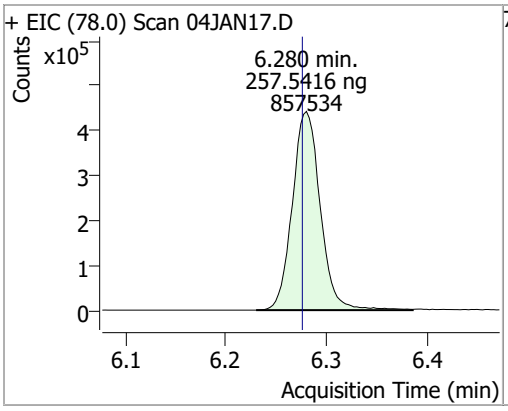
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	258.2324	6.24	0.00	87876	65.0	193.6	166.5	226.5

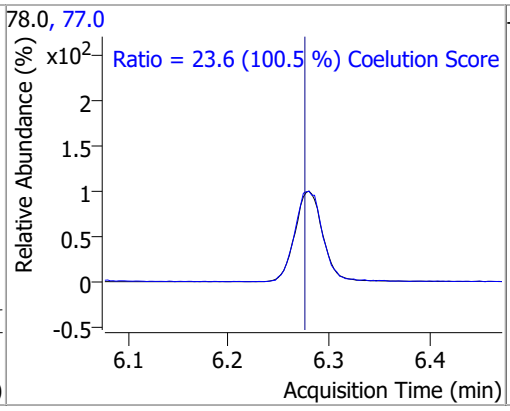


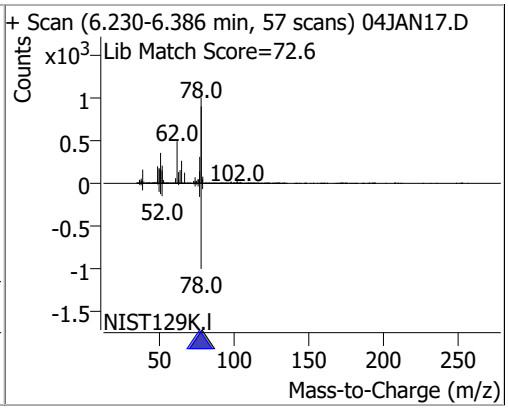




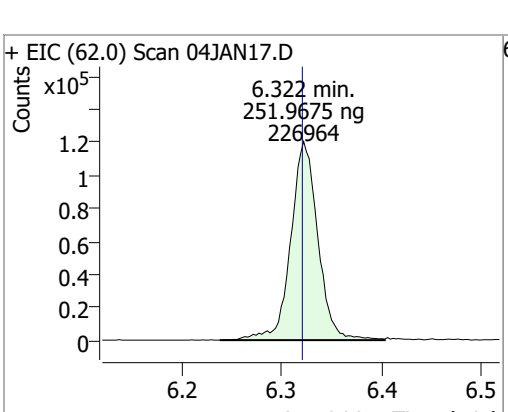
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	257.5416	6.28	0.00	857534	77.0	23.6	0.0	53.5

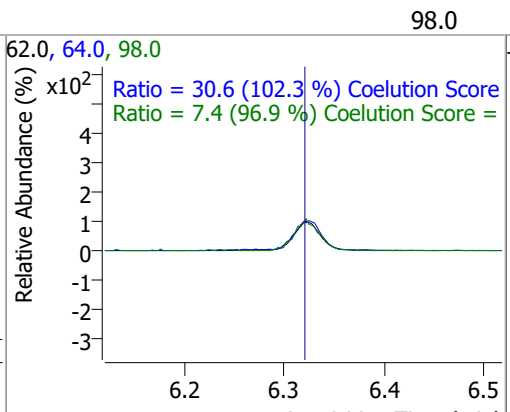


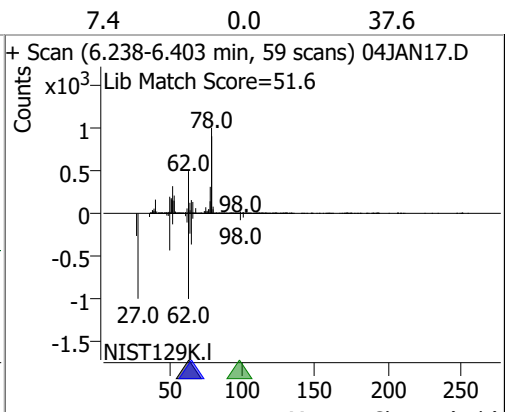




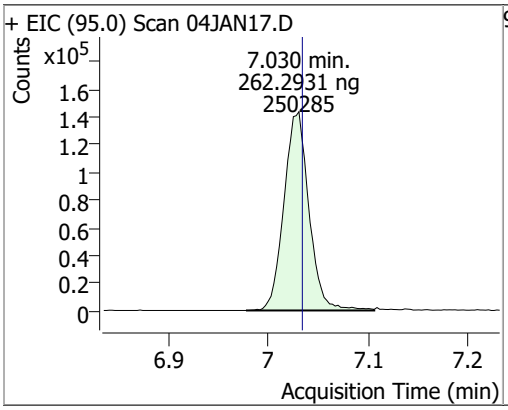
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	251.9675	6.32	0.00	226964	64.0	30.6	0.0	59.9
					98.0	7.4	0.0	37.6

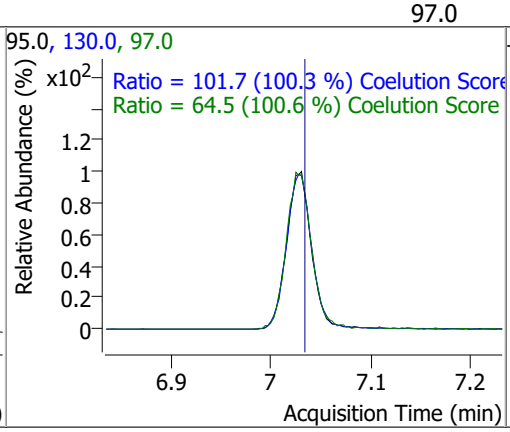


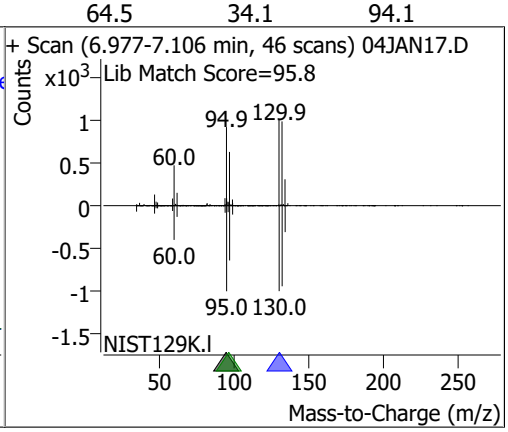




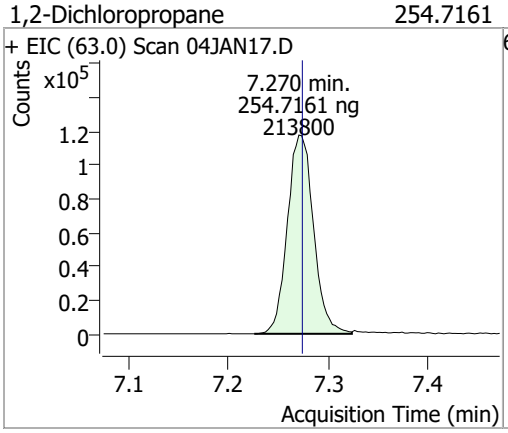
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	262.2931	7.03	0.00	250285	130.0	101.7	71.5	131.5
					97.0	64.5	34.1	94.1

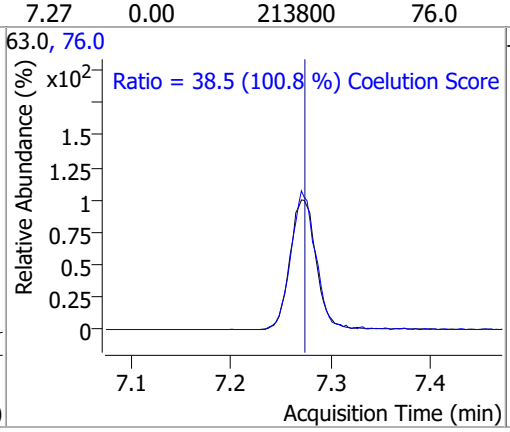


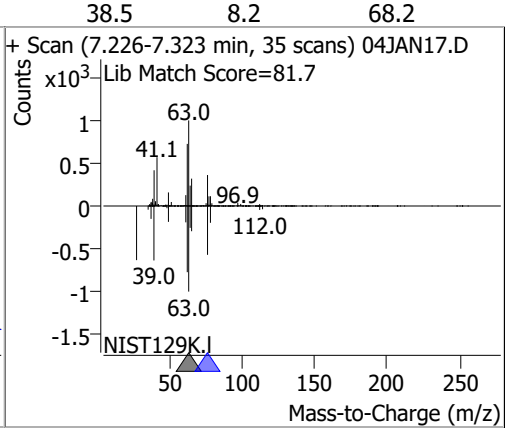




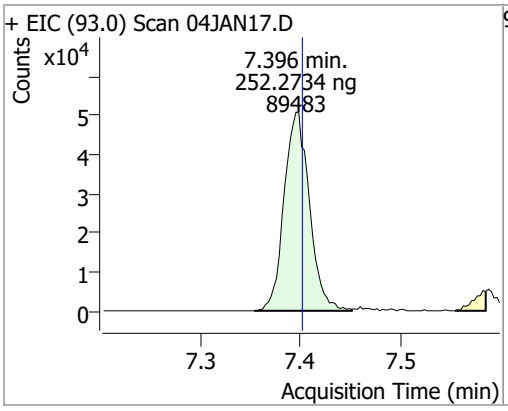
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	254.7161	7.27	0.00	213800	76.0	38.5	8.2	68.2

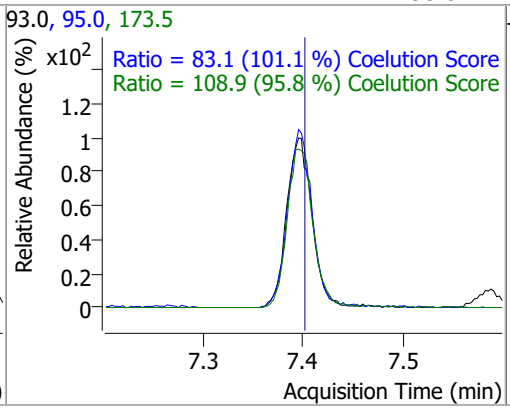


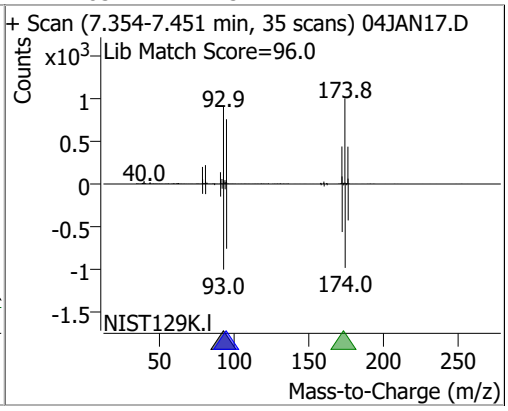




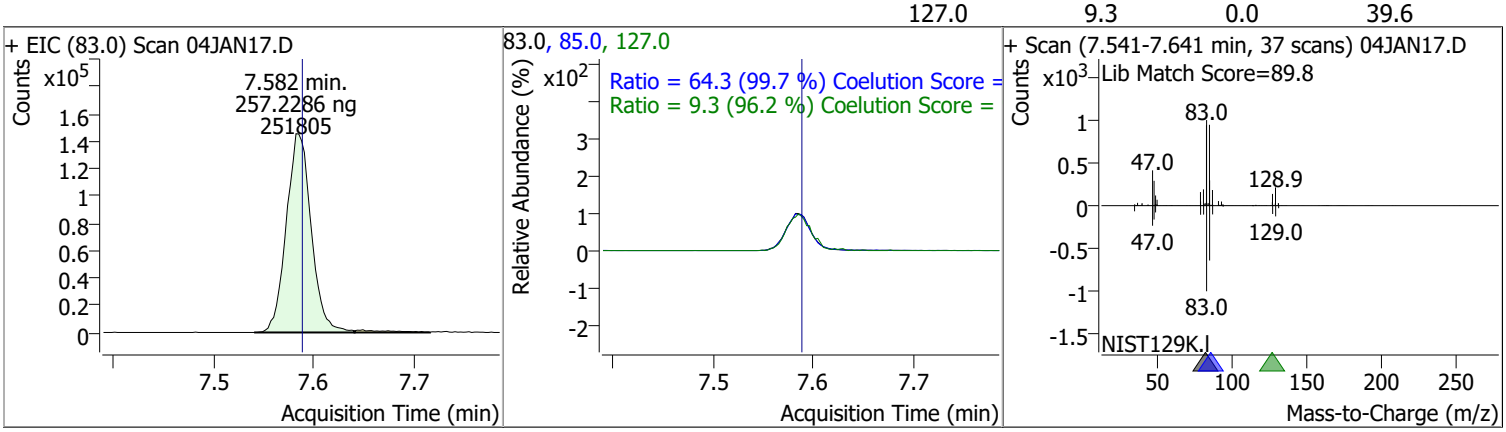
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	252.2734	7.40	0.00	89483	173.5	108.9	83.7	143.7
					95.0	83.1	52.2	112.2



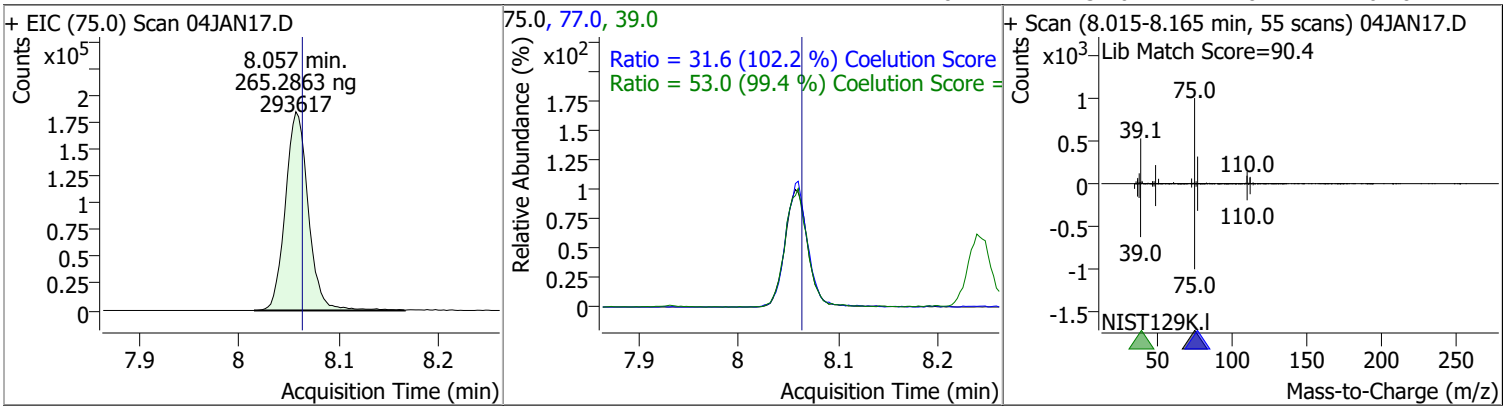




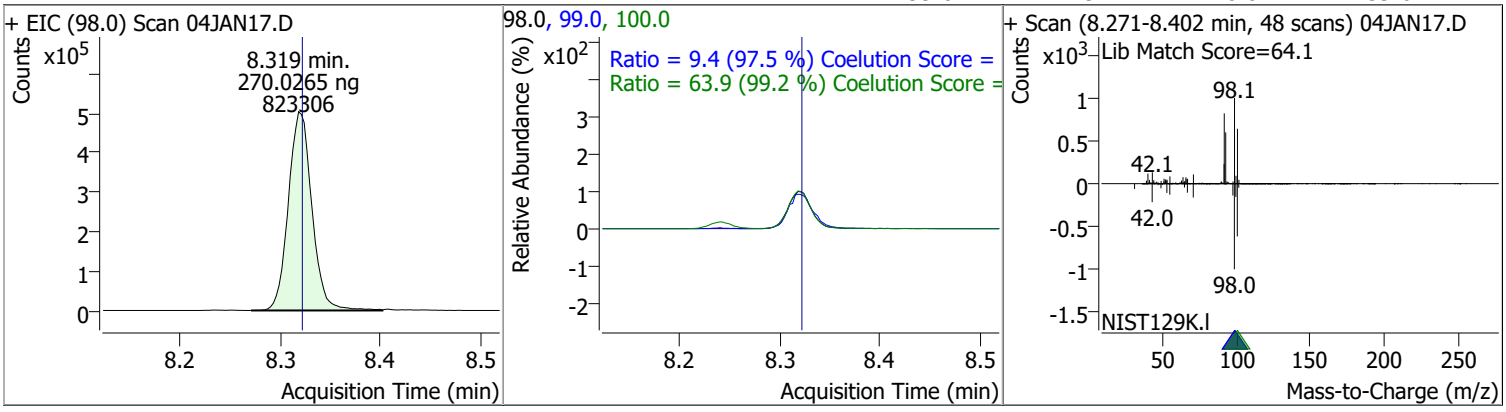
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	257.2286	7.58	0.00	251805	85.0	64.3	34.5	94.5
					127.0	9.3	0.0	39.6



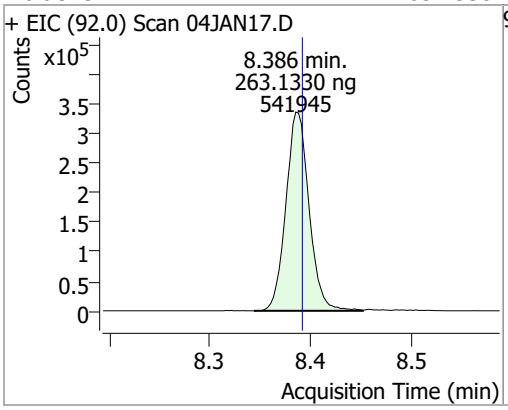
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	265.2863	8.06	0.00	293617	39.0	53.0	23.3	83.3
					77.0	31.6	1.0	61.0

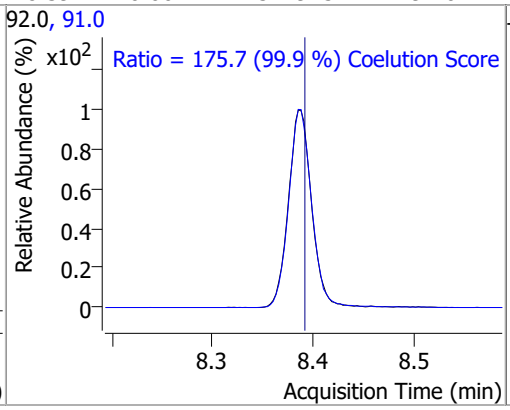


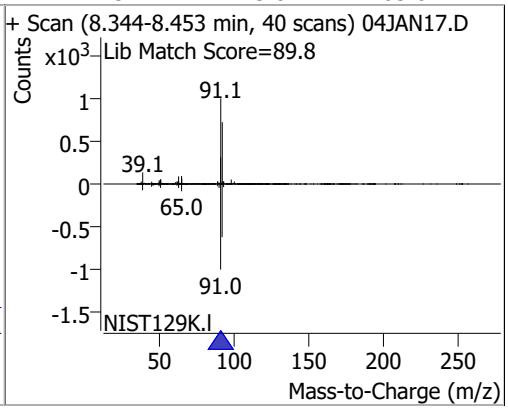
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	270.0265	8.32	0.00	823306	100.0	63.9	34.4	94.4
					99.0	9.4	0.0	39.6



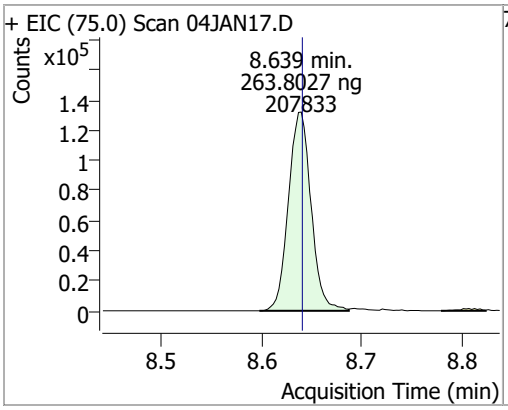
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	263.1330	8.39	0.00	541945	91.0	175.7	145.8	205.8

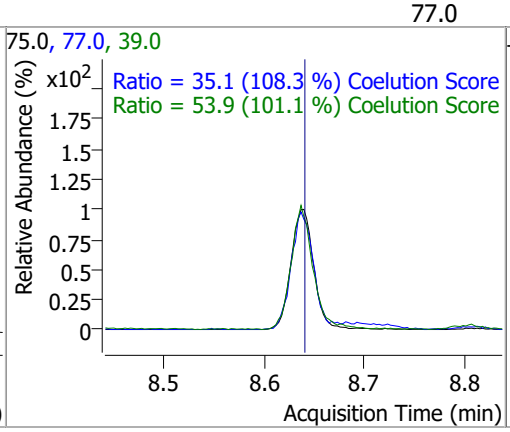


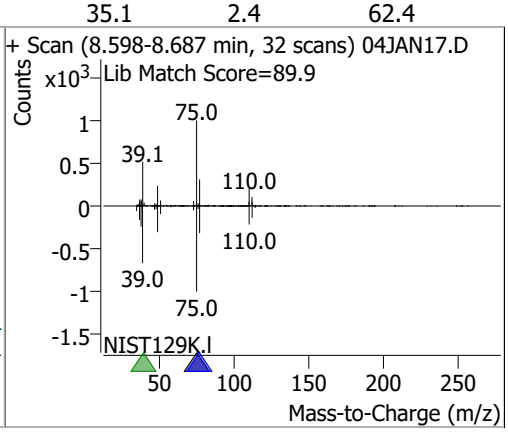




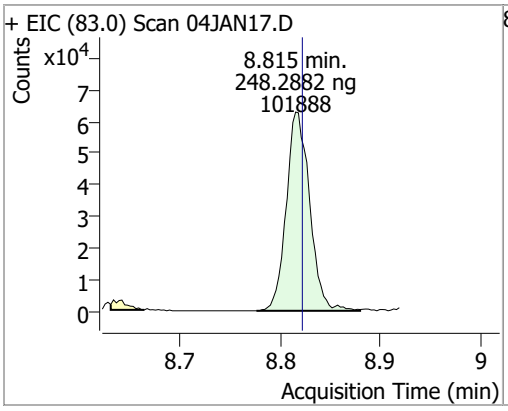
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	263.8027	8.64	0.00	207833	39.0	53.9	23.4	83.4
					77.0	35.1	2.4	62.4

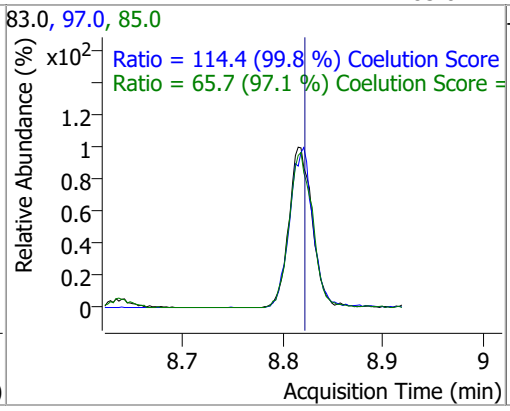


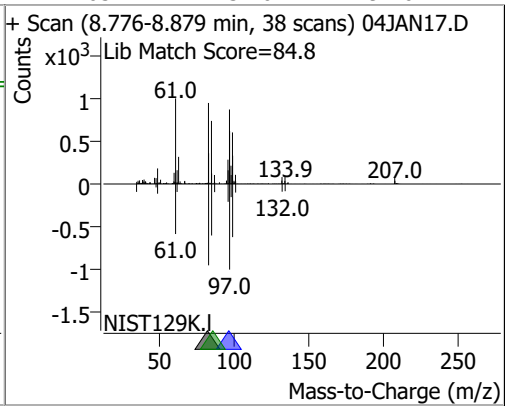




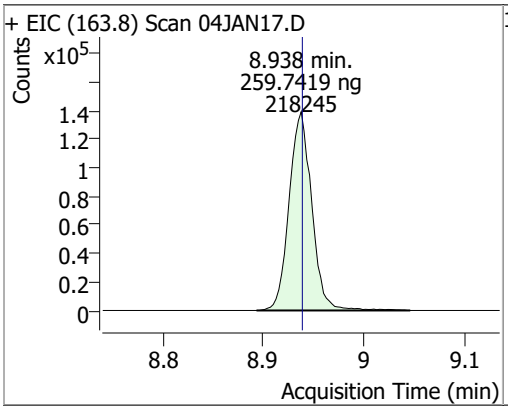
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	248.2882	8.82	0.00	101888	97.0	114.4	84.6	144.6
					85.0	65.7	37.6	97.6

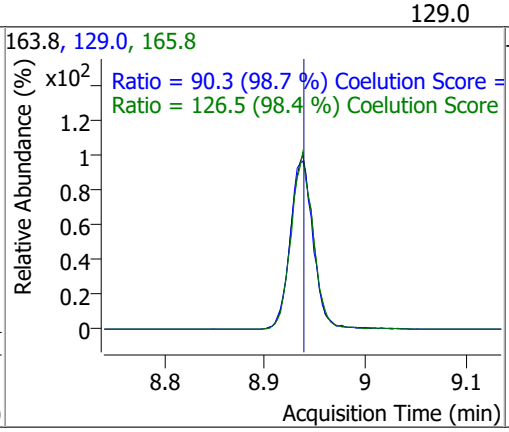


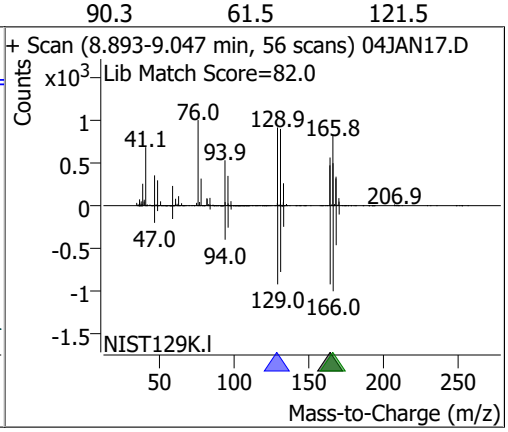




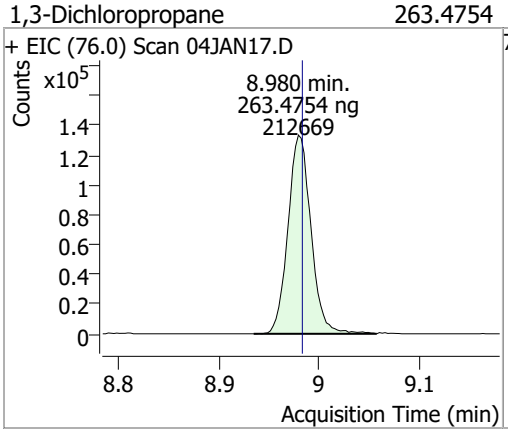
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	259.7419	8.94	0.00	218245	165.8	126.5	98.6	158.6
					129.0	90.3	61.5	121.5

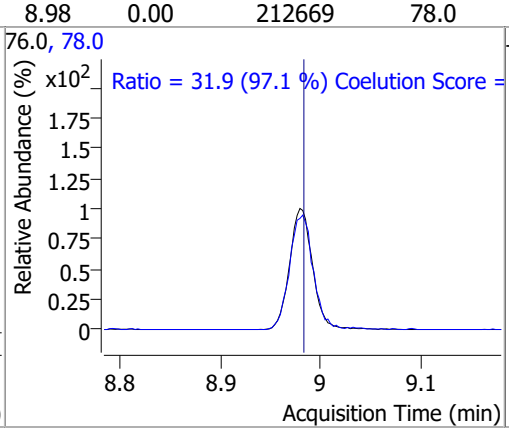


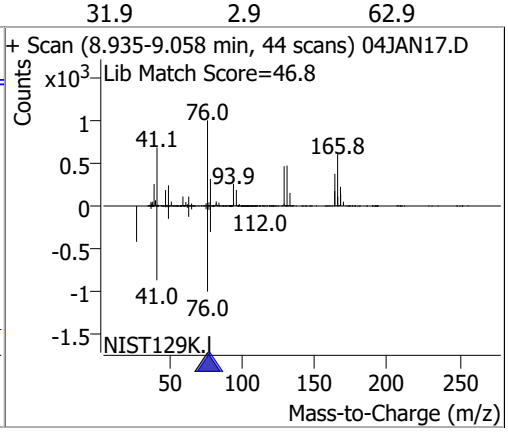




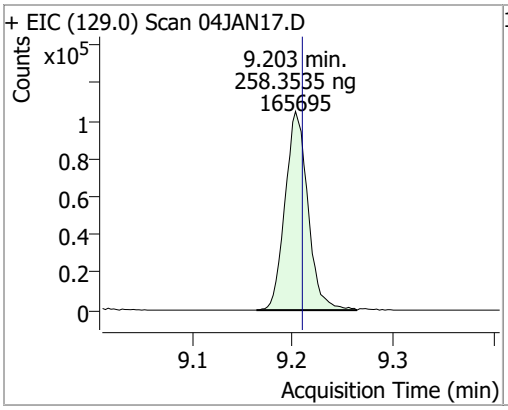
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	263.4754	8.98	0.00	212669	78.0	31.9	2.9	62.9

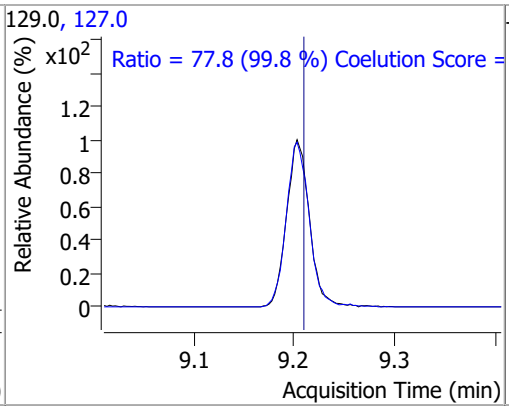


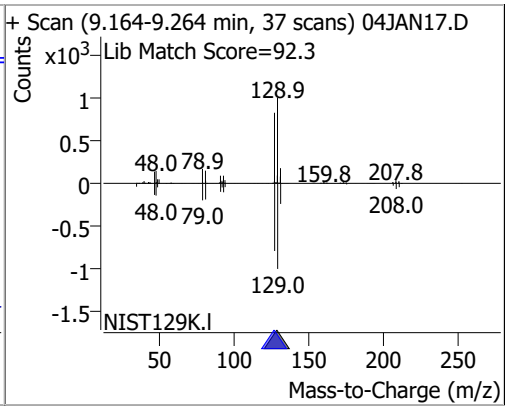




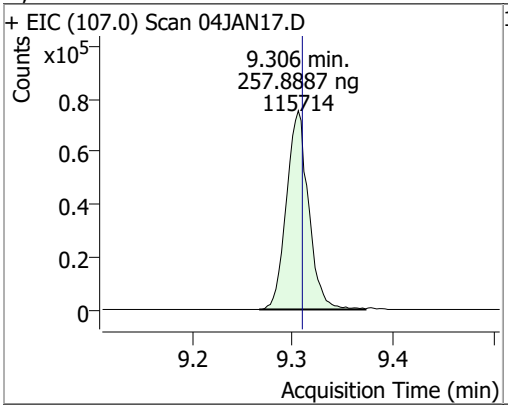
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	258.3535	9.20	0.00	165695	127.0	77.8	48.0	108.0

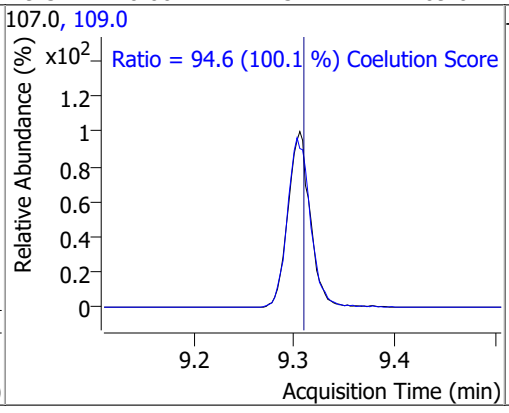


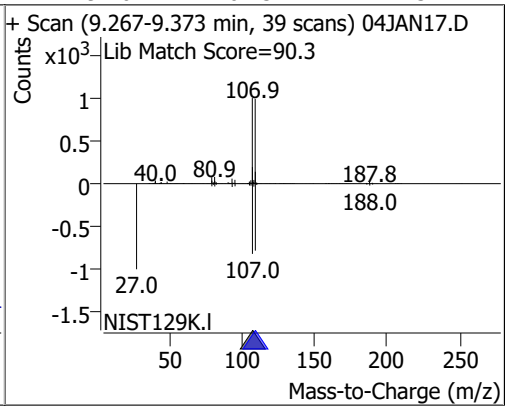




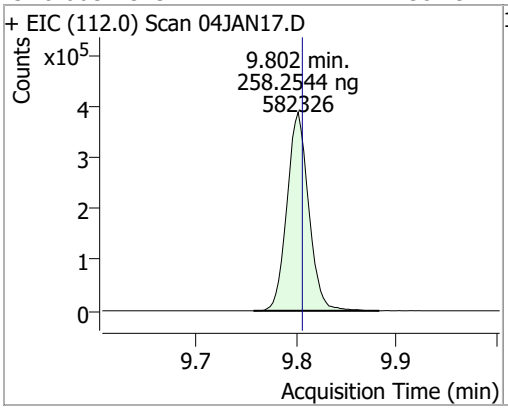
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	257.8887	9.31	0.00	115714	109.0	94.6	64.5	124.5

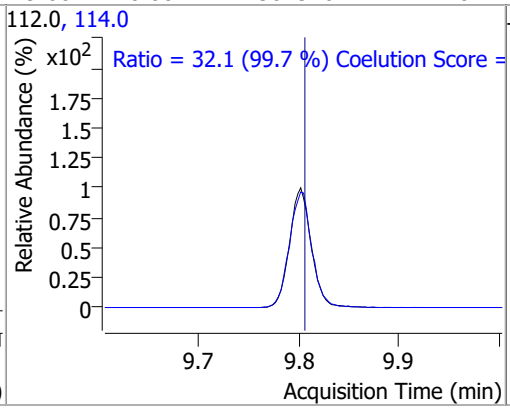


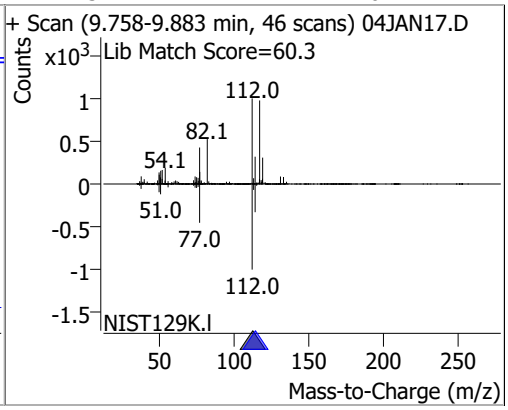




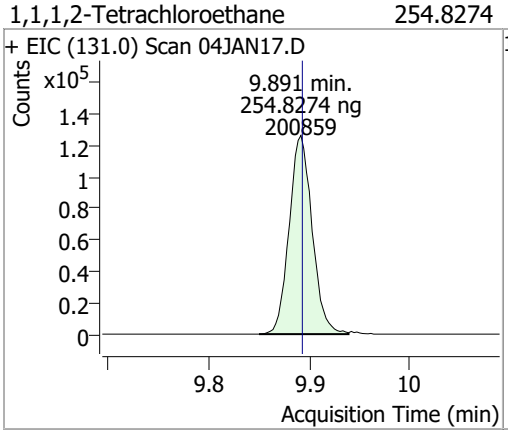
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	258.2544	9.80	0.00	582326	114.0	32.1	2.1	62.1

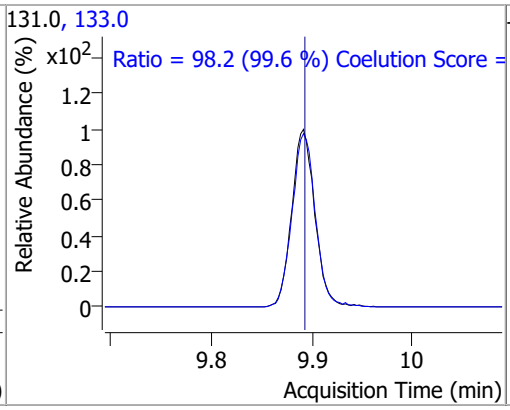


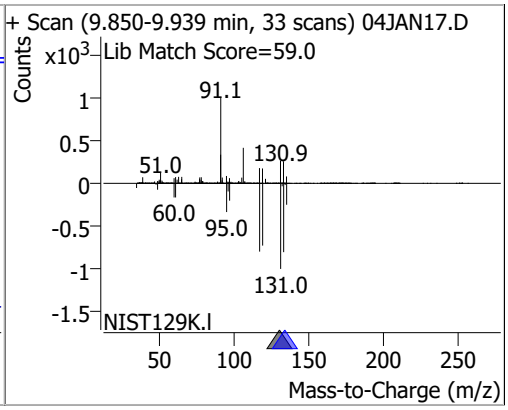




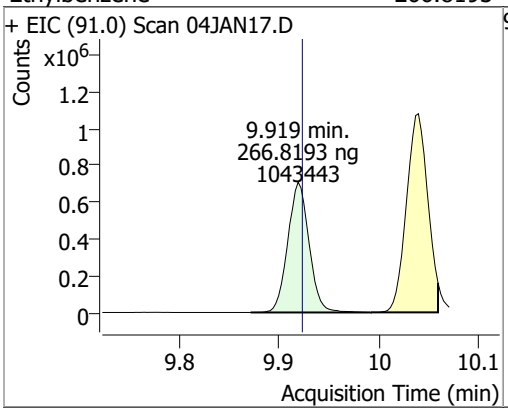
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	254.8274	9.89	0.00	200859	133.0	98.2	68.6	128.6

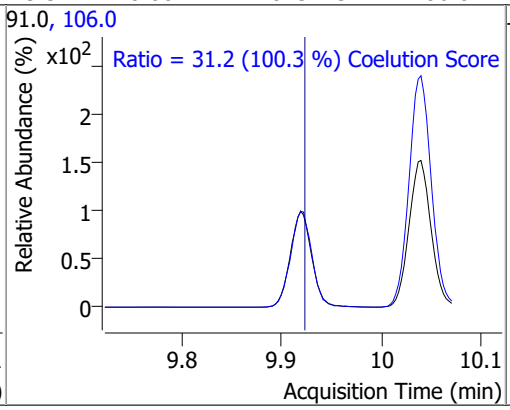


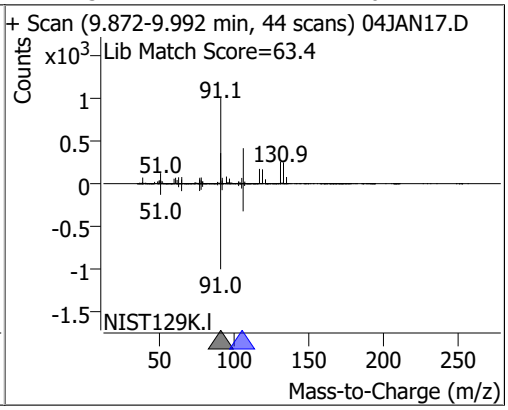




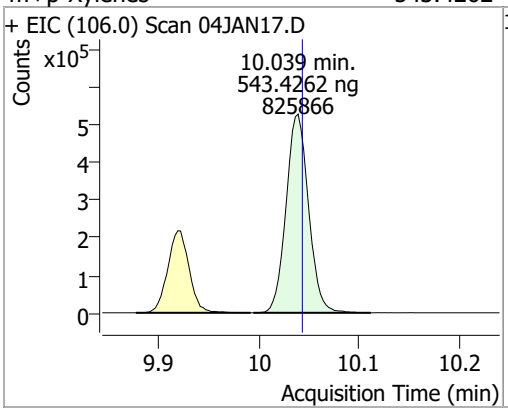
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	266.8193	9.92	0.00	1043443	106.0	31.2	1.1	61.1

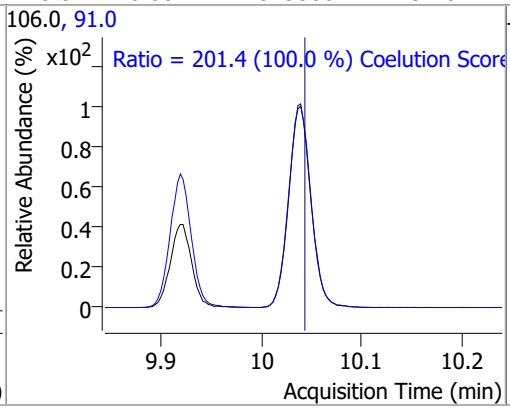


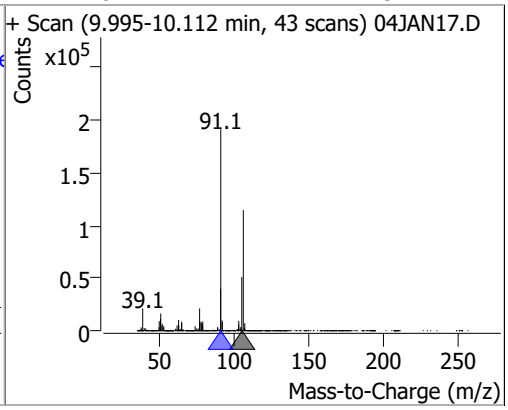




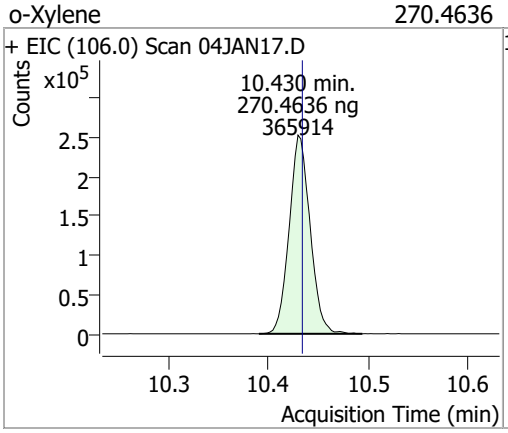
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	543.4262	10.04	0.00	825866	91.0	201.4	171.4	231.4

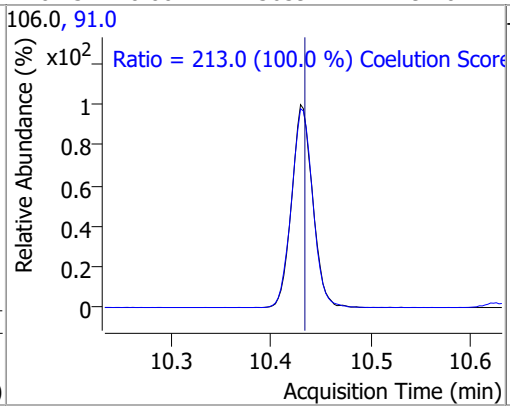


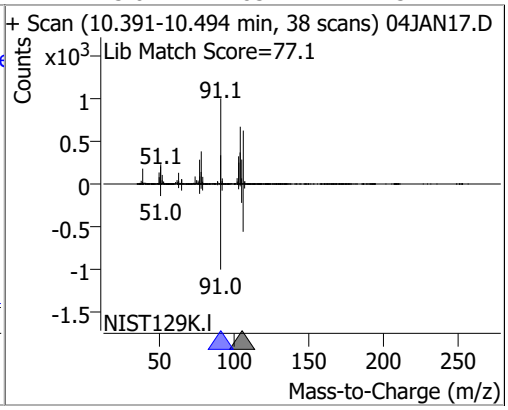




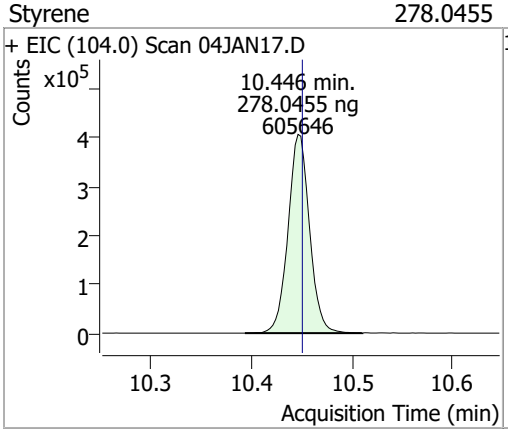
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	270.4636	10.43	0.00	365914	91.0	213.0	183.1	243.1

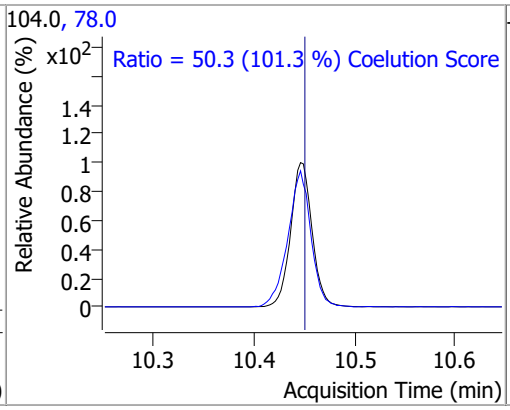


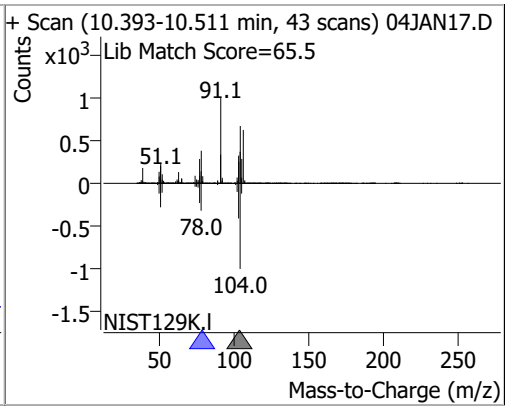




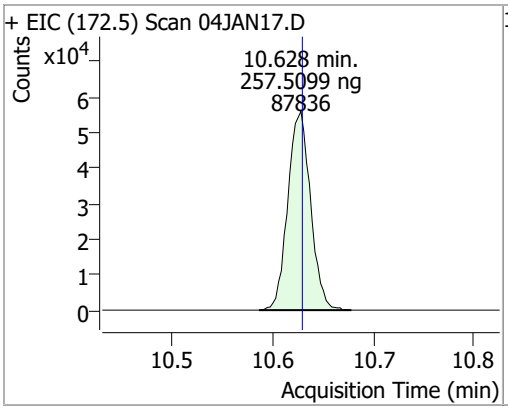
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	278.0455	10.45	0.00	605646	78.0	50.3	19.6	79.6

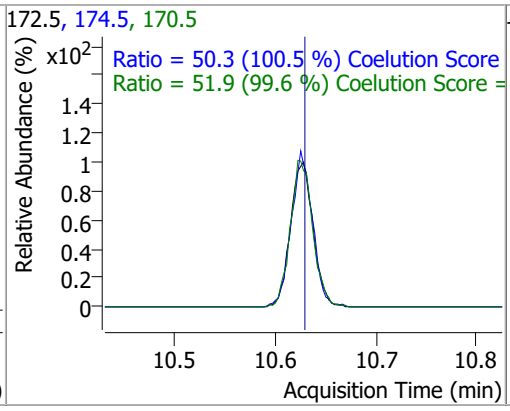


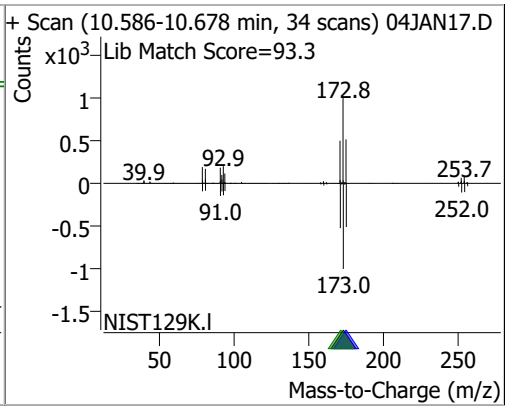




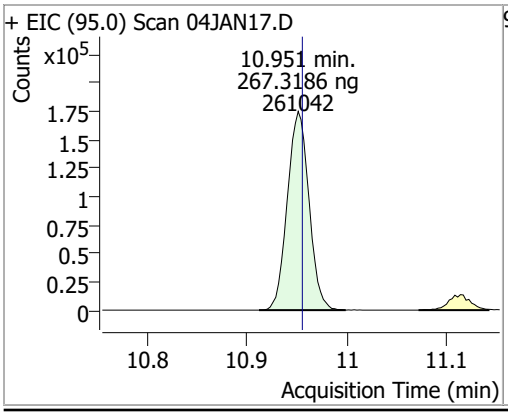
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	257.5099	10.63	0.00	87836	170.5	51.9	22.1	82.1
					174.5	50.3	20.1	80.1

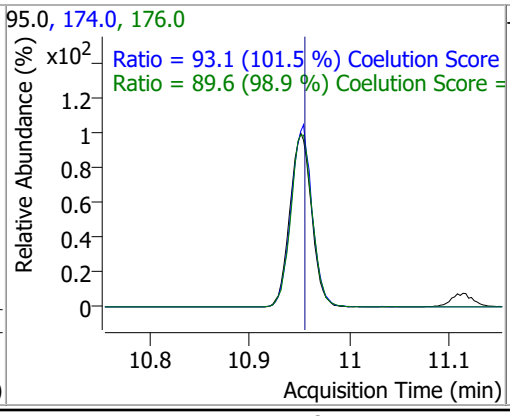


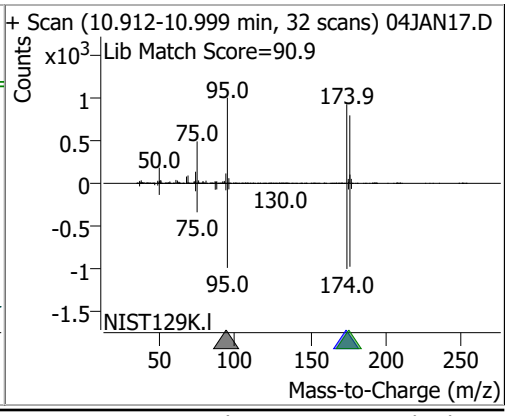




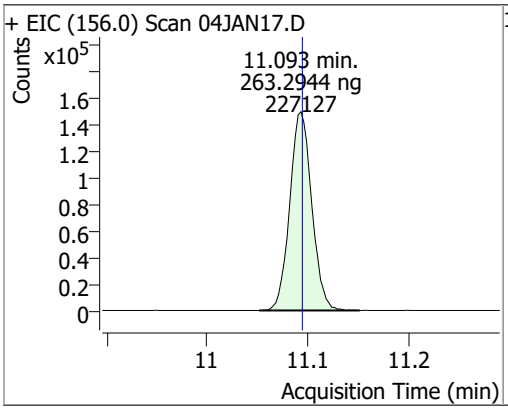
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	267.3186	10.95	0.00	261042	174.0	93.1	61.7	121.7
					176.0	89.6	60.6	120.6

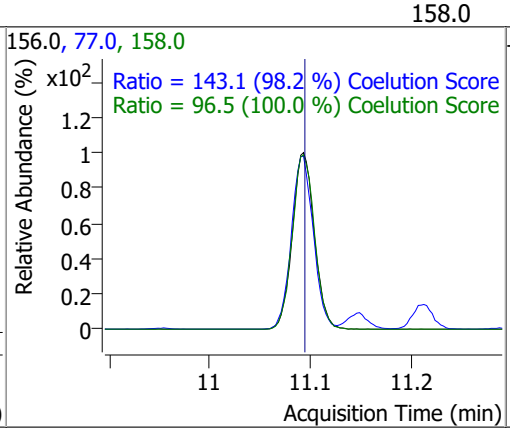


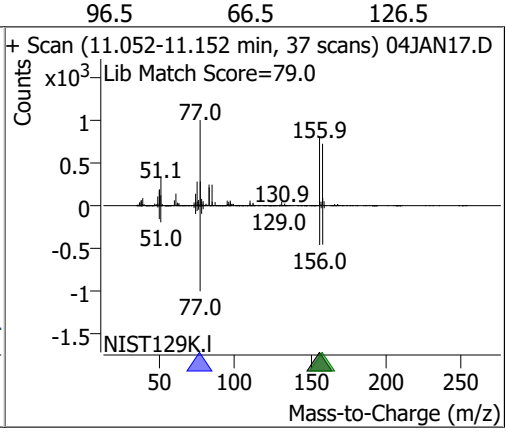




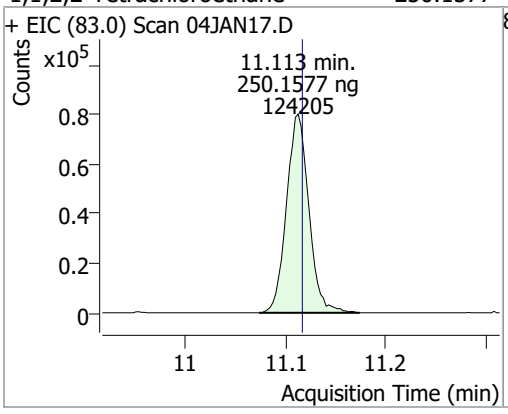
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	263.2944	11.09	0.00	227127	77.0	143.1	115.7	175.7
					158.0	96.5	66.5	126.5

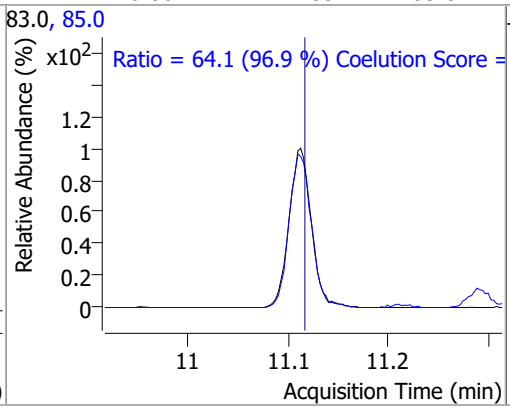


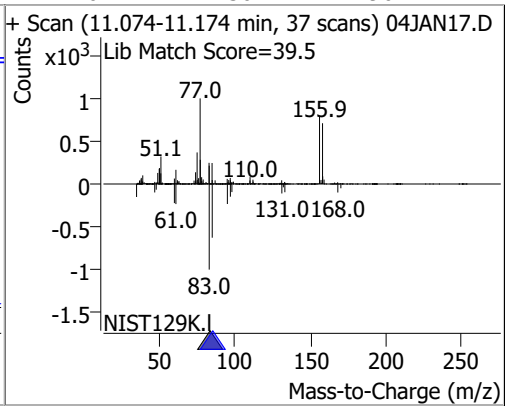




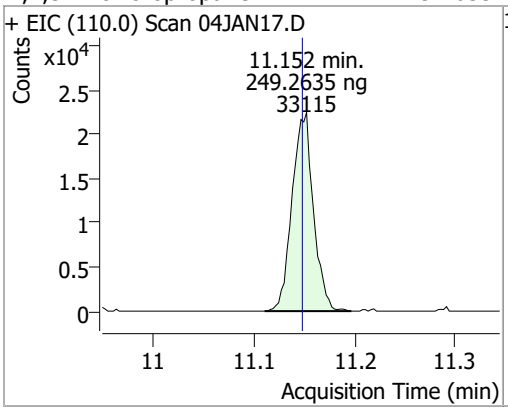
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	250.1577	11.11	0.00	124205	85.0	64.1	36.2	96.2

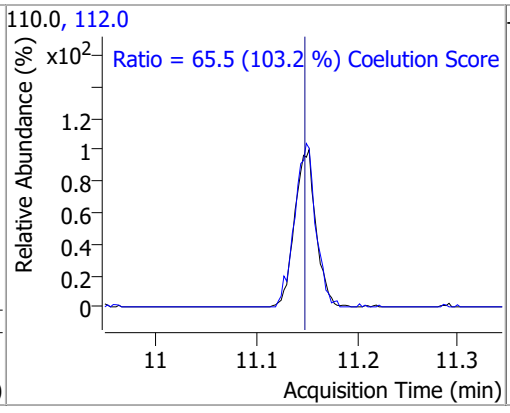


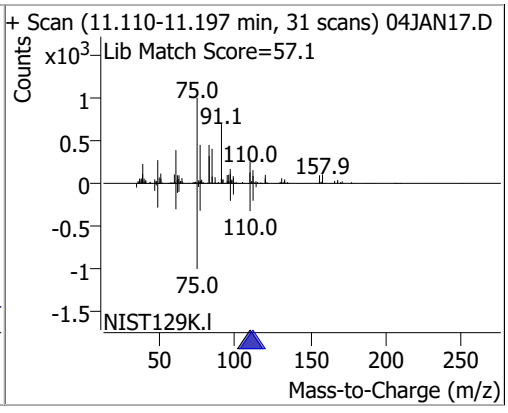




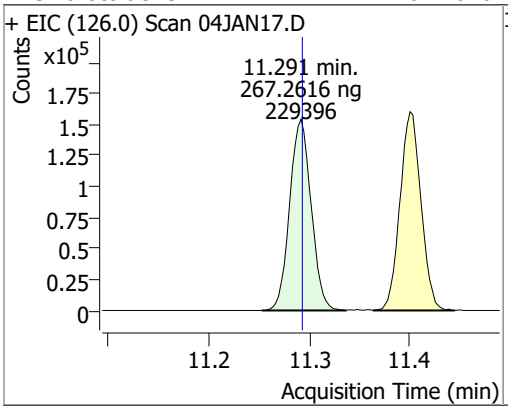
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	249.2635	11.15	0.01	33115	112.0	65.5	33.5	93.5

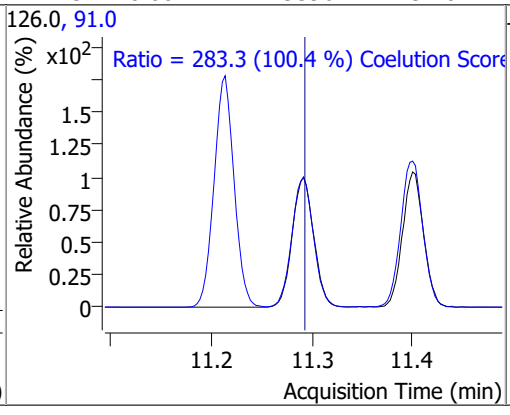


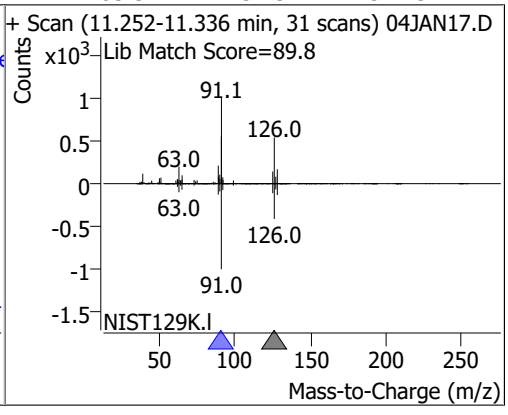




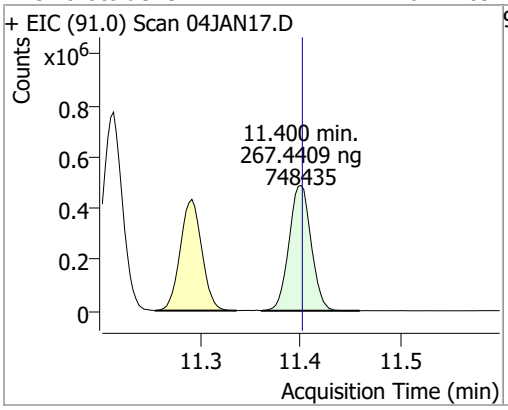
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	267.2616	11.29	0.00	229396	91.0	283.3	252.3	312.3

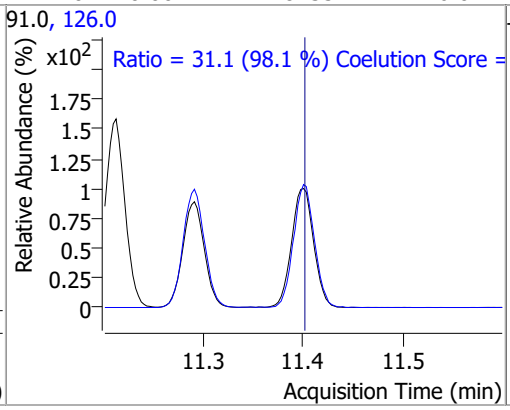


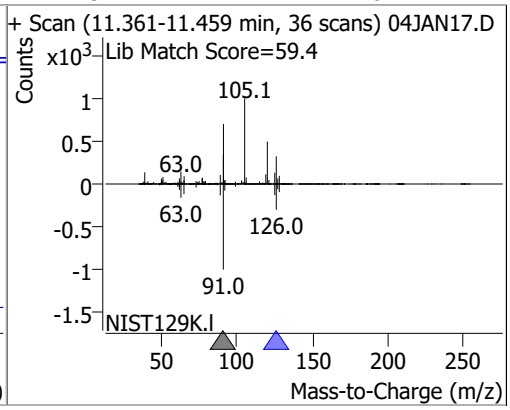




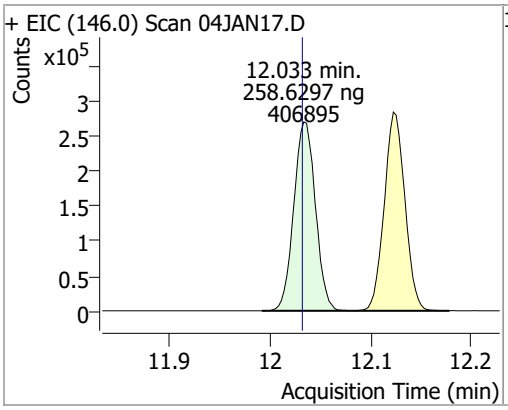
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	267.4409	11.40	0.00	748435	126.0	31.1	1.7	61.7

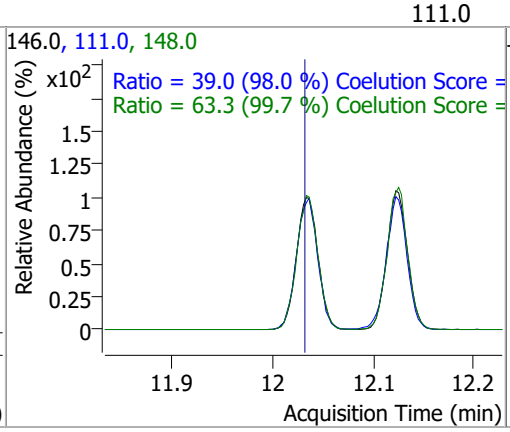


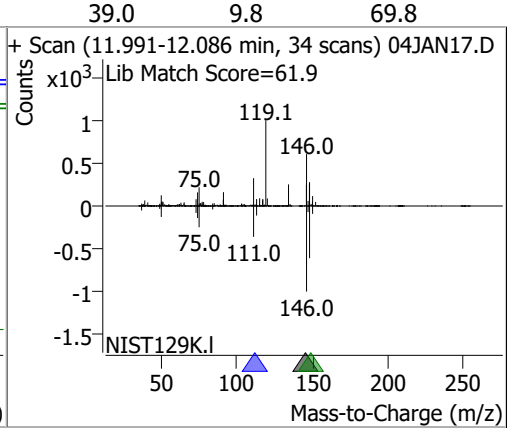




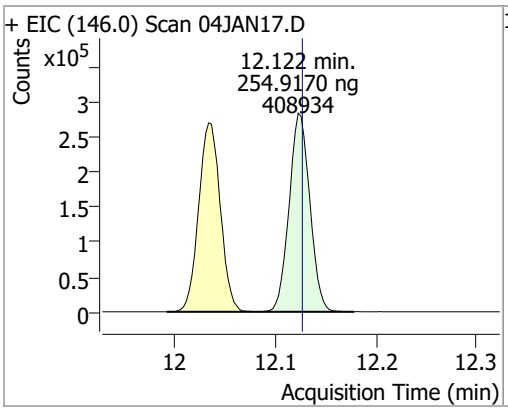
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	258.6297	12.03	0.00	406895	148.0	63.3	33.6	93.6
					111.0	39.0	9.8	69.8

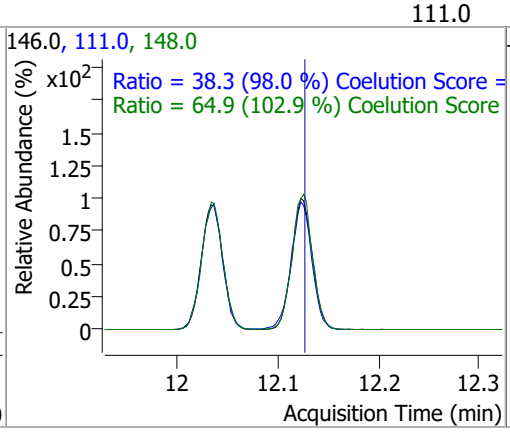


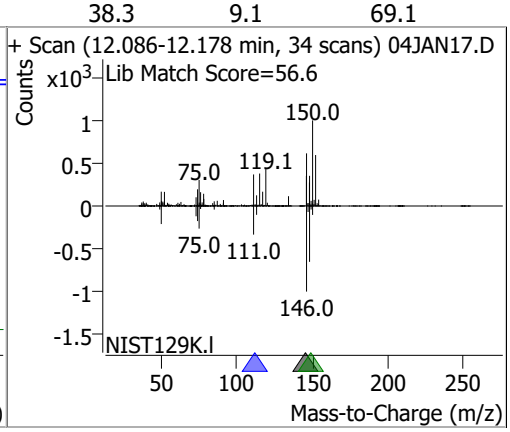




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	254.9170	12.12	0.00	408934	148.0	64.9	33.1	93.1
					111.0	38.3	9.1	69.1

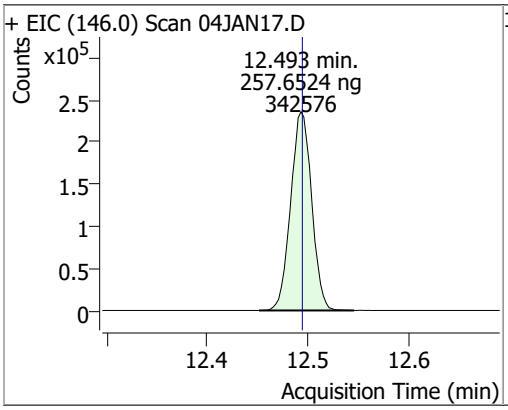


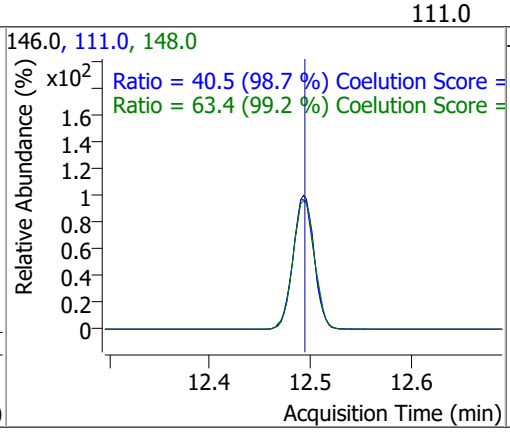


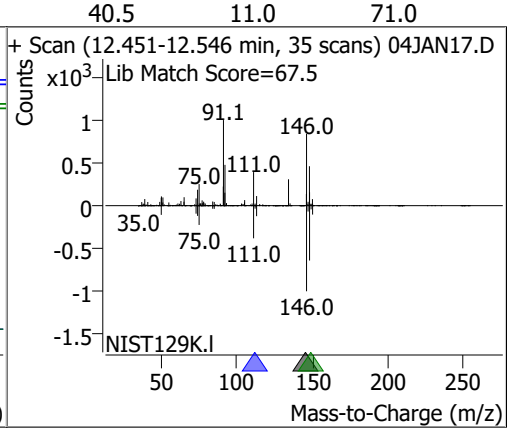


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	257.6524	12.49	0.00	342576	148.0	63.4	33.9	93.9
					111.0	40.5	11.0	71.0







Quantitation Results Report (QT Reviewed)

Data File	04JAN19.D	Operator	MSC
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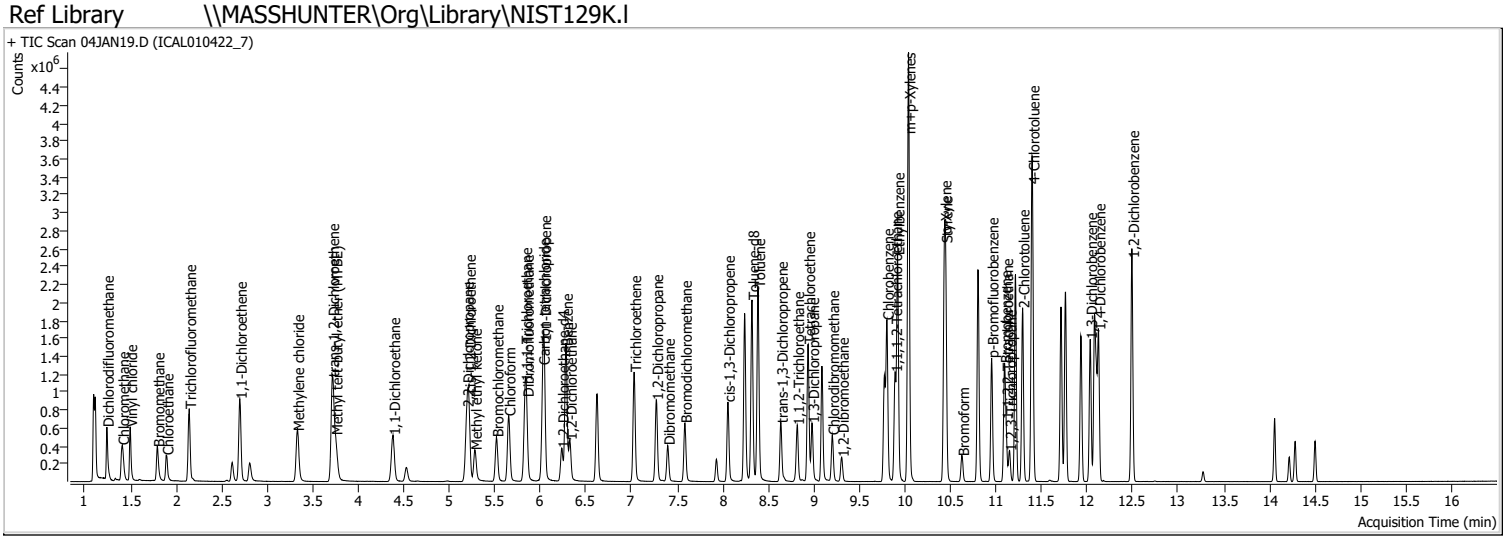
Vial	19	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

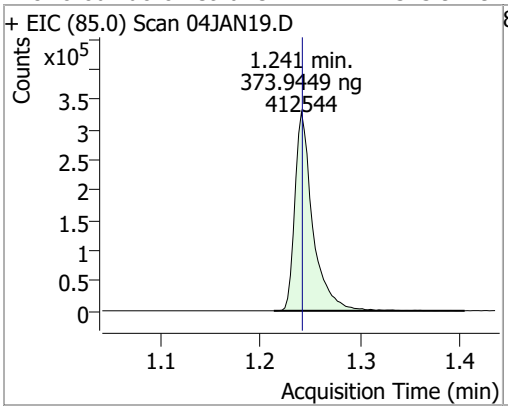


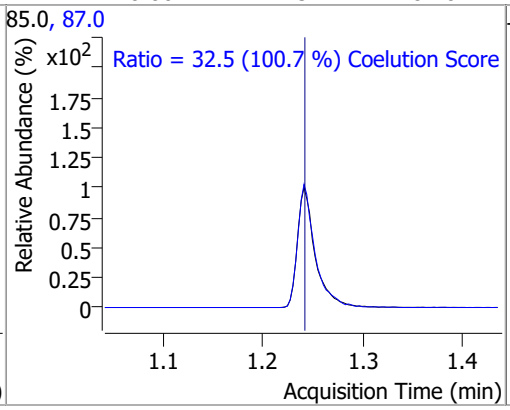
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	841876	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	314668	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	266611	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.848	113.0	305158	384.7503	ng		0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 153.90%		*	
S 1,2-Dichloroethane-d4	6.233	67.0	129608	378.3335	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 151.33%		*	
S Toluene-d8	8.322	98.0	1229775	405.5583	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 162.22%		*	
S p-Bromofluorobenzene	10.951	95.0	385474	394.6566	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 157.86%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.241	85.0	412544	373.9449	ng		100
T Chloromethane	1.409	50.0	471454	352.0836	ng		99
T Vinyl chloride	1.498	62.0	448643	372.3564	ng		95
T Bromomethane	1.796	96.0	207491	385.1259	ng		98
T Chloroethane	1.897	64.0	217393	364.4573	ng		99
T Trichlorofluoromethane	2.145	101.0	555477	371.4290	ng		98
T 1,1-Dichloroethene	2.700	96.0	322557	380.3725	ng		100
T Methylene chloride	3.330	49.0	435116	348.0666	ng		99
T trans-1,2-Dichloroethene	3.715	96.0	325415	376.1367	ng		97
T Methyl tert-butyl ether (MTBE)	3.751	73.0	437439	391.1767	ng		100
T 1,1-Dichloroethane	4.381	63.0	612660	380.4437	ng		99
T 2,2-Dichloropropane	5.190	77.0	446282	369.8436	ng		100
T cis-1,2-Dichloroethene	5.212	96.0	339211	386.7236	ng		97
T Methyl ethyl ketone	5.279	43.0	470653	3961.3410	ng		100
T Bromochloromethane	5.516	128.0	135103	371.8004	ng		99
T Chloroform	5.650	83.0	588080	366.9389	ng		99

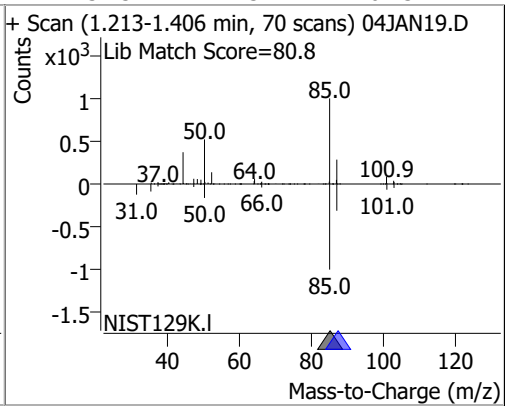
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.831	97.0	580748	386.6625	ng	99
T Carbon tetrachloride	6.024	117.0	572545	386.9014	ng	99
T 1,1-Dichloropropene	6.040	75.0	507157	397.1322	ng	100
T Benzene	6.278	78.0	1293370	385.8526	ng	99
T 1,2-Dichloroethane	6.322	62.0	332775	366.9787	ng	97
T Trichloroethene	7.028	95.0	374370	394.4896	ng	99
T 1,2-Dichloropropane	7.270	63.0	324602	388.8502	ng	98
T Dibromomethane	7.396	93.0	134282	380.6547	ng	96
T Bromodichloromethane	7.585	83.0	375983	386.1940	ng	100
T cis-1,3-Dichloropropene	8.057	75.0	441168	400.7930	ng	99
T Toluene	8.388	92.0	813204	397.0106	ng	100
T trans-1,3-Dichloropropene	8.639	75.0	315063	402.1098	ng	99
T 1,1,2-Trichloroethane	8.818	83.0	152331	373.2534	ng	100
T Tetrachloroethene	8.938	163.8	319950	382.8796	ng	99
T 1,3-Dichloropropane	8.980	76.0	312547	389.3442	ng	99
T Chlorodibromomethane	9.203	129.0	247279	387.6812	ng	99
T 1,2-Dibromoethane	9.306	107.0	168577	377.7698	ng	100
T Chlorobenzene	9.802	112.0	867732	386.9455	ng	100
T 1,1,1,2-Tetrachloroethane	9.892	131.0	307436	392.1859	ng	96
T Ethylbenzene	9.919	91.0	1574219	404.7587	ng	100
T m+p-Xylenes	10.039	106.0	1228570	812.8556	ng	100
T o-Xylene	10.433	106.0	549244	408.2043	ng	100
T Styrene	10.447	104.0	896331	413.7595	ng	99
T Bromoform	10.628	172.5	129038	378.2200	ng	99
T Bromobenzene	11.093	156.0	333431	386.4420	ng	98
T 1,1,2,2-Tetrachloroethane	11.110	83.0	182470	367.4276	ng	98
T 1,2,3-Trichloropropane	11.149	110.0	48325	363.6732	ng	100
T 2-Chlorotoluene	11.291	126.0	336386	391.8269	ng	98
T 4-Chlorotoluene	11.400	91.0	1109221	396.2756	ng	100
T 1,3-Dichlorobenzene	12.033	146.0	603674	383.6225	ng	100
T 1,4-Dichlorobenzene	12.125	146.0	595919	371.3969	ng	98
T 1,2-Dichlorobenzene	12.493	146.0	499147	375.3283	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

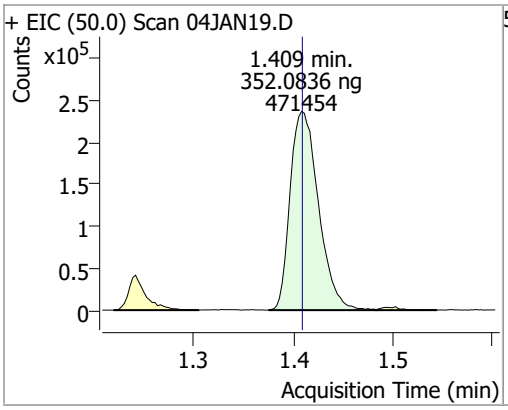
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	373.9449	1.24	0.00	412544	87.0	32.5	2.3	62.3

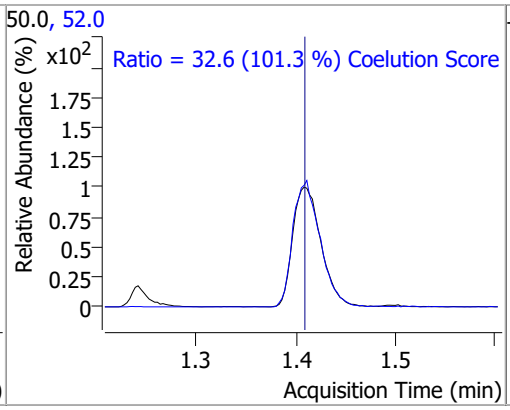


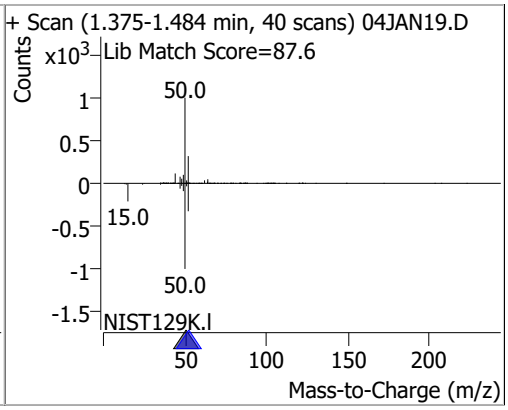




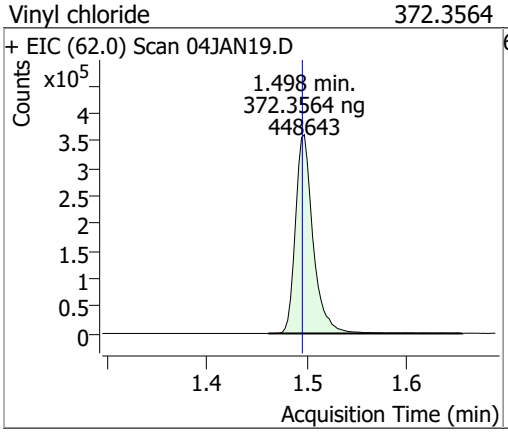
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	352.0836	1.41	0.00	471454	52.0	32.6	2.1	62.1

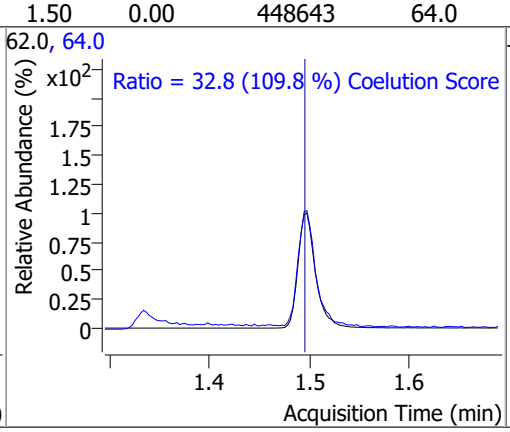


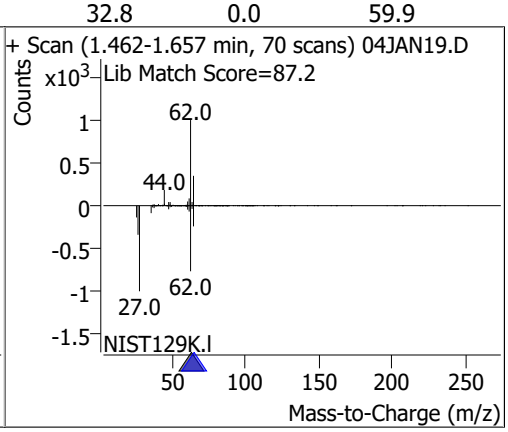




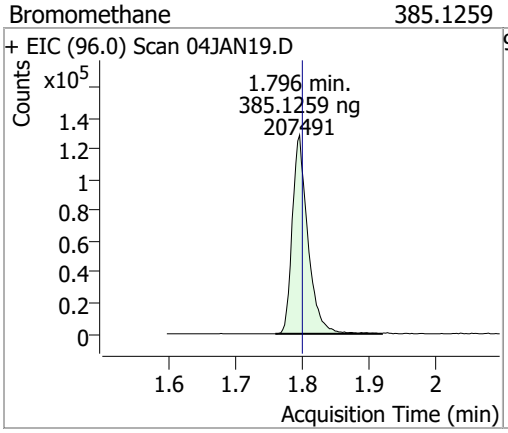
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	372.3564	1.50	0.00	448643	64.0	32.8	0.0	59.9

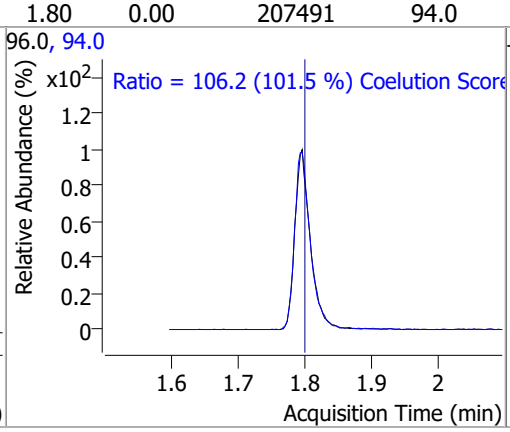


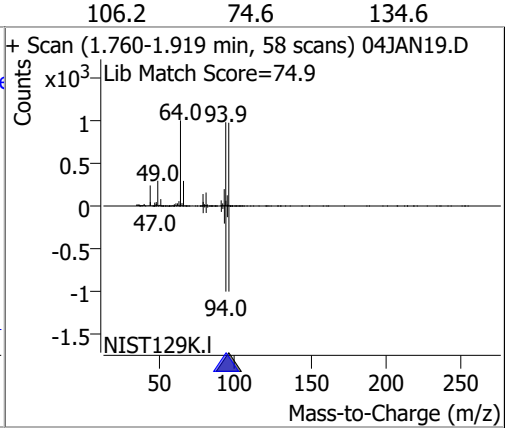




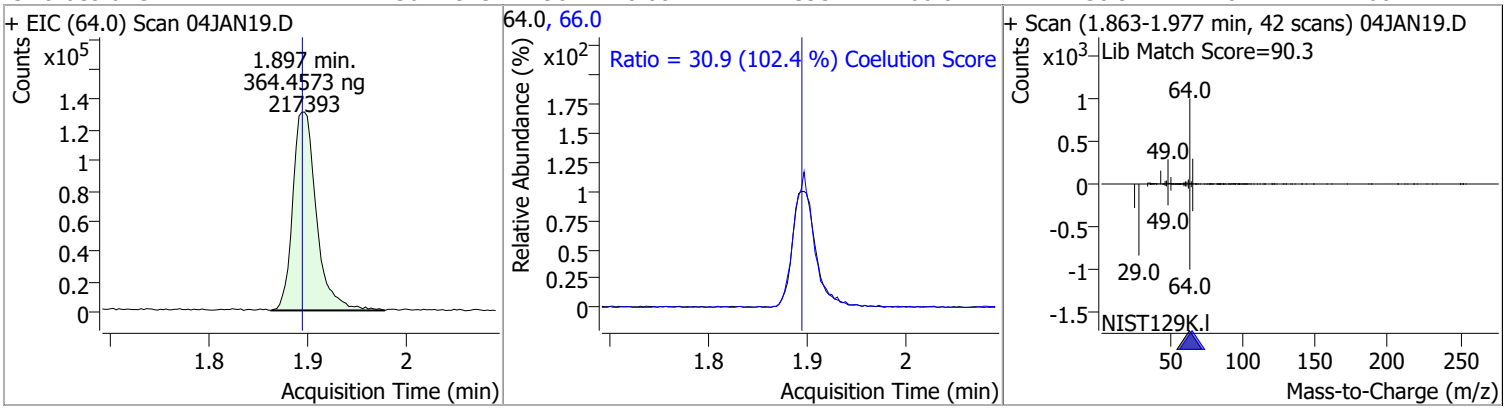
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	385.1259	1.80	0.00	207491	94.0	106.2	74.6	134.6



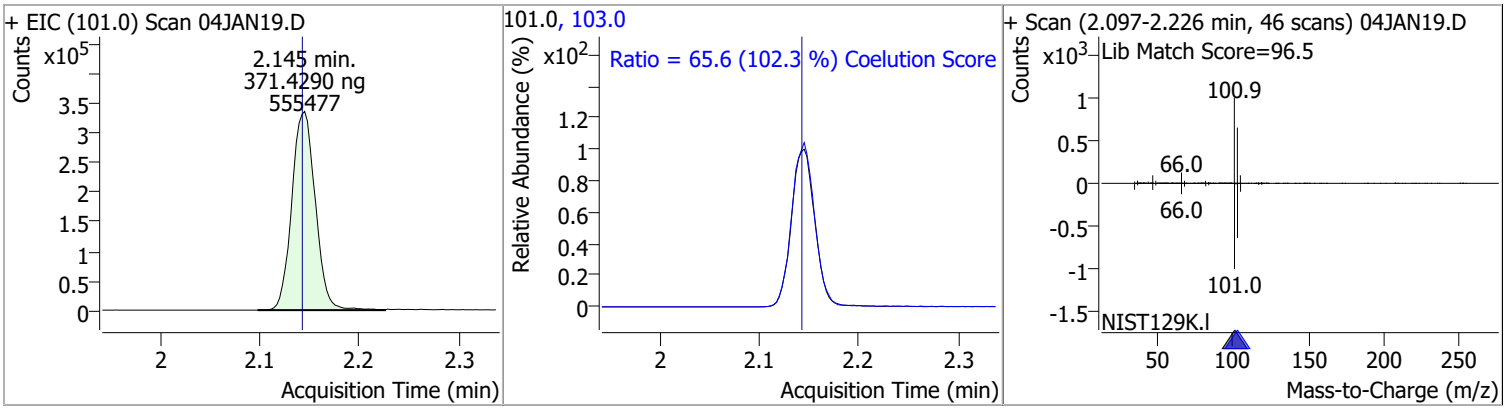




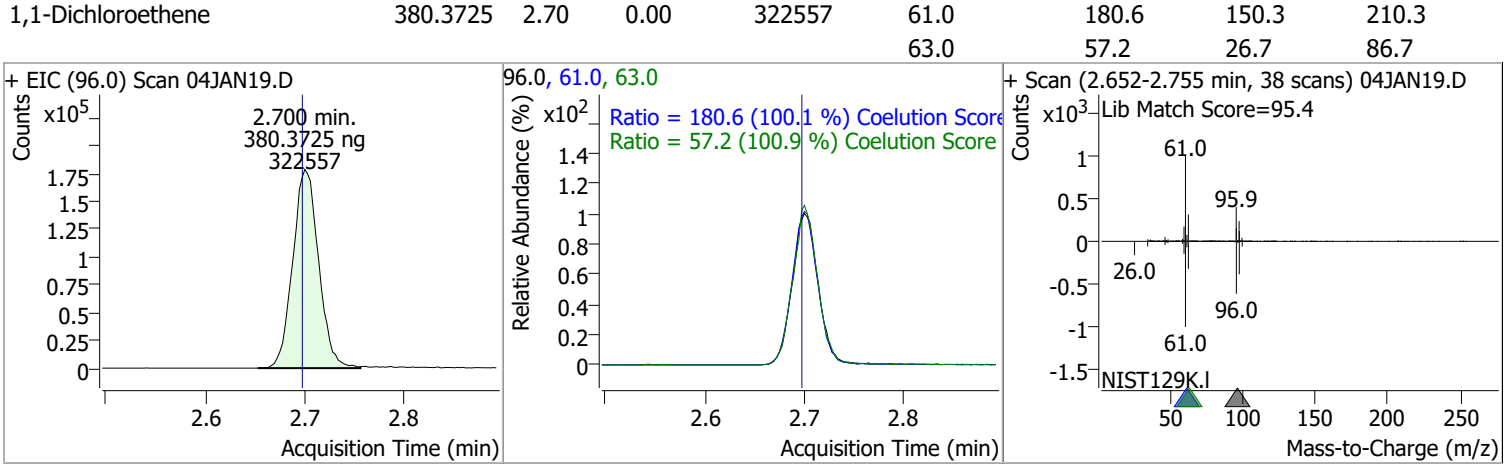
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	364.4573	1.90	0.00	217393	66.0	30.9	0.1	60.1



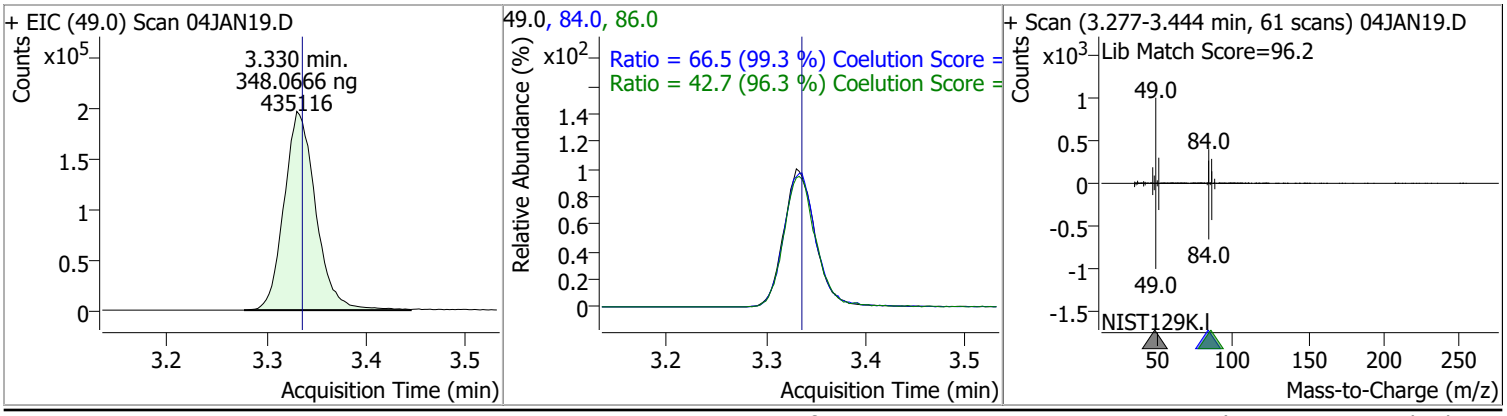
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	371.4290	2.14	0.00	555477	103.0	65.6	34.2	94.2



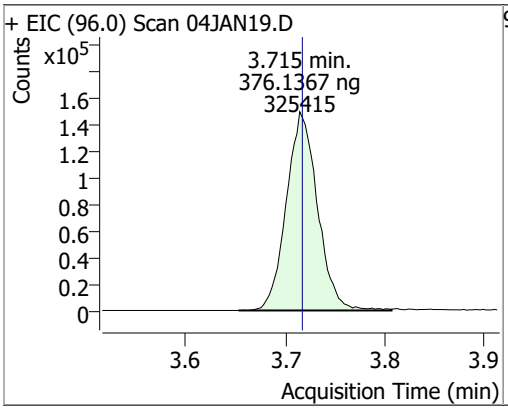
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	380.3725	2.70	0.00	322557	61.0	180.6	150.3	210.3
					63.0	57.2	26.7	86.7

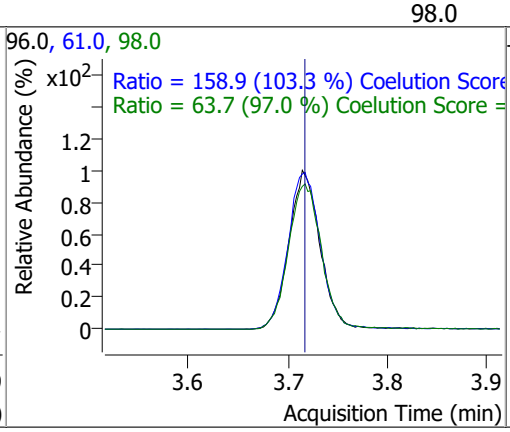


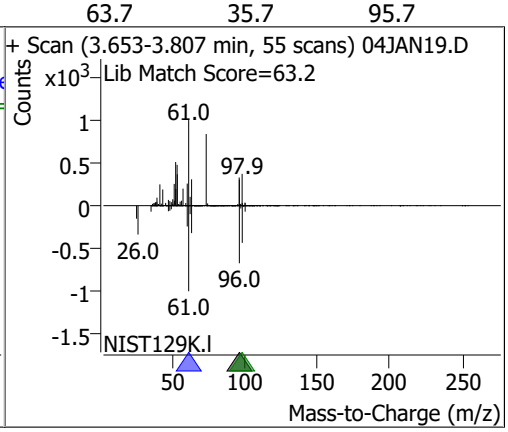
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	348.0666	3.33	-0.01	435116	84.0	66.5	36.9	96.9
					86.0	42.7	14.3	74.3



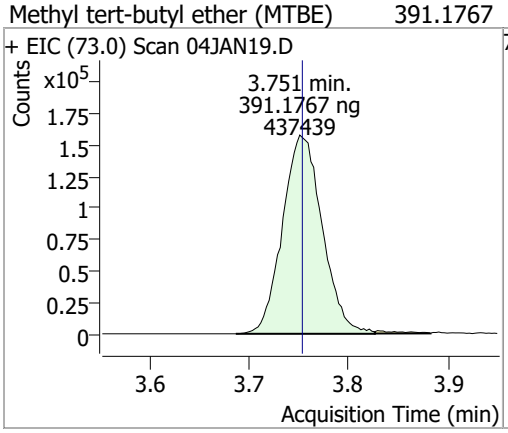
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	376.1367	3.71	0.00	325415	61.0	158.9	123.9	183.9
					98.0	63.7	35.7	95.7

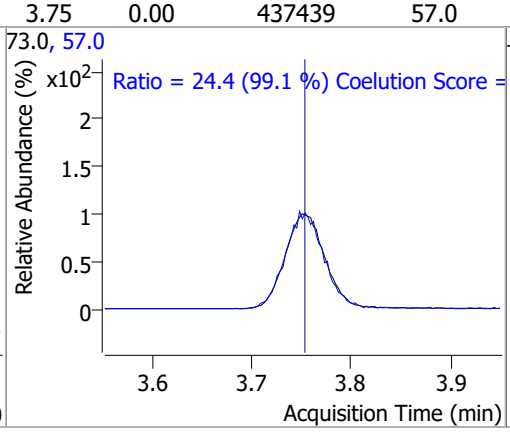


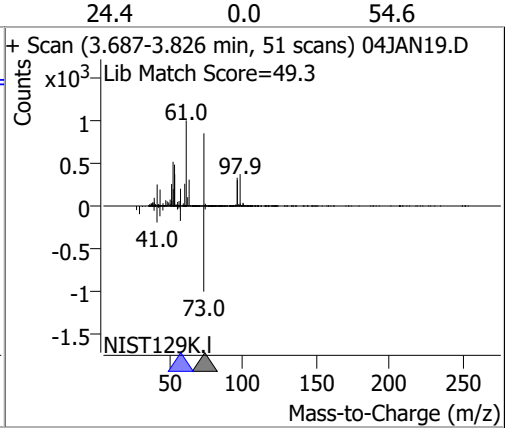




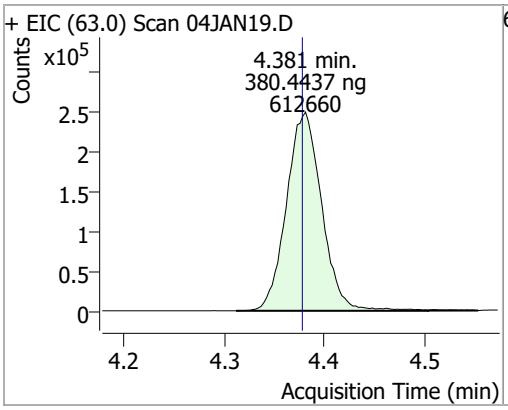
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	391.1767	3.75	0.00	437439	57.0	24.4	0.0	54.6

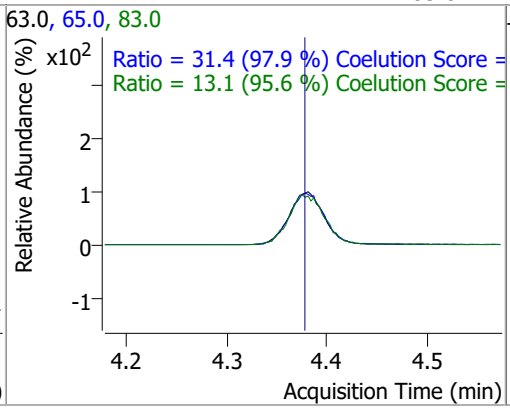


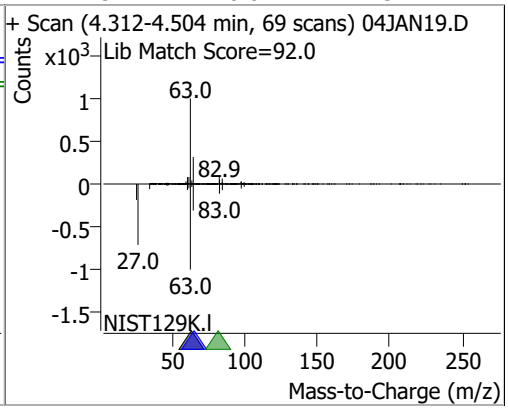




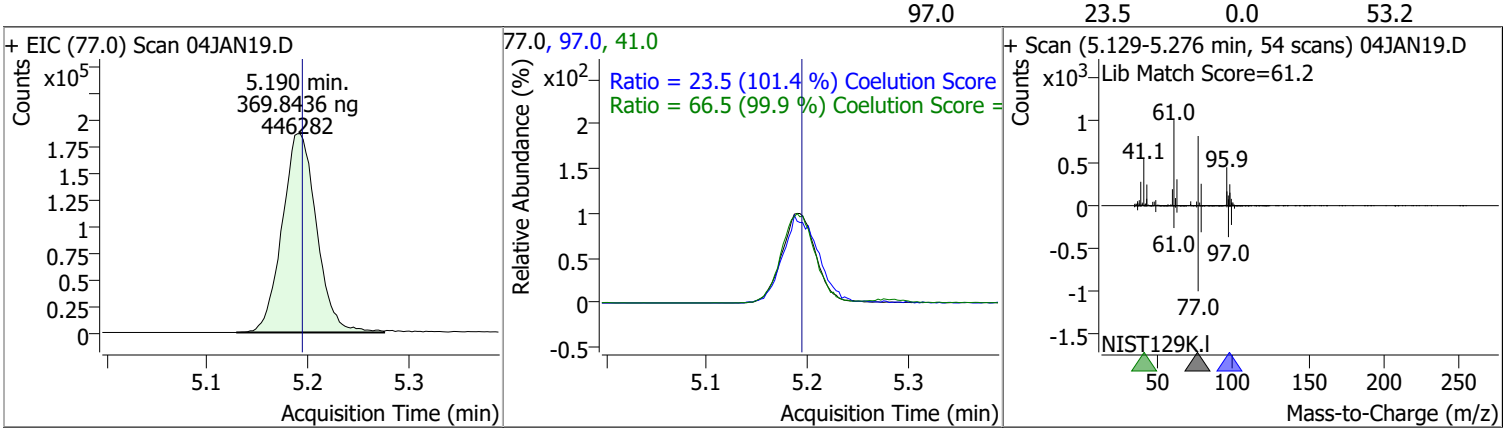
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	380.4437	4.38	0.00	612660	65.0	31.4	2.1	62.1
					83.0	13.1	0.0	43.7



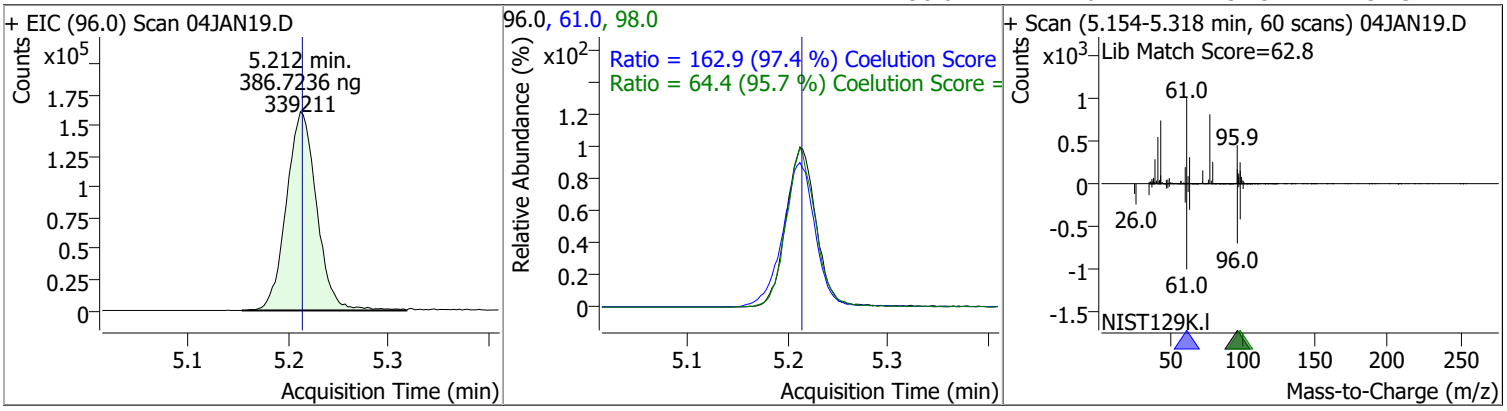




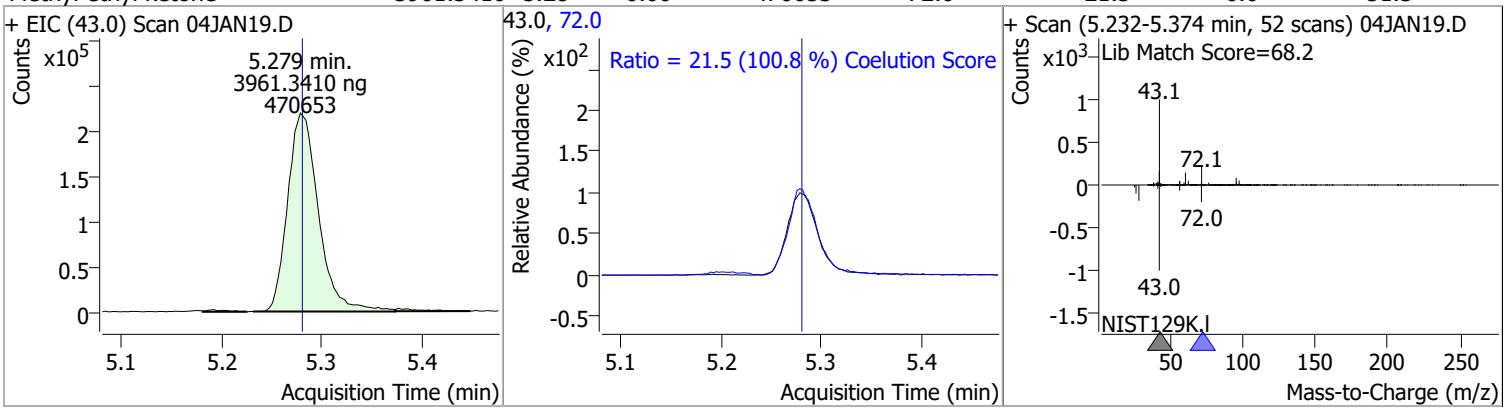
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	369.8436	5.19	-0.01	446282	41.0	66.5	36.5	96.5
					97.0	23.5	0.0	53.2



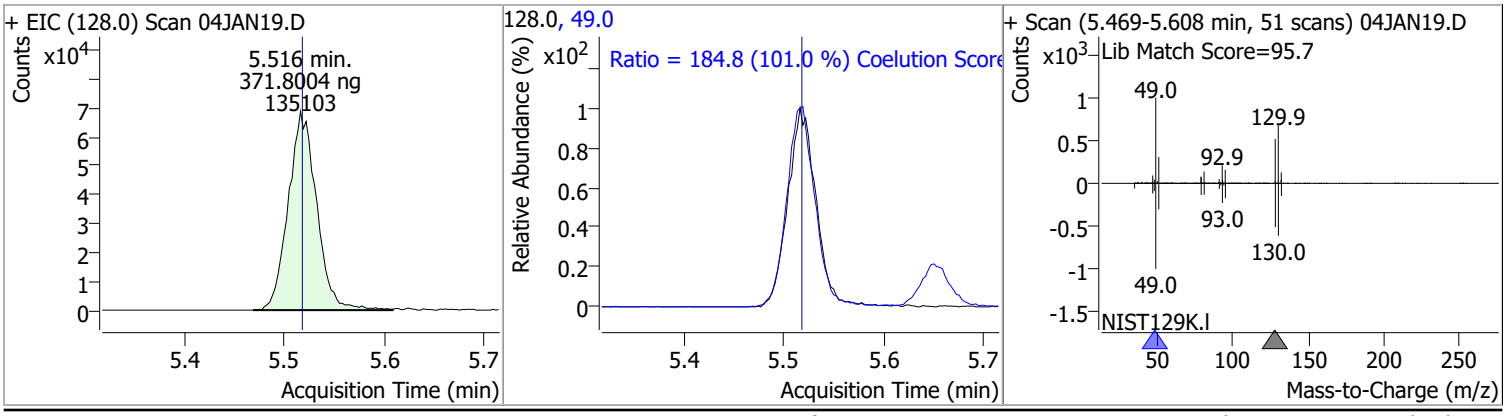
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	386.7236	5.21	0.00	339211	61.0	162.9	137.2	197.2
					98.0	64.4	37.3	97.3



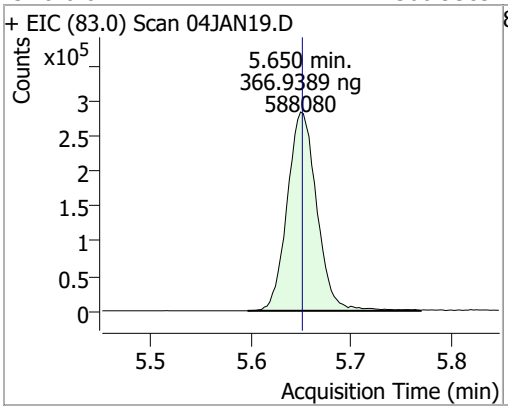
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	3961.3410	5.28	0.00	470653	72.0	21.5	0.0	51.3

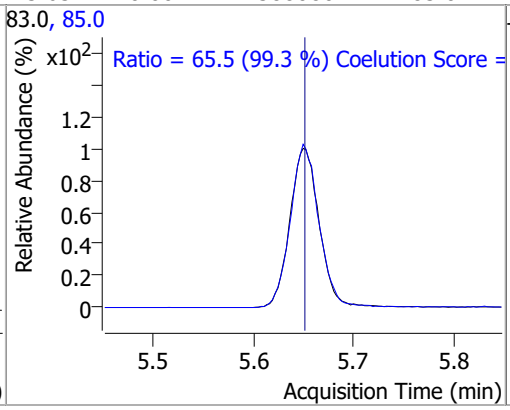


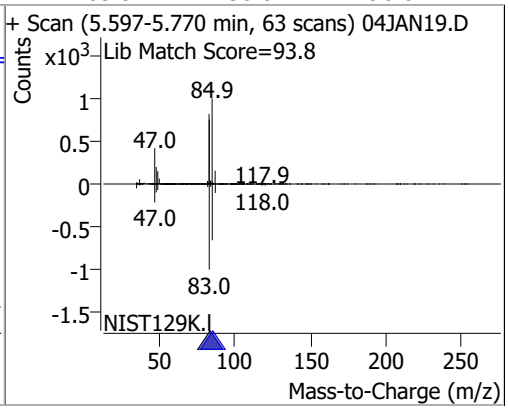
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	371.8004	5.52	0.00	135103	49.0	184.8	152.9	212.9



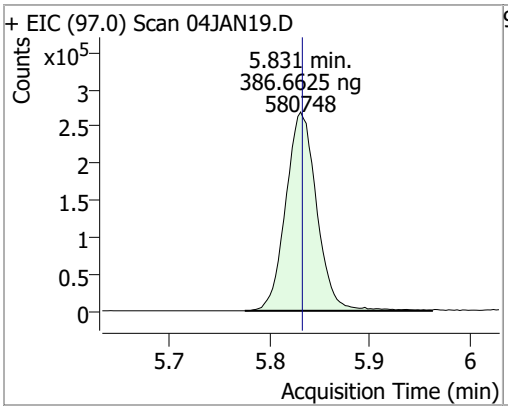
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	366.9389	5.65	0.00	588080	85.0	65.5	36.0	96.0

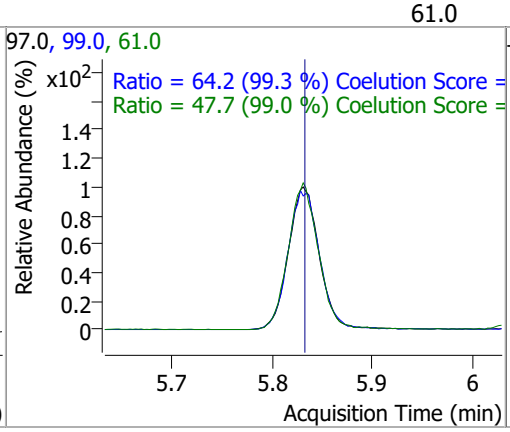


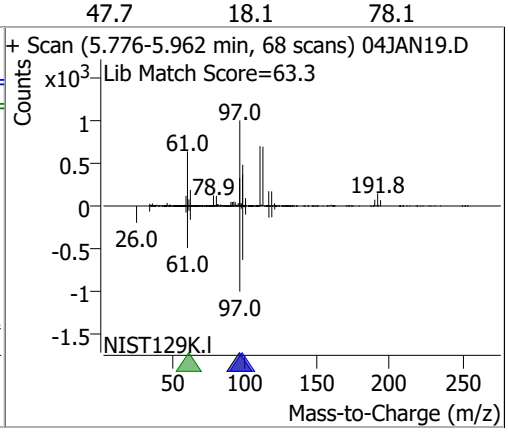




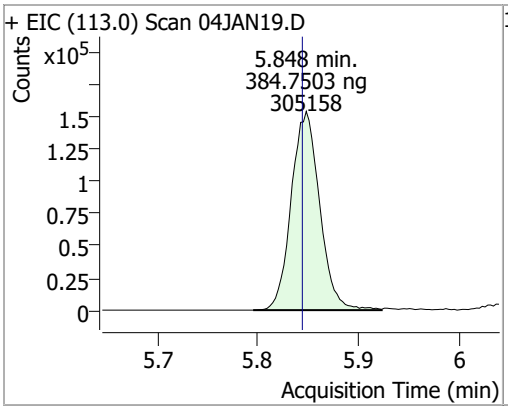
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	386.6625	5.83	0.00	580748	99.0	64.2	34.7	94.7
					61.0	47.7	18.1	78.1

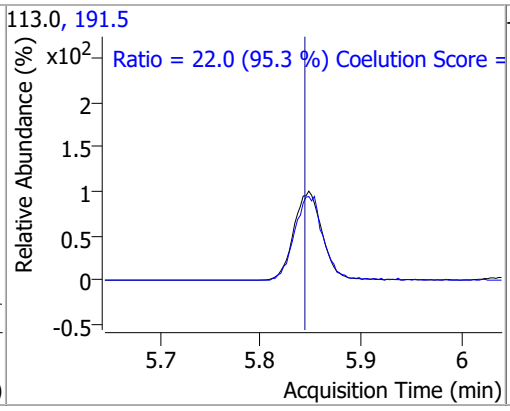


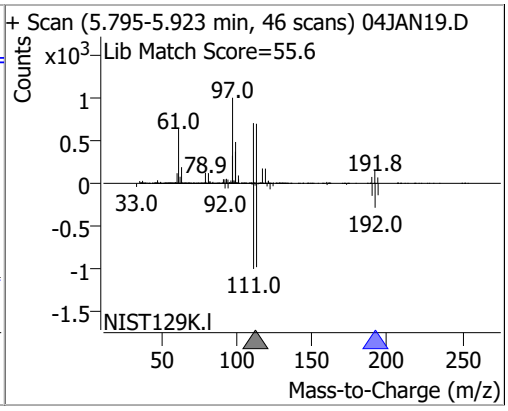




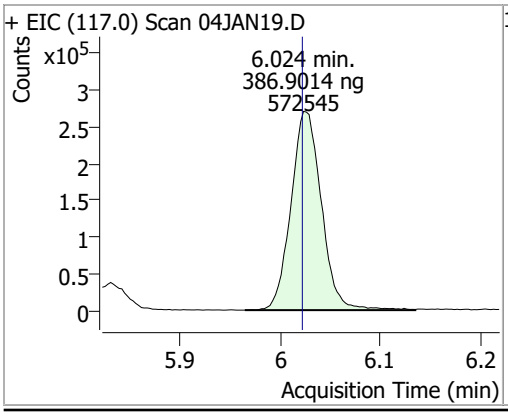
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	384.7503	5.85	0.00	305158	191.5	22.0	0.0	53.1

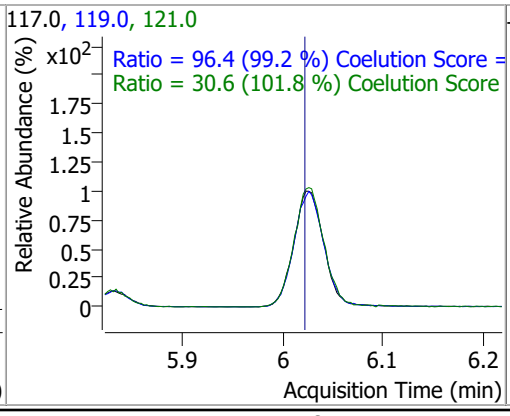


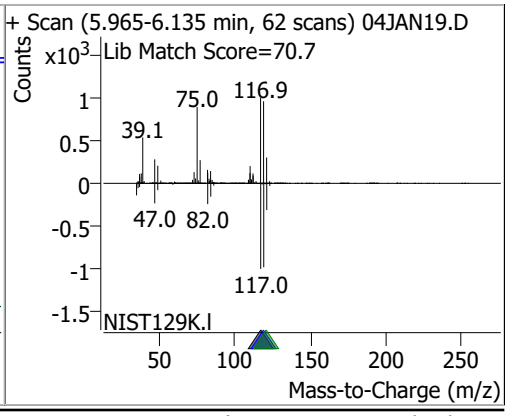




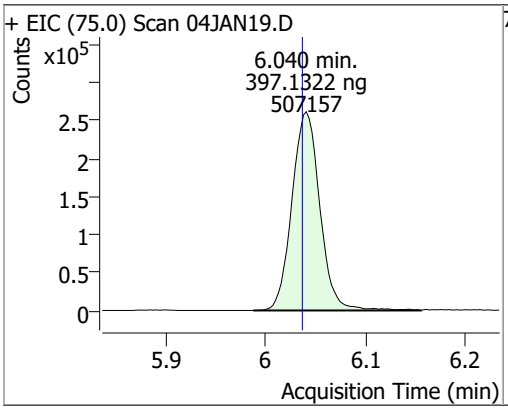
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	386.9014	6.02	0.00	572545	119.0	96.4	67.2	127.2
					121.0	30.6	0.1	60.1

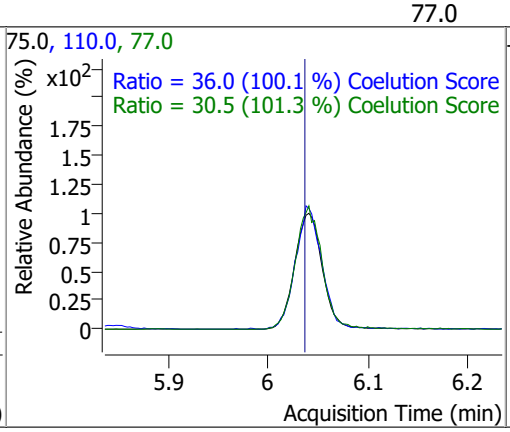


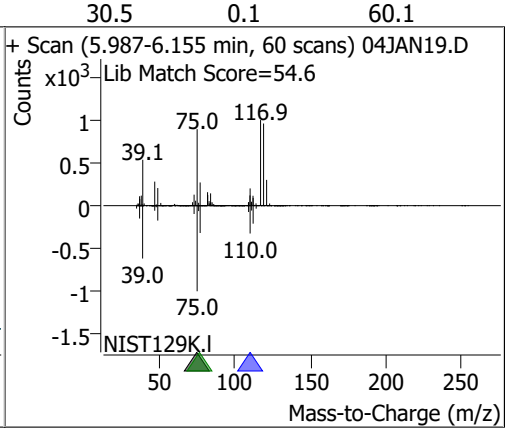




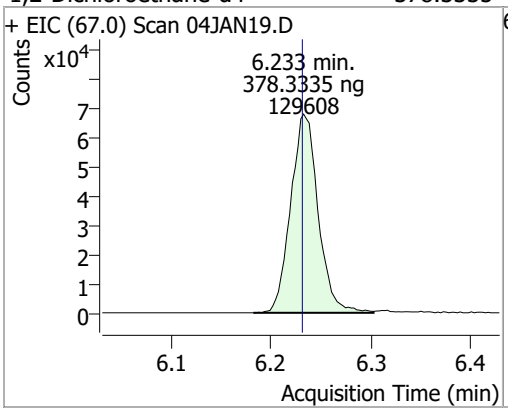
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	397.1322	6.04	0.00	507157	110.0	36.0	5.9	65.9
					77.0	30.5	0.1	60.1

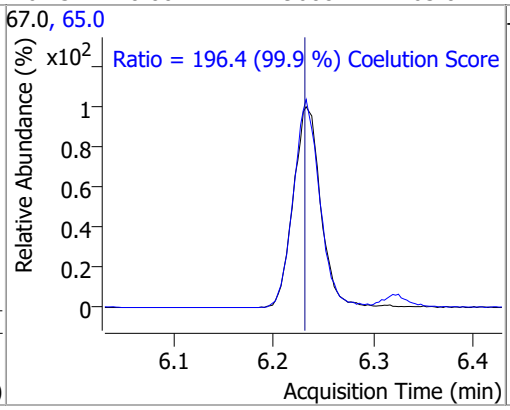


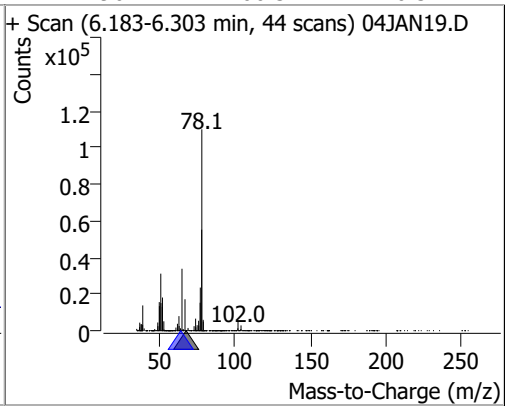




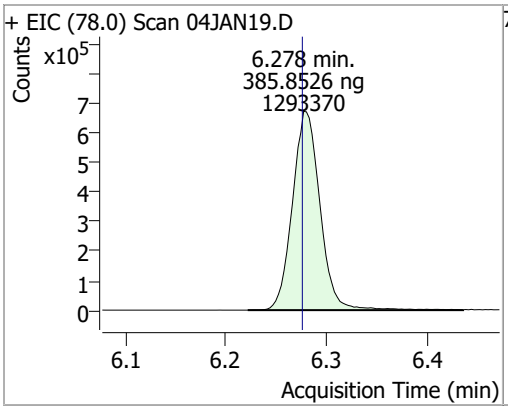
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	378.3335	6.23	0.00	129608	65.0	196.4	166.5	226.5

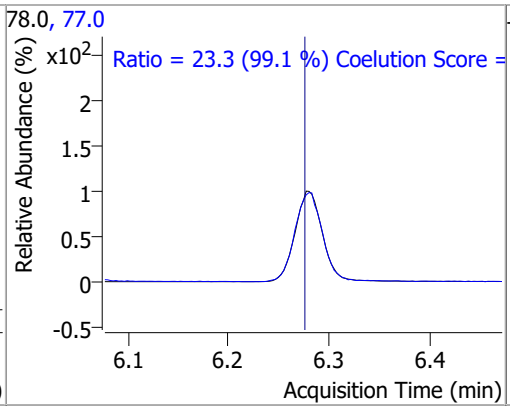


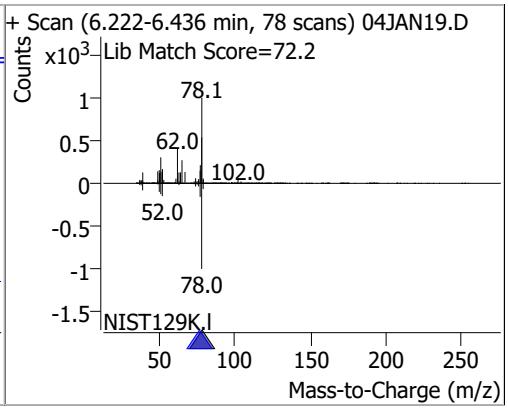




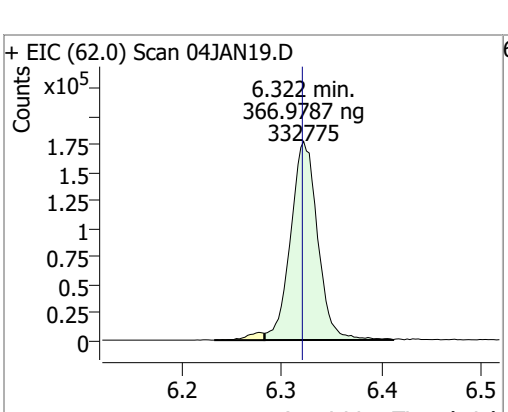
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	385.8526	6.28	0.00	1293370	77.0	23.3	0.0	53.5

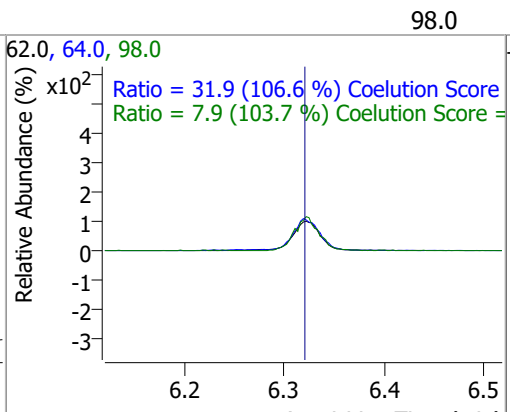


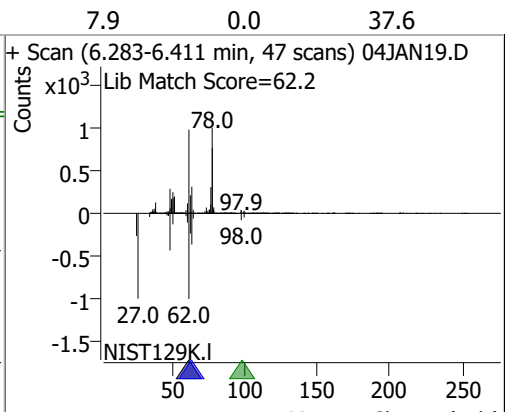




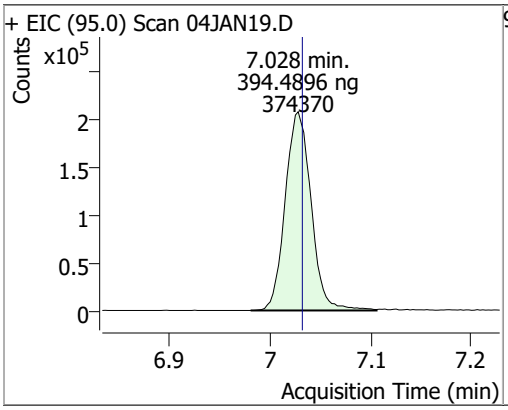
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	366.9787	6.32	0.00	332775	64.0	31.9	0.0	59.9
					98.0	7.9	0.0	37.6

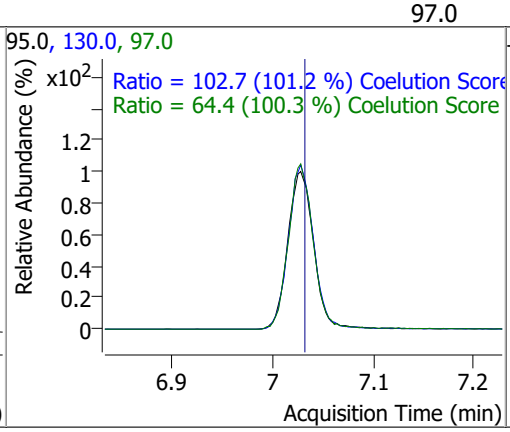


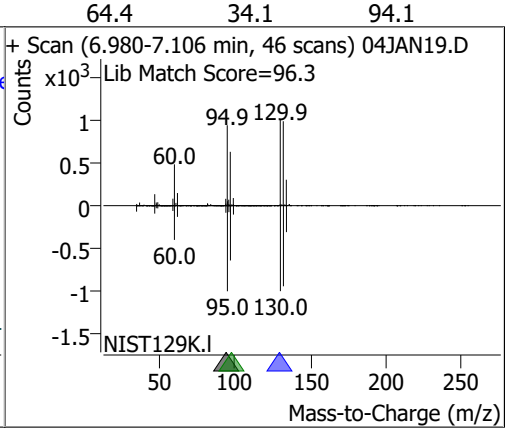




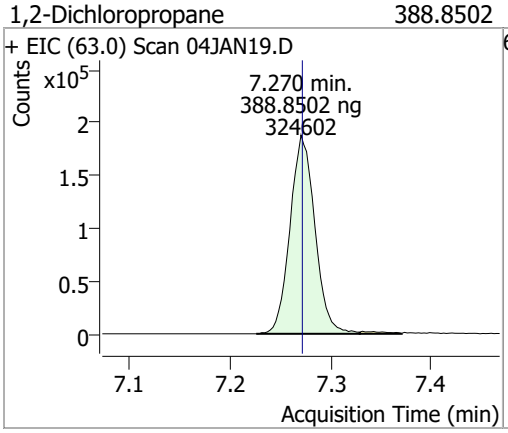
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	394.4896	7.03	0.00	374370	130.0	102.7	71.5	131.5
					97.0	64.4	34.1	94.1

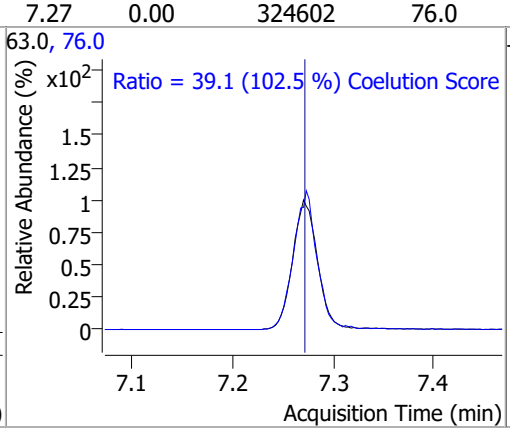


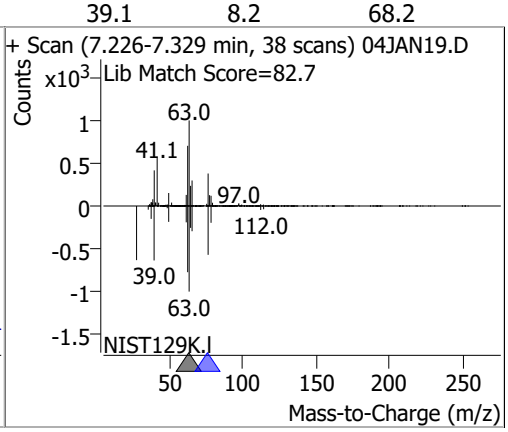




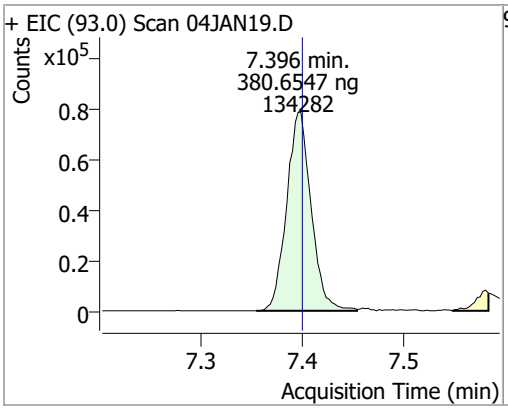
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	388.8502	7.27	0.00	324602	76.0	39.1	8.2	68.2

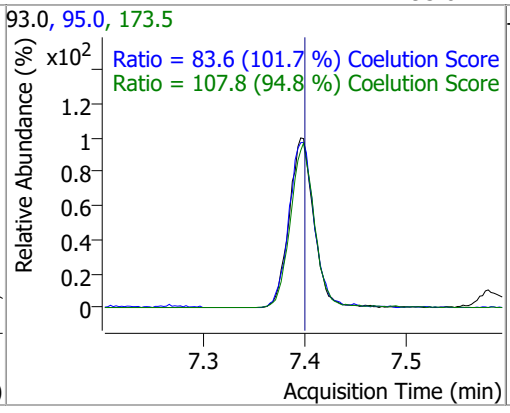


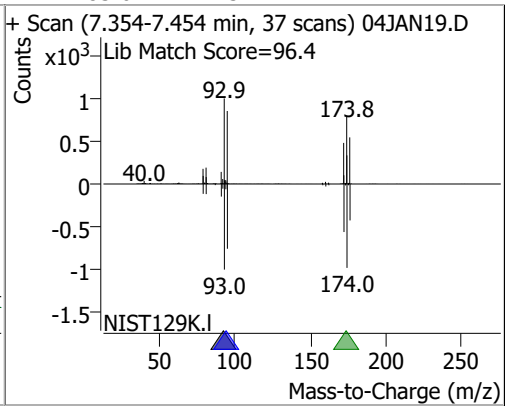




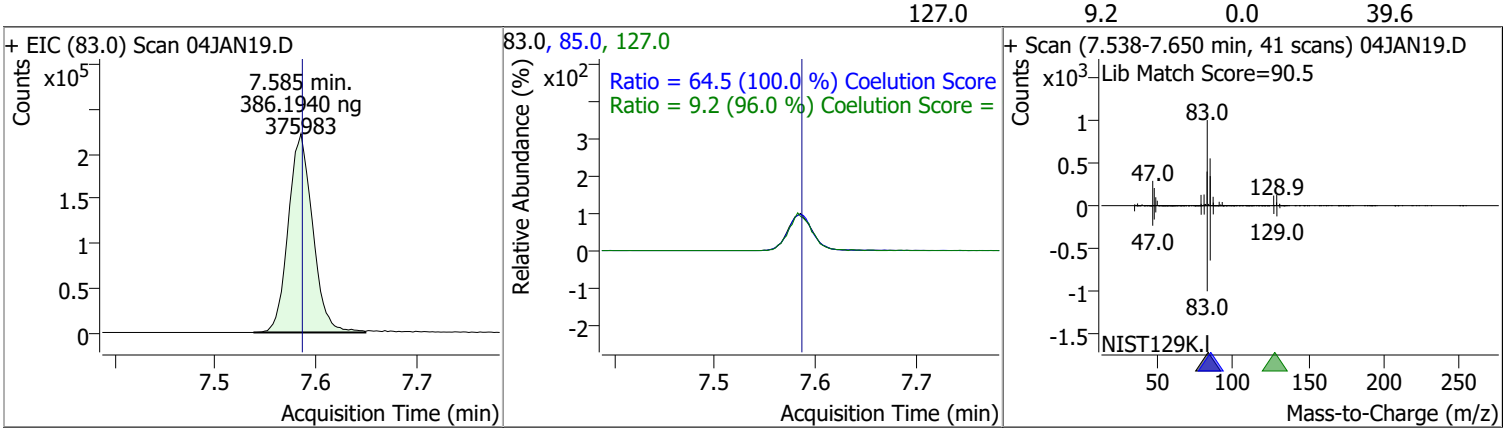
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	380.6547	7.40	0.00	134282	173.5	107.8	83.7	143.7
					95.0	83.6	52.2	112.2



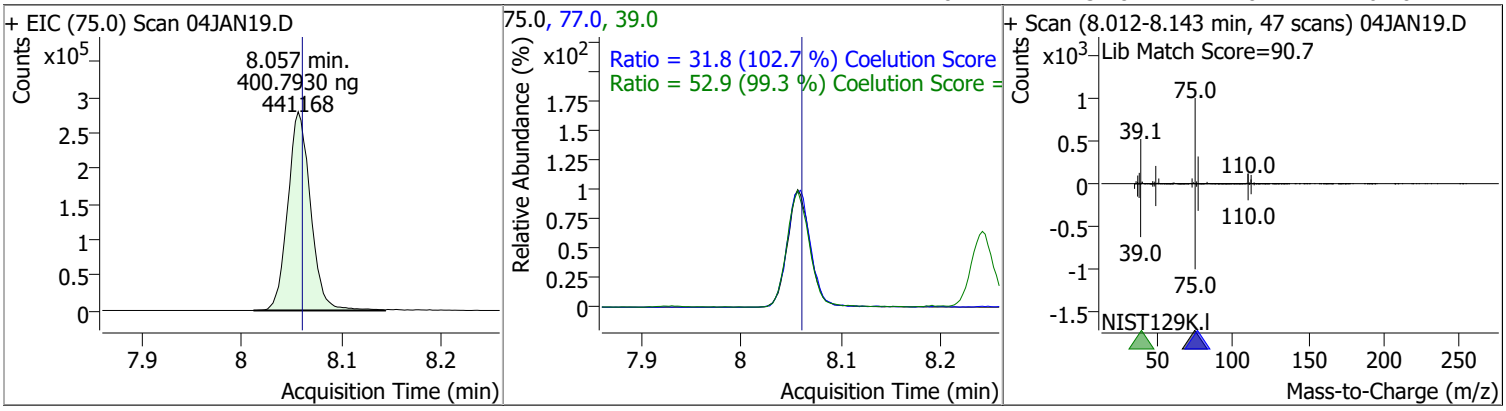




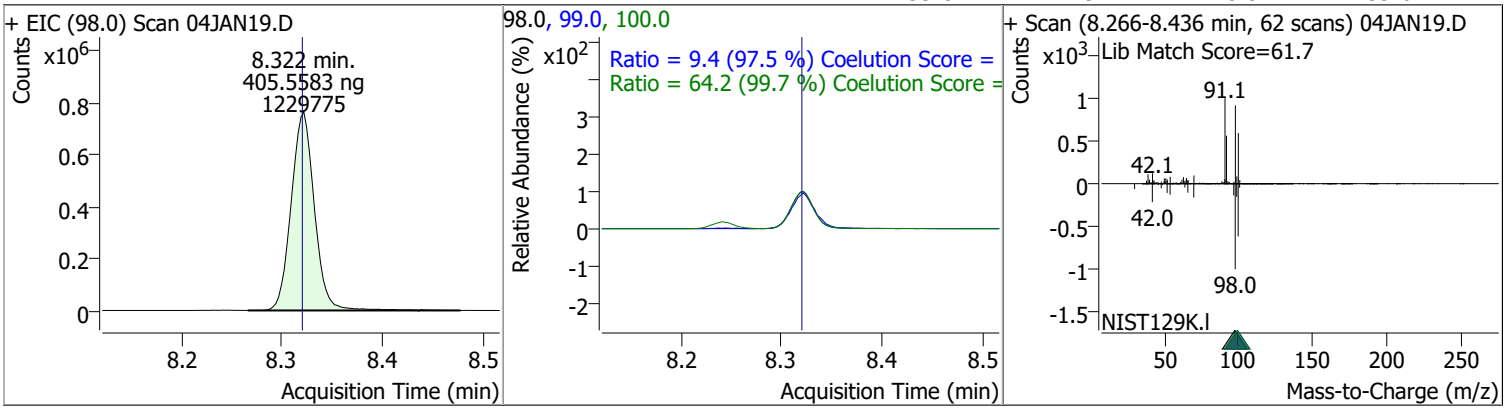
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	386.1940	7.59	0.00	375983	85.0	64.5	34.5	94.5
					127.0	9.2	0.0	39.6



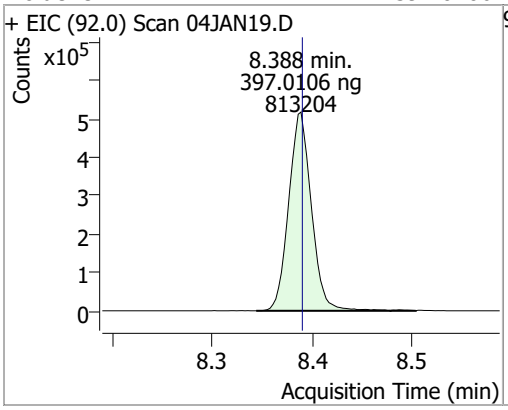
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	400.7930	8.06	0.00	441168	39.0	52.9	23.3	83.3
					77.0	31.8	1.0	61.0

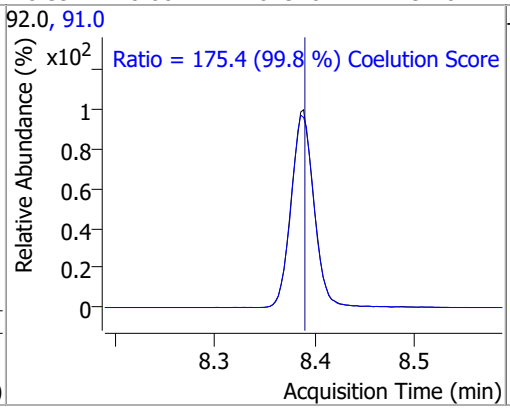


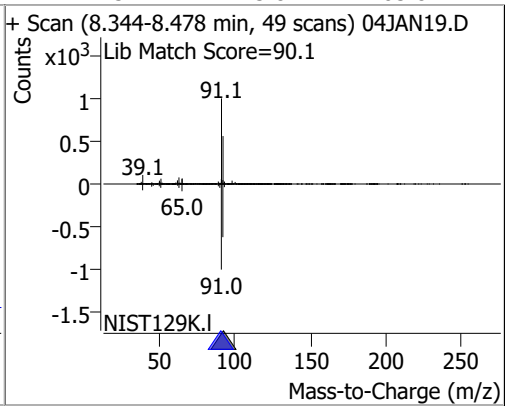
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	405.5583	8.32	0.00	1229775	100.0	64.2	34.4	94.4
					99.0	9.4	0.0	39.6



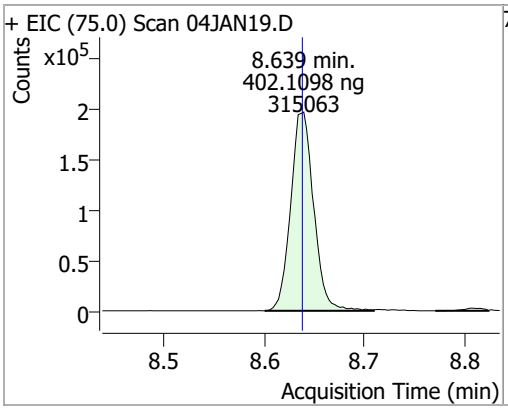
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	397.0106	8.39	0.00	813204	91.0	175.4	145.8	205.8

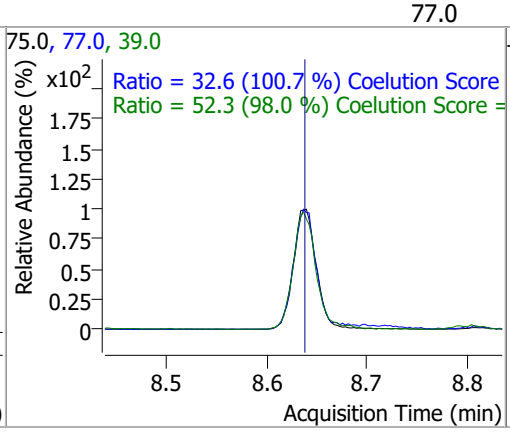


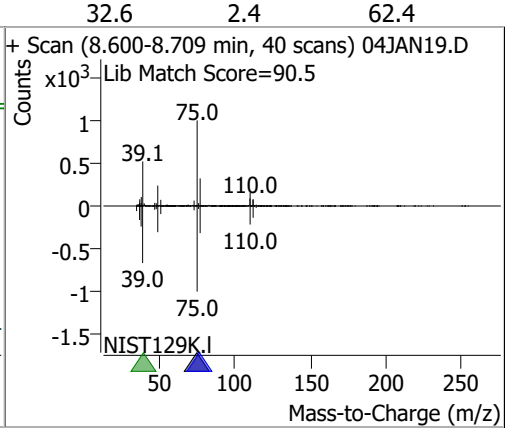




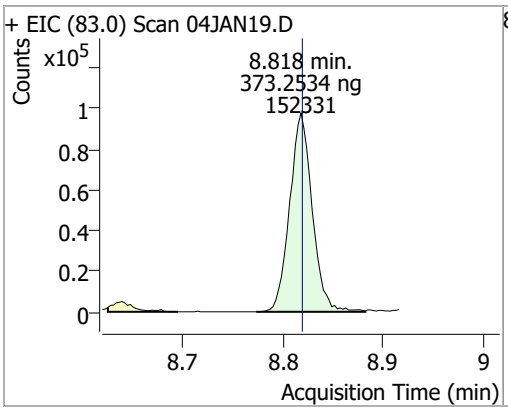
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	402.1098	8.64	0.00	315063	39.0	52.3	23.4	83.4
					77.0	32.6	2.4	62.4

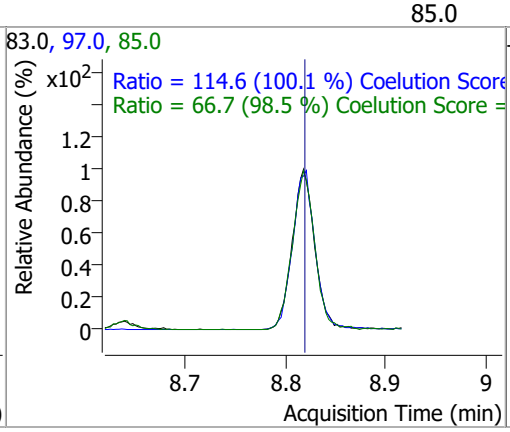


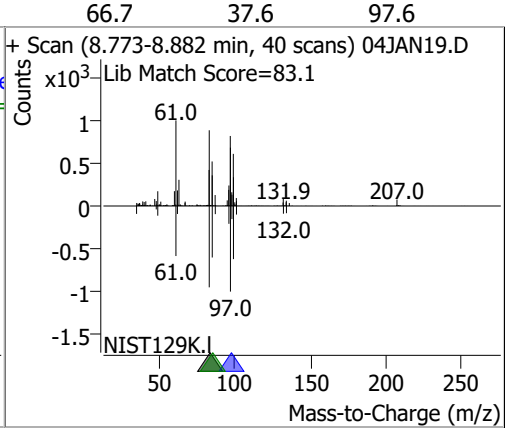




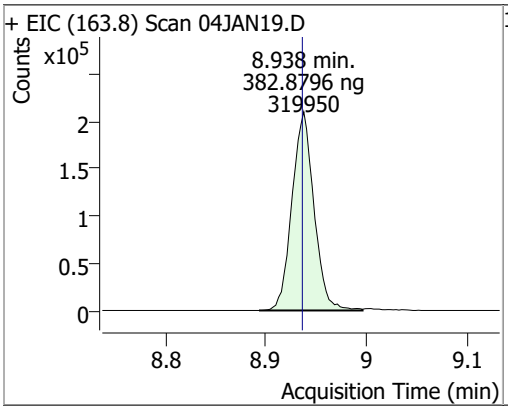
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	373.2534	8.82	0.00	152331	97.0	114.6	84.6	144.6
					85.0	66.7	37.6	97.6

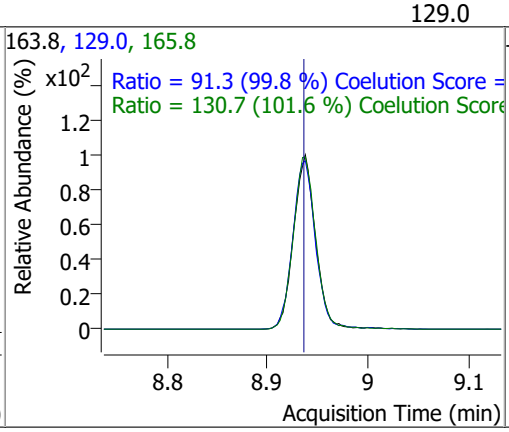


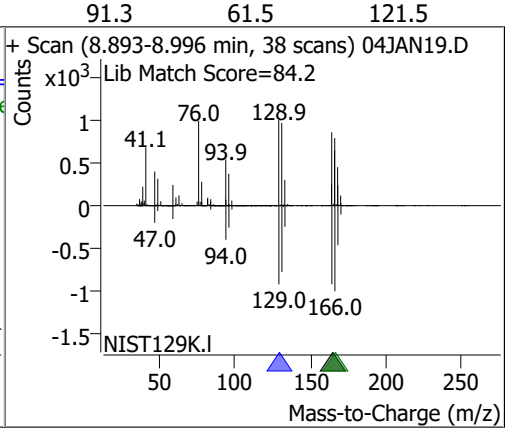




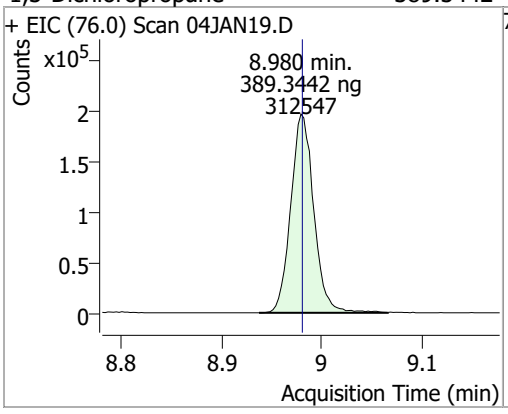
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	382.8796	8.94	0.00	319950	165.8	130.7	98.6	158.6
					129.0	91.3	61.5	121.5

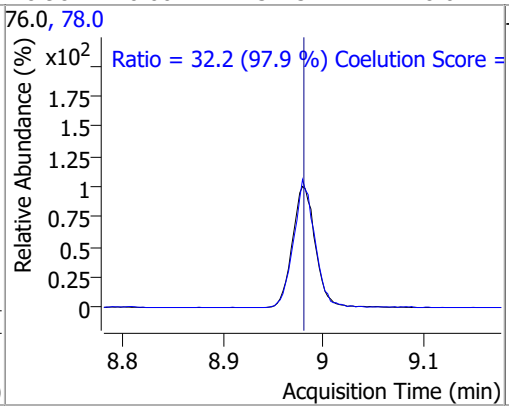


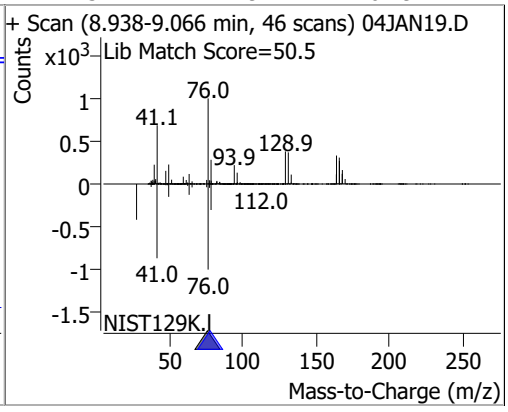




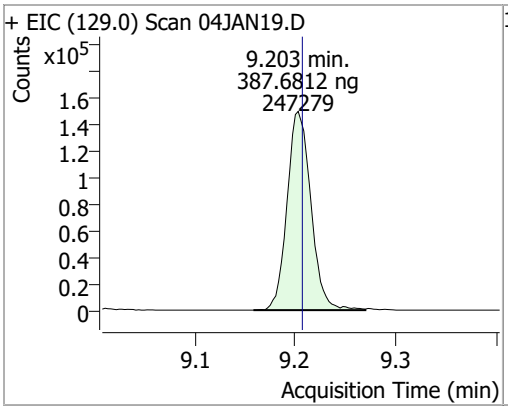
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	389.3442	8.98	0.00	312547	78.0	32.2	2.9	62.9

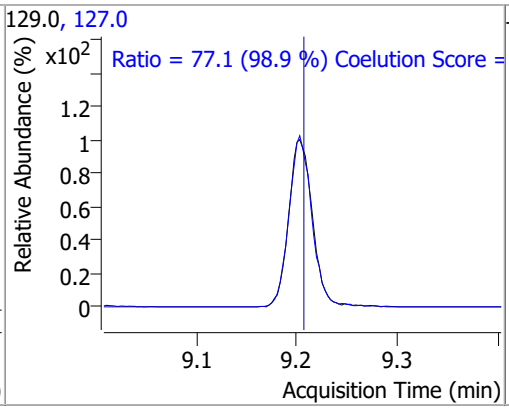


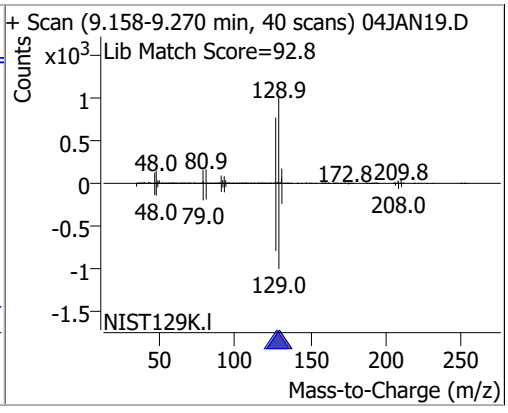




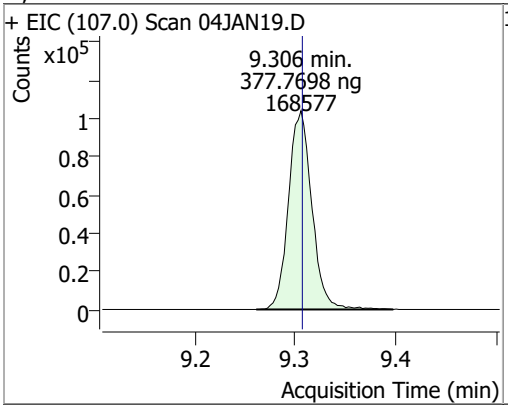
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	387.6812	9.20	0.00	247279	127.0	77.1	48.0	108.0

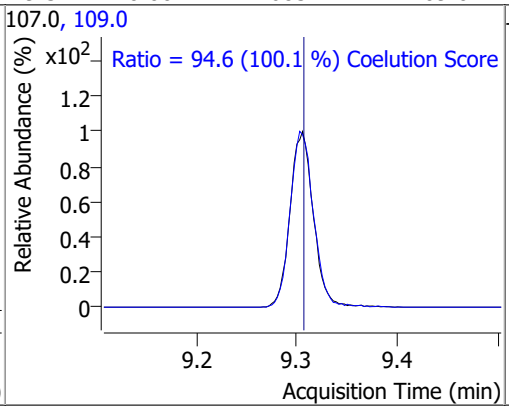


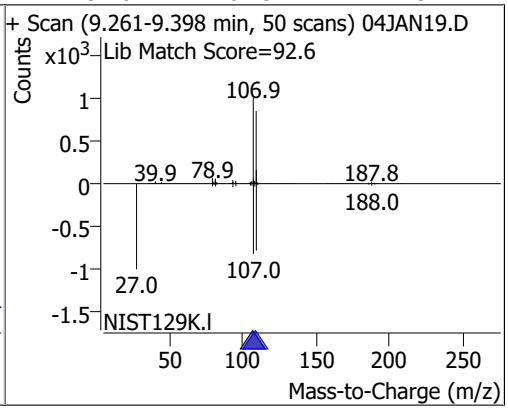




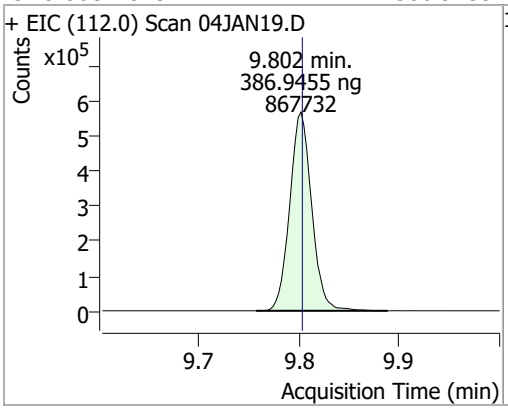
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	377.7698	9.31	0.00	168577	109.0	94.6	64.5	124.5

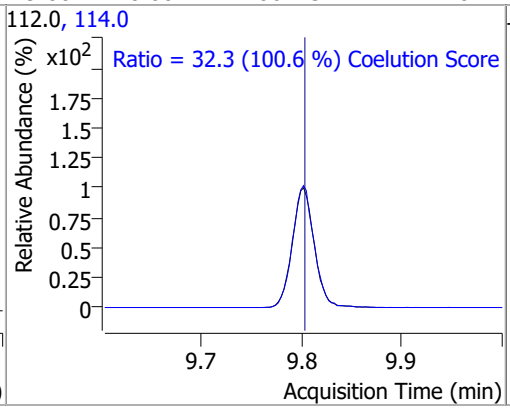


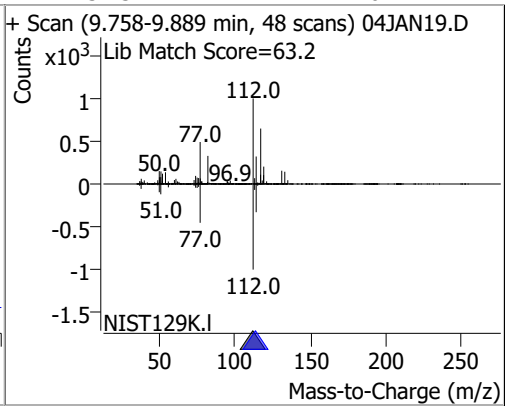




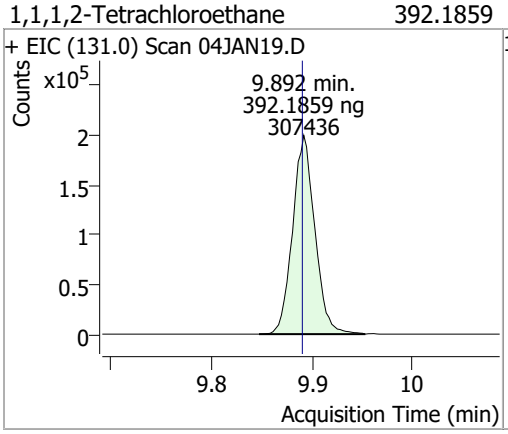
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	386.9455	9.80	0.00	867732	114.0	32.3	2.1	62.1

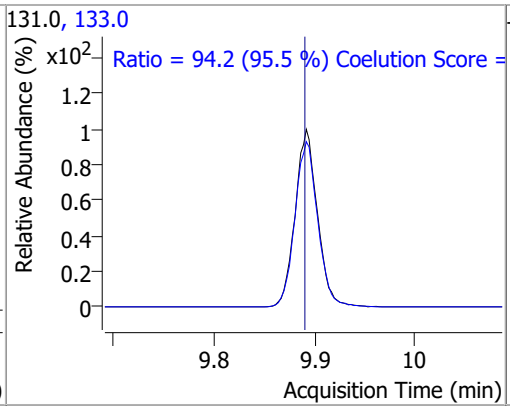


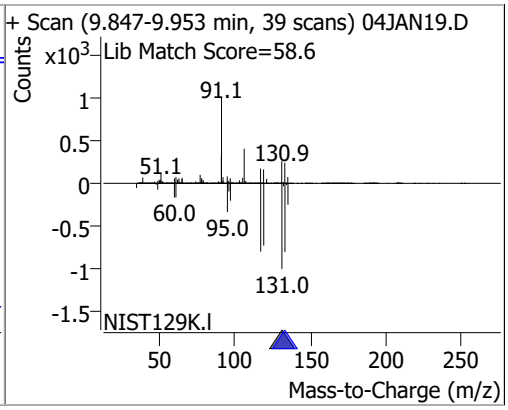




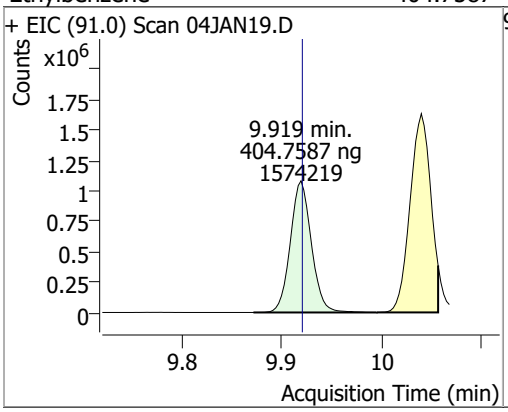
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	392.1859	9.89	0.00	307436	133.0	94.2	68.6	128.6

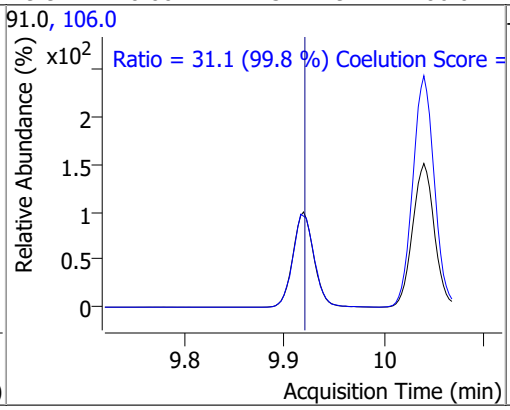


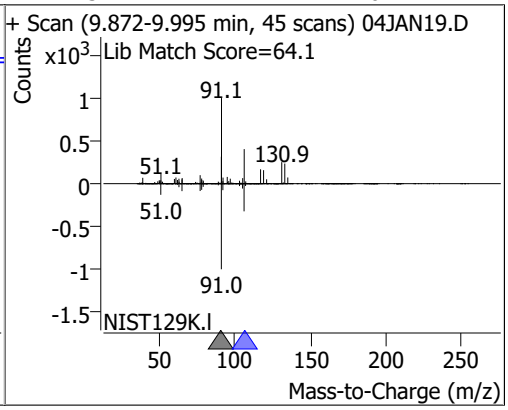




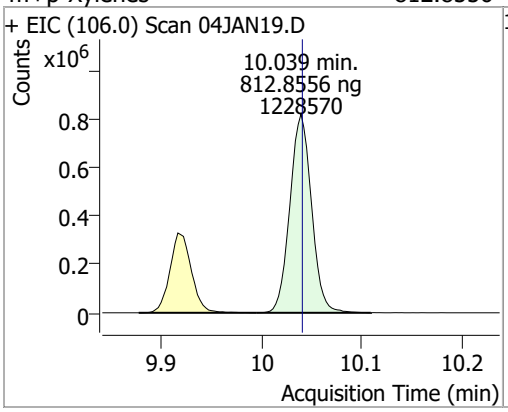
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	404.7587	9.92	0.00	1574219	106.0	31.1	1.1	61.1

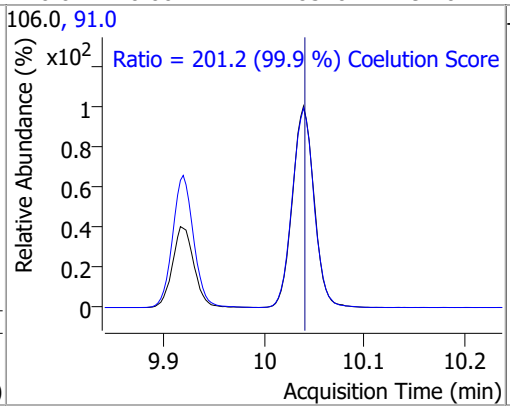


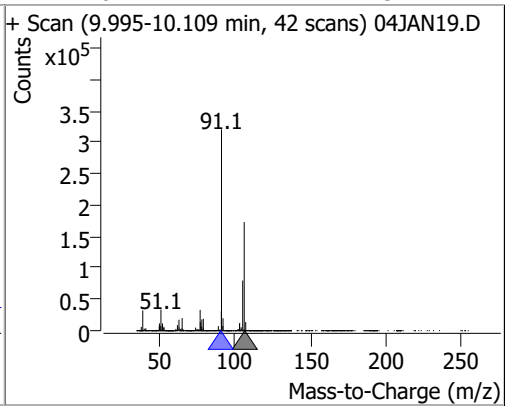




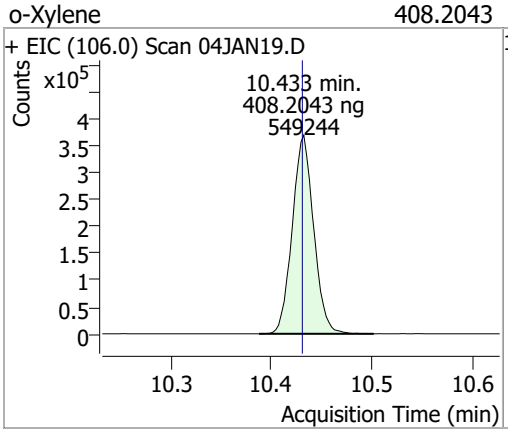
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	812.8556	10.04	0.00	1228570	91.0	201.2	171.4	231.4

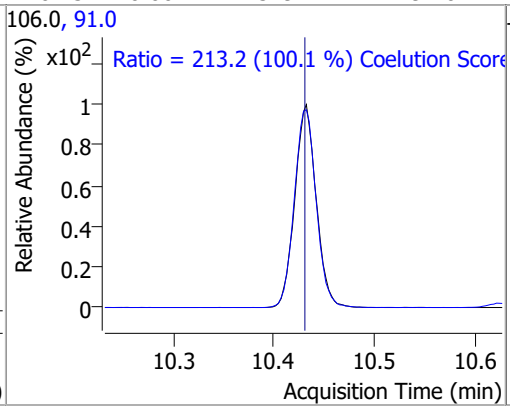


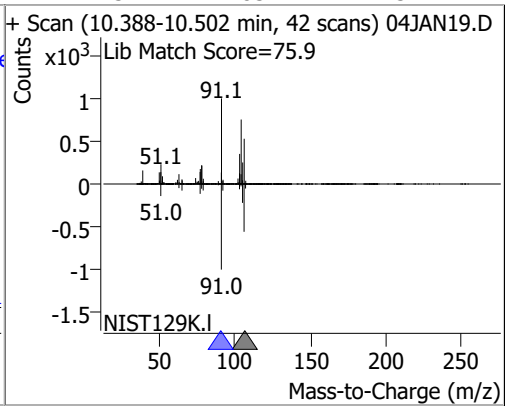




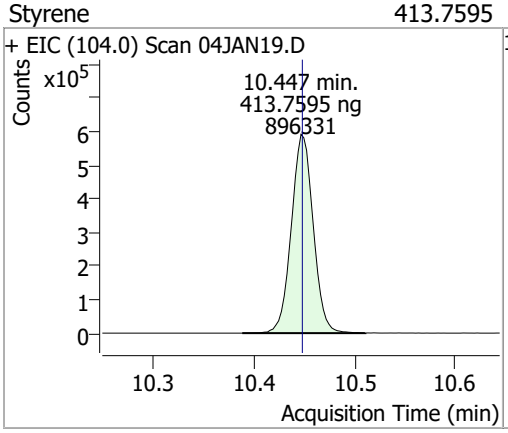
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	408.2043	10.43	0.00	549244	91.0	213.2	183.1	243.1

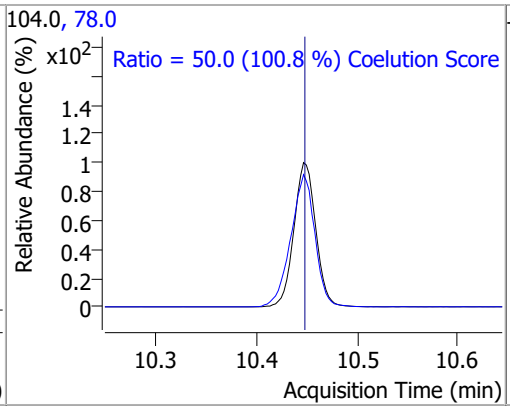


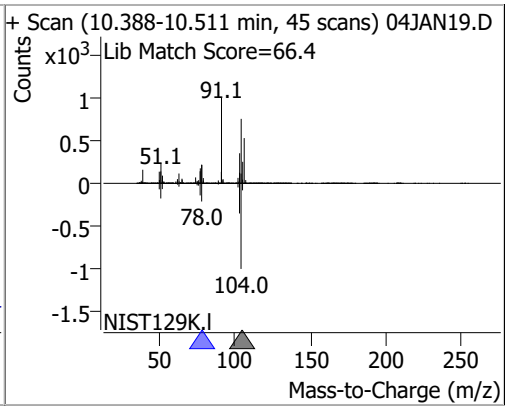




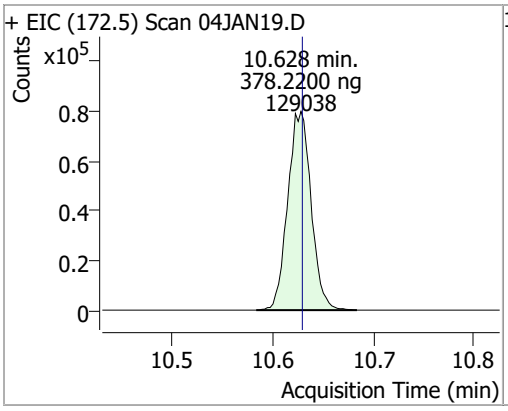
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	413.7595	10.45	0.00	896331	78.0	50.0	19.6	79.6

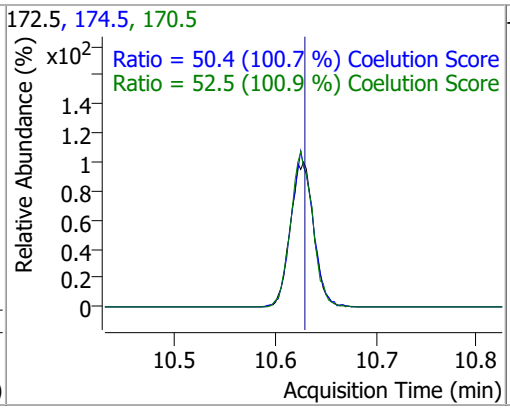


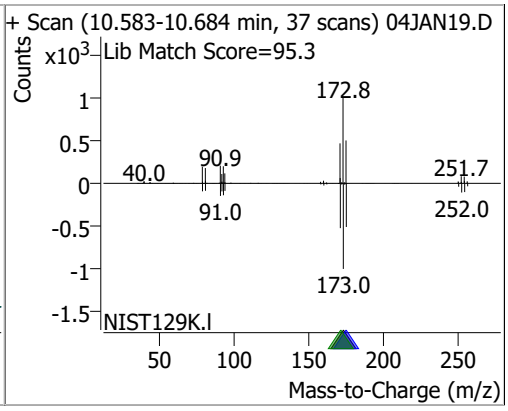




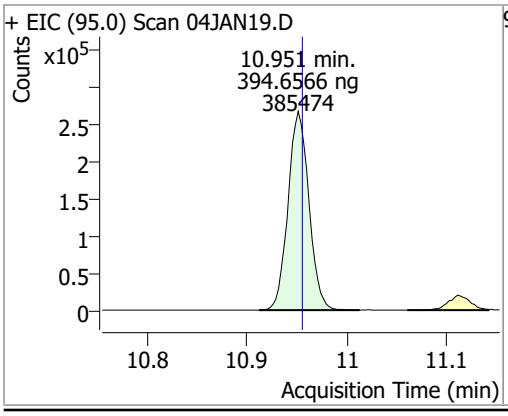
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	378.2200	10.63	0.00	129038	170.5	52.5	22.1	82.1
					174.5	50.4	20.1	80.1

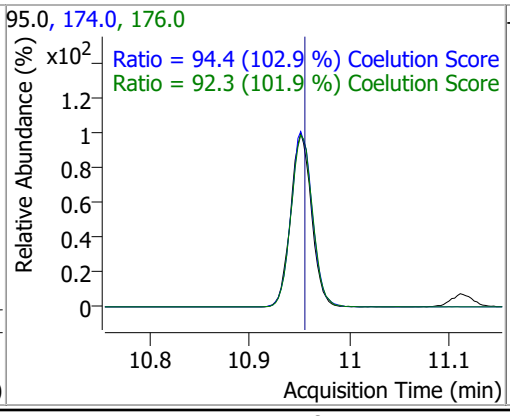


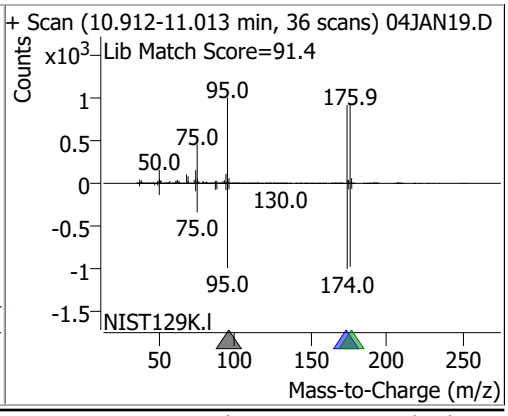




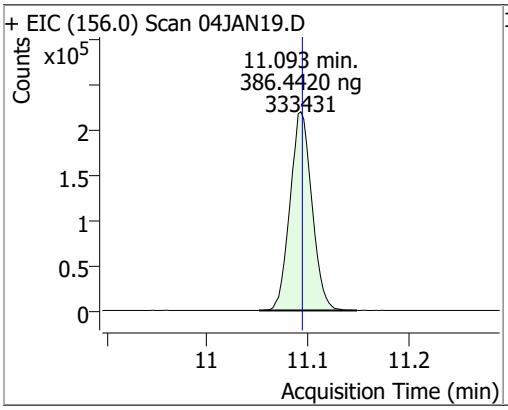
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	394.6566	10.95	0.00	385474	174.0	94.4	61.7	121.7
					176.0	92.3	60.6	120.6

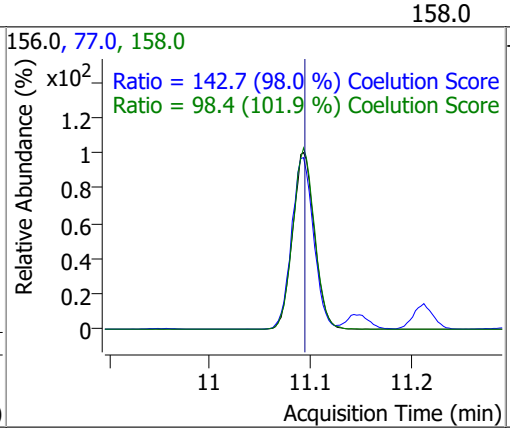


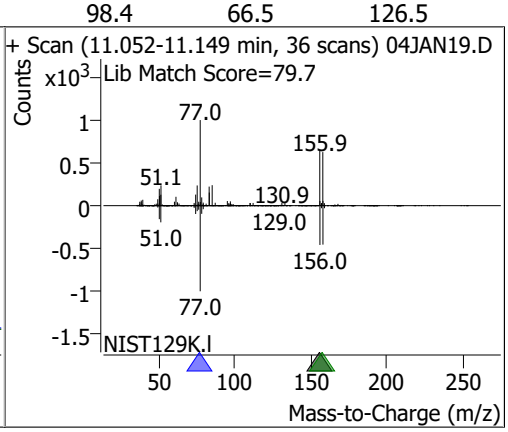




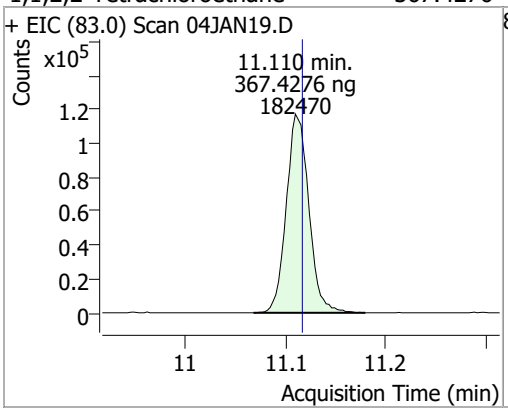
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	386.4420	11.09	0.00	333431	77.0	142.7	115.7	175.7
					158.0	98.4	66.5	126.5

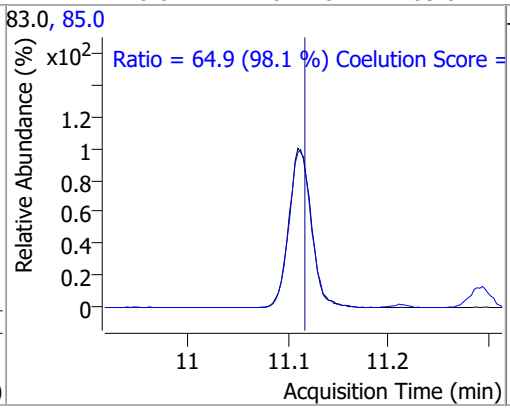


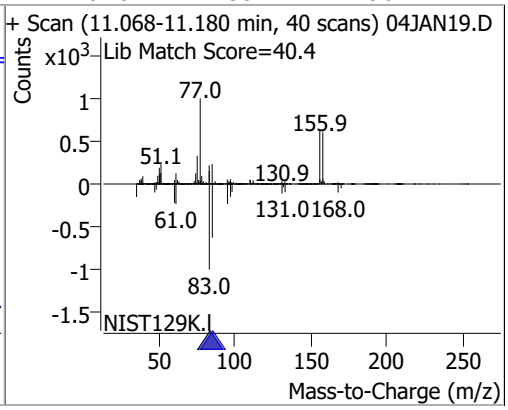




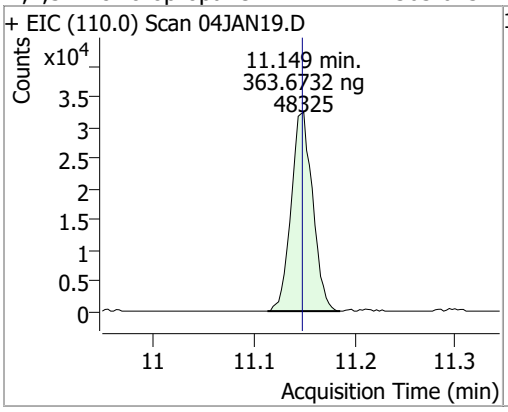
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	367.4276	11.11	-0.01	182470	85.0	64.9	36.2	96.2

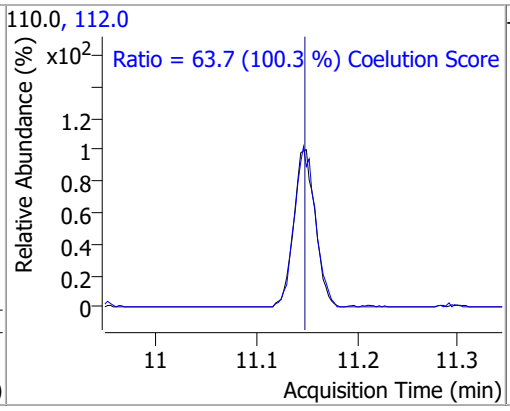


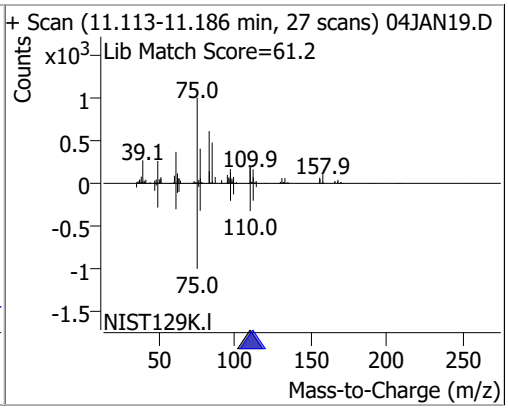




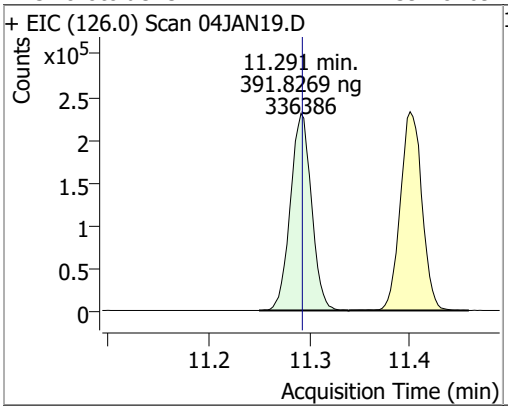
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	363.6732	11.15	0.00	48325	112.0	63.7	33.5	93.5

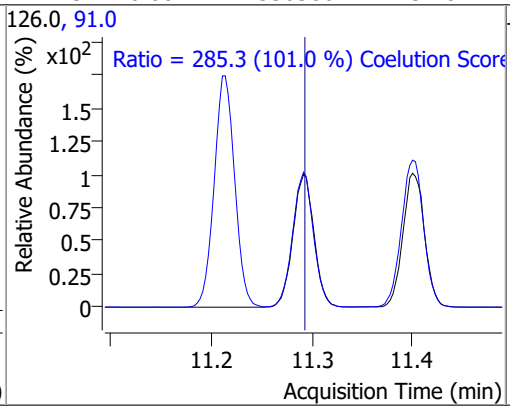


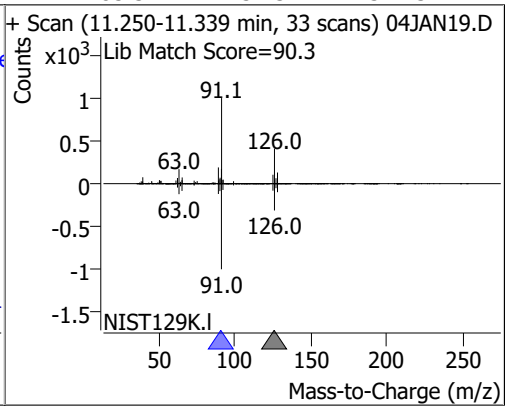




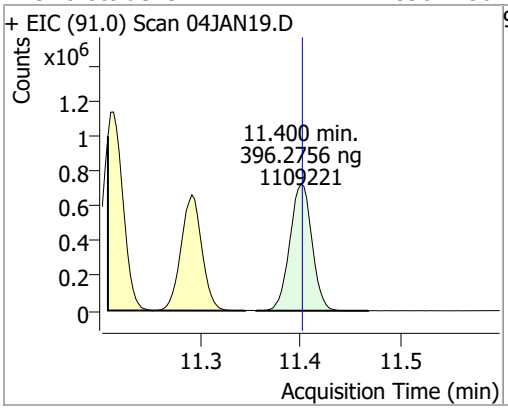
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	391.8269	11.29	0.00	336386	91.0	285.3	252.3	312.3

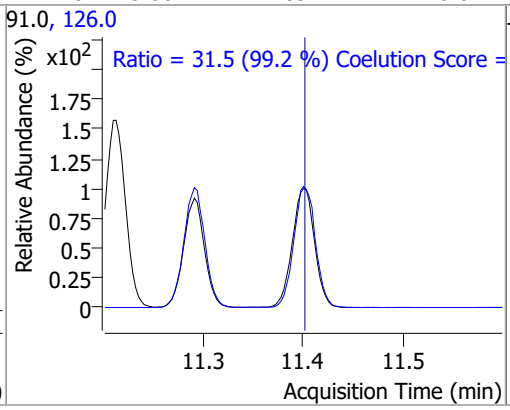


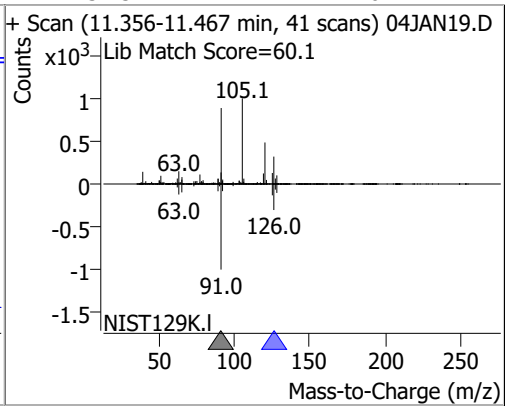




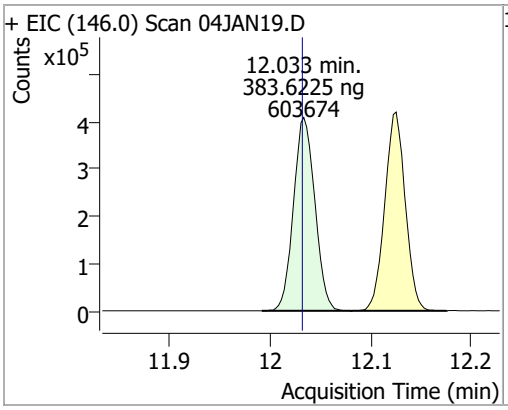
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	396.2756	11.40	0.00	1109221	126.0	31.5	1.7	61.7

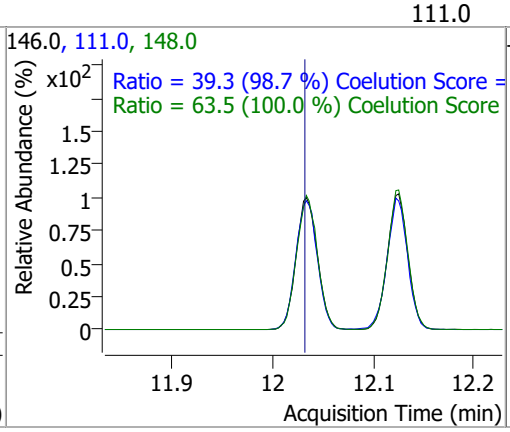


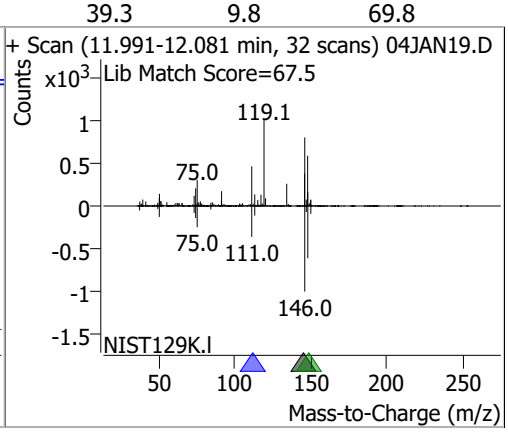




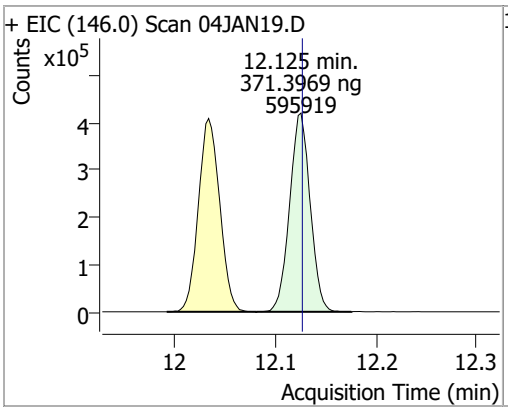
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	383.6225	12.03	0.00	603674	148.0	63.5	33.6	93.6
					111.0	39.3	9.8	69.8

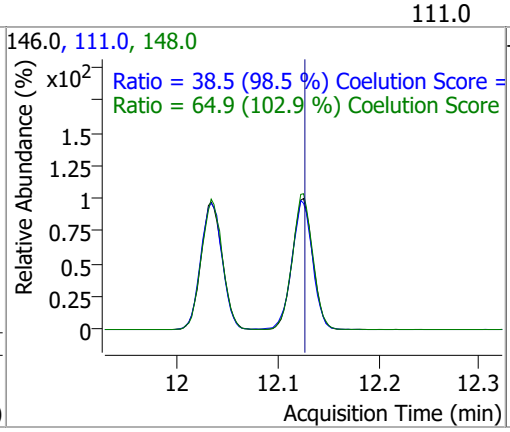


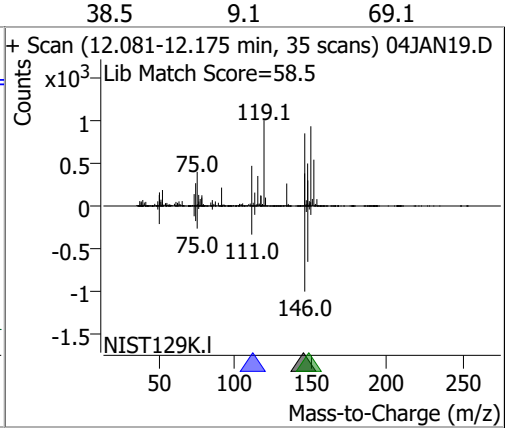




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	371.3969	12.13	0.00	595919	148.0	64.9	33.1	93.1
					111.0	38.5	9.1	69.1

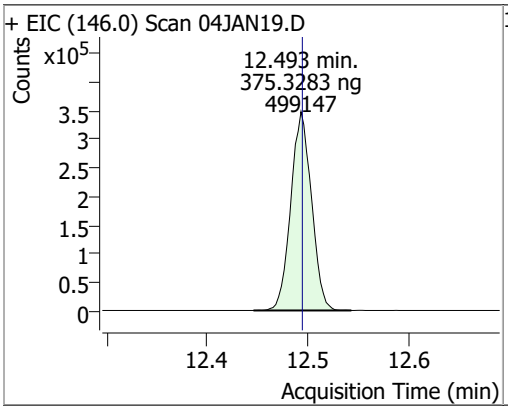


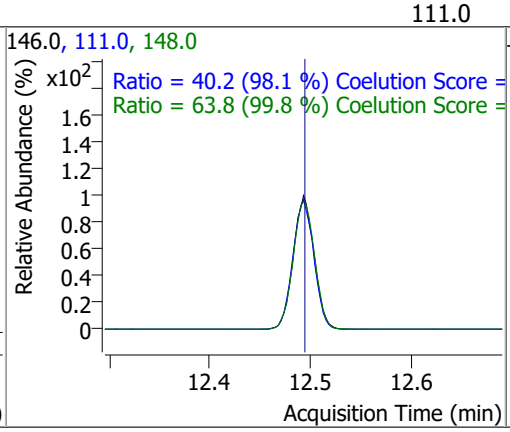


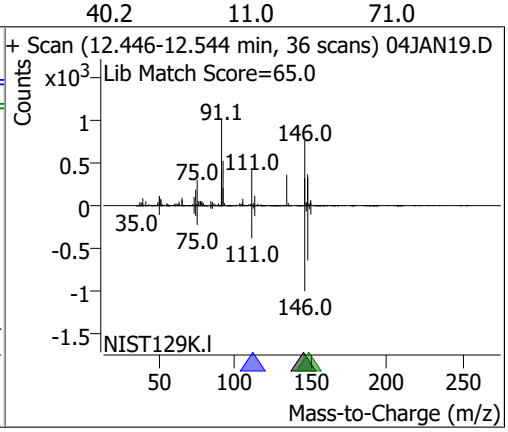


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	375.3283	12.49	0.00	499147	148.0	63.8	33.9	93.9
					111.0	40.2	11.0	71.0







Quantitation Results Report (QT Reviewed)

Data File	04JAN21.D	Operator	MSC
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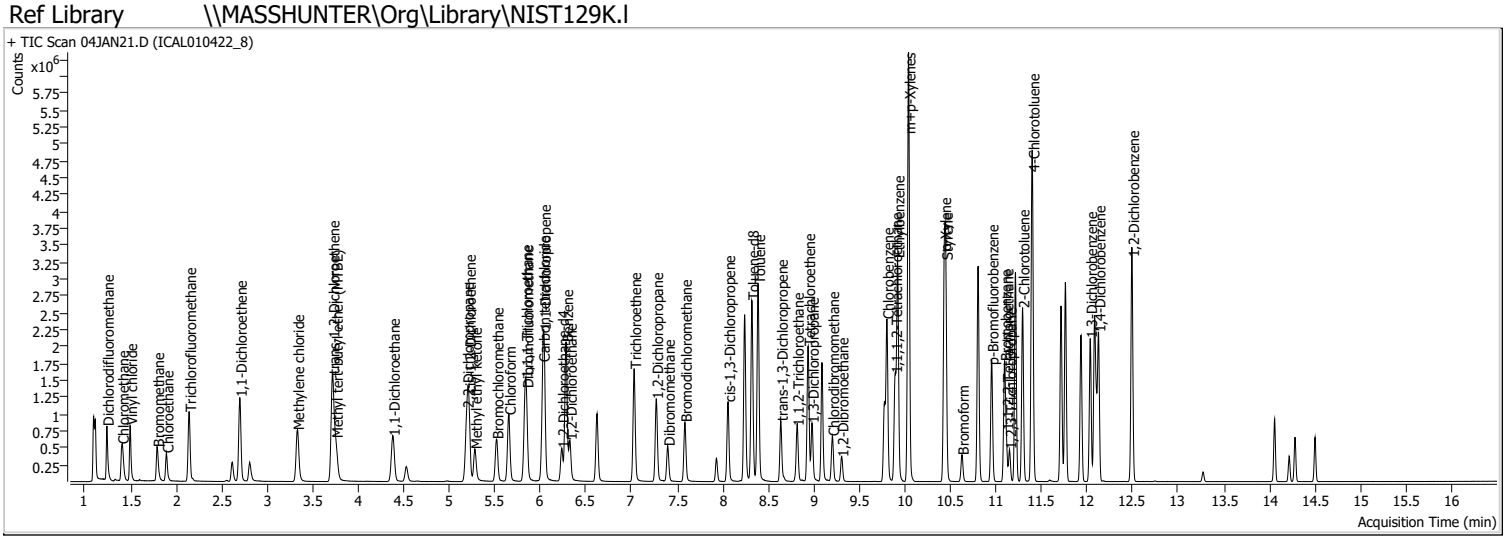
Vial	21	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

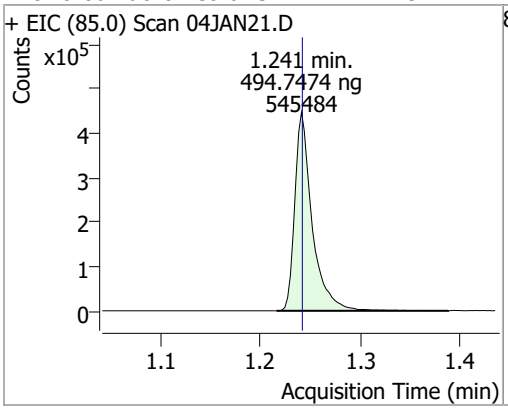


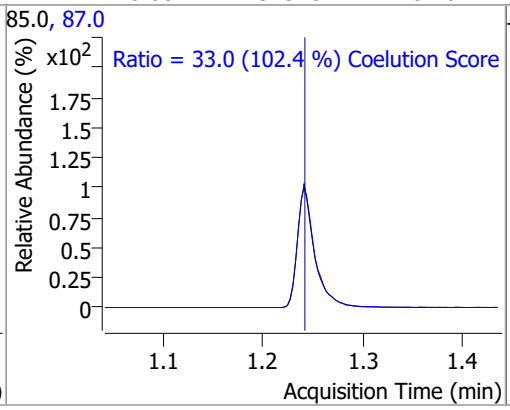
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	841364	250.0000	ng		-0.003
M Chlorobenzene-d5	9.774	82.0	313585	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.103	152.0	262971	250.0000	ng		0.003
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	404568	510.3991	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 204.16%		*	
S 1,2-Dichloroethane-d4	6.233	67.0	174713	510.3080	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 204.12%		*	
S Toluene-d8	8.319	98.0	1644540	544.2136	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 217.69%		*	
S p-Bromofluorobenzene	10.949	95.0	521580	541.3964	ng		-0.006
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 216.56%		*	
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.241	85.0	545484	494.7474	ng		99
T Chloromethane	1.406	50.0	642582	480.1747	ng		100
T Vinyl chloride	1.495	62.0	600092	498.3563	ng		95
T Bromomethane	1.793	96.0	277301	515.0141	ng		98
T Chloroethane	1.894	64.0	287041	481.5143	ng		98
T Trichlorofluoromethane	2.145	101.0	731829	489.6475	ng		99
T 1,1-Dichloroethene	2.700	96.0	436507	515.0603	ng		98
T Methylene chloride	3.330	49.0	583438	466.9993	ng		98
T trans-1,2-Dichloroethene	3.718	96.0	440967	510.0097	ng		98
T Methyl tert-butyl ether (MTBE)	3.754	73.0	584294	522.8187	ng		99
T 1,1-Dichloroethane	4.378	63.0	829359	515.3207	ng		99
T 2,2-Dichloropropane	5.190	77.0	601823	499.0473	ng		98
T cis-1,2-Dichloroethene	5.212	96.0	452377	516.0544	ng		99
T Methyl ethyl ketone	5.279	43.0	632539	5327.1253	ng		99
T Bromochloromethane	5.519	128.0	179618	494.6054	ng		98
T Chloroform	5.653	83.0	783422	489.1221	ng		99

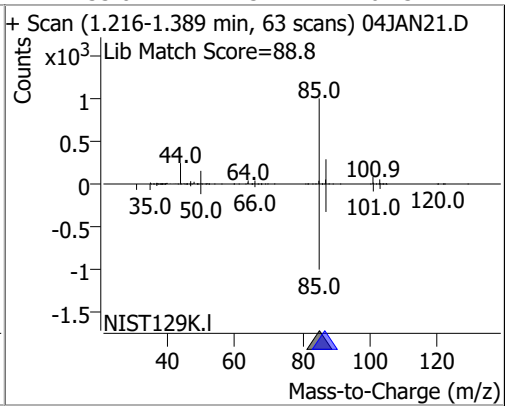
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.834	97.0	778785	518.8312	ng	99
T Carbon tetrachloride	6.024	117.0	770907	521.2630	ng	98
T 1,1-Dichloropropene	6.038	75.0	693669	543.5121	ng	99
T Benzene	6.280	78.0	1714050	511.6658	ng	100
T 1,2-Dichloroethane	6.322	62.0	450739	497.3699	ng	100
T Trichloroethene	7.028	95.0	505400	534.4007	ng	99
T 1,2-Dichloropropane	7.270	63.0	436057	524.1695	ng	100
T Dibromomethane	7.396	93.0	176038	500.7456	ng	98
T Bromodichloromethane	7.585	83.0	502929	518.3718	ng	100
T cis-1,3-Dichloropropene	8.059	75.0	591147	538.9008	ng	99
T Toluene	8.389	92.0	1095161	536.5101	ng	99
T trans-1,3-Dichloropropene	8.637	75.0	416771	533.7551	ng	99
T 1,1,2-Trichloroethane	8.815	83.0	205463	505.1803	ng	97
T Tetrachloroethene	8.938	163.8	428812	514.9255	ng	100
T 1,3-Dichloropropane	8.980	76.0	408993	511.2479	ng	100
T Chlorodibromomethane	9.203	129.0	330813	520.4361	ng	99
T 1,2-Dibromoethane	9.303	107.0	225877	507.9234	ng	99
T Chlorobenzene	9.802	112.0	1153147	515.9957	ng	100
T 1,1,1,2-Tetrachloroethane	9.892	131.0	406450	520.2855	ng	98
T Ethylbenzene	9.919	91.0	2111152	544.6881	ng	100
T m+p-Xylenes	10.039	106.0	1637879	1087.4082	ng	99
T o-Xylene	10.430	106.0	734101	547.4764	ng	100
T Styrene	10.449	104.0	1199879	555.7946	ng	99
T Bromoform	10.625	172.5	175918	522.7660	ng	98
T Bromobenzene	11.094	156.0	439147	516.0104	ng	98
T 1,1,2,2-Tetrachloroethane	11.113	83.0	240837	491.6700	ng	99
T 1,2,3-Trichloropropane	11.146	110.0	64422	491.5229	ng	99
T 2-Chlorotoluene	11.292	126.0	455991	538.4964	ng	99
T 4-Chlorotoluene	11.400	91.0	1468376	531.8471	ng	100
T 1,3-Dichlorobenzene	12.033	146.0	793993	511.5504	ng	100
T 1,4-Dichlorobenzene	12.125	146.0	794954	502.3001	ng	99
T 1,2-Dichlorobenzene	12.493	146.0	664247	506.3871	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

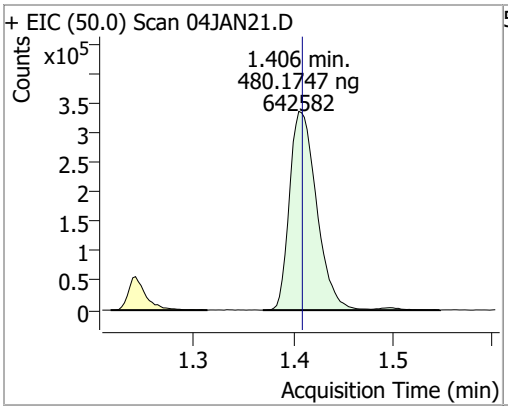
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	494.7474	1.24	0.00	545484	87.0	33.0	2.3	62.3

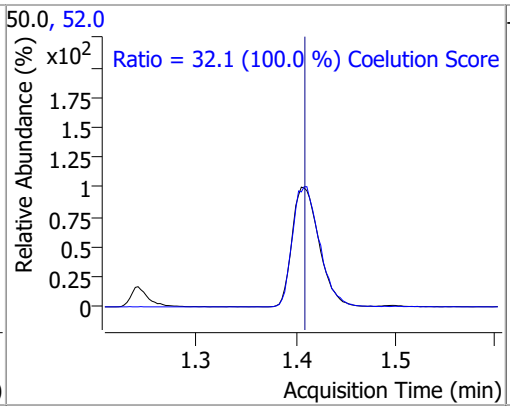


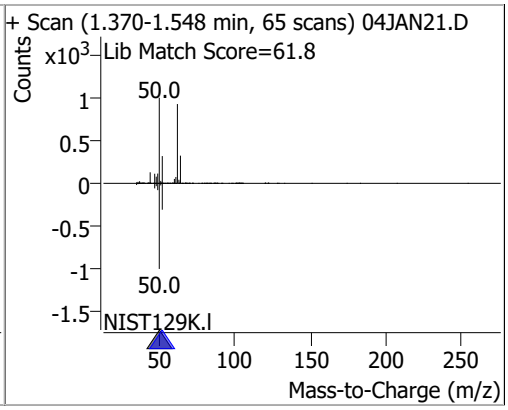




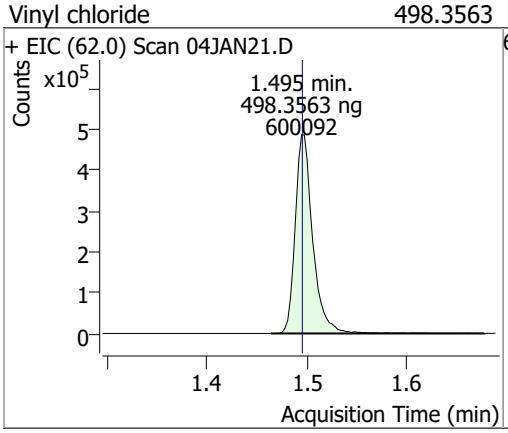
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	480.1747	1.41	0.00	642582	52.0	32.1	2.1	62.1

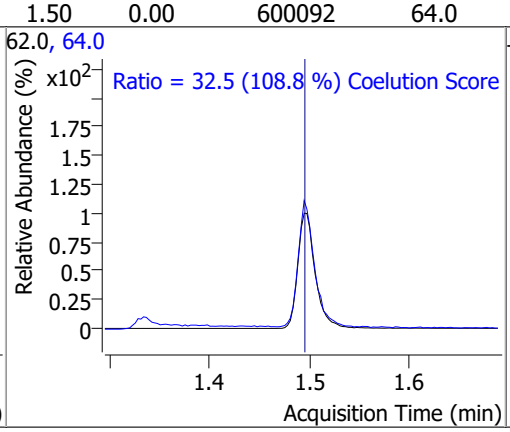


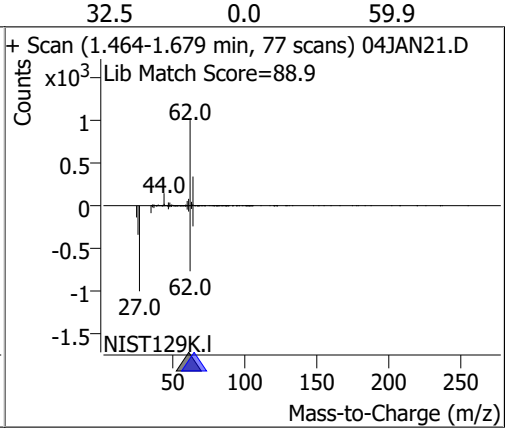




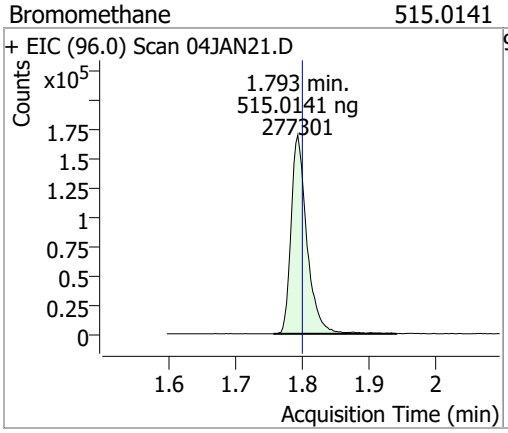
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	498.3563	1.50	0.00	600092	64.0	32.5	0.0	59.9

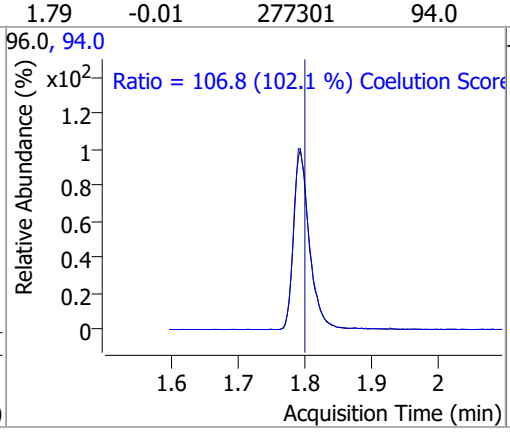


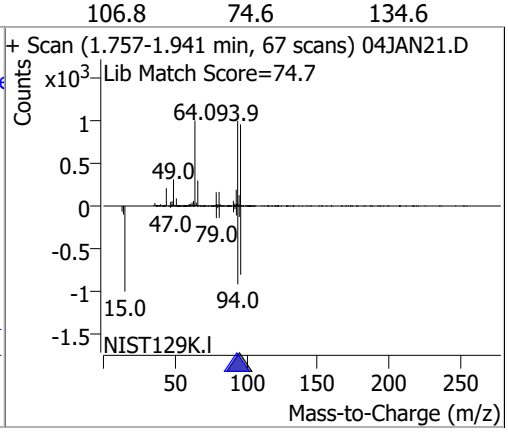




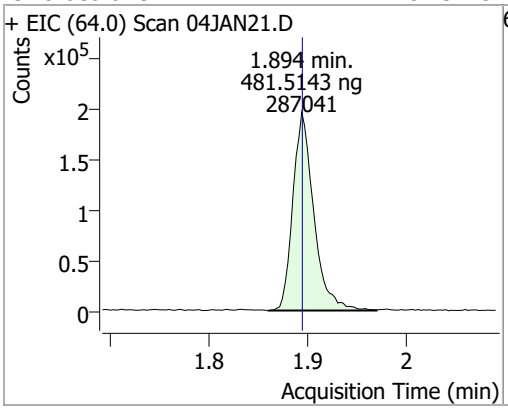
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	515.0141	1.79	-0.01	277301	94.0	106.8	74.6	134.6

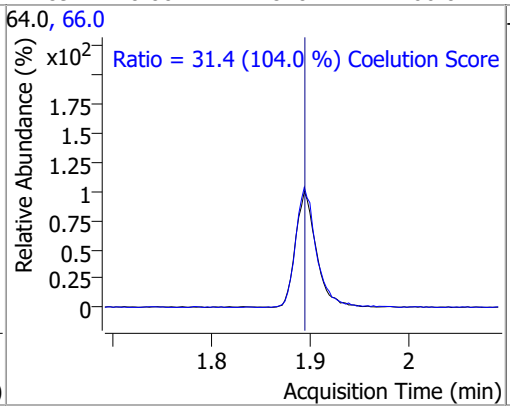


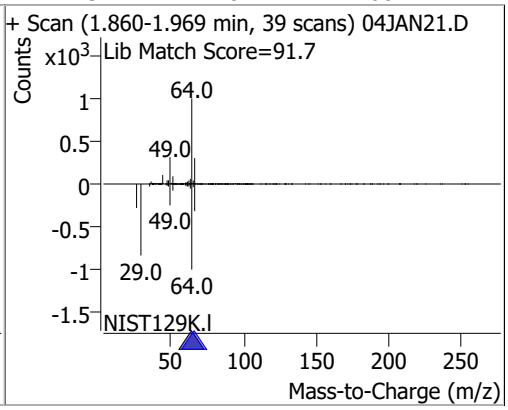




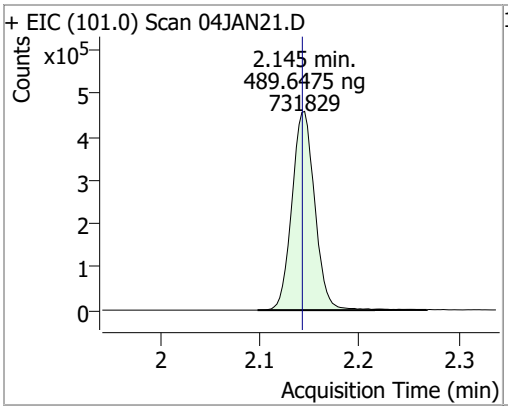
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	481.5143	1.89	0.00	287041	66.0	31.4	0.1	60.1

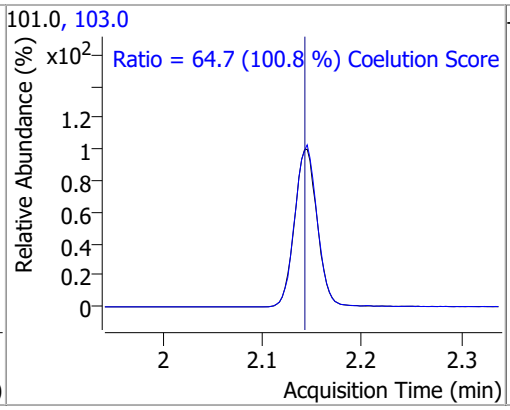


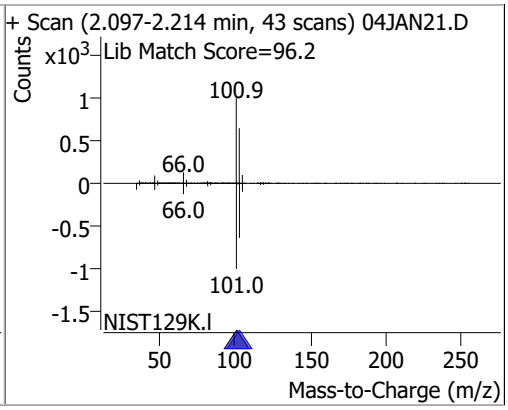




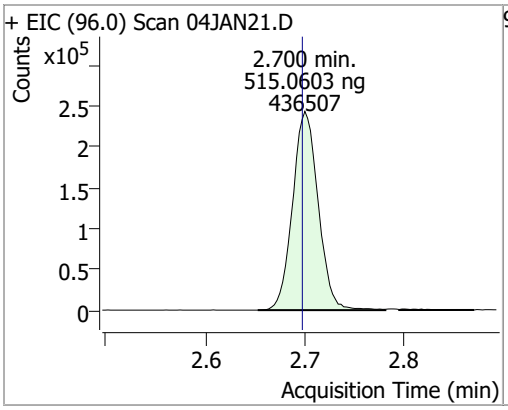
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	489.6475	2.14	0.00	731829	103.0	64.7	34.2	94.2

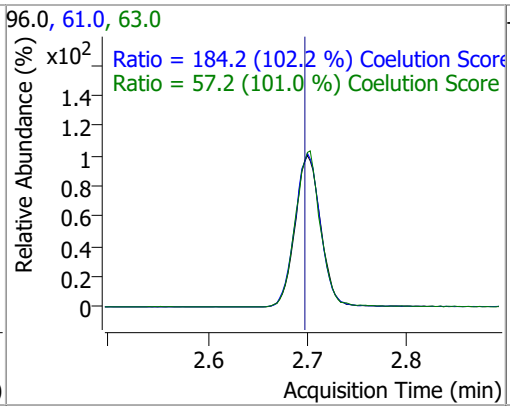


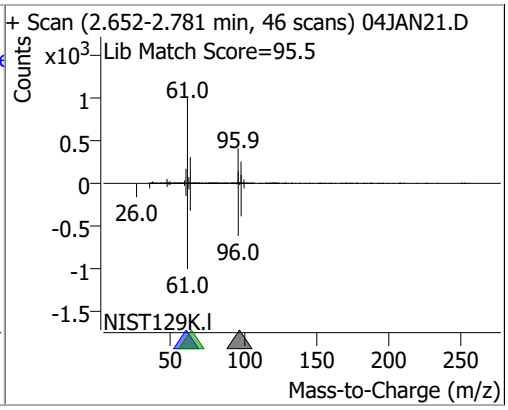




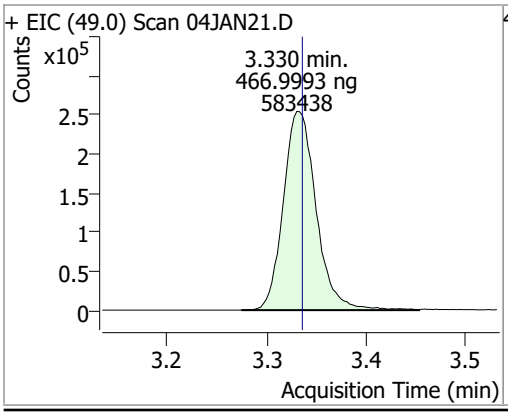
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	515.0603	2.70	0.00	436507	61.0	184.2	150.3	210.3
					63.0	57.2	26.7	86.7

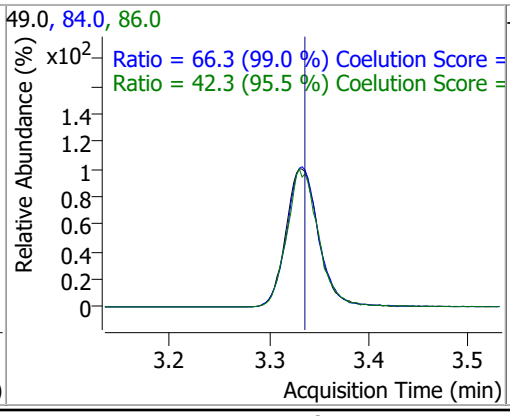


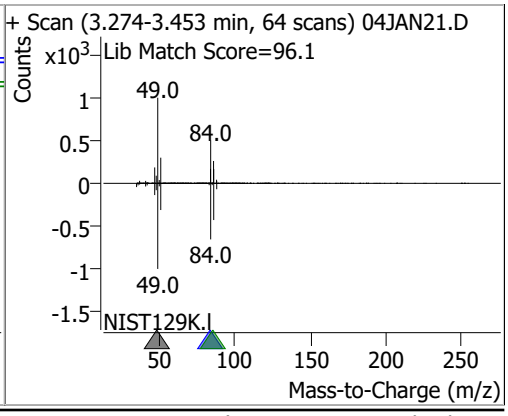




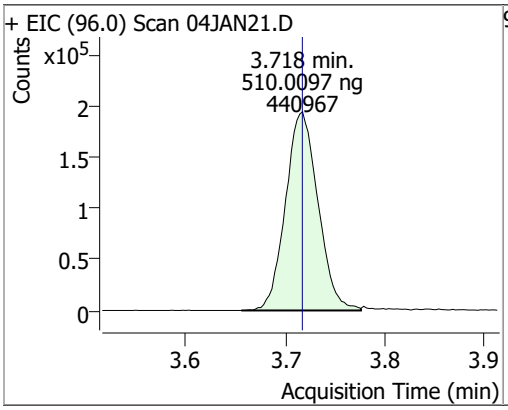
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	466.9993	3.33	-0.01	583438	84.0	66.3	36.9	96.9
					86.0	42.3	14.3	74.3

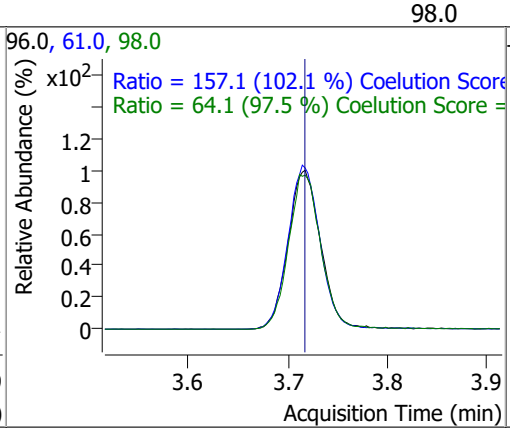


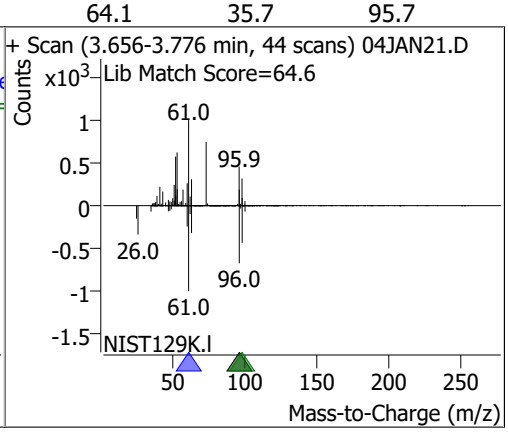




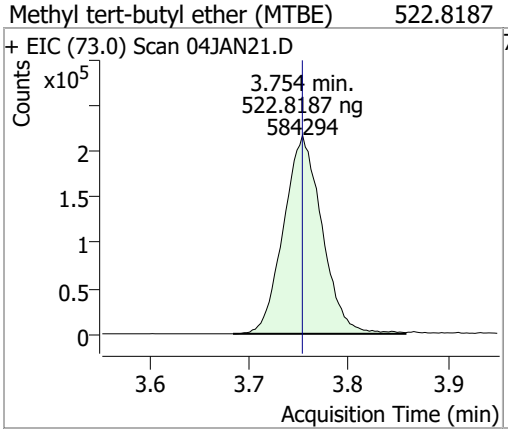
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	510.0097	3.72	0.00	440967	61.0	157.1	123.9	183.9
					98.0	64.1	35.7	95.7

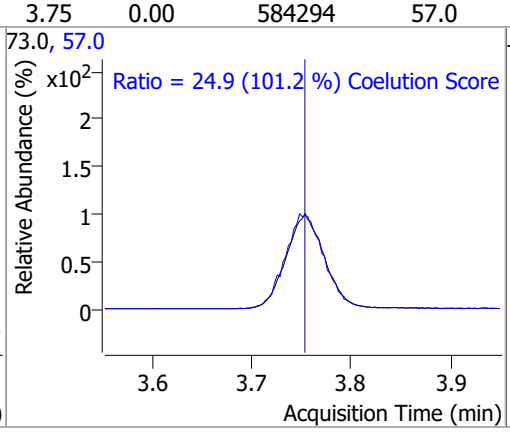


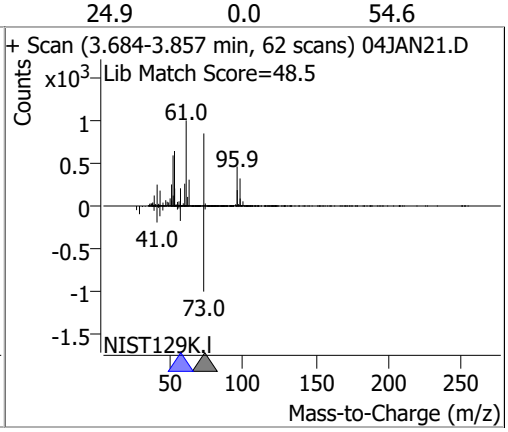




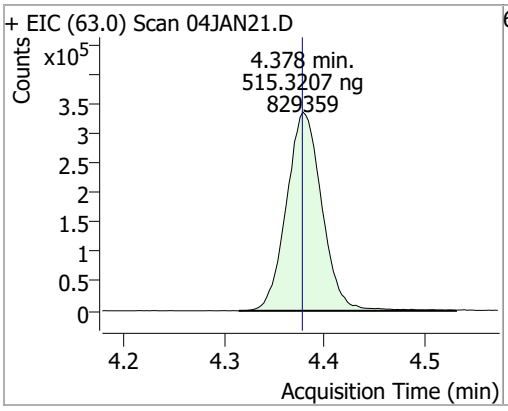
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	522.8187	3.75	0.00	584294	57.0	24.9	0.0	54.6

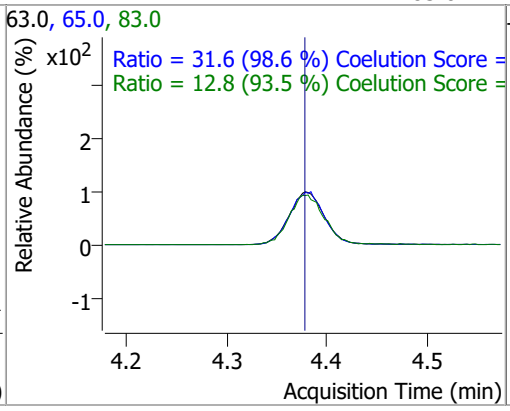


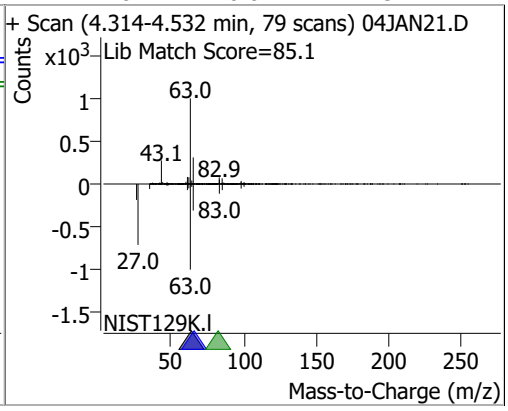




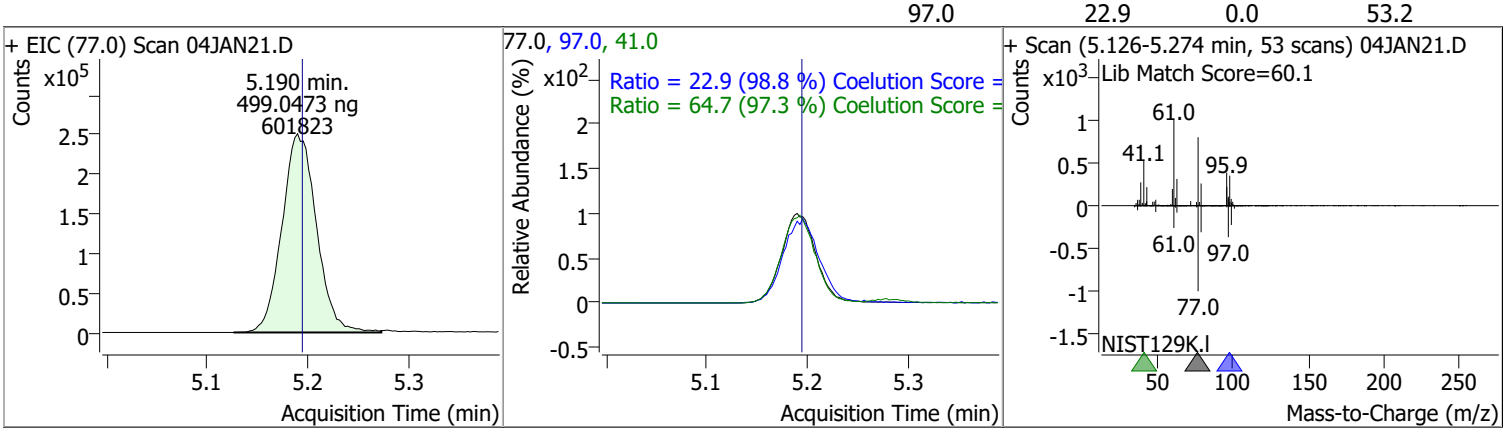
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	515.3207	4.38	0.00	829359	65.0	31.6	2.1	62.1
					83.0	12.8	0.0	43.7



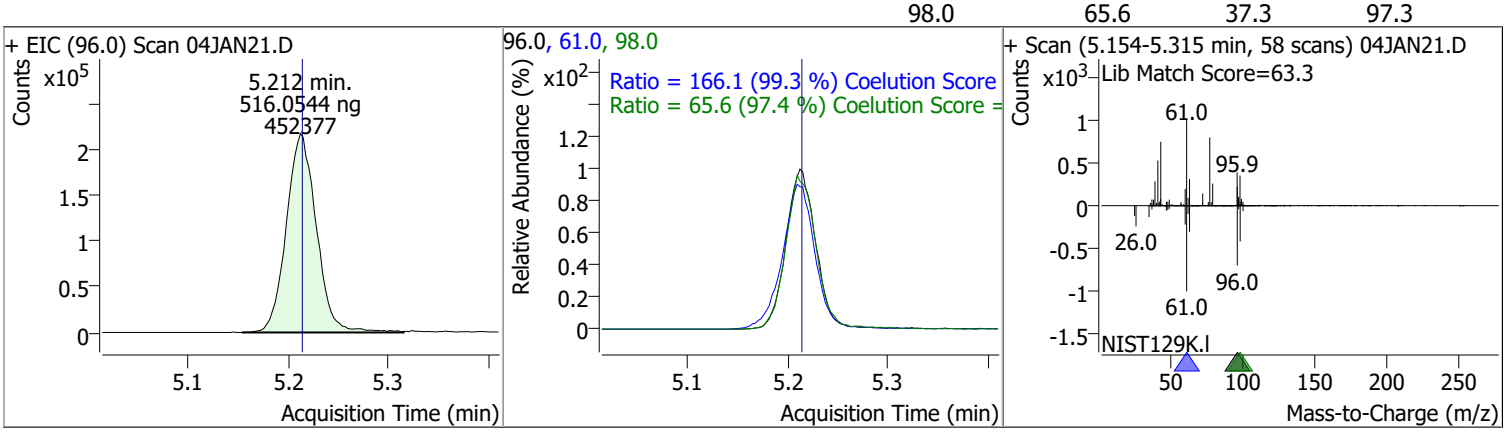




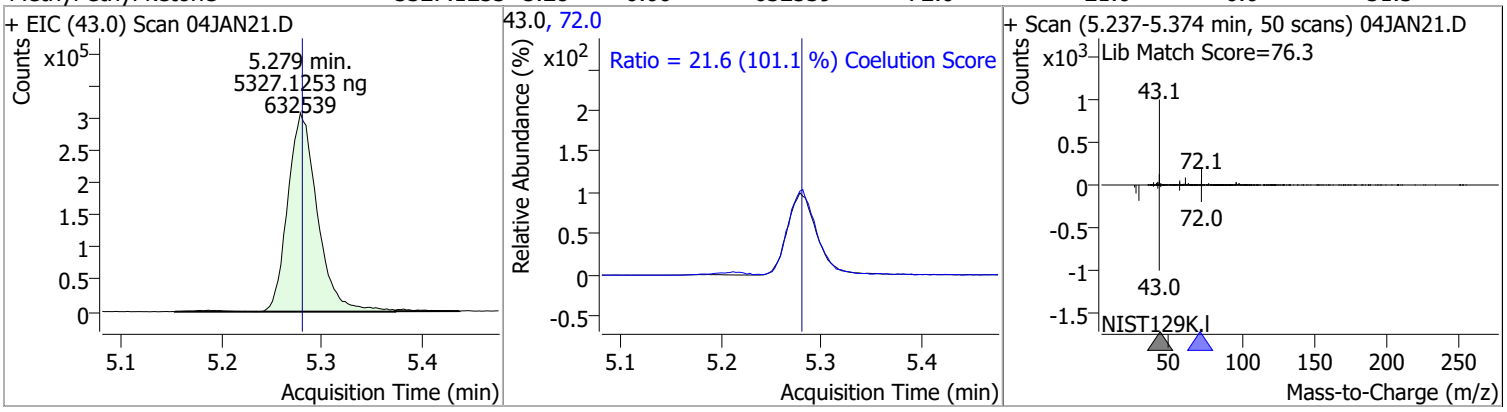
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	499.0473	5.19	-0.01	601823	41.0	64.7	36.5	96.5
					97.0	22.9	0.0	53.2



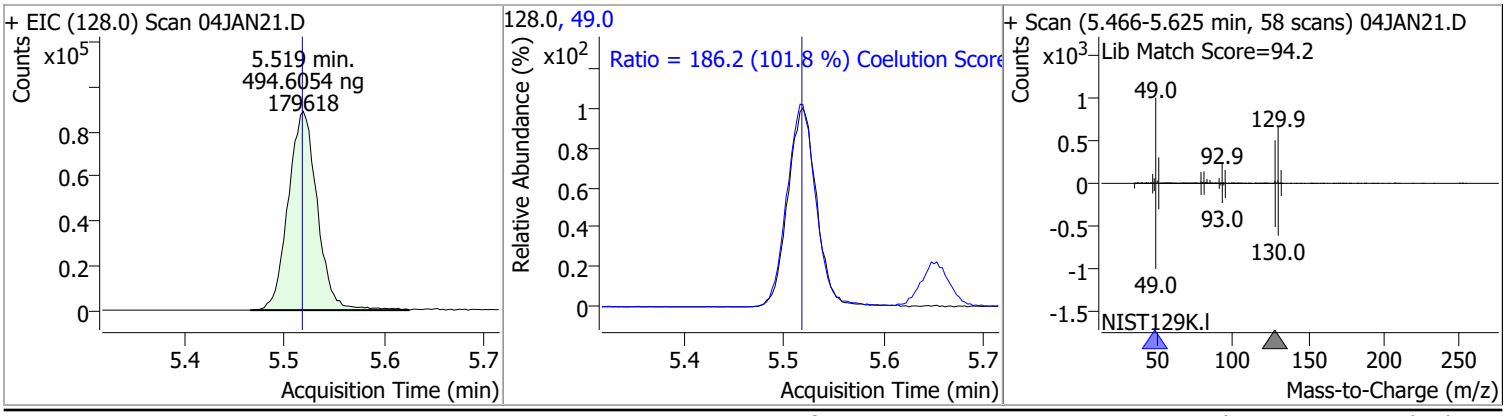
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	516.0544	5.21	0.00	452377	61.0	166.1	137.2	197.2
					98.0	65.6	37.3	97.3



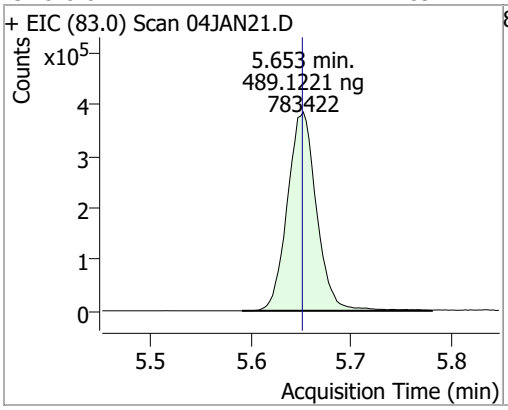
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	5327.1253	5.28	0.00	632539	72.0	21.6	0.0	51.3

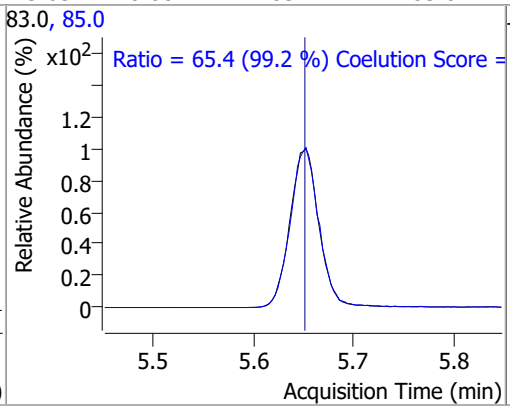


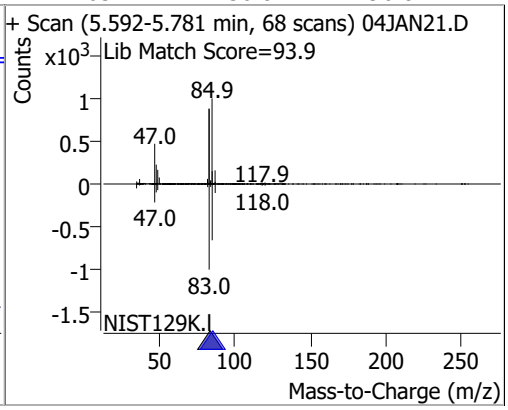
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	494.6054	5.52	0.00	179618	49.0	186.2	152.9	212.9



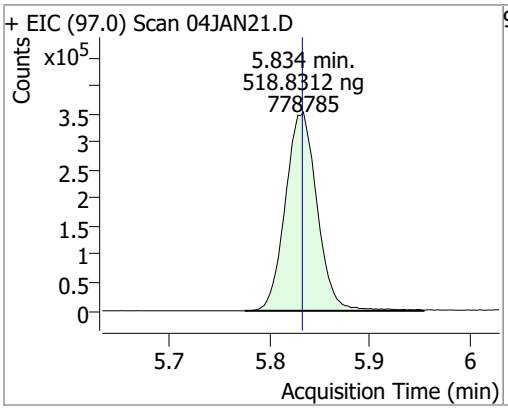
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	489.1221	5.65	0.00	783422	85.0	65.4	36.0	96.0

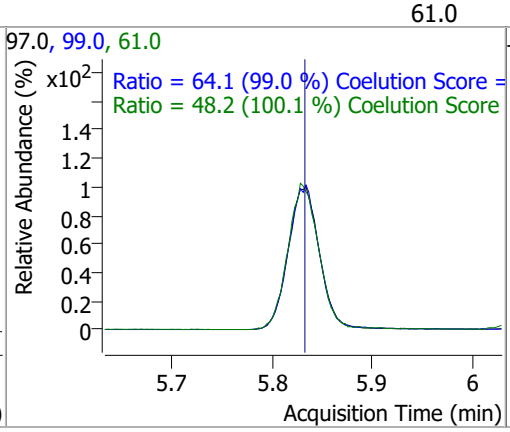


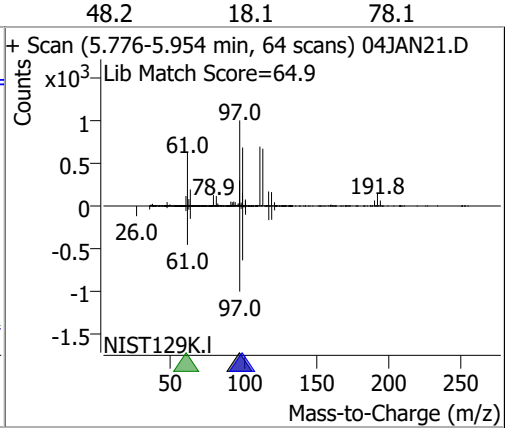




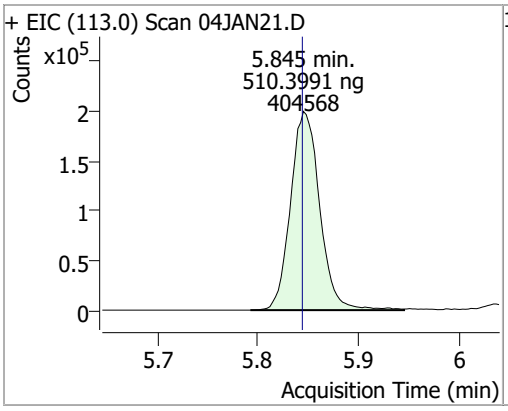
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	518.8312	5.83	0.00	778785	99.0	64.1	34.7	94.7
					61.0	48.2	18.1	78.1

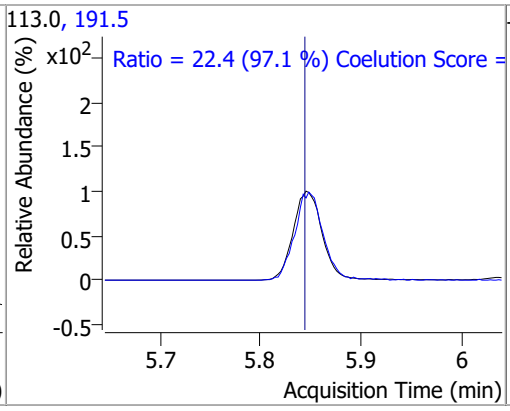


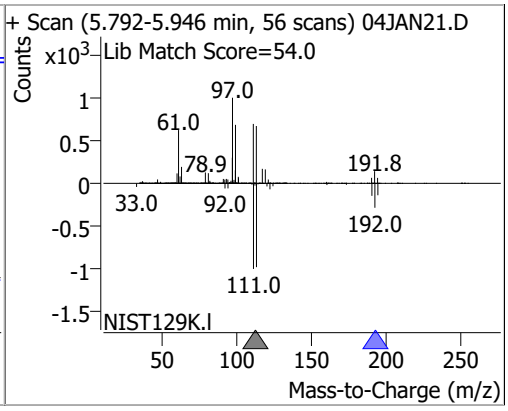




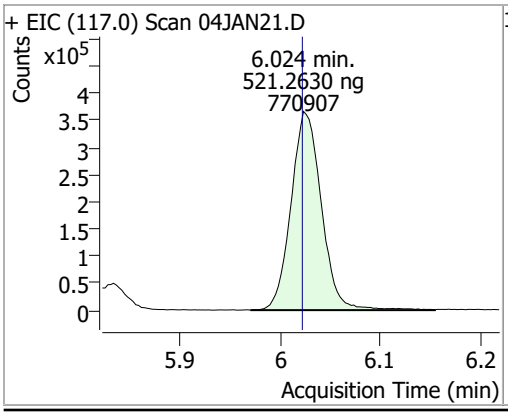
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	510.3991	5.85	0.00	404568	191.5	22.4	0.0	53.1

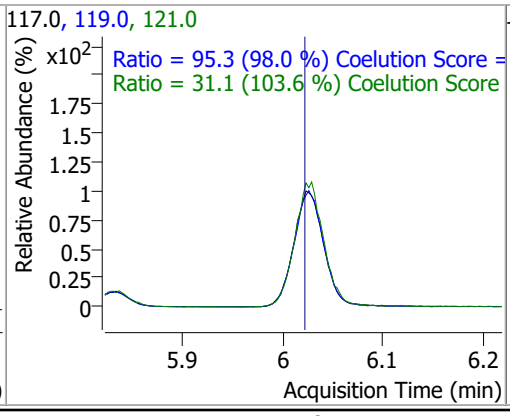


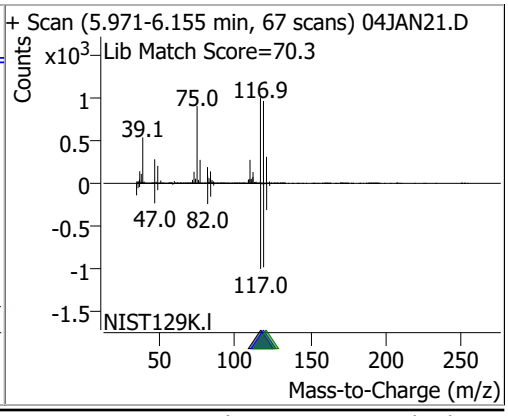




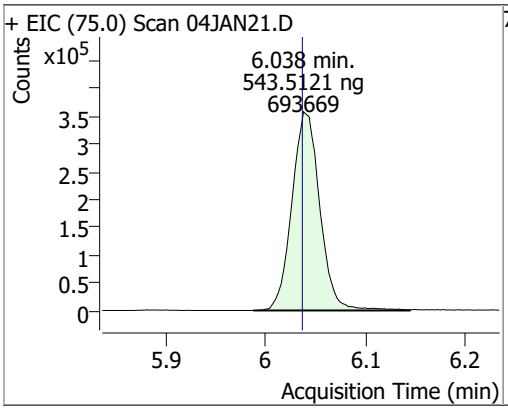
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	521.2630	6.02	0.00	770907	119.0	95.3	67.2	127.2
					121.0	31.1	0.1	60.1

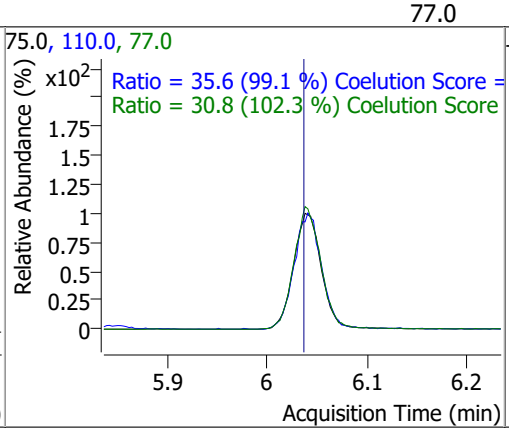


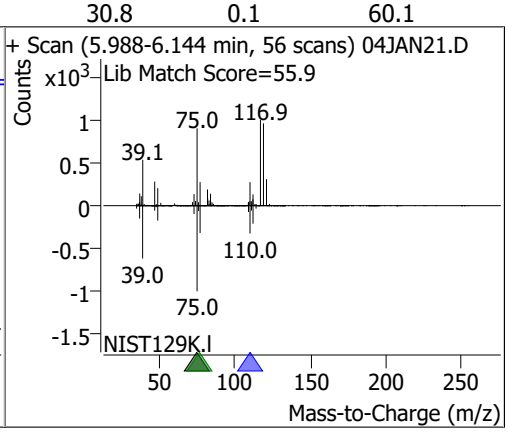




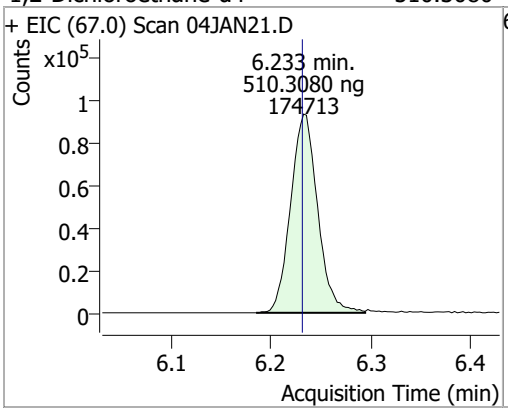
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	543.5121	6.04	0.00	693669	110.0	35.6	5.9	65.9
					77.0	30.8	0.1	60.1

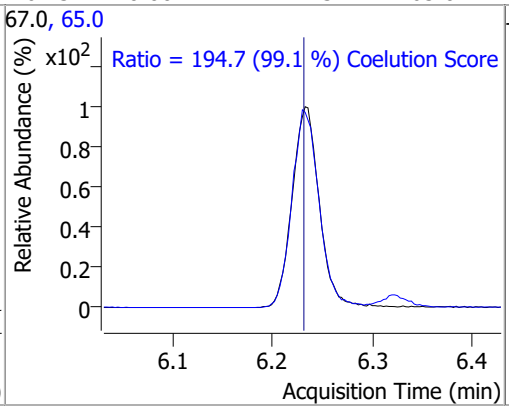


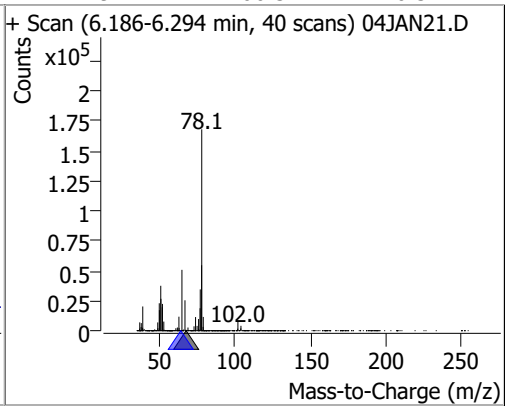




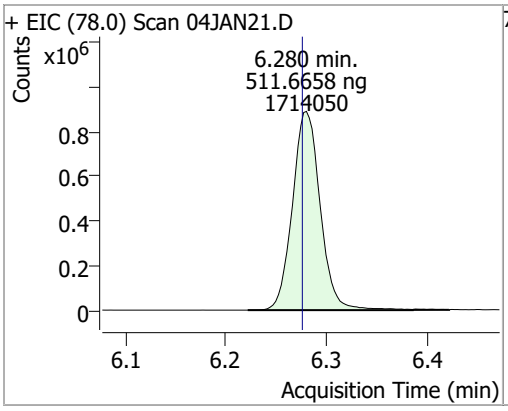
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	510.3080	6.23	0.00	174713	65.0	194.7	166.5	226.5

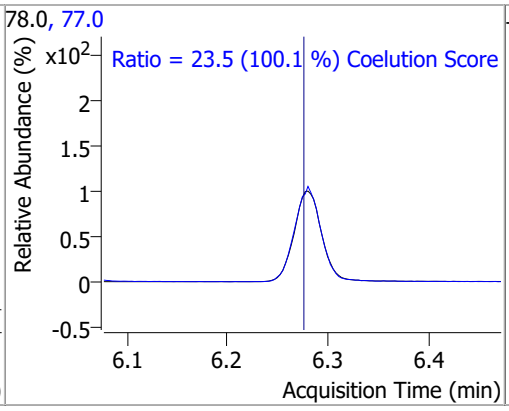


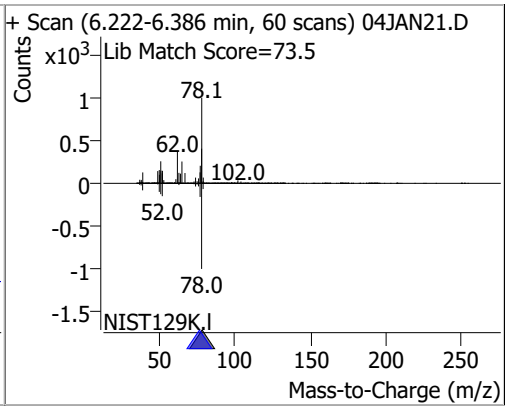




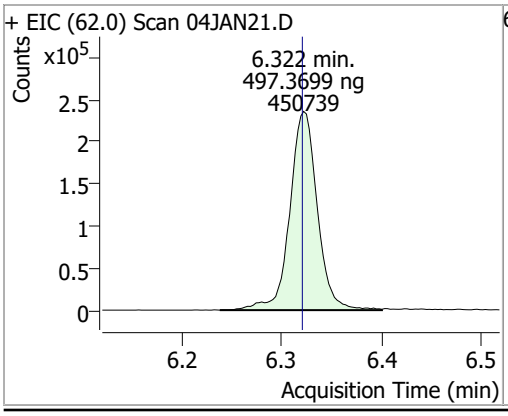
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	511.6658	6.28	0.00	1714050	77.0	23.5	0.0	53.5

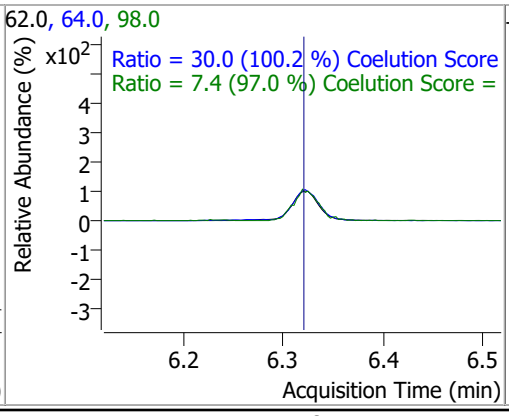


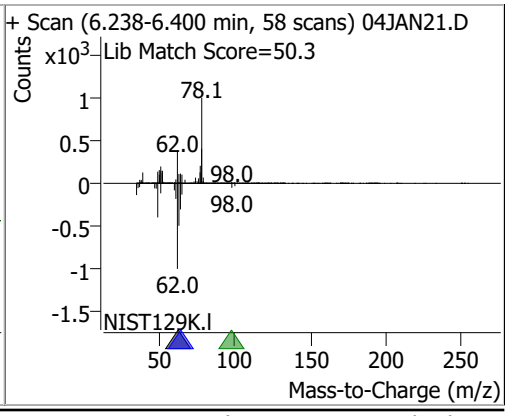




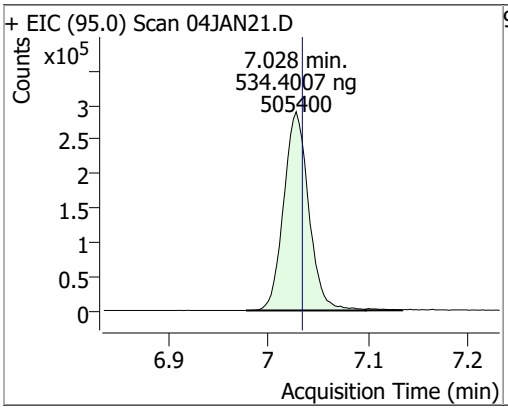
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	497.3699	6.32	0.00	450739	64.0	30.0	0.0	59.9
					98.0	7.4	0.0	37.6

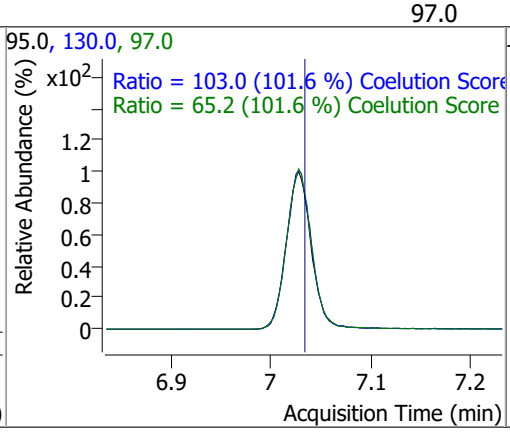


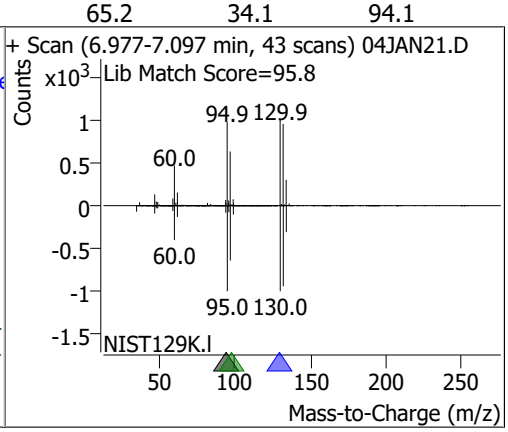




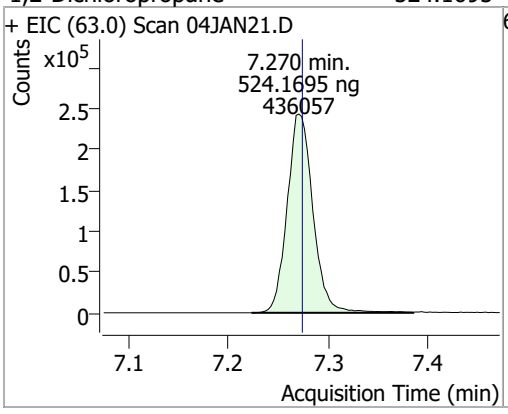
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	534.4007	7.03	0.00	505400	130.0	103.0	71.5	131.5
					97.0	65.2	34.1	94.1

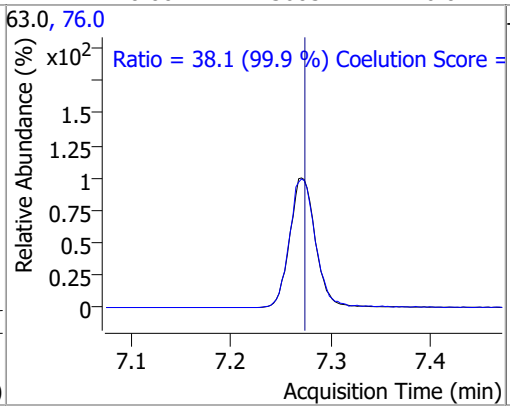


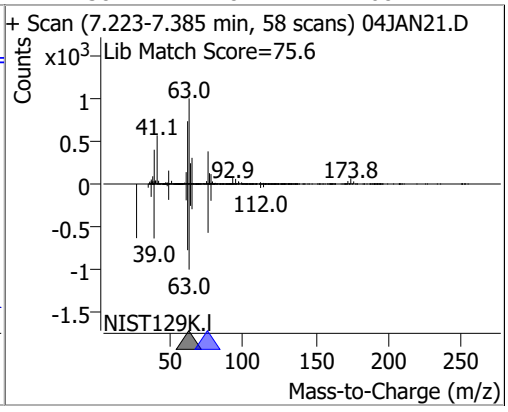




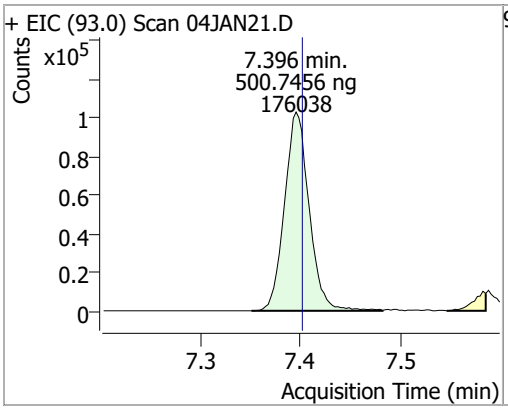
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	524.1695	7.27	0.00	436057	76.0	38.1	8.2	68.2

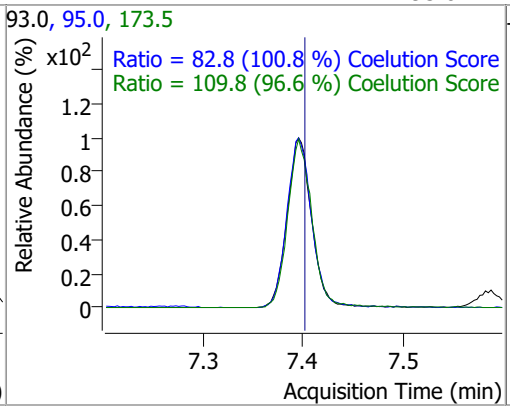


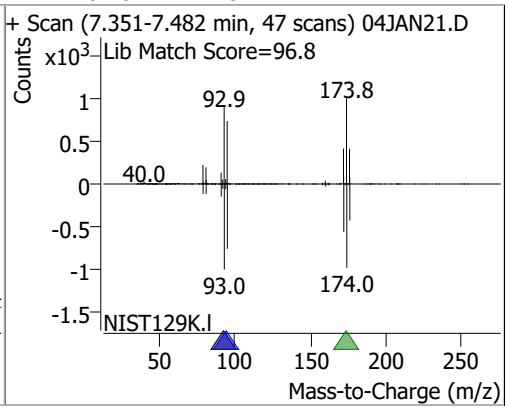




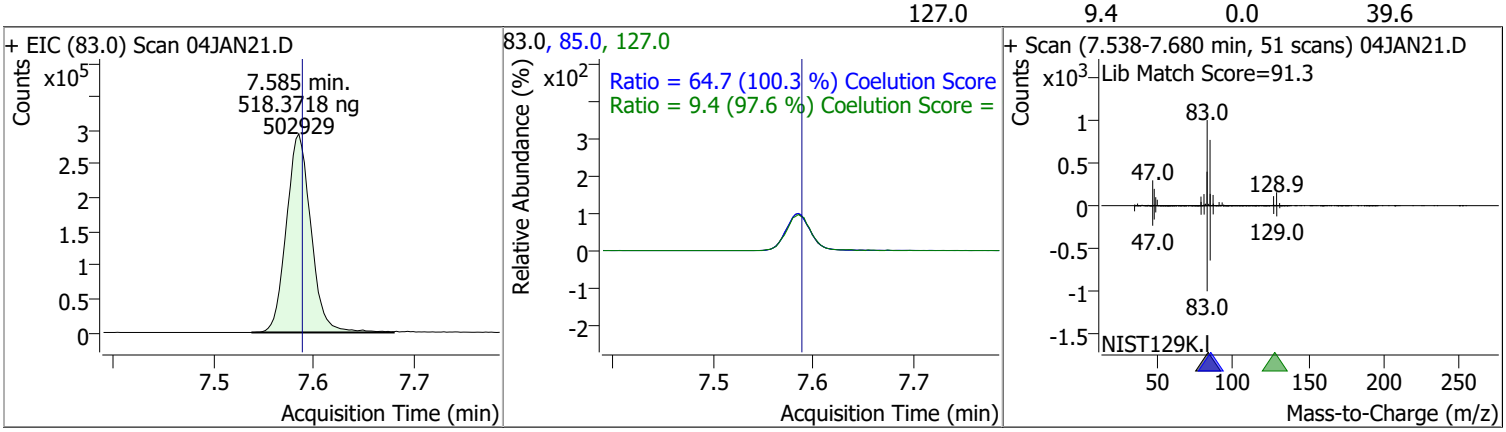
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	500.7456	7.40	0.00	176038	173.5	109.8	83.7	143.7
					95.0	82.8	52.2	112.2



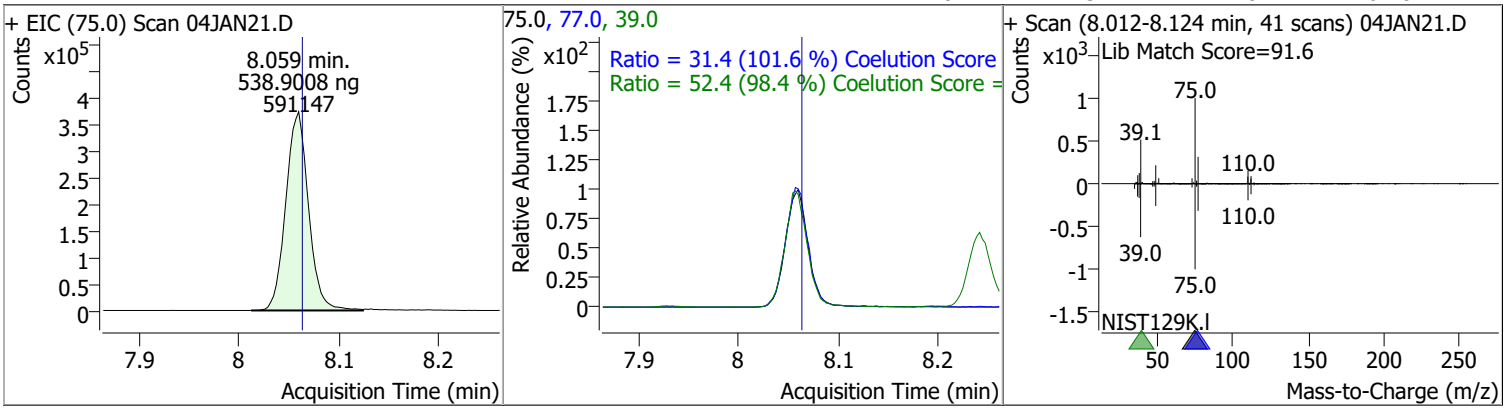




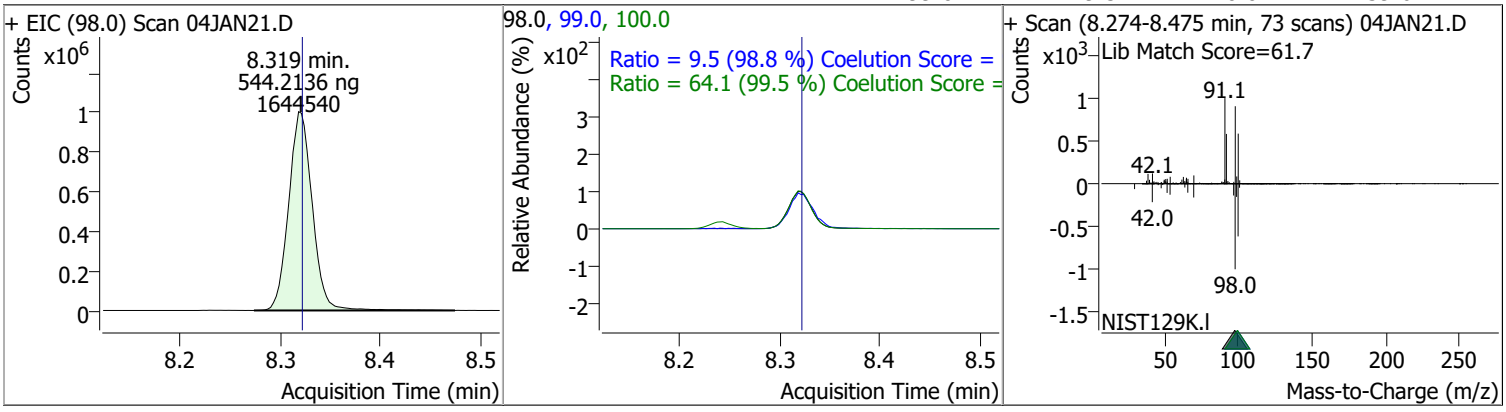
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	518.3718	7.59	0.00	502929	85.0	64.7	34.5	94.5
					127.0	9.4	0.0	39.6



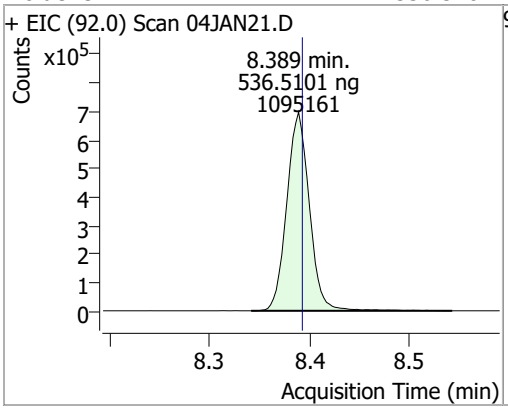
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	538.9008	8.06	0.00	591147	39.0	52.4	23.3	83.3
					77.0	31.4	1.0	61.0

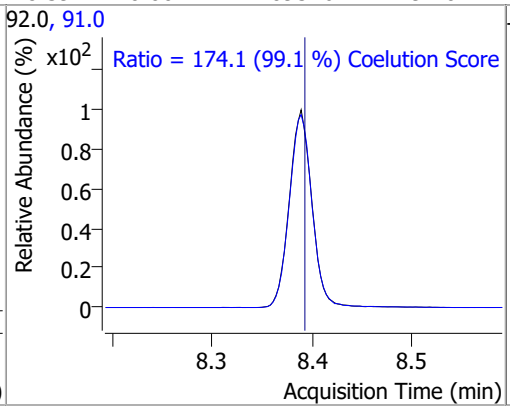


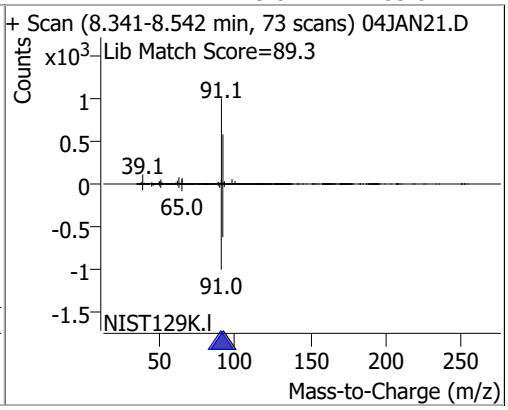
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	544.2136	8.32	0.00	1644540	100.0	64.1	34.4	94.4
					99.0	9.5	0.0	39.6



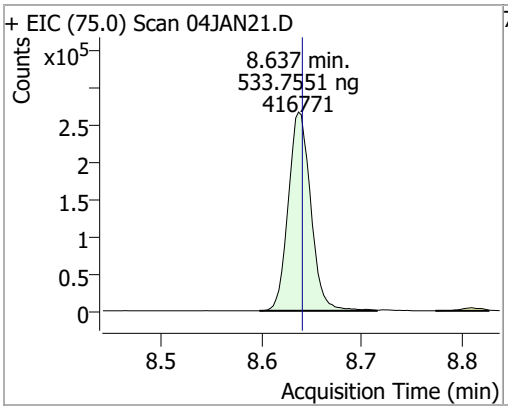
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	536.5101	8.39	0.00	1095161	91.0	174.1	145.8	205.8

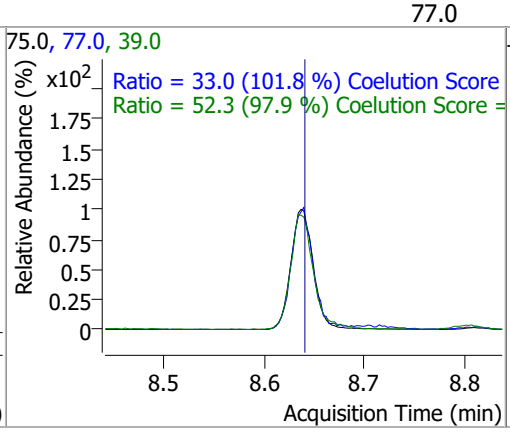


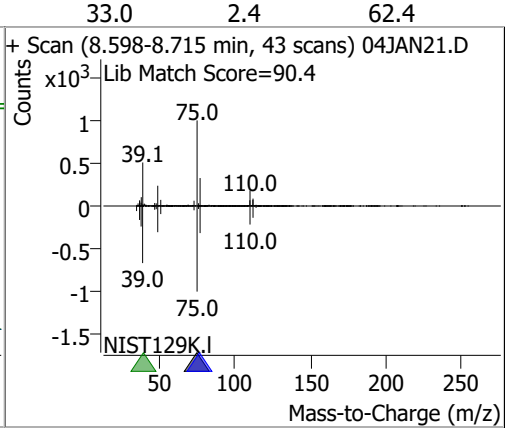




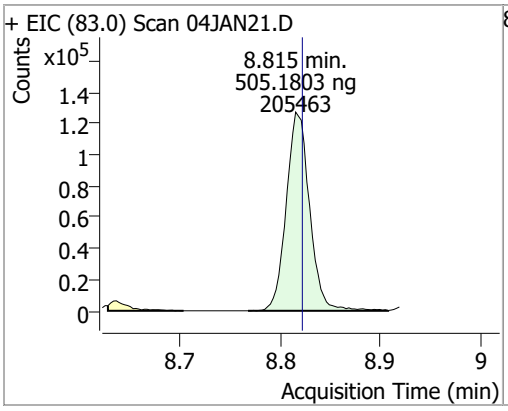
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	533.7551	8.64	0.00	416771	39.0	52.3	23.4	83.4
					77.0	33.0	2.4	62.4

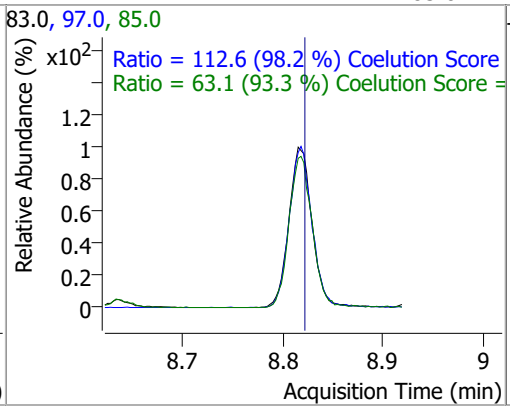


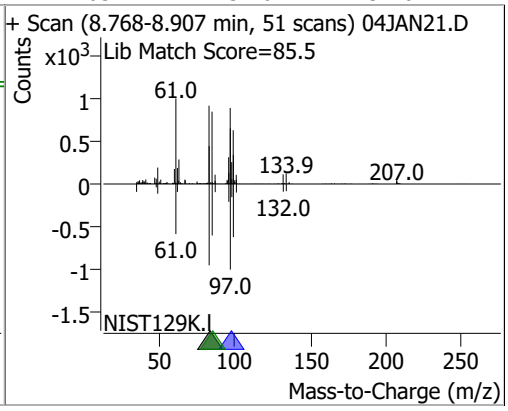




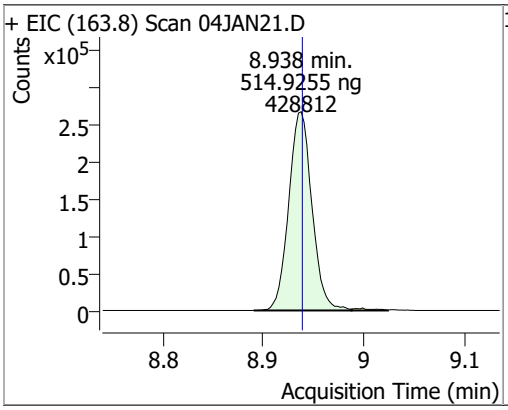
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	505.1803	8.82	0.00	205463	97.0	112.6	84.6	144.6
					85.0	63.1	37.6	97.6

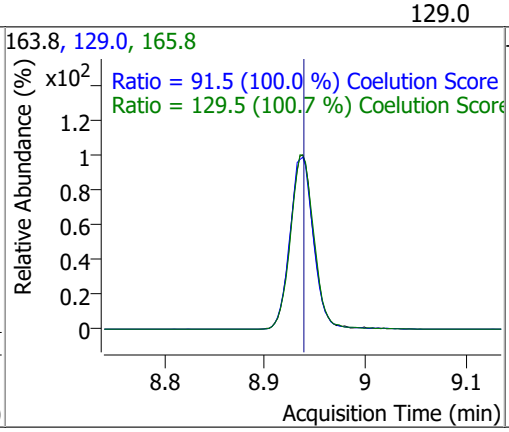


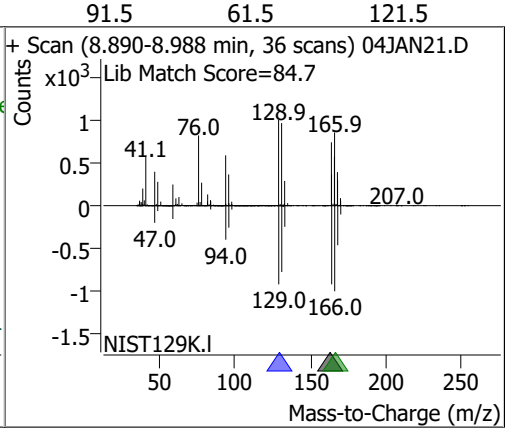




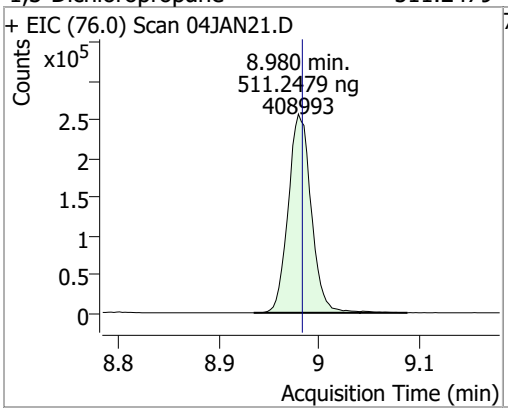
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	514.9255	8.94	0.00	428812	165.8	129.5	98.6	158.6
					129.0	91.5	61.5	121.5

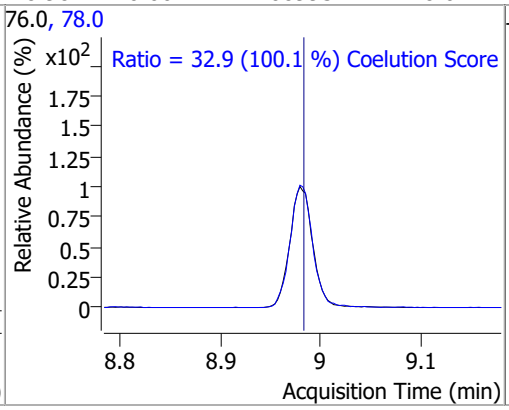


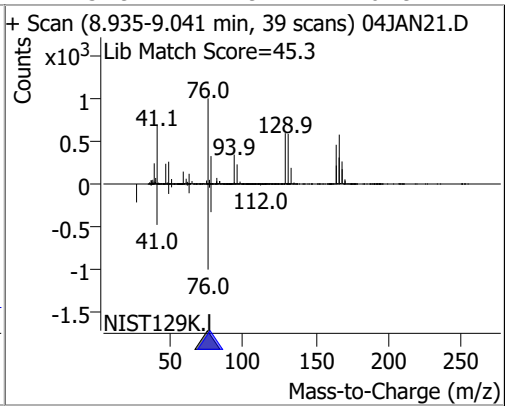




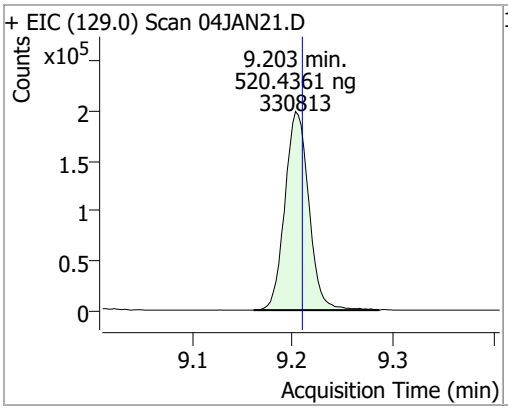
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	511.2479	8.98	0.00	408993	78.0	32.9	2.9	62.9

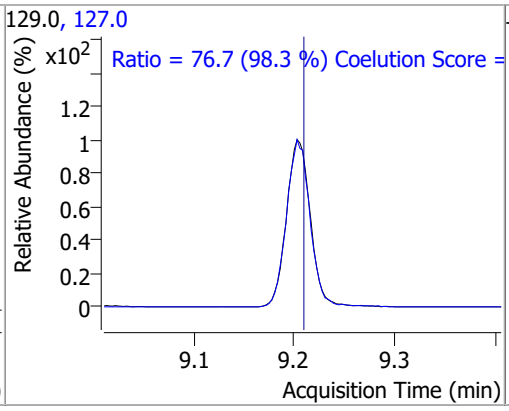


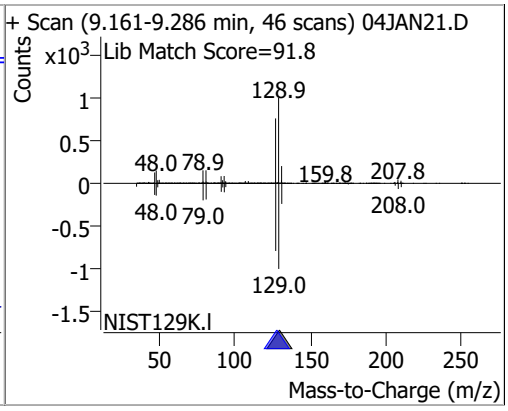




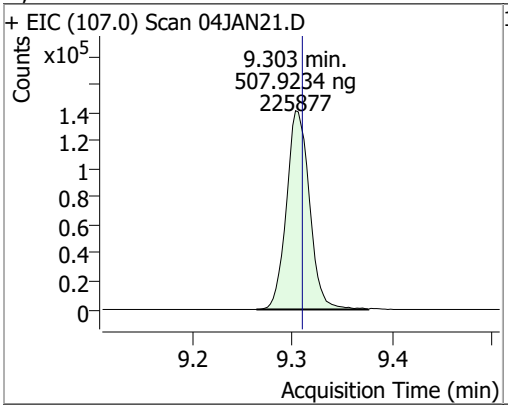
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	520.4361	9.20	0.00	330813	127.0	76.7	48.0	108.0

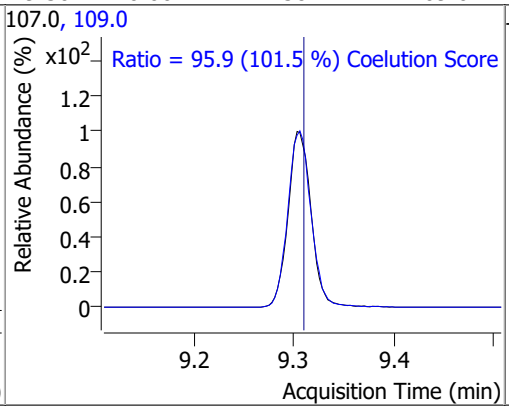


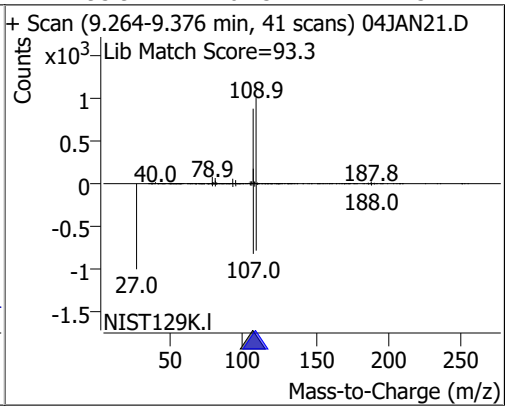




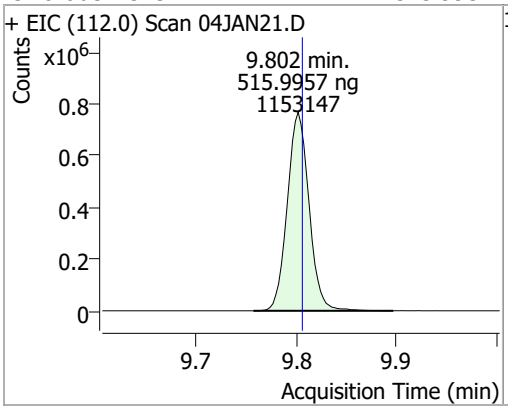
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	507.9234	9.30	0.00	225877	109.0	95.9	64.5	124.5

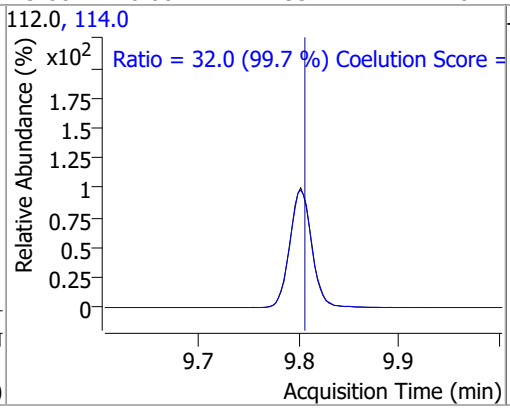


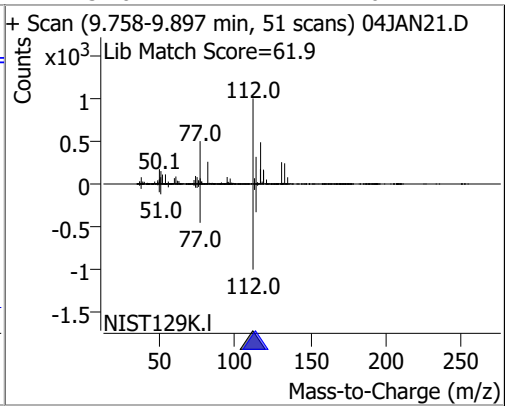




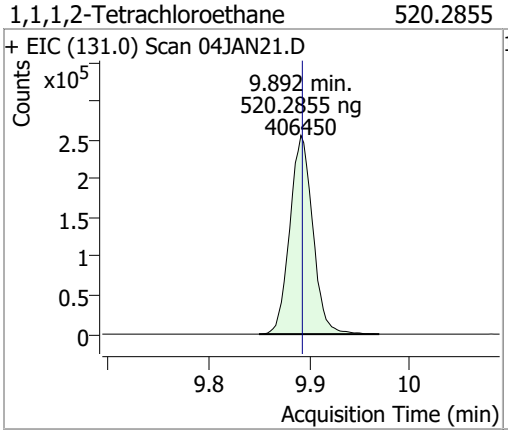
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	515.9957	9.80	0.00	1153147	114.0	32.0	2.1	62.1

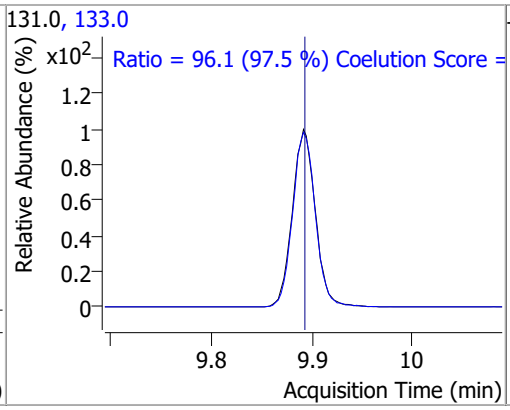


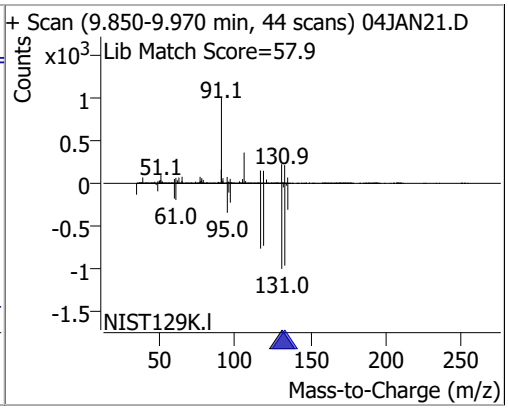




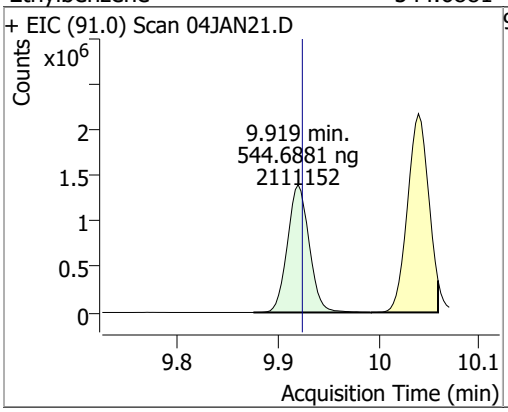
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	520.2855	9.89	0.00	406450	133.0	96.1	68.6	128.6

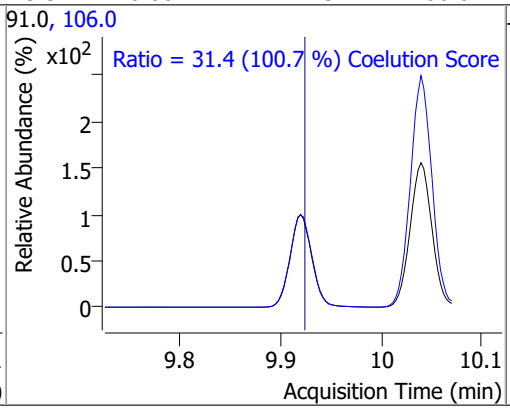


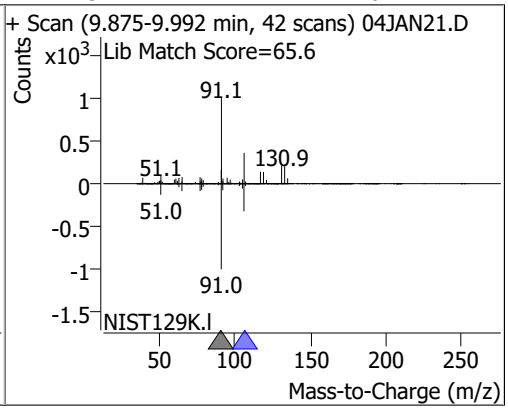




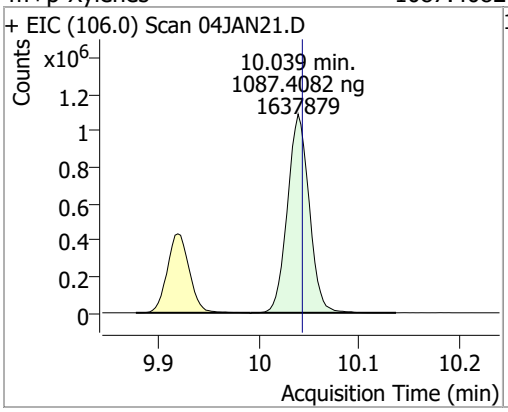
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	544.6881	9.92	0.00	2111152	106.0	31.4	1.1	61.1

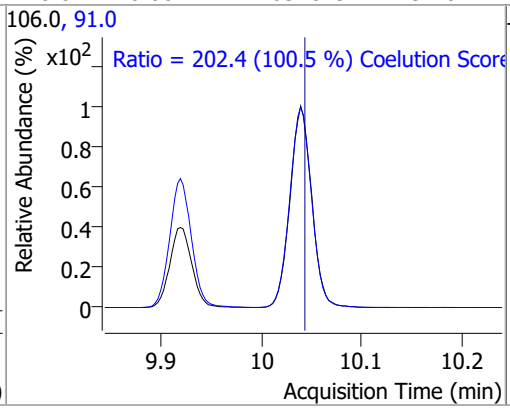


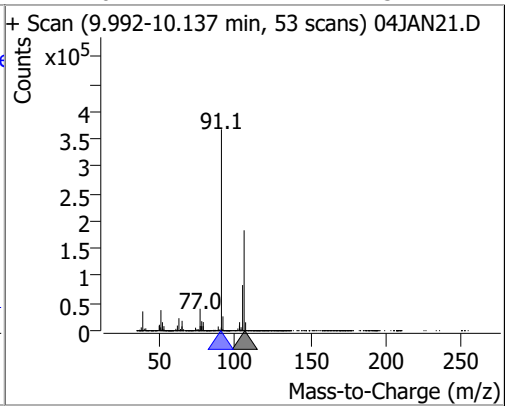




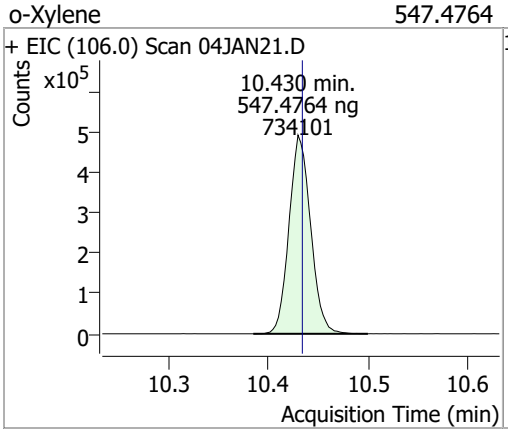
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	1087.4082	10.04	0.00	1637879	91.0	202.4	171.4	231.4

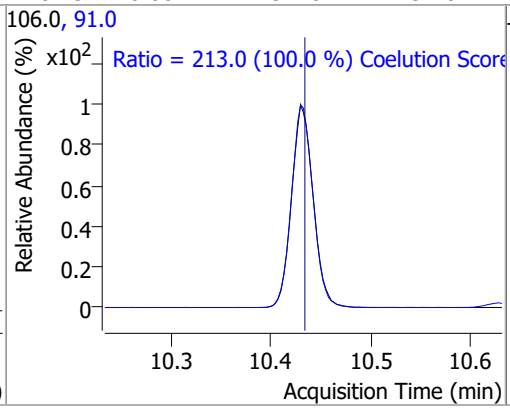


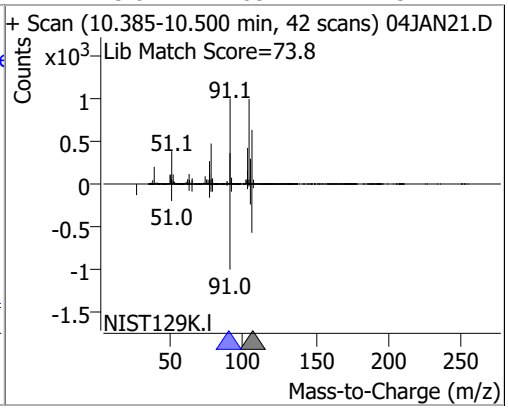




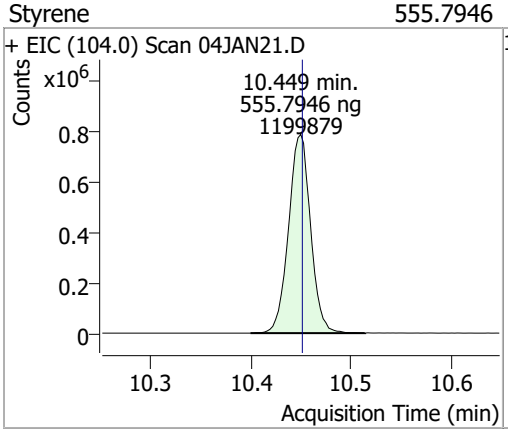
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	547.4764	10.43	0.00	734101	91.0	213.0	183.1	243.1

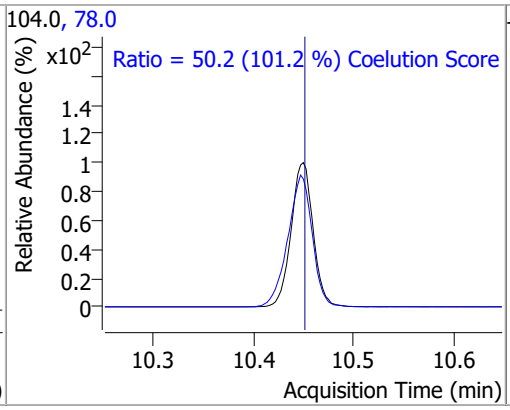


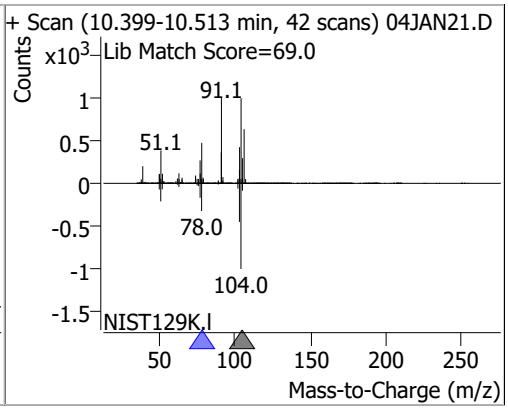




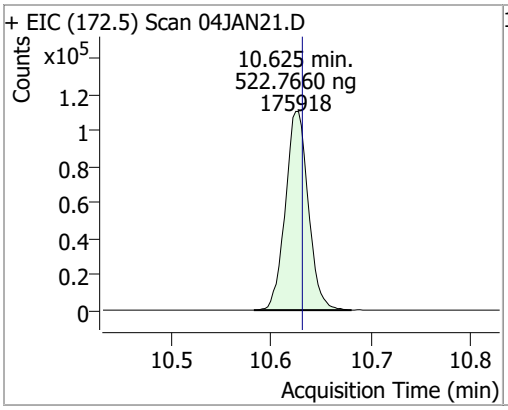
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	555.7946	10.45	0.00	1199879	78.0	50.2	19.6	79.6

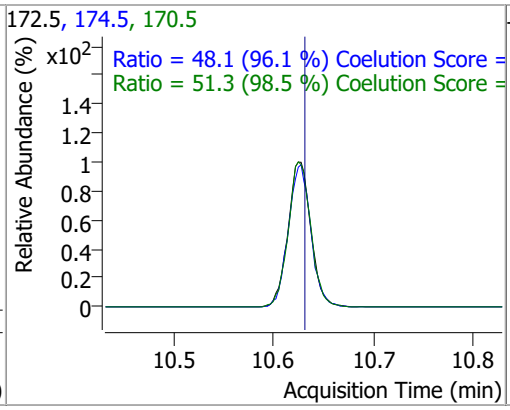


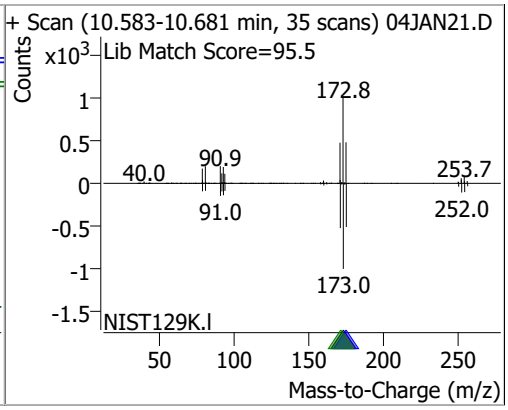




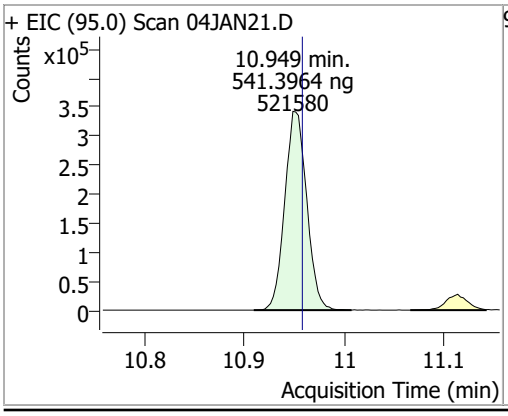
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	522.7660	10.63	0.00	175918	170.5	51.3	22.1	82.1
					174.5	48.1	20.1	80.1

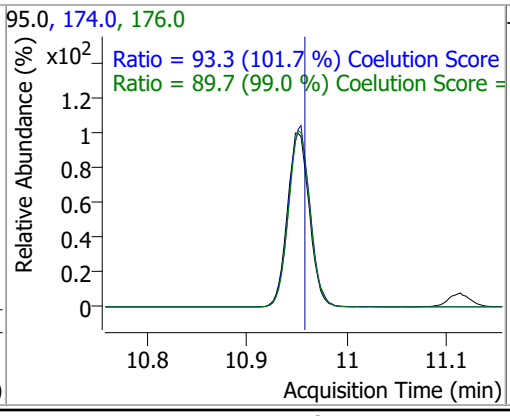


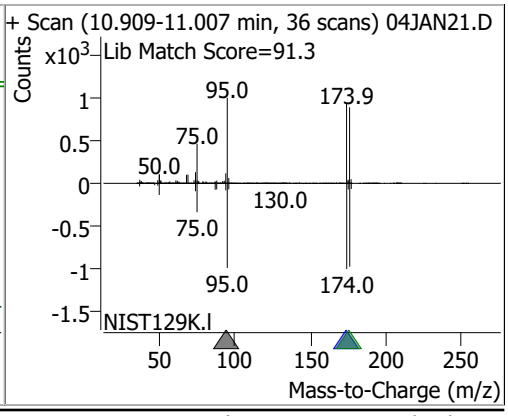




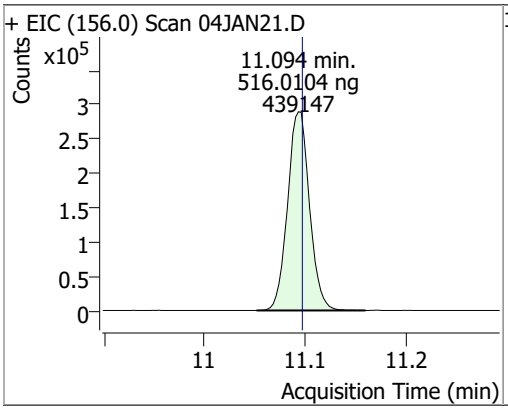
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	541.3964	10.95	-0.01	521580	174.0	93.3	61.7	121.7
					176.0	89.7	60.6	120.6

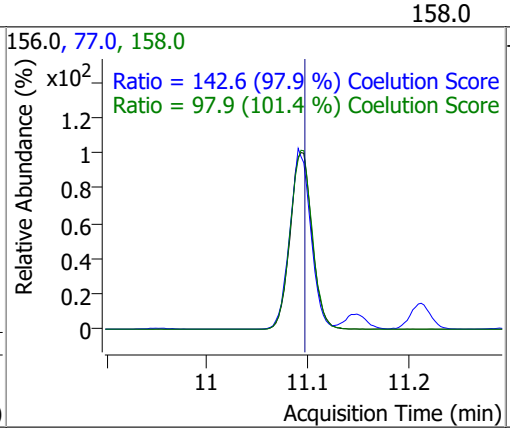


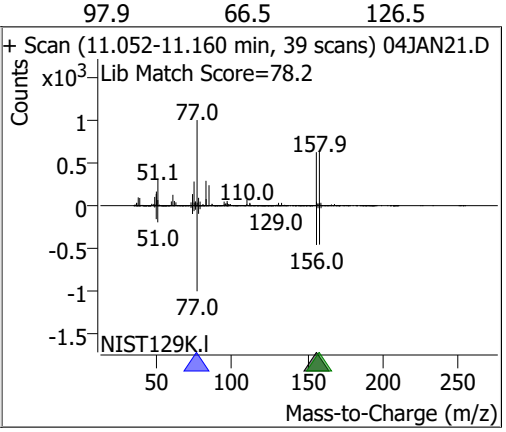




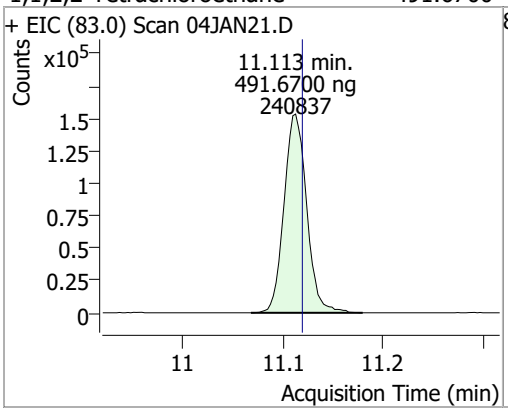
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	516.0104	11.09	0.00	439147	77.0	142.6	115.7	175.7
					158.0	97.9	66.5	126.5

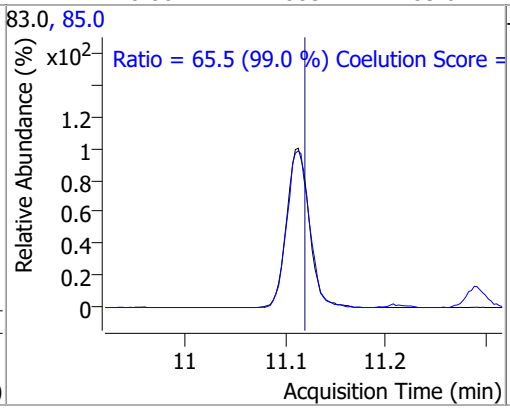


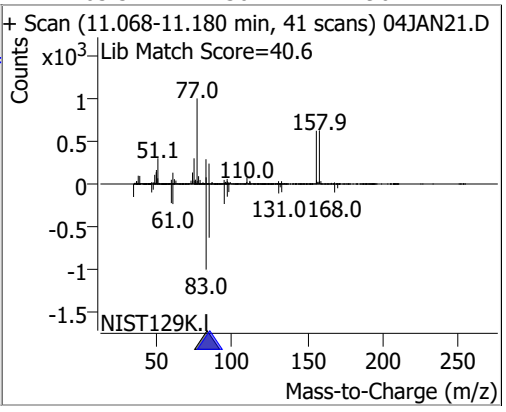




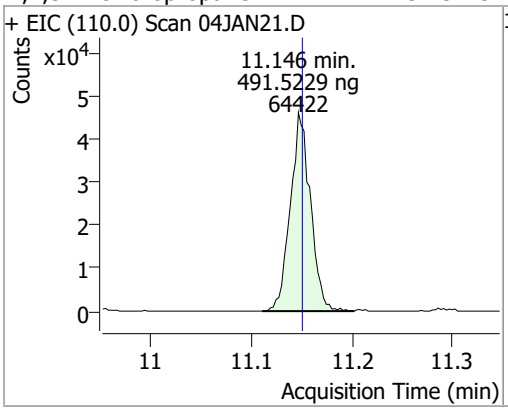
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	491.6700	11.11	0.00	240837	85.0	65.5	36.2	96.2

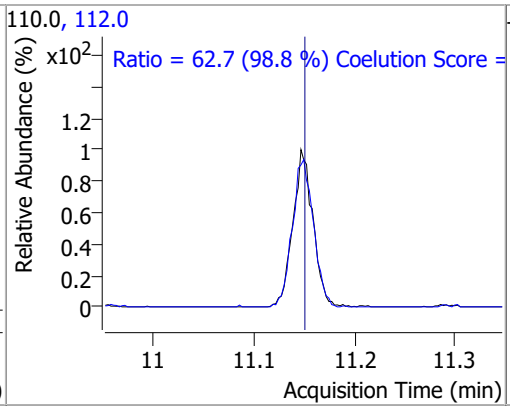


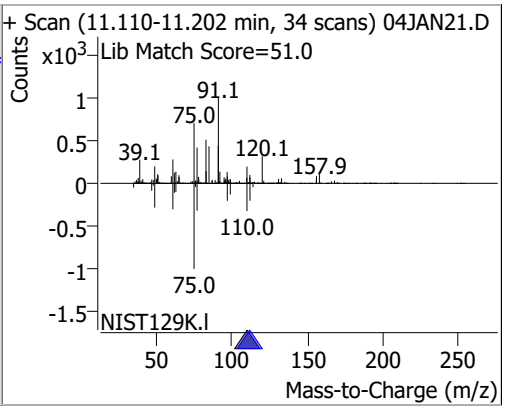




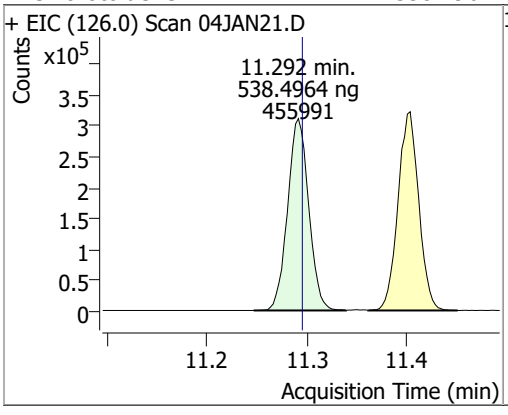
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	491.5229	11.15	0.00	64422	112.0	62.7	33.5	93.5

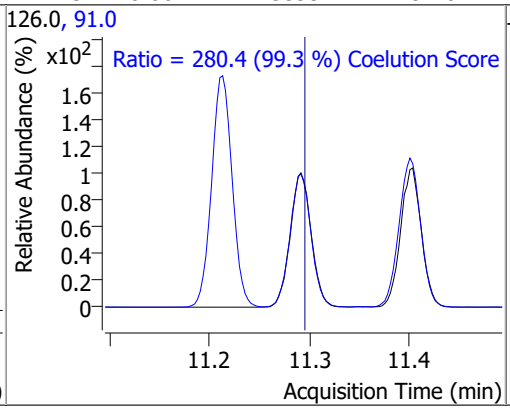


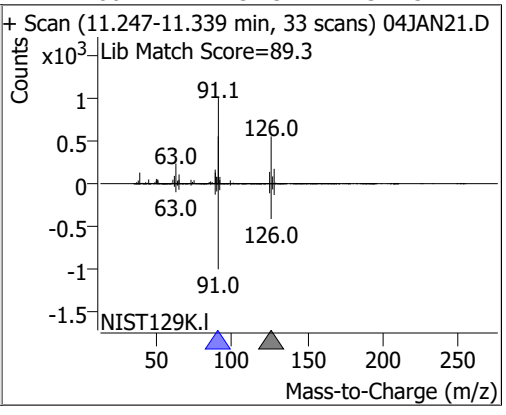




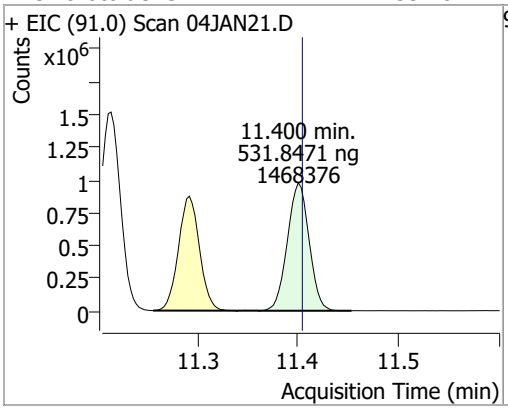
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	538.4964	11.29	0.00	455991	91.0	280.4	252.3	312.3

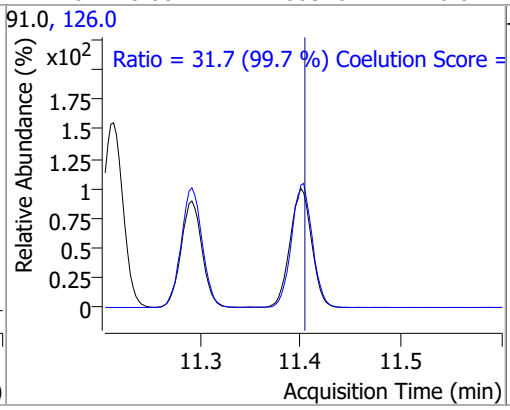


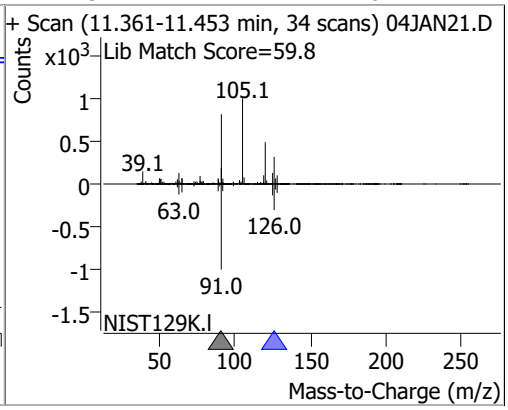




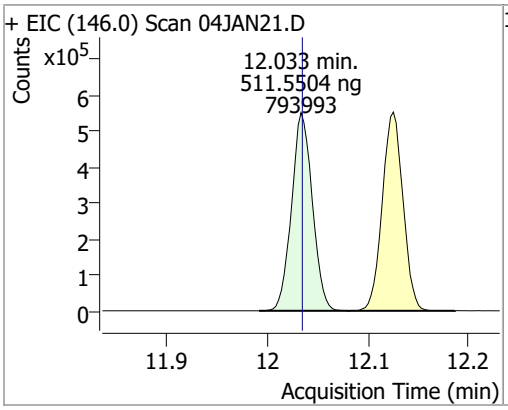
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	531.8471	11.40	0.00	1468376	126.0	31.7	1.7	61.7

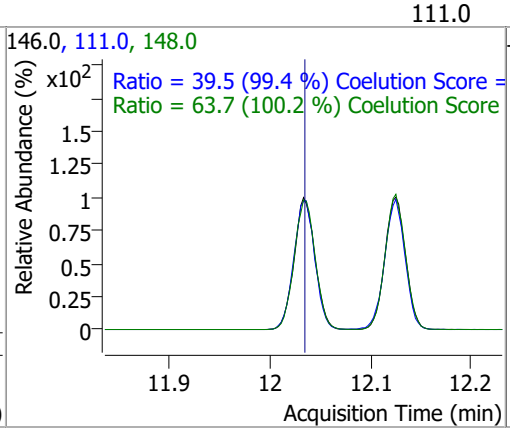


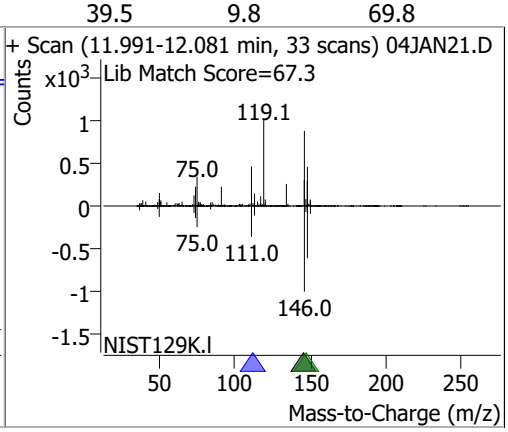




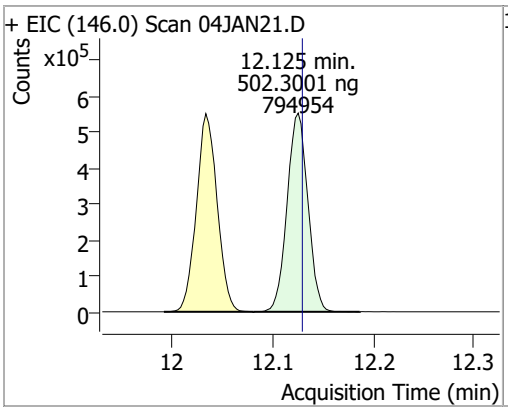
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	511.5504	12.03	0.00	793993	148.0	63.7	33.6	93.6
					111.0	39.5	9.8	69.8

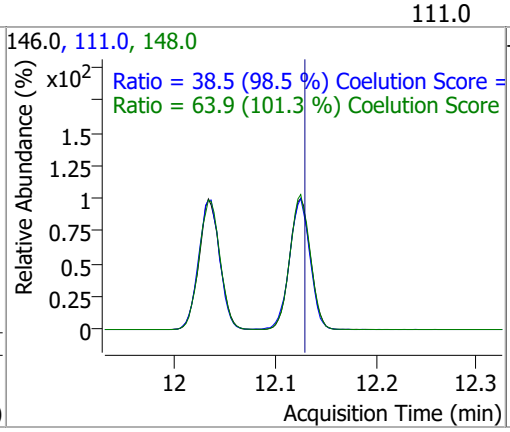


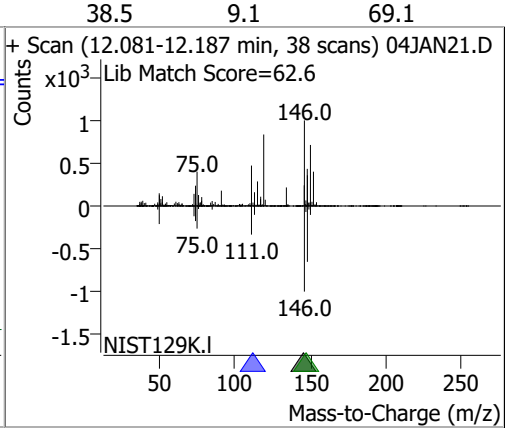




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	502.3001	12.13	0.00	794954	148.0	63.9	33.1	93.1
					111.0	38.5	9.1	69.1

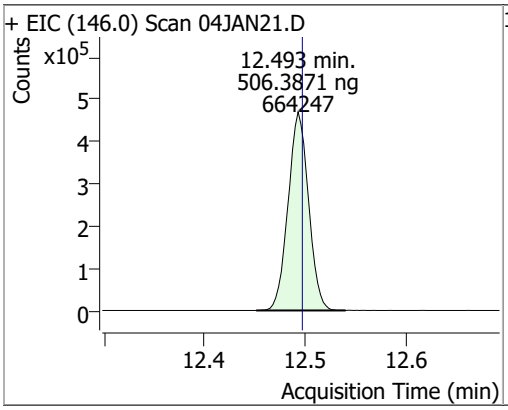


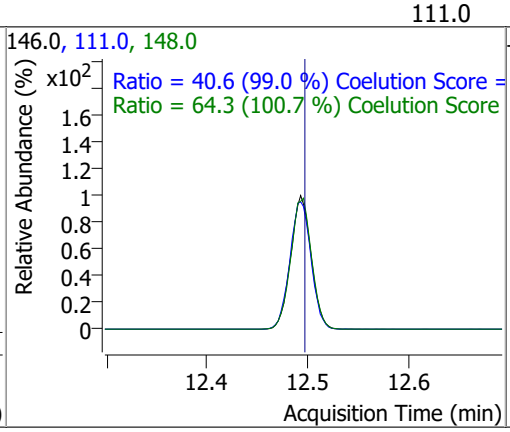


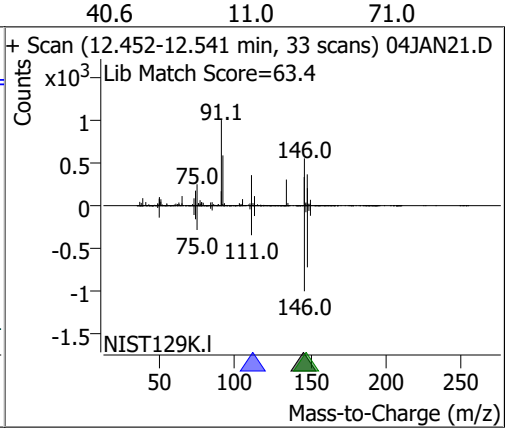


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	506.3871	12.49	0.00	664247	148.0	64.3	33.9	93.9
					111.0	40.6	11.0	71.0







Quantitation Results Report (QT Reviewed)

Data File	04JAN23.D	Operator	MSC
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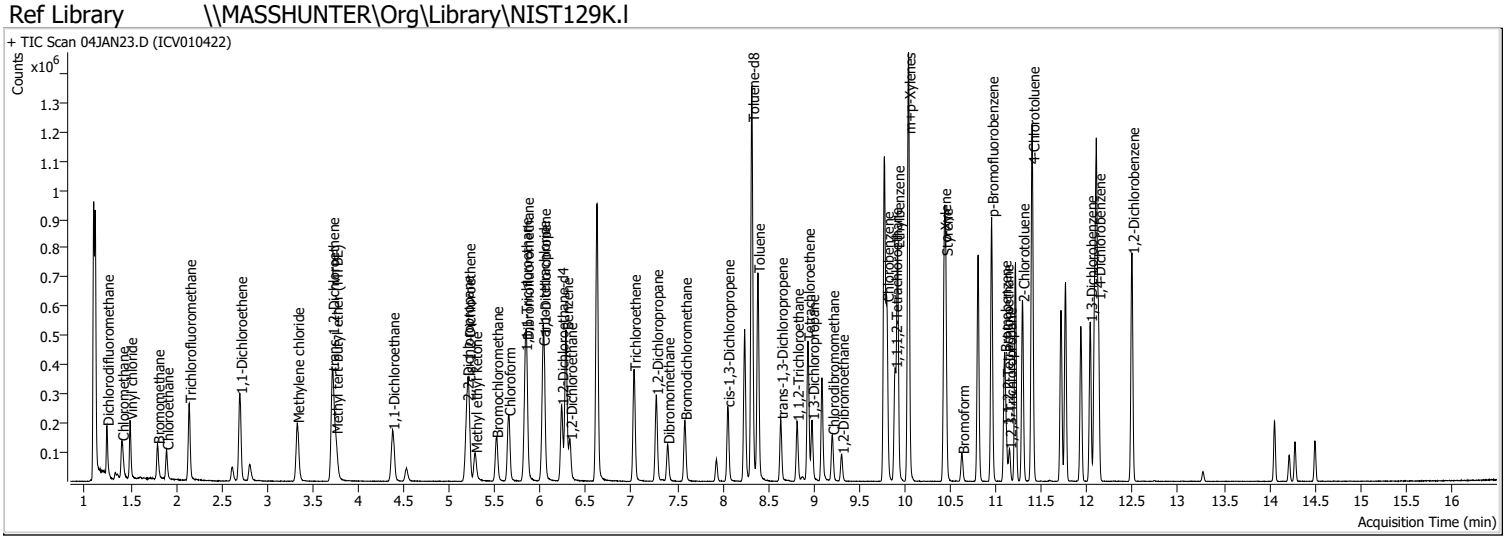
Vial	23	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010422_8260B.batch.bin	Last Calib Update	1/9/2022 8:59:52 PM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

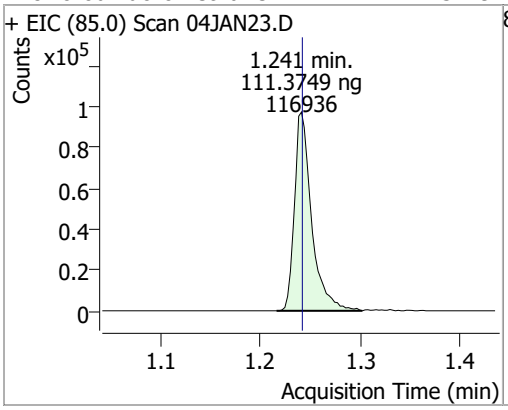


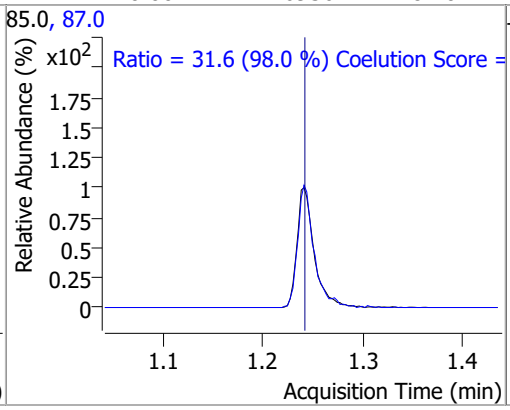
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.621	96.0	801210	250.0000	ng	-0.003
M Chlorobenzene-d5	9.774	82.0	307868	250.0000	ng	0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	255907	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.848	113.0	204707	271.1994	ng	0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 108.48%		
S 1,2-Dichloroethane-d4	6.230	67.0	91382	280.2886	ng	-0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 112.12%		
S Toluene-d8	8.322	98.0	821531	276.9106	ng	0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 110.76%		
S p-Bromofluorobenzene	10.951	95.0	253034	269.8976	ng	-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 107.96%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.241	85.0	116936	111.3749	ng	99
T Chloromethane	1.406	50.0	138617	108.7739	ng	98
T Vinyl chloride	1.495	62.0	137775	120.1518	ng	86
T Bromomethane	1.796	96.0	59947	116.9157	ng	100
T Chloroethane	1.897	64.0	65619	115.5932	ng	99
T Trichlorofluoromethane	2.145	101.0	173333	121.7847	ng	97
T 1,1-Dichloroethene	2.702	96.0	108512	134.4566	ng	100
T Methylene chloride	3.330	49.0	144585	121.5297	ng	98
T trans-1,2-Dichloroethene	3.715	96.0	110909	134.7028	ng	98
T Methyl tert-butyl ether (MTBE)	3.754	73.0	143378	134.7224	ng	99
T 1,1-Dichloroethane	4.376	63.0	208131	135.8030	ng	98
T 2,2-Dichloropropane	5.190	77.0	150902	131.4031	ng	97
T cis-1,2-Dichloroethene	5.209	96.0	108623	130.1231	ng	99
T Methyl ethyl ketone	5.282	43.0	135511	1198.4439	ng	98
T Bromochloromethane	5.513	128.0	42744	123.6009	ng	99
T Chloroform	5.647	83.0	183676	120.4236	ng	100

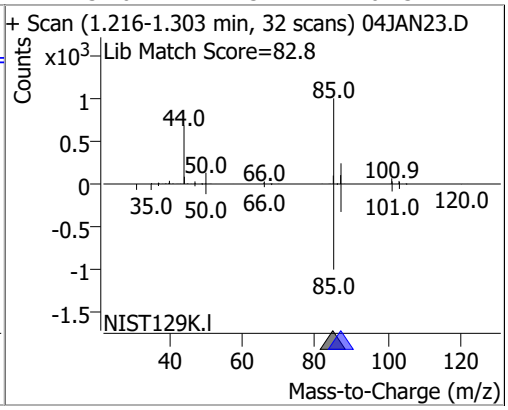
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.831	97.0	183324	128.2524	ng	99
T Carbon tetrachloride	6.027	117.0	181384	128.7928	ng	99
T 1,1-Dichloropropene	6.038	75.0	150930	124.1853	ng	100
T Benzene	6.280	78.0	418900	131.3139	ng	99
T 1,2-Dichloroethane	6.325	62.0	104249	120.7991	ng	95
T Trichloroethene	7.025	95.0	121734	131.1096	ng	99
T 1,2-Dichloropropane	7.270	63.0	102633	125.6626	ng	99
T Dibromomethane	7.393	93.0	43248	125.3047	ng	97
T Bromodichloromethane	7.585	83.0	122757	128.8759	ng	100
T cis-1,3-Dichloropropene	8.054	75.0	130910	121.5561	ng	98
T Toluene	8.386	92.0	264584	132.0244	ng	100
T trans-1,3-Dichloropropene	8.637	75.0	98907	129.0216	ng	97
T 1,1,2-Trichloroethane	8.815	83.0	49128	123.0361	ng	98
T Tetrachloroethene	8.935	163.8	103027	126.0141	ng	99
T 1,3-Dichloropropane	8.980	76.0	95697	121.8442	ng	98
T Chlorodibromomethane	9.203	129.0	78076	125.1103	ng	98
T 1,2-Dibromoethane	9.306	107.0	54259	124.2764	ng	100
T Chlorobenzene	9.802	112.0	288815	131.6352	ng	99
T 1,1,1,2-Tetrachloroethane	9.892	131.0	97148	126.6657	ng	95
T Ethylbenzene	9.917	91.0	501953	131.9113	ng	99
T m+p-Xylenes	10.039	106.0	388558	262.7589	ng	100
T o-Xylene	10.430	106.0	174061	132.2214	ng	98
T Styrene	10.449	104.0	291425	137.4974	ng	98
T Bromoform	10.628	172.5	42560	129.9644	ng	98
T Bromobenzene	11.093	156.0	109054	131.6788	ng	99
T 1,1,2,2-Tetrachloroethane	11.113	83.0	60763	127.4722	ng	98
T 1,2,3-Trichloropropane	11.146	110.0	15682	122.9523	ng	99
T 2-Chlorotoluene	11.291	126.0	108192	131.2948	ng	95
T 4-Chlorotoluene	11.400	91.0	368295	137.0790	ng	100
T 1,3-Dichlorobenzene	12.033	146.0	204088	135.1185	ng	99
T 1,4-Dichlorobenzene	12.122	146.0	200032	129.8812	ng	99
T 1,2-Dichlorobenzene	12.493	146.0	164299	128.7104	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

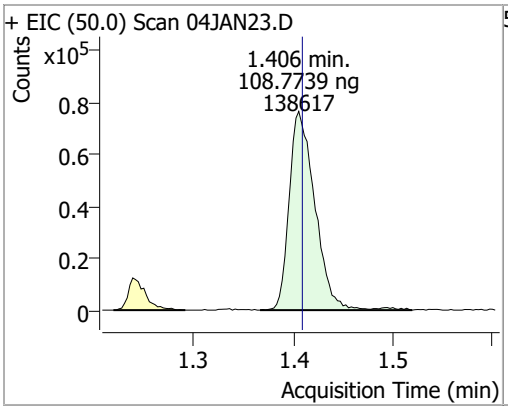
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	111.3749	1.24	0.00	116936	87.0	31.6	2.3	62.3

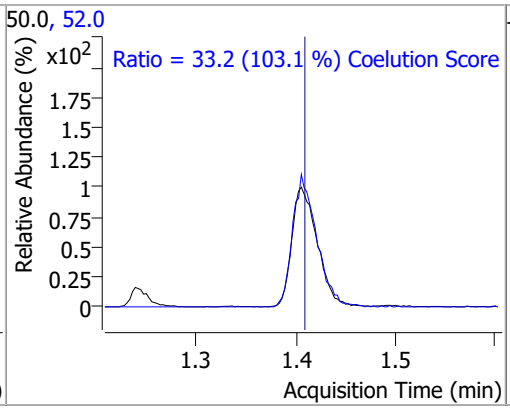


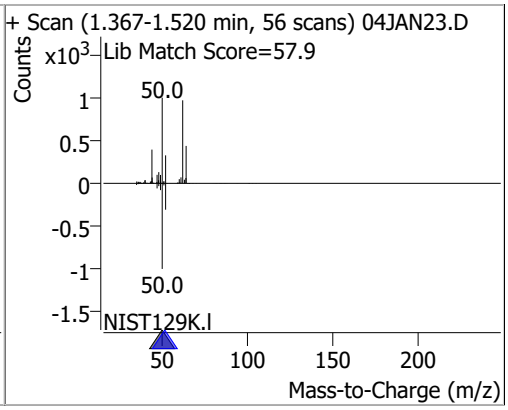




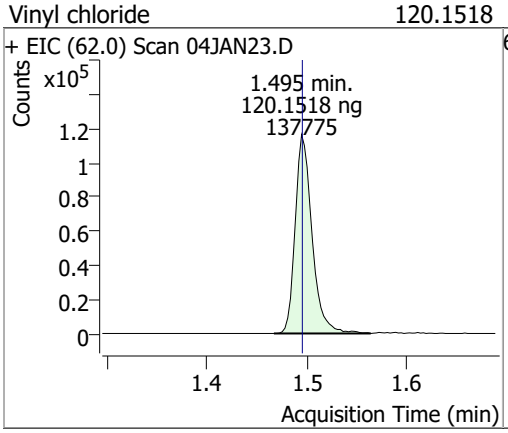
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	108.7739	1.41	0.00	138617	52.0	33.2	2.1	62.1

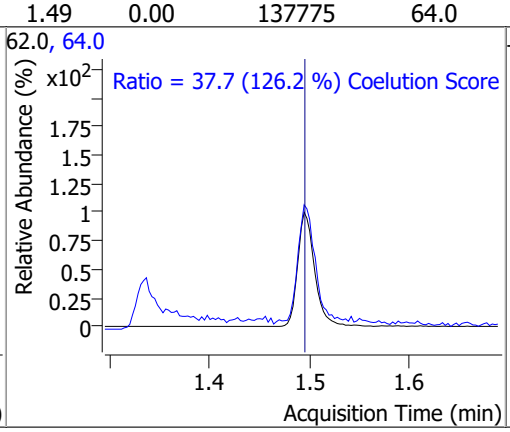


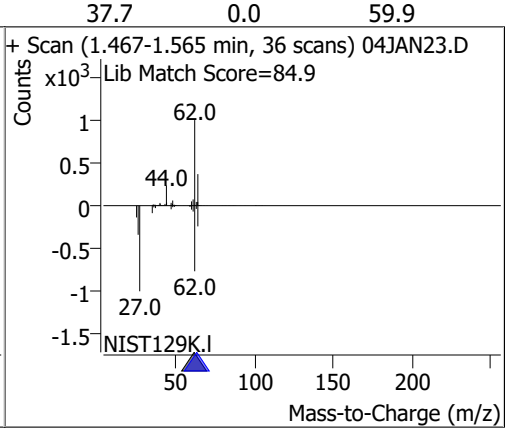




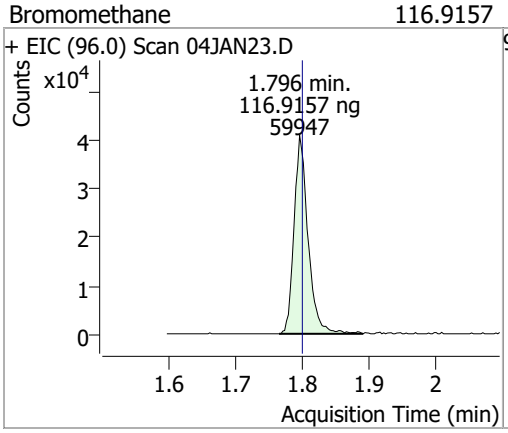
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	120.1518	1.49	0.00	137775	64.0	37.7	0.0	59.9

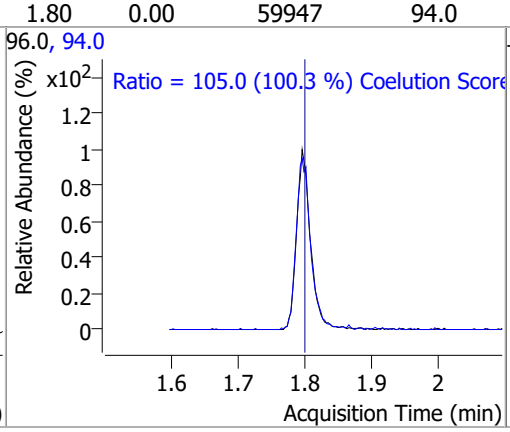


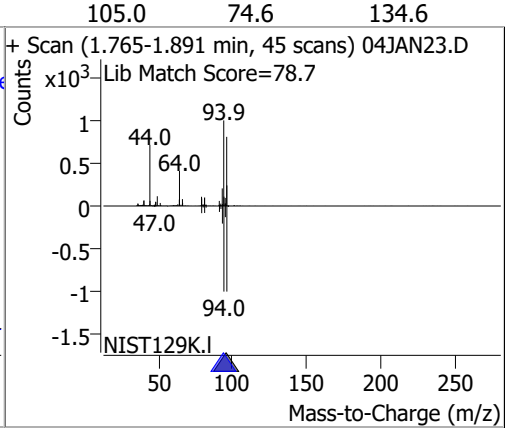




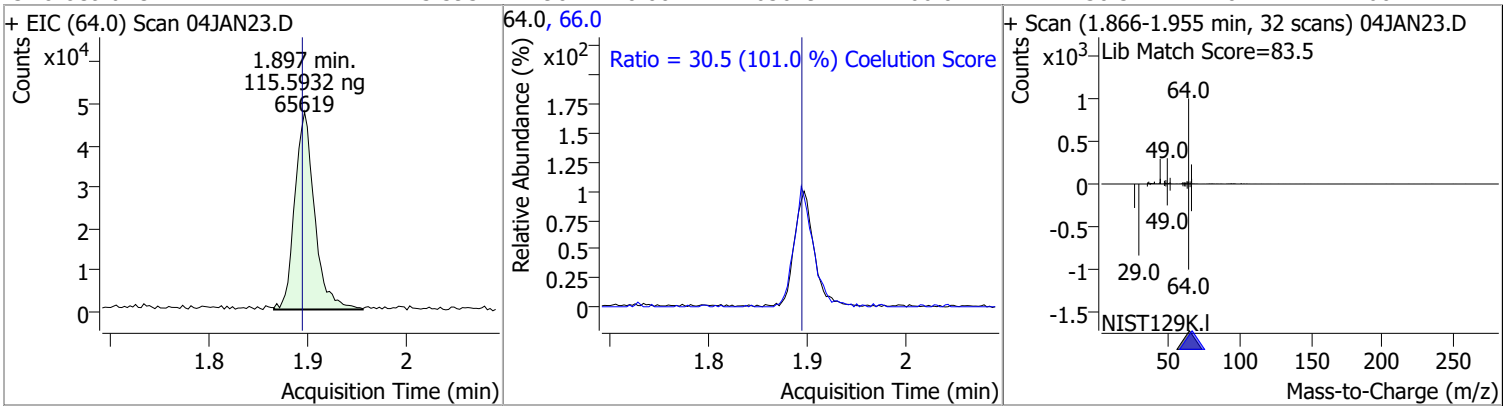
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	116.9157	1.80	0.00	59947	94.0	105.0	74.6	134.6



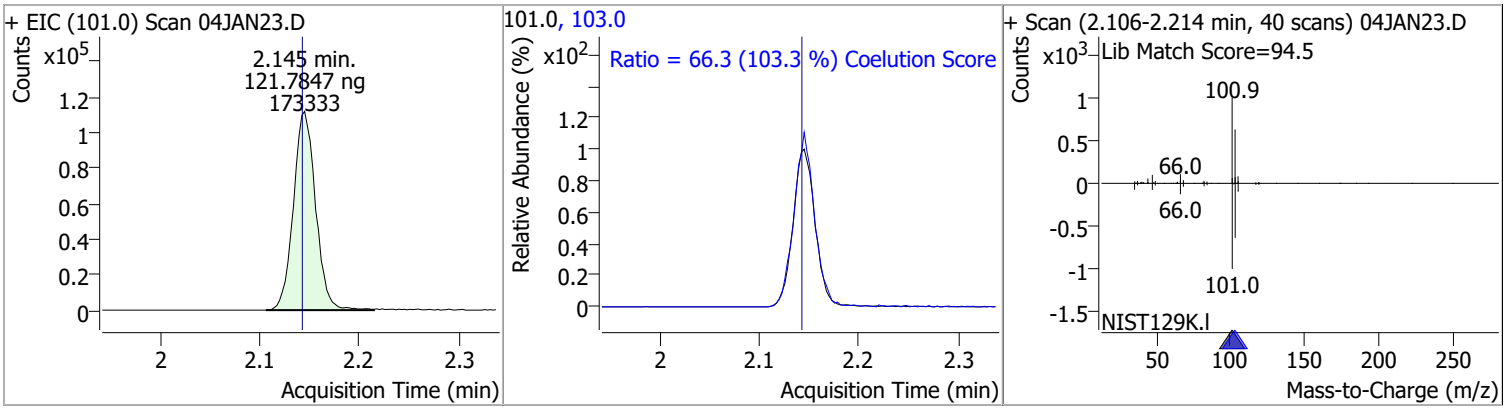




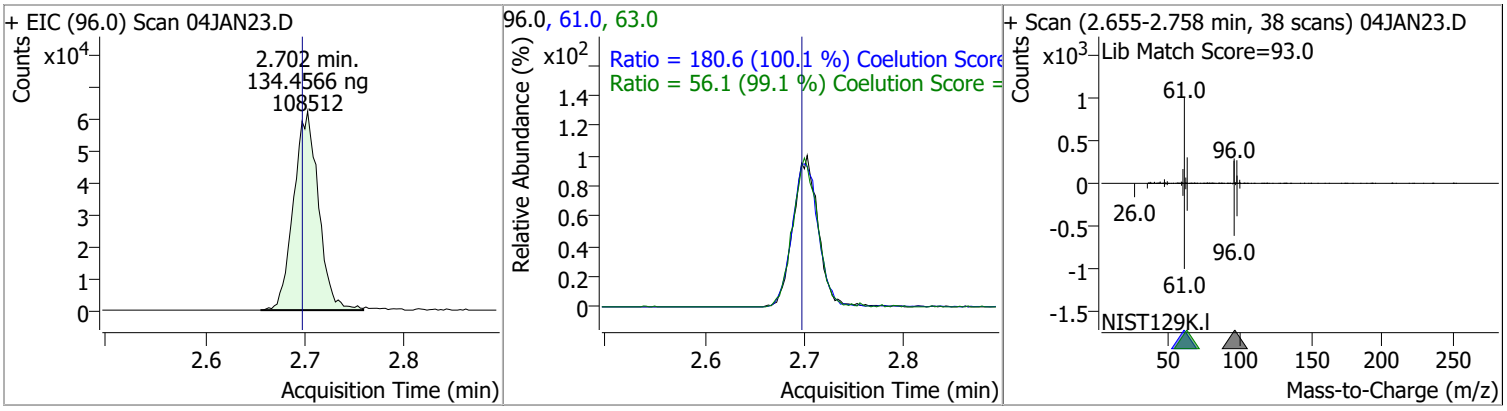
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	115.5932	1.90	0.00	65619	66.0	30.5	0.1	60.1



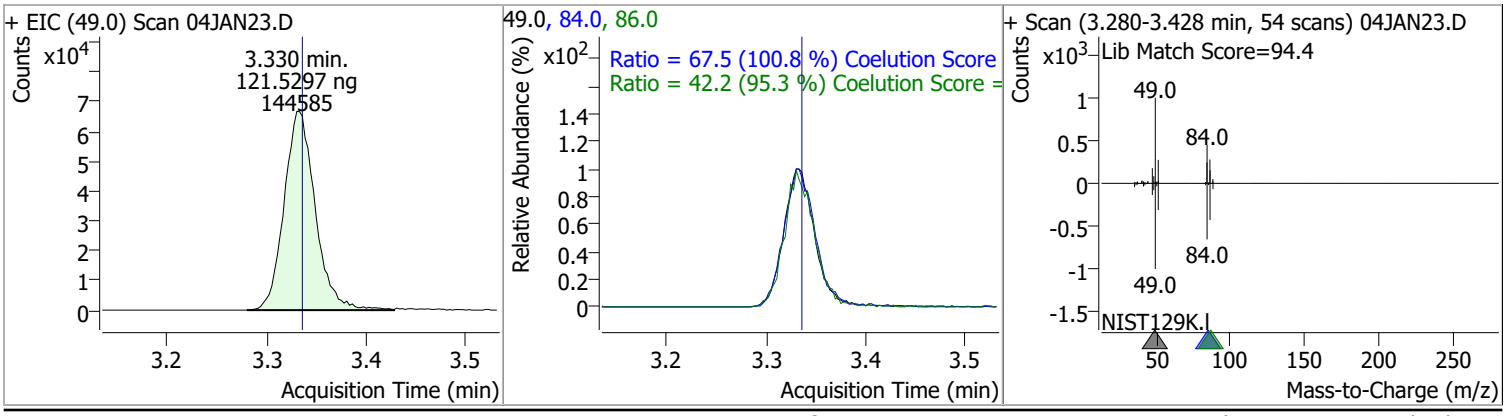
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	121.7847	2.14	0.00	173333	103.0	66.3	34.2	94.2



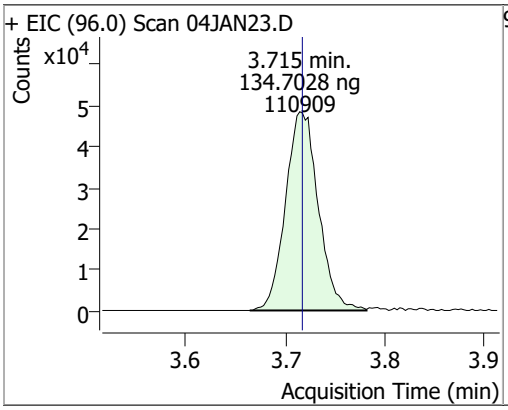
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	134.4566	2.70	0.01	108512	61.0	180.6	150.3	210.3
					63.0	56.1	26.7	86.7

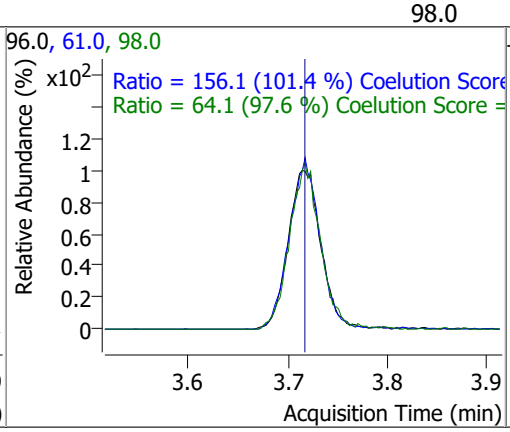


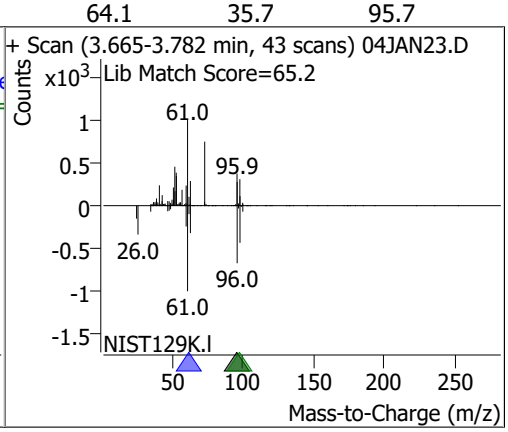
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	121.5297	3.33	-0.01	144585	84.0	67.5	36.9	96.9
					86.0	42.2	14.3	74.3



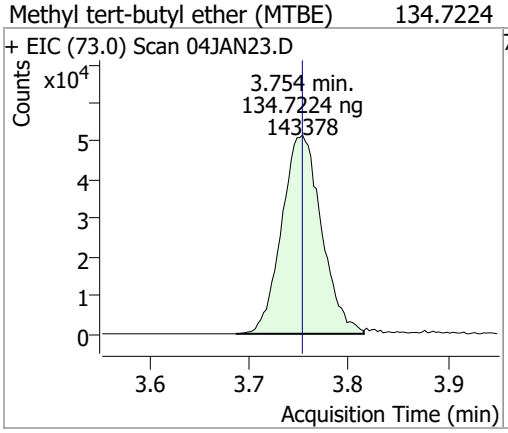
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	134.7028	3.71	0.00	110909	61.0	156.1	123.9	183.9
					98.0	64.1	35.7	95.7

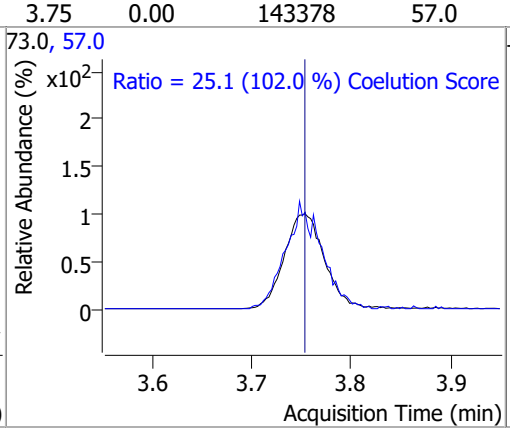


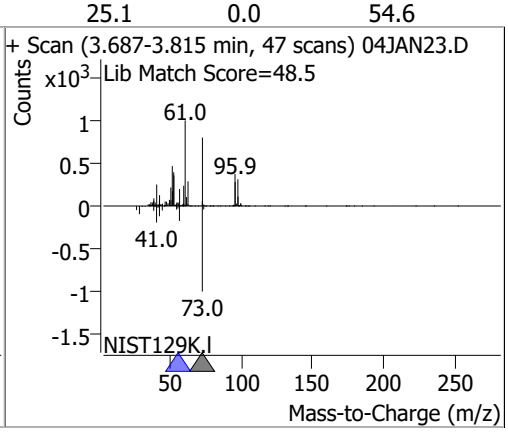




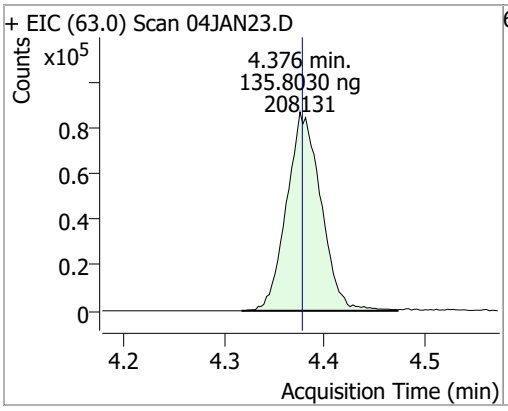
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	134.7224	3.75	0.00	143378	57.0	25.1	0.0	54.6

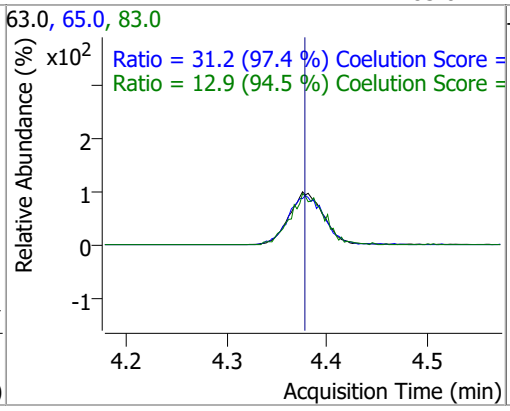


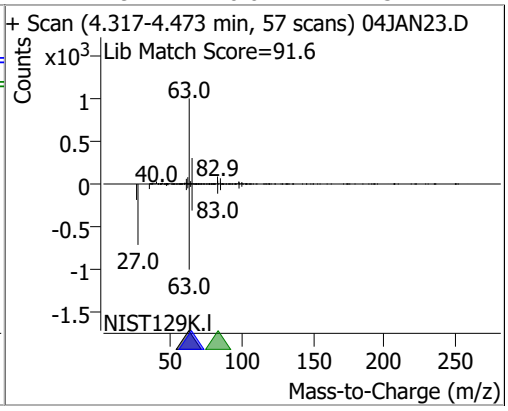




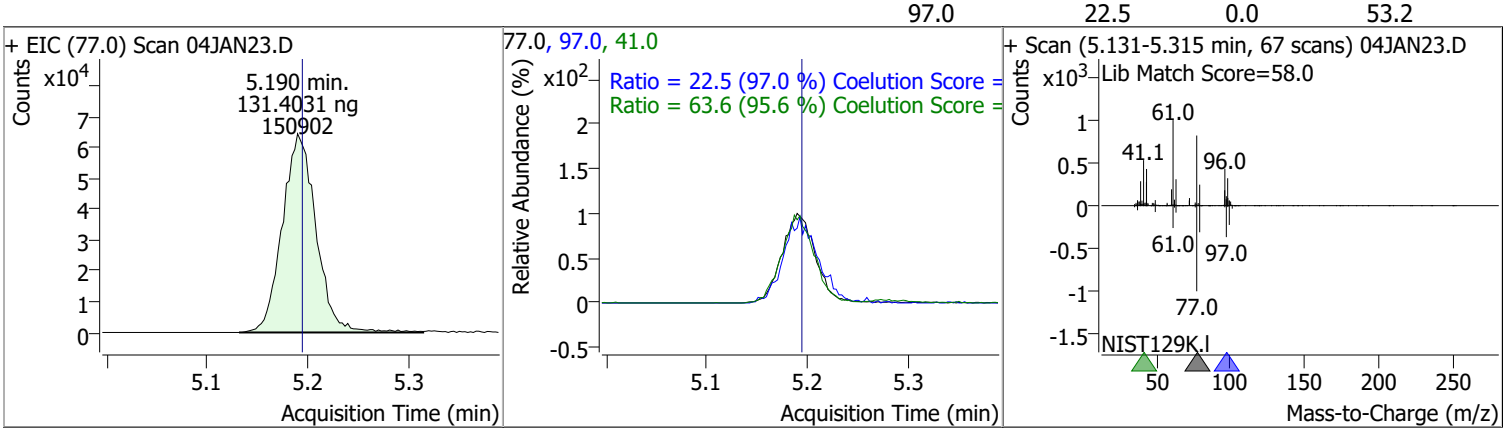
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	135.8030	4.38	0.00	208131	65.0	31.2	2.1	62.1
					83.0	12.9	0.0	43.7



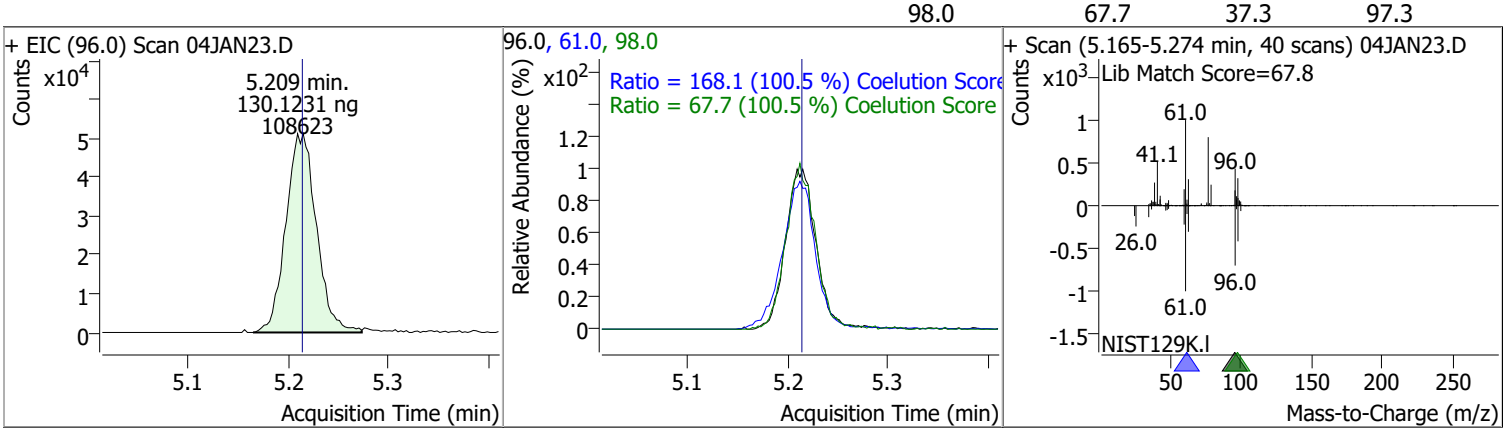




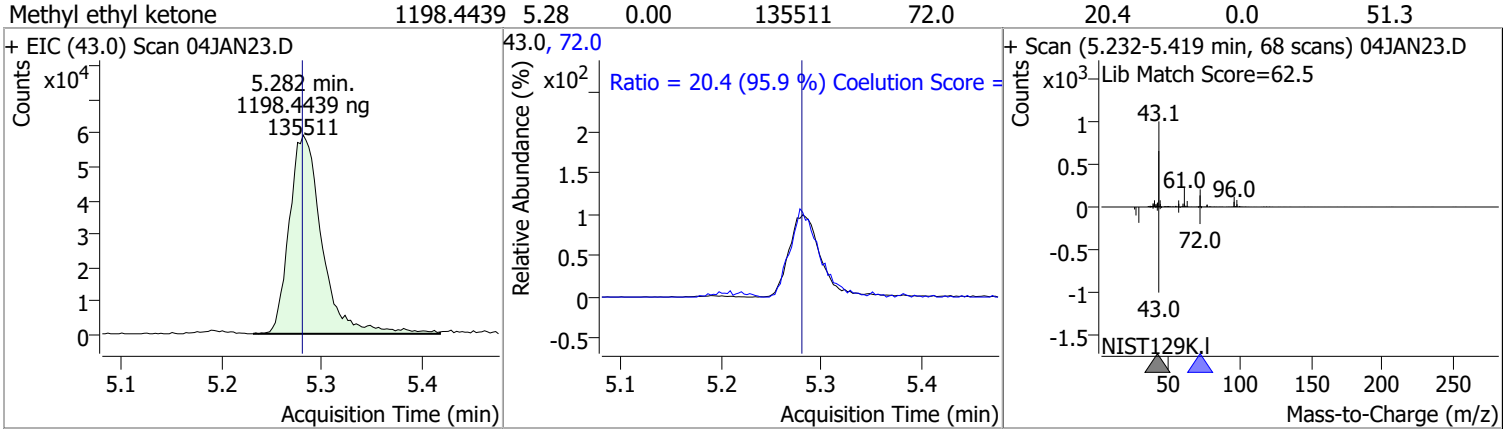
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	131.4031	5.19	-0.01	150902	41.0	63.6	36.5	96.5
					97.0	22.5	0.0	53.2



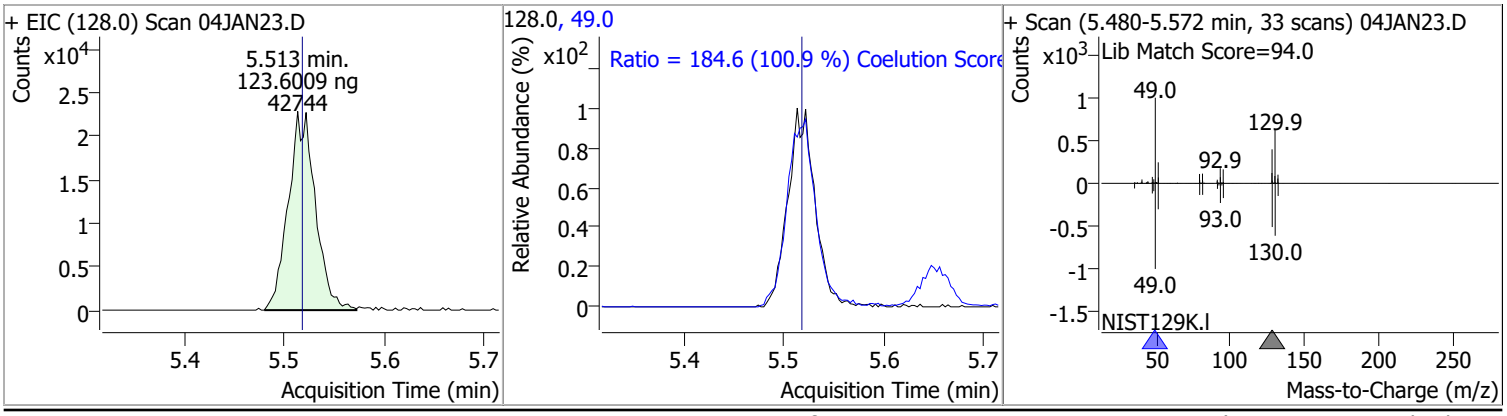
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	130.1231	5.21	-0.01	108623	61.0	168.1	137.2	197.2
					98.0	67.7	37.3	97.3



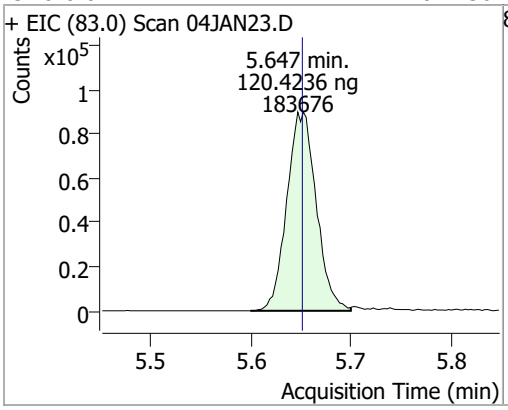
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1198.4439	5.28	0.00	135511	72.0	20.4	0.0	51.3

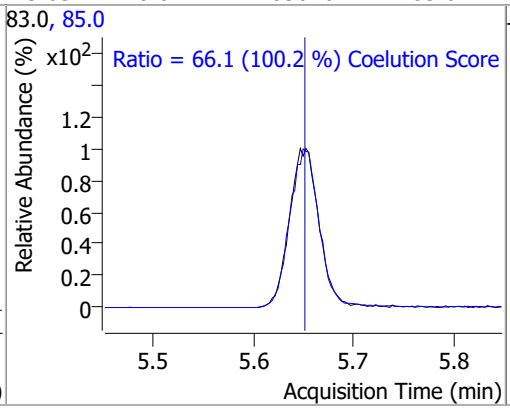


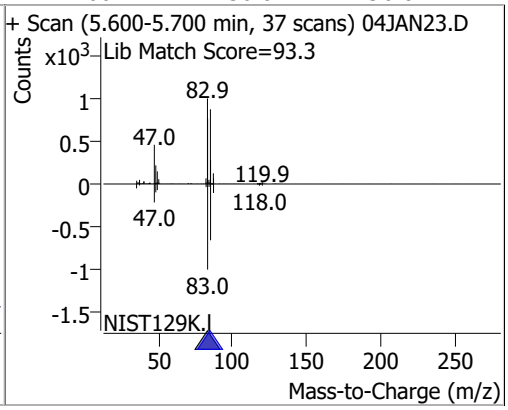
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	123.6009	5.51	-0.01	42744	49.0	184.6	152.9	212.9



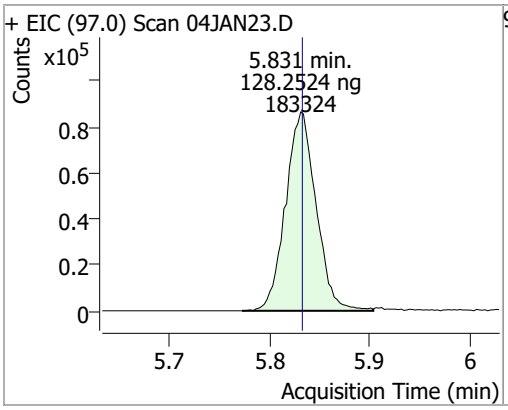
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	120.4236	5.65	-0.01	183676	85.0	66.1	36.0	96.0

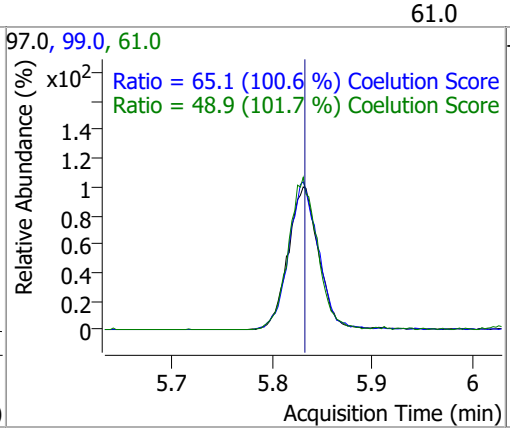


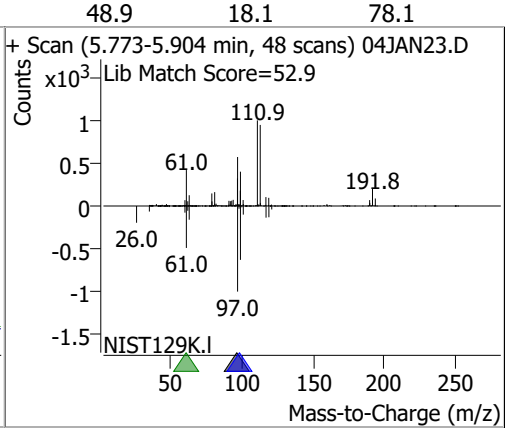




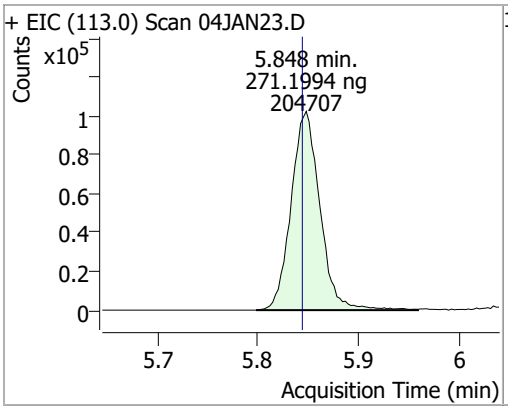
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	128.2524	5.83	0.00	183324	99.0	65.1	34.7	94.7
					61.0	48.9	18.1	78.1

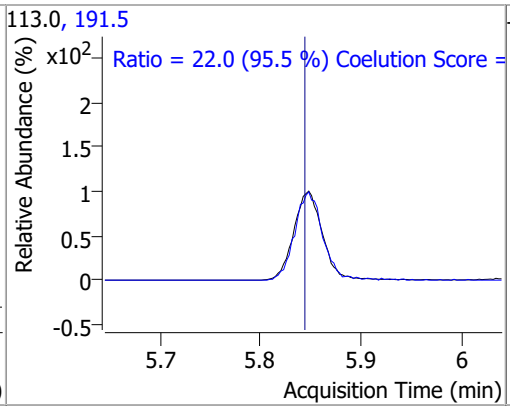


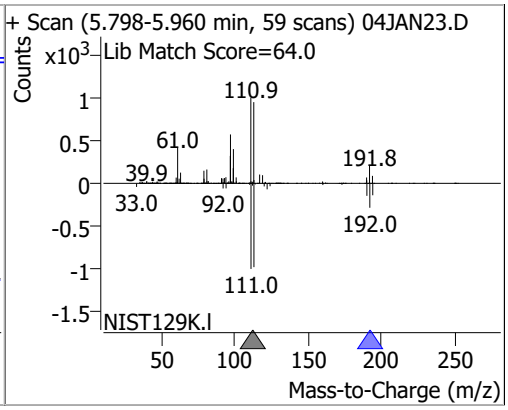




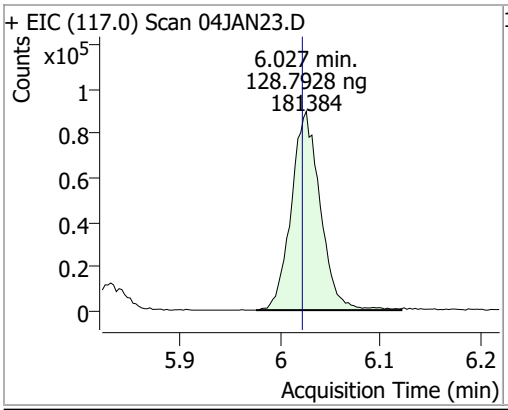
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	271.1994	5.85	0.00	204707	191.5	22.0	0.0	53.1

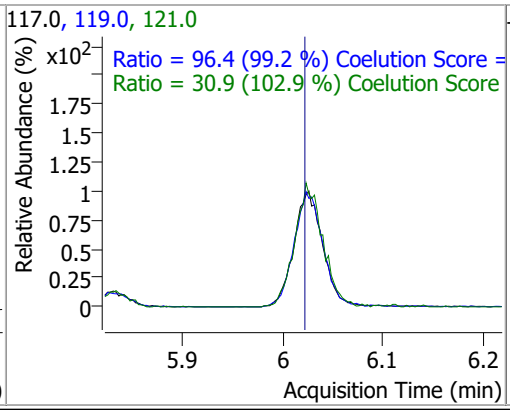


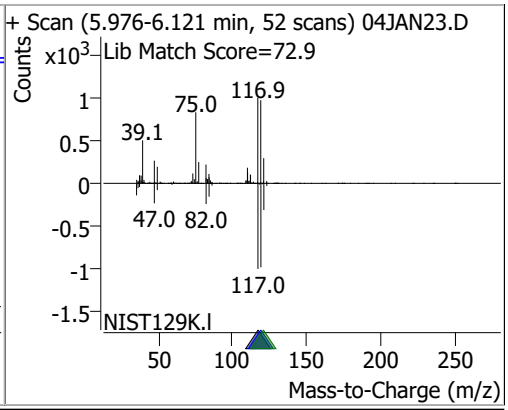




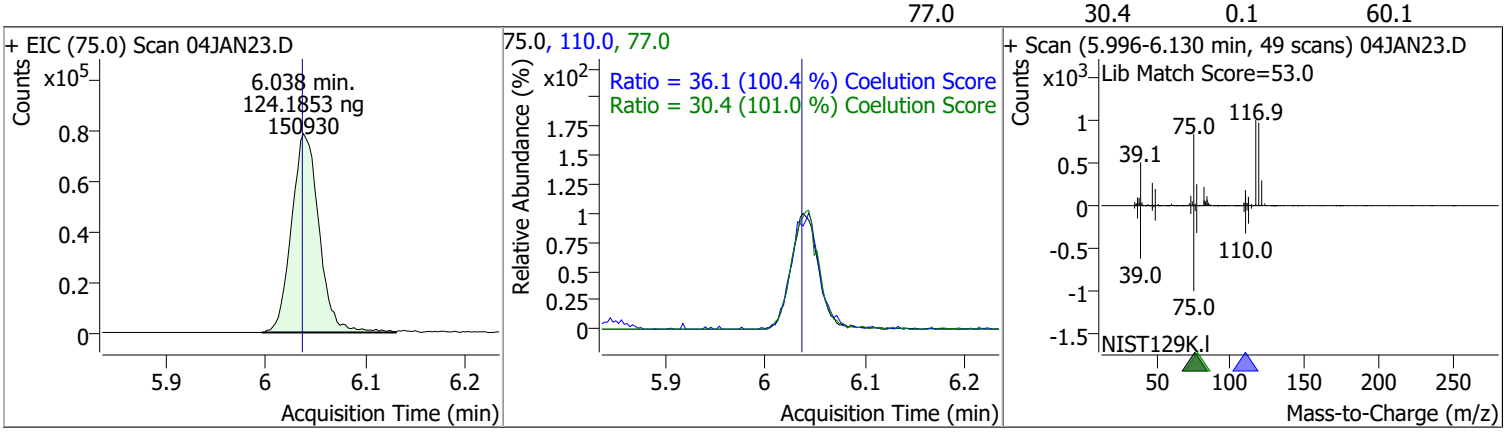
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	128.7928	6.03	0.00	181384	119.0	96.4	67.2	127.2
					121.0	30.9	0.1	60.1



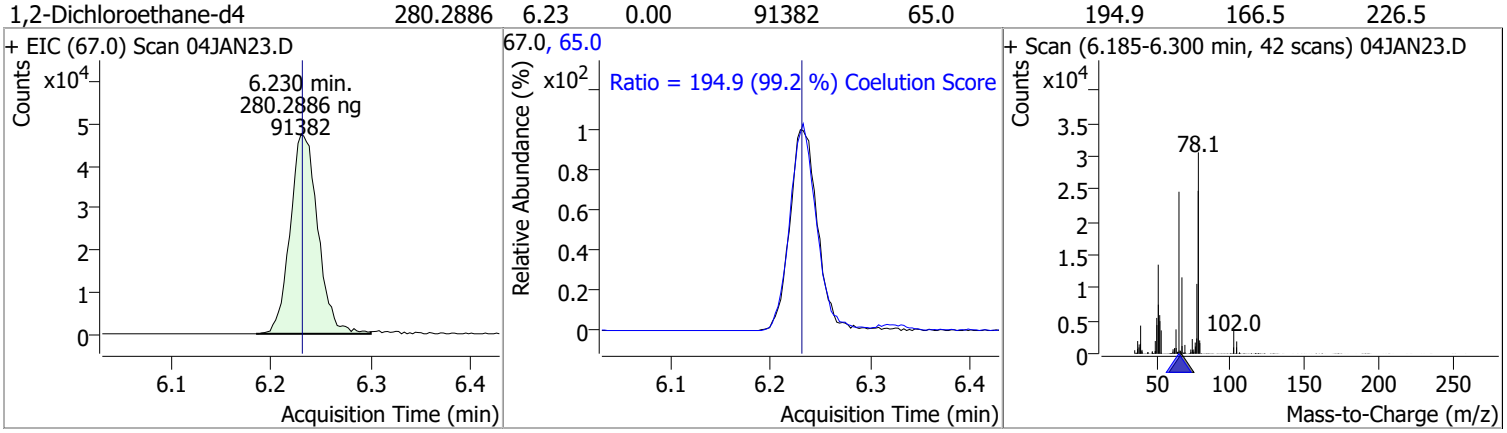




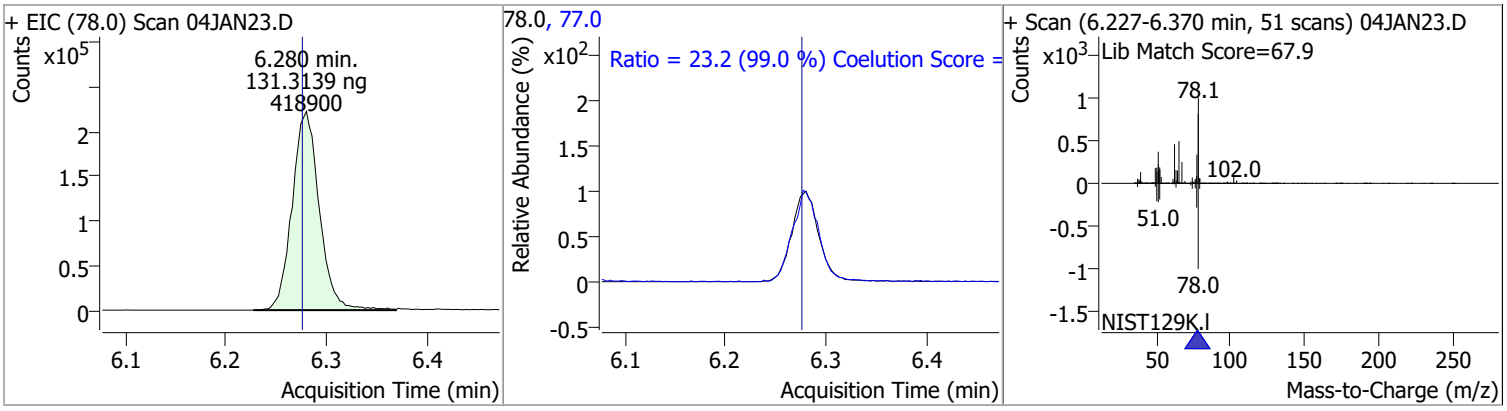
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	124.1853	6.04	0.00	150930	110.0	36.1	5.9	65.9
					77.0	30.4	0.1	60.1



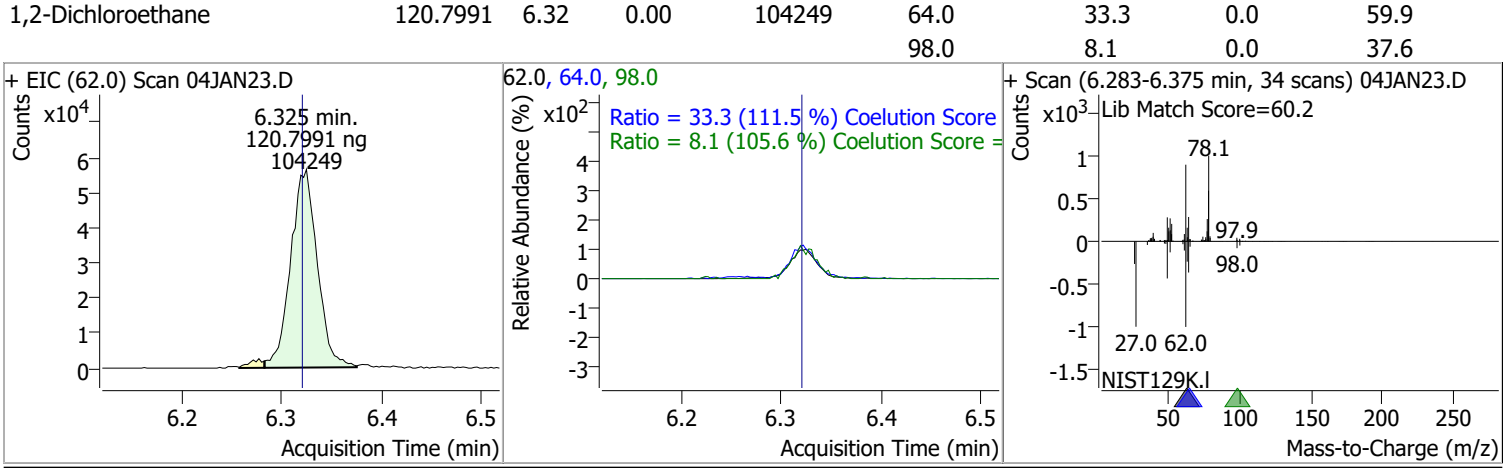
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	280.2886	6.23	0.00	91382	65.0	194.9	166.5	226.5



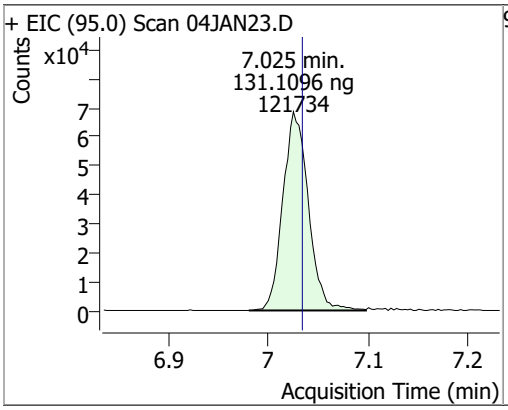
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	131.3139	6.28	0.00	418900	77.0	23.2	0.0	53.5

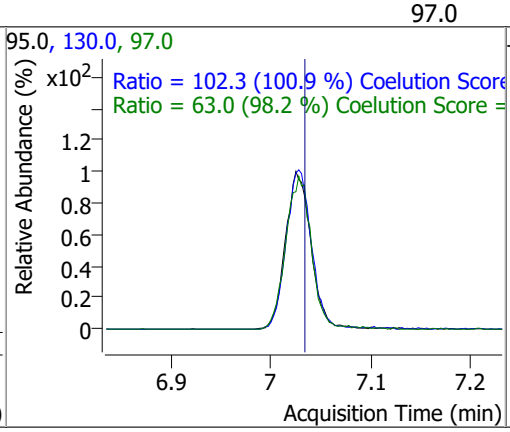


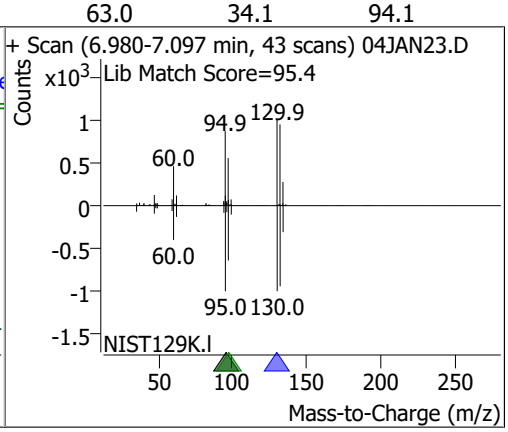
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	120.7991	6.32	0.00	104249	64.0	33.3	0.0	59.9
					98.0	8.1	0.0	37.6



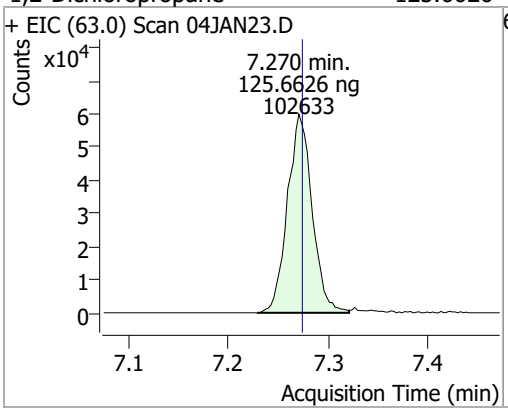
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	131.1096	7.02	-0.01	121734	130.0	102.3	71.5	131.5
					97.0	63.0	34.1	94.1

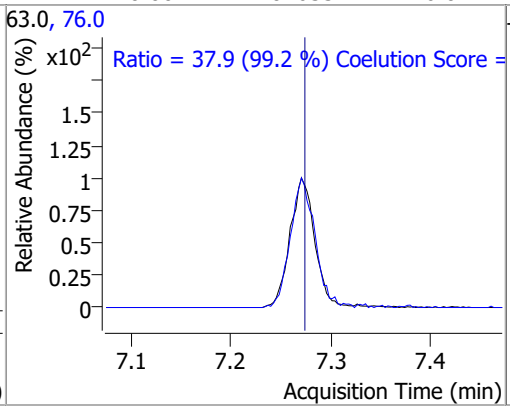


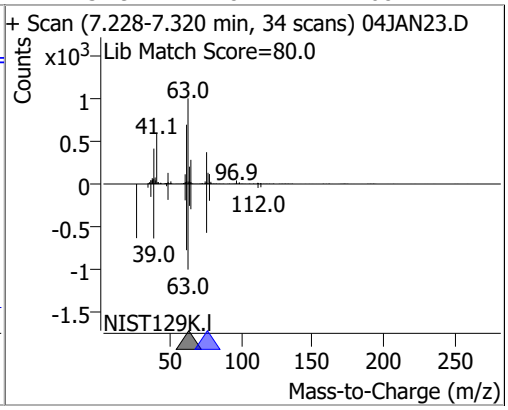




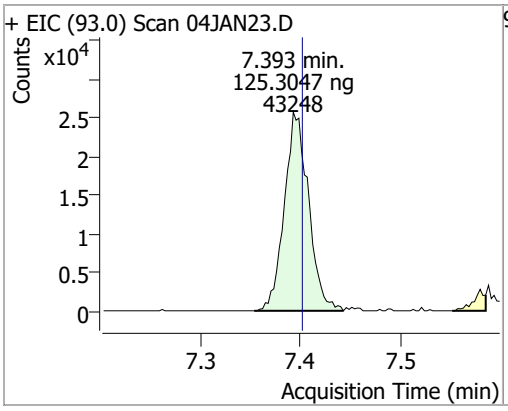
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	125.6626	7.27	0.00	102633	76.0	37.9	8.2	68.2

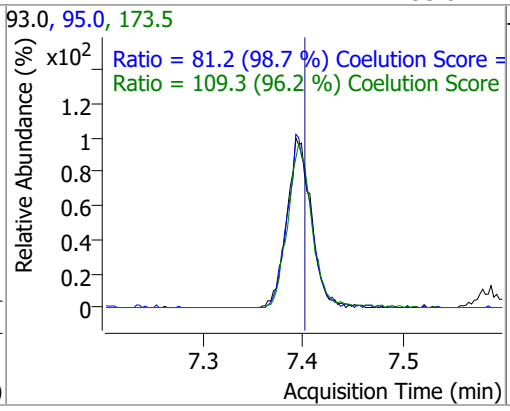


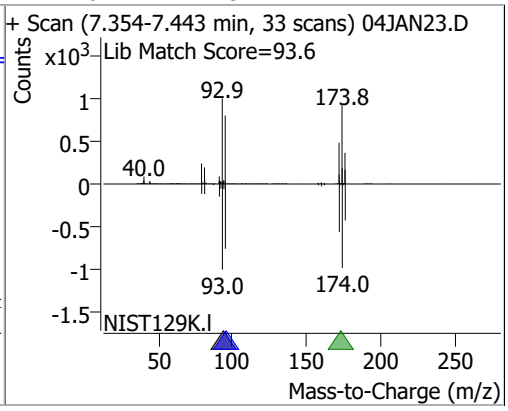




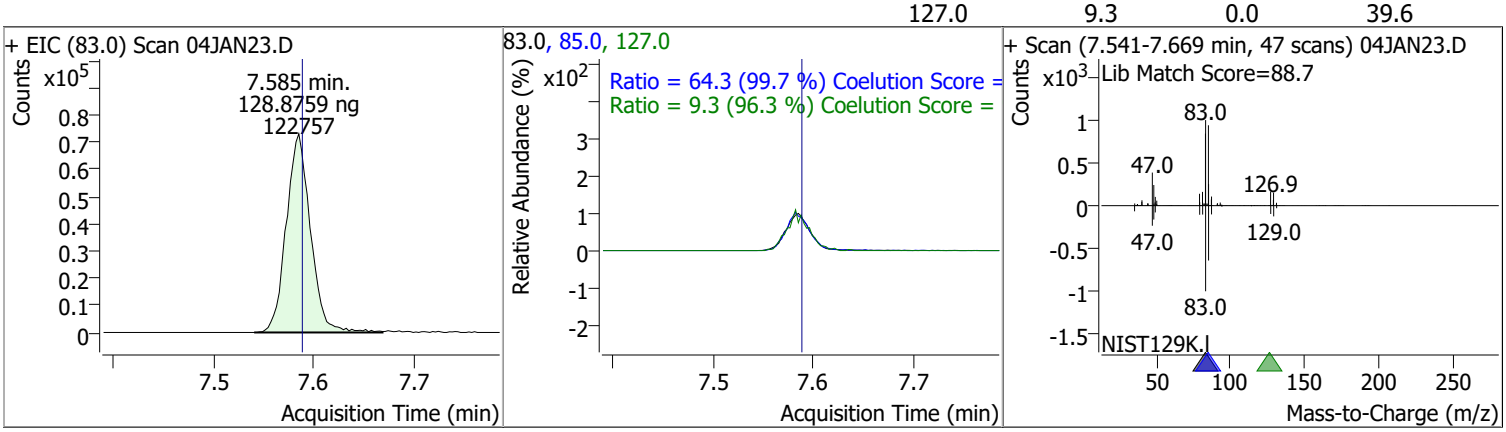
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	125.3047	7.39	-0.01	43248	173.5	109.3	83.7	143.7
					95.0	81.2	52.2	112.2



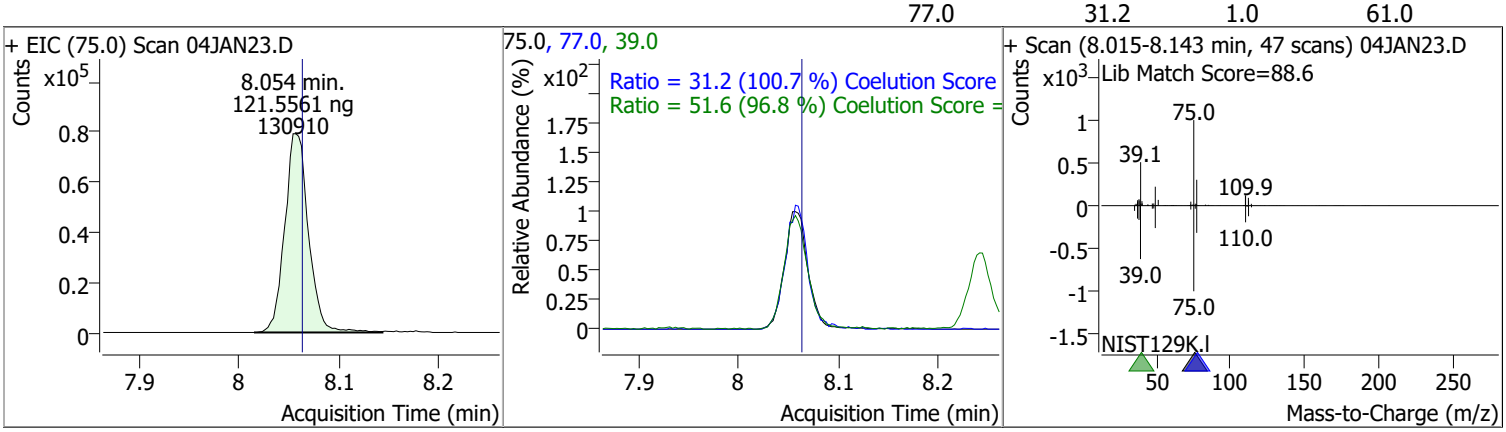




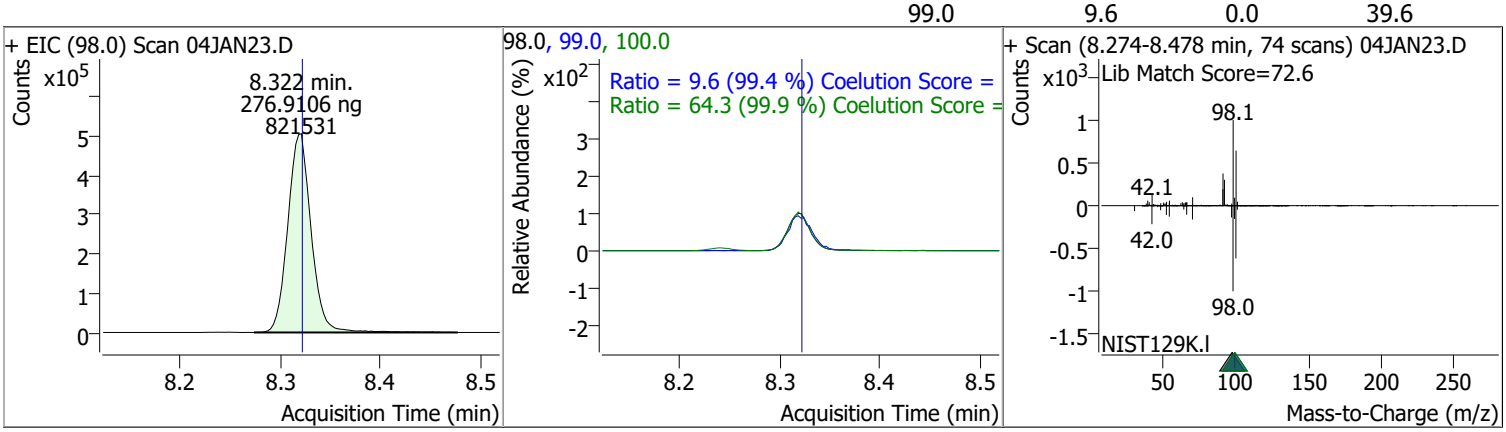
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	128.8759	7.59	0.00	122757	85.0	64.3	34.5	94.5
					127.0	9.3	0.0	39.6



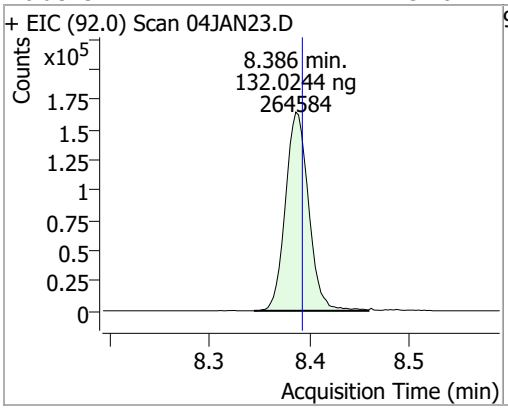
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	121.5561	8.05	-0.01	130910	39.0	51.6	23.3	83.3
					77.0	31.2	1.0	61.0

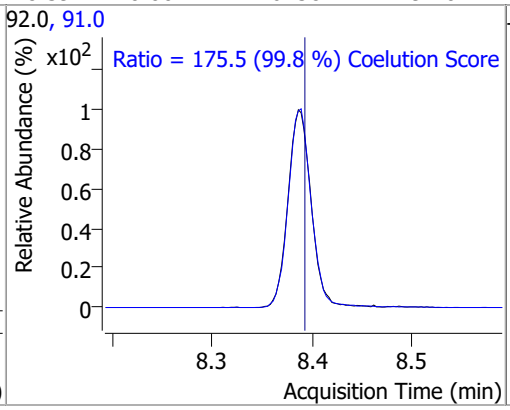


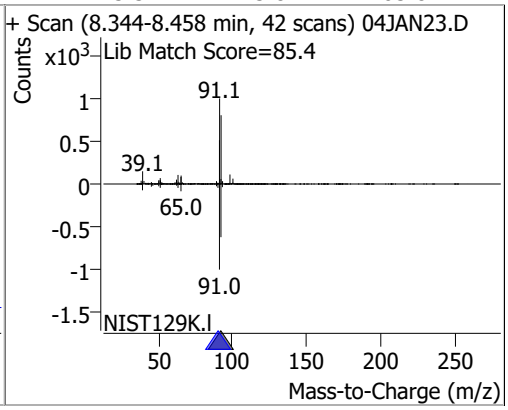
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	276.9106	8.32	0.00	821531	100.0	64.3	34.4	94.4
					99.0	9.6	0.0	39.6



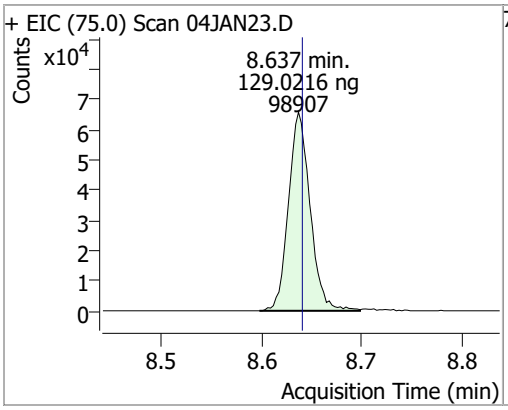
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	132.0244	8.39	0.00	264584	91.0	175.5	145.8	205.8

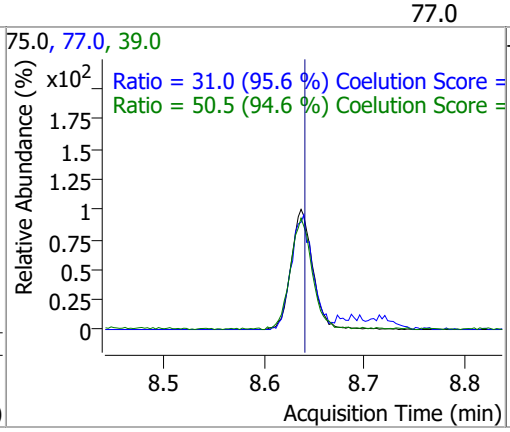


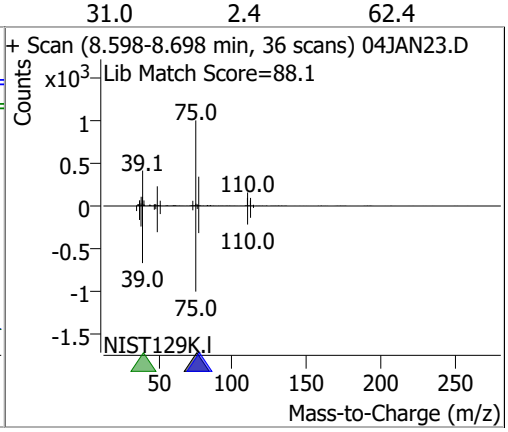




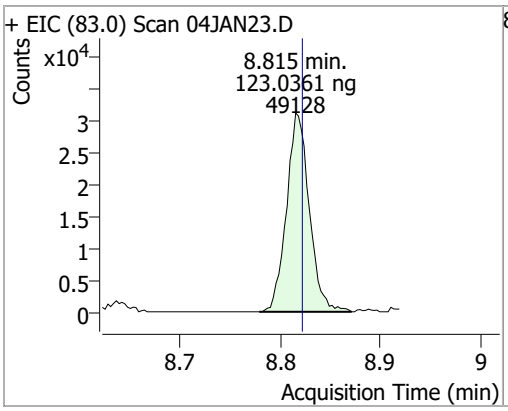
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	129.0216	8.64	0.00	98907	39.0	50.5	23.4	83.4
					77.0	31.0	2.4	62.4

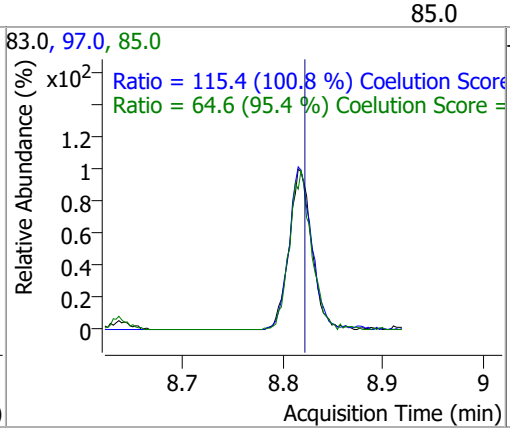


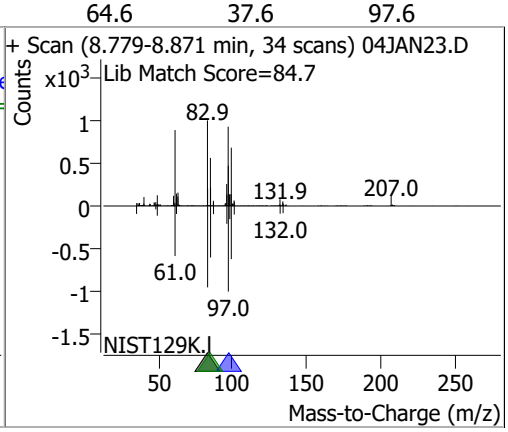




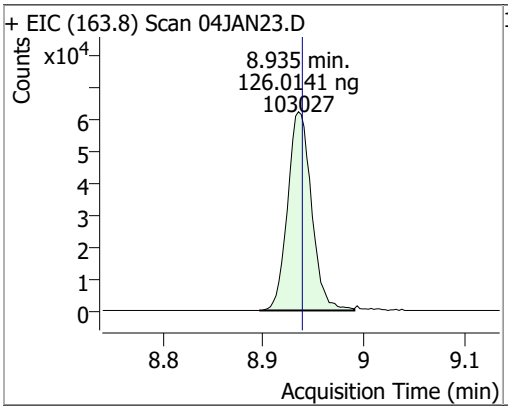
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	123.0361	8.82	0.00	49128	97.0	115.4	84.6	144.6
					85.0	64.6	37.6	97.6

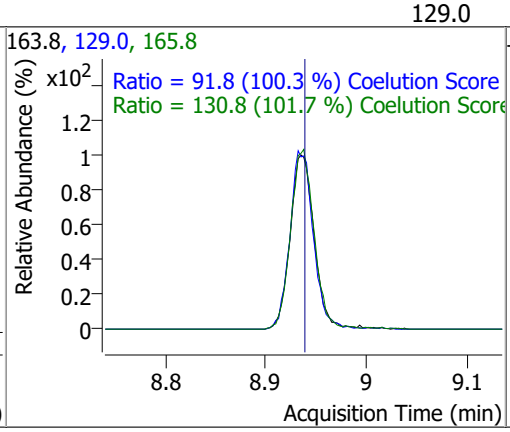


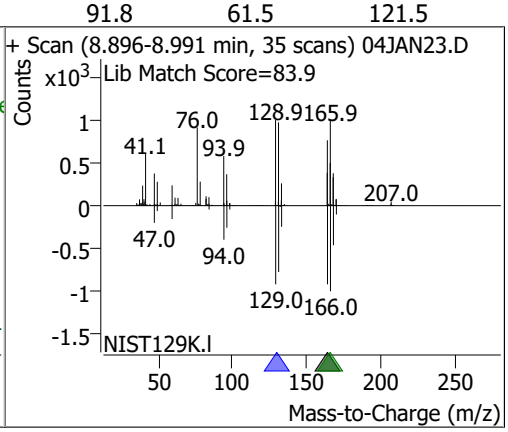




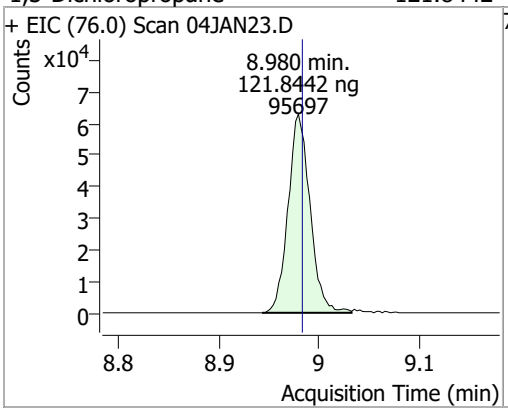
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	126.0141	8.94	0.00	103027	165.8	130.8	98.6	158.6
					129.0	91.8	61.5	121.5

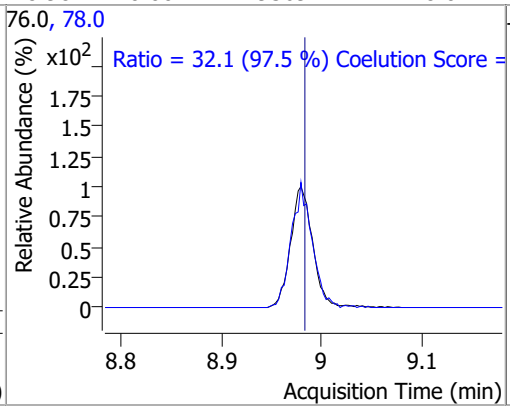


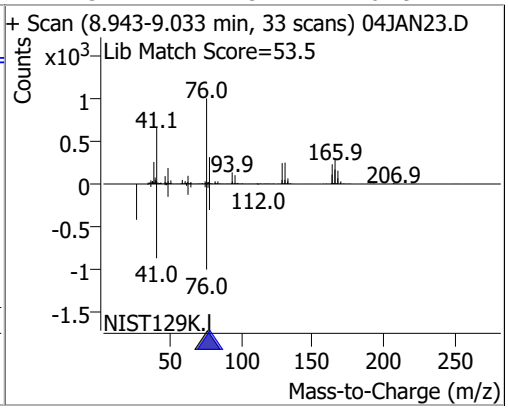




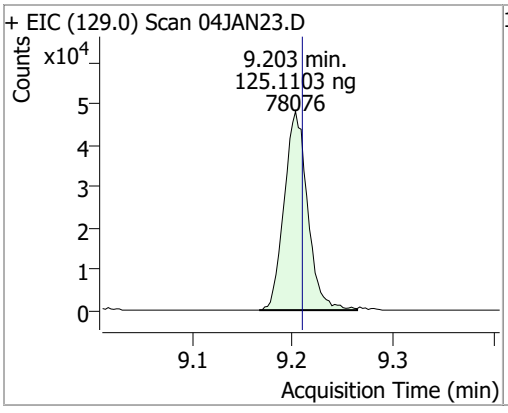
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	121.8442	8.98	0.00	95697	78.0	32.1	2.9	62.9

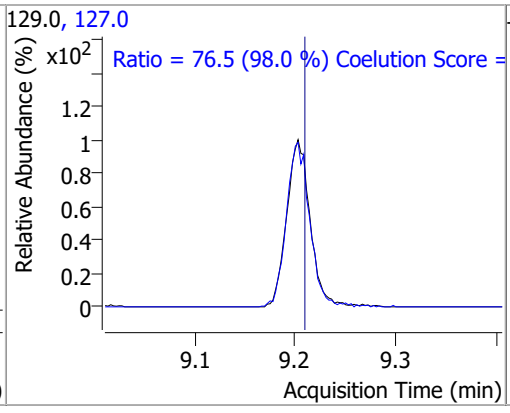


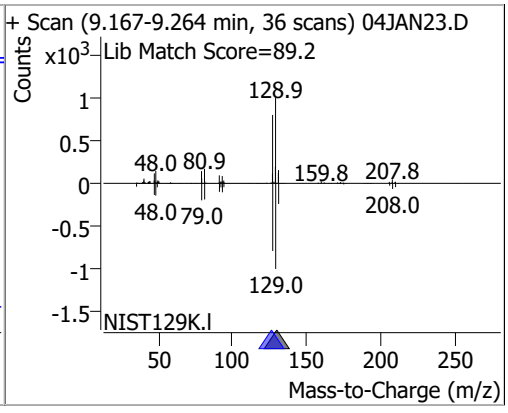




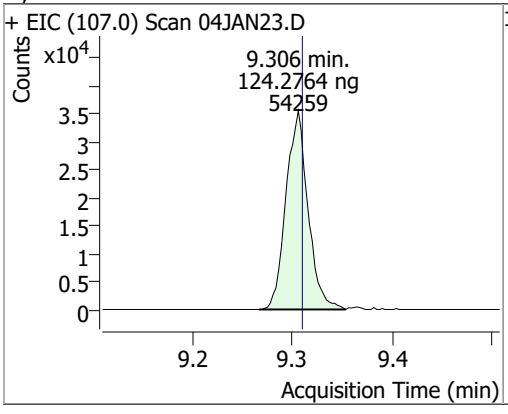
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	125.1103	9.20	0.00	78076	127.0	76.5	48.0	108.0

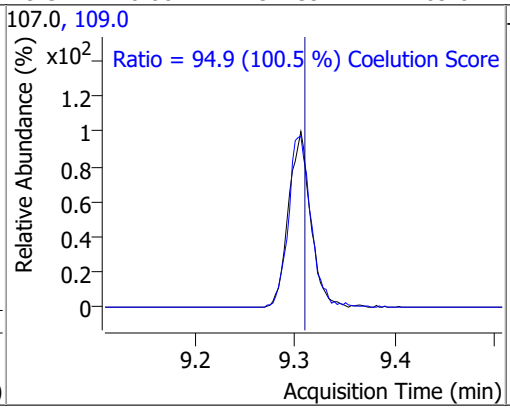


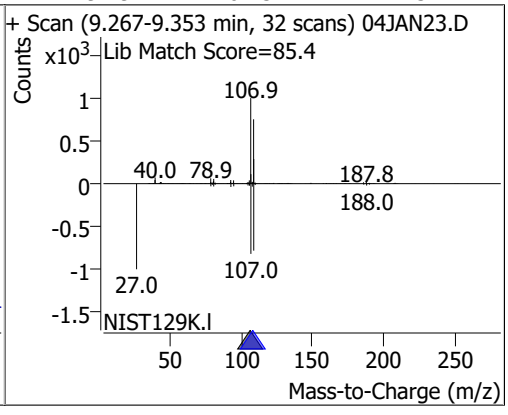




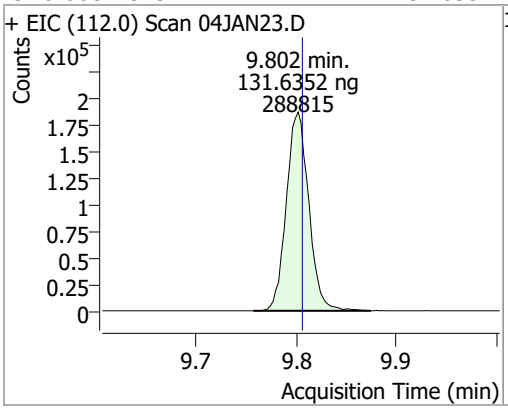
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	124.2764	9.31	0.00	54259	109.0	94.9	64.5	124.5

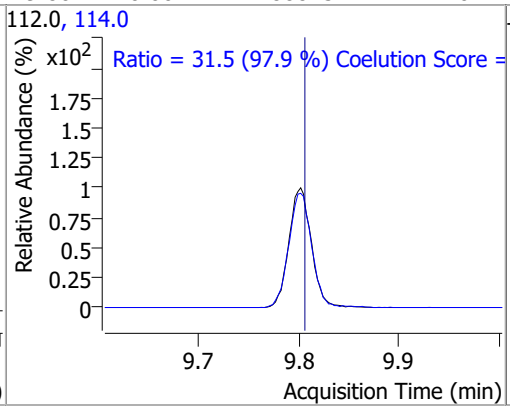


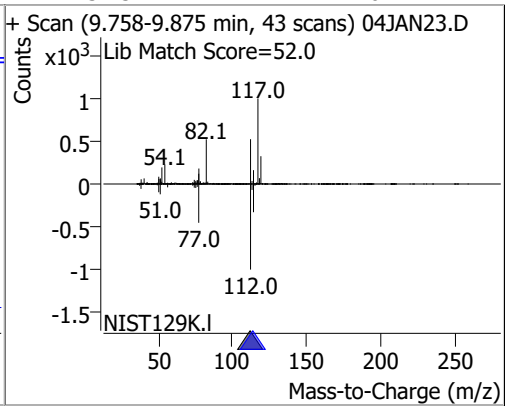




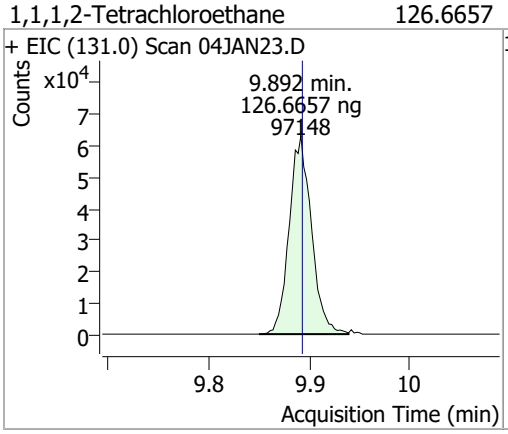
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	131.6352	9.80	0.00	288815	114.0	31.5	2.1	62.1

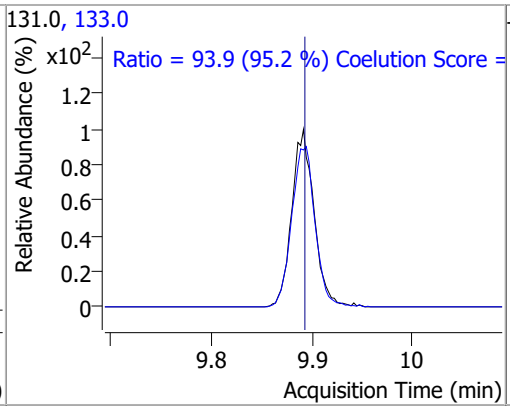


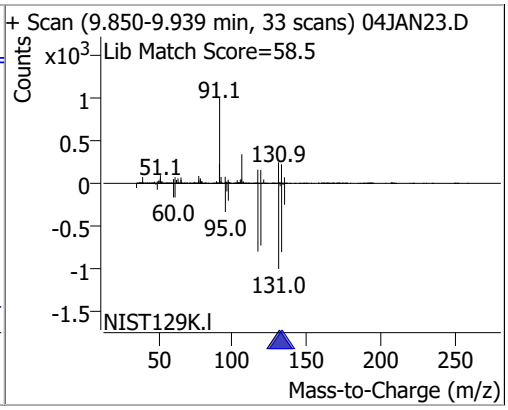




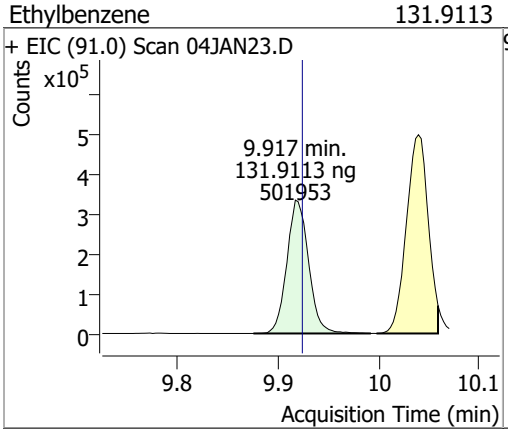
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	126.6657	9.89	0.00	97148	133.0	93.9	68.6	128.6

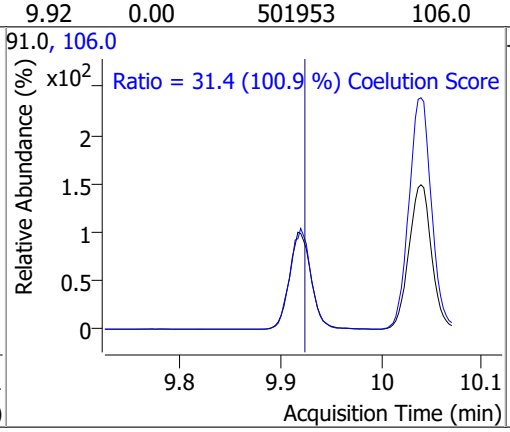


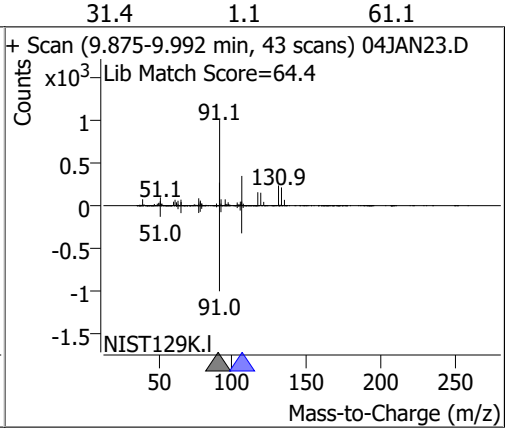




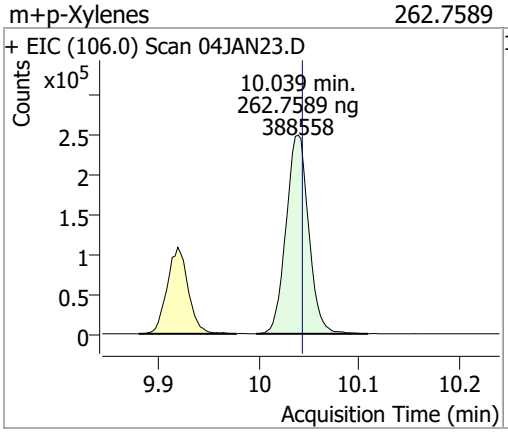
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	131.9113	9.92	0.00	501953	106.0	31.4	1.1	61.1

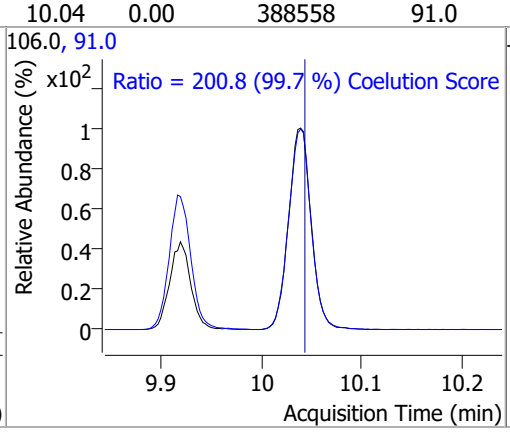


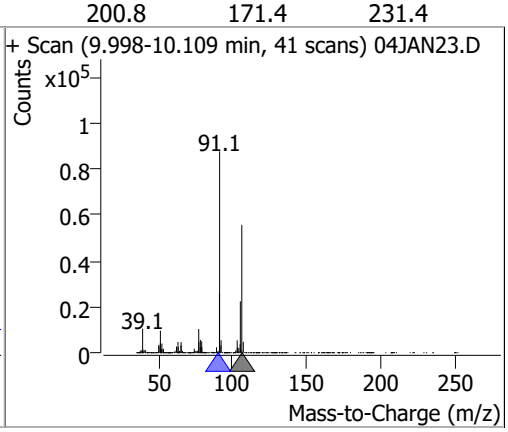




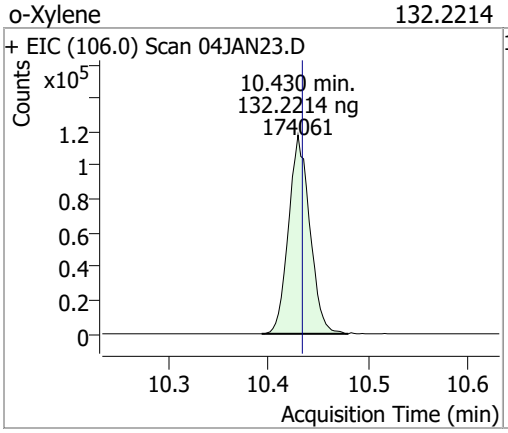
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	262.7589	10.04	0.00	388558	91.0	200.8	171.4	231.4

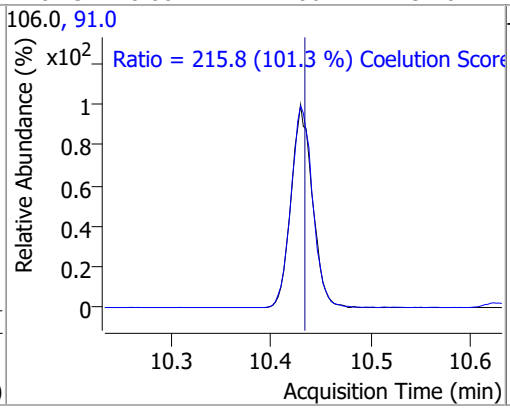


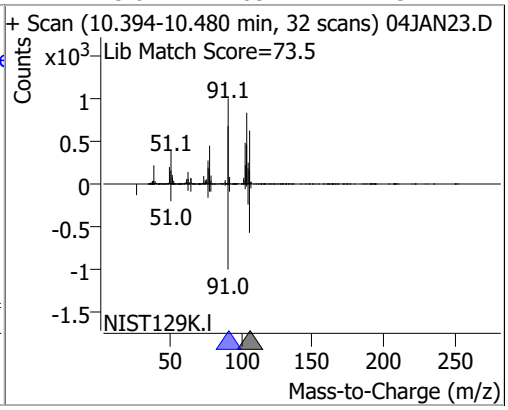




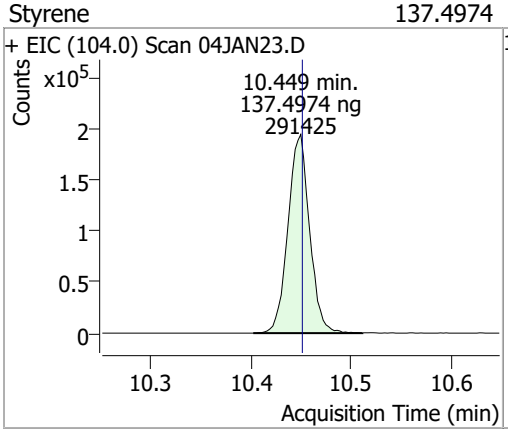
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	132.2214	10.43	0.00	174061	91.0	215.8	183.1	243.1

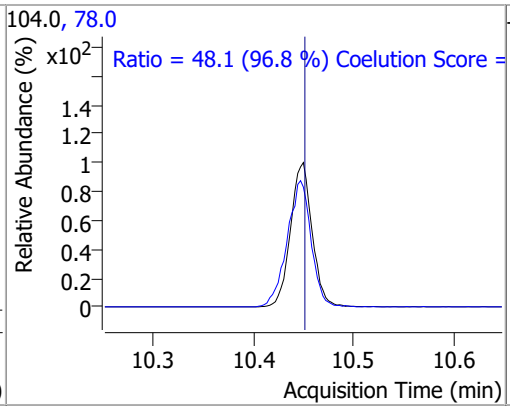


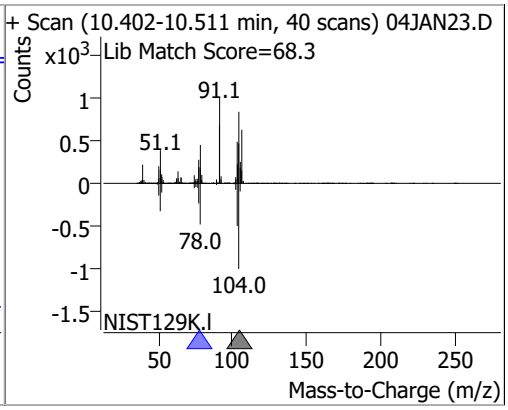




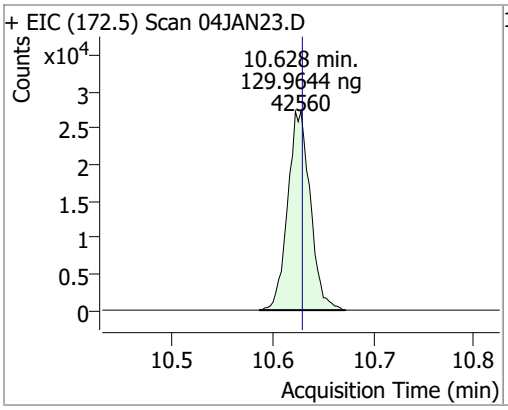
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	137.4974	10.45	0.00	291425	78.0	48.1	19.6	79.6

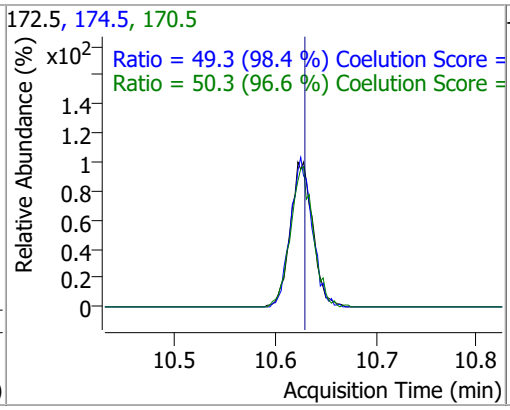


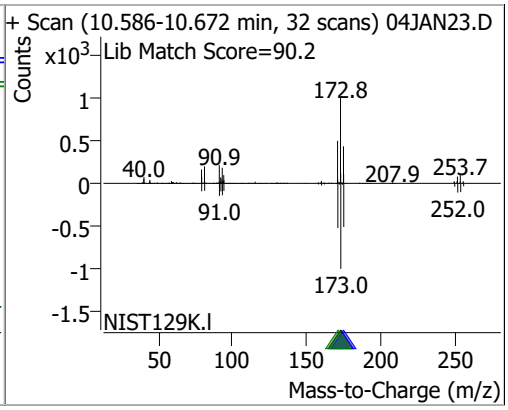




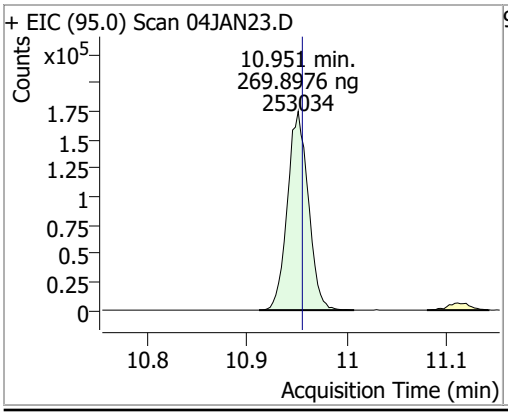
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	129.9644	10.63	0.00	42560	170.5	50.3	22.1	82.1
					174.5	49.3	20.1	80.1

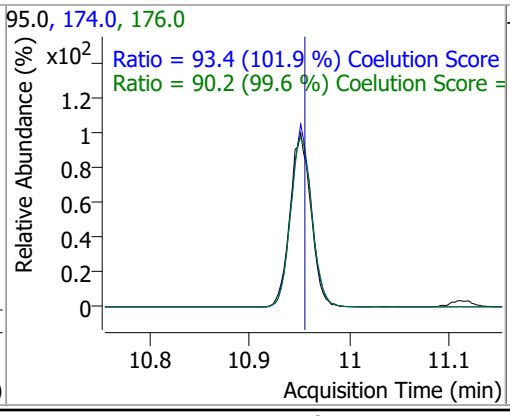


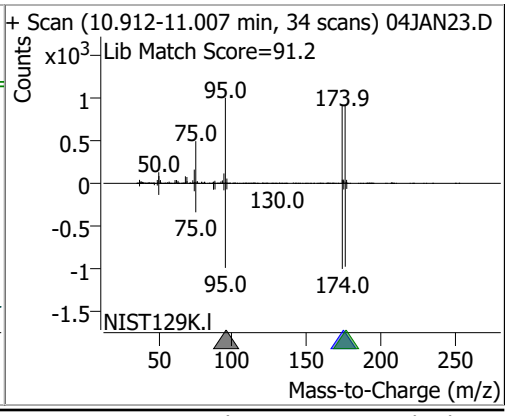




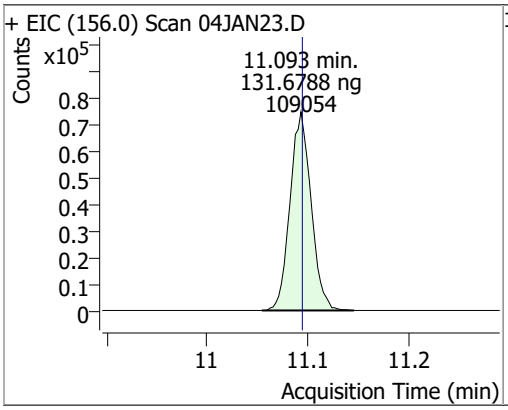
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	269.8976	10.95	0.00	253034	174.0	93.4	61.7	121.7
					176.0	90.2	60.6	120.6

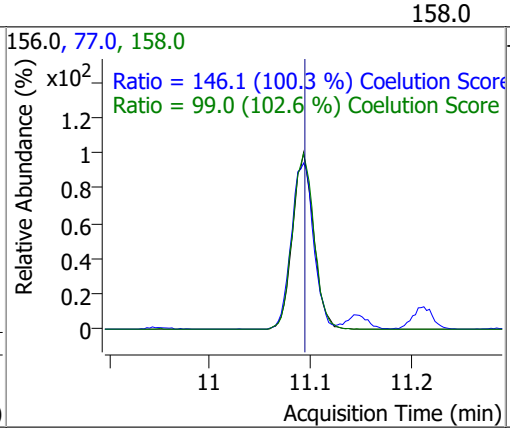


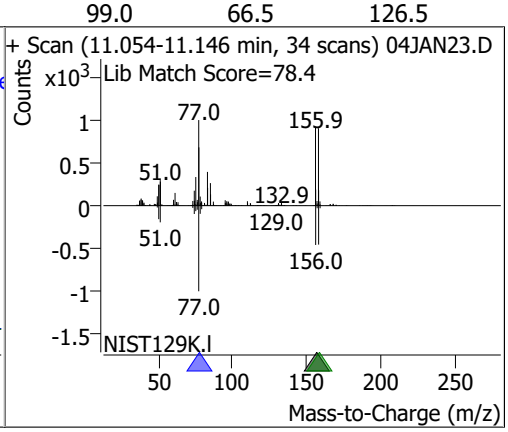




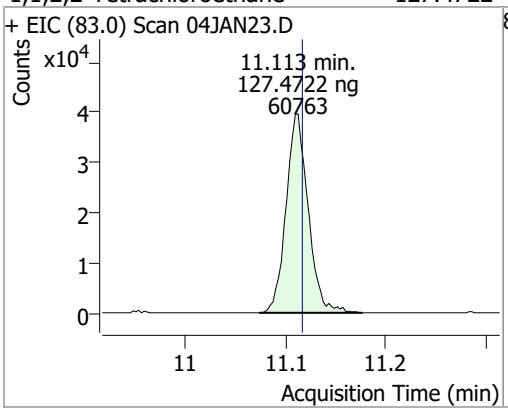
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	131.6788	11.09	0.00	109054	77.0	146.1	115.7	175.7
					158.0	99.0	66.5	126.5

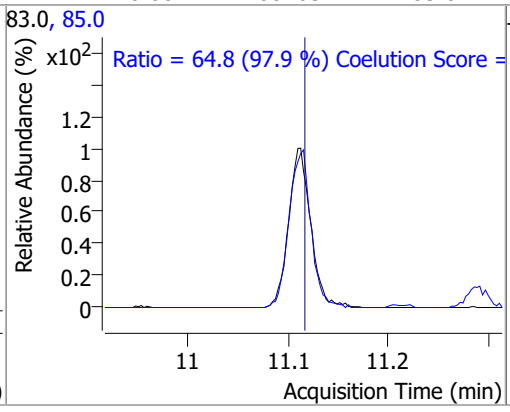


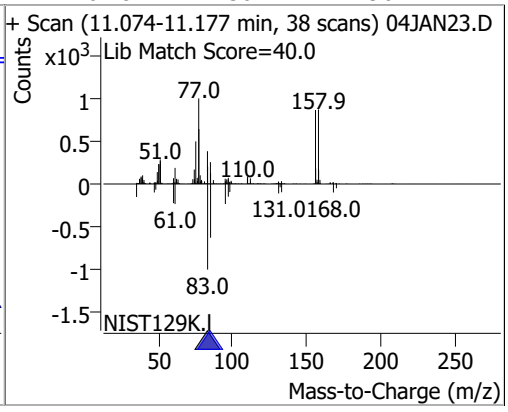




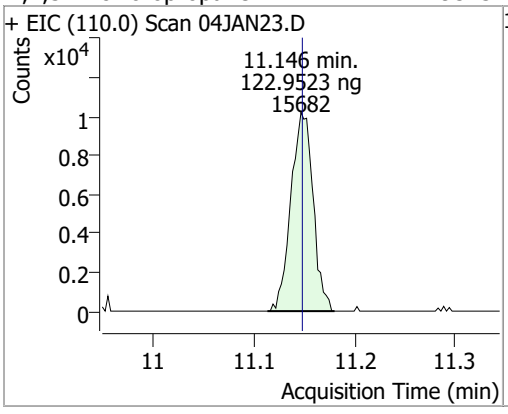
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	127.4722	11.11	0.00	60763	85.0	64.8	36.2	96.2

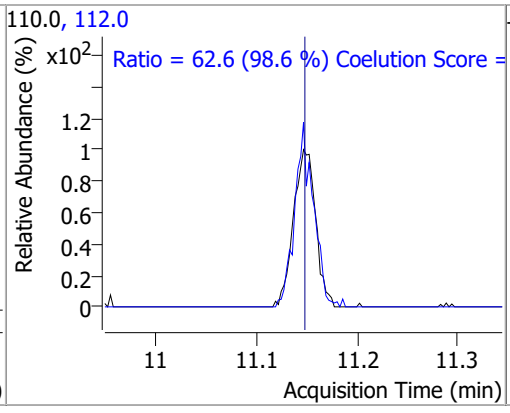


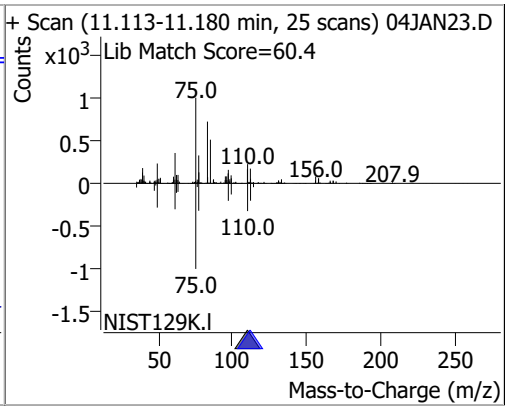




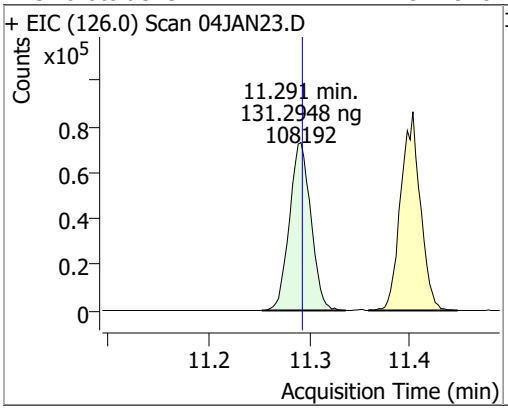
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	122.9523	11.15	0.00	15682	112.0	62.6	33.5	93.5

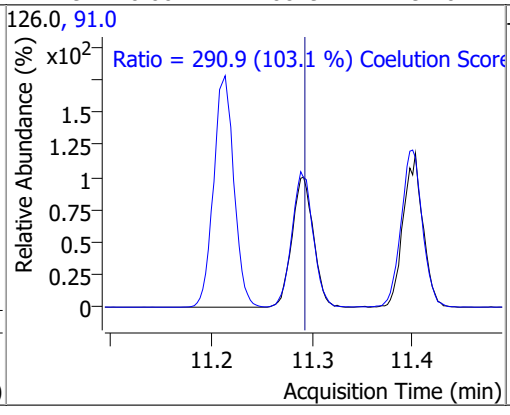


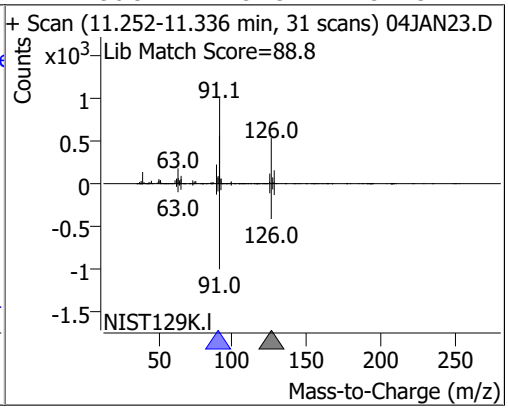




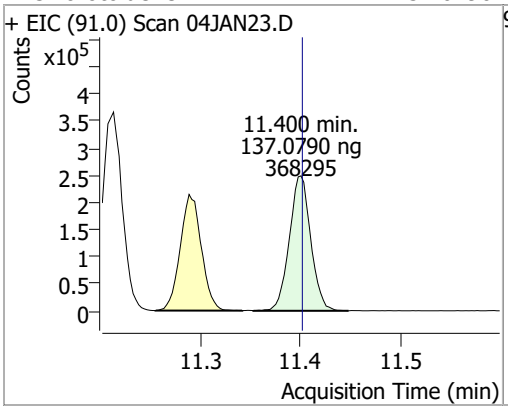
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	131.2948	11.29	0.00	108192	91.0	290.9	252.3	312.3

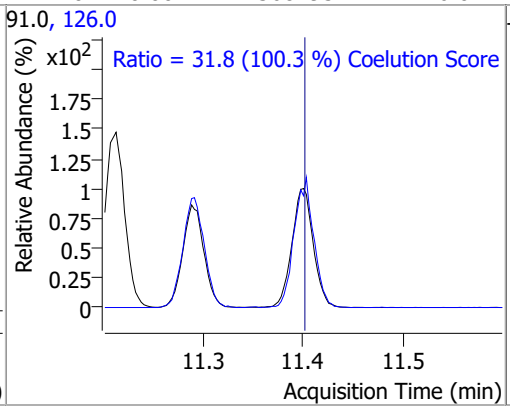


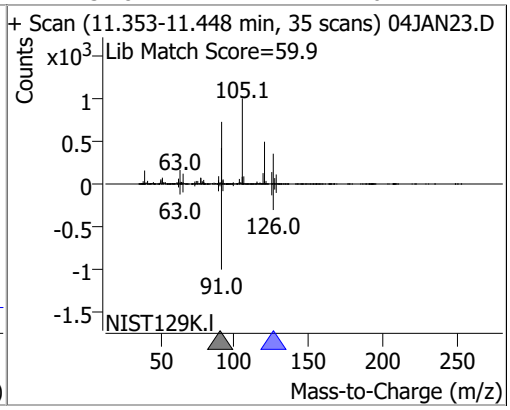




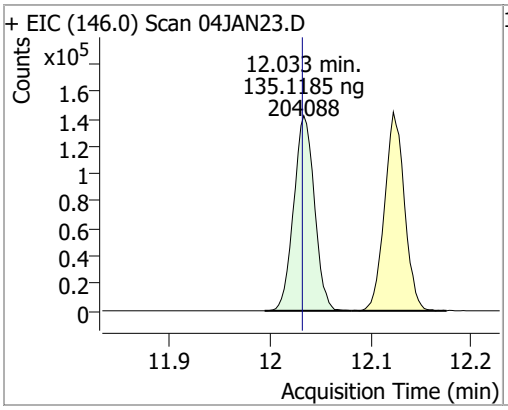
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	137.0790	11.40	0.00	368295	126.0	31.8	1.7	61.7

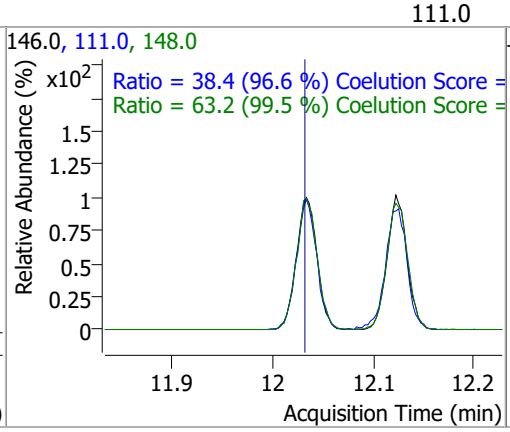


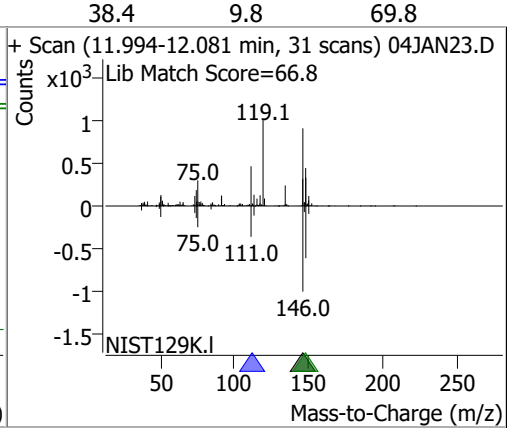




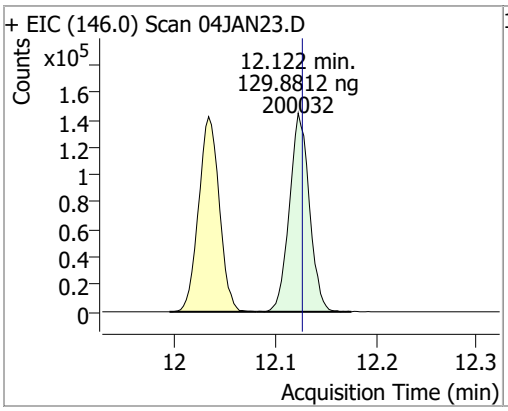
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	135.1185	12.03	0.00	204088	148.0	63.2	33.6	93.6
					111.0	38.4	9.8	69.8

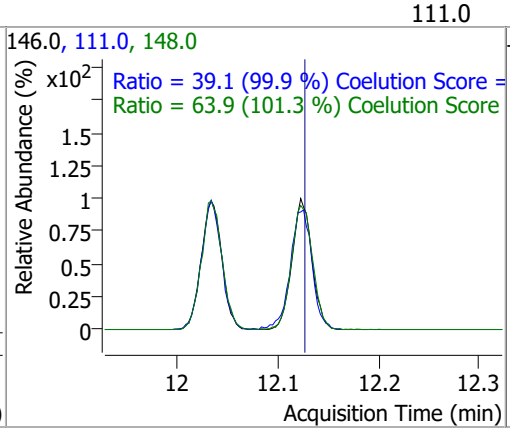


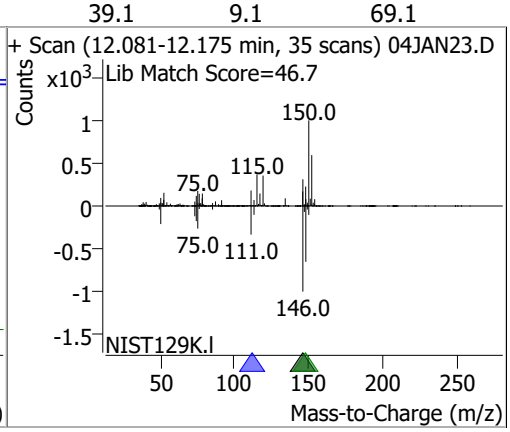




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	129.8812	12.12	0.00	200032	148.0	63.9	33.1	93.1
					111.0	39.1	9.1	69.1

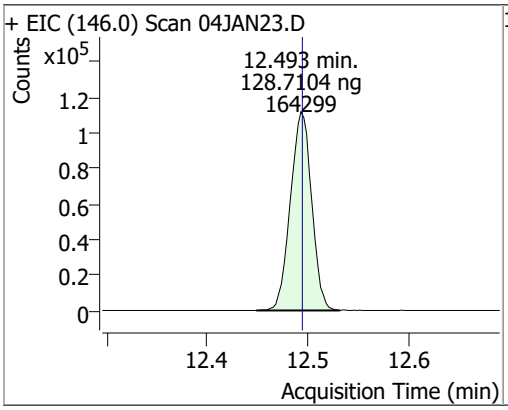


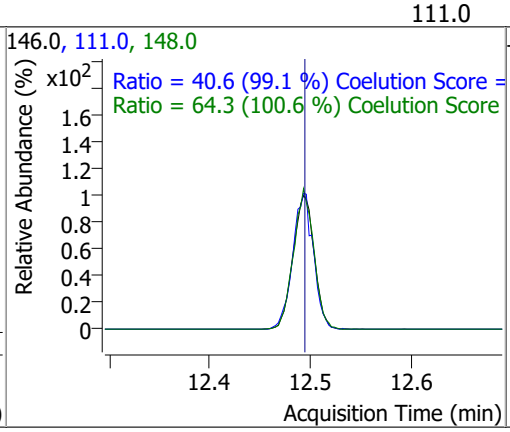


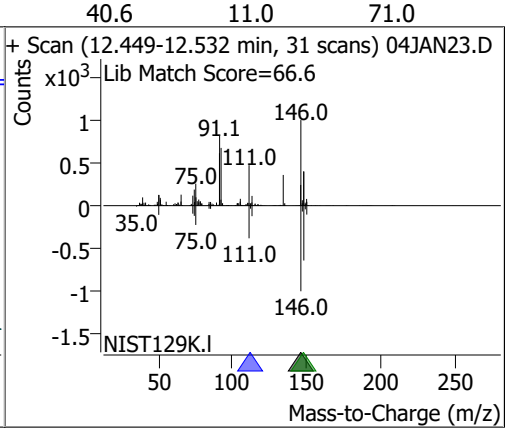


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	128.7104	12.49	0.00	164299	148.0	64.3	33.9	93.9
					111.0	40.6	11.0	71.0







Audit Trail report

Batch name and path: D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin
Quant batch version: 10.0
Quant reporting version: 10.0

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdNewBatchTable	BL2000\mchavez	1/4/2022 10:36:43 AM	Create new batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/4/2022 10:36:56 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN02.D, D:\Org\Data\VOA5975C\VG010422\04JAN01.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 10:37:01 AM	Set SampleType = MatrixBlank for sample 04JAN02.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 10:37:04 AM	Set SampleType = TuneCheck for sample 04JAN02.D; previous value = MatrixBlank			✓	
CmdStartMethodEditing	BL2000\mchavez	1/4/2022 10:52:58 AM	Start method editing			✓	
CmdImportMethodFromBatch	BL2000\mchavez	1/4/2022 10:52:59 AM	Import method from batch D:\Org\Data\VOA5975C\VG010322\VG010322_8260B_2ndRun.batch.bin			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/4/2022 10:53:03 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/4/2022 10:53:03 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/4/2022 10:53:03 AM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/4/2022 10:53:07 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/4/2022 10:54:55 AM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/4/2022 11:14:07 AM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/4/2022 11:14:24 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN03.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 11:14:28 AM	Set SampleType = CC for sample 04JAN03.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 11:14:31 AM	Set LevelName = CC for sample 04JAN03.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/4/2022 11:14:35 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/4/2022 11:15:38 AM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/4/2022 1:02:39 PM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\mchavez	1/4/2022 1:02:58 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN05.D, D:\Org\Data\VOA5975C\VG010422\04JAN04.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 1:03:09 PM	Set SampleType = TuneCheck for sample 04JAN05.D; previous value = Sample			✓	
CmdQuantitate	BL2000\mchavez	1/4/2022 1:03:40 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/4/2022 1:14:26 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/4/2022 1:47:13 PM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/4/2022 1:47:29 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN06.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 1:47:35 PM	Set SampleType = CC for sample 04JAN06.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 1:47:38 PM	Set LevelName = CC for sample 04JAN06.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/4/2022 1:47:44 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/4/2022 1:53:45 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/4/2022 3:05:35 PM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/4/2022 3:06:14 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN08.D, D:\Org\Data\VOA5975C\VG010422\04JAN07.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/4/2022 3:06:26 PM	Set SampleType = TuneCheck for sample 04JAN08.D; previous value = Sample			✓	
CmdSaveBatchTable	BL2000\mchavez	1/4/2022 3:24:14 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/5/2022 8:56:03 AM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\mchavez	1/5/2022 8:58:43 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN28.D, D:\Org\Data\VOA5975C\VG010422\04JAN27.D, D:\Org\Data\VOA5975C\VG010422\04JAN26.D, D:\Org\Data\VOA5975C\VG010422\04JAN25.D, D:\Org\Data\VOA5975C\VG010422\04JAN24.D, D:\Org\Data\VOA5975C\VG010422\04JAN23.D, D:\Org\Data\VOA5975C\VG010422\04JAN22.D, D:\Org\Data\VOA5975C\VG010422\04JAN21.D, D:\Org\Data\VOA5975C\VG010422\04JAN20.D, D:\Org\Data\VOA5975C\VG010422\04JAN19.D, D:\Org\Data\VOA5975C\VG010422\04JAN18.D, D:\Org\Data\VOA5975C\VG010422\04JAN17.D, D:\Org\Data\VOA5975C\VG010422\04JAN16.D, D:\Org\Data\VOA5975C\VG010422\04JAN15.D, D:\Org\Data\VOA5975C\VG010422\04JAN14.D, D:\Org\Data\VOA5975C\VG010422\04JAN13.D, D:\Org\Data\VOA5975C\VG010422\04JAN12.D, D:\Org\Data\VOA5975C\VG010422\04JAN11.D, D:\Org\Data\VOA5975C\VG010422\04JAN10.D, D:\Org\Data\VOA5975C\VG010422\04JAN09.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:26 AM	Set SampleType = Blank for sample 04JAN09.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:31 AM	Set SampleType = Calibration for sample 04JAN10.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:35 AM	Set SampleType = Calibration for sample 04JAN11.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:40 AM	Set SampleType = Calibration for sample 04JAN12.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:45 AM	Set SampleType = Calibration for sample 04JAN13.D; previous value = Sample			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:50 AM	Set SampleType = Calibration for sample 04JAN15.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 8:59:55 AM	Set SampleType = Calibration for sample 04JAN17.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 9:00:00 AM	Set SampleType = Calibration for sample 04JAN19.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 9:00:08 AM	Set SampleType = Calibration for sample 04JAN21.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 9:00:14 AM	Set SampleType = QC for sample 04JAN23.D; previous value = Sample			✓	
CmdSaveBatchTable	BL2000\mchavez	1/5/2022 9:11:16 AM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/5/2022 9:58:34 AM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdQuantitate	BL2000\mchavez	1/5/2022 9:58:40 AM	Quantitate all compounds in all samples				Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Batch quantitation failed ---> Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Sample not validated: Level name is undefined for a Calibration or QC sample. at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.Batch.ValidateBatchMethod() at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.Batch.Quantitate() --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.Batch.Quantitate() at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdQuantitate.QuantitateBatch(Int16 batchId) at Agilent.MassSpectrometry.CommandModel.CommandHistory.Invoke(ICommand cmd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.AppCommandContext._Invoke(ICommand cmd)

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:06:05 AM	Set LevelName = 1 for sample 04JAN10.D; previous value =			✓	
CmdSaveBatchTable	BL2000\mchavez	1/5/2022 10:06:09 AM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/5/2022 10:14:38 AM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:14:47 AM	Set LevelName = 2 for sample 04JAN11.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:14:52 AM	Set LevelName = 3 for sample 04JAN12.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:14:57 AM	Set LevelName = 4 for sample 04JAN13.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:15:05 AM	Set LevelName = 5 for sample 04JAN15.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:15:11 AM	Set LevelName = 6 for sample 04JAN17.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:15:16 AM	Set LevelName = 7 for sample 04JAN19.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:15:21 AM	Set LevelName = 8 for sample 04JAN21.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:15:29 AM	Set LevelName = QC for sample 04JAN23.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:15:36 AM	Set SampleInformation = LCSA for sample 04JAN23.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/5/2022 10:16:01 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:16:49 AM	Manually integrate compound Chloroethane in sample 04JAN10.D, from x, y = 1.874, 1384 to 1.916, 1542, result = 2178; previous integration is from x, y = 1.894, 1143 to 1.933, 1143 and previous response = 2132.			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:16:53 AM	Manually integrate qualifier66.0 of compound Chloroethane in sample 04JAN10.D from x, y = 1.869, 0 to 1.908, 8; result = 781			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:16:54 AM	Manually integrate qualifier 66.0 of compound Chloroethane in sample 04JAN10.D, from x, y = 1.869, 0 to 1.913, 0, result = 824; previous integration is from x, y = 1.869, 0 to 1.908, 8 and previous response = 781.			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:17:08 AM	Manually integrate compound Bromomethane in sample 04JAN10.D from x, y = 1.768, -2 to 1.849, 0; result = 1902			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:17:15 AM	Manually integrate qualifier87.0 of compound Dichlorodifluoromethane in sample 04JAN10.D from x, y = 1.202, 0 to 1.289, 0; result = 1393			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:17:25 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 04JAN10.D from x, y = 1.370, 0 to 1.459, 0; result = 1679			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:17:55 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN10.D, from x, y = 1.492, 6362 to 1.506, 5900, result = 1131; previous integration is from x, y = 1.308, 0 to 1.682, 0 and previous response = 192320.			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:18:11 AM	Manually integrate compound 1,1-Dichloroethene in sample 04JAN10.D from x, y = 2.636, 0 to 2.747, 0; result = 2084			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:18:16 AM	Manually integrate qualifier63.0 of compound 1,1-Dichloroethene in sample 04JAN10.D from x, y = 2.672, 0 to 2.753, 0; result = 1158			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:18:38 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 04JAN10.D from x, y = 3.288, 0 to 3.386, 0; result = 1820			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:19:27 AM	Manually integrate compound trans-1,2-Dichloroethene in sample 04JAN10.D from x, y = 3.673, 0 to 3.762, 0; result = 2146			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:19:31 AM	Manually integrate qualifier98.0 of compound trans-1,2-Dichloroethene in sample 04JAN10.D from x, y = 3.684, 0 to 3.779, 0; result = 1426			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:19:36 AM	Manually integrate compound Methyl tert-butyl ether (MTBE) in sample 04JAN10.D from x, y = 3.698, 0 to 3.832, 0; result = 2717			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:19:38 AM	Manually integrate qualifier57.0 of compound Methyl tert-butyl ether (MTBE) in sample 04JAN10.D from x, y = 3.690, 0 to 3.798, 0; result = 531			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:19:46 AM	Manually integrate compound 2,2-Dichloropropane in sample 04JAN10.D from x, y = 5.137, 0 to 5.279, 0; result = 2930			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:19:49 AM	Manually integrate qualifier97.0 of compound 2,2-Dichloropropane in sample 04JAN10.D from x, y = 5.151, 0 to 5.285, 0; result = 814			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:19:52 AM	Manually integrate qualifier41.0 of compound 2,2-Dichloropropane in sample 04JAN10.D from x, y = 5.151, 0 to 5.257, 0; result = 2246			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:20:11 AM	Manually integrate qualifier65.0 of compound 1,1-Dichloroethane in sample 04JAN10.D from x, y = 4.323, 0 to 4.465, 0; result = 1347			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:20:13 AM	Manually integrate qualifier83.0 of compound 1,1-Dichloroethane in sample 04JAN10.D from x, y = 4.342, 0 to 4.426, 0; result = 227			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:26:02 AM	Manually integrate compound cis-1,2-Dichloroethene in sample 04JAN10.D from x, y = 5.145, 0 to 5.282, 0; result = 2376			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:05 AM	Manually integrate qualifier61.0 of compound cis-1,2-Dichloroethene in sample 04JAN10.D from x, y = 5.151, 0 to 5.274, 0; result = 4139			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:07 AM	Manually integrate qualifier98.0 of compound cis-1,2-Dichloroethene in sample 04JAN10.D from x, y = 5.176, 0 to 5.257, 0; result = 1525			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:12 AM	Manually integrate qualifier72.0 of compound Methyl ethyl ketone in sample 04JAN10.D from x, y = 5.274, 0 to 5.343, 0; result = 435			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:26:18 AM	Manually integrate compound Bromochloromethane in sample 04JAN10.D from x, y = 5.463, 0 to 5.555, 0; result = 807			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:20 AM	Manually integrate qualifier49.0 of compound Bromochloromethane in sample 04JAN10.D from x, y = 5.472, 0 to 5.558, 0; result = 1686			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:23 AM	Manually integrate qualifier85.0 of compound Chloroform in sample 04JAN10.D from x, y = 5.592, 0 to 5.734, 0; result = 2708			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:26:28 AM	Manually integrate compound Dibromofluoromethane in sample 04JAN10.D from x, y = 5.812, 0 to 5.915, 0; result = 2508			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:30 AM	Manually integrate qualifier191.5 of compound Dibromofluoromethane in sample 04JAN10.D from x, y = 5.809, 0 to 5.884, 0; result = 479			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:37 AM	Manually integrate qualifier61.0 of compound 1,1,1-Trichloroethane in sample 04JAN10.D from x, y = 5.790, 0 to 5.890, 0; result = 1705			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:42 AM	Manually integrate qualifier121.0 of compound Carbon tetrachloride in sample 04JAN10.D from x, y = 5.999, 0 to 6.068, 0; result = 903			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:46 AM	Manually integrate qualifier110.0 of compound 1,1-Dichloropropene in sample 04JAN10.D from x, y = 5.993, 0 to 6.074, 0; result = 1122			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:26:48 AM	Manually integrate qualifier77.0 of compound 1,1-Dichloropropene in sample 04JAN10.D from x, y = 6.013, 0 to 6.099, 0; result = 1052			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:26:53 AM	Manually integrate compound 1,2-Dichloroethane-d4 in sample 04JAN10.D from x, y = 6.188, -35 to 6.283, 0; result = 1023			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:26:58 AM	Manually integrate compound 1,2-Dichloroethane-d4 in sample 04JAN10.D, from x, y = 6.197, 0 to 6.283, 0, result = 923; previous integration is from x, y = 6.188, -35 to 6.283, 0 and previous response = 1023.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:00 AM	Manually integrate qualifier65.0 of compound 1,2-Dichloroethane-d4 in sample 04JAN10.D from x, y = 6.199, 0 to 6.275, 0; result = 1927			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:34 AM	Manually integrate qualifier77.0 of compound Benzene in sample 04JAN10.D from x, y = 6.222, 0 to 6.339, 0; result = 1884			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:27:39 AM	Manually integrate compound 1,2-Dichloroethane in sample 04JAN10.D from x, y = 6.269, 0 to 6.386, 0; result = 2415			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:41 AM	Manually integrate qualifier64.0 of compound 1,2-Dichloroethane in sample 04JAN10.D from x, y = 6.280, 0 to 6.378, 0; result = 761			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:43 AM	Manually integrate qualifier98.0 of compound 1,2-Dichloroethane in sample 04JAN10.D from x, y = 6.303, 0 to 6.386, 0; result = 119			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:27:47 AM	Manually integrate compound Trichloroethene in sample 04JAN10.D from x, y = 6.989, 0 to 7.083, 0; result = 2372			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:51 AM	Manually integrate qualifier 130.0 of compound Trichloroethene in sample 04JAN10.D, from x, y = 6.997, 0 to 7.072, 0, result = 2567; previous integration is from x, y = 6.997, 0 to 7.044, 0 and previous response = 2405.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:53 AM	Manually integrate qualifier97.0 of compound Trichloroethene in sample 04JAN10.D from x, y = 6.991, 0 to 7.078, 0; result = 1659			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:27:59 AM	Manually integrate qualifier76.0 of compound 1,2-Dichloropropane in sample 04JAN10.D from x, y = 7.231, 0 to 7.321, 0; result = 733			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:28:03 AM	Manually integrate compound Dibromomethane in sample 04JAN10.D from x, y = 7.357, 0 to 7.424, 0; result = 902			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:06 AM	Manually integrate qualifier95.0 of compound Dibromomethane in sample 04JAN10.D from x, y = 7.357, 0 to 7.454, 0; result = 535			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:09 AM	Manually integrate qualifier173.5 of compound Dibromomethane in sample 04JAN10.D from x, y = 7.360, 0 to 7.429, 0; result = 1002			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:15 AM	Manually integrate qualifier85.0 of compound Bromodichloromethane in sample 04JAN10.D from x, y = 7.546, 0 to 7.633, 0; result = 1631			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:17 AM	Manually integrate qualifier127.0 of compound Bromodichloromethane in sample 04JAN10.D from x, y = 7.569, 0 to 7.633, 0; result = 130			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:23 AM	Manually integrate qualifier77.0 of compound cis-1,3-Dichloropropene in sample 04JAN10.D from x, y = 8.009, 0 to 8.107, 0; result = 922			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:35 AM	Manually integrate qualifier39.0 of compound cis-1,3-Dichloropropene in sample 04JAN10.D from x, y = 8.037, 0 to 8.090, 0; result = 1459			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:40 AM	Manually integrate qualifier99.0 of compound Toluene-d8 in sample 04JAN10.D from x, y = 8.271, 0 to 8.350, 0; result = 699			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:28:47 AM	Manually integrate compound trans-1,3-Dichloropropene in sample 04JAN10.D from x, y = 8.614, 0 to 8.684, 0; result = 1470			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:28:50 AM	Manually integrate qualifier 77.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D, from x, y = 8.601, 0 to 8.656, -10, result = 8023; previous integration is from x, y = 8.656, 57 to 8.709, 135 and previous response = 3575.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:29:03 AM	Manually integrate qualifier 77.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D, from x, y = 8.601, 0 to 8.656, 156, result = 6858; previous integration is from x, y = 8.601, 0 to 8.835, -10 and previous response = 8023.			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/5/2022 10:29:04 AM	Drop baseline for qualifier 77.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D to y = 0, new integration is from x, y = 8.601, 0 to 8.835, 0 and new response = 7954; previous integration is from x, y = 8.601, 0 to 8.835, 156 and previous response = 6858.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:29:09 AM	Manually integrate qualifier 77.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D, from x, y = 8.601, 0 to 8.648, 141, result = 368; previous integration is from x, y = 8.601, 0 to 8.835, 0 and previous response = 7954.			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/5/2022 10:29:12 AM	Drop baseline for qualifier 77.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D to y = 0, new integration is from x, y = 8.601, 0 to 8.648, 0 and new response = 568; previous integration is from x, y = 8.601, 0 to 8.648, 141 and previous response = 368.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:29:19 AM	Manually integrate qualifier39.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D from x, y = 8.614, 0 to 8.648, 26; result = 974			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/5/2022 10:29:21 AM	Drop baseline for qualifier 39.0 of compound trans-1,3-Dichloropropene in sample 04JAN10.D to y = 0, new integration is from x, y = 8.614, 0 to 8.648, 0 and new response = 1000; previous integration is from x, y = 8.614, 0 to 8.648, 26 and previous response = 974.			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:29:29 AM	Manually integrate compound 1,1,2-Trichloroethane in sample 04JAN10.D from x, y = 8.785, 0 to 8.843, 0; result = 960			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:29:31 AM	Manually integrate qualifier97.0 of compound 1,1,2-Trichloroethane in sample 04JAN10.D from x, y = 8.782, 0 to 8.857, 0; result = 1099			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:29:34 AM	Manually integrate qualifier85.0 of compound 1,1,2-Trichloroethane in sample 04JAN10.D from x, y = 8.796, 0 to 8.851, 0; result = 418			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:29:39 AM	Manually integrate compound Tetrachloroethene in sample 04JAN10.D from x, y = 8.899, 0 to 8.983, 0; result = 2105			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:29:43 AM	Manually integrate qualifier165.8 of compound Tetrachloroethene in sample 04JAN10.D from x, y = 8.885, 0 to 9.019, 0; result = 2853			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:29:48 AM	Manually integrate qualifier78.0 of compound 1,3-Dichloropropane in sample 04JAN10.D from x, y = 8.952, 0 to 9.010, 0; result = 452			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:29:52 AM	Manually integrate compound Chlorodibromomethane in sample 04JAN10.D from x, y = 9.169, 0 to 9.256, 0; result = 1468			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:29:54 AM	Manually integrate qualifier127.0 of compound Chlorodibromomethane in sample 04JAN10.D from x, y = 9.175, 0 to 9.242, 0; result = 1140			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:29:58 AM	Manually integrate compound 1,2-Dibromoethane in sample 04JAN10.D from x, y = 9.278, 0 to 9.348, 0; result = 1299			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:30:00 AM	Manually integrate qualifier109.0 of compound 1,2-Dibromoethane in sample 04JAN10.D from x, y = 9.275, 0 to 9.340, 0; result = 1039			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:30:35 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN15.D, from x, y = 1.476, 4348 to 1.540, 3059, result = 44384; previous integration is from x, y = 1.478, 954 to 1.687, 2147 and previous response = 61198.			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:35:03 AM	Set SampleApproved = True for sample 04JAN15.D; previous value = False			✓	
CmdStartMethodEditing	BL2000\mchavez	1/5/2022 10:35:10 AM	Start method editing			✓	
CmdImportMethodFromSample	BL2000\mchavez	1/5/2022 10:35:10 AM	Import method from sample 04JAN15.D			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdUpdateRetentionTimes	BL2000\mchavez	1/5/2022 10:35:27 AM	Update retention time for compound 1,4-Dichlorobenzene; 1,3-Dichlorobenzene; 4-Chlorotoluene; 2-Chlorotoluene; 1,2,3-Trichloropropane; Bromobenzene; 1,1,2,2-Tetrachloroethane; p-Bromofluorobenzene; Bromoform; Styrene; o-Xylene; m+p-Xylenes; Ethylbenzene; 1,1,1,2-Tetrachloroethane; Chlorobenzene; 1,2-Dibromoethane; Chlorodibromomethane; 1,3-Dichloropropane; Tetrachloroethene; 1,1,2-Trichloroethane; trans-1,3-Dichloropropene; Toluene; Toluene-d8; cis-1,3-Dichloropropene; Bromodichloromethane; Dibromomethane; 1,2-Dichloropropane; Trichloroethene; 1,2-Dichloroethane; Benzene; 1,2-Dichloroethane-d4; 1,1-Dichloropropene; Carbon tetrachloride; 1,1,1-Trichloroethane; Dibromofluoromethane; Chloroform; Bromochloromethane; Methyl ethyl ketone; cis-1,2-Dichloroethene; 2,2-Dichloropropane; 1,1-Dichloroethane; Methyl tert-butyl ether (MTBE); trans-1,2-Dichloroethene; Methylene chloride; 1,1-Dichloroethene; Trichlorofluoromethane; Chloroethane; Bromomethane; Vinyl chloride; Chloromethane; Dichlorodifluoromethane; 1,4-Dichlorobenzene-d4; Chlorobenzene-d5; Fluorobenzene; 1,2-Dichlorobenzene;			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdUpdateQualifierRatios	BL2000\mchavez	1/5/2022 10:35:32 AM	Update qualifier ratios for compound 1,4-Dichlorobenzene; Update qualifier ratios for compound 1,3-Dichlorobenzene; Update qualifier ratios for compound 4-Chlorotoluene; Update qualifier ratios for compound 2-Chlorotoluene; Update qualifier ratios for compound 1,2,3-Trichloropropane; Update qualifier ratios for compound Bromobenzene; Update qualifier ratios for compound 1,1,2,2-Tetrachloroethane; Update qualifier ratios for compound p-Bromofluorobenzene; Update qualifier ratios for compound Bromoform; Update qualifier ratios for compound Styrene; Update qualifier ratios for compound o-Xylene; Update qualifier ratios for compound m+p-Xylenes; Update qualifier ratios for compound Ethylbenzene; Update qualifier ratios for compound 1,1,1,2-Tetrachloroethane; Update qualifier ratios for compound Chlorobenzene; Update qualifier ratios for compound 1,2-Dibromoethane; Update qualifier ratios for compound Chlorodibromomethane; Update qualifier ratios for compound 1,3-Dichloropropane; Update qualifier ratios for compound Tetrachloroethene; Update qualifier ratios for compound 1,1,2-Trichloroethane; Update qualifier ratios for compound trans-1,3-Dichloropropene; Update qualifier ratios for compound Toluene; Update qualifier ratios for compound Toluene-d8; Update qualifier ratios for compound cis-1,3-Dichloropropene; Update qualifier ratios for compound Bromodichloromethane; Update qualifier ratios for compound Dibromomethane; Update qualifier ratios for compound 1,2-Dichloropropane; Update qualifier ratios for compound Trichloroethene; Update qualifier ratios for compound 1,2-Dichloroethane; Update qualifier ratios for compound Benzene; Update qualifier ratios for compound 1,2-Dichloroethane-d4; Update qualifier ratios for compound 1,1-Dichloropropene; Update qualifier ratios for compound Carbon tetrachloride; Update qualifier ratios for compound 1,1,1-Trichloroethane; Update qualifier ratios for compound Dibromofluoromethane; Update			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			qualifier ratios for compound Chloroform; Update qualifier ratios for compound Bromochloromethane; Update qualifier ratios for compound Methyl ethyl ketone; Update qualifier ratios for compound cis-1,2-Dichloroethene; Update qualifier ratios for compound 2,2-Dichloropropane; Update qualifier ratios for compound 1,1-Dichloroethane; Update qualifier ratios for compound Methyl tert-butyl ether (MTBE); Update qualifier ratios for compound trans-1,2-Dichloroethene; Update qualifier ratios for compound Methylene chloride; Update qualifier ratios for compound 1,1-Dichloroethene; Update qualifier ratios for compound Trichlorofluoromethane; Update qualifier ratios for compound Chloroethane; Update qualifier ratios for compound Bromomethane; Update qualifier ratios for compound Vinyl chloride; Update qualifier ratios for compound Chloromethane; Update qualifier ratios for compound Dichlorodifluoromethane; Update qualifier ratios for compound 1,4-Dichlorobenzene-d4; Update qualifier ratios for compound Chlorobenzene-d5; Update qualifier ratios for compound Fluorobenzene; Update qualifier ratios for compound 1,2-Dichlorobenzene;				
CmdApplyMethodToAllSamples	BL2000\mchavez	1/5/2022 10:35:43 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/5/2022 10:35:43 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/5/2022 10:35:43 AM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/5/2022 10:36:02 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:37:04 AM	Manually integrate compound 2-Chlorotoluene in sample 04JAN10.D from x, y = 11.241, 0 to 11.353, 0; result = 1844			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:37:15 AM	Manually integrate compound Bromobenzene in sample 04JAN10.D from x, y = 11.049, 0 to 11.127, 0; result = 2024			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:37:18 AM	Manually integrate qualifier158.0 of compound Bromobenzene in sample 04JAN10.D from x, y = 11.063, 0 to 11.152, 0; result = 1934			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:37:23 AM	Manually integrate compound 1,1,2,2-Tetrachloroethane in sample 04JAN10.D from x, y = 11.085, 0 to 11.188, 0; result = 1142			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:37:25 AM	Manually integrate qualifier85.0 of compound 1,1,2,2-Tetrachloroethane in sample 04JAN10.D from x, y = 11.071, 0 to 11.147, 0; result = 834			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:37:35 AM	Manually integrate compound Bromoform in sample 04JAN10.D from x, y = 10.597, 0 to 10.686, 0; result = 708			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:37:36 AM	Manually integrate qualifier174.5 of compound Bromoform in sample 04JAN10.D from x, y = 10.594, 0 to 10.698, 0; result = 258			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:37:38 AM	Manually integrate qualifier170.5 of compound Bromoform in sample 04JAN10.D from x, y = 10.603, 0 to 10.672, 0; result = 339			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:37:45 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,1,2,2-Tetrachloroethane in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:37:53 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 2-Chlorotoluene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:37:55 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 4-Chlorotoluene in sample 04JAN15.D; previous value = True			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:37:58 AM	Manually integrate qualifier126.0 of compound 4-Chlorotoluene in sample 04JAN10.D from x, y = 11.367, 0 to 11.467, 0; result = 1839			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:38:03 AM	Manually integrate qualifier111.0 of compound 1,3-Dichlorobenzene in sample 04JAN10.D from x, y = 12.000, 0 to 12.061, 0; result = 1469			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:05 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,3-Dichlorobenzene in sample 04JAN15.D; previous value = True			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:38:13 AM	Manually integrate qualifier111.0 of compound 1,2-Dichlorobenzene in sample 04JAN10.D from x, y = 12.432, 0 to 12.538, 0; result = 1190			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:15 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,2-Dichlorobenzene in sample 04JAN15.D; previous value = True			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:19 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,4-Dichlorobenzene in sample 04JAN15.D; previous value = True			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:38:25 AM	Manually integrate qualifier148.0 of compound 1,2-Dichlorobenzene in sample 04JAN10.D from x, y = 12.468, 0 to 12.555, 0; result = 1894			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:44 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Styrene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:48 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound o-Xylene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:51 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound m+p-Xylenes in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:38:58 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Ethylbenzene in sample 04JAN15.D; previous value = True			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:40:02 AM	Manually integrate qualifier 106.0 of compound Ethylbenzene in sample 04JAN10.D, from x, y = 9.883, 0 to 9.953, 0, result = 3266; previous integration is from x, y = 9.914, 0 to 9.953, 0 and previous response = 2097.			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 10:40:07 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample 04JAN10.D from x, y = 9.853, 0 to 9.939, 0; result = 1893			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:40:11 AM	Manually integrate qualifier133.0 of compound 1,1,1,2-Tetrachloroethane in sample 04JAN10.D from x, y = 9.841, 0 to 9.931, 0; result = 1911			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 10:40:23 AM	Manually integrate qualifier114.0 of compound Chlorobenzene in sample 04JAN10.D from x, y = 9.783, 18 to 9.844, 0; result = 1827			✓	
CmdManuallyIntegrateDropBaseline	BL2000\mchavez	1/5/2022 10:40:24 AM	Drop baseline for qualifier 114.0 of compound Chlorobenzene in sample 04JAN10.D to y = 0, new integration is from x, y = 9.783, 0 to 9.844, 0 and new response = 1861; previous integration is from x, y = 9.783, 18 to 9.844, 0 and previous response = 1827.			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:41:20 AM	Set SampleApproved = True for sample 04JAN10.D; previous value = False			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:41:35 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN11.D, from x, y = 1.487, 5730 to 1.520, 4652, result = 5622; previous integration is from x, y = 1.311, 0 to 1.679, 0 and previous response = 180129.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:41:38 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN11.D, from x, y = 1.484, 5832 to 1.520, 4652, result = 5972; previous integration is from x, y = 1.487, 5730 to 1.520, 4652 and previous response = 5622.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:41:51 AM	Manually integrate qualifier 63.0 of compound 1,1-Dichloroethene in sample 04JAN11.D, from x, y = 2.663, 0 to 2.736, 0, result = 5268; previous integration is from x, y = 2.697, 0 to 2.736, 0 and previous response = 3135.			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:41:59 AM	Manually integrate compound trans-1,2-Dichloroethene in sample 04JAN11.D, from x, y = 3.678, 0 to 3.765, 0, result = 9821; previous integration is from x, y = 3.678, 0 to 3.718, 0 and previous response = 5041.			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:42:05 AM	Manually integrate compound Methyl tert-butyl ether (MTBE) in sample 04JAN11.D, from x, y = 3.687, 0 to 3.823, 182, result = 11769; previous integration is from x, y = 3.737, 0 to 3.804, 0 and previous response = 10323.			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/5/2022 10:42:08 AM	Drop baseline for compound Methyl tert-butyl ether (MTBE) in sample 04JAN11.D to y = 0, new integration is from x, y = 3.687, 0 to 3.823, 0 and new response = 12515; previous integration is from x, y = 3.687, 0 to 3.823, 182 and previous response = 11769.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:42:12 AM	Manually integrate qualifier 57.0 of compound Methyl tert-butyl ether (MTBE) in sample 04JAN11.D, from x, y = 3.681, 0 to 3.821, 0, result = 3045; previous integration is from x, y = 3.709, 0 to 3.776, 0 and previous response = 2643.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:42:20 AM	Manually integrate qualifier97.0 of compound 2,2-Dichloropropane in sample 04JAN11.D from x, y = 5.140, 0 to 5.240, 0; result = 3733			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:42:32 AM	Manually integrate qualifier191.5 of compound Dibromofluoromethane in sample 04JAN11.D from x, y = 5.809, 0 to 5.918, 0; result = 2020			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:42:45 AM	Manually integrate qualifier 110.0 of compound 1,1-Dichloropropene in sample 04JAN11.D, from x, y = 5.985, 0 to 6.077, 0, result = 5349; previous integration is from x, y = 6.035, 0 to 6.077, 0 and previous response = 3323.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:42:59 AM	Manually integrate qualifier98.0 of compound 1,2-Dichloroethane in sample 04JAN11.D from x, y = 6.275, 0 to 6.381, 0; result = 648			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:43:02 AM	Manually integrate qualifier 64.0 of compound 1,2-Dichloroethane in sample 04JAN11.D, from x, y = 6.266, 0 to 6.361, 0, result = 3587; previous integration is from x, y = 6.317, 0 to 6.361, 0 and previous response = 2405.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:43:11 AM	Manually integrate qualifier127.0 of compound Bromodichloromethane in sample 04JAN11.D from x, y = 7.549, 0 to 7.627, 0; result = 943			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:43:21 AM	Manually integrate compound 1,1,2-Trichloroethane in sample 04JAN11.D, from x, y = 8.768, 0 to 8.882, 0, result = 5090; previous integration is from x, y = 8.818, 0 to 8.851, 0 and previous response = 2437.			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:44:13 AM	Manually integrate compound 1,2,3-Trichloropropane in sample 04JAN11.D from x, y = 11.099, 0 to 11.174, 0; result = 1654			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:44:16 AM	Manually integrate qualifier112.0 of compound 1,2,3-Trichloropropane in sample 04JAN11.D from x, y = 11.096, 0 to 11.191, 0; result = 1059			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:44:25 AM	Set SampleApproved = True for sample 04JAN11.D; previous value = False			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/5/2022 10:44:48 AM	Manually integrate compound 1,2,3-Trichloropropane in sample 04JAN12.D, from x, y = 11.099, 0 to 11.180, 0, result = 3200; previous integration is from x, y = 11.141, 0 to 11.180, 0 and previous response = 2198.			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:45:22 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Chlorodibromomethane in sample 04JAN15.D; previous value = True			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:45:30 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Tetrachloroethene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:45:37 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,1,2-Trichloroethane in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:45:40 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound trans-1,3-Dichloropropene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:45:45 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Toluene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:45:57 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound cis-1,3-Dichloropropene in sample 04JAN15.D; previous value = True			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:46:03 AM	Manually integrate qualifier127.0 of compound Bromodichloromethane in sample 04JAN12.D from x, y = 7.541, 0 to 7.633, 0; result = 2111			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:46:15 AM	Manually integrate qualifier98.0 of compound 1,2-Dichloroethane in sample 04JAN12.D from x, y = 6.278, 83 to 6.370, 0; result = 1440			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/5/2022 10:46:16 AM	Drop baseline for qualifier 98.0 of compound 1,2-Dichloroethane in sample 04JAN12.D to y = 0, new integration is from x, y = 6.278, 0 to 6.370, 0 and new response = 1669; previous integration is from x, y = 6.278, 83 to 6.370, 0 and previous response = 1440.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:46:38 AM	Manually integrate qualifier 97.0 of compound 2,2-Dichloropropane in sample 04JAN12.D, from x, y = 5.143, 0 to 5.254, 0, result = 6975; previous integration is from x, y = 5.187, 0 to 5.229, 0 and previous response = 4210.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:46:55 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN12.D, from x, y = 1.473, 6379 to 1.526, 4265, result = 8175; previous integration is from x, y = 1.305, 0 to 1.676, 0 and previous response = 154800.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:47:03 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN12.D, from x, y = 1.467, 4922 to 1.529, 3954, result = 11779; previous integration is from x, y = 1.473, 6379 to 1.526, 4265 and previous response = 8175.			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:47:11 AM	Set SampleApproved = True for sample 04JAN12.D; previous value = False			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/5/2022 10:47:23 AM	Manually integrate qualifier 64.0 of compound Vinyl chloride in sample 04JAN13.D, from x, y = 1.470, 5496 to 1.517, 5210, result = 15113; previous integration is from x, y = 1.308, 0 to 1.679, 0 and previous response = 148305.			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:47:46 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Chloroform in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:47:57 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,1-Dichloropropene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:48:15 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound Trichloroethene in sample 04JAN15.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/5/2022 10:48:21 AM	Set LevelEnable = False for calibration level 1, levelId = 20 of compound 1,2-Dichloropropane in sample 04JAN15.D; previous value = True			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:48:59 AM	Set SampleApproved = True for sample 04JAN13.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/5/2022 10:49:20 AM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:53:33 AM	Set SampleApproved = True for sample 04JAN17.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:53:45 AM	Set SampleApproved = True for sample 04JAN19.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:53:56 AM	Set SampleApproved = True for sample 04JAN21.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/5/2022 10:55:07 AM	Set SampleApproved = True for sample 04JAN23.D; previous value = False			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdCalibrate	BL2000\mchavez	1/5/2022 10:55:43 AM	Replace level QC with QC sample 04JAN23.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 8 with Calibration sample 04JAN21.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 7 with Calibration sample 04JAN19.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 6 with Calibration sample 04JAN17.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene,				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 5 with Calibration sample 04JAN15.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 4 with Calibration sample 04JAN13.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 3 with Calibration sample 04JAN12.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride,				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 2 with Calibration sample 04JAN11.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene}; Replace level 1 with Calibration sample 04JAN10.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, o-Xylene};				
CmdQuantitate	BL2000\mchavez	1/5/2022 10:55:58 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 11:01:11 AM	Manually integrate compound Methylene chloride in sample 04JAN09.D from x, y = 3.296, 0 to 3.383, 0; result = 1661			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 11:01:17 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 04JAN09.D from x, y = 3.299, 0 to 3.369, 0; result = 1075			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/5/2022 11:01:19 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 04JAN09.D from x, y = 3.294, 0 to 3.413, 0; result = 694			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 11:02:10 AM	Manually integrate compound Vinyl chloride in sample 04JAN09.D from x, y = 1.467, 0 to 1.520, 0; result = 73			✓	
CmdStartMethodEditing	BL2000\mchavez	1/5/2022 11:02:39 AM	Start method editing			✓	
CmdImportMethodFromSample	BL2000\mchavez	1/5/2022 11:02:39 AM	Import method from sample 04JAN09.D			✓	
CmdSetMethodTargetCompoundAttribute	BL2000\mchavez	1/5/2022 11:03:13 AM	Set CurveFit = fitAverageOfResponseFactors for compound Bromomethane; previous value = fitQuadratic			✓	
CmdSetMethodTargetCompoundAttribute	BL2000\mchavez	1/5/2022 11:03:16 AM	Set CurveFitWeight = weightEqual for compound Bromomethane; previous value = weightOneOverX			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/5/2022 11:03:28 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/5/2022 11:03:28 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/5/2022 11:03:29 AM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/5/2022 11:03:49 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 11:05:06 AM	Manually integrate compound Chloroethane in sample 04JAN12.D, from x, y = 1.863, 1400 to 1.922, 1881, result = 14646; previous integration is from x, y = 1.863, 1400 to 1.958, 1400 and previous response = 16843.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/5/2022 11:05:19 AM	Manually integrate compound Chloroethane in sample 04JAN11.D, from x, y = 1.869, 1143 to 1.930, 1702, result = 8052; previous integration is from x, y = 1.869, 1143 to 1.941, 1143 and previous response = 9540.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdCalibrate	BL2000\mchavez	1/5/2022 11:05:53 AM	Replace level QC with QC sample 04JAN23.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 8 with Calibration sample 04JAN21.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 7 with Calibration sample 04JAN19.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 6 with Calibration sample 04JAN17.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 5 with Calibration sample 04JAN15.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 4 with Calibration sample 04JAN13.D for				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 3 with Calibration sample 04JAN12.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 2 with Calibration sample 04JAN11.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane}; Replace level 1 with Calibration sample 04JAN10.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-				

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Chloroethane};				
CmdQuantitate	BL2000\mchavez	1/5/2022 11:06:12 AM	Quantitate all compounds in all samples			✓	
CmdQuantitate	BL2000\mchavez	1/5/2022 11:07:28 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/5/2022 11:07:47 AM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/9/2022 8:45:32 PM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:47:42 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound o-Xylene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:48:01 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:48:47 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound m+p-Xylenes in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:49:03 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:49:35 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound Ethylbenzene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:49:50 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:50:18 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound Styrene in sample 04JAN09.D; previous value = False			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdQuantitate	BL2000\mchavez	1/9/2022 8:50:35 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:50:51 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound Tetrachloroethene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:51:06 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:51:27 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound Toluene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:51:42 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:52:26 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound trans-1,3-Dichloropropene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:52:42 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:52:55 PM	Set LevelEnable = False for calibration level 1, levelId = 30 of compound trans-1,3-Dichloropropene in sample 04JAN09.D; previous value = True			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:53:25 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound Benzene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:53:41 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:54:07 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound Chloroform in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:54:23 PM	Quantitate all compounds in all samples			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:54:39 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound 1,3-Dichlorobenzene in sample 04JAN09.D; previous value = False			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:54:49 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound 1,4-Dichlorobenzene in sample 04JAN09.D; previous value = False			✓	
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:54:54 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound 1,2-Dichlorobenzene in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:55:10 PM	Quantitate all compounds in all samples			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetLevelEnable	BL2000\mchavez	1/9/2022 8:56:06 PM	Set LevelEnable = True for calibration level 1, levelId = 30 of compound 1,2-Dichloroethane in sample 04JAN09.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:56:22 PM	Quantitate all compounds in all samples			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/9/2022 8:56:50 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010422\04JAN15CC.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 8:57:23 PM	Set SampleType = Calibration for sample 04JAN15CC.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 8:57:29 PM	Set LevelName = 5 for sample 04JAN15CC.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 8:57:56 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 8:58:36 PM	Set SampleType = CC for sample 04JAN15CC.D; previous value = Calibration			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 8:58:43 PM	Set LevelName = CC for sample 04JAN15CC.D; previous value = 5			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 8:59:05 PM	Set UserDefined = Reimported midpoint as CC for sample 04JAN15CC.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 8:59:31 PM	Set SampleName = CC010422 for sample 04JAN15CC.D; previous value = ICAL010422_5			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdCalibrate	BL2000\mchavez	1/9/2022 8:59:53 PM	Replace level CC with CC sample 04JAN15CC.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, 1,2-Dichloroethane};			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 9:00:09 PM	Quantitate all compounds in all samples			✓	
CmdStartMethodEditing	BL2000\mchavez	1/9/2022 9:00:22 PM	Start method editing			✓	
CmdImportMethodFromSample	BL2000\mchavez	1/9/2022 9:00:22 PM	Import method from sample 04JAN23.D			✓	
CmdSaveMethodAs	BL2000\mchavez	1/9/2022 9:00:57 PM	Save method to file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL.m			✓	
CmdSaveMethodAs	BL2000\mchavez	1/9/2022 9:02:42 PM	Save method to file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m			✓	
CmdMethodClear	BL2000\mchavez	1/9/2022 9:02:55 PM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/9/2022 9:02:56 PM	End method editing			✓	
CmdStartMethodEditing	BL2000\mchavez	1/9/2022 9:03:07 PM	Start method editing			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportMethodFromFile	BL2000\mchavez	1/9/2022 9:03:07 PM	Import method from file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/9/2022 9:03:19 PM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/9/2022 9:03:19 PM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/9/2022 9:03:20 PM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 9:03:36 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/9/2022 9:04:21 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/10/2022 2:13:24 PM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/10/2022 2:13:38 PM	Set SampleApproved = True for sample 04JAN08.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\mchavez	1/10/2022 2:14:50 PM	Zero out primary peak of compound Chloromethane in sample 04JAN09.D			✓	
CmdZeroOutPeak	BL2000\mchavez	1/10/2022 2:14:56 PM	Zero out primary peak of compound Vinyl chloride in sample 04JAN09.D			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/10/2022 2:15:19 PM	Manually integrate compound Benzene in sample 04JAN09.D from x, y = 6.250, 0 to 6.311, 0; result = 381			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/10/2022 2:15:21 PM	Manually integrate qualifier77.0 of compound Benzene in sample 04JAN09.D from x, y = 6.258, 0 to 6.308, 0; result = 86			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/10/2022 2:16:46 PM	Set SampleApproved = True for sample 04JAN09.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:17:02 PM	Set UserAnnotation = NI for compound Methylene chloride in sample 04JAN09.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:17:09 PM	Set UserAnnotation = NI for compound Benzene in sample 04JAN09.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:18:45 PM	Set UserAnnotation = NI for compound Tetrachloroethene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:18:48 PM	Set UserAnnotation = NI for compound Chlorodibromomethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:18:53 PM	Set UserAnnotation = NI for compound 1,2-Dibromoethane in sample 04JAN10.D; previous value =			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:18:58 PM	Set UserAnnotation = NI for compound 1,1,1,2-Tetrachloroethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:19:02 PM	Set UserAnnotation = NI for compound Bromoform in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:19:08 PM	Set UserAnnotation = NI for compound 1,1,2,2-Tetrachloroethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:19:13 PM	Set UserAnnotation = NI for compound Bromobenzene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:19:30 PM	Set UserAnnotation = NI for compound 2-Chlorotoluene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:20:15 PM	Set UserAnnotation = GT for compound Chloroethane in sample 04JAN11.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:20:34 PM	Set UserAnnotation = LT for compound trans-1,2-Dichloroethene in sample 04JAN11.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:25:39 PM	Set UserAnnotation = LT for compound Methyl tert-butyl ether (MTBE) in sample 04JAN11.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:26:14 PM	Set UserAnnotation = LT for compound 1,1,2-Trichloroethane in sample 04JAN11.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:26:26 PM	Set UserAnnotation = NI for compound 1,2,3-Trichloropropane in sample 04JAN11.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:26:46 PM	Set UserAnnotation = LT for compound 1,2,3-Trichloropropane in sample 04JAN12.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/10/2022 2:26:59 PM	Set UserAnnotation = GT for compound Chloroethane in sample 04JAN12.D; previous value =			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/10/2022 3:00:42 PM	Manually integrate qualifier 148.0 of compound 1,4-Dichlorobenzene in sample 04JAN10.D, from x, y = 12.109, 0 to 12.145, 0, result = 2520; previous integration is from x, y = 12.084, 0 to 12.145, 0 and previous response = 3603.			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/10/2022 3:00:52 PM	Manually integrate qualifier 111.0 of compound 1,4-Dichlorobenzene in sample 04JAN10.D, from x, y = 12.123, 0 to 12.148, 0, result = 1067; previous integration is from x, y = 12.075, 0 to 12.123, 0 and previous response = 3515.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/10/2022 3:01:20 PM	Manually integrate qualifier 91.0 of compound o-Xylene in sample 04JAN10.D, from x, y = 10.405, 0 to 10.483, 0, result = 6062; previous integration is from x, y = 10.405, 0 to 10.466, 0 and previous response = 6034.			✓	
CmdClearManualIntegration	BL2000\mchavez	1/10/2022 3:01:29 PM	Clear manual integration of qualifier 91.0 for compound o-Xylene in sample 04JAN10.D			✓	
CmdSetLevelEnable	BL2000\mchavez	1/10/2022 3:07:00 PM	Set LevelEnable = False for calibration level 1, levelId = 30 of compound o-Xylene in sample 04JAN23.D; previous value = True			✓	
CmdQuantitate	BL2000\mchavez	1/10/2022 3:07:39 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/10/2022 3:12:09 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/11/2022 8:47:03 AM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdStartMethodEditing	BL2000\mchavez	1/11/2022 8:47:16 AM	Start method editing			✓	
CmdImportMethodFromSample	BL2000\mchavez	1/11/2022 8:47:16 AM	Import method from sample 04JAN01.D			✓	
CmdSaveMethodAs	BL2000\mchavez	1/11/2022 8:47:26 AM	Save method to file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/11/2022 8:48:57 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/11/2022 8:48:57 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/11/2022 8:48:58 AM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/11/2022 8:49:22 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/11/2022 8:54:59 AM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	2/28/2022 1:48:49 PM	Open batch D:\Org\Data\VOA5975C\VG010422\VG010422_8260B.batch.bin			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:23 PM	Set UserAnnotation = NI for compound Bromomethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:29 PM	Set UserAnnotation = GT for compound Chloroethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:34 PM	Set UserAnnotation = NI for compound 1,1-Dichloroethene in sample 04JAN10.D; previous value =			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:38 PM	Set UserAnnotation = NI for compound trans-1,2-Dichloroethene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:43 PM	Set UserAnnotation = NI for compound Methyl tert-butyl ether (MTBE) in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:46 PM	Set UserAnnotation = NI for compound 2,2-Dichloropropane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:51 PM	Set UserAnnotation = NI for compound cis-1,2-Dichloroethene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:55 PM	Set UserAnnotation = NI for compound Bromochloromethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:49:59 PM	Set UserAnnotation = NI for compound Dibromofluoromethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:50:03 PM	Set UserAnnotation = NI for compound 1,2-Dichloroethane-d4 in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:50:06 PM	Set UserAnnotation = NI for compound 1,2-Dichloroethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:50:10 PM	Set UserAnnotation = NI for compound Trichloroethene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:50:14 PM	Set UserAnnotation = NI for compound Dibromomethane in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:50:19 PM	Set UserAnnotation = NI for compound trans-1,3-Dichloropropene in sample 04JAN10.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	2/28/2022 1:50:23 PM	Set UserAnnotation = NI for compound 1,1,2-Trichloroethane in sample 04JAN10.D; previous value =			✓	
CmdSaveBatchTable	BL2000\mchavez	2/28/2022 1:54:50 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
GenerateReport	BL2000\mchavez	2/28/2022 1:56:23 PM	Generates report - Method: \\MASSHUNTER\Org\reports\LevelIV_Reports\Calibration\01_Init_Cal+Gen_Calibration+Gen_ResultsSummary.m, Output Path: D:\Org\Data\VOA5975C\VG010422\QuantReports\VG010422_8260B			✓	
CmdStartMethodEditing	BL2000\mchavez	2/28/2022 1:57:20 PM	Start method editing			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportMethodFromFile	BL2000\mchavez	2/28/2022 1:57:21 PM	Import method from file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	2/28/2022 1:57:33 PM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	2/28/2022 1:57:33 PM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	2/28/2022 1:57:33 PM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	2/28/2022 1:57:54 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	2/28/2022 1:59:46 PM	Save batch D:\Org\Data\VOA5975C\VG010422\QuantResults\VG010422_8260B.batch.bin			✓	
GenerateReport	BL2000\mchavez	2/28/2022 2:00:55 PM	Generates report - Method: \\MASSHUNTER\Org\reports\LevelIV_Reports\Calibration\01_Init_Cal+Gen_Calibration+Gen_ResultsSummary.m, Output Path: D:\Org\Data\VOA5975C\VG010422\QuantReports\VG010422_8260B-1			✓	

Energy Laboratories Inc

ANALYTICAL RUN Summary

02-Mar-22

Run ID VOA5975C.I_220107A

Run Start Date:	1/7/2022
Analyst:	Melissa Chavez
Ical:	
Column ID:	
Comments:	

Std ID	Std Name	Std Amount	Std Units	Samp Amount	Samp Units	SampType	Expiration Date
VOCF3517	Internal Standard / Surrogates (INT/SURR)	8.4	ul	42	ml	ALL (TUNE	12/31/2022
VOCF3529B	2nd Source MtBE	1.05	ul	42	ml	LCS, MS, M	1/29/2022
VOCF3546A	Liquids	1.05	ul	42	ml	CCV	1/13/2022
VOCF3549	2nd Source Ketones	1.05	ul	42	ml	LCS, MS, M	1/15/2022
VOCF3550	Ketones	1.05	ul	42	ml	CCV	1/16/2022
VOCF3558B	2nd Source Liquids	1.05	ul	42	ml	LCS, MS, M	2/27/2022
VOCF3559A	MtBE	1.05	ul	42	ml	CCV	1/27/2022
VOCF3562A	Gases	1.05	ul	42	ml	CCV	1/10/2022
VOCF3566A	2nd Source Gases	1.05	ul	42	ml	LCS, MS, M	1/11/2022

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973448	07JAN02_D_TU	VOC-8260-BFB	TUNE	)A5975C\VG010´	1/7/2022 9:03:00	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
173, % of mass 174	A	%	0	0		100	0	0	0	0	0	0%	0	1.99	0%	
174, % of mass 95	A	%	85.7	85.7		100	0	0	0	0	0	86%	50	99.99	0%	
175, % of mass 174	A	%	8.3	8.3		100	0	0	0	0	0	8%	5	9	0%	
176, % of mass 174	A	%	99.3	99.3		100	0	0	0	0	0	99%	95	101	0%	
177, % of mass 176	A	%	7.7	7.7		100	0	0	0	0	0	8%	5	9	0%	
50, % of mass 95	A	%	19.2	19.2		100	0	0	0	0	0	19%	15	40	0%	
75, % of mass 95	A	%	51.3	51.3		100	0	0	0	0	0	51%	30	60	0%	
95, Base Peak	A	%	100	100		100	0	0	0	0	0	100%	0	100	0%	
96, % of mass 95	A	%	6.8	6.8		100	0	0	0	0	0	7%	5	9	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973449	CCV010722_	VOC-8260-W-Q	CCV	)A5975C\VG010	1/7/2022 9:43:06	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	118.08544	4.7234176		5	0	0	0.101	0.5	500	94%	80	120	0%	
1,1,1-Trichloroethane	A	ug/L	125.09168	5.0036672		5	0	0	0.131	0.5	500	100%	80	120	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	130.59813	5.2239252		5	0	0	0.0872	0.5	500	104%	80	120	0%	
1,1,2-Trichloroethane	A	ug/L	123.52129	4.9408516		5	0	0	0.108	0.5	500	99%	80	120	0%	
1,1-Dichloroethane	A	ug/L	127.83306	5.1133224		5	0	0	0.135	0.5	500	102%	80	120	0%	
1,1-Dichloroethene	A	ug/L	121.47259	4.8589036		5	0	0	0.141	0.5	500	97%	80	120	0%	
1,1-Dichloropropene	A	ug/L	125.60688	5.0242752		5	0	0	0.083	0.5	500	100%	80	120	0%	
1,2,3-Trichloropropane	A	ug/L	129.90227	5.1960908		5	0	0	0.235	0.5	500	104%	80	120	0%	
1,2-Dibromoethane	A	ug/L	125.22993	5.0091972		5	0	0	0.0916	0.5	500	100%	80	120	0%	
1,2-Dichlorobenzene	A	ug/L	124.08023	4.9632092		5	0	0	0.0746	0.5	500	99%	80	120	0%	
1,2-Dichloroethane	A	ug/L	123.7158	4.948632		5	0	0	0.116	0.5	500	99%	80	120	0%	
1,2-Dichloropropane	A	ug/L	127.33164	5.0932656		5	0	0	0.0847	0.5	500	102%	80	120	0%	
1,3-Dichlorobenzene	A	ug/L	124.85744	4.9942976		5	0	0	0.0803	0.5	500	100%	80	120	0%	
1,3-Dichloropropane	A	ug/L	124.55548	4.9822192		5	0	0	0.0791	0.5	500	100%	80	120	0%	
1,4-Dichlorobenzene	A	ug/L	126.09813	5.0439252		5	0	0	0.0858	0.5	500	101%	80	120	0%	
2,2-Dichloropropane	A	ug/L	130.08225	5.20329		5	0	0	0.186	0.5	500	104%	80	120	0%	
2-Chlorotoluene	A	ug/L	129.2634	5.170536		5	0	0	0.0876	0.5	500	103%	80	120	0%	
4-Chlorotoluene	A	ug/L	132.69103	5.3076412		5	0	0	0.0728	0.5	500	106%	80	120	0%	
Benzene	A	ug/L	127.07276	5.0829104		5	0	0	0.0914	0.5	500	102%	80	120	0%	
Bromobenzene	A	ug/L	130.11606	5.2046424		5	0	0	0.0831	0.5	500	104%	80	120	0%	
Bromochloromethane	A	ug/L	128.60244	5.1440976		5	0	0	0.141	0.5	500	103%	80	120	0%	
Bromodichloromethane	A	ug/L	123.9145	4.95658		5	0	0	0.12	0.5	500	99%	80	120	0%	
Bromoform	A	ug/L	128.72304	5.1489216		5	0	0	0.119	0.5	500	103%	80	120	0%	
Bromomethane	A	ug/L	132.28115	5.291246		5	0	0	0.253	0.5	500	106%	80	120	0%	
Carbon tetrachloride	A	ug/L	120.27745	4.811098		5	0	0	0.143	0.5	500	96%	80	120	0%	
Chlorobenzene	A	ug/L	123.25086	4.9300344		5	0	0	0.0914	0.5	500	99%	80	120	0%	
Chlorodibromomethane	A	ug/L	124.96785	4.998714		5	0	0	0.0841	0.5	500	100%	80	120	0%	
Chloroethane	A	ug/L	118.04495	4.721798		5	0	0	0.169	0.5	500	94%	80	120	0%	
Chloroform	A	ug/L	121.57152	4.8628608		5	0	0	0.0789	0.5	500	97%	80	120	0%	
Chloromethane	A	ug/L	116.39918	4.6559672		5	0	0	0.162	0.5	500	93%	80	120	0%	
cis-1,2-Dichloroethene	A	ug/L	128.70764	5.1483056		5	0	0	0.108	0.5	500	103%	80	120	0%	
cis-1,3-Dichloropropene	A	ug/L	122.22136	4.8888544		5	0	0	0.073	0.5	500	98%	80	120	0%	
Dibromomethane	A	ug/L	121.6934	4.867736		5	0	0	0.147	0.5	500	97%	80	120	0%	
Dichlorodifluoromethane	A	ug/L	114.48502	4.5794008		5	0	0	0.175	0.5	500	92%	80	120	0%	
Ethylbenzene	A	ug/L	123.92738	4.9570952		5	0	0	0.0836	0.5	500	99%	80	120	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973449	CCV010722_	VOC-8260-W-Q	CCV	)A5975C\VG010´	1/7/2022 9:43:06	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	252.2653	10.090612		10	0	0	0.15	0.5	1000	101%	80	120	0%	
Methyl ethyl ketone	A	ug/L	1257.21492	50.2885968		50	0	0	1.77	10	5000	101%	80	120	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	127.82414	5.1129656		5	0	0	0.101	0.5	500	102%	80	120	0%	
Methylene chloride	A	ug/L	119.35881	4.7743524		5	0	0	0.338	0.5	500	95%	80	120	0%	
o-Xylene	A	ug/L	124.69512	4.9878048		5	0	0	0.0604	0.5	500	100%	80	120	0%	
Styrene	A	ug/L	129.82914	5.1931656		5	0	0	0.067	0.5	500	104%	80	120	0%	
Tetrachloroethene	A	ug/L	117.83369	4.7133476		5	0	0	0.0671	0.5	500	94%	80	120	0%	
Toluene	A	ug/L	127.1937	5.087748		5	0	0	0.0679	0.5	500	102%	80	120	0%	
trans-1,2-Dichloroethene	A	ug/L	125.43198	5.0172792		5	0	0	0.125	0.5	500	100%	80	120	0%	
trans-1,3-Dichloropropene	A	ug/L	127.53024	5.1012096		5	0	0	0.0846	0.5	500	102%	80	120	0%	
Trichloroethene	A	ug/L	120.79095	4.831638		5	0	0	0.0993	0.5	500	97%	80	120	0%	
Trichlorofluoromethane	A	ug/L	112.41757	4.4967028		5	0	0	0.134	0.5	500	90%	80	120	0%	
Vinyl chloride	A	ug/L	121.48241	4.8592964		5	0	0	0.153	0.5	500	97%	80	120	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	376.96042	15.0784168		15	0	0	0.0604	0.5	1500	101%	80	120	0%	
1,2-Dichloroethane-d4	S	ug/L	280.59167	11.2236668		10	0	0	0.229	0.5	500	112%	80	120	0%	
Dibromofluoromethane	S	ug/L	279.46855	11.178742		10	0	0	0.129	0.5	500	112%	80	120	0%	
p-Bromofluorobenzene	S	ug/L	276.60604	11.0642416		10	0	0	0.149	0.5	500	111%	80	120	0%	
Toluene-d8	S	ug/L	269.54488	10.7817952		10	0	0	0.23	0.5	500	108%	80	120	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973450	LCS010722_	VOC-8260-W-Q	LCS-DOD	)A5975C\VG010´	1/7/2022 10:50:3	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	125.15027	5.0060108		5	0	0	0.101	0.5	500	100%	78	124	0%	
1,1,1-Trichloroethane	A	ug/L	122.15451	4.8861804		5	0	0	0.131	0.5	500	98%	74	131	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	123.11919	4.9247676		5	0	0	0.0872	0.5	500	98%	71	121	0%	
1,1,2-Trichloroethane	A	ug/L	125.1597	5.006388		5	0	0	0.108	0.5	500	100%	80	119	0%	
1,1-Dichloroethane	A	ug/L	131.07885	5.243154		5	0	0	0.135	0.5	500	105%	77	125	0%	
1,1-Dichloroethene	A	ug/L	133.65174	5.3460696		5	0	0	0.141	0.5	500	107%	71	131	0%	
1,1-Dichloropropene	A	ug/L	118.08273	4.7233092		5	0	0	0.083	0.5	500	94%	79	125	0%	
1,2,3-Trichloropropane	A	ug/L	125.17636	5.0070544		5	0	0	0.235	0.5	500	100%	73	125	0%	
1,2-Dibromoethane	A	ug/L	123.15191	4.9260764		5	0	0	0.0916	0.5	500	99%	78	122	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist				
14973450	LCS010722_	VOC-8260-W-Q	LCS-DOD	)A5975C\VG010	1/7/2022 10:50:3	1	R373037		0	0					

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	125.55956	5.0223824		5	0	0	0.0746	0.5	500	100%	80	119	0%	
1,2-Dichloroethane	A	ug/L	121.9792	4.879168		5	0	0	0.116	0.5	500	98%	73	128	0%	
1,2-Dichloropropane	A	ug/L	124.07441	4.9629764		5	0	0	0.0847	0.5	500	99%	78	122	0%	
1,3-Dichlorobenzene	A	ug/L	131.94939	5.2779756		5	0	0	0.0803	0.5	500	106%	80	119	0%	
1,3-Dichloropropane	A	ug/L	121.38496	4.8553984		5	0	0	0.0791	0.5	500	97%	80	119	0%	
1,4-Dichlorobenzene	A	ug/L	128.63551	5.1454204		5	0	0	0.0858	0.5	500	103%	79	118	0%	
2,2-Dichloropropane	A	ug/L	128.77441	5.1509764		5	0	0	0.186	0.5	500	103%	60	139	0%	
2-Chlorotoluene	A	ug/L	130.04655	5.201862		5	0	0	0.0876	0.5	500	104%	79	122	0%	
4-Chlorotoluene	A	ug/L	134.9812	5.399248		5	0	0	0.0728	0.5	500	108%	78	122	0%	
Benzene	A	ug/L	127.6093	5.104372		5	0	0	0.0914	0.5	500	102%	79	120	0%	
Bromobenzene	A	ug/L	131.56488	5.2625952		5	0	0	0.0831	0.5	500	105%	80	120	0%	
Bromochloromethane	A	ug/L	120.19576	4.8078304		5	0	0	0.141	0.5	500	96%	78	123	0%	
Bromodichloromethane	A	ug/L	127.5739	5.102956		5	0	0	0.12	0.5	500	102%	79	125	0%	
Bromoform	A	ug/L	131.15559	5.2462236		5	0	0	0.119	0.5	500	105%	66	130	0%	
Bromomethane	A	ug/L	120.90262	4.8361048		5	0	0	0.253	0.5	500	97%	53	141	0%	
Carbon tetrachloride	A	ug/L	117.63797	4.7055188		5	0	0	0.143	0.5	500	94%	72	136	0%	
Chlorobenzene	A	ug/L	125.64774	5.0259096		5	0	0	0.0914	0.5	500	101%	82	118	0%	
Chlorodibromomethane	A	ug/L	122.73184	4.9092736		5	0	0	0.0841	0.5	500	98%	74	126	0%	
Chloroethane	A	ug/L	109.6085	4.38434		5	0	0	0.169	0.5	500	88%	60	138	0%	
Chloroform	A	ug/L	119.37452	4.7749808		5	0	0	0.0789	0.5	500	95%	79	124	0%	
Chloromethane	A	ug/L	108.69779	4.3479116		5	0	0	0.162	0.5	500	87%	50	139	0%	
cis-1,2-Dichloroethene	A	ug/L	129.68356	5.1873424		5	0	0	0.108	0.5	500	104%	78	123	0%	
cis-1,3-Dichloropropene	A	ug/L	117.25682	4.6902728		5	0	0	0.073	0.5	500	94%	75	124	0%	
Dibromomethane	A	ug/L	124.14929	4.9659716		5	0	0	0.147	0.5	500	99%	79	123	0%	
Dichlorodifluoromethane	A	ug/L	103.63824	4.1455296		5	0	0	0.175	0.5	500	83%	32	152	0%	
Ethylbenzene	A	ug/L	125.94524	5.0378096		5	0	0	0.0836	0.5	500	101%	79	121	0%	
m+p-Xylenes	A	ug/L	250.90789	10.0363156		10	0	0	0.15	0.5	1000	100%	80	121	0%	
Methyl ethyl ketone	A	ug/L	1138.31283	45.5325132		50	0	0	1.77	10	5000	91%	56	143	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	132.80502	5.3122008		5	0	0	0.101	0.5	500	106%	71	124	0%	
Methylene chloride	A	ug/L	118.10926	4.7243704		5	0	0	0.338	0.5	500	94%	74	124	0%	
o-Xylene	A	ug/L	129.8994	5.195976		5	0	0	0.0604	0.5	500	104%	78	122	0%	
Styrene	A	ug/L	132.51502	5.3006008		5	0	0	0.067	0.5	500	106%	78	123	0%	
Tetrachloroethene	A	ug/L	118.88372	4.7553488		5	0	0	0.0671	0.5	500	95%	74	129	0%	
Toluene	A	ug/L	129.15194	5.1660776		5	0	0	0.0679	0.5	500	103%	80	121	0%	
trans-1,2-Dichloroethene	A	ug/L	129.83201	5.1932804		5	0	0	0.125	0.5	500	104%	75	124	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973450	LCS010722_	VOC-8260-W-Q	LCS-DOD	)A5975C\VG010´	1/7/2022 10:50:3	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
trans-1,3-Dichloropropene	A	ug/L	131.25712	5.2502848		5	0	0	0.0846	0.5	500	105%	73	127	0%	
Trichloroethene	A	ug/L	125.89894	5.0359576		5	0	0	0.0993	0.5	500	101%	79	123	0%	
Trichlorofluoromethane	A	ug/L	112.75149	4.5100596		5	0	0	0.134	0.5	500	90%	65	141	0%	
Vinyl chloride	A	ug/L	119.23499	4.7693996		5	0	0	0.153	0.5	500	95%	58	137	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	380.80729	15.2322916		15	0	0	0.0604	0.5	1500	102%	79	121	0%	
1,2-Dichloroethane-d4	S	ug/L	282.91531	11.3166124		10	0	0	0.229	0.5	500	113%	81	118	0%	
Dibromofluoromethane	S	ug/L	270.45155	10.818062		10	0	0	0.129	0.5	500	108%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	271.98175	10.87927		10	0	0	0.149	0.5	500	109%	85	114	0%	
Toluene-d8	S	ug/L	271.05199	10.8420796		10	0	0	0.23	0.5	500	108%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973453	MBLK010722_	VOC-8260-W-Q	MBLK	)A5975C\VG010´	1/7/2022 11:45:0	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	0.5	500	0%	0	0	0%	
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	0.5	500	0%	0	0	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	0.5	500	0%	0	0	0%	
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	0.5	500	0%	0	0	0%	
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	0.5	500	0%	0	0	0%	
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	0.5	500	0%	0	0	0%	
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	0.5	500	0%	0	0	0%	
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	0.5	500	0%	0	0	0%	
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	0.5	500	0%	0	0	0%	
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	0.5	500	0%	0	0	0%	
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	0.5	500	0%	0	0	0%	
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	0.5	500	0%	0	0	0%	
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	0.5	500	0%	0	0	0%	
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	0.5	500	0%	0	0	0%	
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	0.5	500	0%	0	0	0%	
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	0.5	500	0%	0	0	0%	
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	0.5	500	0%	0	0	0%	
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	0.5	500	0%	0	0	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973453	MBLK010722_	VOC-8260-W-Q	MBLK	)A5975C\VG010	1/7/2022 11:45:0	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Benzene	A	ug/L	0	0		0	0	0	0.0914	0.5	500	0%	0	0	0%	
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	0.5	500	0%	0	0	0%	
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	0.5	500	0%	0	0	0%	
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	0.5	500	0%	0	0	0%	
Bromoform	A	ug/L	0	0		0	0	0	0.119	0.5	500	0%	0	0	0%	
Bromomethane	A	ug/L	0	0		0	0	0	0.253	0.5	500	0%	0	0	0%	
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	0.5	500	0%	0	0	0%	
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	0.5	500	0%	0	0	0%	
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	0.5	500	0%	0	0	0%	
Chloroethane	A	ug/L	0	0		0	0	0	0.169	0.5	500	0%	0	0	0%	
Chloroform	A	ug/L	0	0		0	0	0	0.0789	0.5	500	0%	0	0	0%	
Chloromethane	A	ug/L	0	0		0	0	0	0.162	0.5	500	0%	0	0	0%	
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	0.5	500	0%	0	0	0%	
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	0.5	500	0%	0	0	0%	
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	0.5	500	0%	0	0	0%	
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	0.5	500	0%	0	0	0%	
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	0.5	500	0%	0	0	0%	
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	0.5	1000	0%	0	0	0%	
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	10	5000	0%	0	0	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	0.5	500	0%	0	0	0%	
Methylene chloride	A	ug/L	1.28894	0		0	0	0	0.338	0.5	500	0%	0	0	0%	
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	0.5	500	0%	0	0	0%	
Styrene	A	ug/L	0	0		0	0	0	0.067	0.5	500	0%	0	0	0%	
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	0.5	500	0%	0	0	0%	
Toluene	A	ug/L	0	0		0	0	0	0.0679	0.5	500	0%	0	0	0%	
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	0.5	500	0%	0	0	0%	
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	0.5	500	0%	0	0	0%	
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	0.5	500	0%	0	0	0%	
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	0.5	500	0%	0	0	0%	
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	0.5	500	0%	0	0	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	0.5	1500	0%	0	0	0%	
1,2-Dichloroethane-d4	S	ug/L	279.14576	11.1658304		10	0	0	0.229	0.5	500	112%	81	118	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973453	MBLK010722_	VOC-8260-W-Q	MBLK	)A5975C\VG010´	1/7/2022 11:45:0	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Dibromofluoromethane	S	ug/L	270.76663	10.8306652		10	0	0	0.129	0.5	500	108%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	274.98941	10.9995764		10	0	0	0.149	0.5	500	110%	85	114	0%	
Toluene-d8	S	ug/L	264.37748	10.5750992		10	0	0	0.23	0.5	500	106%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973457	B22010214-002	VOC-8260-W-S	SAMP	)A5975C\VG010´	1/7/2022 12:24:5	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0.048	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	1	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973457	B22010214-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 12:24:5		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Chloroethane	A	ug/L	0	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	0.14554	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	1.44449	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0.10634	0		0	0	0	0.0836	1	500	0%	0	0	0%	U
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	1.98693	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	1	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	1.99774	0.0799096		0	0	0	0.0679	1	500	0%	0	0	0%	J
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	1	0	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	314.75492	12.5901968		10	0	0	0.229	1	500	126%	81	118	0%	S
Dibromofluoromethane	S	ug/L	292.59389	11.7037556		10	0	0	0.129	1	500	117%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	283.55508	11.3422032		10	0	0	0.149	1	500	113%	85	114	0%	
Toluene-d8	S	ug/L	260.0307	10.401228		10	0	0	0.23	1	500	104%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973458	B22010219-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 12:52:1		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist				
14973458	B22010219-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 12:52:1		1	R373037		0	0					

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0.06481	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	1	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	1	500	0%	0	0	0%	U
Chloroethane	A	ug/L	0	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	0	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	1.90607	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0.12145	0		0	0	0	0.0836	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973458	B22010219-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 12:52:1		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	1.833	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	1	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	1.80909	0.0723636		0	0	0	0.0679	1	500	0%	0	0	0%	J
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	1	0	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	289.53136	11.5812544		10	0	0	0.229	1	500	116%	81	118	0%	
Dibromofluoromethane	S	ug/L	275.67606	11.0270424		10	0	0	0.129	1	500	110%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	279.27018	11.1708072		10	0	0	0.149	1	500	112%	85	114	0%	
Toluene-d8	S	ug/L	263.89535	10.555814		10	0	0	0.23	1	500	106%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973459	B22010262-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 1:19:35		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973459	B22010262-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 1:19:35		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	1	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	1	500	0%	0	0	0%	U
Chloroethane	A	ug/L	0	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	0.24661	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	0.80639	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	1	500	0%	0	0	0%	U
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	2.20608	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	1	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	2.19497	0.0877988		0	0	0	0.0679	1	500	0%	0	0	0%	J
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973459	B22010262-002	VOC-8260-W-S	SAMP	)A5975C\VG010´	1/7/2022 1:19:35	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	1	0	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	281.55972	11.2623888		10	0	0	0.229	1	500	113%	81	118	0%	
Dibromofluoromethane	S	ug/L	277.07768	11.0831072		10	0	0	0.129	1	500	111%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	270.77107	10.8308428		10	0	0	0.149	1	500	108%	85	114	0%	
Toluene-d8	S	ug/L	271.91357	10.8765428		10	0	0	0.23	1	500	109%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973460	B22010214-002	VOC-8260-W-S	SAMP	)A5975C\VG010´	1/7/2022 1:46:47	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973461	B22010219-001	VOC-8260-W-S	SAMP	)A5975C\VG010´	1/7/2022 2:14:04	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0.6175	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973461	B22010219-001	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 2:14:04		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0.40048	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0.43706	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	1	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	1	500	0%	0	0	0%	U
Chloroethane	A	ug/L	2.80665	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	0.50573	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	1.01509	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	1	500	0%	0	0	0%	U
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	1.07273	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	1	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	0.34358	0		0	0	0	0.0679	1	500	0%	0	0	0%	UT
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973461	B22010219-001	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 2:14:04		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	1	0	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	295.01742	11.8006968		10	0	0	0.229	1	500	118%	81	118	0%	
Dibromofluoromethane	S	ug/L	277.84253	11.1137012		10	0	0	0.129	1	500	111%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	270.69927	10.8279708		10	0	0	0.149	1	500	108%	85	114	0%	
Toluene-d8	S	ug/L	269.72701	10.7890804		10	0	0	0.23	1	500	108%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973462	B22010213-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 2:41:20		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0.12888	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973462	B22010213-002	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 2:41:20		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Bromodichloromethane	A	ug/L	0.66597	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	6.9642	0.278568		0	0	0	0.119	1	500	0%	0	0	0%	J
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	2.48888	0.0995552		0	0	0	0.0841	1	500	0%	0	0	0%	J
Chloroethane	A	ug/L	0	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	1.08961	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	0.67305	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0.31217	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0.36732	0		0	0	0	0.0836	1	500	0%	0	0	0%	U
m+p-Xylenes	A	ug/L	0.88259	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	5.01792	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	0	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	2.80675	0.11227		0	0	0	0.0604	1	500	0%	0	0	0%	J
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	0.23877	0		0	0	0	0.0679	1	500	0%	0	0	0%	UT
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	3.68934	0.11227		0	0	0	0.0604	1	0	0%	0	0	0%	J
1,2-Dichloroethane-d4	S	ug/L	291.79466	11.6717864		10	0	0	0.229	1	500	117%	81	118	0%	
Dibromofluoromethane	S	ug/L	278.83419	11.1533676		10	0	0	0.129	1	500	112%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	270.38071	10.8152284		10	0	0	0.149	1	500	108%	85	114	0%	
Toluene-d8	S	ug/L	261.60968	10.4643872		10	0	0	0.23	1	500	105%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973463	B22010214-001	VOC-8260-W-S	SAMP	)A5975C\VG010	1/7/2022 3:35:59	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	1	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	1	500	0%	0	0	0%	U
Chloroethane	A	ug/L	0	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	0	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	1.7362	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973463	B22010214-001	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 3:35:59		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	0.70004	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	1	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	0	0		0	0	0	0.0679	1	500	0%	0	0	0%	UT
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	1	0	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	287.08913	11.4835652		10	0	0	0.229	1	500	115%	81	118	0%	
Dibromofluoromethane	S	ug/L	279.63706	11.1854824		10	0	0	0.129	1	500	112%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	276.52735	11.061094		10	0	0	0.149	1	500	111%	85	114	0%	
Toluene-d8	S	ug/L	267.8901	10.715604		10	0	0	0.23	1	500	107%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973464	B22010262-001	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 4:03:14		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	1	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	1	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0	0		0	0	0	0.135	1	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	1	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	1	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973464	B22010262-001	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 4:03:14		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	1	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	1	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	1	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	1	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0	0		0	0	0	0.0791	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	1	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	1	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	1	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	1	500	0%	0	0	0%	U
Benzene	A	ug/L	0.22658	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	1	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	1	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	1	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	1	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	1	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	1	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	1	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	1	500	0%	0	0	0%	U
Chloroethane	A	ug/L	0	0		0	0	0	0.169	1	500	0%	0	0	0%	U
Chloroform	A	ug/L	0.38011	0		0	0	0	0.0789	1	500	0%	0	0	0%	U
Chloromethane	A	ug/L	0.4748	0		0	0	0	0.162	1	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	1	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	1	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	1	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	1.09038	0		0	0	0	0.175	1	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	1	500	0%	0	0	0%	U
m+p-Xylenes	A	ug/L	0.33148	0		0	0	0	0.15	1	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	20	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	1	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	0.52611	0		0	0	0	0.338	1	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	1	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	1	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	1	500	0%	0	0	0%	U
Toluene	A	ug/L	0.35025	0		0	0	0	0.0679	1	500	0%	0	0	0%	UT
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	1	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973464	B22010262-001	VOC-8260-W-S	SAMP	)A5975C\VG010´1/7/2022 4:03:14		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	1	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	1	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	1	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	1	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	0	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0.1	0	0	0%	0	0	0%	
Xylenes, Total	M	ug/L	0.33148	0		0	0	0	0.0604	1	0	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	286.13131	11.4452524		10	0	0	0.229	1	500	114%	81	118	0%	
Dibromofluoromethane	S	ug/L	273.56212	10.9424848		10	0	0	0.129	1	500	109%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	269.84089	10.7936356		10	0	0	0.149	1	500	108%	85	114	0%	
Toluene-d8	S	ug/L	270.53246	10.8212984		10	0	0	0.23	1	500	108%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973525	B22010219-001	VOC-8260-W-Q	SAMP	)A5975C\VG010´1/7/2022 2:14:04		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.101	0.5	500	0%	0	0	0%	U
1,1,1-Trichloroethane	A	ug/L	0	0		0	0	0	0.131	0.5	500	0%	0	0	0%	U
1,1,2,2-Tetrachloroethane	A	ug/L	0	0		0	0	0	0.0872	0.5	500	0%	0	0	0%	U
1,1,2-Trichloroethane	A	ug/L	0	0		0	0	0	0.108	0.5	500	0%	0	0	0%	U
1,1-Dichloroethane	A	ug/L	0.6175	0		0	0	0	0.135	0.5	500	0%	0	0	0%	U
1,1-Dichloroethene	A	ug/L	0	0		0	0	0	0.141	0.5	500	0%	0	0	0%	U
1,1-Dichloropropene	A	ug/L	0	0		0	0	0	0.083	0.5	500	0%	0	0	0%	U
1,2,3-Trichloropropane	A	ug/L	0	0		0	0	0	0.235	0.5	500	0%	0	0	0%	U
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0916	0.5	500	0%	0	0	0%	U
1,2-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0746	0.5	500	0%	0	0	0%	U
1,2-Dichloroethane	A	ug/L	0	0		0	0	0	0.116	0.5	500	0%	0	0	0%	U
1,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.0847	0.5	500	0%	0	0	0%	U
1,3-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0803	0.5	500	0%	0	0	0%	U
1,3-Dichloropropane	A	ug/L	0.40048	0		0	0	0	0.0791	0.5	500	0%	0	0	0%	U
1,4-Dichlorobenzene	A	ug/L	0	0		0	0	0	0.0858	0.5	500	0%	0	0	0%	U
2,2-Dichloropropane	A	ug/L	0	0		0	0	0	0.186	0.5	500	0%	0	0	0%	U
2-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0876	0.5	500	0%	0	0	0%	U
4-Chlorotoluene	A	ug/L	0	0		0	0	0	0.0728	0.5	500	0%	0	0	0%	U

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973525	B22010219-001	VOC-8260-W-Q	SAMP	)A5975C\VG010´1/7/2022 2:14:04		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Benzene	A	ug/L	0.43706	0		0	0	0	0.0914	0.5	500	0%	0	0	0%	U
Bromobenzene	A	ug/L	0	0		0	0	0	0.0831	0.5	500	0%	0	0	0%	U
Bromochloromethane	A	ug/L	0	0		0	0	0	0.141	0.5	500	0%	0	0	0%	U
Bromodichloromethane	A	ug/L	0	0		0	0	0	0.12	0.5	500	0%	0	0	0%	U
Bromoform	A	ug/L	0	0		0	0	0	0.119	0.5	500	0%	0	0	0%	U
Bromomethane	A	ug/L	0	0		0	0	0	0.253	0.5	500	0%	0	0	0%	U
Carbon tetrachloride	A	ug/L	0	0		0	0	0	0.143	0.5	500	0%	0	0	0%	U
Chlorobenzene	A	ug/L	0	0		0	0	0	0.0914	0.5	500	0%	0	0	0%	U
Chlorodibromomethane	A	ug/L	0	0		0	0	0	0.0841	0.5	500	0%	0	0	0%	U
Chloroethane	A	ug/L	2.80665	0		0	0	0	0.169	0.5	500	0%	0	0	0%	U
Chloroform	A	ug/L	0.50573	0		0	0	0	0.0789	0.5	500	0%	0	0	0%	U
Chloromethane	A	ug/L	1.01509	0		0	0	0	0.162	0.5	500	0%	0	0	0%	U
cis-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.108	0.5	500	0%	0	0	0%	U
cis-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.073	0.5	500	0%	0	0	0%	U
Dibromomethane	A	ug/L	0	0		0	0	0	0.147	0.5	500	0%	0	0	0%	U
Dichlorodifluoromethane	A	ug/L	0	0		0	0	0	0.175	0.5	500	0%	0	0	0%	U
Ethylbenzene	A	ug/L	0	0		0	0	0	0.0836	0.5	500	0%	0	0	0%	U
m+p-Xylenes	A	ug/L	0	0		0	0	0	0.15	0.5	1000	0%	0	0	0%	U
Methyl ethyl ketone	A	ug/L	0	0		0	0	0	1.77	10	5000	0%	0	0	0%	U
Methyl tert-butyl ether (MTBE)	A	ug/L	0	0		0	0	0	0.101	0.5	500	0%	0	0	0%	U
Methylene chloride	A	ug/L	1.07273	0		0	0	0	0.338	0.5	500	0%	0	0	0%	U
o-Xylene	A	ug/L	0	0		0	0	0	0.0604	0.5	500	0%	0	0	0%	U
Styrene	A	ug/L	0	0		0	0	0	0.067	0.5	500	0%	0	0	0%	U
Tetrachloroethene	A	ug/L	0	0		0	0	0	0.0671	0.5	500	0%	0	0	0%	U
Toluene	A	ug/L	0.34358	0		0	0	0	0.0679	0.5	500	0%	0	0	0%	U
trans-1,2-Dichloroethene	A	ug/L	0	0		0	0	0	0.125	0.5	500	0%	0	0	0%	U
trans-1,3-Dichloropropene	A	ug/L	0	0		0	0	0	0.0846	0.5	500	0%	0	0	0%	U
Trichloroethene	A	ug/L	0	0		0	0	0	0.0993	0.5	500	0%	0	0	0%	U
Trichlorofluoromethane	A	ug/L	0	0		0	0	0	0.134	0.5	500	0%	0	0	0%	U
Vinyl chloride	A	ug/L	0	0		0	0	0	0.153	0.5	500	0%	0	0	0%	U
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	0	0		0	0	0	0.0604	0.5	1500	0%	0	0	0%	U
1,2-Dichloroethane-d4	S	ug/L	295.01742	11.8006968		10	0	0	0.229	0.5	500	118%	81	118	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973525	B22010219-001	VOC-8260-W-Q	SAMP	)A5975C\VG010´1/7/2022 2:14:04		1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Dibromofluoromethane	S	ug/L	277.84253	11.1137012		10	0	0	0.129	0.5	500	111%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	270.69927	10.8279708		10	0	0	0.149	0.5	500	108%	85	114	0%	
Toluene-d8	S	ug/L	269.72701	10.7890804		10	0	0	0.23	0.5	500	108%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973526	B22010219-001	VOC-8260-W-Q	MS-DOD	)A5975C\VG010´1/7/2022 5:07:38		1	R373037		1E+07	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	123.97204	4.9588816		5	0	0	0.101	0.5	500	99%	78	124	0%	
1,1,1-Trichloroethane	A	ug/L	124.76854	4.9907416		5	0	0	0.131	0.5	500	100%	74	131	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	131.23289	5.2493156		5	0	0	0.0872	0.5	500	105%	71	121	0%	
1,1,2-Trichloroethane	A	ug/L	123.55913	4.9423652		5	0	0	0.108	0.5	500	99%	80	119	0%	
1,1-Dichloroethane	A	ug/L	135.5713	5.422852		5	0	0	0.135	0.5	500	108%	77	125	0%	
1,1-Dichloroethene	A	ug/L	131.07964	5.2431856		5	0	0	0.141	0.5	500	105%	71	131	0%	
1,1-Dichloropropene	A	ug/L	121.41514	4.8566056		5	0	0	0.083	0.5	500	97%	79	125	0%	
1,2,3-Trichloropropane	A	ug/L	123.82462	4.9529848		5	0	0	0.235	0.5	500	99%	73	125	0%	
1,2-Dibromoethane	A	ug/L	124.83894	4.9935576		5	0	0	0.0916	0.5	500	100%	78	122	0%	
1,2-Dichlorobenzene	A	ug/L	125.95905	5.038362		5	0	0	0.0746	0.5	500	101%	80	119	0%	
1,2-Dichloroethane	A	ug/L	127.86267	5.1145068		5	0	0	0.116	0.5	500	102%	73	128	0%	
1,2-Dichloropropane	A	ug/L	125.3364	5.013456		5	0	0	0.0847	0.5	500	100%	78	122	0%	
1,3-Dichlorobenzene	A	ug/L	131.39699	5.2558796		5	0	0	0.0803	0.5	500	105%	80	119	0%	
1,3-Dichloropropane	A	ug/L	126.75683	5.0702732		5	0	0	0.0791	0.5	500	101%	80	119	0%	
1,4-Dichlorobenzene	A	ug/L	129.89641	5.1958564		5	0	0	0.0858	0.5	500	104%	79	118	0%	
2,2-Dichloropropane	A	ug/L	131.81307	5.2725228		5	0	0	0.186	0.5	500	105%	60	139	0%	
2-Chlorotoluene	A	ug/L	131.31335	5.252534		5	0	0	0.0876	0.5	500	105%	79	122	0%	
4-Chlorotoluene	A	ug/L	133.3048	5.332192		5	0	0	0.0728	0.5	500	107%	78	122	0%	
Benzene	A	ug/L	130.23088	5.2092352		5	0	0	0.0914	0.5	500	104%	79	120	0%	
Bromobenzene	A	ug/L	133.61494	5.3445976		5	0	0	0.0831	0.5	500	107%	80	120	0%	
Bromochloromethane	A	ug/L	124.44772	4.9779088		5	0	0	0.141	0.5	500	100%	78	123	0%	
Bromodichloromethane	A	ug/L	128.24178	5.1296712		5	0	0	0.12	0.5	500	103%	79	125	0%	
Bromoform	A	ug/L	129.70364	5.1881456		5	0	0	0.119	0.5	500	104%	66	130	0%	
Bromomethane	A	ug/L	120.98596	4.8394384		5	0	0	0.253	0.5	500	97%	53	141	0%	
Carbon tetrachloride	A	ug/L	121.69468	4.8677872		5	0	0	0.143	0.5	500	97%	72	136	0%	
Chlorobenzene	A	ug/L	128.27202	5.1308808		5	0	0	0.0914	0.5	500	103%	82	118	0%	
Chlorodibromomethane	A	ug/L	125.15946	5.0063784		5	0	0	0.0841	0.5	500	100%	74	126	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973526	B22010219-001	VOC-8260-W-Q	MS-DOD	)A5975C\VG010´	1/7/2022 5:07:38	1	R373037		1E+07	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
Chloroethane	A	ug/L	141.01823	5.6407292		5	0	0	0.169	0.5	500	113%	60	138	0%	
Chloroform	A	ug/L	121.47415	4.858966		5	0	0	0.0789	0.5	500	97%	79	124	0%	
Chloromethane	A	ug/L	111.80488	4.4721952		5	0	0	0.162	0.5	500	89%	50	139	0%	
cis-1,2-Dichloroethene	A	ug/L	131.27992	5.2511968		5	0	0	0.108	0.5	500	105%	78	123	0%	
cis-1,3-Dichloropropene	A	ug/L	121.0494	4.841976		5	0	0	0.073	0.5	500	97%	75	124	0%	
Dibromomethane	A	ug/L	125.77977	5.0311908		5	0	0	0.147	0.5	500	101%	79	123	0%	
Dichlorodifluoromethane	A	ug/L	107.33342	4.2933368		5	0	0	0.175	0.5	500	86%	32	152	0%	
Ethylbenzene	A	ug/L	127.05741	5.0822964		5	0	0	0.0836	0.5	500	102%	79	121	0%	
m+p-Xylenes	A	ug/L	253.24827	10.1299308		10	0	0	0.15	0.5	1000	101%	80	121	0%	
Methyl ethyl ketone	A	ug/L	1256.63414	50.2653656		50	0	0	1.77	10	5000	101%	56	143	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	140.12523	5.6050092		5	0	0	0.101	0.5	500	112%	71	124	0%	
Methylene chloride	A	ug/L	121.94927	4.8779708		5	0	0	0.338	0.5	500	98%	74	124	0%	
o-Xylene	A	ug/L	130.73381	5.2293524		5	0	0	0.0604	0.5	500	105%	78	122	0%	
Styrene	A	ug/L	132.42019	5.2968076		5	0	0	0.067	0.5	500	106%	78	123	0%	
Tetrachloroethene	A	ug/L	122.39033	4.8956132		5	0	0	0.0671	0.5	500	98%	74	129	0%	
Toluene	A	ug/L	129.27156	5.1708624		5	0	0	0.0679	0.5	500	103%	80	121	0%	
trans-1,2-Dichloroethene	A	ug/L	129.83227	5.1932908		5	0	0	0.125	0.5	500	104%	75	124	0%	
trans-1,3-Dichloropropene	A	ug/L	132.48323	5.2993292		5	0	0	0.0846	0.5	500	106%	73	127	0%	
Trichloroethene	A	ug/L	125.4471	5.017884		5	0	0	0.0993	0.5	500	100%	79	123	0%	
Trichlorofluoromethane	A	ug/L	124.51499	4.9805996		5	0	0	0.134	0.5	500	100%	65	141	0%	
Vinyl chloride	A	ug/L	122.60319	4.9041276		5	0	0	0.153	0.5	500	98%	58	137	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	383.98208	15.3592832		15	0	0	0.0604	0.5	1500	102%	79	121	0%	
1,2-Dichloroethane-d4	S	ug/L	277.78361	11.1113444		10	0	0	0.229	0.5	500	111%	81	118	0%	
Dibromofluoromethane	S	ug/L	273.52362	10.9409448		10	0	0	0.129	0.5	500	109%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	274.05376	10.9621504		10	0	0	0.149	0.5	500	110%	85	114	0%	
Toluene-d8	S	ug/L	271.7626	10.870504		10	0	0	0.23	0.5	500	109%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973527	B22010219-001	VOC-8260-W-Q	MSD-DOD	)A5975C\VG010´	1/7/2022 5:34:50	1	R373037		1E+07	1E+07	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973527	B22010219-001	VOC-8260-W-Q	MSD-DOD	)A5975C\VG010	1/7/2022 5:34:50	1	R373037		1E+07	1E+07	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	126.97081	5.0788324		5	0	4.9588816	0.101	0.5	500	102%	78	124	2%	
1,1,1-Trichloroethane	A	ug/L	129.67471	5.1869884		5	0	4.9907416	0.131	0.5	500	104%	74	131	4%	
1,1,2,2-Tetrachloroethane	A	ug/L	137.05211	5.4820844		5	0	5.2493156	0.0872	0.5	500	110%	71	121	4%	
1,1,2-Trichloroethane	A	ug/L	128.35043	5.1340172		5	0	4.9423652	0.108	0.5	500	103%	80	119	4%	
1,1-Dichloroethane	A	ug/L	139.44565	5.577826		5	0	5.422852	0.135	0.5	500	112%	77	125	3%	
1,1-Dichloroethene	A	ug/L	134.54038	5.3816152		5	0	5.2431856	0.141	0.5	500	108%	71	131	3%	
1,1-Dichloropropene	A	ug/L	125.60877	5.0243508		5	0	4.8566056	0.083	0.5	500	100%	79	125	3%	
1,2,3-Trichloropropane	A	ug/L	134.57211	5.3828844		5	0	4.9529848	0.235	0.5	500	108%	73	125	8%	
1,2-Dibromoethane	A	ug/L	131.63801	5.2655204		5	0	4.9935576	0.0916	0.5	500	105%	78	122	5%	
1,2-Dichlorobenzene	A	ug/L	132.96807	5.3187228		5	0	5.038362	0.0746	0.5	500	106%	80	119	5%	
1,2-Dichloroethane	A	ug/L	127.77295	5.110918		5	0	5.1145068	0.116	0.5	500	102%	73	128	0%	
1,2-Dichloropropane	A	ug/L	133.30768	5.3323072		5	0	5.013456	0.0847	0.5	500	107%	78	122	6%	
1,3-Dichlorobenzene	A	ug/L	137.9499	5.517996		5	0	5.2558796	0.0803	0.5	500	110%	80	119	5%	
1,3-Dichloropropane	A	ug/L	132.70403	5.3081612		5	0	5.0702732	0.0791	0.5	500	106%	80	119	5%	
1,4-Dichlorobenzene	A	ug/L	137.42865	5.497146		5	0	5.1958564	0.0858	0.5	500	110%	79	118	6%	
2,2-Dichloropropane	A	ug/L	141.24186	5.6496744		5	0	5.2725228	0.186	0.5	500	113%	60	139	7%	
2-Chlorotoluene	A	ug/L	141.2126	5.648504		5	0	5.252534	0.0876	0.5	500	113%	79	122	7%	
4-Chlorotoluene	A	ug/L	142.04727	5.6818908		5	0	5.332192	0.0728	0.5	500	114%	78	122	6%	
Benzene	A	ug/L	133.07465	5.322986		5	0	5.2092352	0.0914	0.5	500	106%	79	120	2%	
Bromobenzene	A	ug/L	142.54012	5.7016048		5	0	5.3445976	0.0831	0.5	500	114%	80	120	6%	
Bromochloromethane	A	ug/L	127.87548	5.1150192		5	0	4.9779088	0.141	0.5	500	102%	78	123	3%	
Bromodichloromethane	A	ug/L	135.38706	5.4154824		5	0	5.1296712	0.12	0.5	500	108%	79	125	5%	
Bromoform	A	ug/L	139.78015	5.591206		5	0	5.1881456	0.119	0.5	500	112%	66	130	7%	
Bromomethane	A	ug/L	119.87937	4.7951748		5	0	4.8394384	0.253	0.5	500	96%	53	141	1%	
Carbon tetrachloride	A	ug/L	124.65019	4.9860076		5	0	4.8677872	0.143	0.5	500	100%	72	136	2%	
Chlorobenzene	A	ug/L	135.94651	5.4378604		5	0	5.1308808	0.0914	0.5	500	109%	82	118	6%	
Chlorodibromomethane	A	ug/L	127.72526	5.1090104		5	0	5.0063784	0.0841	0.5	500	102%	74	126	2%	
Chloroethane	A	ug/L	121.58436	4.8633744		5	0	5.6407292	0.169	0.5	500	97%	60	138	15%	
Chloroform	A	ug/L	126.24502	5.0498008		5	0	4.858966	0.0789	0.5	500	101%	79	124	4%	
Chloromethane	A	ug/L	117.71804	4.7087216		5	0	4.4721952	0.162	0.5	500	94%	50	139	5%	
cis-1,2-Dichloroethene	A	ug/L	133.90986	5.3563944		5	0	5.2511968	0.108	0.5	500	107%	78	123	2%	
cis-1,3-Dichloropropene	A	ug/L	127.54766	5.1019064		5	0	4.841976	0.073	0.5	500	102%	75	124	5%	
Dibromomethane	A	ug/L	129.42028	5.1768112		5	0	5.0311908	0.147	0.5	500	104%	79	123	3%	
Dichlorodifluoromethane	A	ug/L	110.63166	4.4252664		5	0	4.2933368	0.175	0.5	500	89%	32	152	3%	
Ethylbenzene	A	ug/L	135.43417	5.4173668		5	0	5.0822964	0.0836	0.5	500	108%	79	121	6%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973527	B22010219-001	VOC-8260-W-Q	MSD-DOD	)A5975C\VG010´	1/7/2022 5:34:50	1	R373037		1E+07	1E+07	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
m+p-Xylenes	A	ug/L	265.23121	10.6092484		10	0	10.129931	0.15	0.5	1000	106%	80	121	5%	
Methyl ethyl ketone	A	ug/L	1218.10205	48.724082		50	0	50.265366	1.77	10	5000	97%	56	143	3%	
Methyl tert-butyl ether (MTBE)	A	ug/L	142.97082	5.7188328		5	0	5.6050092	0.101	0.5	500	114%	71	124	2%	
Methylene chloride	A	ug/L	125.73719	5.0294876		5	0	4.8779708	0.338	0.5	500	101%	74	124	3%	
o-Xylene	A	ug/L	139.35978	5.5743912		5	0	5.2293524	0.0604	0.5	500	111%	78	122	6%	
Styrene	A	ug/L	141.51056	5.6604224		5	0	5.2968076	0.067	0.5	500	113%	78	123	7%	
Tetrachloroethene	A	ug/L	132.08546	5.2834184		5	0	4.8956132	0.0671	0.5	500	106%	74	129	8%	
Toluene	A	ug/L	138.2772	5.531088		5	0	5.1708624	0.0679	0.5	500	111%	80	121	7%	
trans-1,2-Dichloroethene	A	ug/L	138.08523	5.5234092		5	0	5.1932908	0.125	0.5	500	110%	75	124	6%	
trans-1,3-Dichloropropene	A	ug/L	135.69869	5.4279476		5	0	5.2993292	0.0846	0.5	500	109%	73	127	2%	
Trichloroethene	A	ug/L	133.54359	5.3417436		5	0	5.017884	0.0993	0.5	500	107%	79	123	6%	
Trichlorofluoromethane	A	ug/L	120.34118	4.8136472		5	0	4.9805996	0.134	0.5	500	96%	65	141	3%	
Vinyl chloride	A	ug/L	125.03335	5.001334		5	0	4.9041276	0.153	0.5	500	100%	58	137	2%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	404.59099	16.1836396		15	0	15.359283	0.0604	0.5	1500	108%	79	121	5%	
1,2-Dichloroethane-d4	S	ug/L	272.54518	10.9018072		10	0	0	0.229	0.5	500	109%	81	118	0%	
Dibromofluoromethane	S	ug/L	268.6027	10.744108		10	0	0	0.129	0.5	500	107%	80	119	0%	
p-Bromofluorobenzene	S	ug/L	271.17267	10.8469068		10	0	0	0.149	0.5	500	108%	85	114	0%	
Toluene-d8	S	ug/L	276.39197	11.0556788		10	0	0	0.23	0.5	500	111%	89	112	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973528	CCV010722_Cl	VOC-8260-W-Q	CCV	)A5975C\VG010´	1/7/2022 6:29:09	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,1,1,2-Tetrachloroethane	A	ug/L	116.51857	4.6607428		5	0	0	0.101	0.5	500	93%	50	150	0%	
1,1,1-Trichloroethane	A	ug/L	118.9139	4.756556		5	0	0	0.131	0.5	500	95%	50	150	0%	
1,1,2,2-Tetrachloroethane	A	ug/L	121.30049	4.8520196		5	0	0	0.0872	0.5	500	97%	50	150	0%	
1,1,2-Trichloroethane	A	ug/L	116.98429	4.6793716		5	0	0	0.108	0.5	500	94%	50	150	0%	
1,1-Dichloroethane	A	ug/L	125.05691	5.0022764		5	0	0	0.135	0.5	500	100%	50	150	0%	
1,1-Dichloroethene	A	ug/L	118.78652	4.7514608		5	0	0	0.141	0.5	500	95%	50	150	0%	
1,1-Dichloropropene	A	ug/L	119.866	4.79464		5	0	0	0.083	0.5	500	96%	50	150	0%	
1,2,3-Trichloropropane	A	ug/L	114.97648	4.5990592		5	0	0	0.235	0.5	500	92%	50	150	0%	
1,2-Dibromoethane	A	ug/L	120.00893	4.8003572		5	0	0	0.0916	0.5	500	96%	50	150	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973528	CCV010722_Cl	VOC-8260-W-Q	CCV	)A5975C\VG010	1/7/2022 6:29:09	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dichlorobenzene	A	ug/L	118.73276	4.7493104		5	0	0	0.0746	0.5	500	95%	50	150	0%	
1,2-Dichloroethane	A	ug/L	115.09774	4.6039096		5	0	0	0.116	0.5	500	92%	50	150	0%	
1,2-Dichloropropane	A	ug/L	123.78967	4.9515868		5	0	0	0.0847	0.5	500	99%	50	150	0%	
1,3-Dichlorobenzene	A	ug/L	121.39871	4.8559484		5	0	0	0.0803	0.5	500	97%	50	150	0%	
1,3-Dichloropropane	A	ug/L	121.54909	4.8619636		5	0	0	0.0791	0.5	500	97%	50	150	0%	
1,4-Dichlorobenzene	A	ug/L	120.7629	4.830516		5	0	0	0.0858	0.5	500	97%	50	150	0%	
2,2-Dichloropropane	A	ug/L	121.75209	4.8700836		5	0	0	0.186	0.5	500	97%	50	150	0%	
2-Chlorotoluene	A	ug/L	123.99093	4.9596372		5	0	0	0.0876	0.5	500	99%	50	150	0%	
4-Chlorotoluene	A	ug/L	126.3413	5.053652		5	0	0	0.0728	0.5	500	101%	50	150	0%	
Benzene	A	ug/L	122.52703	4.9010812		5	0	0	0.0914	0.5	500	98%	50	150	0%	
Bromobenzene	A	ug/L	124.73788	4.9895152		5	0	0	0.0831	0.5	500	100%	50	150	0%	
Bromochloromethane	A	ug/L	115.32869	4.6131476		5	0	0	0.141	0.5	500	92%	50	150	0%	
Bromodichloromethane	A	ug/L	118.25683	4.7302732		5	0	0	0.12	0.5	500	95%	50	150	0%	
Bromoform	A	ug/L	120.83294	4.8333176		5	0	0	0.119	0.5	500	97%	50	150	0%	
Bromomethane	A	ug/L	124.27001	4.9708004		5	0	0	0.253	0.5	500	99%	50	150	0%	
Carbon tetrachloride	A	ug/L	117.6182	4.704728		5	0	0	0.143	0.5	500	94%	50	150	0%	
Chlorobenzene	A	ug/L	121.07577	4.8430308		5	0	0	0.0914	0.5	500	97%	50	150	0%	
Chlorodibromomethane	A	ug/L	118.19991	4.7279964		5	0	0	0.0841	0.5	500	95%	50	150	0%	
Chloroethane	A	ug/L	121.35074	4.8540296		5	0	0	0.169	0.5	500	97%	50	150	0%	
Chloroform	A	ug/L	118.39682	4.7358728		5	0	0	0.0789	0.5	500	95%	50	150	0%	
Chloromethane	A	ug/L	114.27651	4.5710604		5	0	0	0.162	0.5	500	91%	50	150	0%	
cis-1,2-Dichloroethene	A	ug/L	122.15956	4.8863824		5	0	0	0.108	0.5	500	98%	50	150	0%	
cis-1,3-Dichloropropene	A	ug/L	119.11853	4.7647412		5	0	0	0.073	0.5	500	95%	50	150	0%	
Dibromomethane	A	ug/L	120.85967	4.8343868		5	0	0	0.147	0.5	500	97%	50	150	0%	
Dichlorodifluoromethane	A	ug/L	114.70512	4.5882048		5	0	0	0.175	0.5	500	92%	50	150	0%	
Ethylbenzene	A	ug/L	121.63922	4.8655688		5	0	0	0.0836	0.5	500	97%	50	150	0%	
m+p-Xylenes	A	ug/L	248.2876	9.931504		10	0	0	0.15	0.5	1000	99%	50	150	0%	
Methyl ethyl ketone	A	ug/L	1168.59215	46.743686		50	0	0	1.77	10	5000	93%	50	150	0%	
Methyl tert-butyl ether (MTBE)	A	ug/L	119.22551	4.7690204		5	0	0	0.101	0.5	500	95%	50	150	0%	
Methylene chloride	A	ug/L	114.81968	4.5927872		5	0	0	0.338	0.5	500	92%	50	150	0%	
o-Xylene	A	ug/L	123.16365	4.926546		5	0	0	0.0604	0.5	500	99%	50	150	0%	
Styrene	A	ug/L	126.27154	5.0508616		5	0	0	0.067	0.5	500	101%	50	150	0%	
Tetrachloroethene	A	ug/L	119.752	4.79008		5	0	0	0.0671	0.5	500	96%	50	150	0%	
Toluene	A	ug/L	123.86607	4.9546428		5	0	0	0.0679	0.5	500	99%	50	150	0%	
trans-1,2-Dichloroethene	A	ug/L	120.86833	4.8347332		5	0	0	0.125	0.5	500	97%	50	150	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
14973528	CCV010722_Cl	VOC-8260-W-Q	CCV	)A5975C\VG010	1/7/2022 6:29:09	1	R373037		0	0	

Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
trans-1,3-Dichloropropene	A	ug/L	124.4317	4.977268		5	0	0	0.0846	0.5	500	100%	50	150	0%	
Trichloroethene	A	ug/L	122.28717	4.8914868		5	0	0	0.0993	0.5	500	98%	50	150	0%	
Trichlorofluoromethane	A	ug/L	105.41691	4.2166764		5	0	0	0.134	0.5	500	84%	50	150	0%	
Vinyl chloride	A	ug/L	117.75408	4.7101632		5	0	0	0.153	0.5	500	94%	50	150	0%	
1,4-Dichlorobenzene-d4	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Chlorobenzene-d5	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Fluorobenzene	I	ug/L	250	10		0	0	0	0	0	500	0%	0	0	0%	
Xylenes, Total	M	ug/L	371.45125	14.85805		15	0	0	0.0604	0.5	1500	99%	50	150	0%	
1,2-Dichloroethane-d4	S	ug/L	273.91948	10.9567792		10	0	0	0.229	0.5	500	110%	50	150	0%	
Dibromofluoromethane	S	ug/L	267.46235	10.698494		10	0	0	0.129	0.5	500	107%	50	150	0%	
p-Bromofluorobenzene	S	ug/L	268.41426	10.7365704		10	0	0	0.149	0.5	500	107%	50	150	0%	
Toluene-d8	S	ug/L	272.36346	10.8945384		10	0	0	0.23	0.5	500	109%	50	150	0%	

DATAFILE HEADERS FROM C:\MSDCHEM\1\DATA\VG010722

```
Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN01.D
Sample Name    : PRIMER
Operator       : MSC
Date injected  :  7 Jan 2022   8:36 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 1
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN02.D
Sample Name    : BFB010722_
Operator       : MSC
Date injected  :  7 Jan 2022   9:03 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 2
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN03.D
Sample Name    : CCV010722_
Operator       : MSC
Date injected  :  7 Jan 2022   9:43 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 3
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN04.D
Sample Name    : LCS010722_
Operator       : MSC
Date injected  :  7 Jan 2022  10:50 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 4
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN05.D
```

```
Sample Name    : BLK
Operator       : MSC
Date injected  :  7 Jan 2022  11:17 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 5
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN06.D
Sample Name    : MBLK010722_
Operator       : MSC
Date injected  :  7 Jan 2022  11:45 am
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 6
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN07.D
Sample Name    : B22010214-002A
Operator       : MSC
Date injected  :  7 Jan 2022  12:24 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 7
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN08.D
Sample Name    : B22010219-002A
Operator       : MSC
Date injected  :  7 Jan 2022  12:52 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 8
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN09.D
Sample Name    : B22010226-002A
Operator       : MSC
Date injected  :  7 Jan 2022   1:19 pm
```

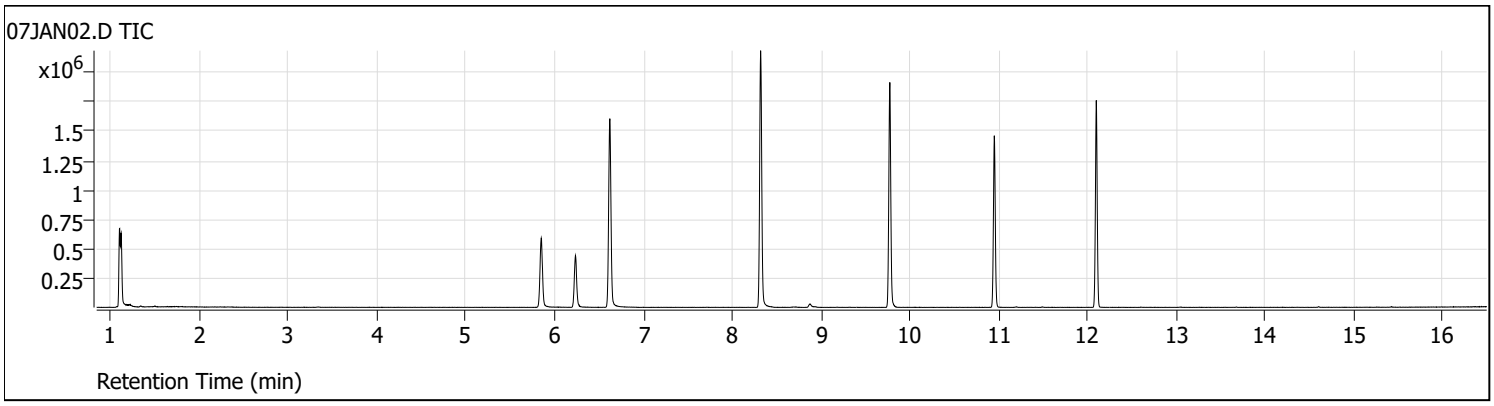
```
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 9
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN10.D
Sample Name    : B22010214-002A
Operator       : MSC
Date injected  :  7 Jan 2022   1:46 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 10
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN11.D
Sample Name    : B22010219-001F
Operator       : MSC
Date injected  :  7 Jan 2022   2:14 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 11
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN12.D
Sample Name    : B22010213-002C
Operator       : MSC
Date injected  :  7 Jan 2022   2:41 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 12
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN13.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  7 Jan 2022   3:08 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
```

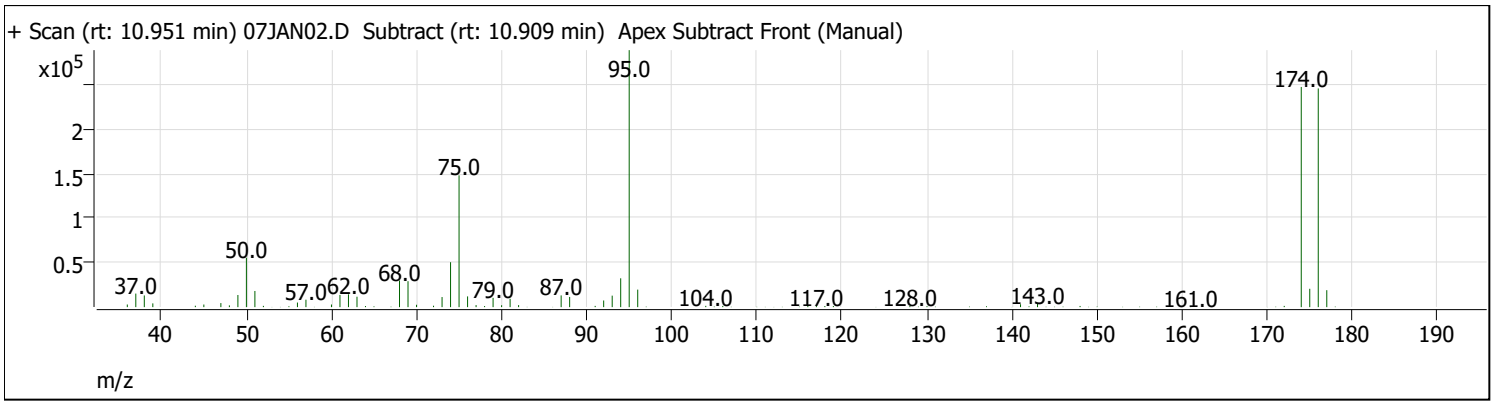
```
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 13
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN14.D
Sample Name    : B22010214-001F
Operator       : MSC
Date injected  :  7 Jan 2022   3:35 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 14
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN15.D
Sample Name    : B22010262-001F
Operator       : MSC
Date injected  :  7 Jan 2022   4:03 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 15
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN16.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  7 Jan 2022   4:40 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.839
End Time       : 16.498
Vial Number    : 16
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN17.D
Sample Name    : B22010219-001FMS
Operator       : MSC
Date injected  :  7 Jan 2022   5:07 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 17
```

```
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN18.D
Sample Name    : B22010219-001FMSD
Operator       : MSC
Date injected  :  7 Jan 2022   5:34 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 18
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN19.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  7 Jan 2022   6:01 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 19
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN20.D
Sample Name    : CCV010722_Closing
Operator       : MSC
Date injected  :  7 Jan 2022   6:29 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 20
---------------------------------------------------------------

Data file Name : C:\MSDCHEM\1\DATA\VG010722\07JAN21.D
Sample Name    : BLK
Operator       : MSC
Date injected  :  7 Jan 2022   6:56 pm
Instrument     : VOA5975C
Method used    : 5975CACQF
No of spectra  : 5616
Start Time     : 0.840
End Time       : 16.498
Vial Number    : 21
---------------------------------------------------------------
```

Tune Evaluation Report

Data Path:	D:\Org\Data\VOA5975C\VG010722\07JAN02.D
Acq on:	1/7/2022 9:03:26 AM
Operator:	MSC
Sample:	BFB010722_
Inst Name:	VOA5975C
ALS Vial:	2
Method:	\\MASSHUNTER\Org\Data\Methods\BFBapex.m





Target Mass	Rel. To Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Pass/Fail
50	95	15	40	19.2	55384	Pass
75	95	30	60	51.3	147840	Pass
95	95	100	100	100.0	288256	Pass
96	95	5	9	6.8	19656	Pass
173	174	0	2	0.0	0	Pass
174	95	50	100	85.7	246976	Pass
175	174	5	9	8.3	20592	Pass
176	174	95	101	99.3	245312	Pass
177	176	5	9	7.7	19000	Pass

Continuing Calibration Report

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Batch Name D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin
Method File \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m
Daily CC D:\Org\Data\VOA5975C\VG01072207JAN03.D

Level name	Injection Time	Calibration Files
1	1/4/2022 3:33:04 PM	D:\Org\Data\VOA5975C\VG010422\04JAN10.D
2	1/4/2022 4:00:35 PM	D:\Org\Data\VOA5975C\VG010422\04JAN11.D
3	1/4/2022 4:28:05 PM	D:\Org\Data\VOA5975C\VG010422\04JAN12.D
4	1/4/2022 4:55:32 PM	D:\Org\Data\VOA5975C\VG010422\04JAN13.D
5	1/4/2022 5:50:25 PM	D:\Org\Data\VOA5975C\VG010422\04JAN15.D
6	1/4/2022 6:45:10 PM	D:\Org\Data\VOA5975C\VG010422\04JAN17.D
7	1/4/2022 7:39:45 PM	D:\Org\Data\VOA5975C\VG010422\04JAN19.D
8	1/4/2022 8:34:31 PM	D:\Org\Data\VOA5975C\VG010422\04JAN21.D
CC	1/7/2022 9:43:06 AM	D:\Org\Data\VOA5975C\VG010722\07JAN03.D <======

ISTD Compound:	Avg Resp	Mid Resp	CC Resp	Area%	A/M
Fluorobenzene	805964	778120	868396	111.60	M
Chlorobenzene-d5	305684	300356	338774	112.79	M
1,4-Dichlorobenzene-d4	252451	248636	271840	109.33	M

Target Compound	AvgRF/R2	CC RF	Exp. Conc	Calc. Conc	%Dev	Area%	Curve Fit
Fluorobenzene	----------ISTD----------						
Dichlorodifluoromethane	0.3276	0.3000	125.00	114.49	8.41	94.45	Avg RF
Chloromethane	0.3976	0.3703	125.00	116.40	6.88	100.11	Avg RF
Vinyl chloride	0.3578	0.3477	125.00	121.48	2.81	101.77	Avg RF
Bromomethane	0.1600	0.1693	125.00	132.28	-5.82	112.81	Avg RF
Chloroethane	0.1771	0.1673	125.00	118.04	5.56	101.69	Avg RF
Trichlorofluoromethane	0.4441	0.3994	125.00	112.42	10.07	91.85	Avg RF
1,1-Dichloroethene	0.2518	0.2447	125.00	121.47	2.82	106.85	Avg RF
Methylene chloride	0.3712	0.3545	125.00	119.36	4.51	113.78	Avg RF
trans-1,2-Dichloroethene	0.2569	0.2578	125.00	125.43	-0.35	111.48	Avg RF
Methyl tert-butyl ether (MTBE)	0.3321	0.3396	125.00	127.82	-2.26	106.02	Avg RF
1,1-Dichloroethane	0.4782	0.4891	125.00	127.83	-2.27	114.13	Avg RF
2,2-Dichloropropane	0.3583	0.3729	125.00	130.08	-4.07	115.94	Avg RF
cis-1,2-Dichloroethene	0.2605	0.2682	125.00	128.71	-2.97	116.38	Avg RF
Methyl ethyl ketone	0.0353	0.0355 #	1250.00	1257.21	-0.58	114.36	Avg RF
Bromochloromethane	0.1079	0.1110	125.00	128.60	-2.88	114.86	Avg RF
Chloroform	0.4759	0.4629	125.00	121.57	2.74	111.88	Avg RF
1,1,1-Trichloroethane	0.4460	0.4463	125.00	125.09	-0.07	111.25	Avg RF
Dibromofluoromethane	0.2355	0.2633	250.00	279.47	-11.79	256.01	Avg RF
Carbon tetrachloride	0.4394	0.4228	125.00	120.28	3.78	106.17	Avg RF
1,1-Dichloropropene	0.3792	0.3811	125.00	125.61	-0.49	110.56	Avg RF
1,2-Dichloroethane-d4	0.1017	0.1142	250.00	280.59	-12.24	253.68	Avg RF
Benzene	0.9954	1.0119	125.00	127.07	-1.66	114.58	Avg RF
1,2-Dichloroethane	0.2693	0.2665	125.00	123.72	1.03	110.36	Avg RF
Chlorobenzene-d5	----------ISTD----------						
Trichloroethene	0.7540	0.7286	125.00	120.79	3.37	108.14	Avg RF
1,2-Dichloropropane	0.6632	0.6756	125.00	127.33	-1.87	115.37	Avg RF
Dibromomethane	0.2803	0.2729	125.00	121.69	2.65	113.76	Avg RF
Bromodichloromethane	0.7735	0.7668	125.00	123.91	0.87	112.29	Avg RF
cis-1,3-Dichloropropene	0.8745	0.8551	125.00	122.22	2.22	111.92	Avg RF
Toluene-d8	2.4091	2.5975	250.00	269.54	-7.82	245.67	Avg RF
Toluene	1.6274	1.6559	125.00	127.19	-1.75	114.62	Avg RF
trans-1,3-Dichloropropene	0.6225	0.6351	125.00	127.53	-2.02	116.03	Avg RF
1,1,2-Trichloroethane	0.3242	0.3204	125.00	123.52	1.18	116.28	Avg RF
Tetrachloroethene	0.6639	0.6258	125.00	117.83	5.73	108.63	Avg RF

Target Compound	AvgRF/R2	CC RF	Exp. Conc	Calc. Conc	%Dev	Area%	Curve Fit
1,3-Dichloropropane	0.6378	0.6355	125.00	124.56	0.36	111.92	Avg RF
Chlorodibromomethane	0.5068	0.5066	125.00	124.97	0.03	114.40	Avg RF
1,2-Dibromoethane	0.3545	0.3552	125.00	125.23	-0.18	116.09	Avg RF
Chlorobenzene	1.7817	1.7567	125.00	123.25	1.40	112.88	Avg RF
1,1,1,2-Tetrachloroethane	0.6228	0.5884	125.00	118.09	5.53	109.64	Avg RF
Ethylbenzene	3.0900	3.0635	125.00	123.93	0.86	111.80	Avg RF
m+p-Xylenes	1.2008	1.2117	250.00	252.27	-0.91	111.42	Avg RF
o-Xylene	1.0690	1.0664	125.00	124.70	0.24	111.84	Avg RF
Styrene	1.7211	1.7876	125.00	129.83	-3.86	112.83	Avg RF
1,4-Dichlorobenzene-d4	----------ISTD----------						
Bromoform	0.3199	0.3294	125.00	128.72	-2.98	114.33	Avg RF
p-Bromofluorobenzene	0.9159	1.0133	250.00	276.61	-10.64	241.07	Avg RF
Bromobenzene	0.8091	0.8422	125.00	130.12	-4.09	111.93	Avg RF
1,1,2,2-Tetrachloroethane	0.4657	0.4865	125.00	130.60	-4.48	116.10	Avg RF
1,2,3-Trichloropropane	0.1246	0.1295	125.00	129.90	-3.92	53.15	Avg RF
2-Chlorotoluene	0.8050	0.8325	125.00	129.26	-3.41	110.47	Avg RF
4-Chlorotoluene	2.6247	2.7862	125.00	132.69	-6.15	112.66	Avg RF
1,3-Dichlorobenzene	1.4756	1.4739	125.00	124.86	0.11	109.23	Avg RF
1,4-Dichlorobenzene	1.5046	1.5178	125.00	126.10	-0.88	109.13	Avg RF
1,2-Dichlorobenzene	1.2470	1.2379	125.00	124.08	0.74	110.48	Avg RF

A -- against Average; M -- against Mid Point; P -- against Previous CC in the Method;

Continuing Calibration Report

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Batch Name D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin
Method File \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m
Daily CC D:\Org\Data\VOA5975C\VG01072207JAN20.D

Level name	Injection Time	Calibration Files
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5	1/4/2022 5:50:25 PM	D:\Org\Data\VOA5975C\VG010422\04JAN15.D
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CC	1/7/2022 6:29:09 PM	D:\Org\Data\VOA5975C\VG010722\07JAN20.D <======

ISTD Compound:	Avg Resp	Mid Resp	CC Resp	Area%	A/M
Fluorobenzene	805964	778120	875212	112.48	M
Chlorobenzene-d5	305684	300356	332712	110.77	M
1,4-Dichlorobenzene-d4	252451	248636	273694	110.08	M

Target Compound	AvgRF/R2	CC RF	Exp. Conc	Calc. Conc	%Dev	Area%	Curve Fit
Fluorobenzene	----------ISTD----------						
Dichlorodifluoromethane	0.3276	0.3006	125.00	114.71	8.24	95.38	Avg RF
Chloromethane	0.3976	0.3635	125.00	114.28	8.58	99.05	Avg RF
Vinyl chloride	0.3578	0.3371	125.00	117.75	5.80	99.42	Avg RF
Bromomethane	0.1600	0.1591	125.00	124.27	0.58	106.81	Avg RF
Chloroethane	0.1771	0.1720	125.00	121.35	2.92	105.36	Avg RF
Trichlorofluoromethane	0.4441	0.3745	125.00	105.42	15.67	86.81	Avg RF
1,1-Dichloroethene	0.2518	0.2393	125.00	118.79	4.97	105.31	Avg RF
Methylene chloride	0.3712	0.3410	125.00	114.82	8.14	110.31	Avg RF
trans-1,2-Dichloroethene	0.2569	0.2484	125.00	120.87	3.31	108.27	Avg RF
Methyl tert-butyl ether (MTBE)	0.3321	0.3167	125.00	119.23	4.62	99.67	Avg RF
1,1-Dichloroethane	0.4782	0.4784	125.00	125.06	-0.05	112.53	Avg RF
2,2-Dichloropropane	0.3583	0.3490	125.00	121.75	2.60	109.36	Avg RF
cis-1,2-Dichloroethene	0.2605	0.2546	125.00	122.16	2.27	111.33	Avg RF
Methyl ethyl ketone	0.0353	0.0330 #	1250.00	1168.59	6.51	107.13	Avg RF
Bromochloromethane	0.1079	0.0996 #	125.00	115.33	7.74	103.81	Avg RF
Chloroform	0.4759	0.4508	125.00	118.40	5.28	109.81	Avg RF
1,1,1-Trichloroethane	0.4460	0.4243	125.00	118.91	4.87	106.58	Avg RF
Dibromofluoromethane	0.2355	0.2520	250.00	267.46	-6.98	246.94	Avg RF
Carbon tetrachloride	0.4394	0.4135	125.00	117.62	5.91	104.64	Avg RF
1,1-Dichloropropene	0.3792	0.3637	125.00	119.87	4.11	106.34	Avg RF
1,2-Dichloroethane-d4	0.1017	0.1115	250.00	273.92	-9.57	249.59	Avg RF
Benzene	0.9954	0.9757	125.00	122.53	1.98	111.34	Avg RF
1,2-Dichloroethane	0.2693	0.2479	125.00	115.10	7.92	103.48	Avg RF
Chlorobenzene-d5	----------ISTD----------						
Trichloroethene	0.7540	0.7376	125.00	122.29	2.17	107.52	Avg RF
1,2-Dichloropropane	0.6632	0.6568	125.00	123.79	0.97	110.16	Avg RF
Dibromomethane	0.2803	0.2710	125.00	120.86	3.31	110.96	Avg RF
Bromodichloromethane	0.7735	0.7318	125.00	118.26	5.39	105.25	Avg RF
cis-1,3-Dichloropropene	0.8745	0.8334	125.00	119.12	4.71	107.12	Avg RF
Toluene-d8	2.4091	2.6246	250.00	272.36	-8.95	243.80	Avg RF
Toluene	1.6274	1.6126	125.00	123.87	0.91	109.63	Avg RF
trans-1,3-Dichloropropene	0.6225	0.6197	125.00	124.43	0.45	111.18	Avg RF
1,1,2-Trichloroethane	0.3242	0.3035	125.00	116.98	6.41	108.16	Avg RF
Tetrachloroethene	0.6639	0.6360	125.00	119.75	4.20	108.42	Avg RF

Target Compound	AvgRF/R2	CC RF	Exp. Conc	Calc. Conc	%Dev	Area%	Curve Fit
1,3-Dichloropropane	0.6378	0.6202	125.00	121.55	2.76	107.26	Avg RF
Chlorodibromomethane	0.5068	0.4792	125.00	118.20	5.44	106.27	Avg RF
1,2-Dibromoethane	0.3545	0.3404	125.00	120.01	3.99	109.26	Avg RF
Chlorobenzene	1.7817	1.7257	125.00	121.08	3.14	108.90	Avg RF
1,1,1,2-Tetrachloroethane	0.6228	0.5805	125.00	116.52	6.79	106.25	Avg RF
Ethylbenzene	3.0900	3.0069	125.00	121.64	2.69	107.77	Avg RF
m+p-Xylenes	1.2008	1.1926	250.00	248.29	0.68	107.70	Avg RF
o-Xylene	1.0690	1.0533	125.00	123.16	1.47	108.49	Avg RF
Styrene	1.7211	1.7386	125.00	126.27	-1.02	107.77	Avg RF
1,4-Dichlorobenzene-d4	----------ISTD----------						
Bromoform	0.3199	0.3093	125.00	120.83	3.33	108.06	Avg RF
p-Bromofluorobenzene	0.9159	0.9833	250.00	268.41	-7.37	235.53	Avg RF
Bromobenzene	0.8091	0.8074	125.00	124.74	0.21	108.04	Avg RF
1,1,2,2-Tetrachloroethane	0.4657	0.4519	125.00	121.30	2.96	108.57	Avg RF
1,2,3-Trichloropropane	0.1246	0.1146	125.00	114.98	8.02	47.36	Avg RF
2-Chlorotoluene	0.8050	0.7985	125.00	123.99	0.81	106.69	Avg RF
4-Chlorotoluene	2.6247	2.6529	125.00	126.34	-1.07	108.00	Avg RF
1,3-Dichlorobenzene	1.4756	1.4331	125.00	121.40	2.88	106.93	Avg RF
1,4-Dichlorobenzene	1.5046	1.4536	125.00	120.76	3.39	105.22	Avg RF
1,2-Dichlorobenzene	1.2470	1.1845	125.00	118.73	5.01	106.44	Avg RF

A -- against Average; M -- against Mid Point; P -- against Previous CC in the Method;

Quantitation Results Report (QT Reviewed)

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Data File	07JAN03.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 9:43:06 AM
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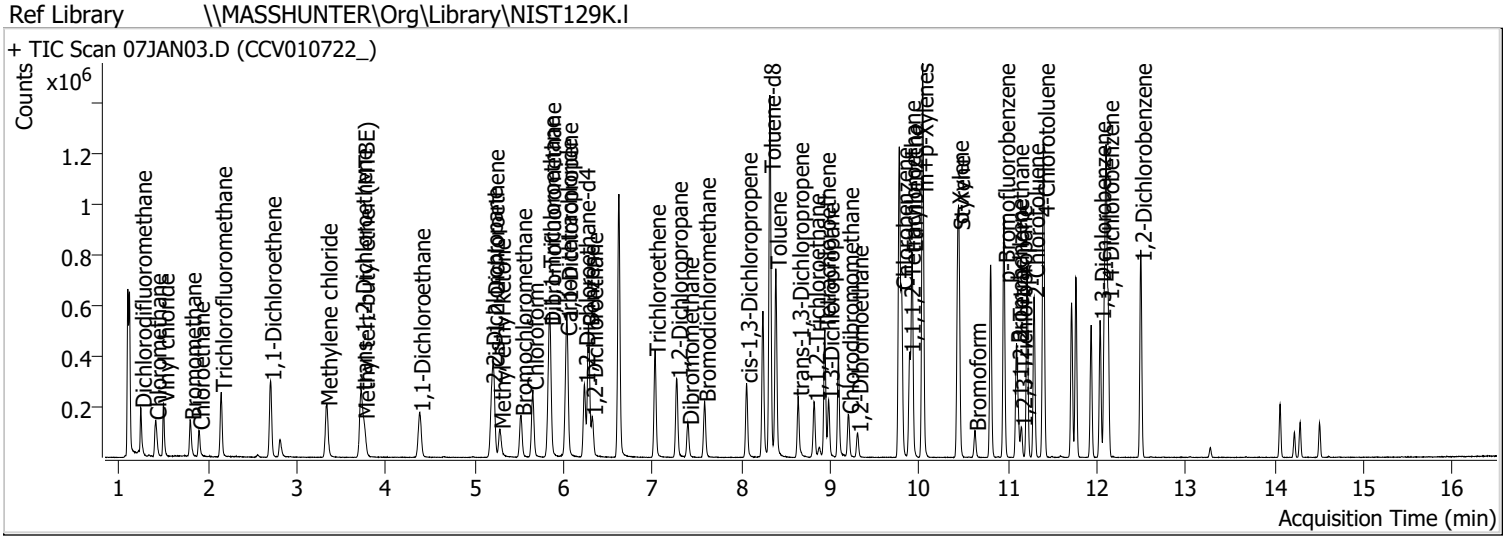
Vial	3	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

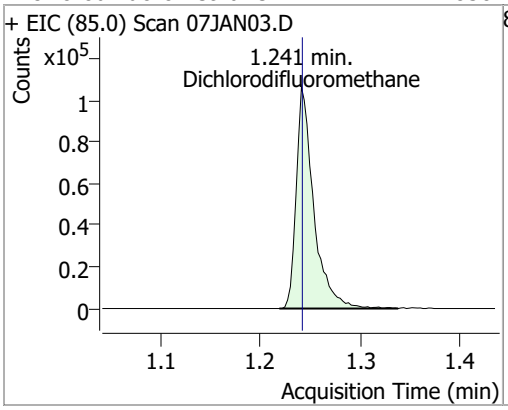


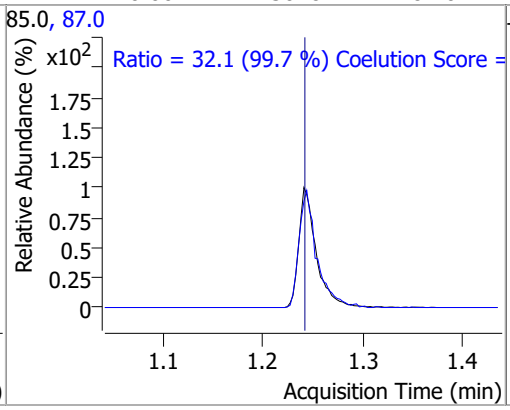
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.620	96.0	868396	250.0000	ng	-0.003
M Chlorobenzene-d5	9.774	82.0	338774	250.0000	ng	0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	271840	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.848	113.0	228638	279.4685	ng	0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 111.79%		
S 1,2-Dichloroethane-d4	6.233	67.0	99152	280.5917	ng	0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 112.24%		
S Toluene-d8	8.319	98.0	879956	269.5449	ng	0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 107.82%		
S p-Bromofluorobenzene	10.954	95.0	275469	276.6060	ng	0.000
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 110.64%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.241	85.0	130281	114.4850	ng	100
T Chloromethane	1.408	50.0	160773	116.3992	ng	100
T Vinyl chloride	1.498	62.0	150982	121.4824	ng	97
T Bromomethane	1.802	96.0	73513	132.2812	ng	97
T Chloroethane	1.897	64.0	72630	118.0450	ng	99
T Trichlorofluoromethane	2.147	101.0	173418	112.4176	ng	100
T 1,1-Dichloroethene	2.702	96.0	106254	121.4726	ng	97
T Methylene chloride	3.330	49.0	153910	119.3588	ng	99
T trans-1,2-Dichloroethene	3.720	96.0	111936	125.4320	ng	98
T Methyl tert-butyl ether (MTBE)	3.751	73.0	147444	127.8241	ng	98
T 1,1-Dichloroethane	4.378	63.0	212345	127.8331	ng	98
T 2,2-Dichloropropane	5.190	77.0	161912	130.0823	ng	99
T cis-1,2-Dichloroethene	5.209	96.0	116451	128.7076	ng	97
T Methyl ethyl ketone	5.282	43.0	154077	1257.2149	ng	98
T Bromochloromethane	5.522	128.0	48203	128.6024	ng	99
T Chloroform	5.653	83.0	200976	121.5715	ng	99

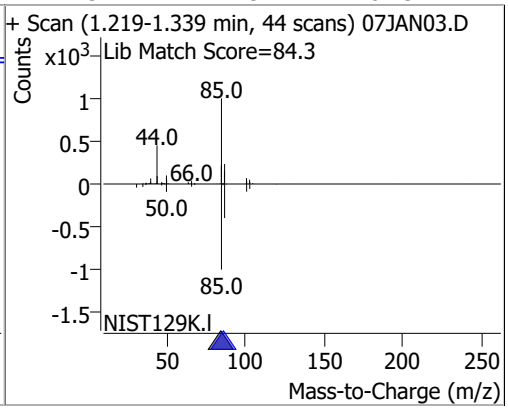
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.834	97.0	193800	125.0917	ng	98
T Carbon tetrachloride	6.026	117.0	183596	120.2774	ng	99
T 1,1-Dichloropropene	6.038	75.0	165459	125.6069	ng	99
T Benzene	6.280	78.0	439363	127.0728	ng	100
T 1,2-Dichloroethane	6.319	62.0	115719	123.7158	ng	97
T Trichloroethene	7.025	95.0	123412	120.7910	ng	97
T 1,2-Dichloropropane	7.270	63.0	114436	127.3316	ng	98
T Dibromomethane	7.401	93.0	46218	121.6934	ng	97
T Bromodichloromethane	7.585	83.0	129880	123.9145	ng	98
T cis-1,3-Dichloropropene	8.057	75.0	144840	122.2214	ng	98
T Toluene	8.386	92.0	280492	127.1937	ng	98
T trans-1,3-Dichloropropene	8.637	75.0	107578	127.5302	ng	98
T 1,1,2-Trichloroethane	8.818	83.0	54273	123.5213	ng	95
T Tetrachloroethene	8.938	163.8	106010	117.8337	ng	99
T 1,3-Dichloropropane	8.980	76.0	107647	124.5555	ng	99
T Chlorodibromomethane	9.200	129.0	85816	124.9678	ng	96
T 1,2-Dibromoethane	9.306	107.0	60164	125.2299	ng	99
T Chlorobenzene	9.802	112.0	297566	123.2509	ng	99
T 1,1,1,2-Tetrachloroethane	9.892	131.0	99659	118.0854	ng	100
T Ethylbenzene	9.919	91.0	518912	123.9274	ng	100
T m+p-Xylenes	10.037	106.0	410489	252.2653	ng	100
T o-Xylene	10.433	106.0	180632	124.6951	ng	99
T Styrene	10.446	104.0	302796	129.8291	ng	99
T Bromoform	10.628	172.5	44778	128.7230	ng	99
T Bromobenzene	11.093	156.0	114469	130.1161	ng	98
T 1,1,2,2-Tetrachloroethane	11.110	83.0	66129	130.5981	ng	96
T 1,2,3-Trichloropropane	11.149	110.0	17600	129.9023	ng	100
T 2-Chlorotoluene	11.289	126.0	113150	129.2634	ng	99
T 4-Chlorotoluene	11.400	91.0	378702	132.6910	ng	99
T 1,3-Dichlorobenzene	12.033	146.0	200331	124.8574	ng	99
T 1,4-Dichlorobenzene	12.125	146.0	206297	126.0981	ng	99
T 1,2-Dichlorobenzene	12.493	146.0	168250	124.0802	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

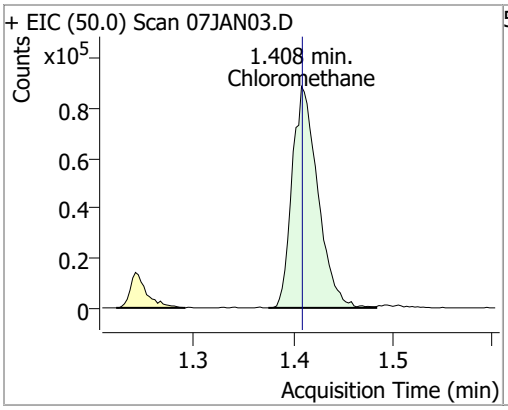
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	114.4850	1.24	0.00	130281	87.0	32.1	2.3	62.3

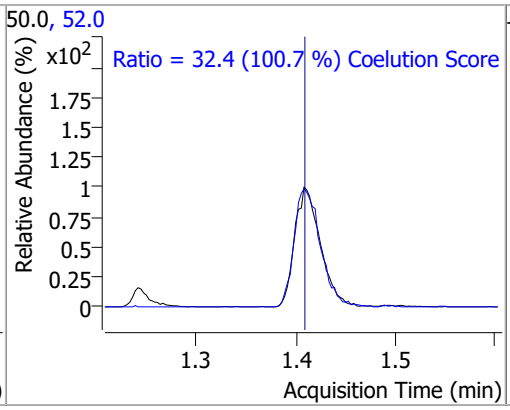


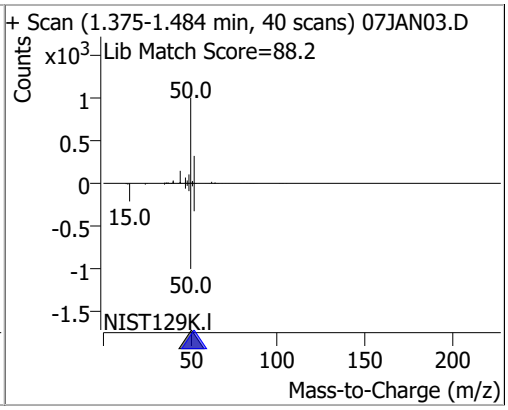




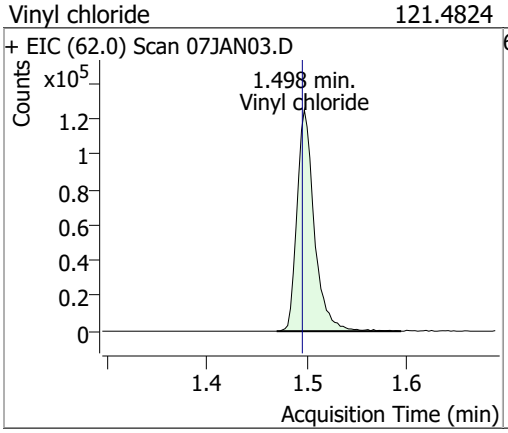
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	116.3992	1.41	0.00	160773	52.0	32.4	2.1	62.1

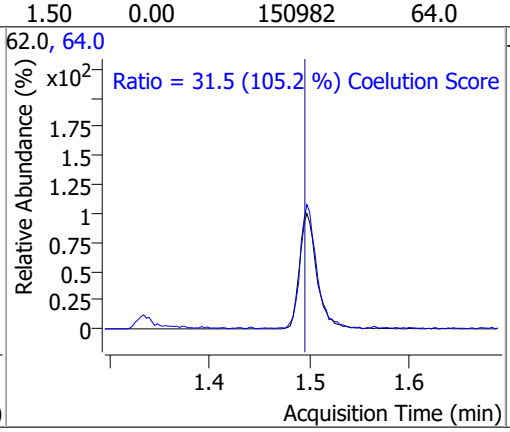


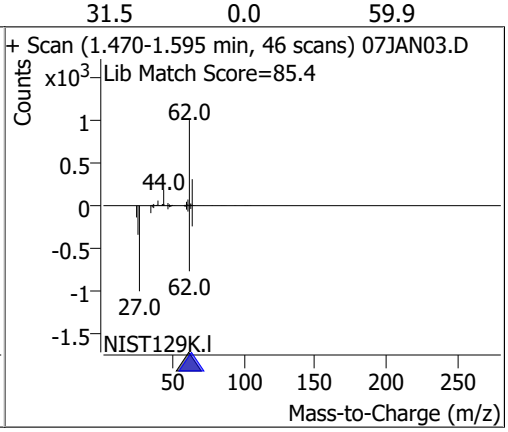




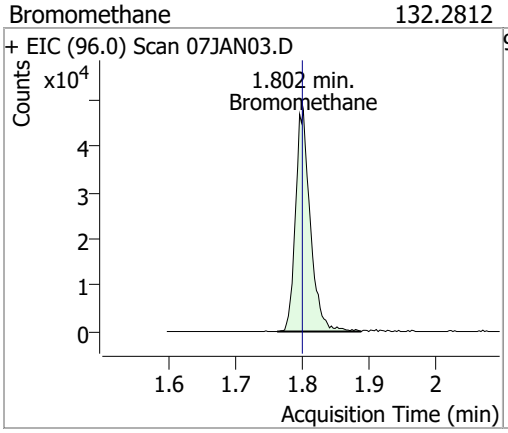
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	121.4824	1.50	0.00	150982	64.0	31.5	0.0	59.9

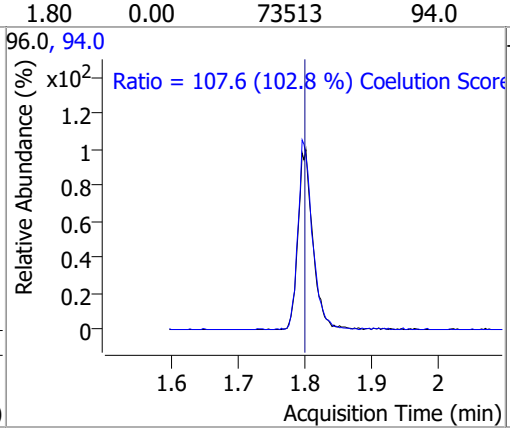


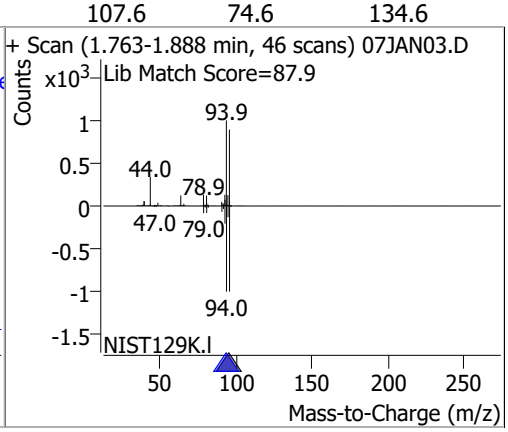




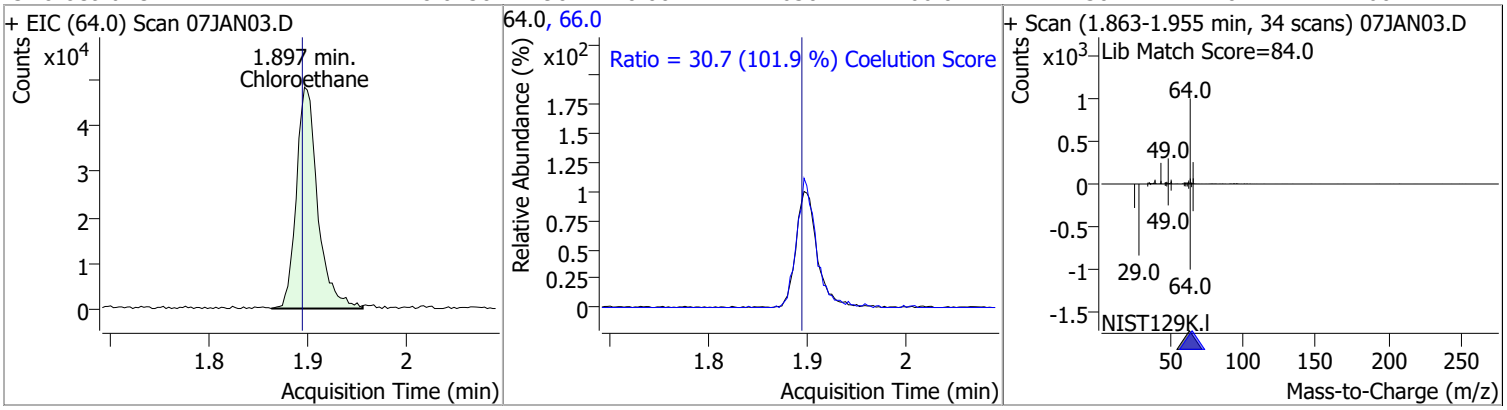
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	132.2812	1.80	0.00	73513	94.0	107.6	74.6	134.6



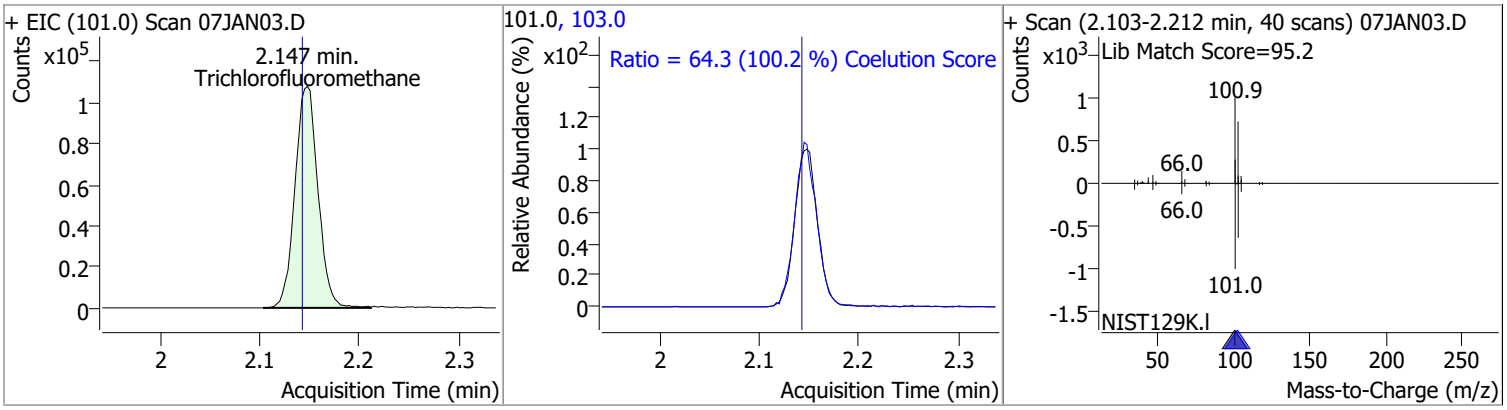




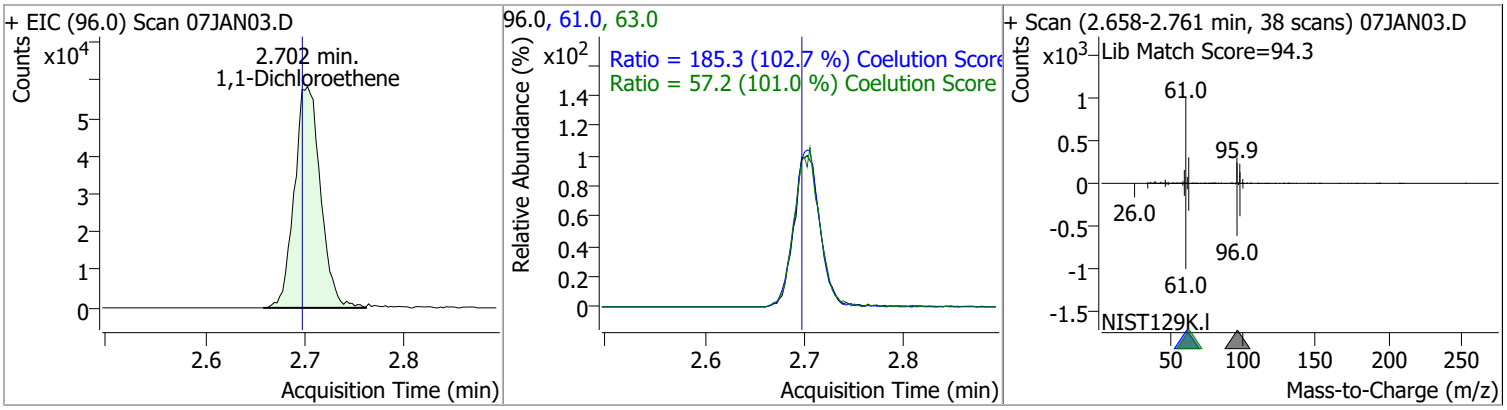
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	118.0450	1.90	0.00	72630	66.0	30.7	0.1	60.1



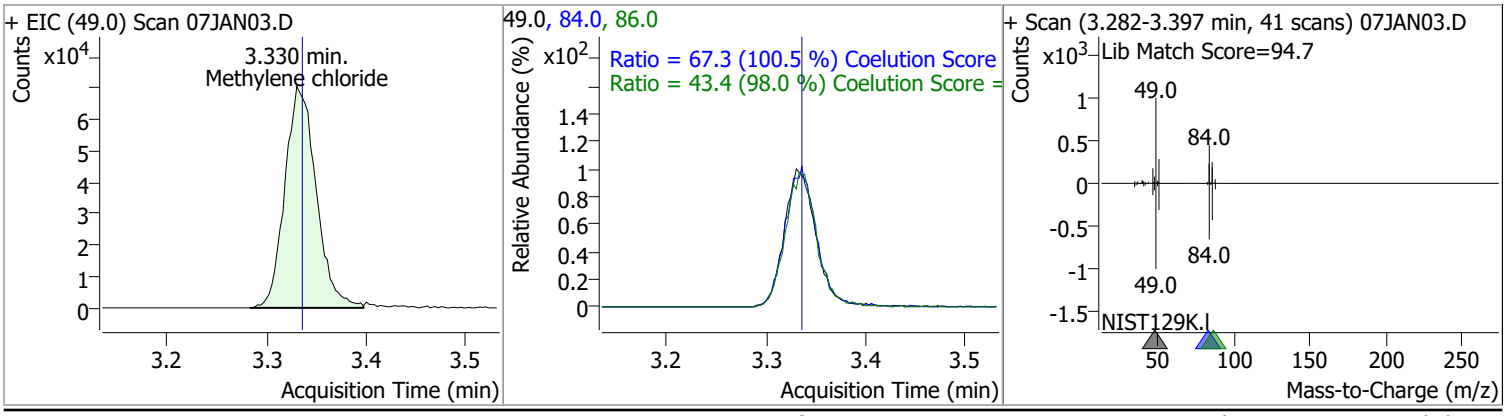
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	112.4176	2.15	0.01	173418	103.0	64.3	34.2	94.2



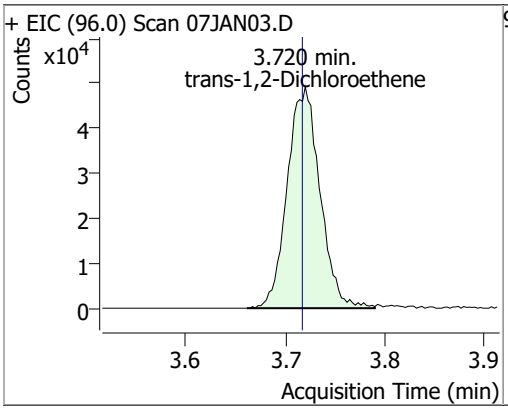
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	121.4726	2.70	0.01	106254	61.0	185.3	150.3	210.3
					63.0	57.2	26.7	86.7

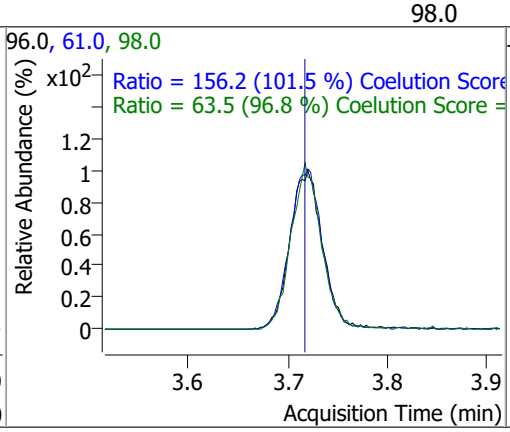


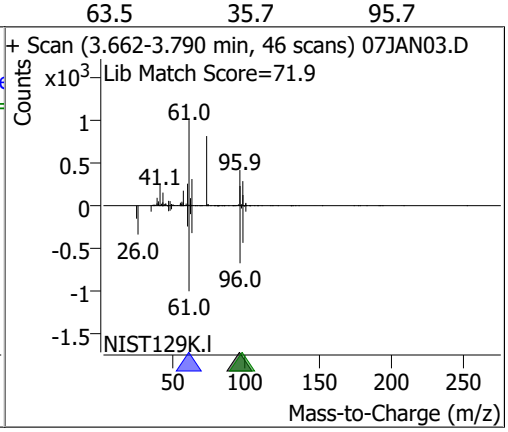
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	119.3588	3.33	-0.01	153910	84.0	67.3	36.9	96.9
					86.0	43.4	14.3	74.3



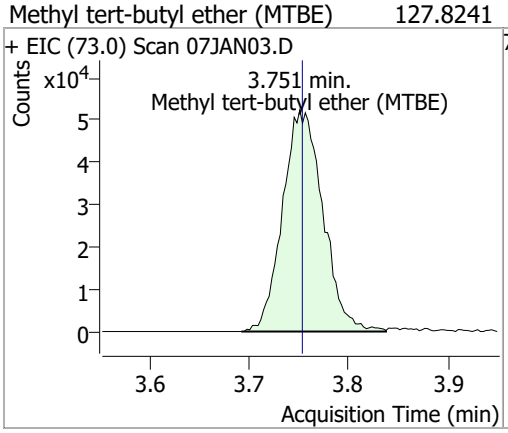
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	125.4320	3.72	0.00	111936	61.0	156.2	123.9	183.9
					98.0	63.5	35.7	95.7

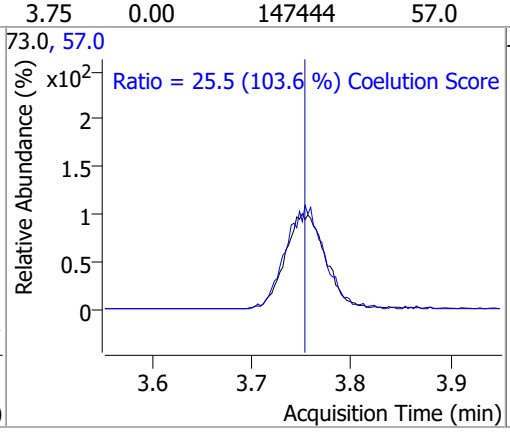


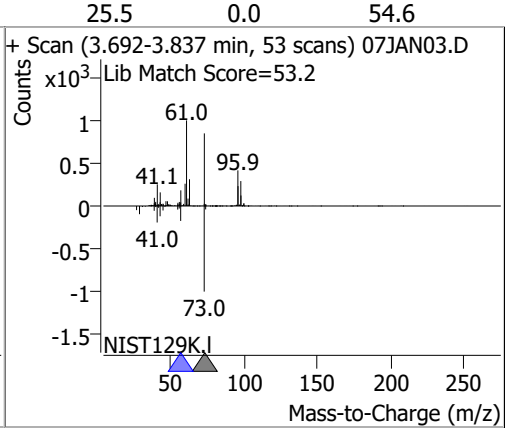




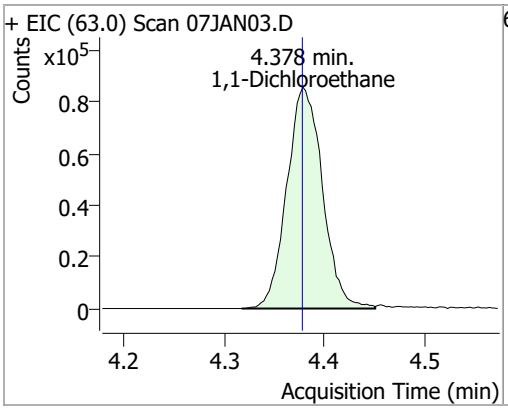
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	127.8241	3.75	0.00	147444	57.0	25.5	0.0	54.6

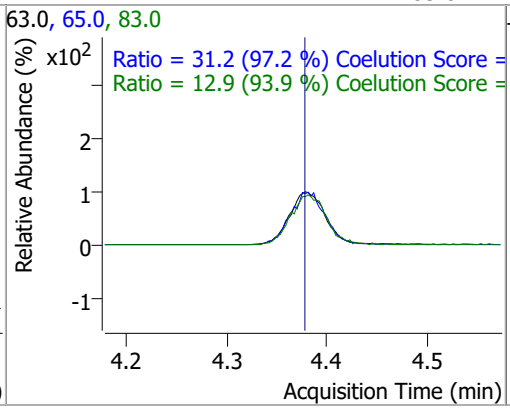


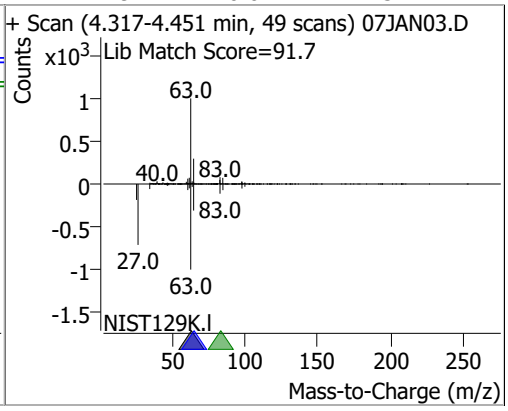




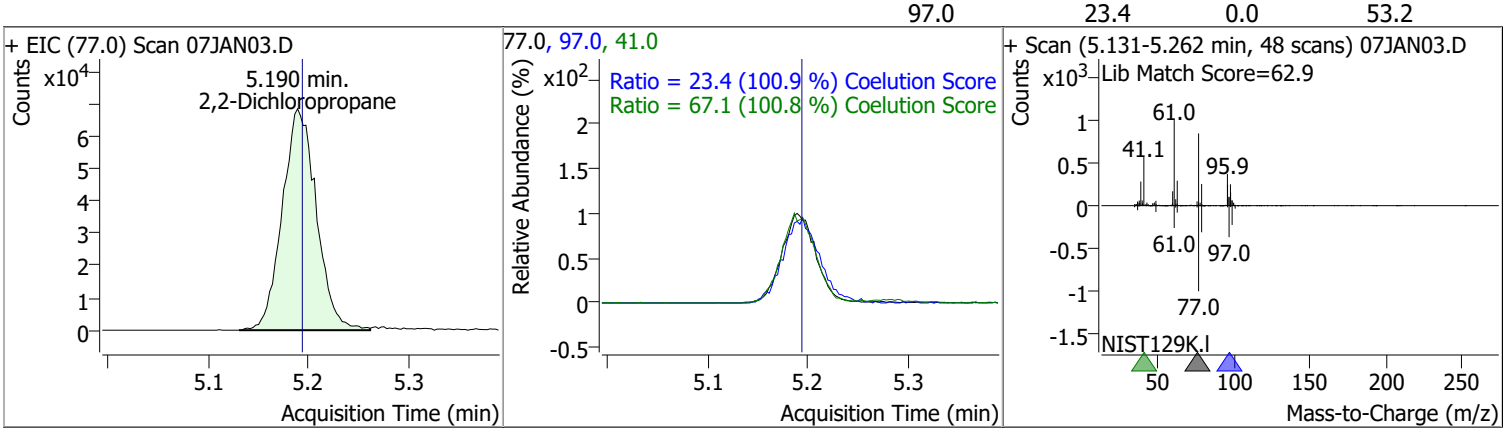
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	127.8331	4.38	0.00	212345	65.0	31.2	2.1	62.1
					83.0	12.9	0.0	43.7



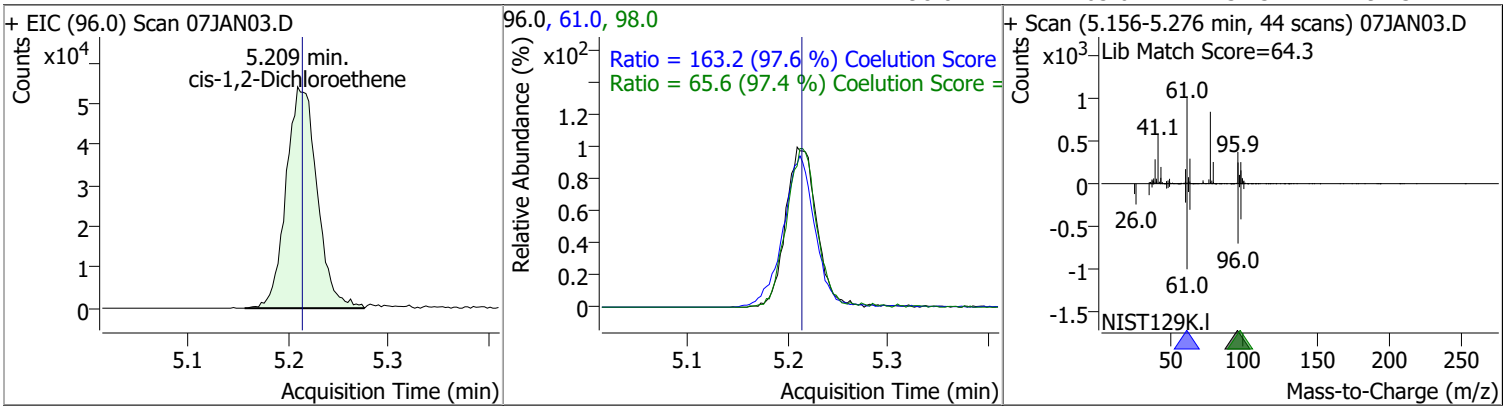




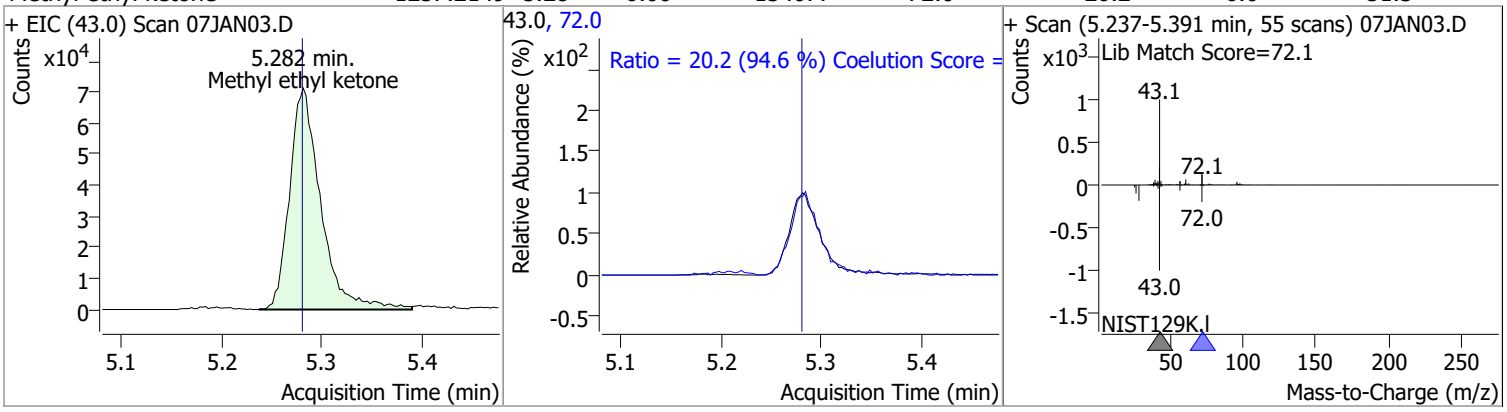
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	130.0823	5.19	-0.01	161912	41.0	67.1	36.5	96.5
					97.0	23.4	0.0	53.2



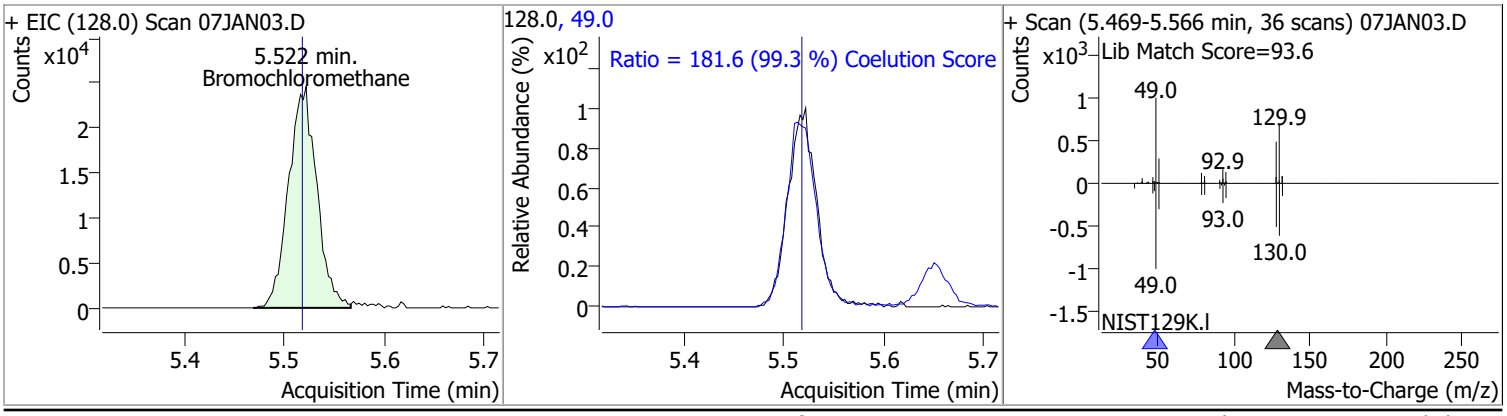
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	128.7076	5.21	-0.01	116451	61.0	163.2	137.2	197.2
					98.0	65.6	37.3	97.3



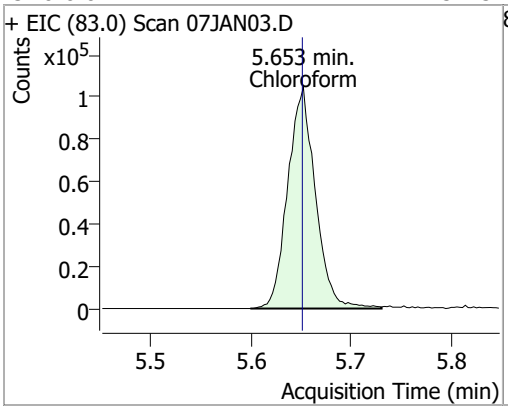
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1257.2149	5.28	0.00	154077	72.0	20.2	0.0	51.3

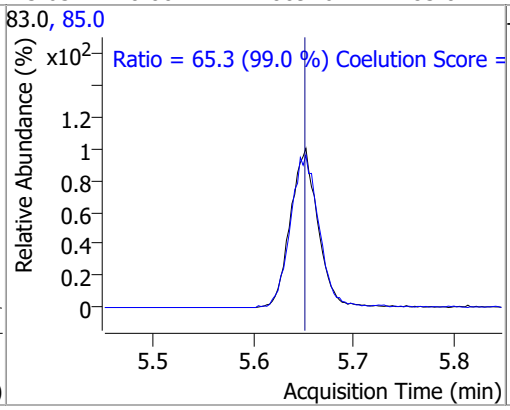


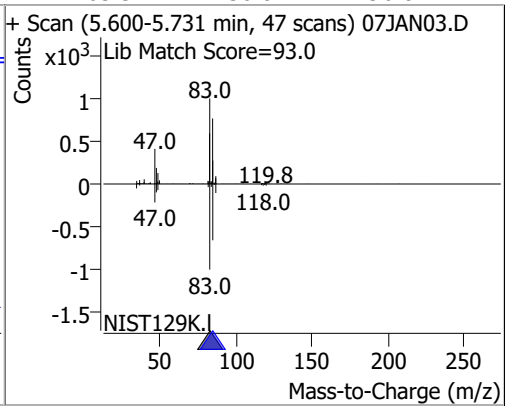
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	128.6024	5.52	0.00	48203	49.0	181.6	152.9	212.9



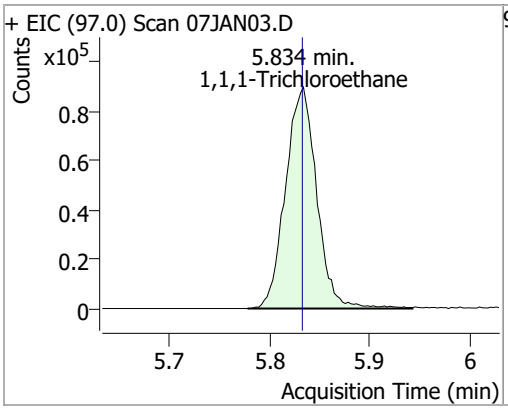
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	121.5715	5.65	0.00	200976	85.0	65.3	36.0	96.0

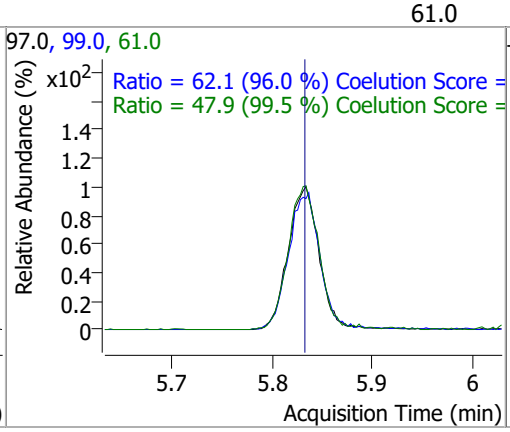


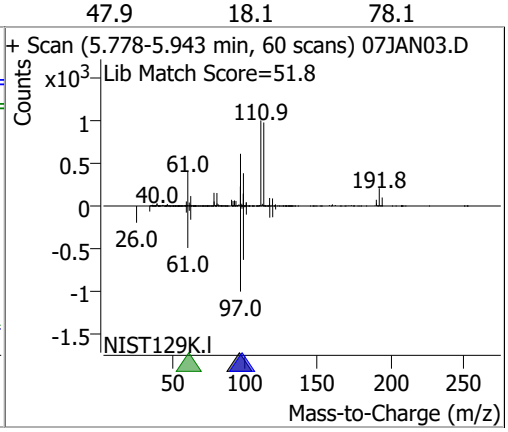




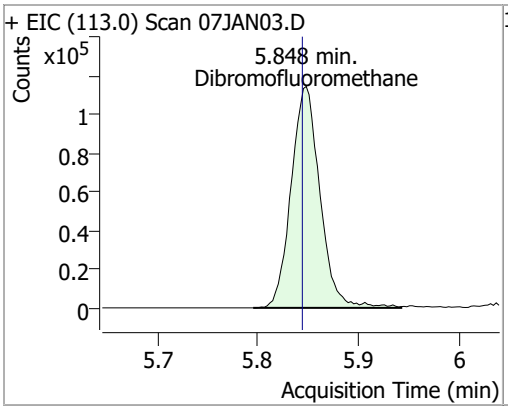
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	125.0917	5.83	0.00	193800	99.0	62.1	34.7	94.7
					61.0	47.9	18.1	78.1

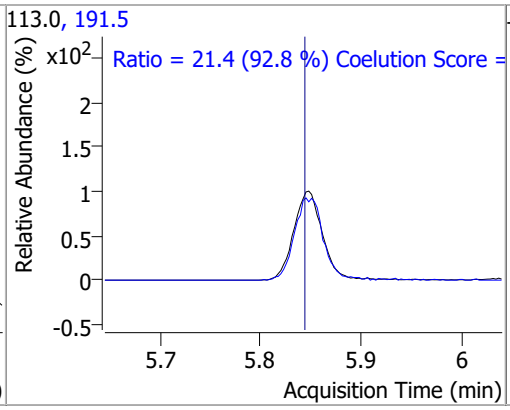


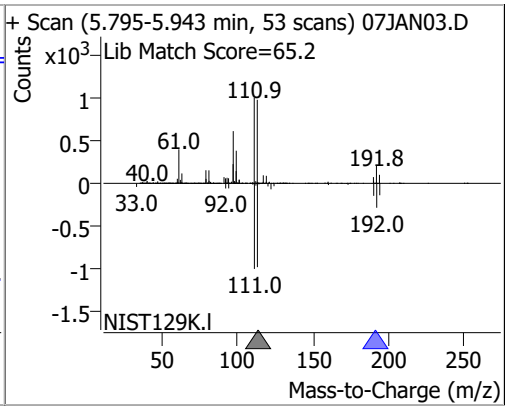




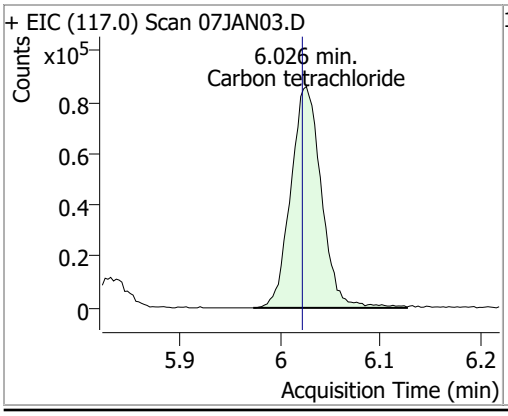
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	279.4685	5.85	0.00	228638	191.5	21.4	0.0	53.1

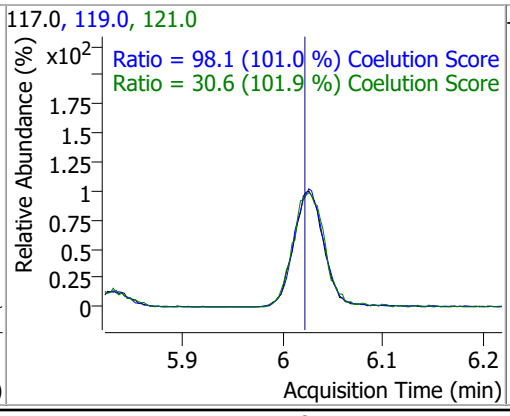


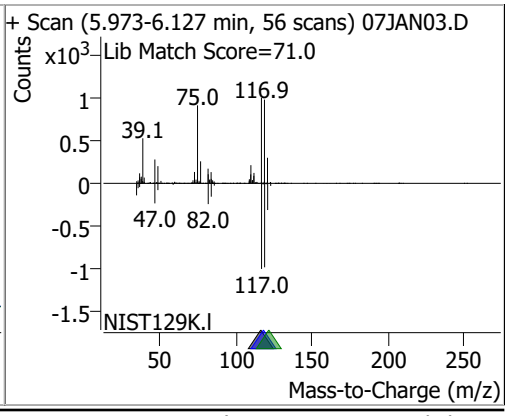




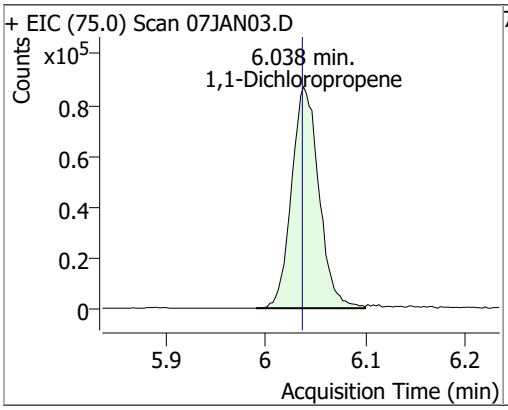
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	120.2774	6.03	0.00	183596	119.0	98.1	67.2	127.2
					121.0	30.6	0.1	60.1

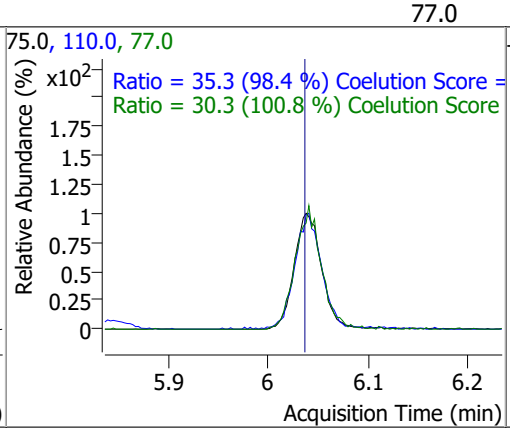


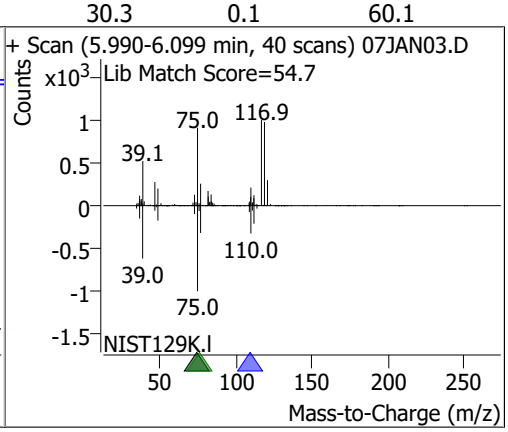




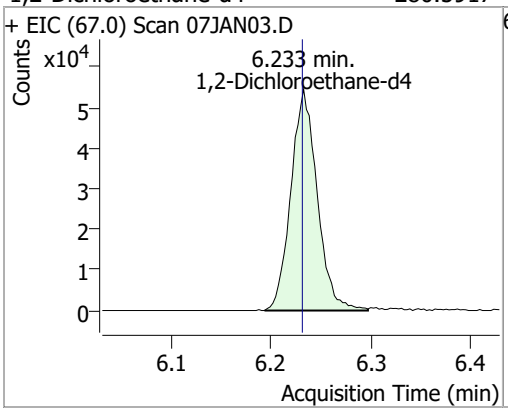
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	125.6069	6.04	0.00	165459	110.0	35.3	5.9	65.9
					77.0	30.3	0.1	60.1

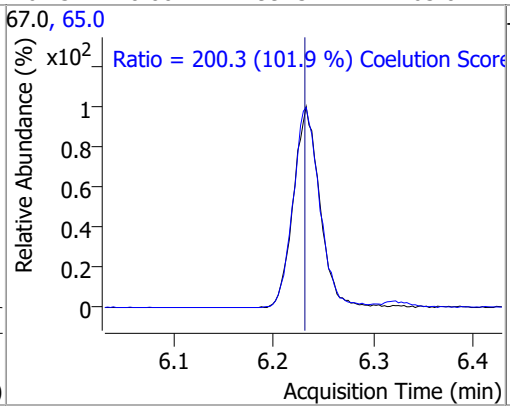


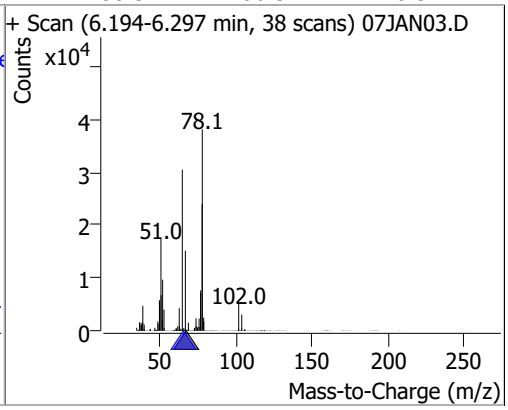




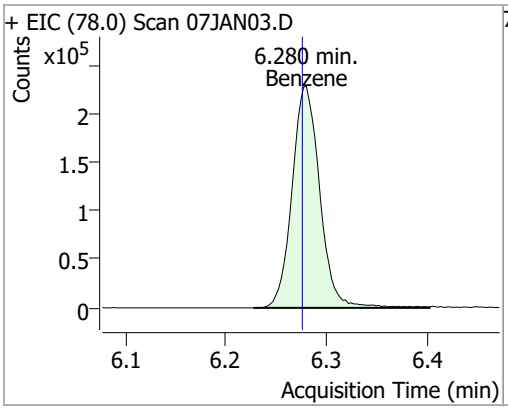
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	280.5917	6.23	0.00	99152	65.0	200.3	166.5	226.5

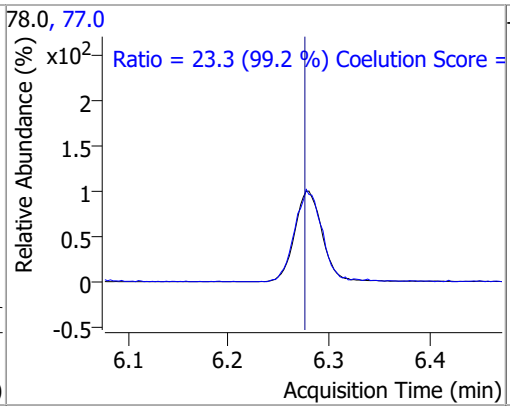


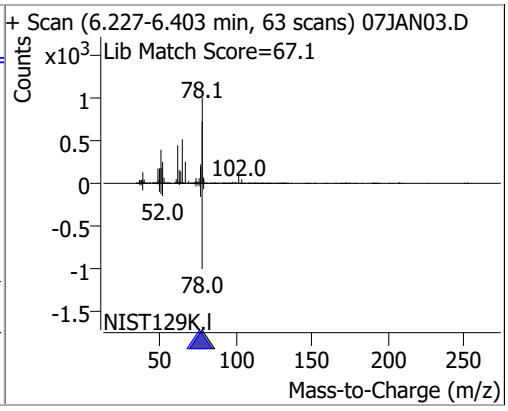




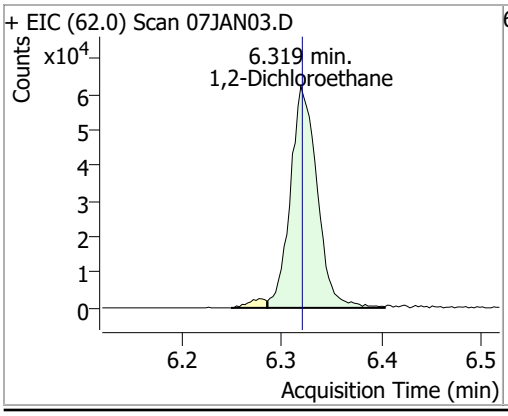
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	127.0728	6.28	0.00	439363	77.0	23.3	0.0	53.5

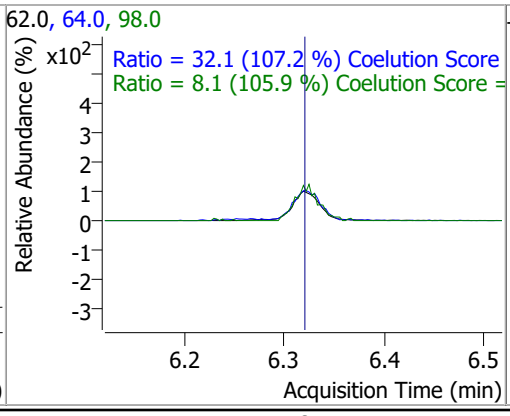


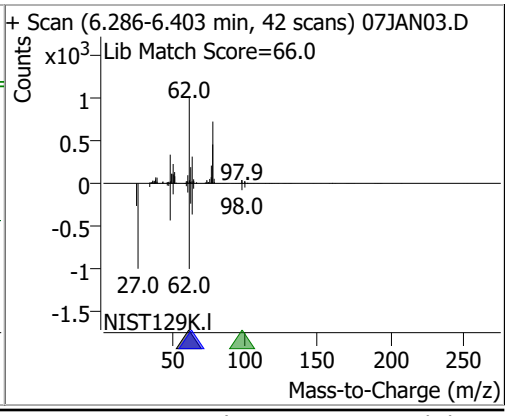




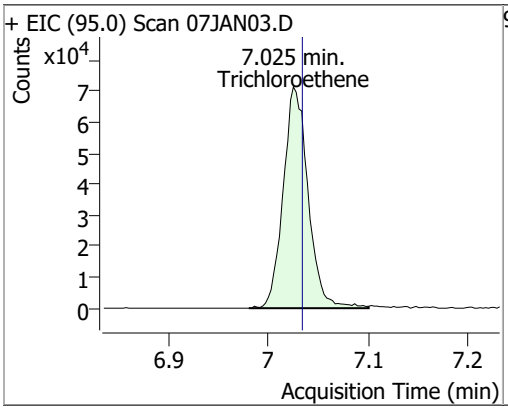
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	123.7158	6.32	0.00	115719	64.0	32.1	0.0	59.9
					98.0	8.1	0.0	37.6

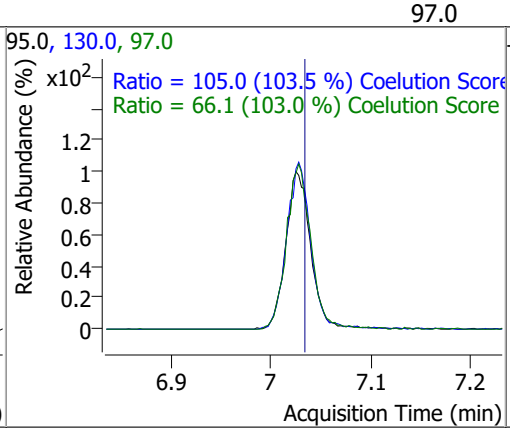


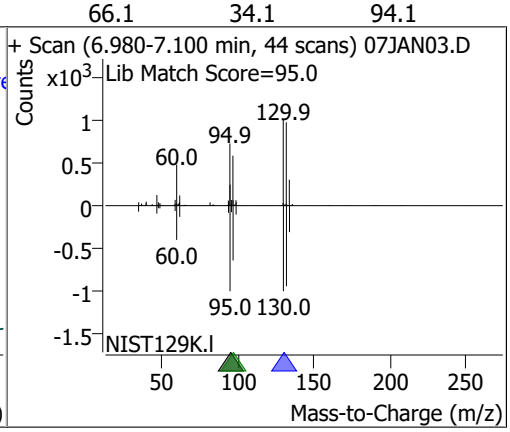




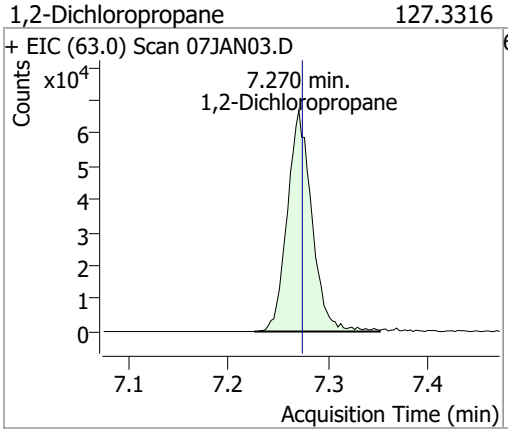
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	120.7910	7.02	-0.01	123412	130.0	105.0	71.5	131.5
					97.0	66.1	34.1	94.1

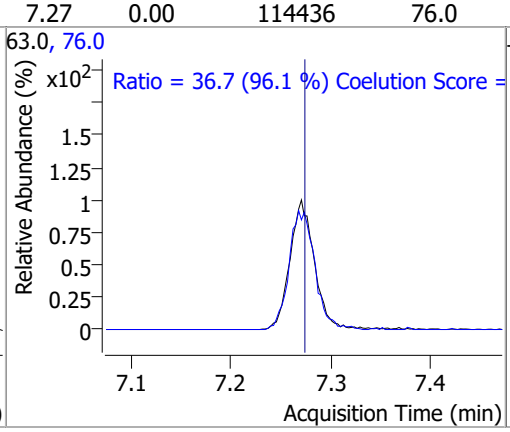


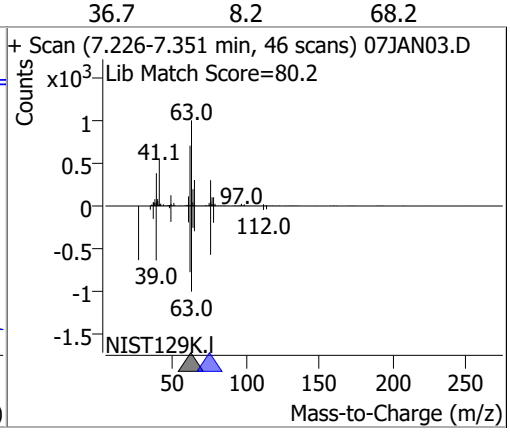




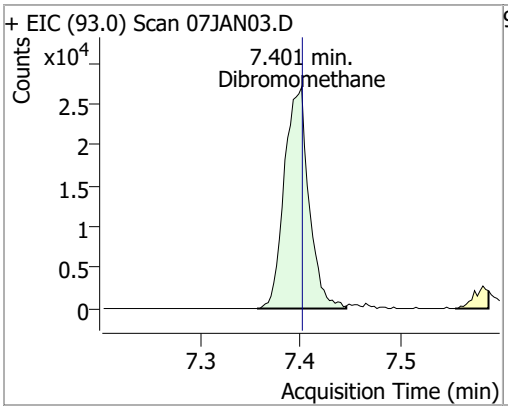
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	127.3316	7.27	0.00	114436	76.0	36.7	8.2	68.2

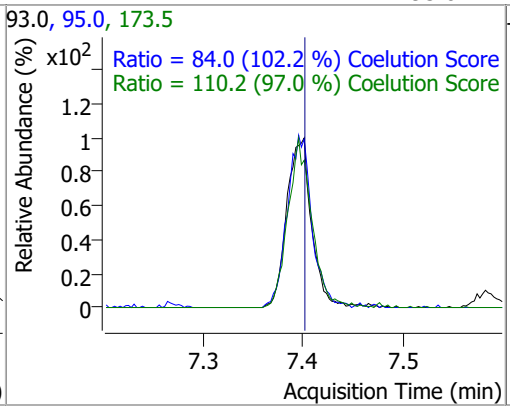


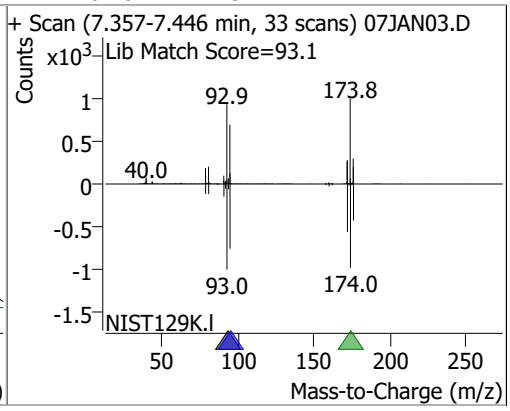




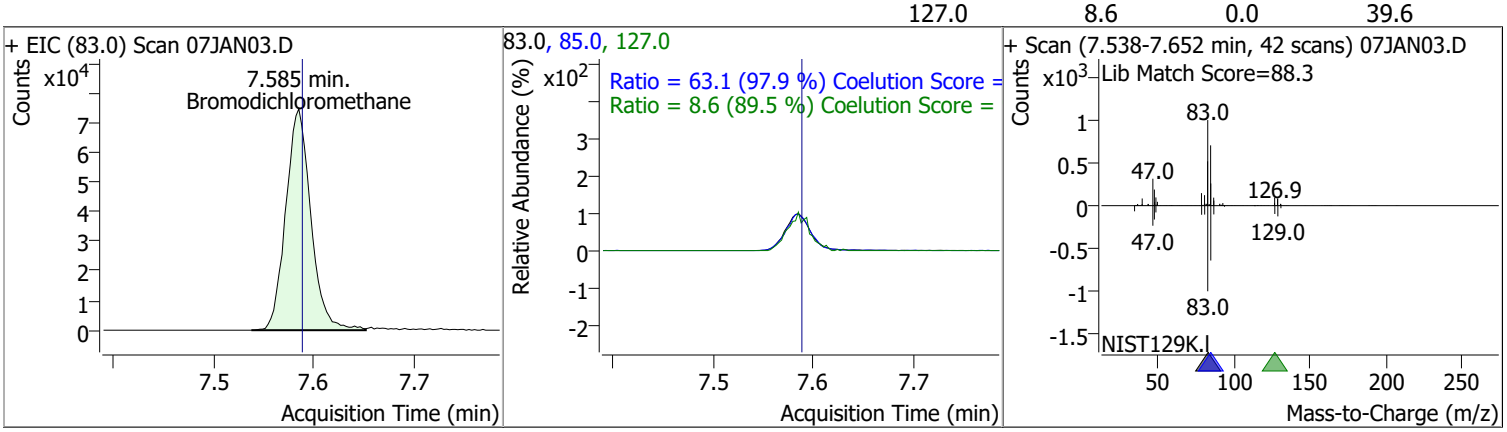
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	121.6934	7.40	0.00	46218	173.5	110.2	83.7	143.7
					95.0	84.0	52.2	112.2



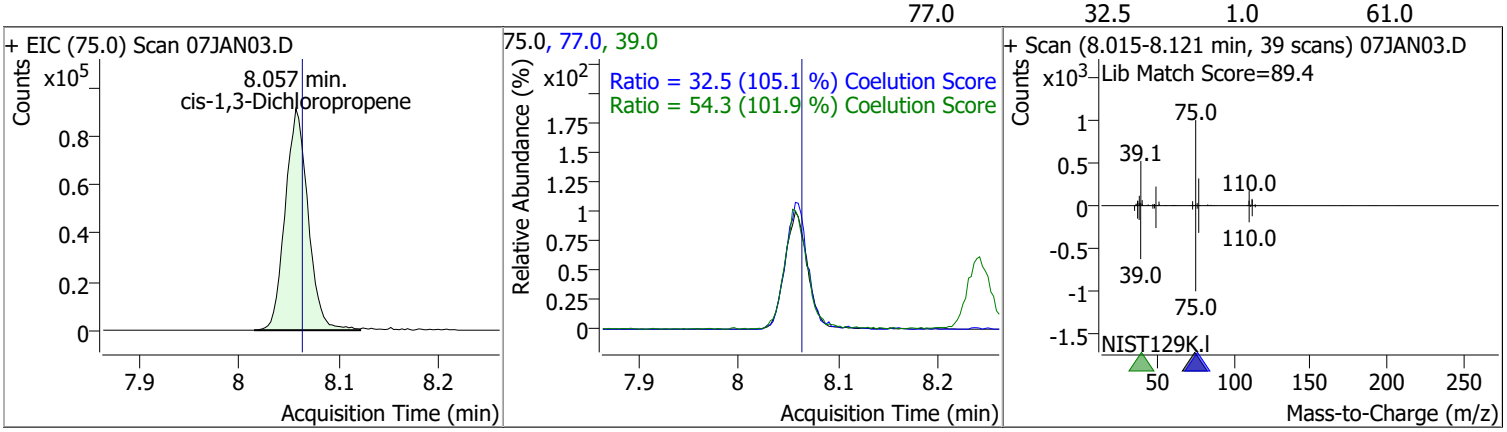




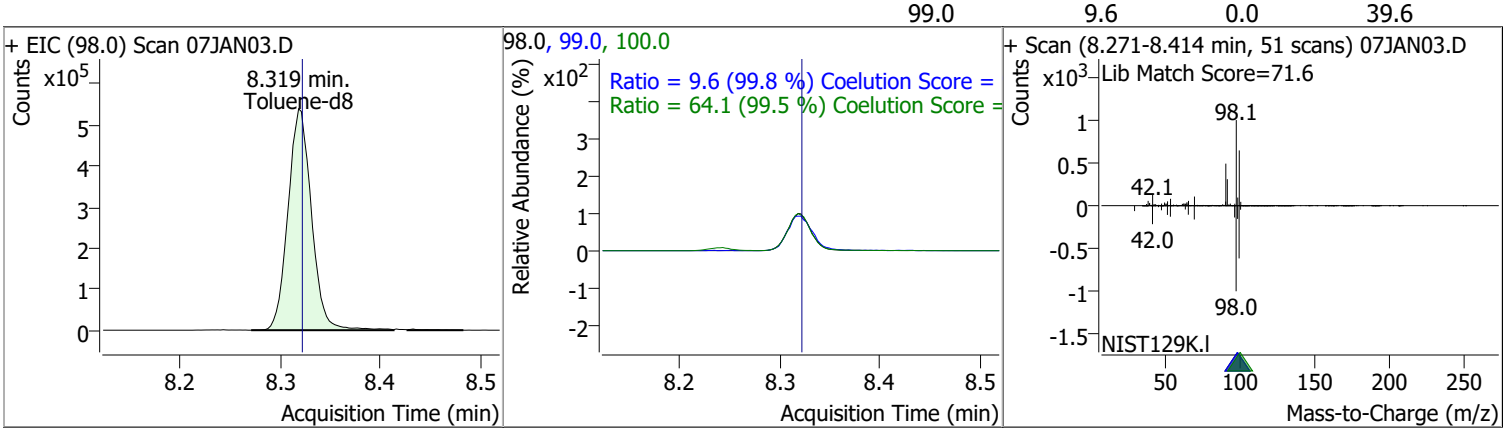
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	123.9145	7.59	0.00	129880	85.0	63.1	34.5	94.5
					127.0	8.6	0.0	39.6



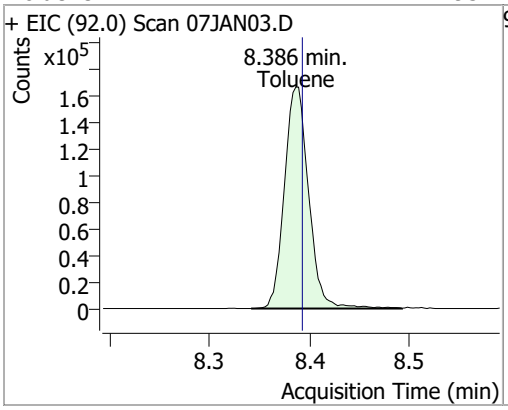
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	122.2214	8.06	0.00	144840	39.0	54.3	23.3	83.3
					77.0	32.5	1.0	61.0

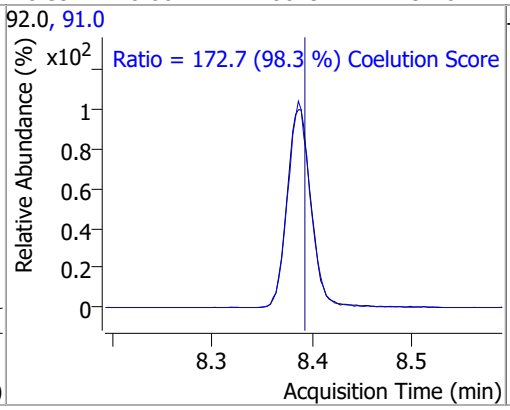


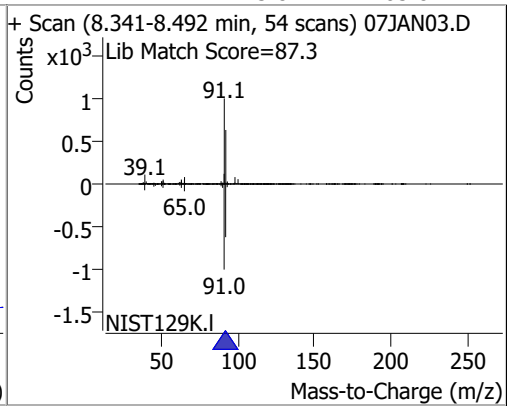
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	269.5449	8.32	0.00	879956	100.0	64.1	34.4	94.4
					99.0	9.6	0.0	39.6



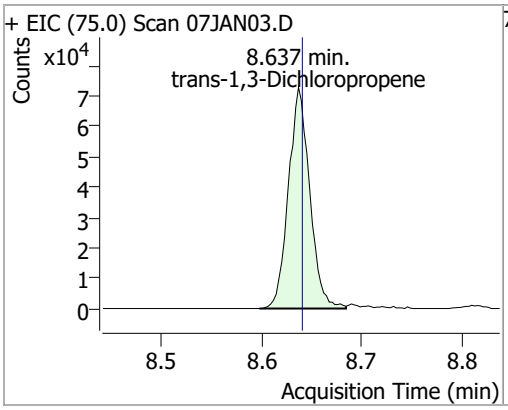
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	127.1937	8.39	0.00	280492	91.0	172.7	145.8	205.8

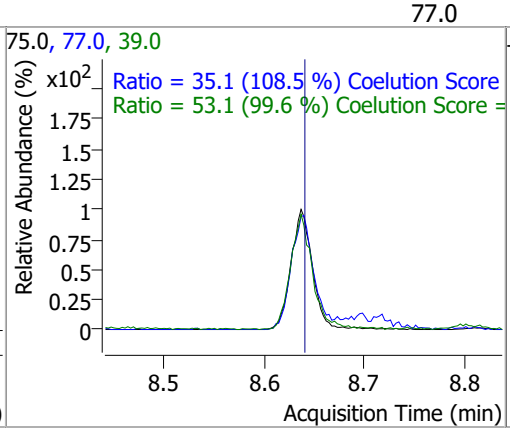


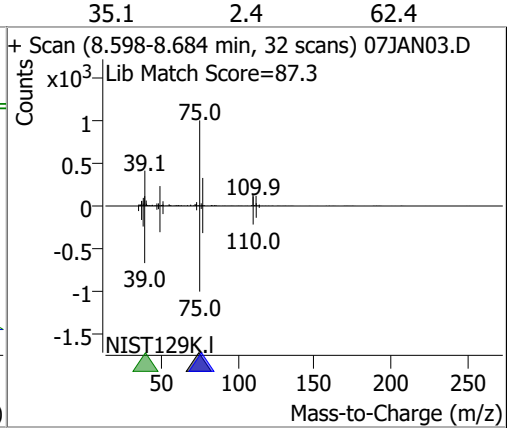




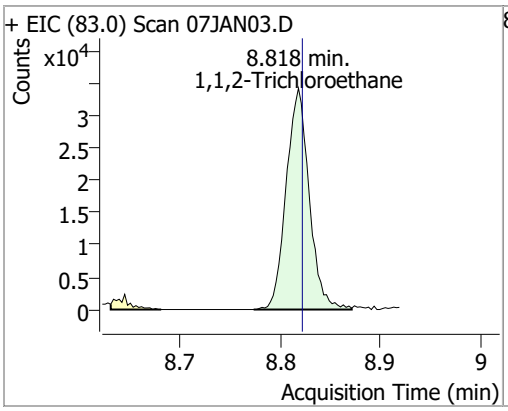
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	127.5302	8.64	0.00	107578	39.0	53.1	23.4	83.4
					77.0	35.1	2.4	62.4

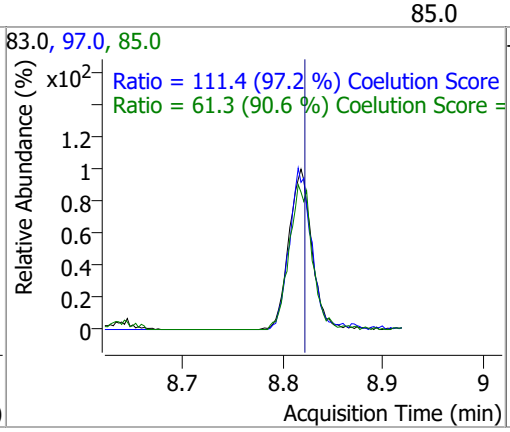


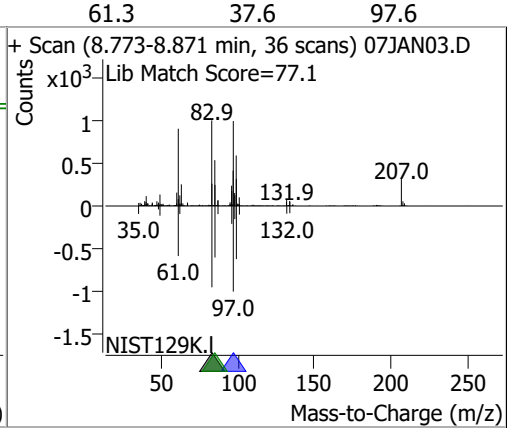




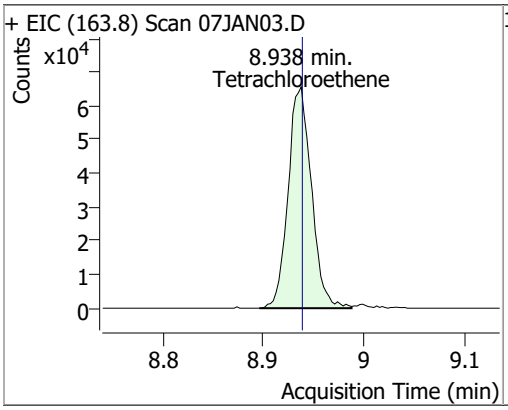
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	123.5213	8.82	0.00	54273	97.0	111.4	84.6	144.6
					85.0	61.3	37.6	97.6

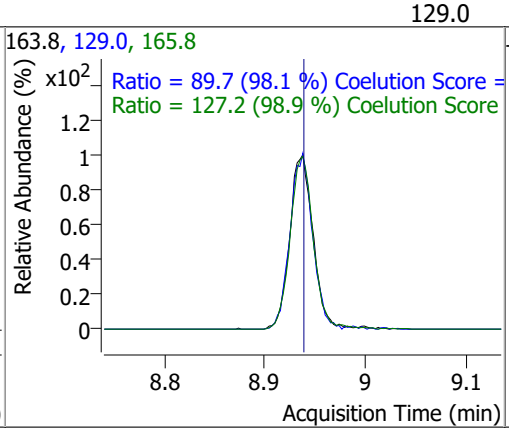


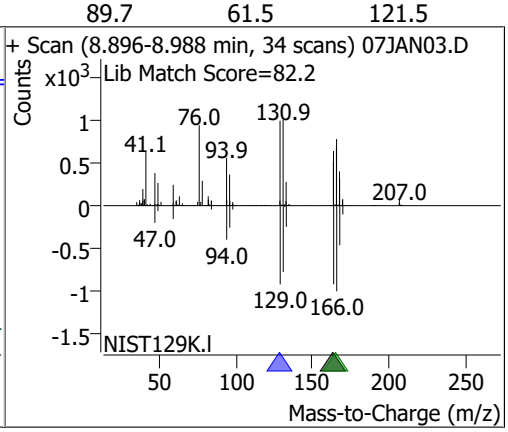




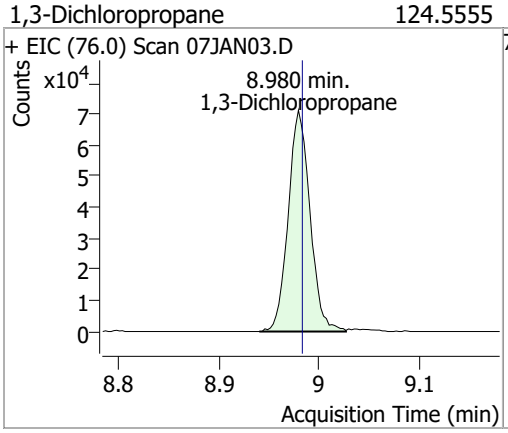
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	117.8337	8.94	0.00	106010	165.8	127.2	98.6	158.6
					129.0	89.7	61.5	121.5

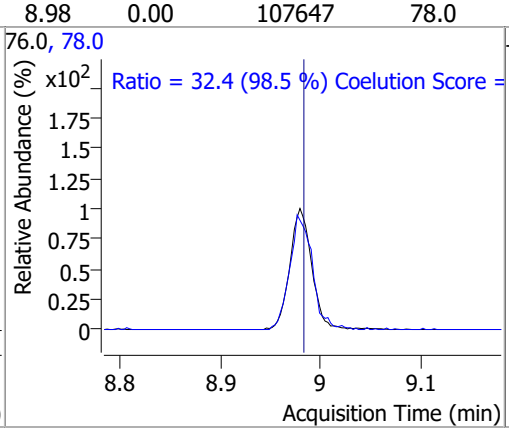


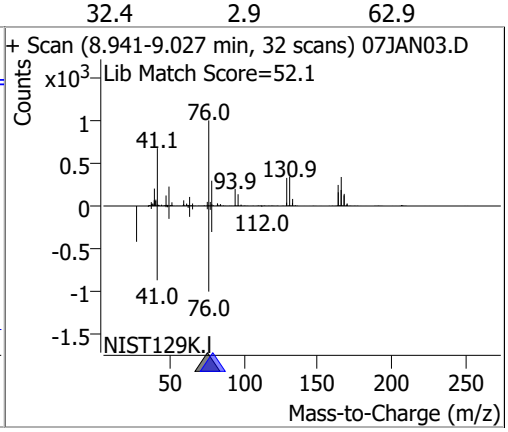




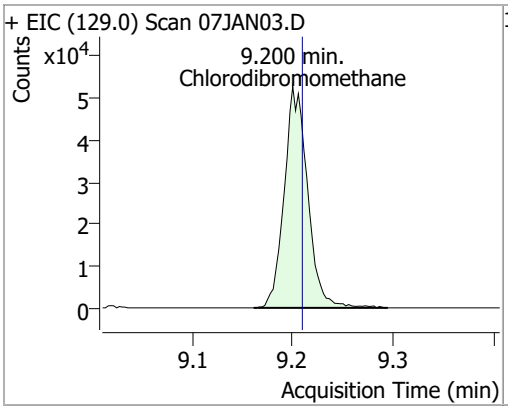
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	124.5555	8.98	0.00	107647	78.0	32.4	2.9	62.9

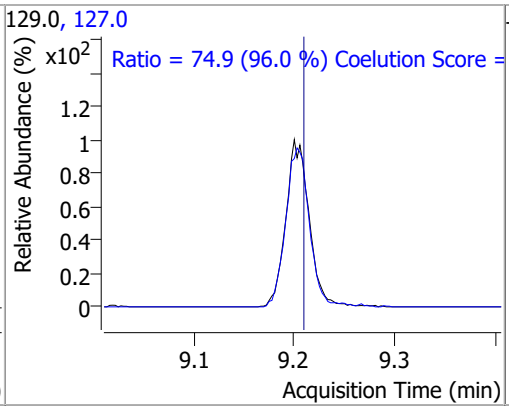


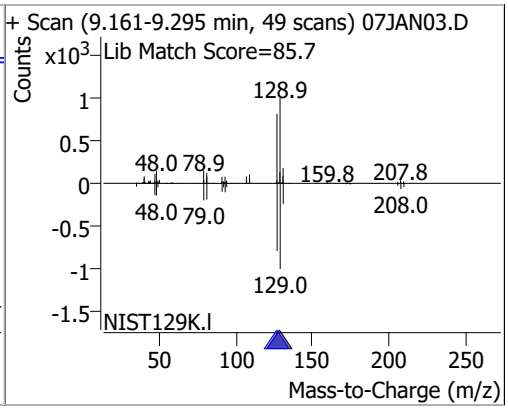




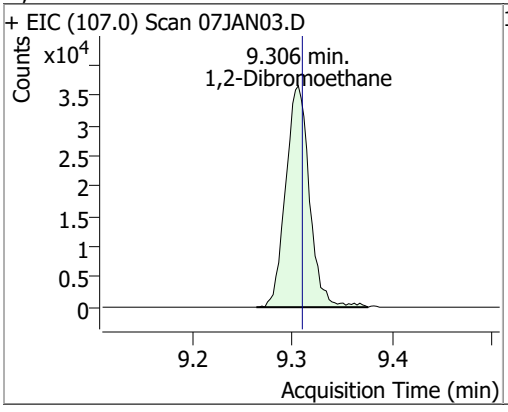
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	124.9678	9.20	-0.01	85816	127.0	74.9	48.0	108.0

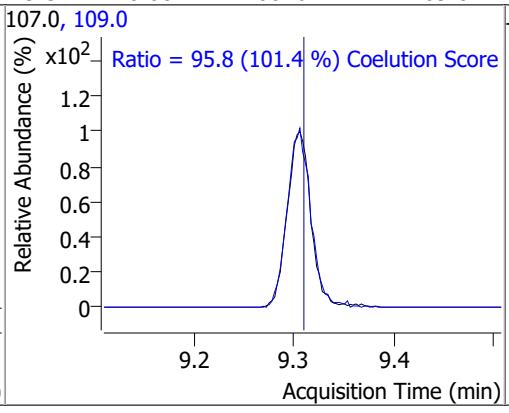


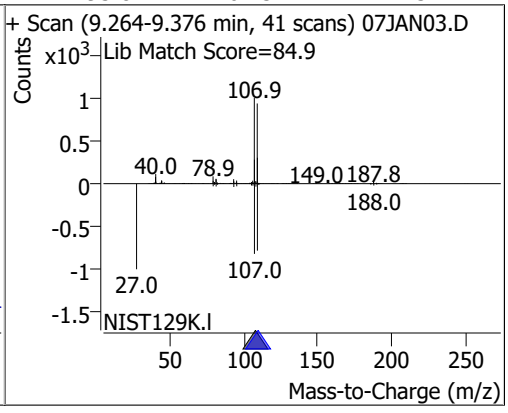




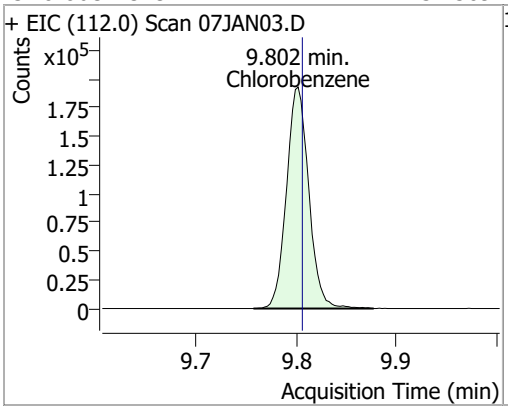
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	125.2299	9.31	0.00	60164	109.0	95.8	64.5	124.5

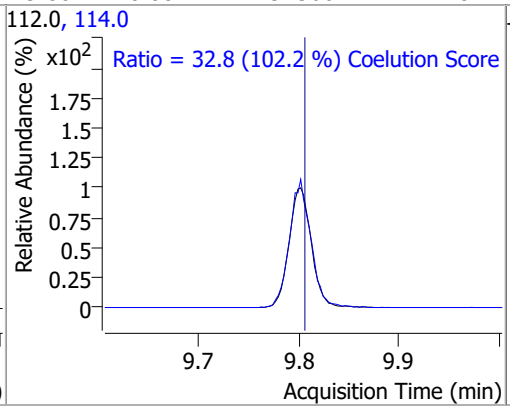


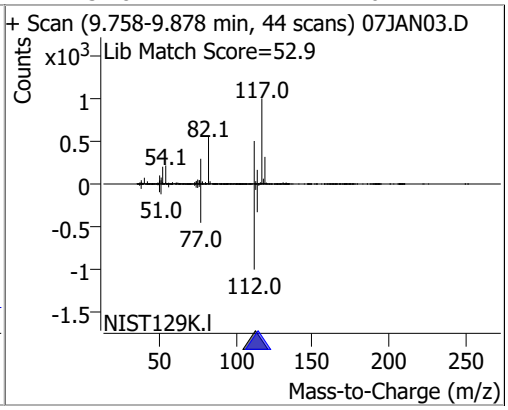




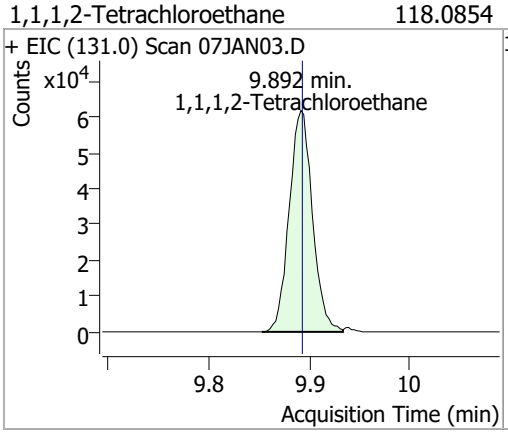
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	123.2509	9.80	0.00	297566	114.0	32.8	2.1	62.1

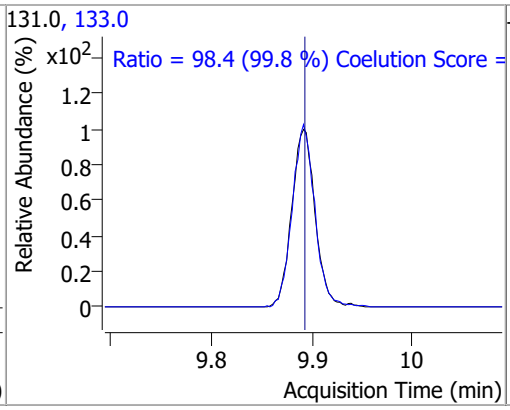


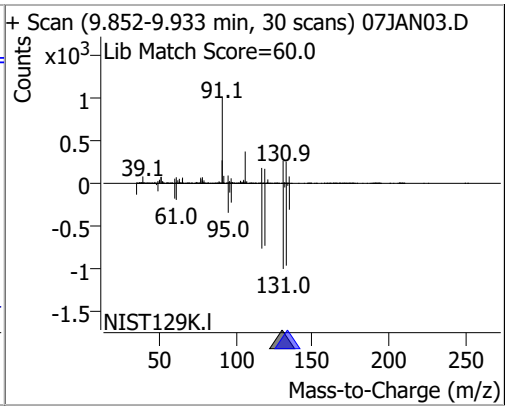




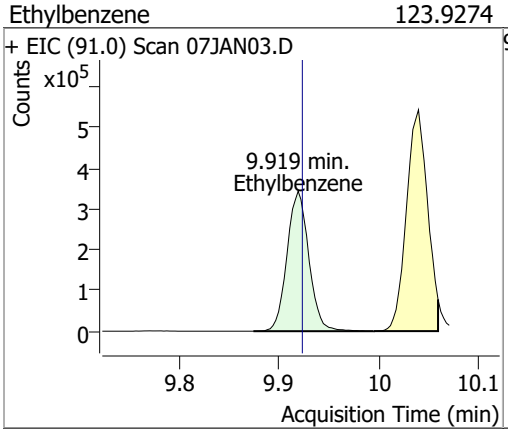
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	118.0854	9.89	0.00	99659	133.0	98.4	68.6	128.6

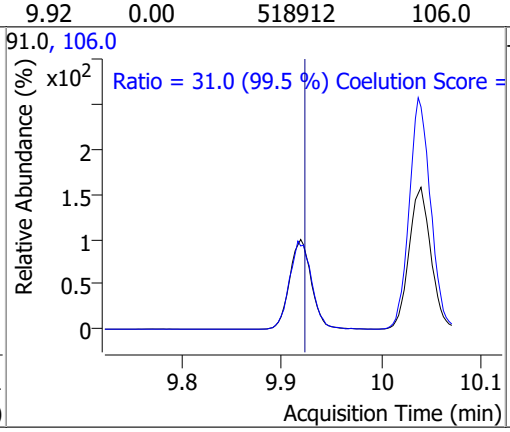


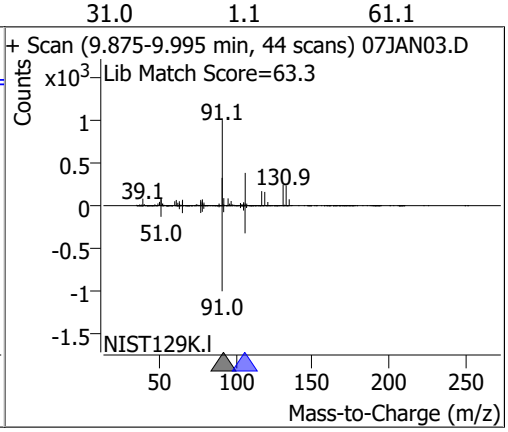




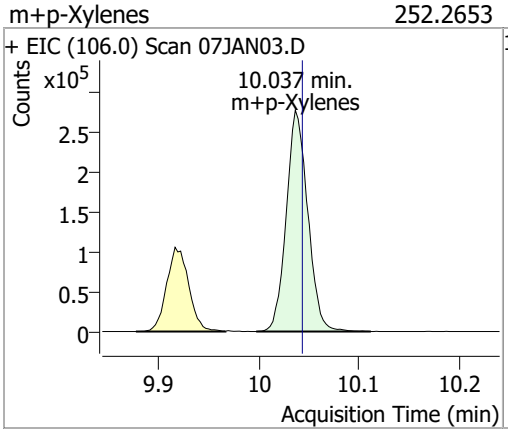
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	123.9274	9.92	0.00	518912	106.0	31.0	1.1	61.1

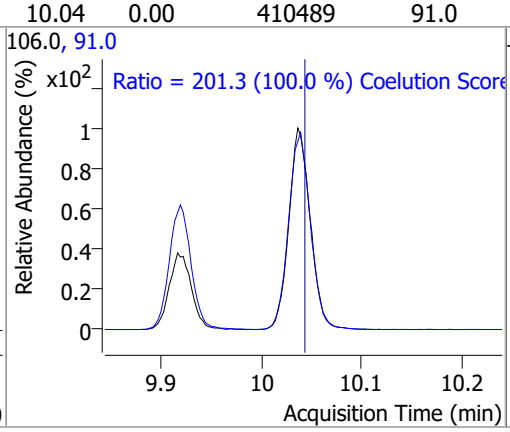


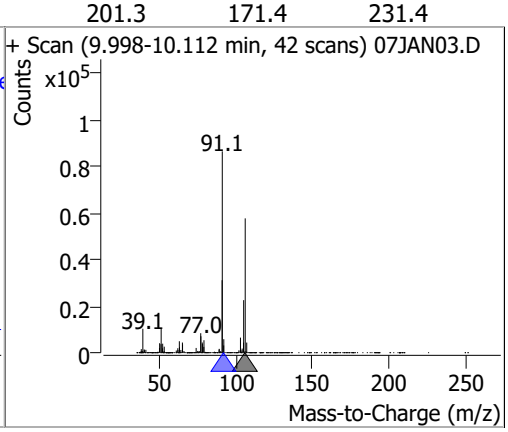




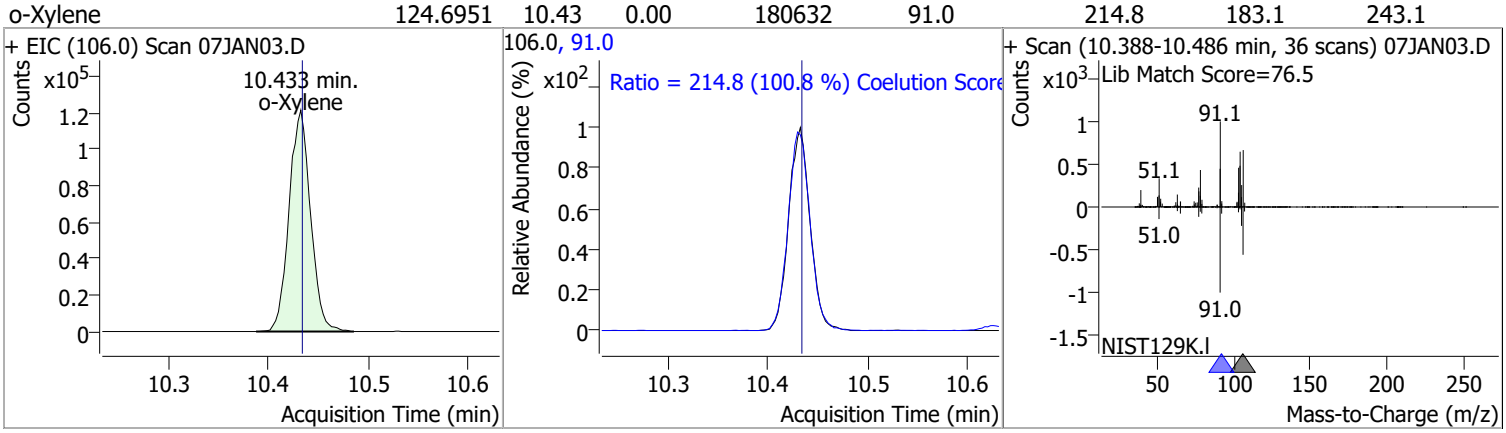
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	252.2653	10.04	0.00	410489	91.0	201.3	171.4	231.4



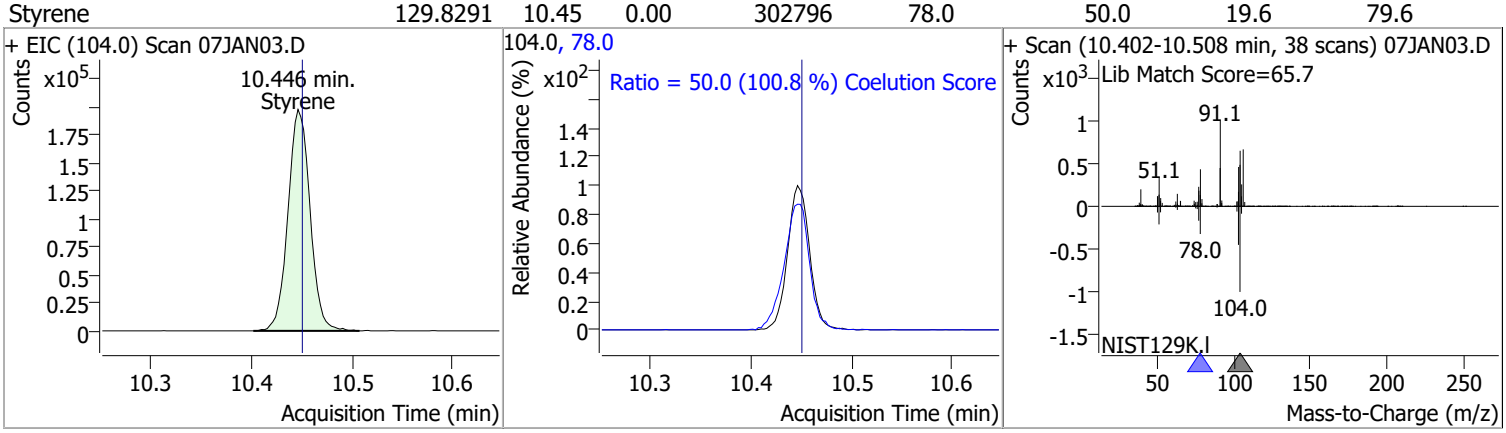




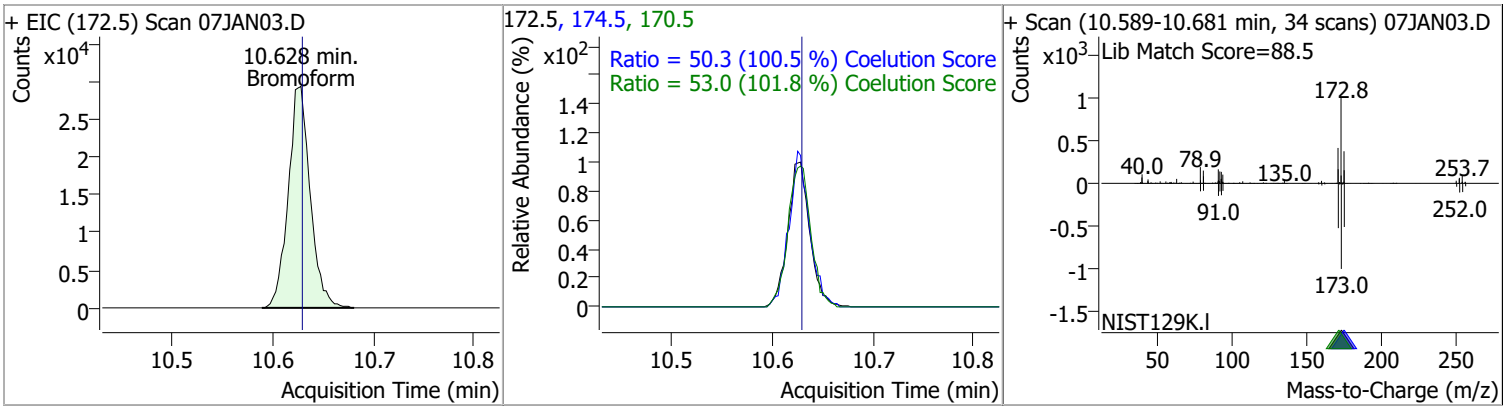
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	124.6951	10.43	0.00	180632	91.0	214.8	183.1	243.1



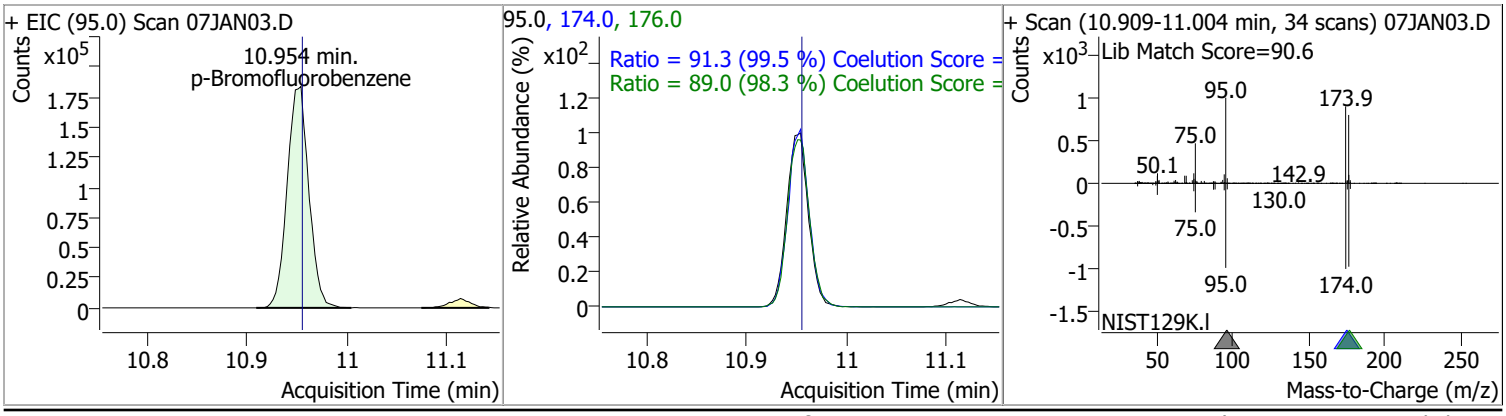
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	129.8291	10.45	0.00	302796	78.0	50.0	19.6	79.6



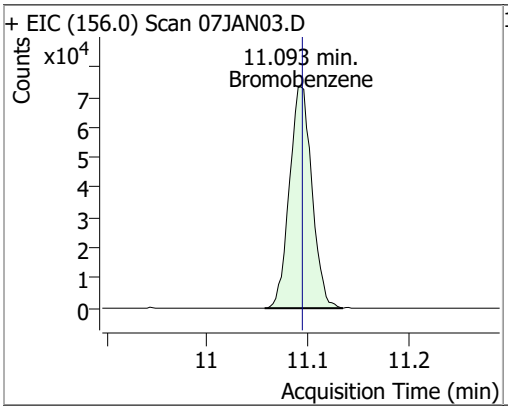
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	128.7230	10.63	0.00	44778	170.5	53.0	22.1	82.1
					174.5	50.3	20.1	80.1

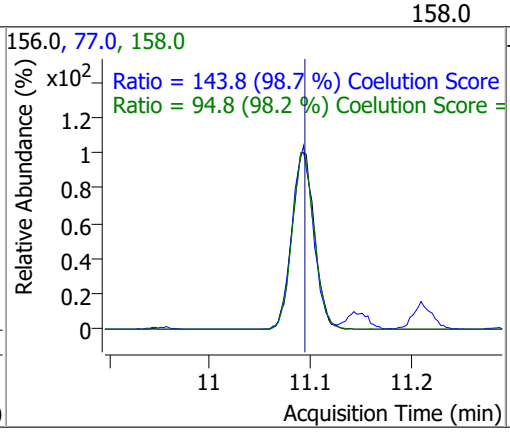


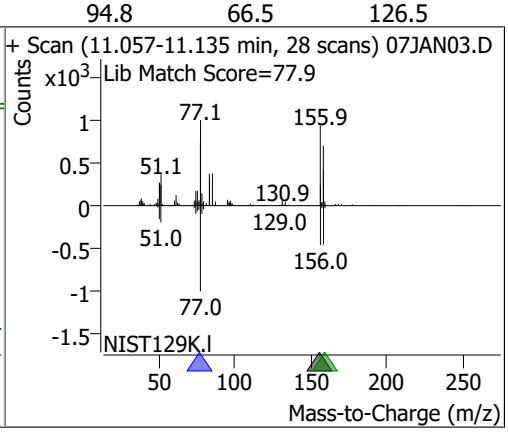
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	276.6060	10.95	0.00	275469	174.0	91.3	61.7	121.7
					176.0	89.0	60.6	120.6



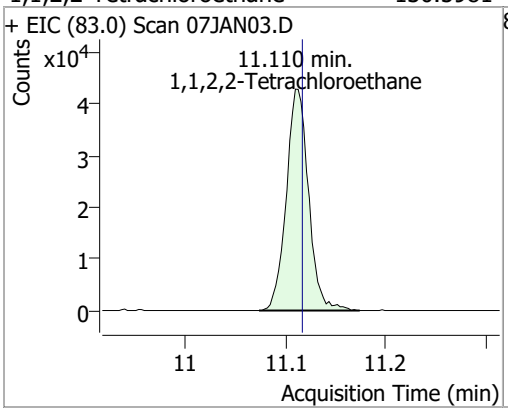
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	130.1161	11.09	0.00	114469	77.0	143.8	115.7	175.7
					158.0	94.8	66.5	126.5

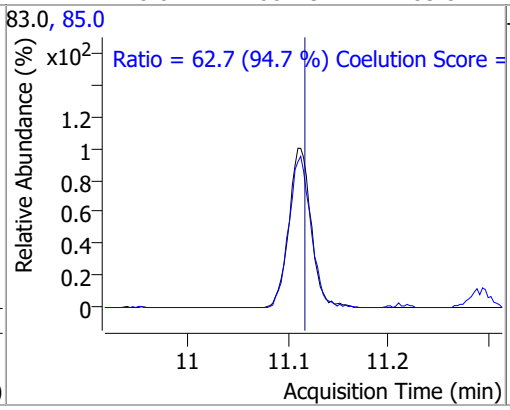


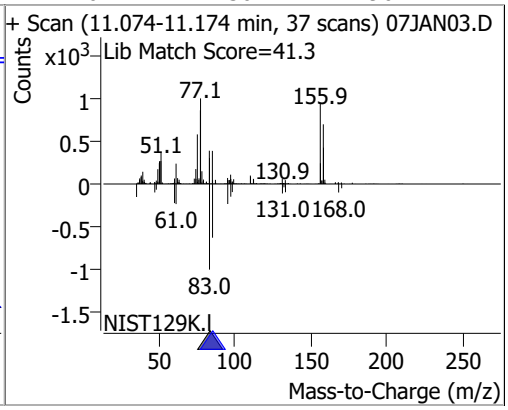




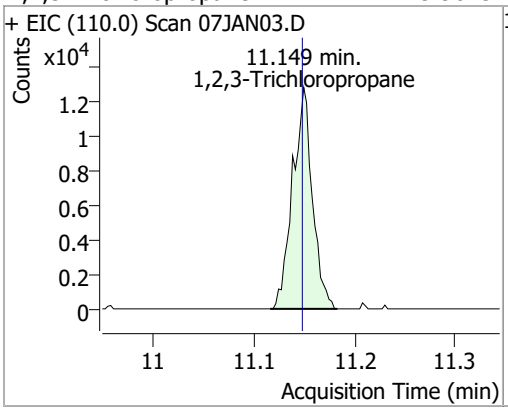
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	130.5981	11.11	-0.01	66129	85.0	62.7	36.2	96.2

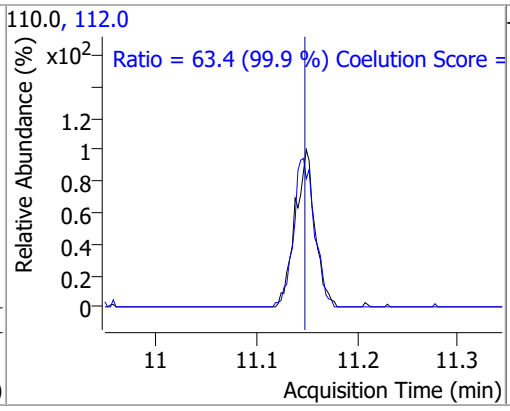


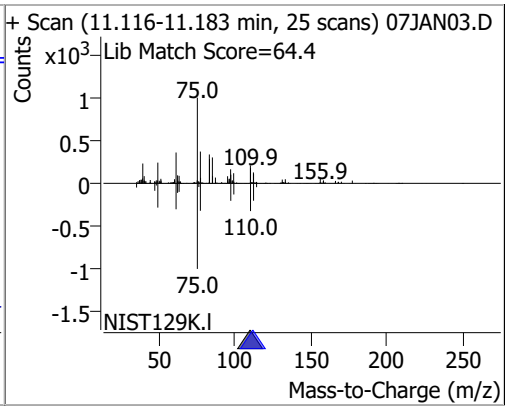




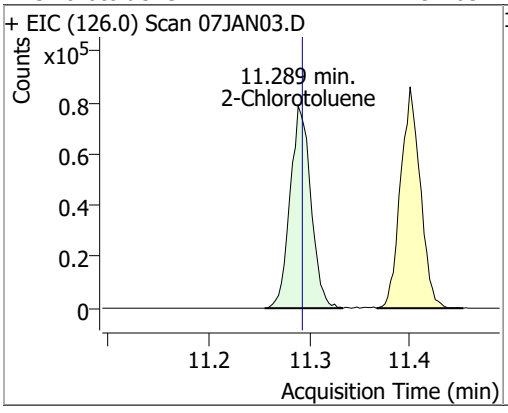
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	129.9023	11.15	0.00	17600	112.0	63.4	33.5	93.5

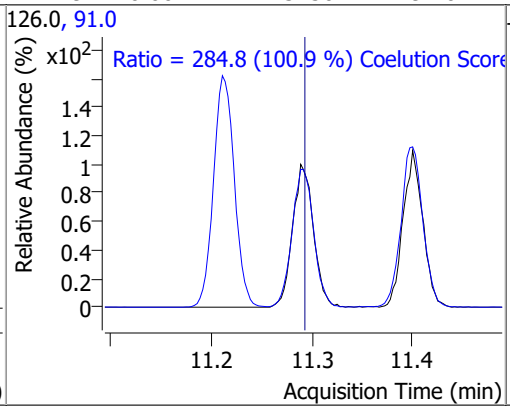


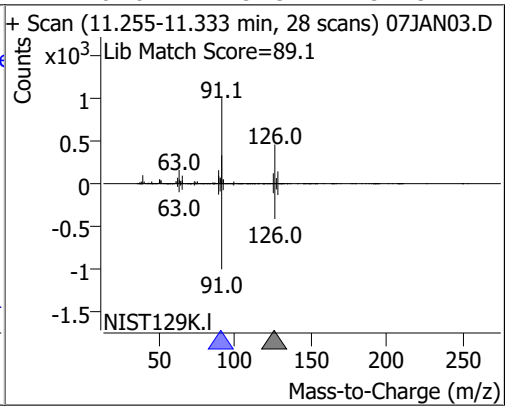




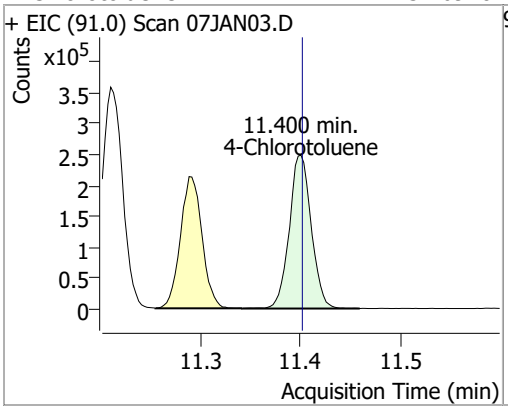
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	129.2634	11.29	0.00	113150	91.0	284.8	252.3	312.3

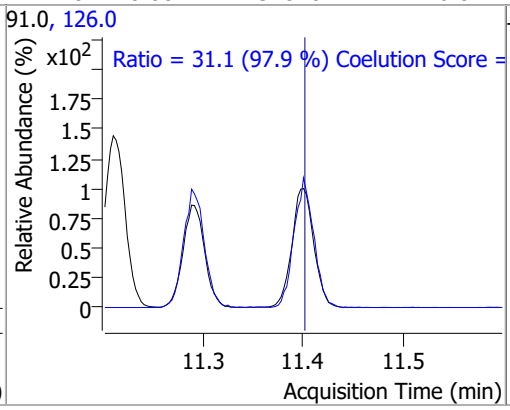


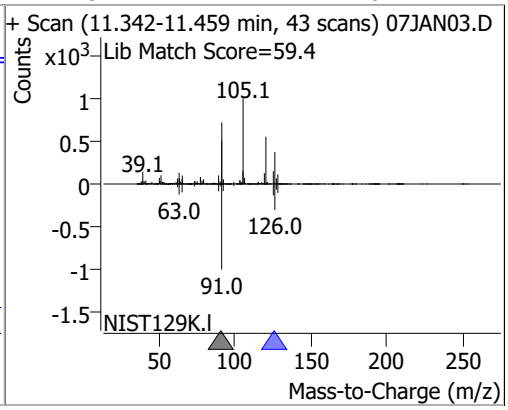




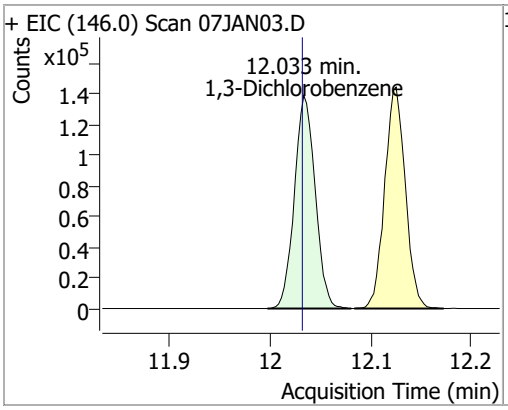
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	132.6910	11.40	0.00	378702	126.0	31.1	1.7	61.7

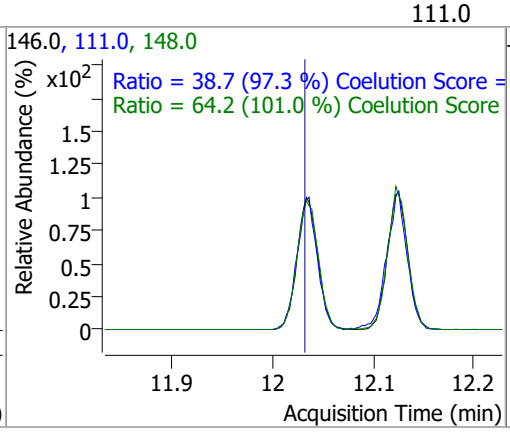


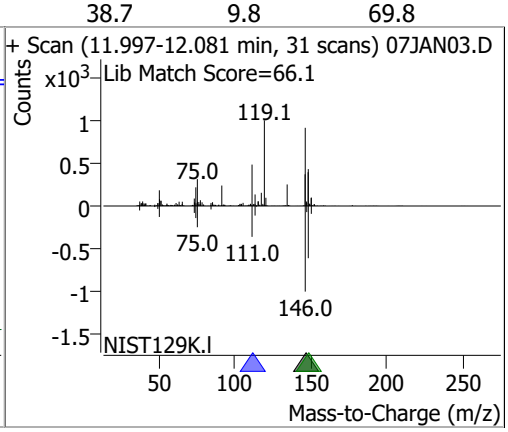




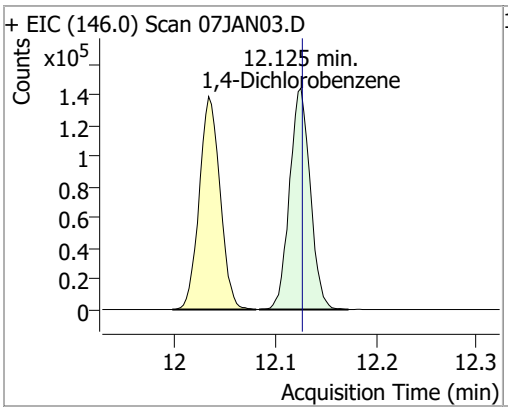
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	124.8574	12.03	0.00	200331	148.0	64.2	33.6	93.6
					111.0	38.7	9.8	69.8

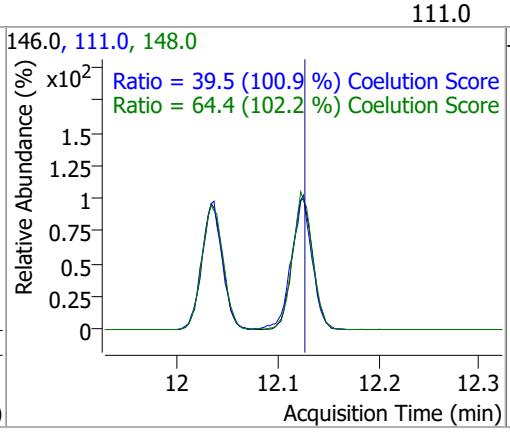


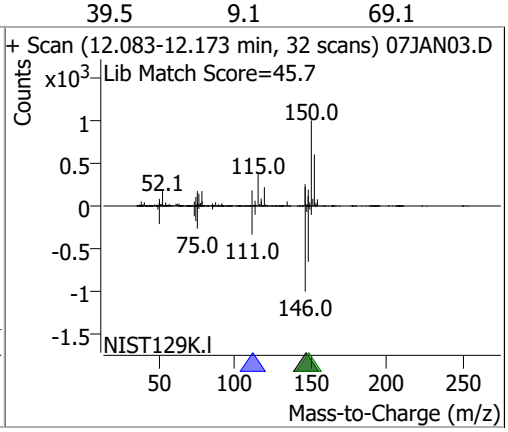




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	126.0981	12.13	0.00	206297	148.0	64.4	33.1	93.1
					111.0	39.5	9.1	69.1

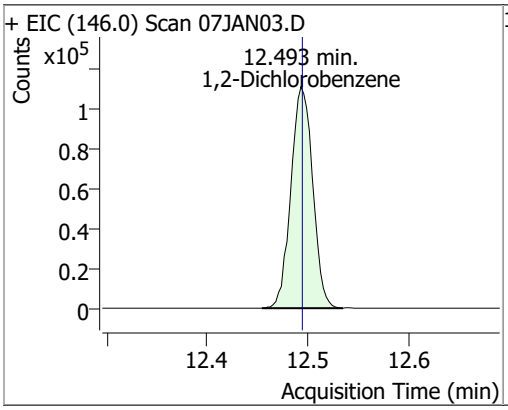


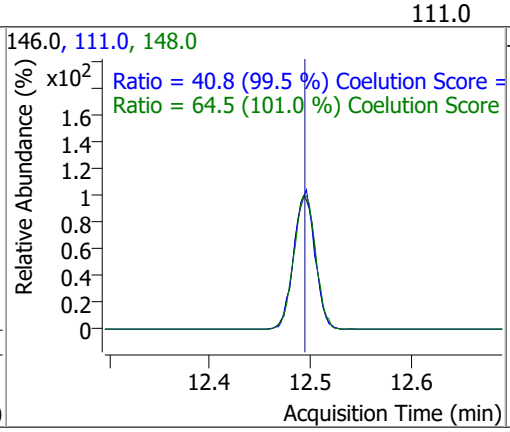


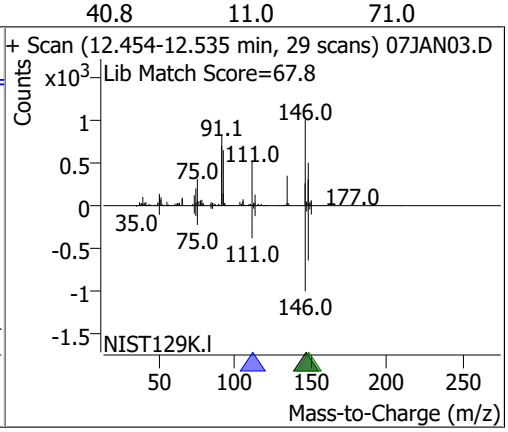


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	124.0802	12.49	0.00	168250	148.0	64.5	33.9	93.9
					111.0	40.8	11.0	71.0







Quantitation Results Report (QT Reviewed)

Data File	07JAN04.D	Operator	MSC
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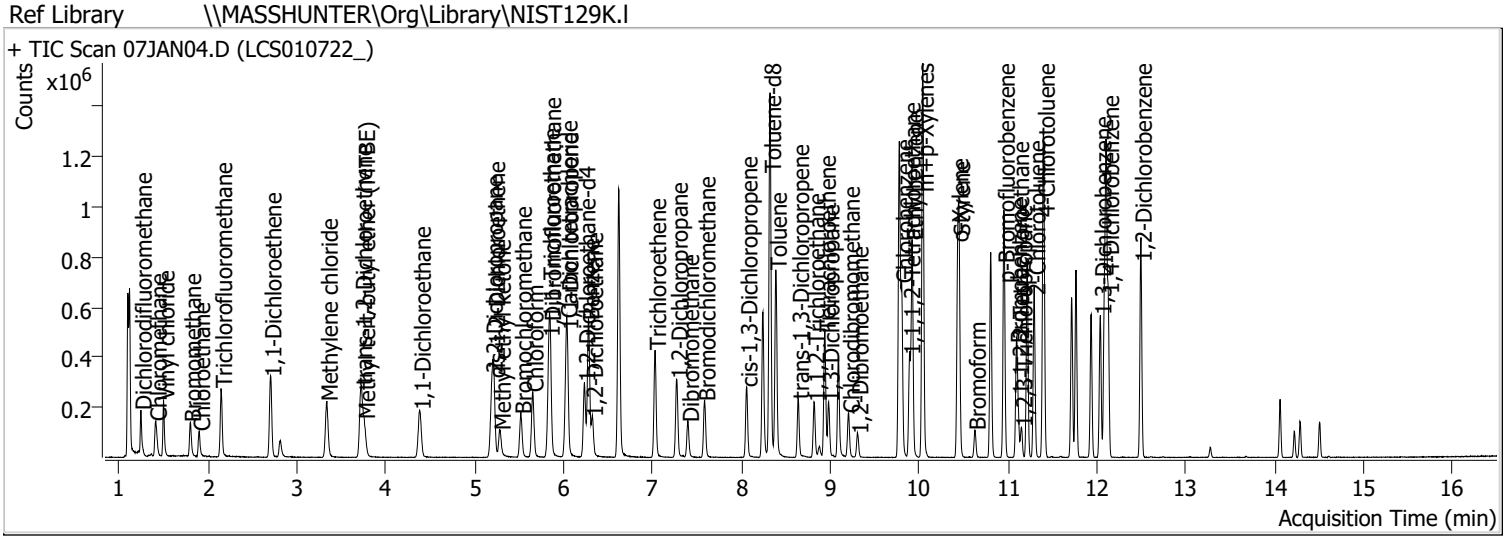
Vial	4	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

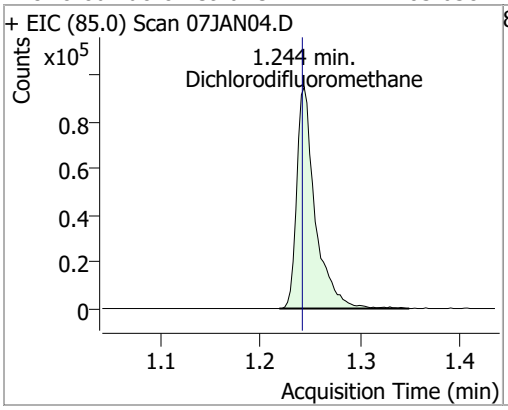


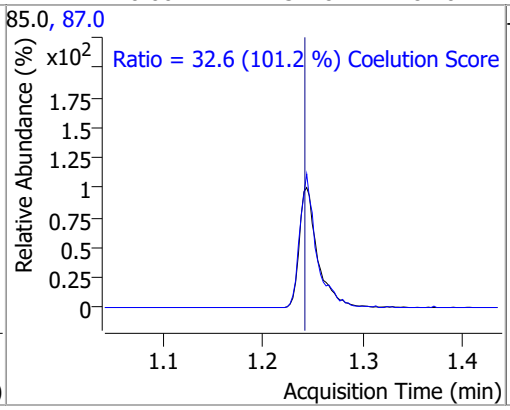
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.620	96.0	907718	250.0000	ng	-0.003
M Chlorobenzene-d5	9.774	82.0	346586	250.0000	ng	0.003
M 1,4-Dichlorobenzene-d4	12.103	152.0	280949	250.0000	ng	0.003
System Monitoring Compounds						
S Dibromofluoromethane	5.848	113.0	231280	270.4515	ng	0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 108.18%		
S 1,2-Dichloroethane-d4	6.233	67.0	104500	282.9153	ng	0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 113.17%		
S Toluene-d8	8.321	98.0	905281	271.0520	ng	0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 108.42%		
S p-Bromofluorobenzene	10.951	95.0	279940	271.9817	ng	-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 108.79%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.244	85.0	123278	103.6382	ng	99
T Chloromethane	1.411	50.0	156934	108.6978	ng	99
T Vinyl chloride	1.498	62.0	154899	119.2350	ng	97
T Bromomethane	1.799	96.0	70232	120.9026	ng	99
T Chloroethane	1.899	64.0	70493	109.6085	ng	97
T Trichlorofluoromethane	2.145	101.0	181809	112.7515	ng	99
T 1,1-Dichloroethene	2.700	96.0	122201	133.6517	ng	96
T Methylene chloride	3.335	49.0	159195	118.1093	ng	99
T trans-1,2-Dichloroethene	3.720	96.0	121109	129.8320	ng	97
T Methyl tert-butyl ether (MTBE)	3.756	73.0	160126	132.8050	ng	99
T 1,1-Dichloroethane	4.381	63.0	227596	131.0789	ng	100
T 2,2-Dichloropropane	5.193	77.0	167542	128.7744	ng	100
T cis-1,2-Dichloroethene	5.218	96.0	122647	129.6836	ng	98
T Methyl ethyl ketone	5.282	43.0	145822	1138.3128	ng	98
T Bromochloromethane	5.519	128.0	47092	120.1958	ng	93
T Chloroform	5.650	83.0	206280	119.3745	ng	100

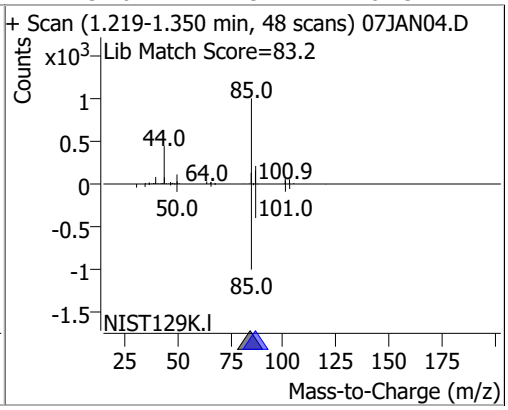
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.831	97.0	197819	122.1545	ng	99
T Carbon tetrachloride	6.032	117.0	187698	117.6380	ng	99
T 1,1-Dichloropropene	6.040	75.0	162591	118.0827	ng	99
T Benzene	6.280	78.0	461197	127.6093	ng	100
T 1,2-Dichloroethane	6.319	62.0	119261	121.9792	ng	100
T Trichloroethene	7.028	95.0	131597	125.8989	ng	99
T 1,2-Dichloropropane	7.270	63.0	114080	124.0744	ng	99
T Dibromomethane	7.398	93.0	48238	124.1493	ng	97
T Bromodichloromethane	7.582	83.0	136799	127.5739	ng	99
T cis-1,3-Dichloropropene	8.057	75.0	142161	117.2568	ng	99
T Toluene	8.386	92.0	291378	129.1519	ng	99
T trans-1,3-Dichloropropene	8.637	75.0	113275	131.2571	ng	95
T 1,1,2-Trichloroethane	8.815	83.0	56261	125.1597	ng	99
T Tetrachloroethene	8.935	163.8	109421	118.8837	ng	100
T 1,3-Dichloropropane	8.980	76.0	107326	121.3850	ng	99
T Chlorodibromomethane	9.205	129.0	86224	122.7318	ng	99
T 1,2-Dibromoethane	9.306	107.0	60530	123.1519	ng	100
T Chlorobenzene	9.802	112.0	310348	125.6477	ng	99
T 1,1,1,2-Tetrachloroethane	9.889	131.0	108057	125.1503	ng	94
T Ethylbenzene	9.919	91.0	539522	125.9452	ng	99
T m+p-Xylenes	10.039	106.0	417695	250.9079	ng	99
T o-Xylene	10.430	106.0	192510	129.8994	ng	100
T Styrene	10.449	104.0	316187	132.5150	ng	100
T Bromoform	10.625	172.5	47153	131.1556	ng	99
T Bromobenzene	11.091	156.0	119622	131.5649	ng	100
T 1,1,2,2-Tetrachloroethane	11.110	83.0	64431	123.1192	ng	98
T 1,2,3-Trichloropropane	11.149	110.0	17528	125.1764	ng	99
T 2-Chlorotoluene	11.294	126.0	117650	130.0465	ng	97
T 4-Chlorotoluene	11.400	91.0	398147	134.9812	ng	100
T 1,3-Dichlorobenzene	12.036	146.0	218804	131.9494	ng	99
T 1,4-Dichlorobenzene	12.125	146.0	217500	128.6355	ng	99
T 1,2-Dichlorobenzene	12.493	146.0	175961	125.5596	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

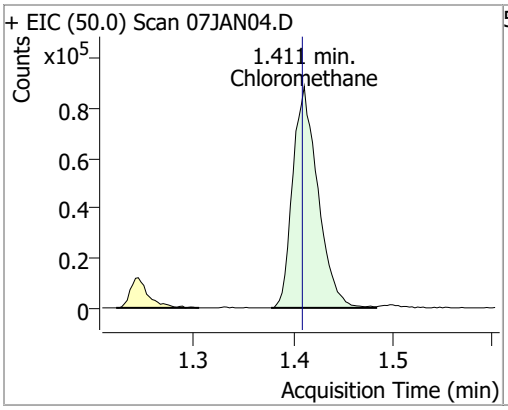
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	103.6382	1.24	0.00	123278	87.0	32.6	2.3	62.3

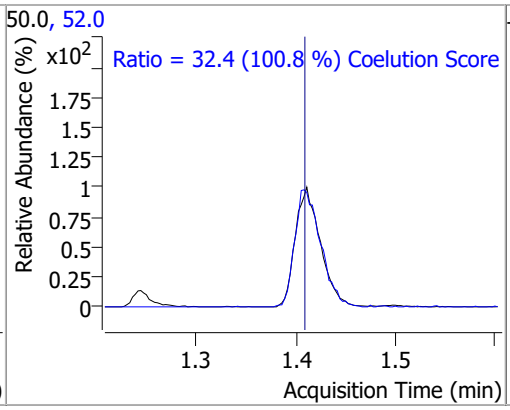


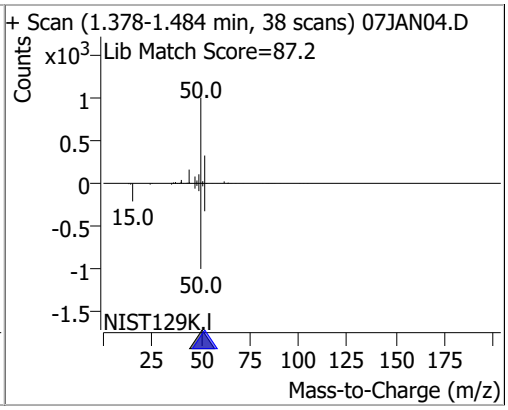




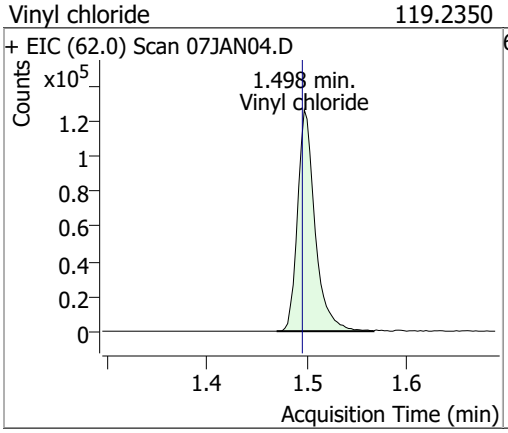
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	108.6978	1.41	0.00	156934	52.0	32.4	2.1	62.1

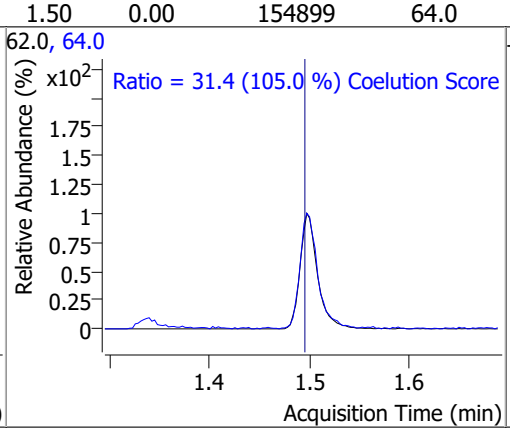


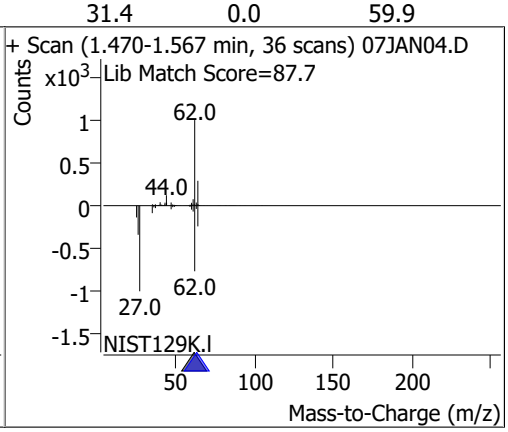




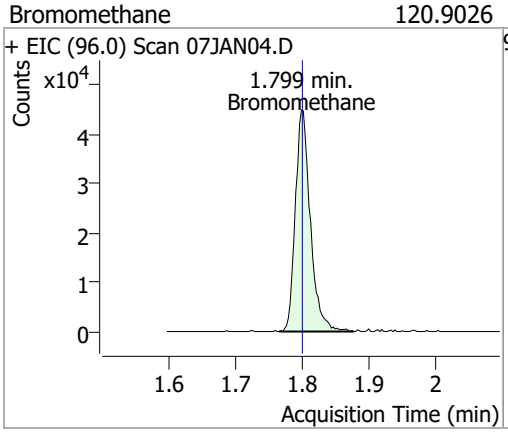
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	119.2350	1.50	0.00	154899	64.0	31.4	0.0	59.9

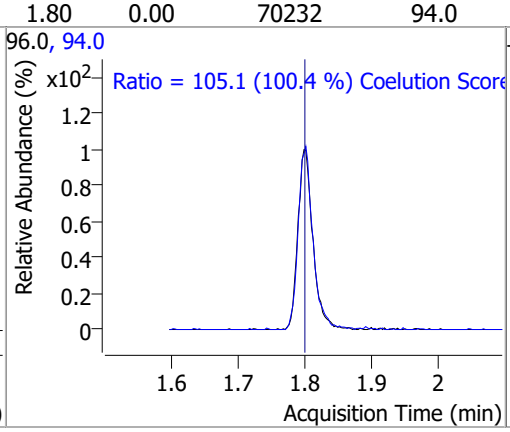


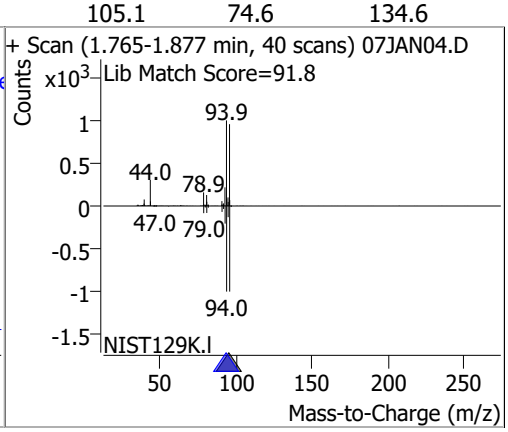




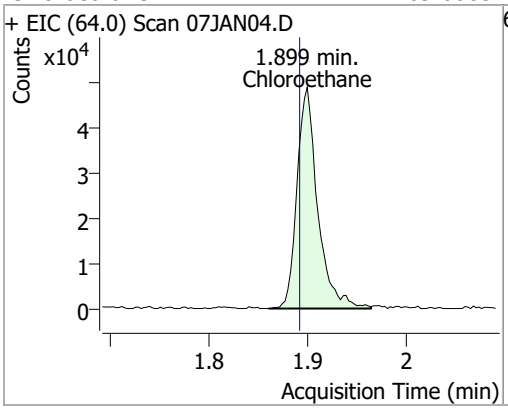
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	120.9026	1.80	0.00	70232	94.0	105.1	74.6	134.6

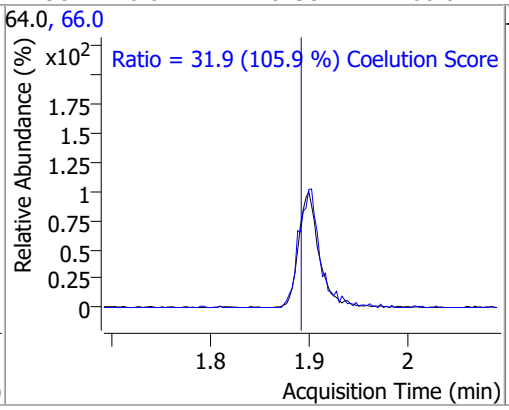


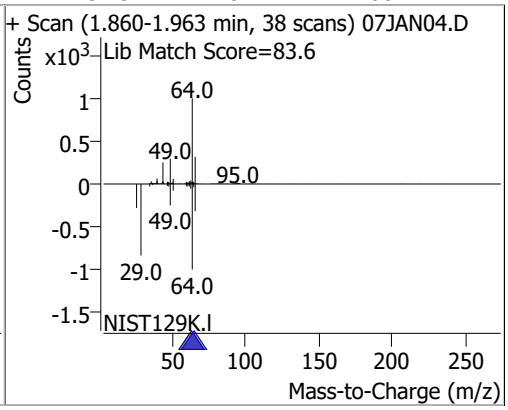




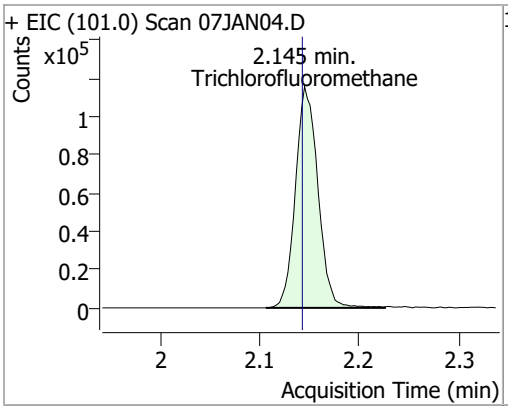
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	109.6085	1.90	0.01	70493	66.0	31.9	0.1	60.1

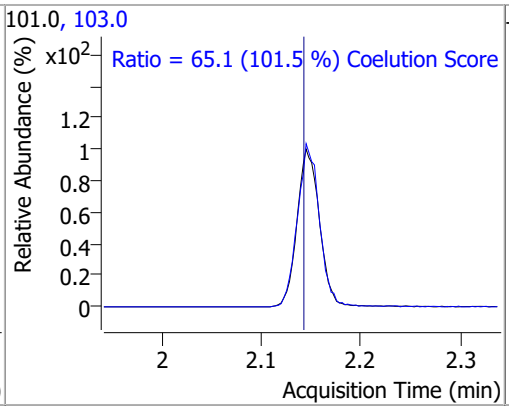


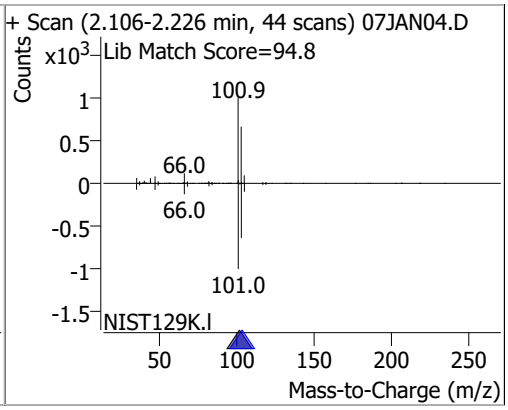




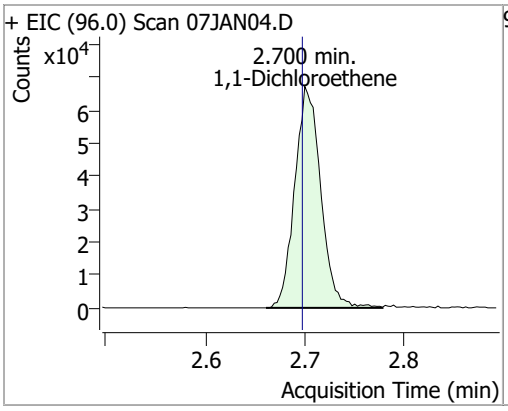
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	112.7515	2.14	0.00	181809	103.0	65.1	34.2	94.2

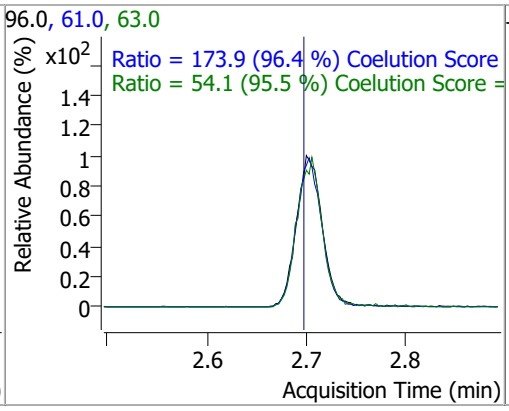


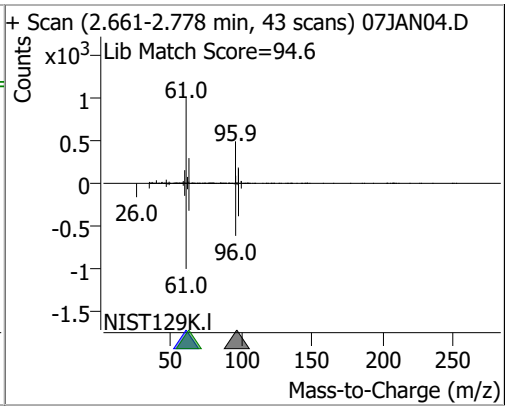




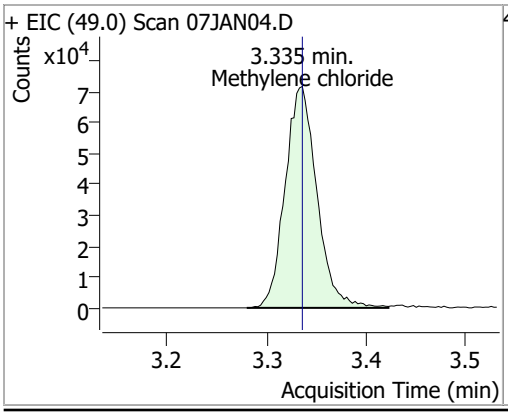
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	133.6517	2.70	0.00	122201	61.0	173.9	150.3	210.3
					63.0	54.1	26.7	86.7

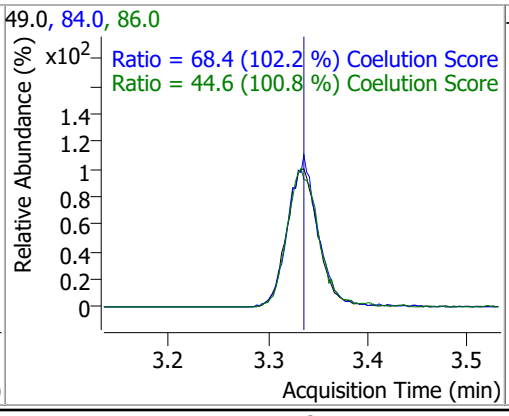


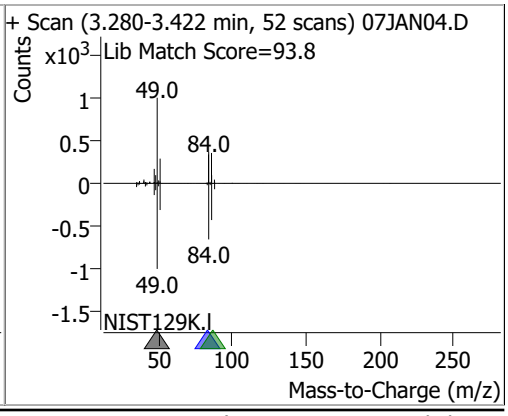




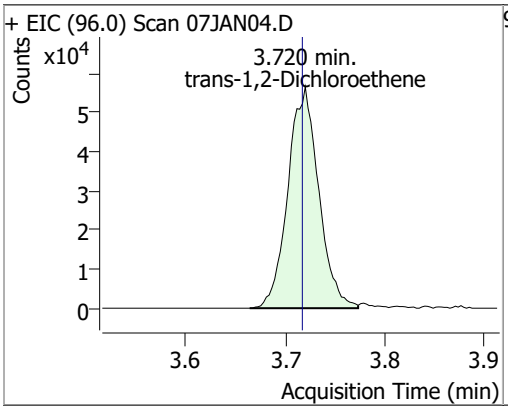
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	118.1093	3.34	0.00	159195	84.0	68.4	36.9	96.9
					86.0	44.6	14.3	74.3

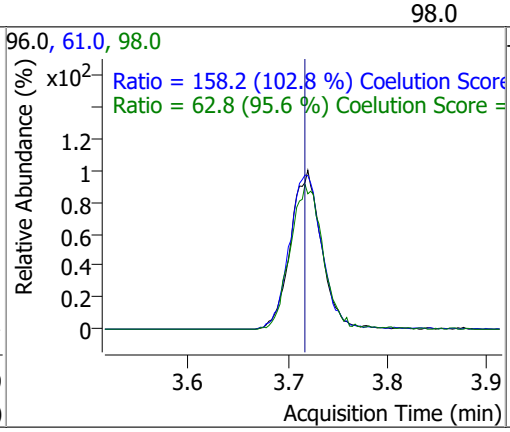


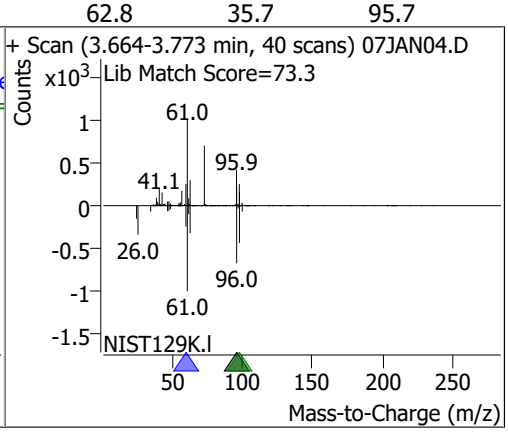




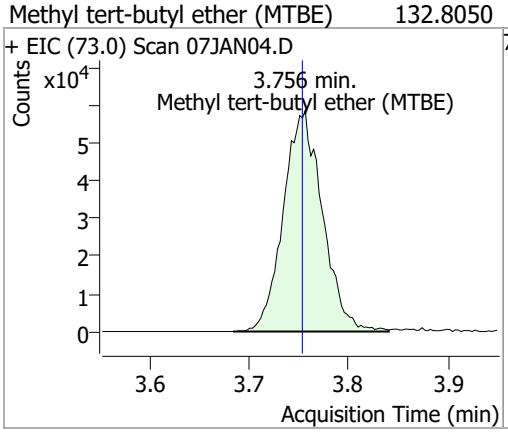
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	129.8320	3.72	0.00	121109	61.0	158.2	123.9	183.9
					98.0	62.8	35.7	95.7

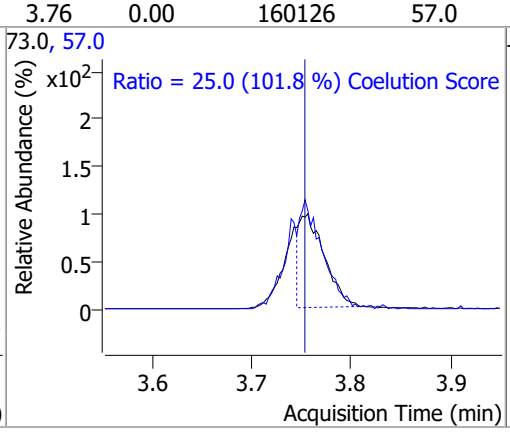


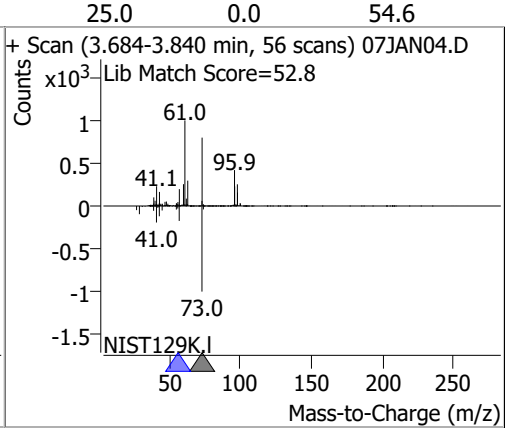




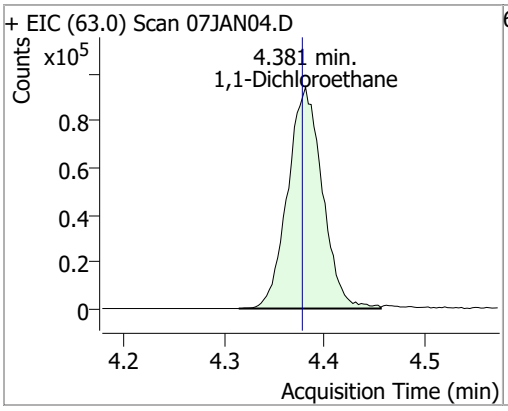
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	132.8050	3.76	0.00	160126	57.0	25.0	0.0	54.6

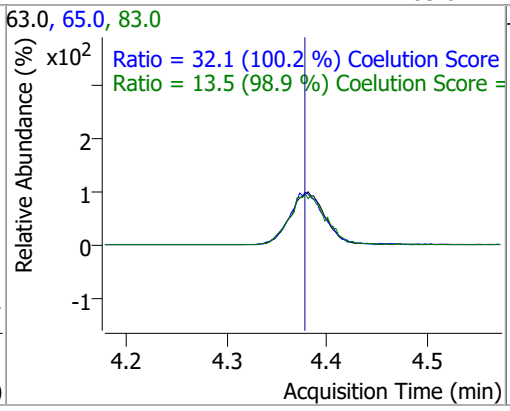


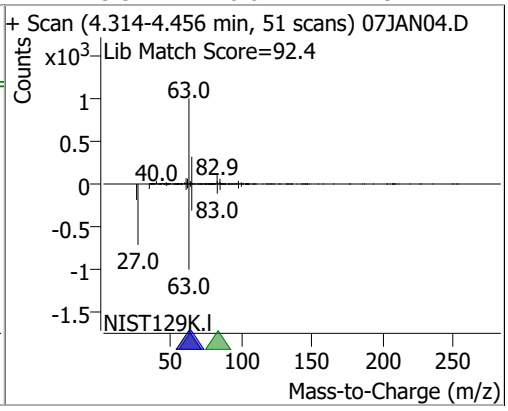




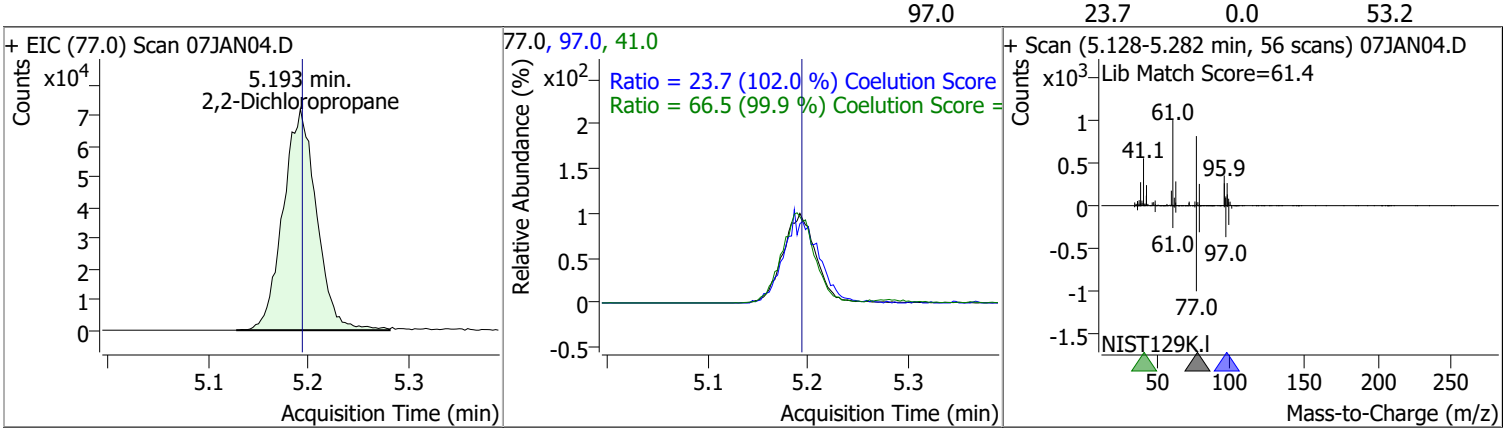
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	131.0789	4.38	0.00	227596	65.0	32.1	2.1	62.1
					83.0	13.5	0.0	43.7



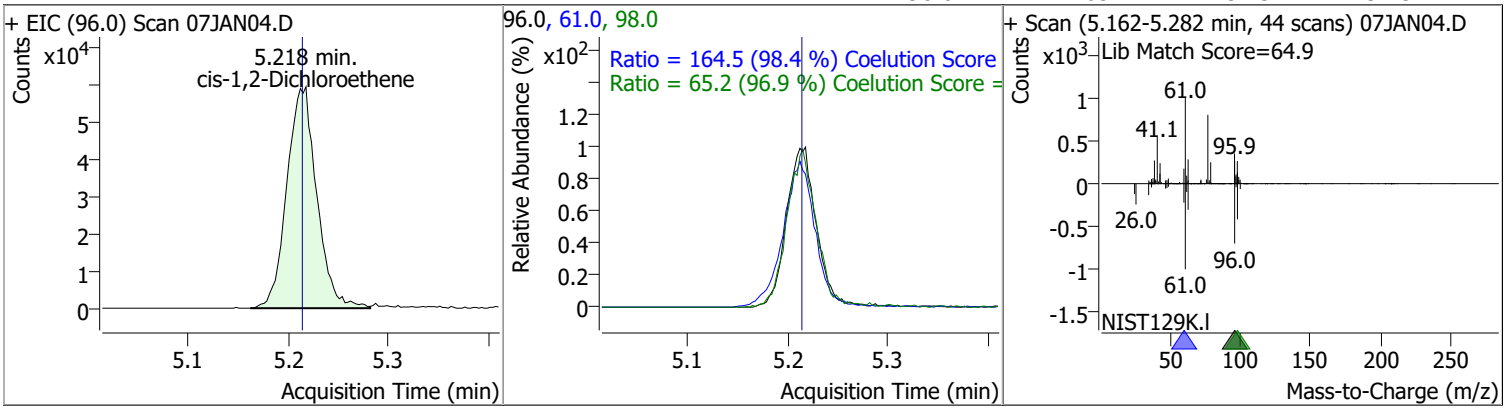




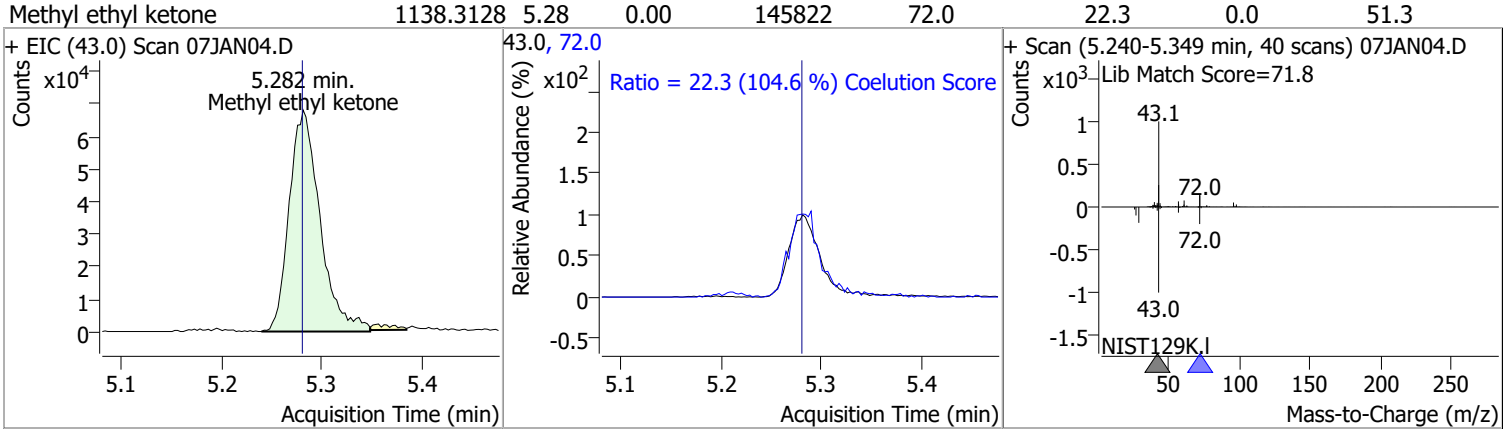
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	128.7744	5.19	0.00	167542	41.0	66.5	36.5	96.5
					97.0	23.7	0.0	53.2



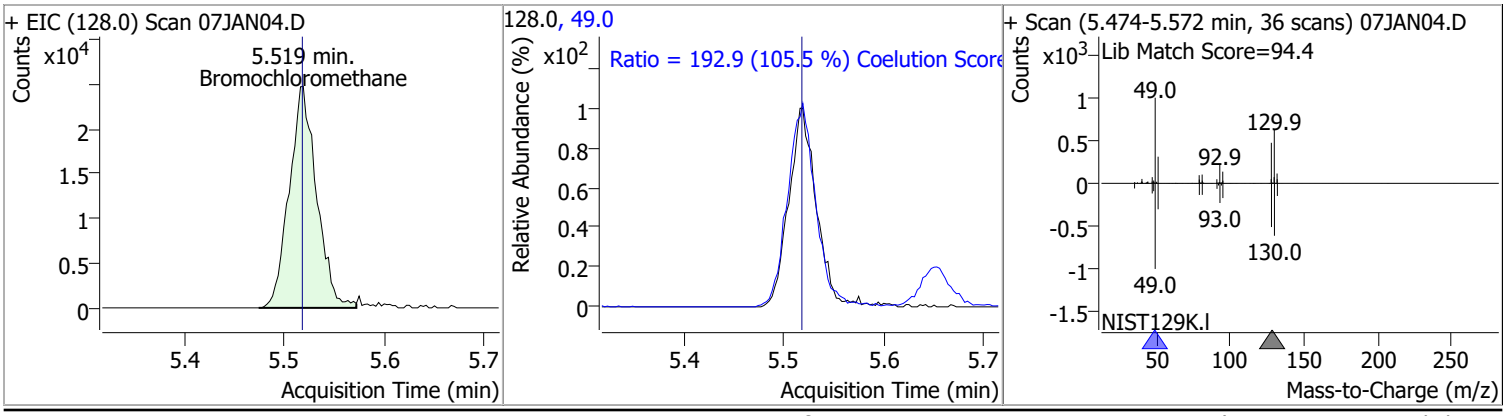
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	129.6836	5.22	0.00	122647	61.0	164.5	137.2	197.2
					98.0	65.2	37.3	97.3



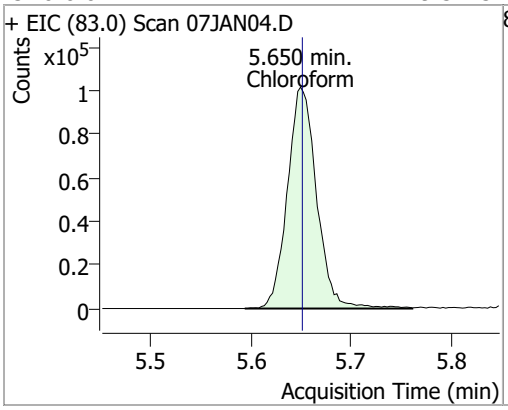
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1138.3128	5.28	0.00	145822	72.0	22.3	0.0	51.3

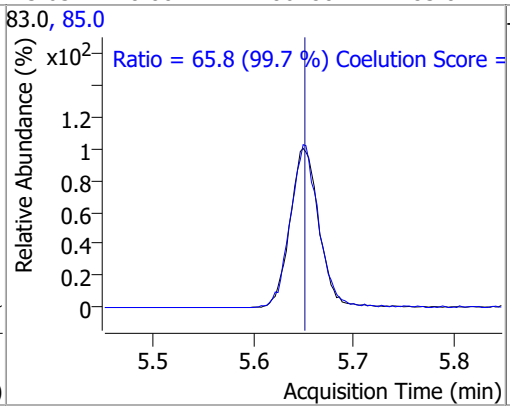


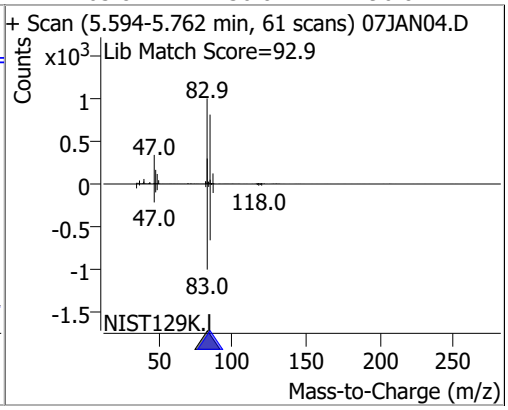
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	120.1958	5.52	0.00	47092	49.0	192.9	152.9	212.9



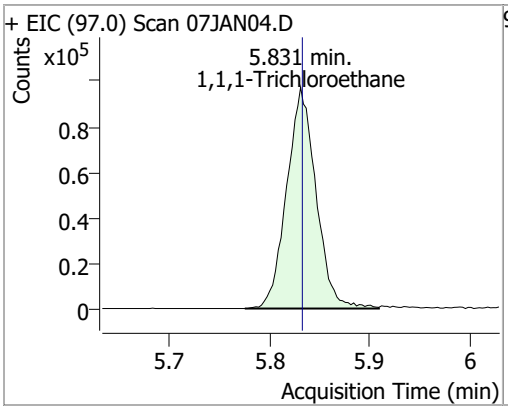
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	119.3745	5.65	0.00	206280	85.0	65.8	36.0	96.0

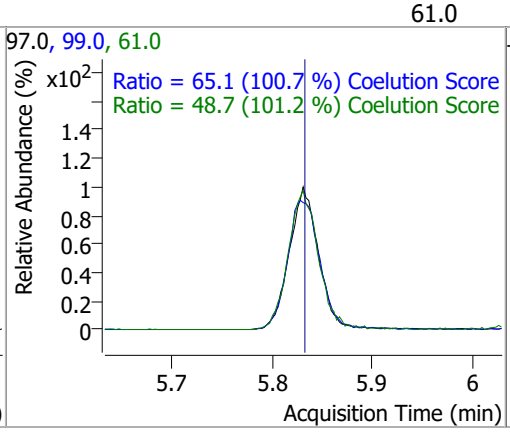


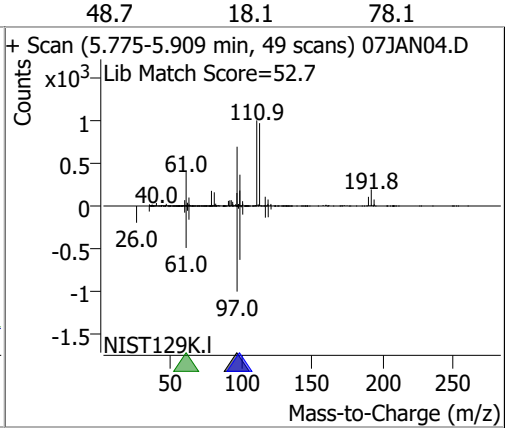




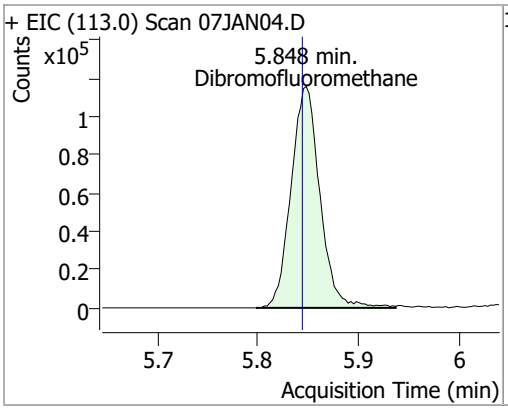
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	122.1545	5.83	0.00	197819	99.0	65.1	34.7	94.7
					61.0	48.7	18.1	78.1

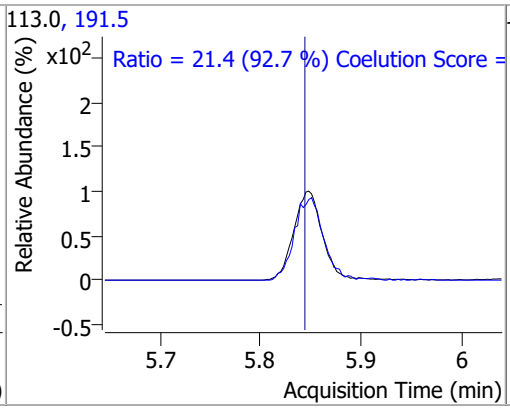


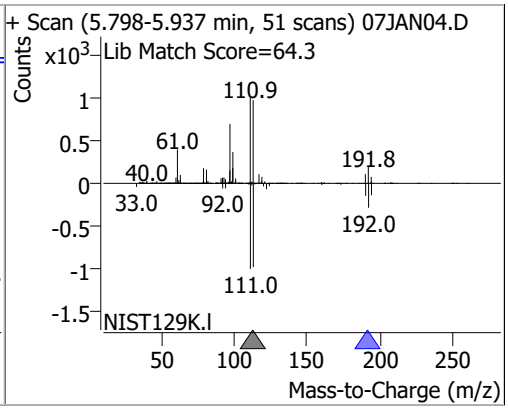




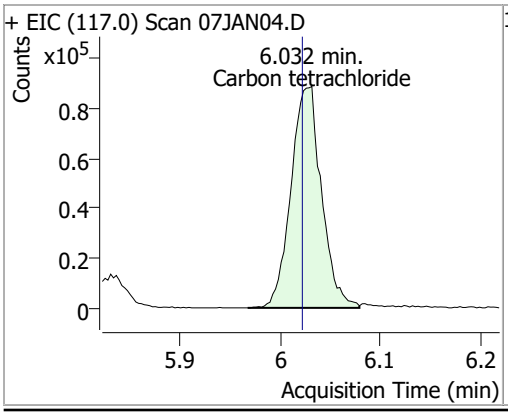
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	270.4515	5.85	0.00	231280	191.5	21.4	0.0	53.1

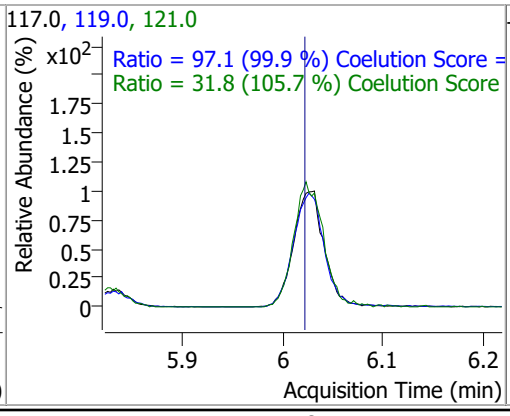


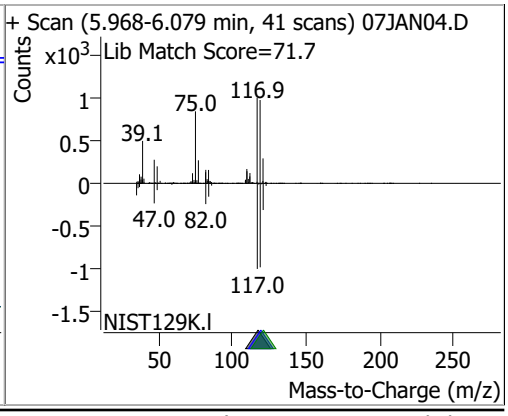




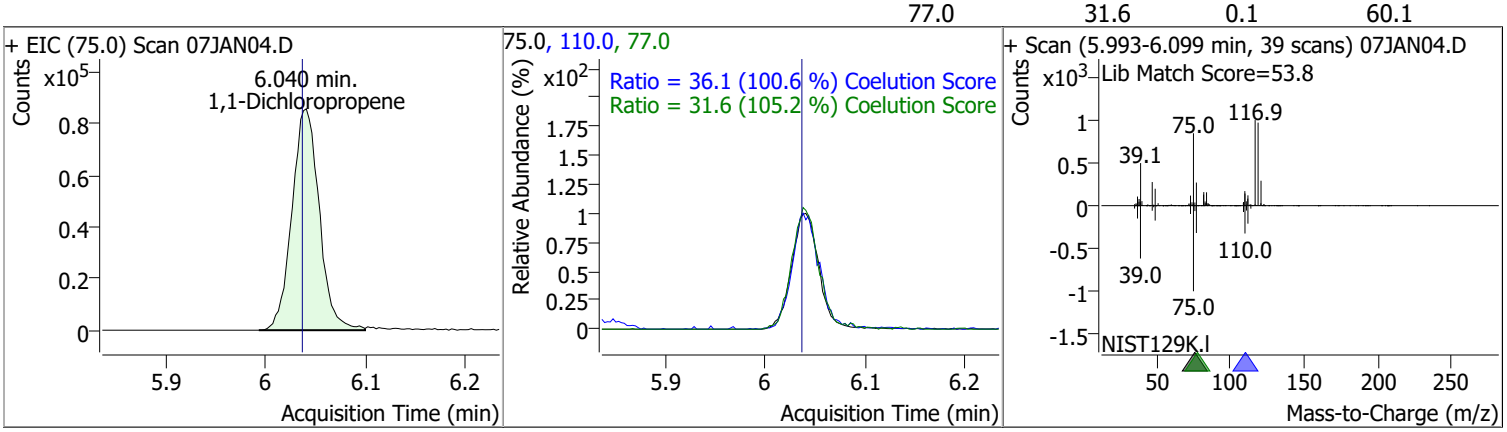
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	117.6380	6.03	0.01	187698	119.0	97.1	67.2	127.2
					121.0	31.8	0.1	60.1



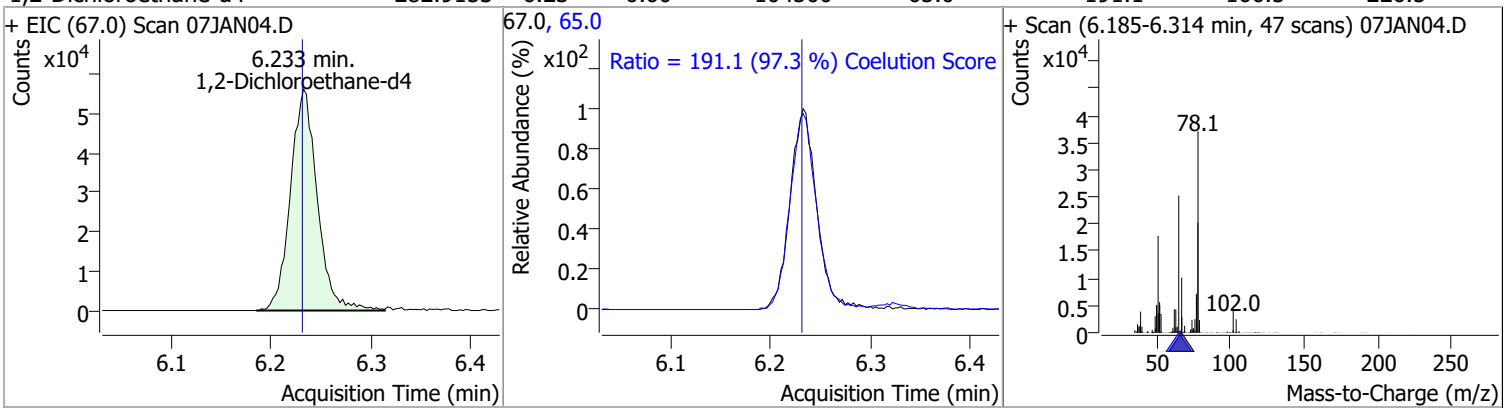




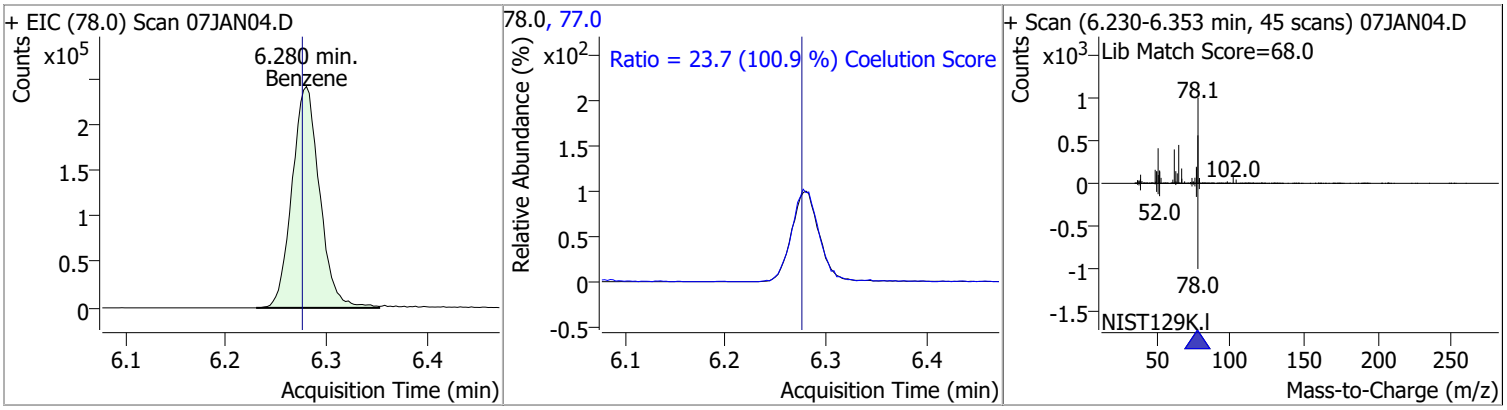
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	118.0827	6.04	0.00	162591	110.0	36.1	5.9	65.9
					77.0	31.6	0.1	60.1



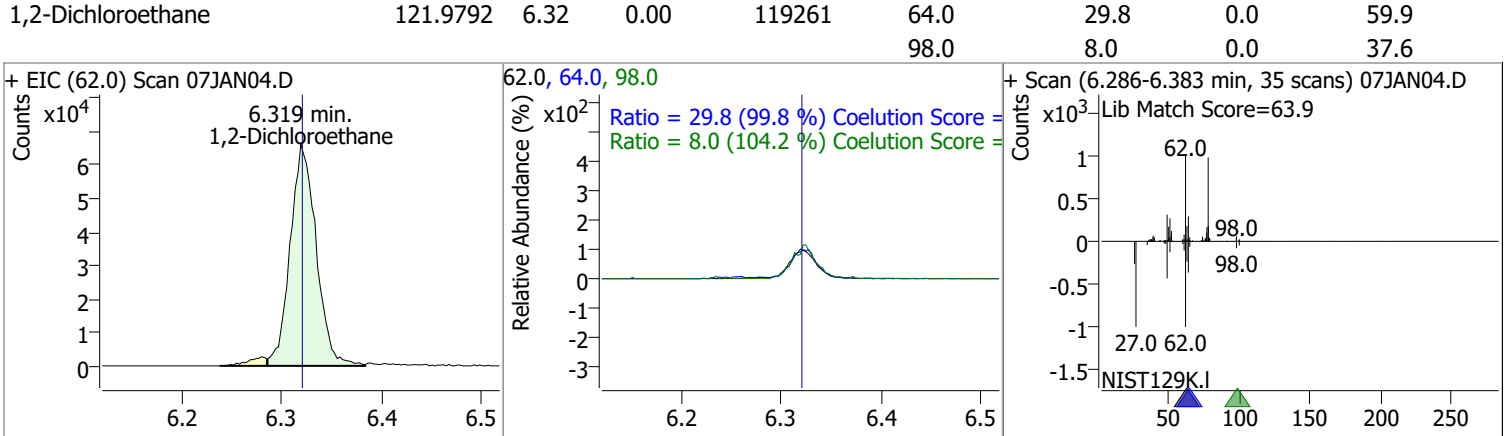
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	282.9153	6.23	0.00	104500	65.0	191.1	166.5	226.5



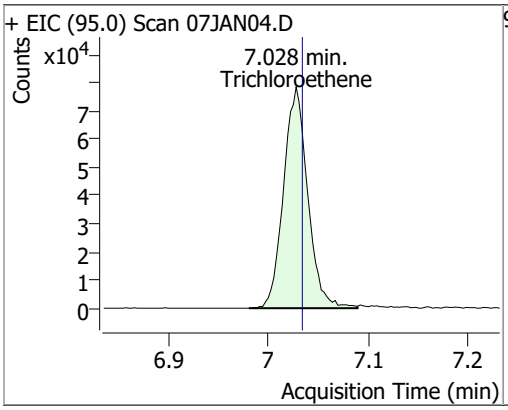
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	127.6093	6.28	0.00	461197	77.0	23.7	0.0	53.5

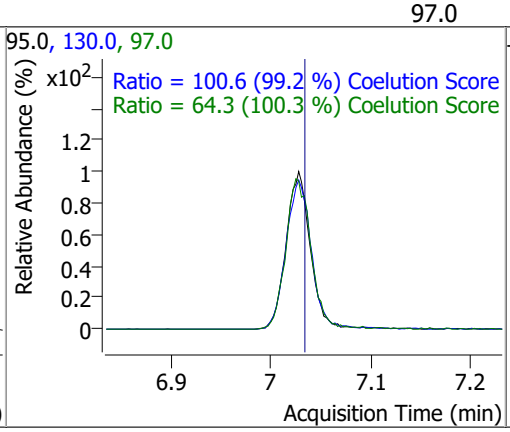


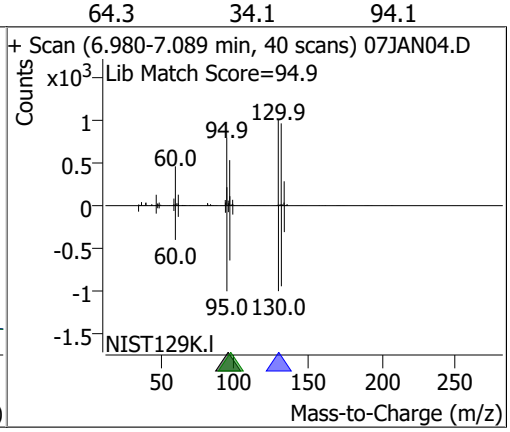
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	121.9792	6.32	0.00	119261	64.0	29.8	0.0	59.9
					98.0	8.0	0.0	37.6



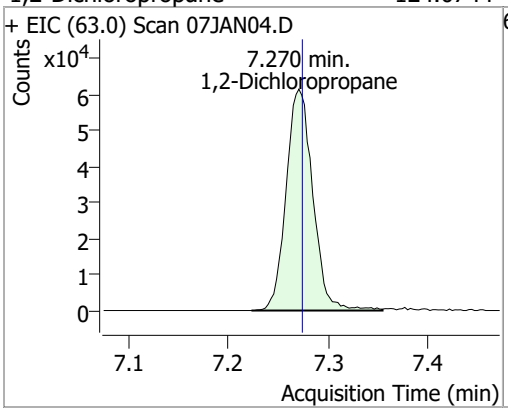
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	125.8989	7.03	0.00	131597	130.0	100.6	71.5	131.5
					97.0	64.3	34.1	94.1

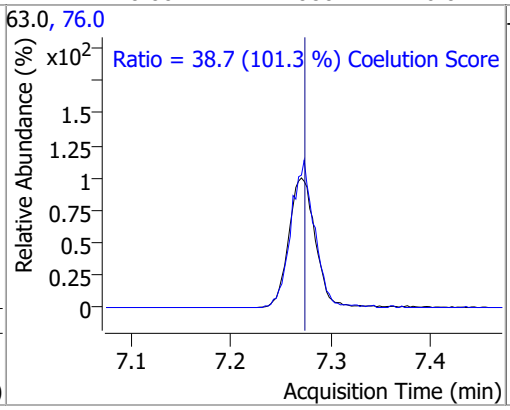


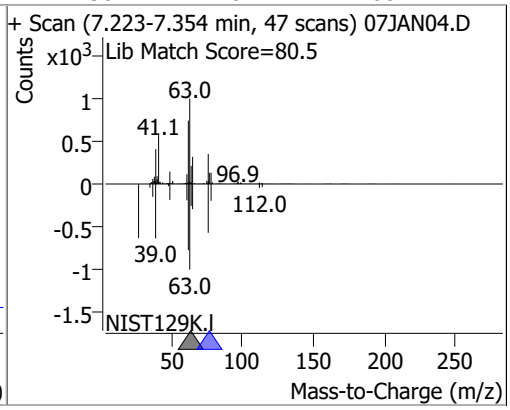




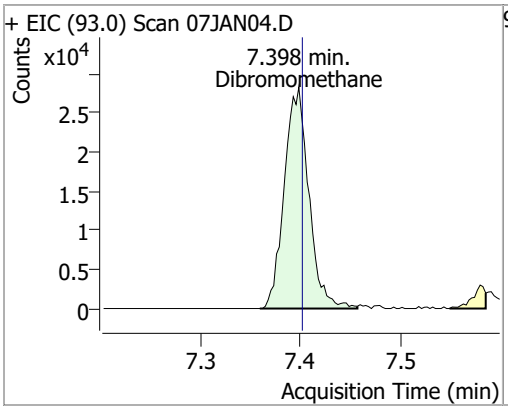
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	124.0744	7.27	0.00	114080	76.0	38.7	8.2	68.2

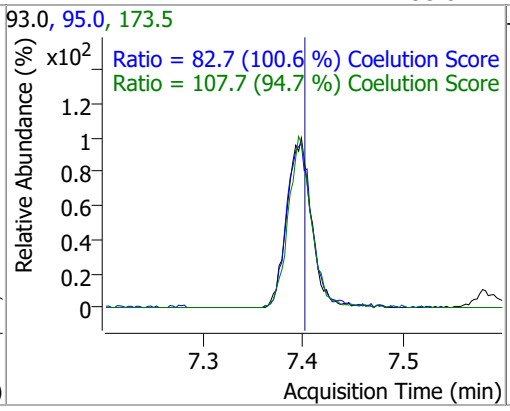


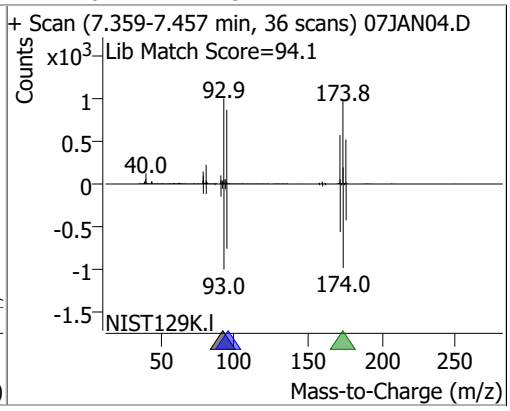




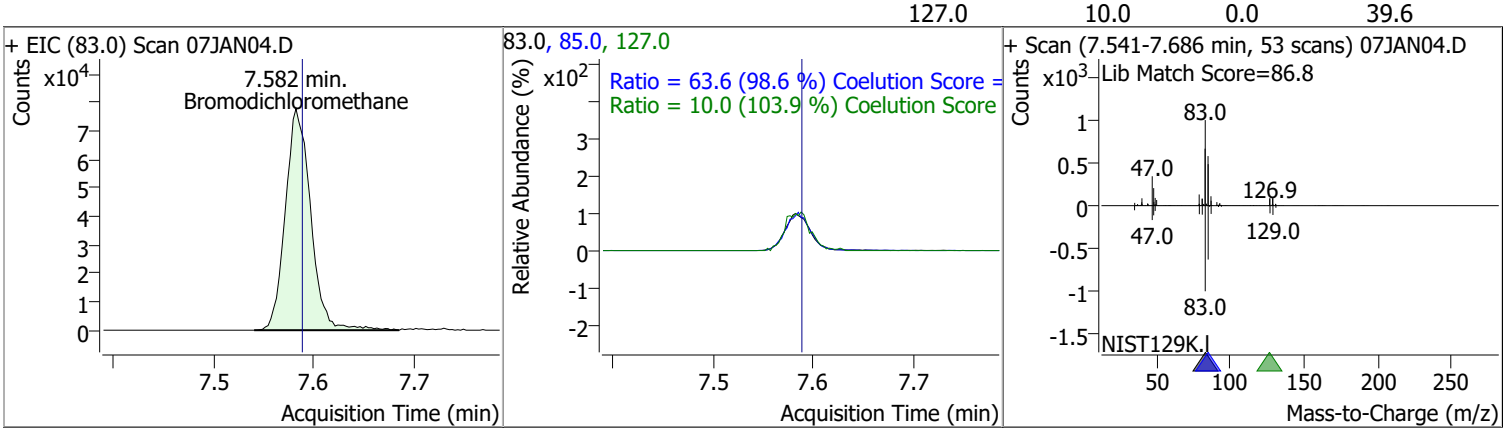
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	124.1493	7.40	0.00	48238	173.5	107.7	83.7	143.7
					95.0	82.7	52.2	112.2



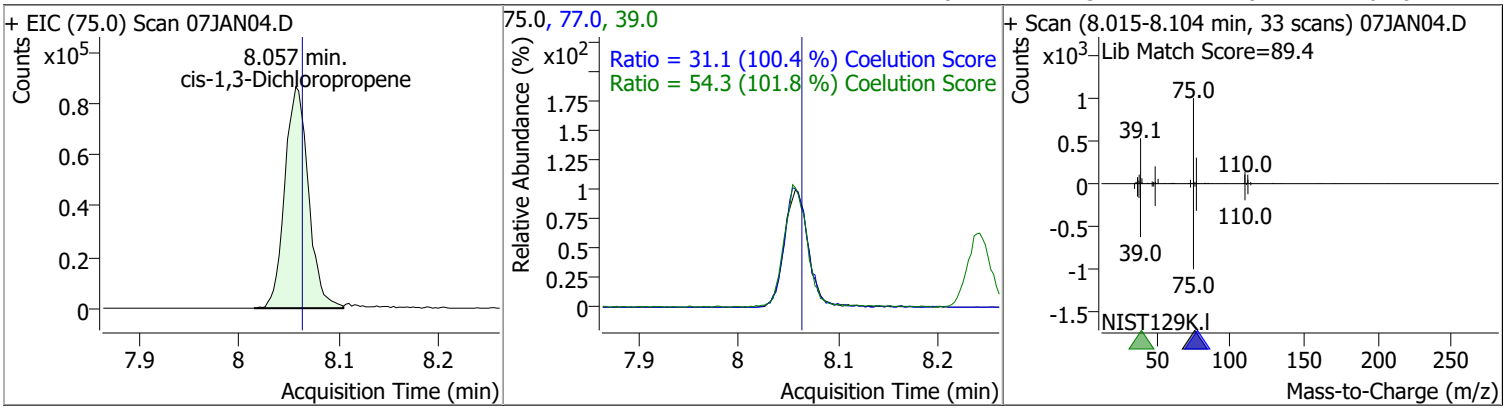




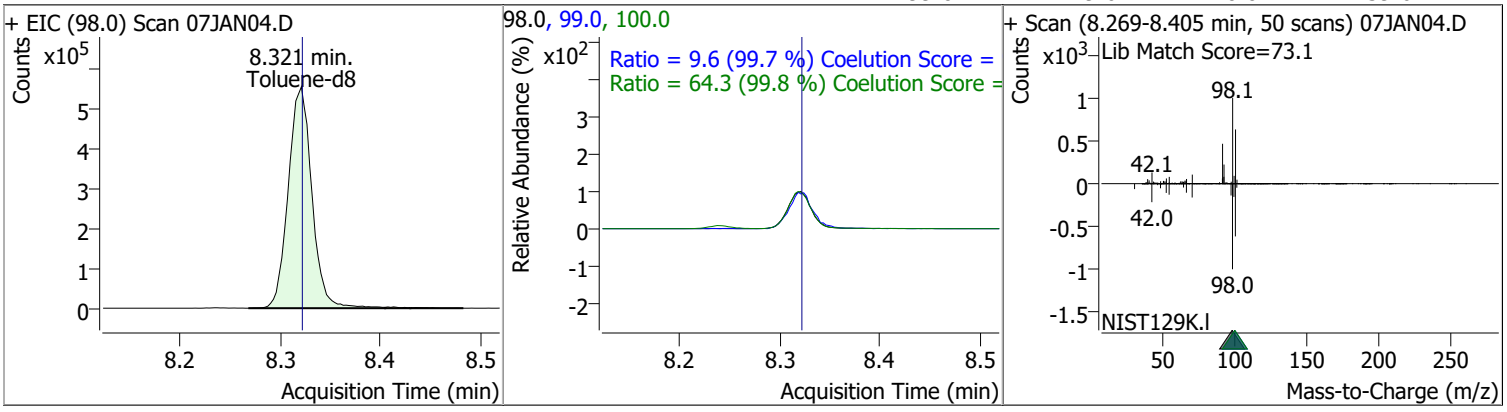
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	127.5739	7.58	0.00	136799	85.0	63.6	34.5	94.5
					127.0	10.0	0.0	39.6



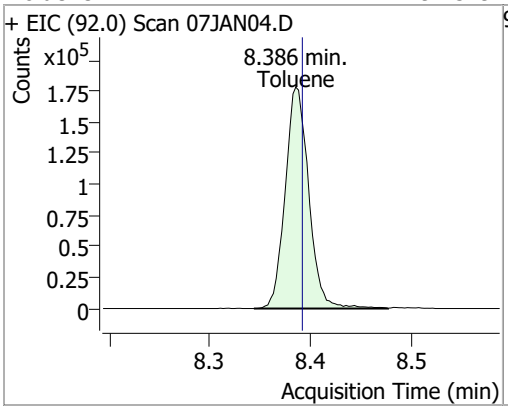
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	117.2568	8.06	0.00	142161	39.0	54.3	23.3	83.3
					77.0	31.1	1.0	61.0

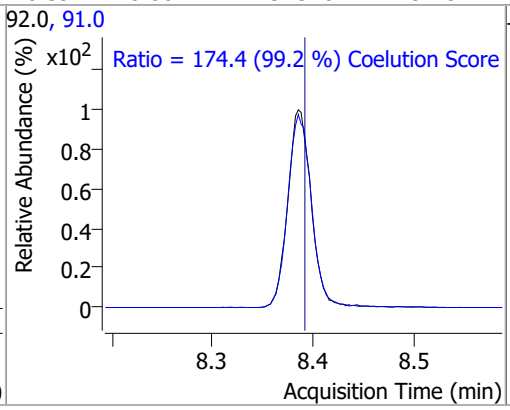


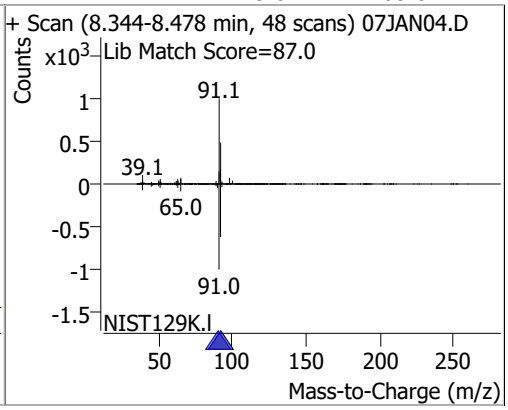
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	271.0520	8.32	0.00	905281	100.0	64.3	34.4	94.4
					99.0	9.6	0.0	39.6



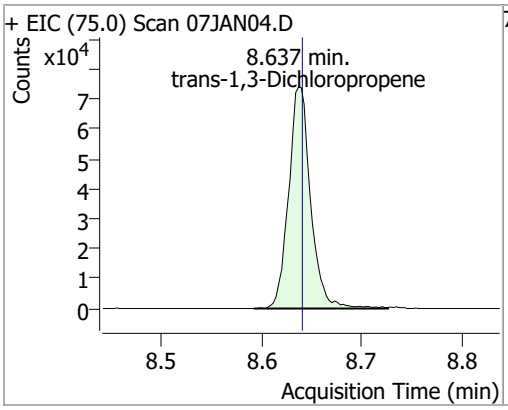
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	129.1519	8.39	0.00	291378	91.0	174.4	145.8	205.8

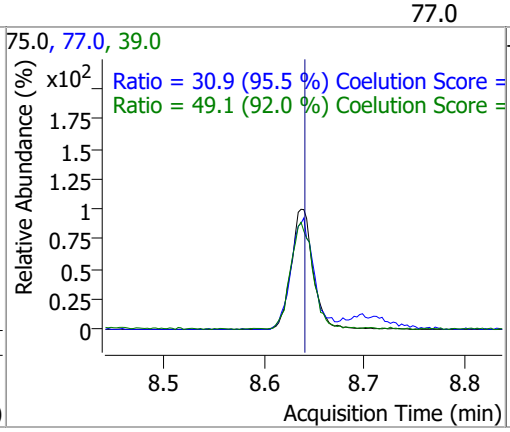


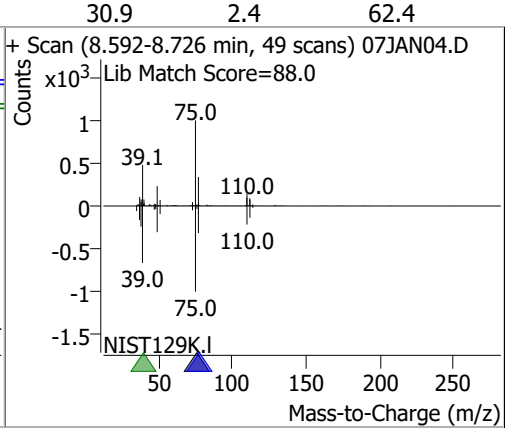




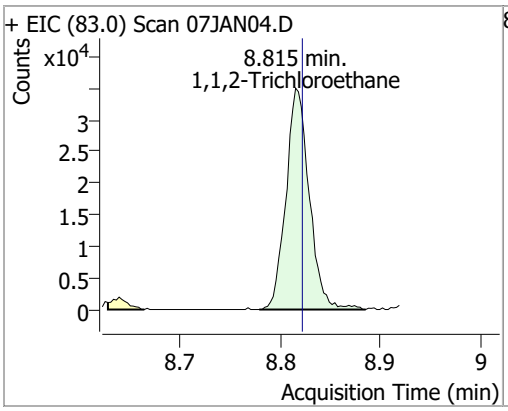
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	131.2571	8.64	0.00	113275	39.0	49.1	23.4	83.4
					77.0	30.9	2.4	62.4

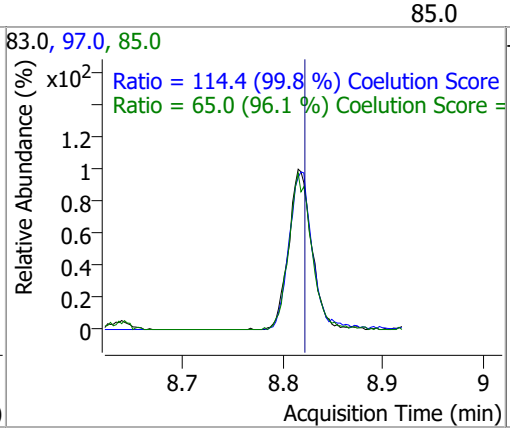


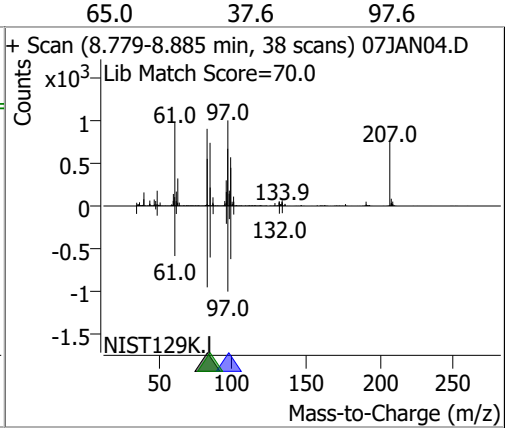




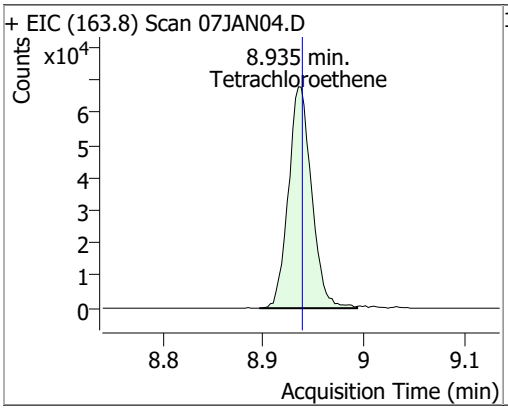
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	125.1597	8.82	0.00	56261	97.0	114.4	84.6	144.6
					85.0	65.0	37.6	97.6

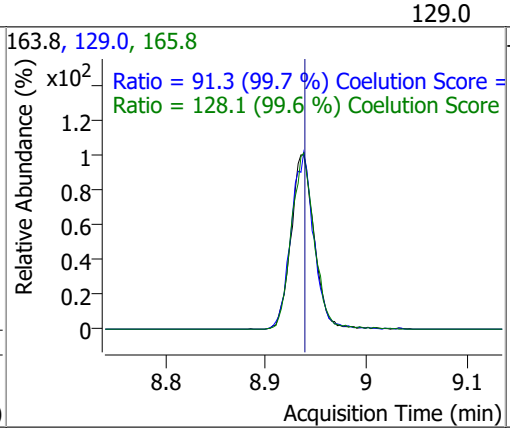


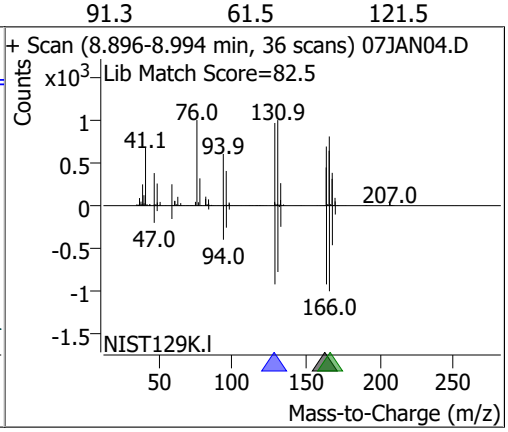




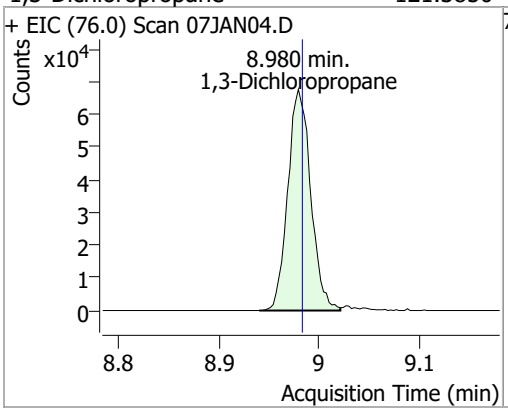
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	118.8837	8.93	0.00	109421	165.8	128.1	98.6	158.6
					129.0	91.3	61.5	121.5

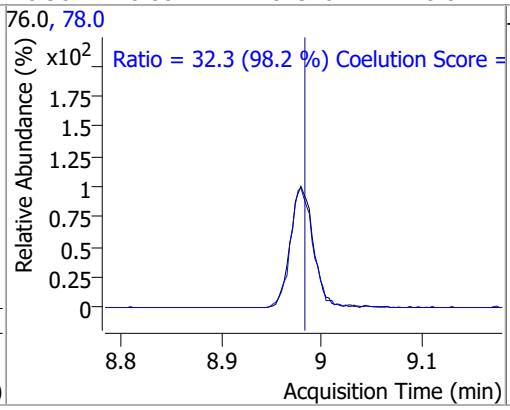


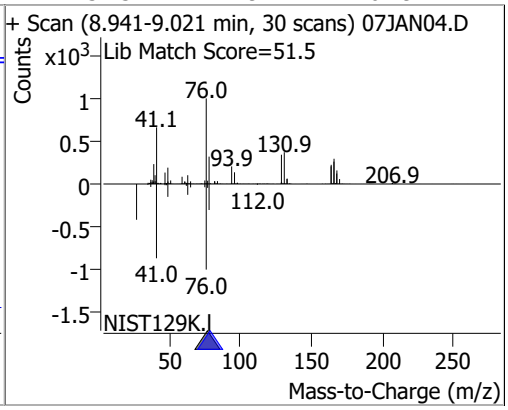




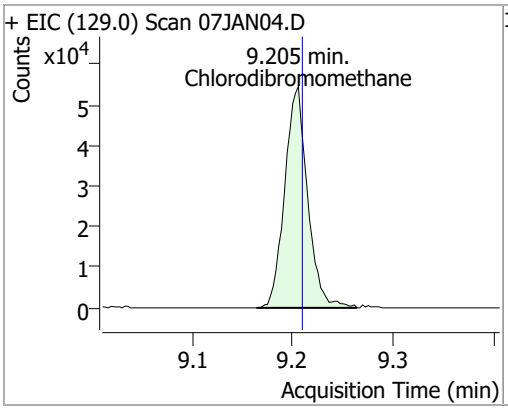
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	121.3850	8.98	0.00	107326	78.0	32.3	2.9	62.9

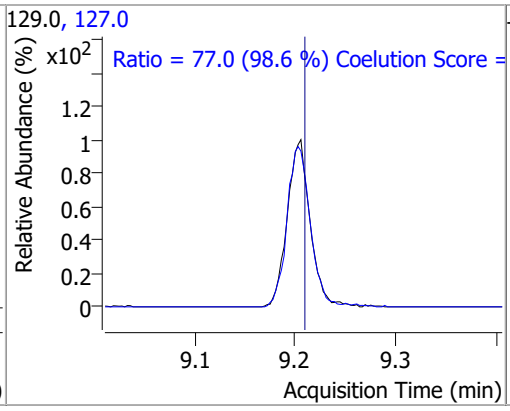


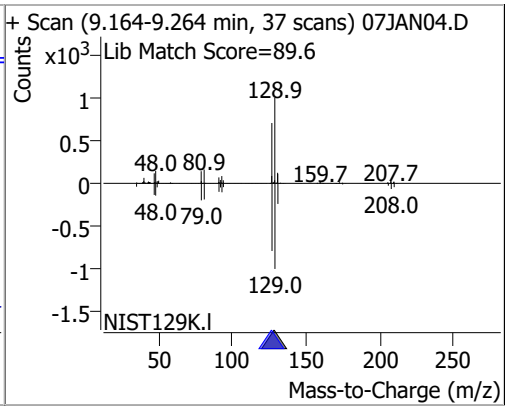




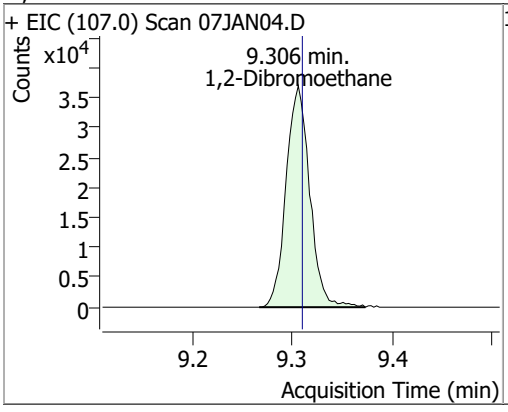
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	122.7318	9.21	0.00	86224	127.0	77.0	48.0	108.0

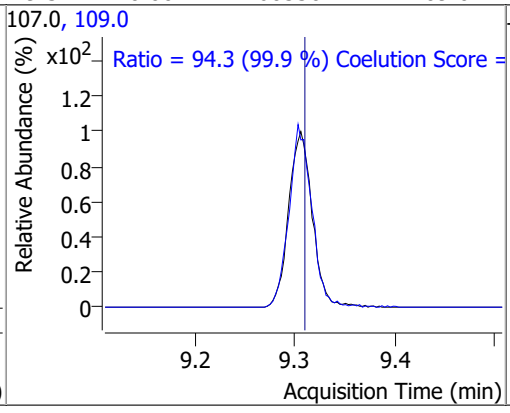


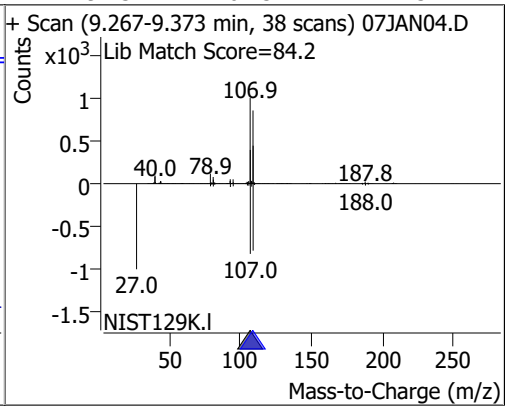




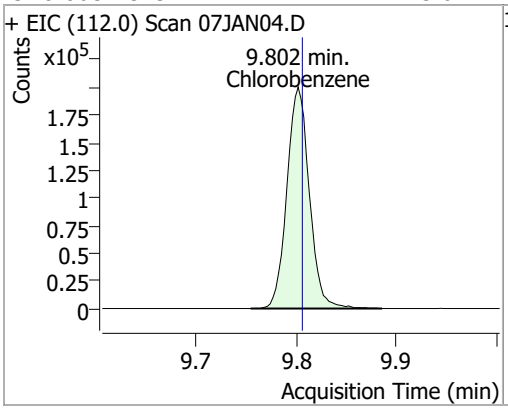
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	123.1519	9.31	0.00	60530	109.0	94.3	64.5	124.5

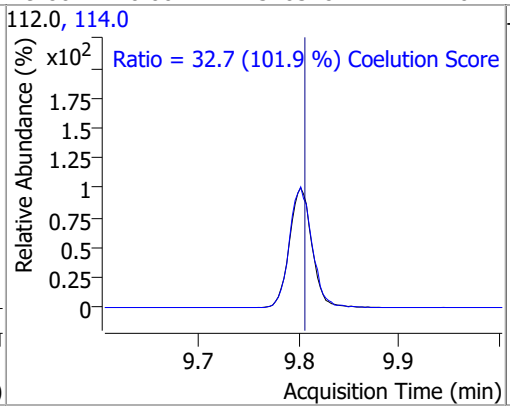


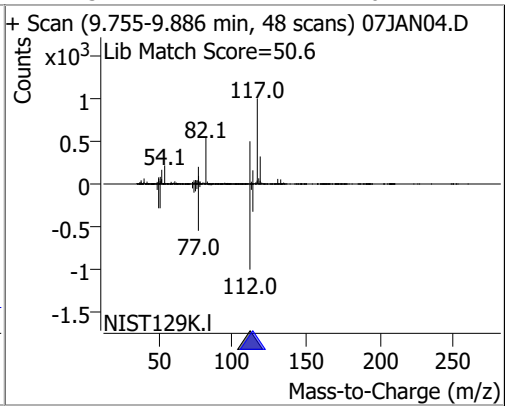




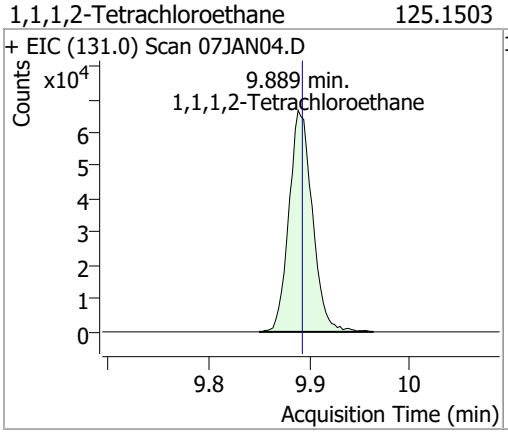
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	125.6477	9.80	0.00	310348	114.0	32.7	2.1	62.1

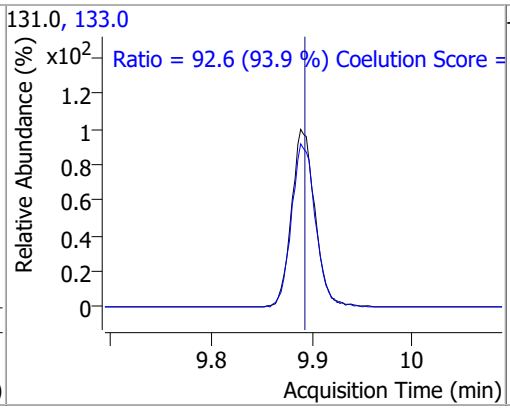


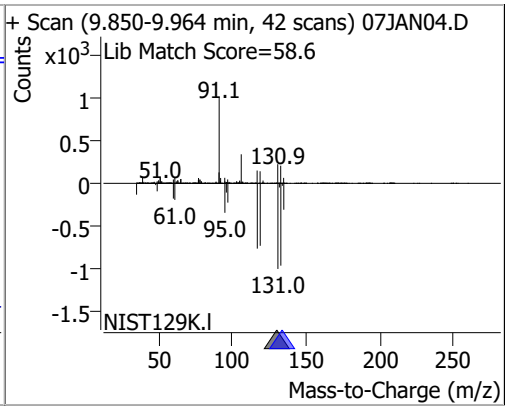




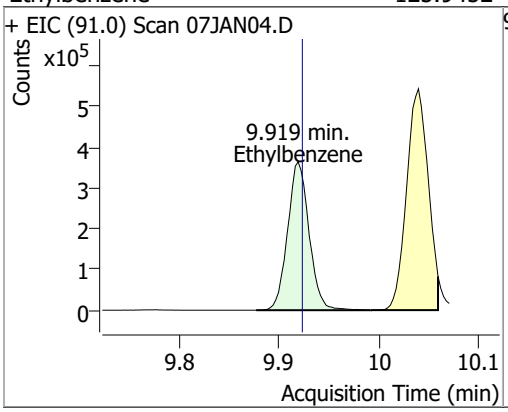
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	125.1503	9.89	0.00	108057	133.0	92.6	68.6	128.6

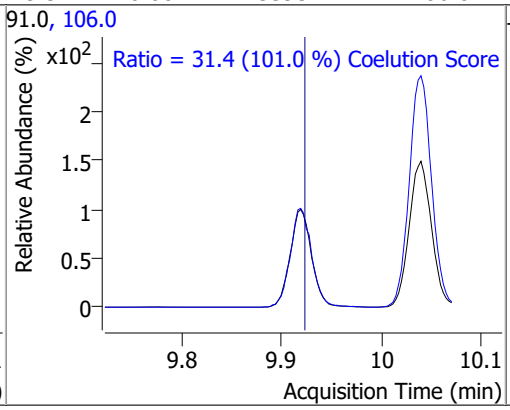


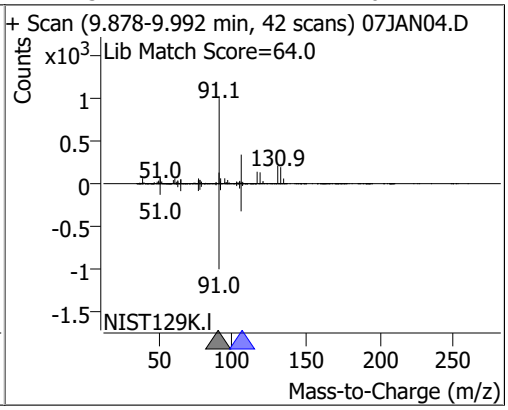




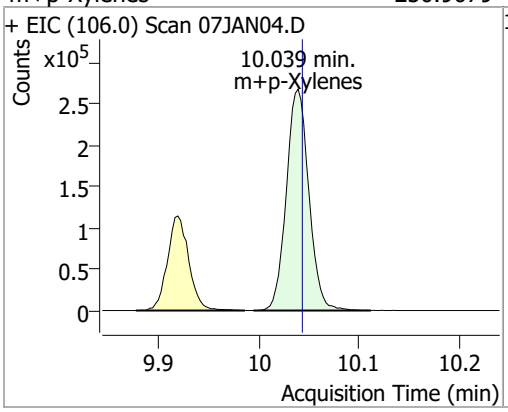
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	125.9452	9.92	0.00	539522	106.0	31.4	1.1	61.1

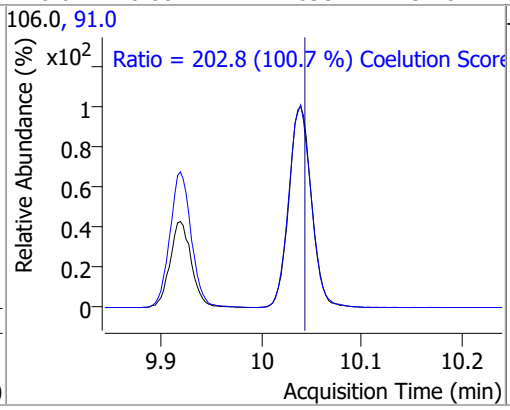


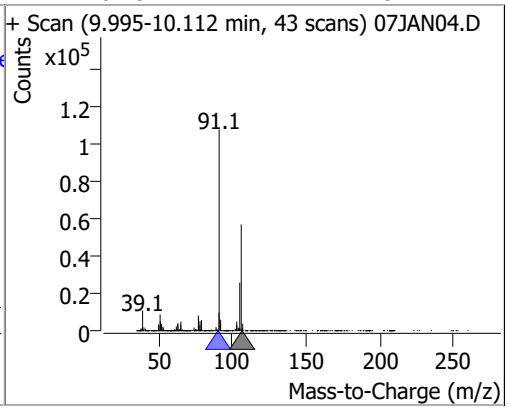




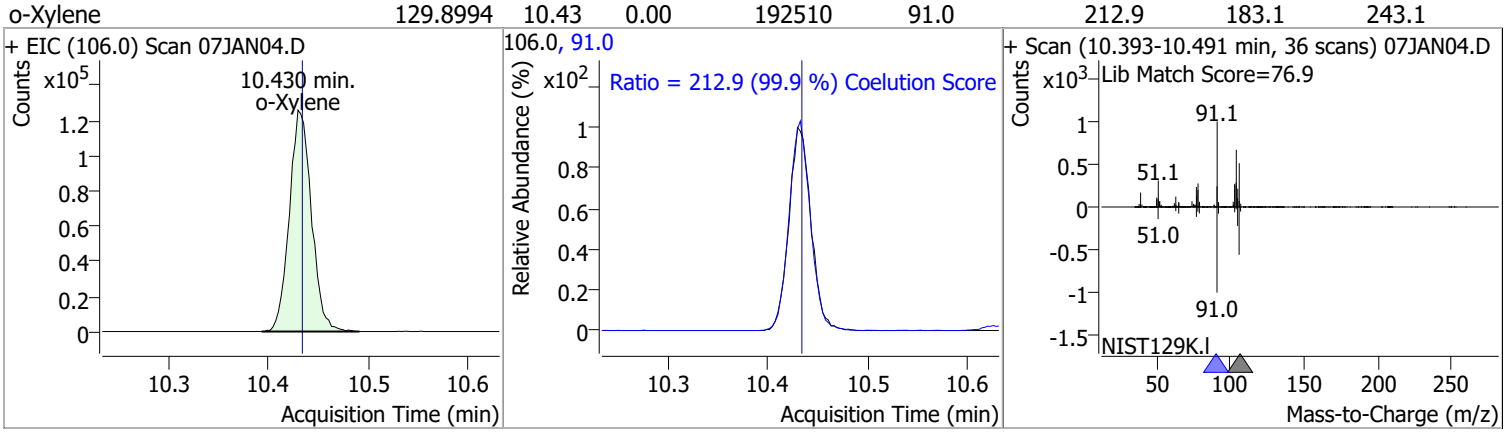
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	250.9079	10.04	0.00	417695	91.0	202.8	171.4	231.4



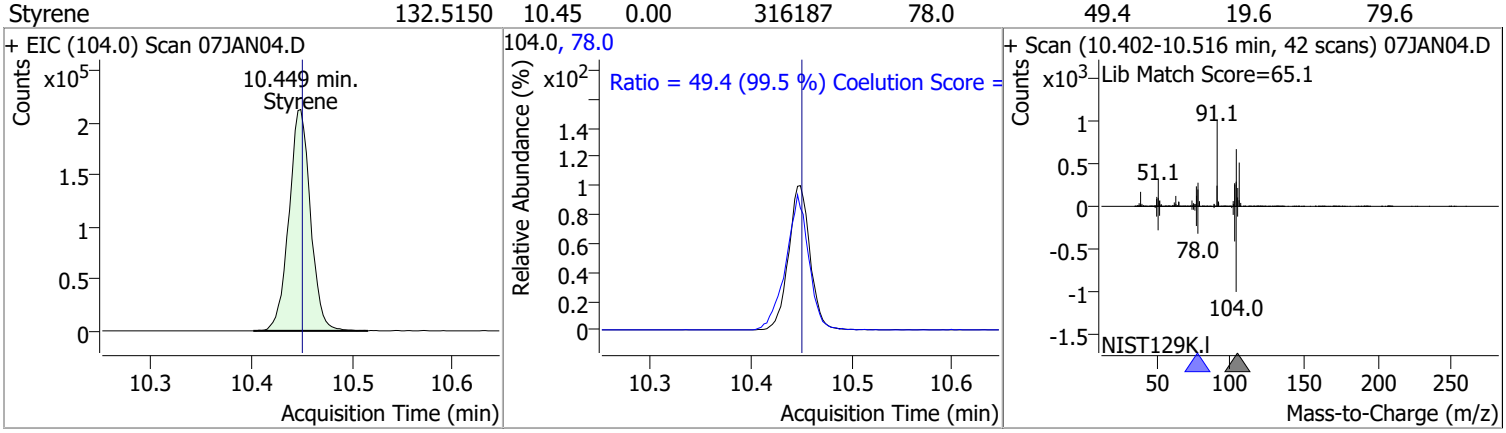




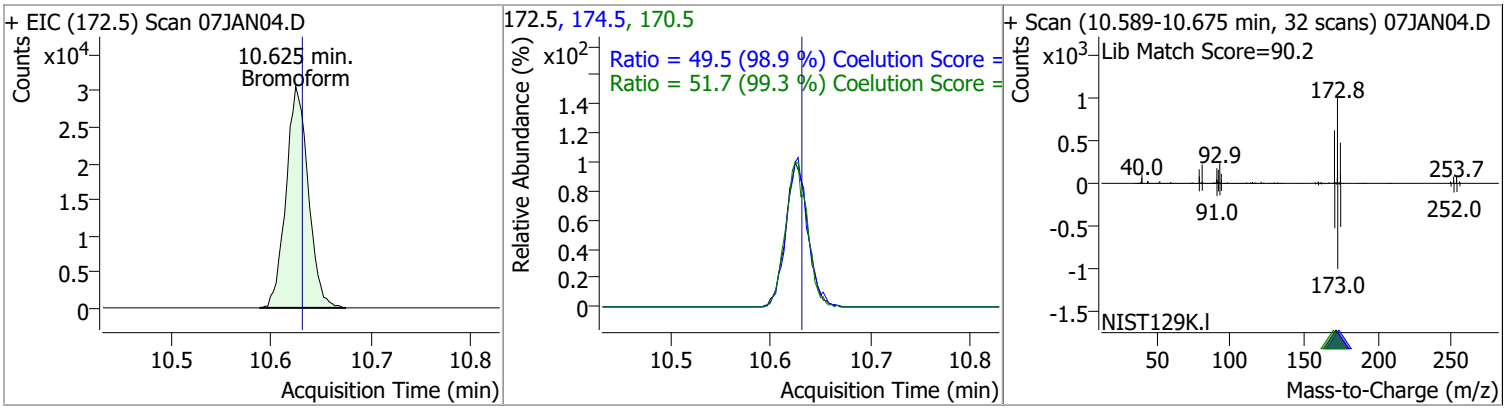
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	129.8994	10.43	0.00	192510	91.0	212.9	183.1	243.1



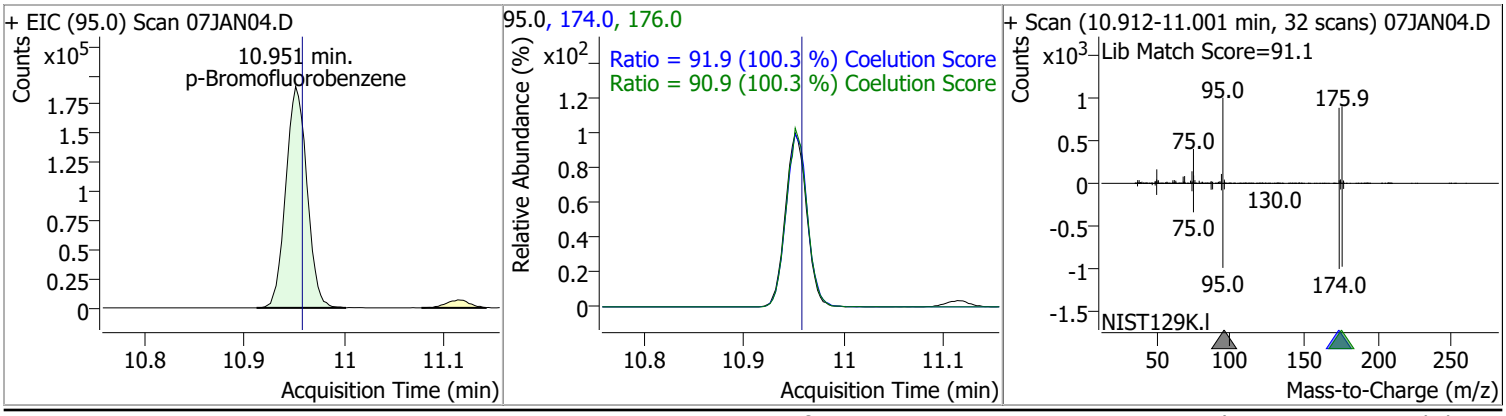
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	132.5150	10.45	0.00	316187	78.0	49.4	19.6	79.6



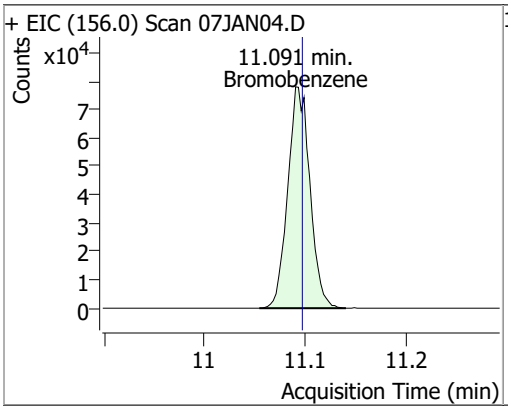
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	131.1556	10.62	0.00	47153	170.5	51.7	22.1	82.1
					174.5	49.5	20.1	80.1

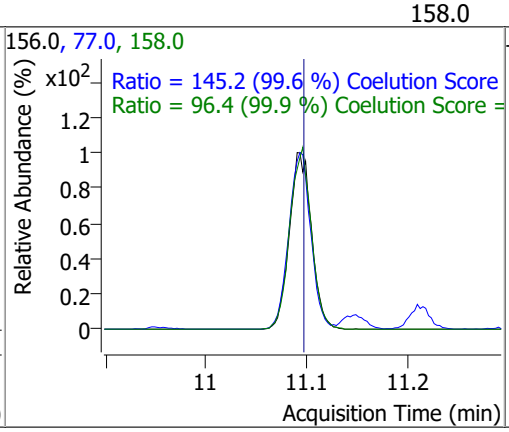


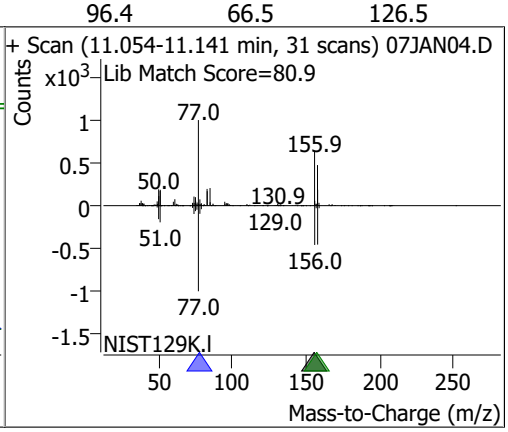
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	271.9817	10.95	0.00	279940	174.0	91.9	61.7	121.7
					176.0	90.9	60.6	120.6



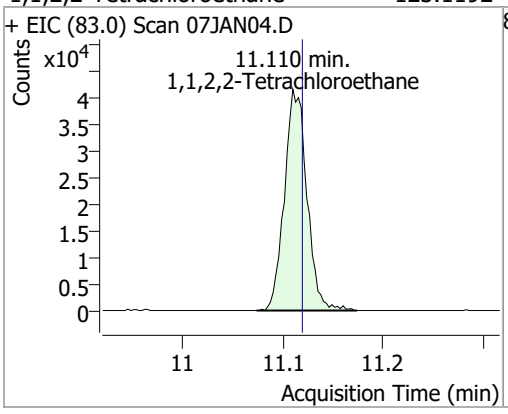
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	131.5649	11.09	0.00	119622	77.0	145.2	115.7	175.7
					158.0	96.4	66.5	126.5

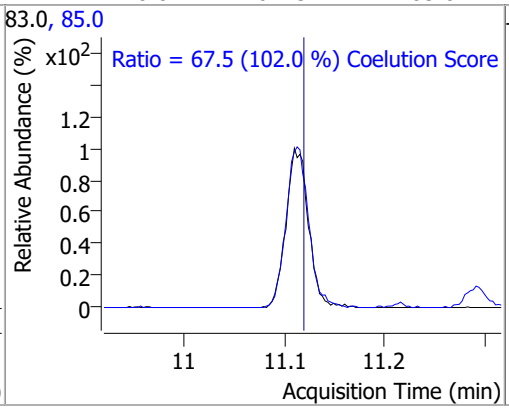


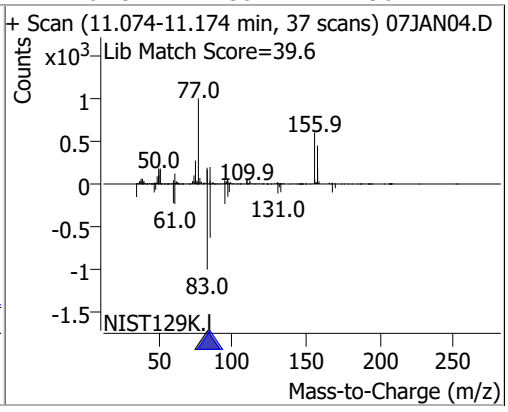




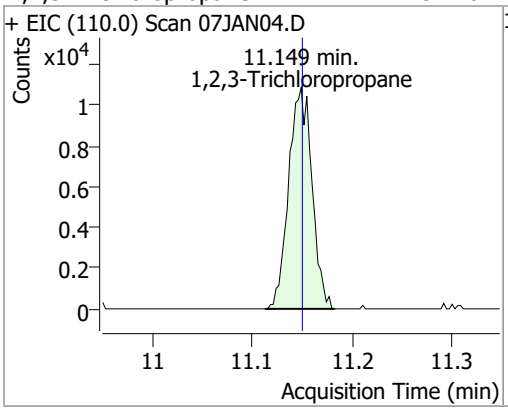
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	123.1192	11.11	-0.01	64431	85.0	67.5	36.2	96.2

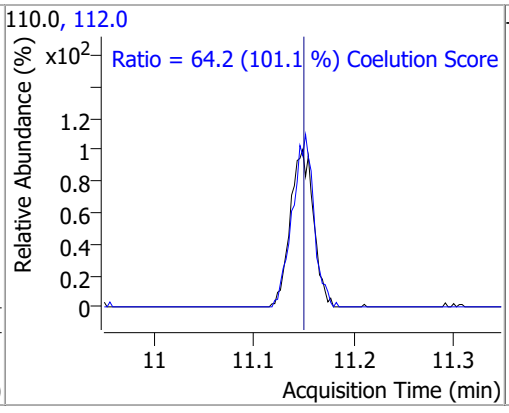


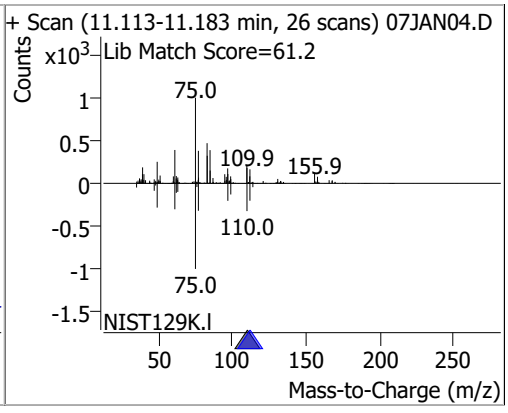




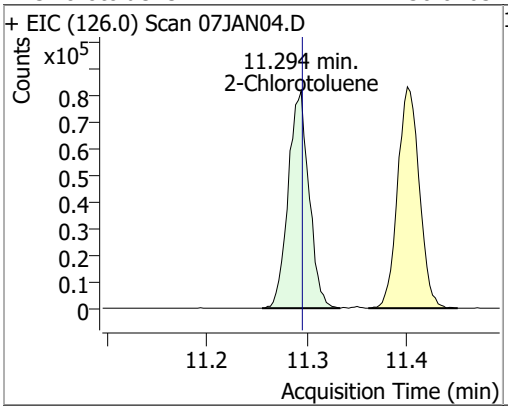
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	125.1764	11.15	0.00	17528	112.0	64.2	33.5	93.5

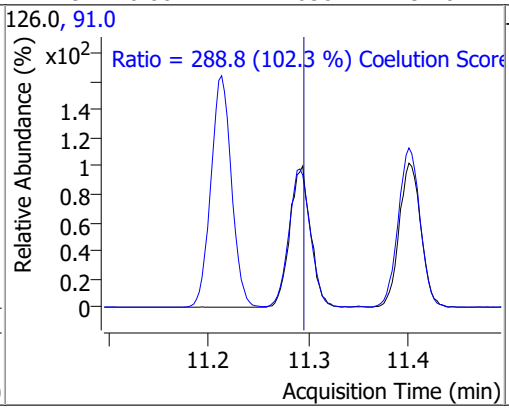


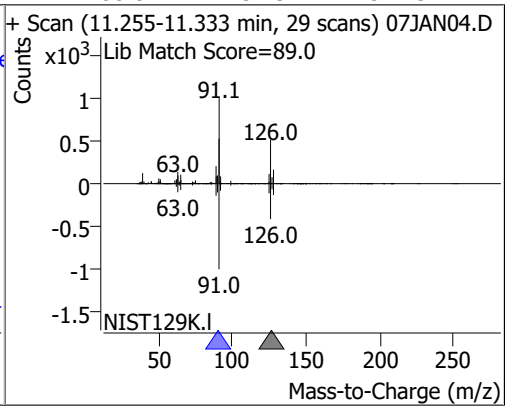




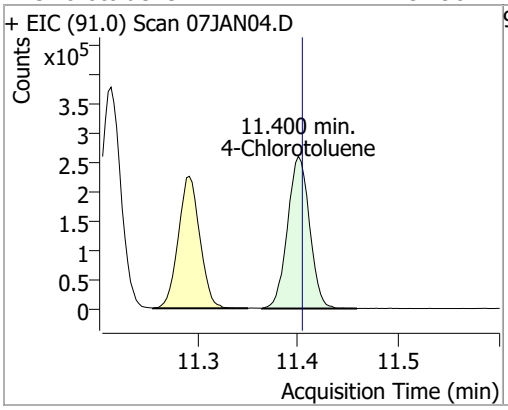
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	130.0465	11.29	0.00	117650	91.0	288.8	252.3	312.3

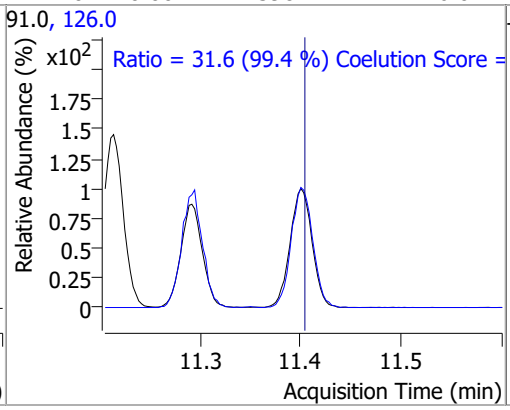


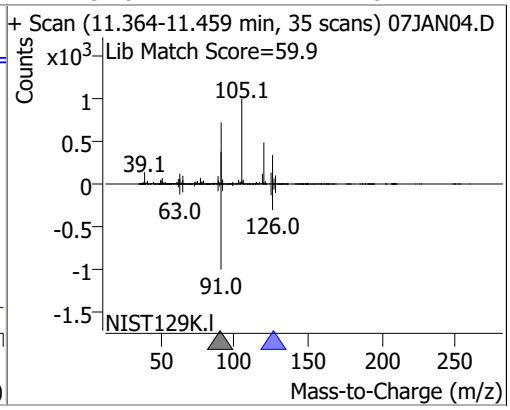




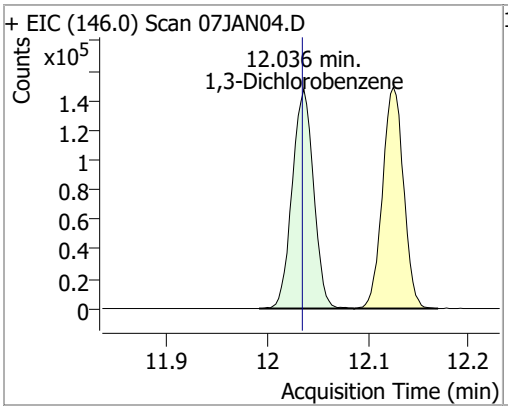
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	134.9812	11.40	0.00	398147	126.0	31.6	1.7	61.7

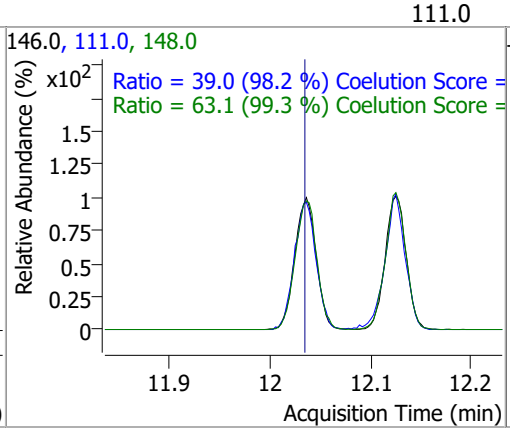


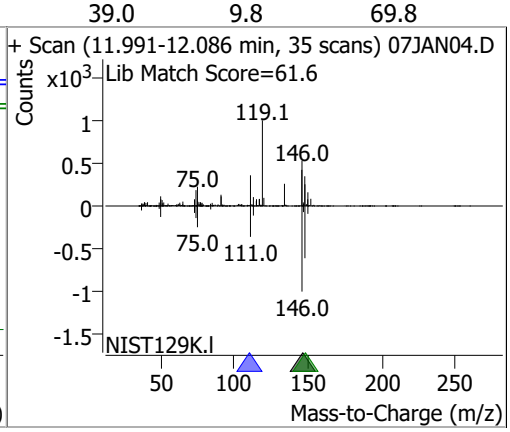




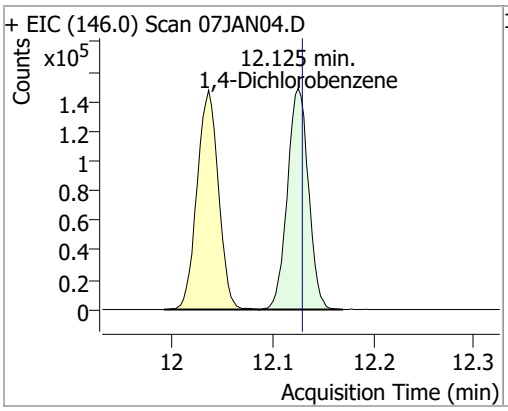
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	131.9494	12.04	0.01	218804	148.0	63.1	33.6	93.6
					111.0	39.0	9.8	69.8

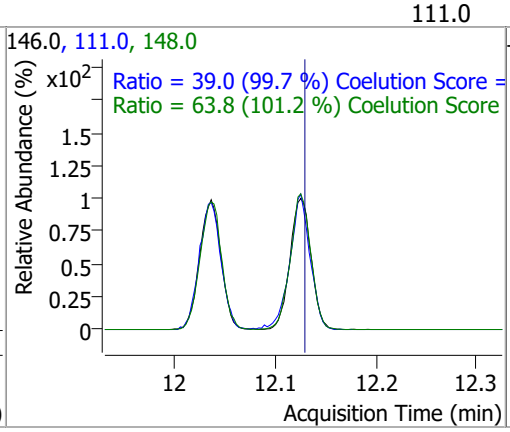


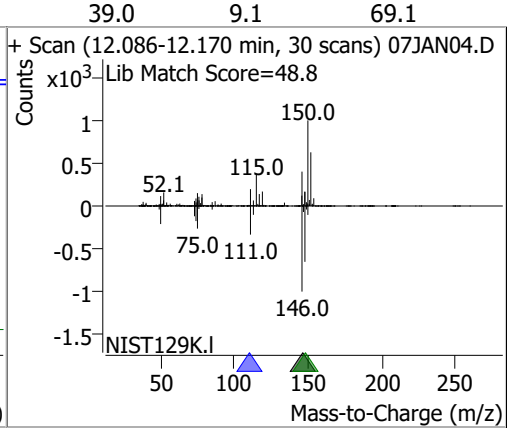




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	128.6355	12.13	0.00	217500	148.0	63.8	33.1	93.1
					111.0	39.0	9.1	69.1

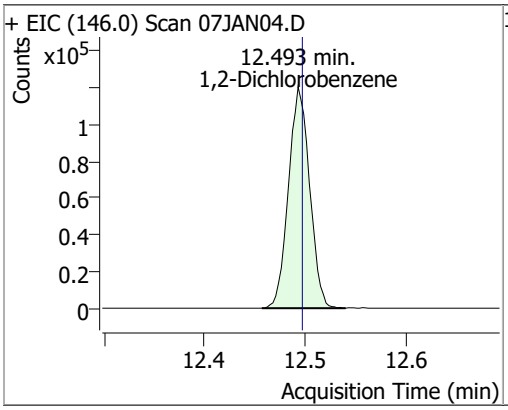


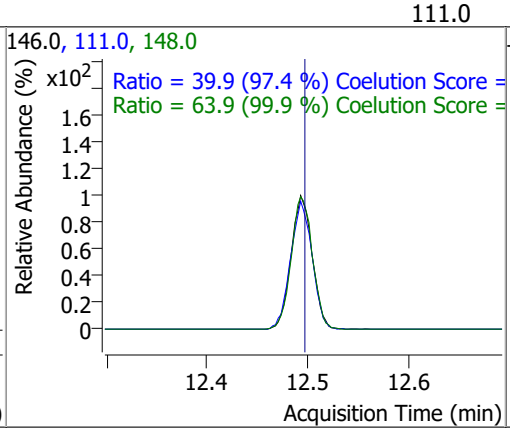


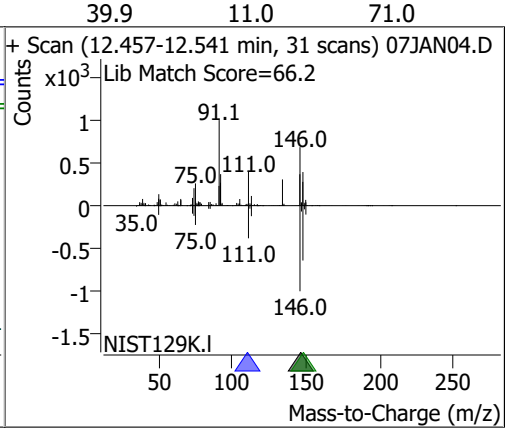


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	125.5596	12.49	0.00	175961	148.0	63.9	33.9	93.9
					111.0	39.9	11.0	71.0







Quantitation Results Report (QT Reviewed)

Data File	07JAN06.D	Operator	MSC
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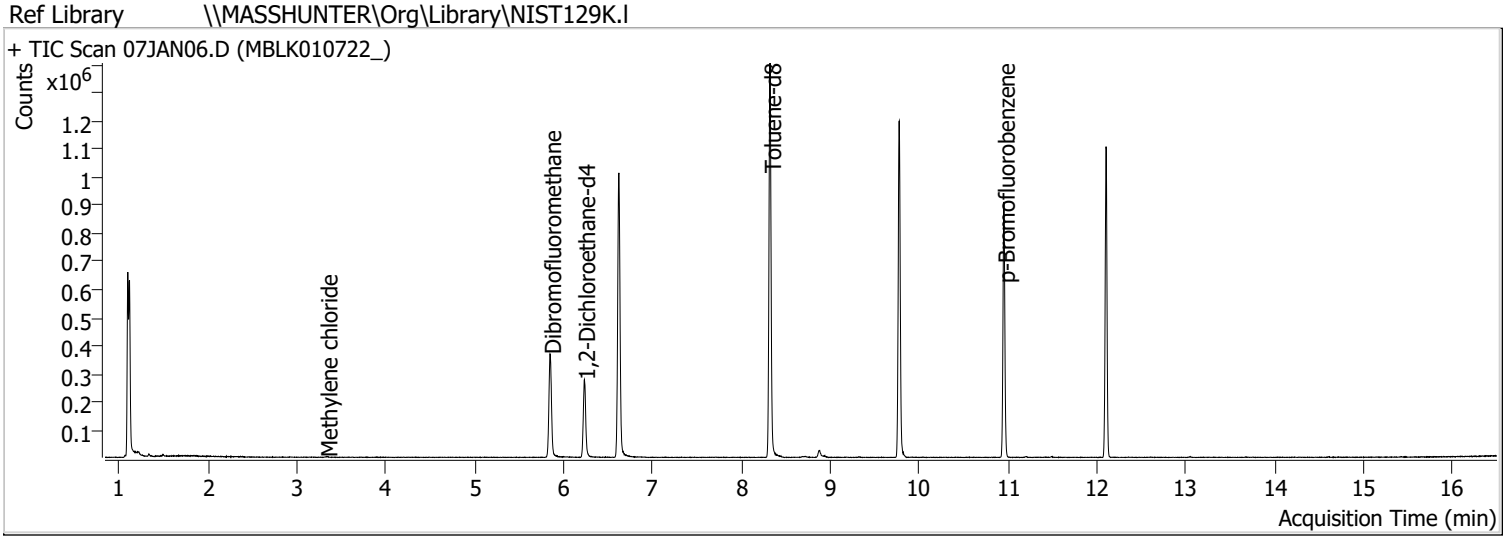
Sample Name	MBLK010722_	Instrument	VOA5975C
Vial	6	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

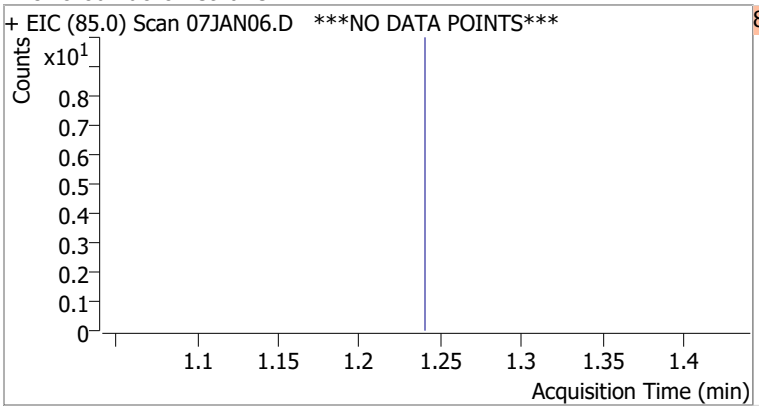


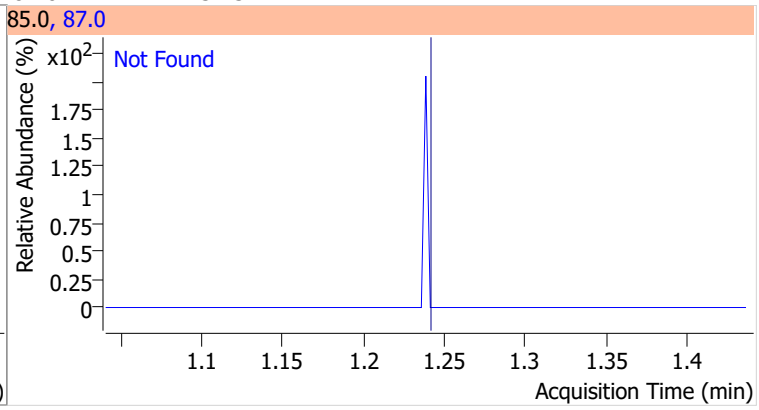
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.620	96.0	869267	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	339487	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.103	152.0	257432	250.0000	ng		0.003
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	221741	270.7666	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 108.31%			
S 1,2-Dichloroethane-d4	6.230	67.0	98740	279.1458	ng		-0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 111.66%			
S Toluene-d8	8.321	98.0	864903	264.3775	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 105.75%			
S p-Bromofluorobenzene	10.951	95.0	259344	274.9894	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 110.00%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	0.000		0	N.D.			
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.330	49.0	1664	1.2889	ng	m	84
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	0.000		0	N.D.			

Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.		
T Carbon tetrachloride	0.000		0	N.D.		
T 1,1-Dichloropropene	0.000		0	N.D.		
T Benzene	0.000		0	N.D.		
T 1,2-Dichloroethane	0.000		0	N.D.		
T Trichloroethene	0.000		0	N.D.		
T 1,2-Dichloropropane	0.000		0	N.D.		
T Dibromomethane	0.000		0	N.D.		
T Bromodichloromethane	0.000		0	N.D.		
T cis-1,3-Dichloropropene	0.000		0	N.D.		
T Toluene	0.000		0	N.D.		
T trans-1,3-Dichloropropene	0.000		0	N.D.		
T 1,1,2-Trichloroethane	0.000		0	N.D.		
T Tetrachloroethene	0.000		0	N.D.		
T 1,3-Dichloropropane	0.000		0	N.D.		
T Chlorodibromomethane	0.000		0	N.D.		
T 1,2-Dibromoethane	0.000		0	N.D.		
T Chlorobenzene	0.000		0	N.D.		
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.		
T Ethylbenzene	0.000		0	N.D.		
T m+p-Xylenes	0.000		0	N.D.		
T o-Xylene	0.000		0	N.D.		
T Styrene	0.000		0	N.D.		
T Bromoform	0.000		0	N.D.		
T Bromobenzene	0.000		0	N.D.		
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
T 1,2,3-Trichloropropane	0.000		0	N.D.		
T 2-Chlorotoluene	0.000		0	N.D.		
T 4-Chlorotoluene	0.000		0	N.D.		
T 1,3-Dichlorobenzene	0.000		0	N.D.		
T 1,4-Dichlorobenzene	0.000		0	N.D.		
T 1,2-Dichlorobenzene	0.000		0	N.D.		

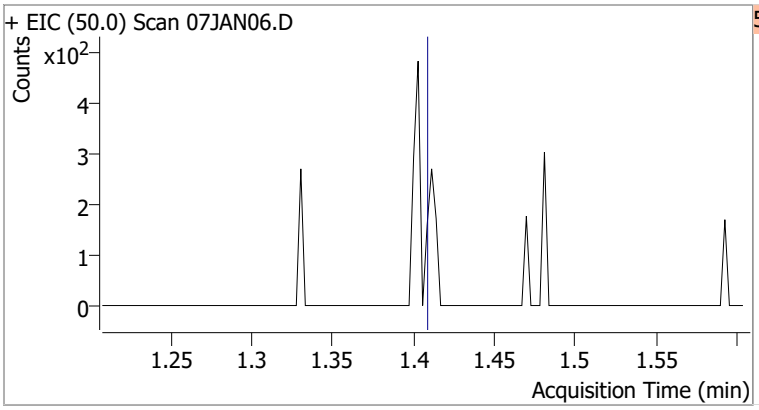
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

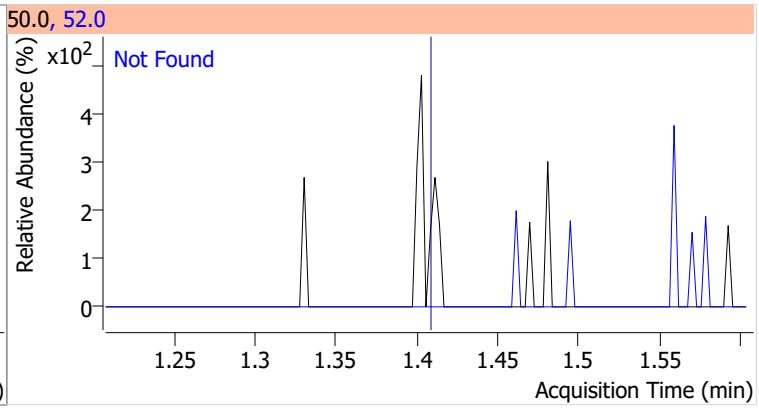
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3



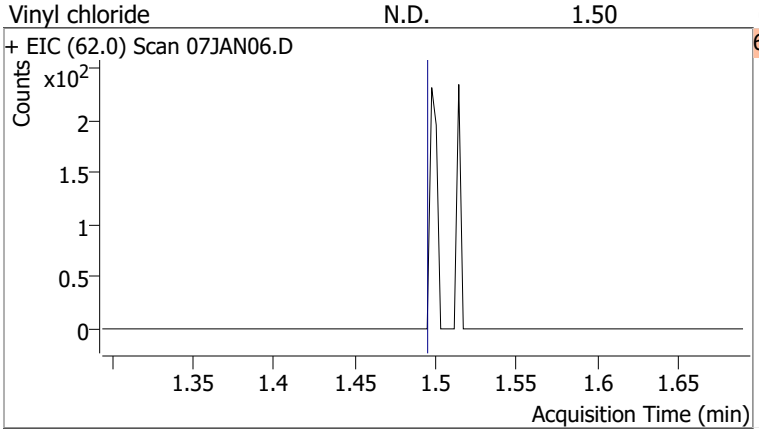


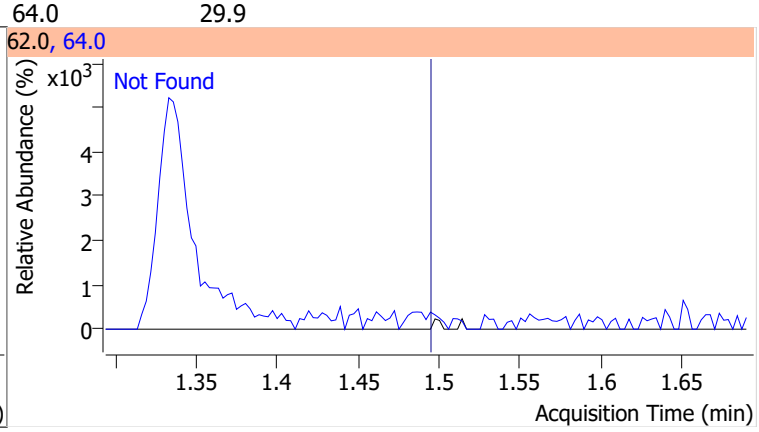
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloromethane	N.D.	1.41	52.0	32.1



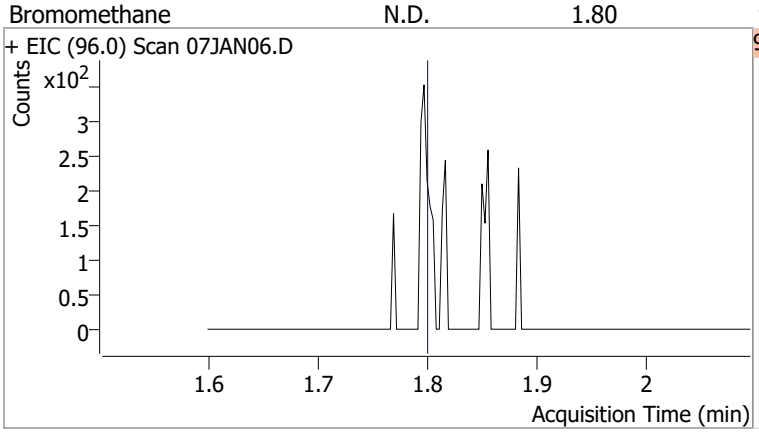


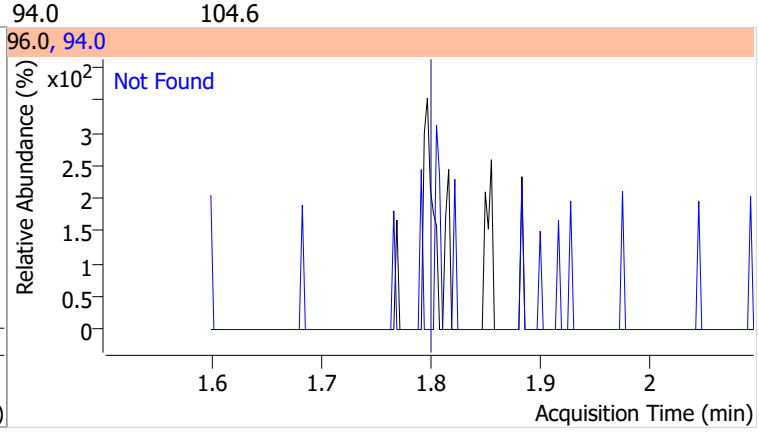
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



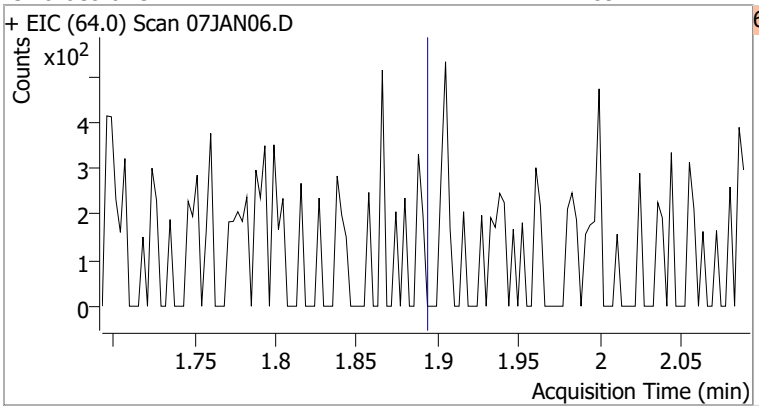


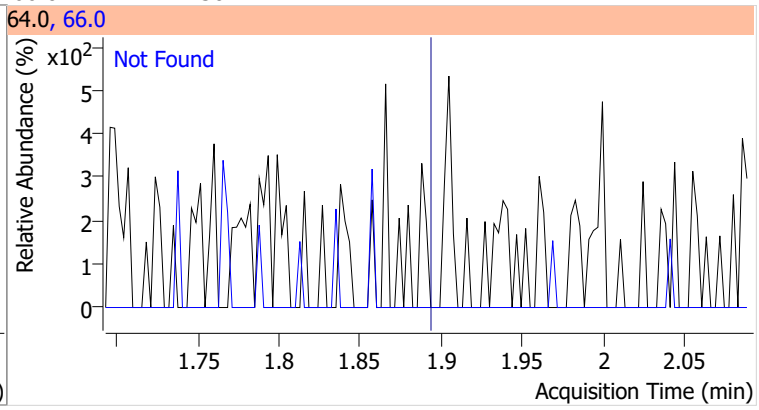
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



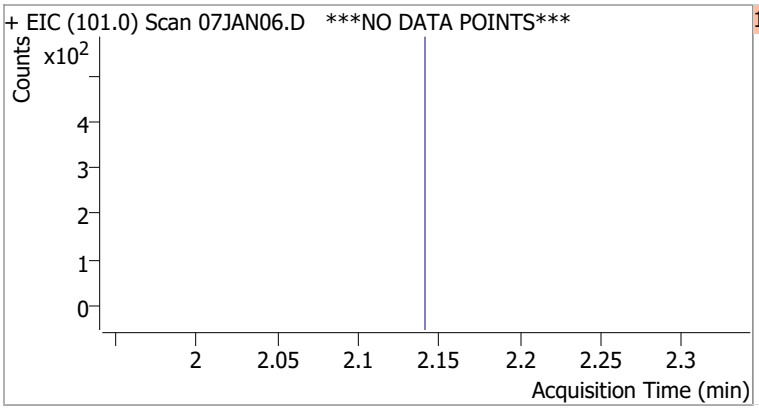


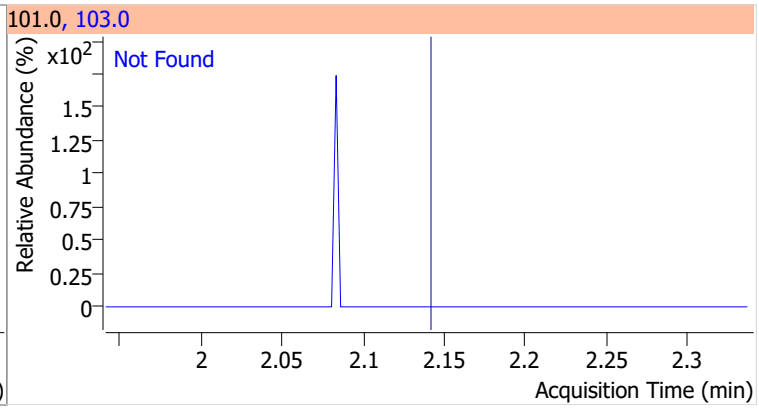
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



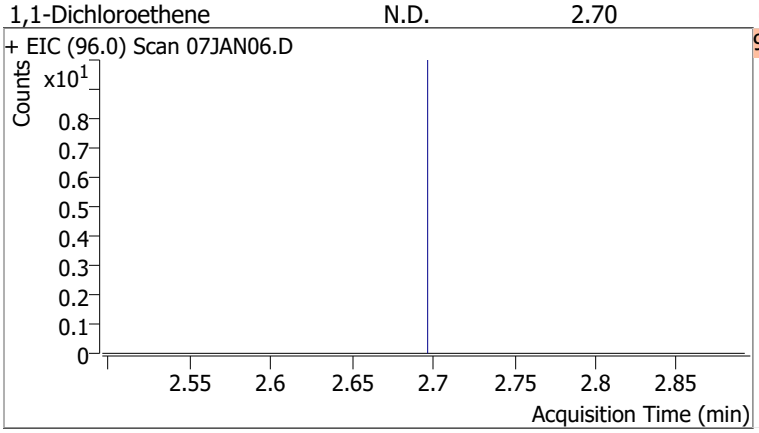


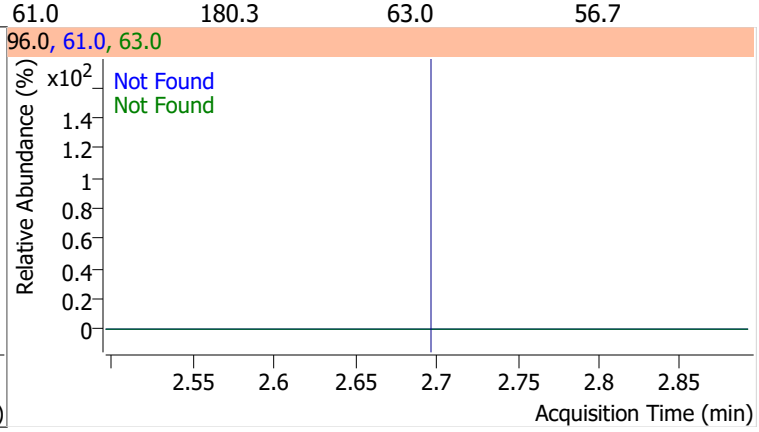
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



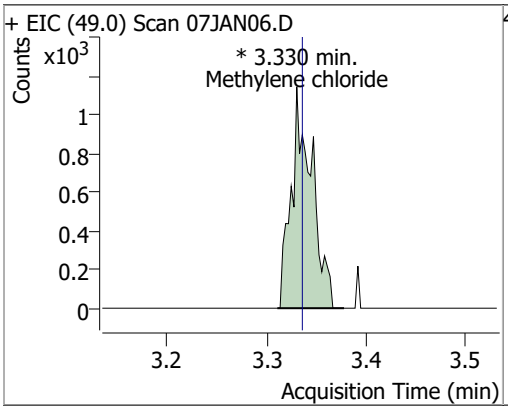


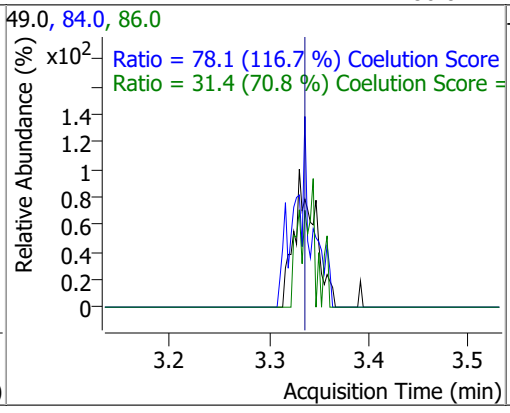
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

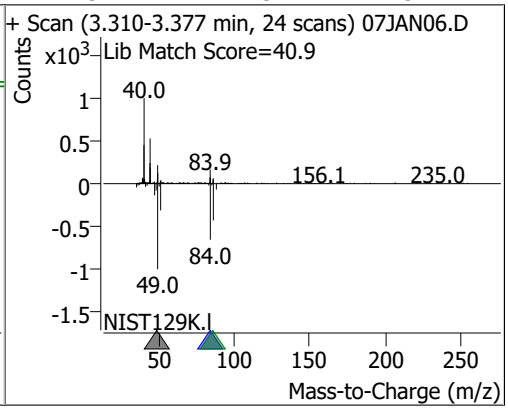




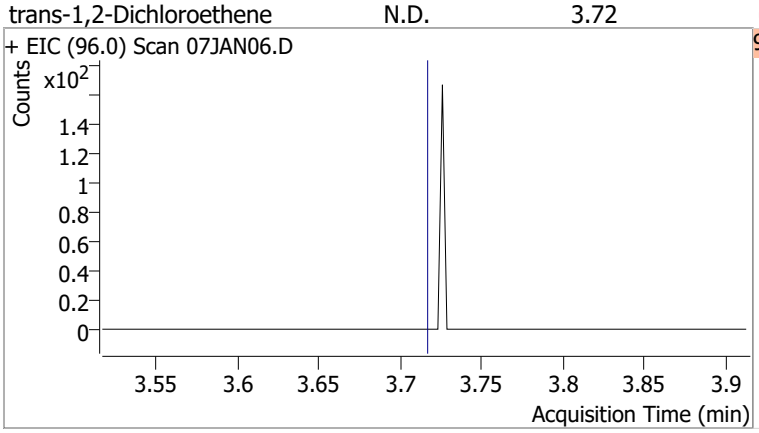
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	1.2889	3.33	-0.01	1664 (m)	84.0	78.1	36.9	96.9
					86.0	31.4	14.3	74.3

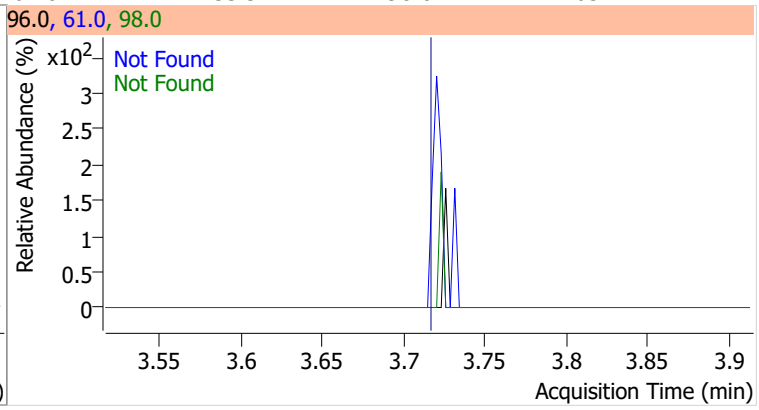




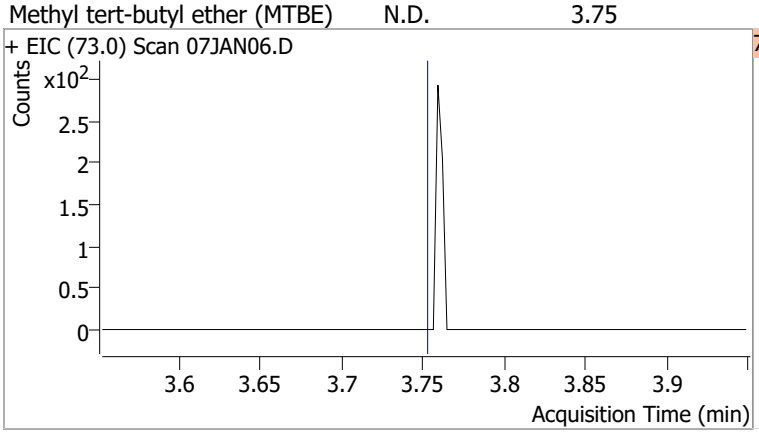


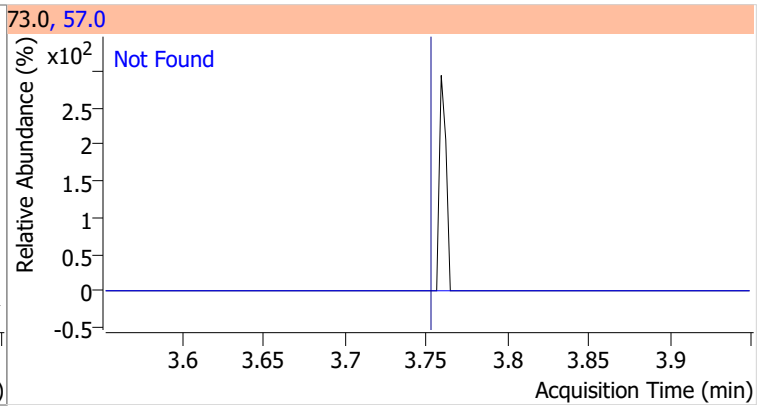
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



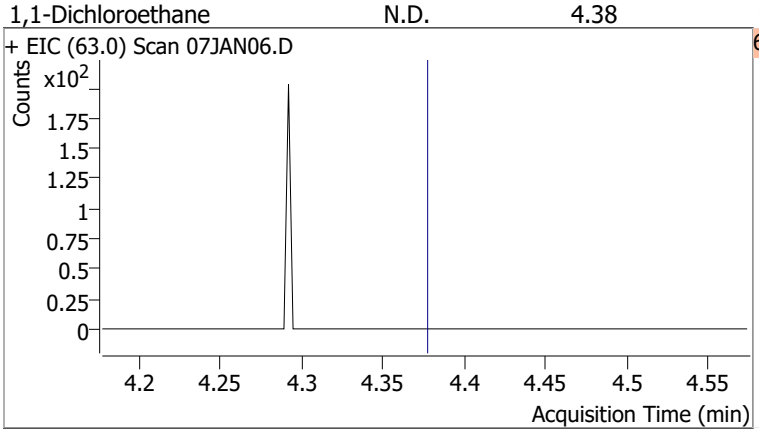


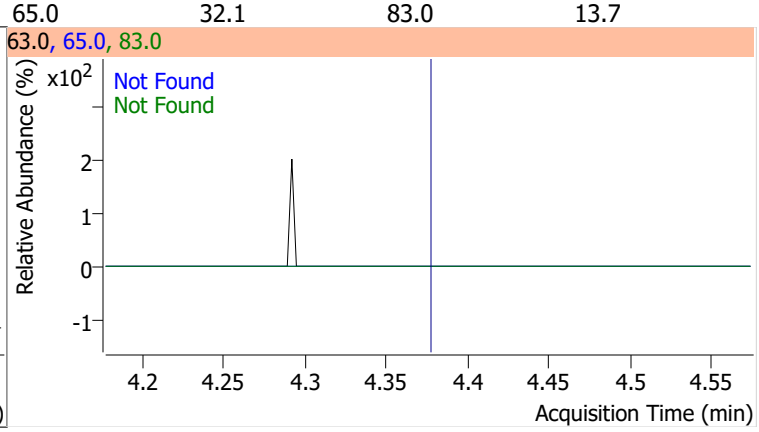
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



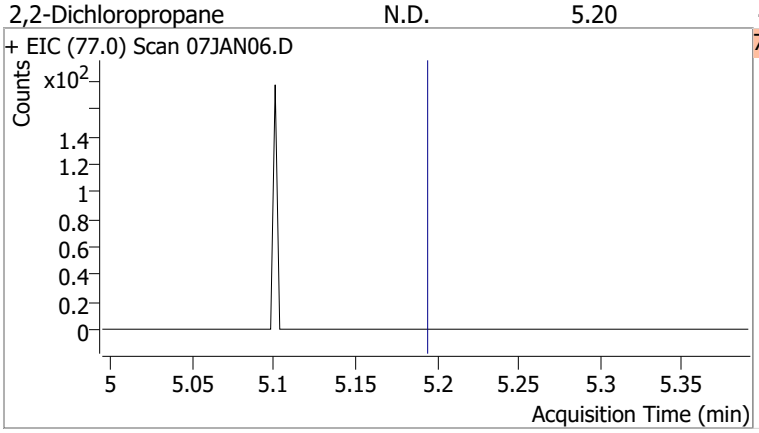


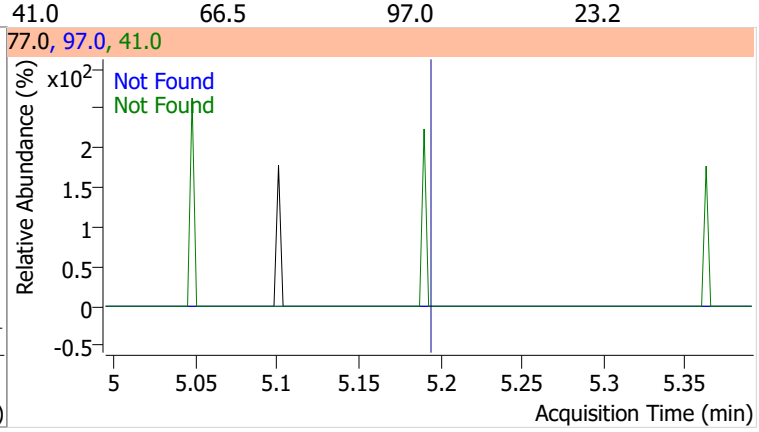
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



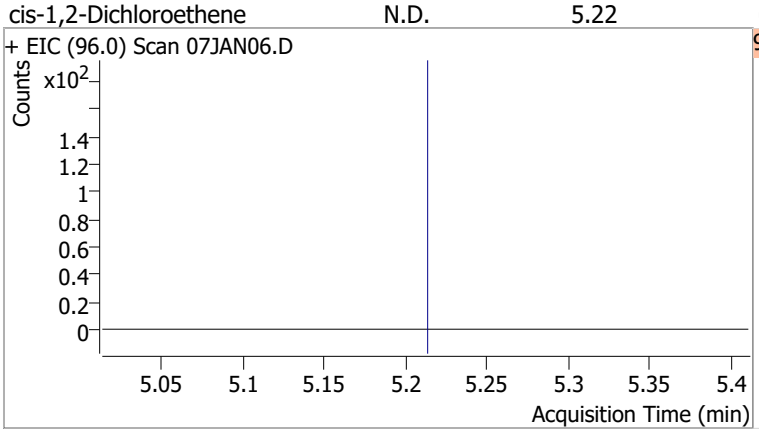


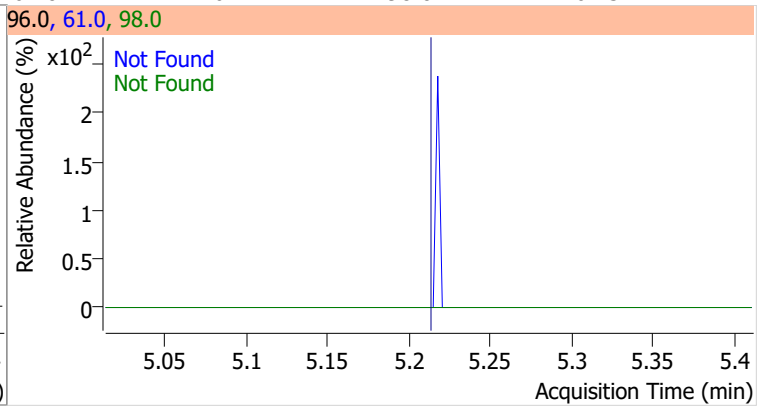
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



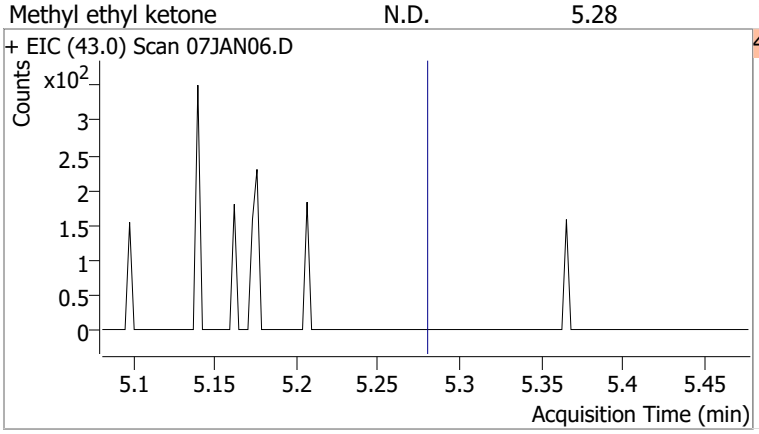


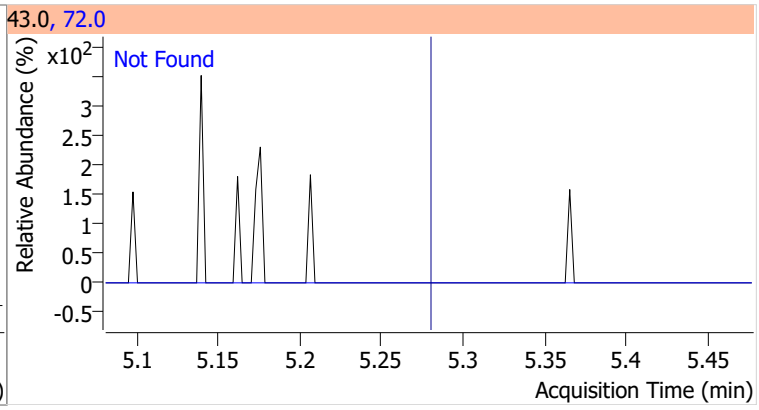
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



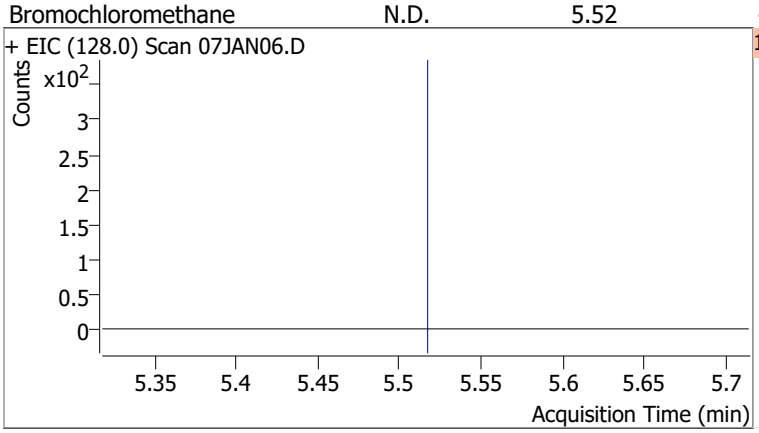


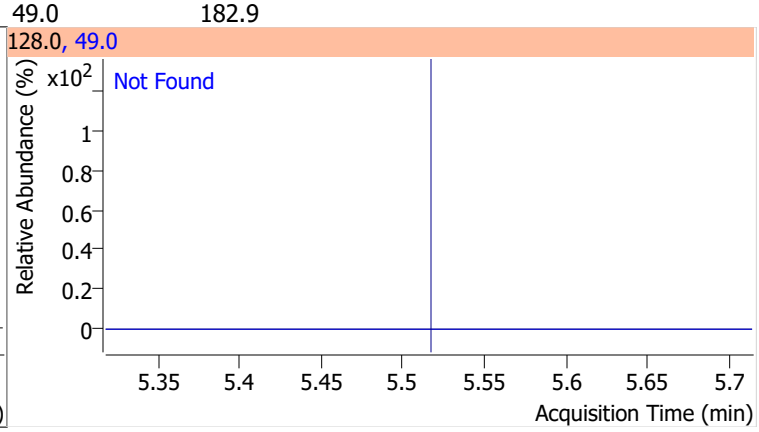
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



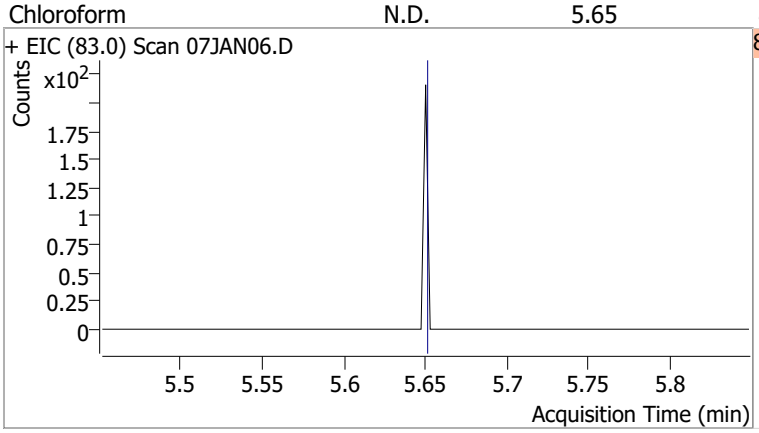


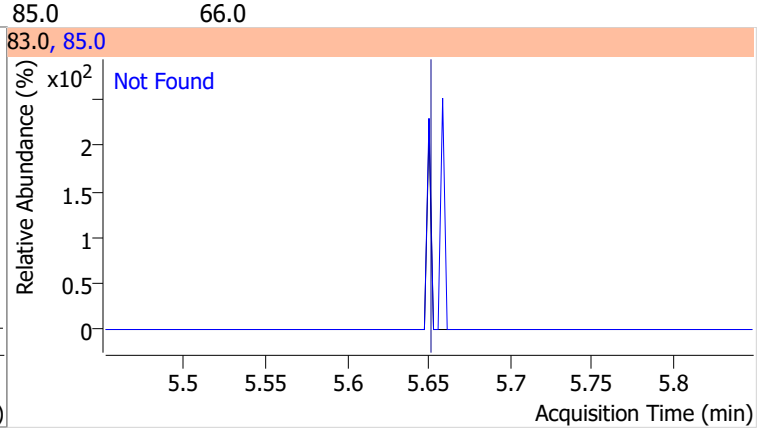
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9



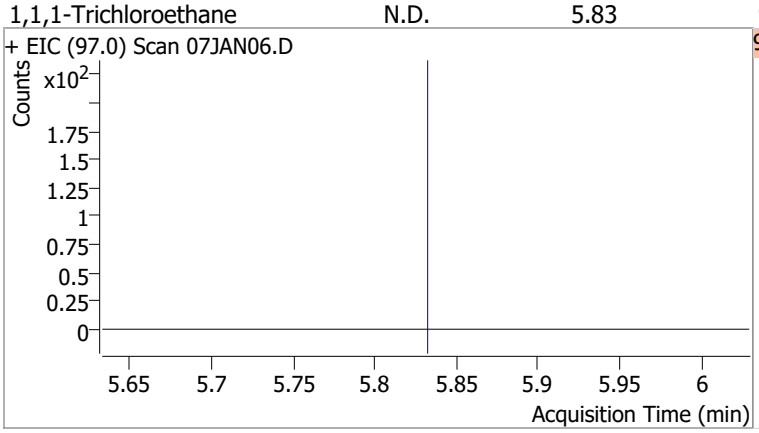


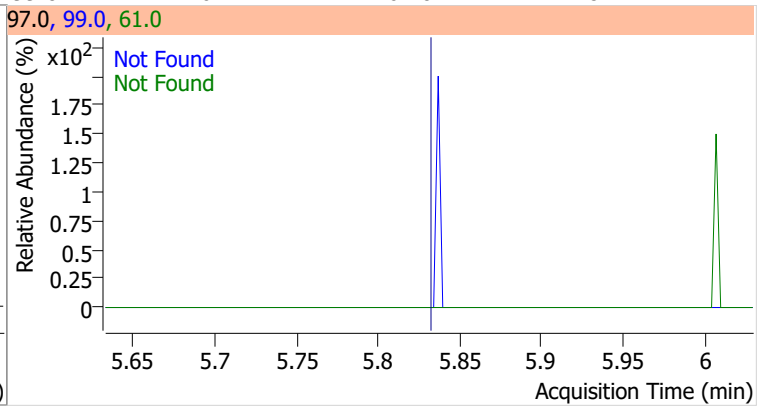
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroform	N.D.	5.65	85.0	66.0



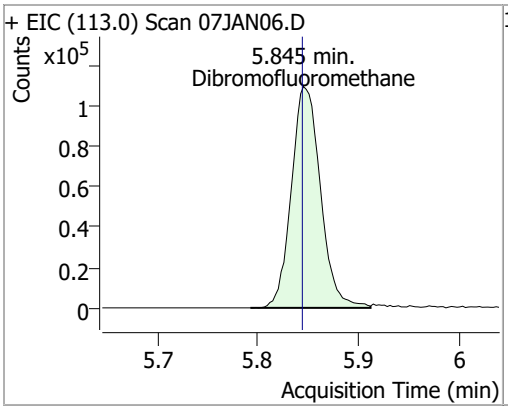


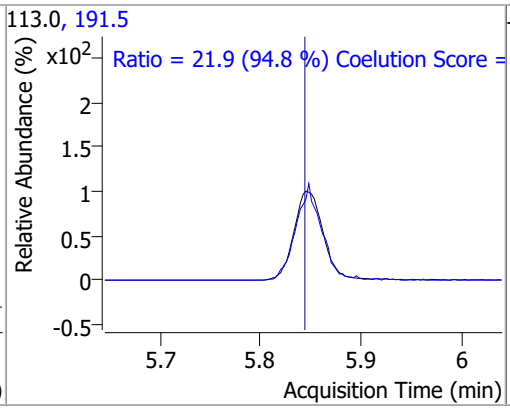
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

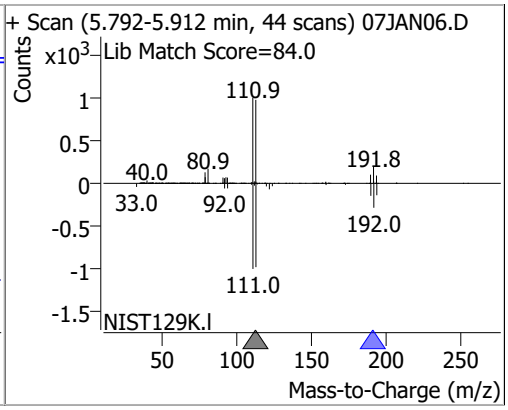




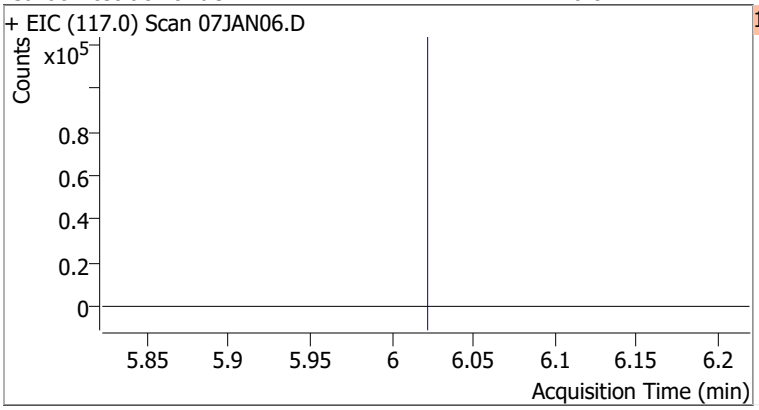
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	270.7666	5.85	0.00	221741	191.5	21.9	0.0	53.1

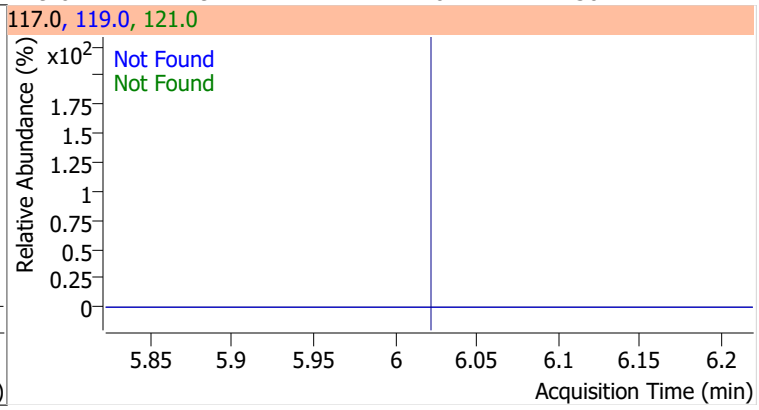




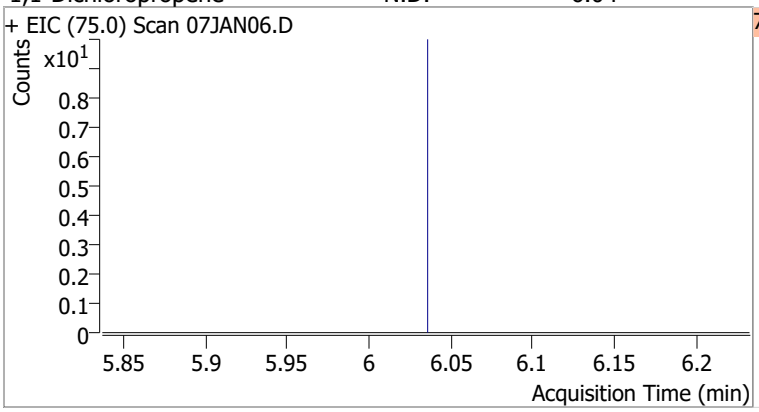


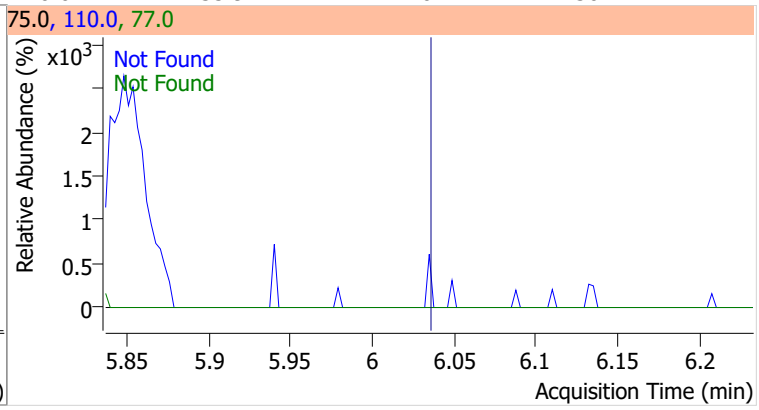
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



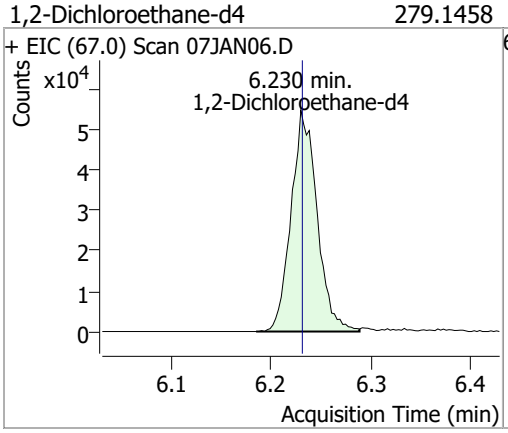


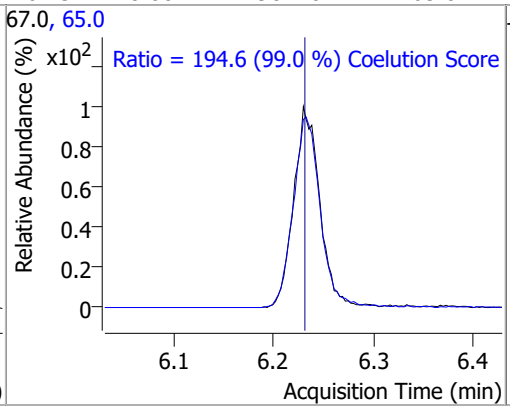
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

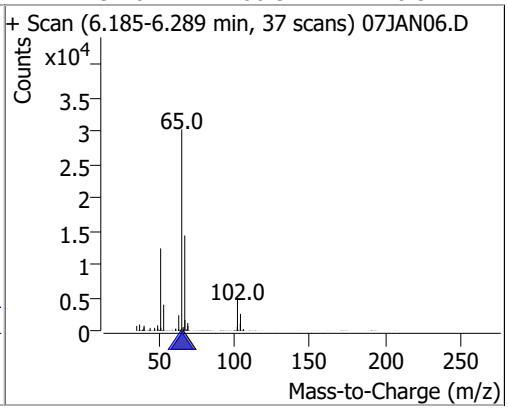




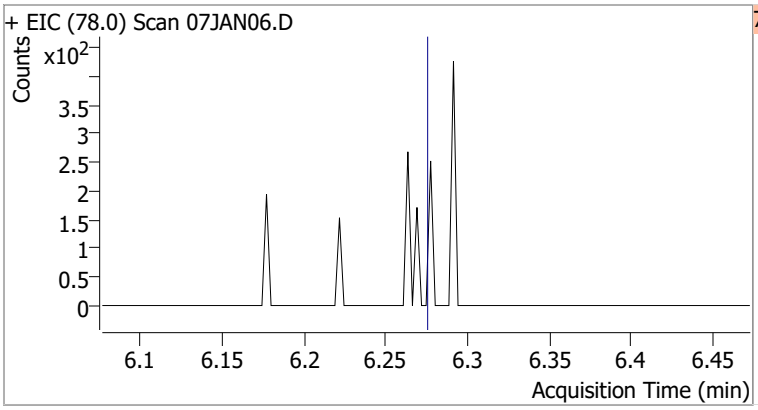
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	279.1458	6.23	0.00	98740	65.0	194.6	166.5	226.5

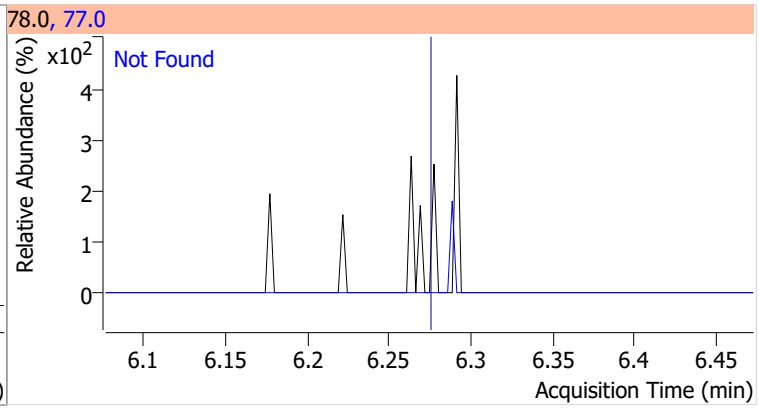




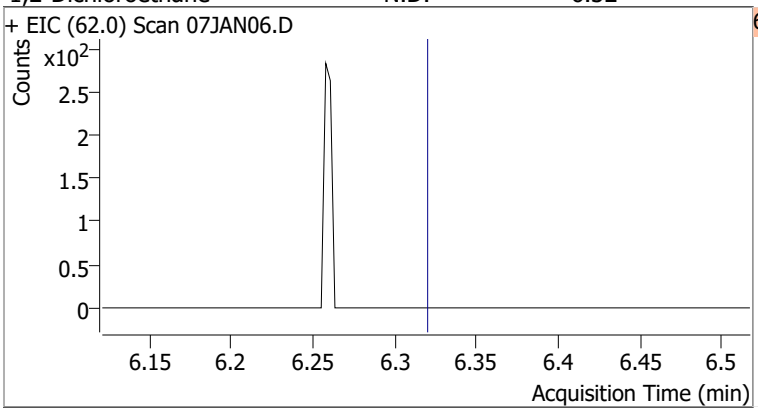


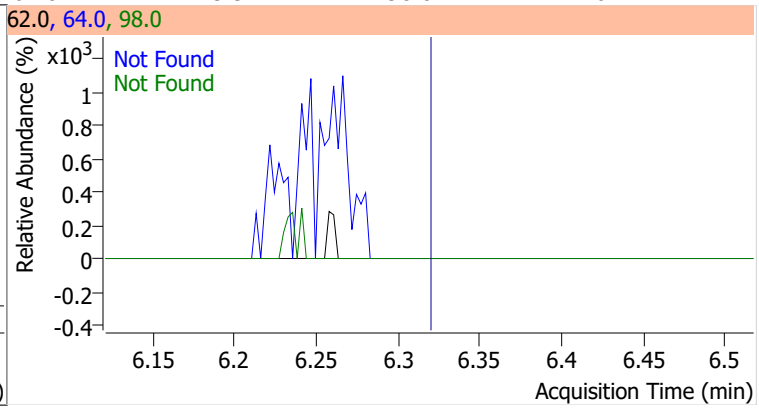
Compound	Conc.	Exp RT	QIon	Exp Ratio
Benzene	N.D.	6.28	77.0	23.5



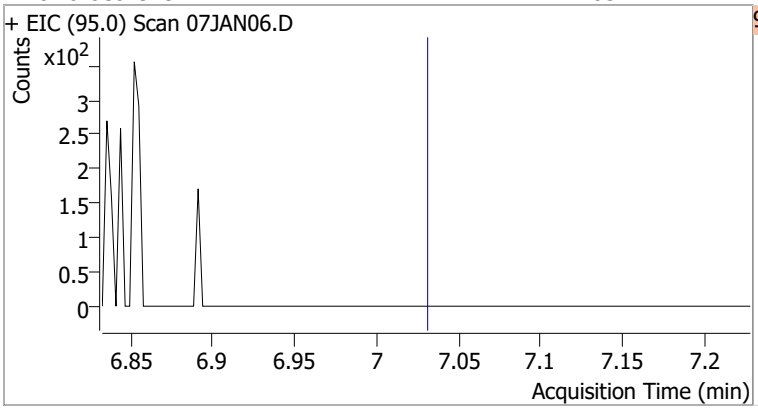


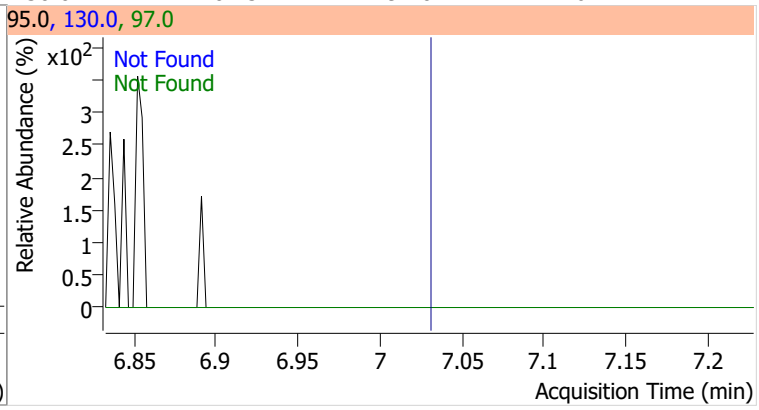
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



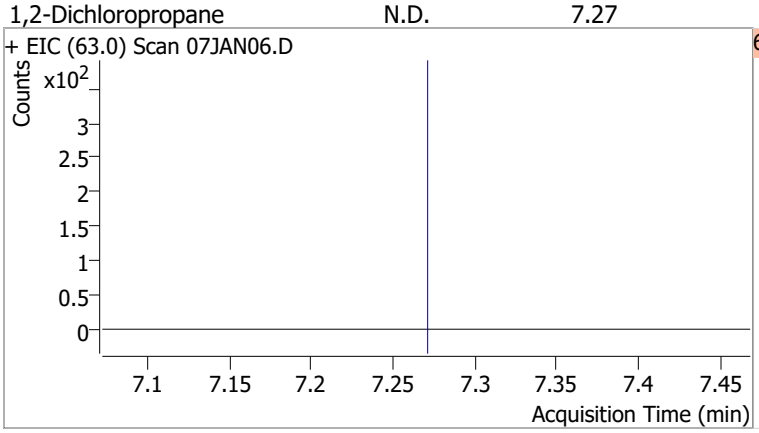


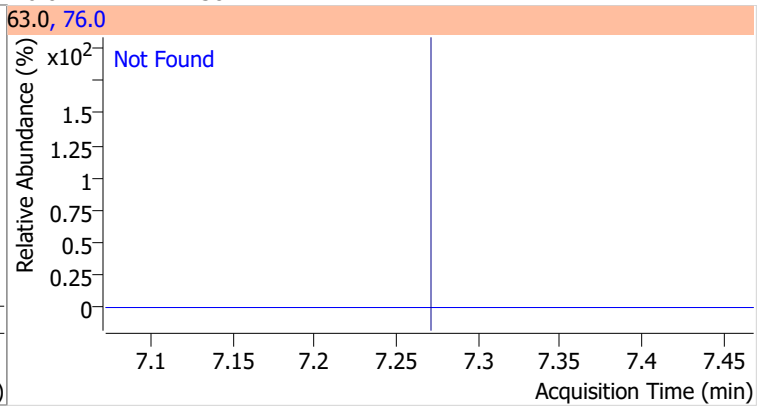
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



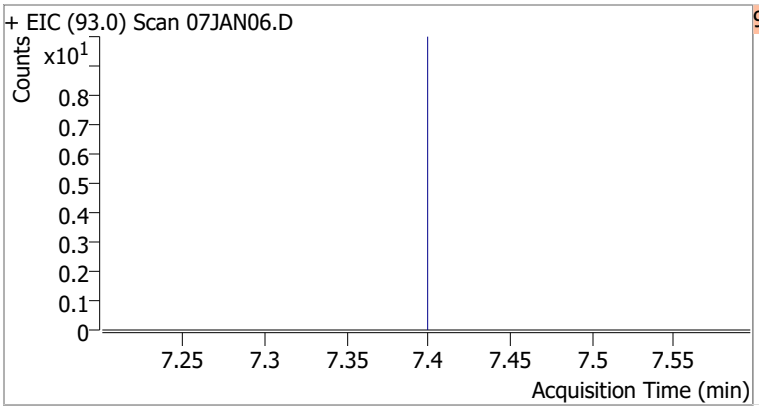


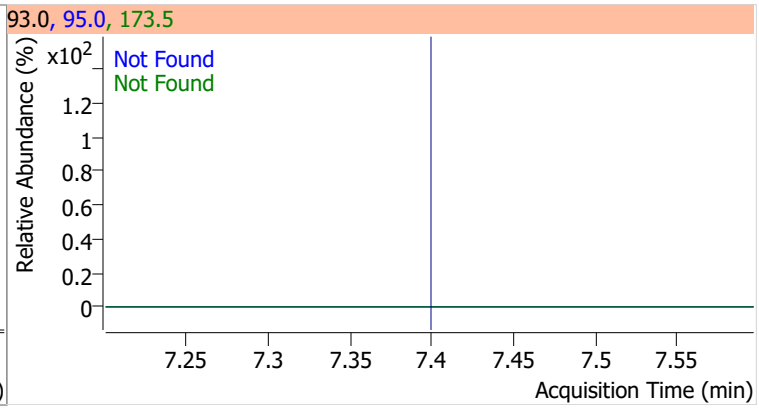
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



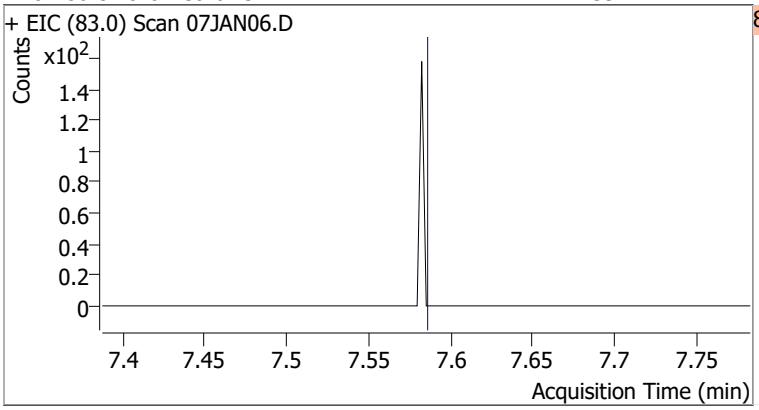


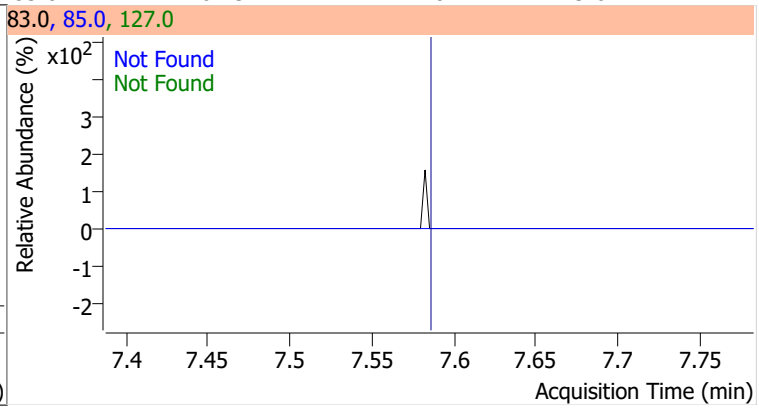
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



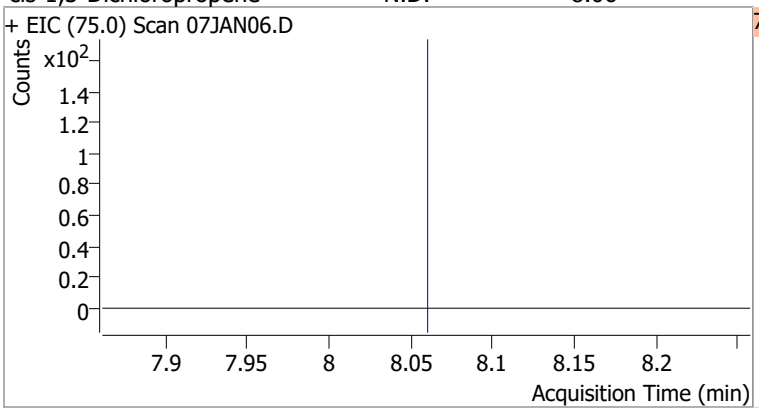


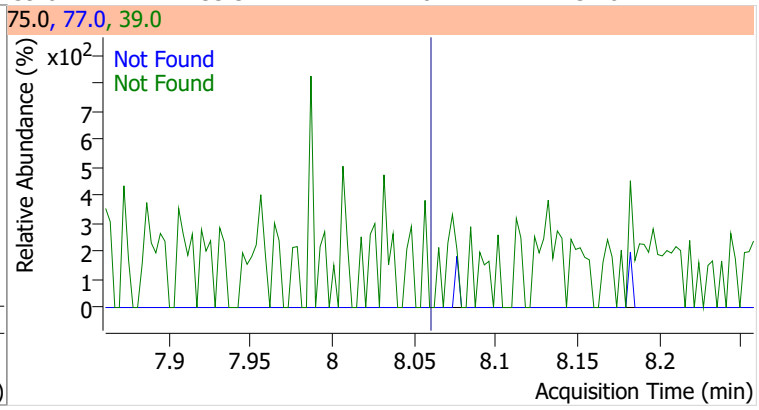
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



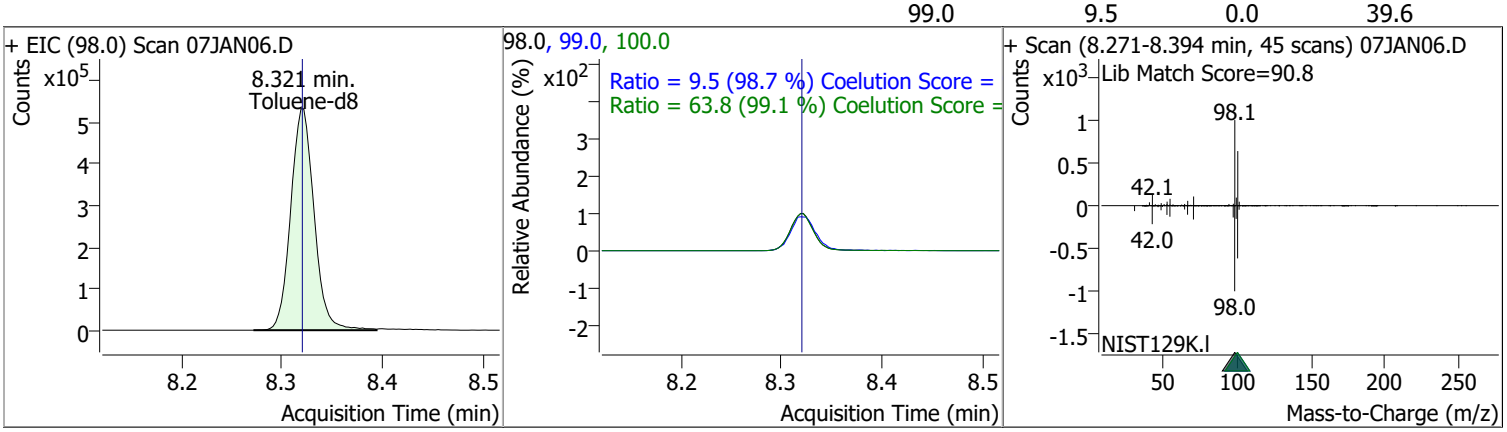


Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

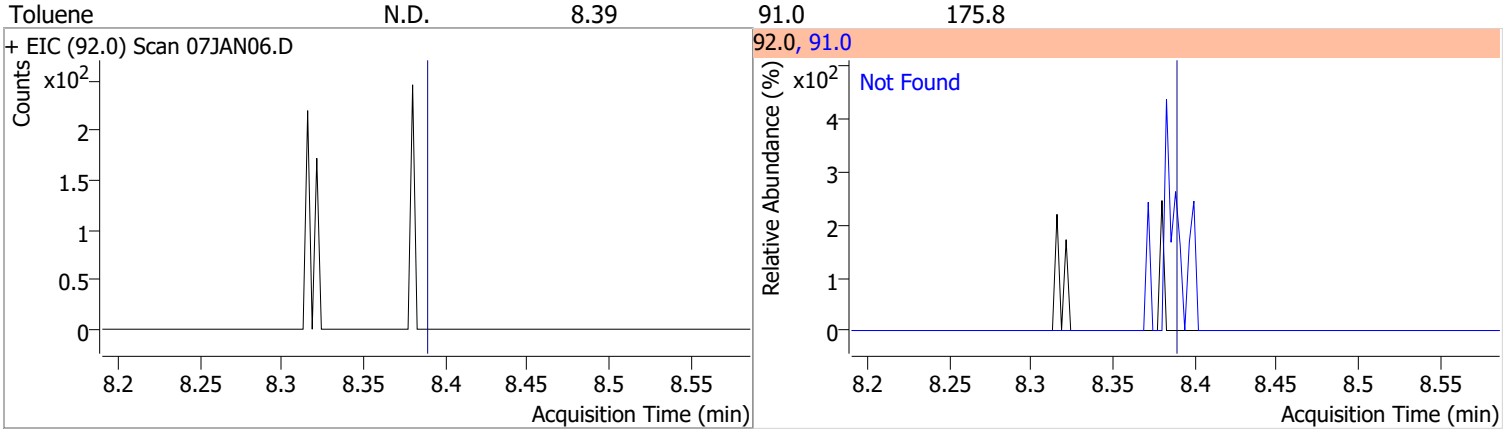




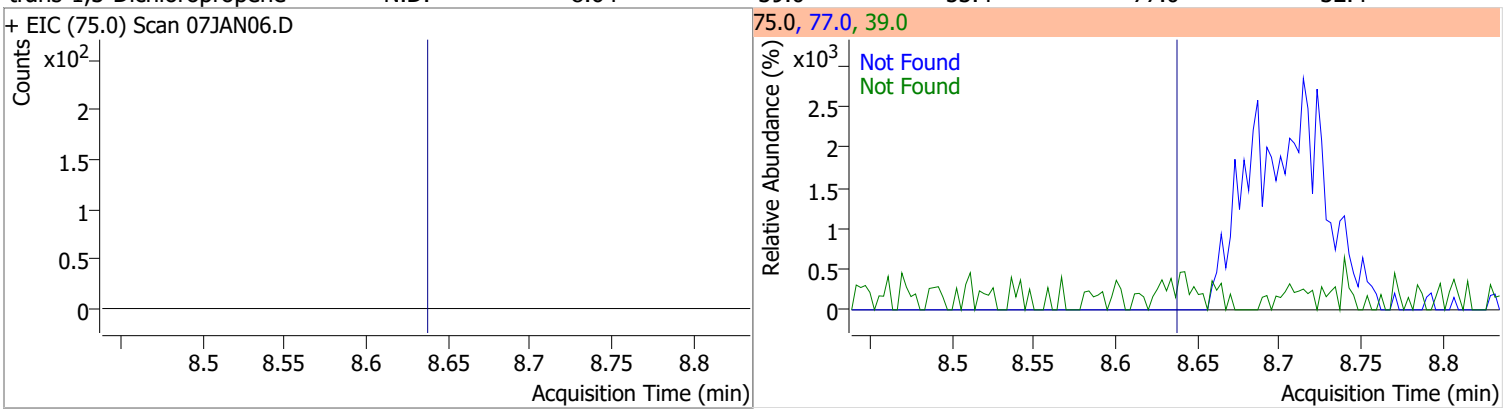
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	264.3775	8.32	0.00	864903	100.0	63.8	34.4	94.4
					99.0	9.5	0.0	39.6



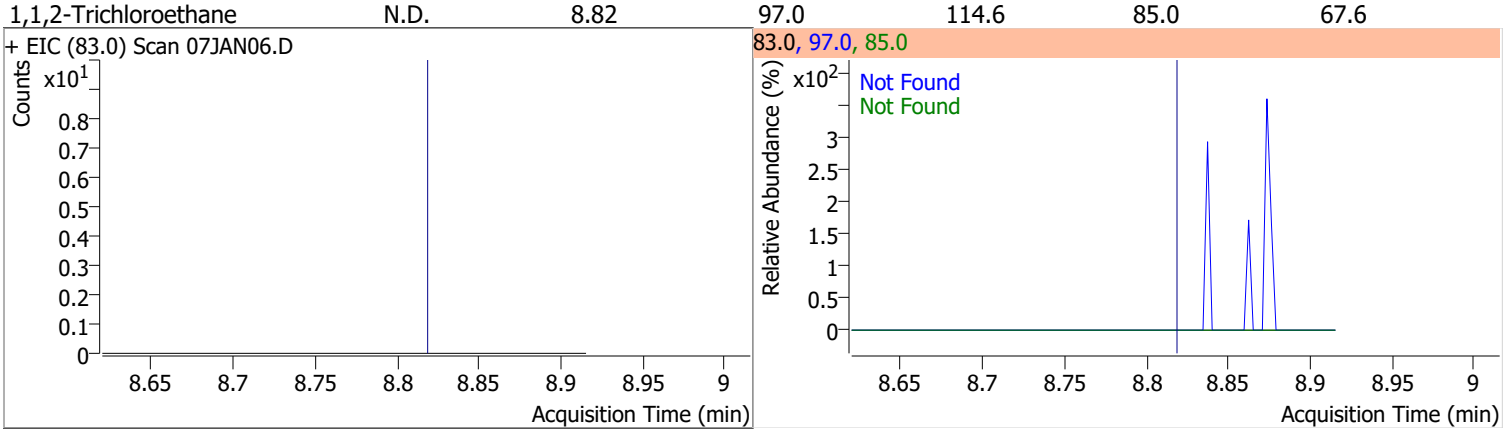
Compound	Conc.	Exp RT	QIon	Exp Ratio
Toluene	N.D.	8.39	91.0	175.8



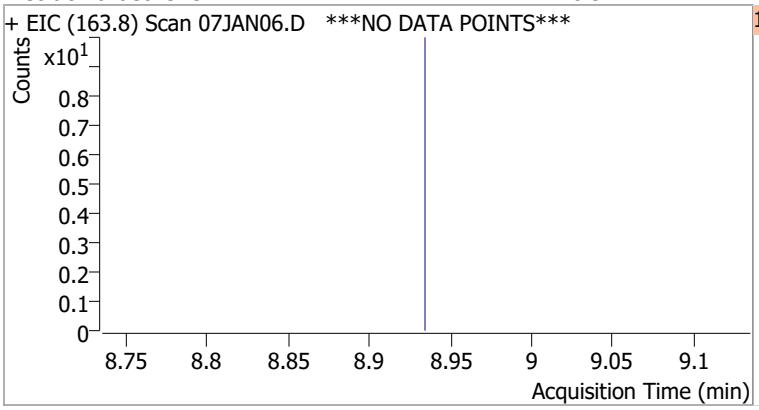
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4

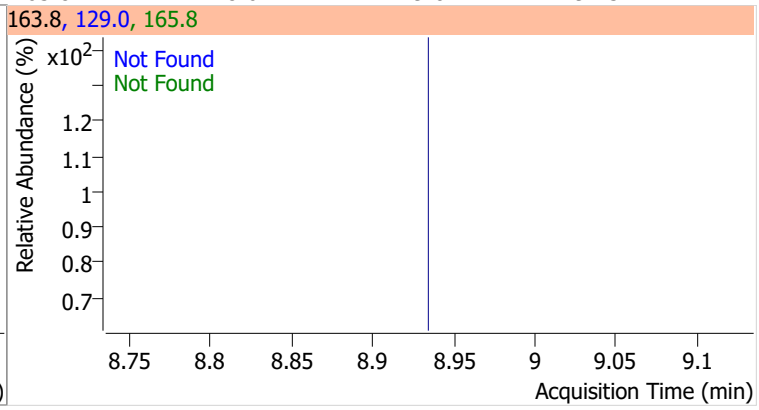


Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6

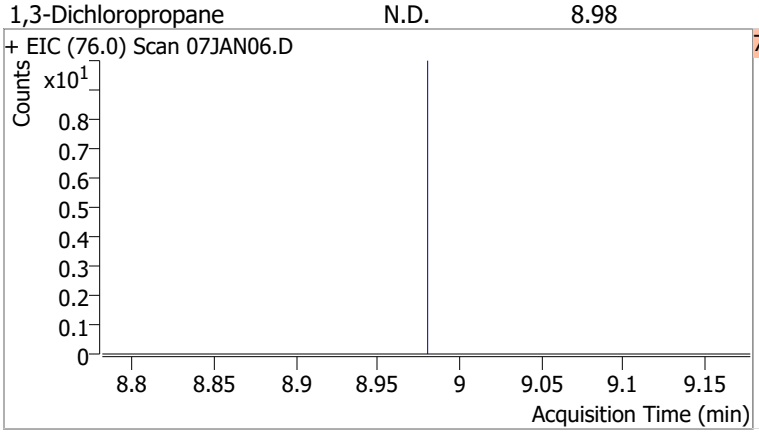


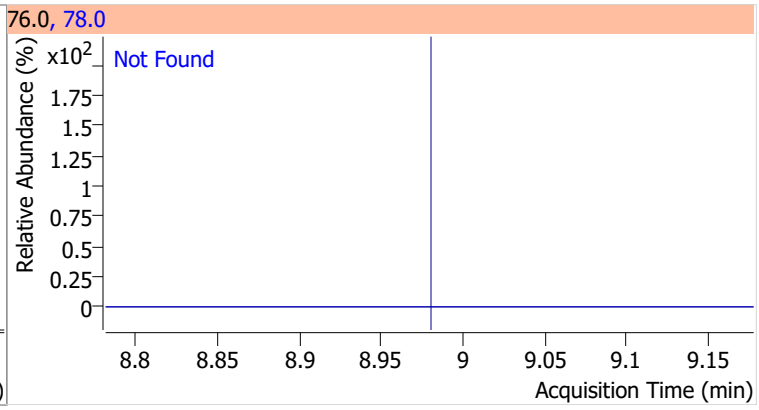
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



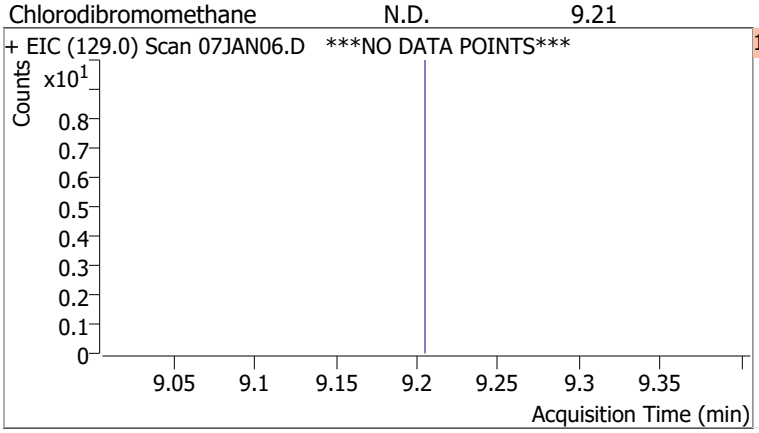


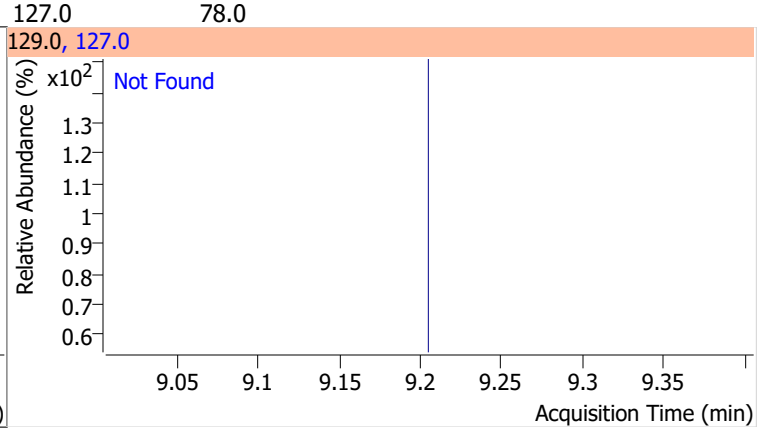
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



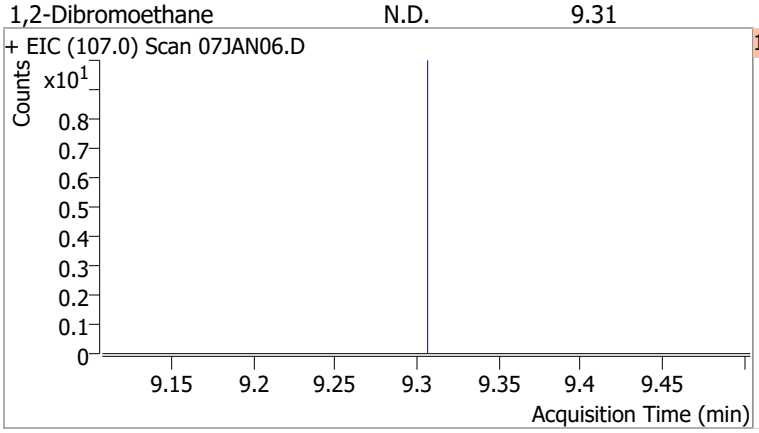


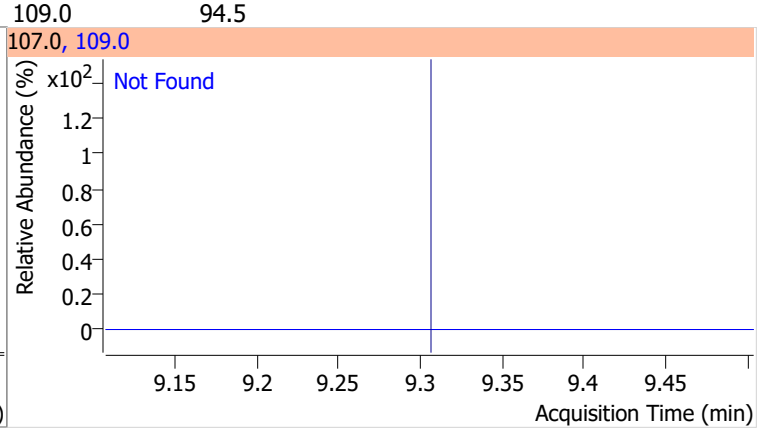
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



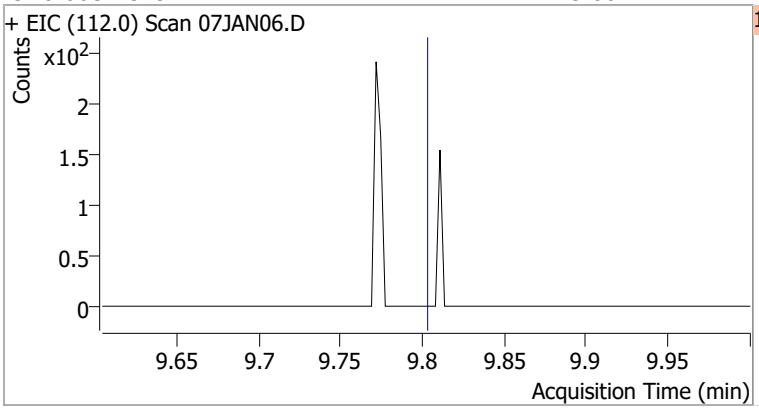


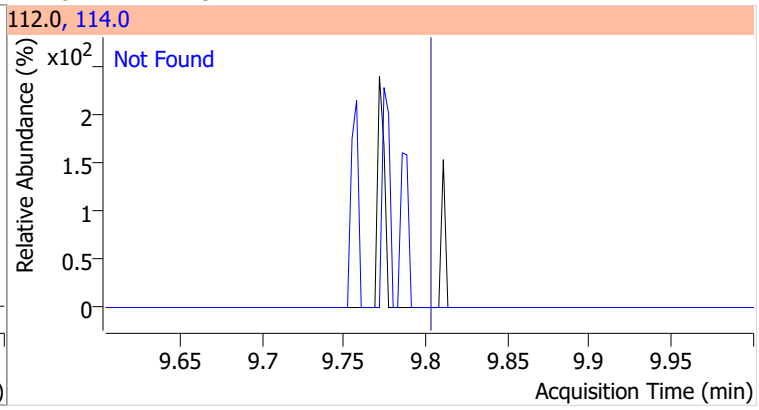
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



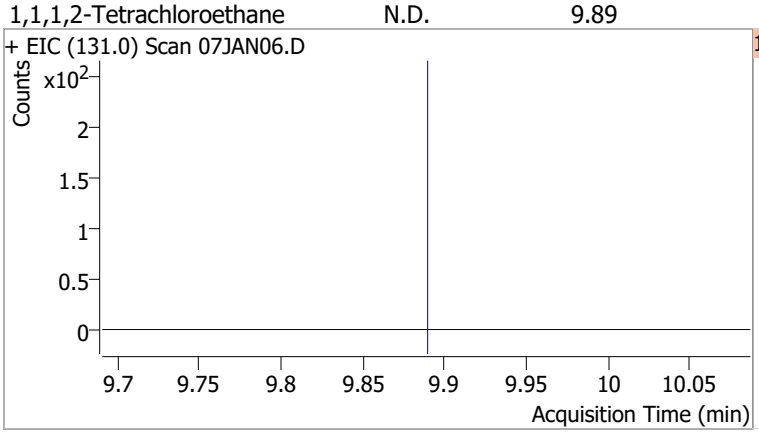


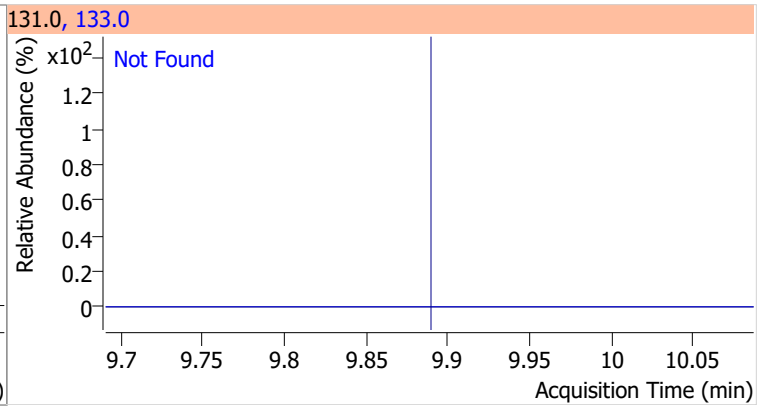
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



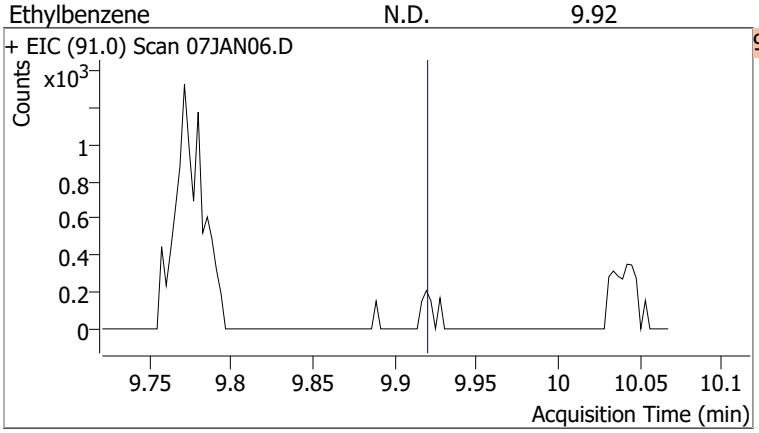


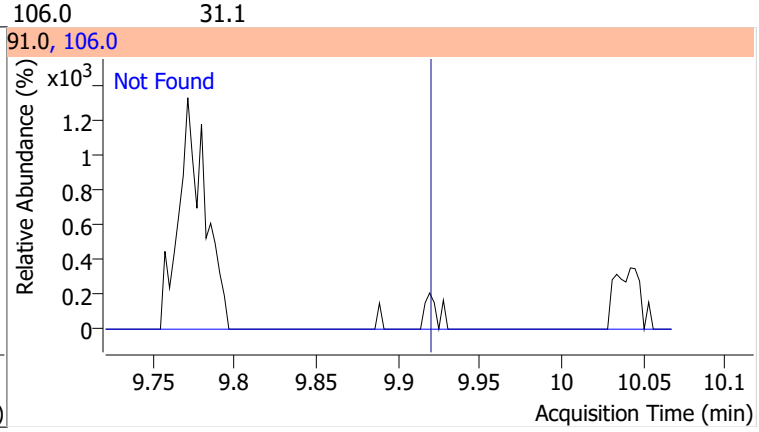
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



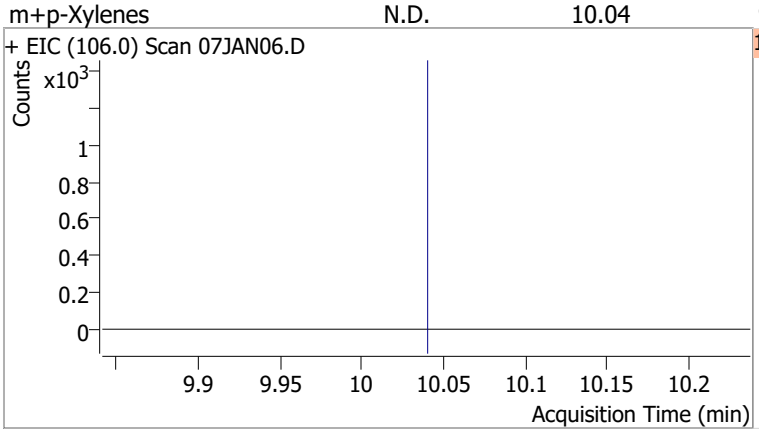


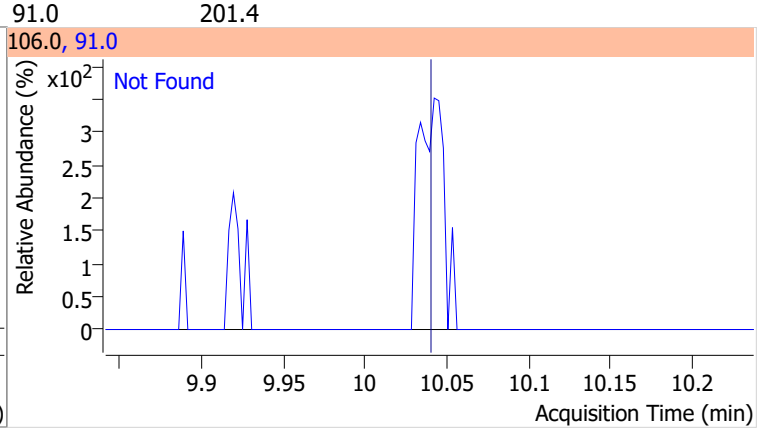
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1



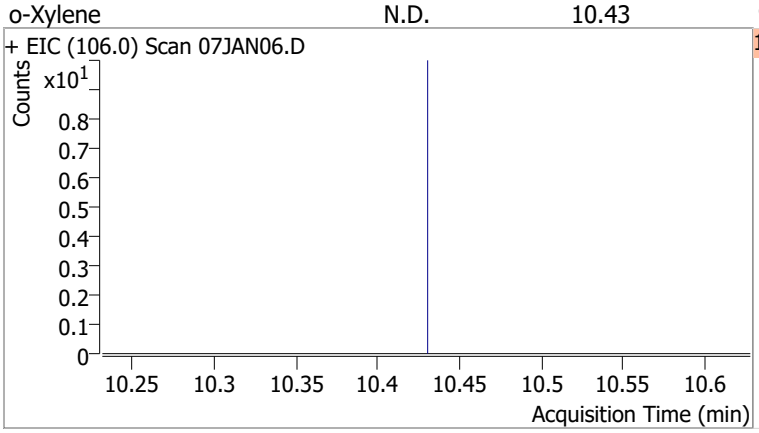


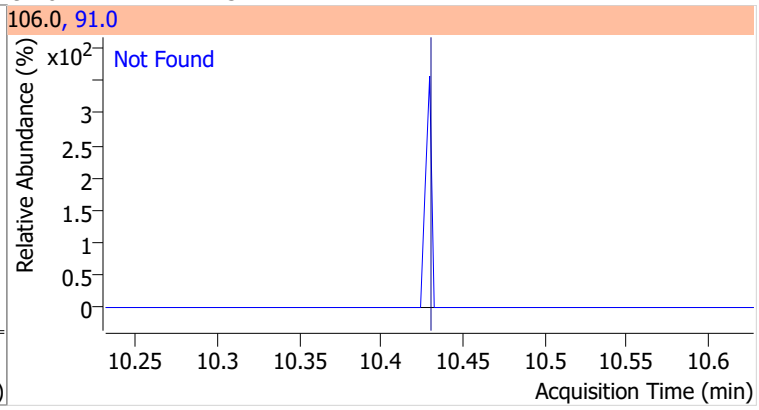
Compound	Conc.	Exp RT	QIon	Exp Ratio
m+p-Xylenes	N.D.	10.04	91.0	201.4



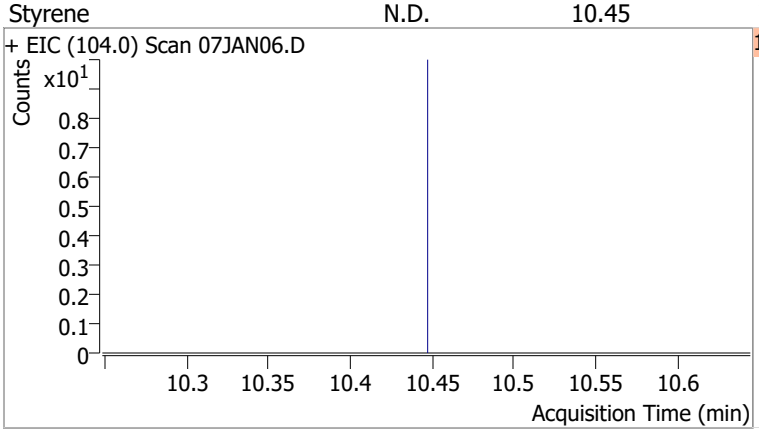


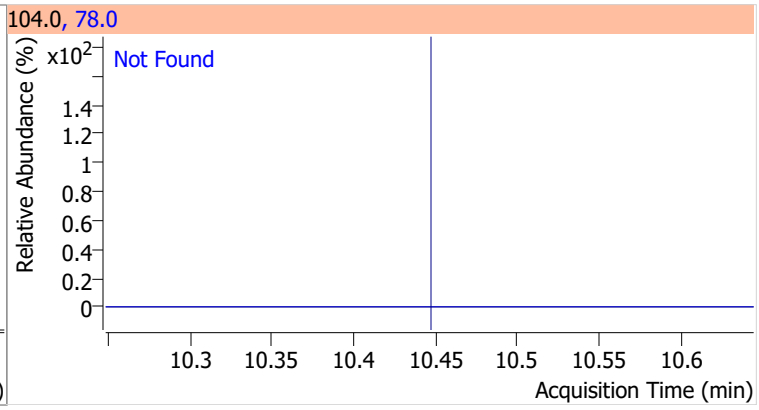
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



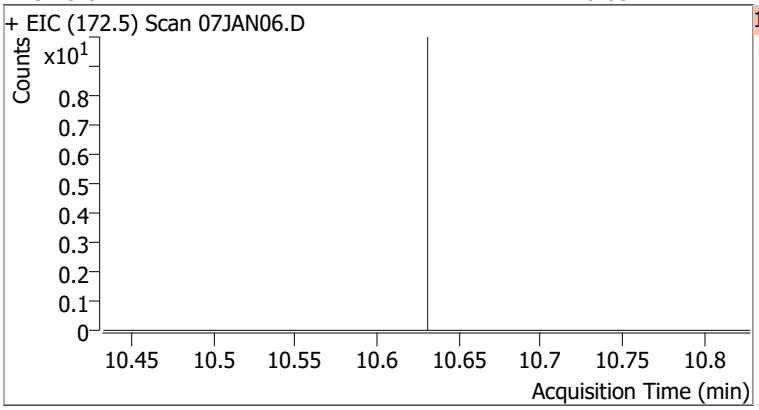


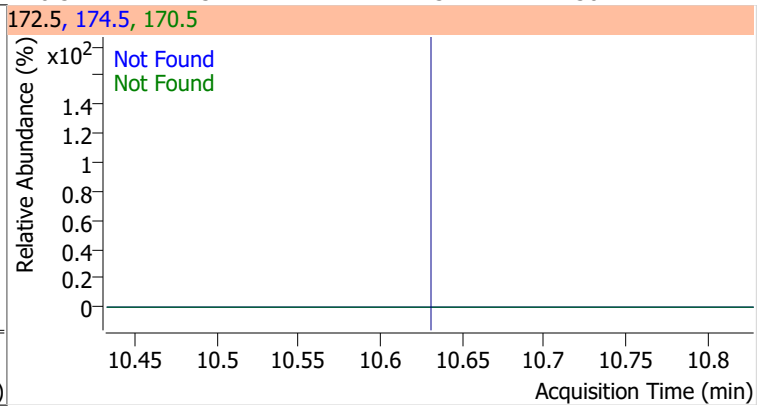
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



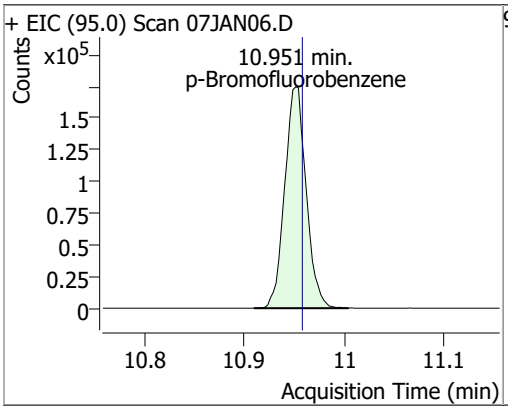


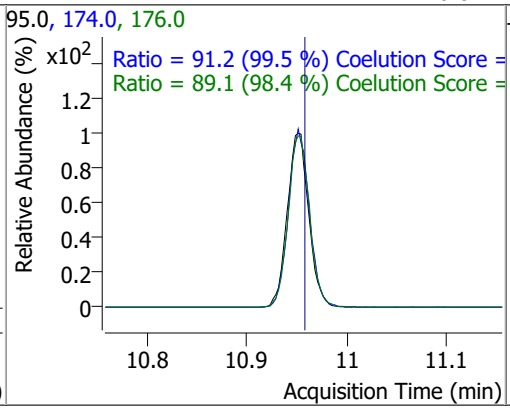
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

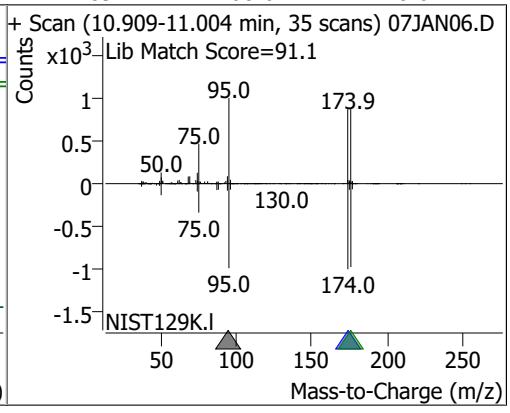




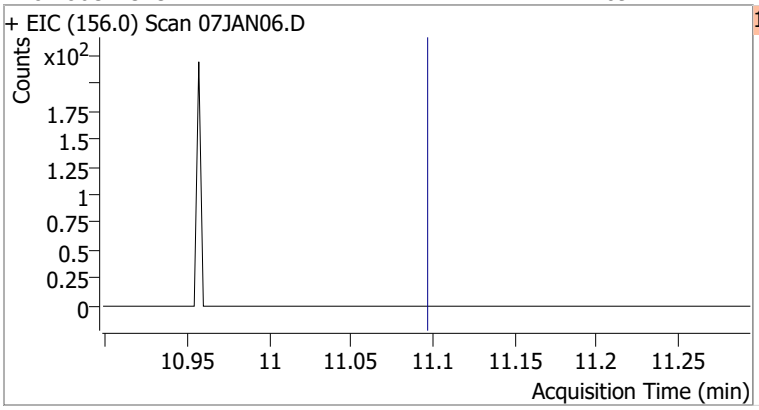
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	274.9894	10.95	0.00	259344	174.0	91.2	61.7	121.7
					176.0	89.1	60.6	120.6

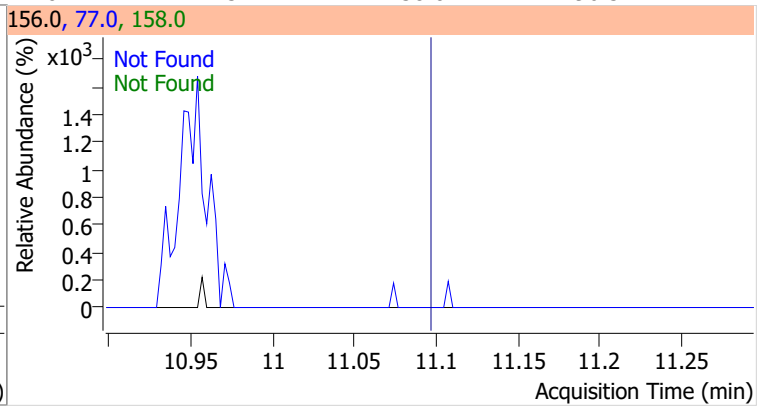




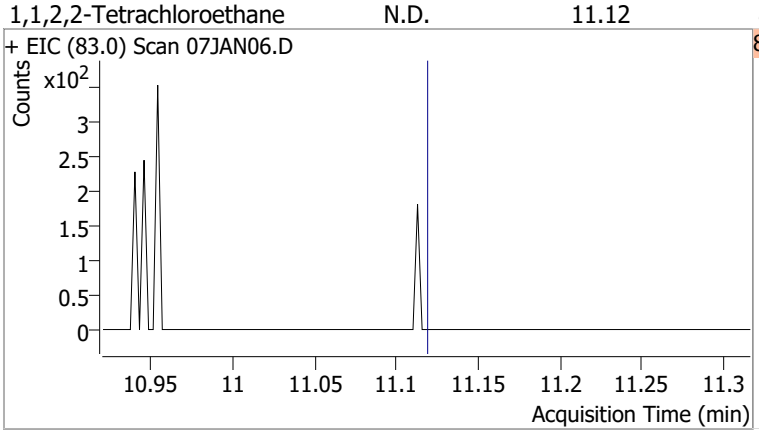


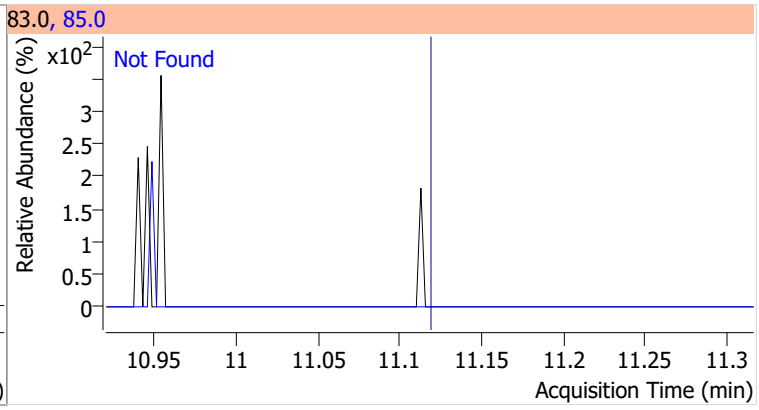
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



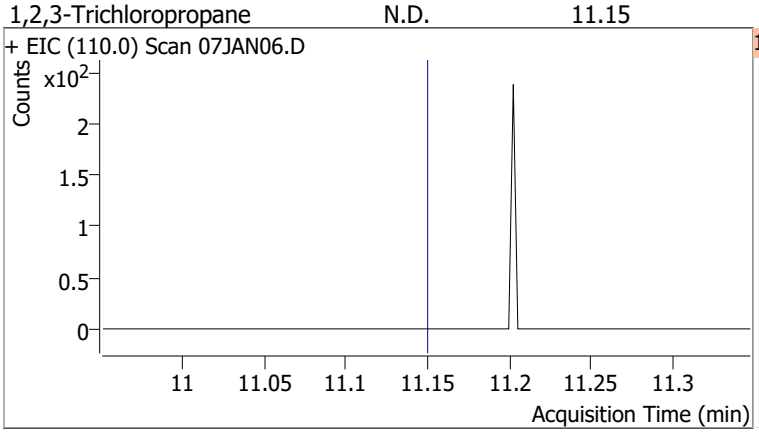


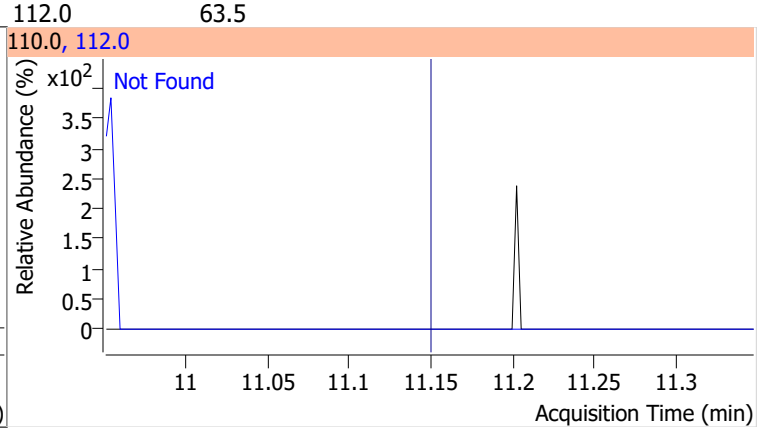
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



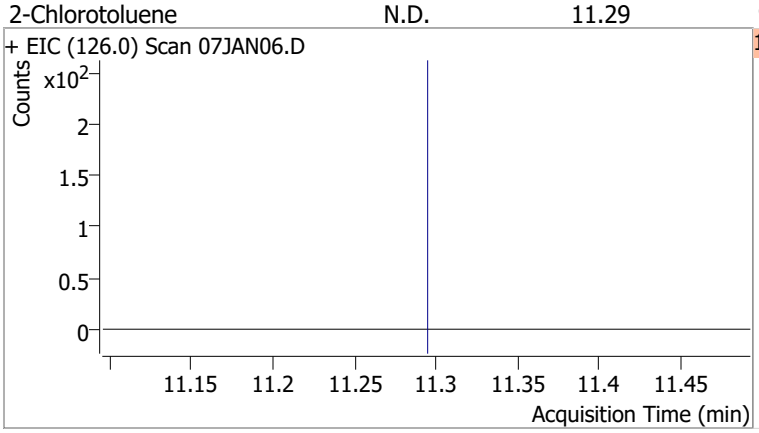


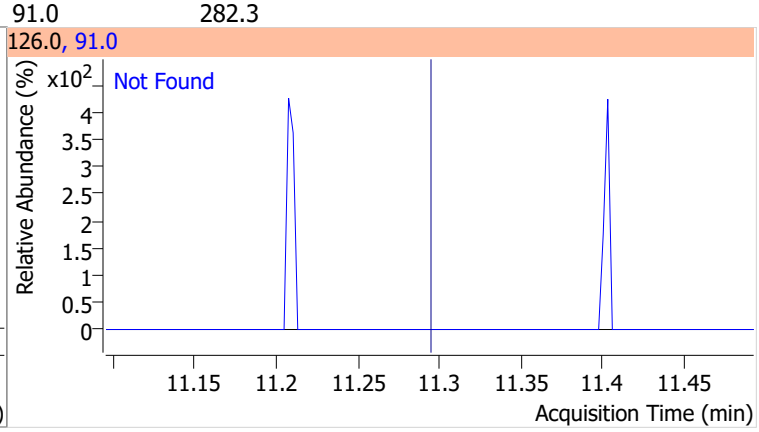
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



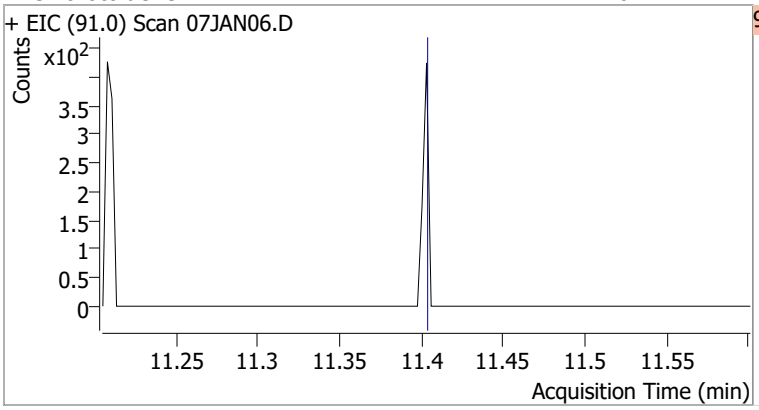


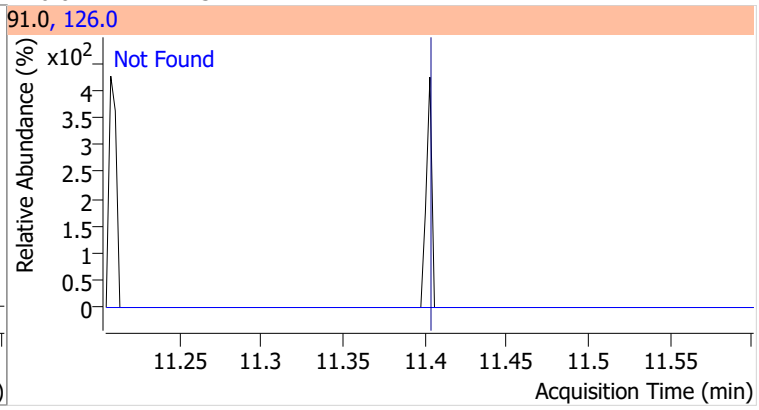
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



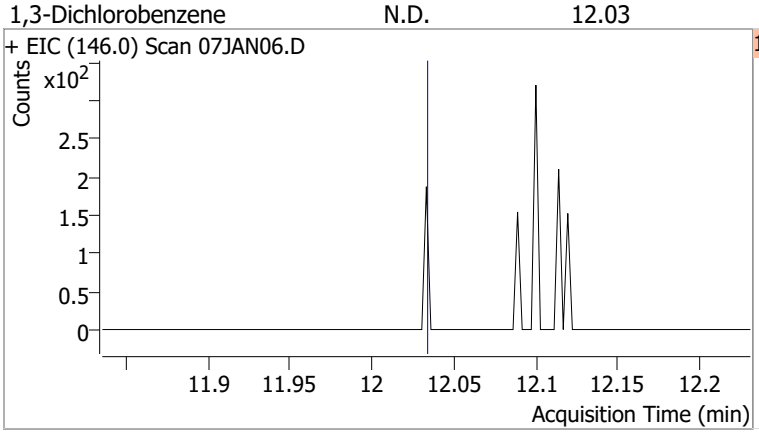


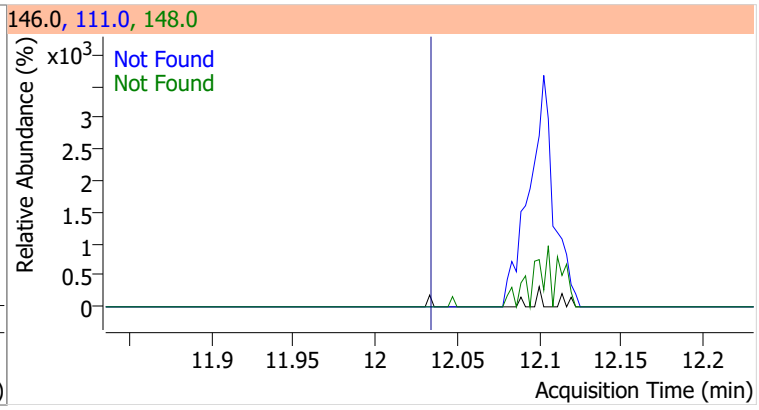
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



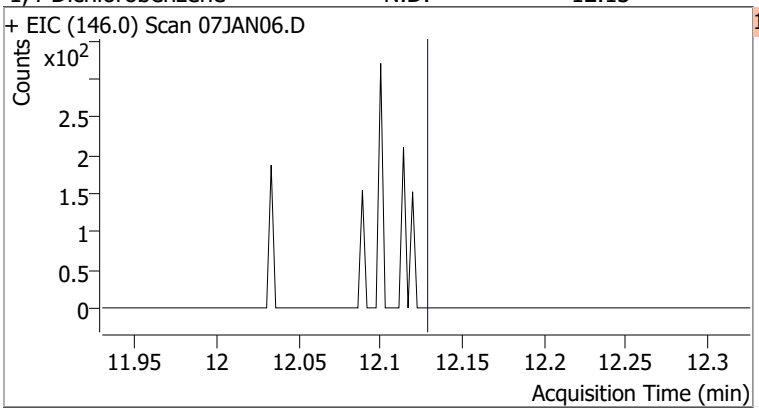


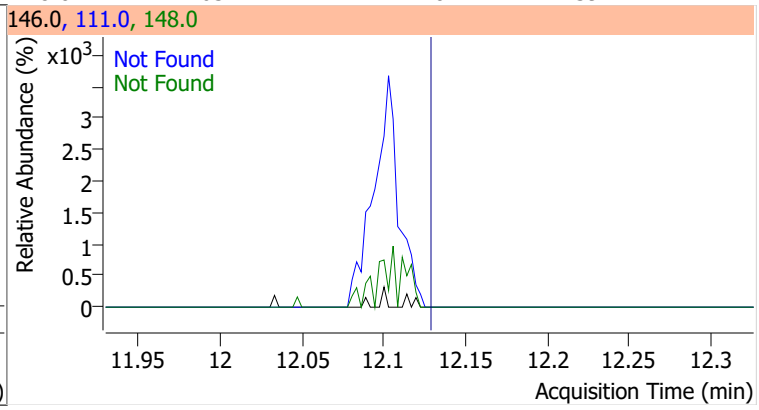
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



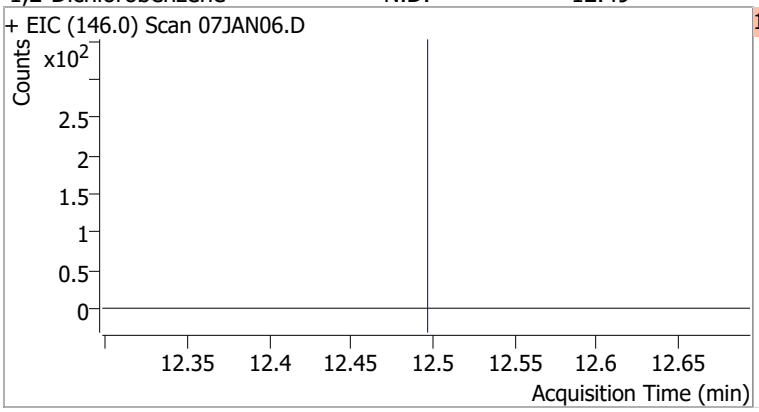


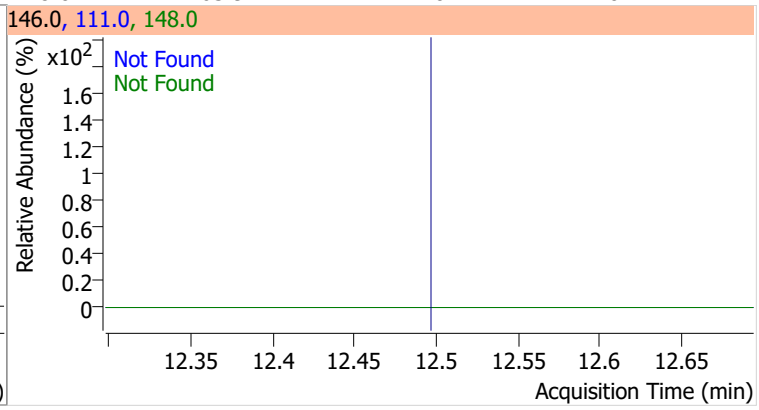
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





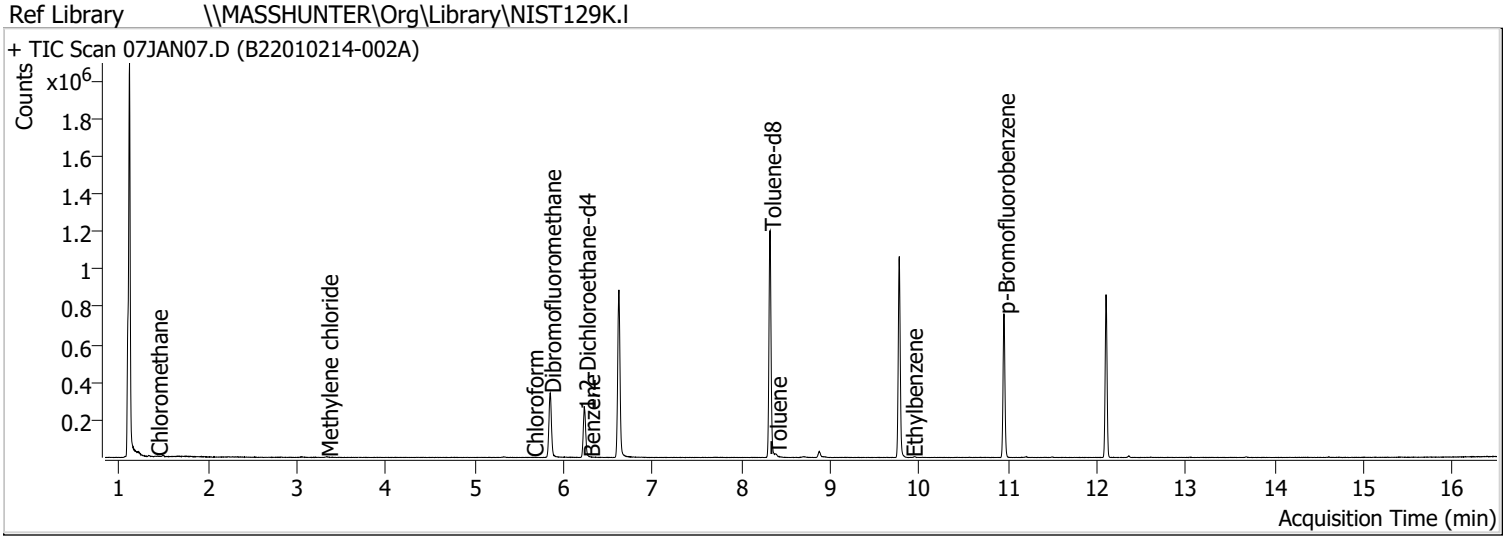
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN07.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 12:24:50 PM
Sample Name	B22010214-002A	Instrument	VOA5975C
Vial	7	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

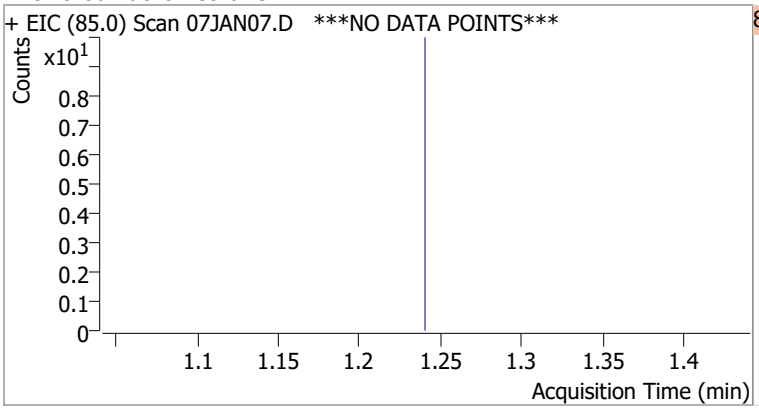


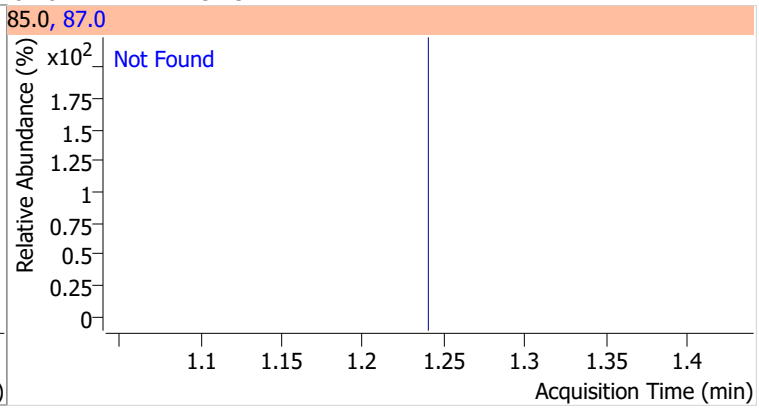
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.620	96.0	753037	250.0000	ng		-0.003
M Chlorobenzene-d5	9.774	82.0	295673	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	207588	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.848	113.0	207577	292.5939	ng		0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 117.04%			
S 1,2-Dichloroethane-d4	6.236	67.0	96449	314.7549	ng		0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 125.90%		*	
S Toluene-d8	8.321	98.0	740894	260.0307	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 104.01%			
S p-Bromofluorobenzene	10.951	95.0	215644	283.5551	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 113.42%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.422	50.0	1730	1.4445	ng	m	98
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.338	49.0	2222	1.9869	ng	m	91
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	5.647	83.0	209	0.1455	ng	m	80

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.286	78.0	144	0.0480	ng	m	89
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	0.000		0	N.D.			
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.386	92.0	3845	1.9977	ng		100
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	0.000		0	N.D.			
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	9.917	91.0	389	0.1063	ng	m	61
T m+p-Xylenes	10.045	106.0	0		ng	md	1
T o-Xylene	0.000		0	N.D.			
T Styrene	0.000		0	N.D.			
T Bromoform	0.000		0	N.D.			
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

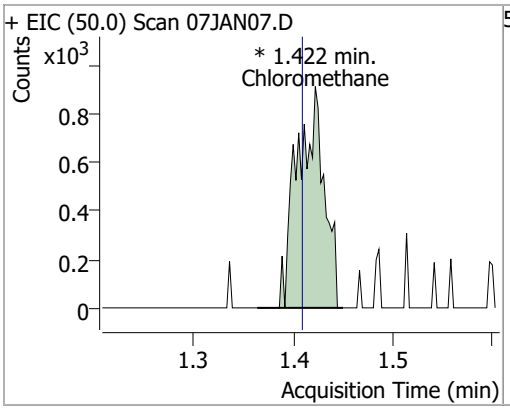
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

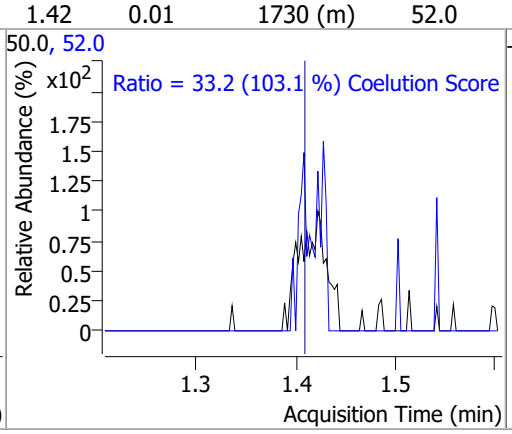
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

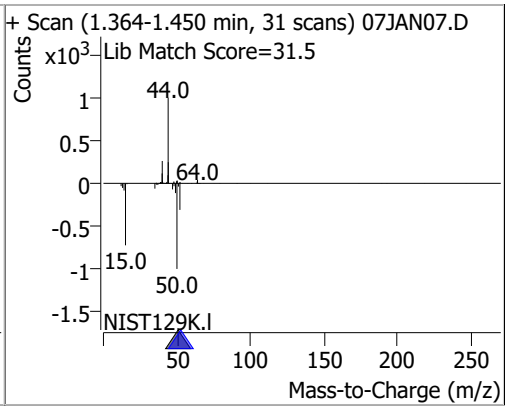




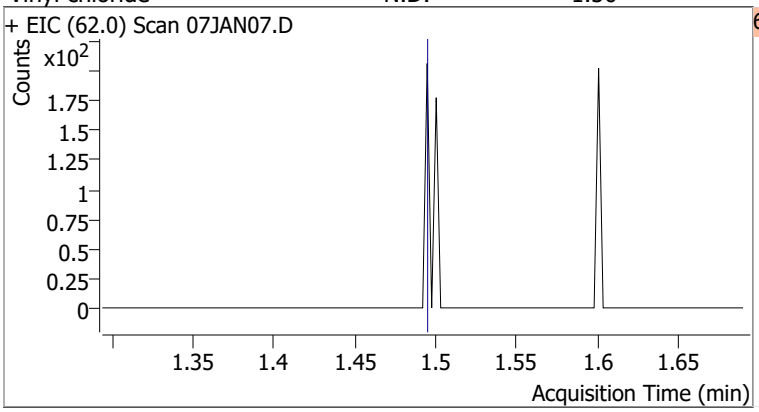
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	1.4445	1.42	0.01	1730 (m)	52.0	33.2	2.1	62.1

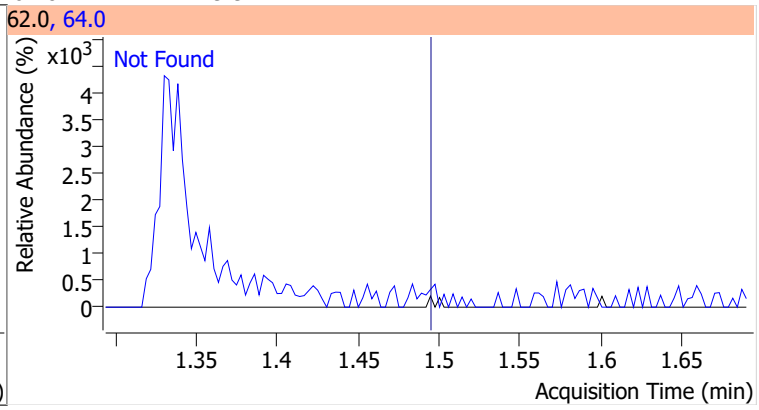




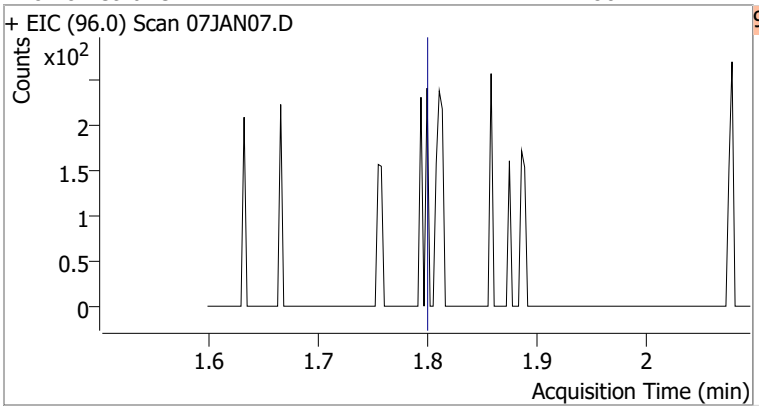


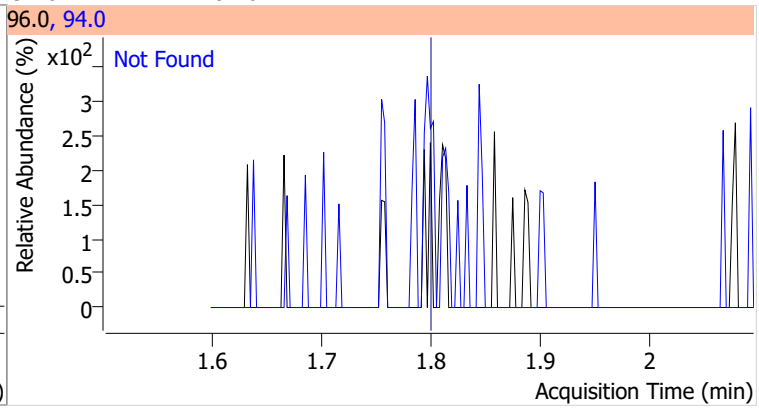
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



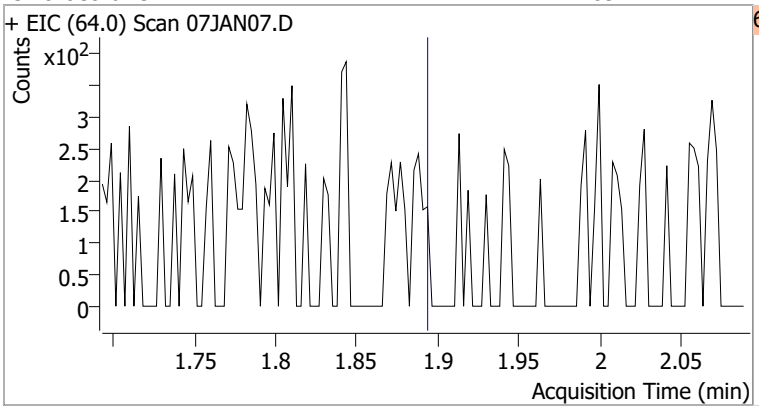


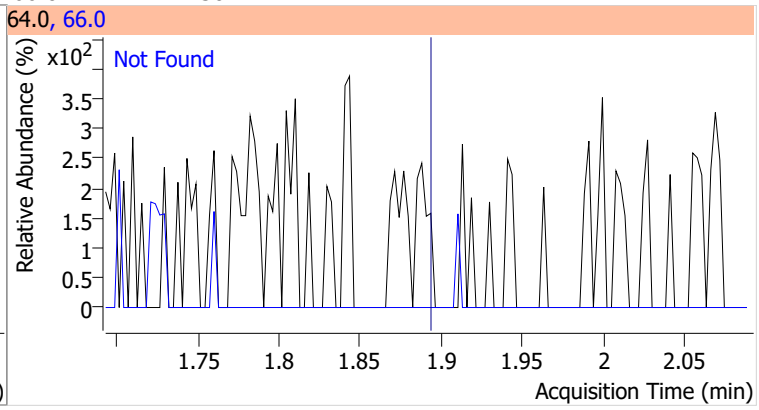
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



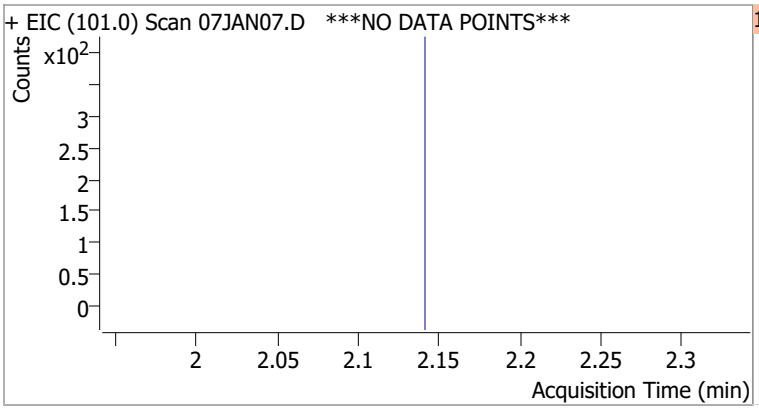


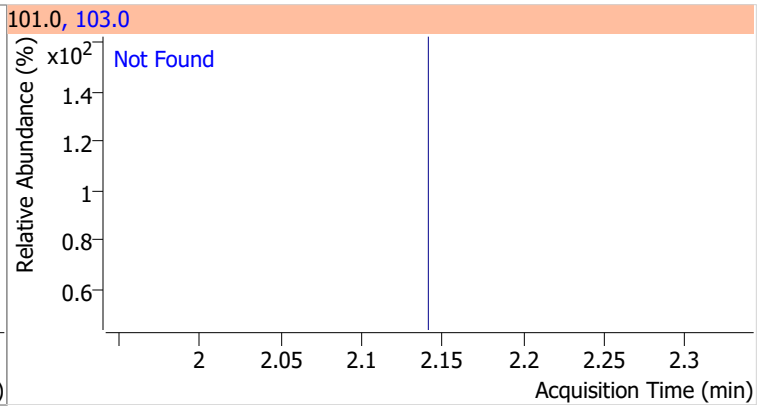
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



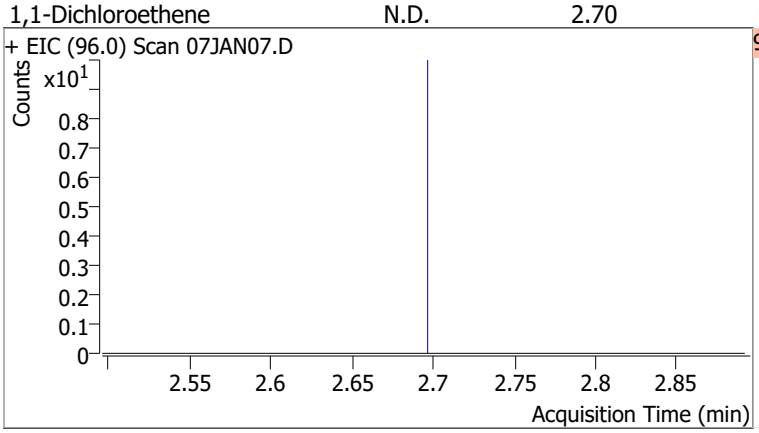


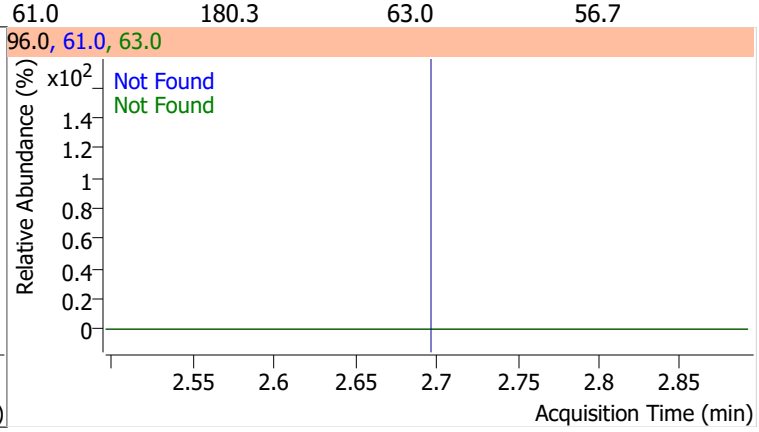
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



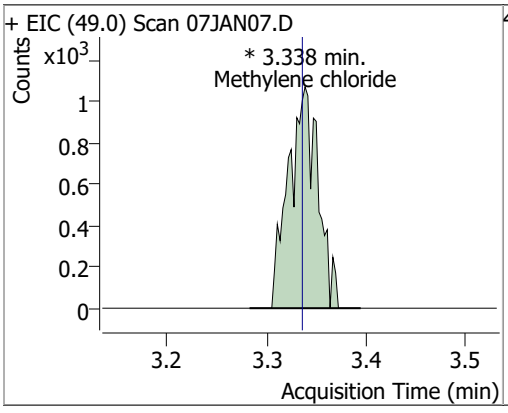


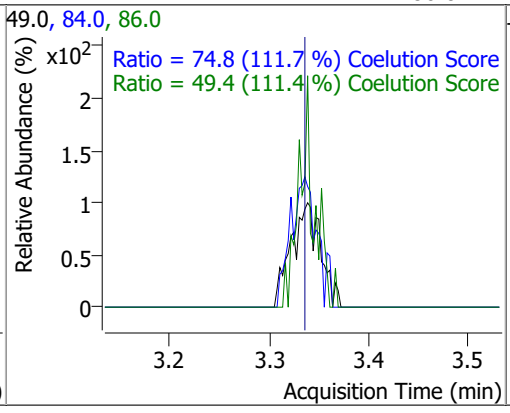
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

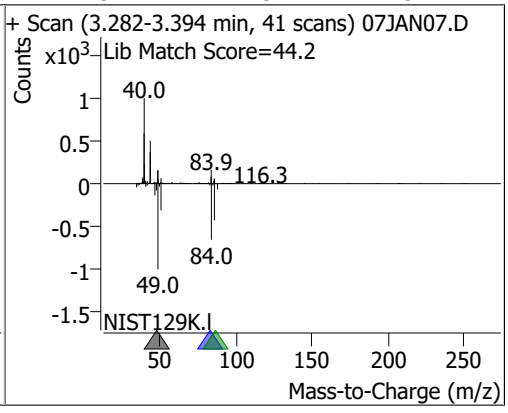




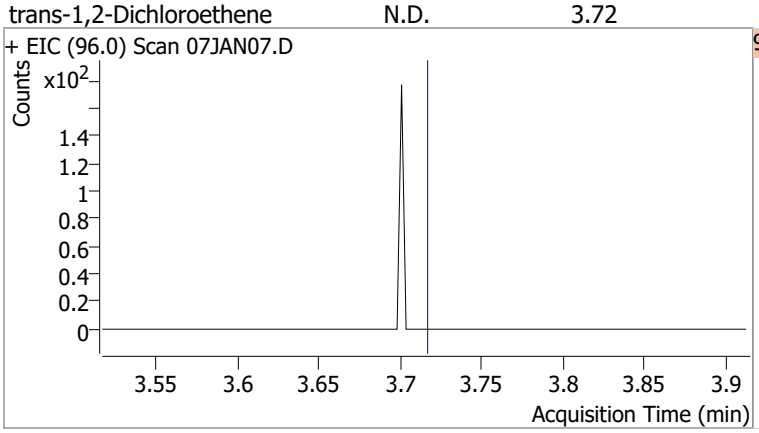
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	1.9869	3.34	0.00	2222 (m)	84.0	74.8	36.9	96.9
					86.0	49.4	14.3	74.3

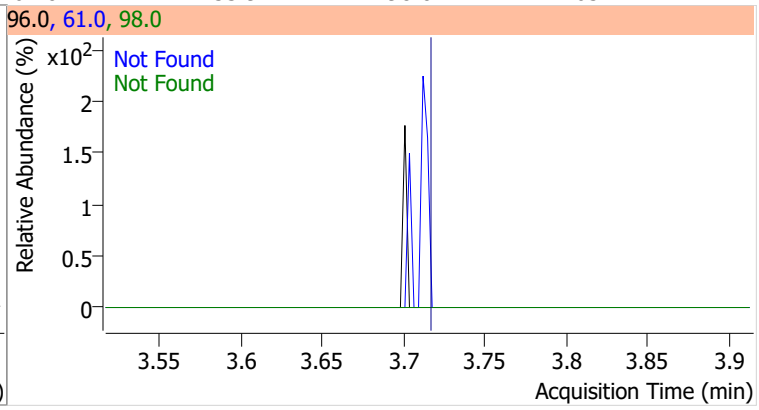




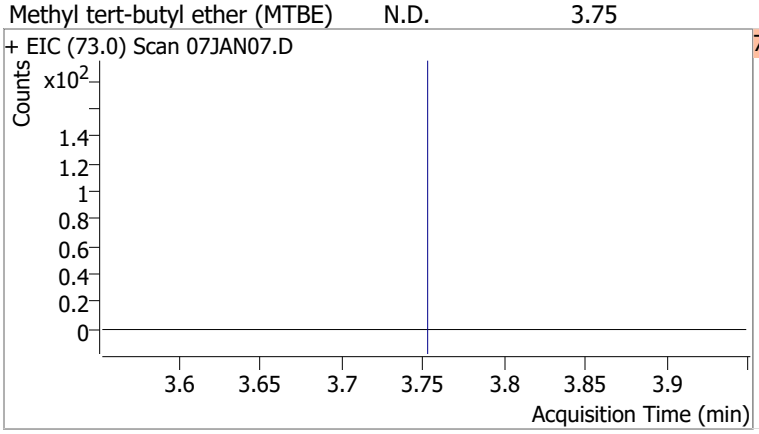


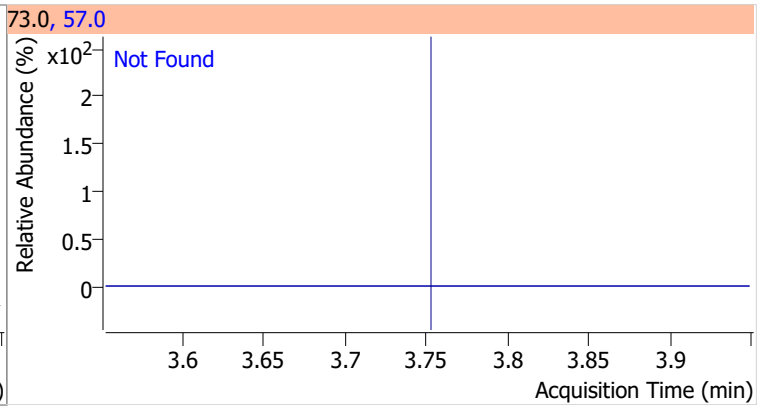
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



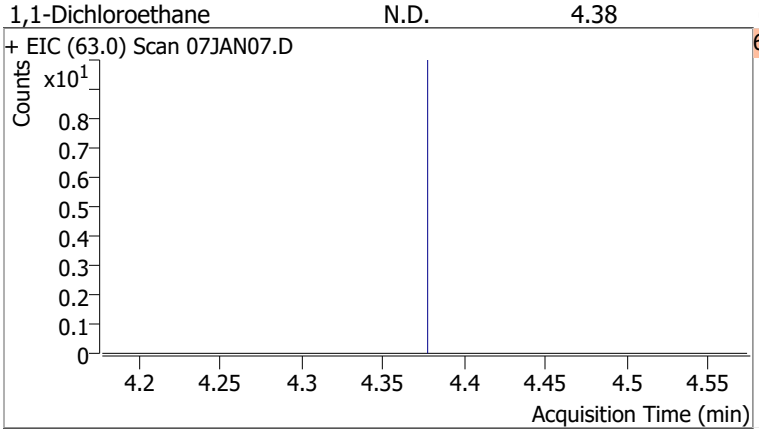


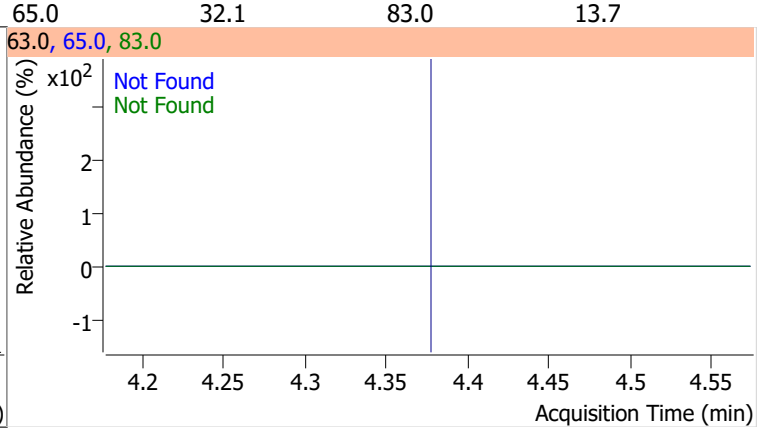
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



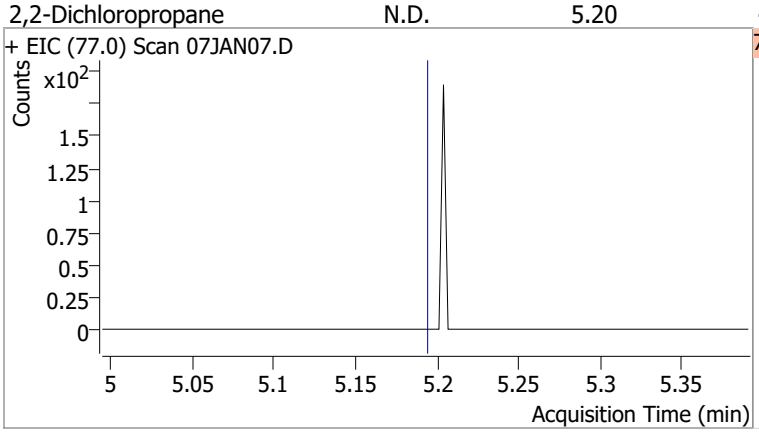


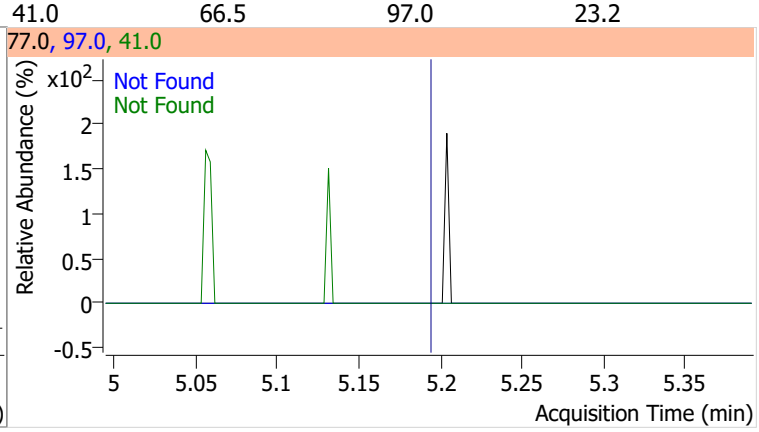
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



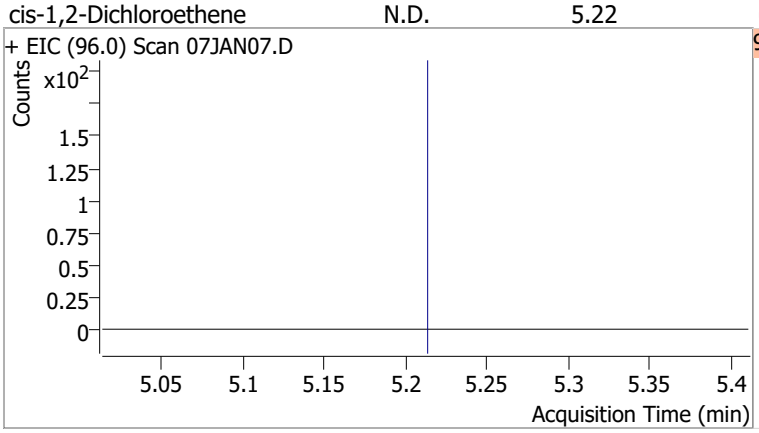


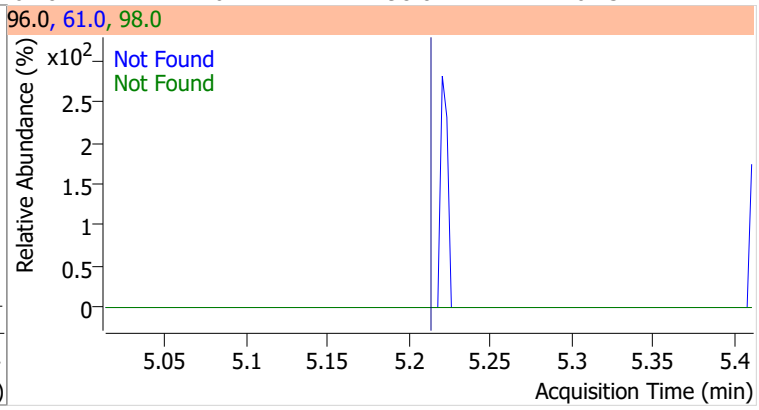
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



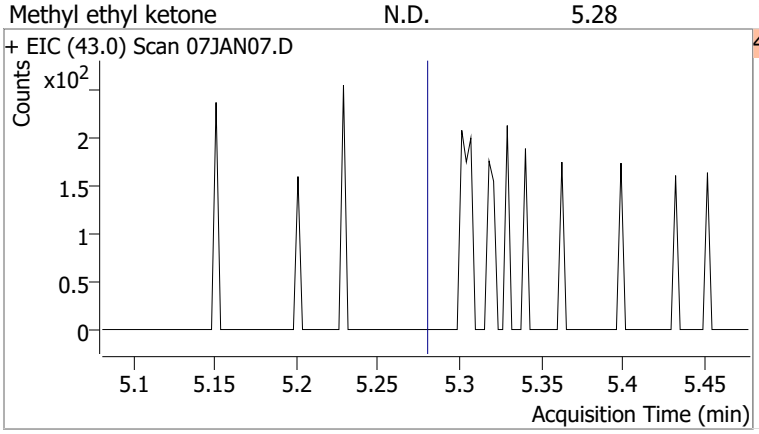


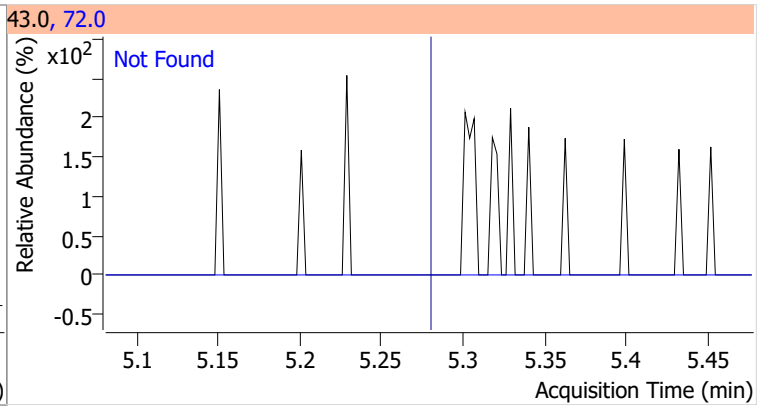
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



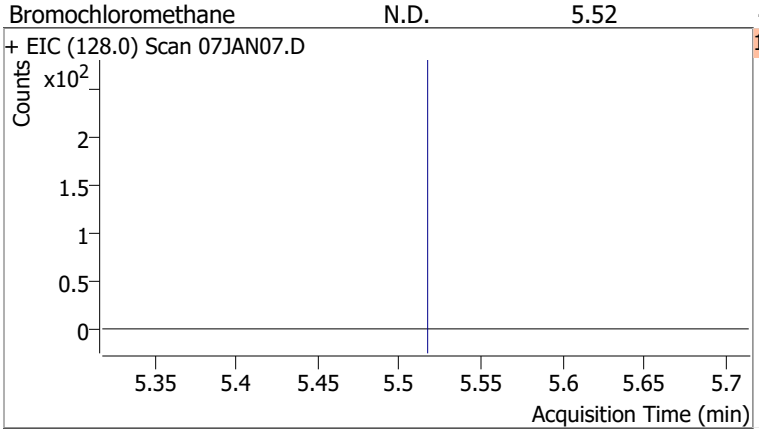


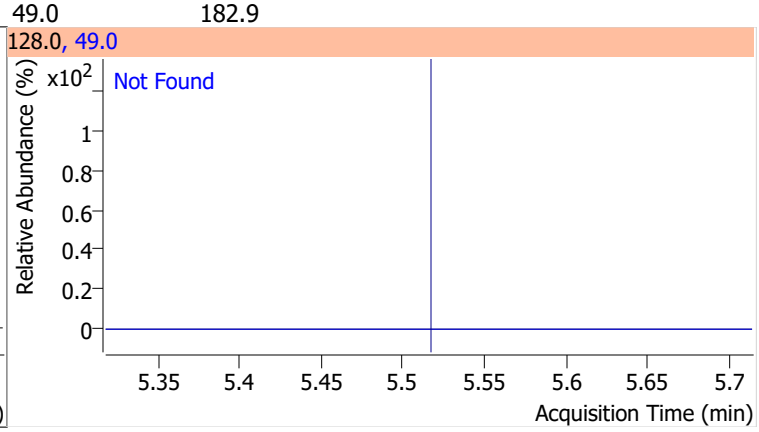
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



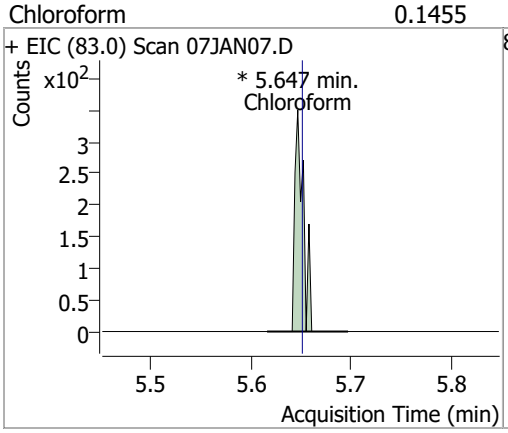


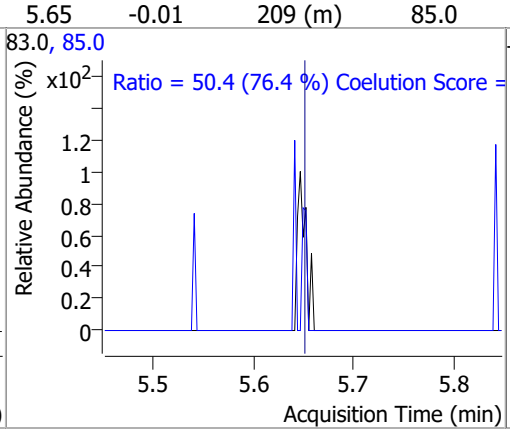
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9

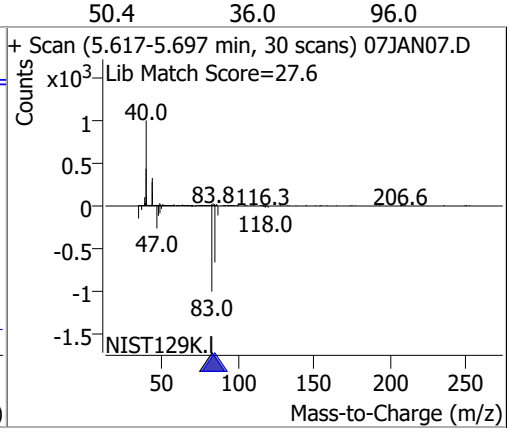




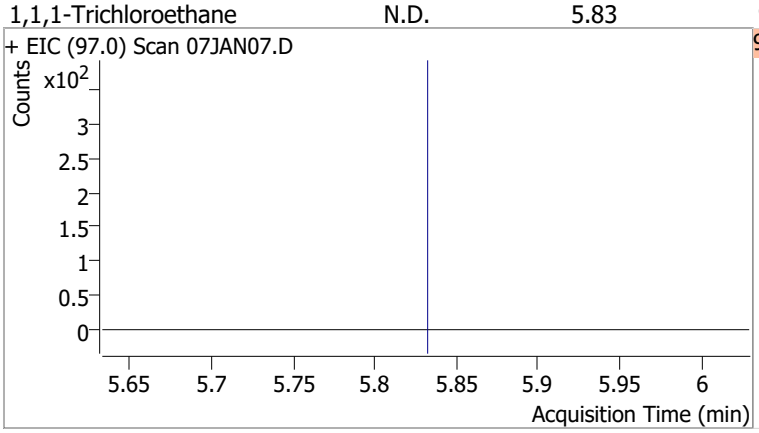
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	0.1455	5.65	-0.01	209 (m)	85.0	50.4	36.0	96.0

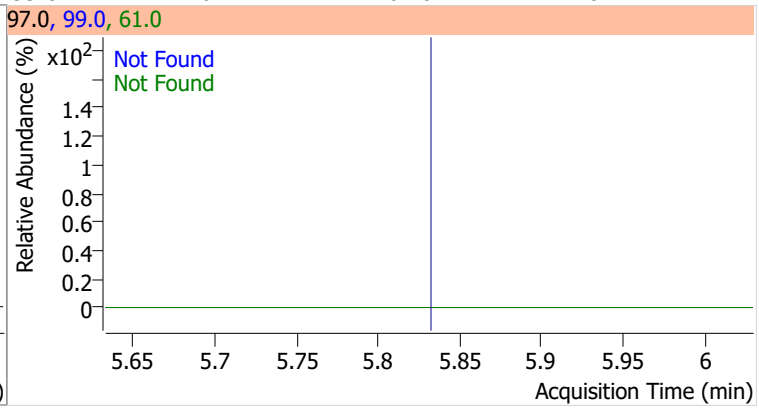




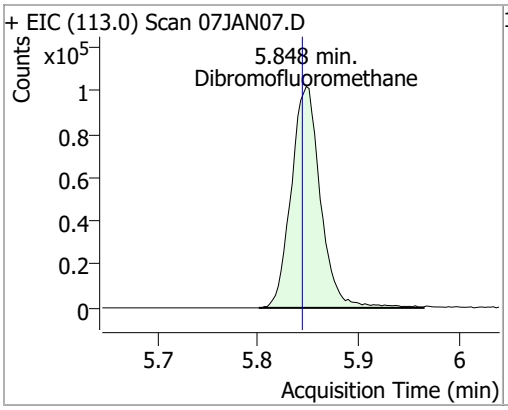


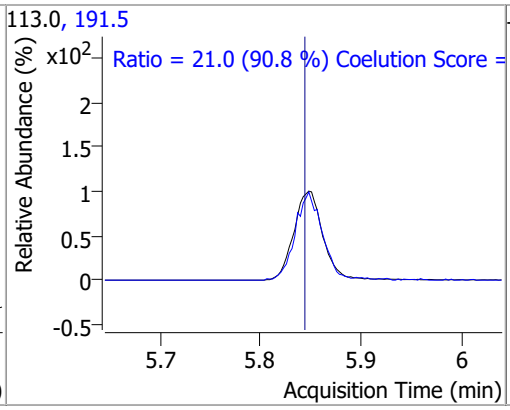
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

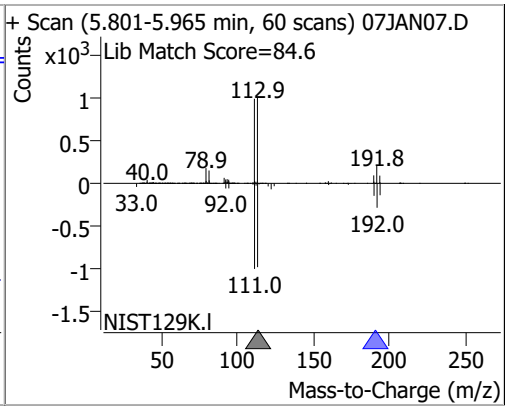




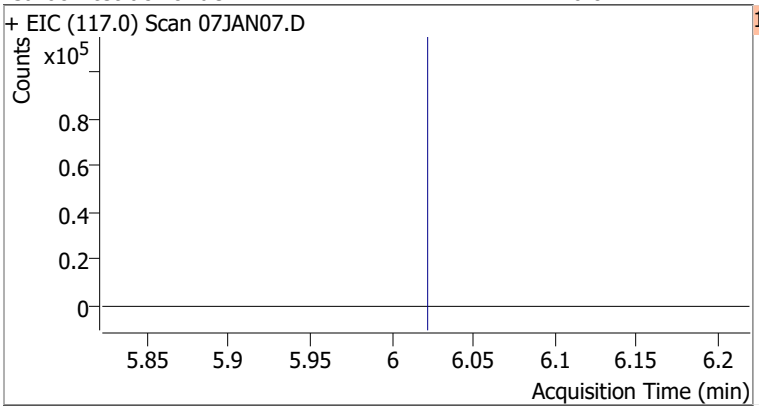
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	292.5939	5.85	0.00	207577	191.5	21.0	0.0	53.1

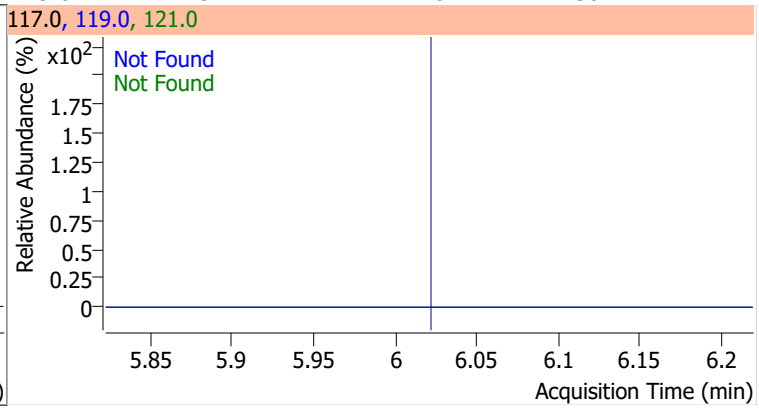




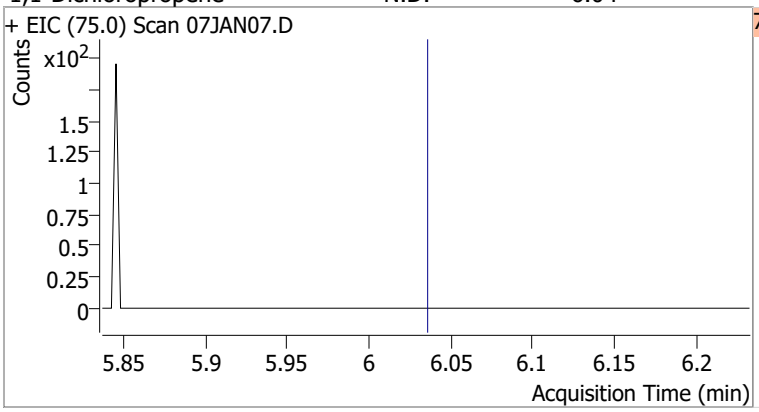


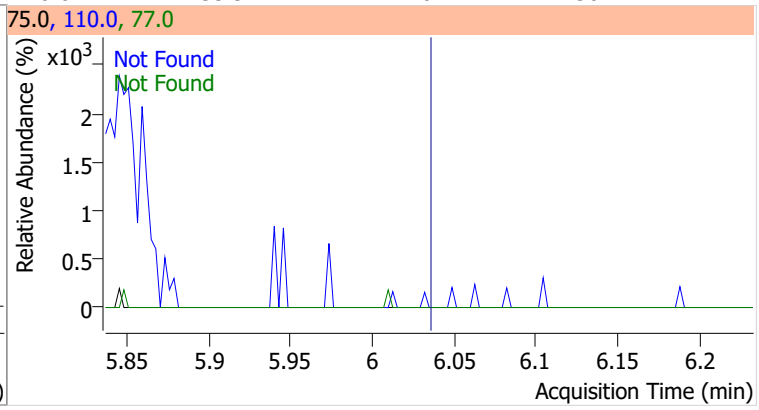
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



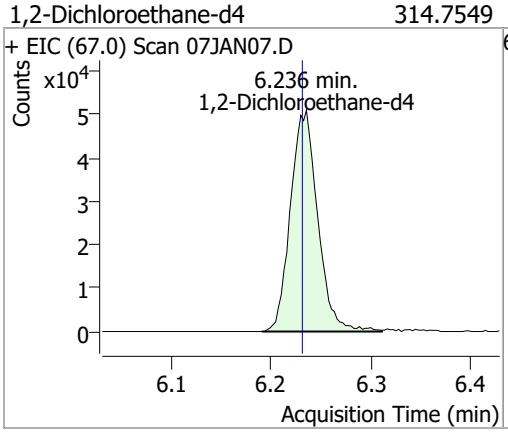


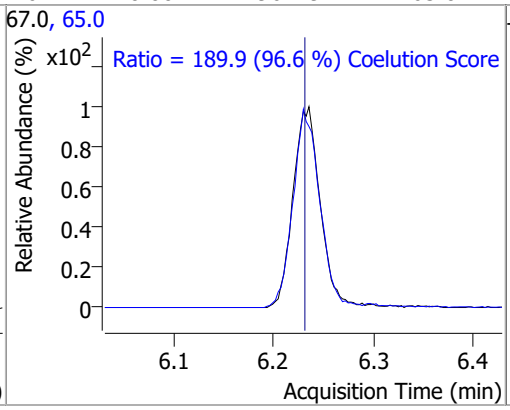
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

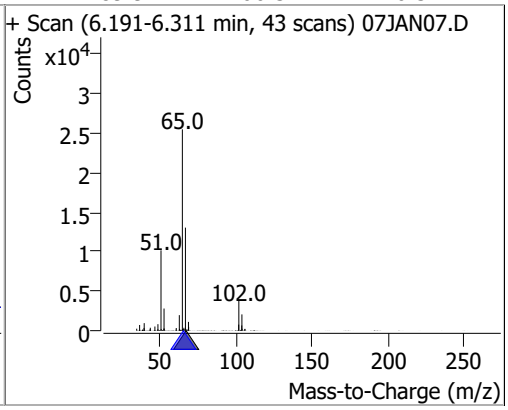




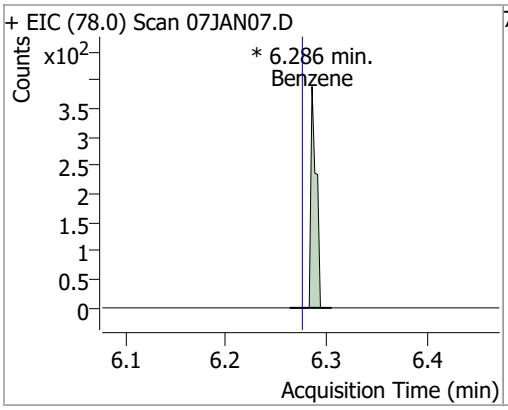
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	314.7549	6.24	0.00	96449	65.0	189.9	166.5	226.5

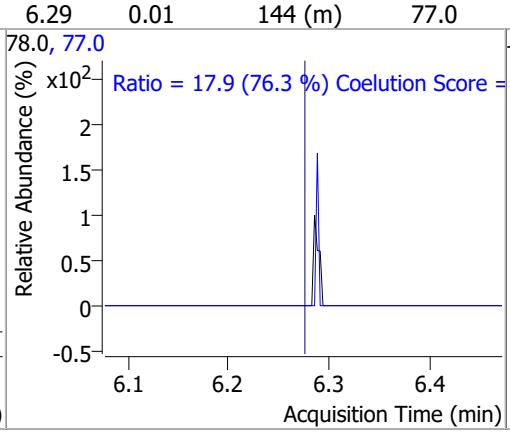


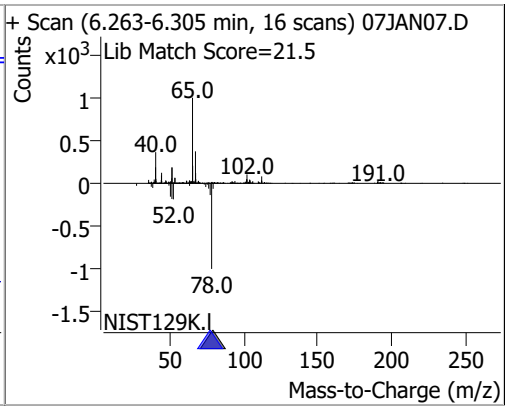




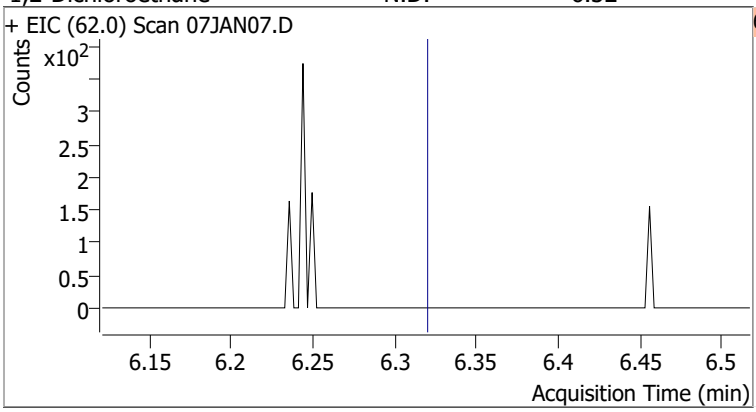
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.0480	6.29	0.01	144 (m)	77.0	17.9	0.0	53.5

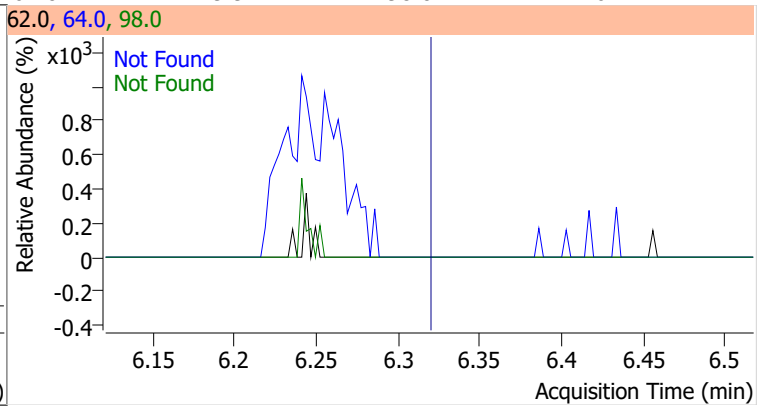




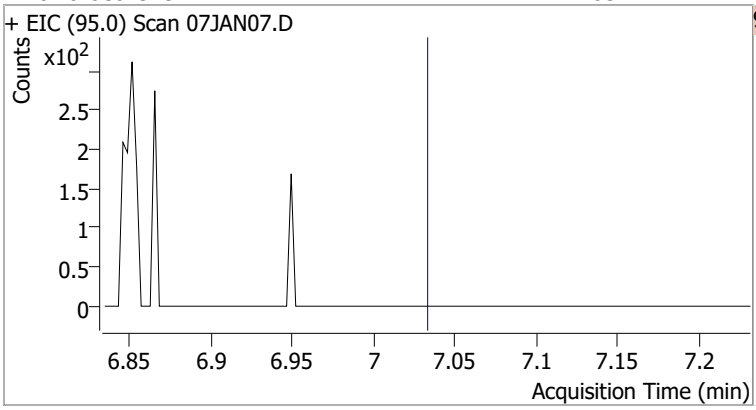


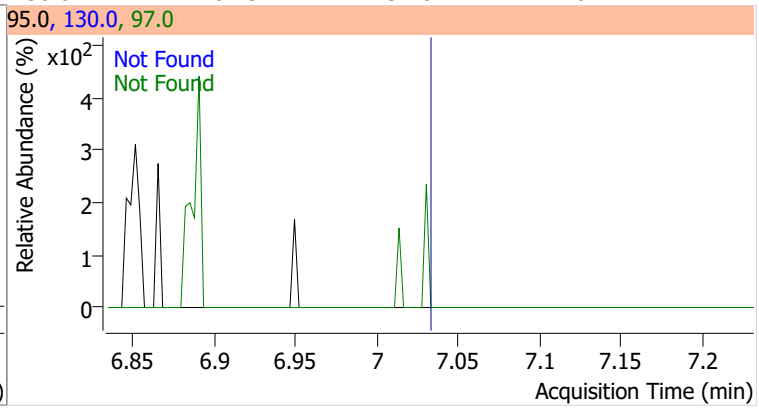
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



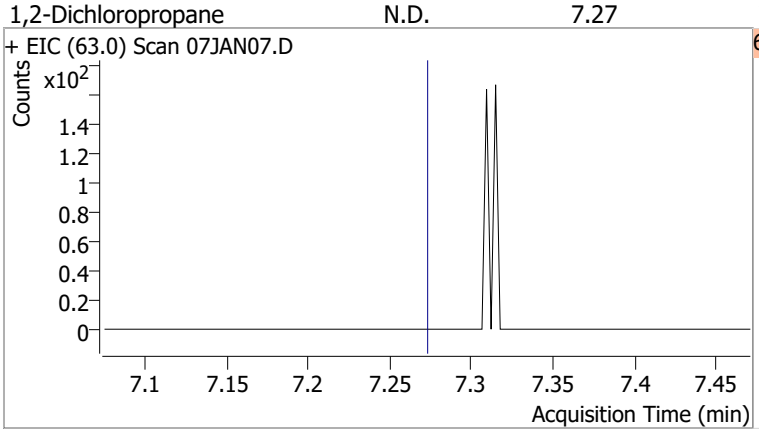


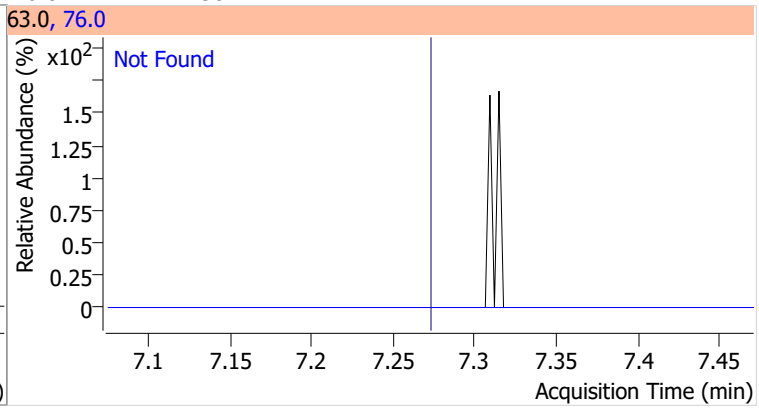
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



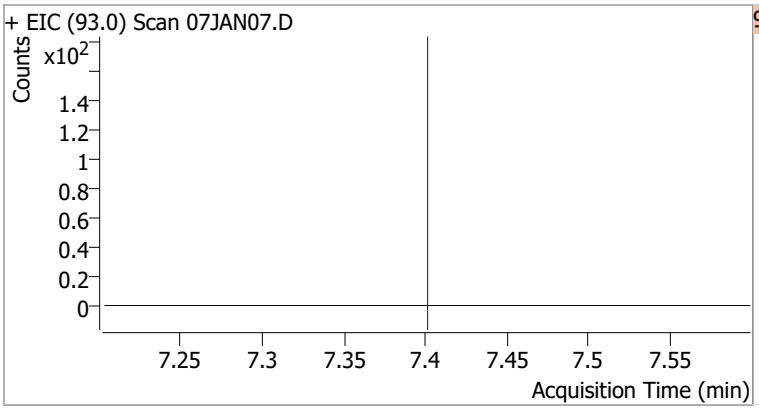


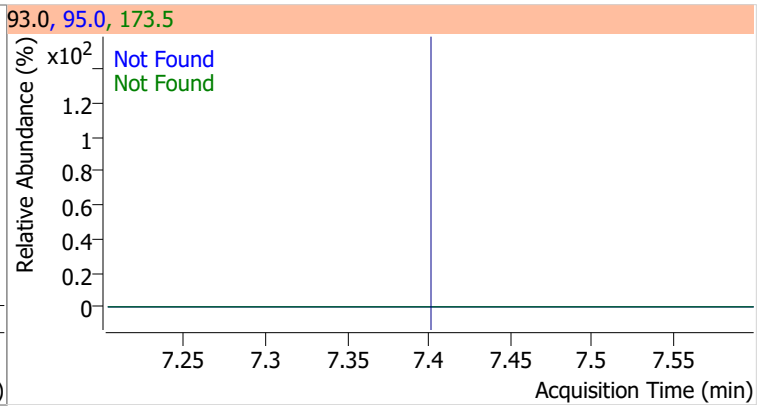
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



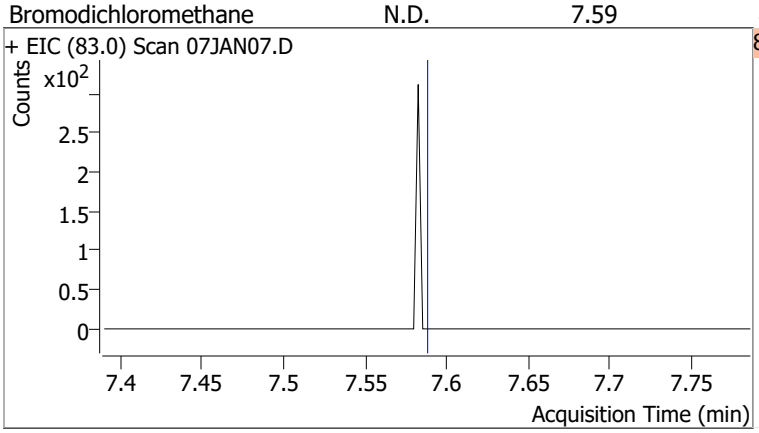


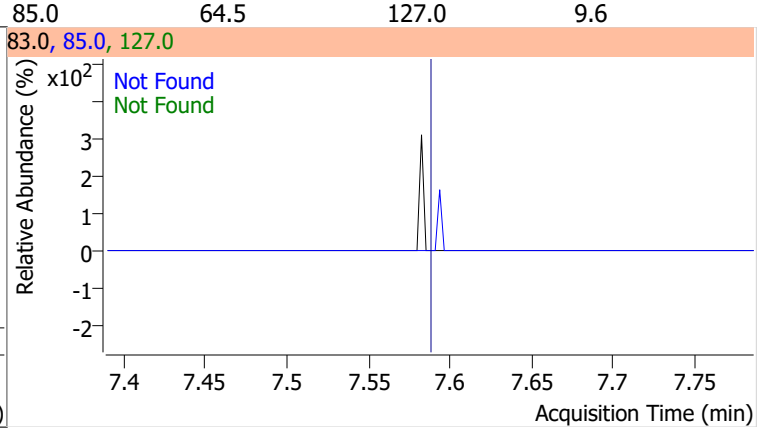
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



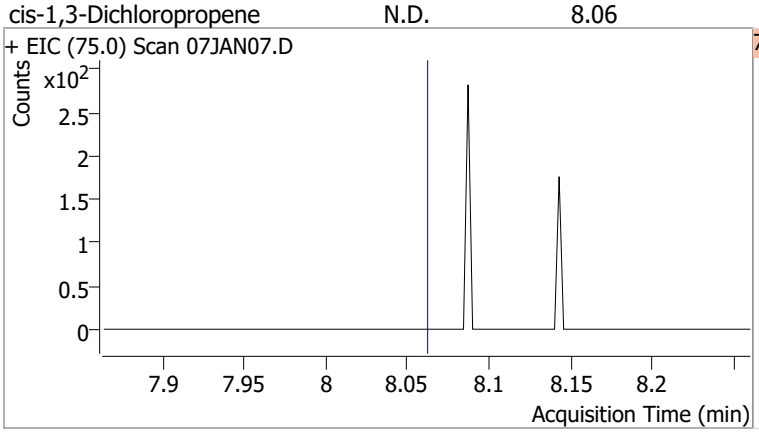


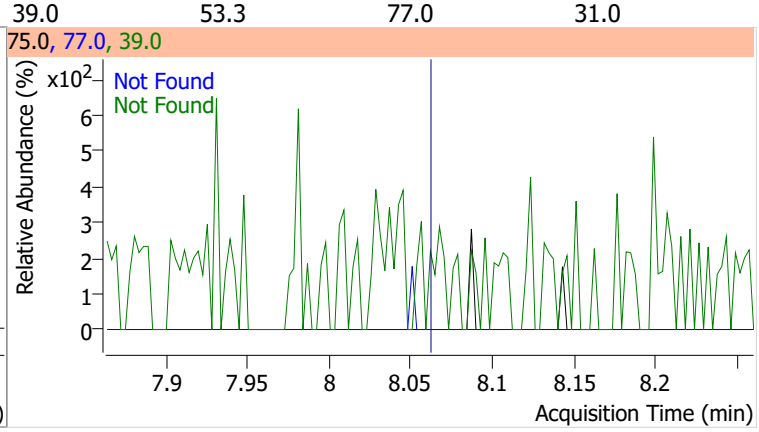
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



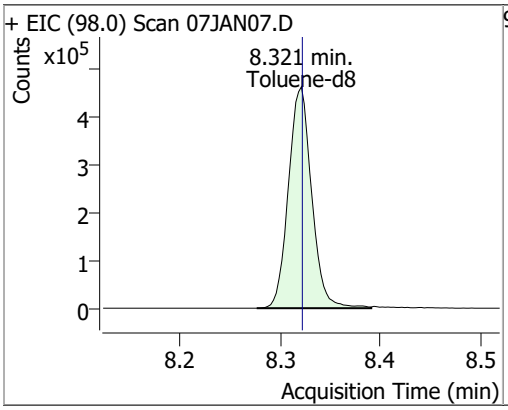


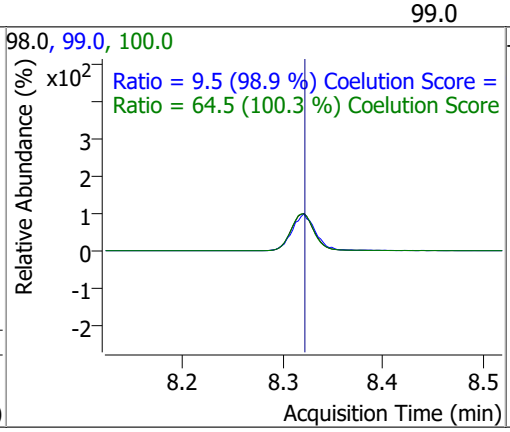
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

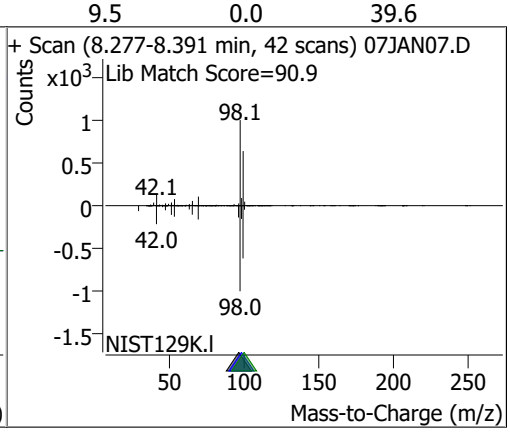




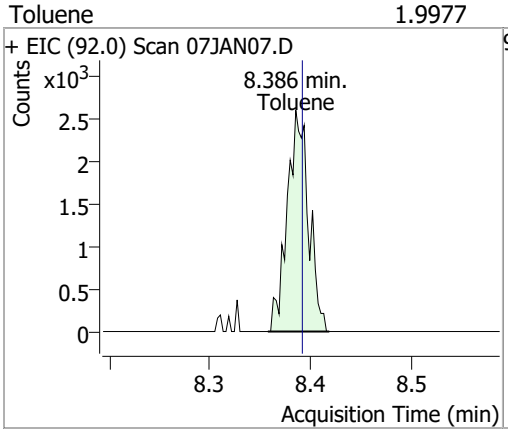
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	260.0307	8.32	0.00	740894	100.0	64.5	34.4	94.4
					99.0	9.5	0.0	39.6

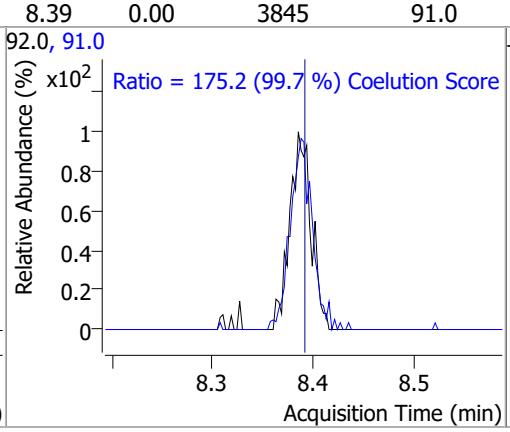


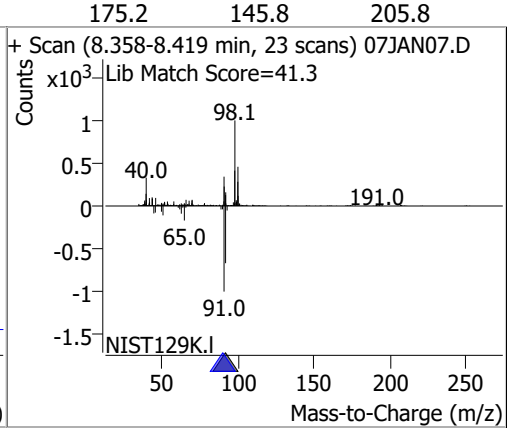




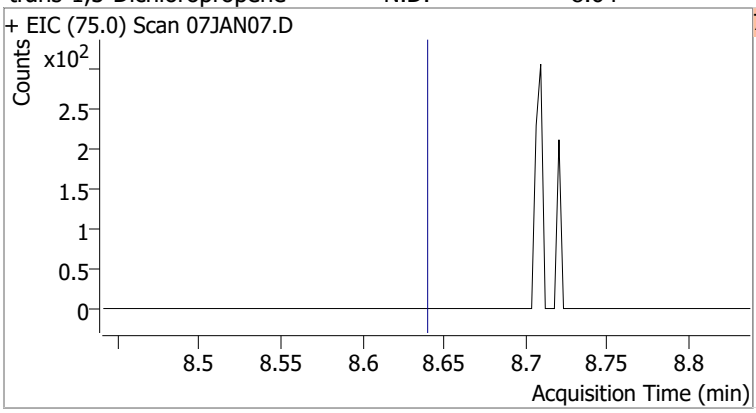
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	1.9977	8.39	0.00	3845	91.0	175.2	145.8	205.8

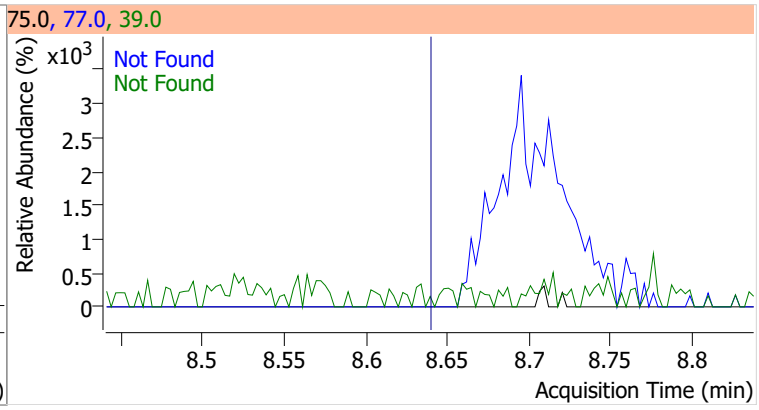




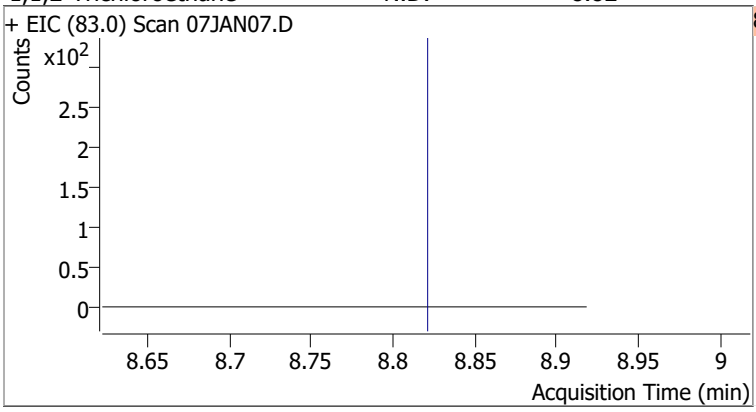


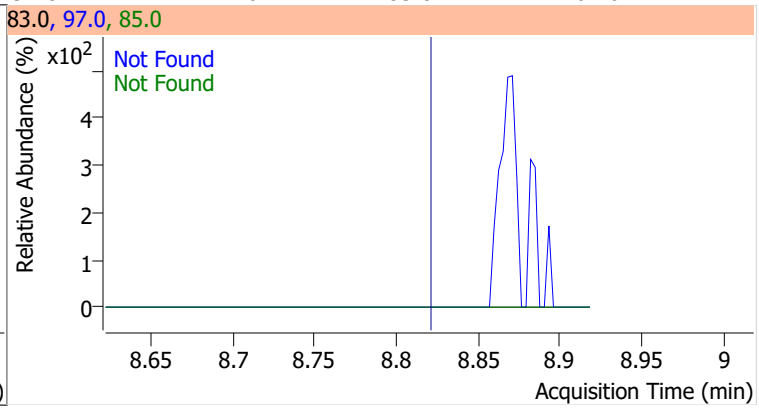
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



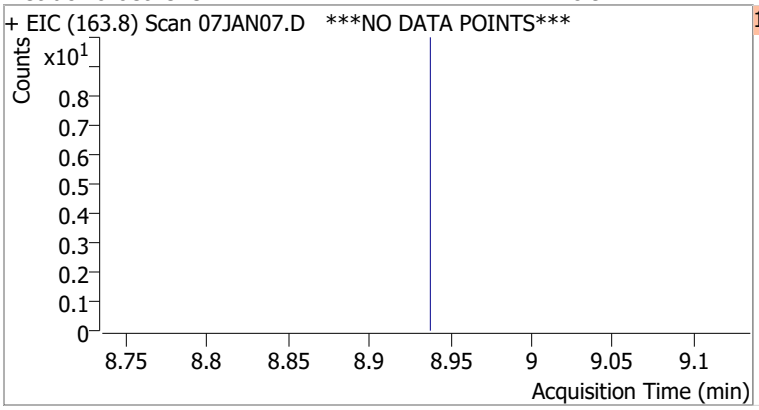


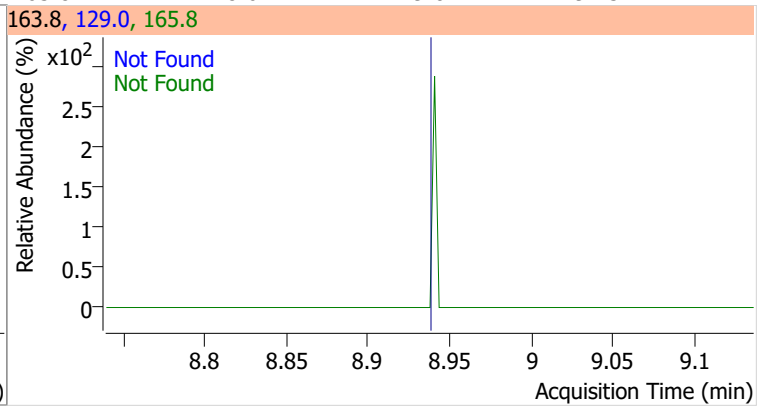
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



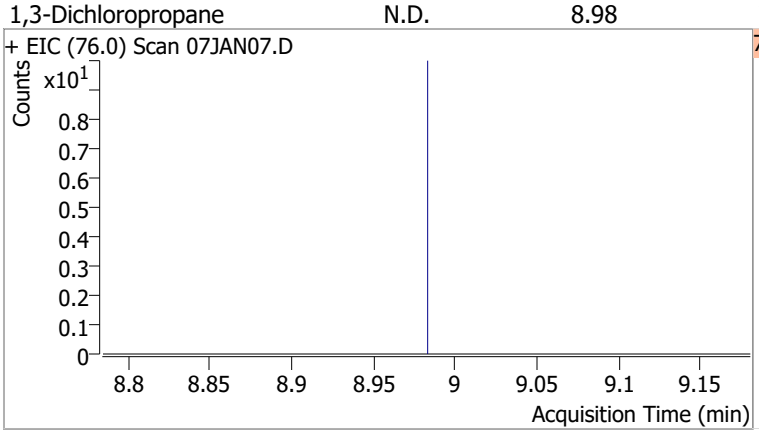


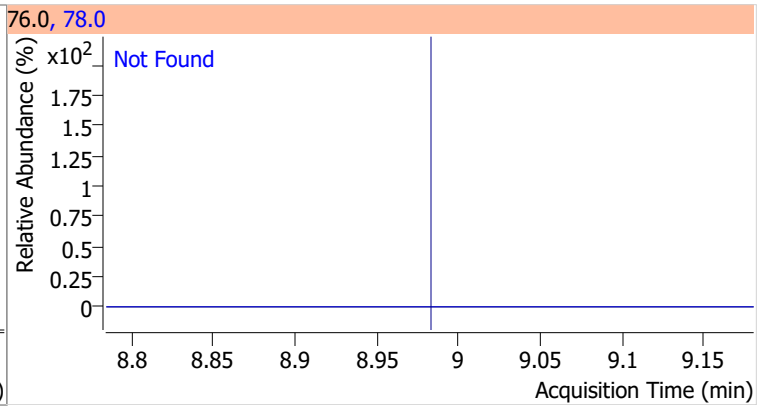
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



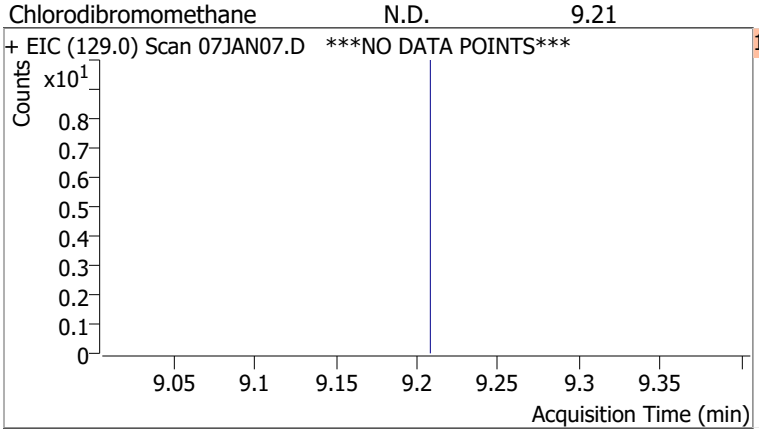


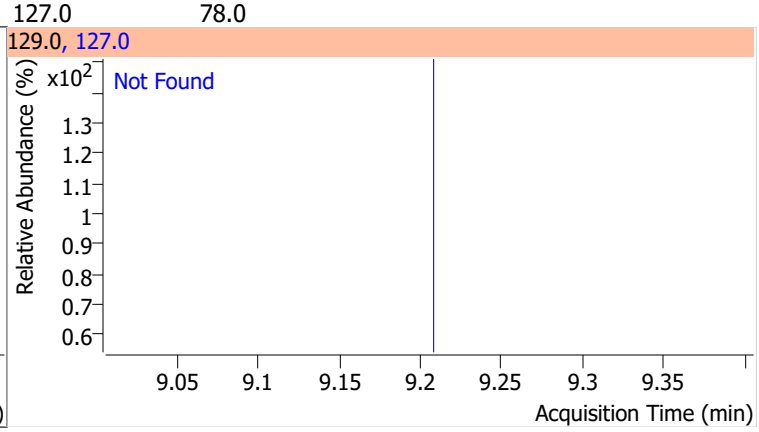
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



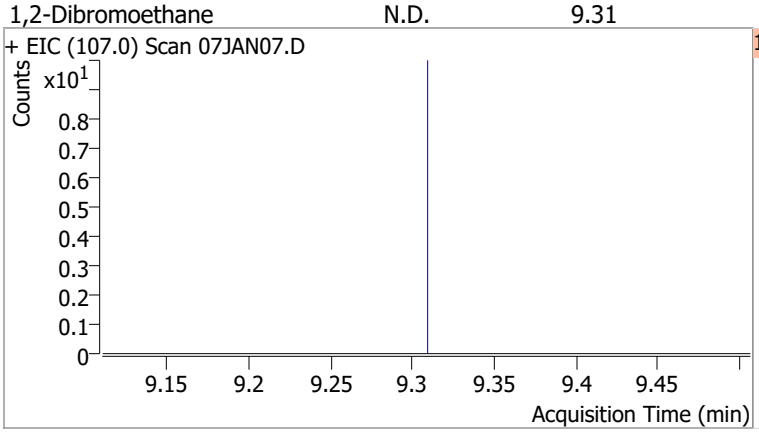


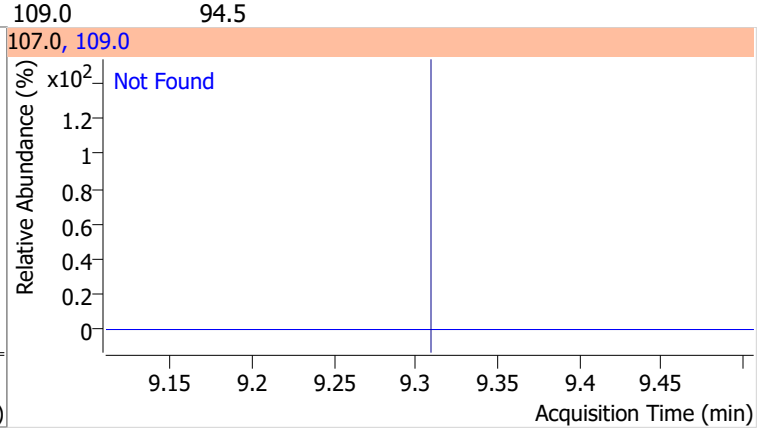
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



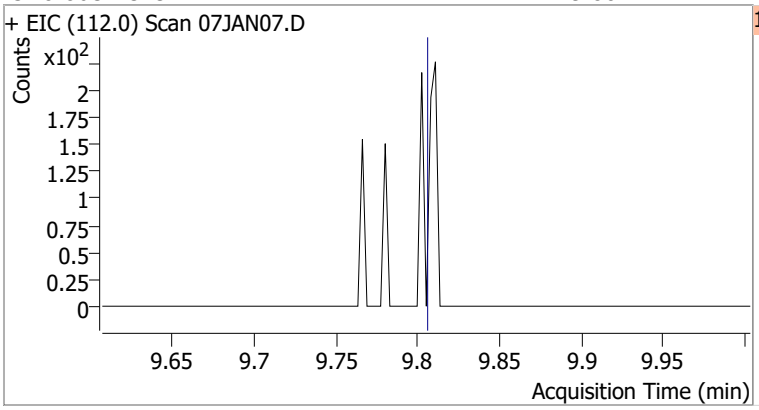


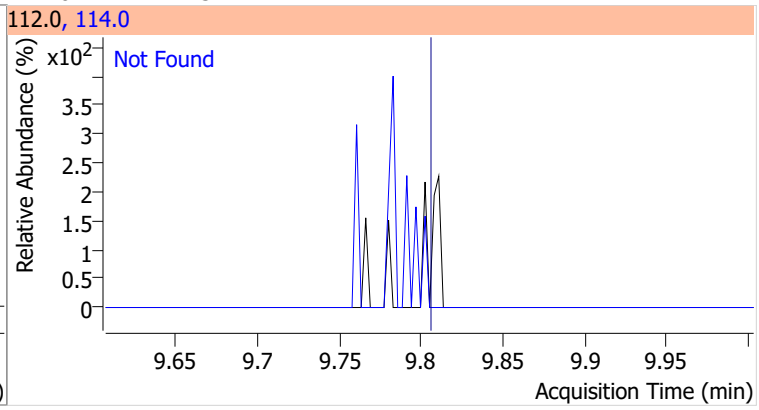
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



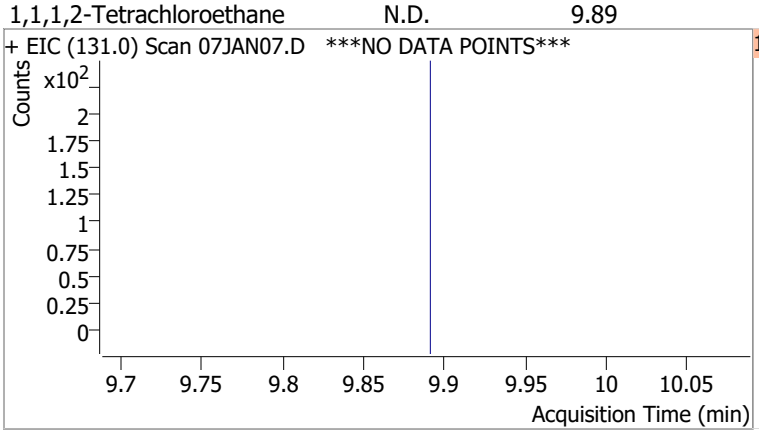


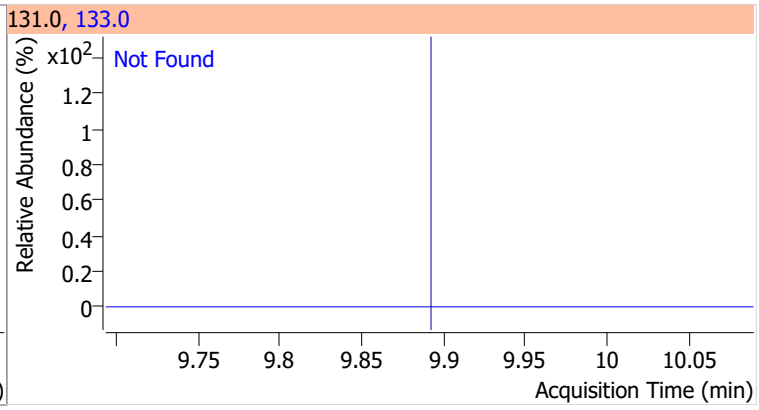
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



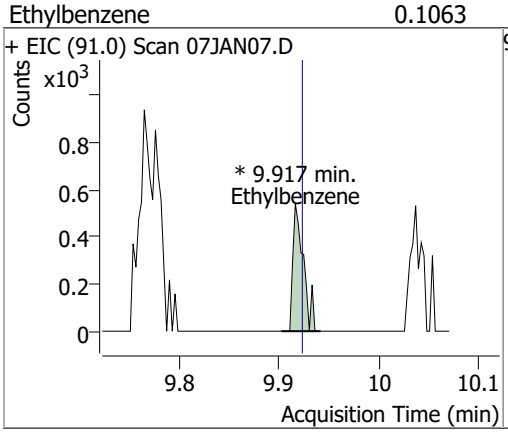


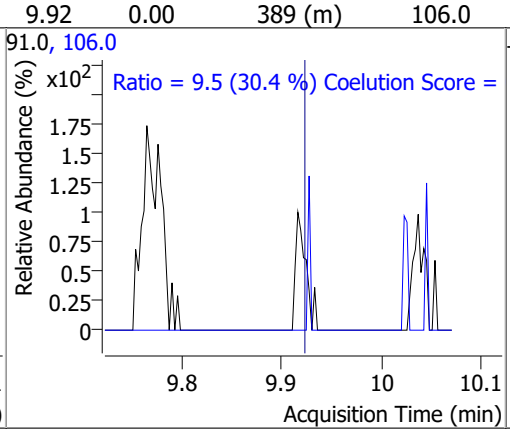
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6

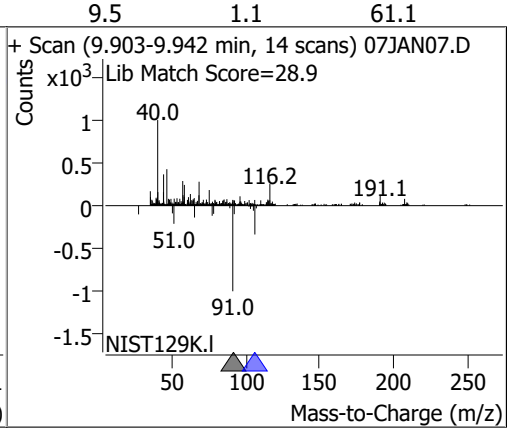




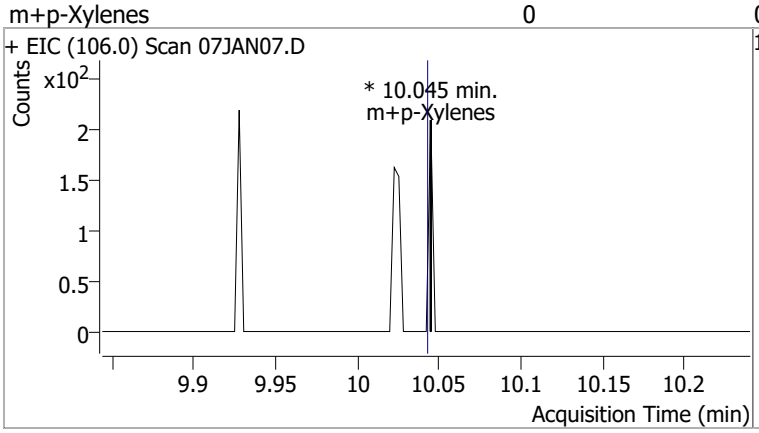
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	0.1063	9.92	0.00	389 (m)	106.0	9.5	1.1	61.1

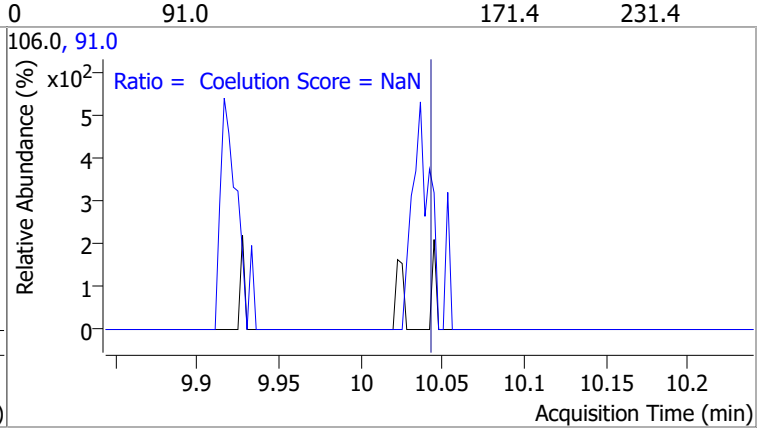




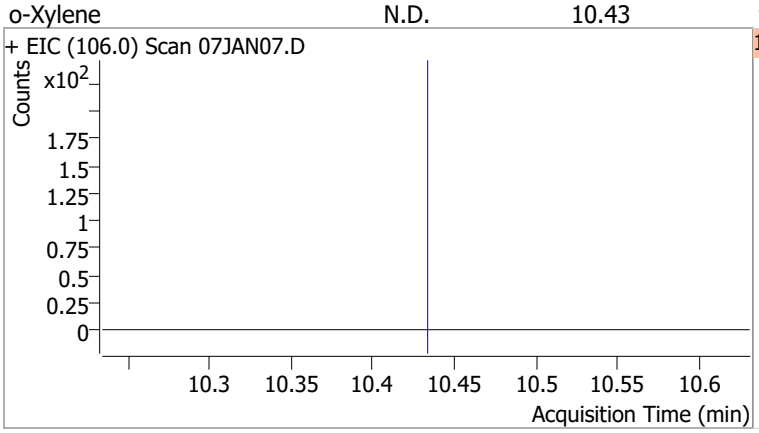


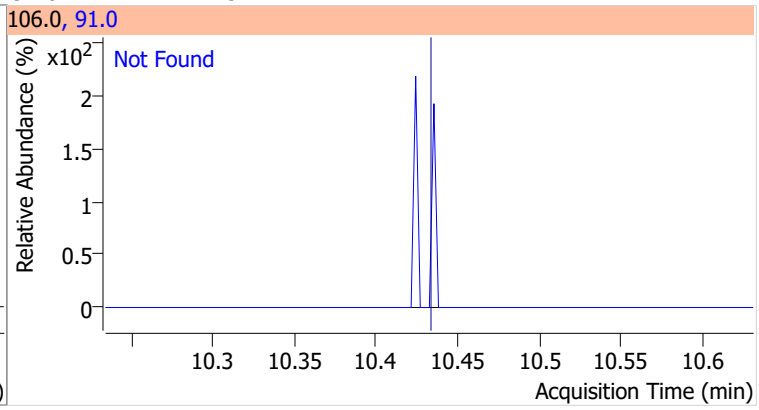
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes		0		0	91.0		171.4	231.4



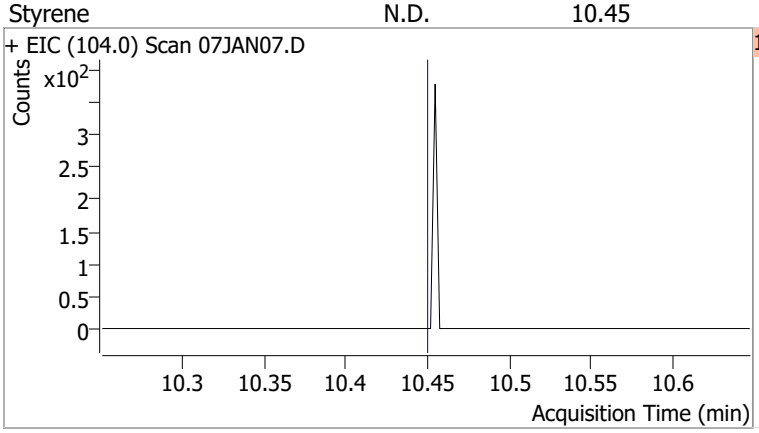


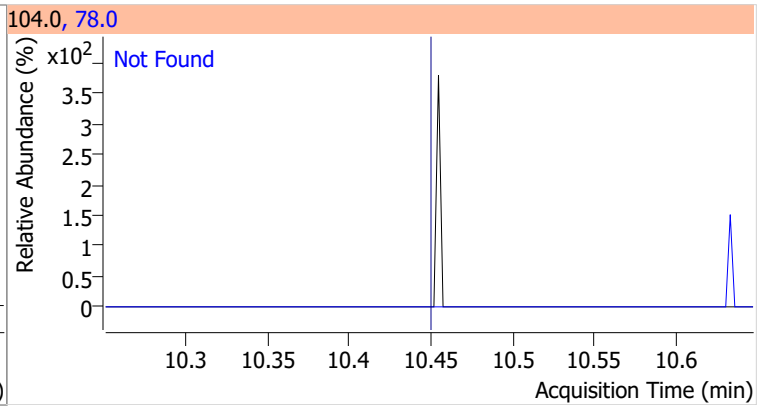
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



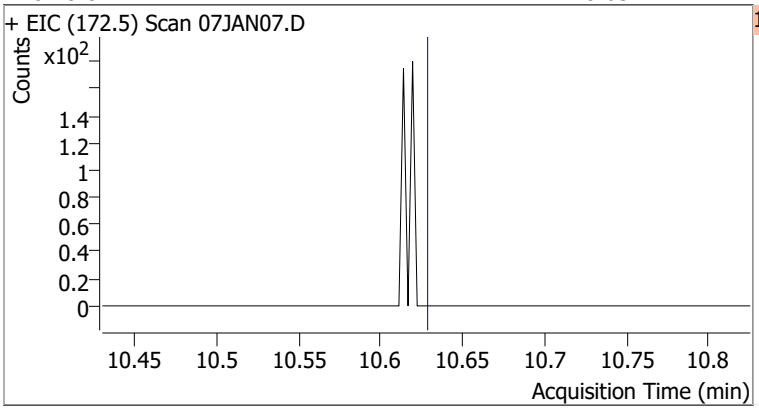


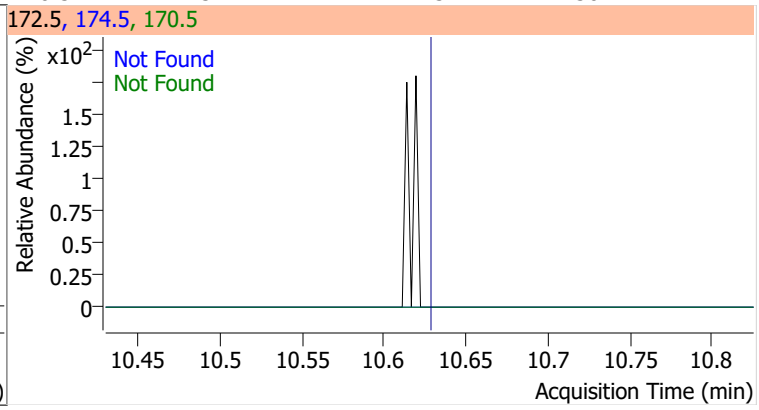
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



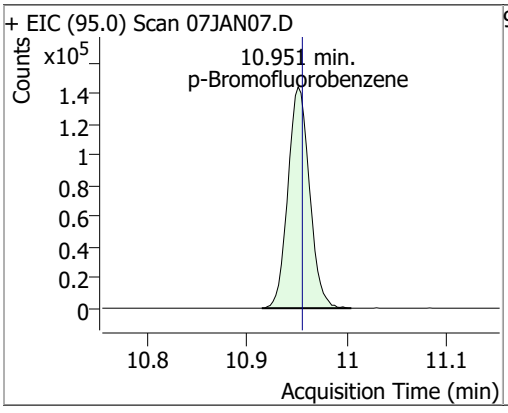


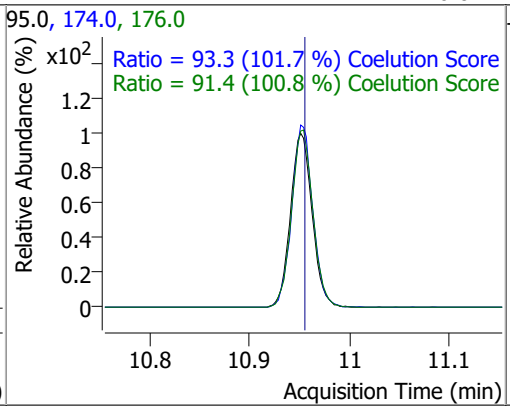
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

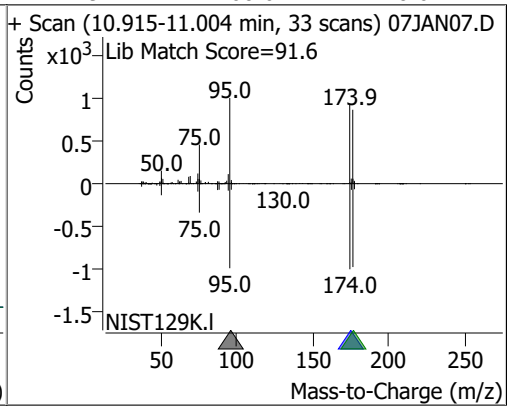




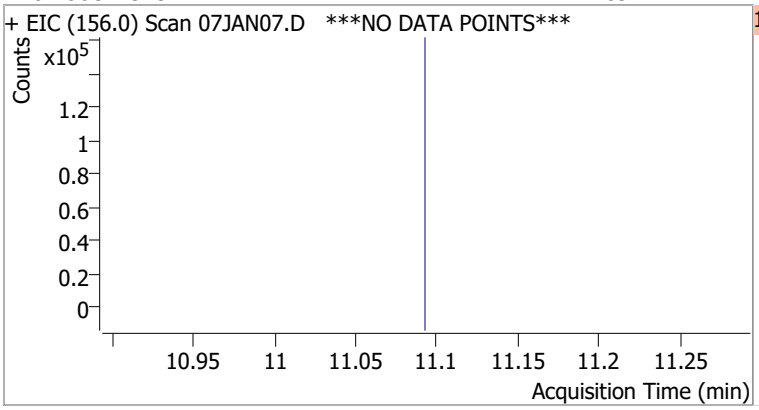
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	283.5551	10.95	0.00	215644	174.0	93.3	61.7	121.7
					176.0	91.4	60.6	120.6

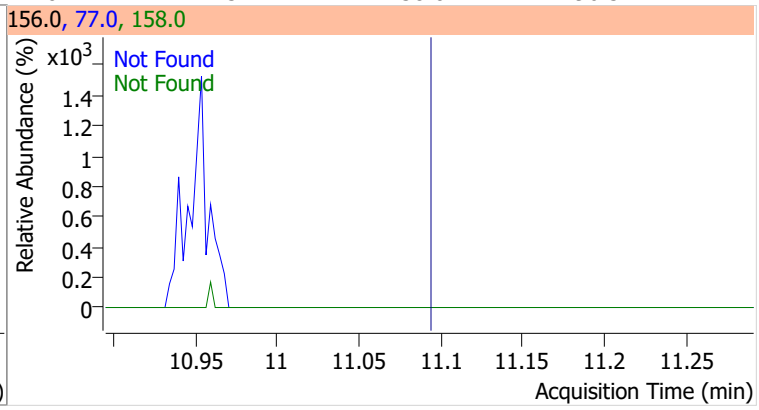




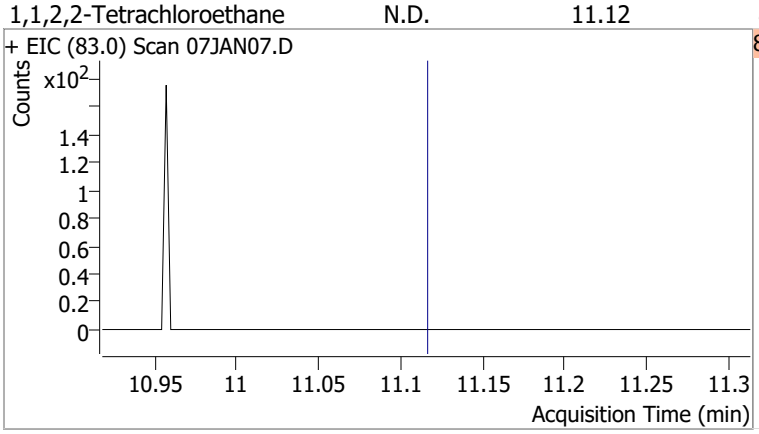


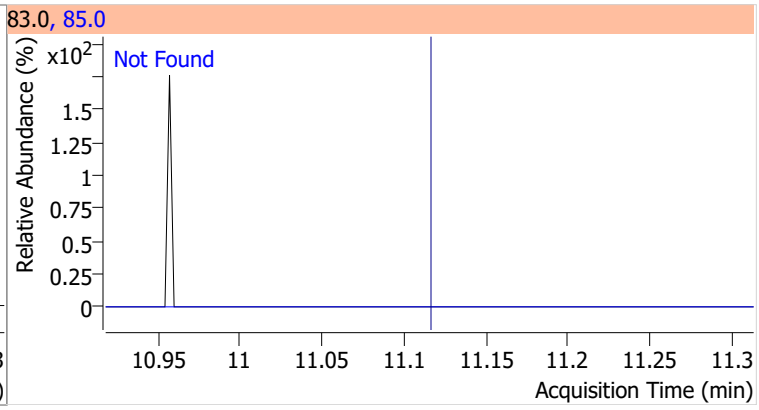
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



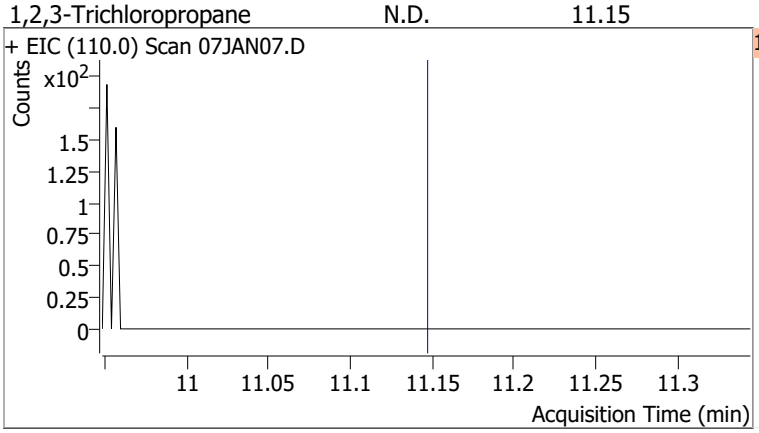


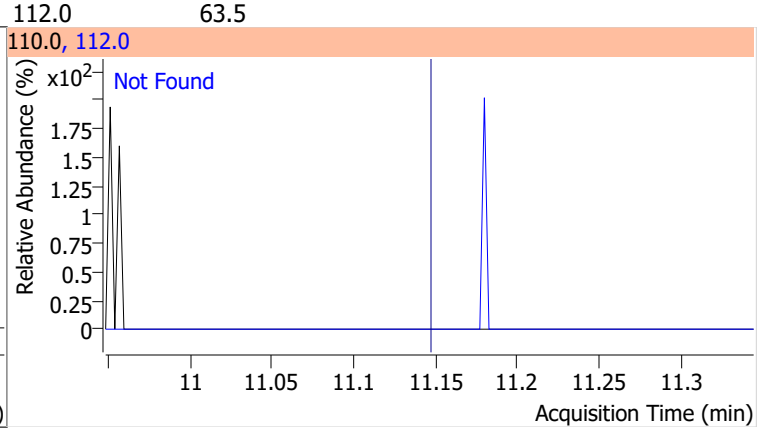
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



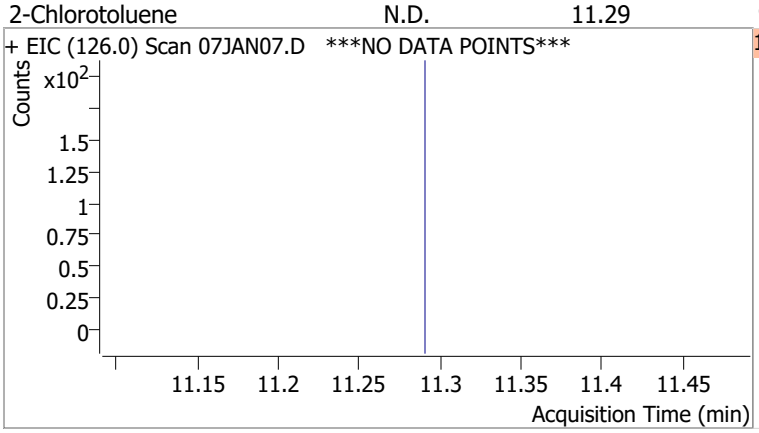


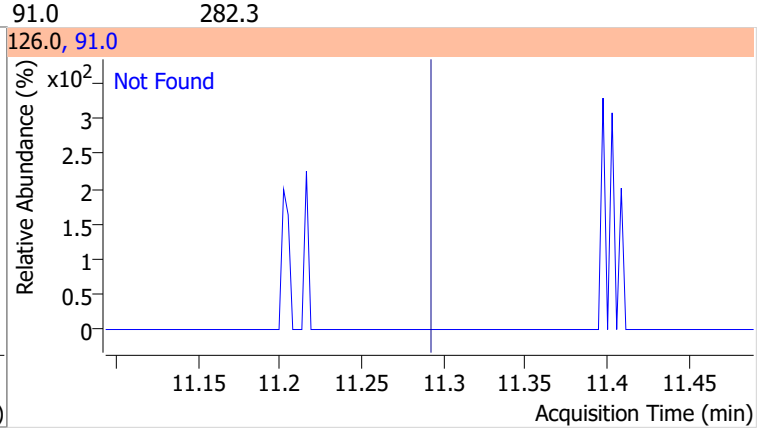
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



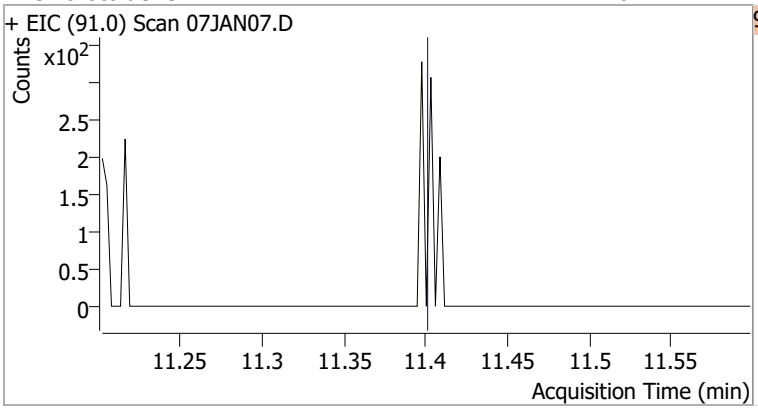


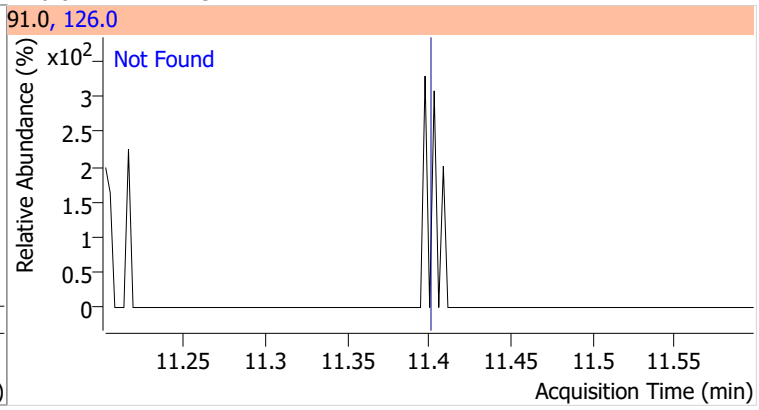
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



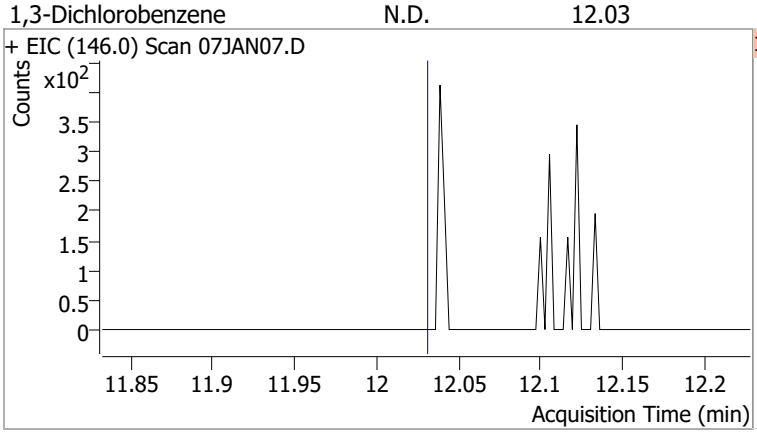


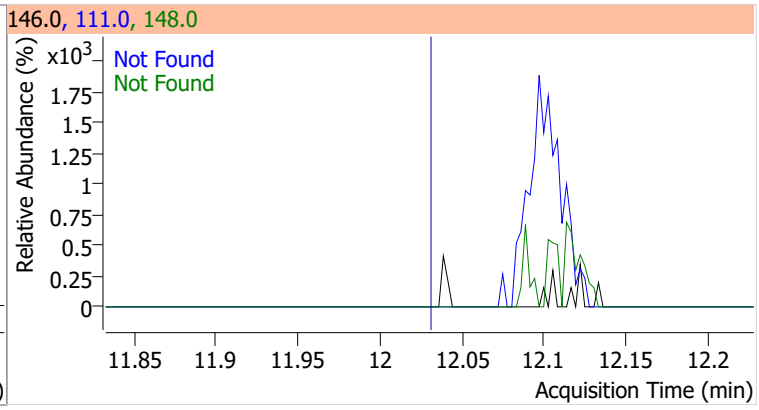
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



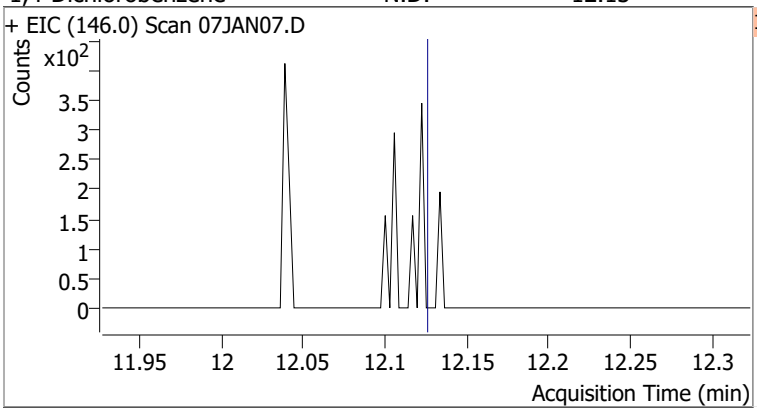


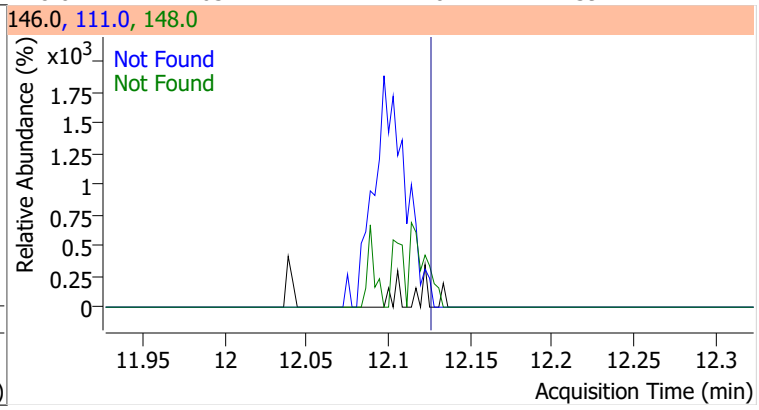
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



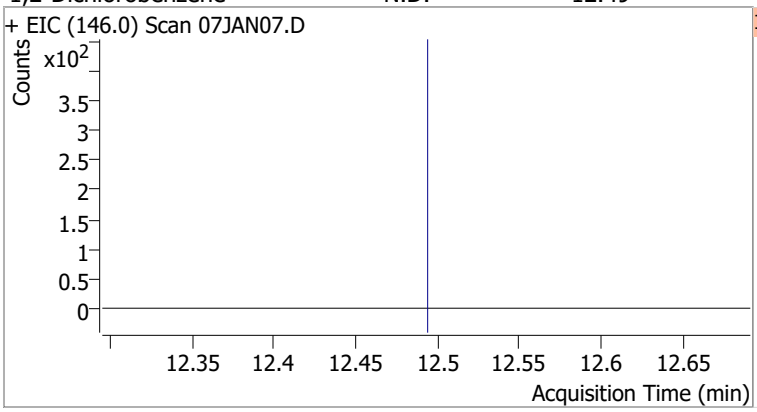


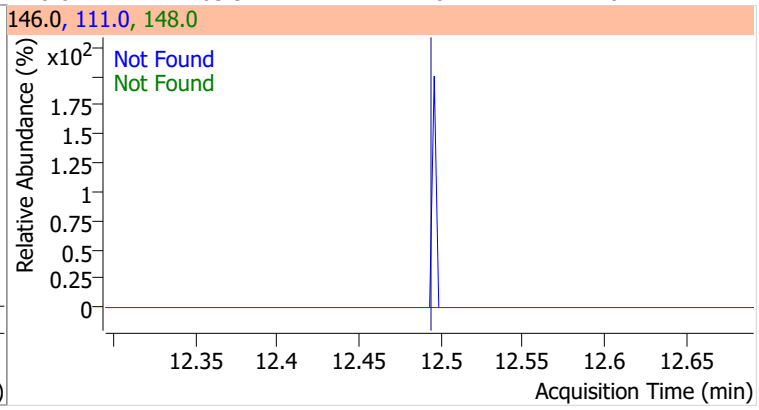
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





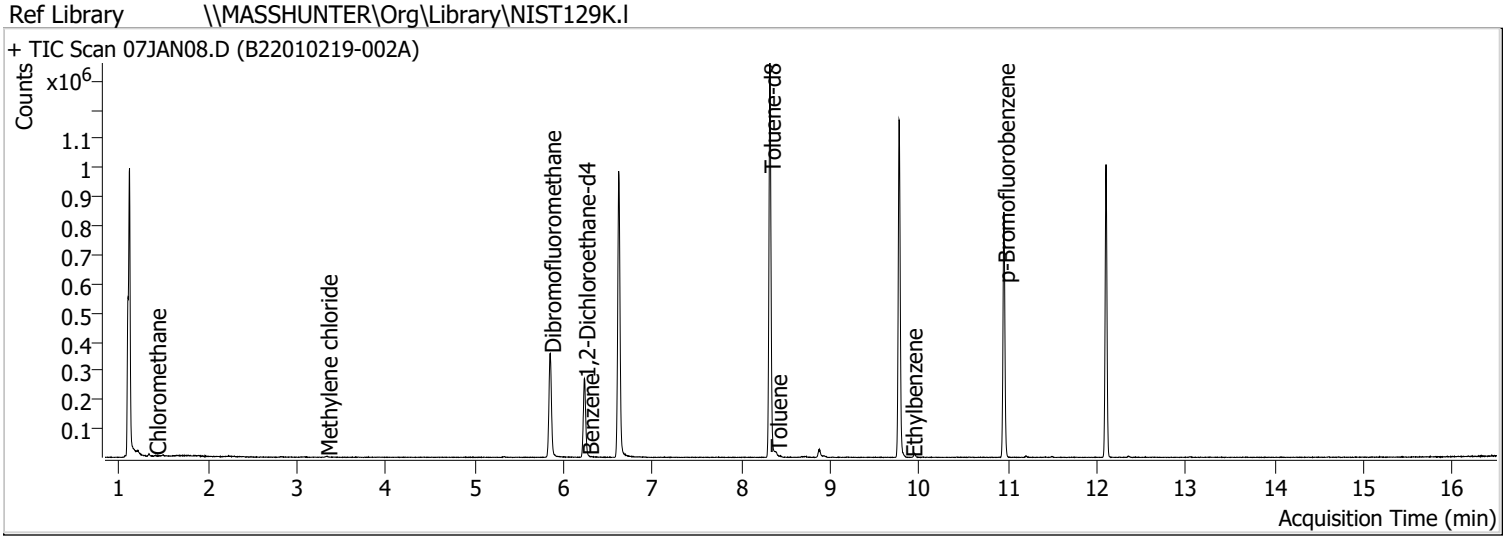
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN08.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 12:52:15 PM
Sample Name	B22010219-002A	Instrument	VOA5975C
Vial	8	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

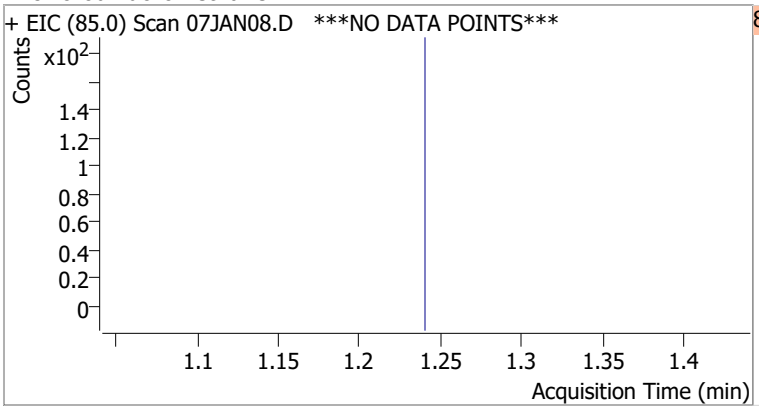


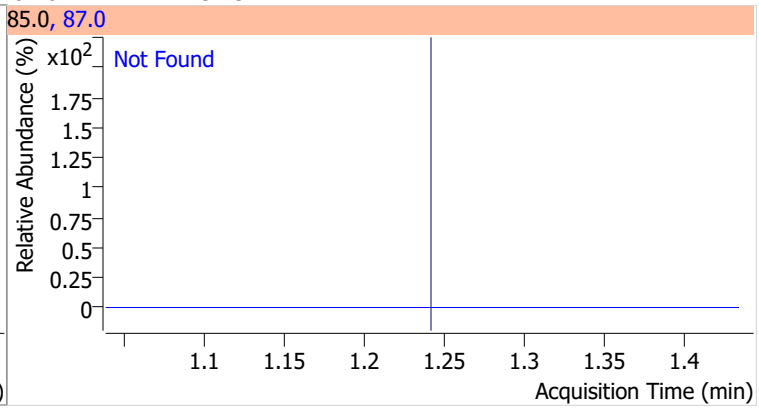
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.618	96.0	819676	250.0000	ng	-0.006
M Chlorobenzene-d5	9.772	82.0	322006	250.0000	ng	0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	235354	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.848	113.0	212882	275.6761	ng	0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 110.27%		
S 1,2-Dichloroethane-d4	6.233	67.0	96571	289.5314	ng	0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 115.81%		
S Toluene-d8	8.319	98.0	818871	263.8953	ng	0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 105.56%		
S p-Bromofluorobenzene	10.951	95.0	240793	279.2702	ng	-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 111.71%		
Target Compounds						**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.		
T Chloromethane	1.406	50.0	2485	1.9061	ng	96
T Vinyl chloride	0.000		0	N.D.		
T Bromomethane	0.000		0	N.D.		
T Chloroethane	0.000		0	N.D.		
T Trichlorofluoromethane	0.000		0	N.D.		
T 1,1-Dichloroethene	0.000		0	N.D.		
T Methylene chloride	3.333	49.0	2231	1.8330	ng	97
T trans-1,2-Dichloroethene	0.000		0	N.D.		
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.		
T 1,1-Dichloroethane	0.000		0	N.D.		
T 2,2-Dichloropropane	0.000		0	N.D.		
T cis-1,2-Dichloroethene	0.000		0	N.D.		
T Methyl ethyl ketone	0.000		0	N.D.		
T Bromochloromethane	0.000		0	N.D.		
T Chloroform	0.000		0	N.D.		

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.275	78.0	212	0.0648	ng	m	67
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	0.000		0	N.D.			
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.388	92.0	3792	1.8091	ng		95
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	0.000		0	N.D.			
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	9.911	91.0	483	0.1214	ng	m	53
T m+p-Xylenes	10.037	106.0	0		ng	md	1
T o-Xylene	0.000		0	N.D.			
T Styrene	0.000		0	N.D.			
T Bromoform	0.000		0	N.D.			
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

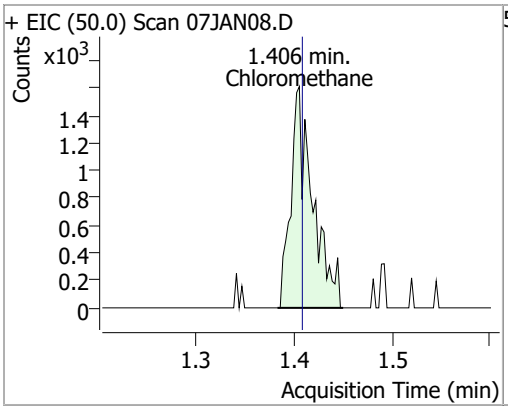
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

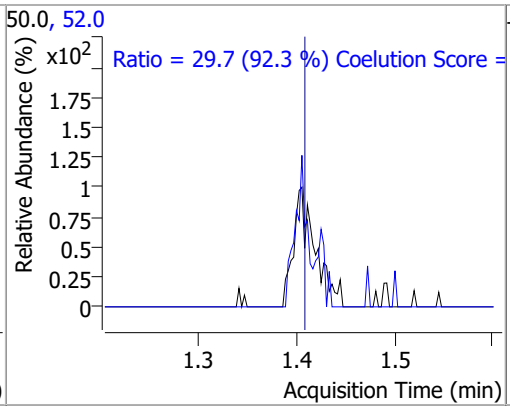
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

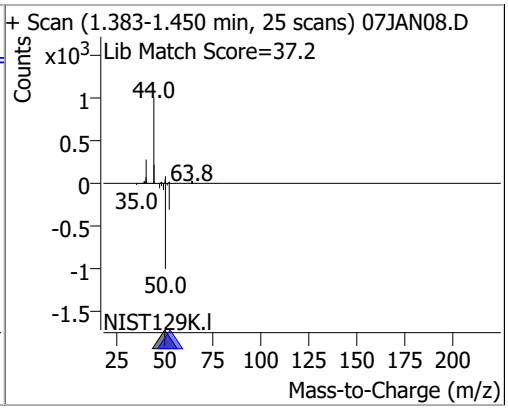




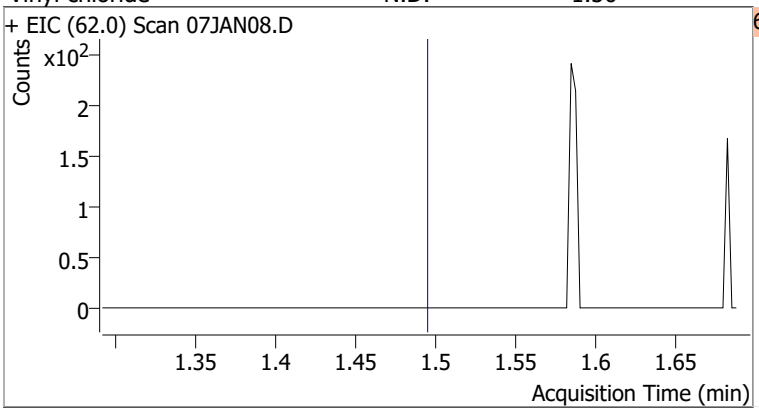
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	1.9061	1.41	0.00	2485	52.0	29.7	2.1	62.1

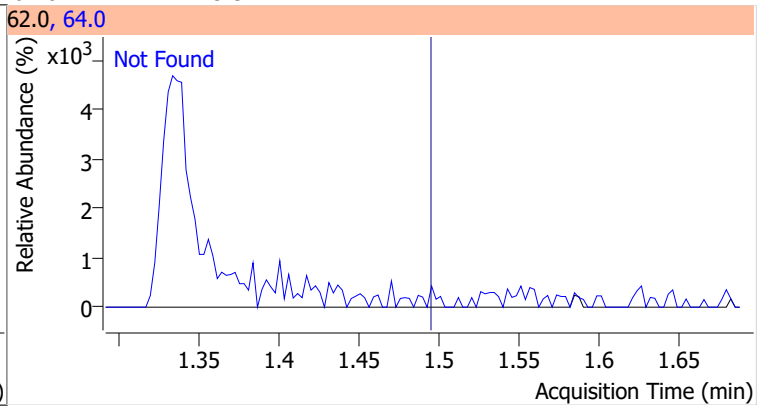




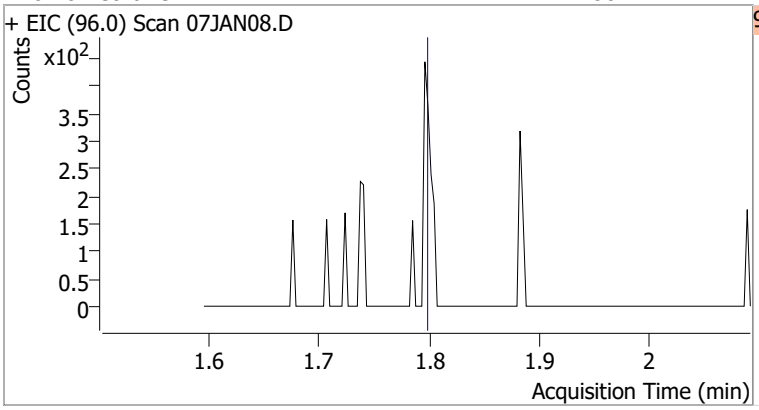


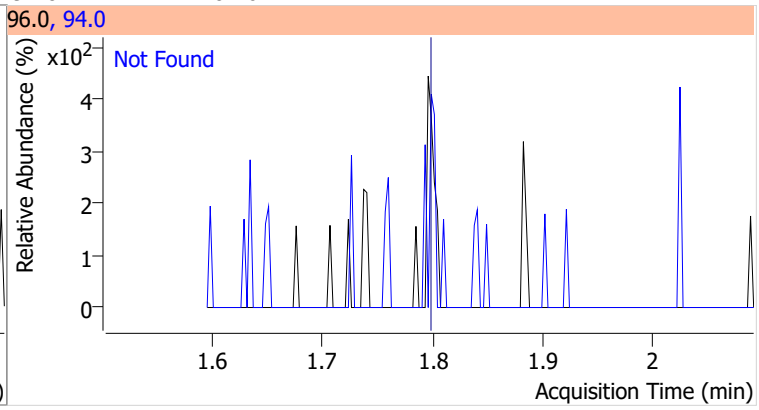
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



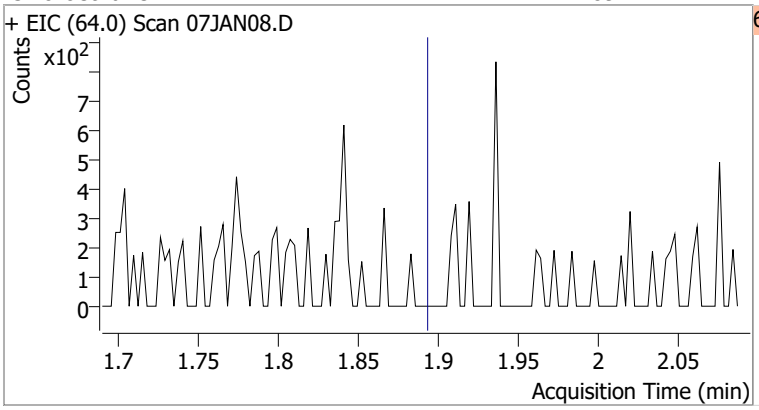


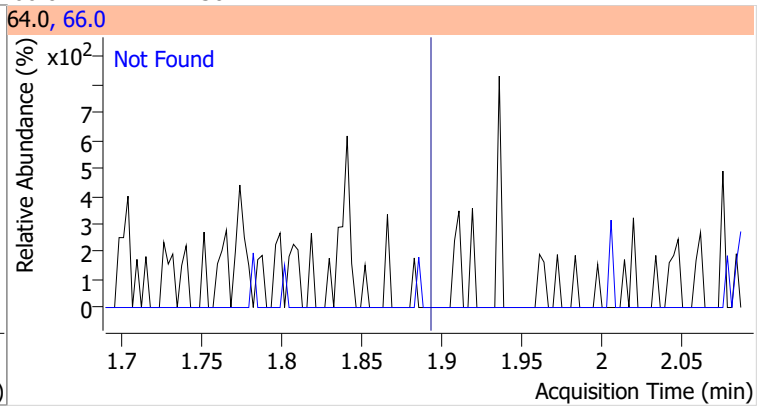
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



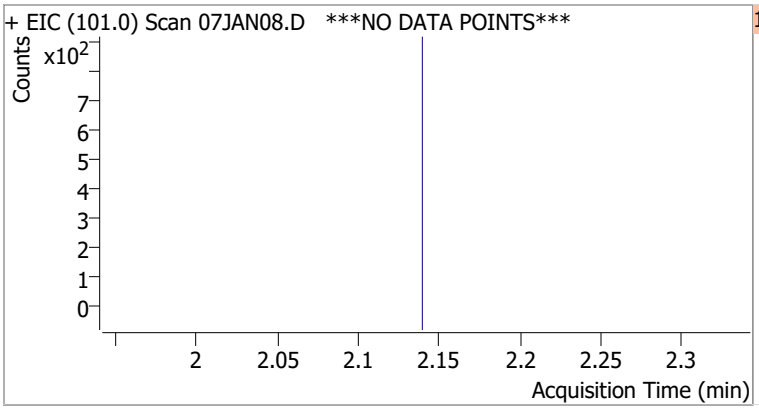


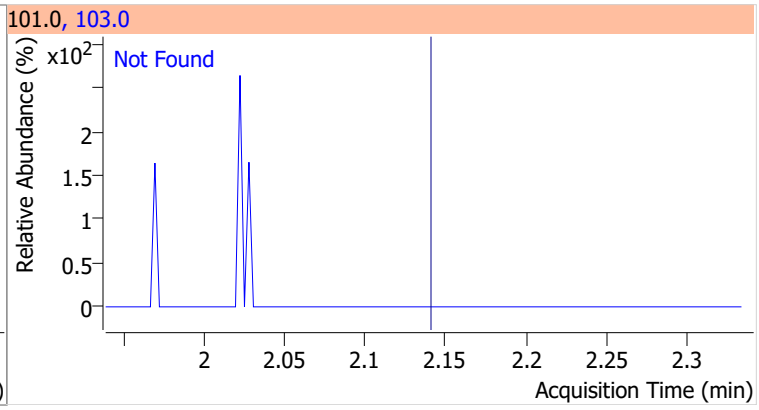
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



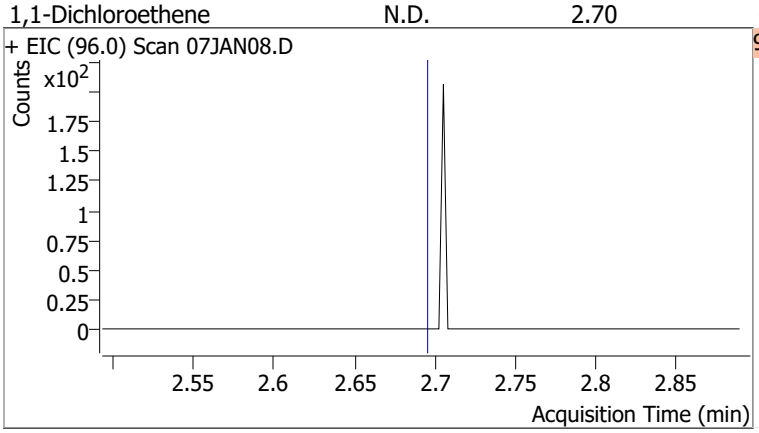


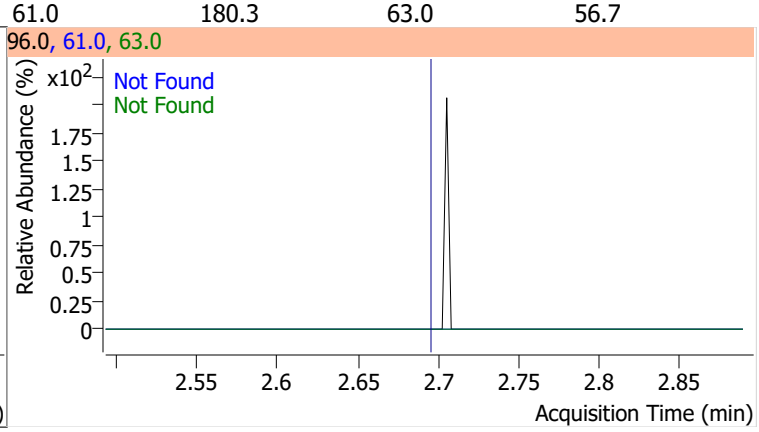
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



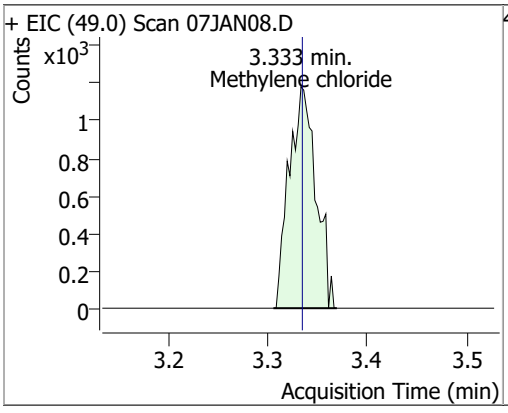


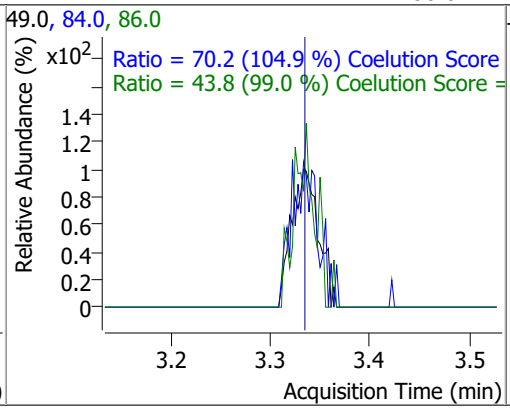
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

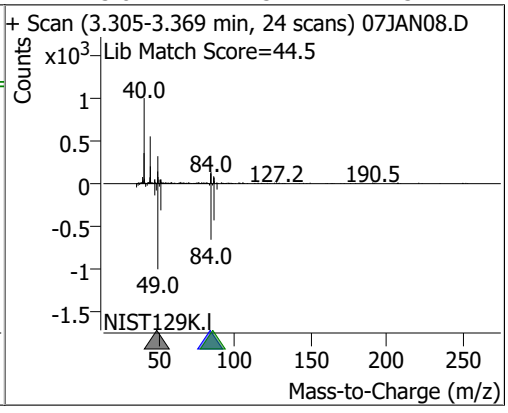




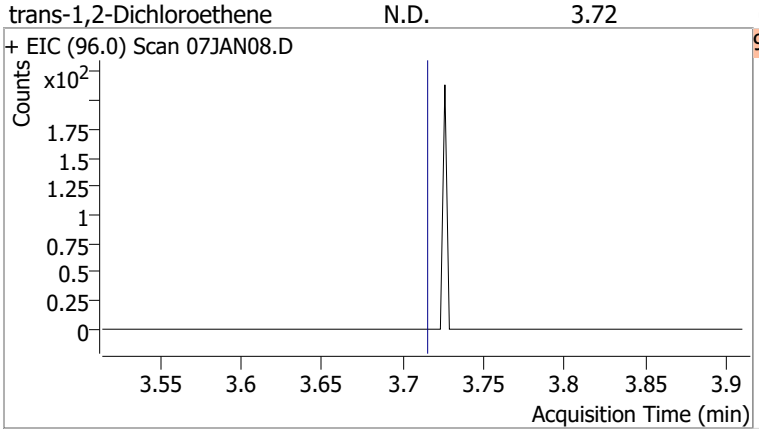
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	1.8330	3.33	0.00	2231	84.0	70.2	36.9	96.9
					86.0	43.8	14.3	74.3

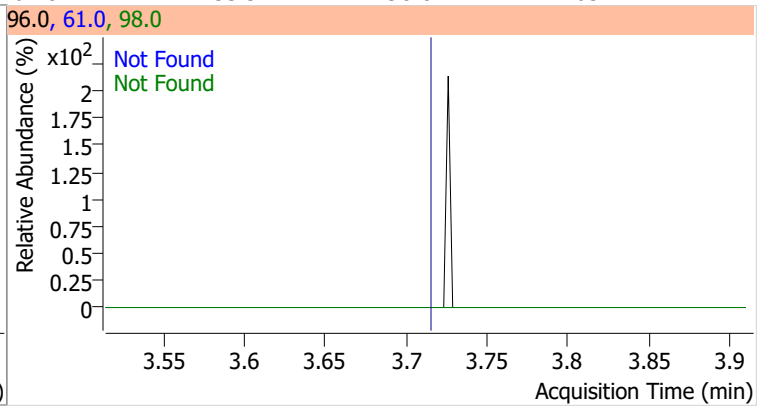




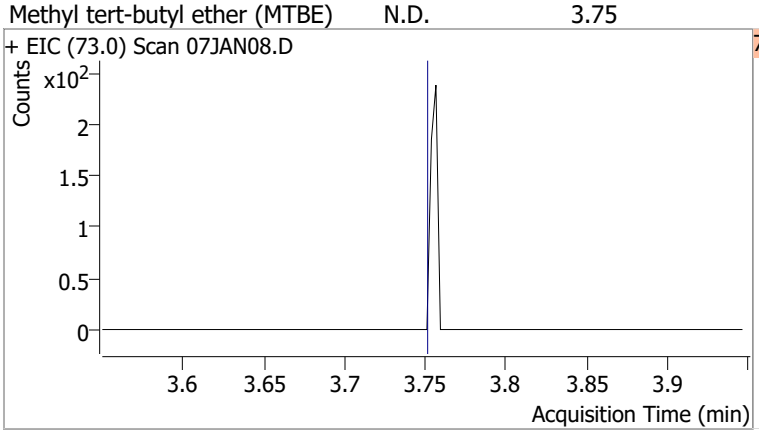


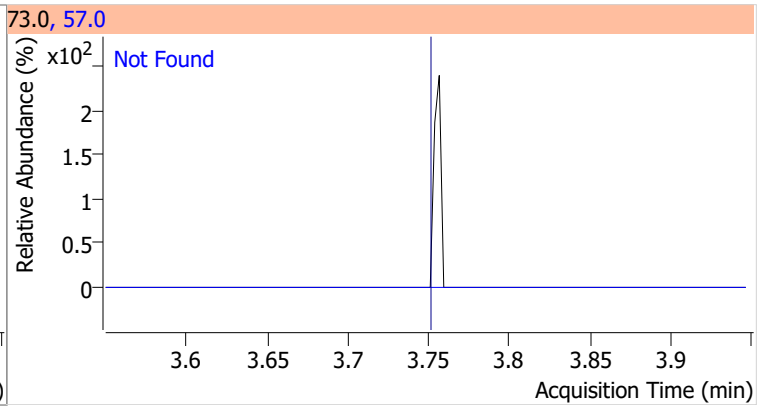
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



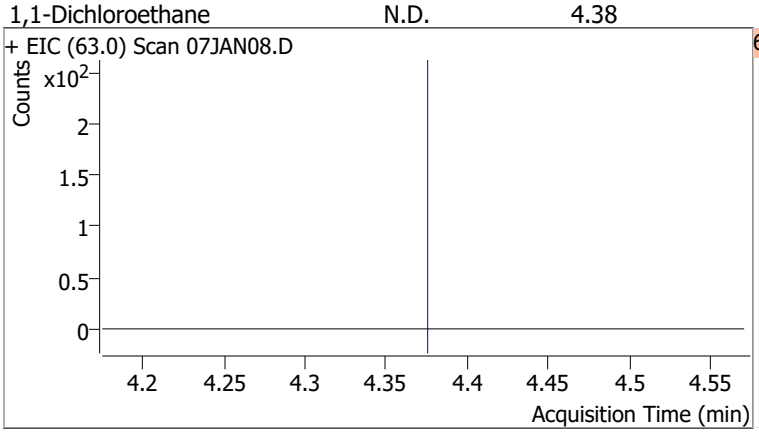


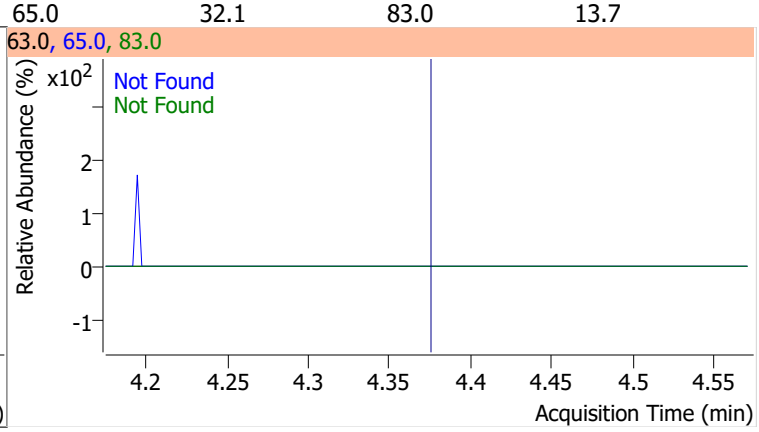
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



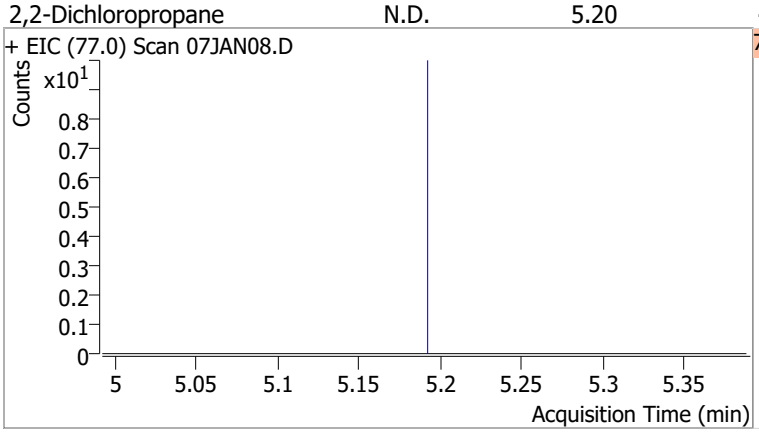


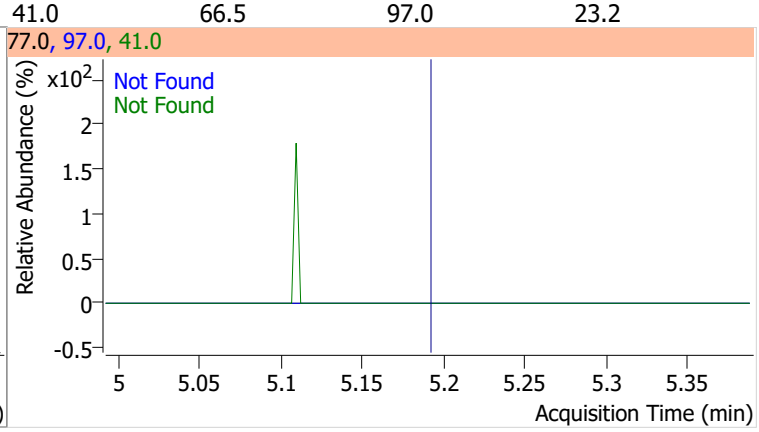
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



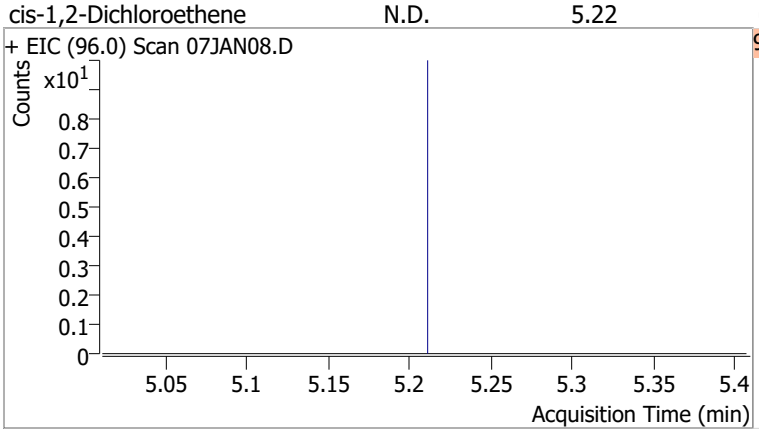


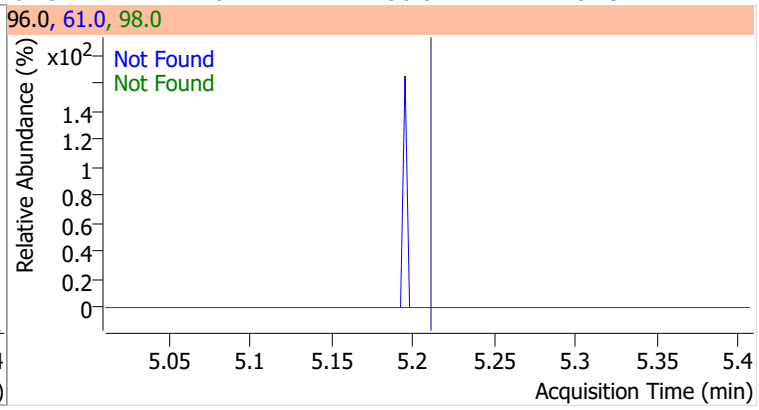
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



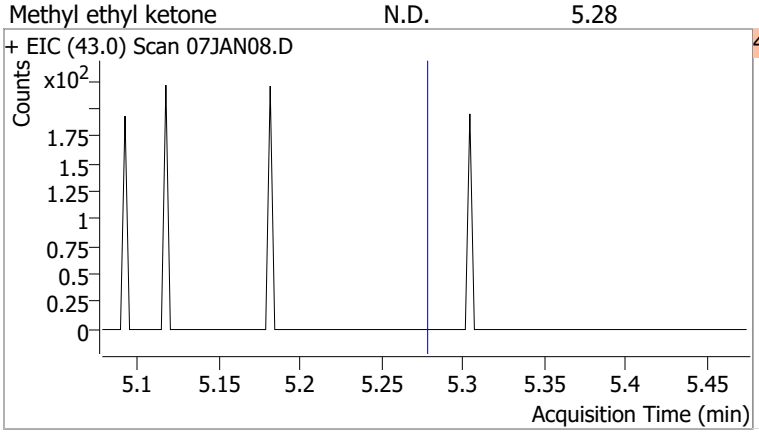


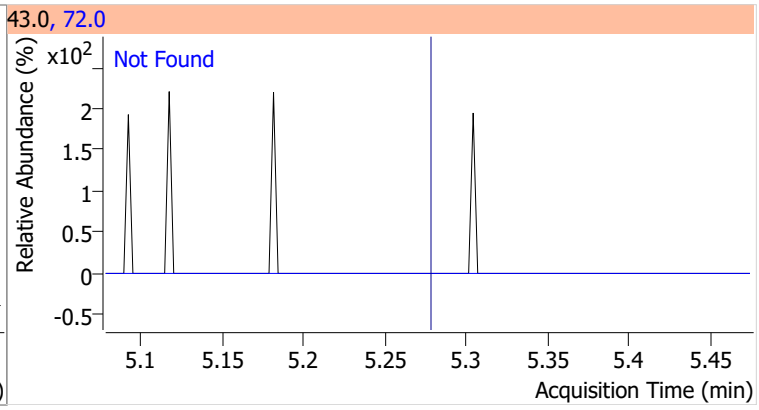
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



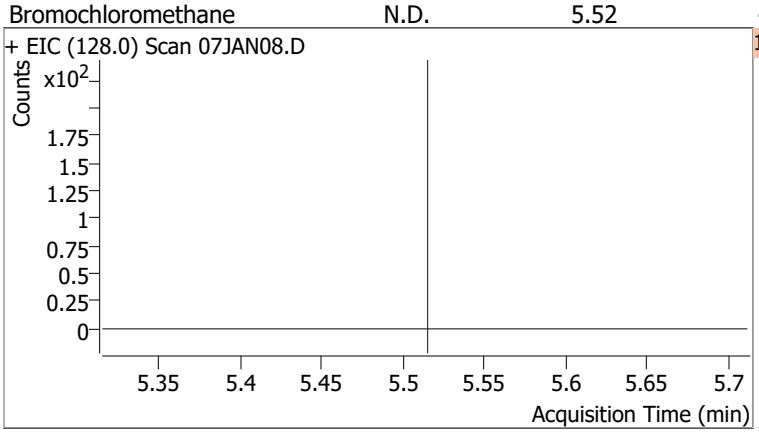


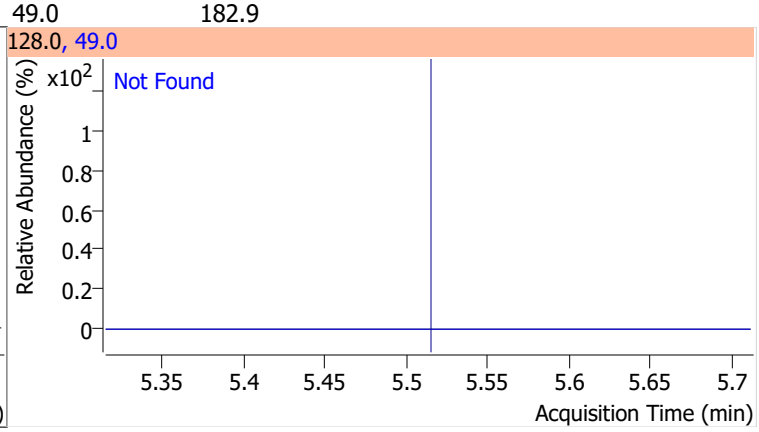
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



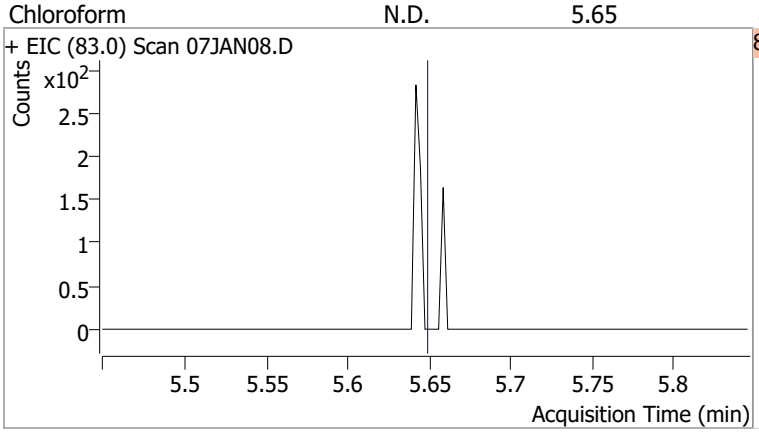


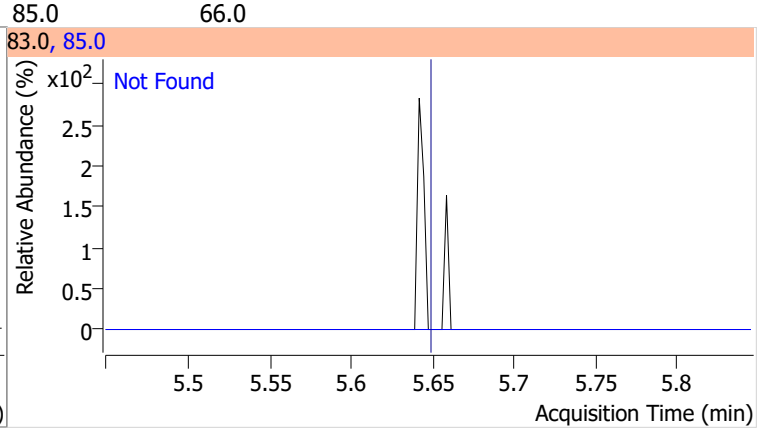
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9



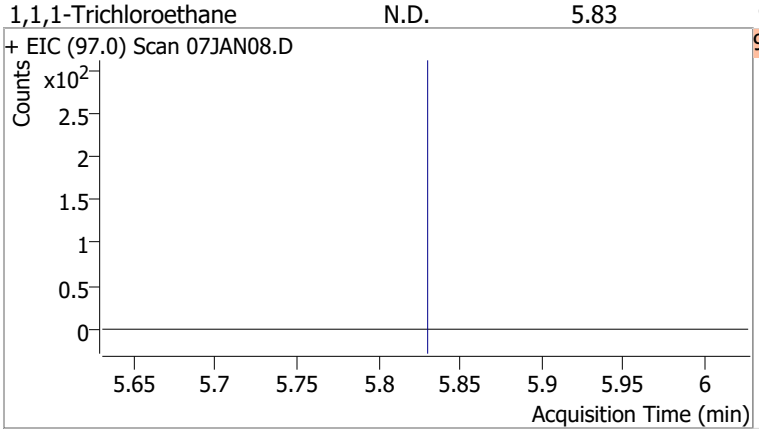


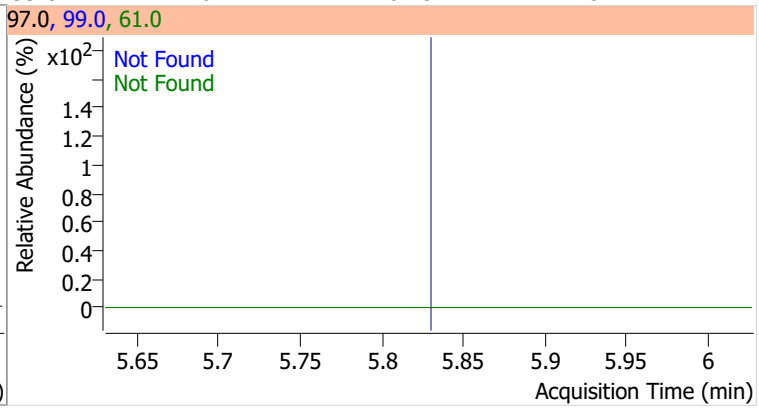
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroform	N.D.	5.65	85.0	66.0



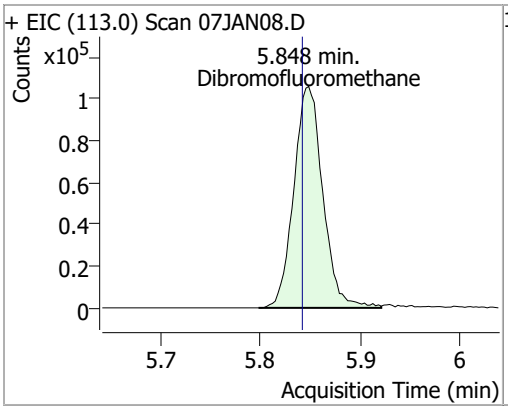


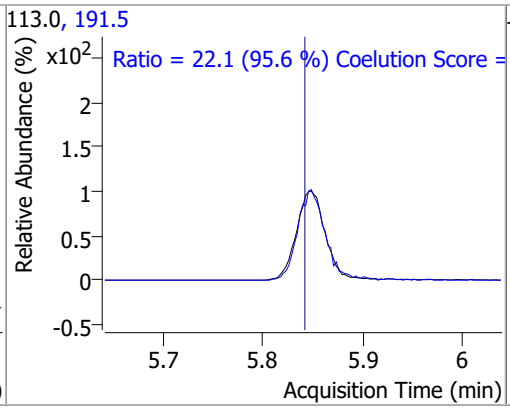
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

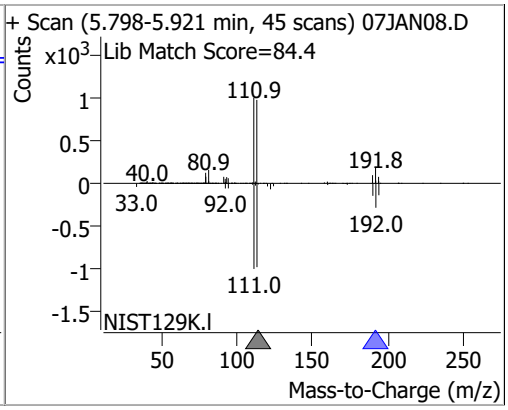




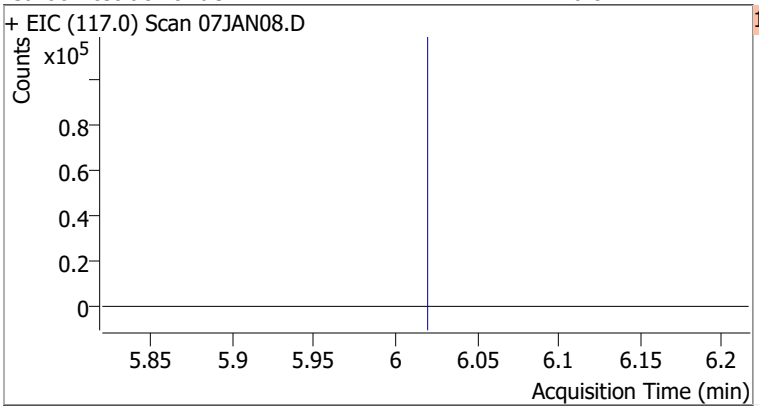
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	275.6761	5.85	0.00	212882	191.5	22.1	0.0	53.1

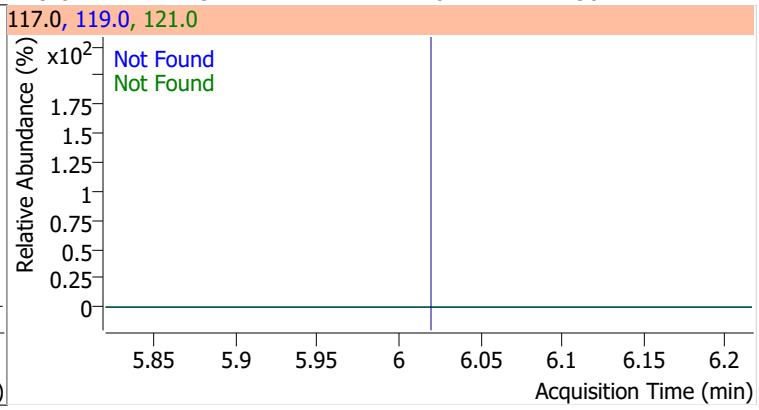




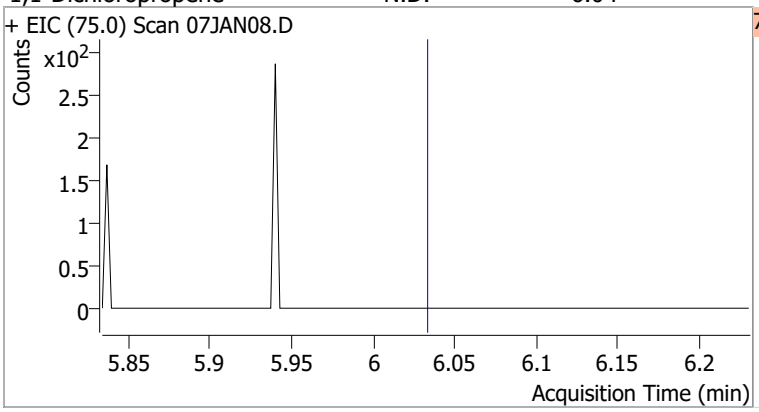


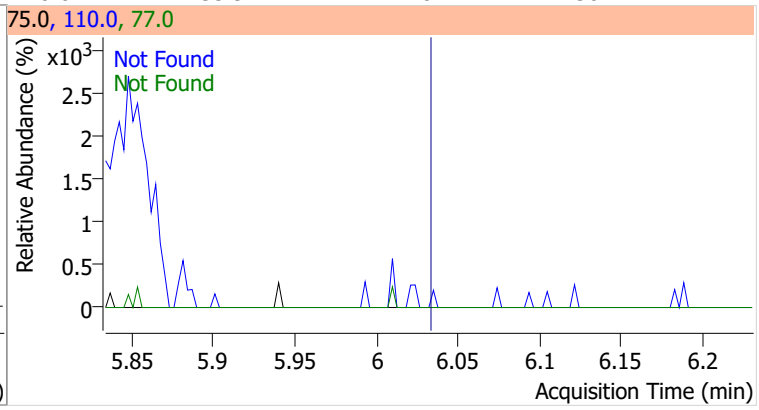
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



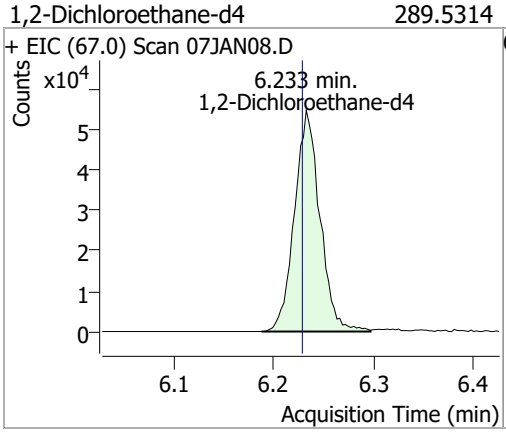


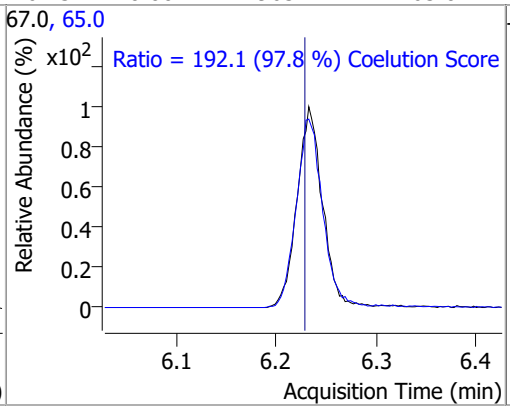
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

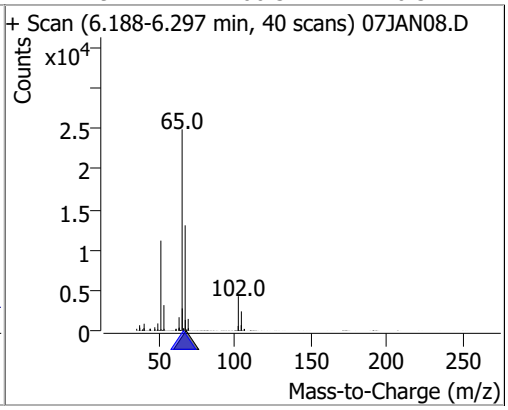




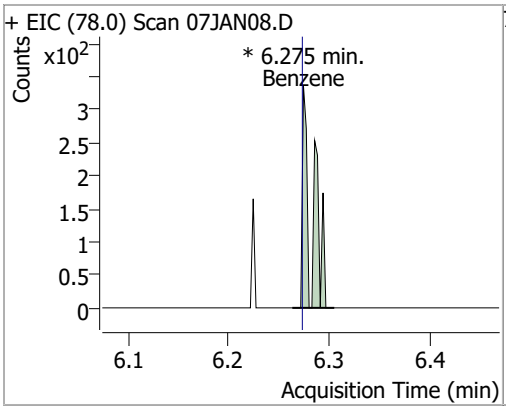
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	289.5314	6.23	0.00	96571	65.0	192.1	166.5	226.5

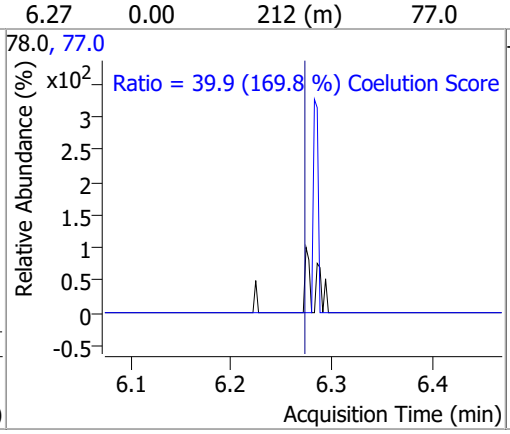


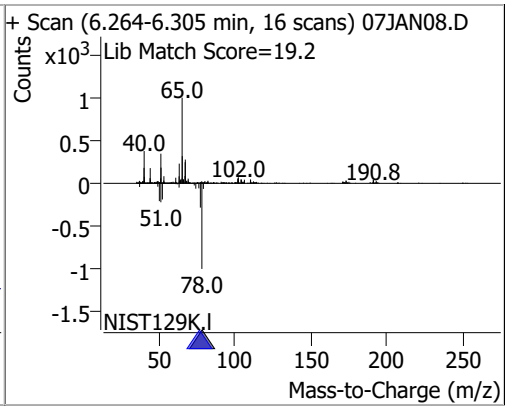




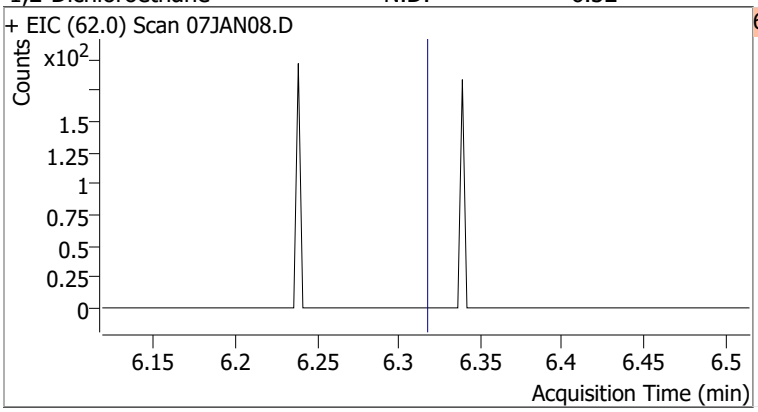
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.0648	6.27	0.00	212 (m)	77.0	39.9	0.0	53.5

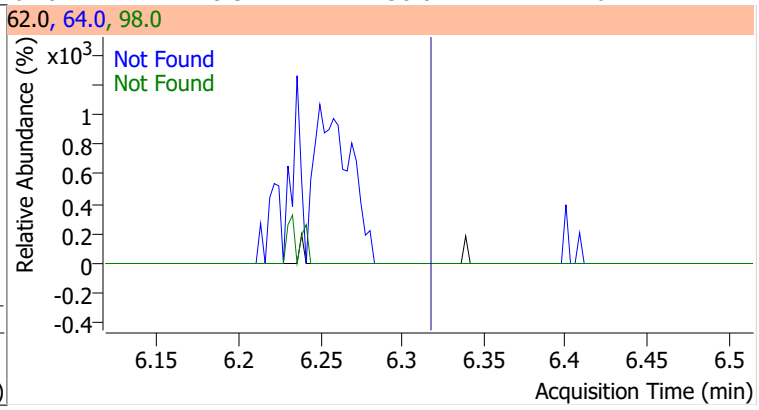




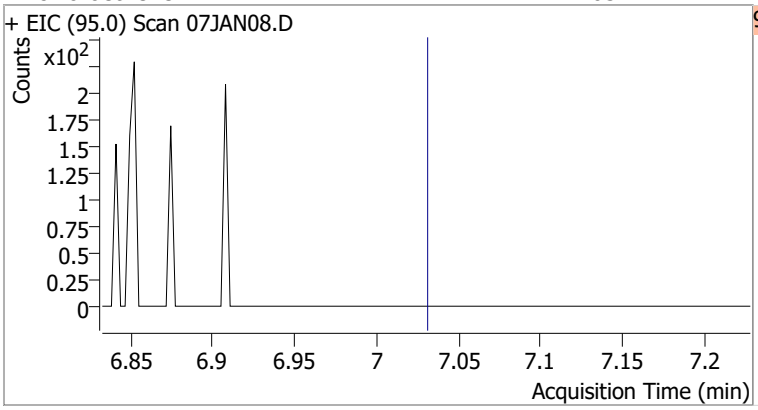


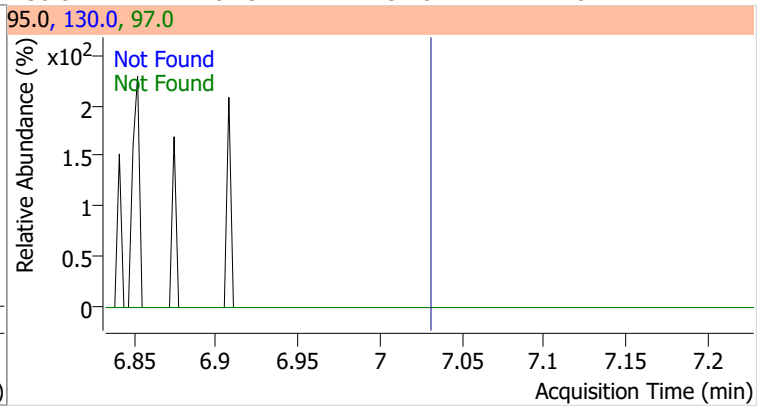
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



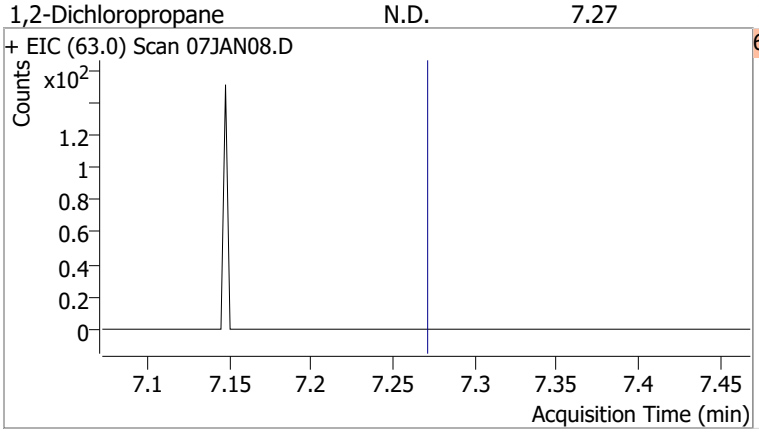


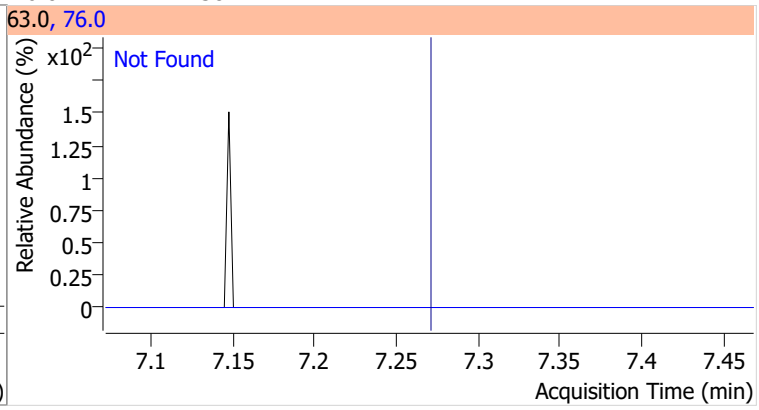
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



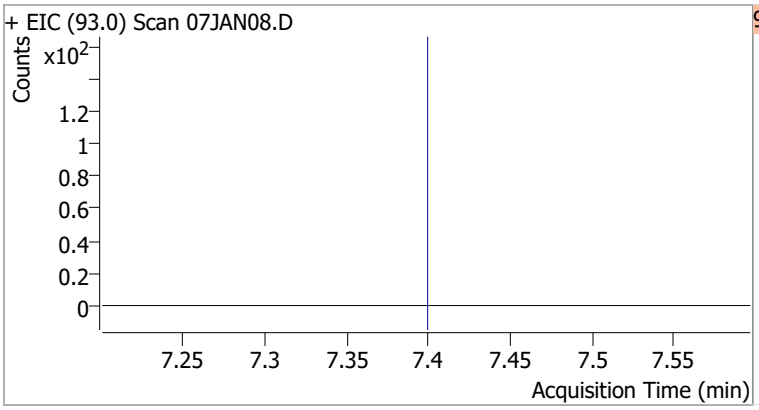


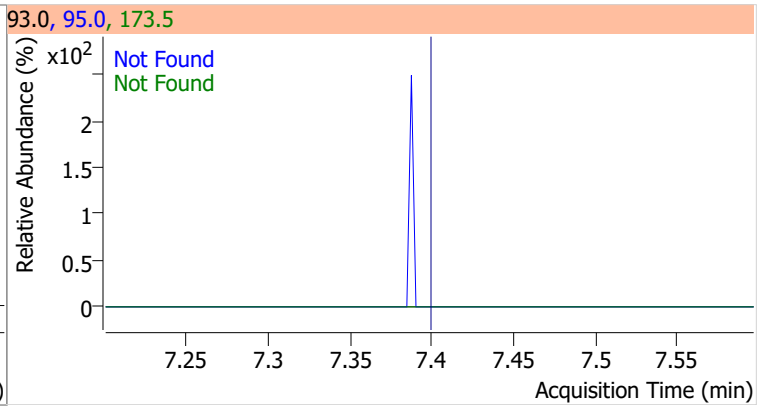
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



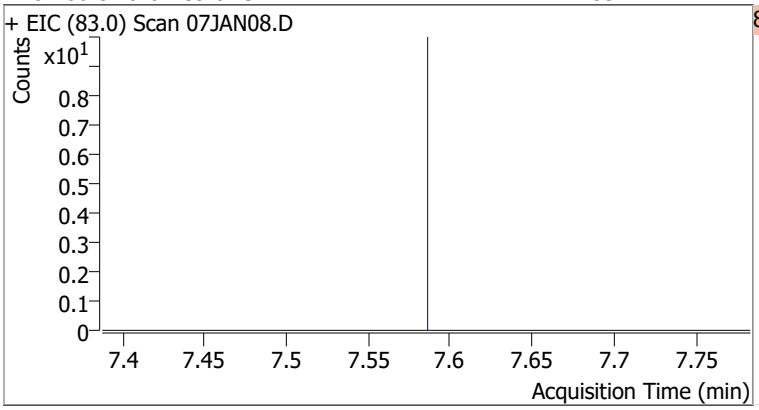


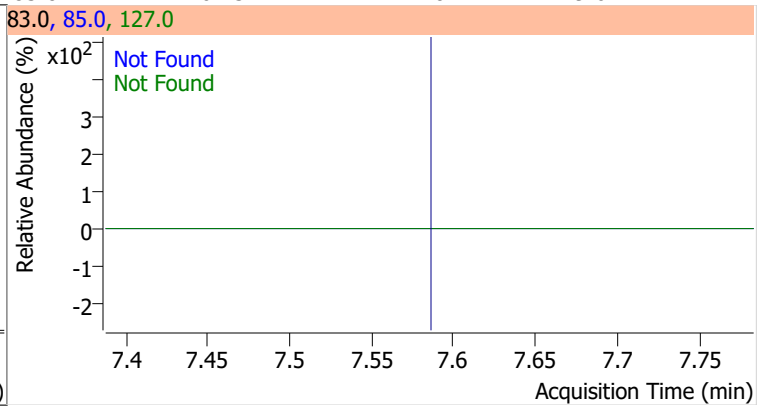
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



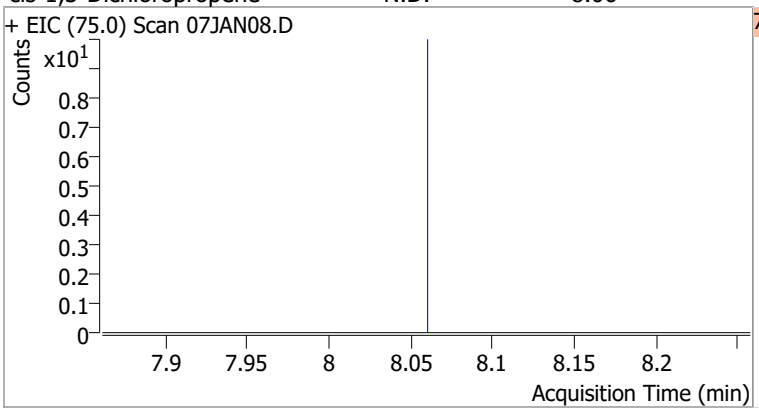


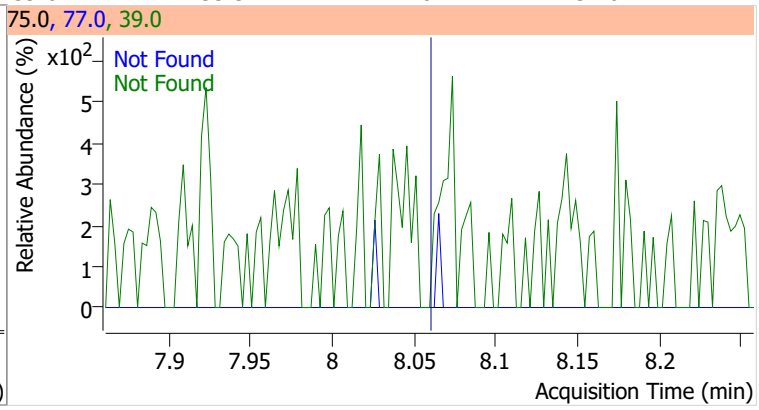
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



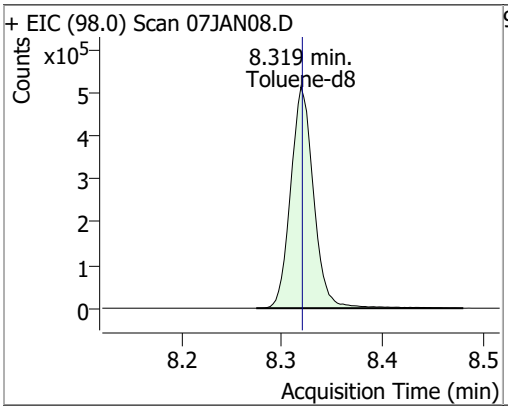


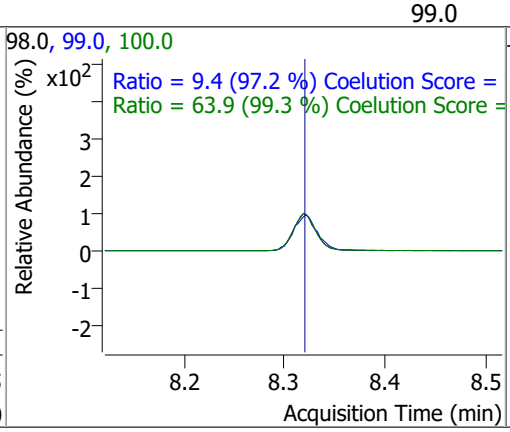
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

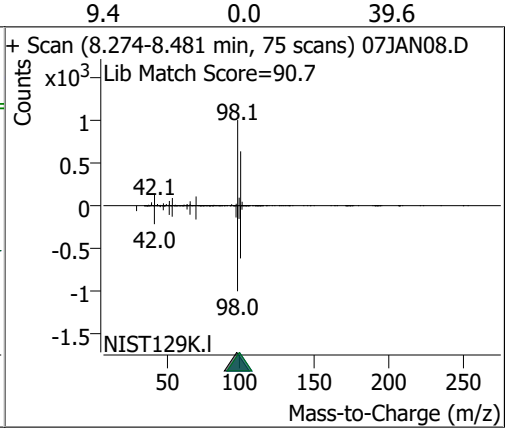




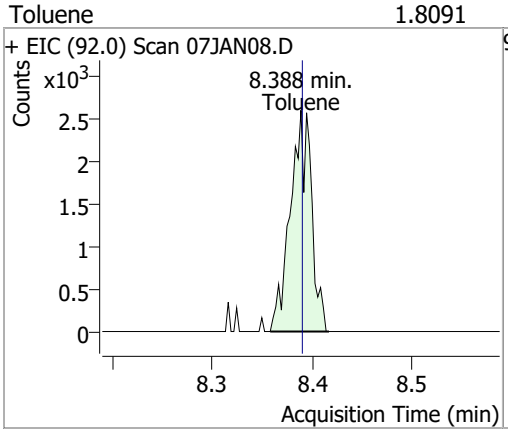
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	263.8953	8.32	0.00	818871	100.0	63.9	34.4	94.4
					99.0	9.4	0.0	39.6

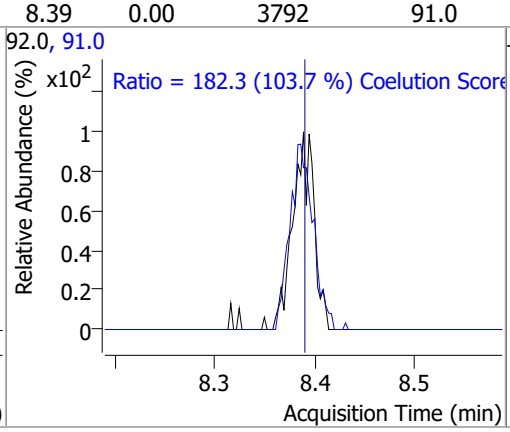


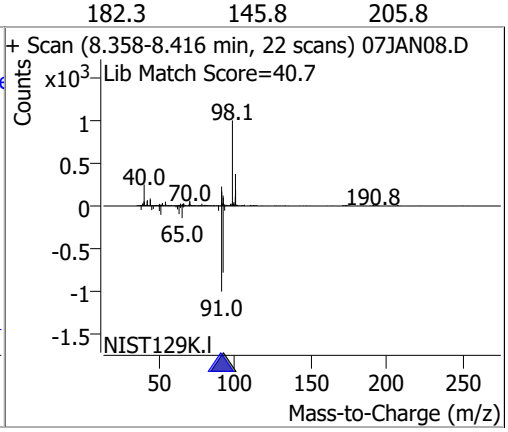




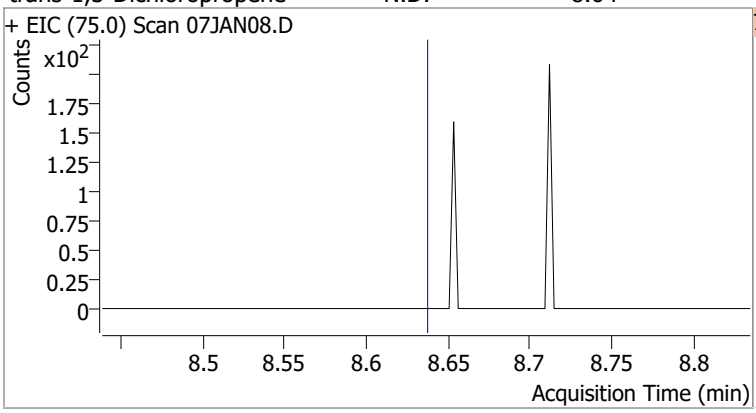
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	1.8091	8.39	0.00	3792	91.0	182.3	145.8	205.8

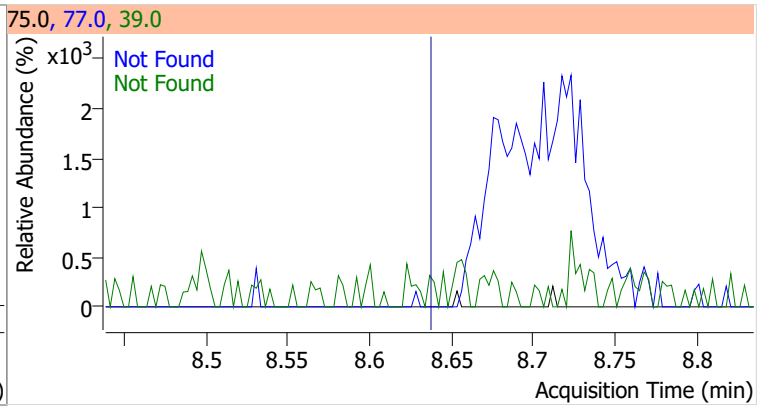




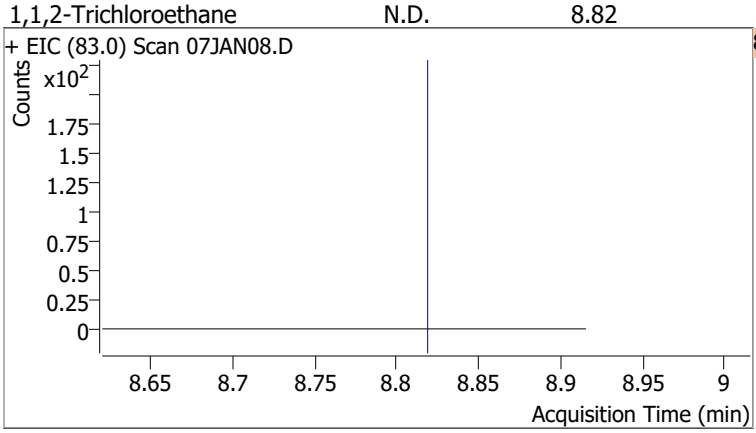


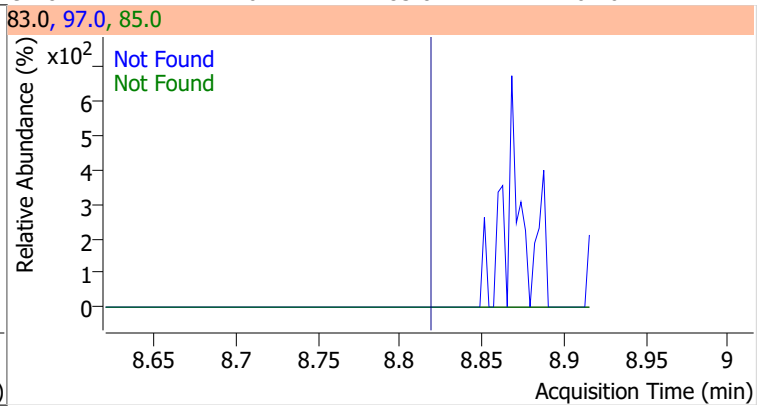
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



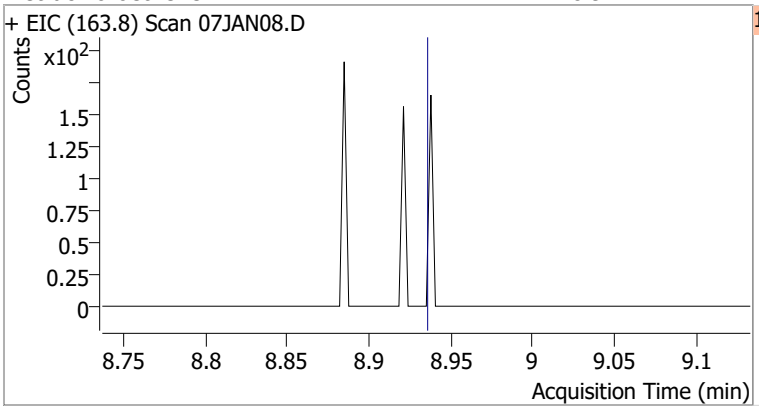


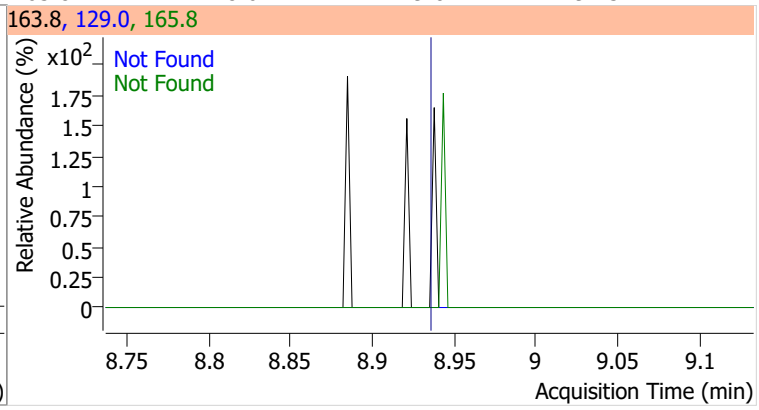
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



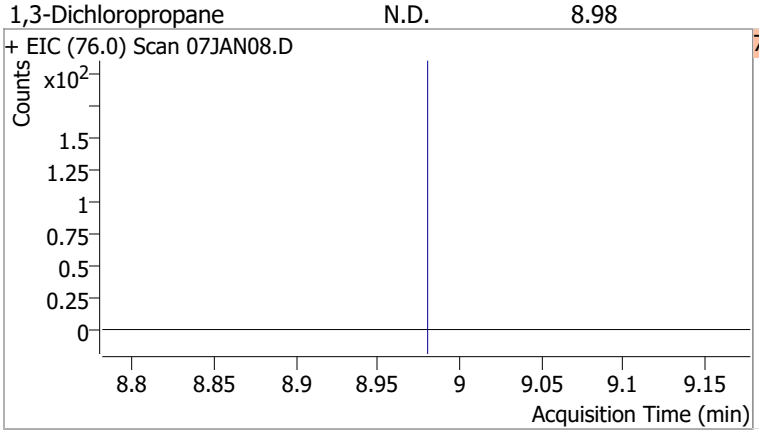


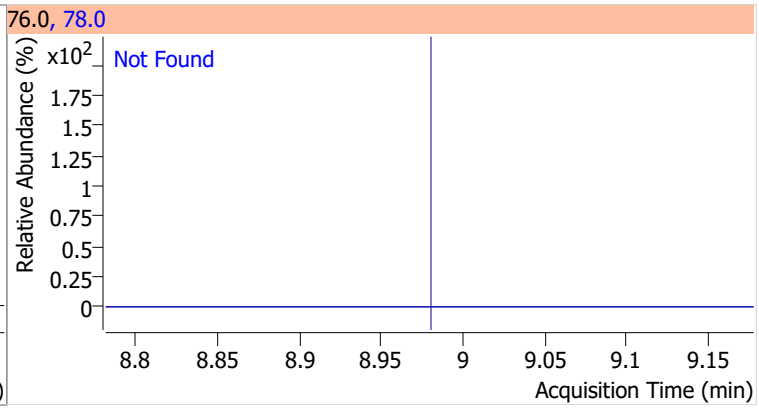
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



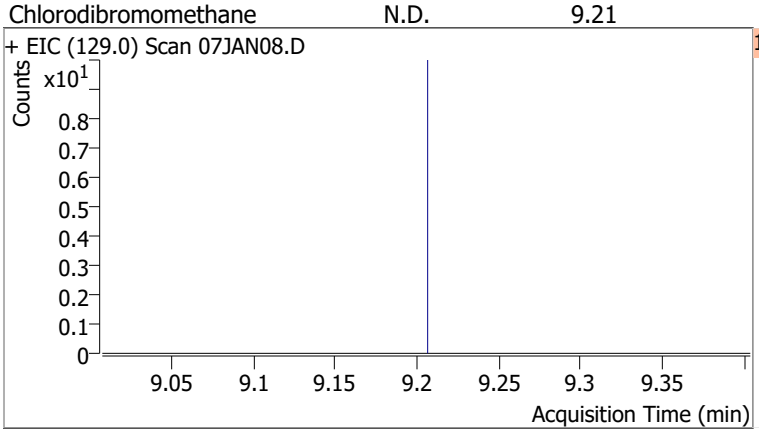


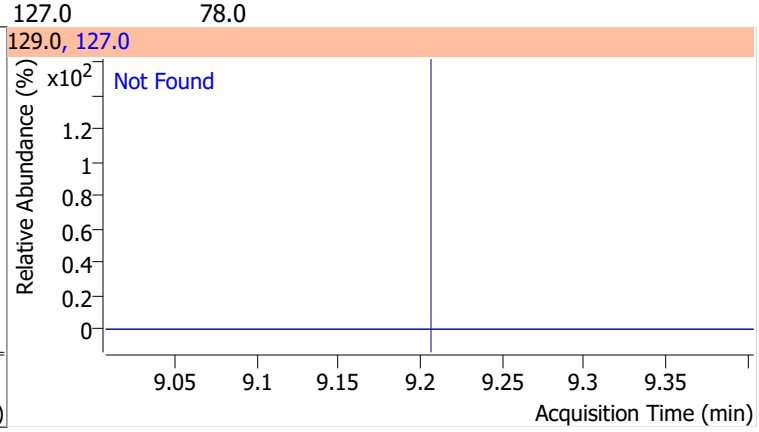
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



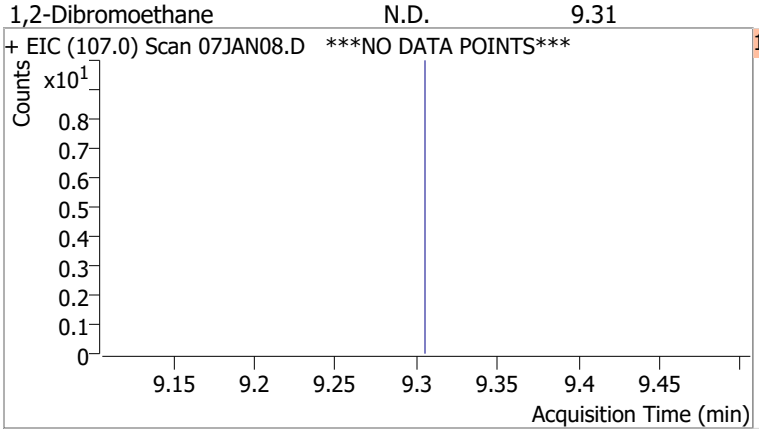


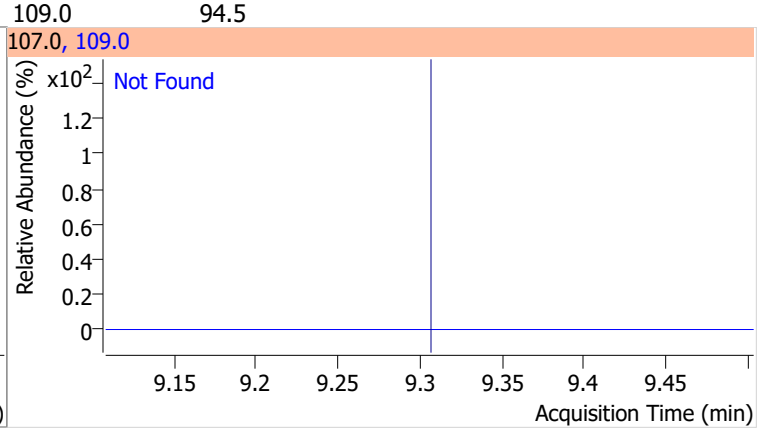
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



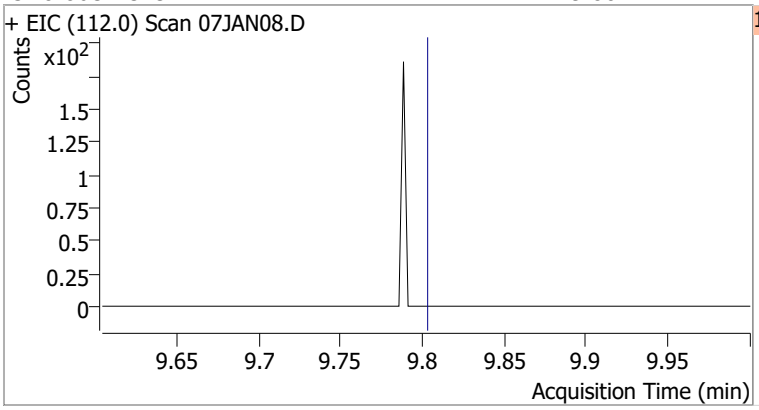


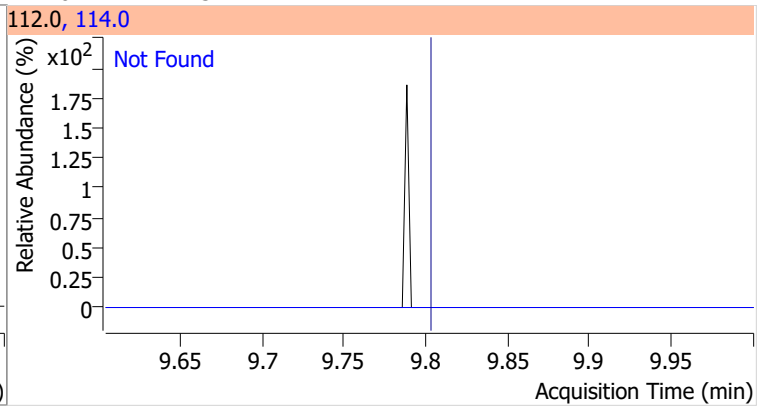
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



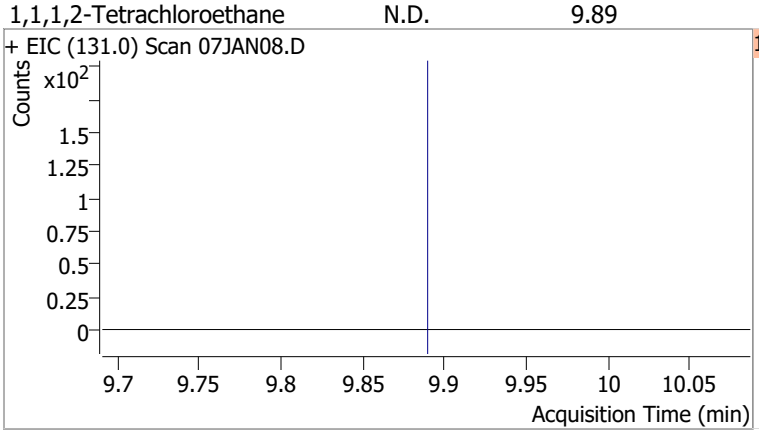


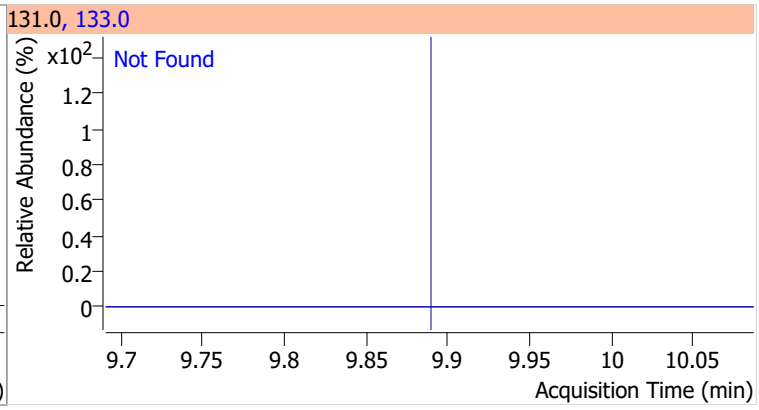
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



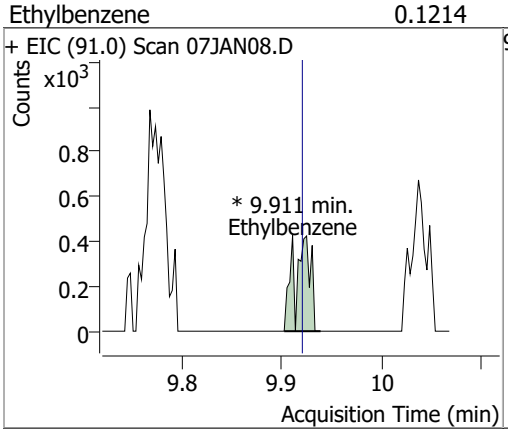


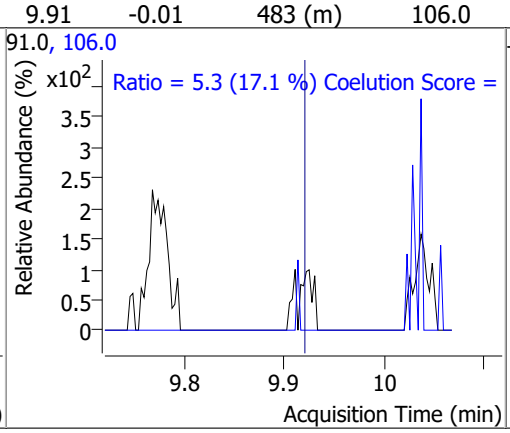
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6

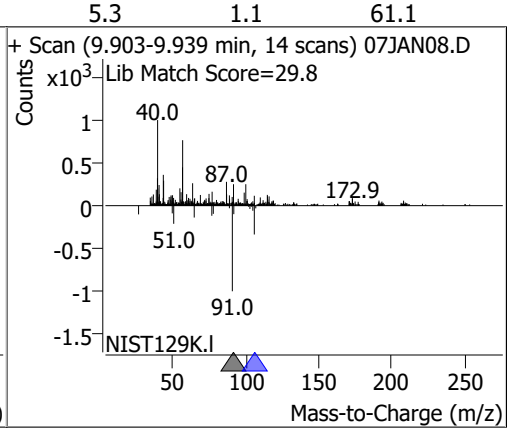




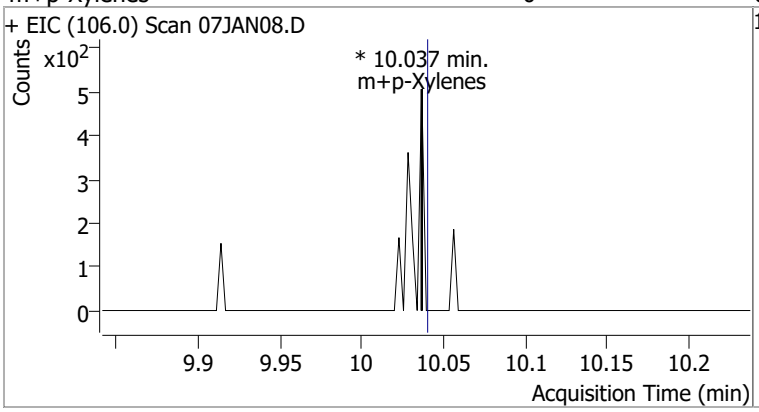
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	0.1214	9.91	-0.01	483 (m)	106.0	5.3	1.1	61.1

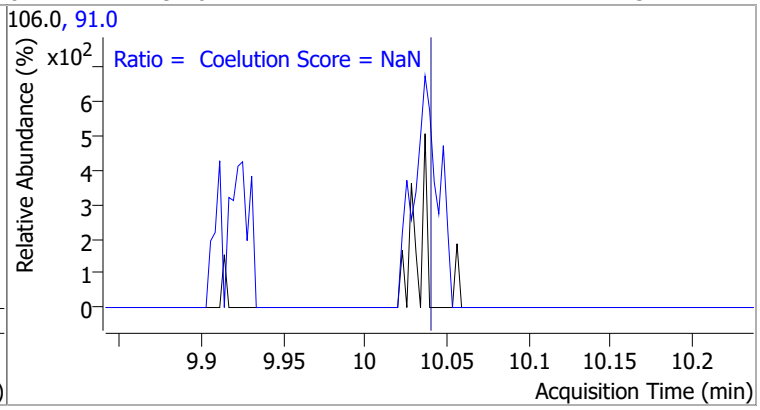




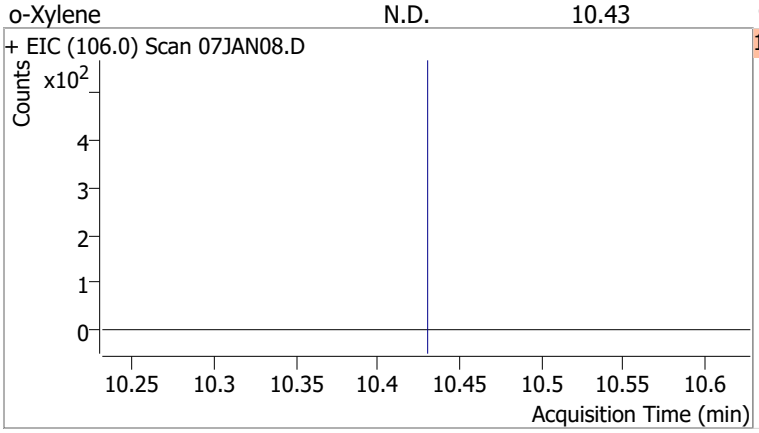


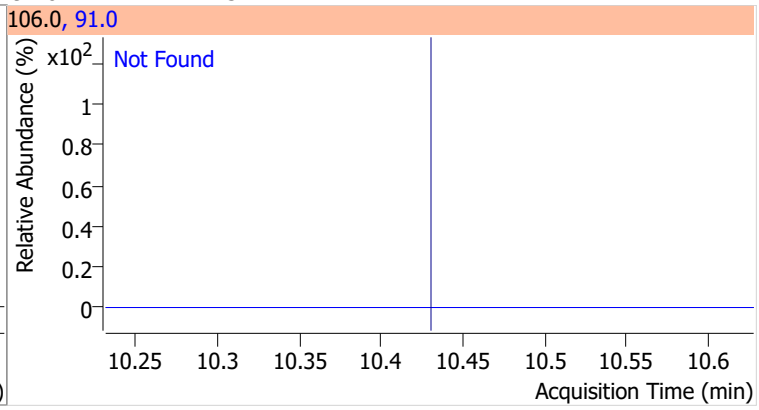
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes		0		0	91.0		171.4	231.4



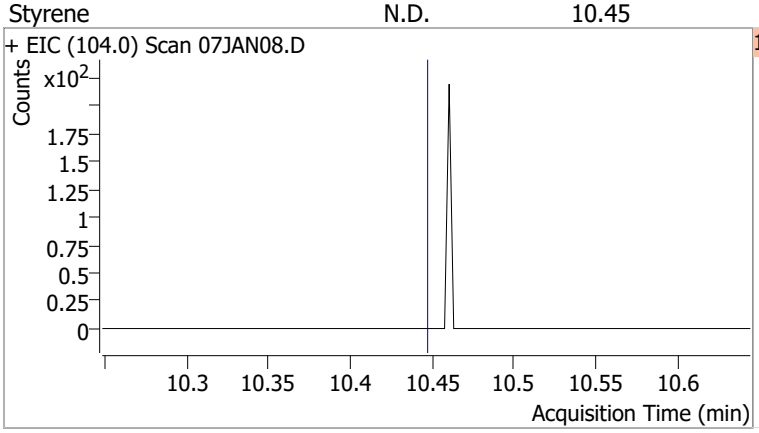


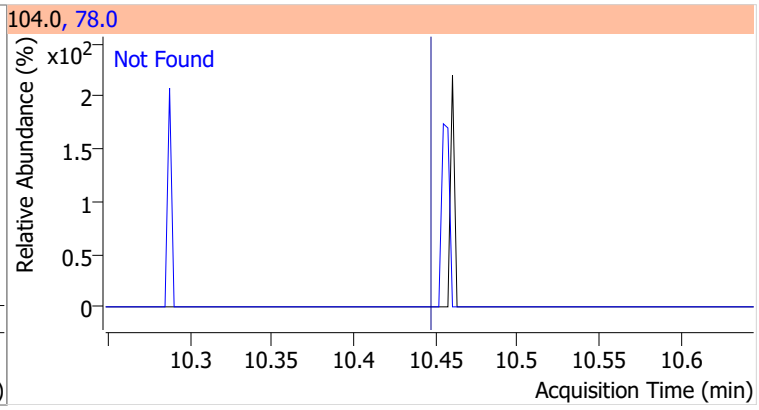
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



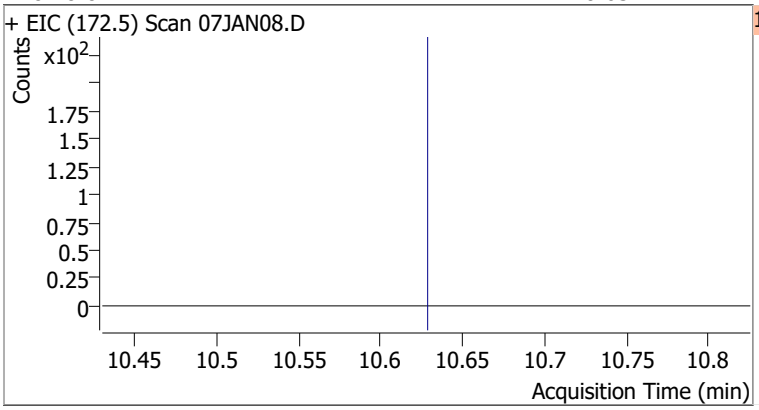


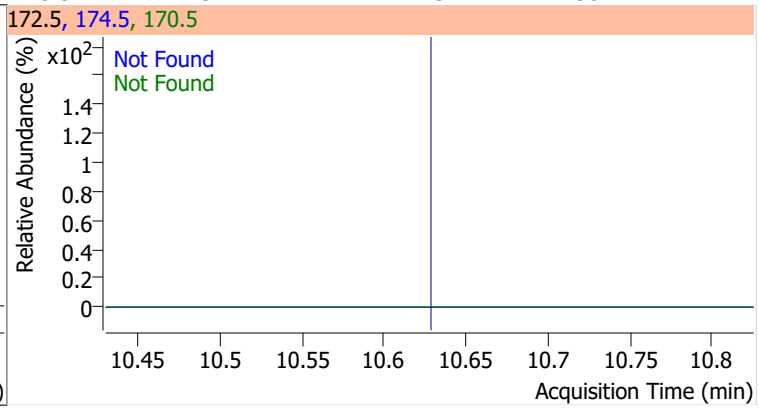
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



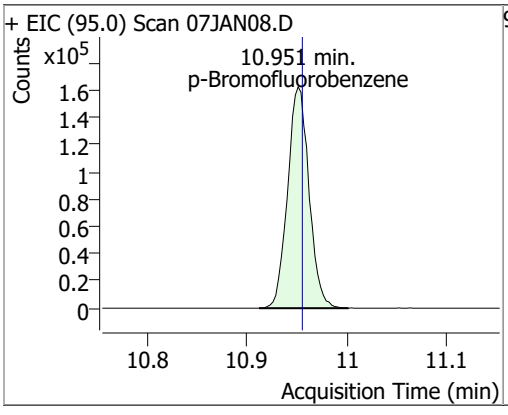


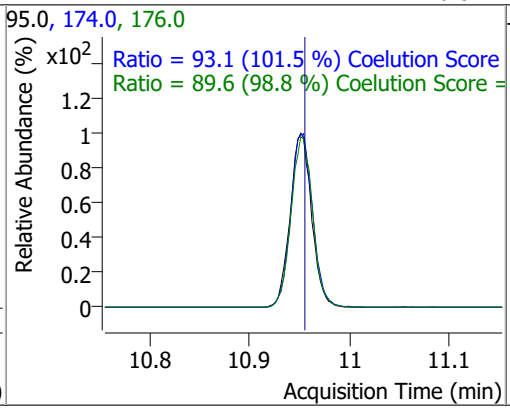
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

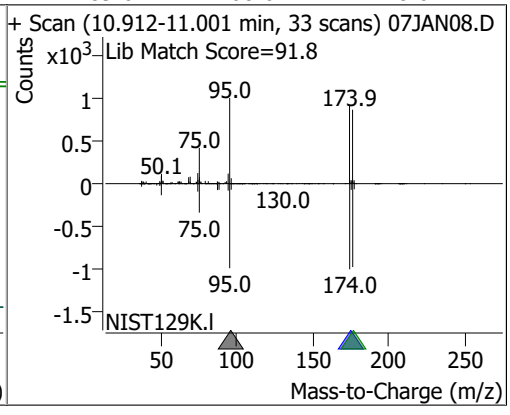




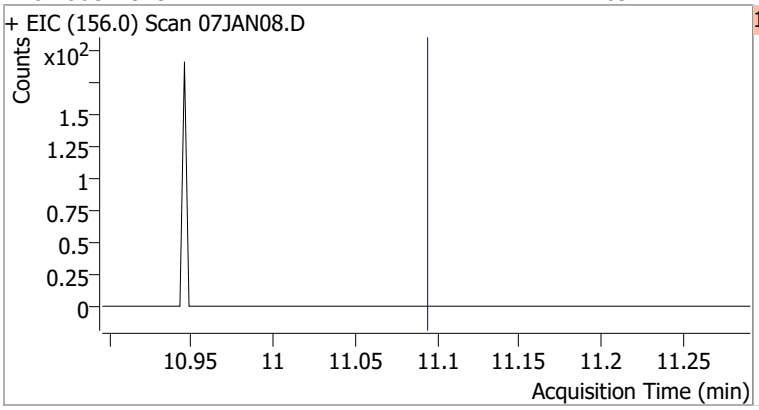
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	279.2702	10.95	0.00	240793	174.0	93.1	61.7	121.7
					176.0	89.6	60.6	120.6

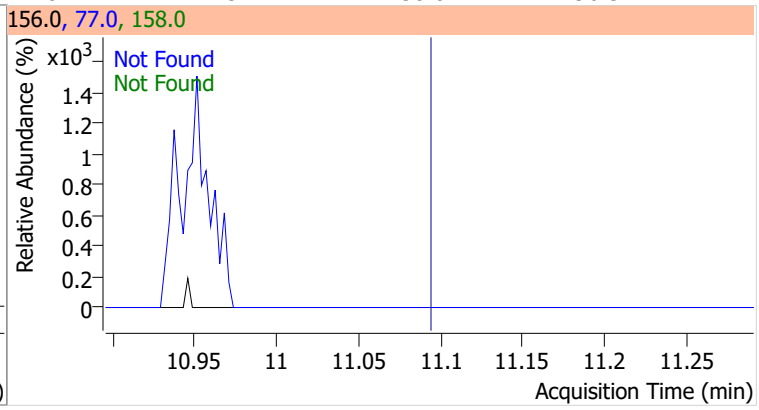




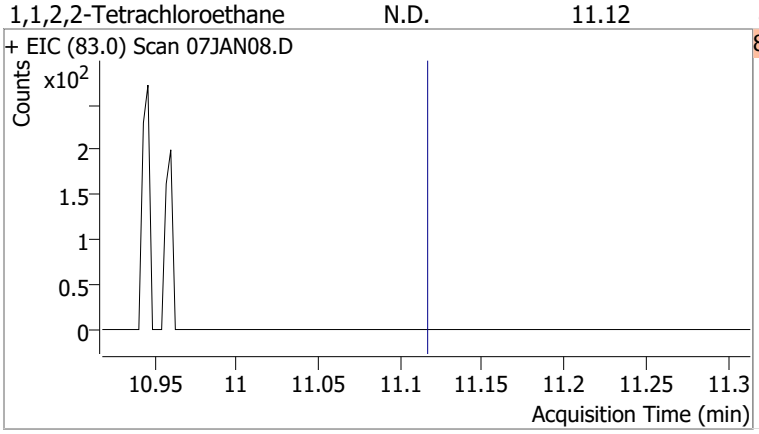


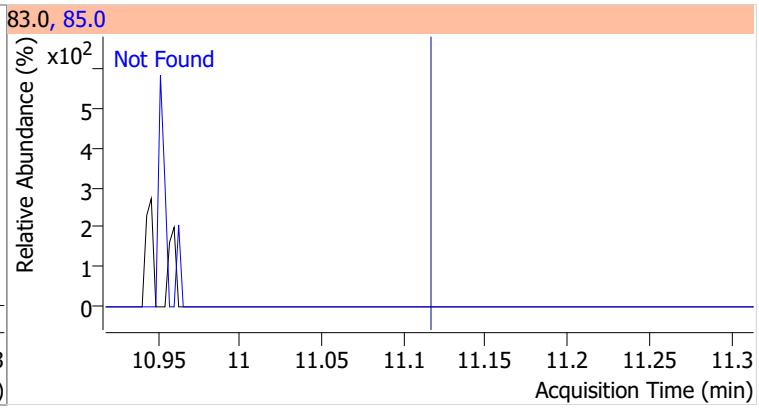
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



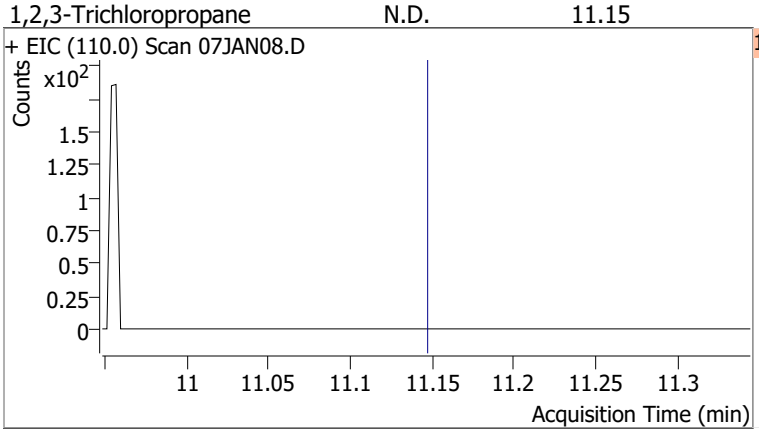


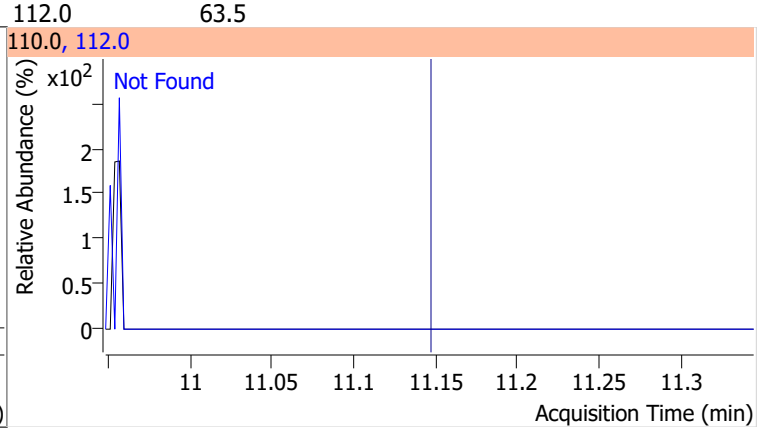
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



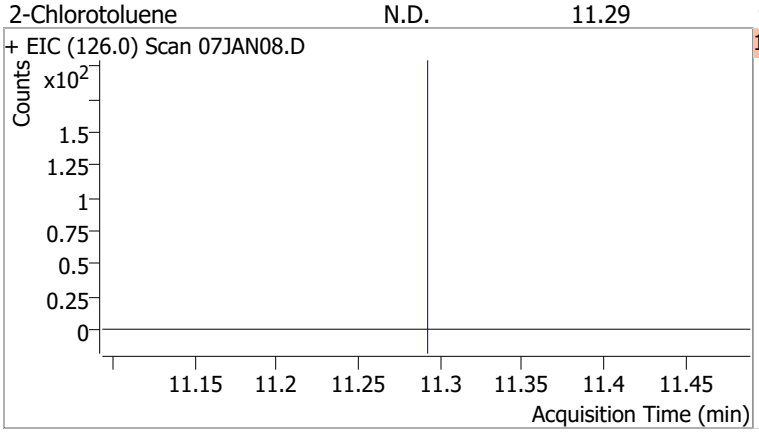


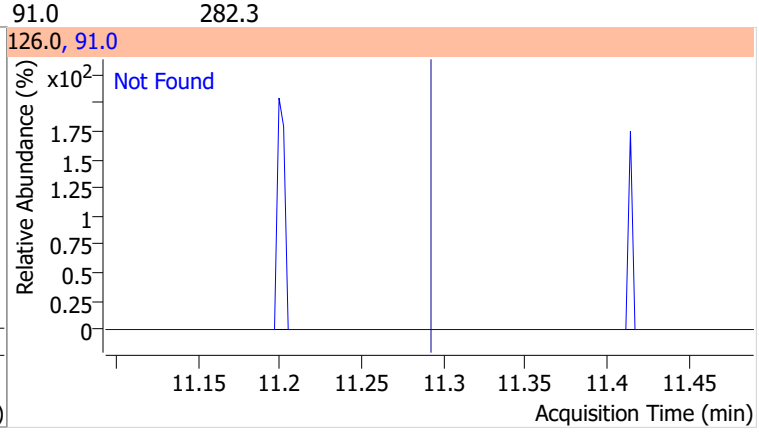
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



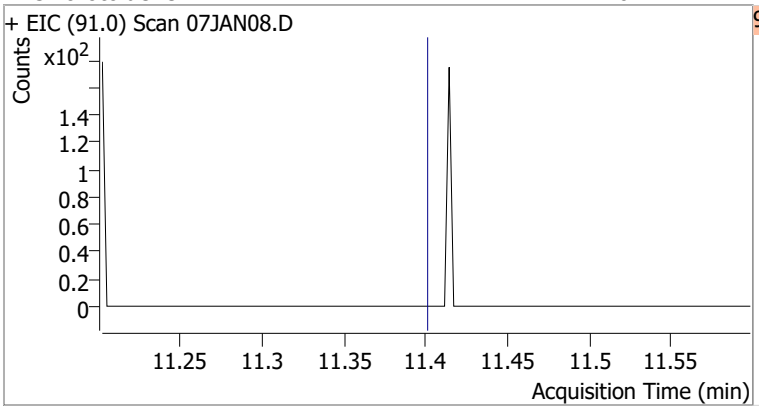


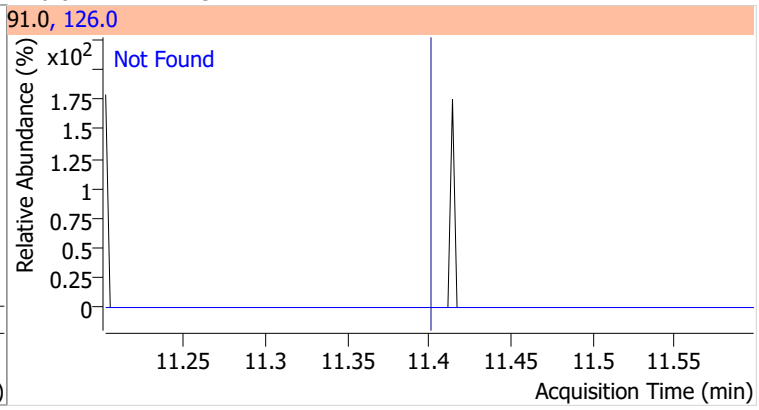
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



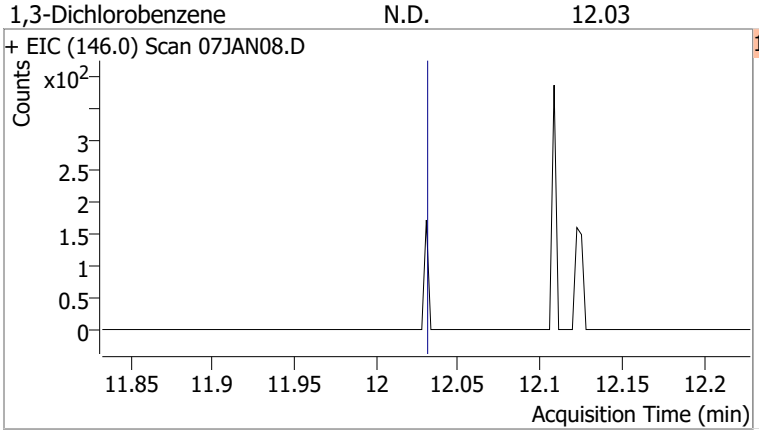


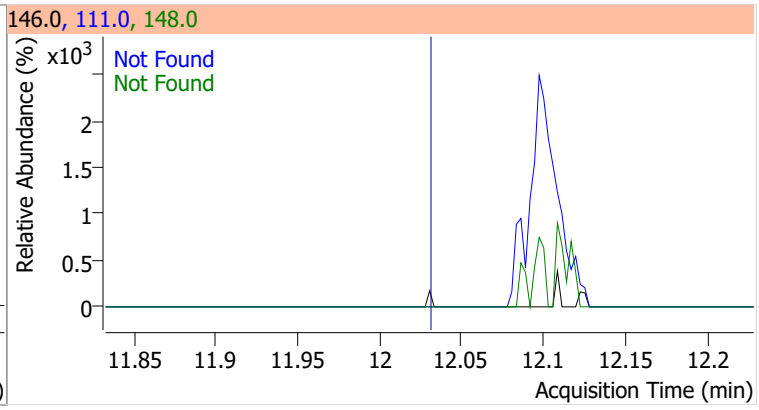
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



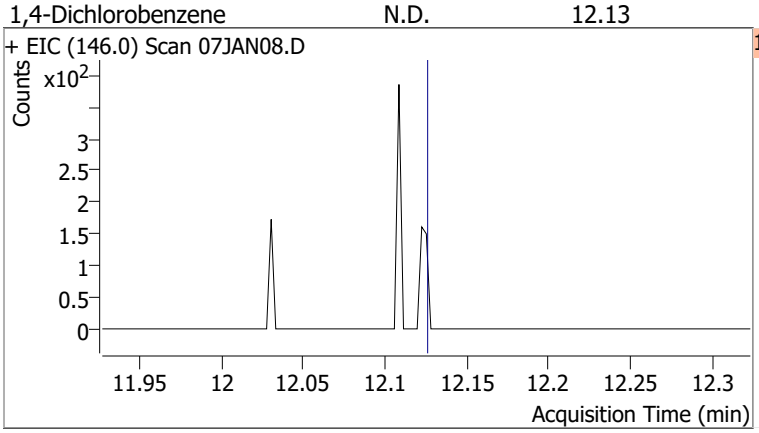


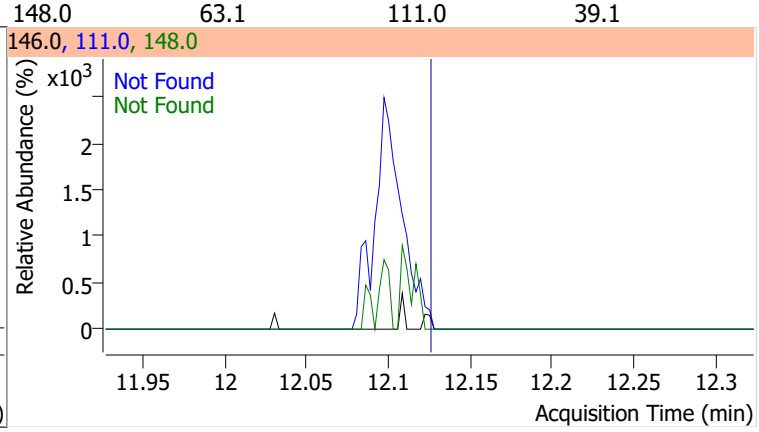
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



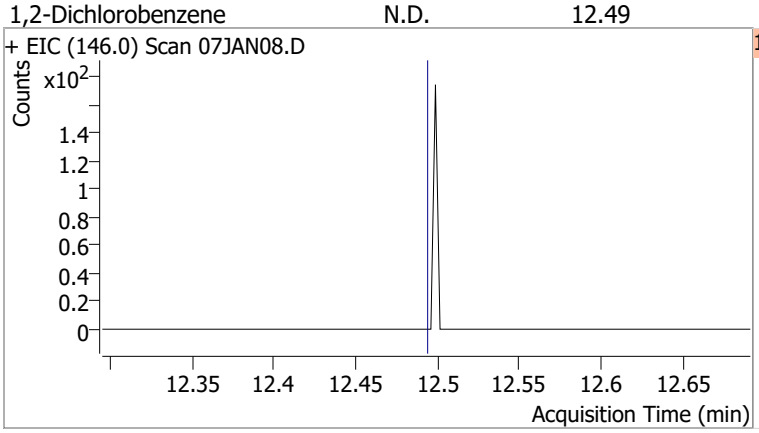


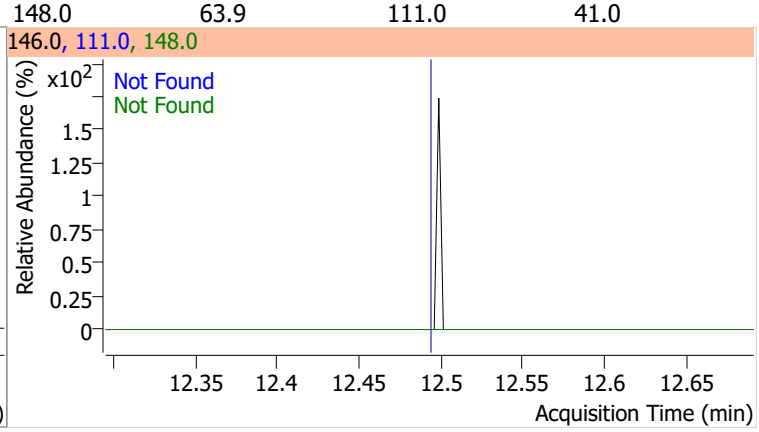
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





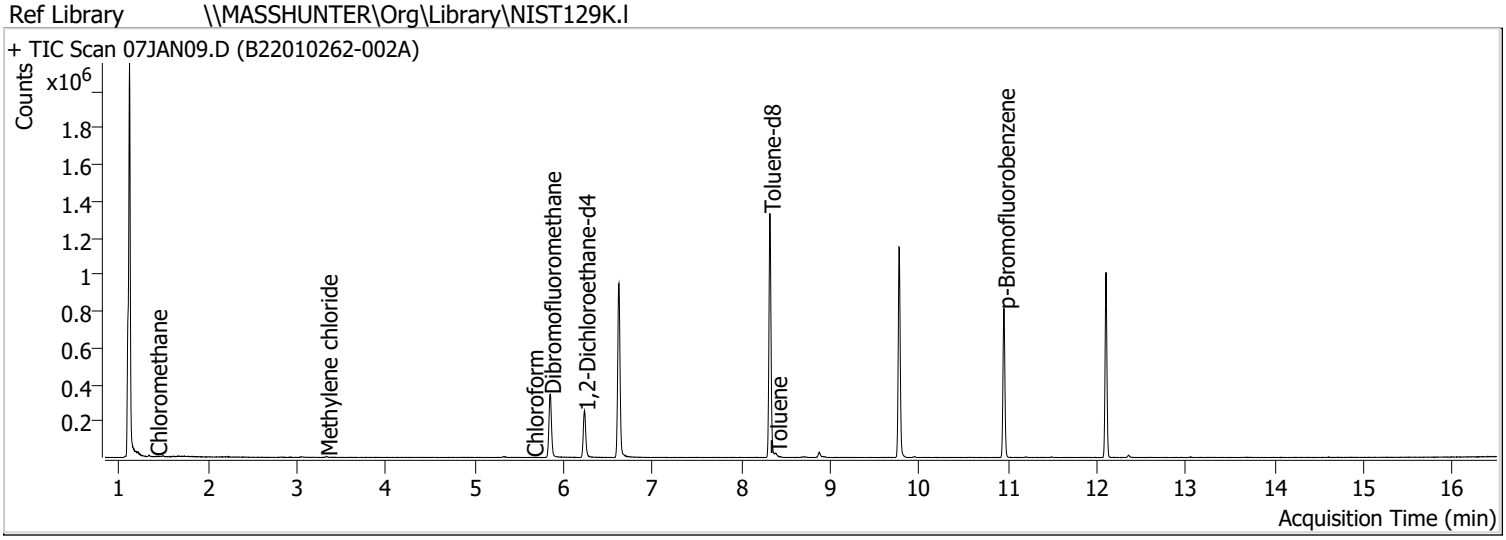
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN09.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 1:19:35 PM
Sample Name	B22010262-002A	Instrument	VOA5975C
Vial	9	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

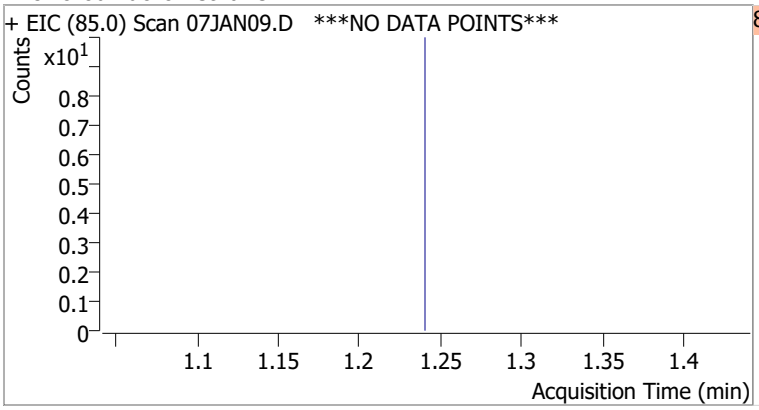


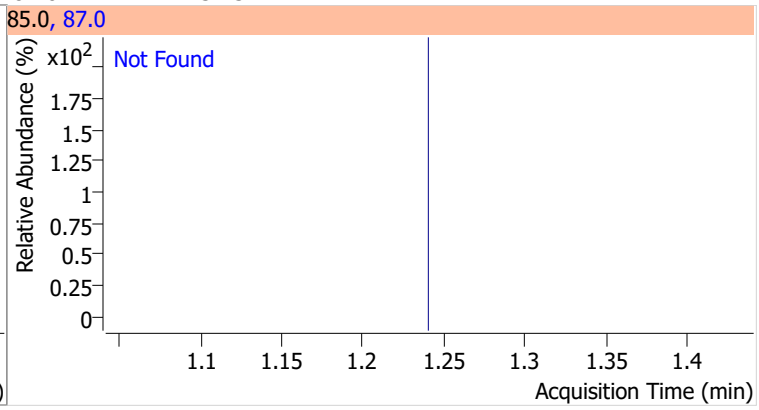
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	811488	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	311589	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	236584	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	211827	277.0777	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 110.83%			
S 1,2-Dichloroethane-d4	6.233	67.0	92974	281.5597	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 112.62%			
S Toluene-d8	8.319	98.0	816456	271.9136	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 108.77%			
S p-Bromofluorobenzene	10.951	95.0	234685	270.7711	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 108.31%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.414	50.0	1041	0.8064	ng	m	85
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.330	49.0	2658	2.2061	ng	m	98
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	5.650	83.0	381	0.2466	ng	m	99

Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.		
T Carbon tetrachloride	0.000		0	N.D.		
T 1,1-Dichloropropene	0.000		0	N.D.		
T Benzene	0.000		0	N.D.		
T 1,2-Dichloroethane	0.000		0	N.D.		
T Trichloroethene	0.000		0	N.D.		
T 1,2-Dichloropropane	0.000		0	N.D.		
T Dibromomethane	0.000		0	N.D.		
T Bromodichloromethane	0.000		0	N.D.		
T cis-1,3-Dichloropropene	0.000		0	N.D.		
T Toluene	8.391	92.0	4452	2.1950	ng	99
T trans-1,3-Dichloropropene	0.000		0	N.D.		
T 1,1,2-Trichloroethane	0.000		0	N.D.		
T Tetrachloroethene	0.000		0	N.D.		
T 1,3-Dichloropropane	0.000		0	N.D.		
T Chlorodibromomethane	0.000		0	N.D.		
T 1,2-Dibromoethane	0.000		0	N.D.		
T Chlorobenzene	0.000		0	N.D.		
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.		
T Ethylbenzene	0.000		0	N.D.		
T m+p-Xylenes	0.000		0	N.D.		
T o-Xylene	0.000		0	N.D.		
T Styrene	0.000		0	N.D.		
T Bromoform	0.000		0	N.D.		
T Bromobenzene	0.000		0	N.D.		
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
T 1,2,3-Trichloropropane	0.000		0	N.D.		
T 2-Chlorotoluene	0.000		0	N.D.		
T 4-Chlorotoluene	0.000		0	N.D.		
T 1,3-Dichlorobenzene	0.000		0	N.D.		
T 1,4-Dichlorobenzene	0.000		0	N.D.		
T 1,2-Dichlorobenzene	0.000		0	N.D.		

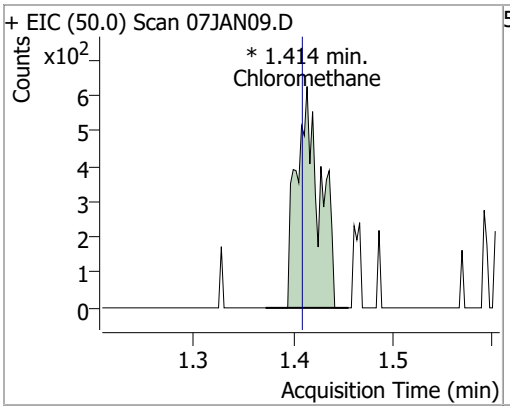
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

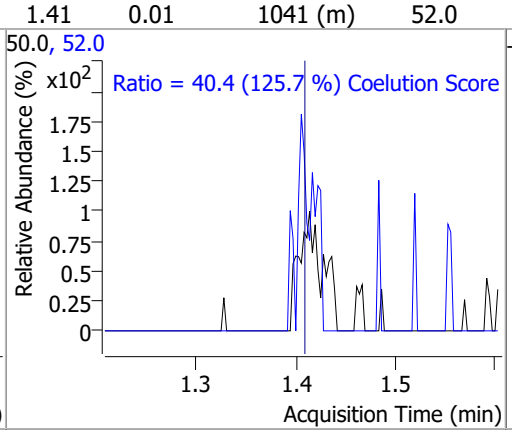
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

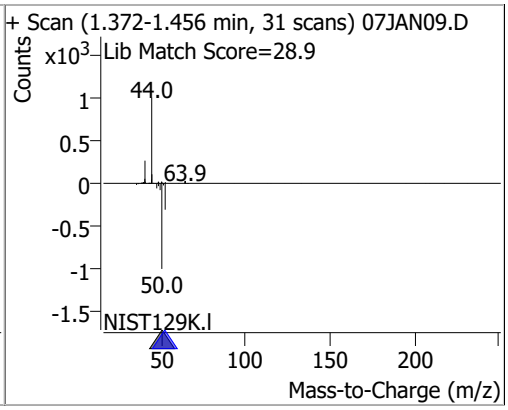




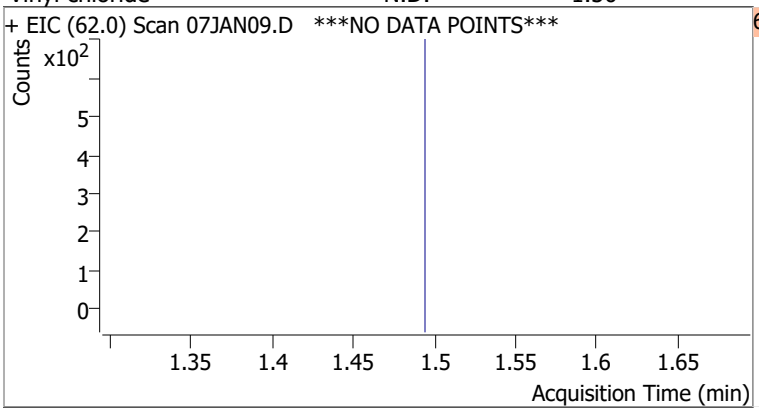
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	0.8064	1.41	0.01	1041 (m)	52.0	40.4	2.1	62.1

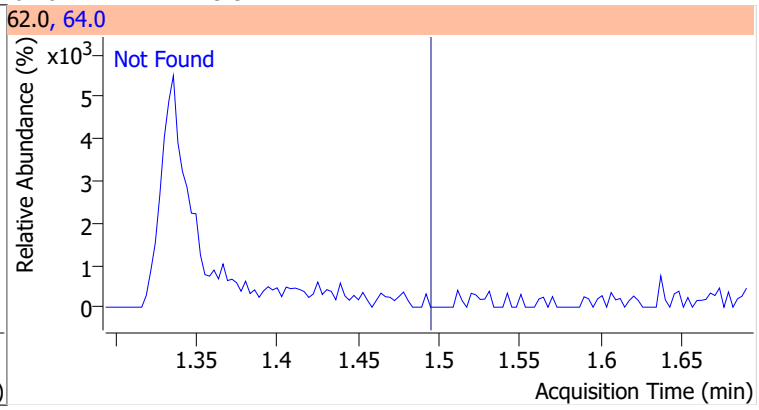




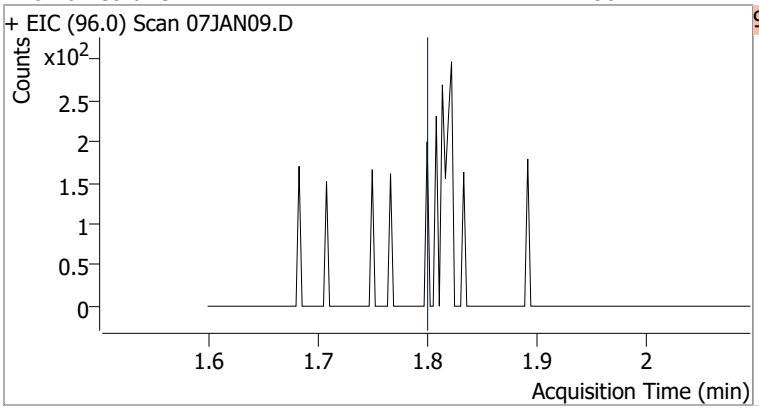


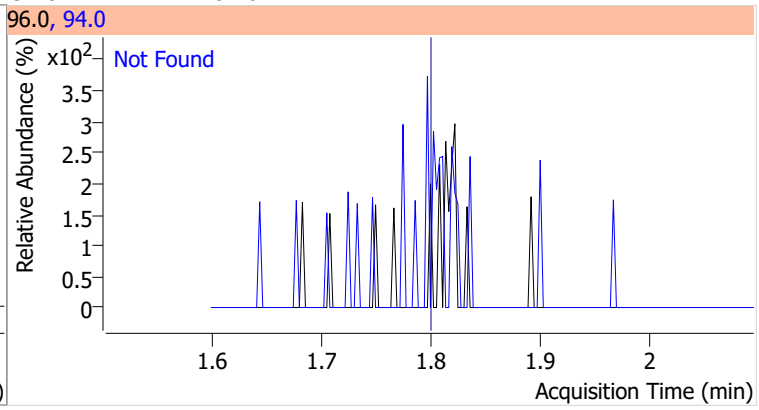
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



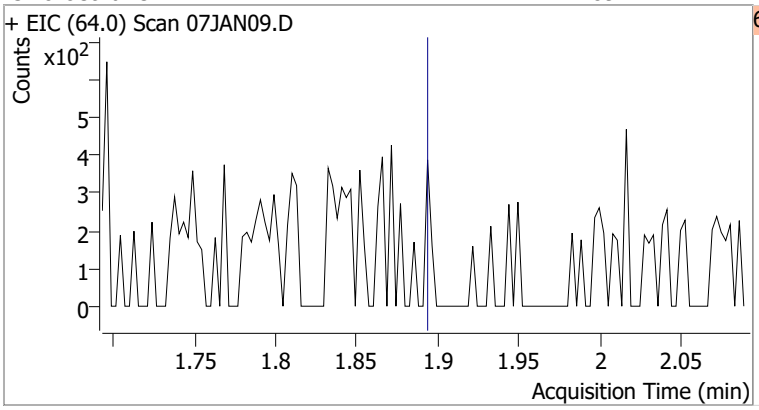


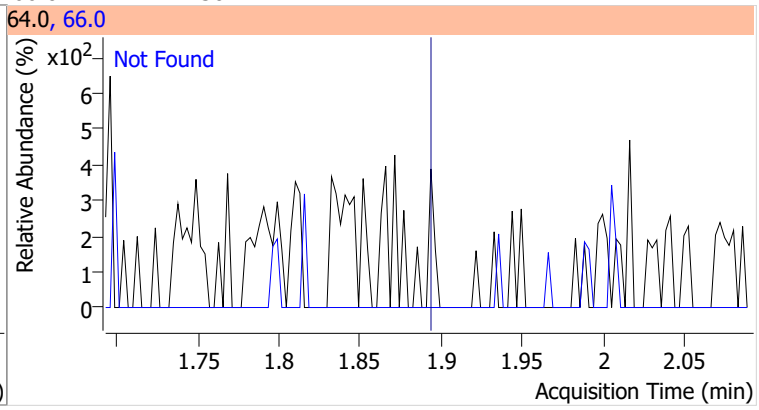
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



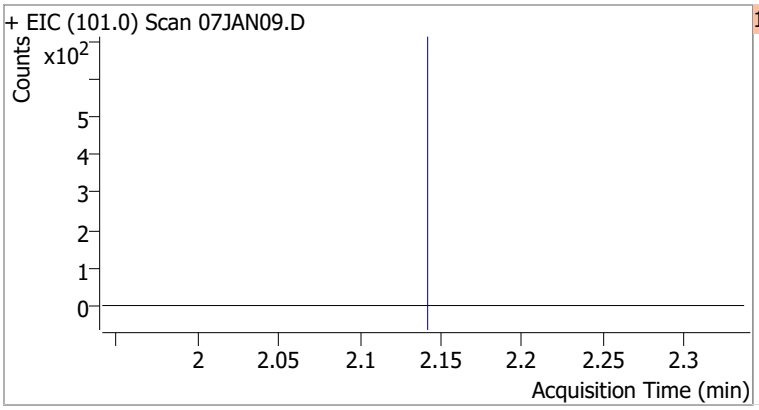


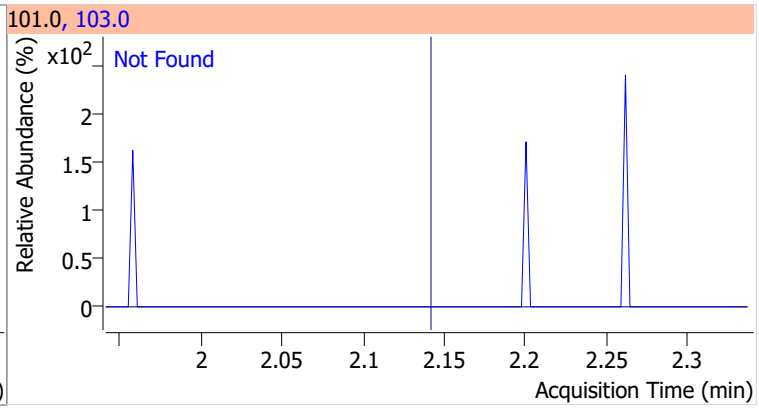
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



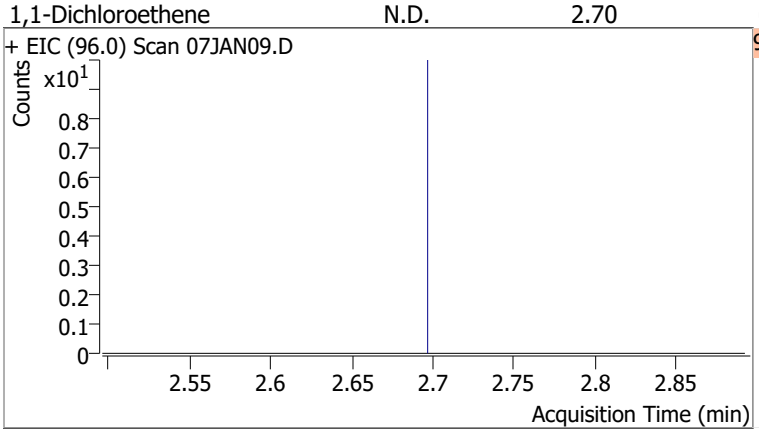


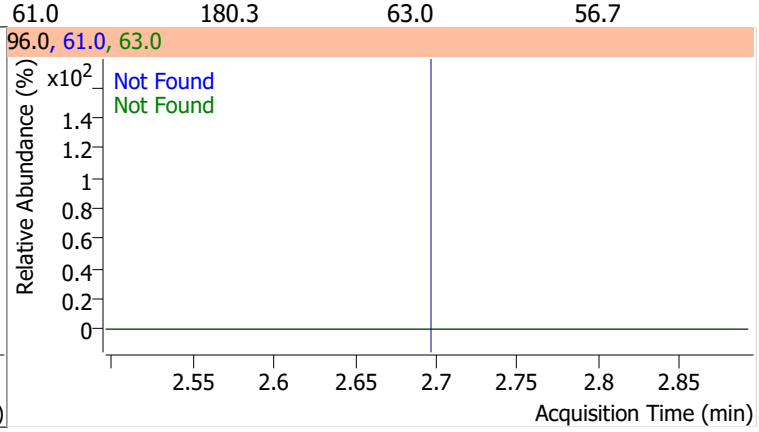
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



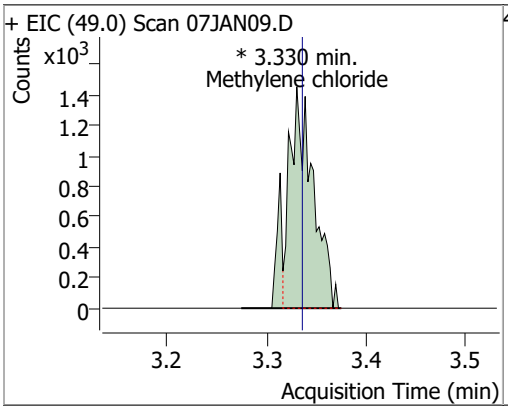


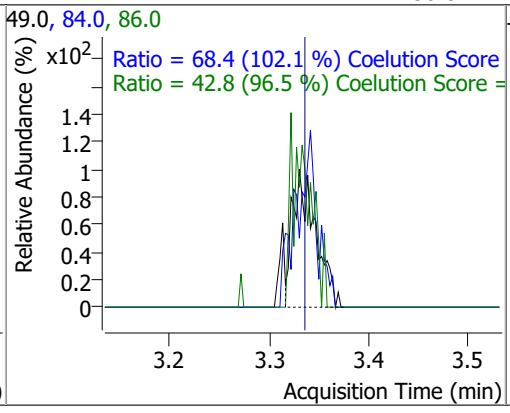
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

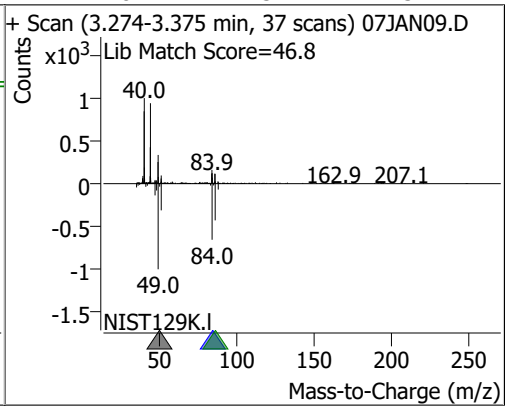




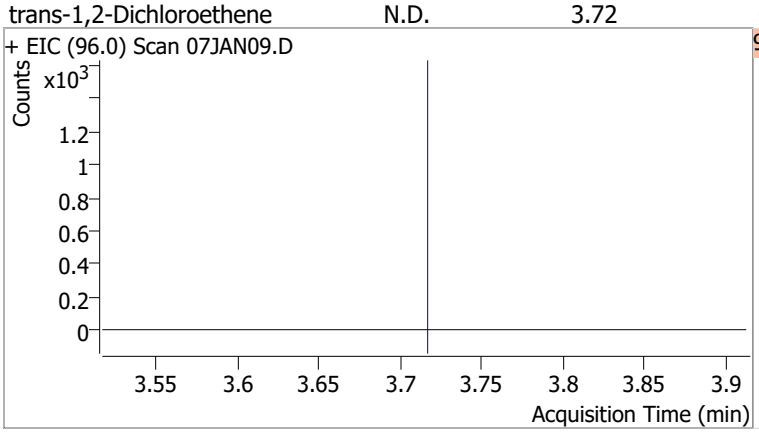
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	2.2061	3.33	-0.01	2658 (m)	84.0	68.4	36.9	96.9
					86.0	42.8	14.3	74.3

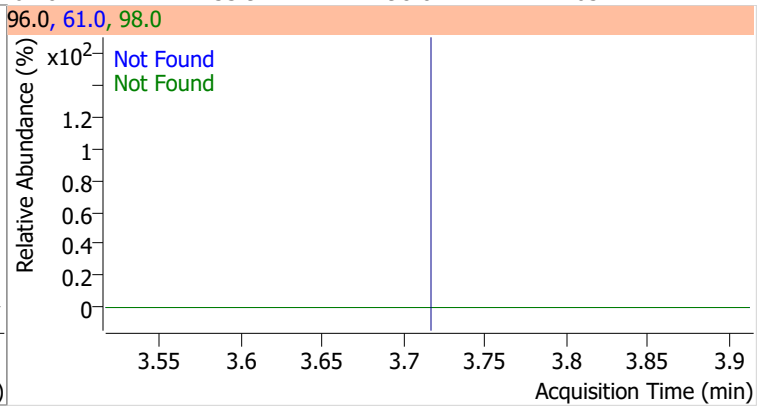




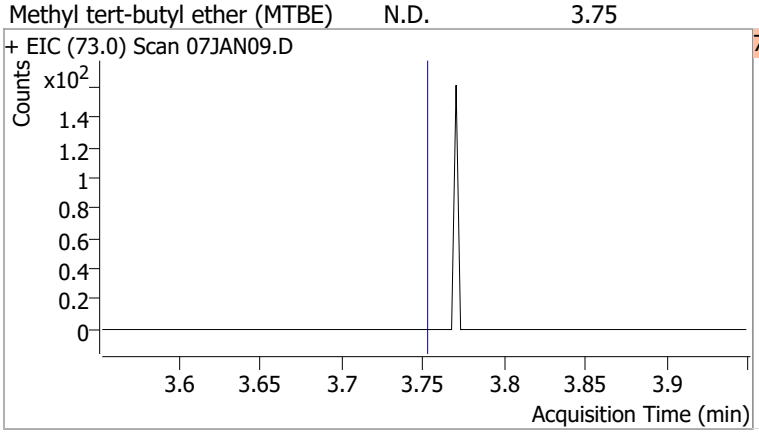


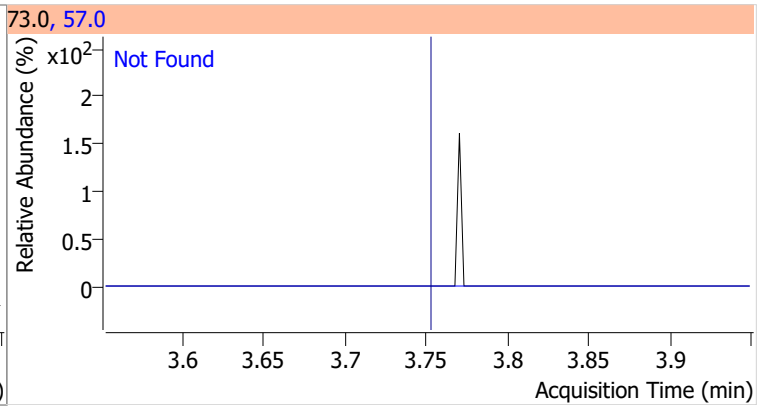
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



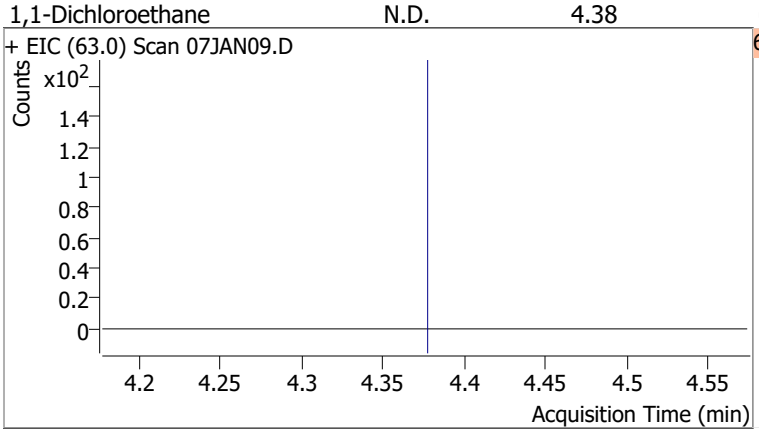


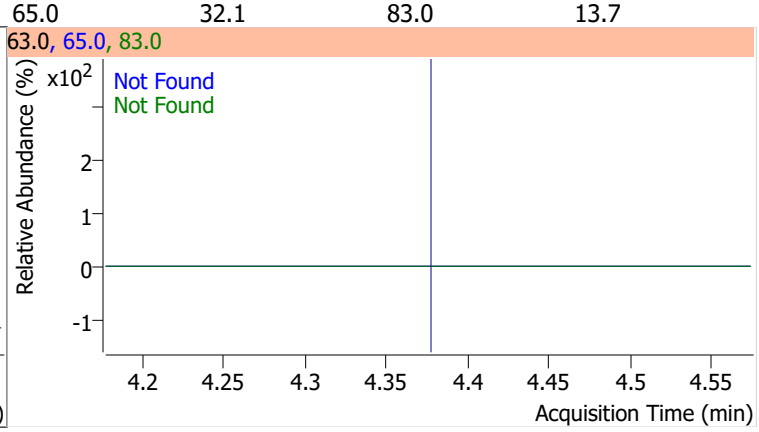
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



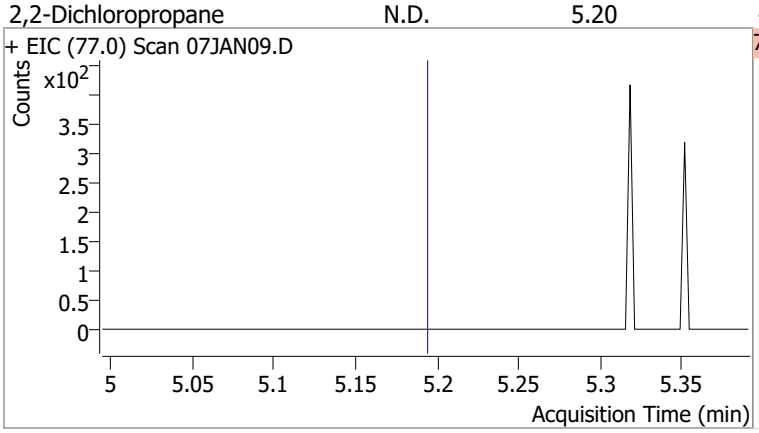


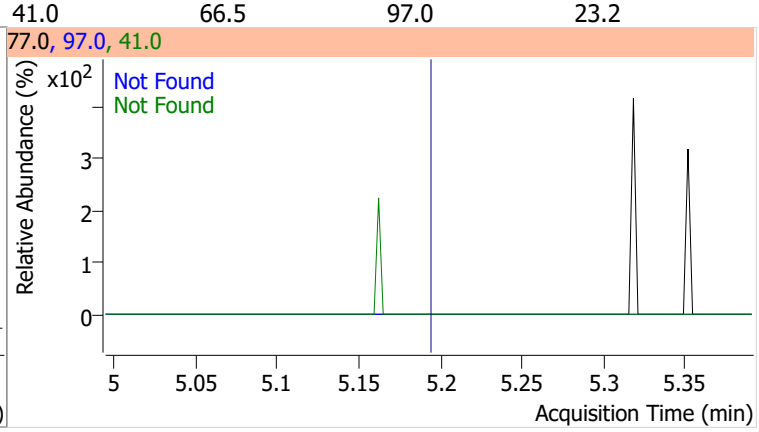
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



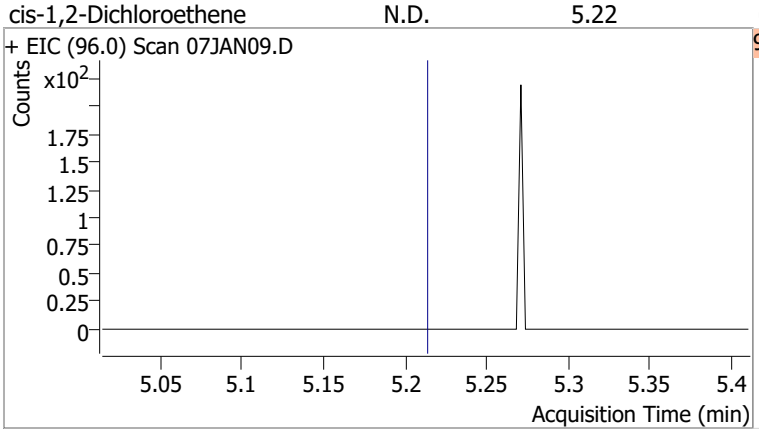


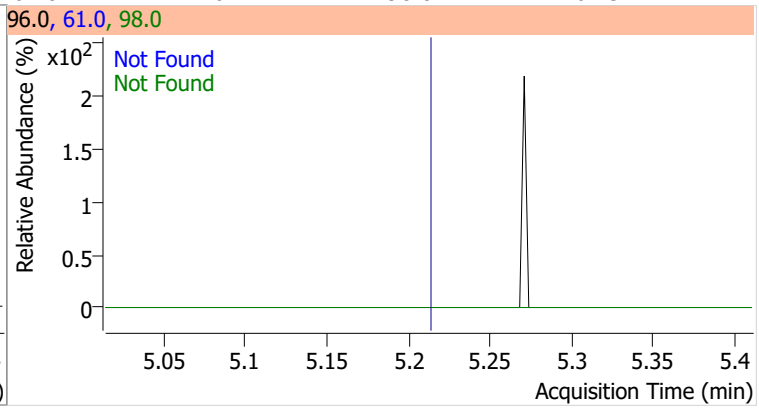
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



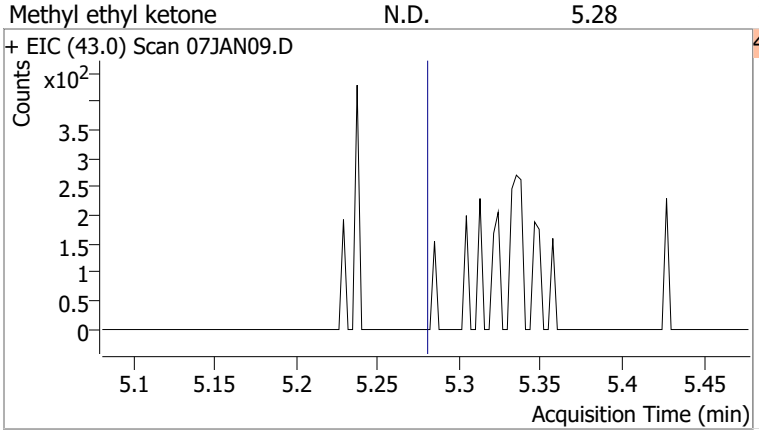


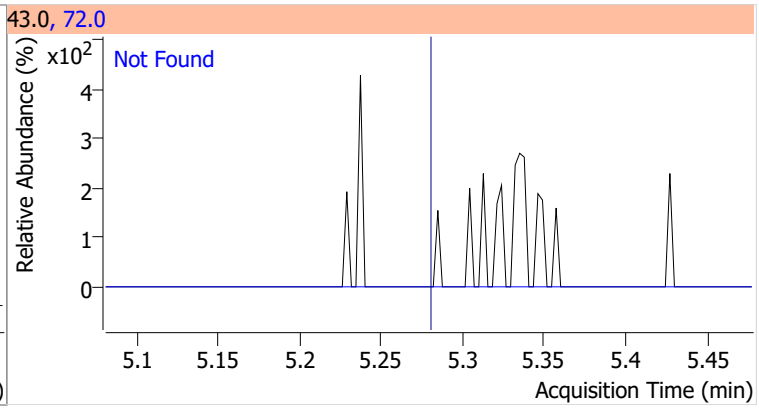
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



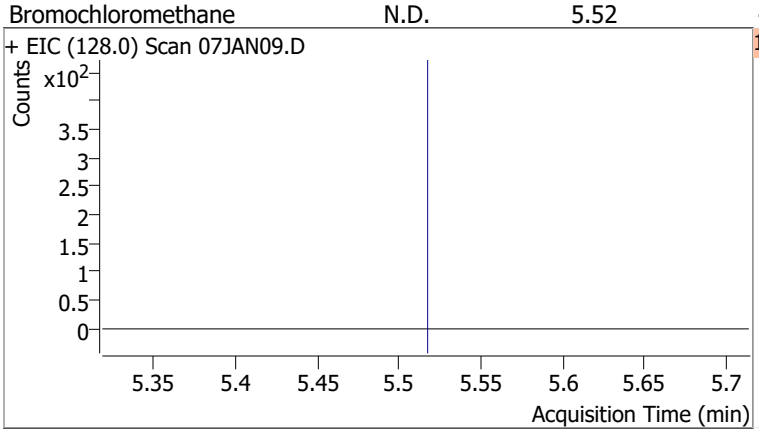


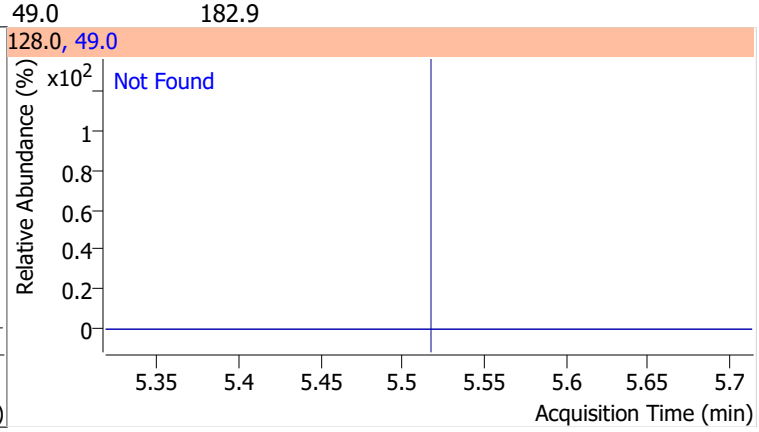
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



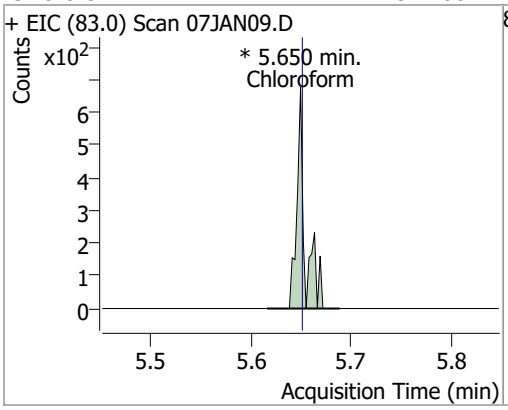


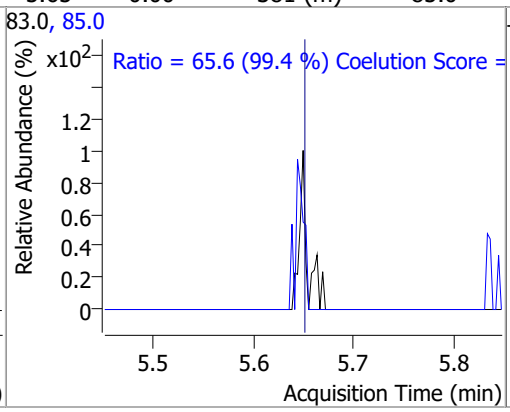
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9

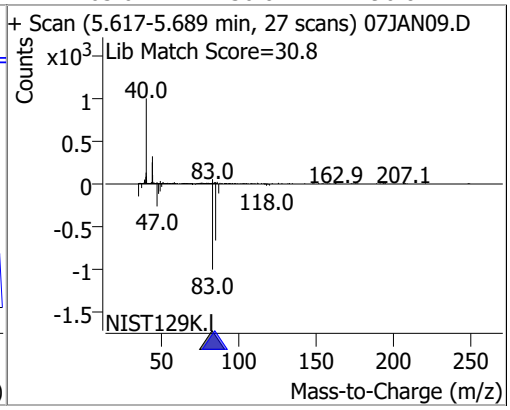




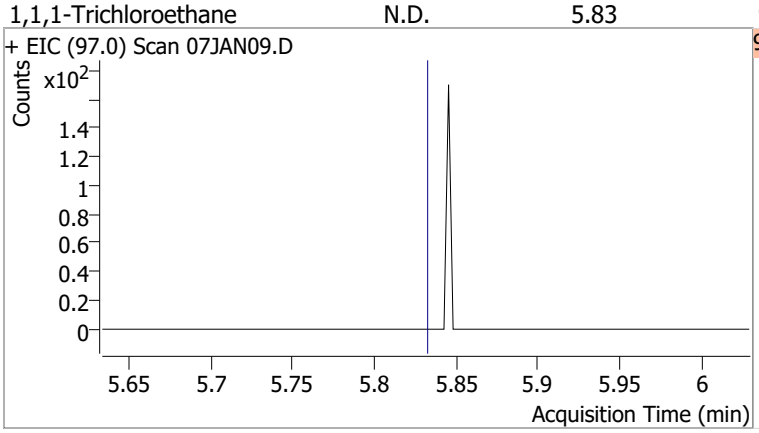
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	0.2466	5.65	0.00	381 (m)	85.0	65.6	36.0	96.0

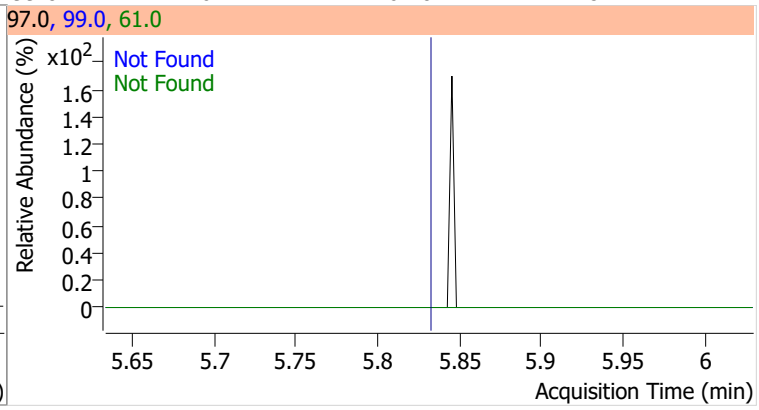




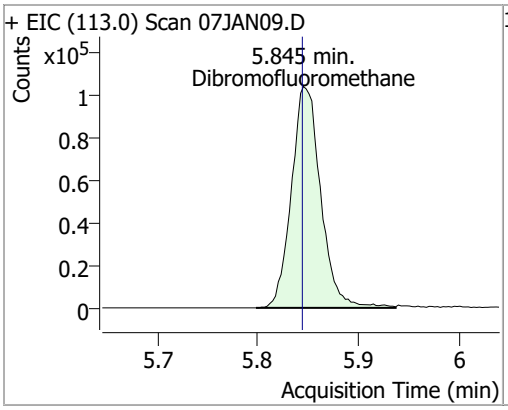


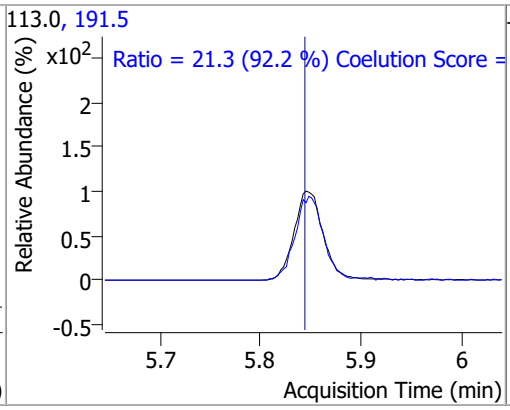
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

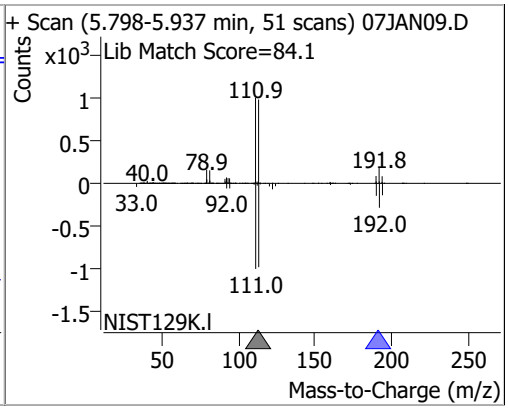




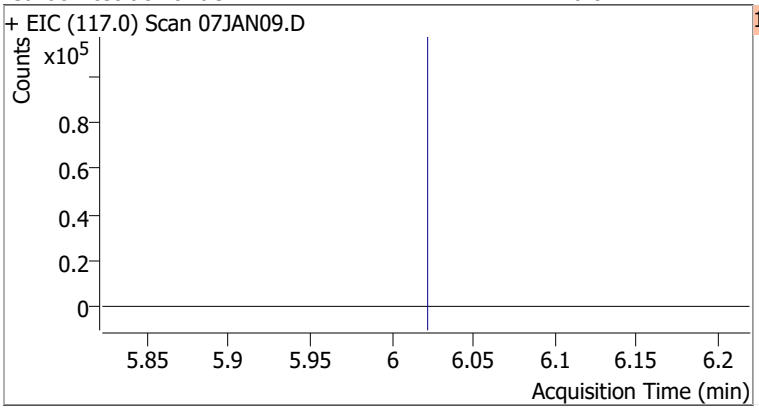
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	277.0777	5.85	0.00	211827	191.5	21.3	0.0	53.1

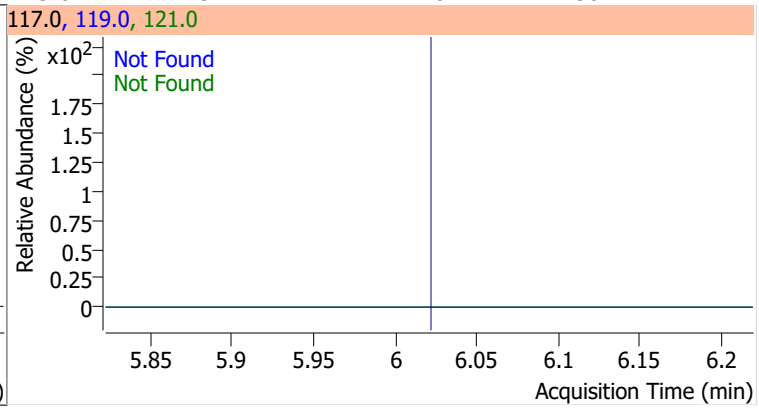




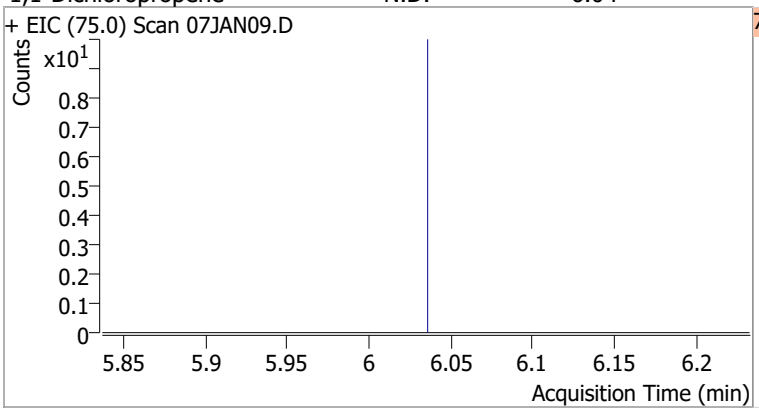


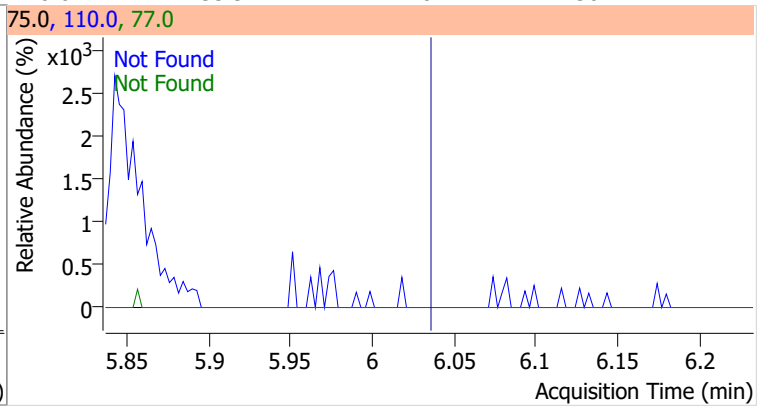
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



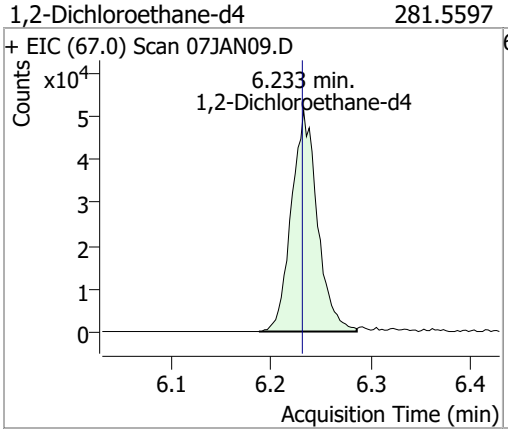


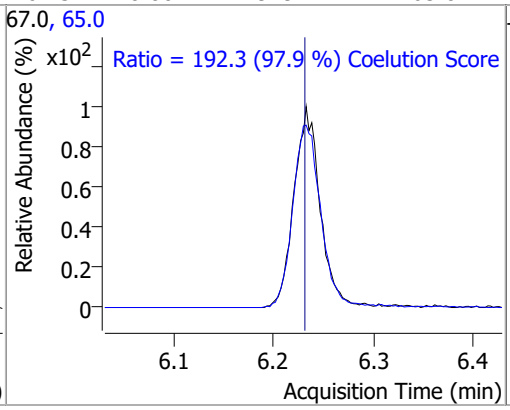
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

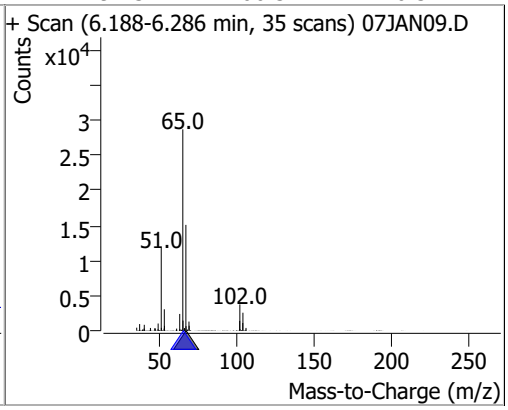




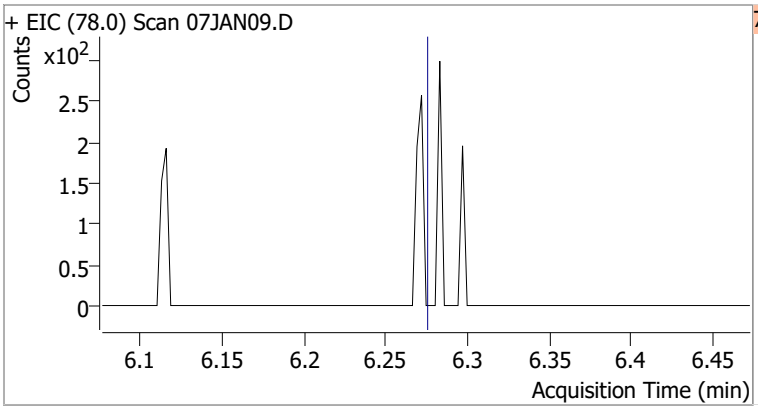
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	281.5597	6.23	0.00	92974	65.0	192.3	166.5	226.5

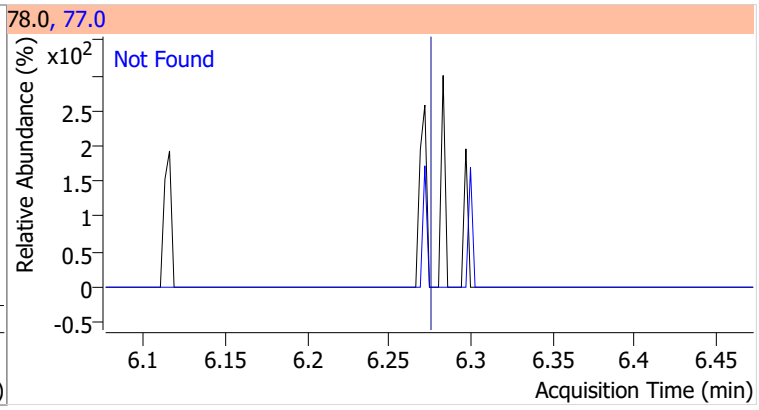




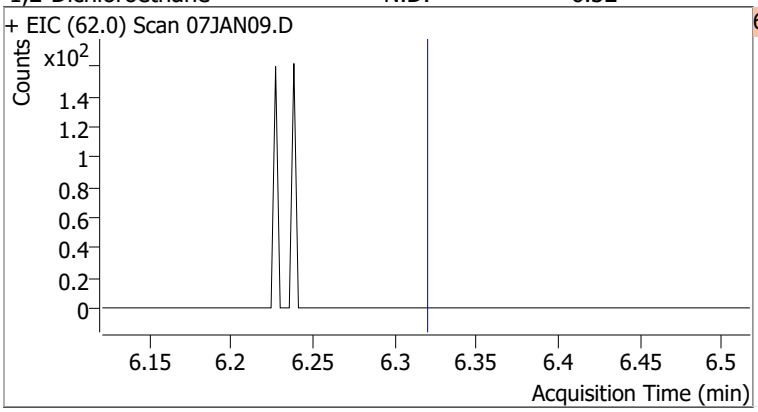


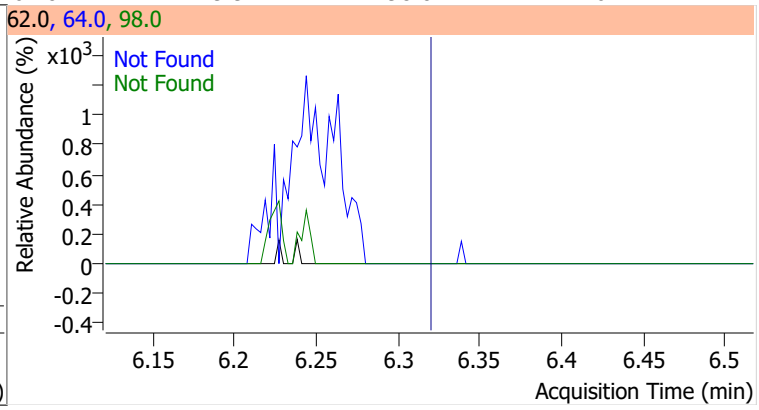
Compound	Conc.	Exp RT	QIon	Exp Ratio
Benzene	N.D.	6.28	77.0	23.5



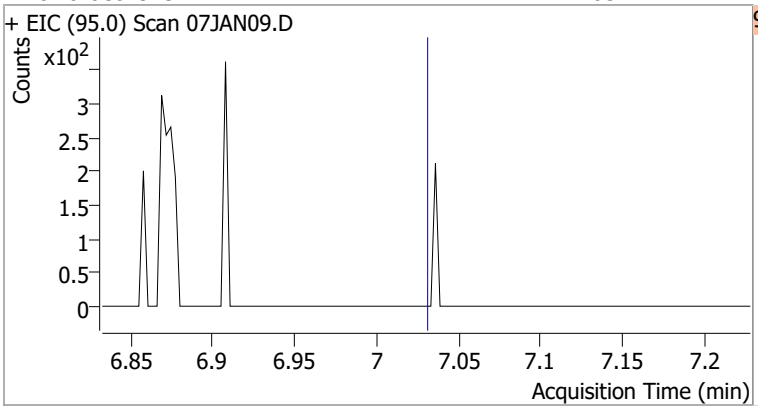


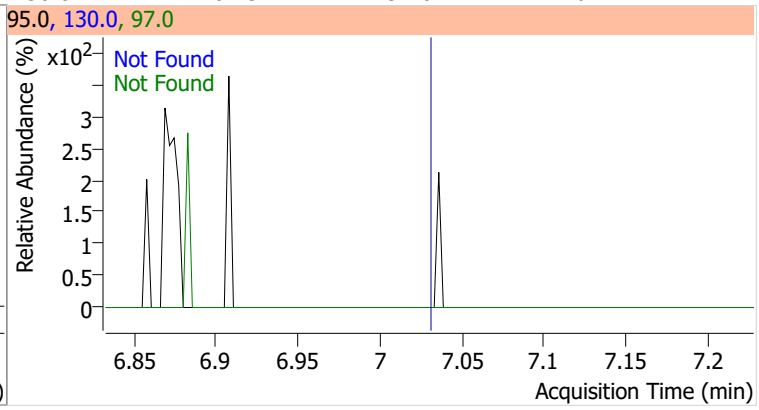
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



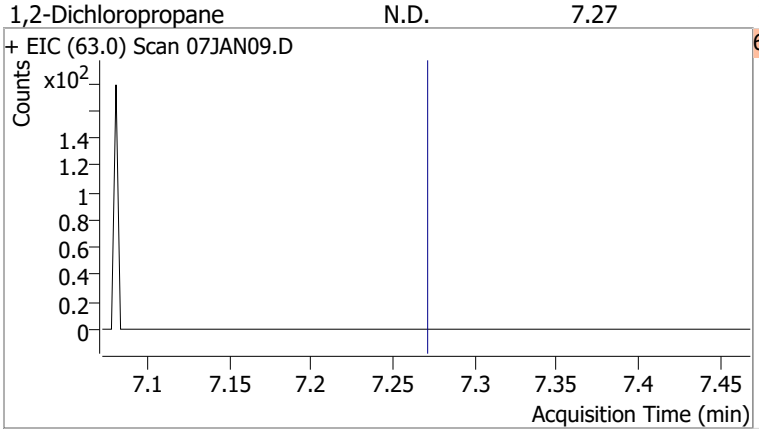


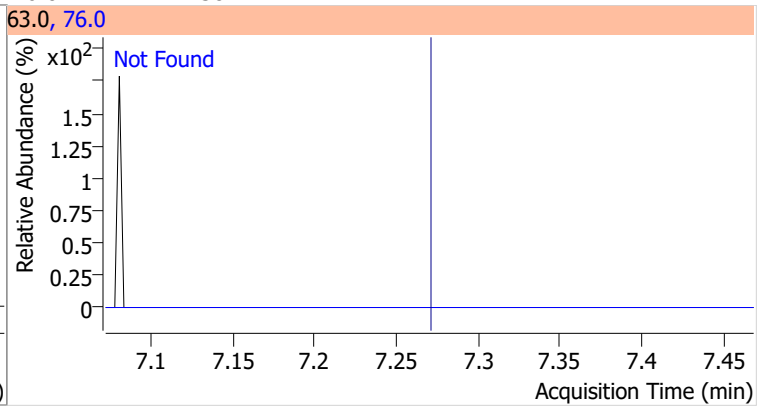
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



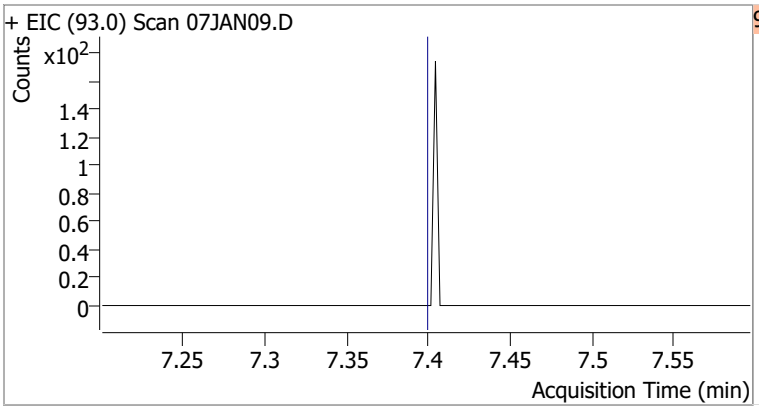


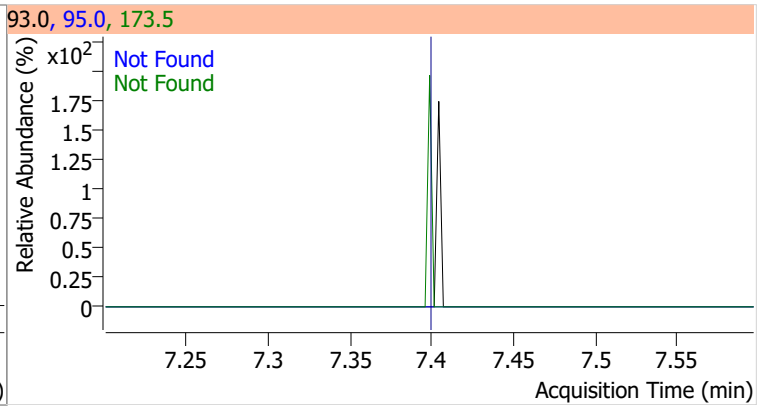
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



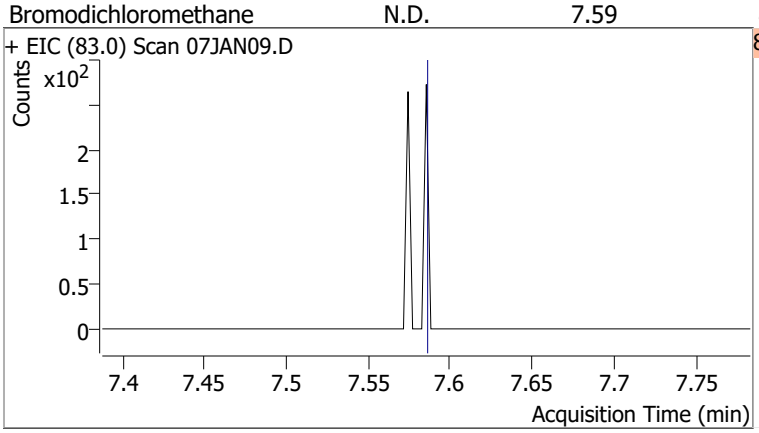


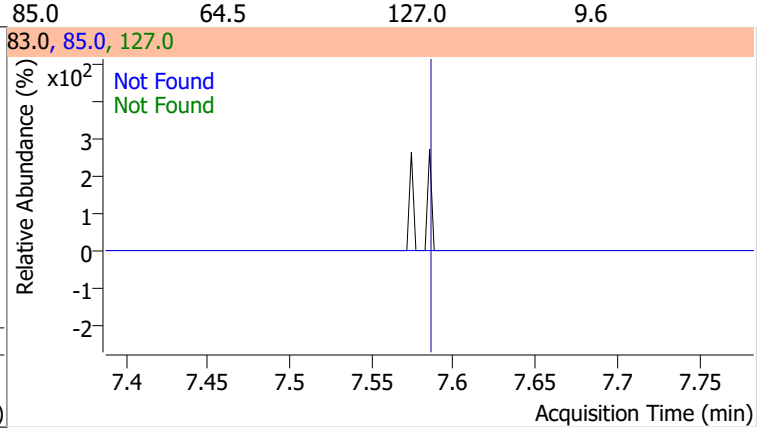
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



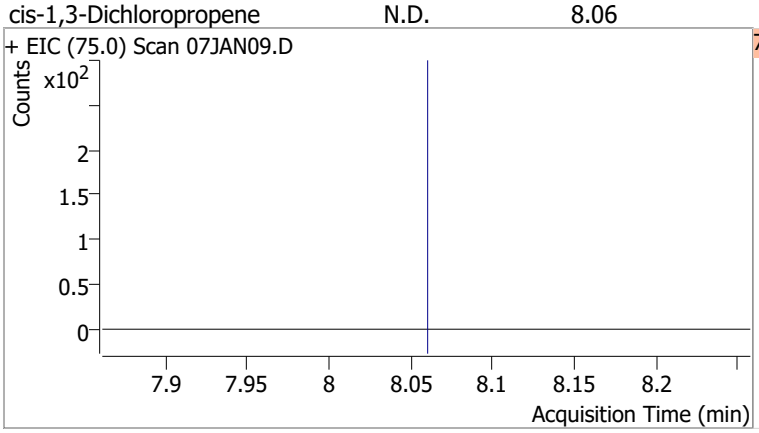


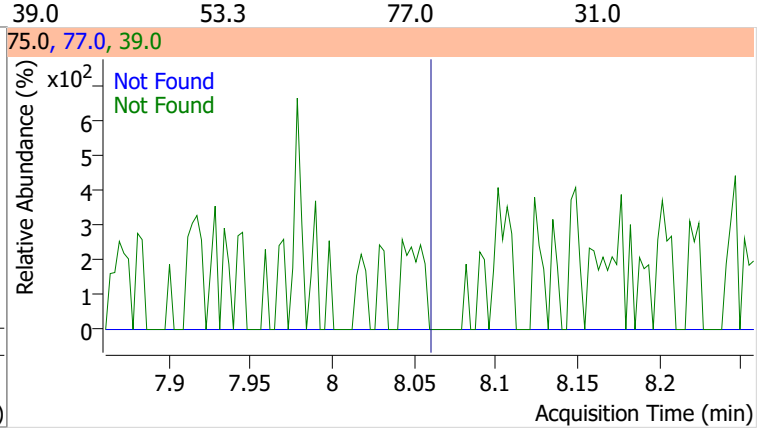
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



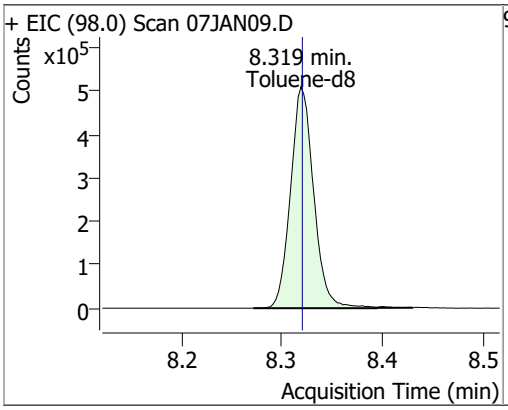


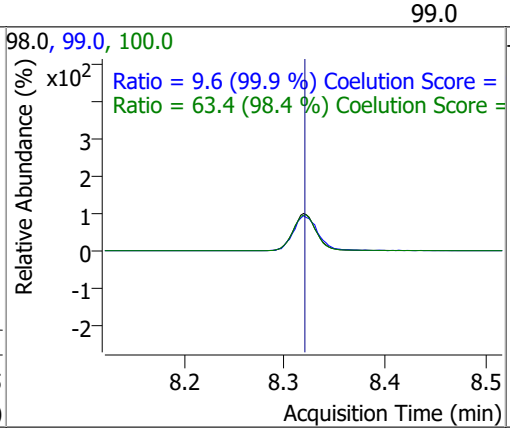
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

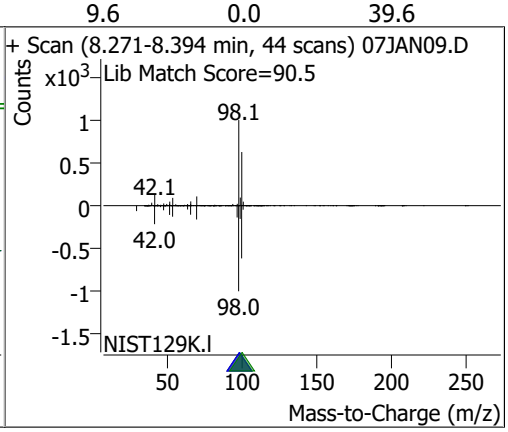




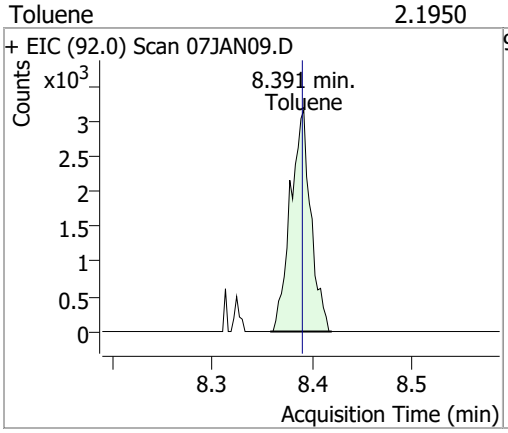
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	271.9136	8.32	0.00	816456	100.0	63.4	34.4	94.4
					99.0	9.6	0.0	39.6

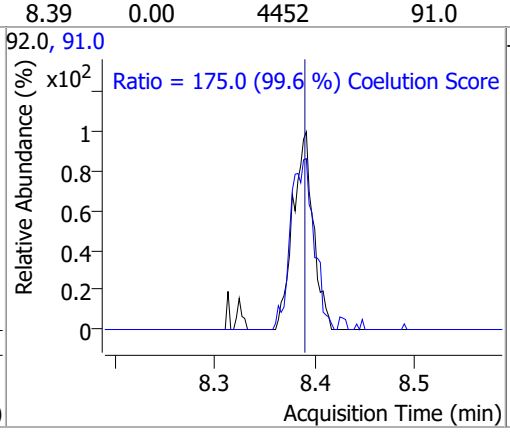


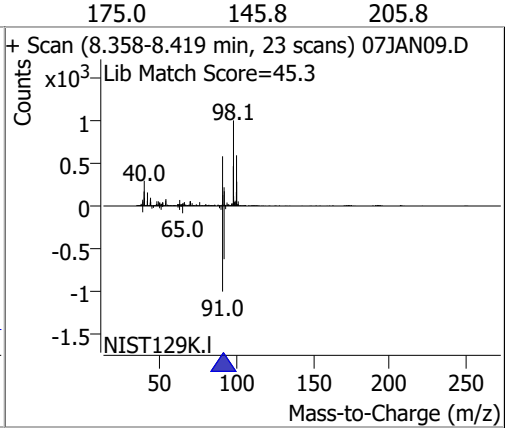




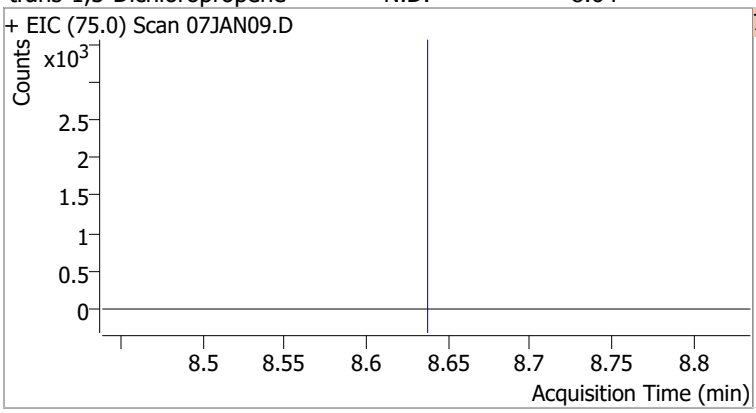
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	2.1950	8.39	0.00	4452	91.0	175.0	145.8	205.8

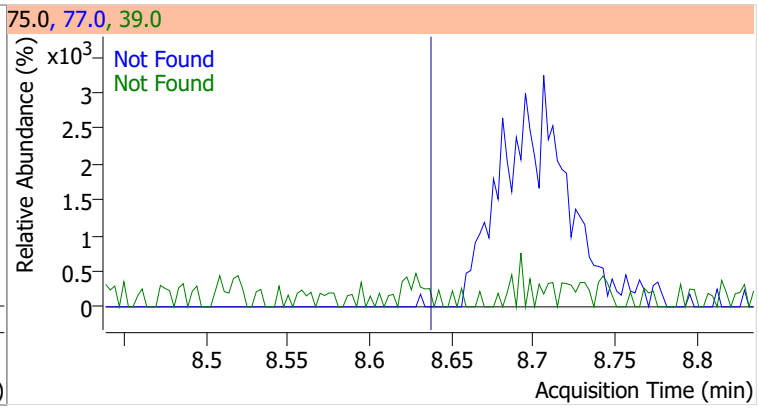




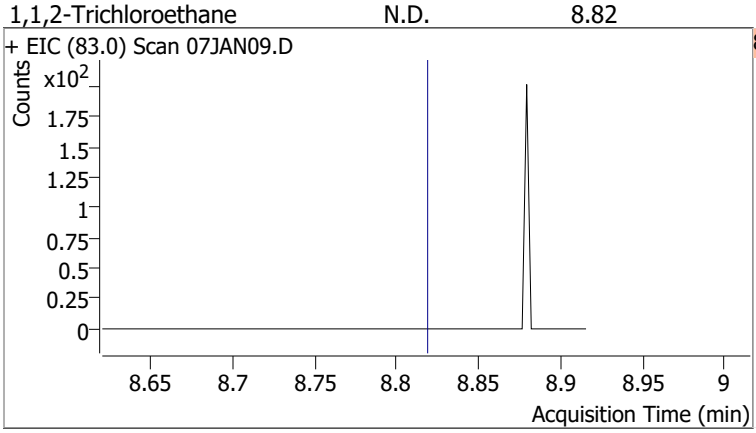


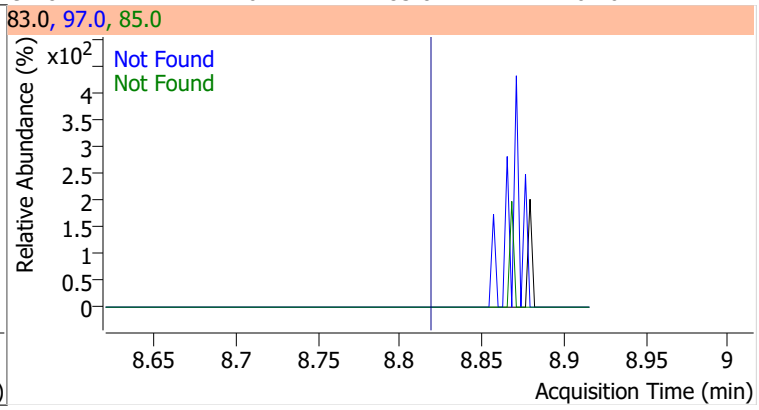
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



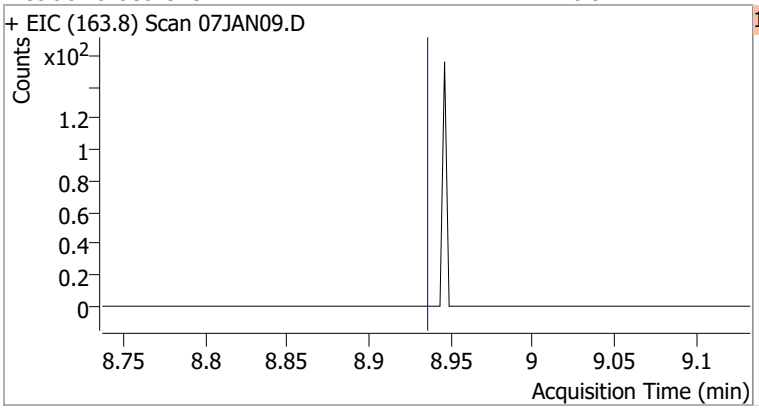


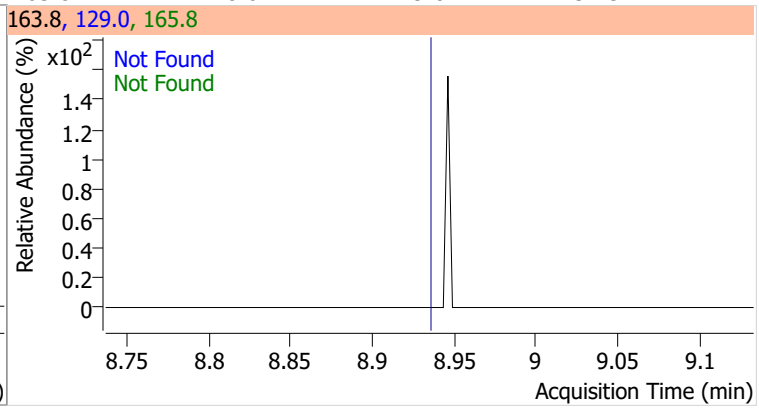
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



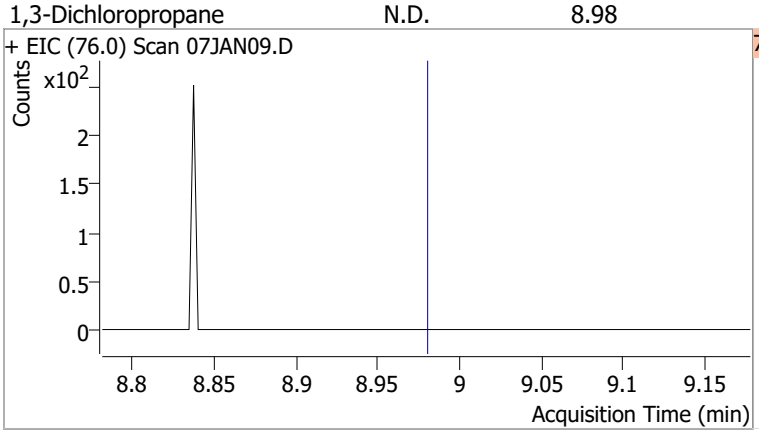


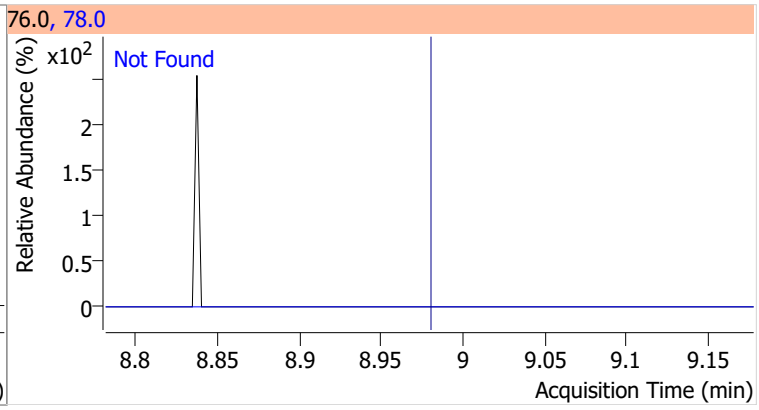
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



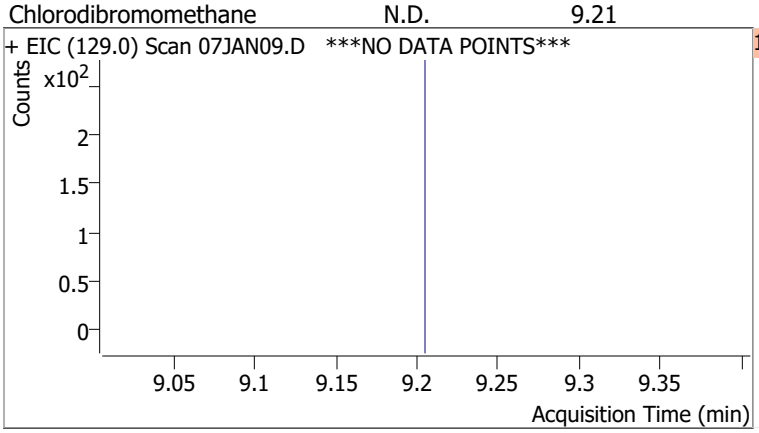


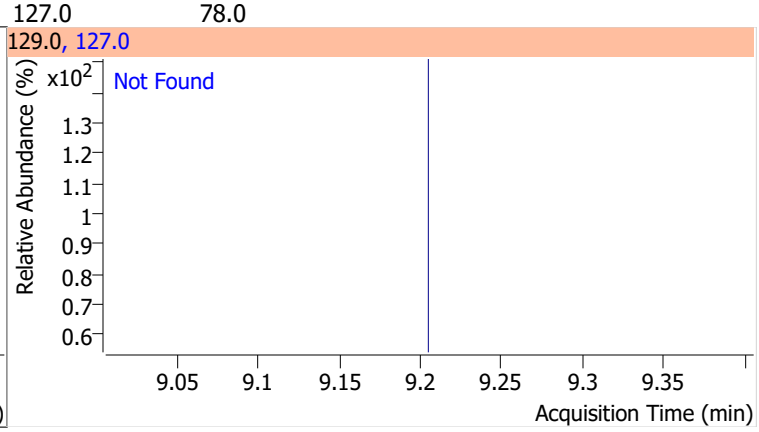
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



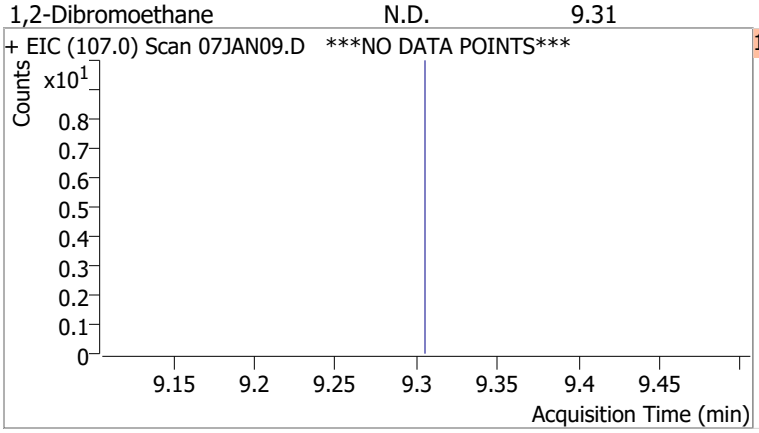


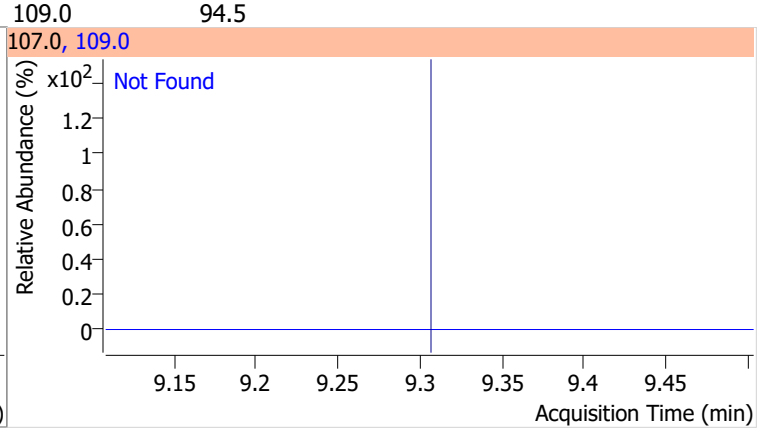
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



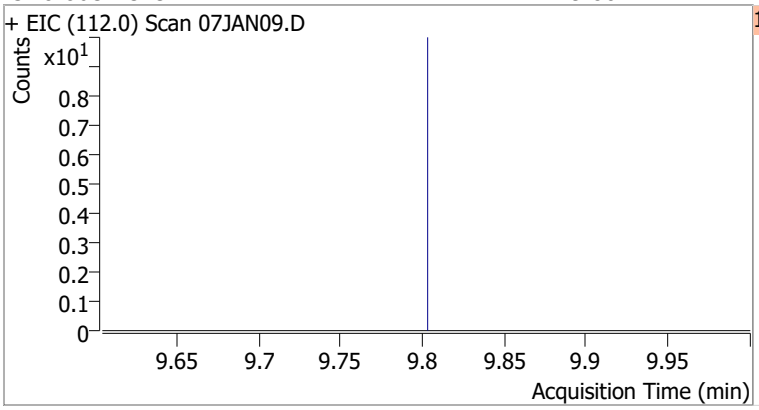


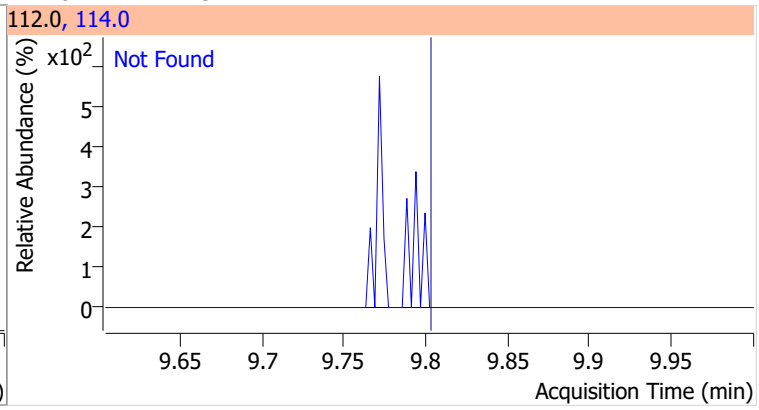
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



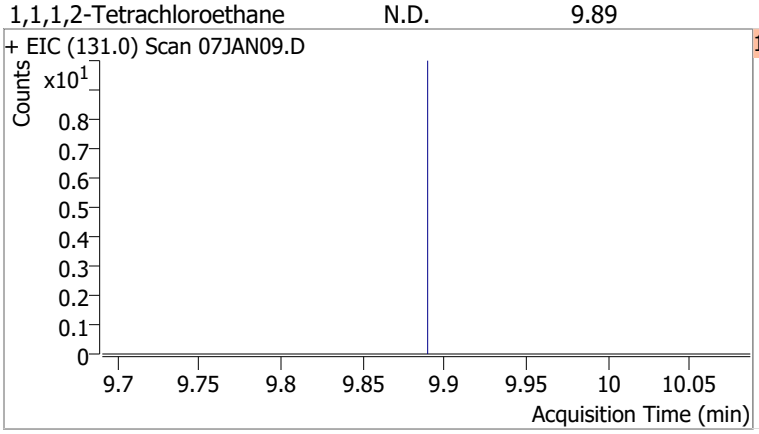


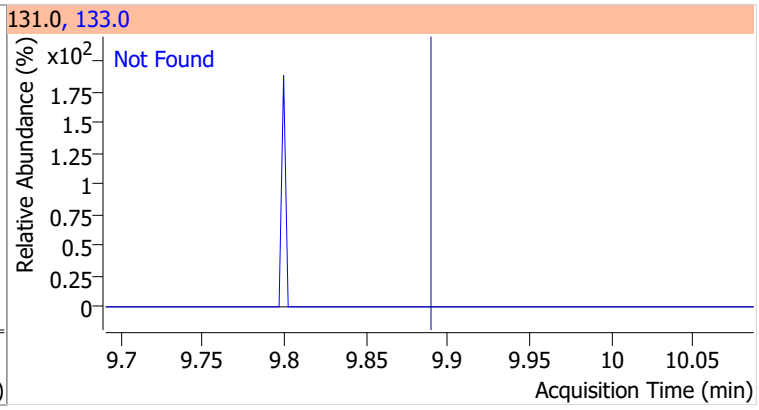
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



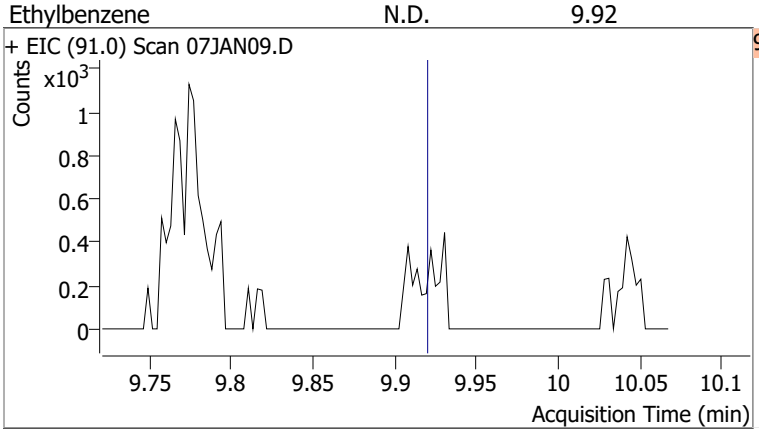


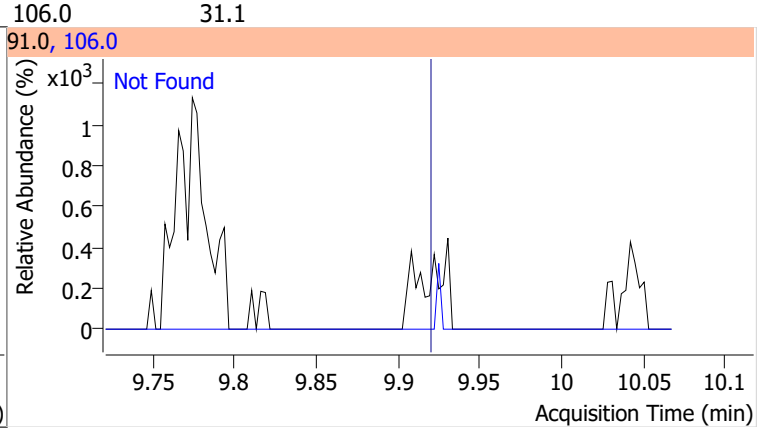
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



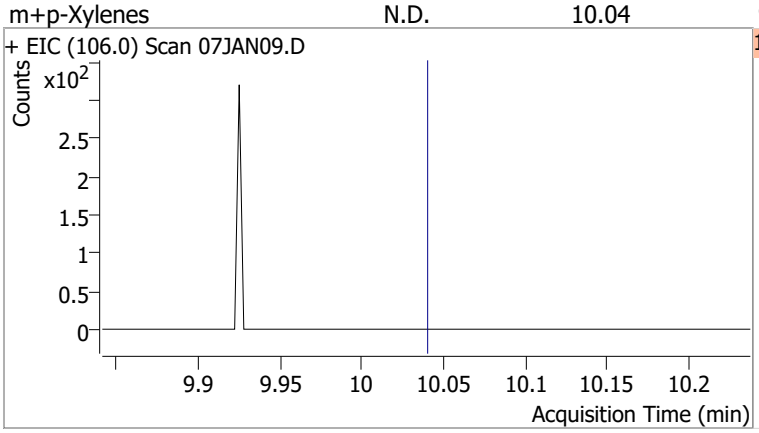


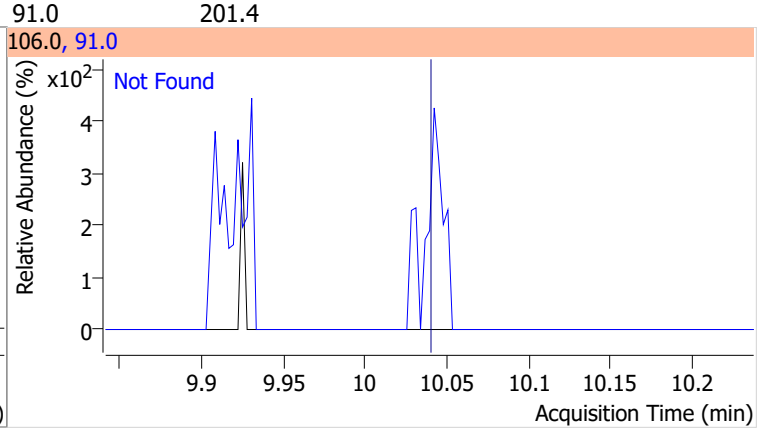
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1



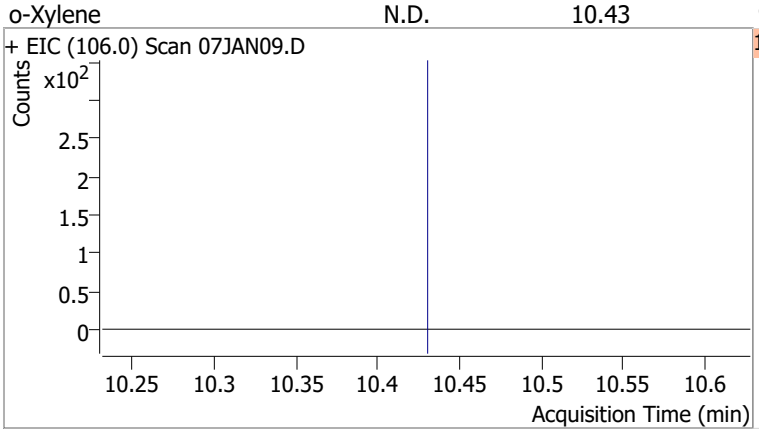


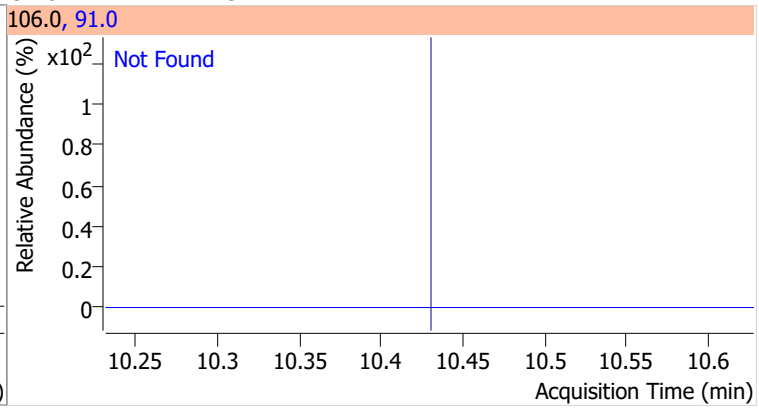
Compound	Conc.	Exp RT	QIon	Exp Ratio
m+p-Xylenes	N.D.	10.04	91.0	201.4



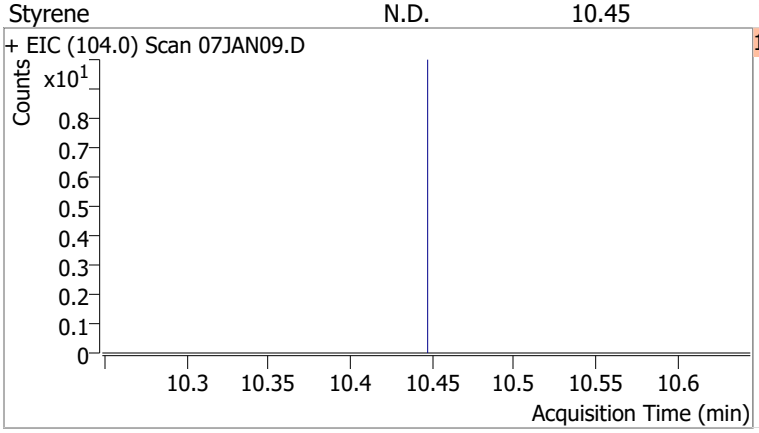


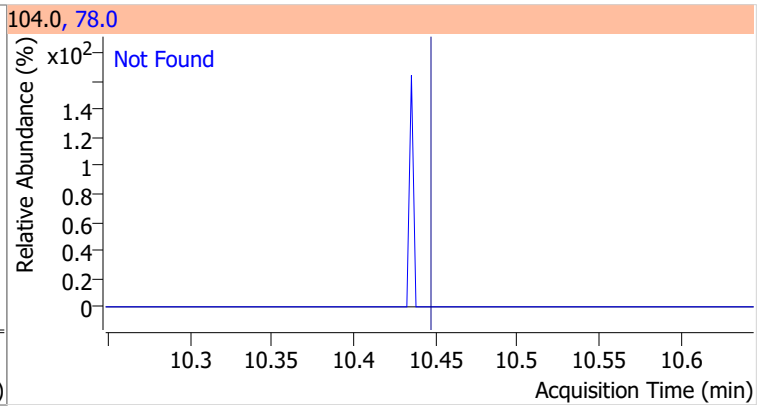
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



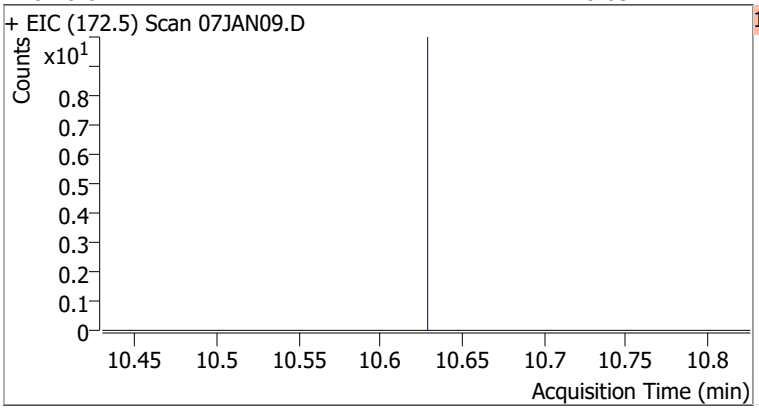


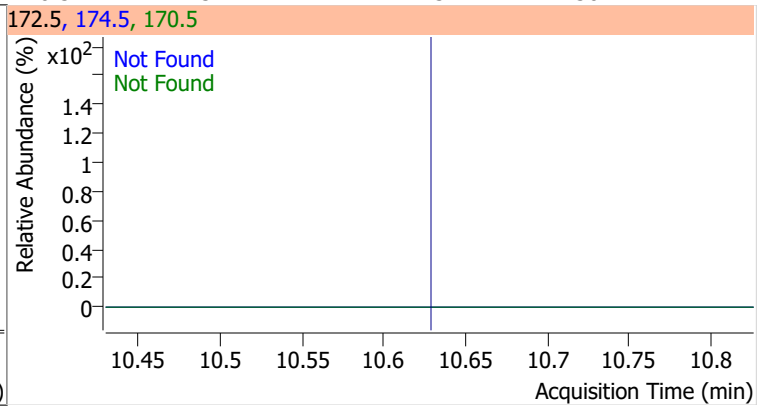
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



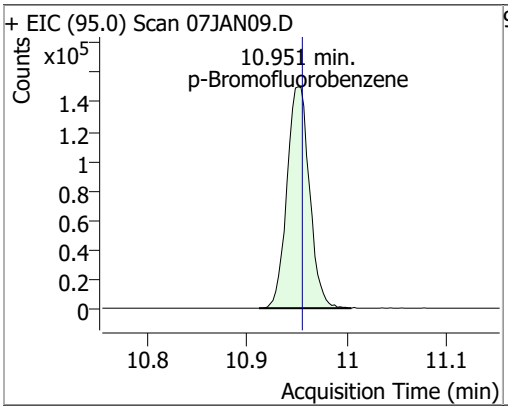


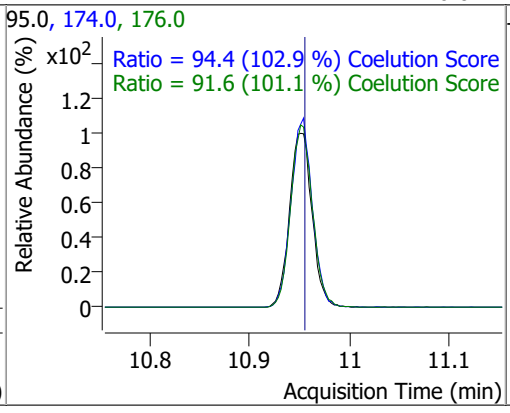
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

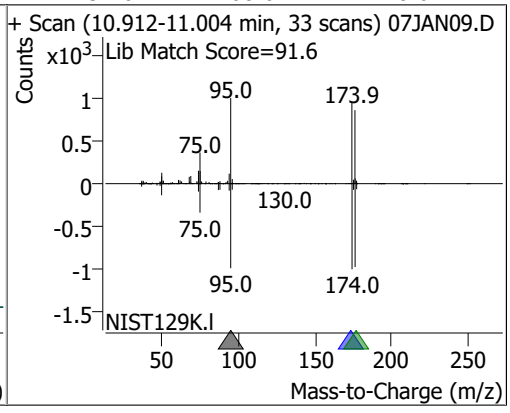




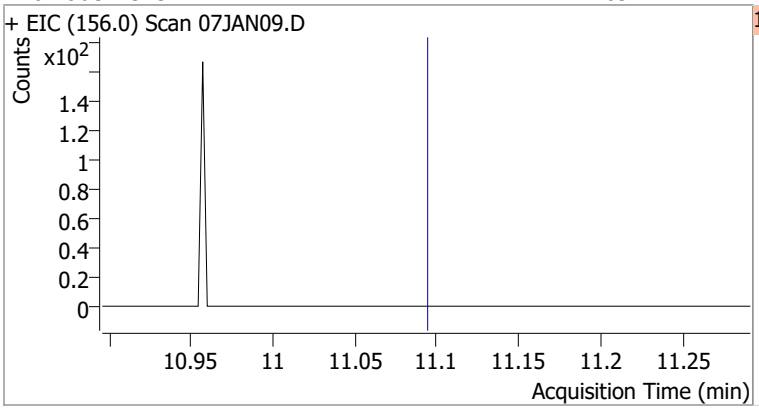
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	270.7711	10.95	0.00	234685	174.0	94.4	61.7	121.7
					176.0	91.6	60.6	120.6

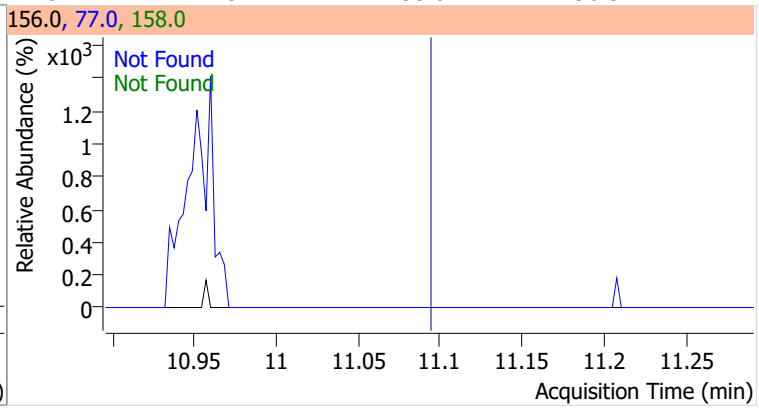




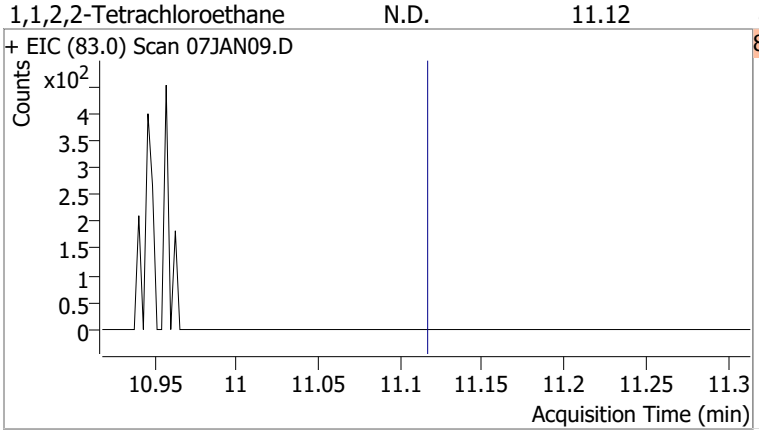


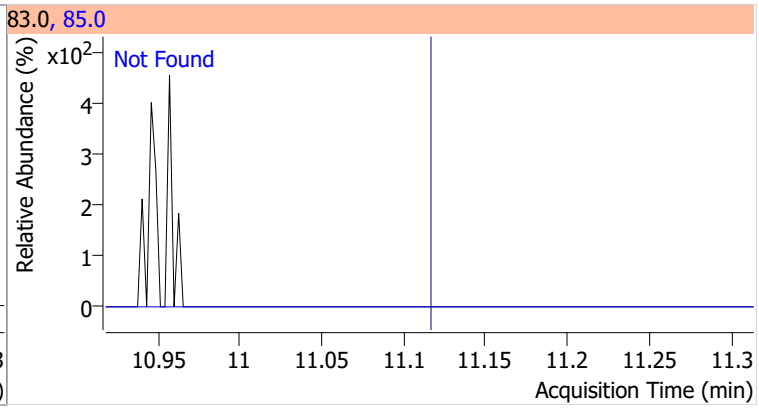
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



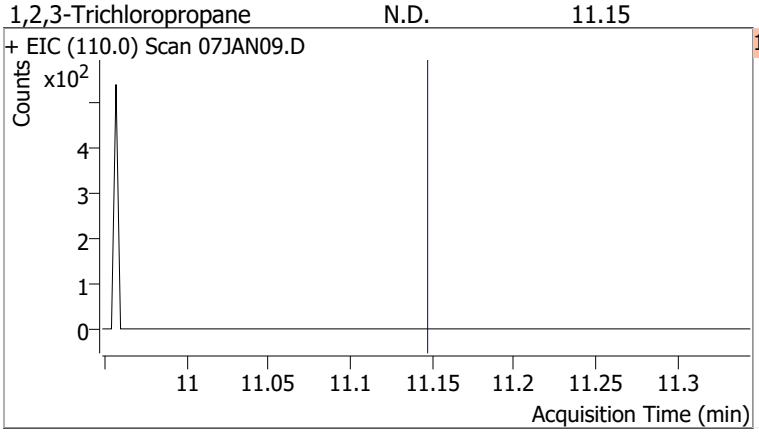


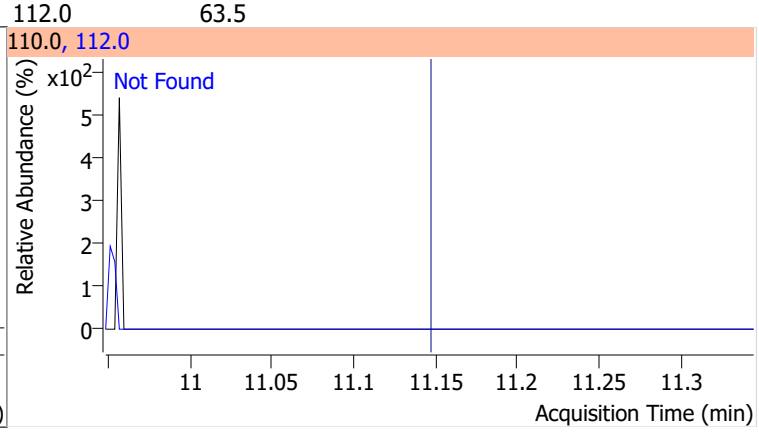
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



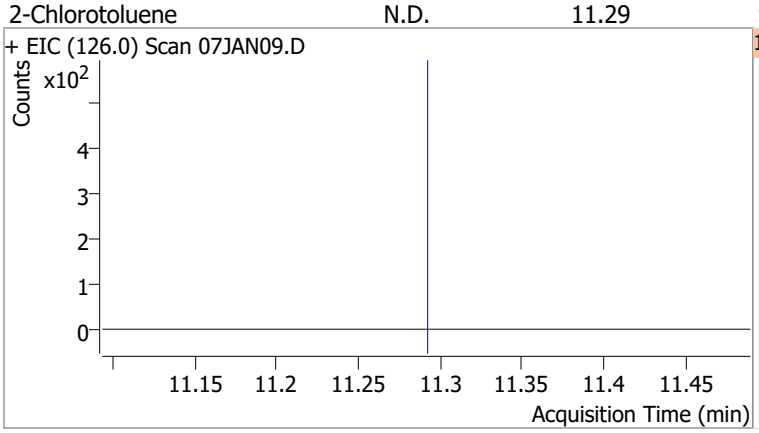


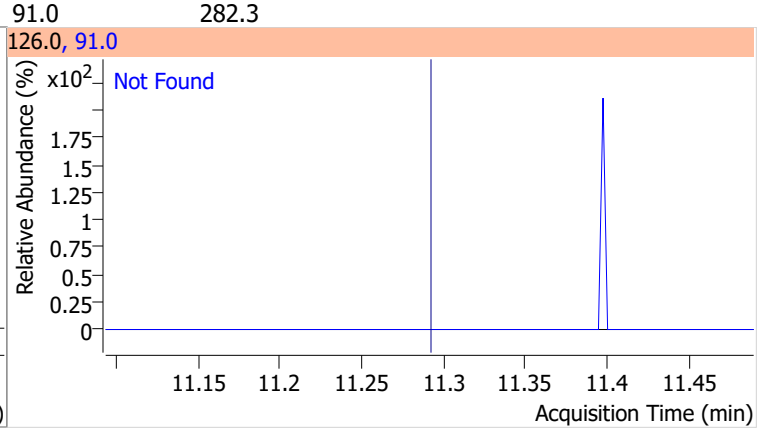
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



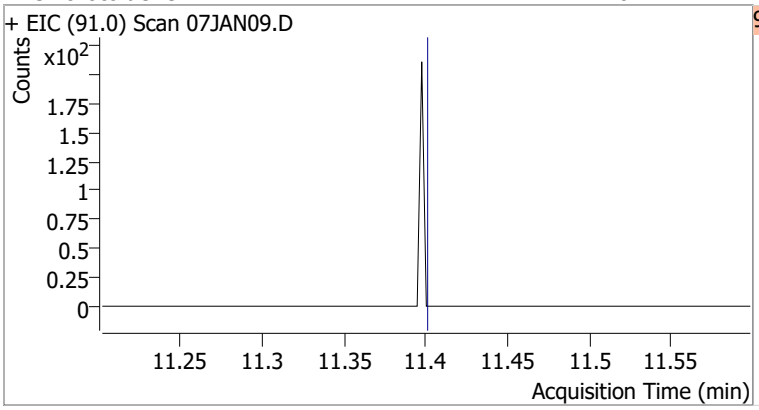


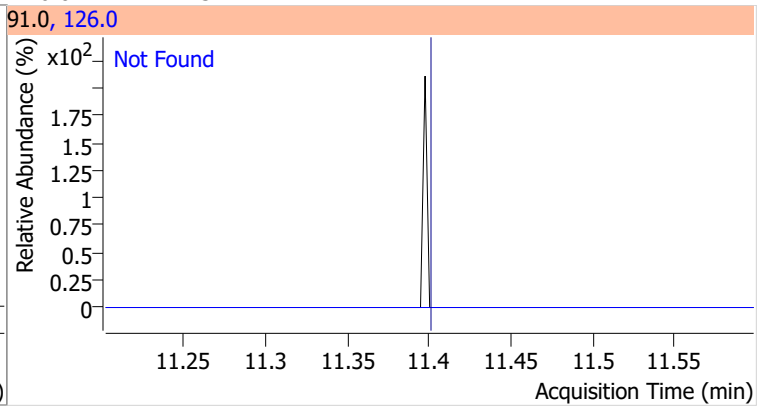
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



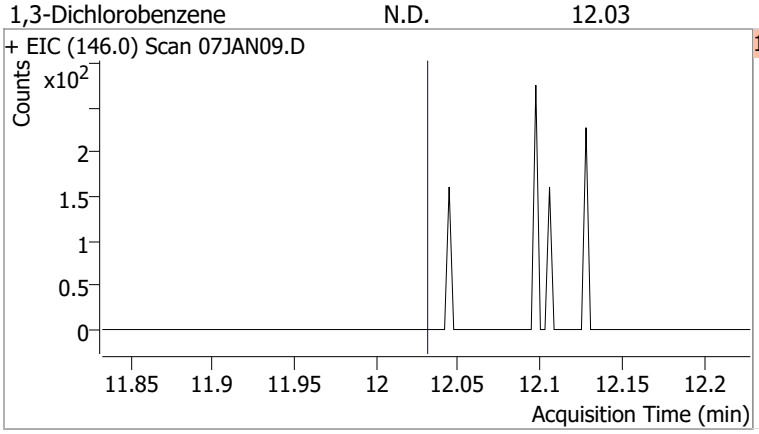


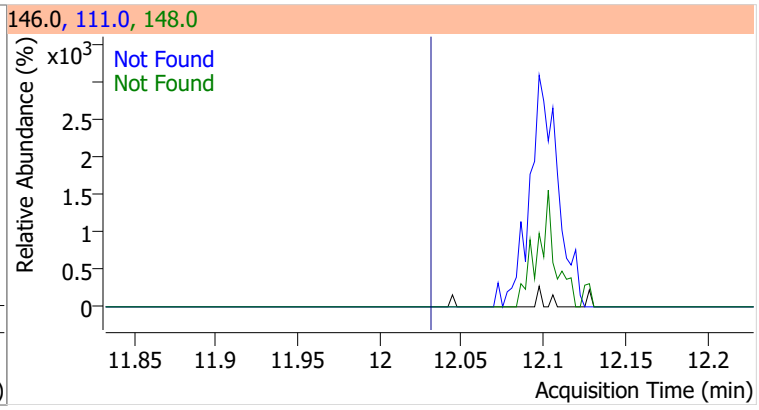
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



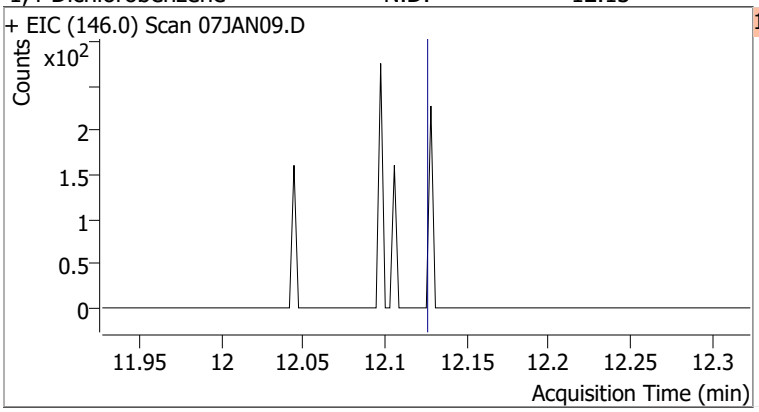


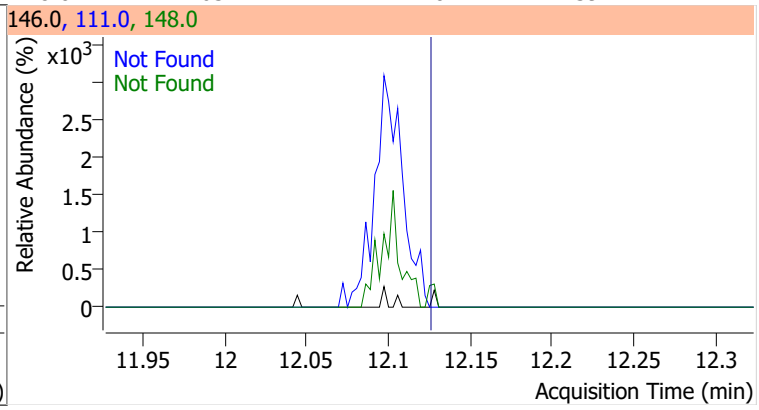
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



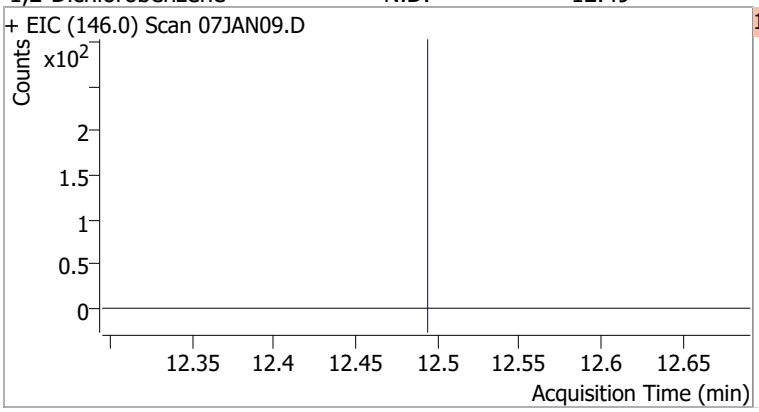


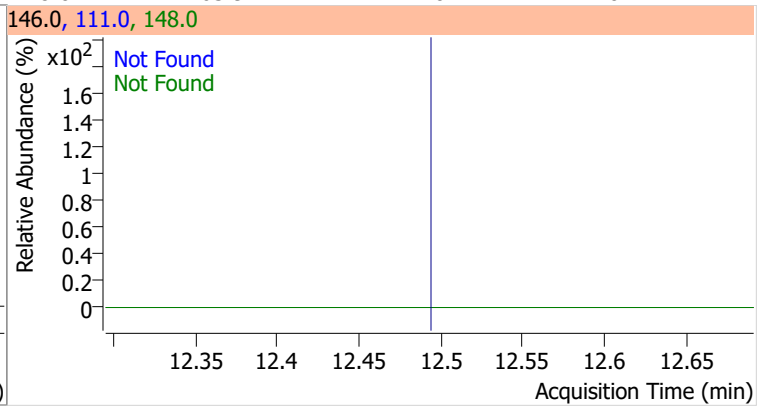
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





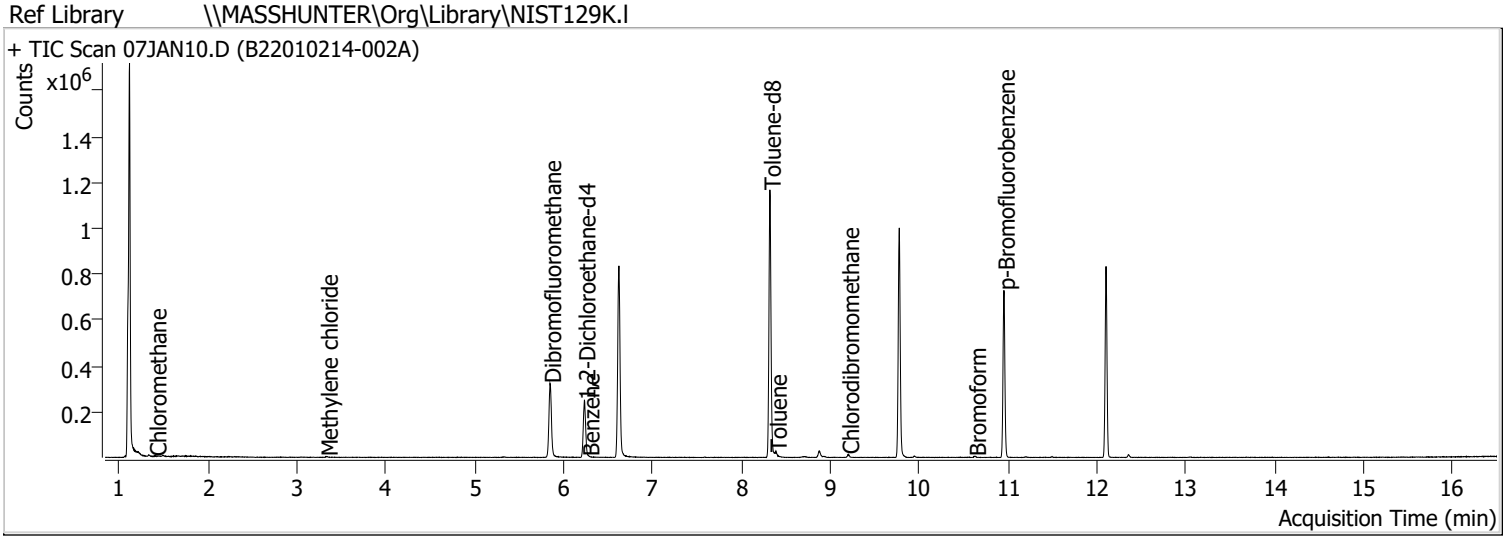
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN10.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 1:46:47 PM
Sample Name	B22010214-002A	Instrument	VOA5975C
Vial	10	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

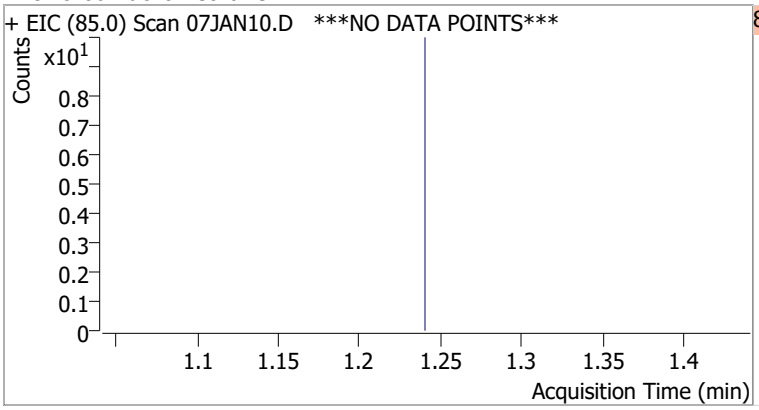


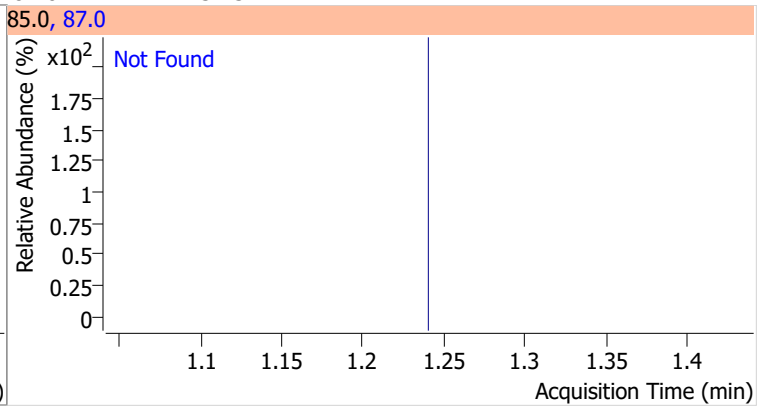
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	710622	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	276679	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	197130	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	189885	283.6314	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 113.45%			
S 1,2-Dichloroethane-d4	6.233	67.0	88230	305.1186	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 122.05%		*	
S Toluene-d8	8.319	98.0	699981	262.5368	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 105.01%			
S p-Bromofluorobenzene	10.951	95.0	203968	282.4305	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 112.97%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.409	50.0	2332	2.0632	ng		99
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.336	49.0	1975	1.8718	ng	m	96
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	5.656	83.0	0		ng	md	1

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.278	78.0	229	0.0810	ng	m	81
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	7.580	83.0	0		ng	md	1
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.386	92.0	5263	2.9222	ng		91
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	9.200	129.0	5690	10.1448	ng	m	97
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	0.000		0	N.D.			
T m+p-Xylenes	10.037	106.0	0		ng	md	1
T o-Xylene	0.000		0	N.D.			
T Styrene	0.000		0	N.D.			
T Bromoform	10.628	172.5	2536	10.0531	ng		95
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

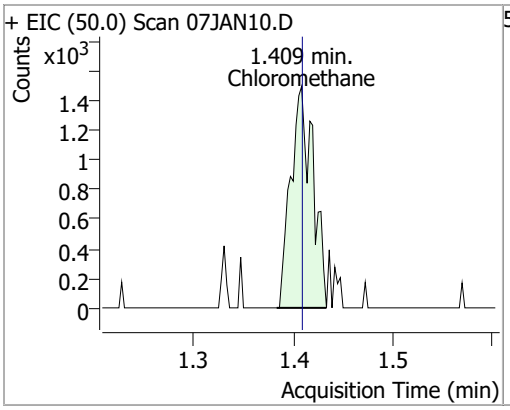
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

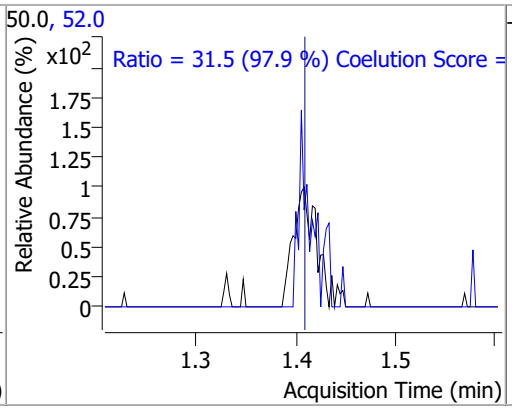
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

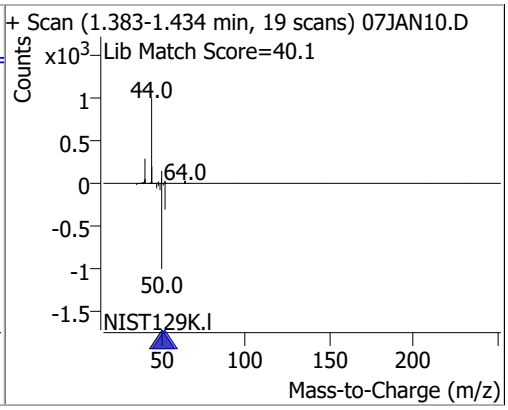




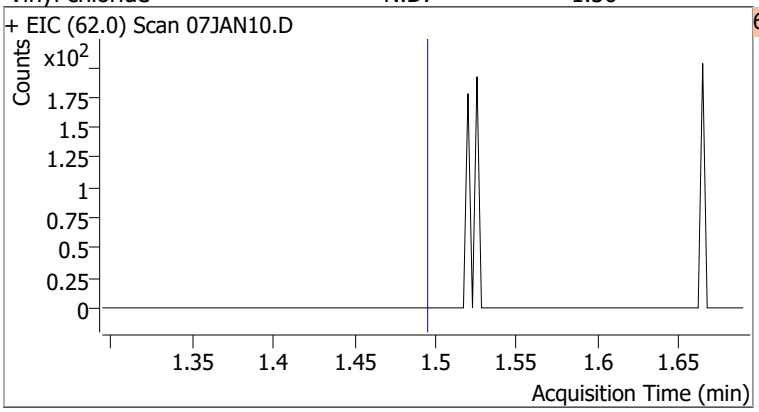
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	2.0632	1.41	0.00	2332	52.0	31.5	2.1	62.1

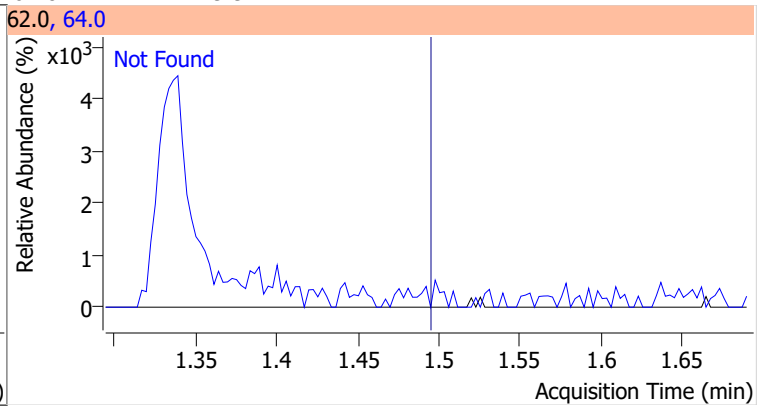




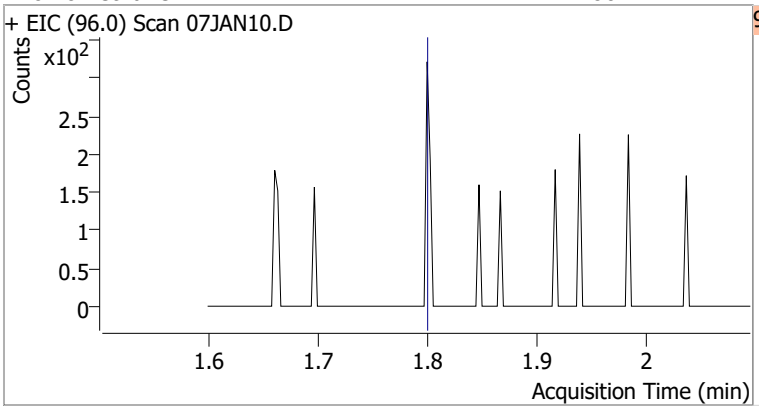


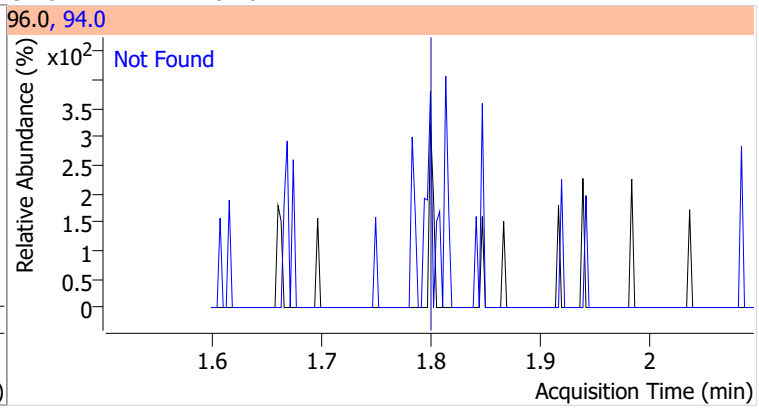
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



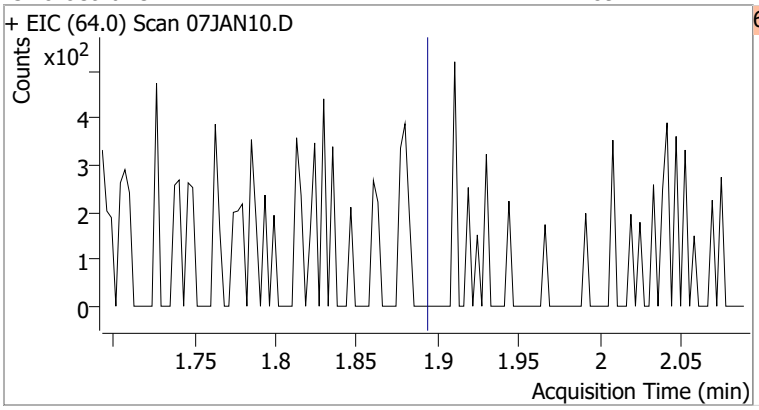


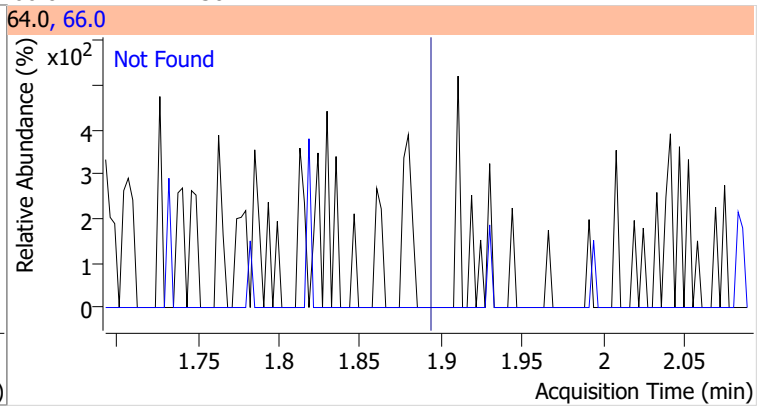
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



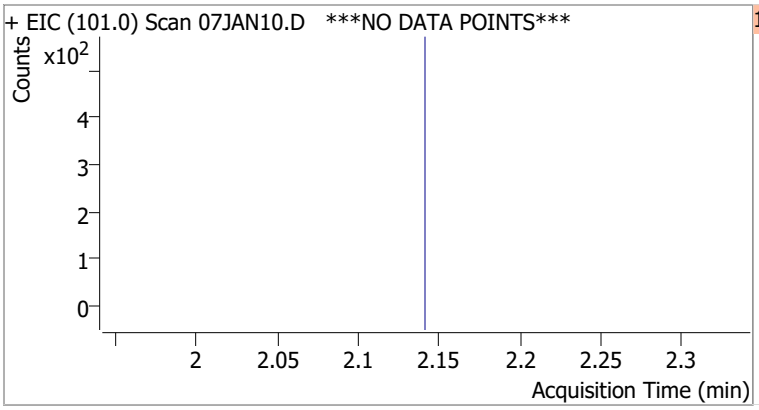


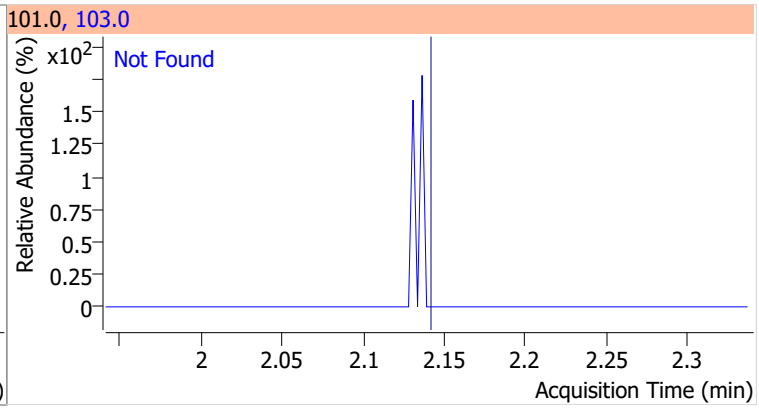
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



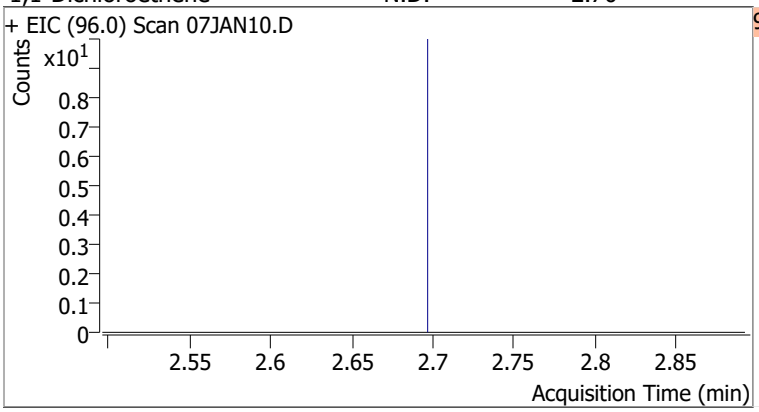


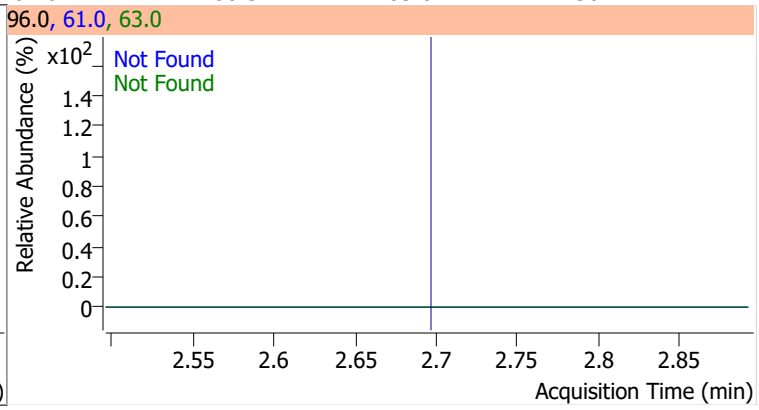
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



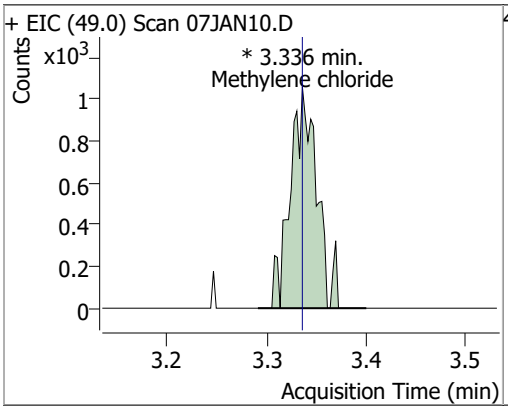


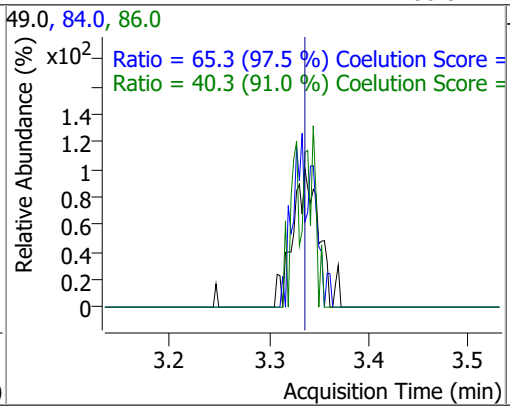
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

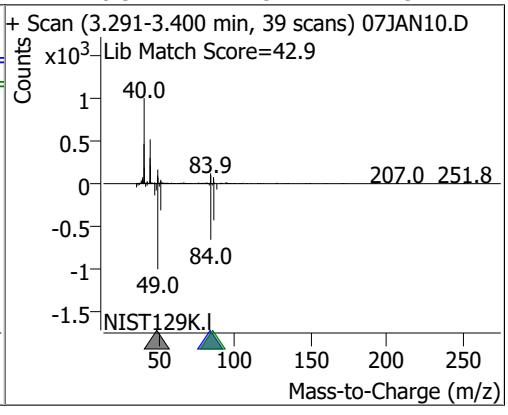




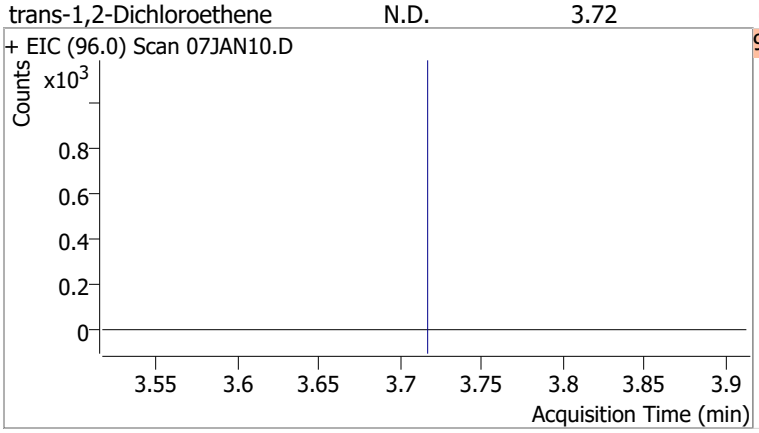
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	1.8718	3.34	0.00	1975 (m)	84.0	65.3	36.9	96.9
					86.0	40.3	14.3	74.3

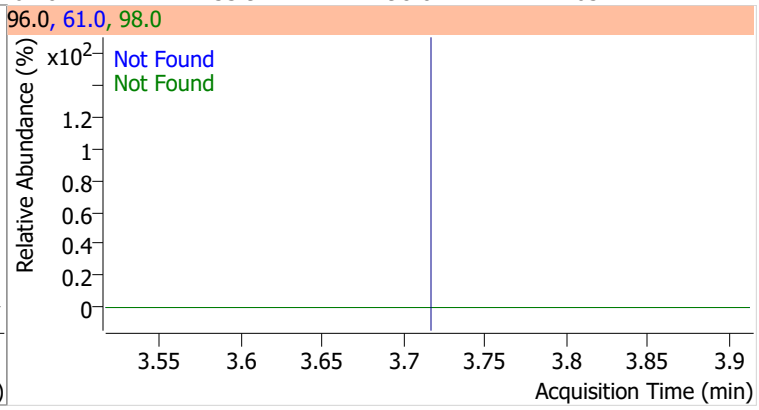




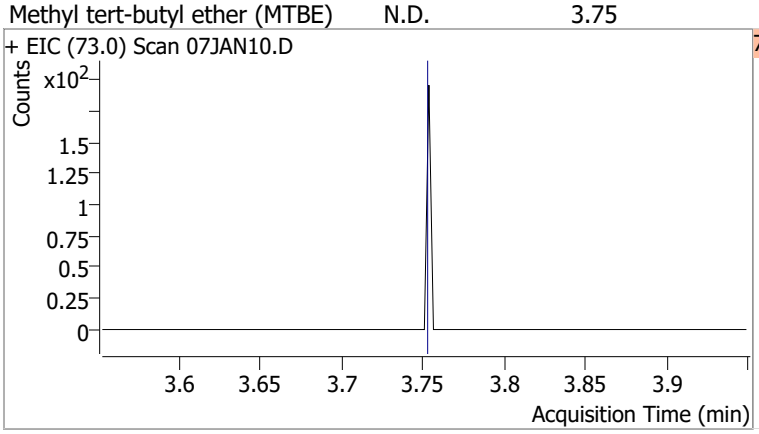


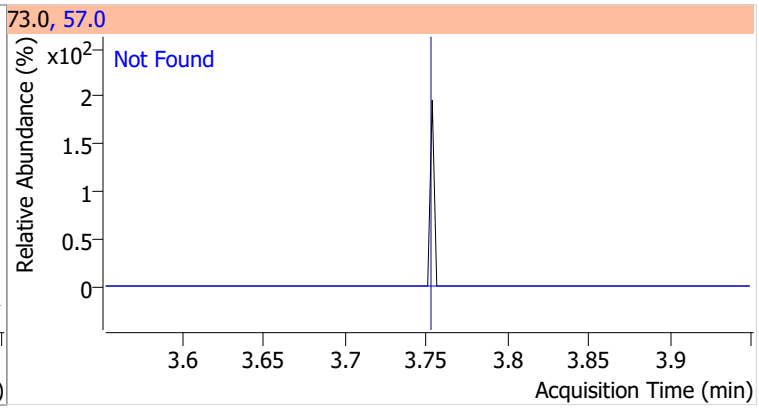
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



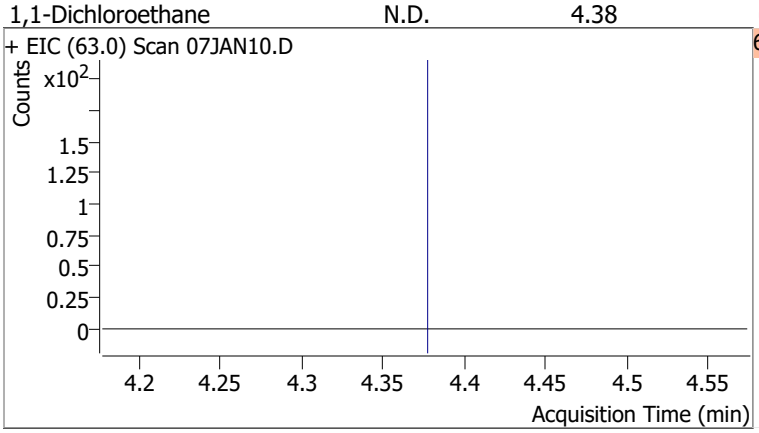


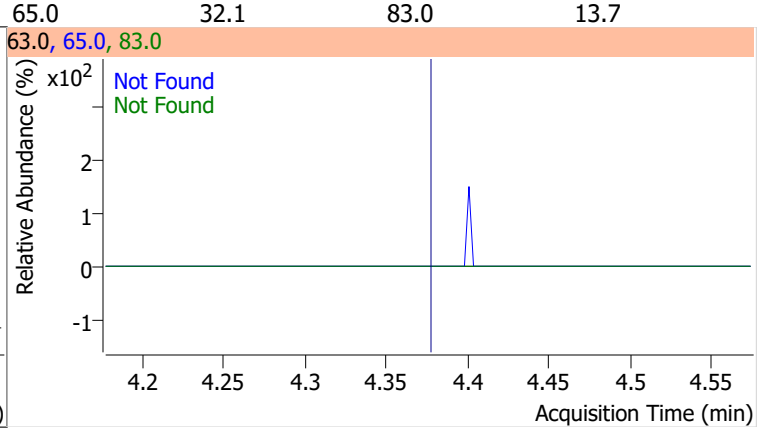
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



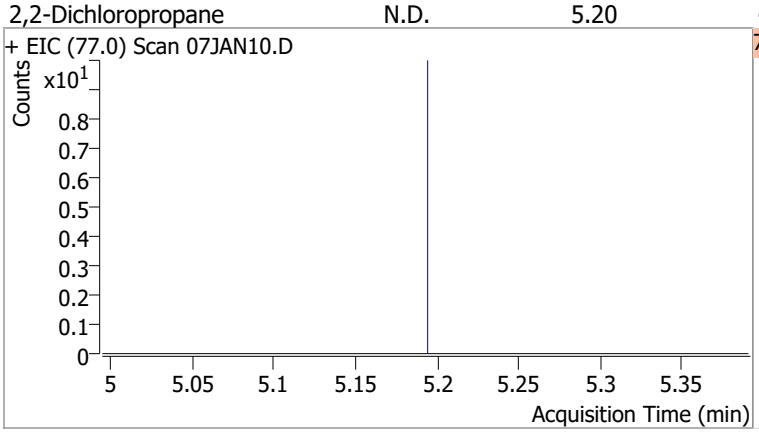


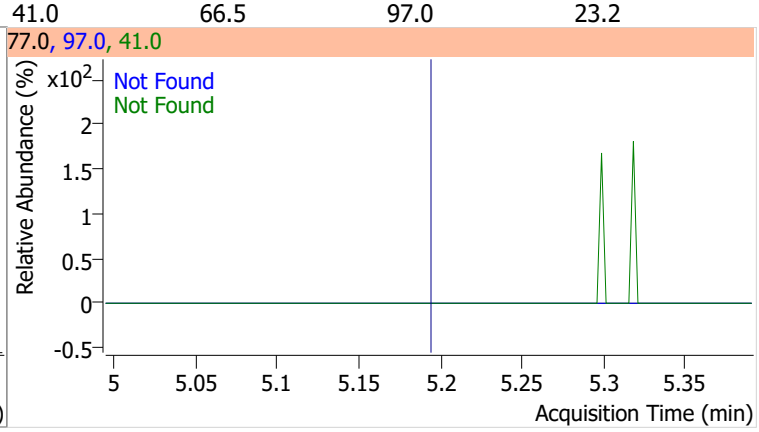
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



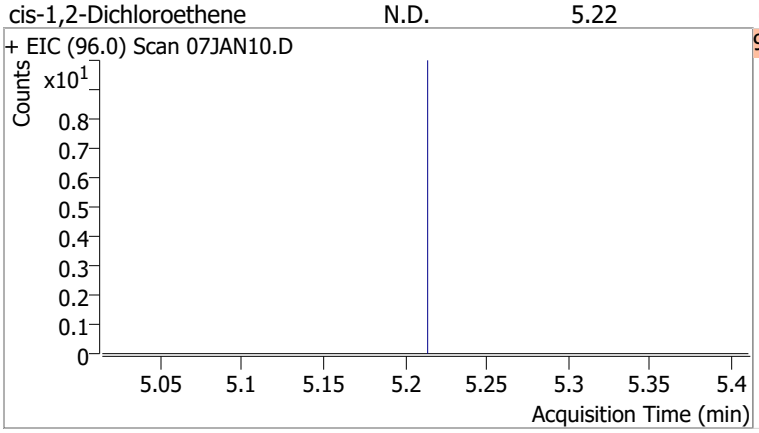


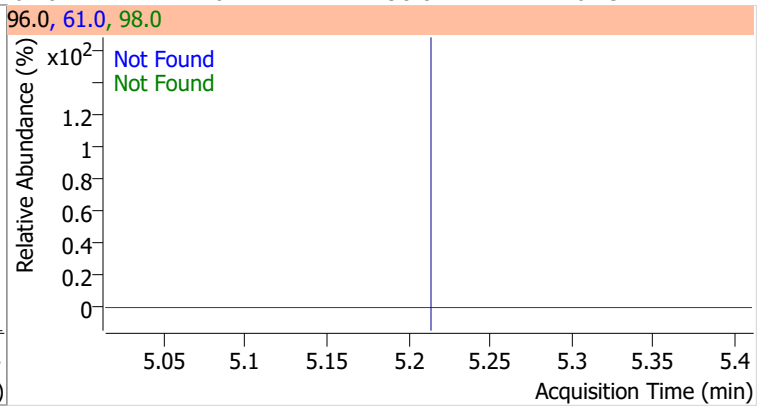
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



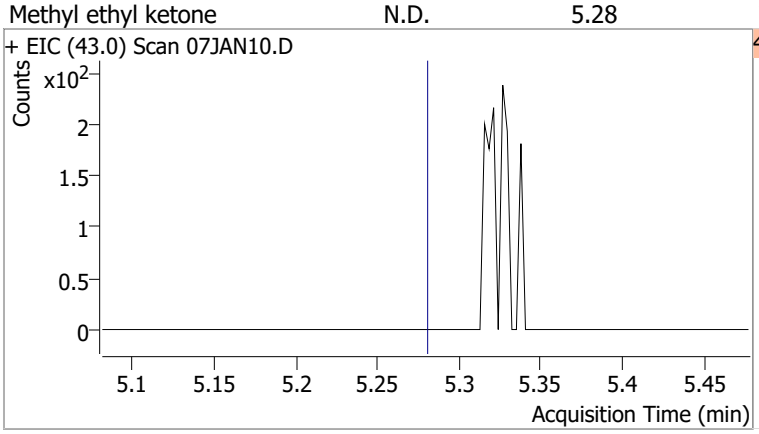


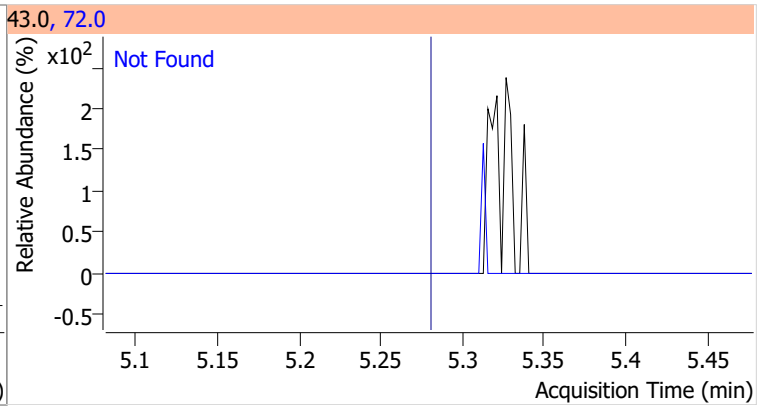
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



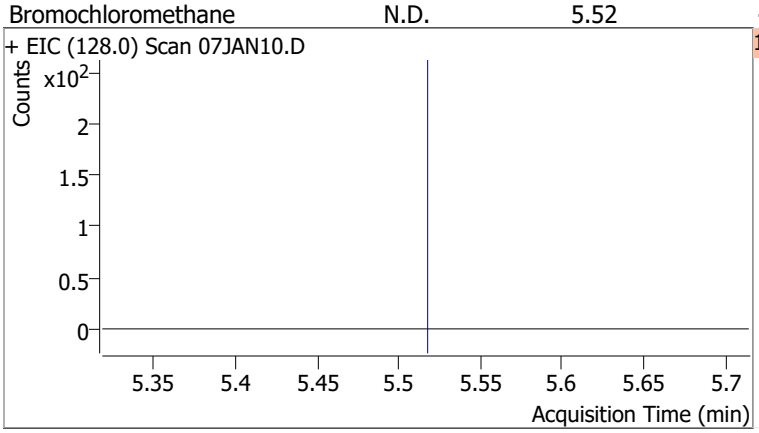


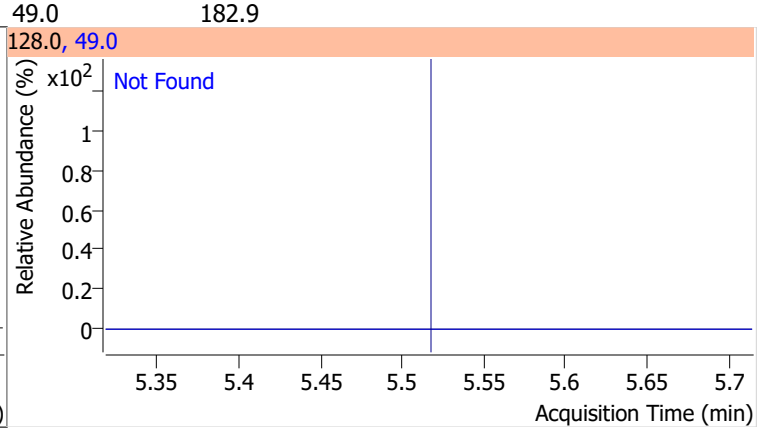
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



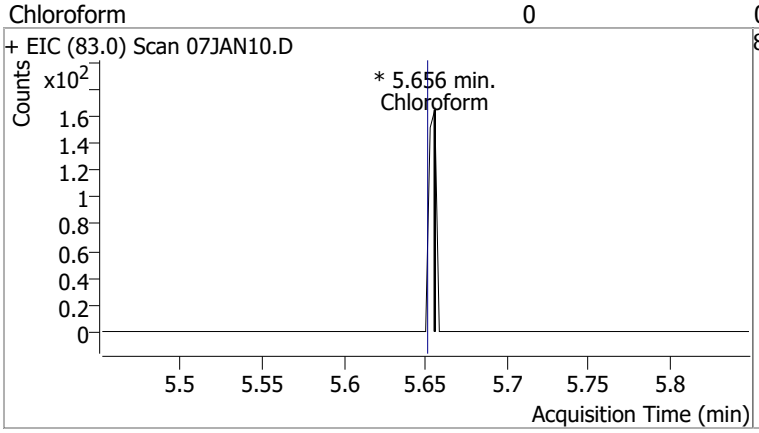


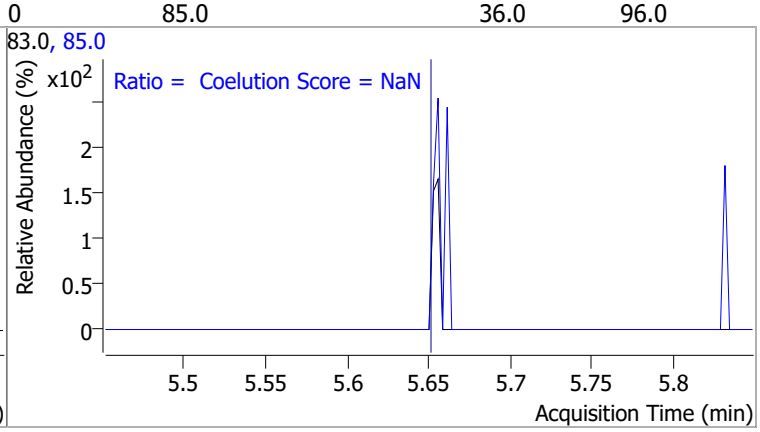
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9



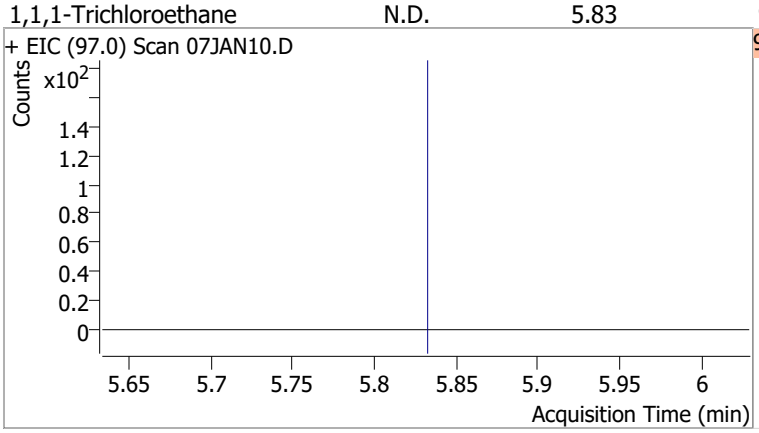


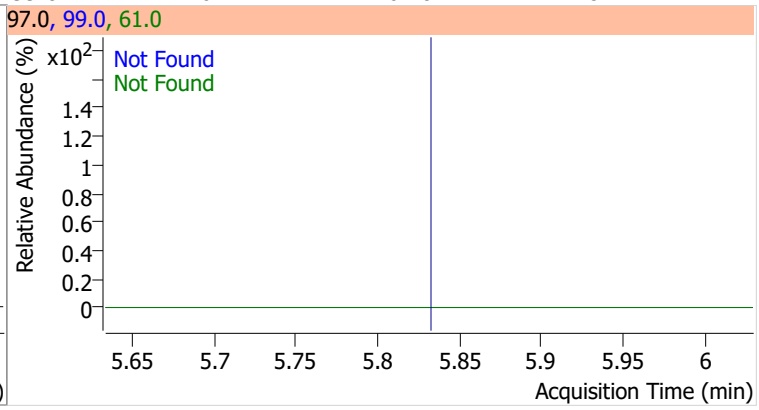
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform		0		0	85.0		36.0	96.0



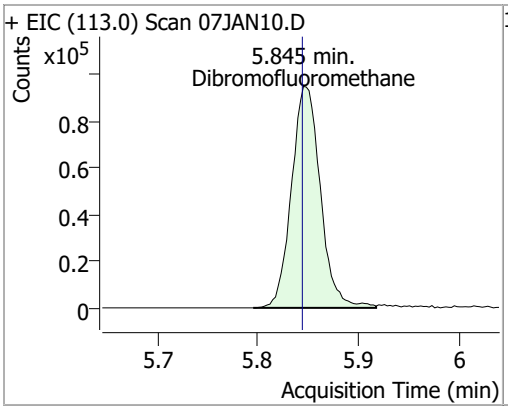


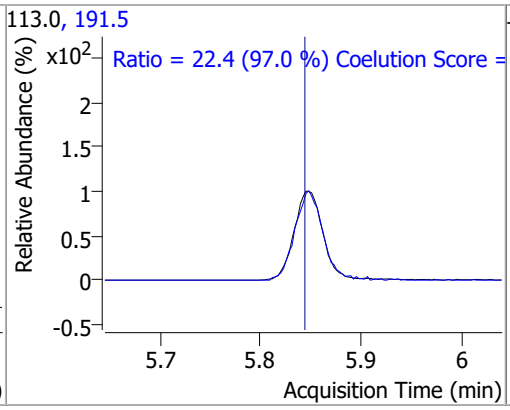
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

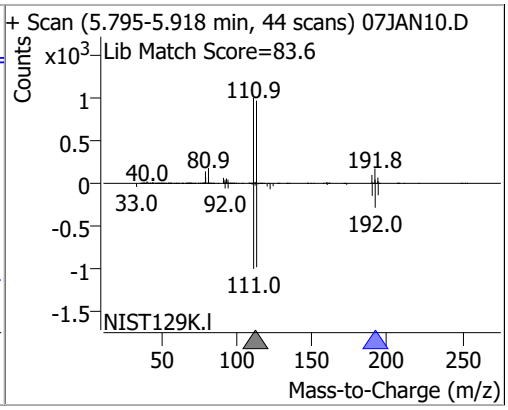




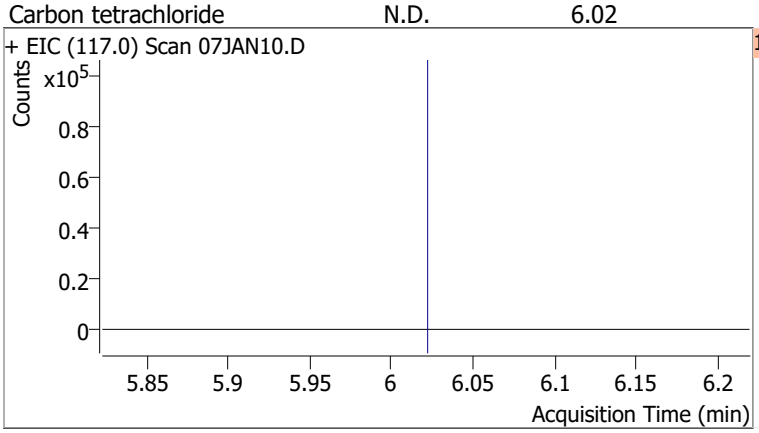
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	283.6314	5.85	0.00	189885	191.5	22.4	0.0	53.1

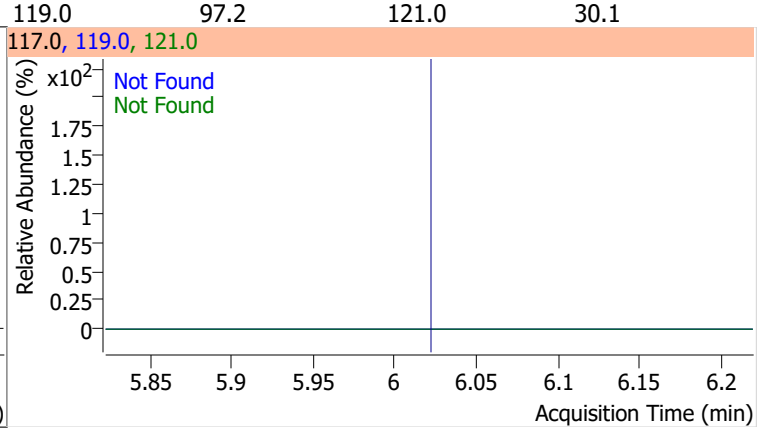




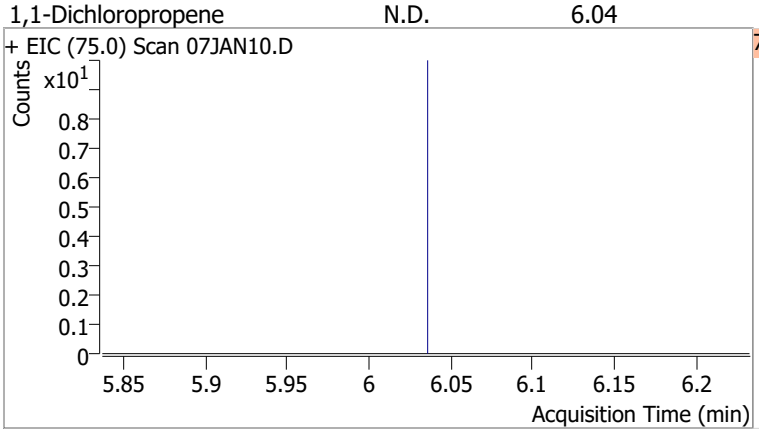


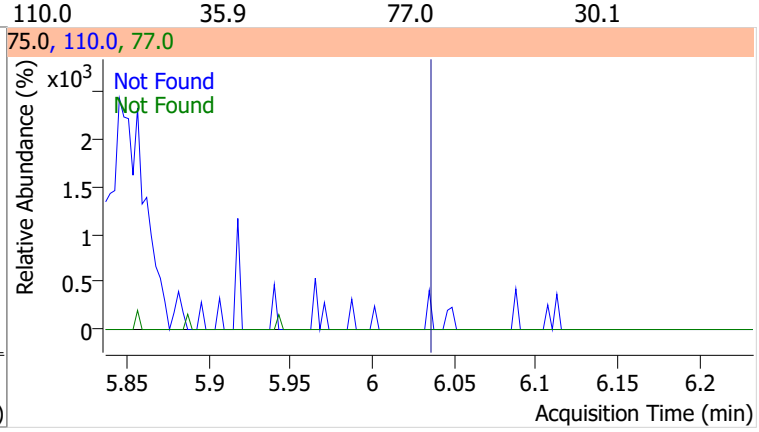
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



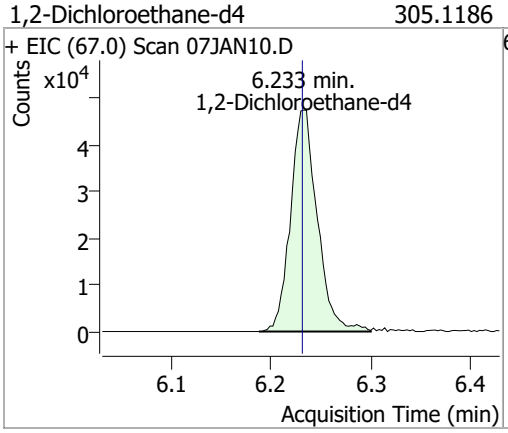


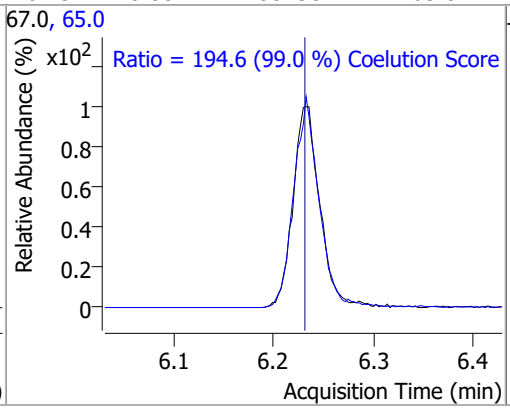
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

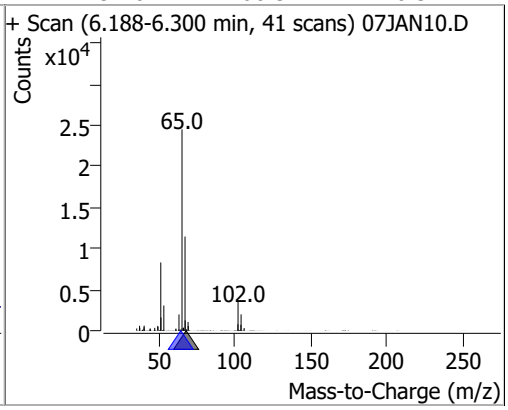




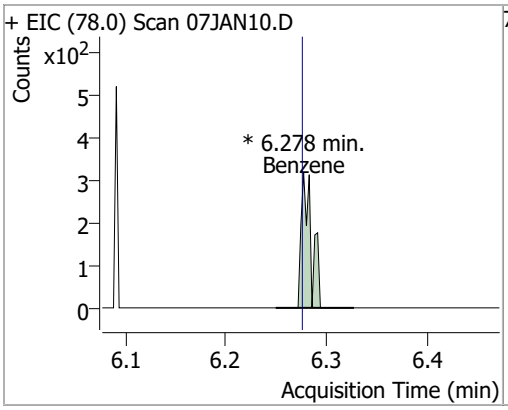
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	305.1186	6.23	0.00	88230	65.0	194.6	166.5	226.5

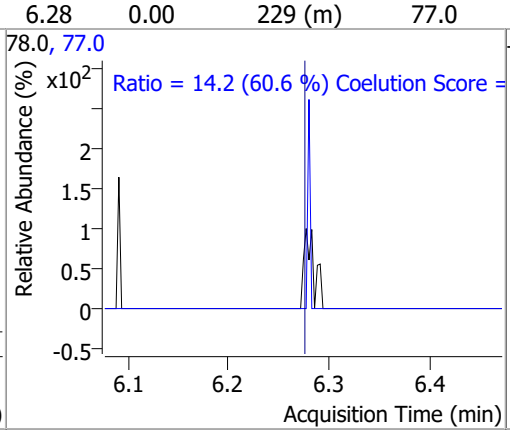


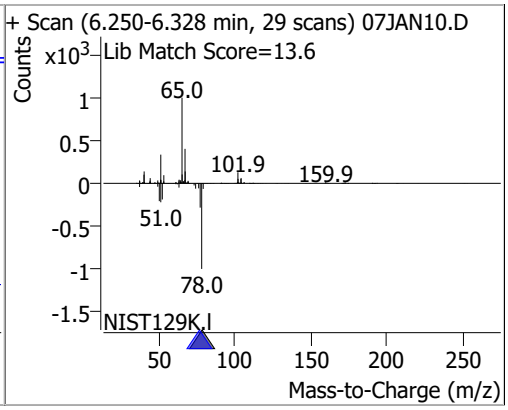




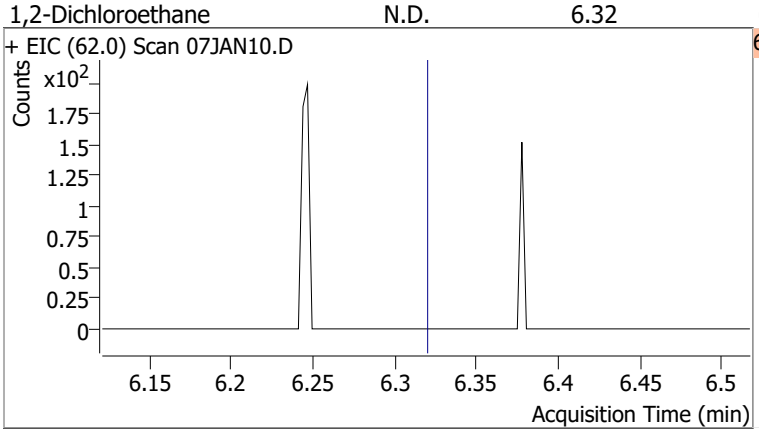
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.0810	6.28	0.00	229 (m)	77.0	14.2	0.0	53.5

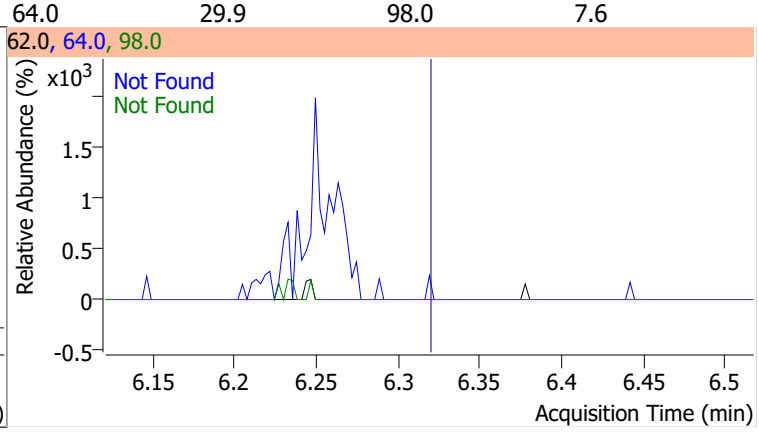




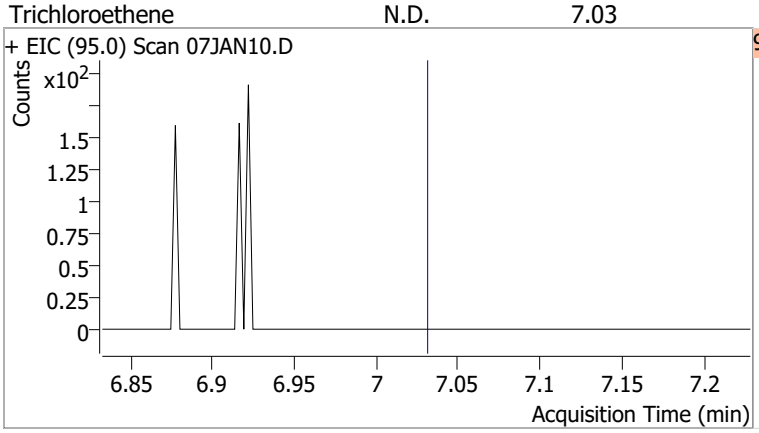


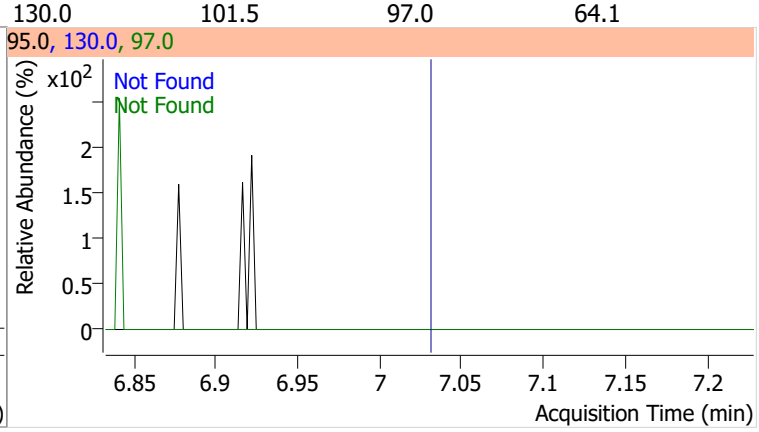
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



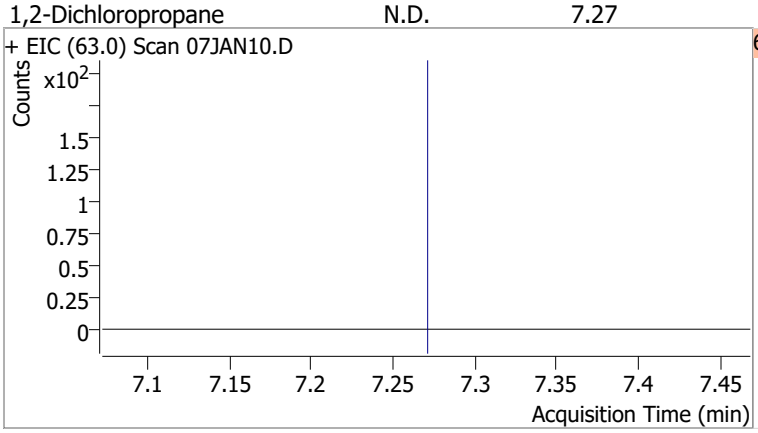


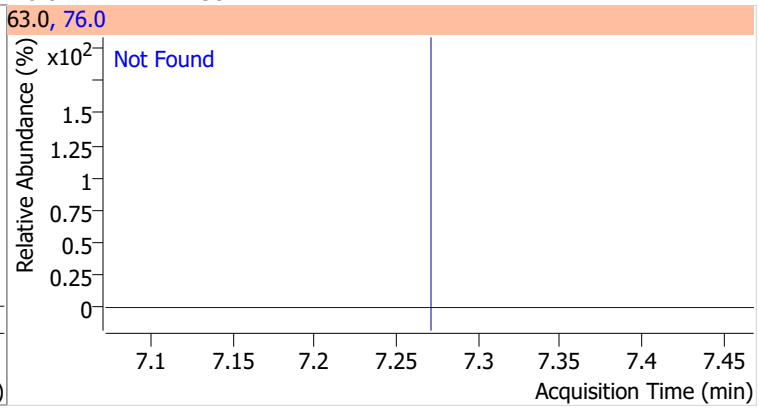
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



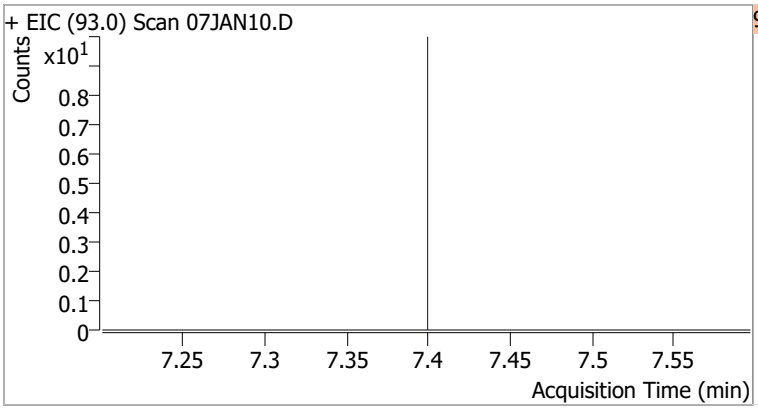


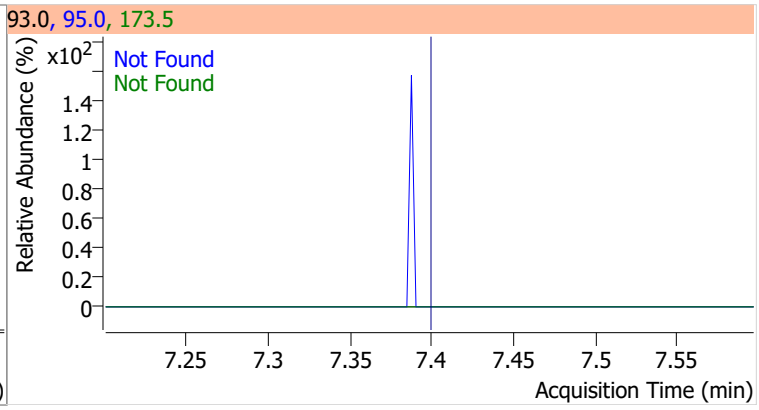
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



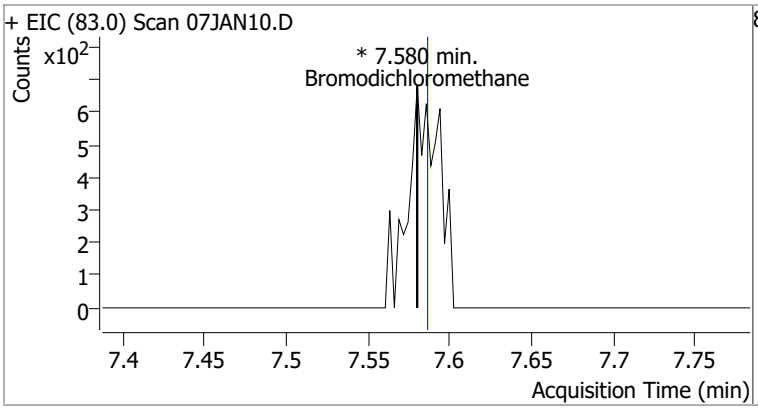


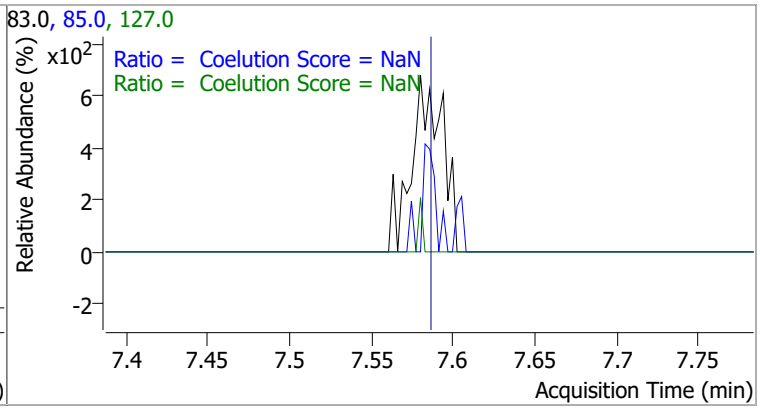
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



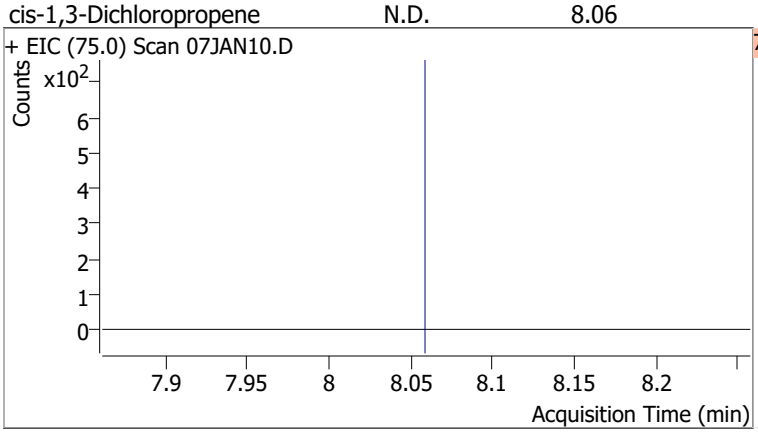


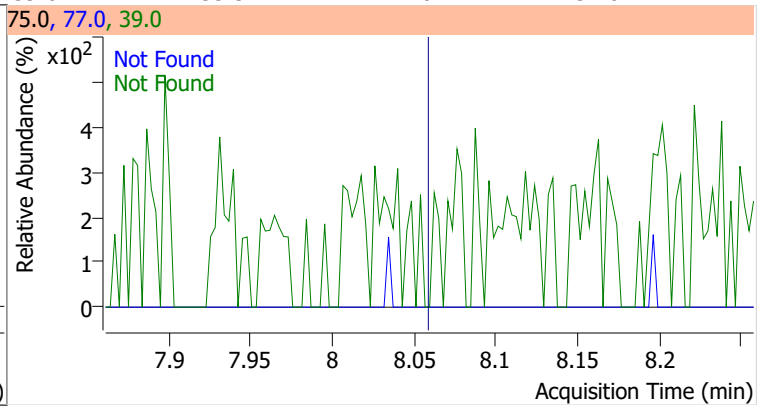
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane		0		0	85.0		34.5	94.5
					127.0		0.0	39.6



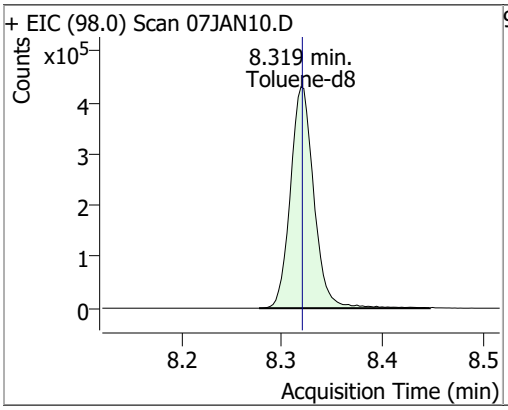


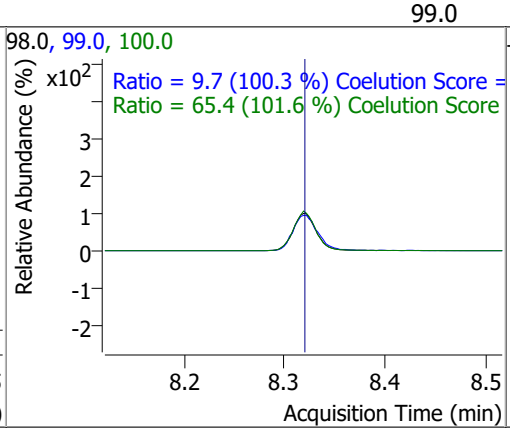
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

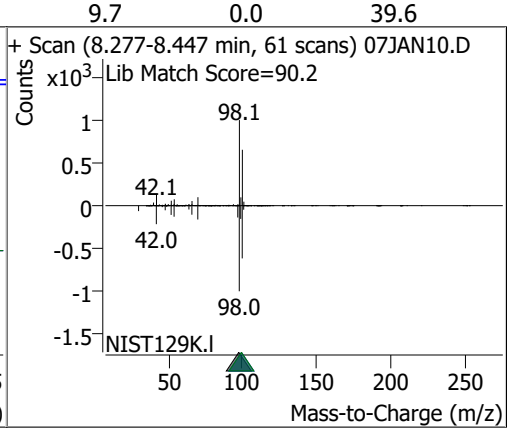




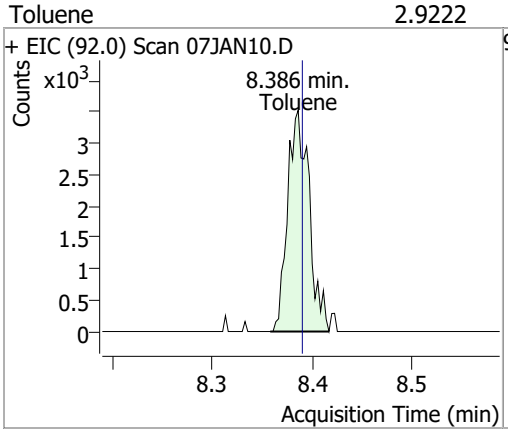
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	262.5368	8.32	0.00	699981	100.0	65.4	34.4	94.4
					99.0	9.7	0.0	39.6

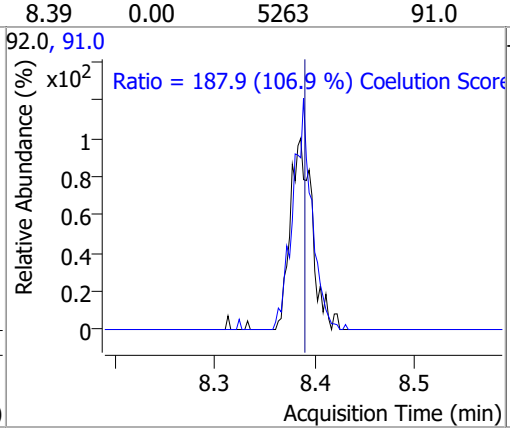


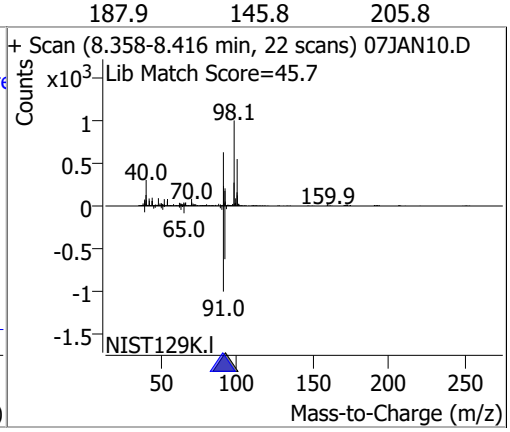




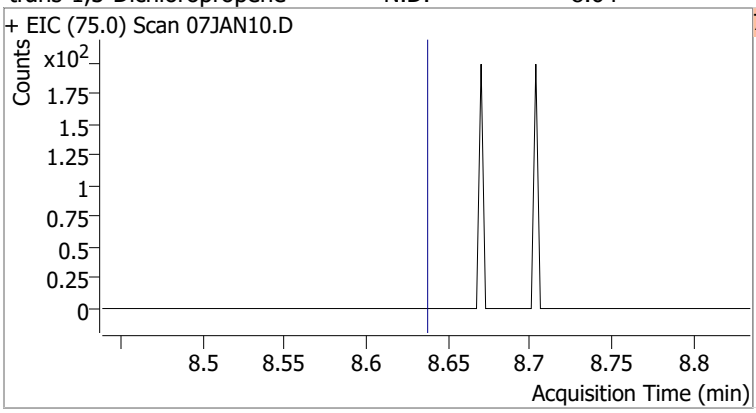
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	2.9222	8.39	0.00	5263	91.0	187.9	145.8	205.8

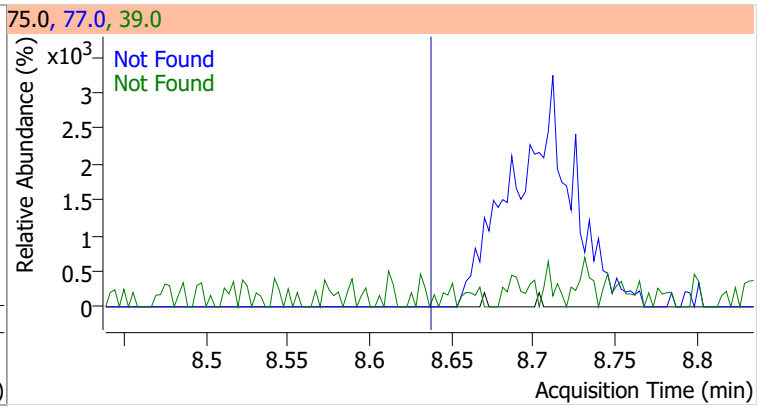




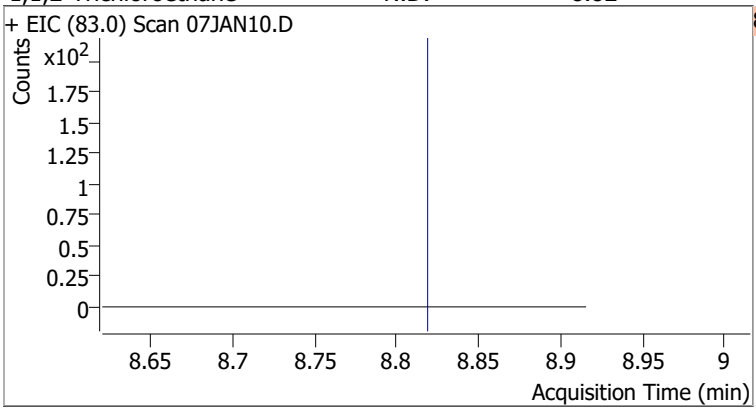


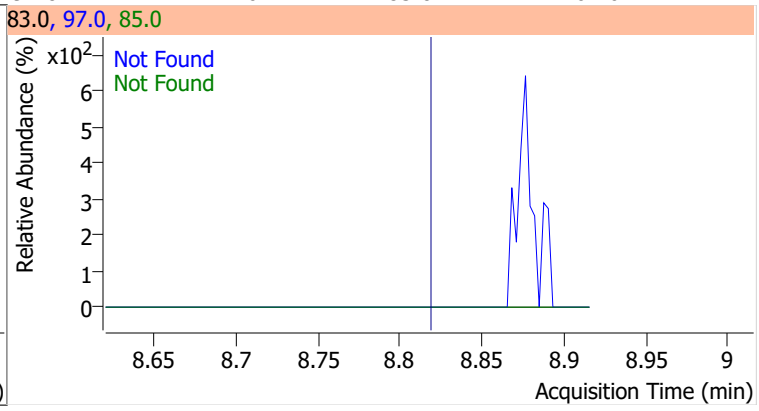
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



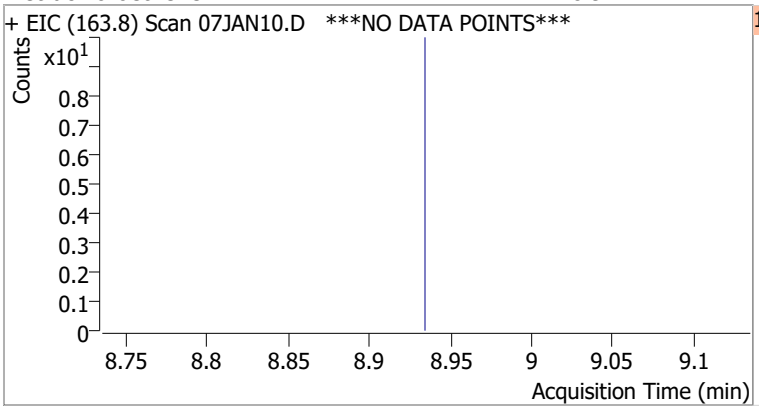


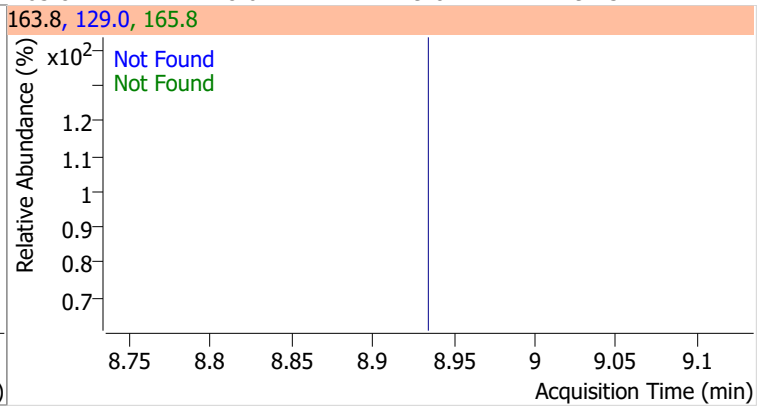
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



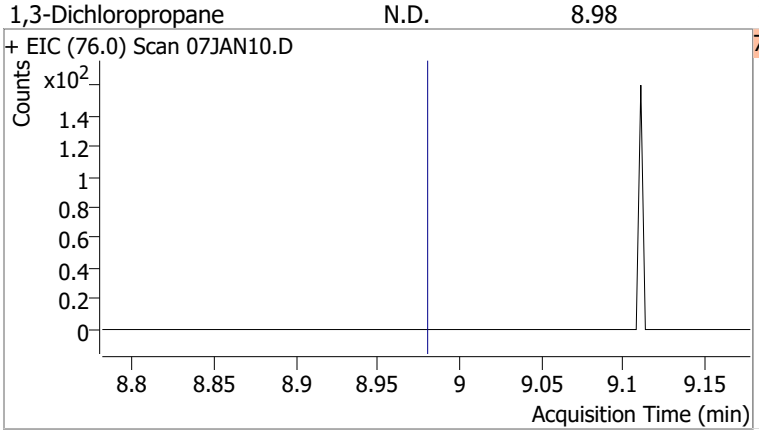


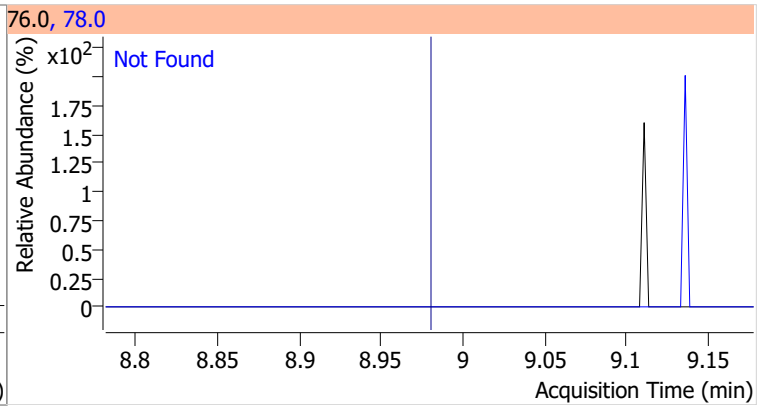
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



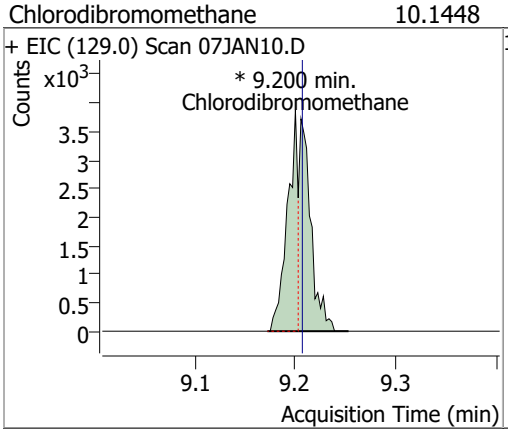


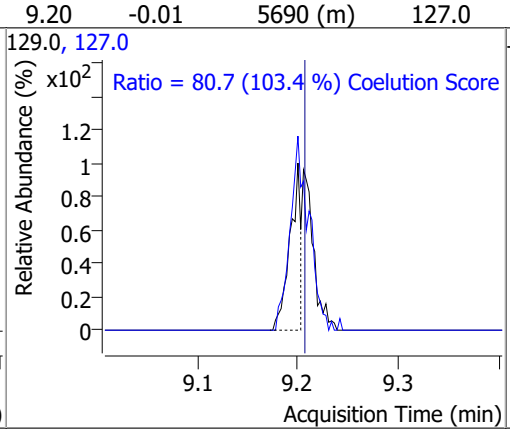
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9

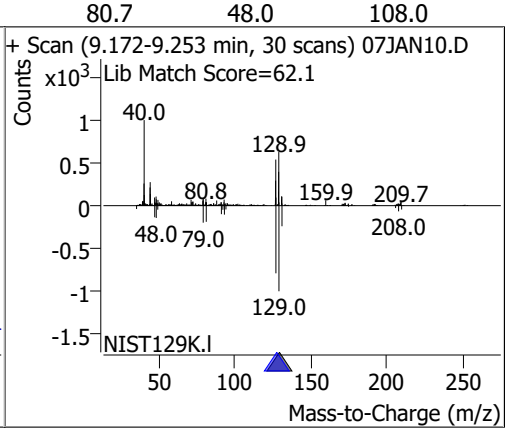




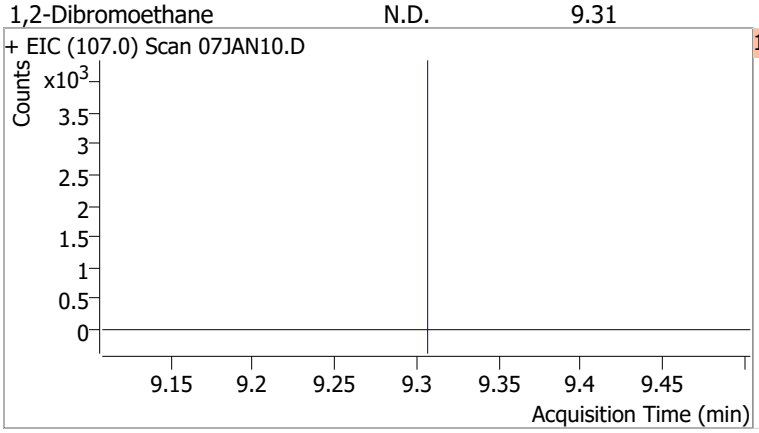
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	10.1448	9.20	-0.01	5690 (m)	127.0	80.7	48.0	108.0

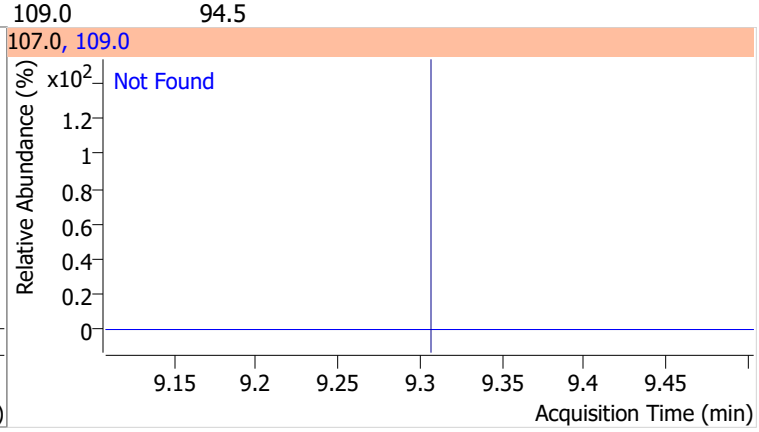




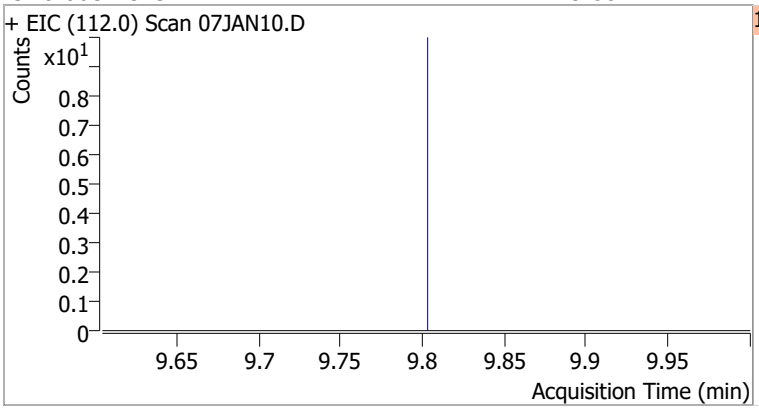


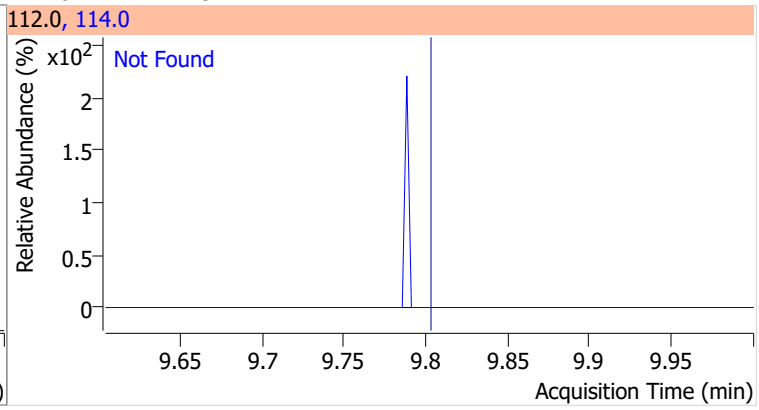
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



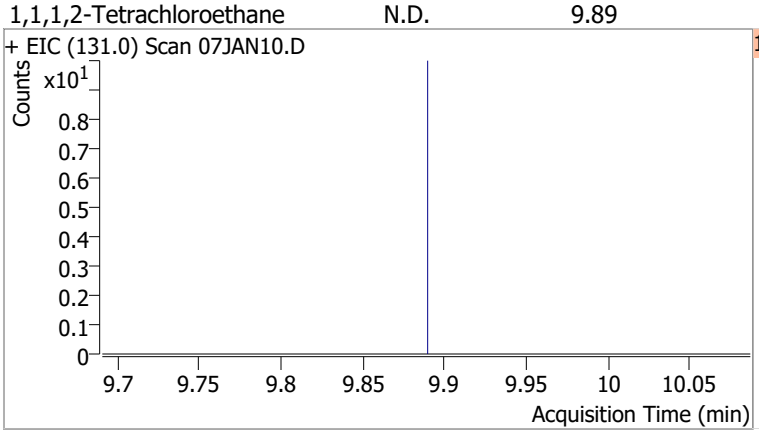


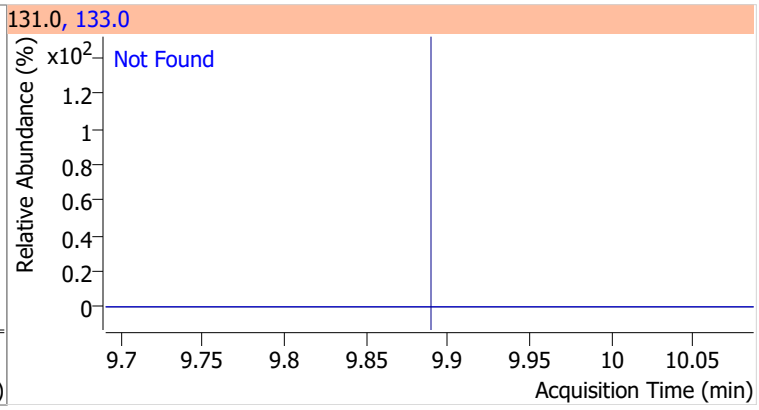
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



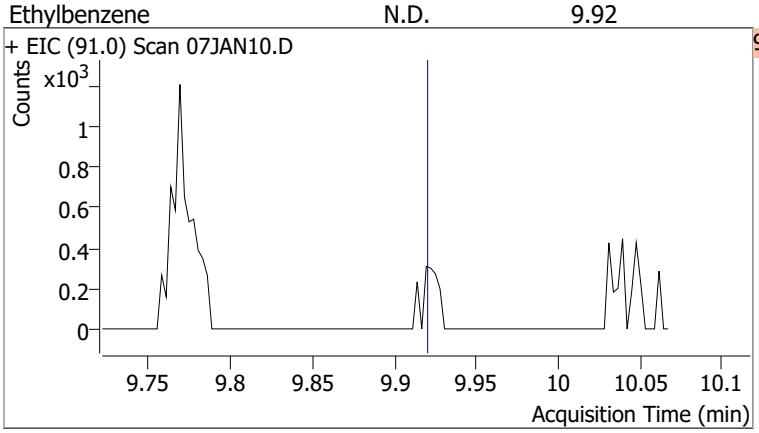


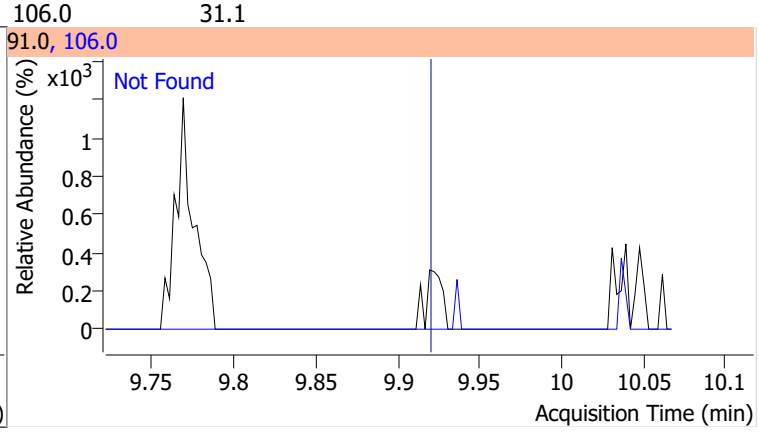
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



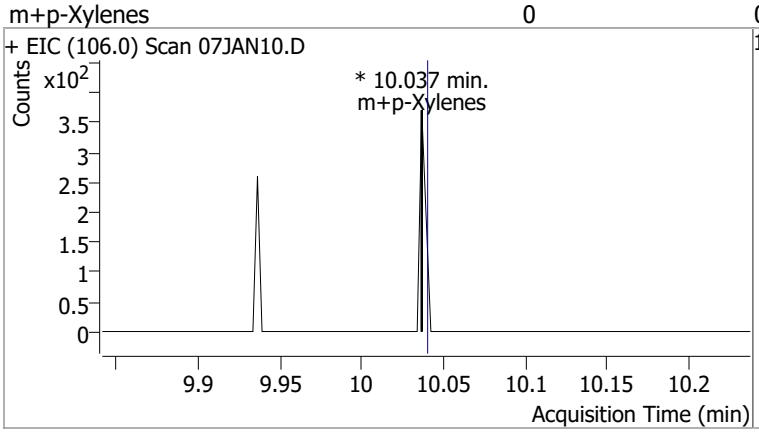


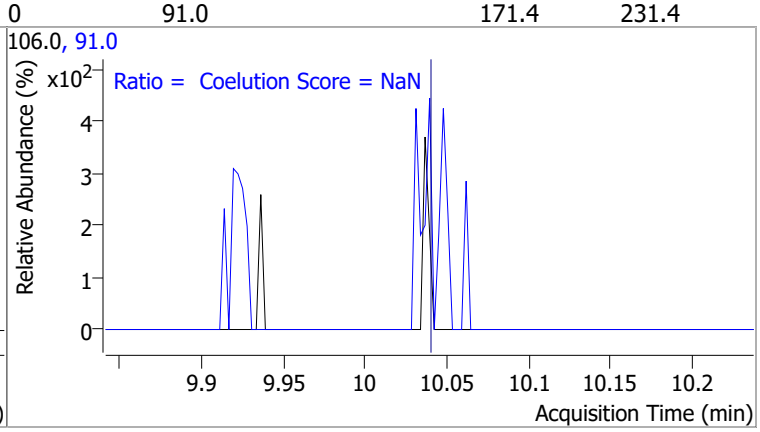
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1



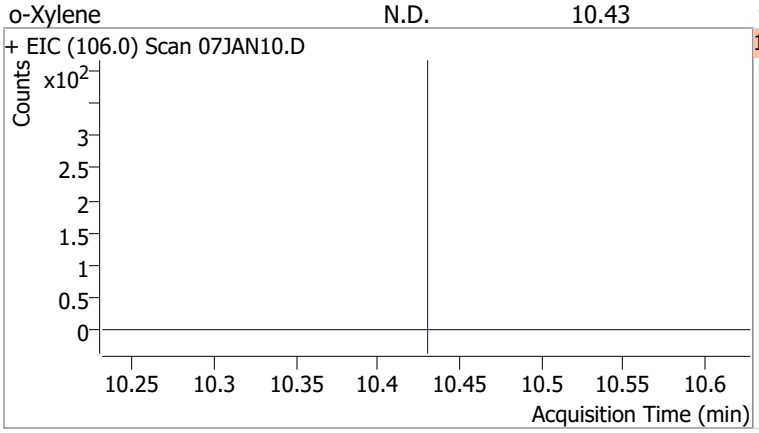


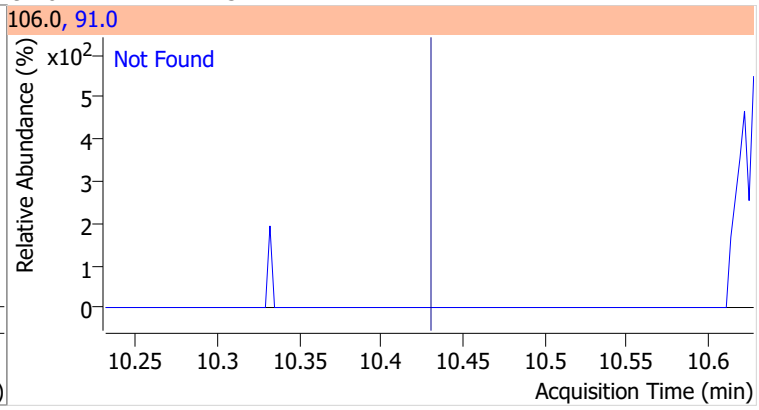
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes		0		0	91.0		171.4	231.4



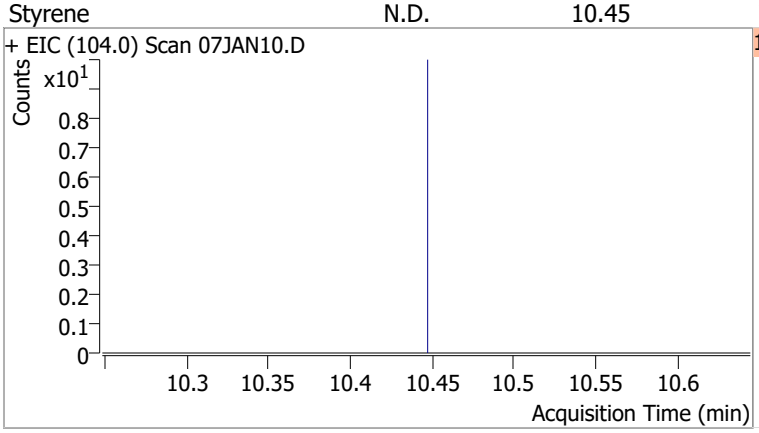


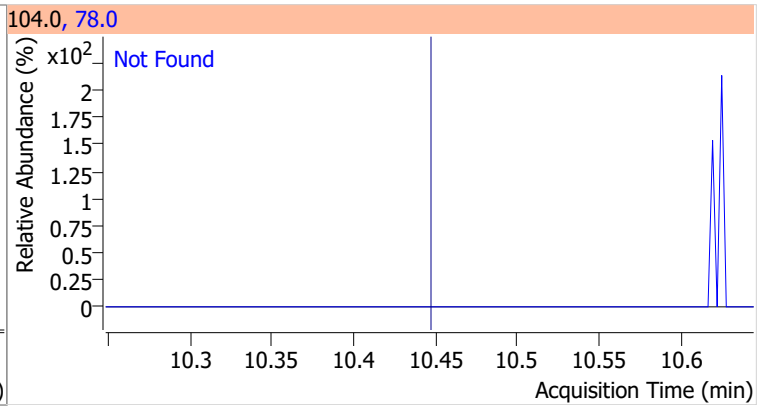
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



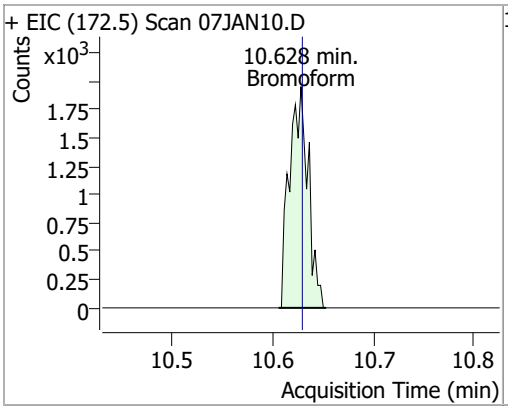


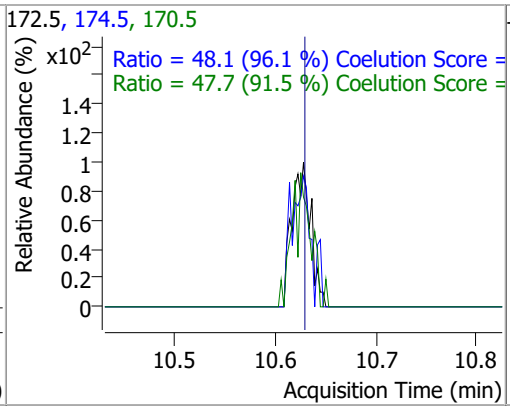
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6

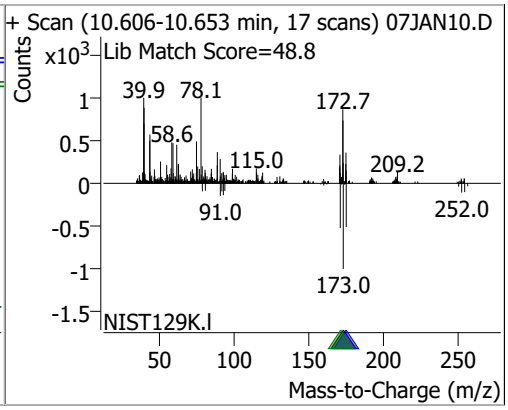




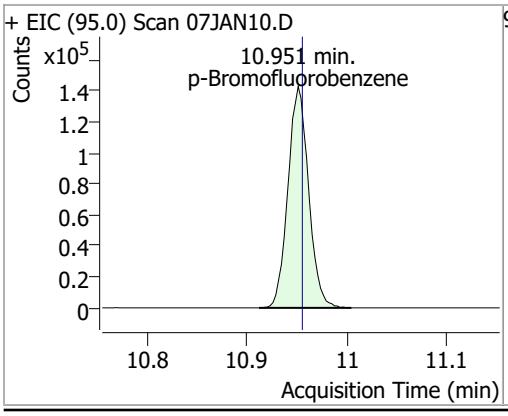
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	10.0531	10.63	0.00	2536	170.5	47.7	22.1	82.1
					174.5	48.1	20.1	80.1

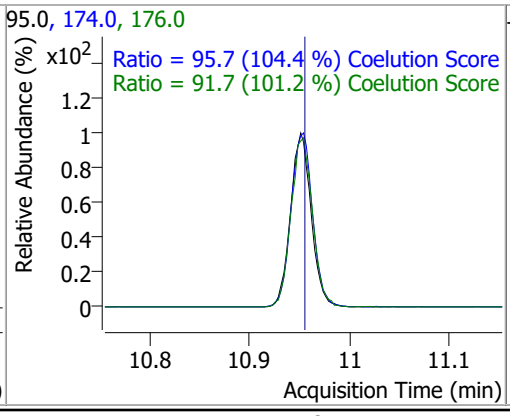


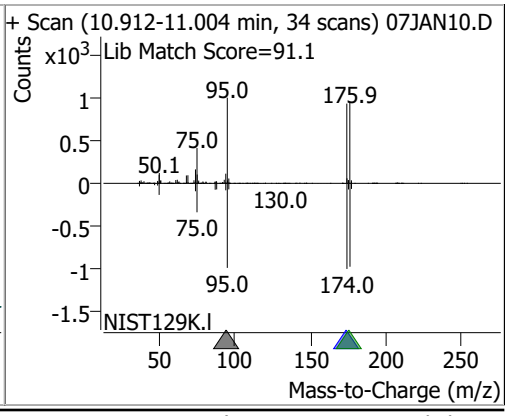




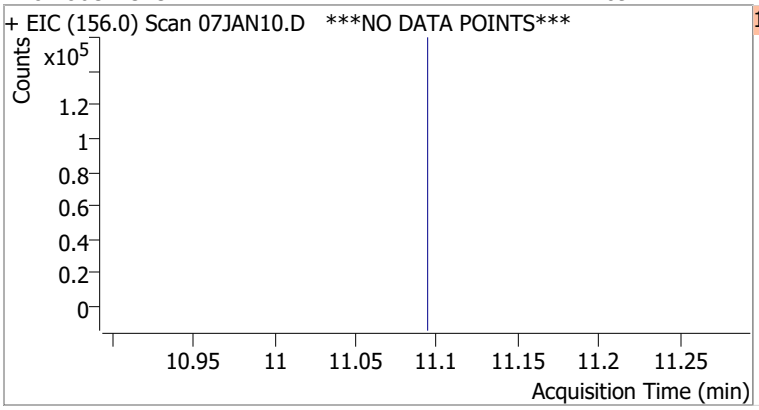
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	282.4305	10.95	0.00	203968	174.0	95.7	61.7	121.7
					176.0	91.7	60.6	120.6

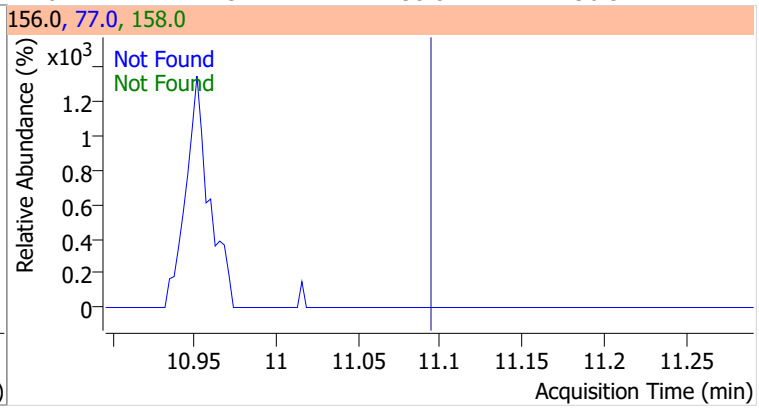




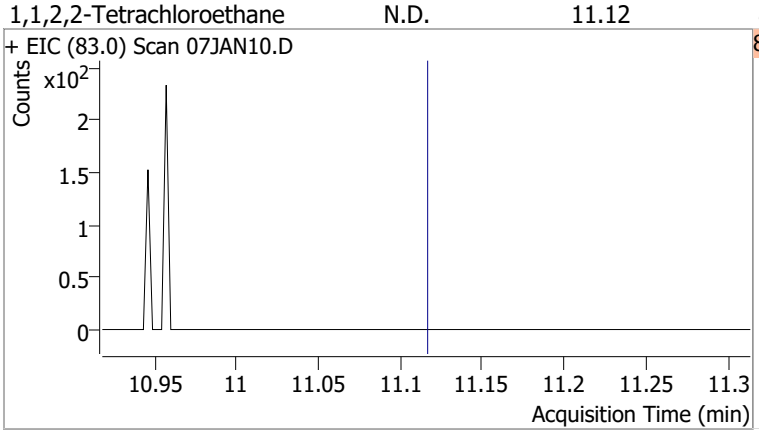


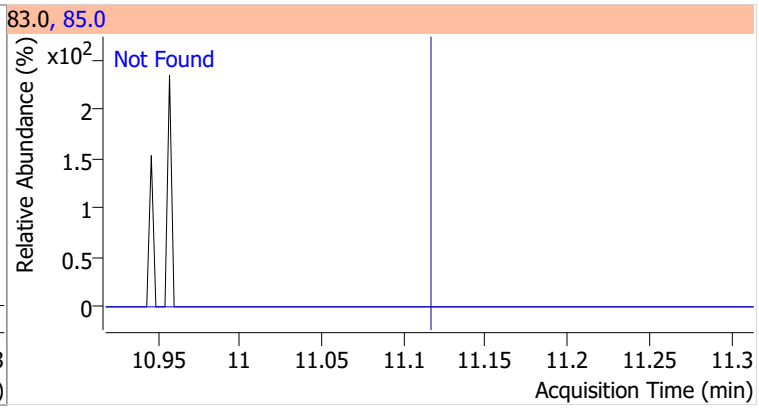
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



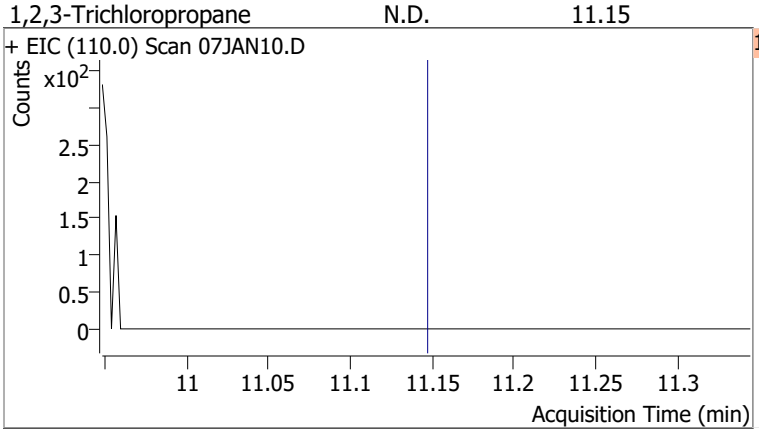


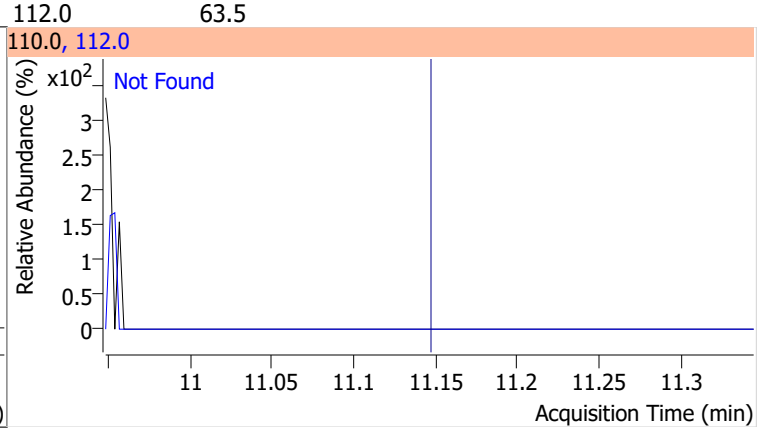
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



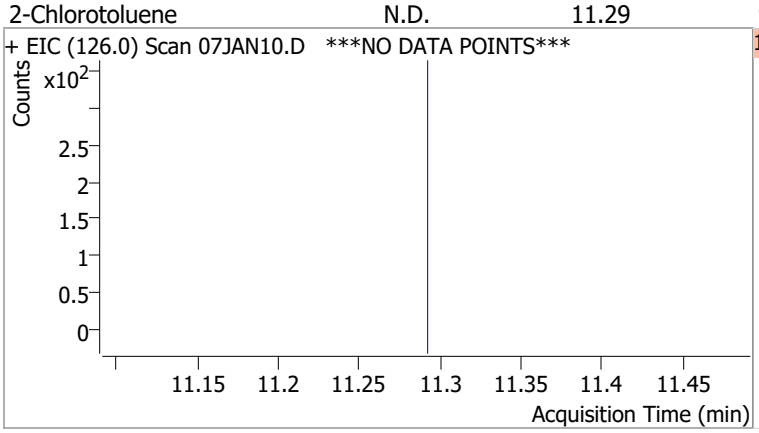


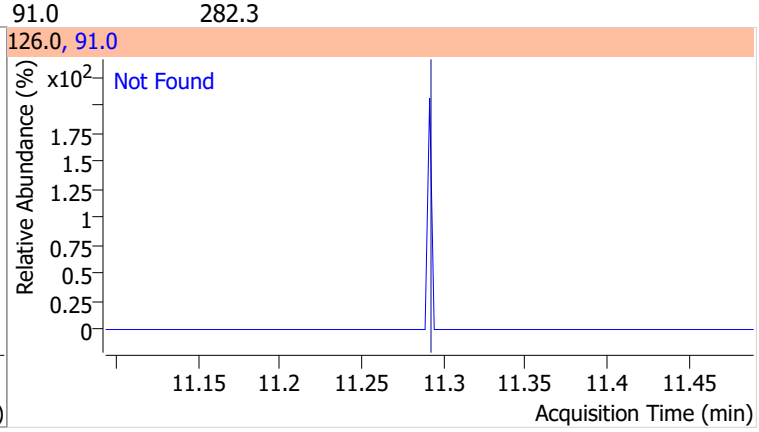
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



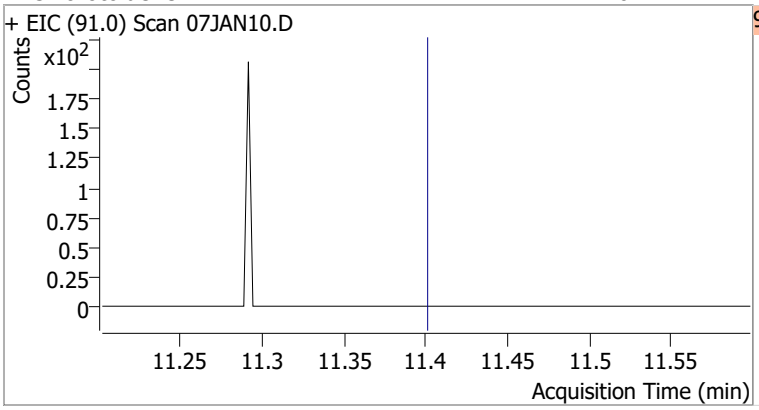


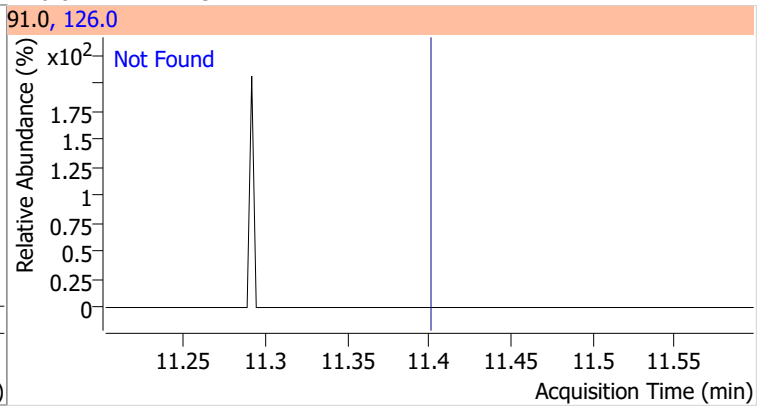
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



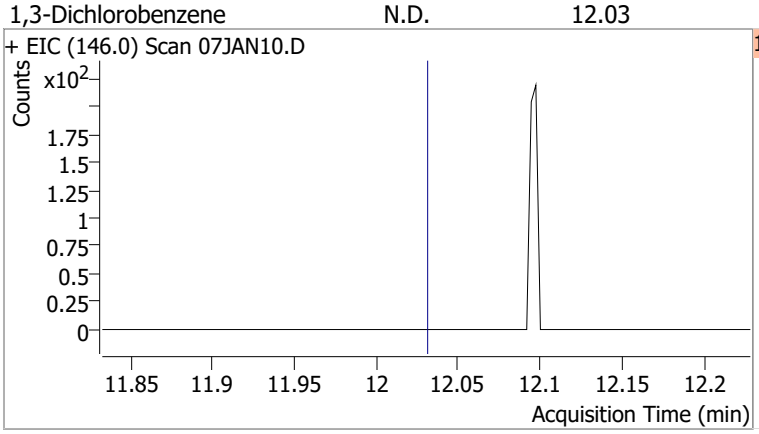


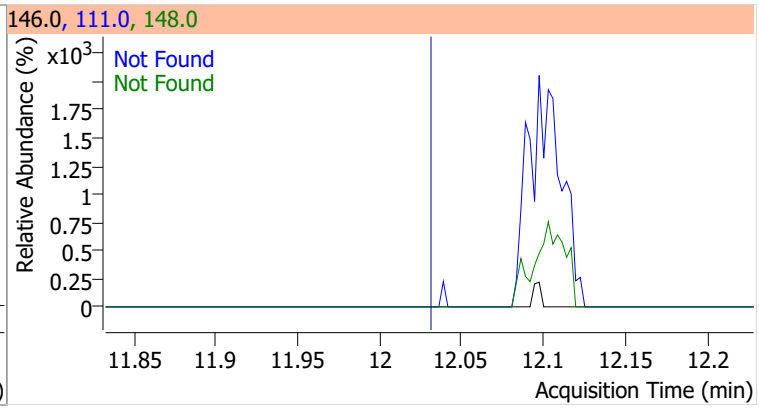
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



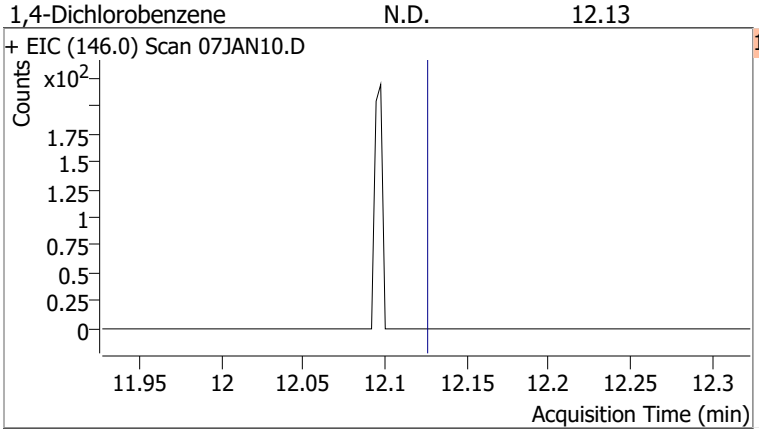


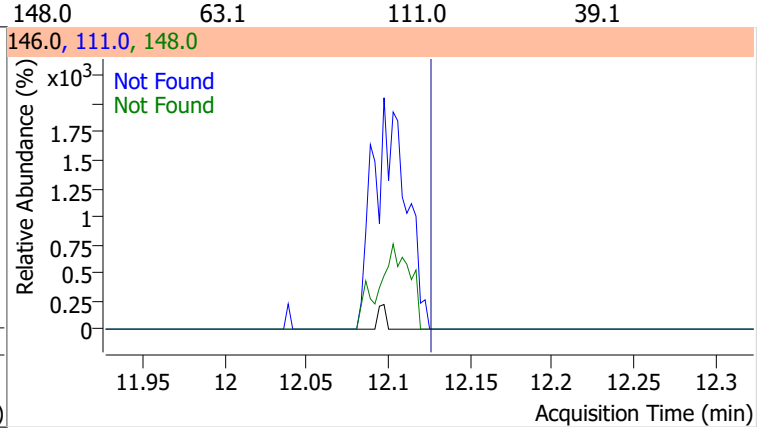
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



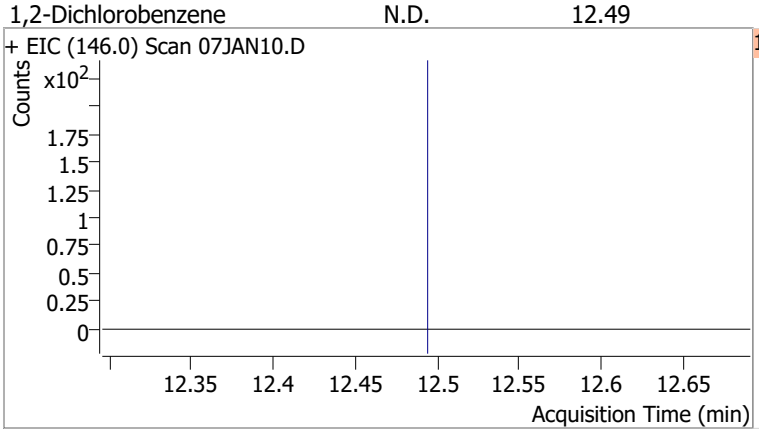


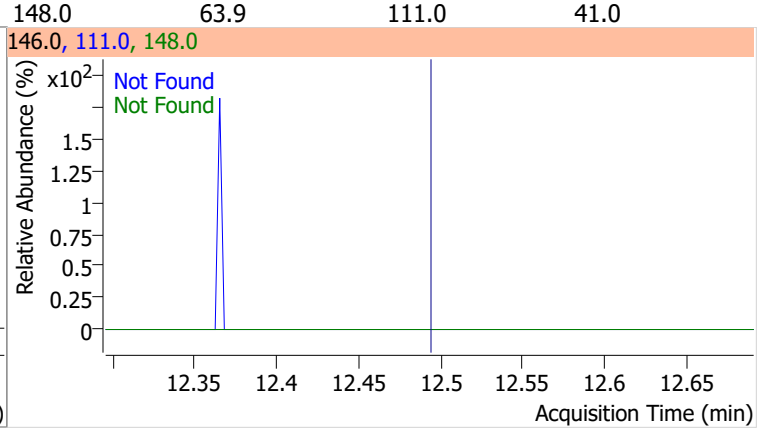
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





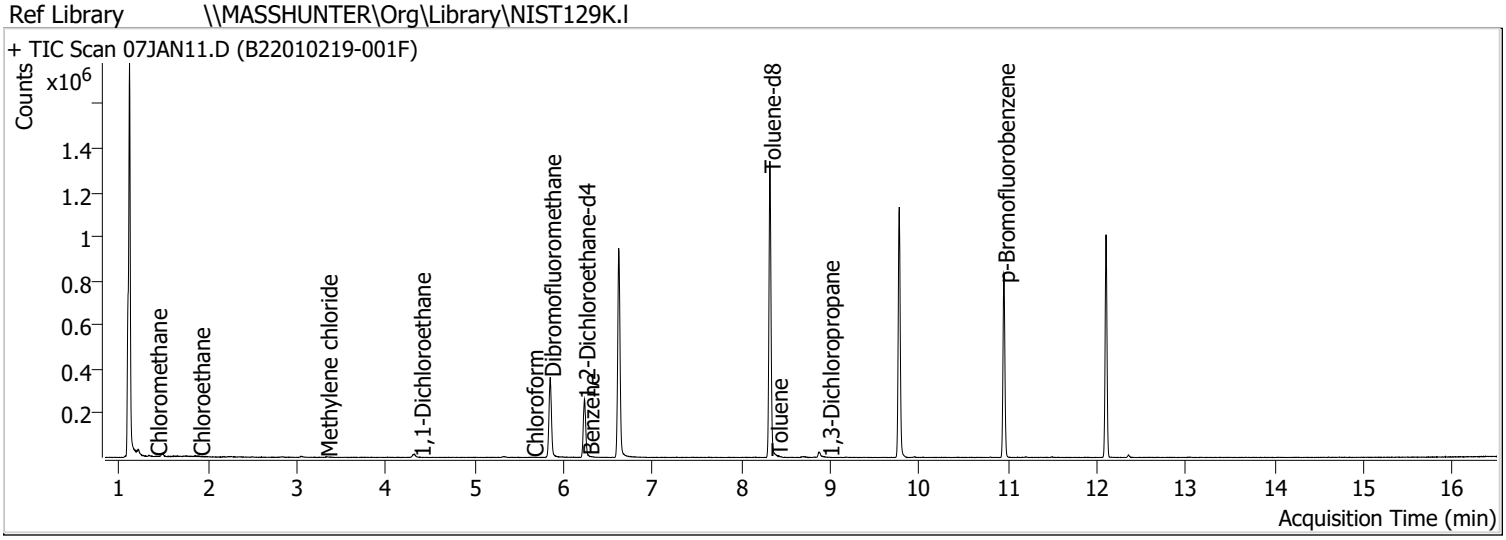
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN11.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 2:14:04 PM
Sample Name	B22010219-001F	Instrument	VOA5975C
Vial	11	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

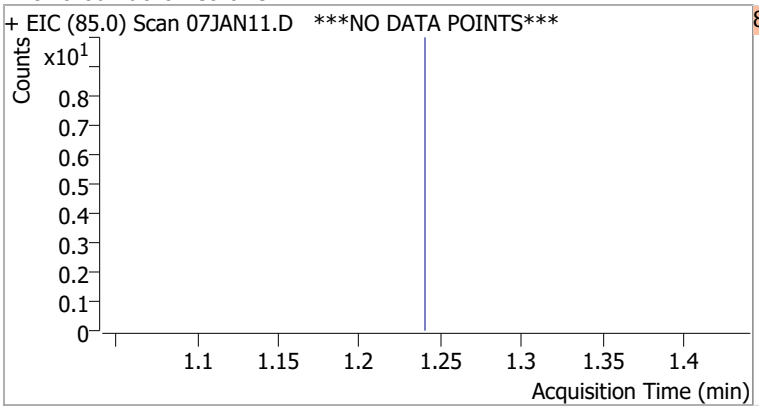


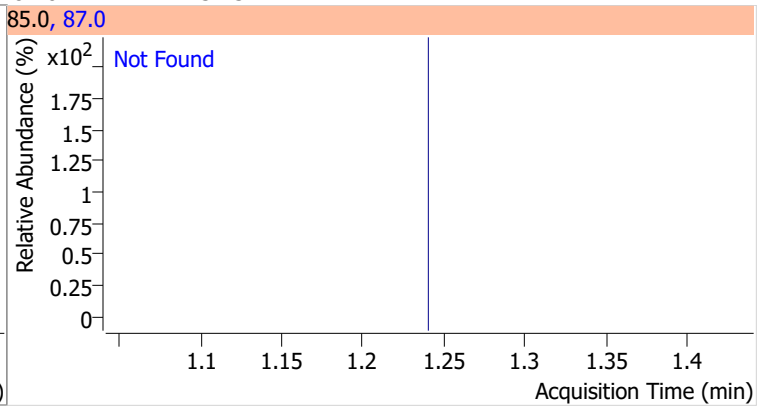
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.621	96.0	806641	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	312339	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	239572	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.848	113.0	211143	277.8425	ng		0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 111.14%			
S 1,2-Dichloroethane-d4	6.233	67.0	96836	295.0174	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 118.01%		*	
S Toluene-d8	8.319	98.0	811840	269.7270	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 107.89%			
S p-Bromofluorobenzene	10.951	95.0	237586	270.6993	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 108.28%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.417	50.0	1302	1.0151	ng	m	84
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	1.899	64.0	1604	2.8067	ng	m	94
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.333	49.0	1285	1.0727	ng	m	83
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	4.381	63.0	953	0.6175	ng	m	76
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	5.650	83.0	777	0.5057	ng	m	77

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.278	78.0	1404	0.4371	ng	m	75
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	0.000		0	N.D.			
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.397	92.0	699	0.3436	ng	m	84
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	8.812	83.0	0		ng	md	1
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	8.977	76.0	319	0.4005	ng	m	81
T Chlorodibromomethane	0.000		0	N.D.			
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	0.000		0	N.D.			
T m+p-Xylenes	10.039	106.0	0		ng	md	1
T o-Xylene	10.430	106.0	0		ng	md	1
T Styrene	0.000		0	N.D.			
T Bromoform	0.000		0	N.D.			
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

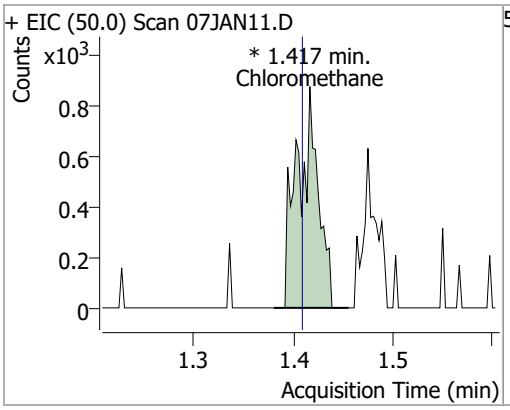
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

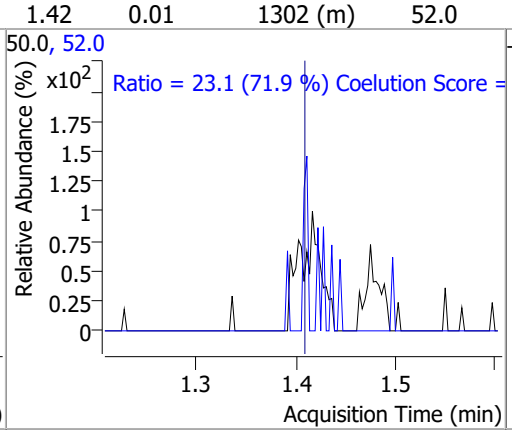
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

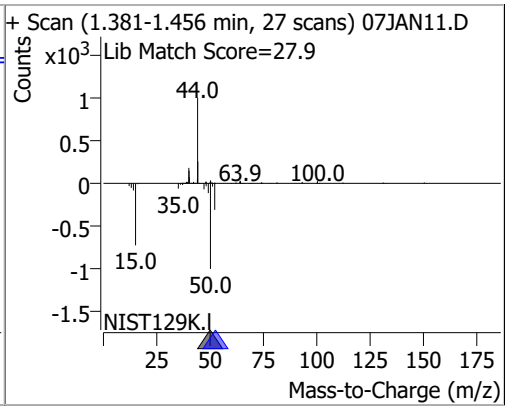




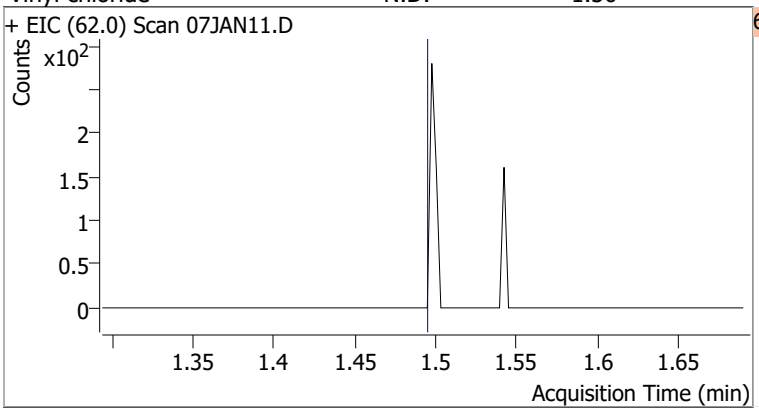
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	1.0151	1.42	0.01	1302 (m)	52.0	23.1	2.1	62.1

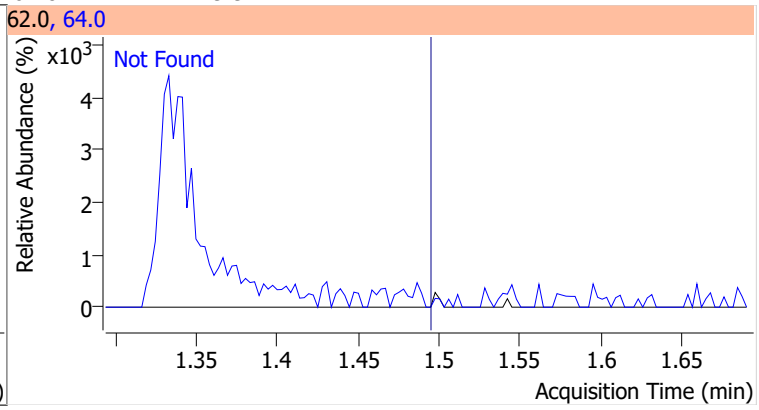




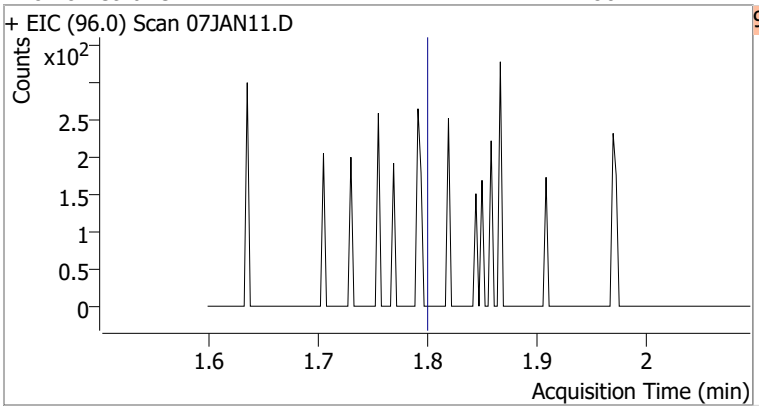


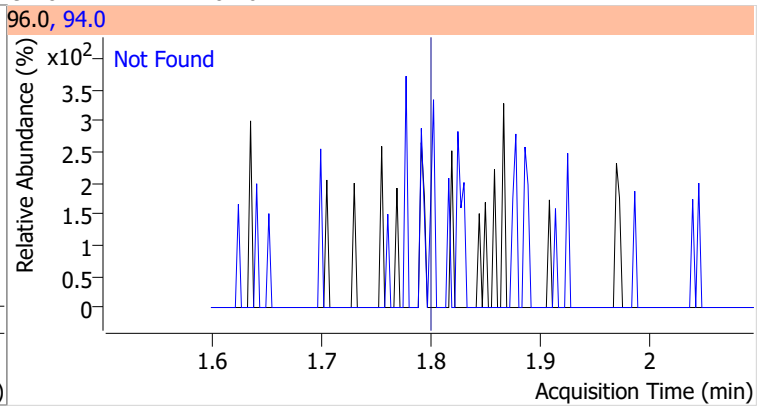
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



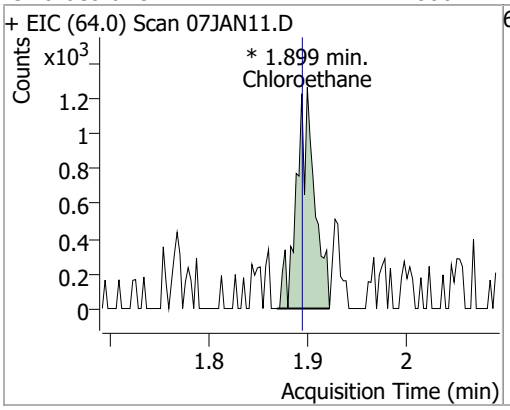


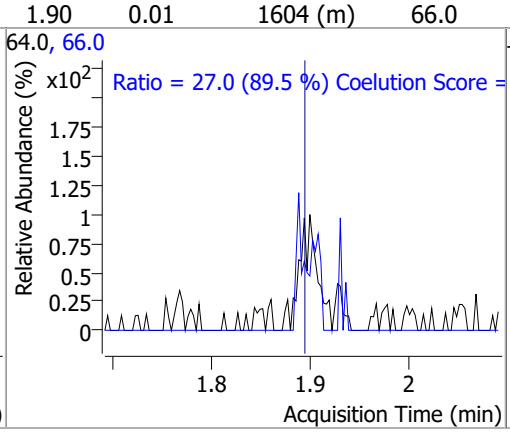
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6

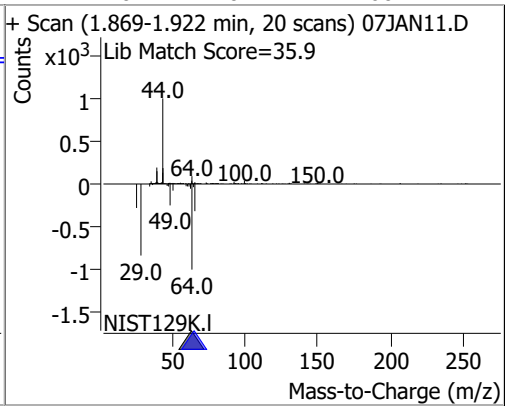




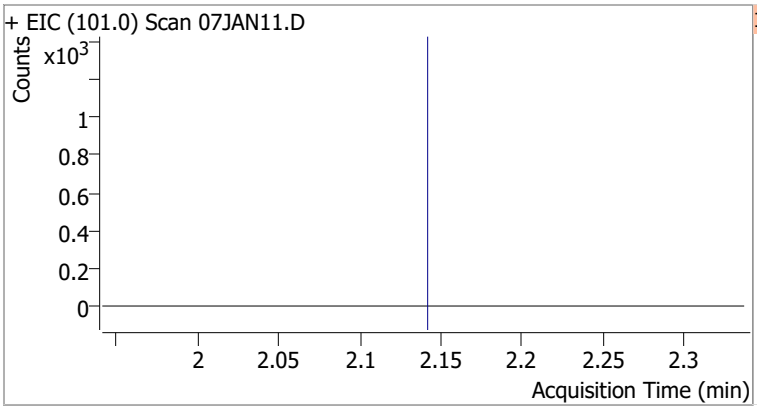
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	2.8067	1.90	0.01	1604 (m)	66.0	27.0	0.1	60.1

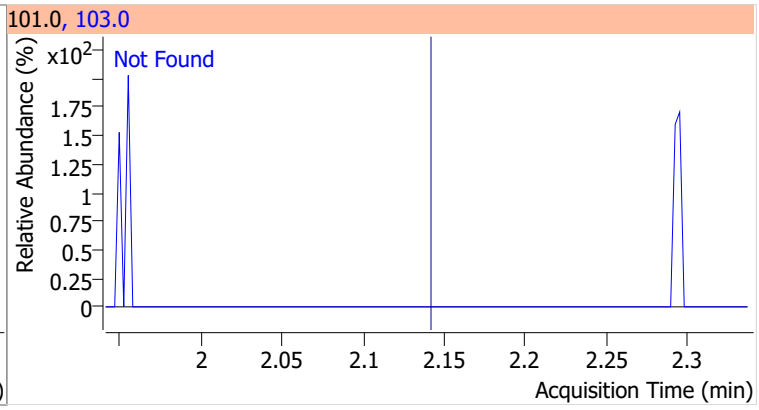




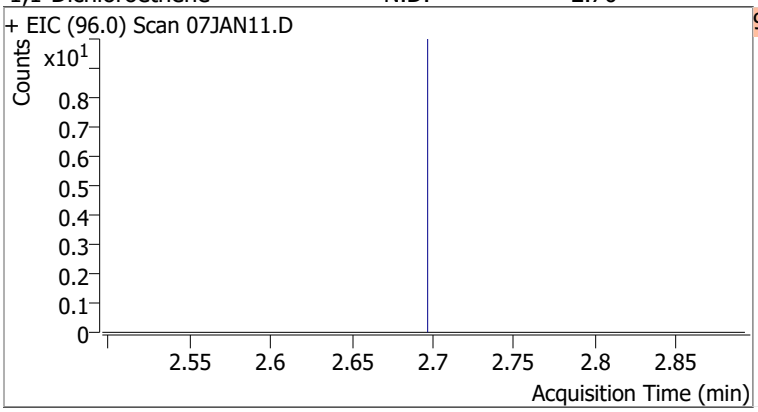


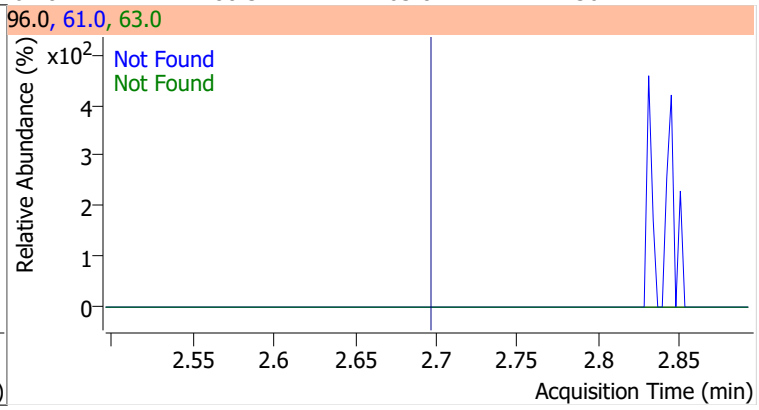
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



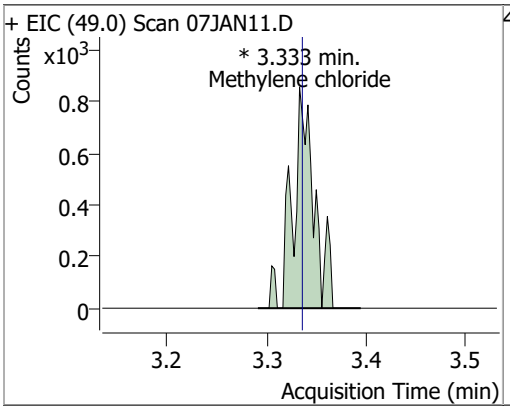


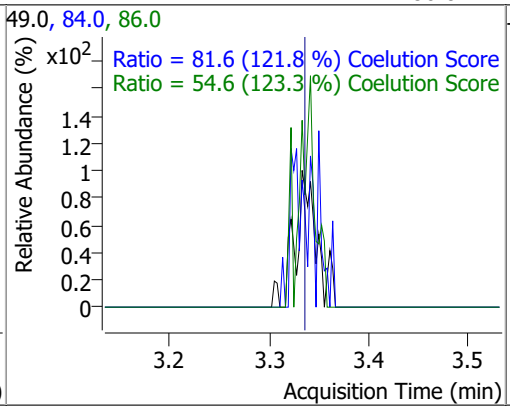
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

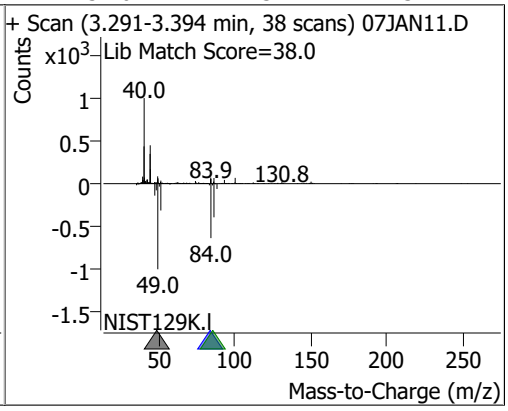




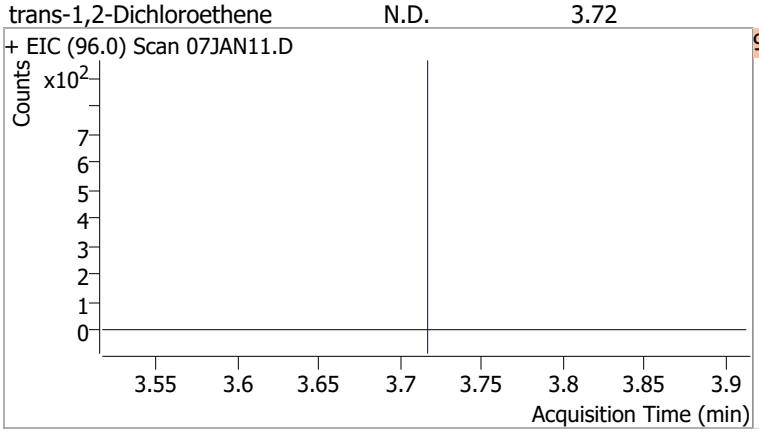
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	1.0727	3.33	0.00	1285 (m)	84.0	81.6	36.9	96.9
					86.0	54.6	14.3	74.3

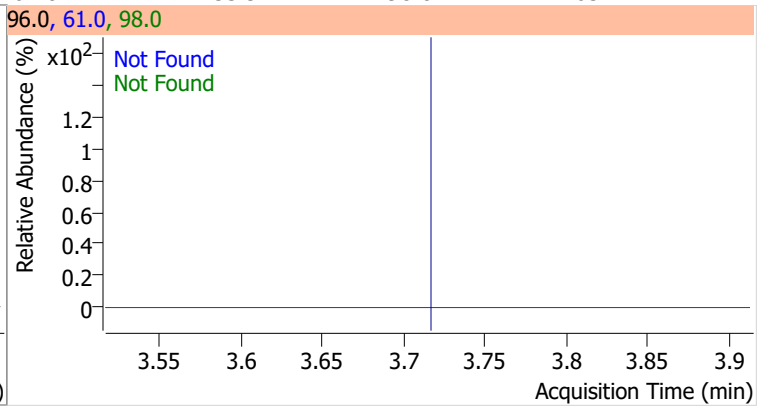




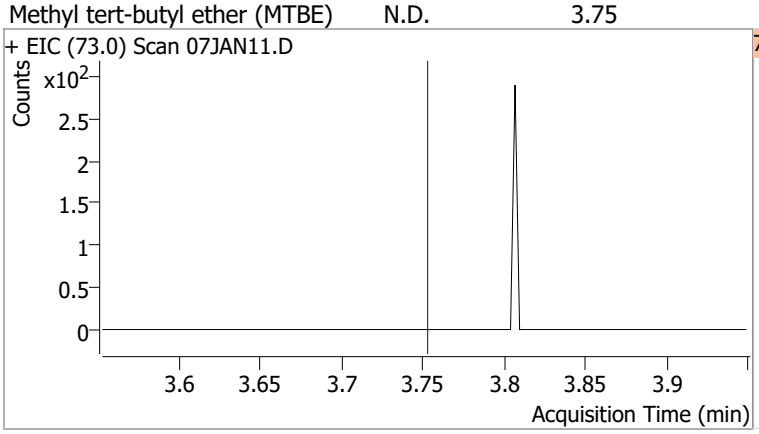


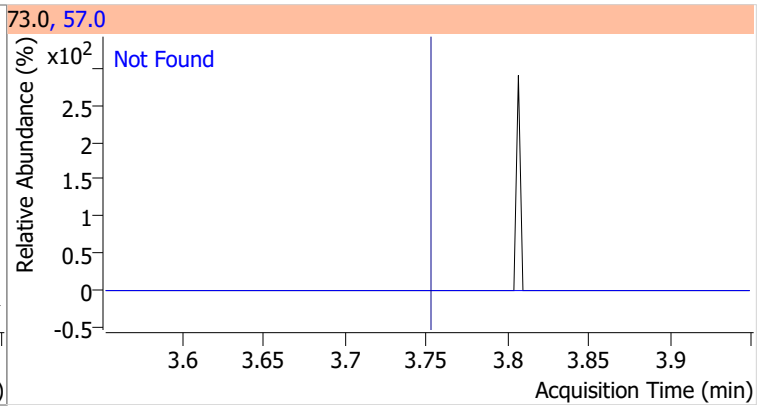
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



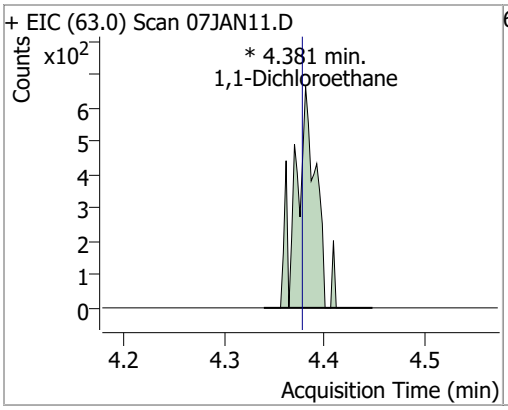


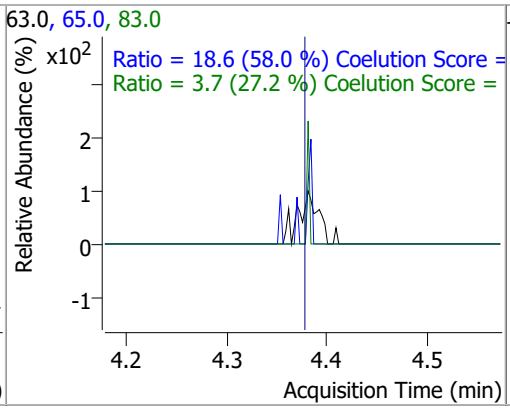
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6

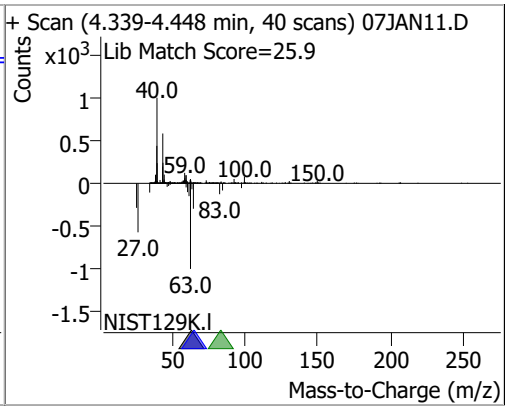




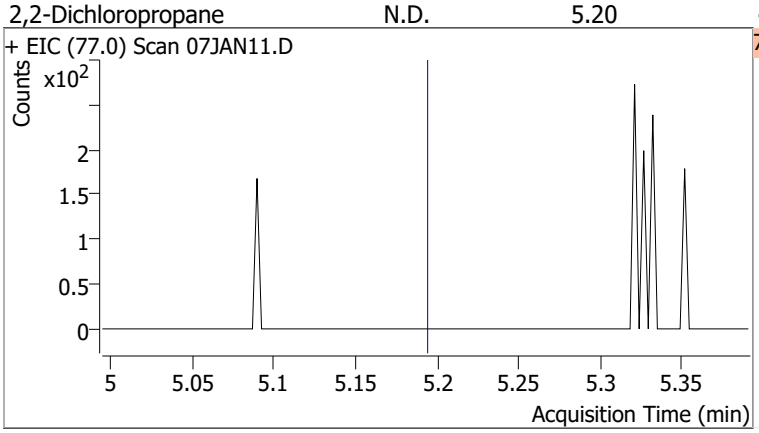
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	0.6175	4.38	0.00	953 (m)	65.0	18.6	2.1	62.1
					83.0	3.7	0.0	43.7

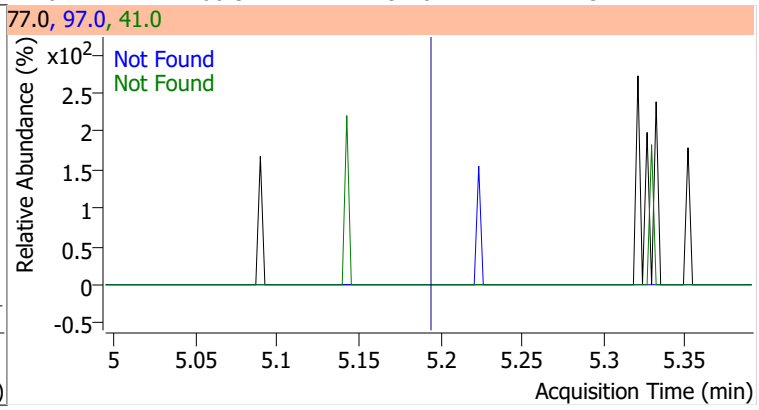




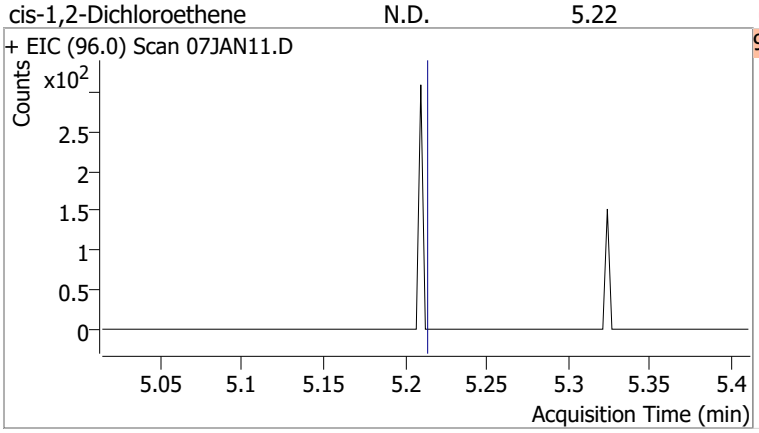


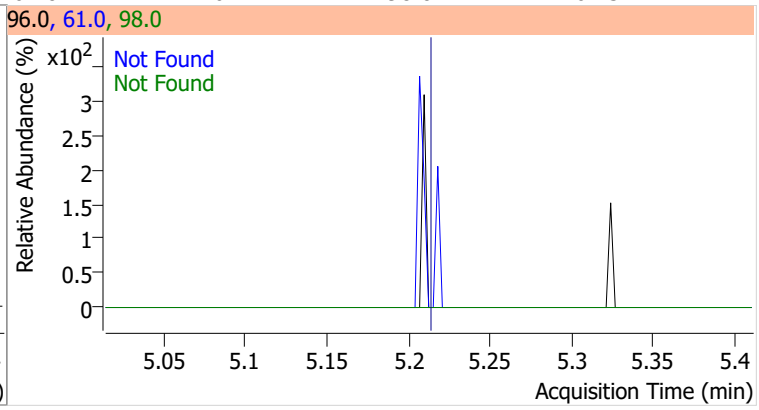
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



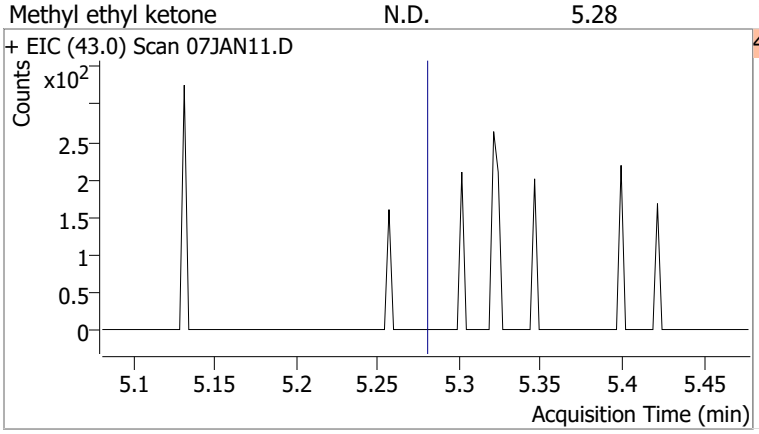


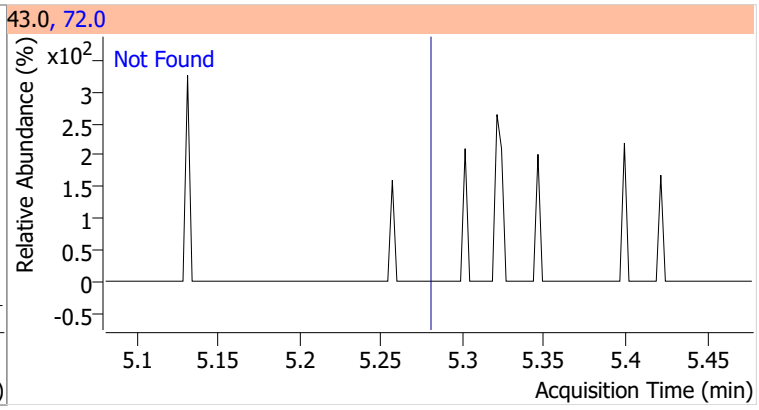
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



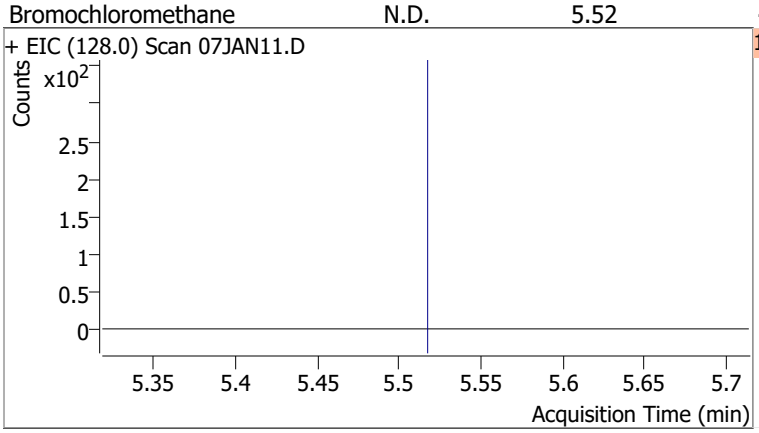


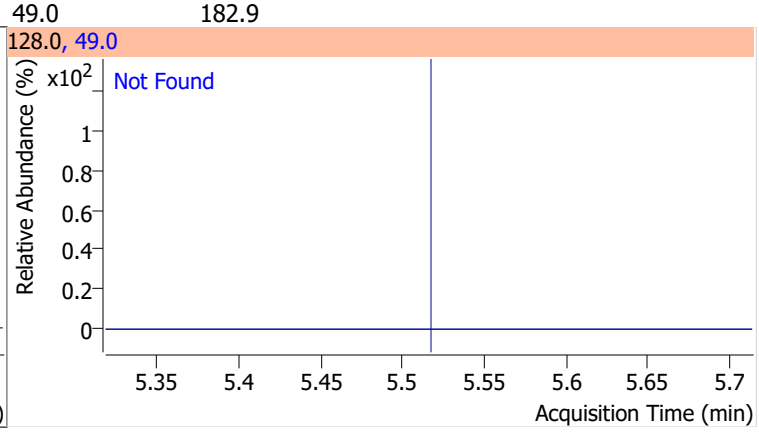
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



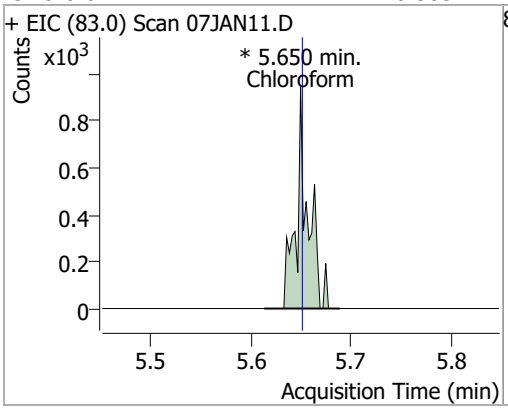


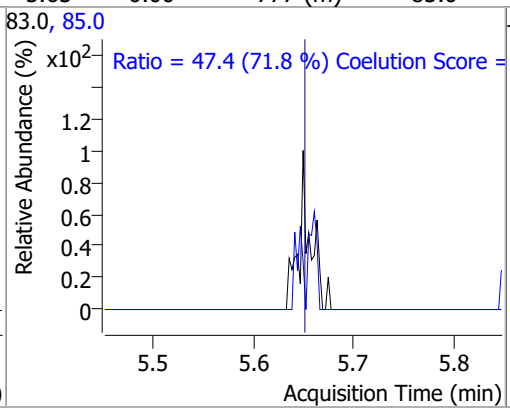
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9

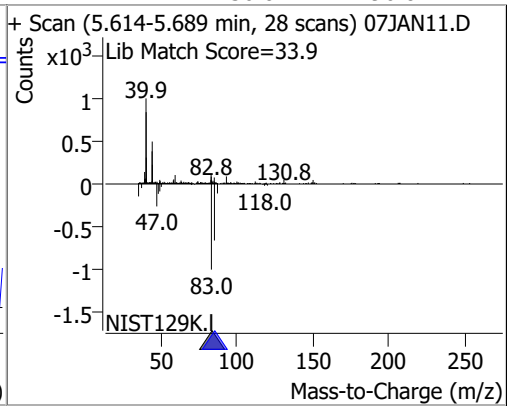




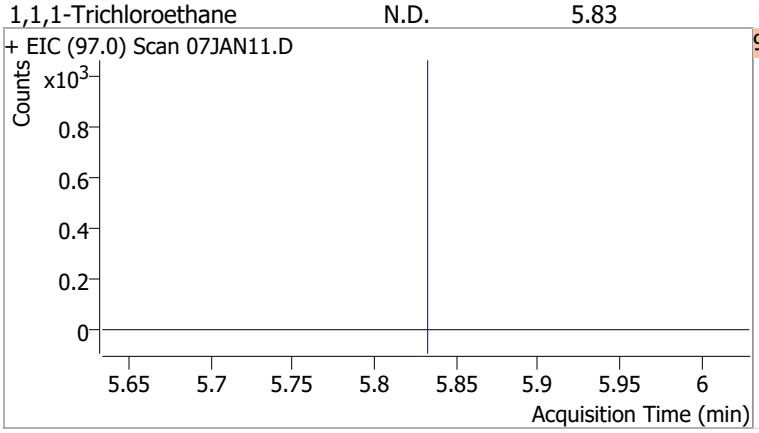
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	0.5057	5.65	0.00	777 (m)	85.0	47.4	36.0	96.0

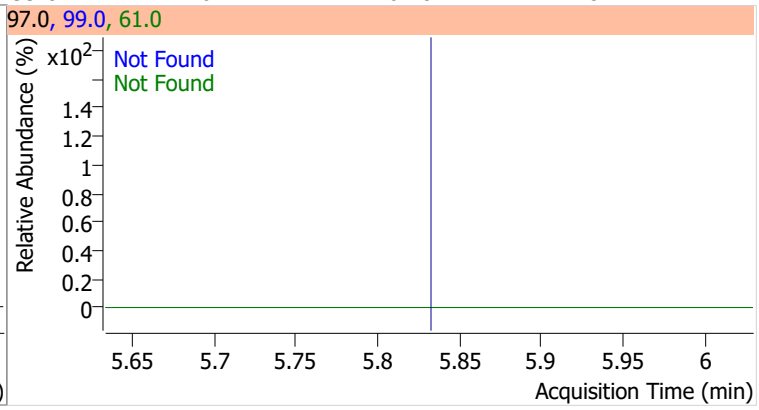




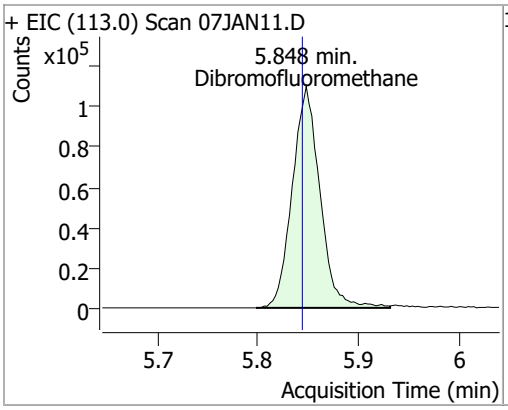


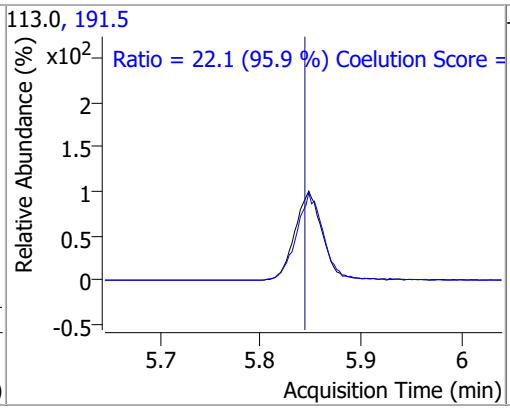
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

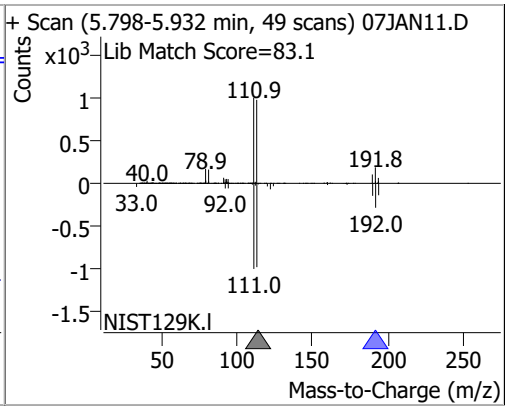




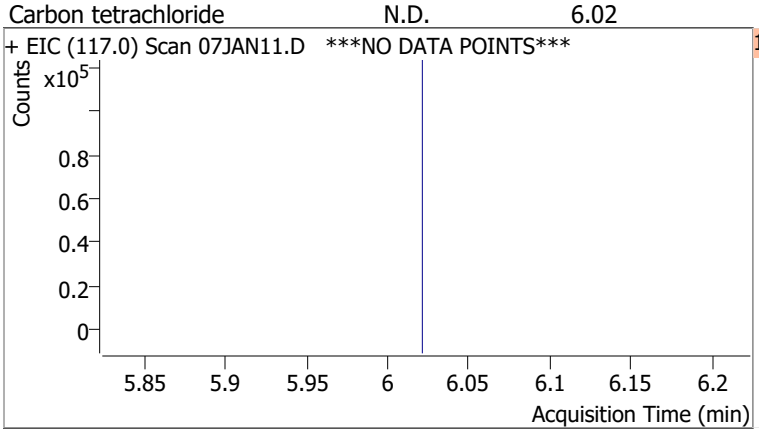
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	277.8425	5.85	0.00	211143	191.5	22.1	0.0	53.1

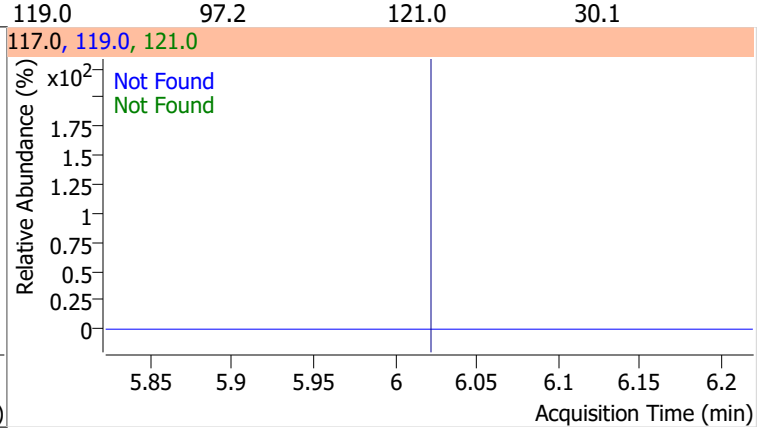




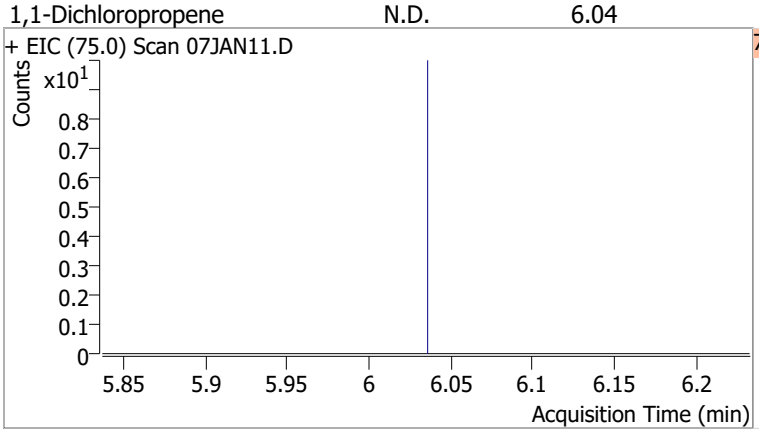


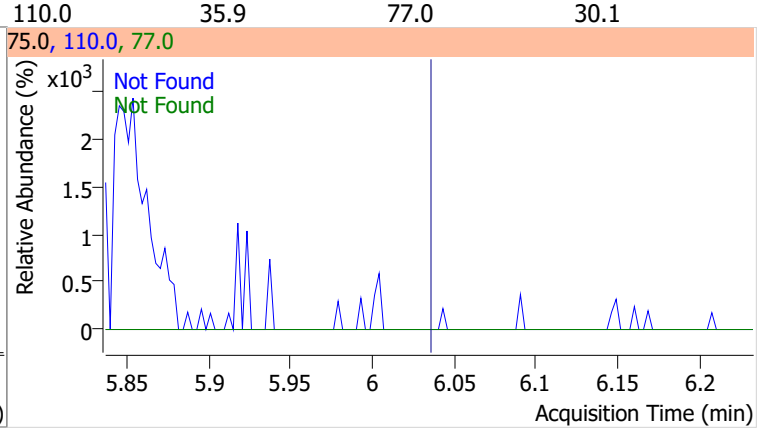
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



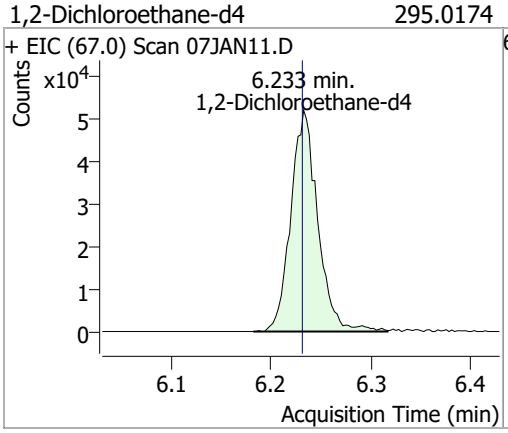


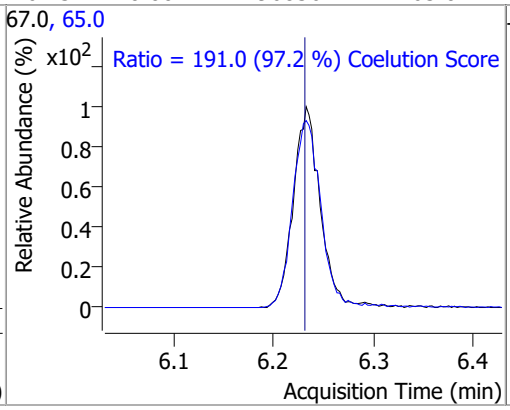
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

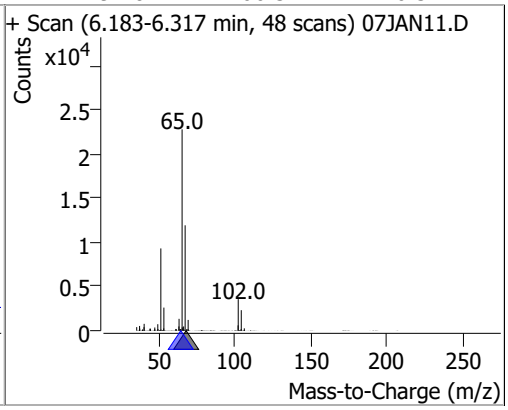




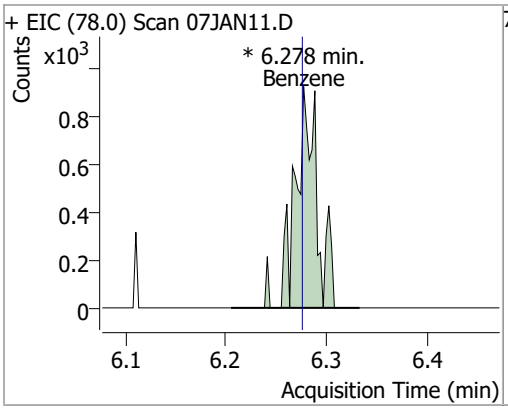
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	295.0174	6.23	0.00	96836	65.0	191.0	166.5	226.5

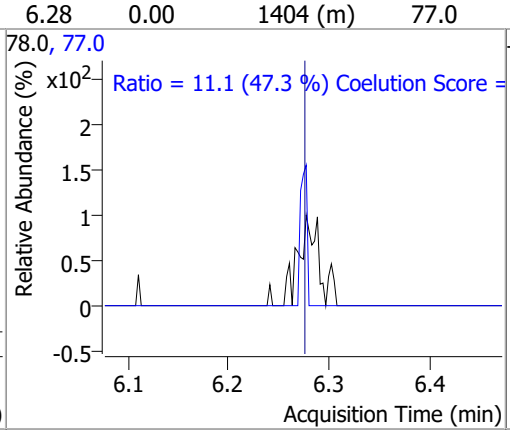


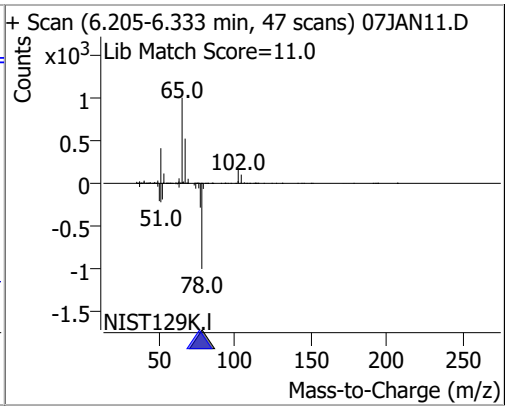




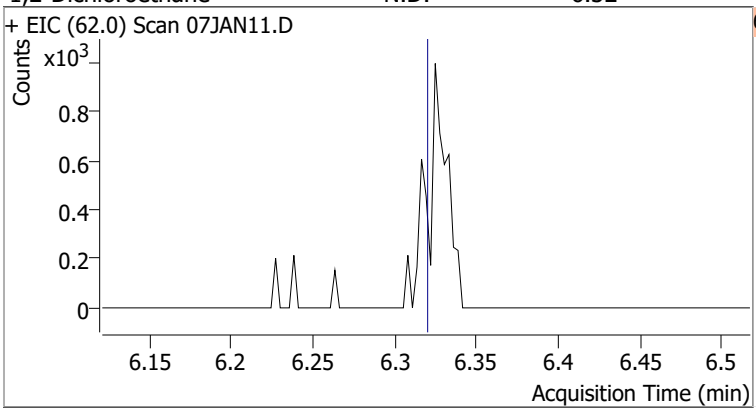
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.4371	6.28	0.00	1404 (m)	77.0	11.1	0.0	53.5

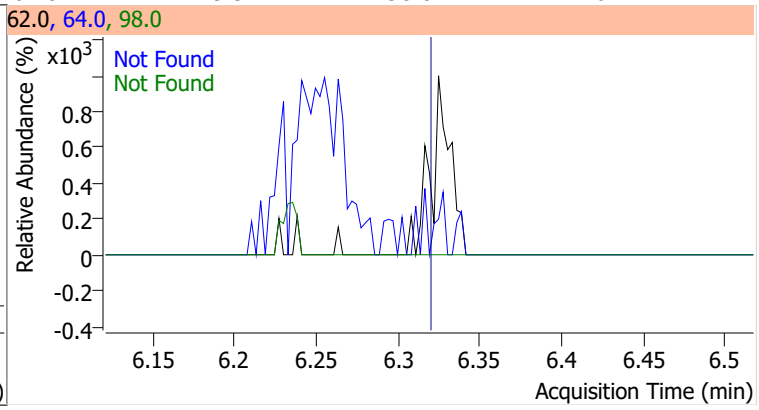




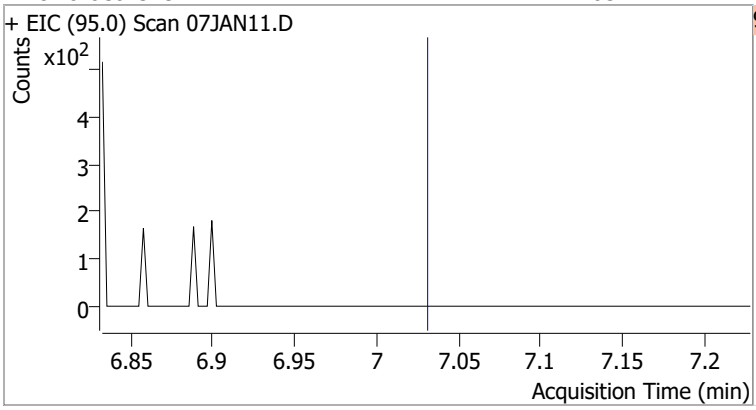


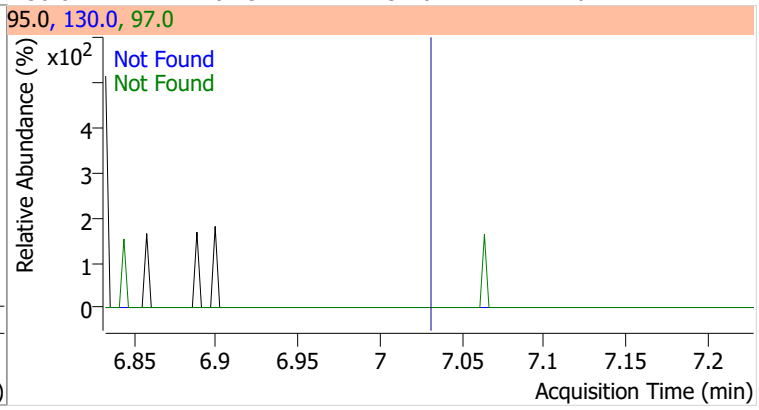
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



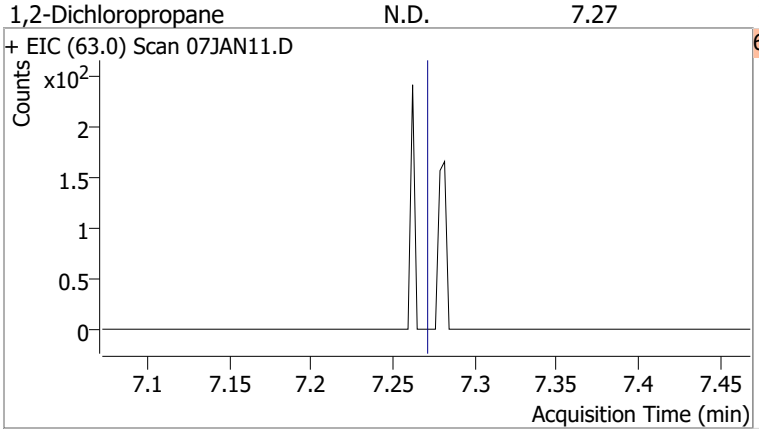


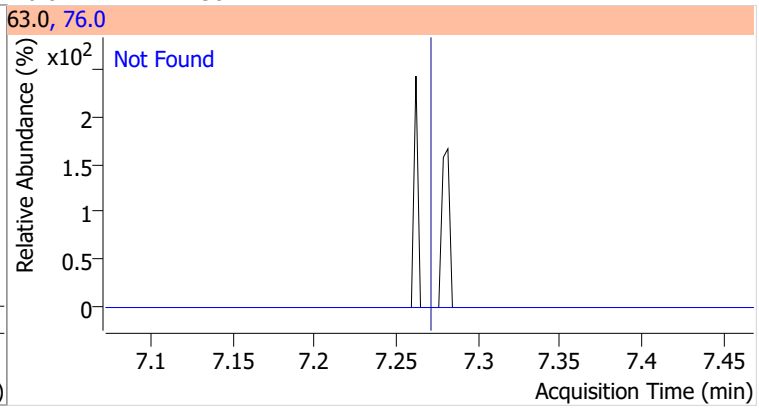
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



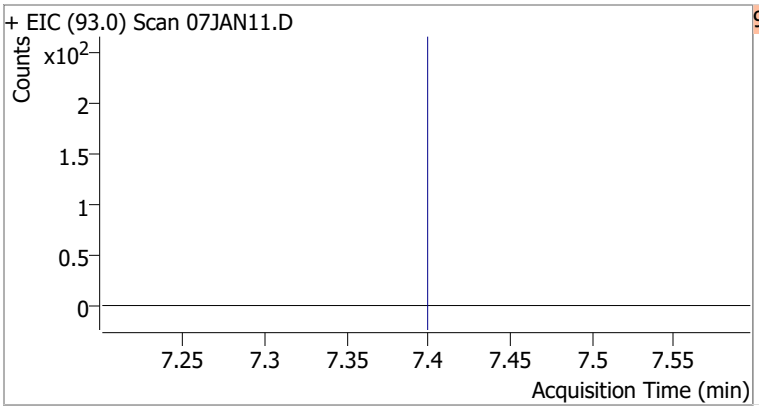


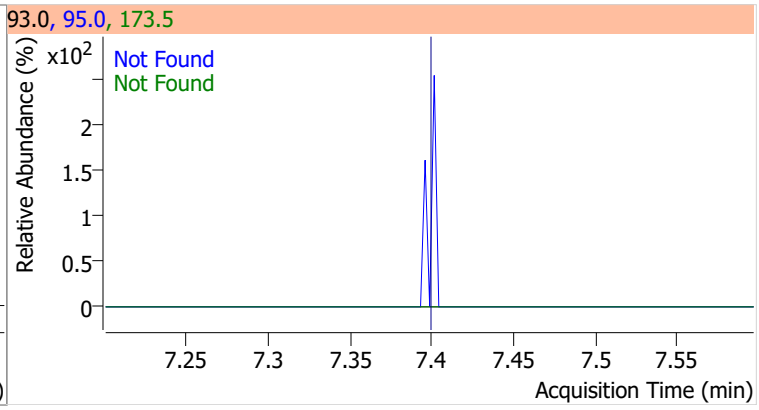
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



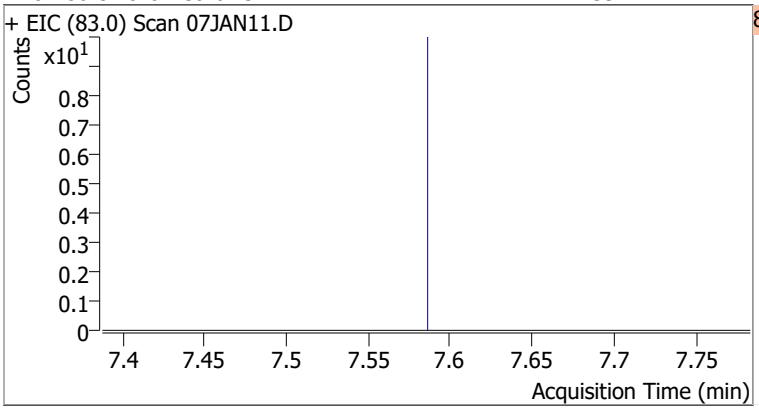


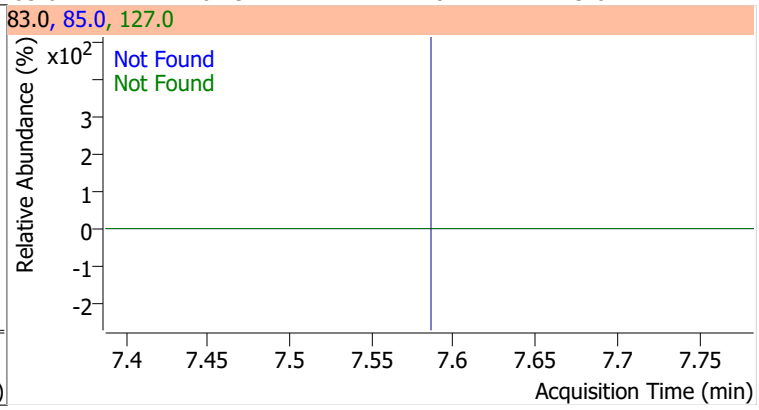
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



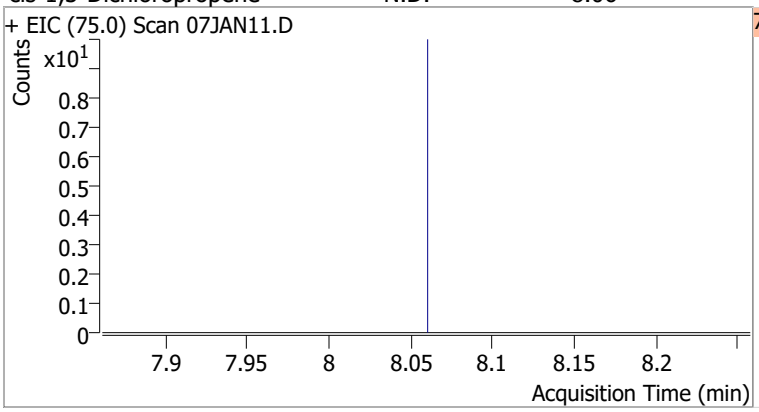


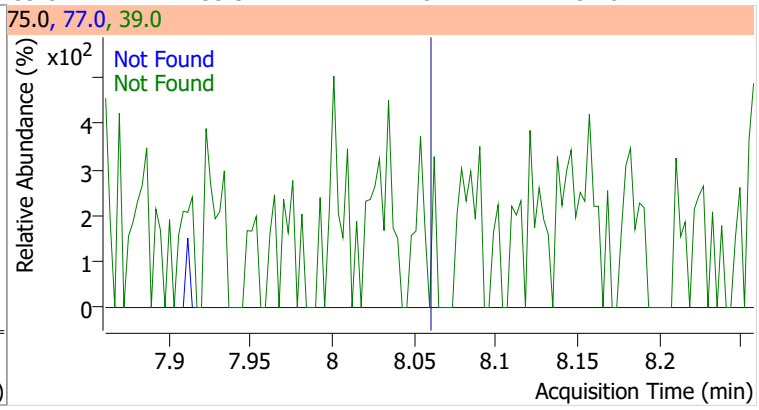
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



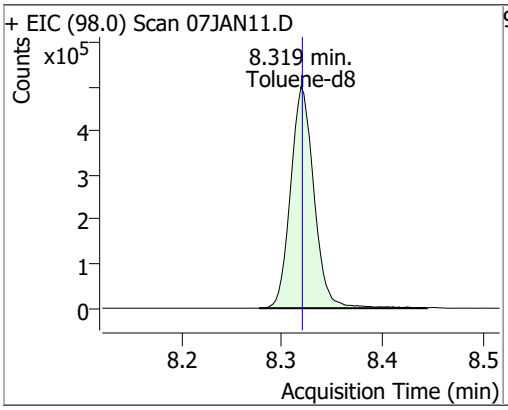


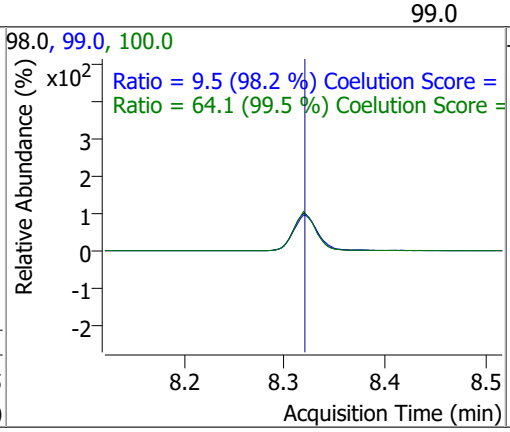
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

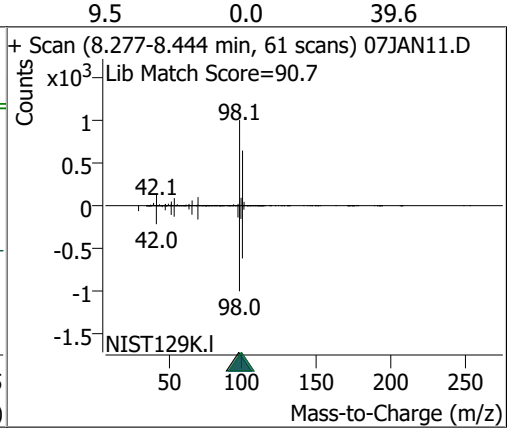




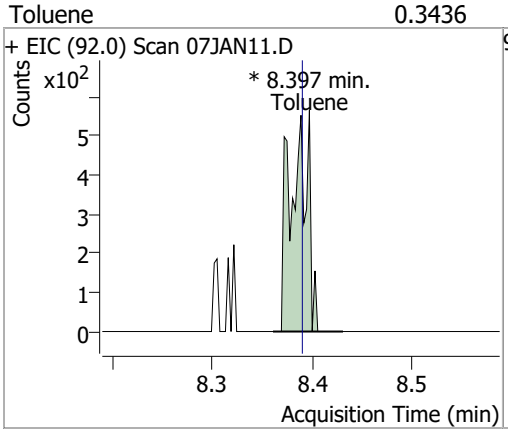
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	269.7270	8.32	0.00	811840	100.0	64.1	34.4	94.4
					99.0	9.5	0.0	39.6

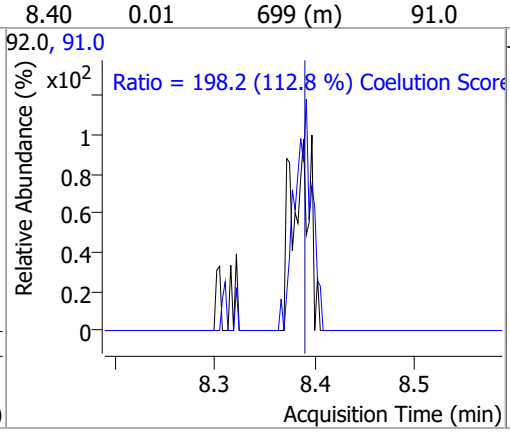


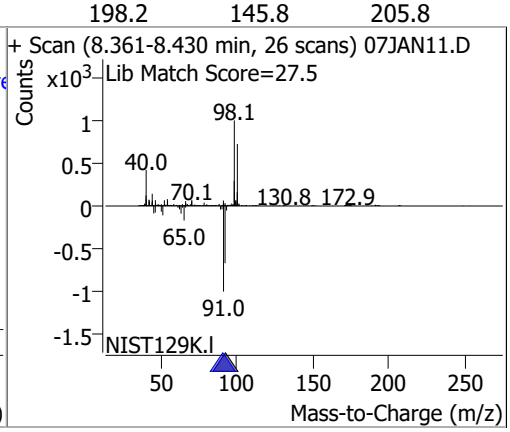




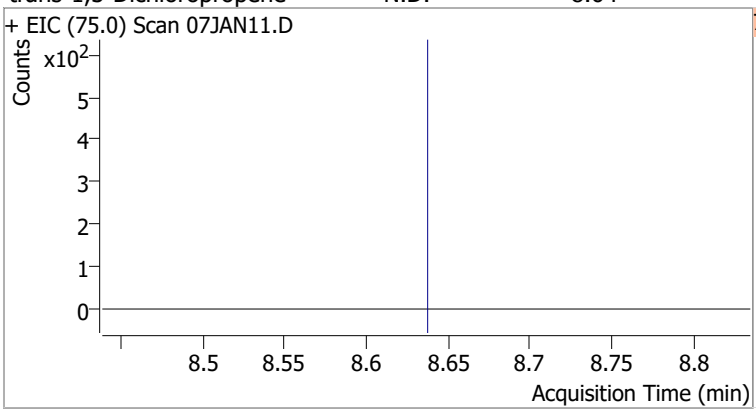
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	0.3436	8.40	0.01	699 (m)	91.0	198.2	145.8	205.8

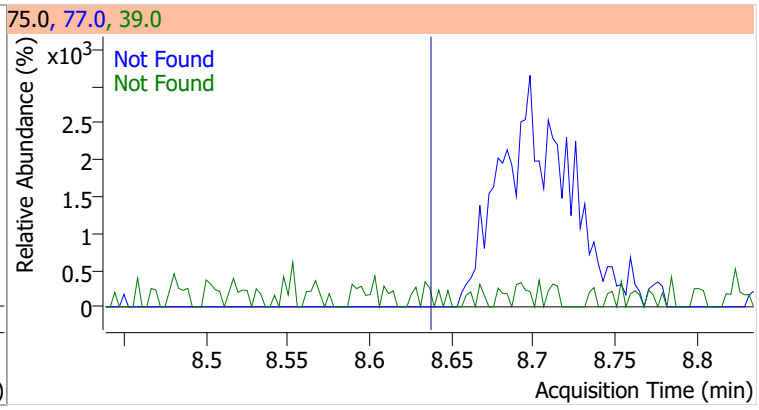




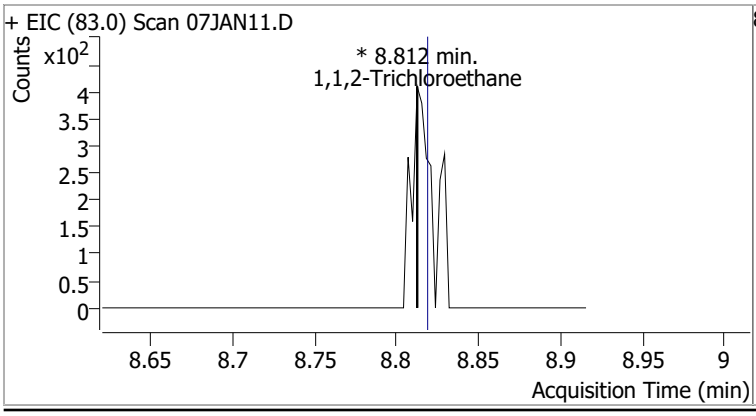


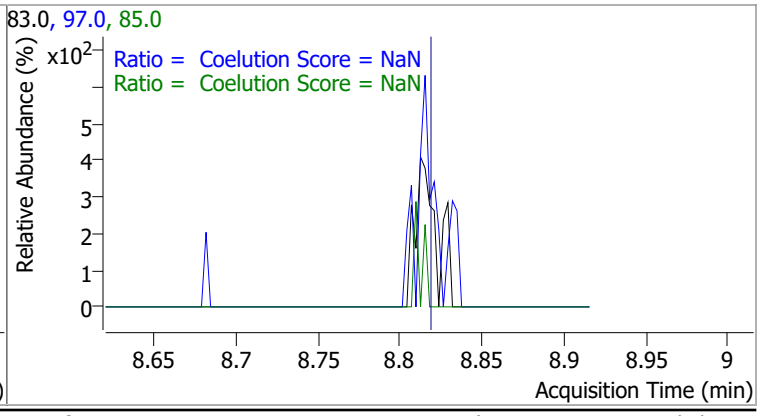
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



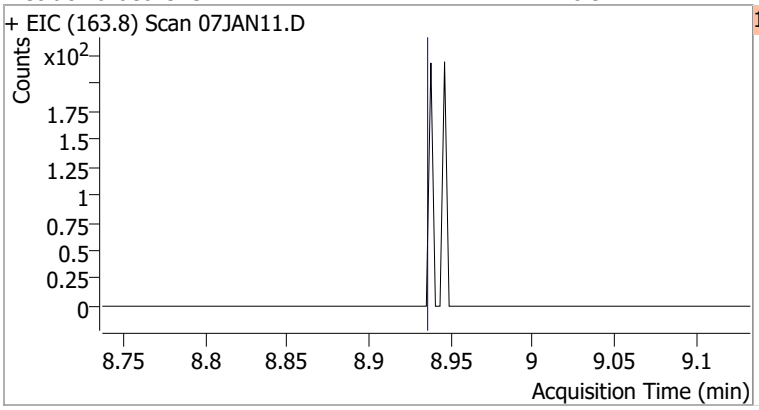


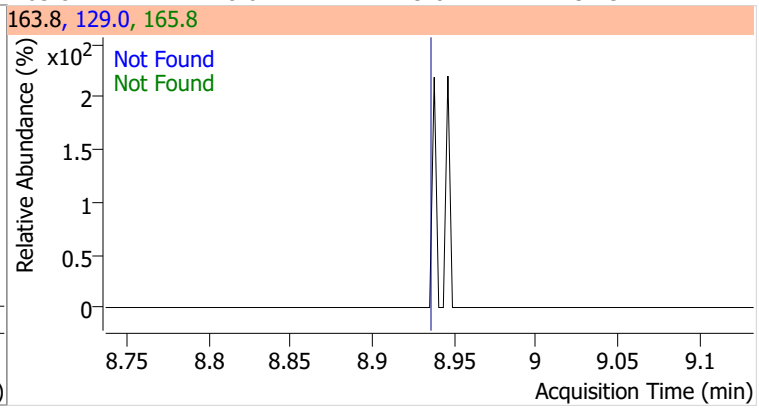
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane		0		0	97.0		84.6	144.6
					85.0		37.6	97.6



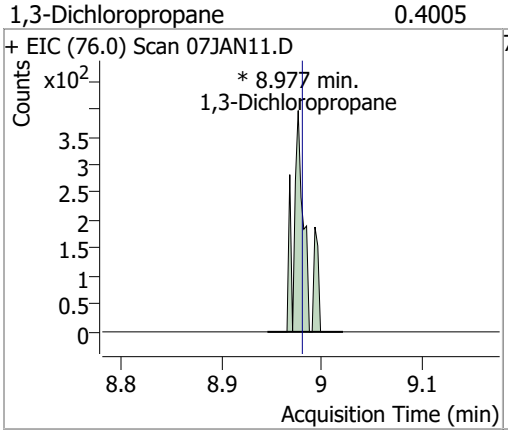


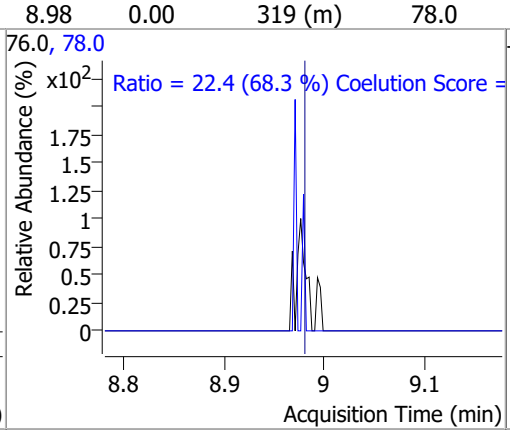
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5

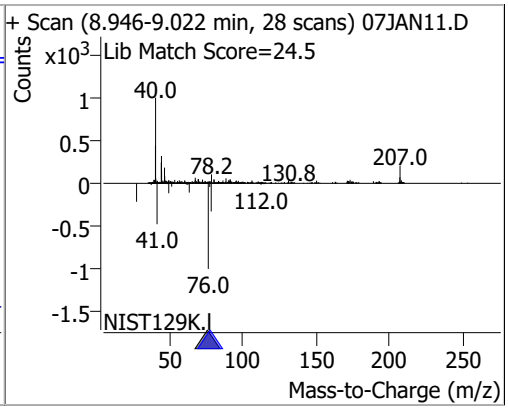




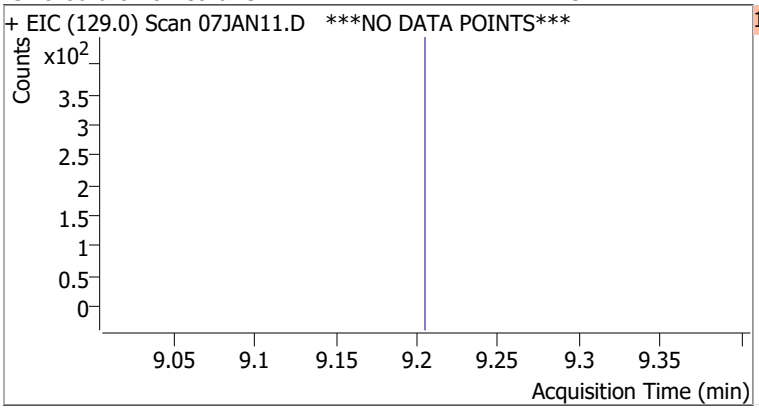
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	0.4005	8.98	0.00	319 (m)	78.0	22.4	2.9	62.9

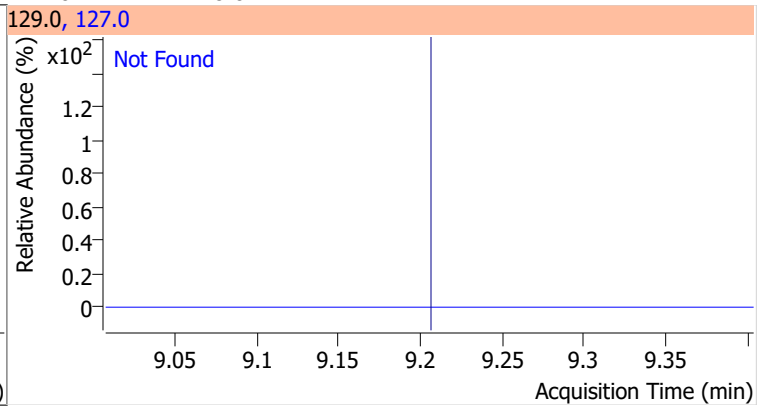




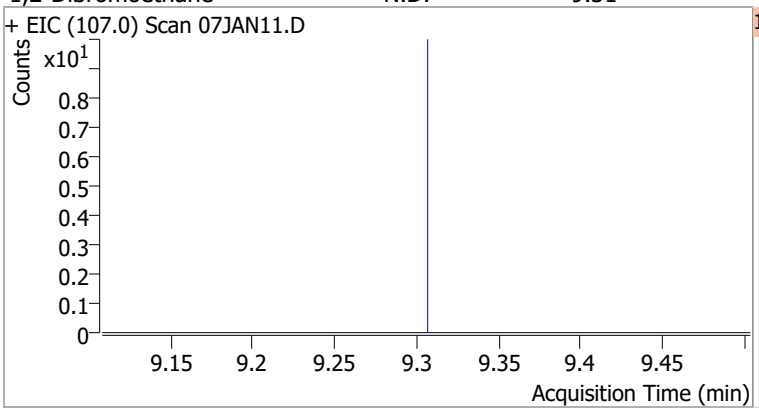


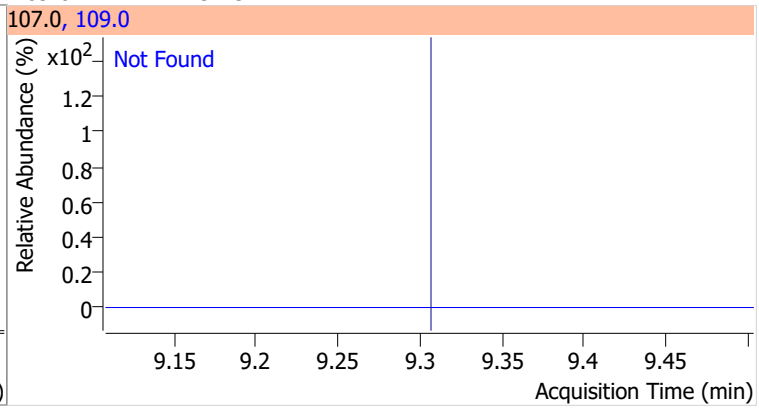
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



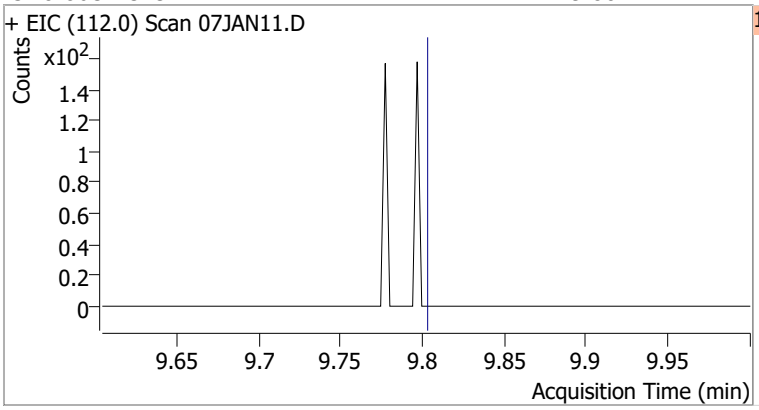


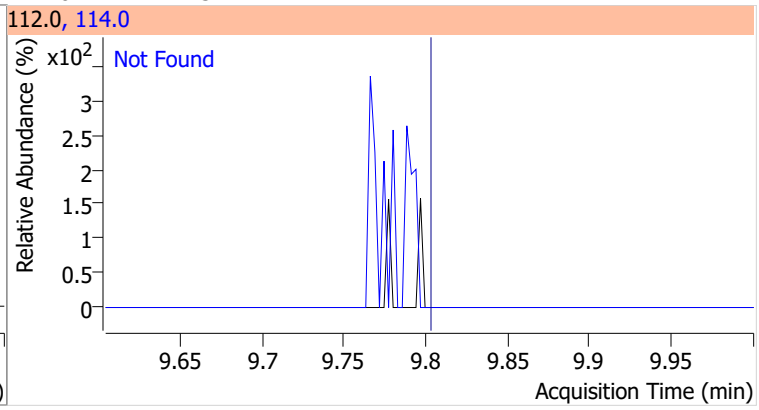
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



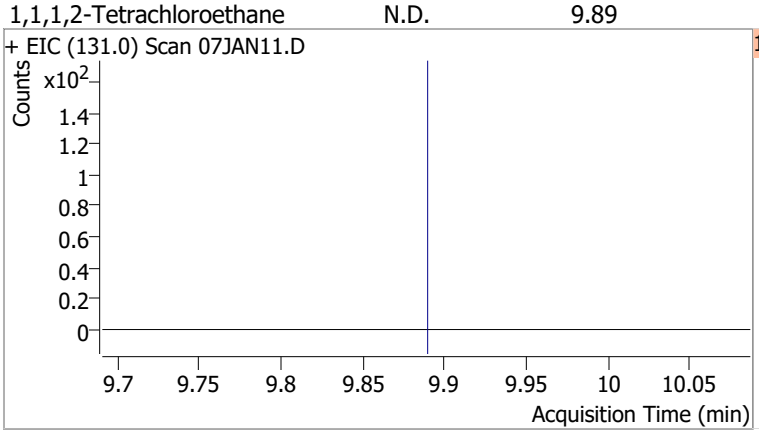


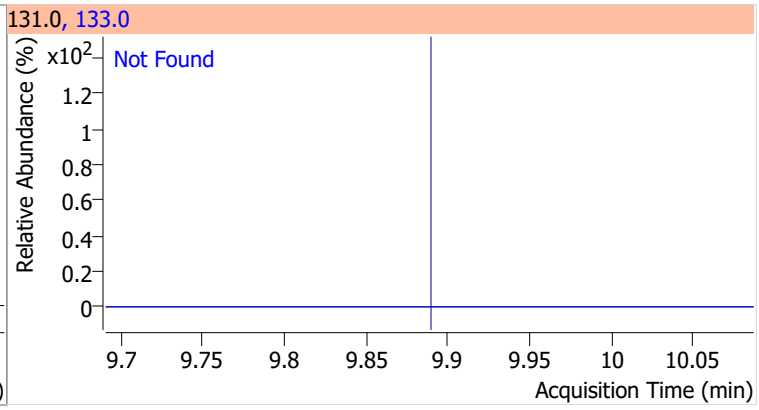
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



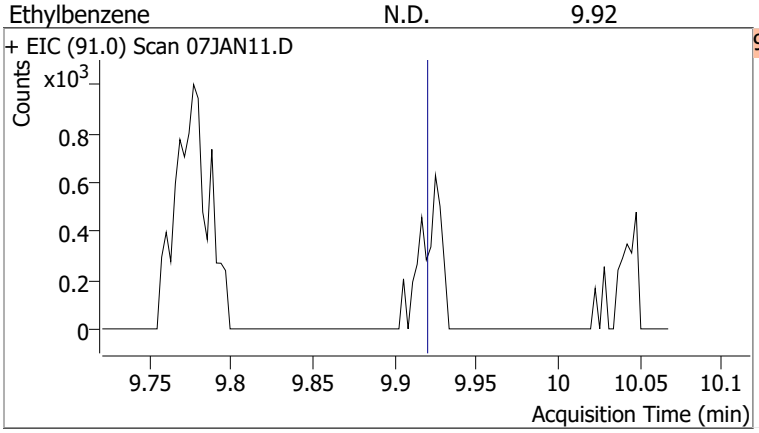


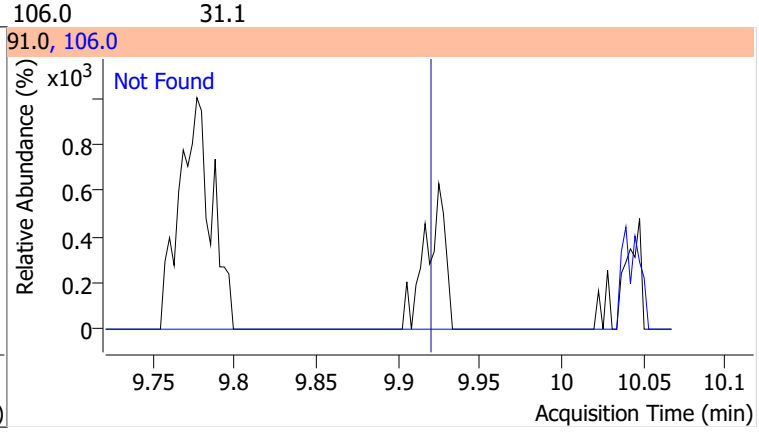
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



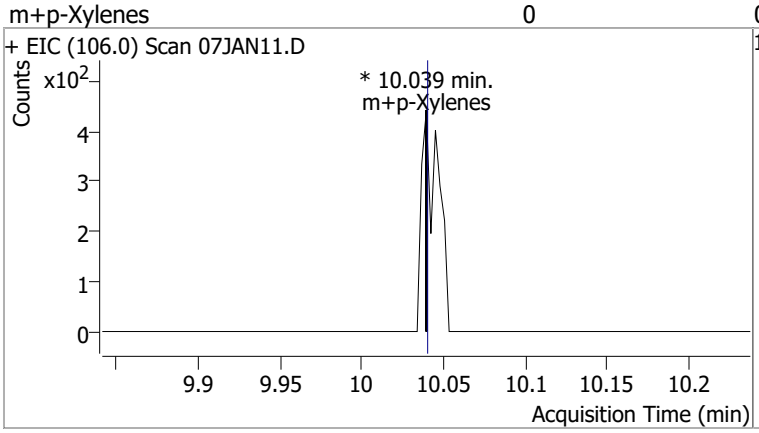


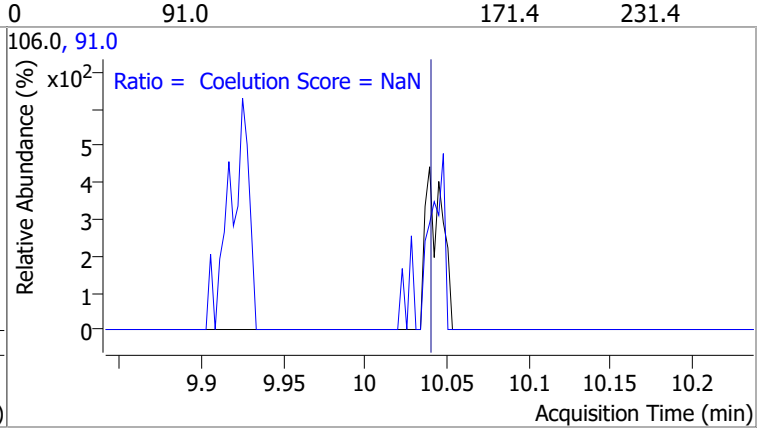
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1



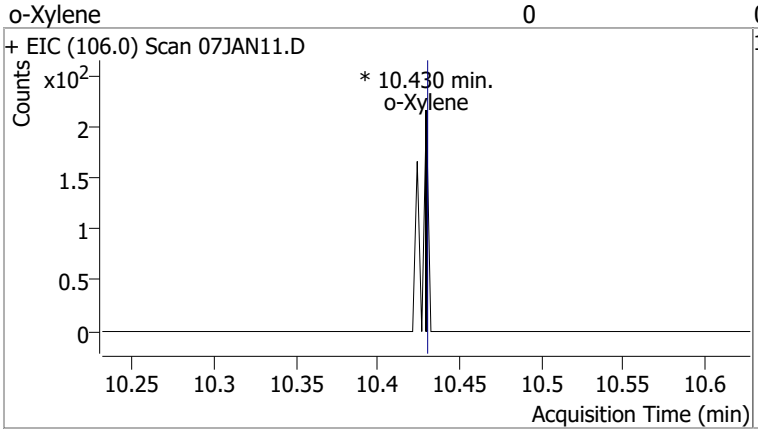


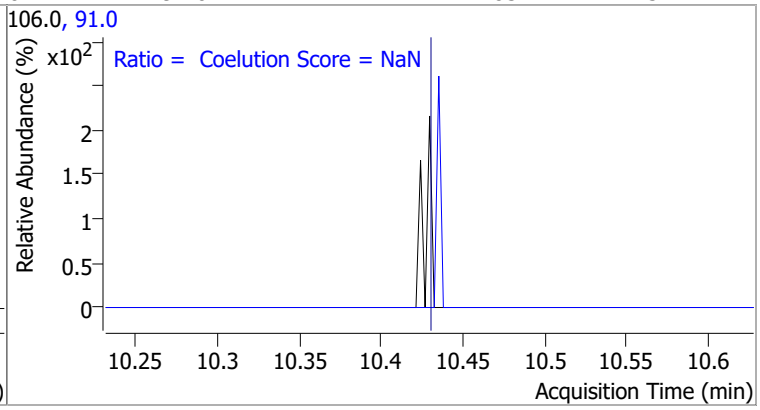
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes		0		0	91.0		171.4	231.4



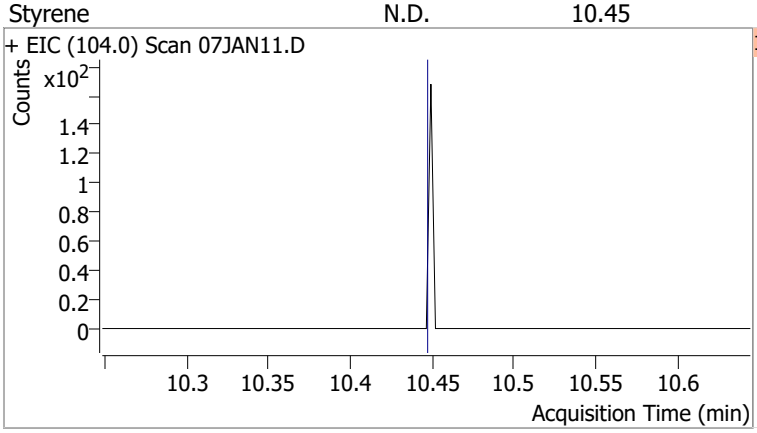


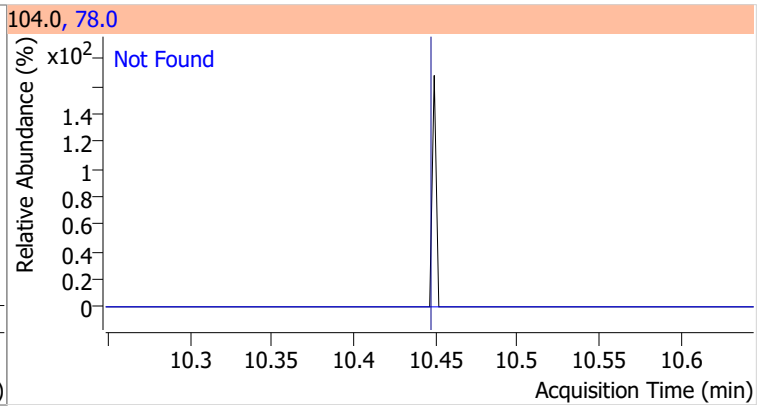
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene		0		0	91.0		183.1	243.1



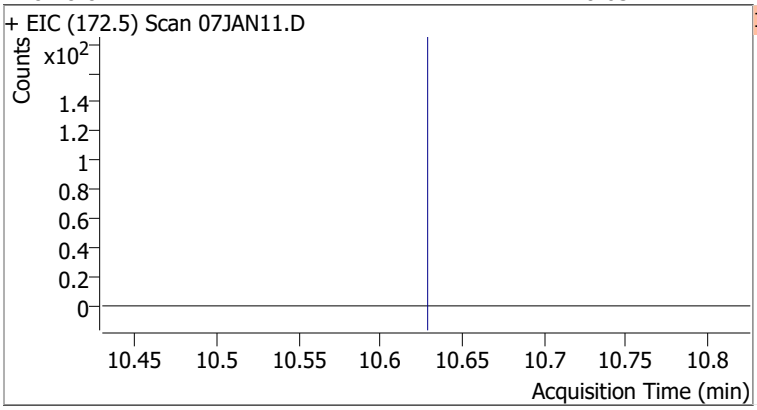


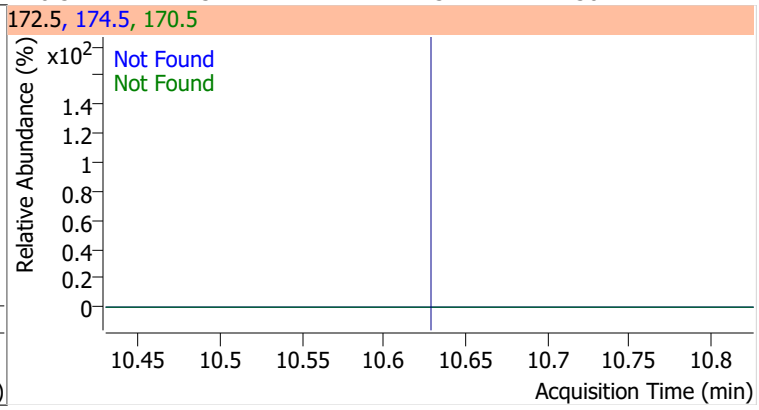
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



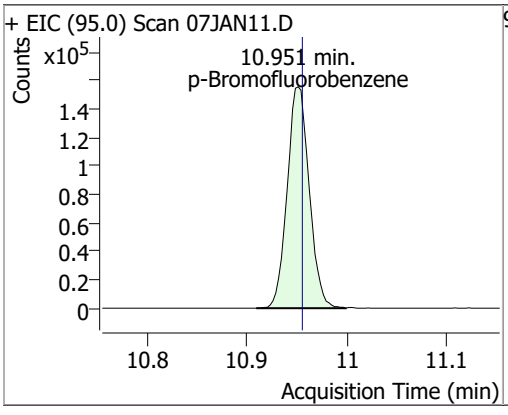


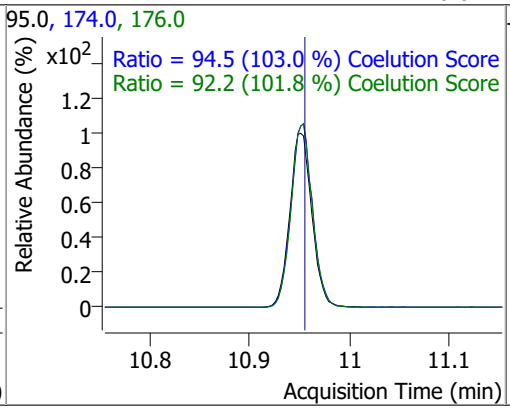
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

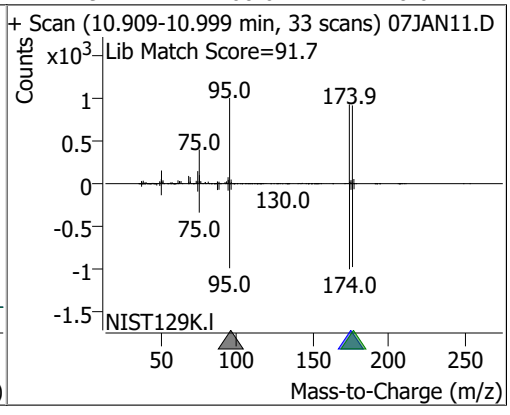




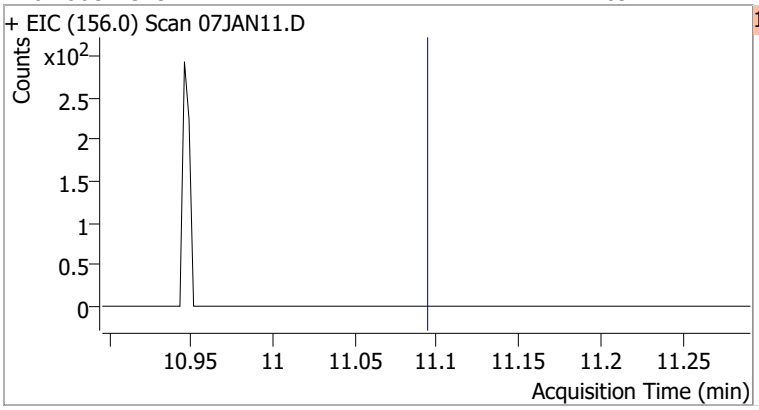
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	270.6993	10.95	0.00	237586	174.0	94.5	61.7	121.7
					176.0	92.2	60.6	120.6

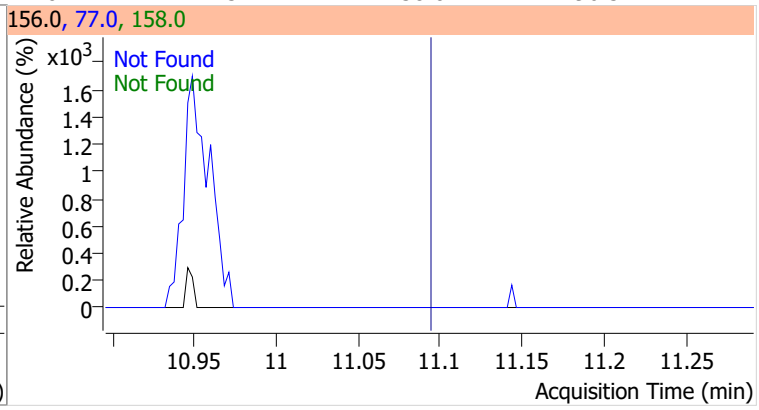




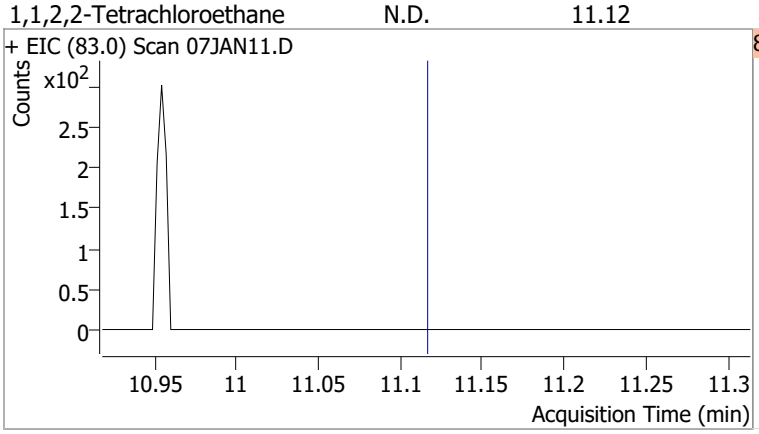


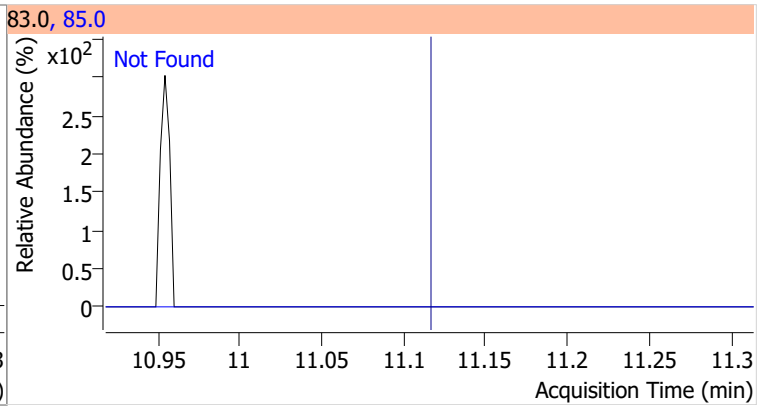
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



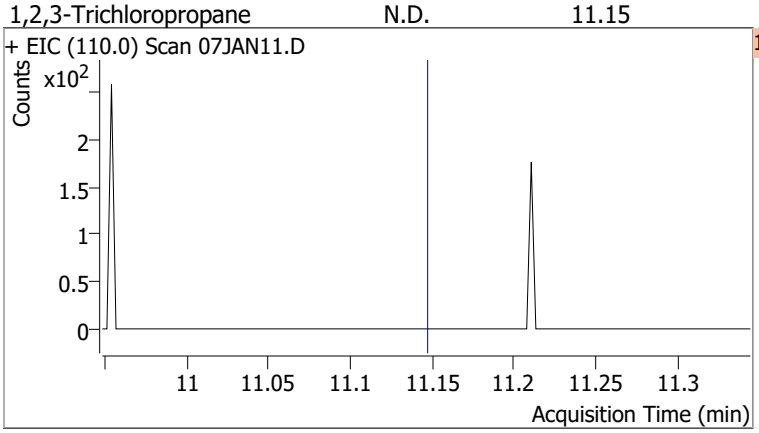


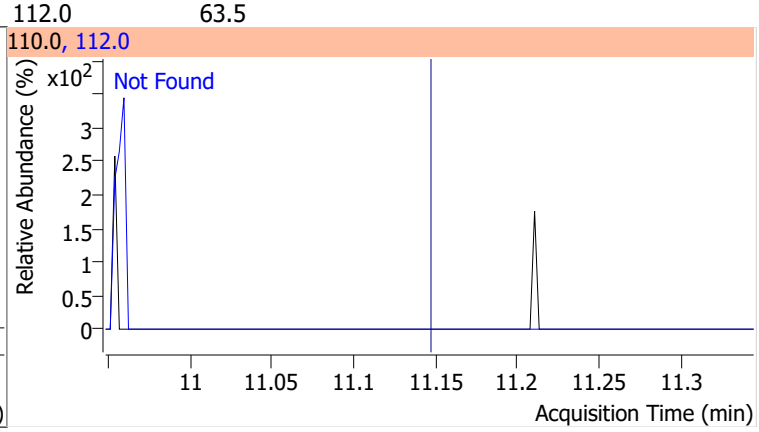
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



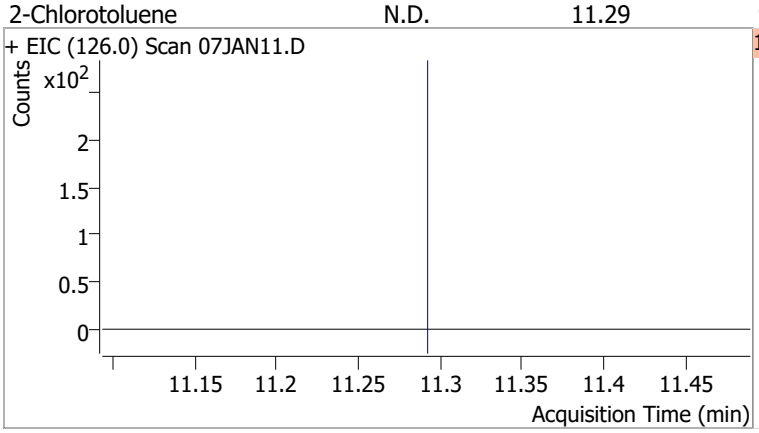


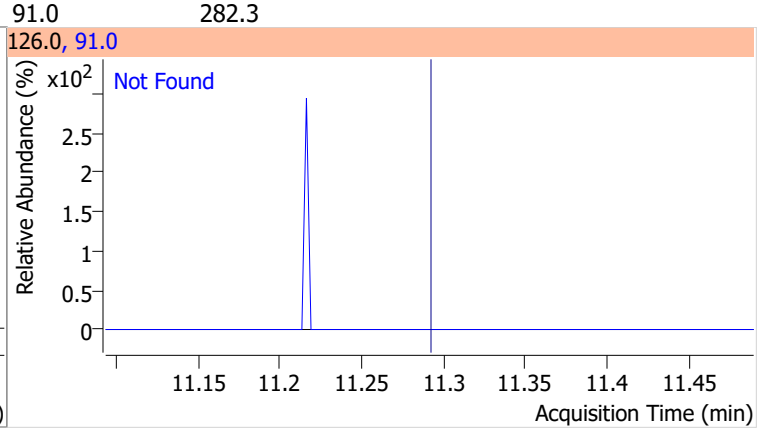
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



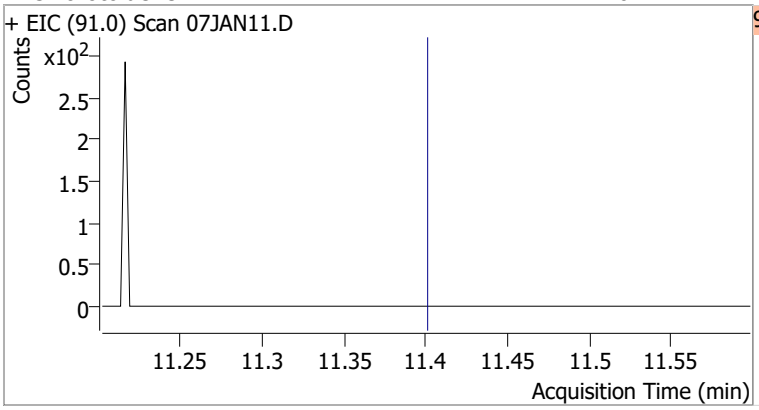


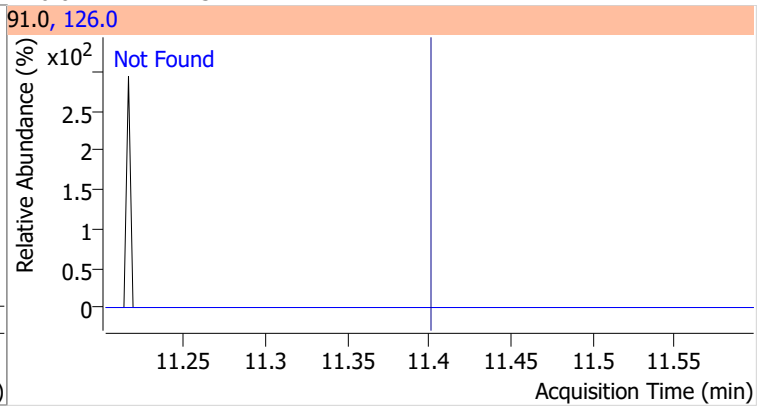
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



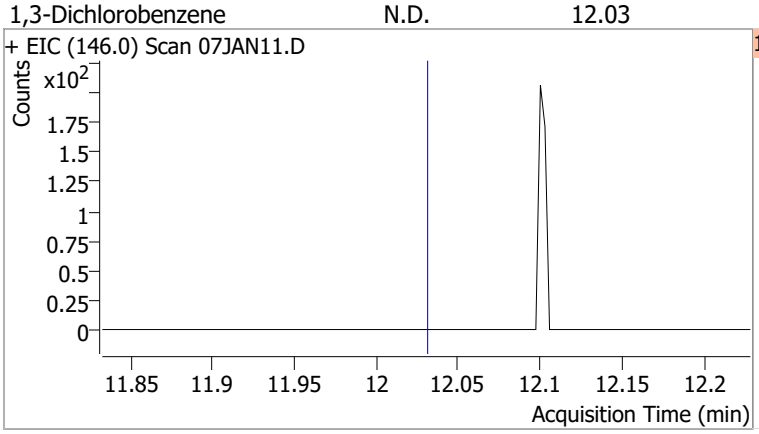


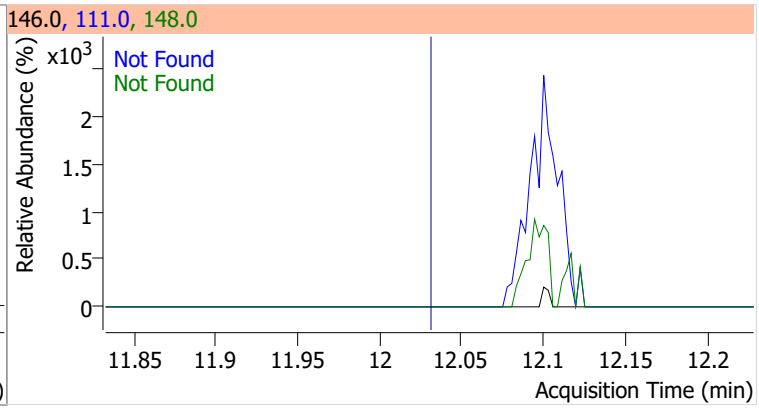
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



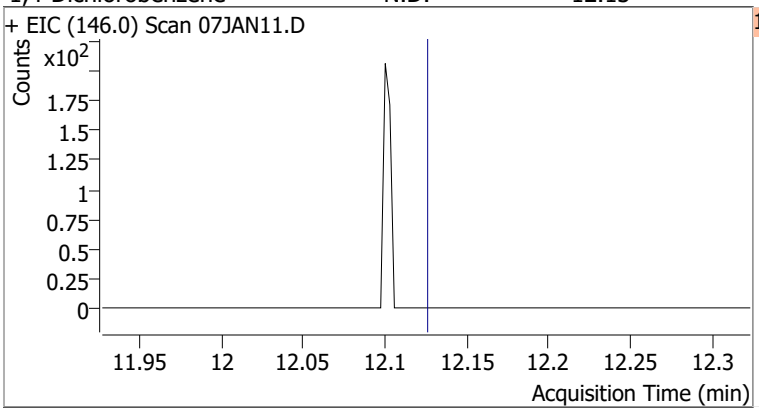


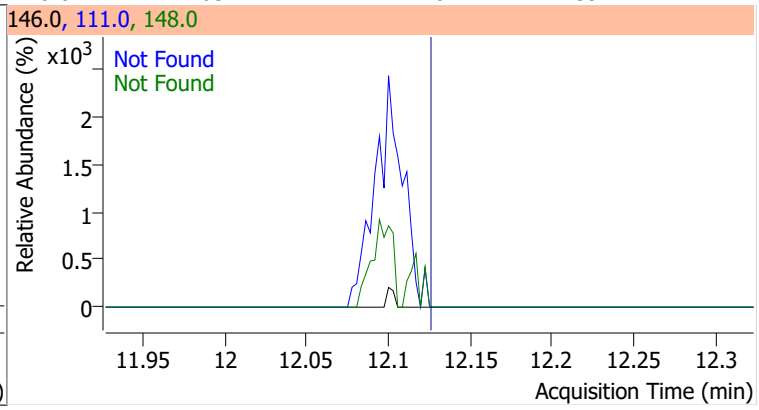
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



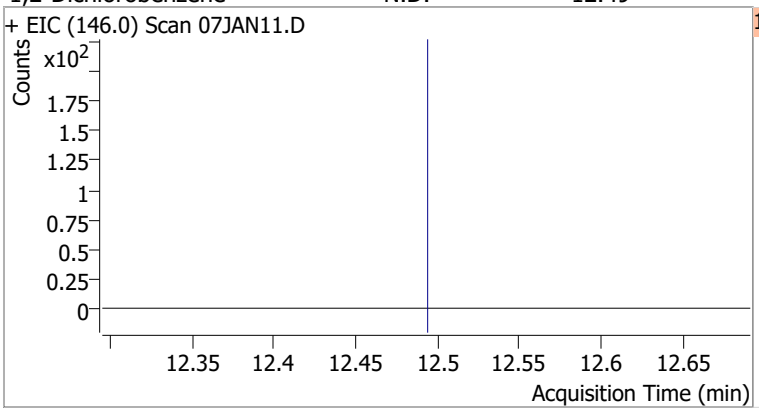


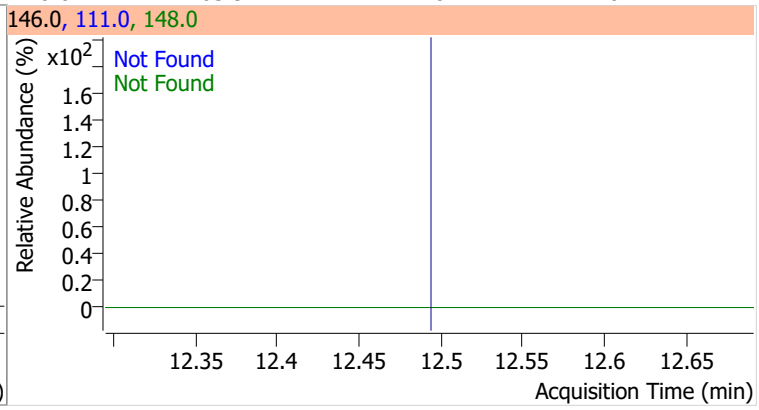
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





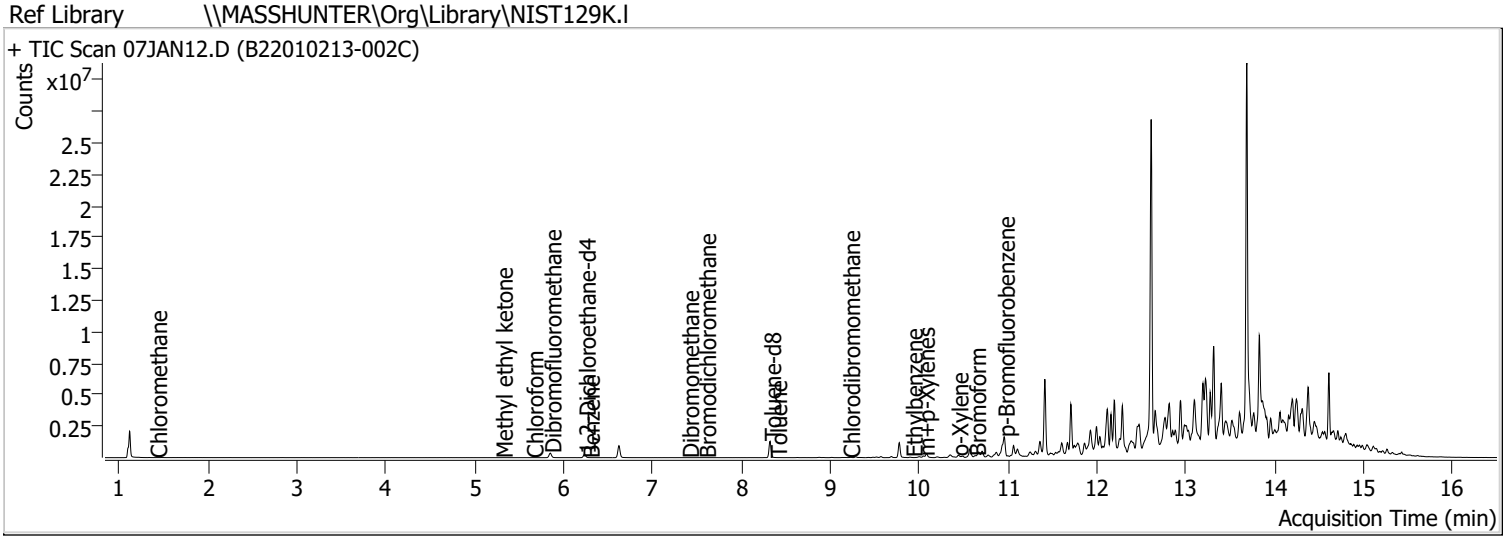
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN12.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 2:41:20 PM
Sample Name	B22010213-002C	Instrument	VOA5975C
Vial	12	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

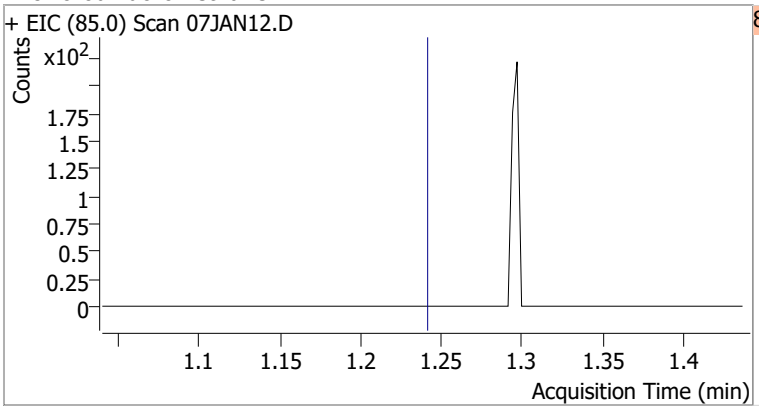


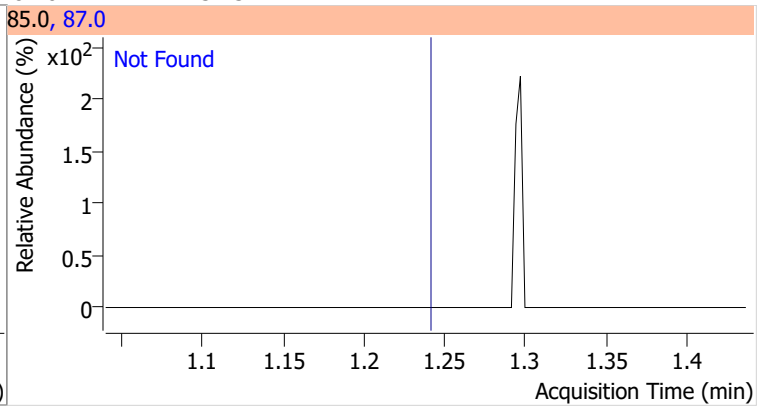
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.620	96.0	797906	250.0000	ng		-0.003
M Chlorobenzene-d5	9.774	82.0	317458	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	260553	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.851	113.0	209602	278.8342	ng		0.005
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 111.53%			
S 1,2-Dichloroethane-d4	6.236	67.0	94741	291.7947	ng		0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 116.72%			
S Toluene-d8	8.321	98.0	800313	261.6097	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 104.64%			
S p-Bromofluorobenzene	10.951	95.0	258089	270.3807	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 108.15%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.417	50.0	854	0.6730	ng	m	63
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	0.000		0	N.D.			
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	5.304	43.0	565	5.0179	ng	m	97
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	5.647	83.0	1655	1.0896	ng	m	89

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.280	78.0	409	0.1289	ng	m	74
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	7.398	93.0	111	0.3122	ng	m	86
T Bromodichloromethane	7.591	83.0	654	0.6660	ng	m	93
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.388	92.0	493	0.2388	ng	m	97
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	9.211	129.0	1602	2.4889	ng	m	80
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	9.917	91.0	1441	0.3673	ng	m	62
T m+p-Xylenes	10.050	106.0	1346	0.8826	ng	m	82
T o-Xylene	10.430	106.0	3810	2.8067	ng		82
T Styrene	0.000		0	N.D.			
T Bromoform	10.628	172.5	2322	6.9642	ng		98
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	11.099	83.0	0		ng	md	1
T 1,2,3-Trichloropropane	11.146	110.0	0		ng	md	1
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	11.392	91.0	0		ng	md	1
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

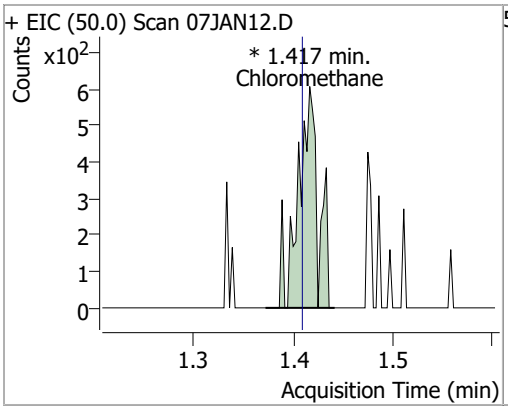
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

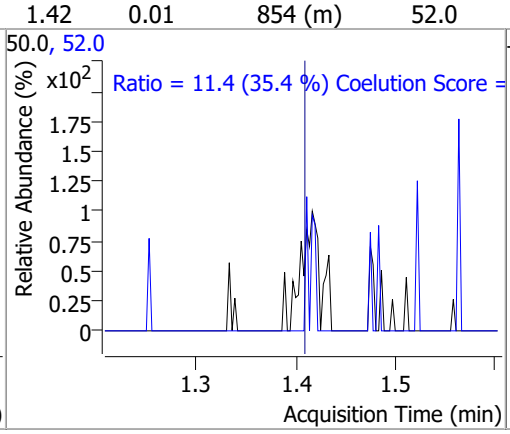
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

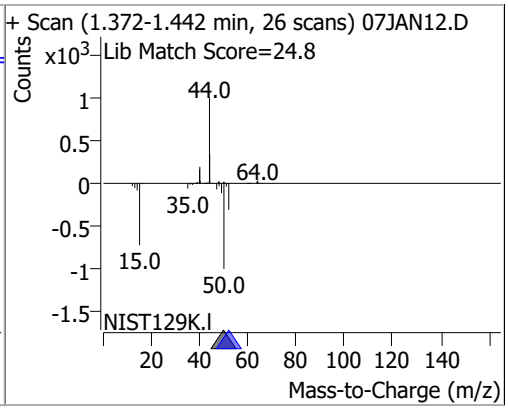




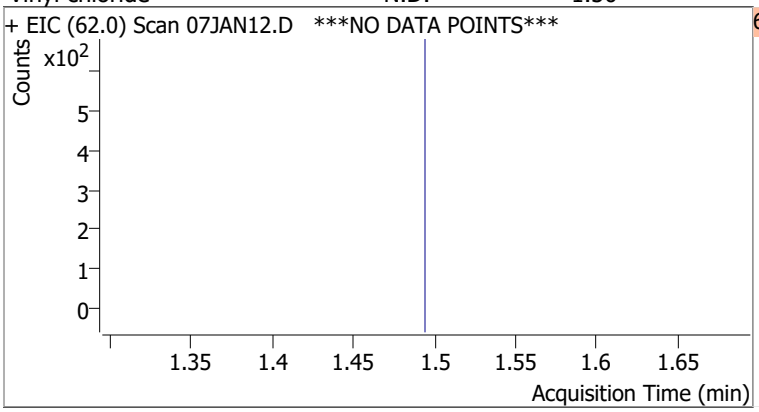
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	0.6730	1.42	0.01	854 (m)	52.0	11.4	2.1	62.1

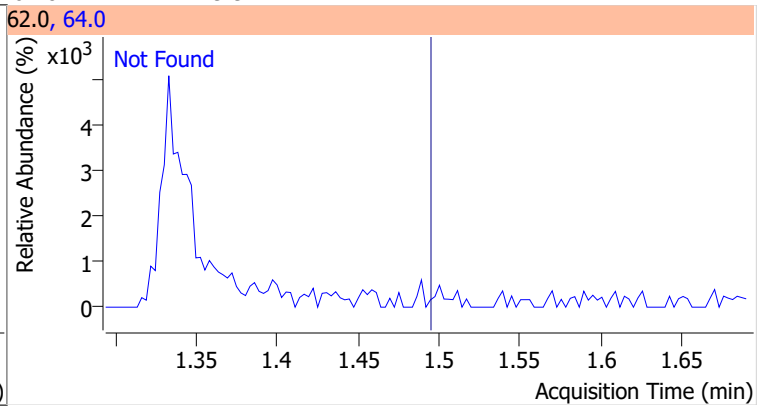




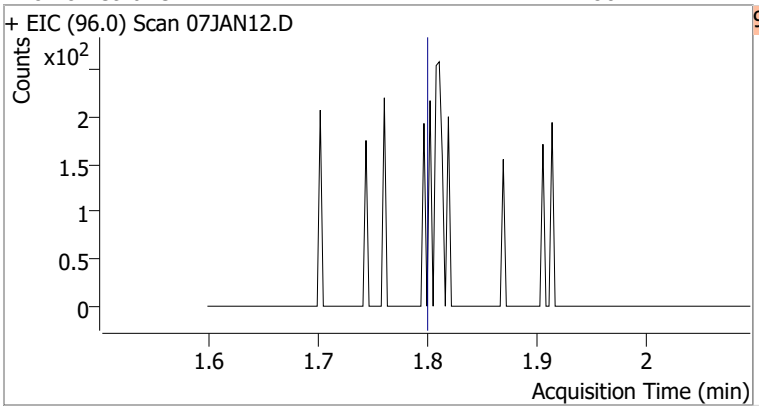


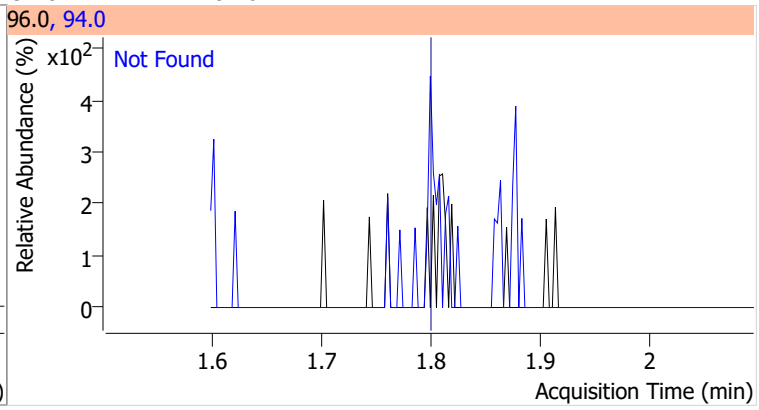
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



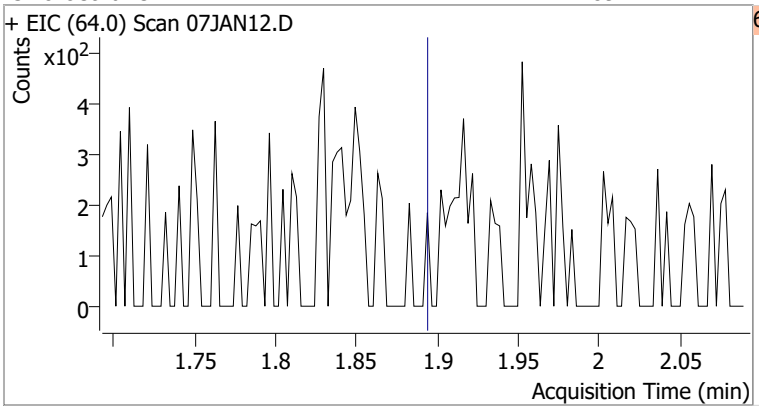


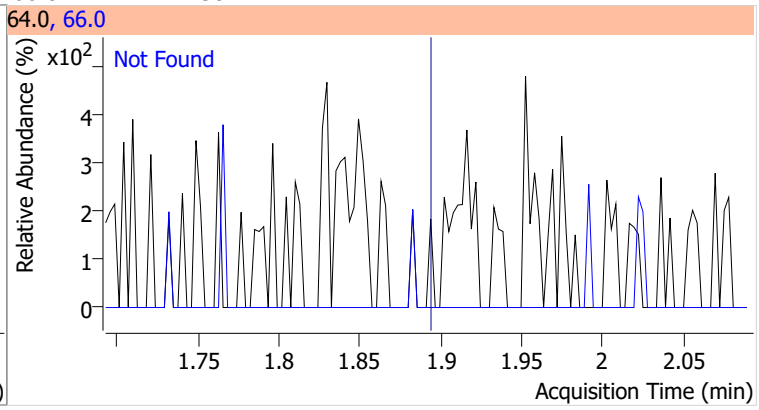
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



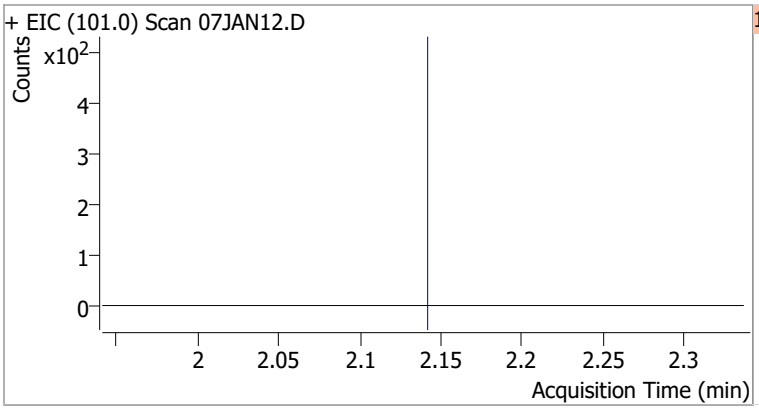


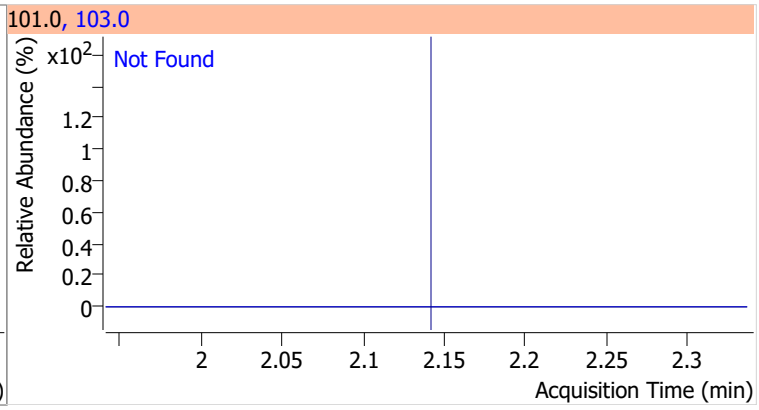
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



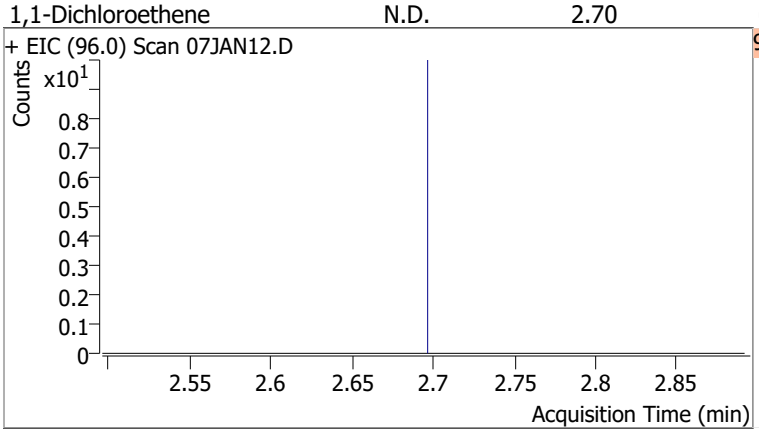


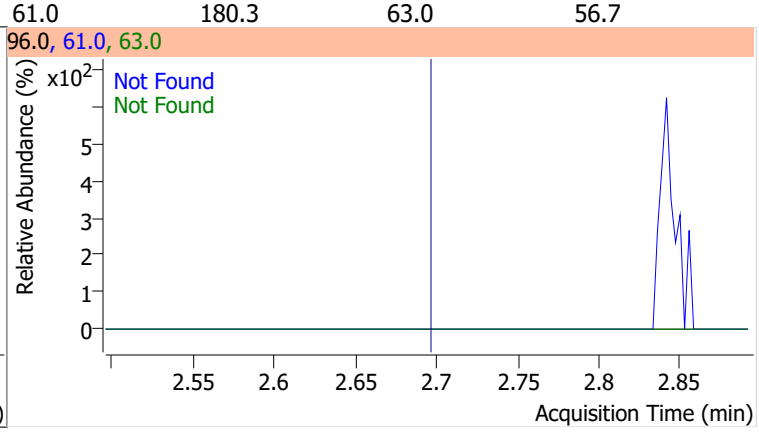
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



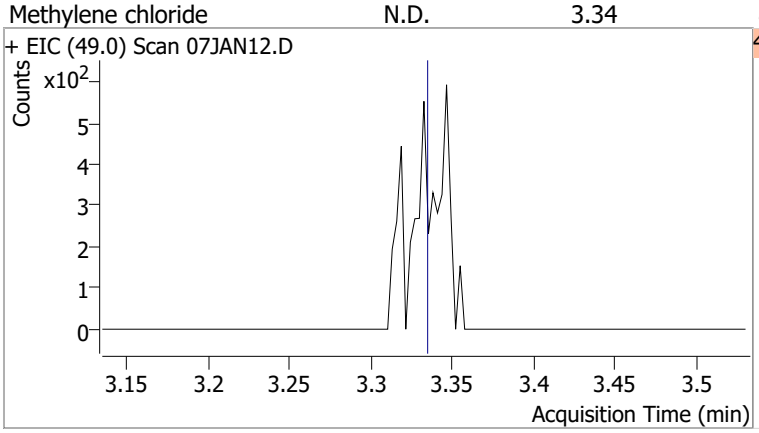


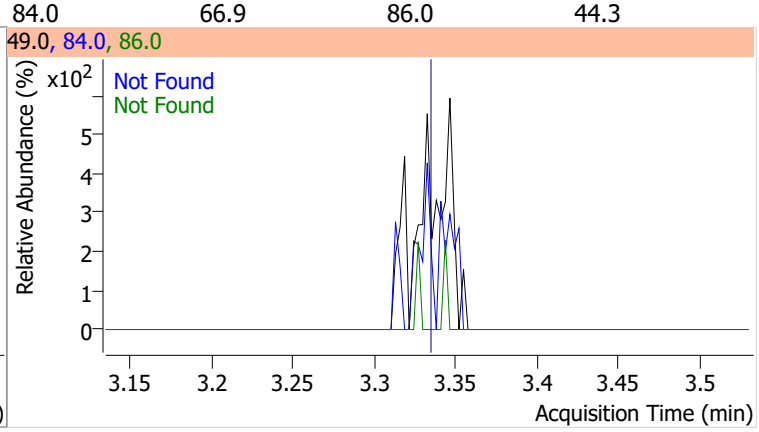
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7



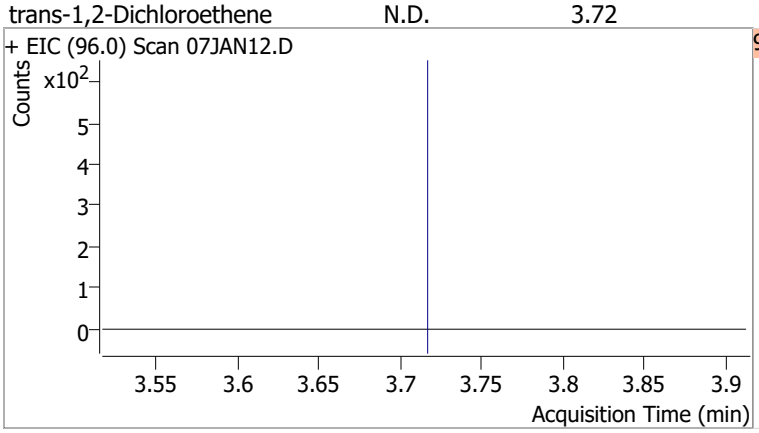


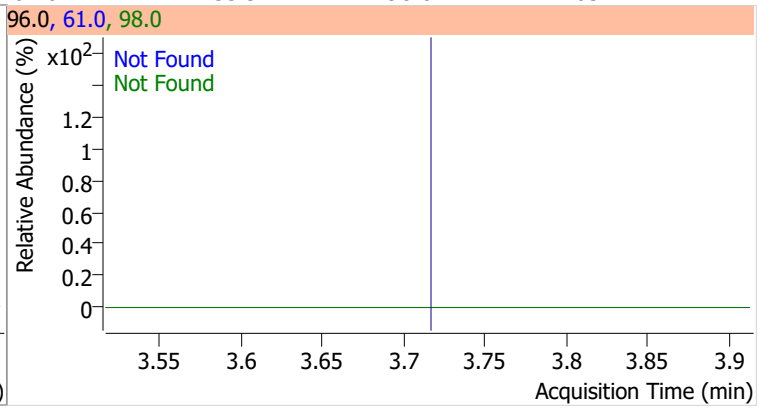
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Methylene chloride	N.D.	3.34	84.0	66.9	86.0	44.3



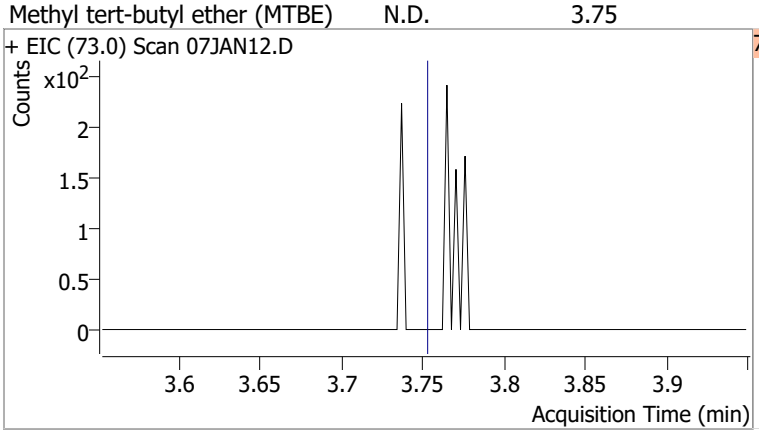


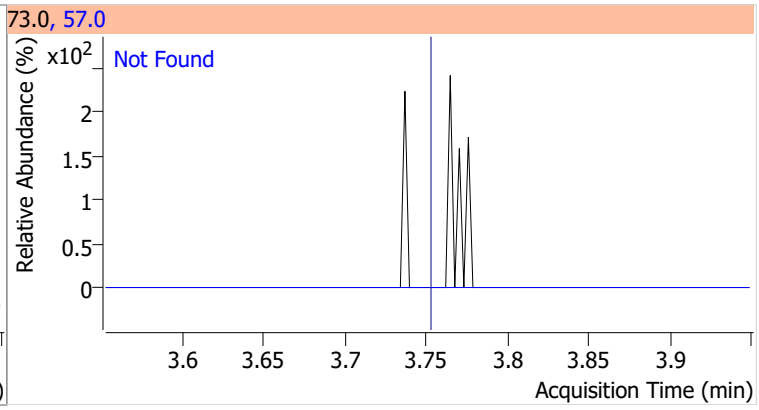
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



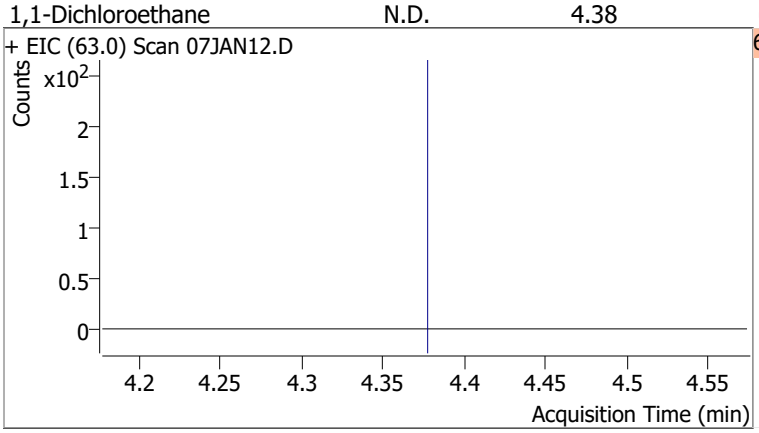


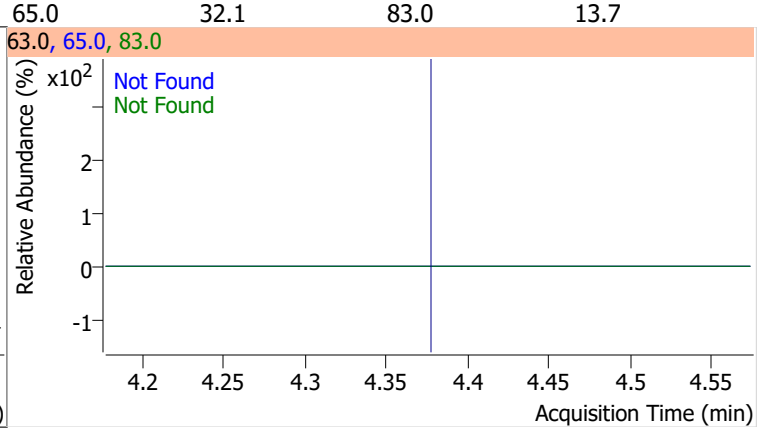
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



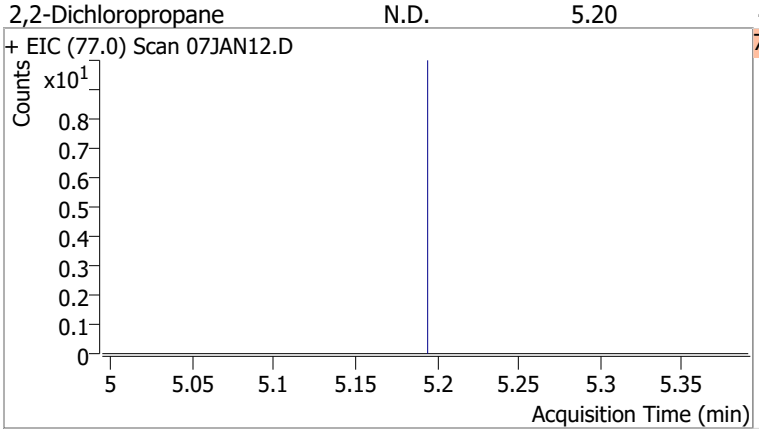


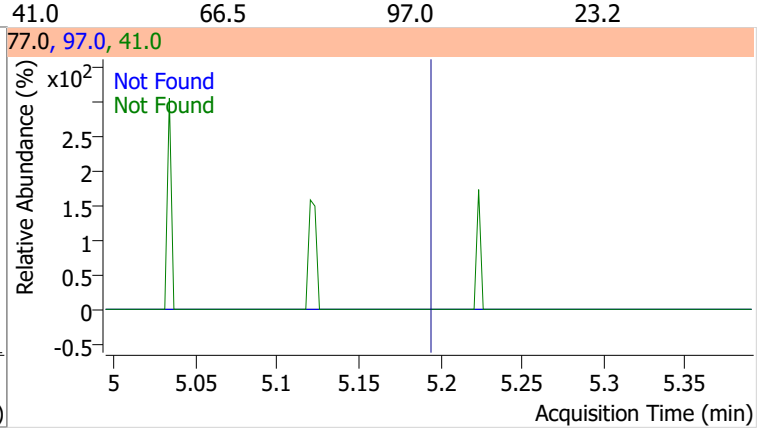
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



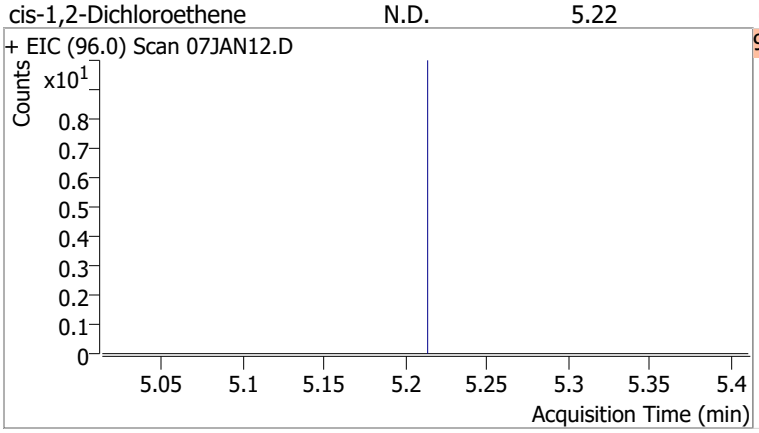


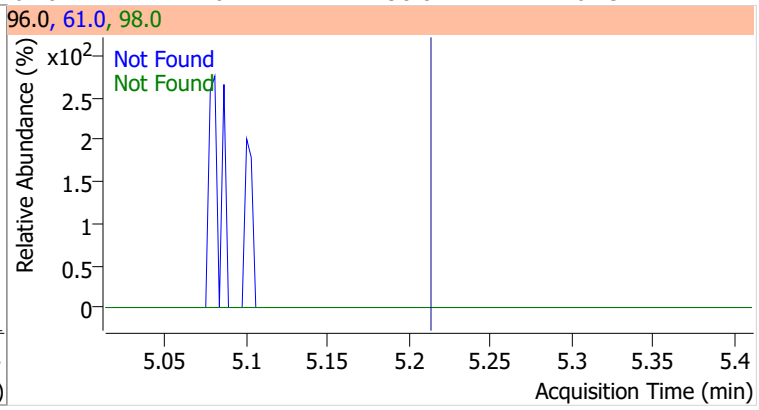
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



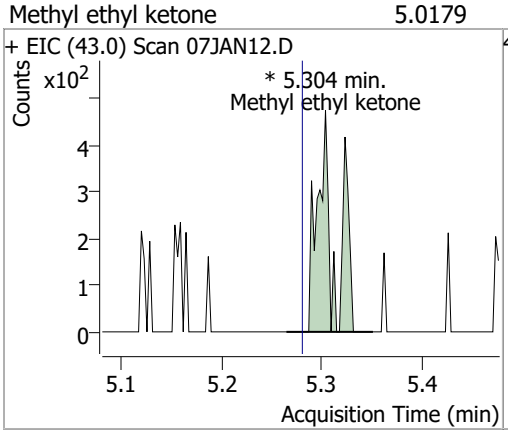


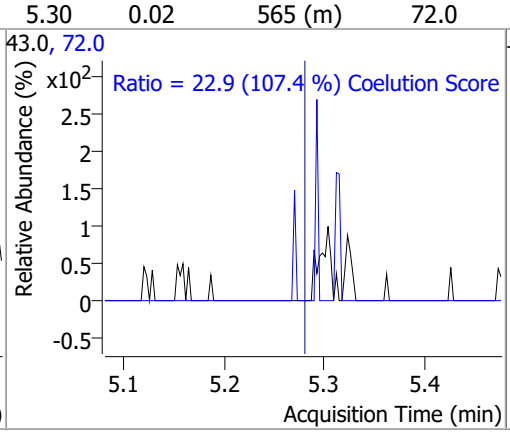
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3

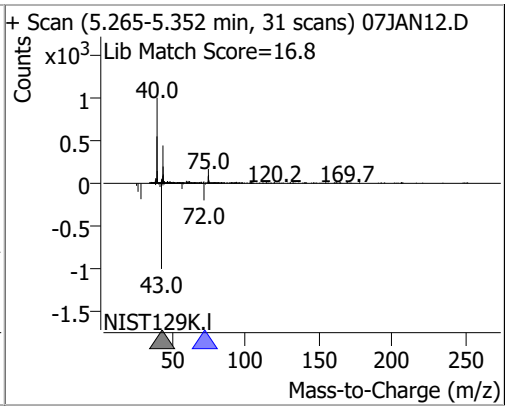




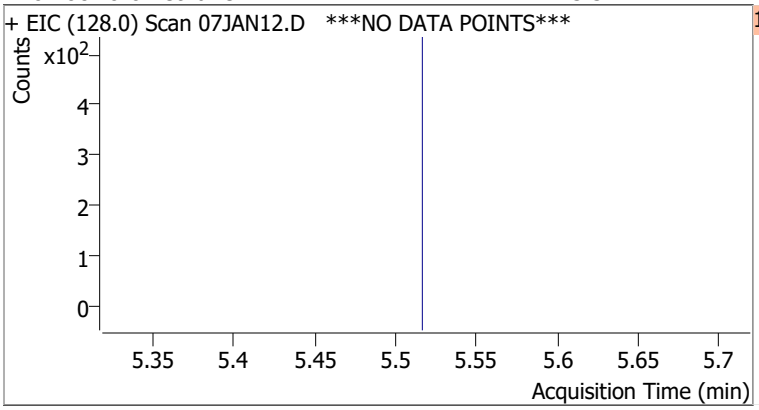
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	5.0179	5.30	0.02	565 (m)	72.0	22.9	0.0	51.3

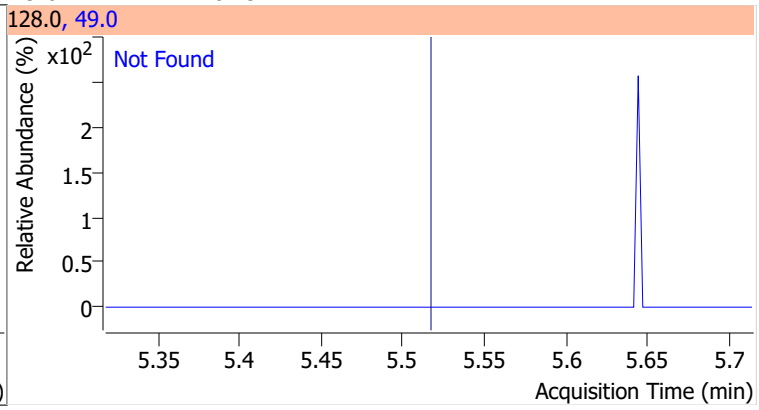




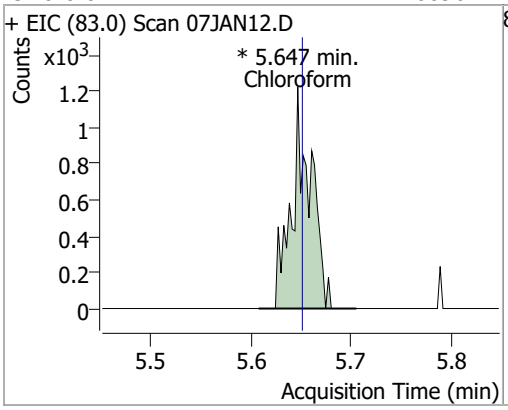


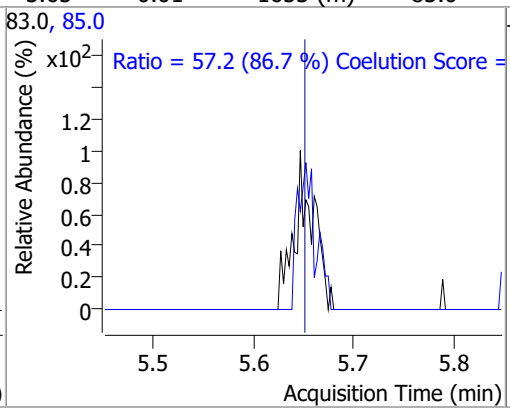
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9

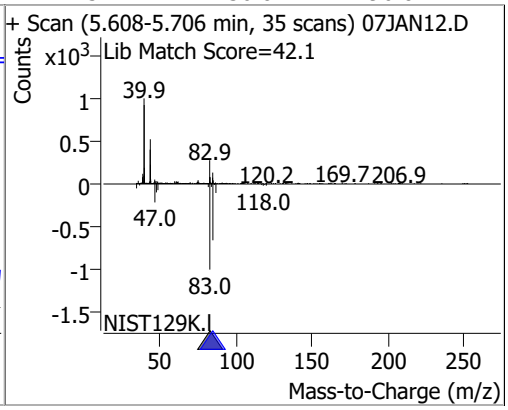




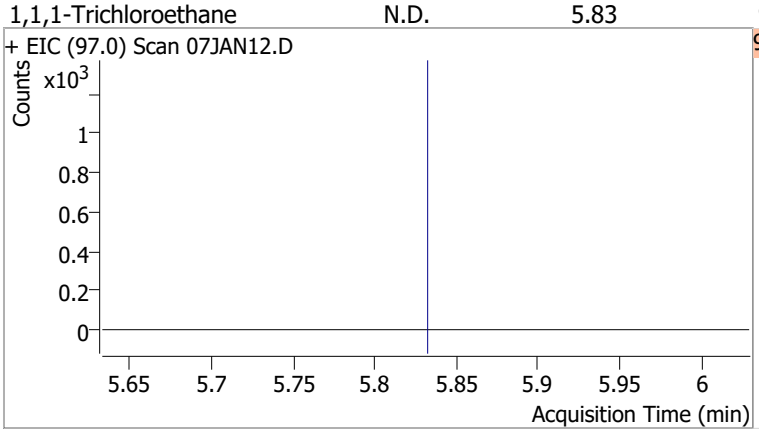
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	1.0896	5.65	-0.01	1655 (m)	85.0	57.2	36.0	96.0

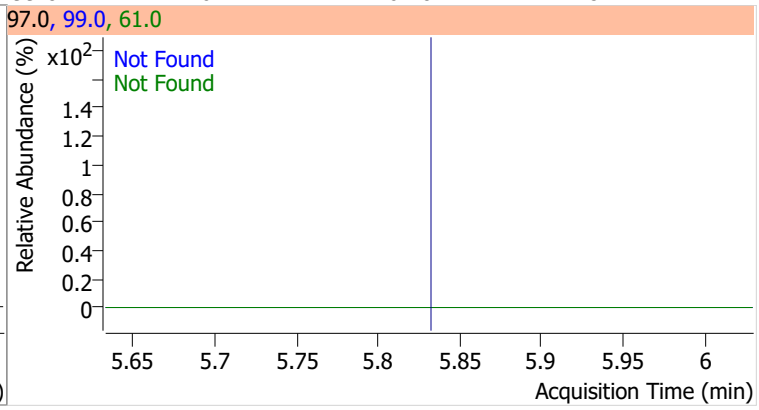




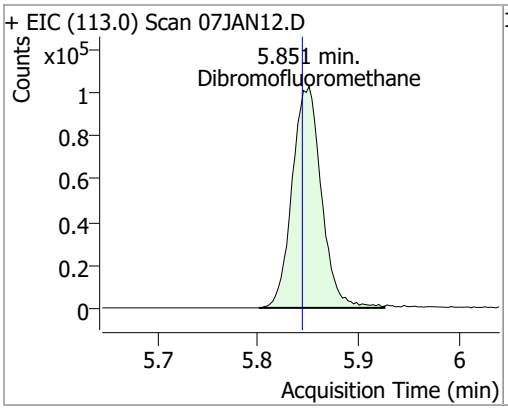


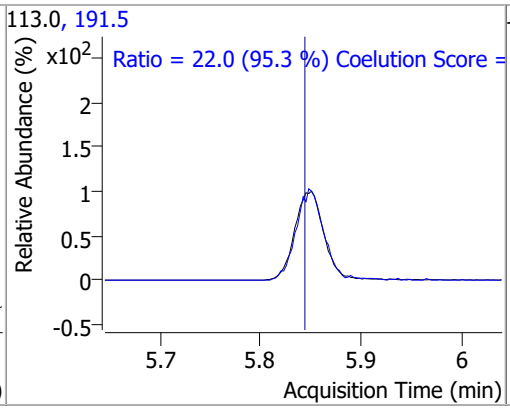
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

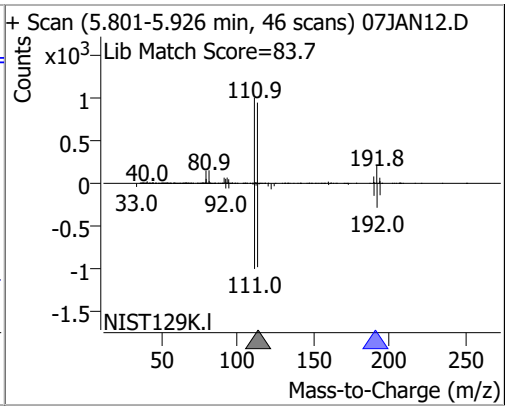




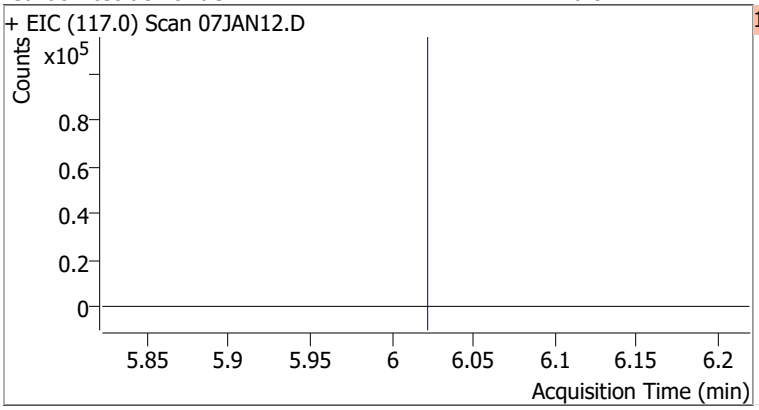
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	278.8342	5.85	0.01	209602	191.5	22.0	0.0	53.1

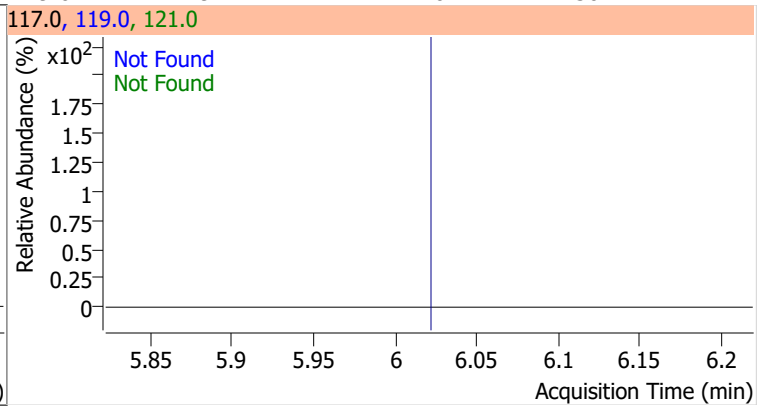




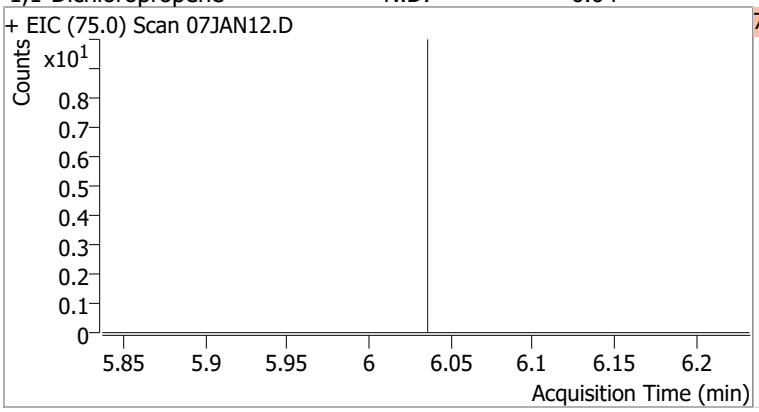


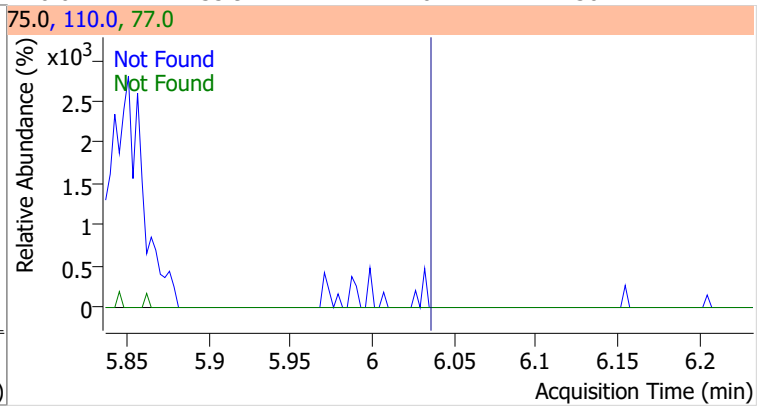
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



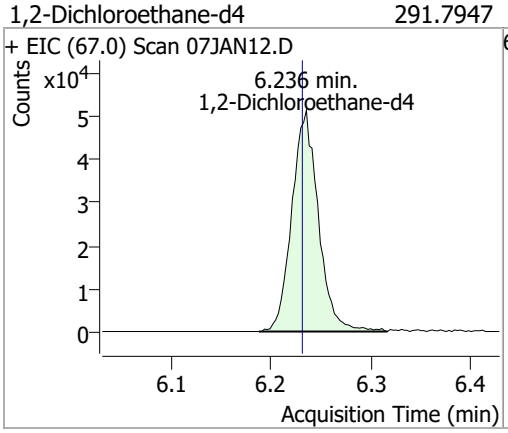


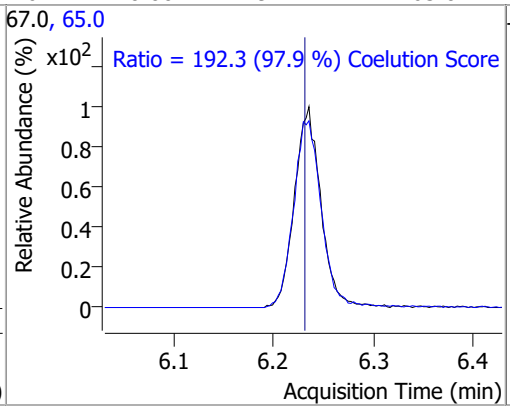
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

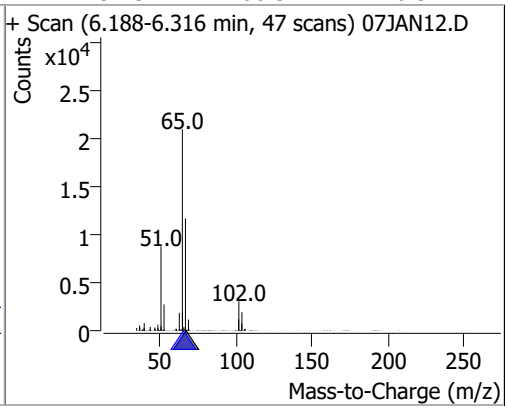




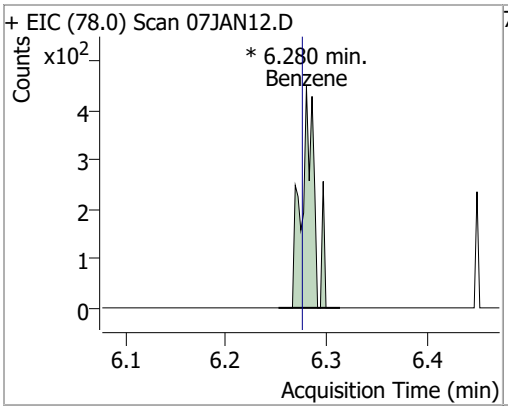
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	291.7947	6.24	0.00	94741	65.0	192.3	166.5	226.5

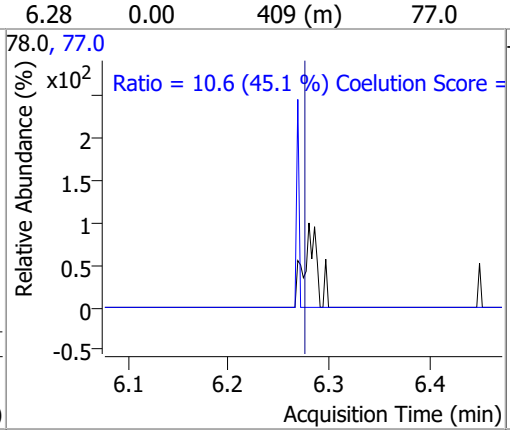


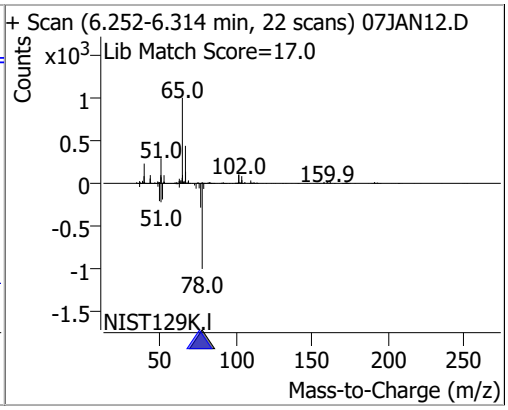




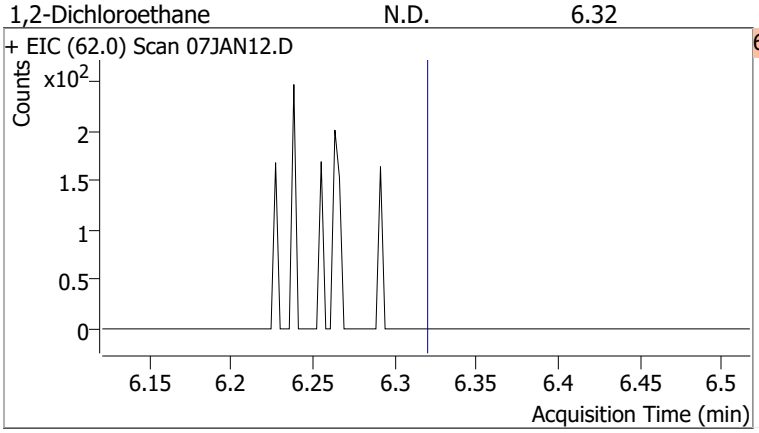
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.1289	6.28	0.00	409 (m)	77.0	10.6	0.0	53.5

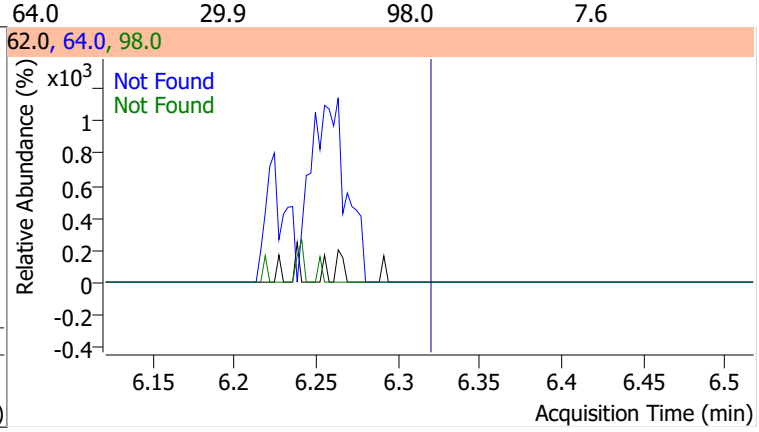




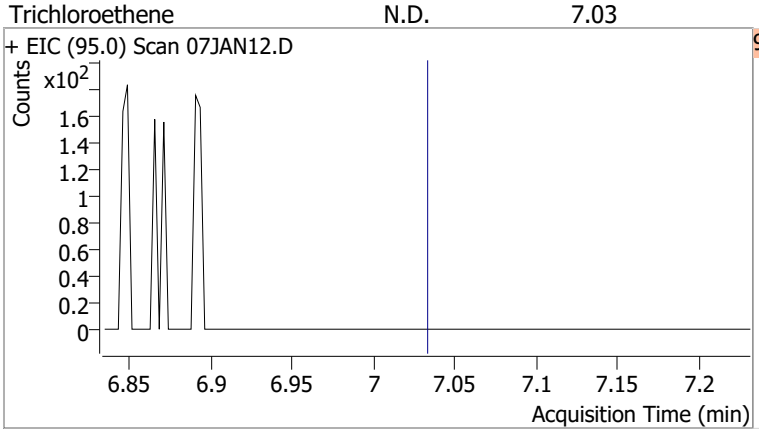


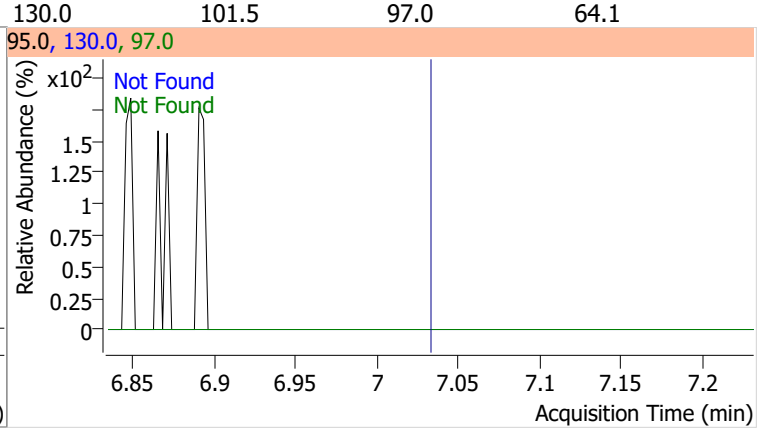
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



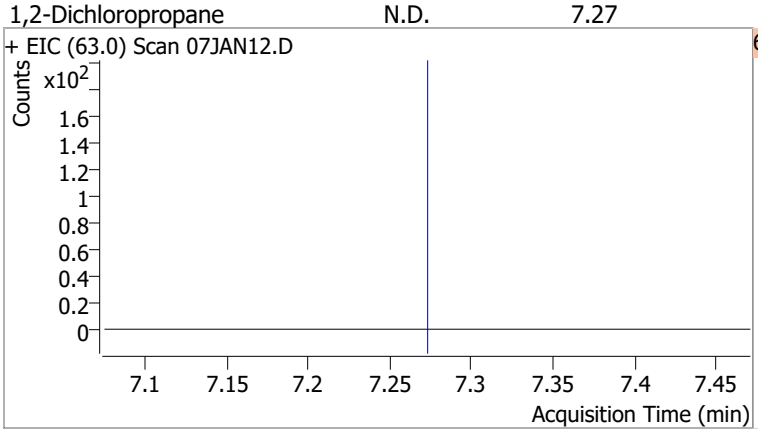


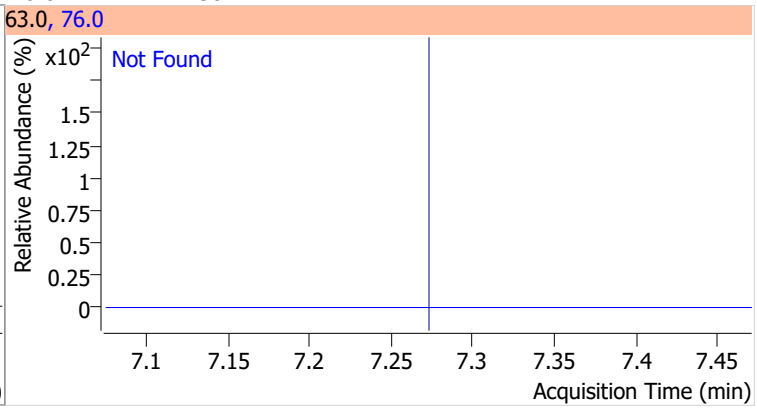
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



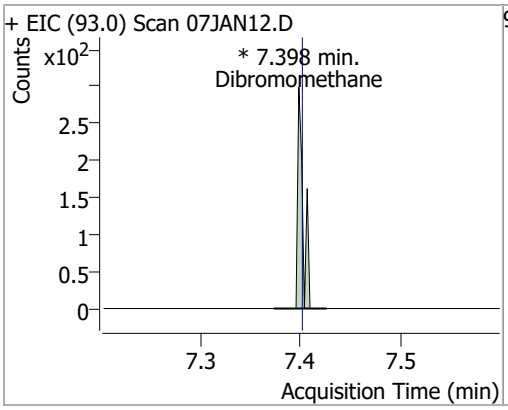


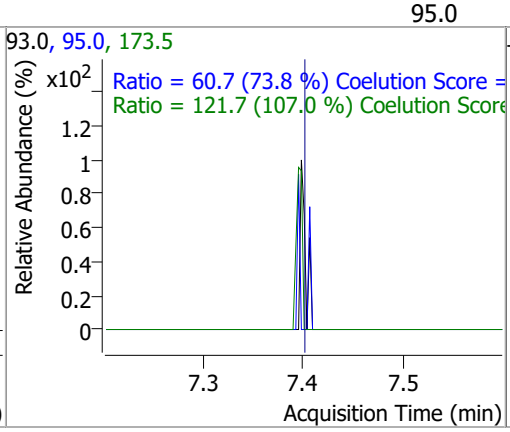
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2

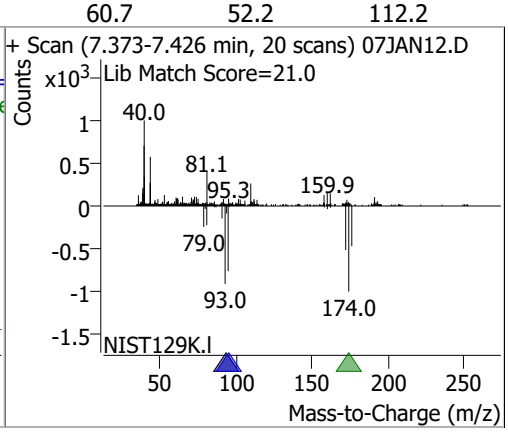




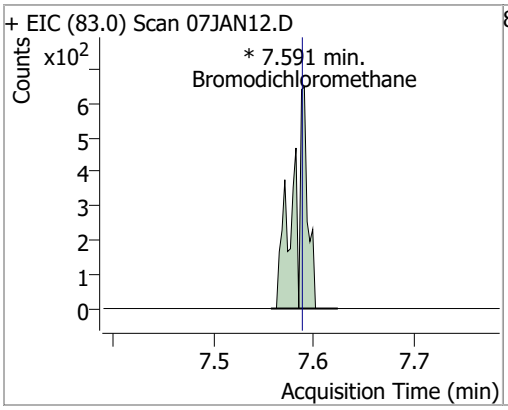
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	0.3122	7.40	0.00	111 (m)	173.5	121.7	83.7	143.7
					95.0	60.7	52.2	112.2

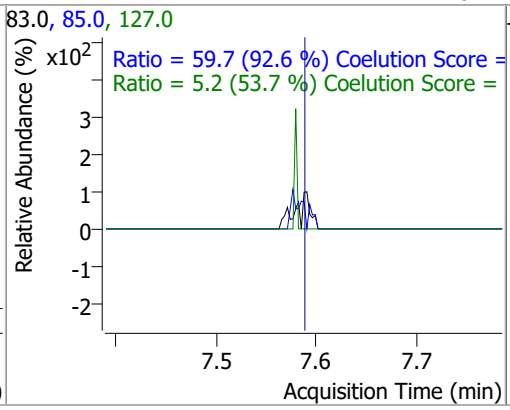


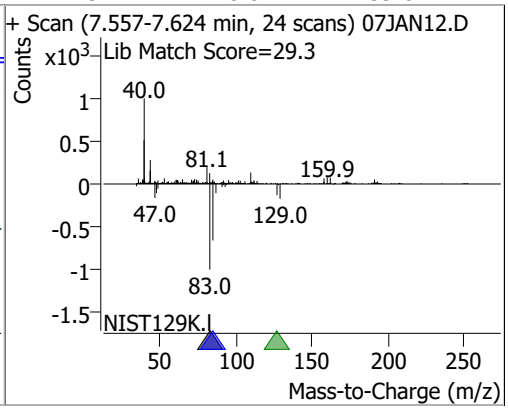




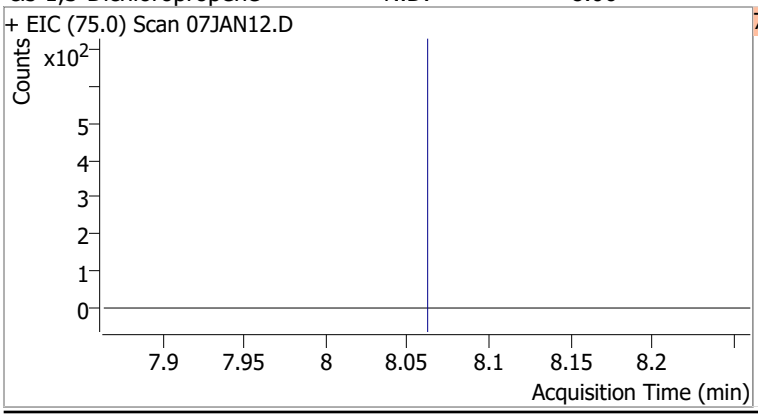
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	0.6660	7.59	0.01	654 (m)	85.0	59.7	34.5	94.5
					127.0	5.2	0.0	39.6

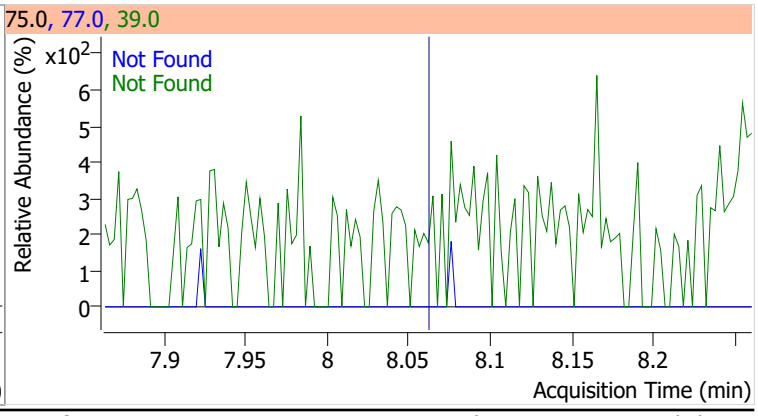




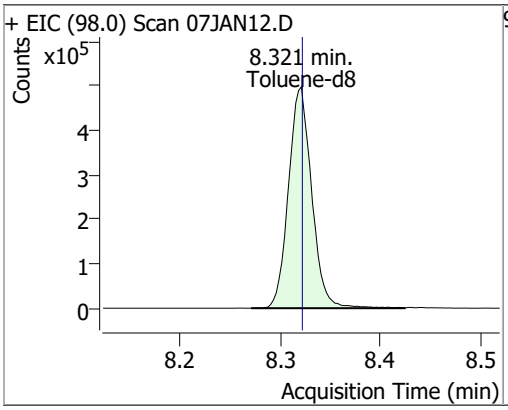


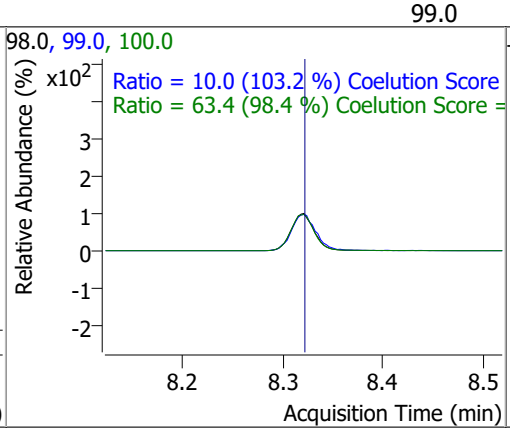
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

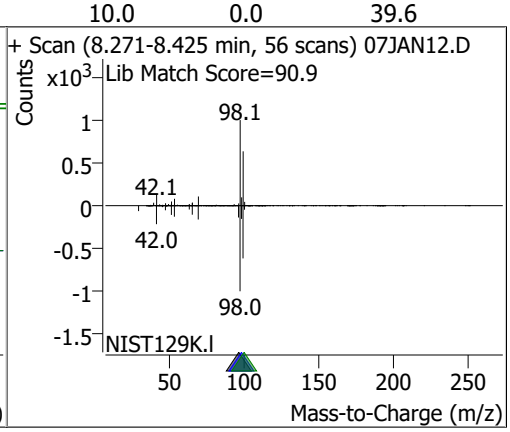




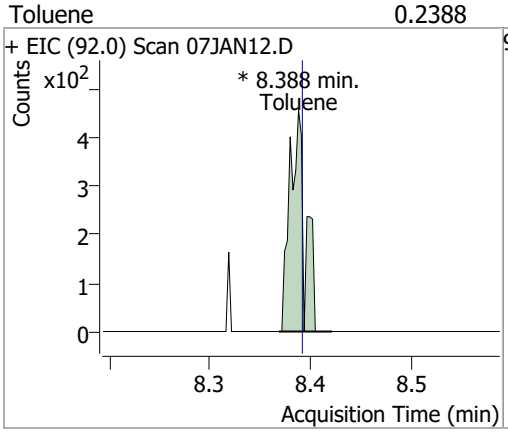
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	261.6097	8.32	0.00	800313	100.0	63.4	34.4	94.4
					99.0	10.0	0.0	39.6

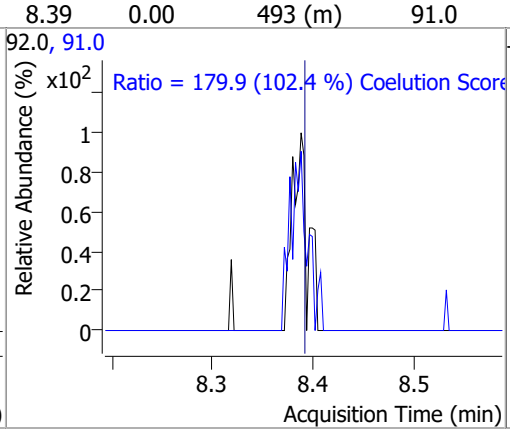


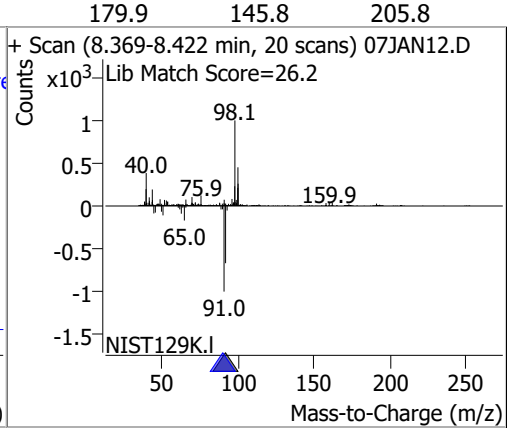




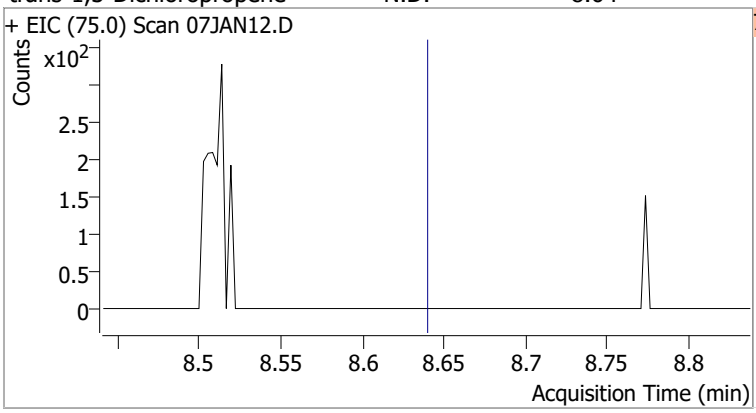
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	0.2388	8.39	0.00	493 (m)	91.0	179.9	145.8	205.8

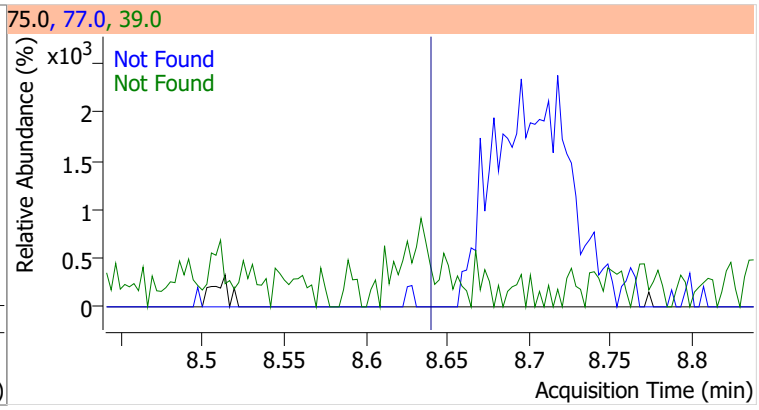




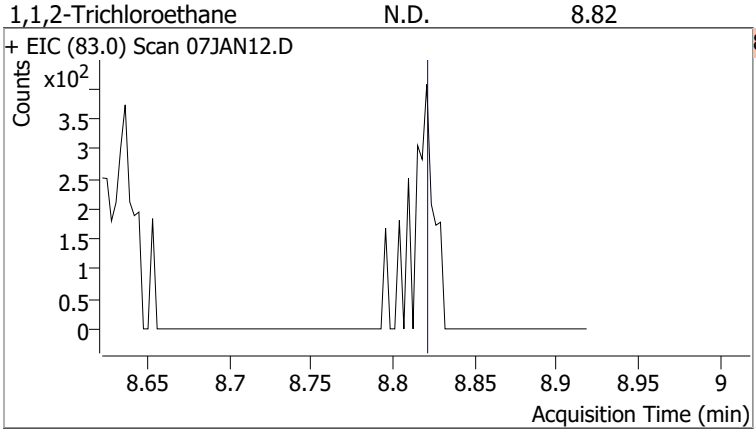


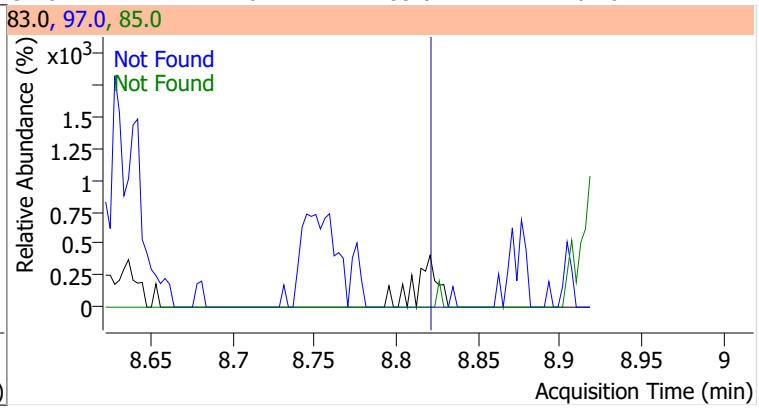
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



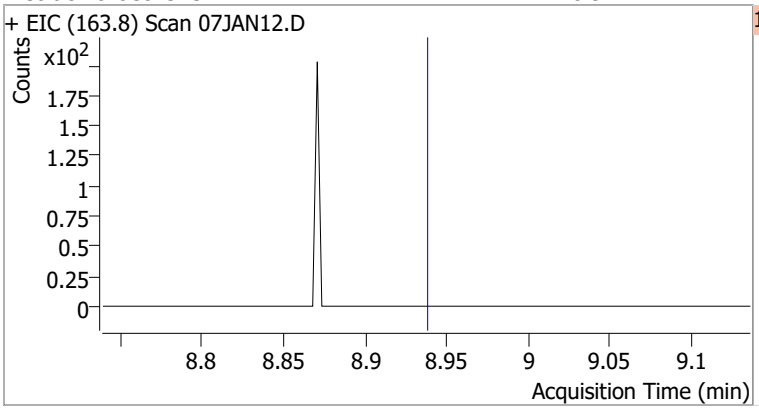


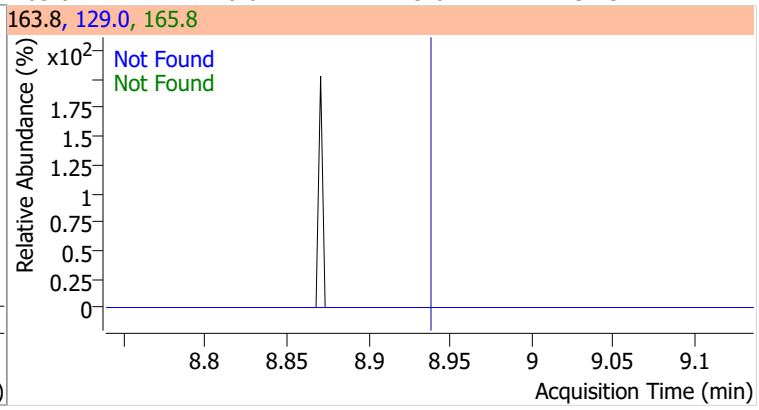
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



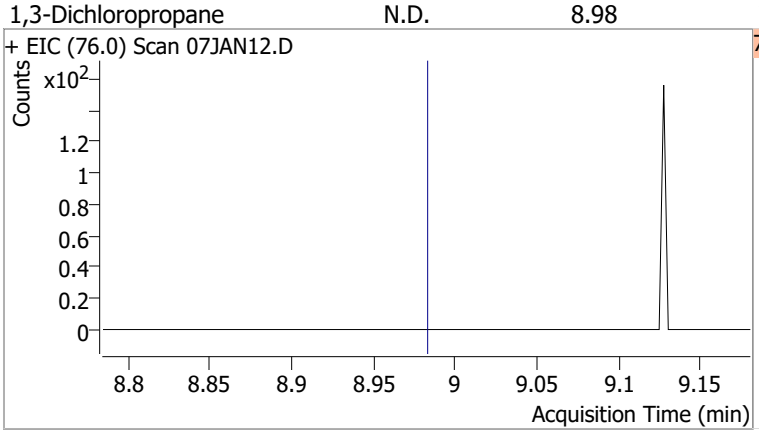


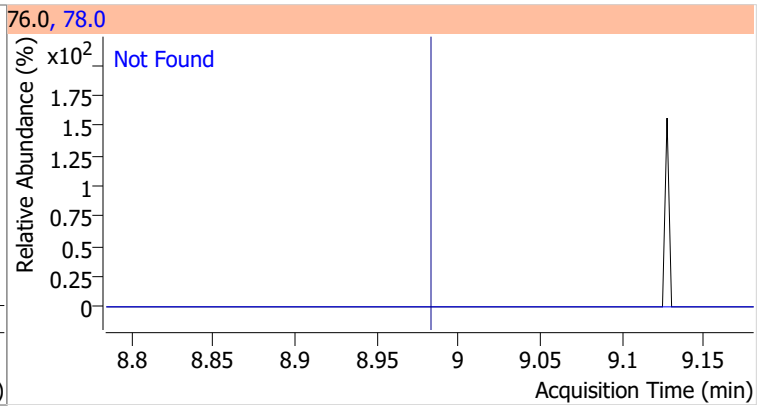
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



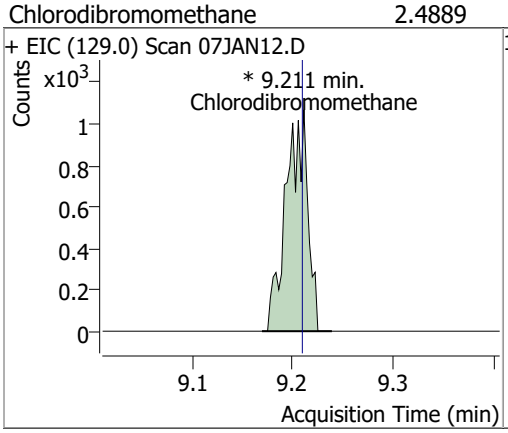


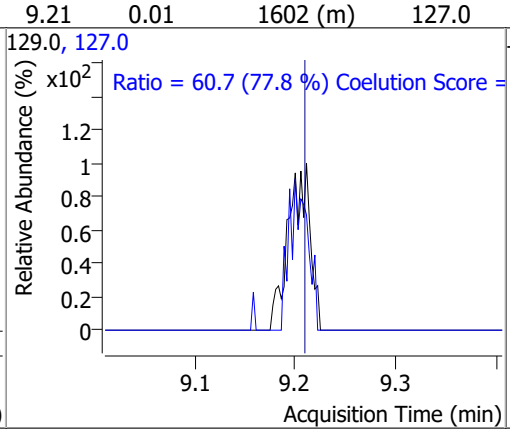
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9

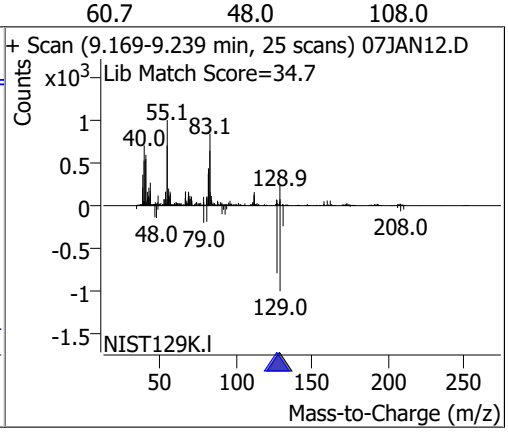




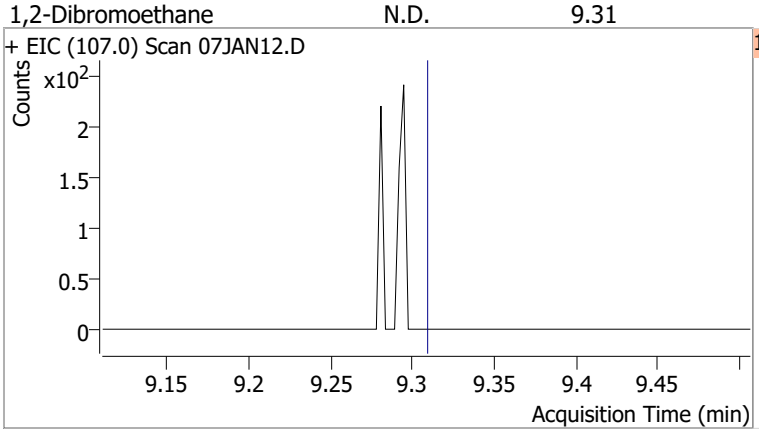
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	2.4889	9.21	0.01	1602 (m)	127.0	60.7	48.0	108.0

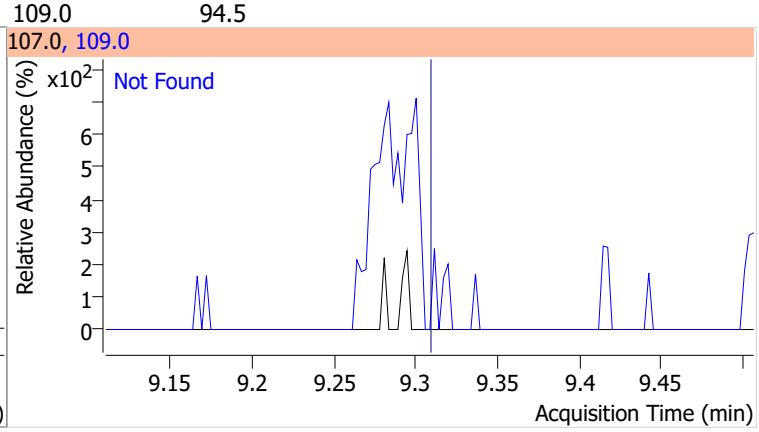




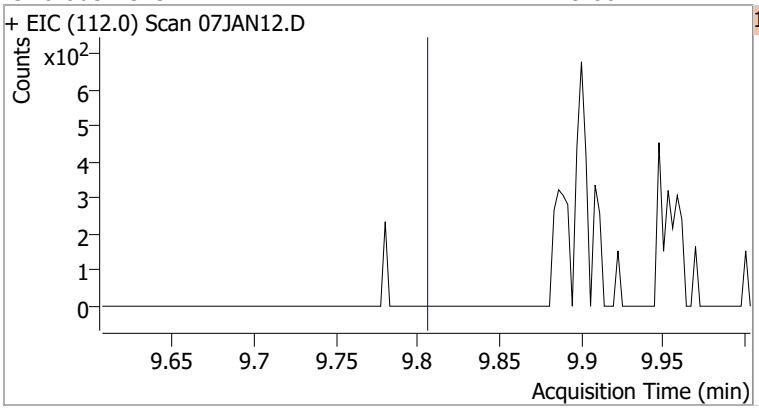


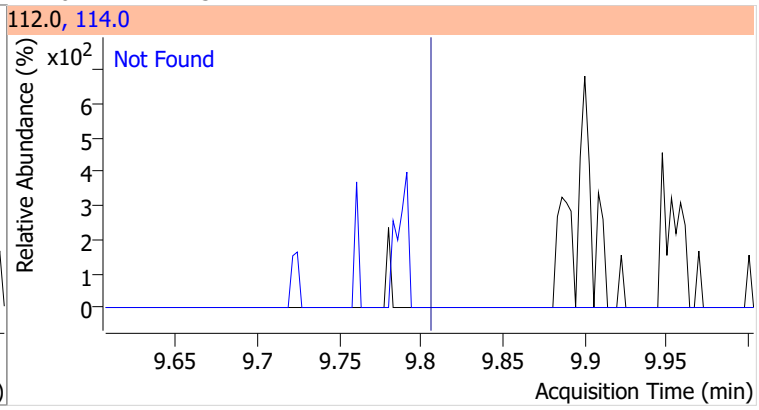
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



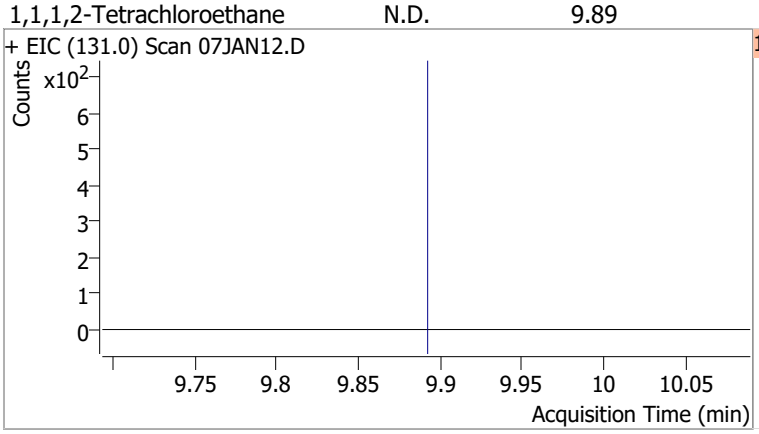


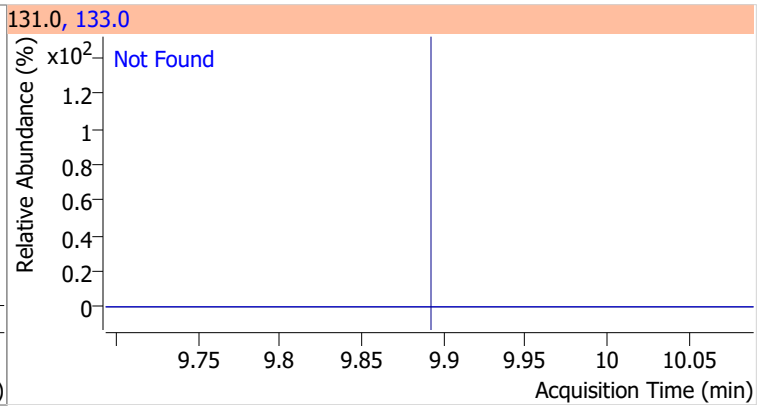
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



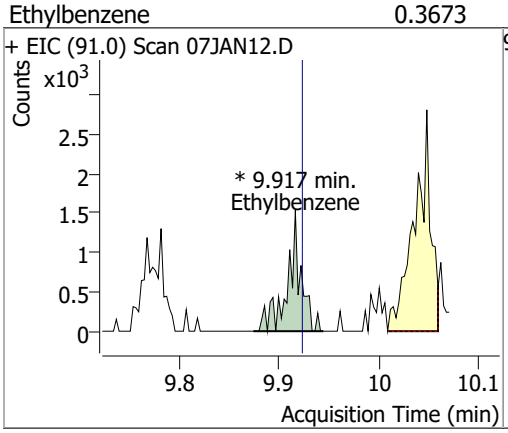


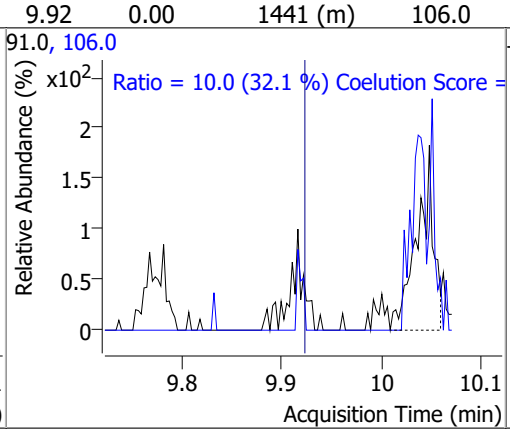
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6

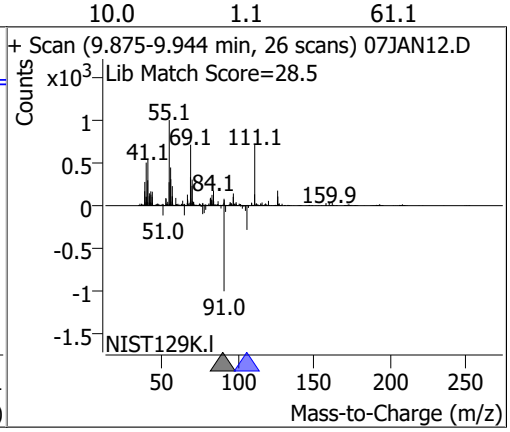




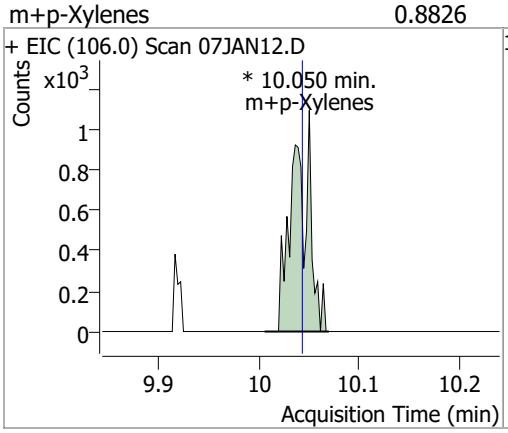
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	0.3673	9.92	0.00	1441 (m)	106.0	10.0	1.1	61.1

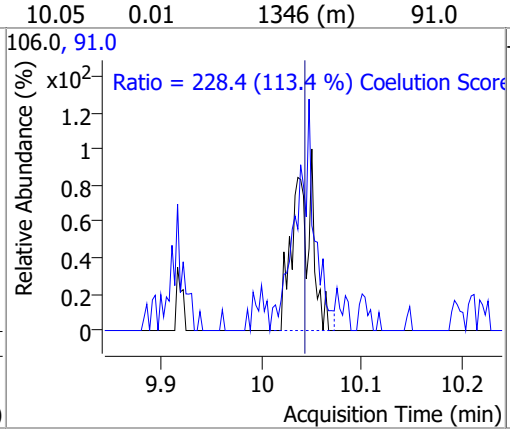


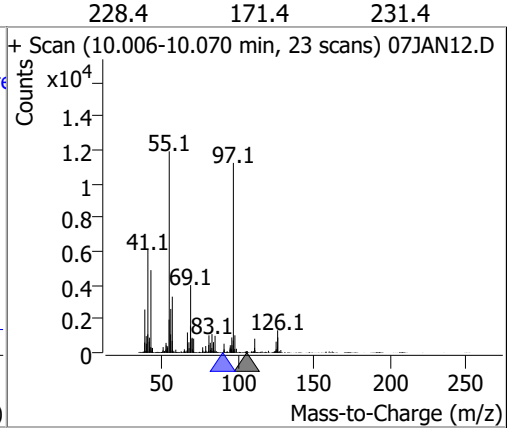




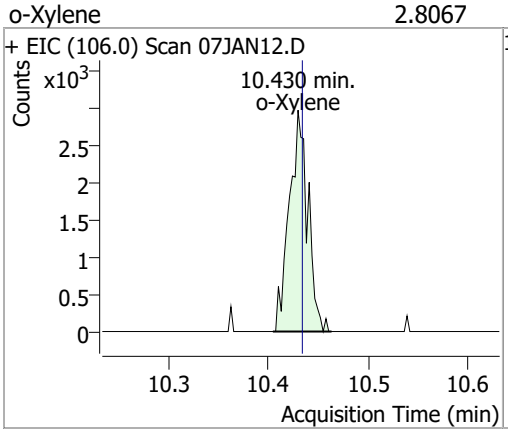
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	0.8826	10.05	0.01	1346 (m)	91.0	228.4	171.4	231.4

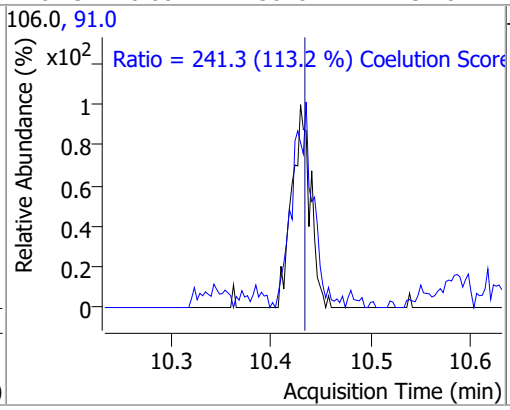


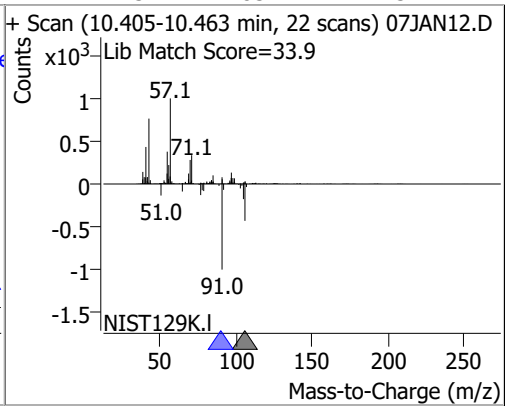




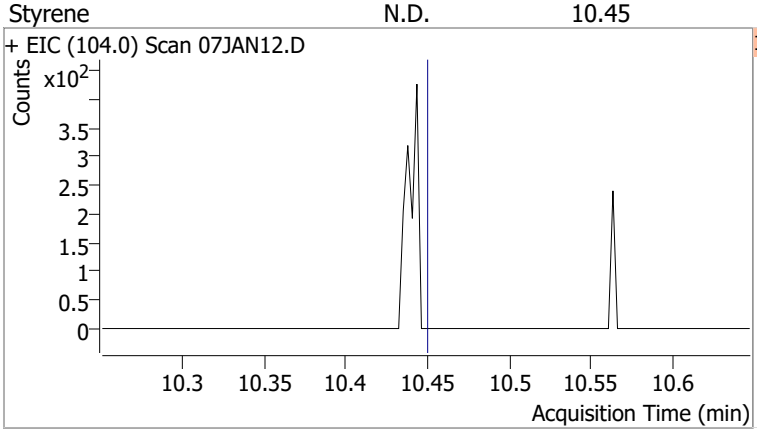
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	2.8067	10.43	0.00	3810	91.0	241.3	183.1	243.1

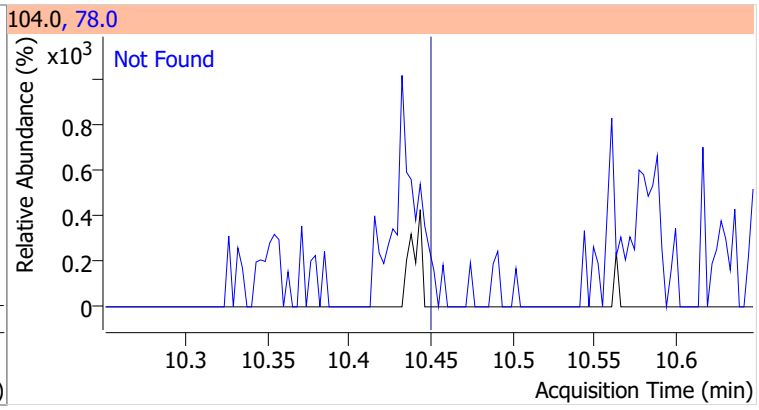




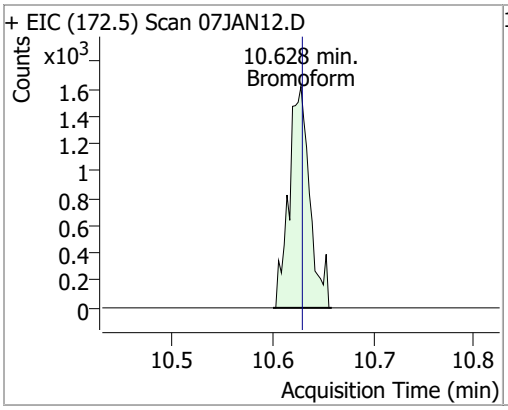


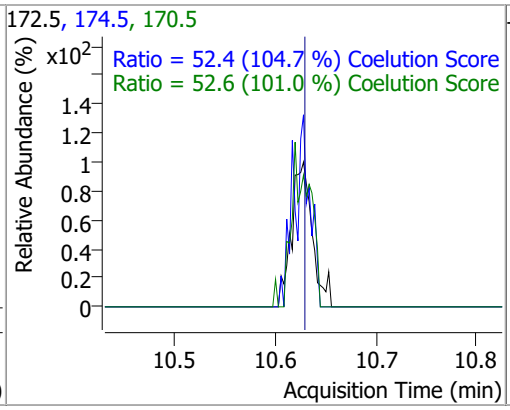
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6

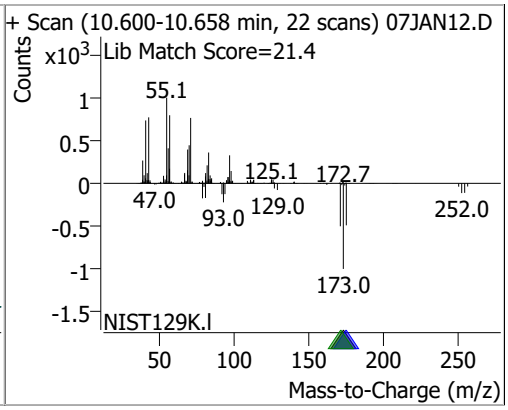




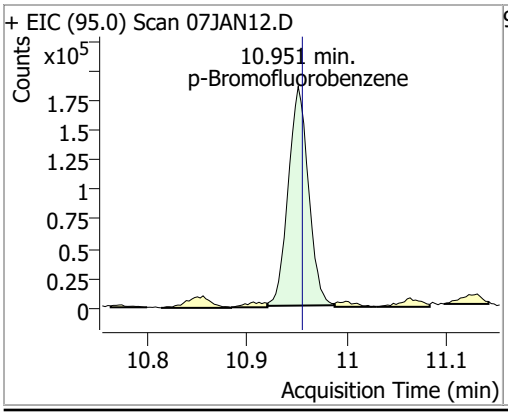
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	6.9642	10.63	0.00	2322	170.5	52.6	22.1	82.1
					174.5	52.4	20.1	80.1

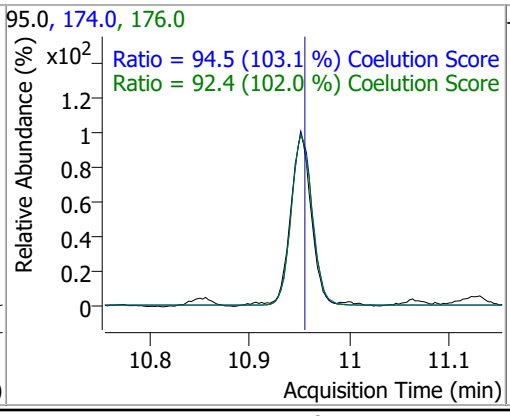


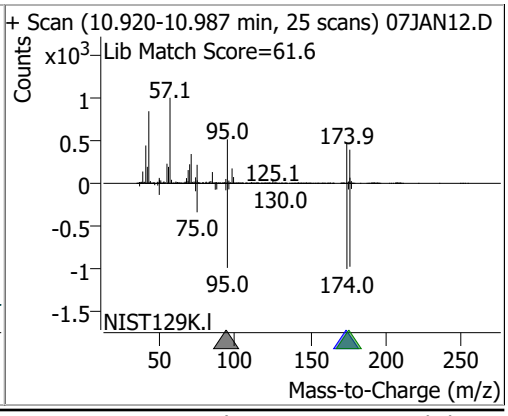




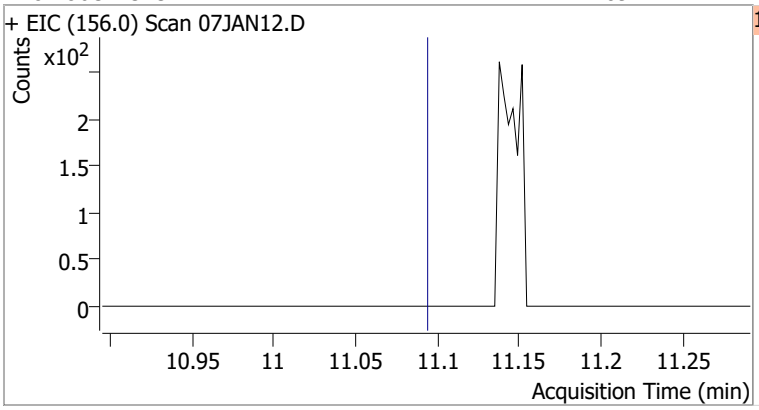
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	270.3807	10.95	0.00	258089	174.0	94.5	61.7	121.7
					176.0	92.4	60.6	120.6

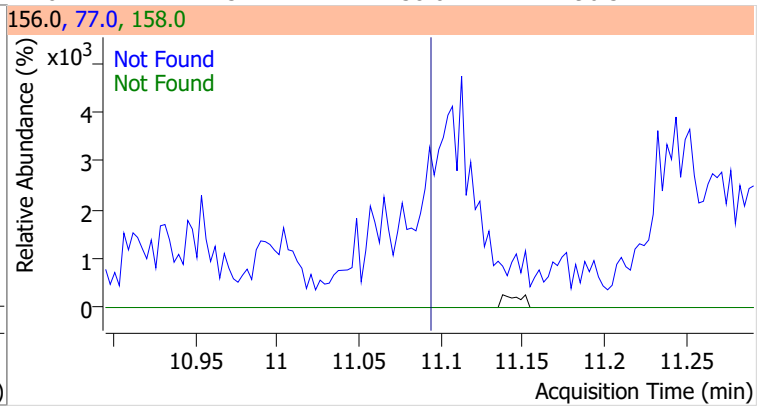




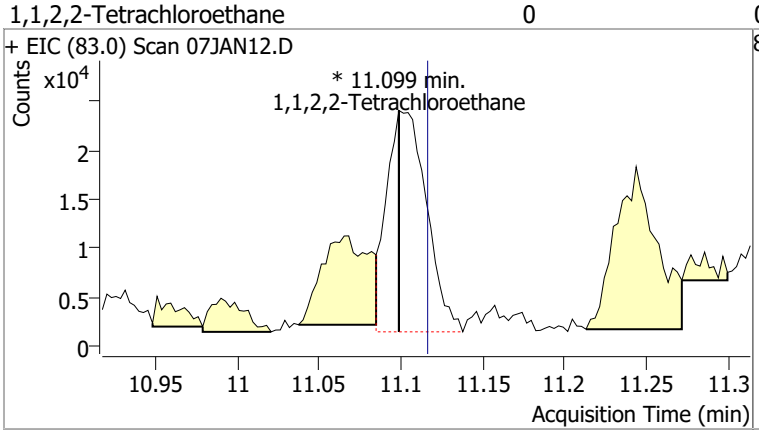


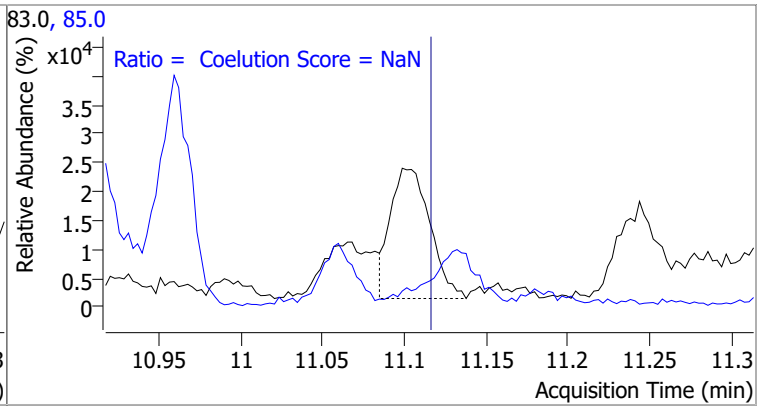
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



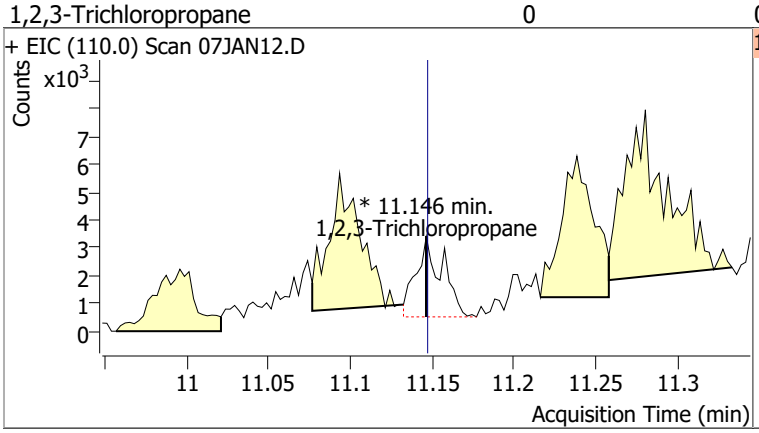


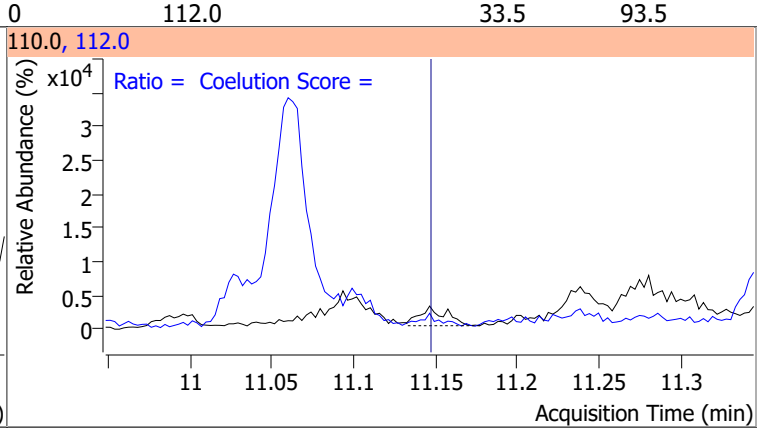
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane		0		0	85.0		36.2	96.2



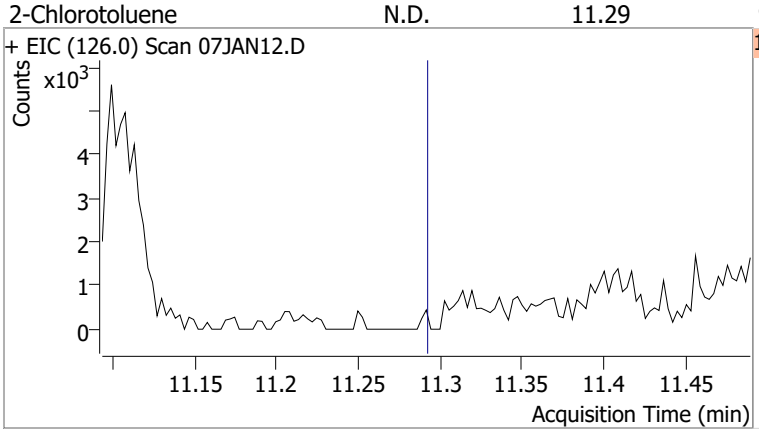


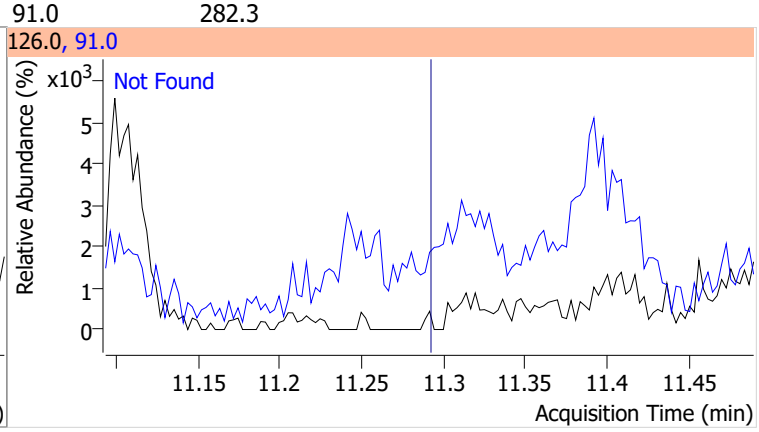
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane		0		0	112.0		33.5	93.5



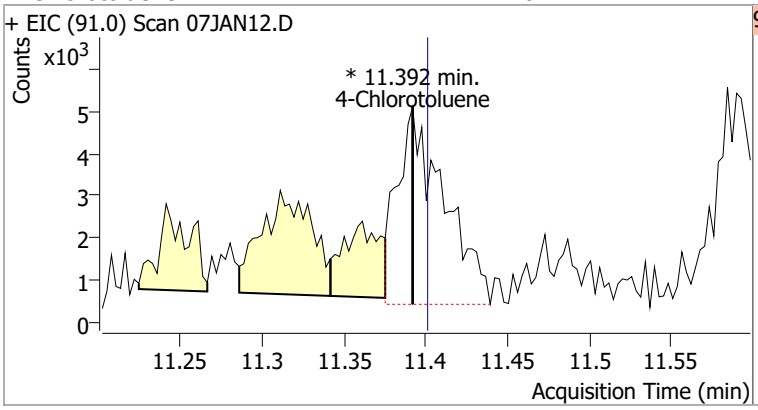


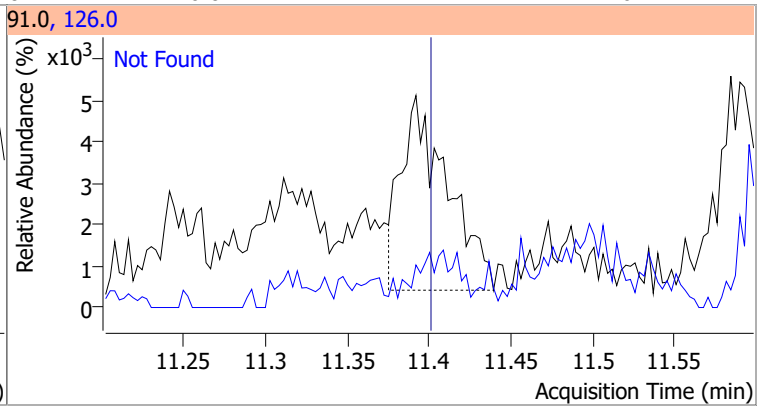
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



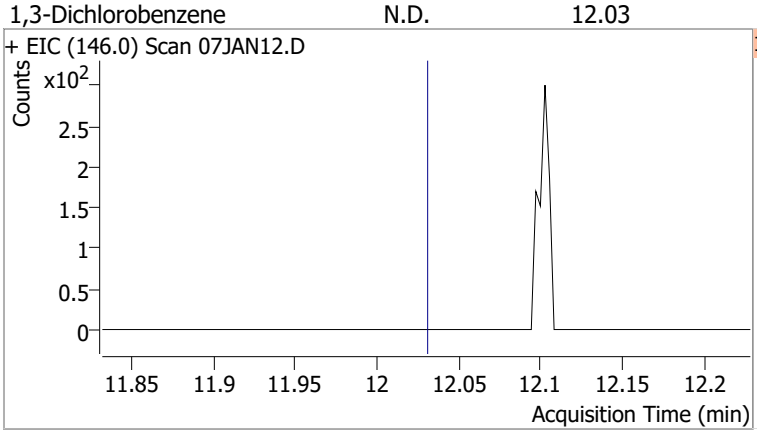


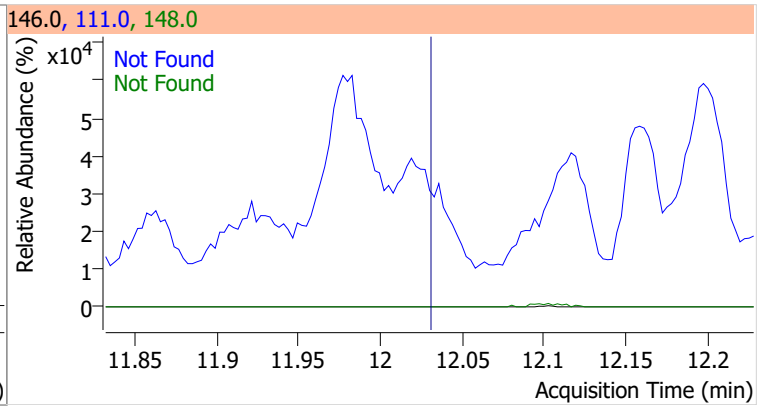
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene		0		0	126.0		1.7	61.7



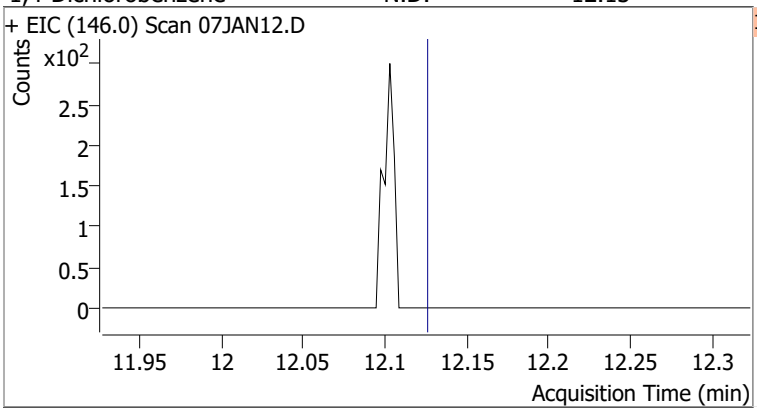


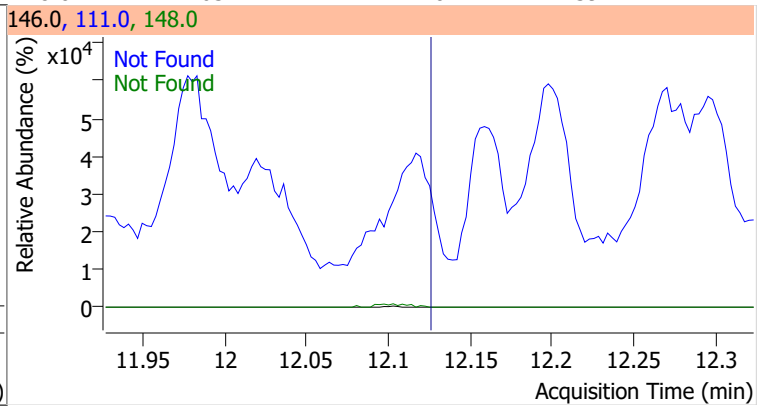
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



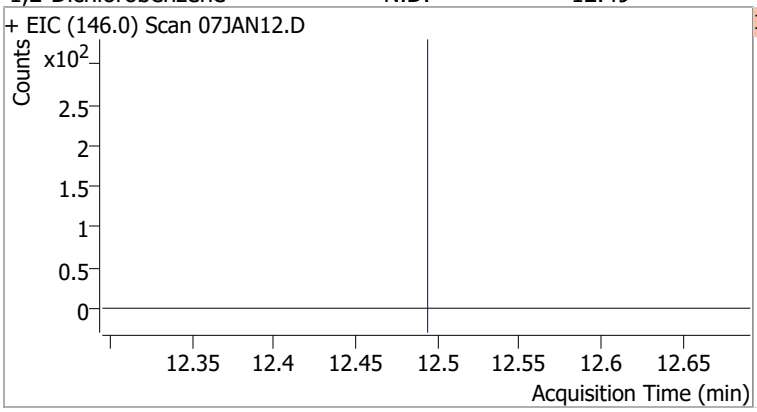


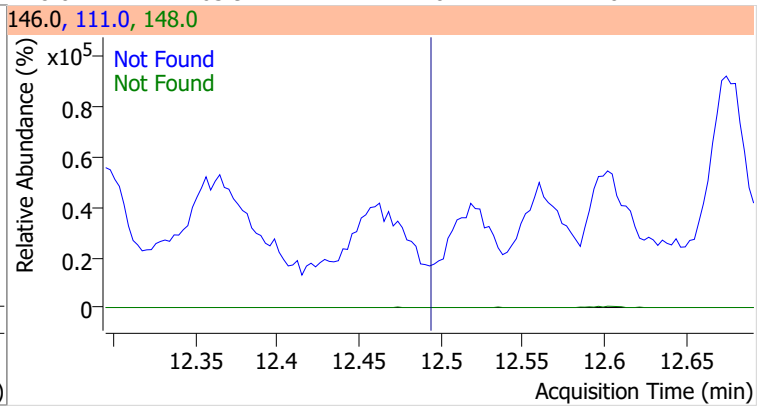
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





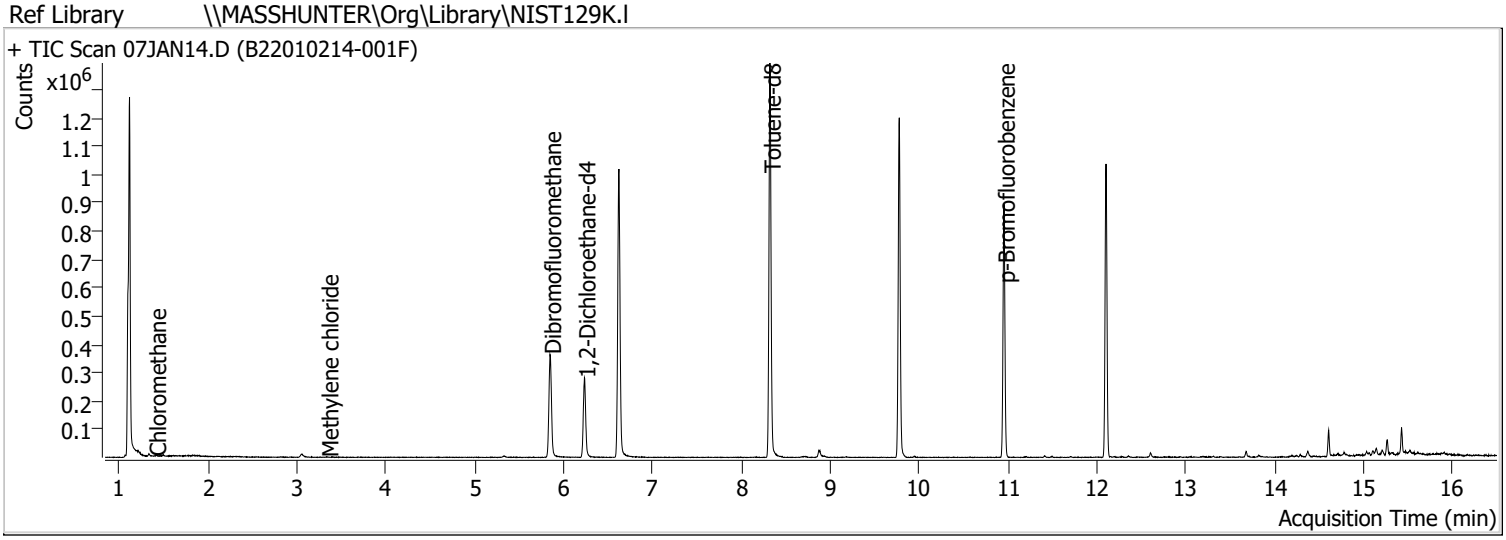
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN14.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 3:35:59 PM
Sample Name	B22010214-001F	Instrument	VOA5975C
Vial	14	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

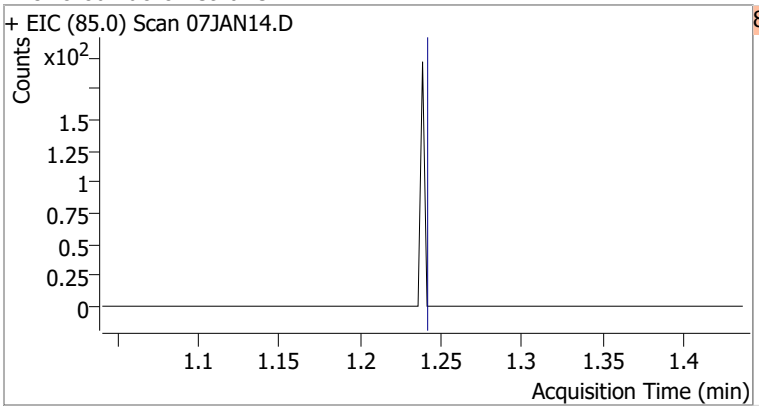


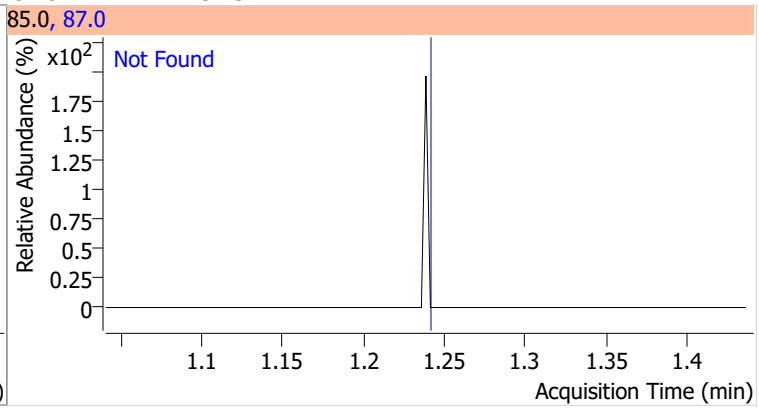
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.620	96.0	850985	250.0000	ng		-0.003
M Chlorobenzene-d5	9.772	82.0	331791	250.0000	ng		0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	251886	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.845	113.0	224189	279.6371	ng		0.000
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 111.85%			
S 1,2-Dichloroethane-d4	6.233	67.0	99414	287.0891	ng		0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 114.84%			
S Toluene-d8	8.319	98.0	856527	267.8901	ng		0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 107.16%			
S p-Bromofluorobenzene	10.951	95.0	255176	276.5273	ng		-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 110.61%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	0.000		0	N.D.			
T Chloromethane	1.406	50.0	2350	1.7362	ng		98
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.344	49.0	885	0.7000	ng	m	82
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	0.000		0	N.D.			

Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	0.000		0	N.D.			
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	0.000		0	N.D.			
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.377	92.0	0		ng	md	1
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	0.000		0	N.D.			
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	0.000		0	N.D.			
T m+p-Xylenes	0.000		0	N.D.			
T o-Xylene	0.000		0	N.D.			
T Styrene	0.000		0	N.D.			
T Bromoform	0.000		0	N.D.			
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

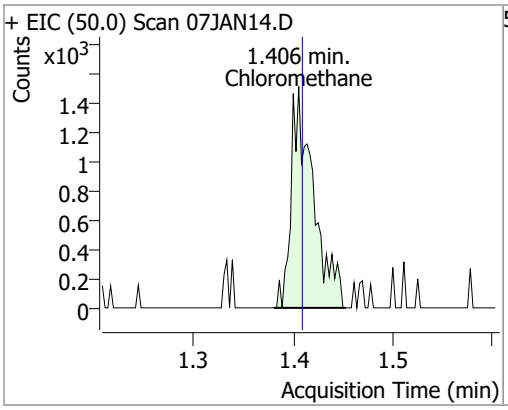
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

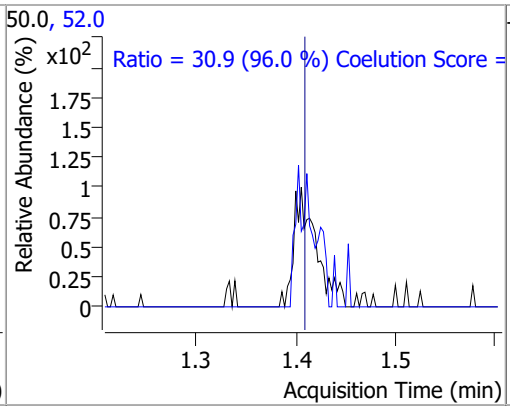
Compound	Conc.	Exp RT	QIon	Exp Ratio
Dichlorodifluoromethane	N.D.	1.24	87.0	32.3

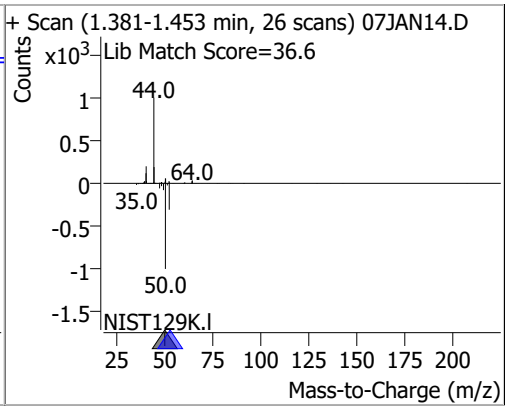




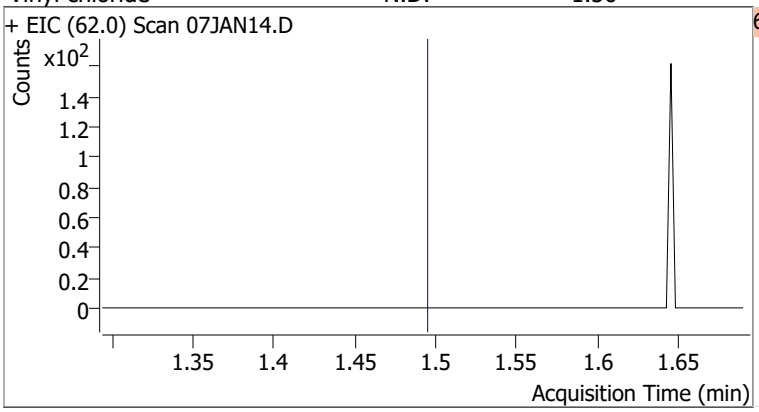
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	1.7362	1.41	0.00	2350	52.0	30.9	2.1	62.1

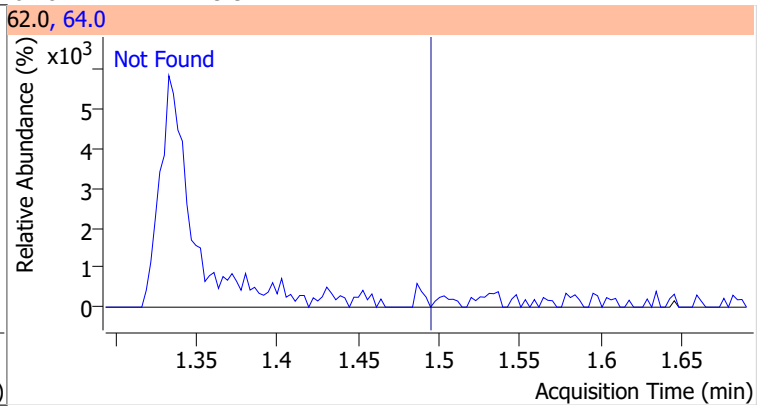




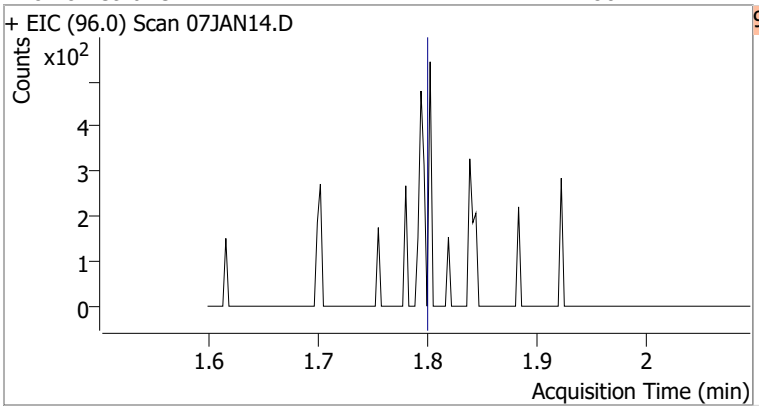


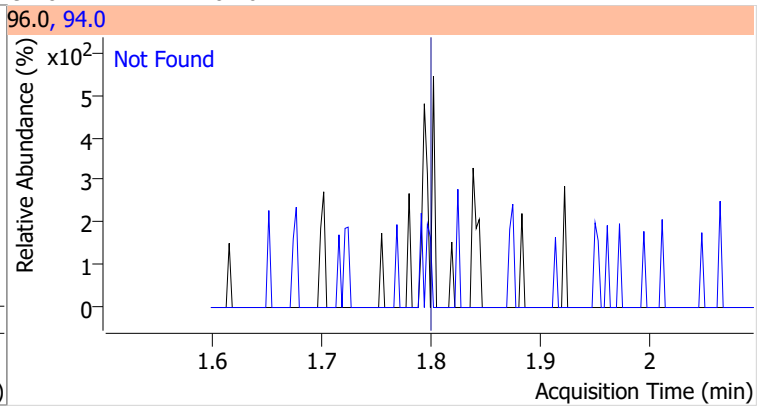
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



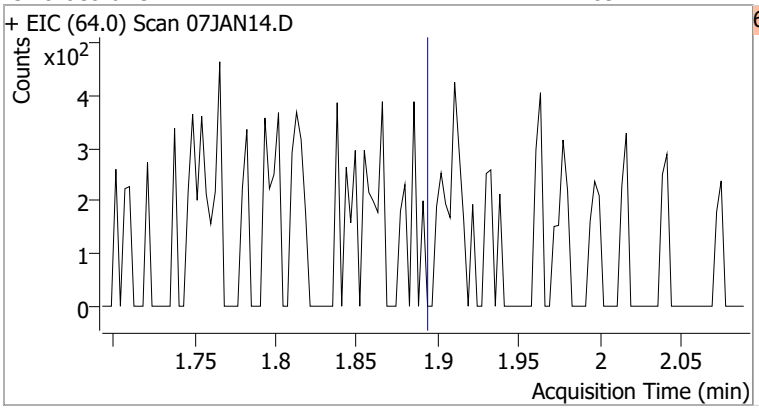


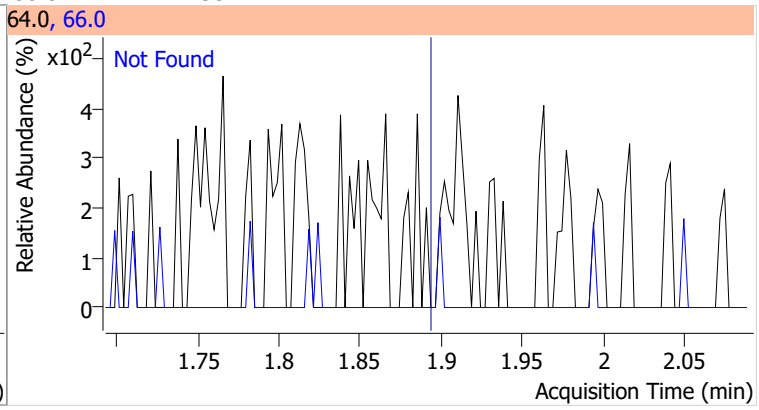
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



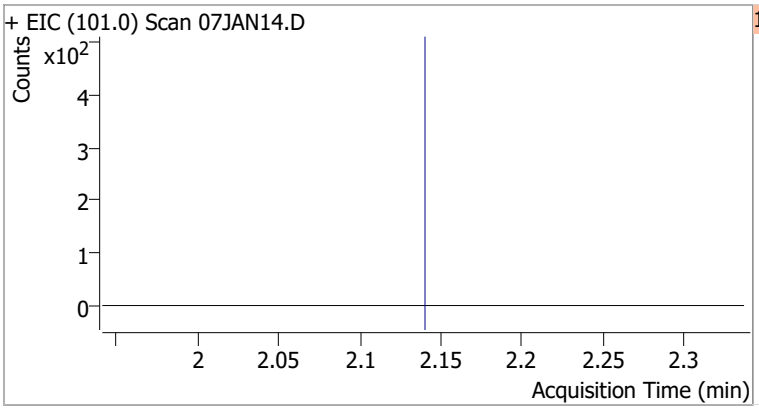


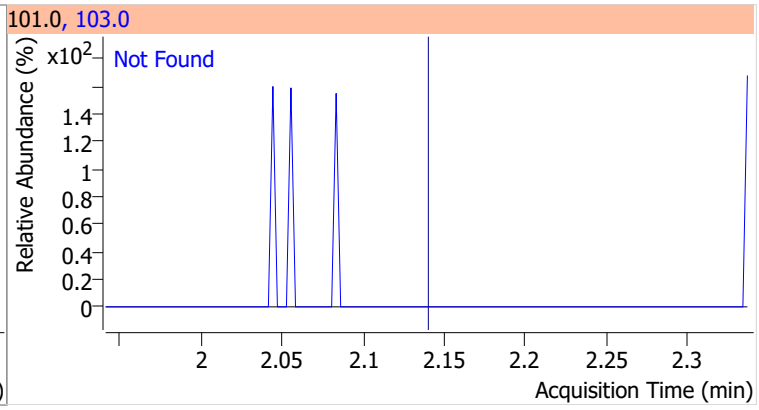
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



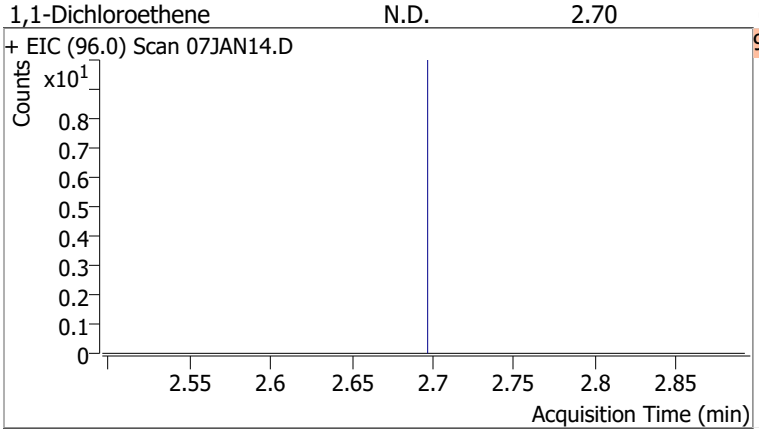


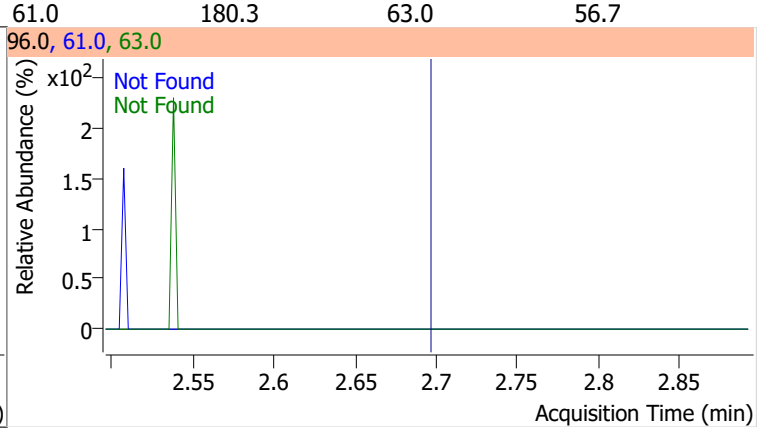
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



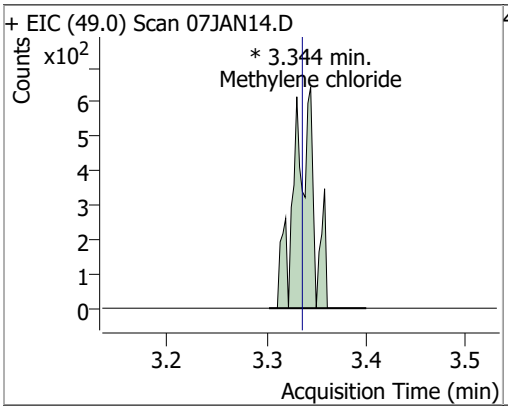


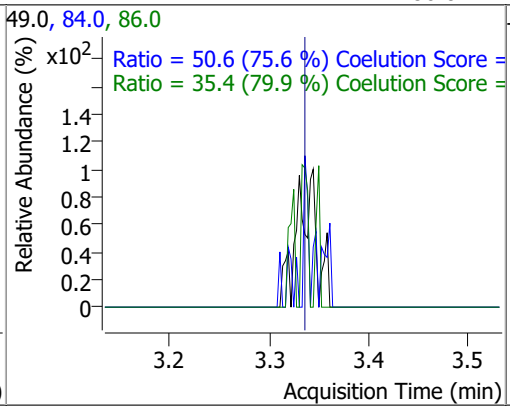
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

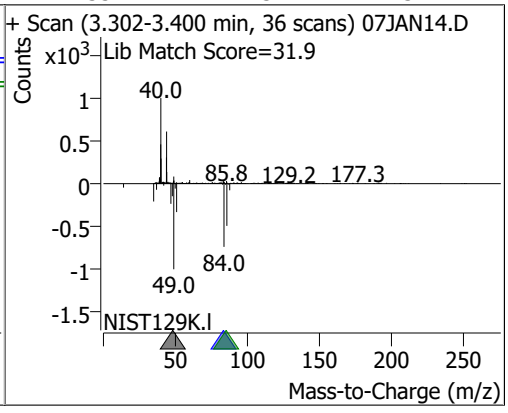




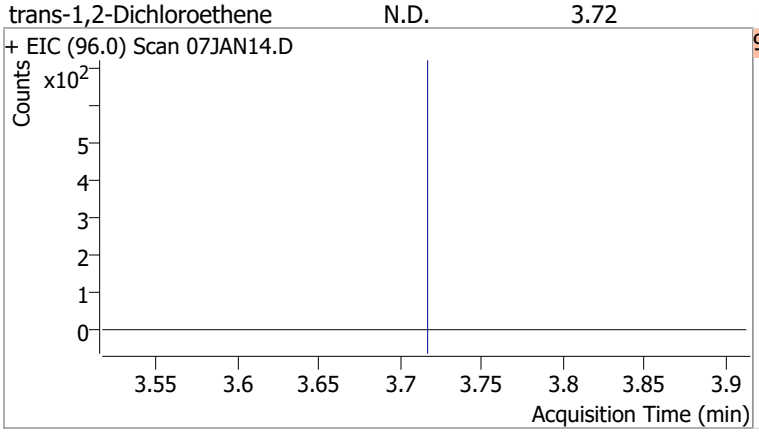
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	0.7000	3.34	0.01	885 (m)	84.0	50.6	36.9	96.9
					86.0	35.4	14.3	74.3

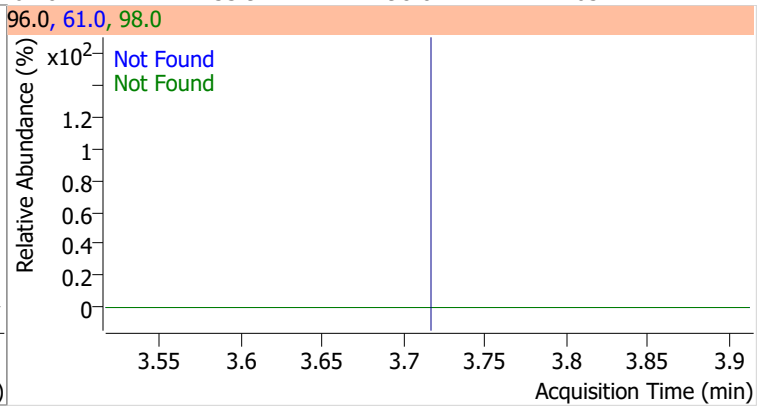




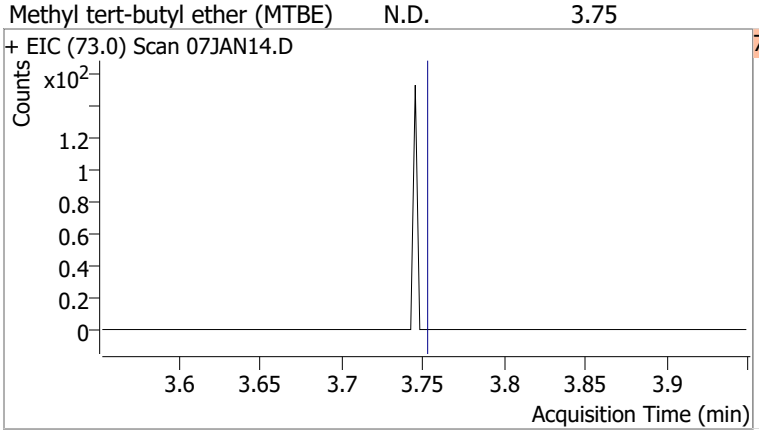


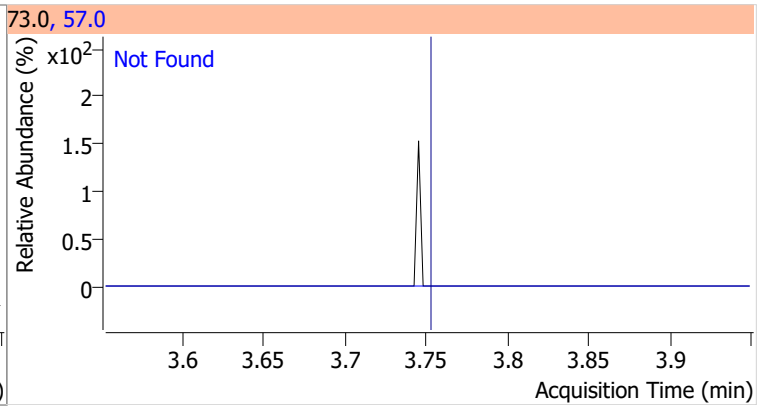
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



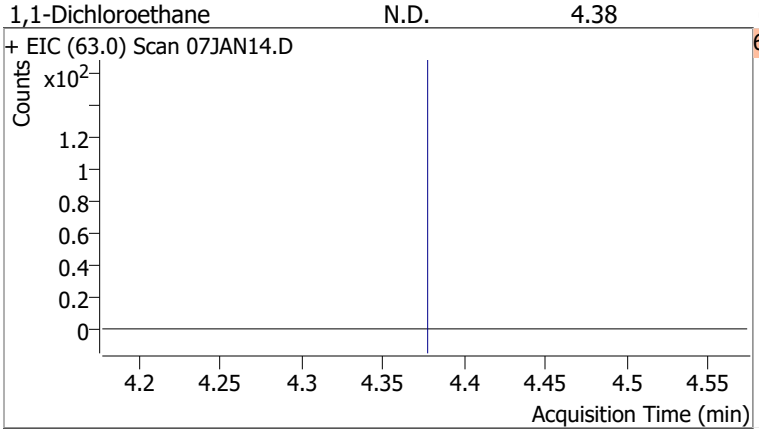


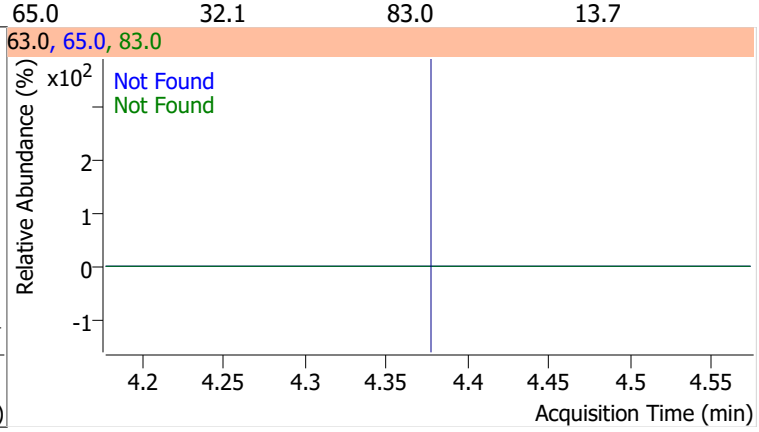
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



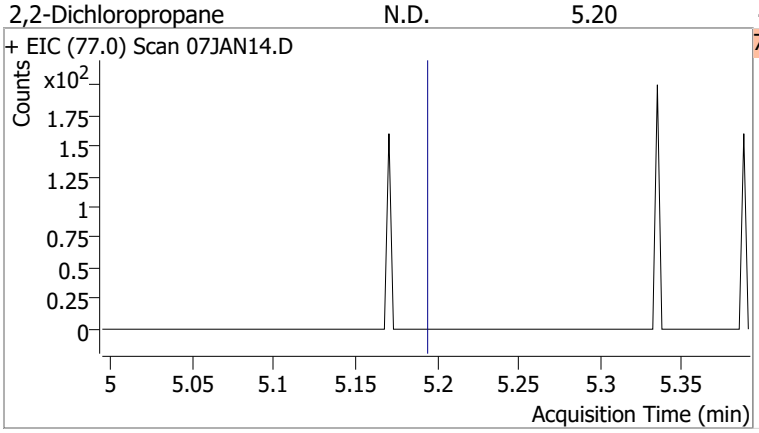


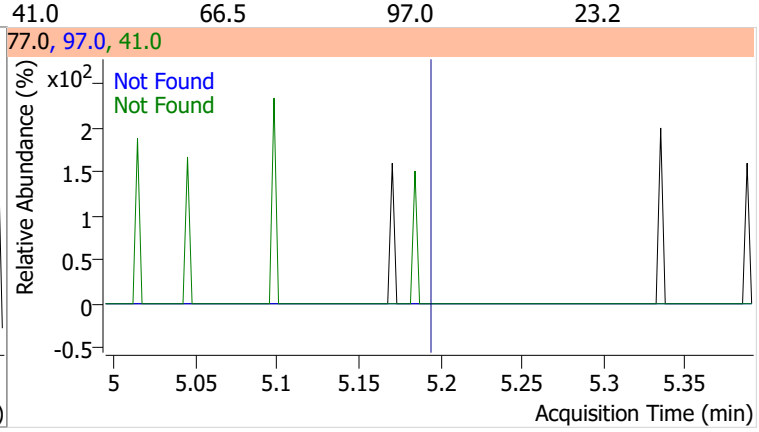
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



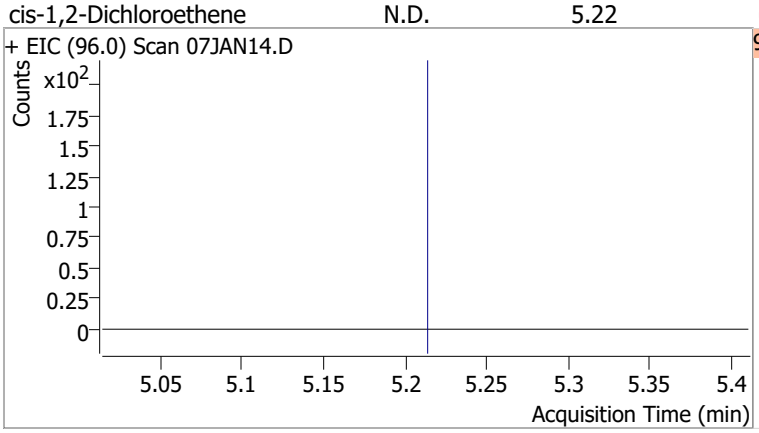


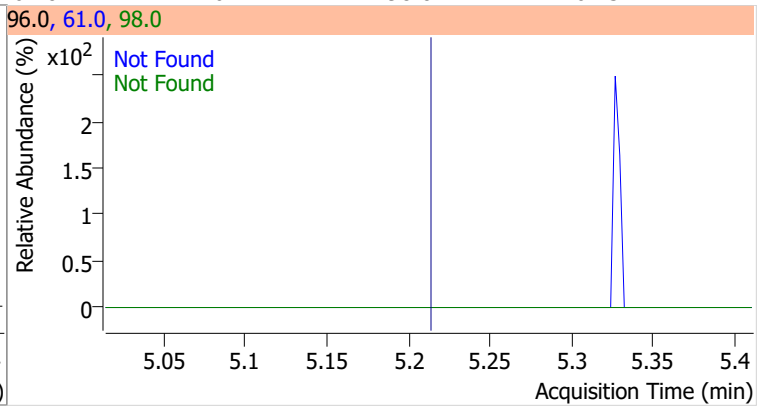
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



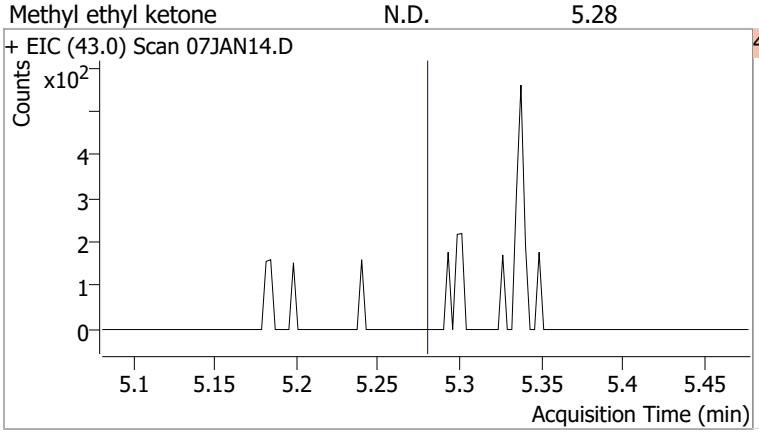


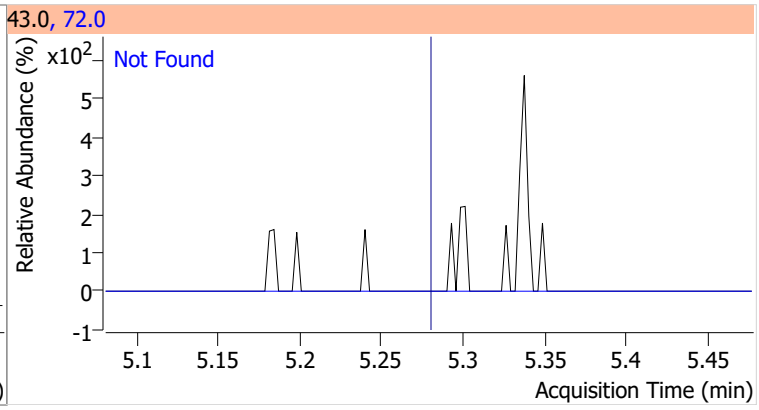
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



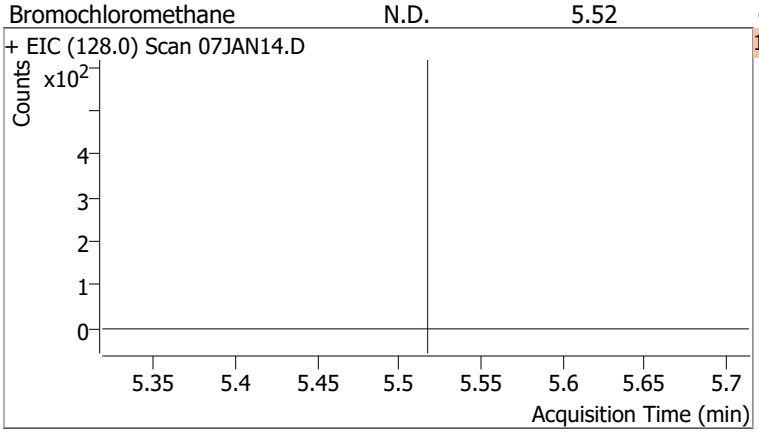


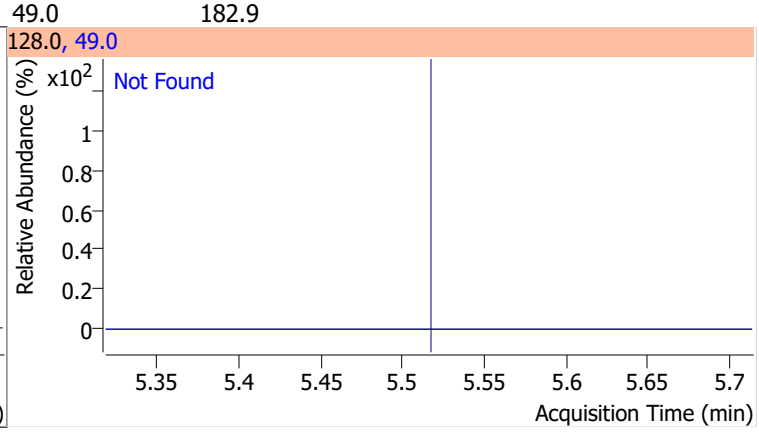
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



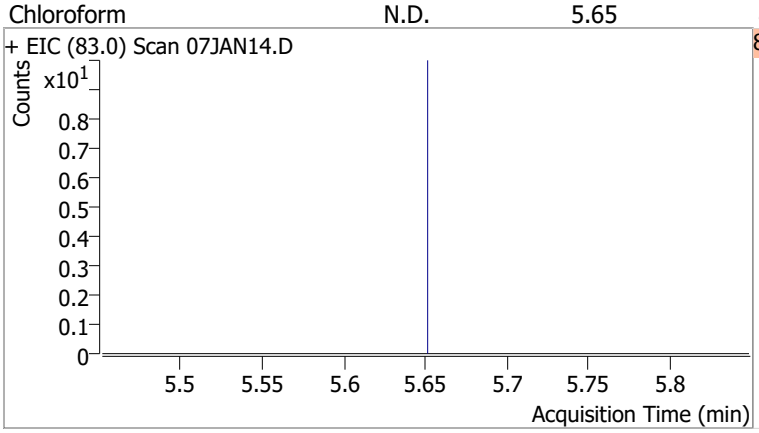


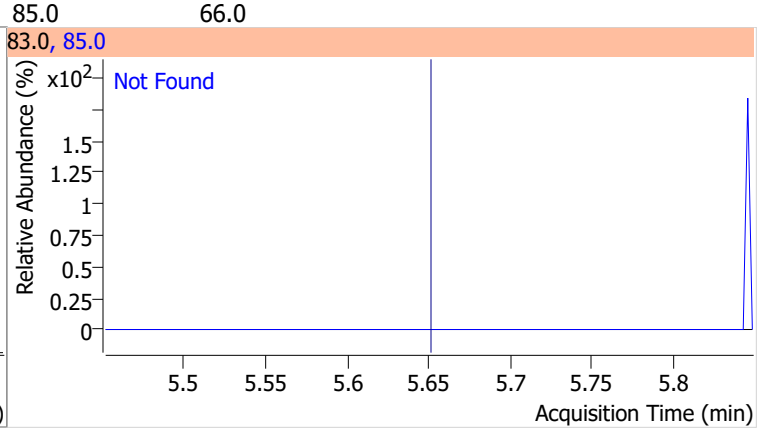
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9



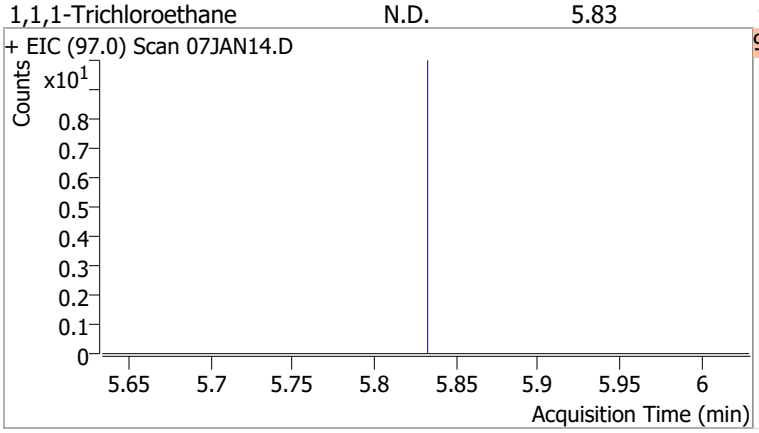


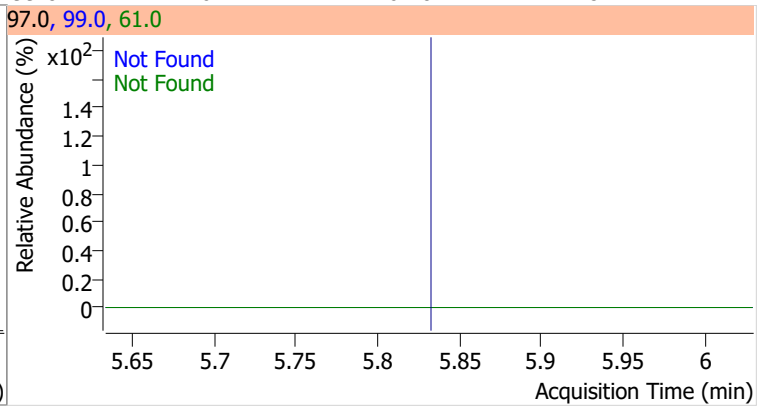
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroform	N.D.	5.65	85.0	66.0



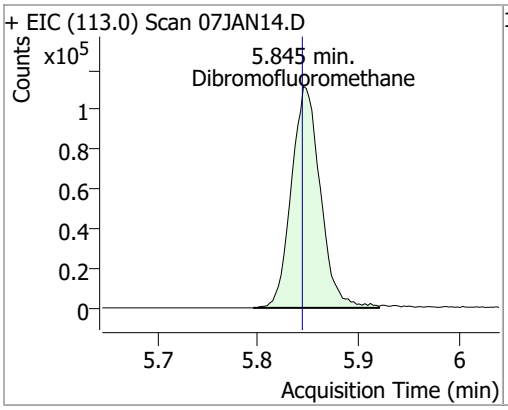


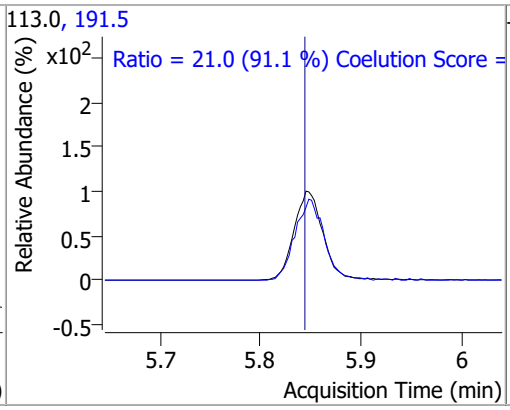
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

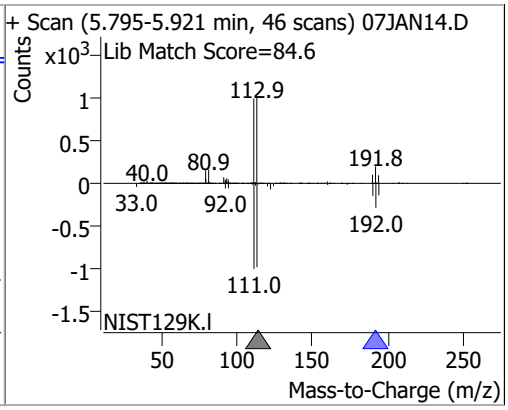




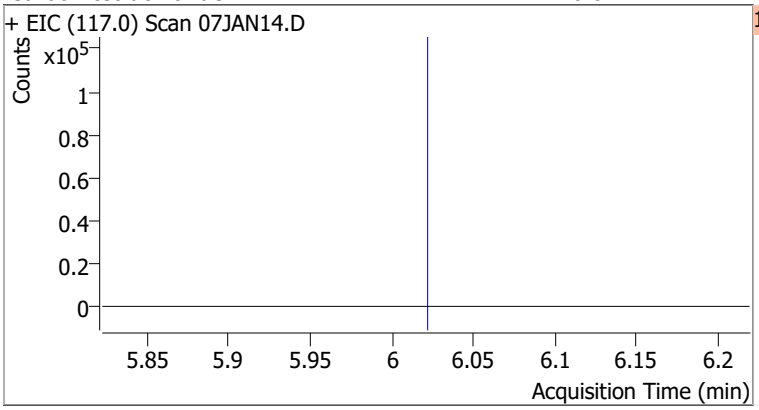
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	279.6371	5.85	0.00	224189	191.5	21.0	0.0	53.1

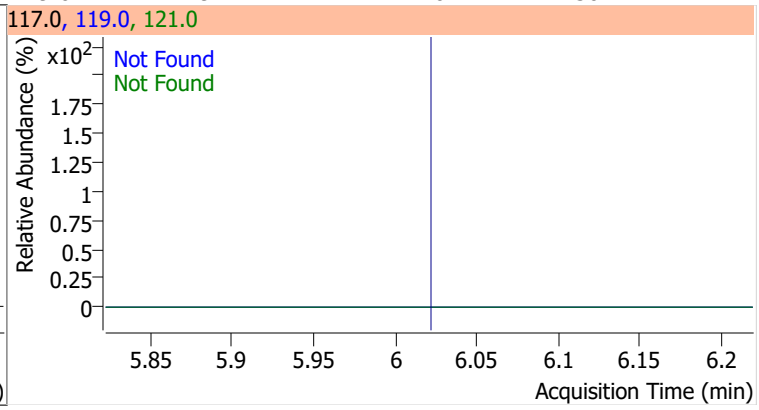




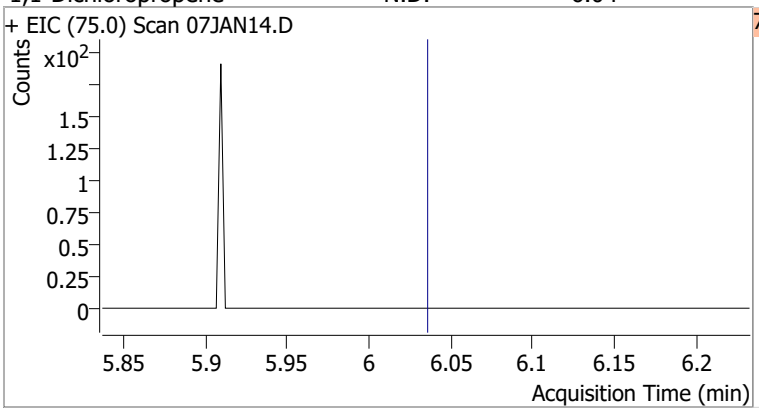


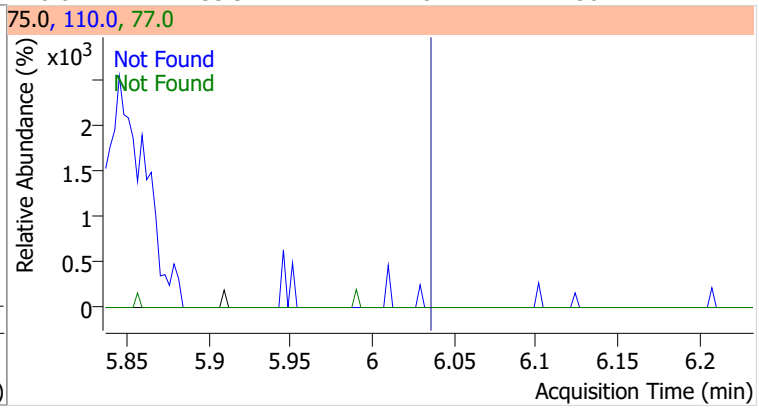
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



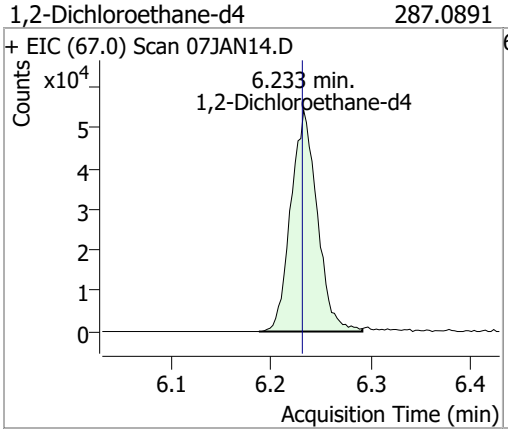


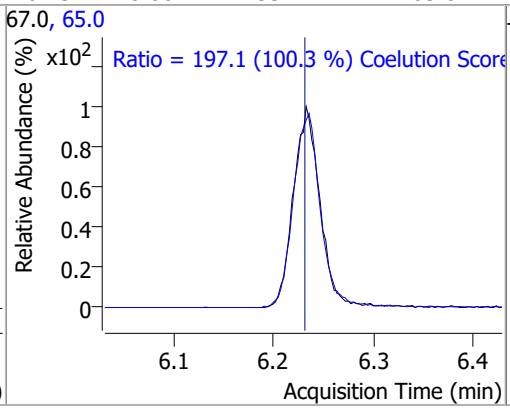
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

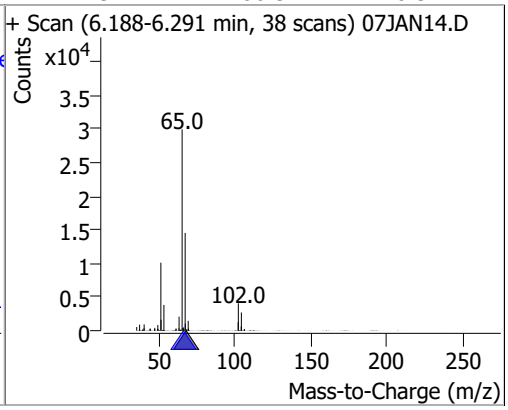




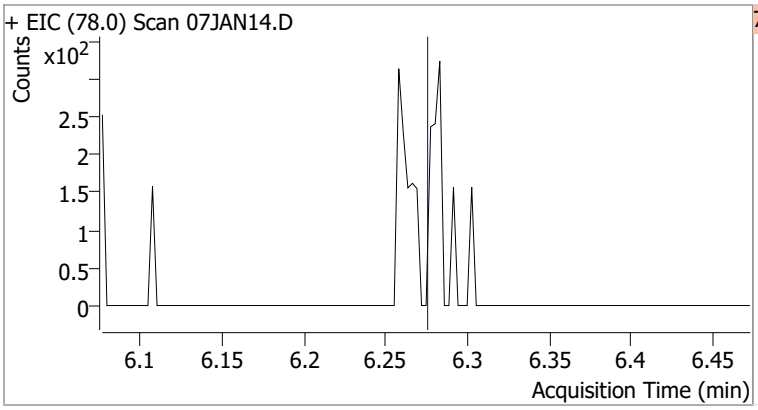
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	287.0891	6.23	0.00	99414	65.0	197.1	166.5	226.5

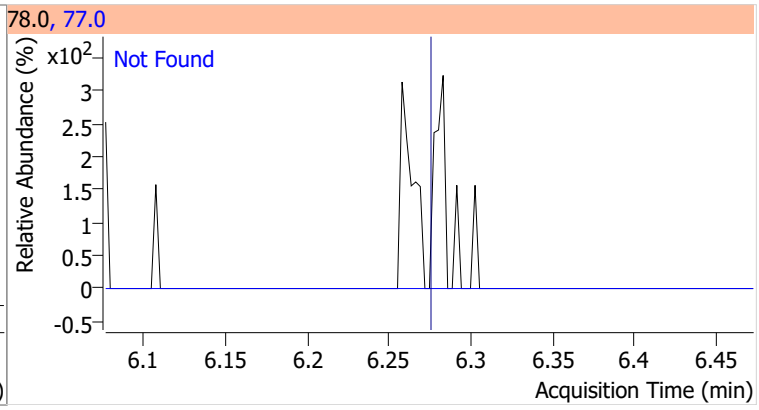




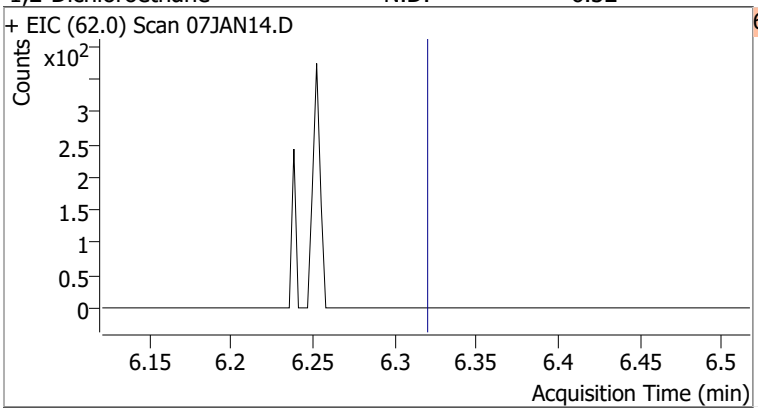


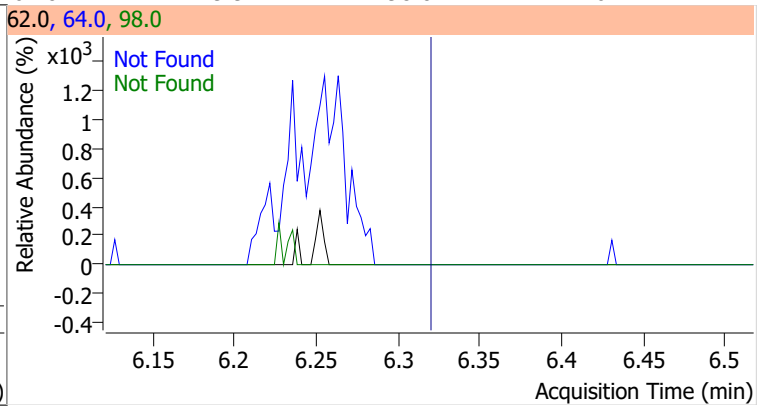
Compound	Conc.	Exp RT	QIon	Exp Ratio
Benzene	N.D.	6.28	77.0	23.5



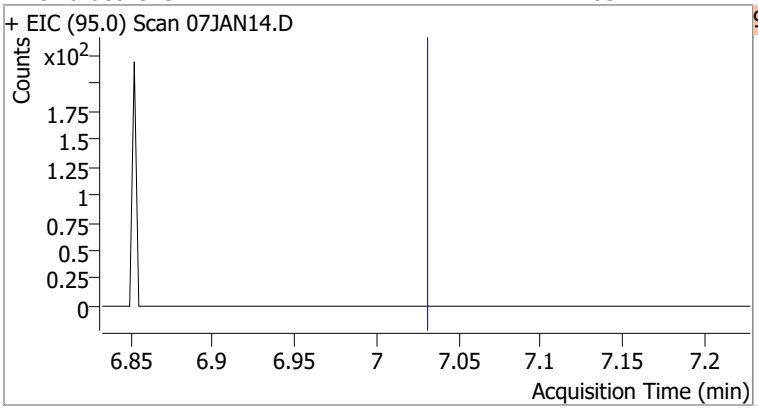


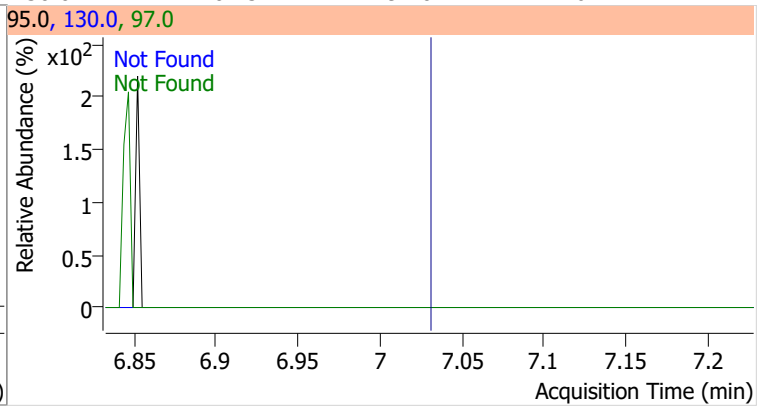
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



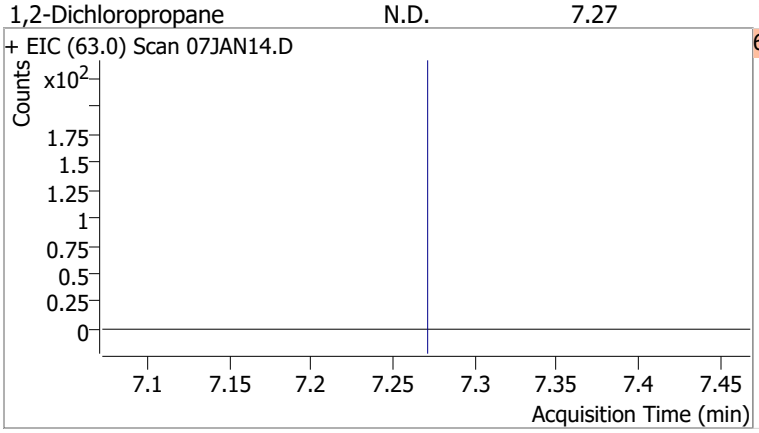


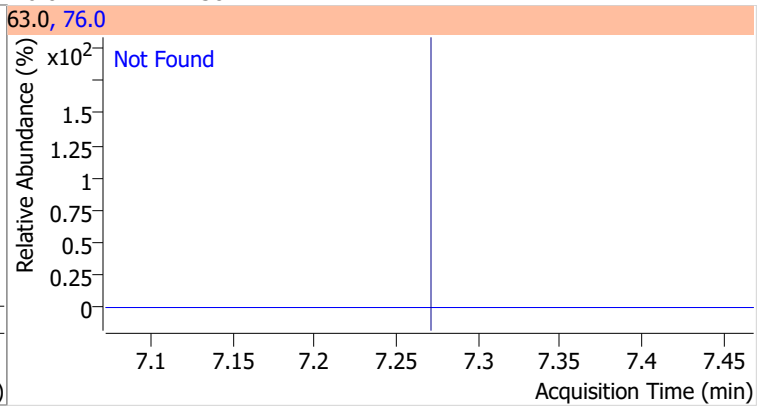
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



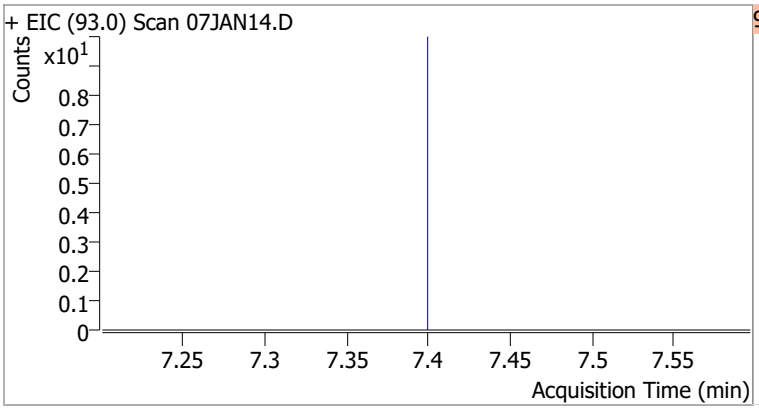


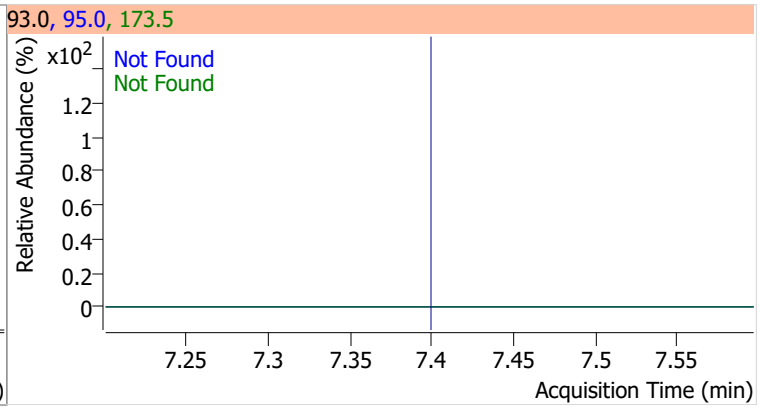
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



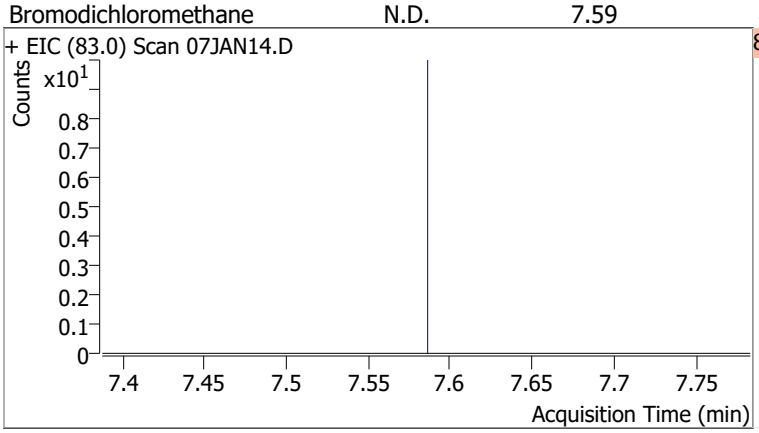


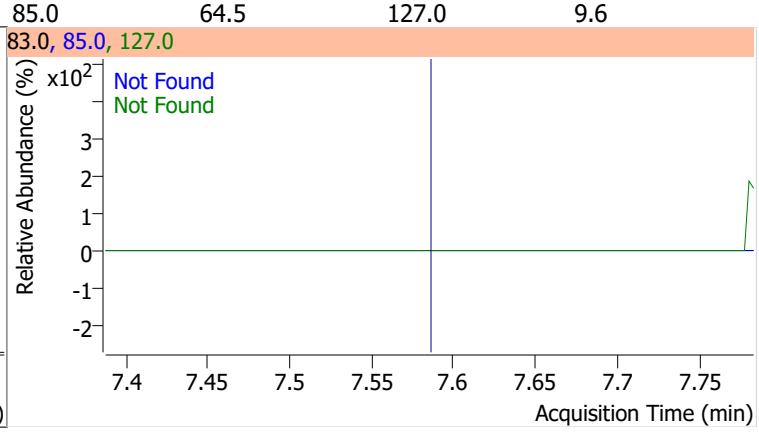
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



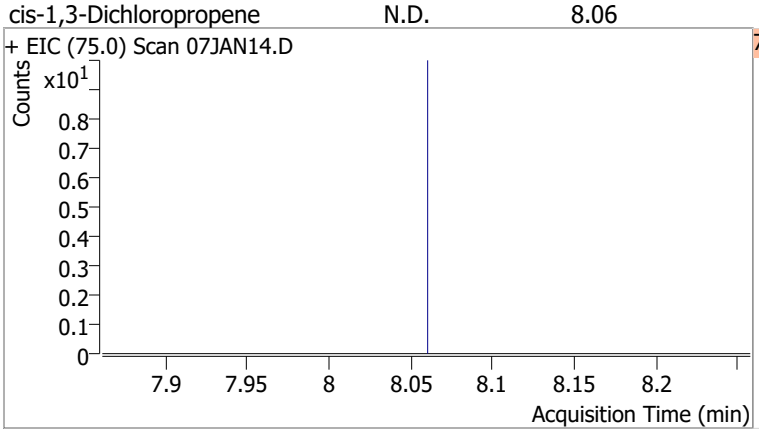


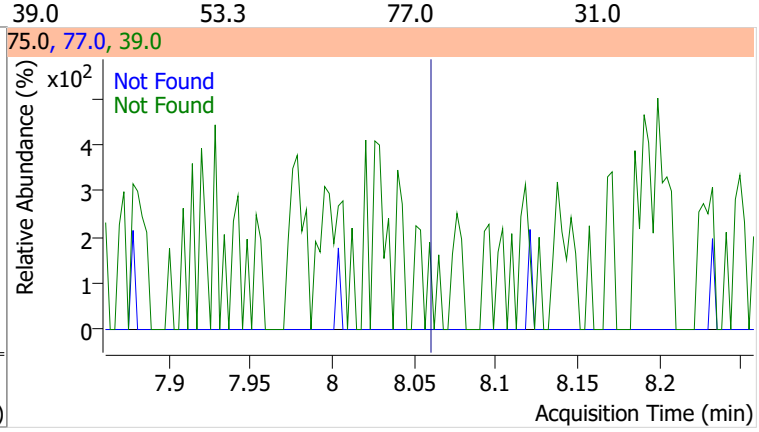
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



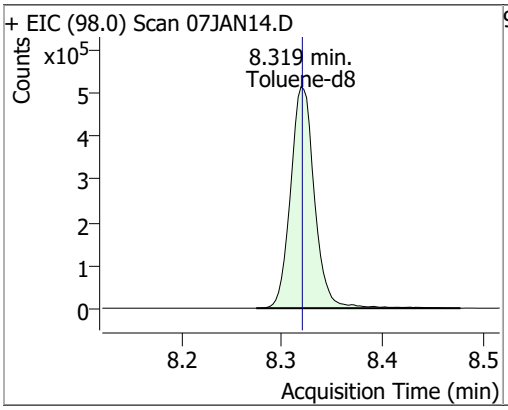


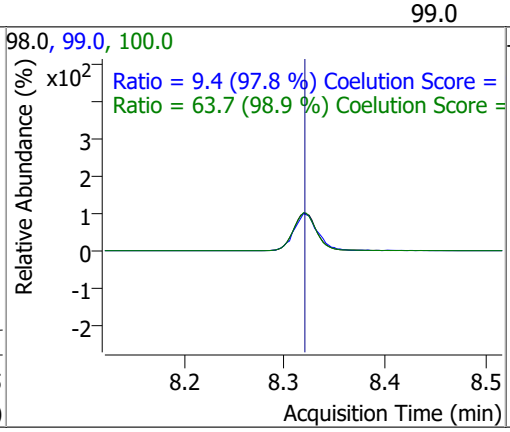
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

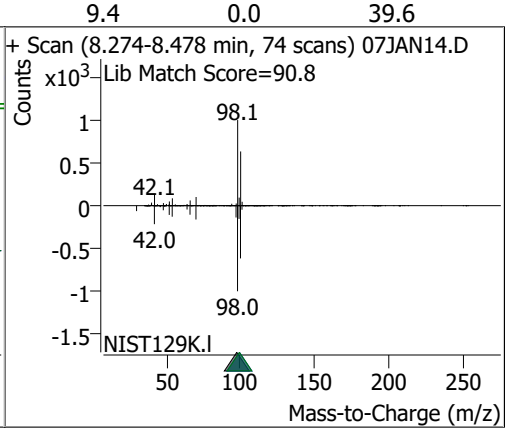




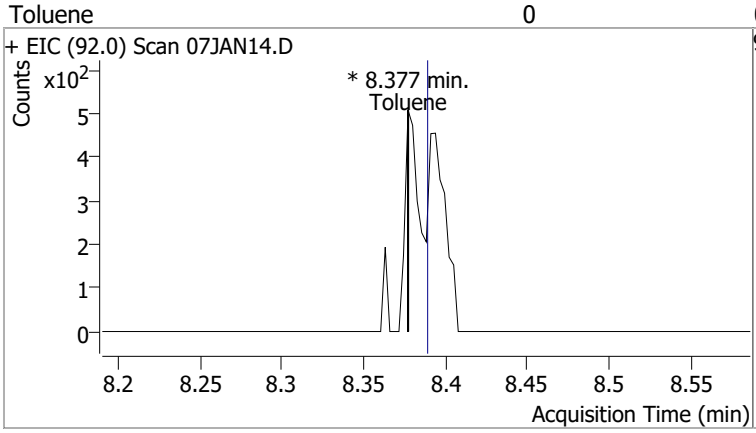
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	267.8901	8.32	0.00	856527	100.0	63.7	34.4	94.4
					99.0	9.4	0.0	39.6

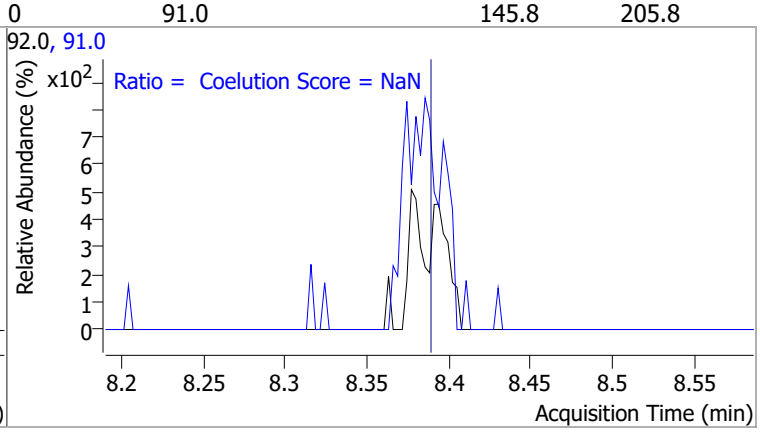




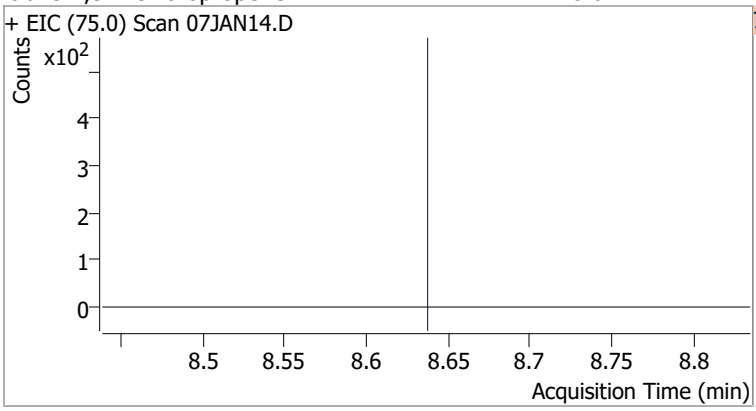


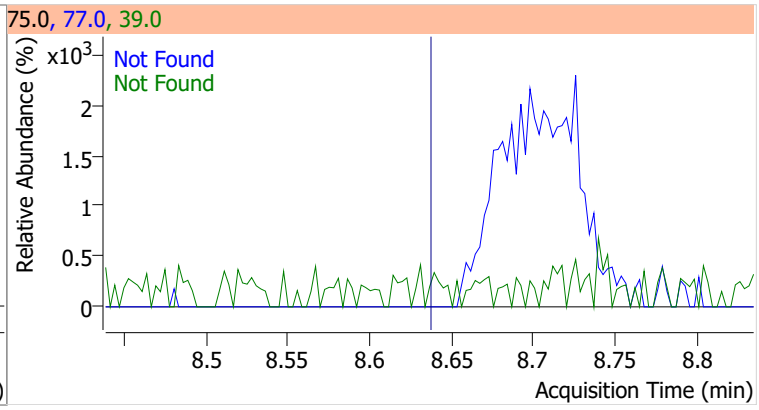
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene		0		0	91.0		145.8	205.8



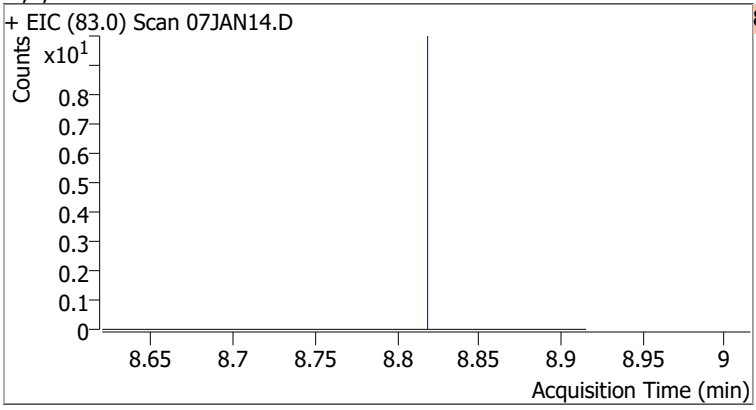


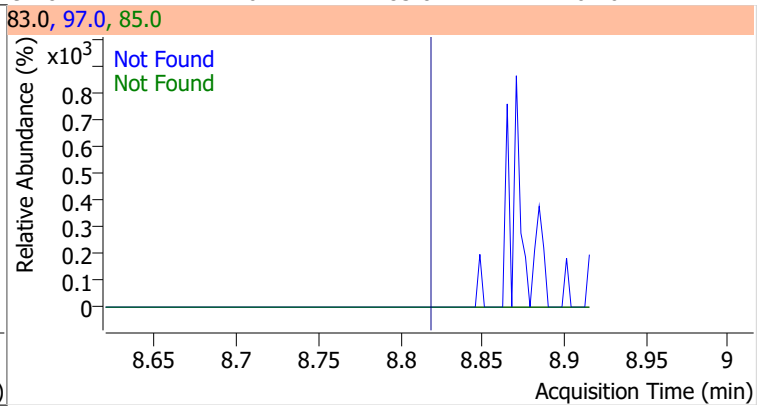
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



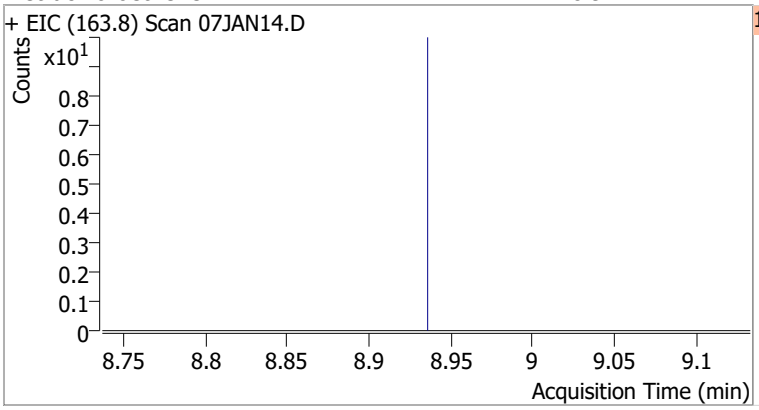


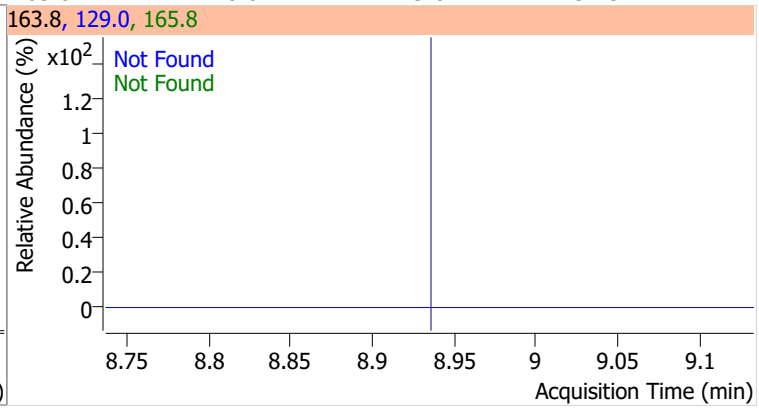
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



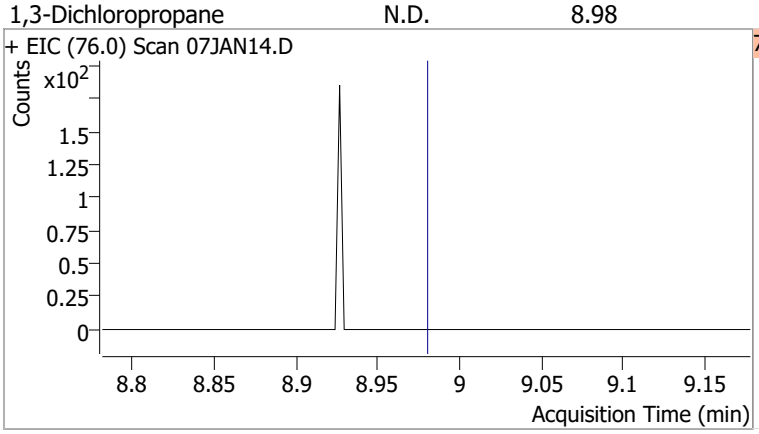


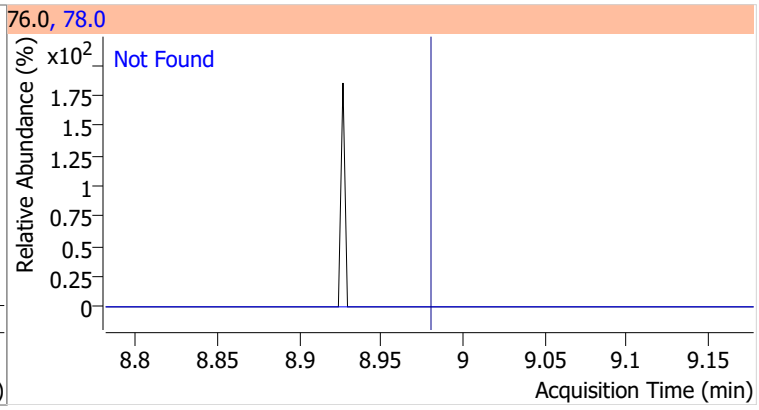
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



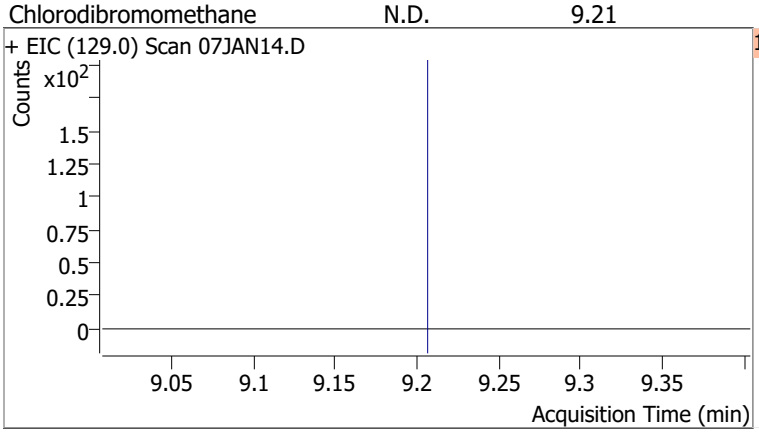


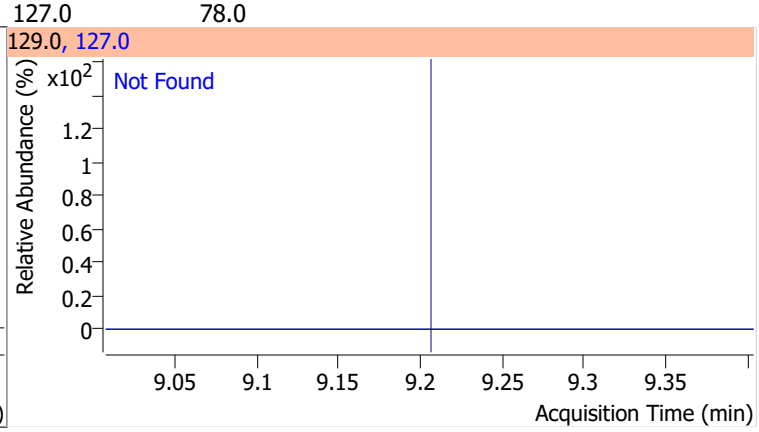
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



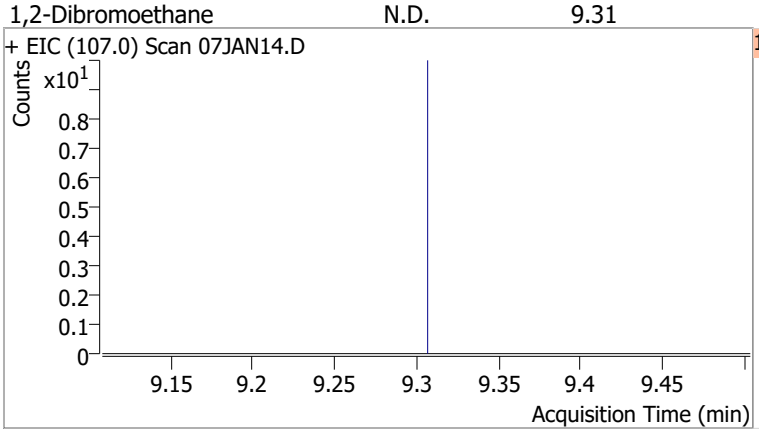


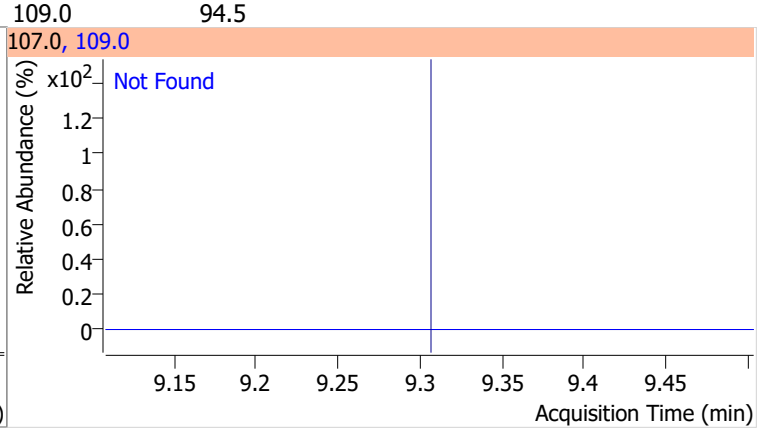
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



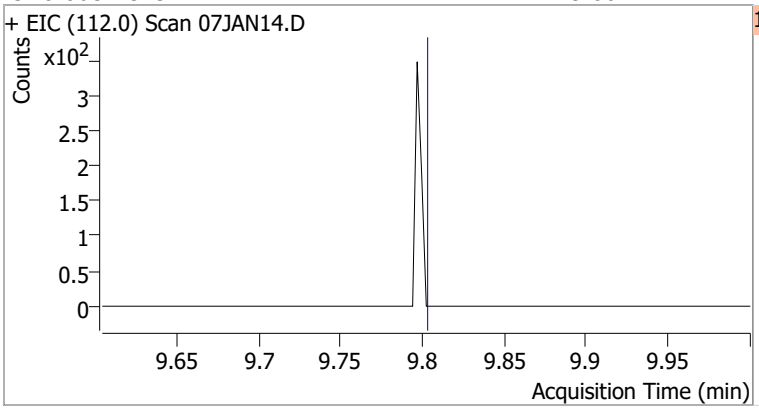


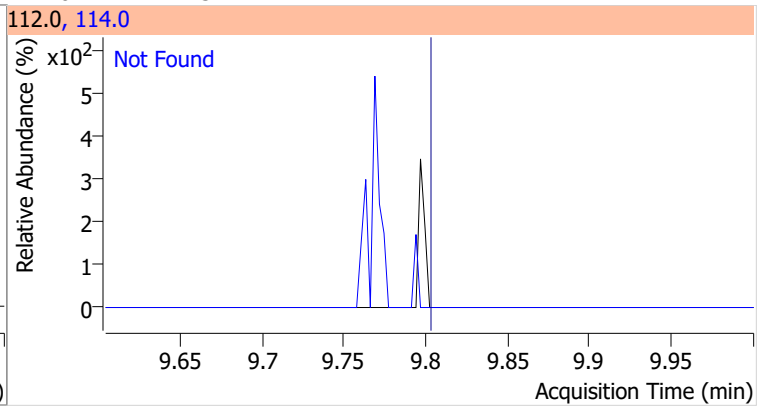
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



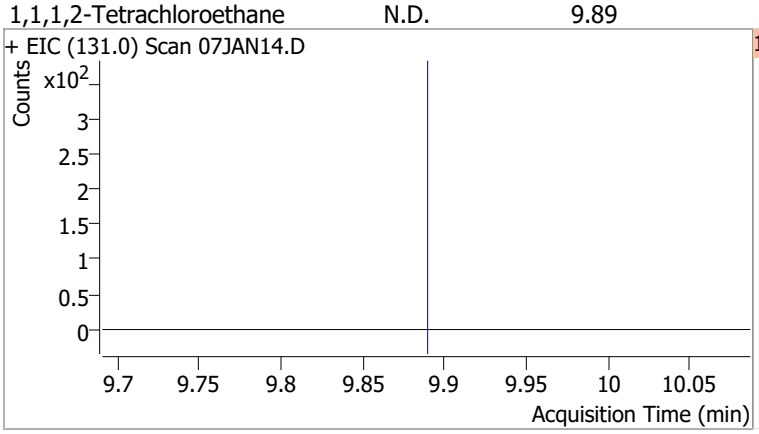


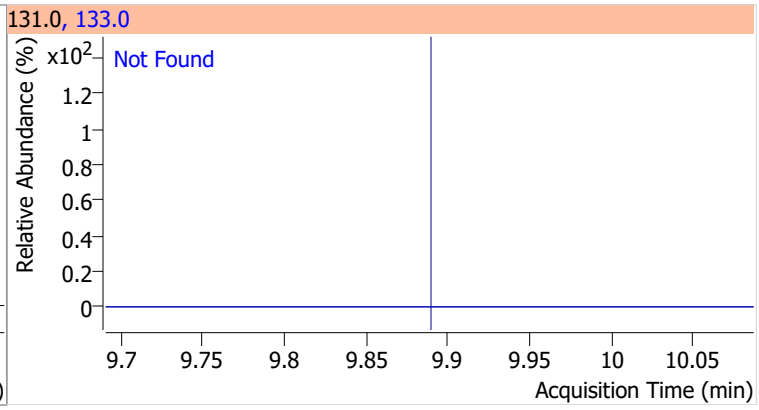
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



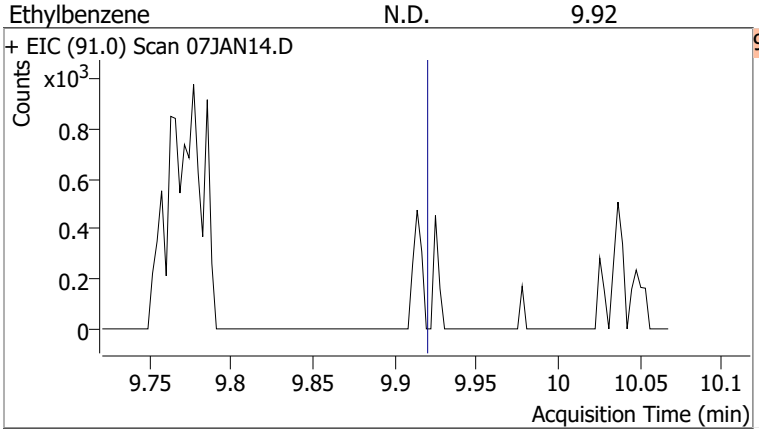


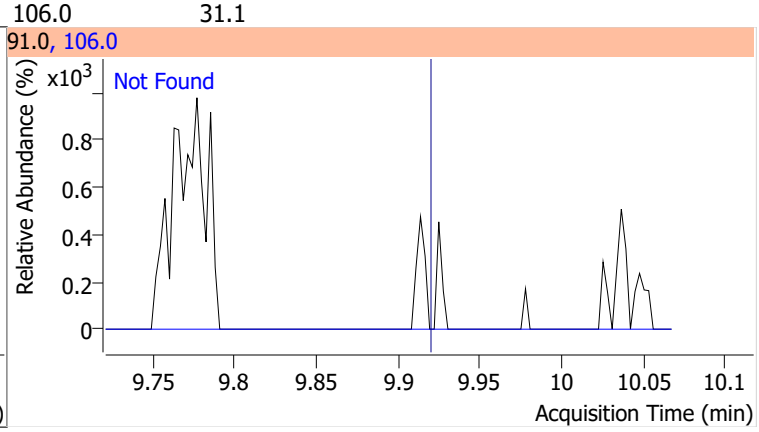
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



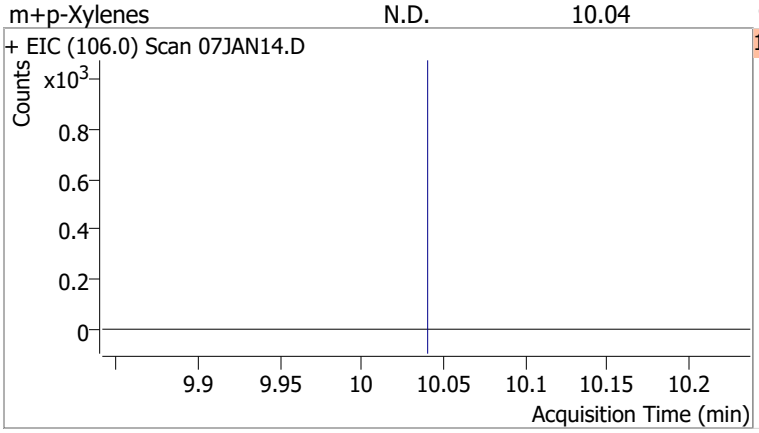


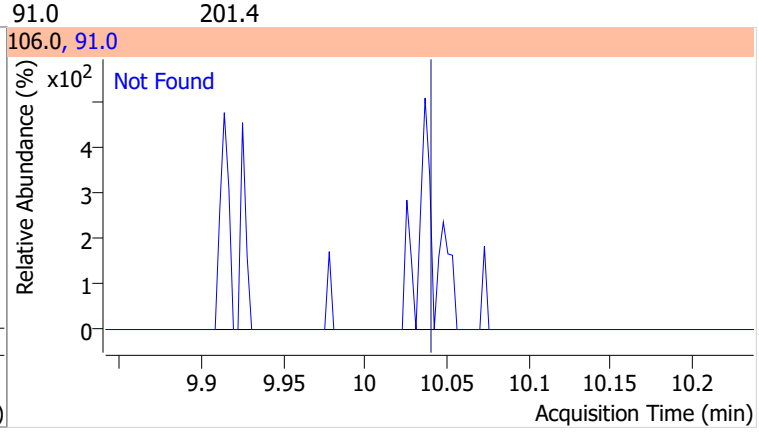
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1



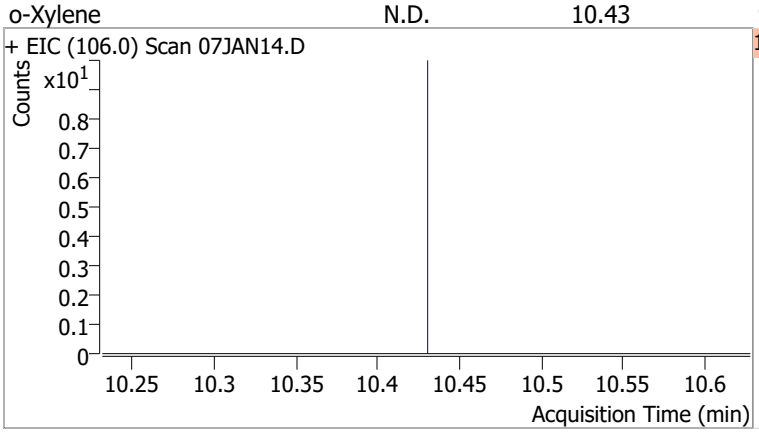


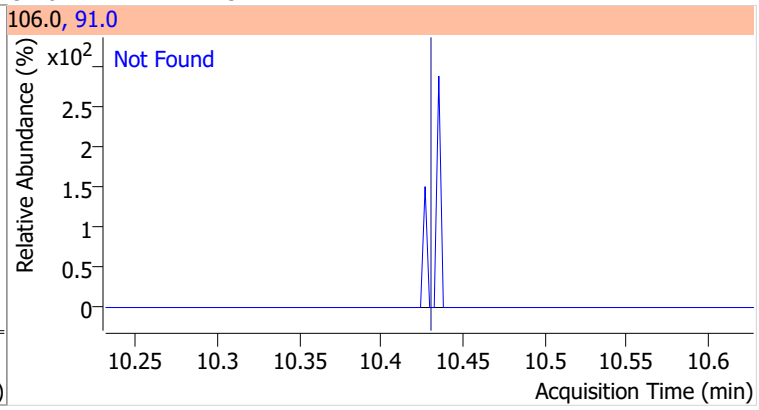
Compound	Conc.	Exp RT	QIon	Exp Ratio
m+p-Xylenes	N.D.	10.04	91.0	201.4



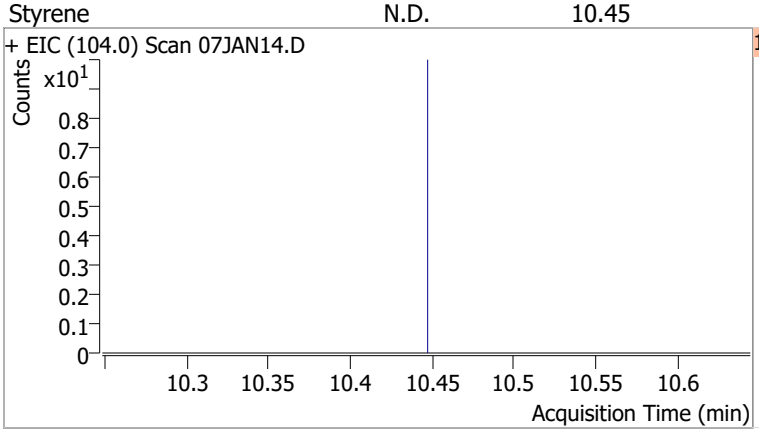


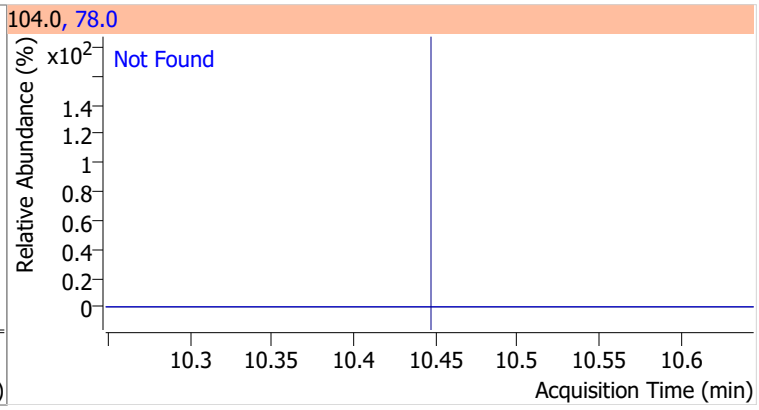
Compound	Conc.	Exp RT	QIon	Exp Ratio
o-Xylene	N.D.	10.43	91.0	213.1



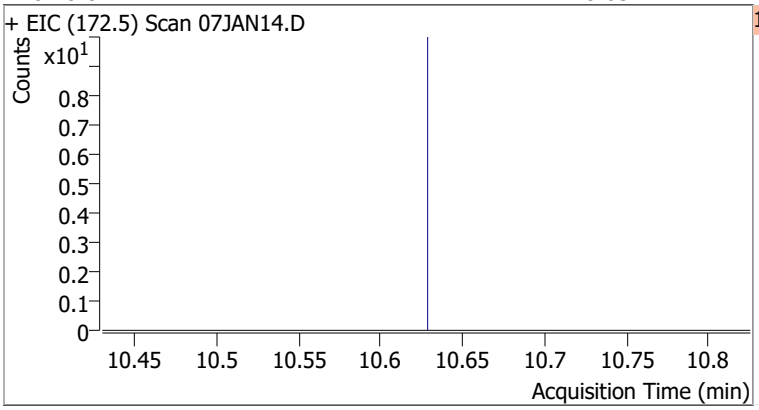


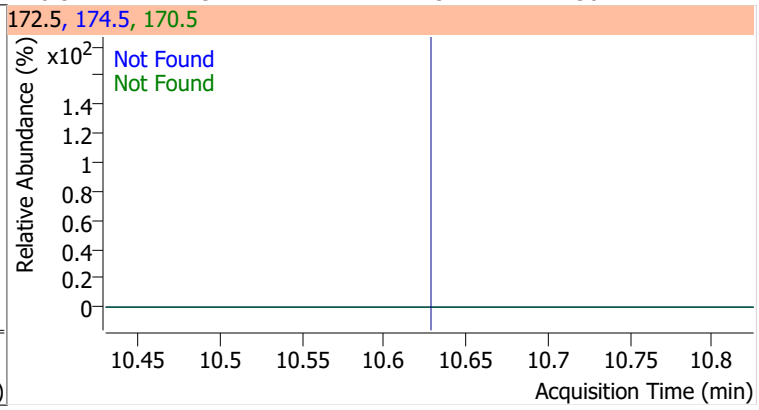
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



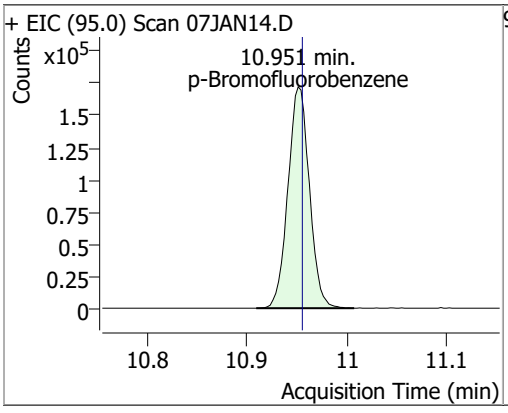


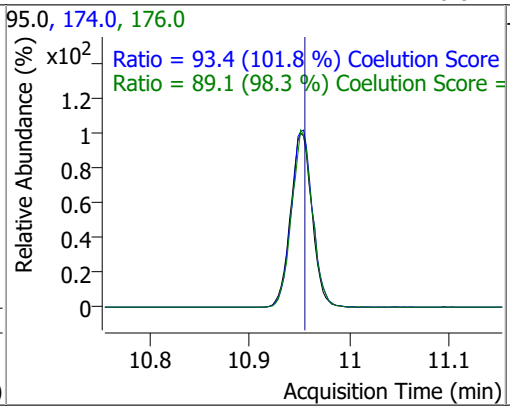
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

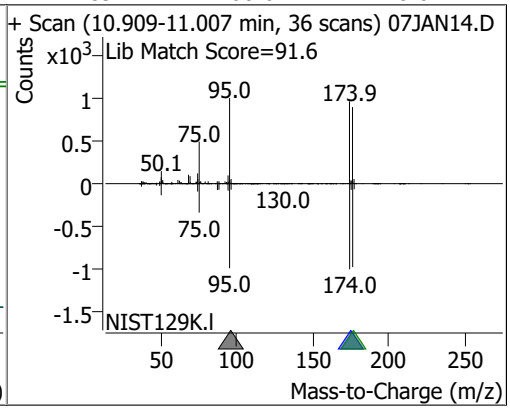




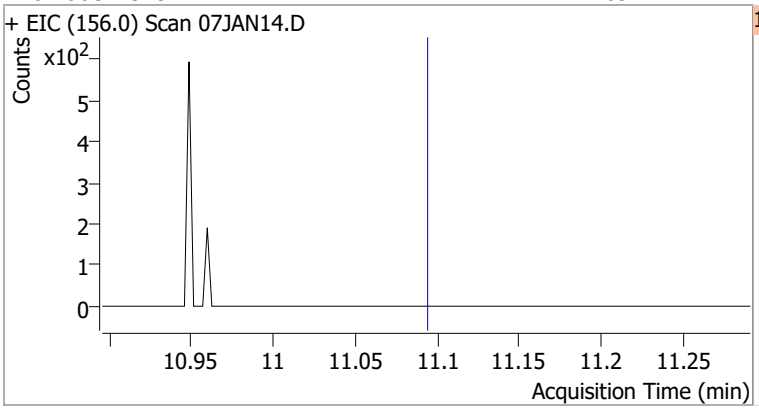
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	276.5273	10.95	0.00	255176	174.0	93.4	61.7	121.7
					176.0	89.1	60.6	120.6

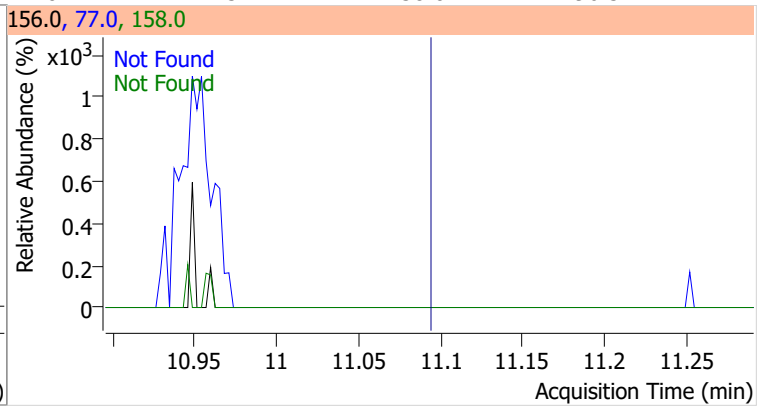




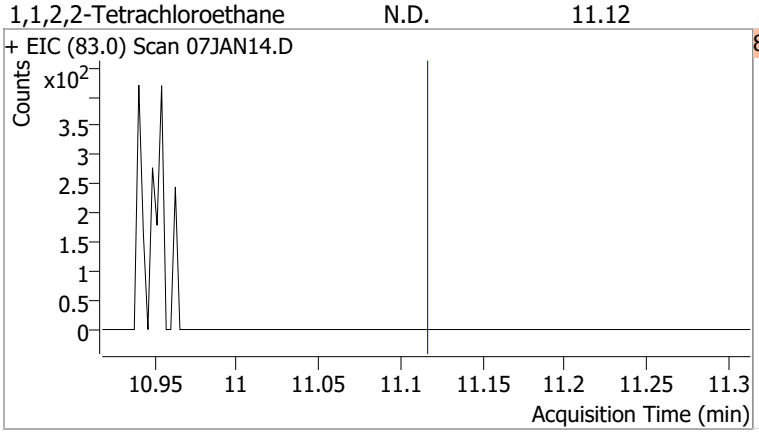


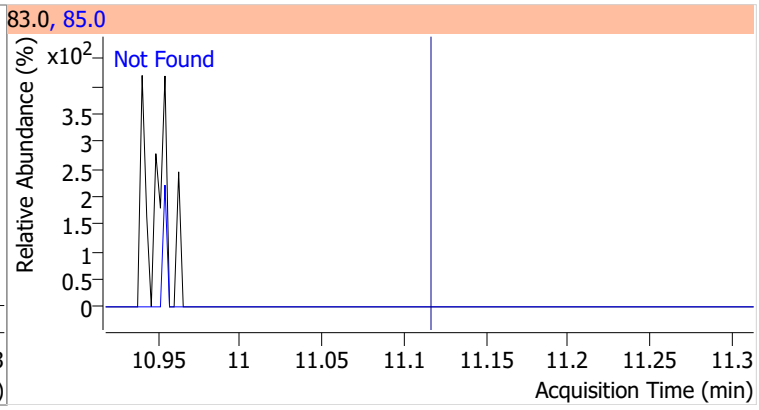
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



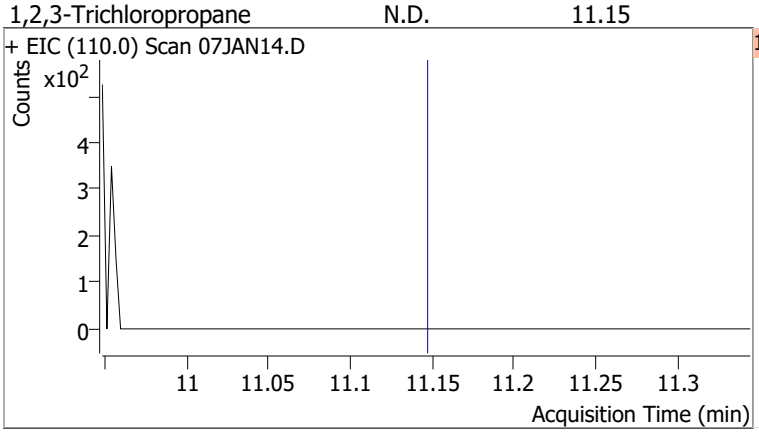


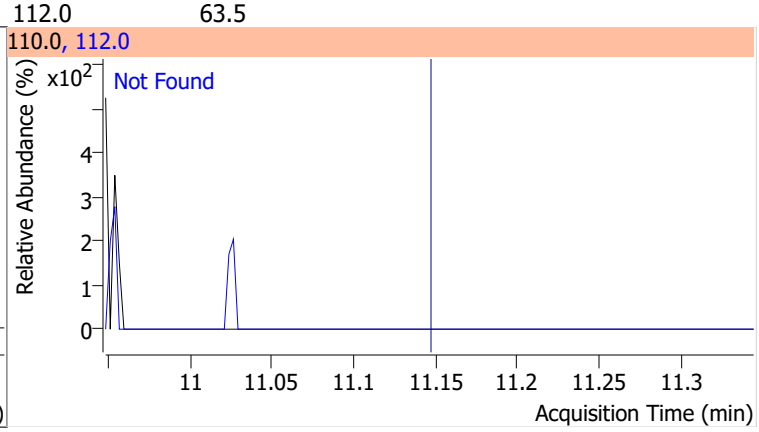
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



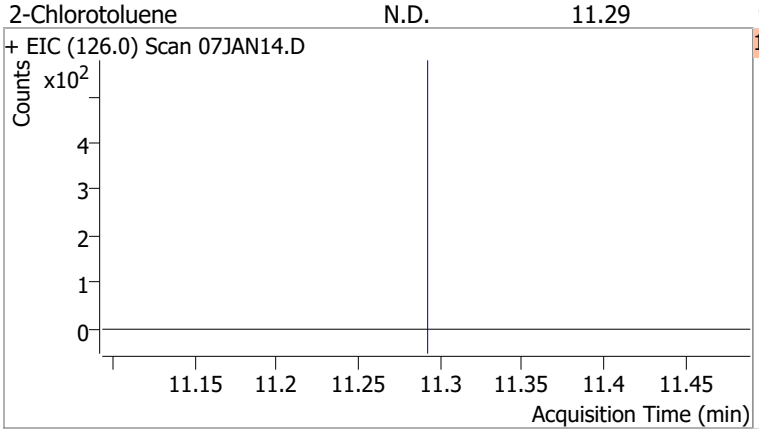


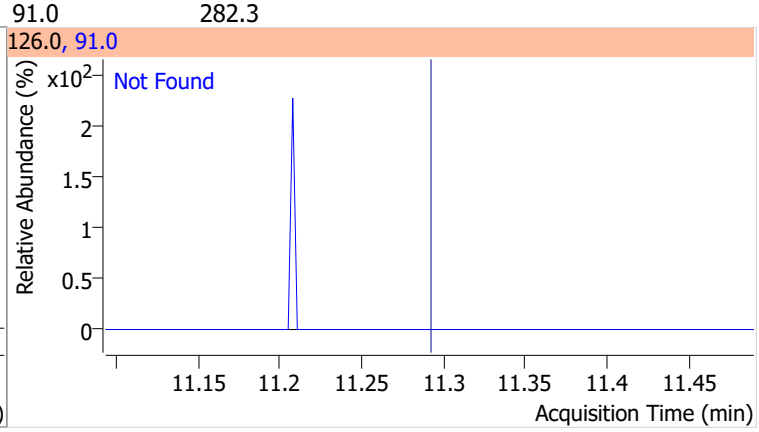
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



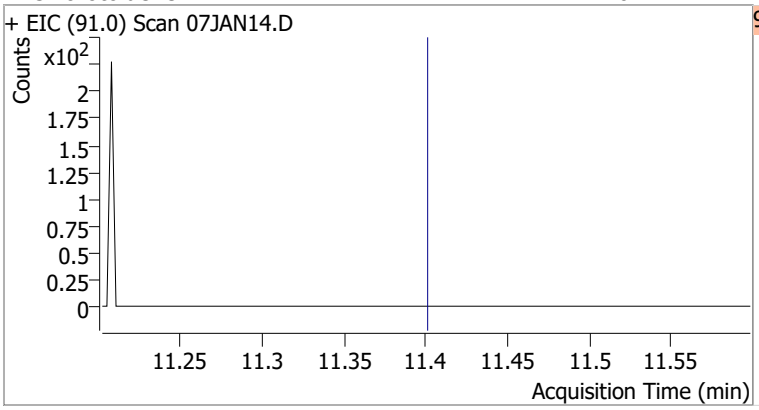


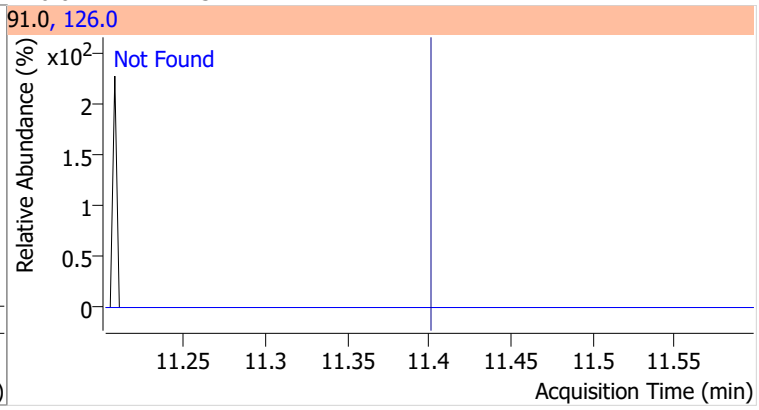
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



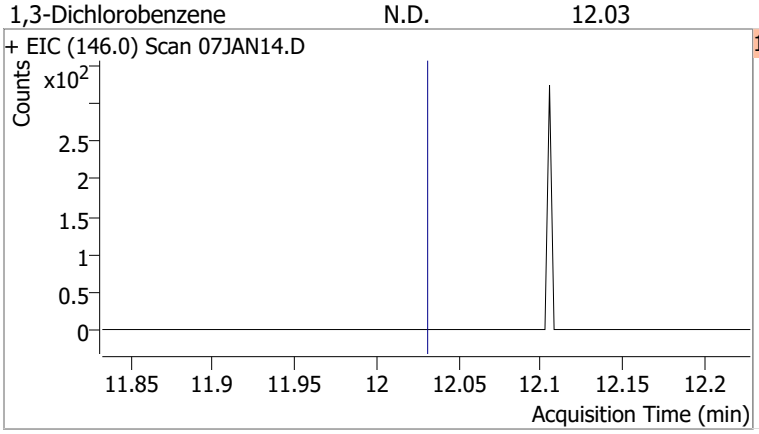


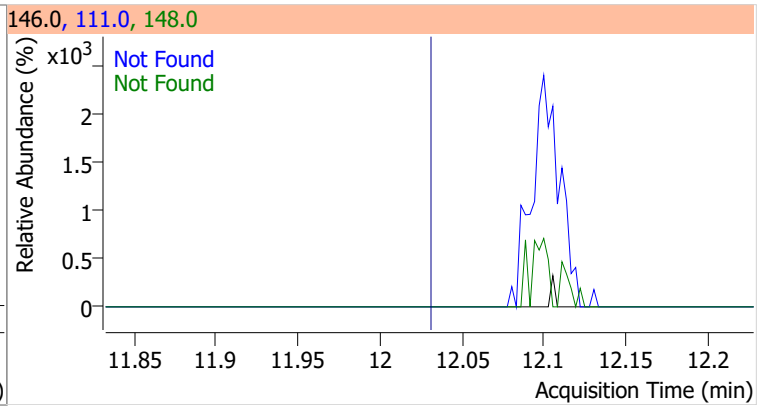
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



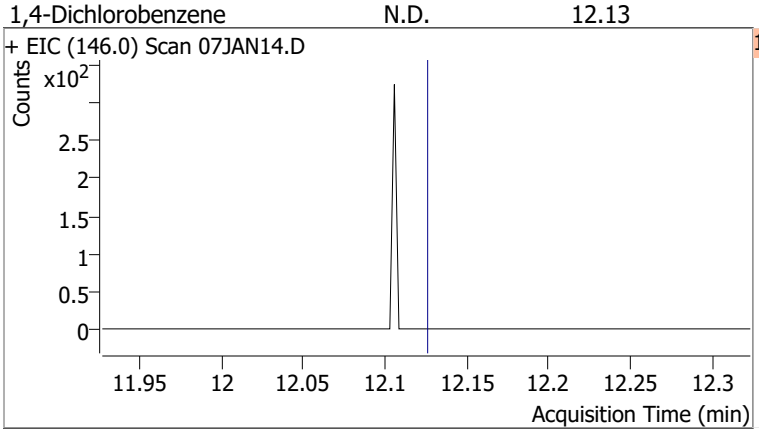


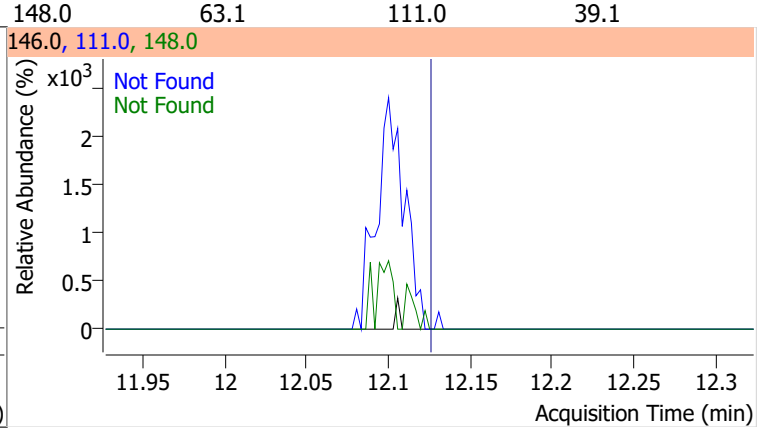
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



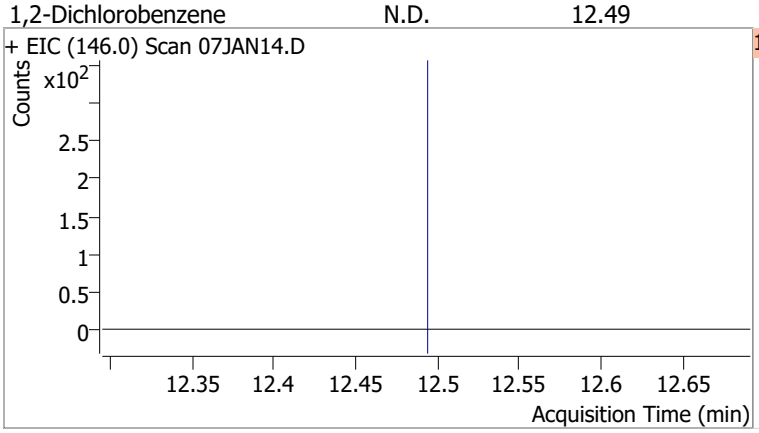


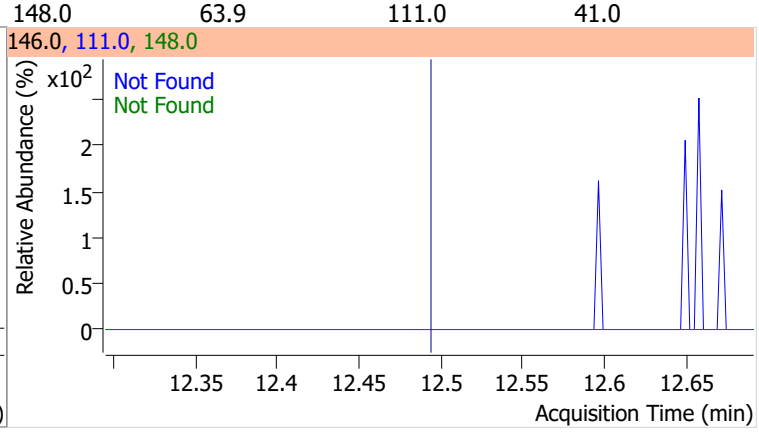
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





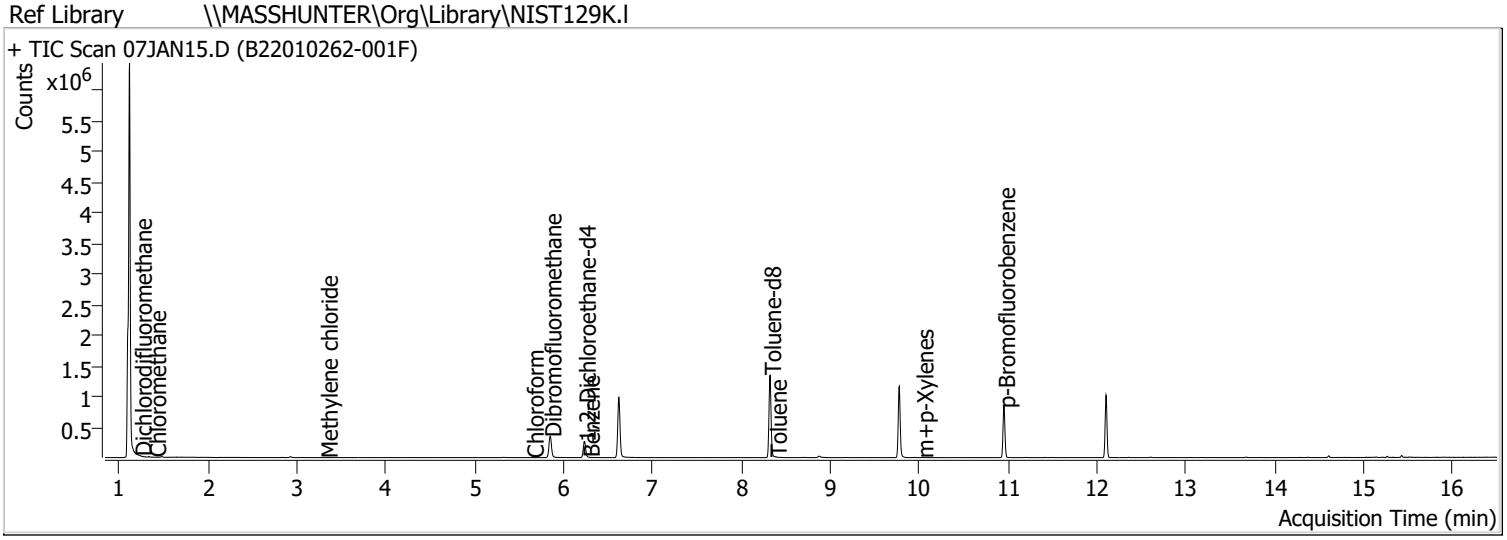
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN15.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 4:03:14 PM
Sample Name	B22010262-001F	Instrument	VOA5975C
Vial	15	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

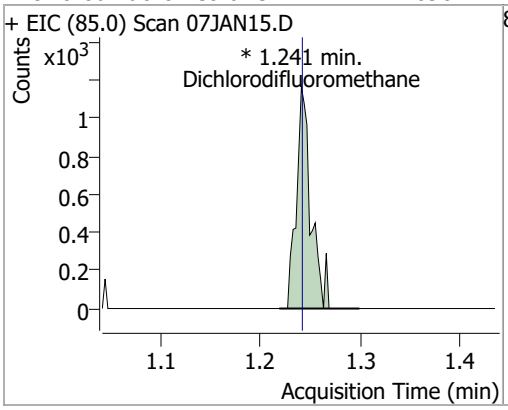


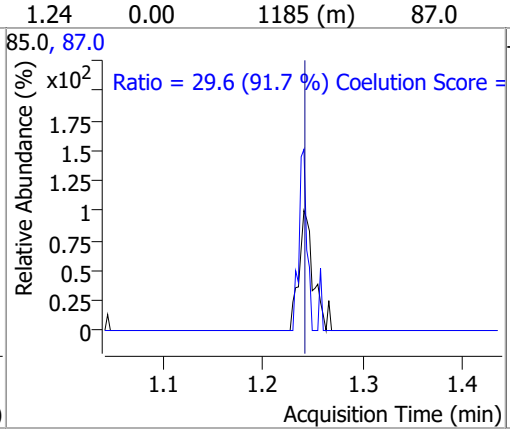
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
Internal Standards							
M Fluorobenzene	6.620	96.0	829021	250.0000	ng		-0.003
M Chlorobenzene-d5	9.774	82.0	317136	250.0000	ng		0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	243125	250.0000	ng		0.000
System Monitoring Compounds							
S Dibromofluoromethane	5.851	113.0	213658	273.5621	ng		0.005
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 109.42%			
S 1,2-Dichloroethane-d4	6.236	67.0	96525	286.1313	ng		0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 114.45%			
S Toluene-d8	8.321	98.0	826770	270.5325	ng		0.003
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 108.21%			
S p-Bromofluorobenzene	10.954	95.0	240345	269.8409	ng		0.000
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 107.94%			
Target Compounds							**QValue**
T Dichlorodifluoromethane	1.241	85.0	1185	1.0904	ng	m	95
T Chloromethane	1.406	50.0	626	0.4748	ng	m	60
T Vinyl chloride	0.000		0	N.D.			
T Bromomethane	0.000		0	N.D.			
T Chloroethane	0.000		0	N.D.			
T Trichlorofluoromethane	0.000		0	N.D.			
T 1,1-Dichloroethene	0.000		0	N.D.			
T Methylene chloride	3.335	49.0	648	0.5261	ng	m	98
T trans-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl tert-butyl ether (MTBE)	0.000		0	N.D.			
T 1,1-Dichloroethane	0.000		0	N.D.			
T 2,2-Dichloropropane	0.000		0	N.D.			
T cis-1,2-Dichloroethene	0.000		0	N.D.			
T Methyl ethyl ketone	0.000		0	N.D.			
T Bromochloromethane	0.000		0	N.D.			
T Chloroform	5.653	83.0	600	0.3801	ng	#m	35

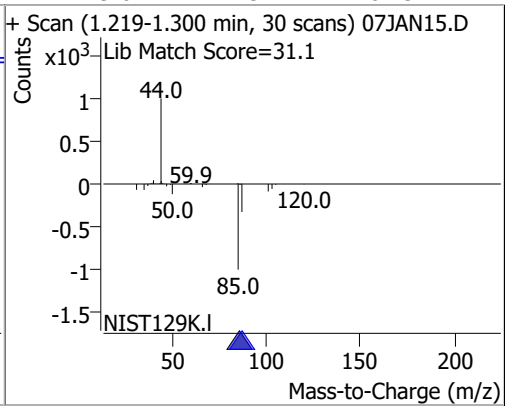
Compound	RT	QIon	Resp.	Conc.	Units		Dev(Min)
T 1,1,1-Trichloroethane	0.000		0	N.D.			
T Carbon tetrachloride	0.000		0	N.D.			
T 1,1-Dichloropropene	0.000		0	N.D.			
T Benzene	6.283	78.0	748	0.2266	ng	m	81
T 1,2-Dichloroethane	0.000		0	N.D.			
T Trichloroethene	0.000		0	N.D.			
T 1,2-Dichloropropane	0.000		0	N.D.			
T Dibromomethane	0.000		0	N.D.			
T Bromodichloromethane	0.000		0	N.D.			
T cis-1,3-Dichloropropene	0.000		0	N.D.			
T Toluene	8.386	92.0	723	0.3502	ng	m	100
T trans-1,3-Dichloropropene	0.000		0	N.D.			
T 1,1,2-Trichloroethane	0.000		0	N.D.			
T Tetrachloroethene	0.000		0	N.D.			
T 1,3-Dichloropropane	0.000		0	N.D.			
T Chlorodibromomethane	0.000		0	N.D.			
T 1,2-Dibromoethane	0.000		0	N.D.			
T Chlorobenzene	0.000		0	N.D.			
T 1,1,1,2-Tetrachloroethane	0.000		0	N.D.			
T Ethylbenzene	0.000		0	N.D.			
T m+p-Xylenes	10.036	106.0	505	0.3315	ng	m	87
T o-Xylene	10.427	106.0	0		ng	md	1
T Styrene	0.000		0	N.D.			
T Bromoform	0.000		0	N.D.			
T Bromobenzene	0.000		0	N.D.			
T 1,1,2,2-Tetrachloroethane	0.000		0	N.D.			
T 1,2,3-Trichloropropane	0.000		0	N.D.			
T 2-Chlorotoluene	0.000		0	N.D.			
T 4-Chlorotoluene	0.000		0	N.D.			
T 1,3-Dichlorobenzene	0.000		0	N.D.			
T 1,4-Dichlorobenzene	0.000		0	N.D.			
T 1,2-Dichlorobenzene	0.000		0	N.D.			

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

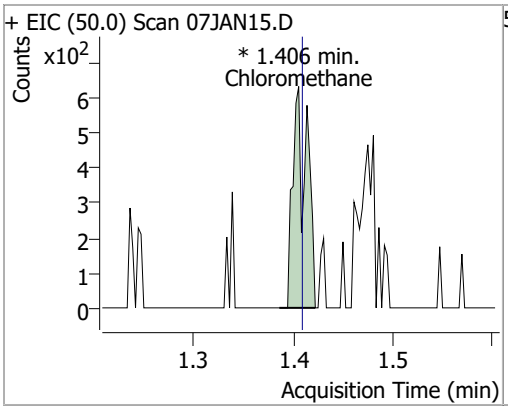
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	1.0904	1.24	0.00	1185 (m)	87.0	29.6	2.3	62.3

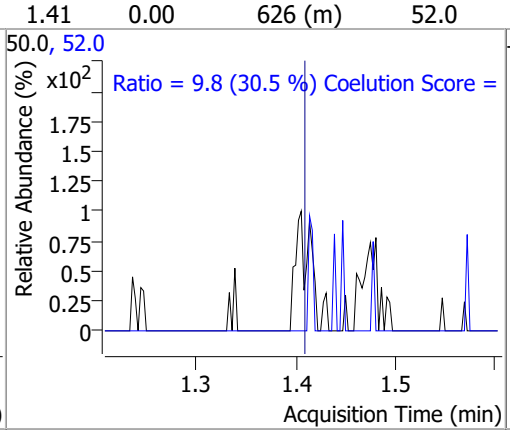


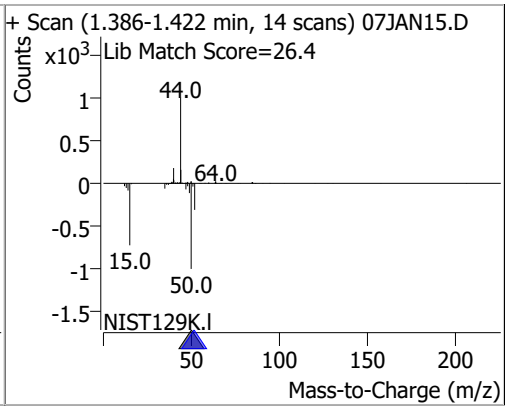




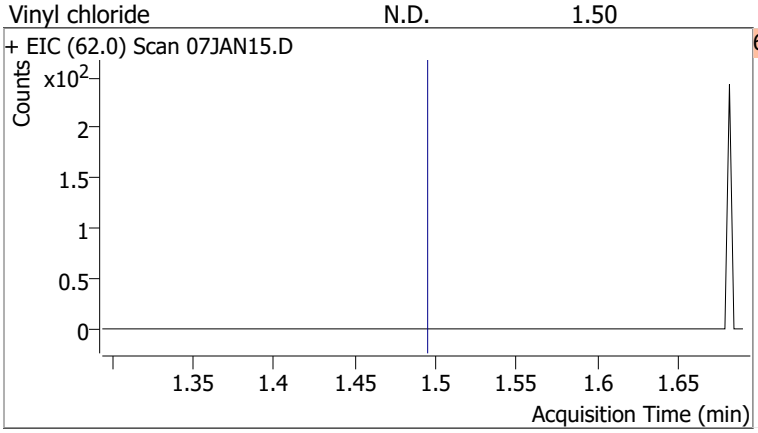
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	0.4748	1.41	0.00	626 (m)	52.0	9.8	2.1	62.1

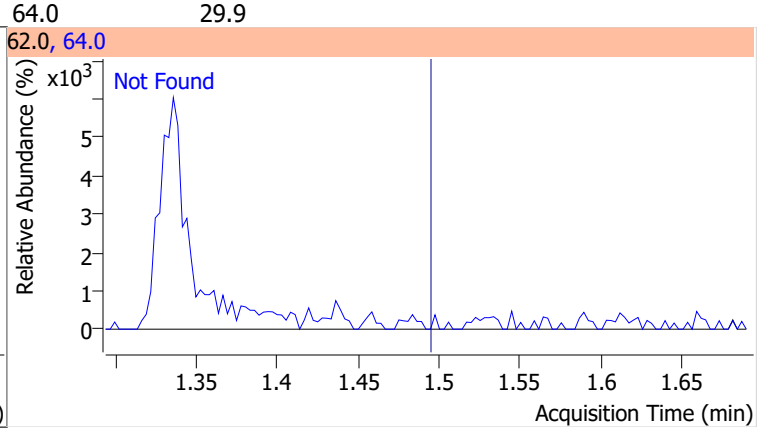




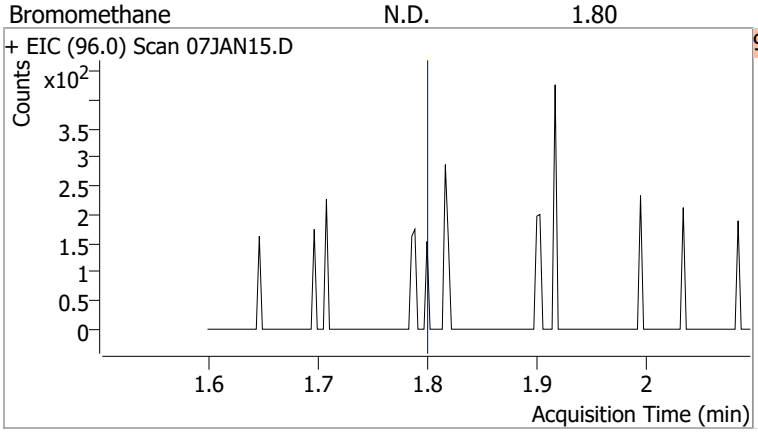


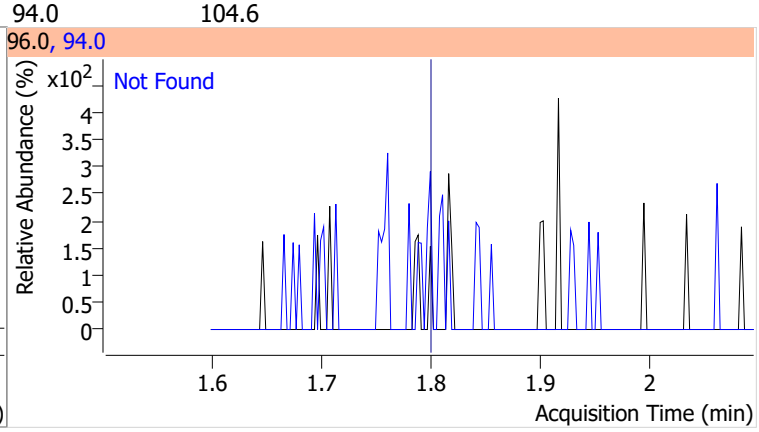
Compound	Conc.	Exp RT	QIon	Exp Ratio
Vinyl chloride	N.D.	1.50	64.0	29.9



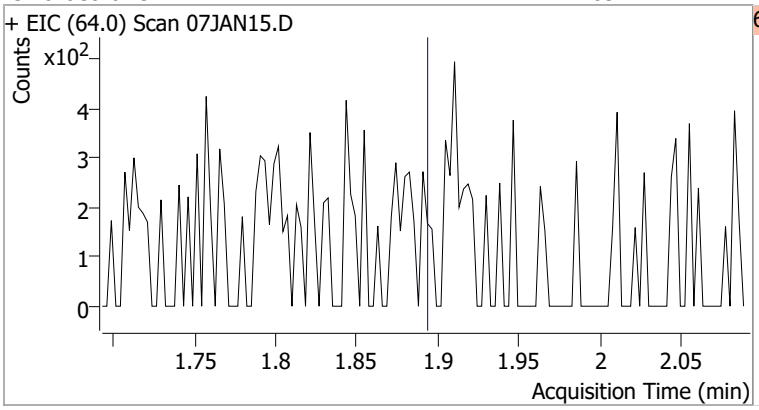


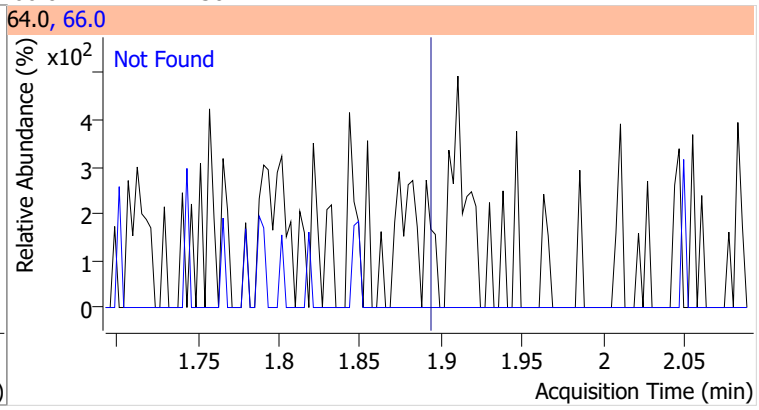
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromomethane	N.D.	1.80	94.0	104.6



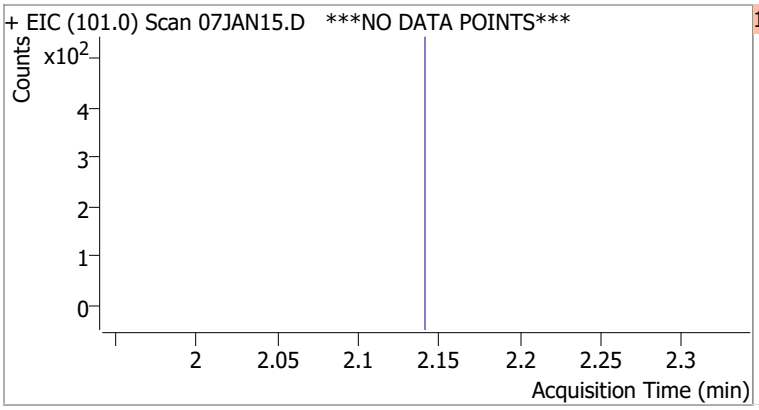


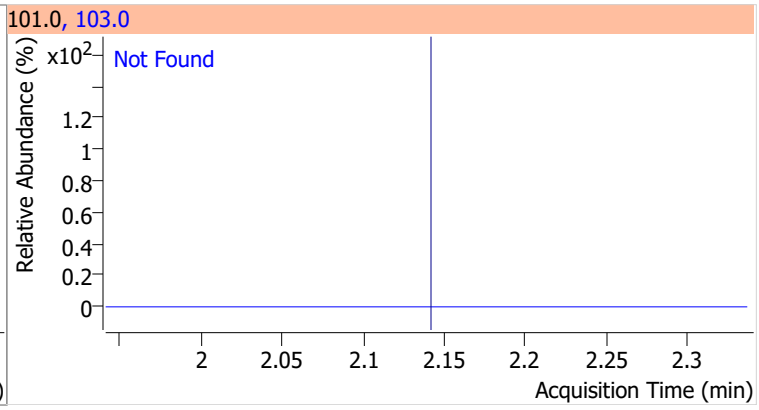
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chloroethane	N.D.	1.89	66.0	30.1



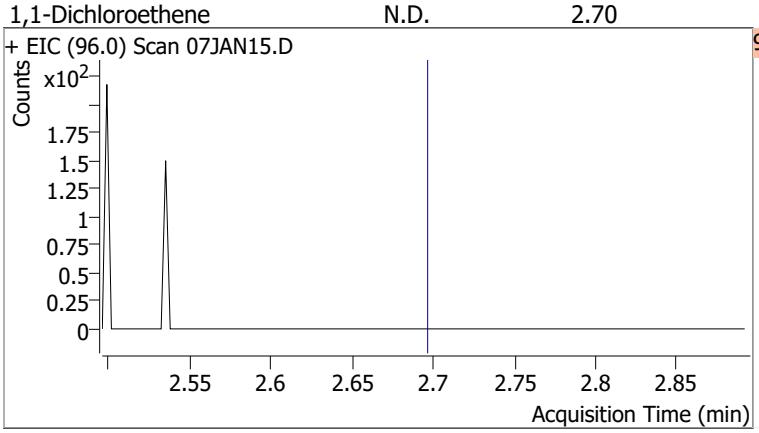


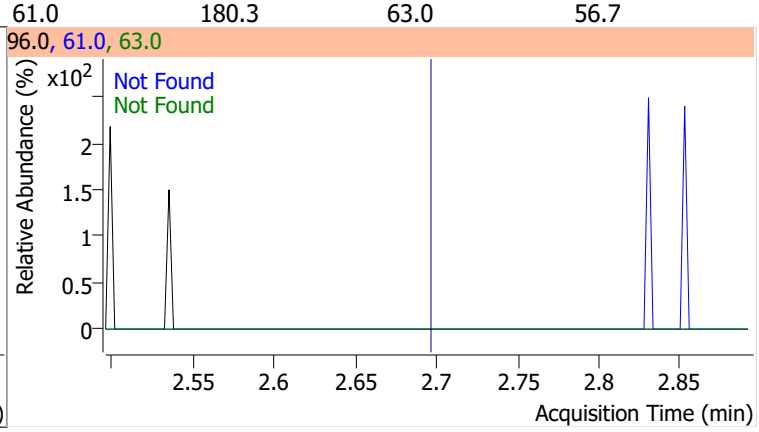
Compound	Conc.	Exp RT	QIon	Exp Ratio
Trichlorofluoromethane	N.D.	2.14	103.0	64.2



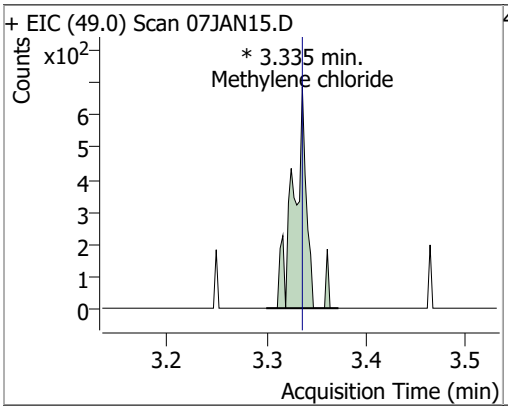


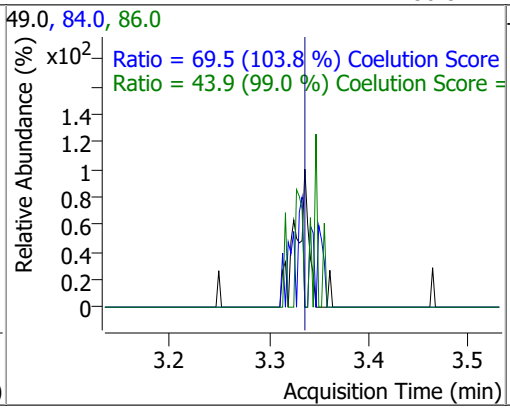
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethene	N.D.	2.70	61.0	180.3	63.0	56.7

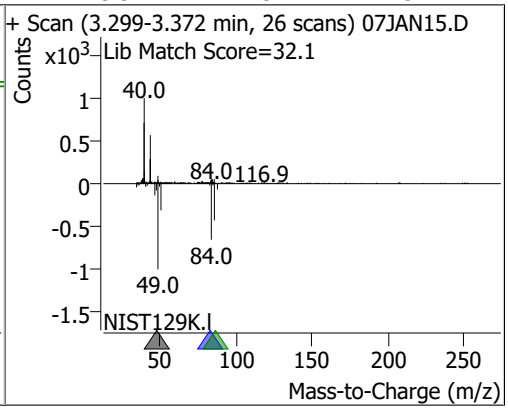




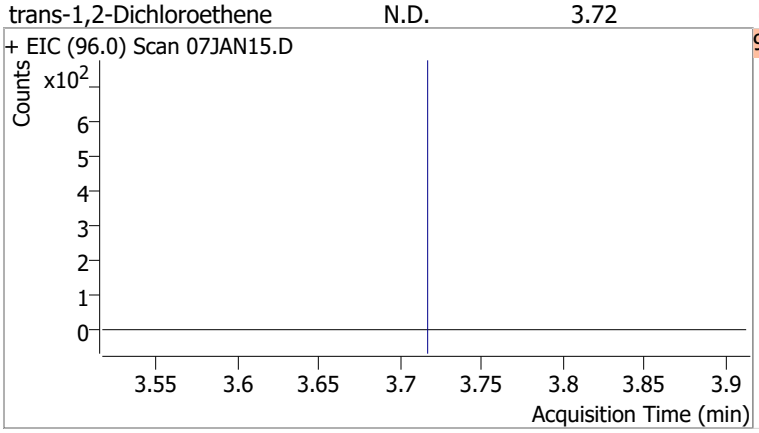
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	0.5261	3.34	0.00	648 (m)	84.0	69.5	36.9	96.9
					86.0	43.9	14.3	74.3

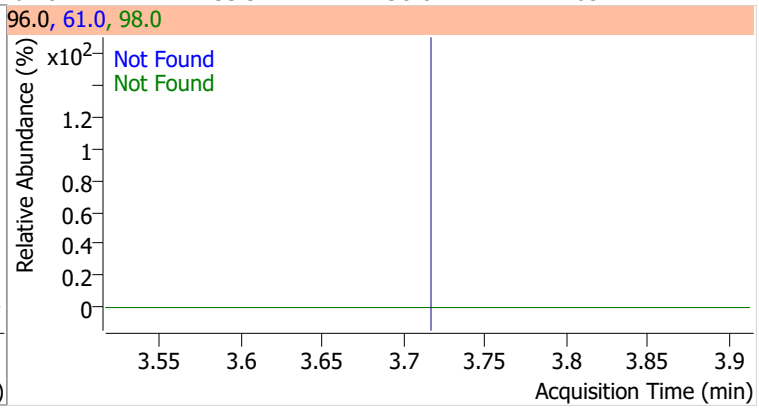




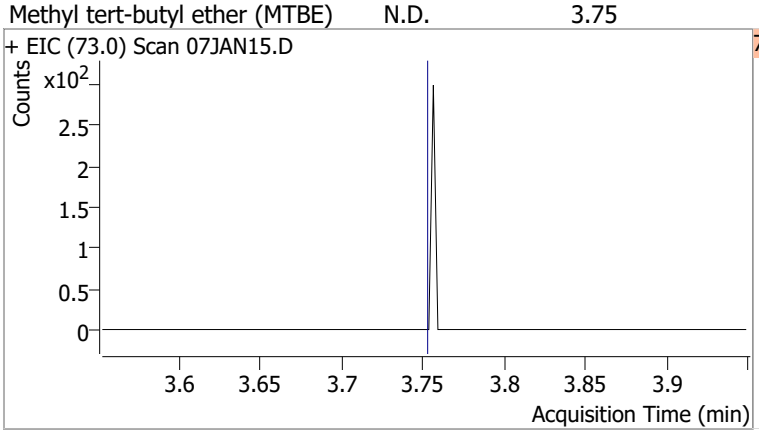


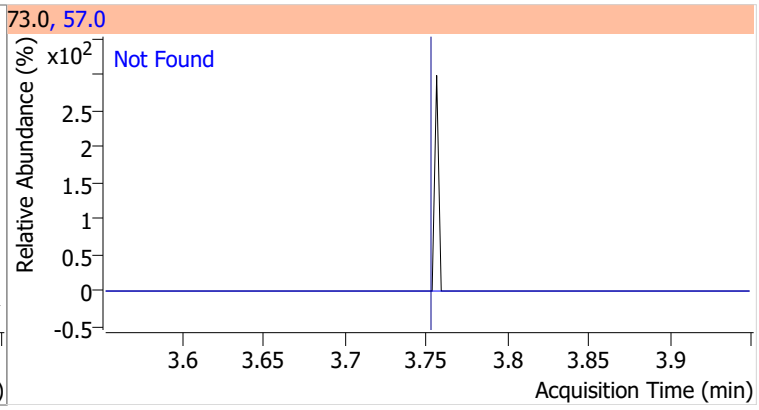
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,2-Dichloroethene	N.D.	3.72	61.0	153.9	98.0	65.7



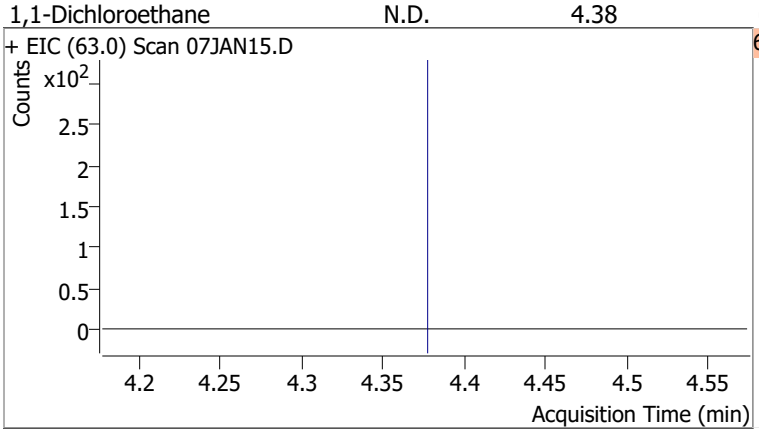


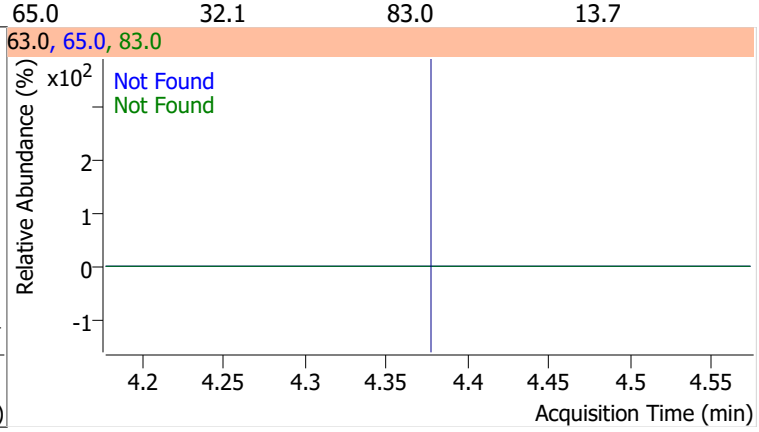
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl tert-butyl ether (MTBE)	N.D.	3.75	57.0	24.6



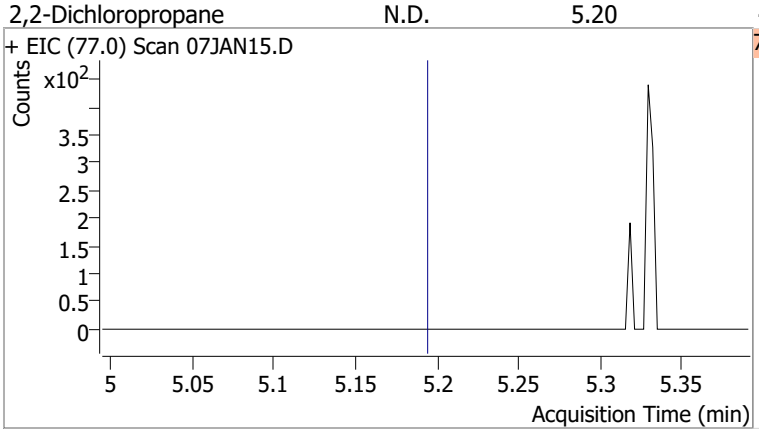


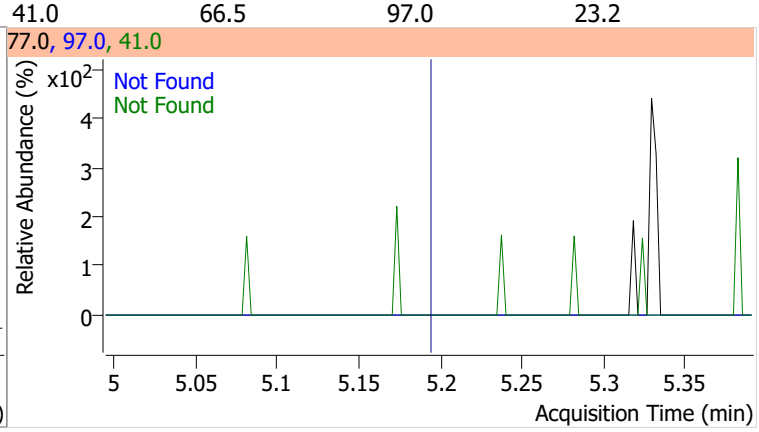
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloroethane	N.D.	4.38	65.0	32.1	83.0	13.7



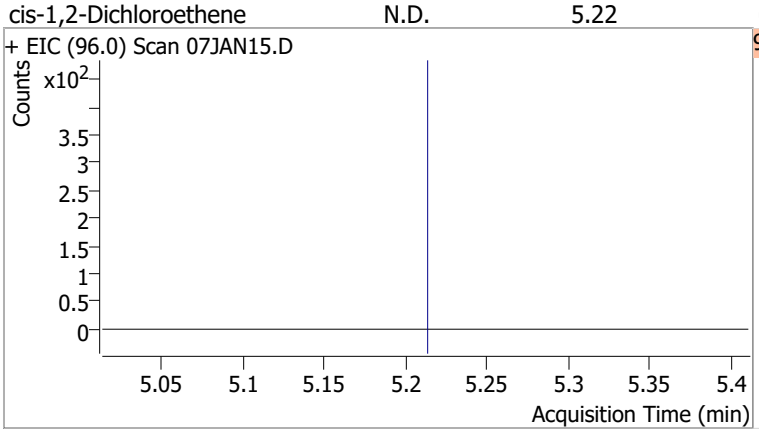


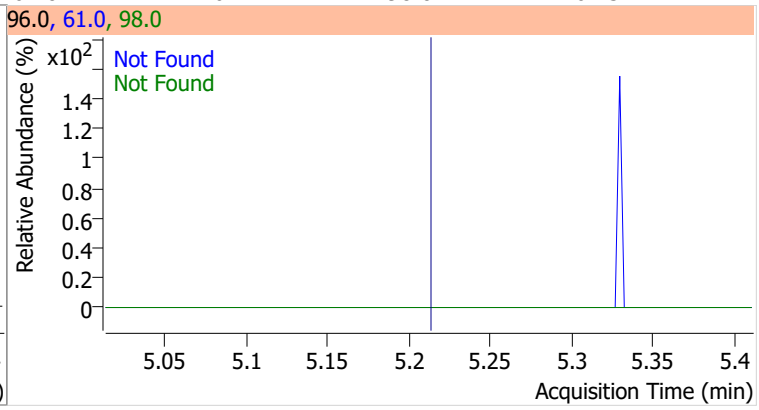
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
2,2-Dichloropropane	N.D.	5.20	41.0	66.5	97.0	23.2



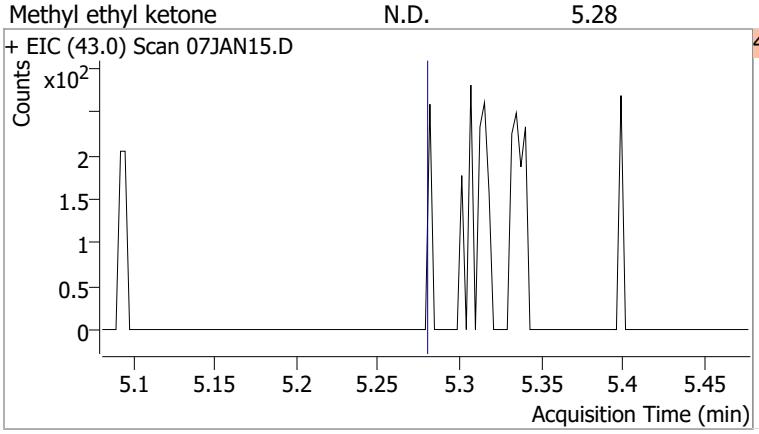


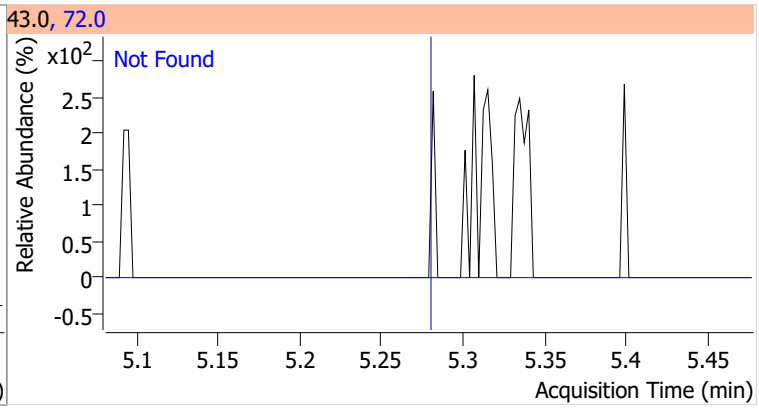
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,2-Dichloroethene	N.D.	5.22	61.0	167.2	98.0	67.3



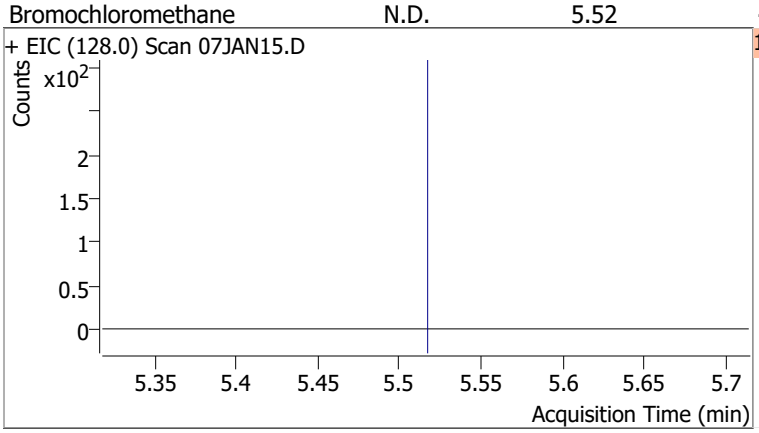


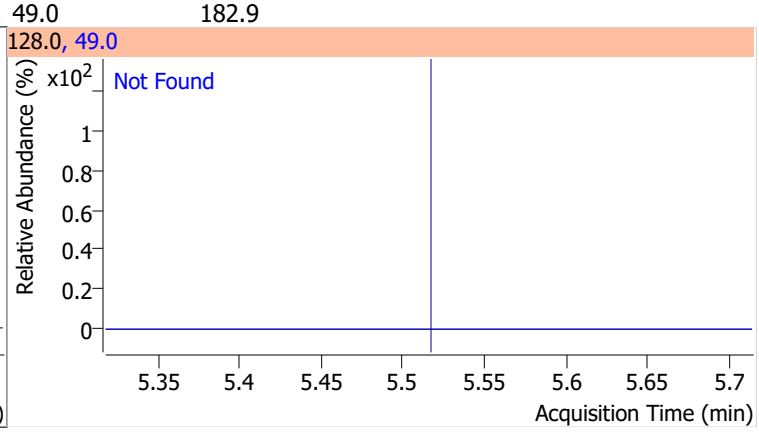
Compound	Conc.	Exp RT	QIon	Exp Ratio
Methyl ethyl ketone	N.D.	5.28	72.0	21.3



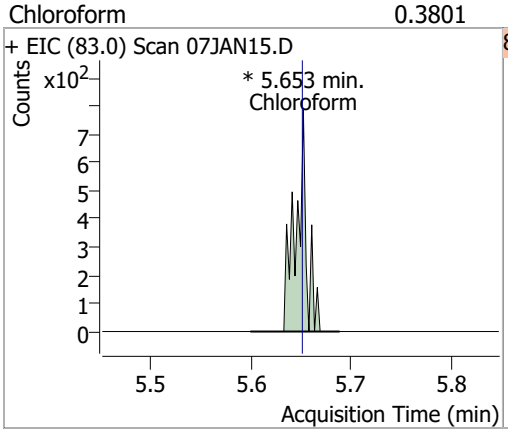


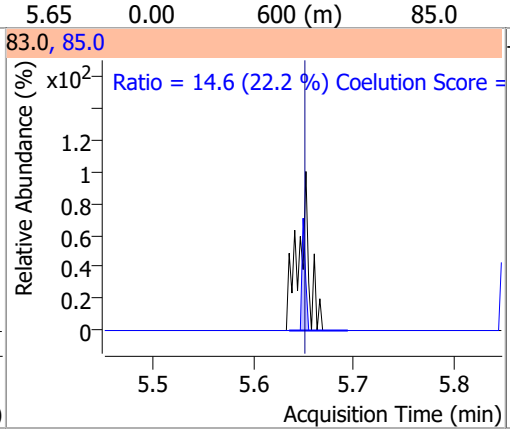
Compound	Conc.	Exp RT	QIon	Exp Ratio
Bromochloromethane	N.D.	5.52	49.0	182.9

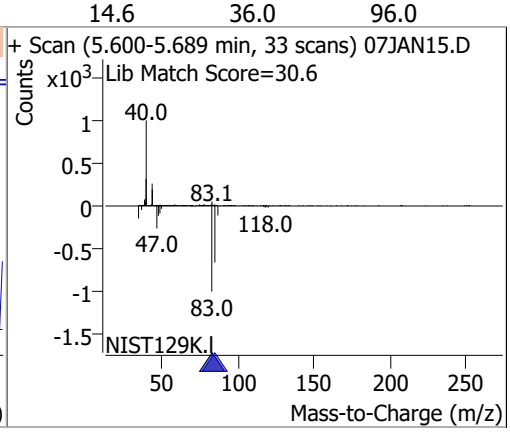




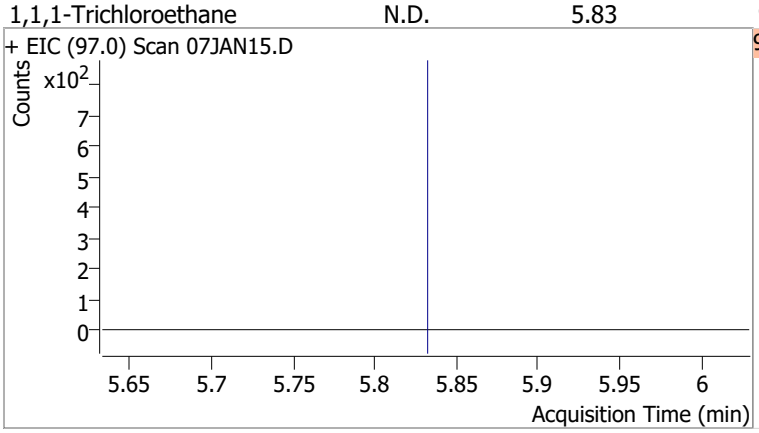
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	0.3801	5.65	0.00	600 (m)	85.0	14.6	36.0	96.0

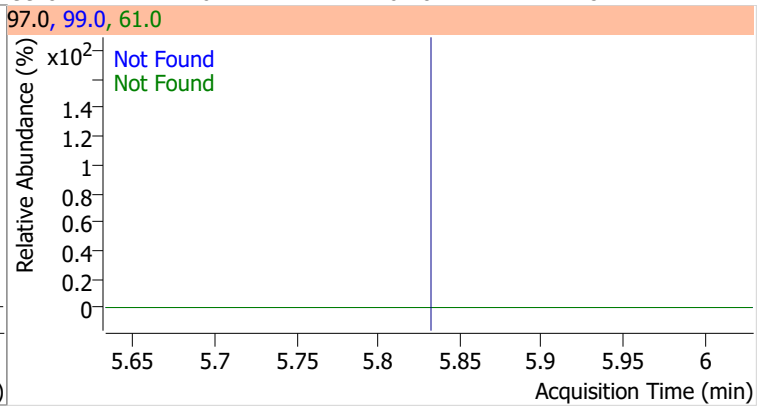




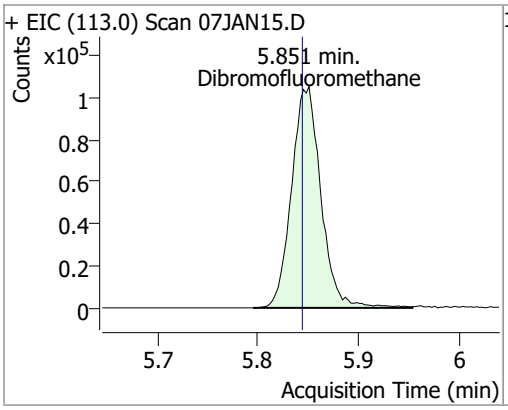


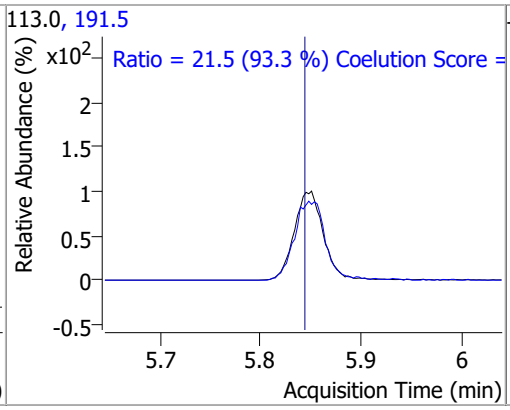
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,1-Trichloroethane	N.D.	5.83	99.0	64.7	61.0	48.1

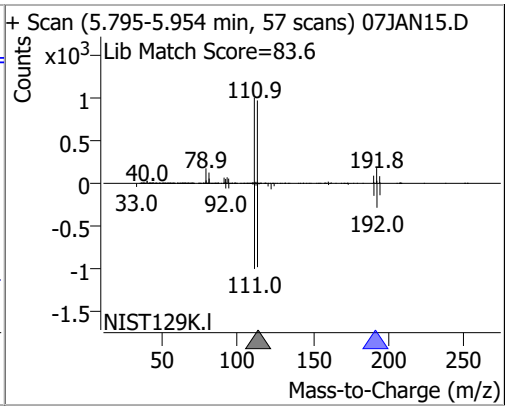




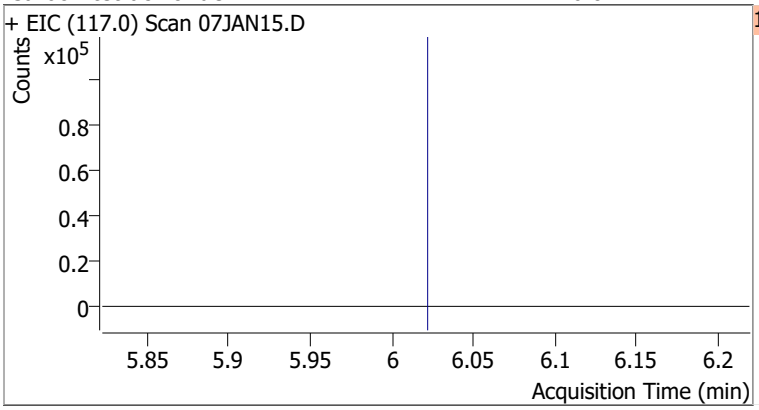
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	273.5621	5.85	0.01	213658	191.5	21.5	0.0	53.1

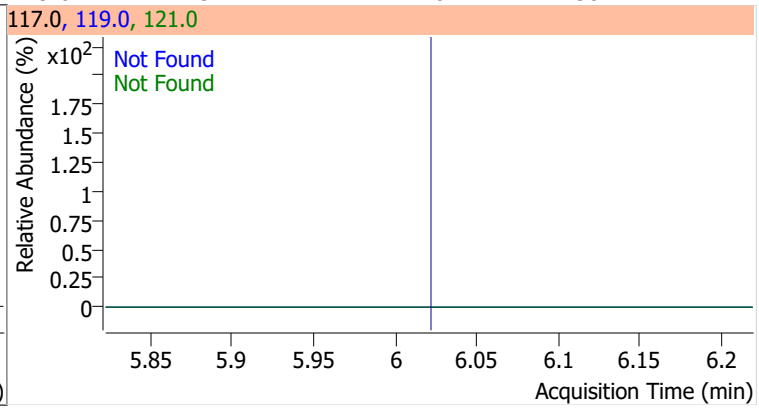




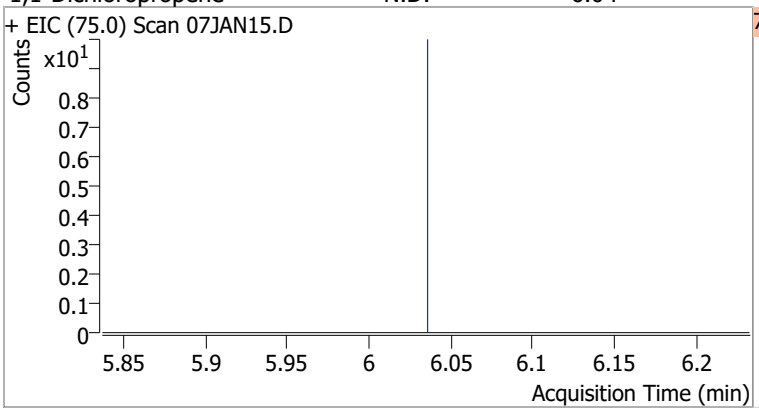


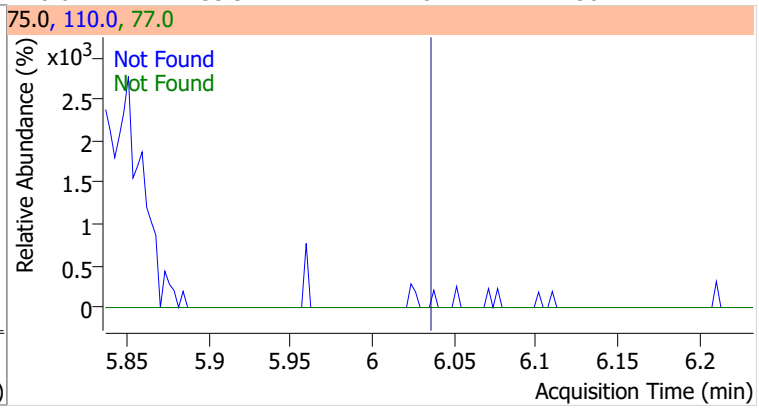
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Carbon tetrachloride	N.D.	6.02	119.0	97.2	121.0	30.1



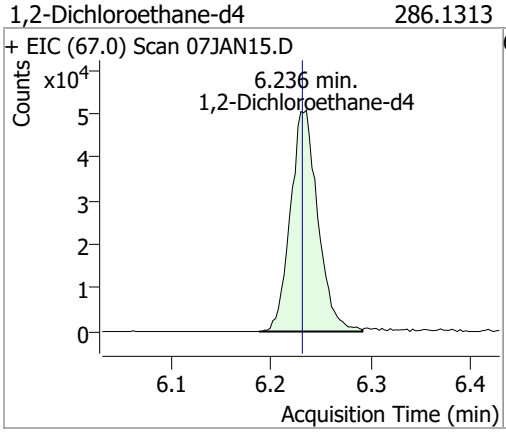


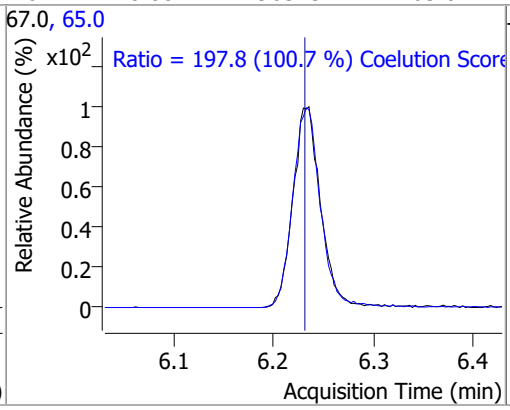
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1-Dichloropropene	N.D.	6.04	110.0	35.9	77.0	30.1

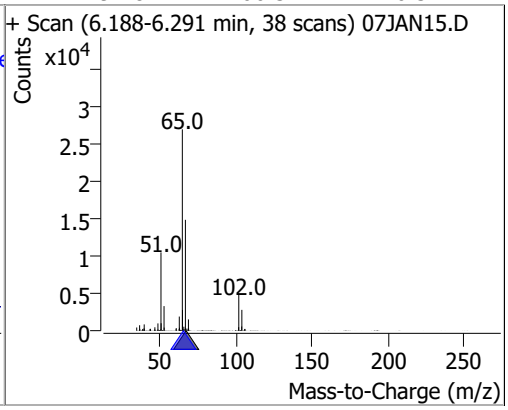




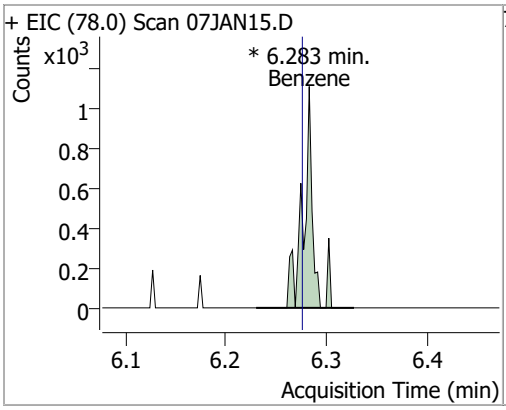
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	286.1313	6.24	0.00	96525	65.0	197.8	166.5	226.5

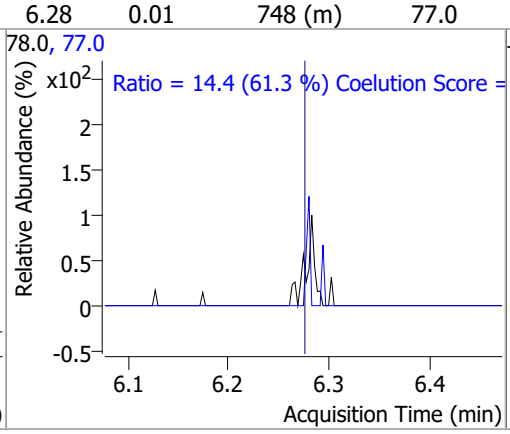


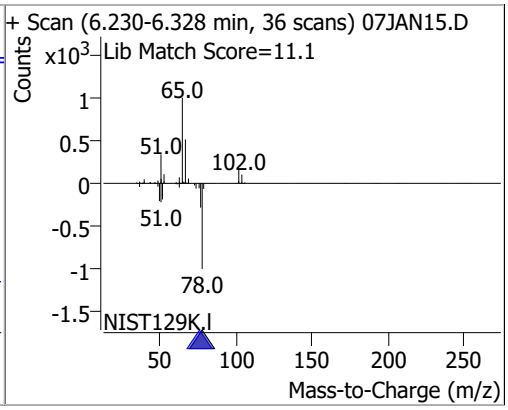




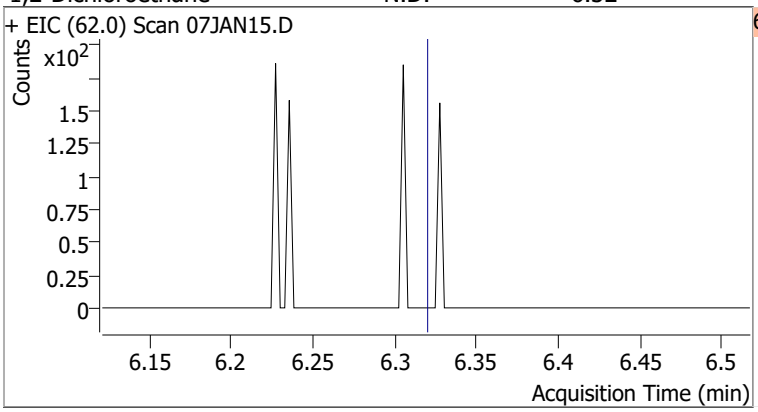
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	0.2266	6.28	0.01	748 (m)	77.0	14.4	0.0	53.5

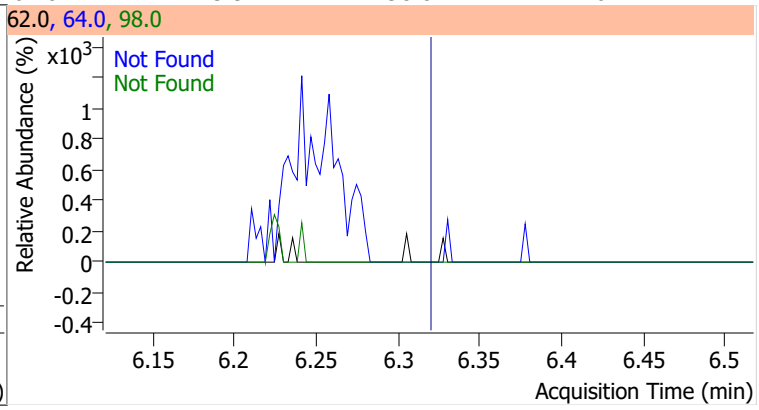




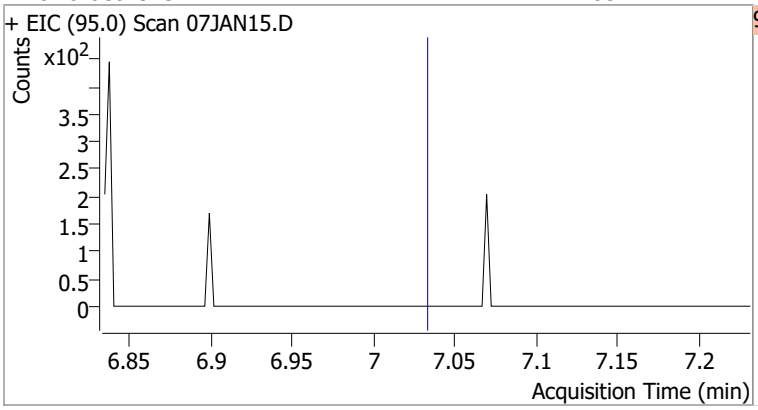


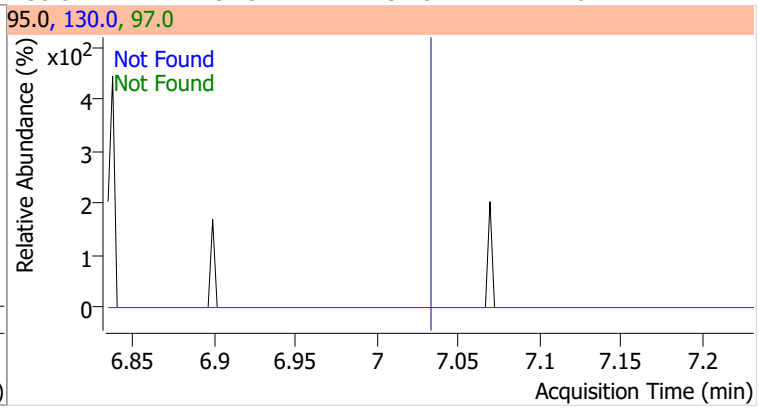
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichloroethane	N.D.	6.32	64.0	29.9	98.0	7.6



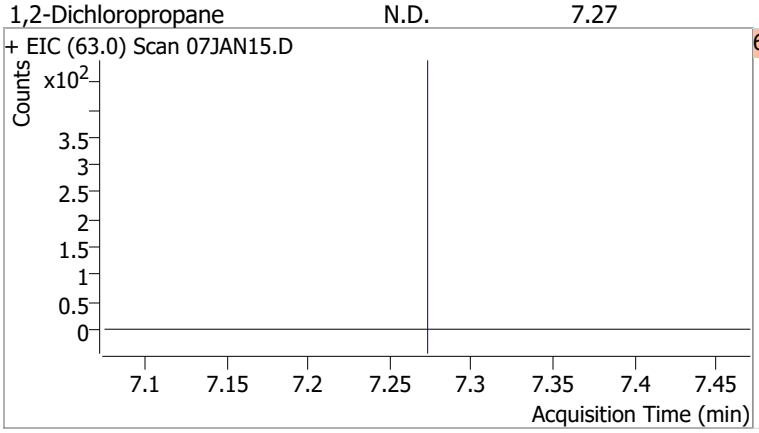


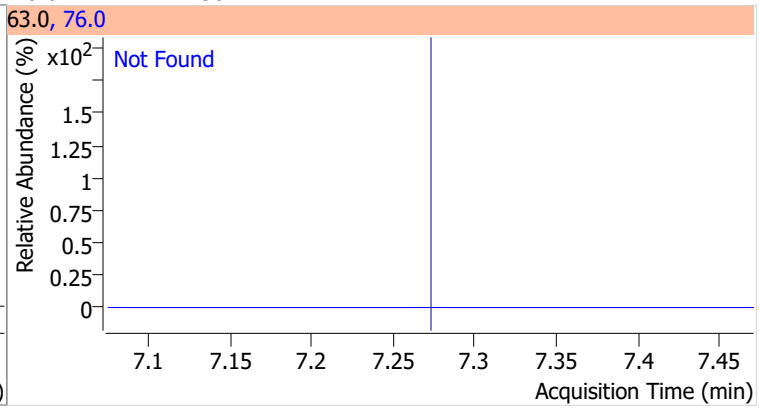
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Trichloroethene	N.D.	7.03	130.0	101.5	97.0	64.1



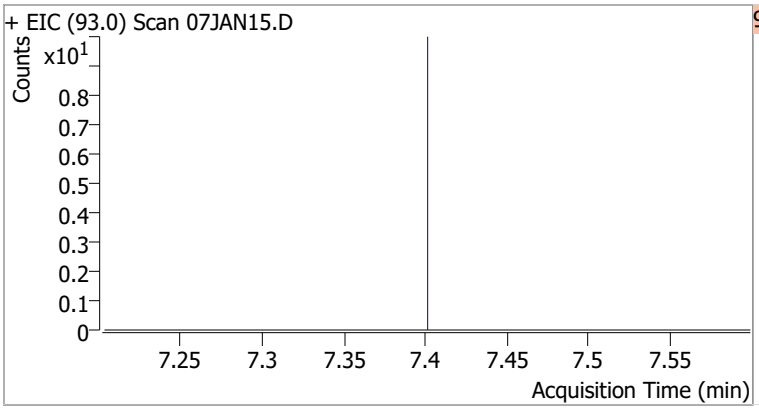


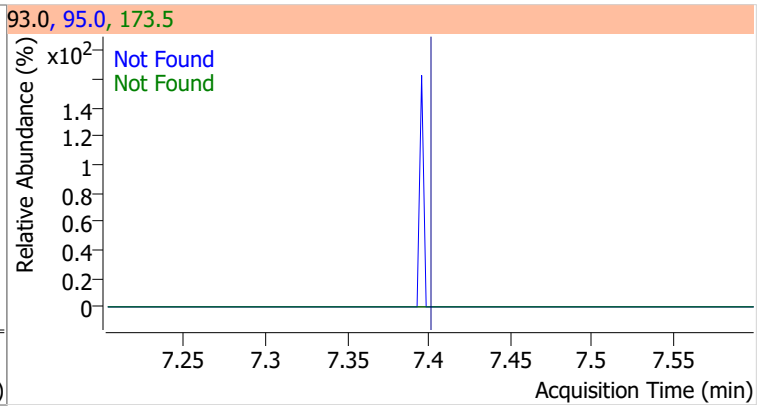
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dichloropropane	N.D.	7.27	76.0	38.2



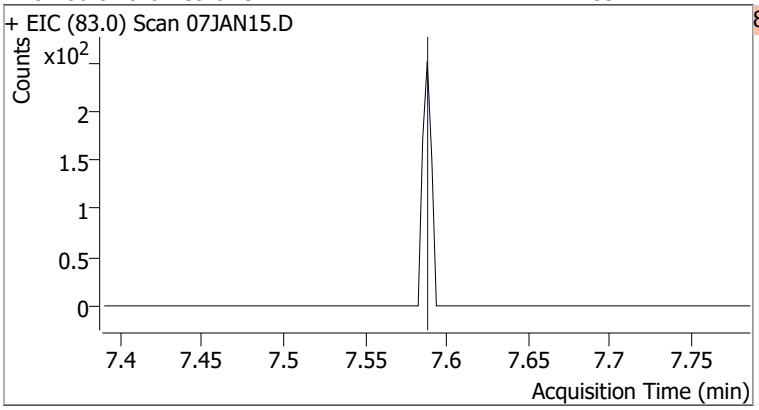


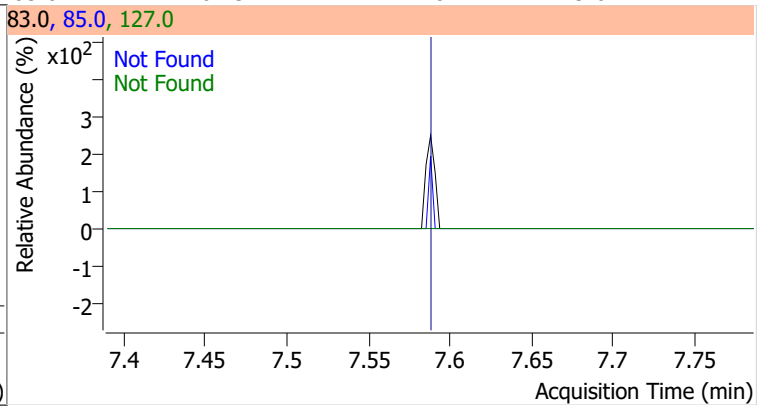
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Dibromomethane	N.D.	7.40	173.5	113.7	95.0	82.2



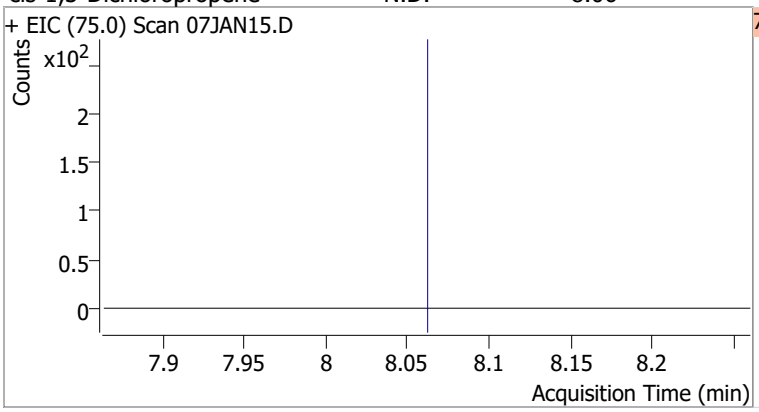


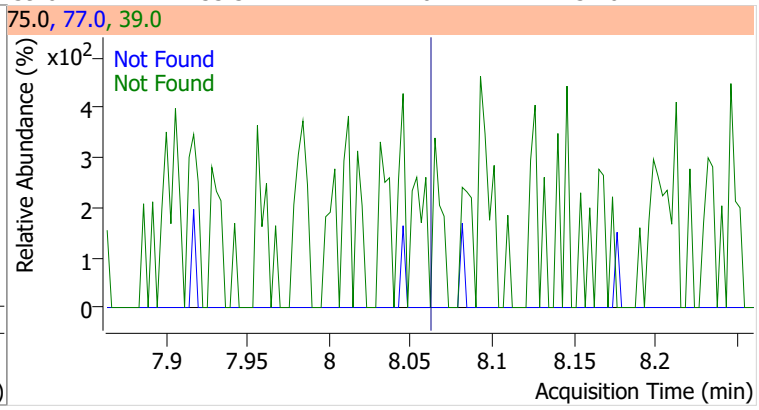
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromodichloromethane	N.D.	7.59	85.0	64.5	127.0	9.6



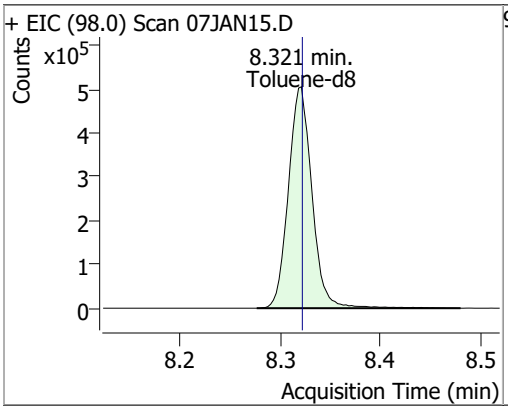


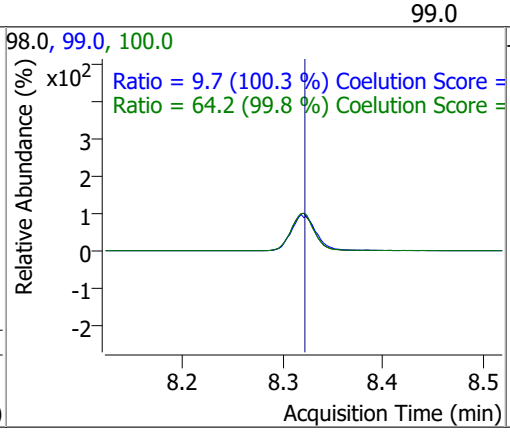
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
cis-1,3-Dichloropropene	N.D.	8.06	39.0	53.3	77.0	31.0

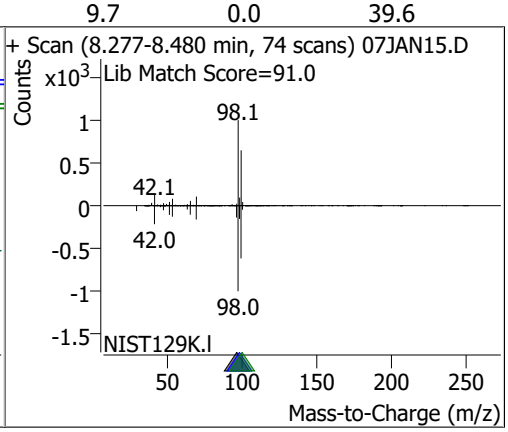




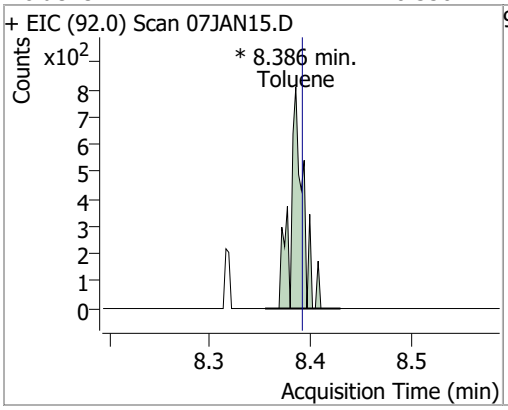
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	270.5325	8.32	0.00	826770	100.0	64.2	34.4	94.4
					99.0	9.7	0.0	39.6

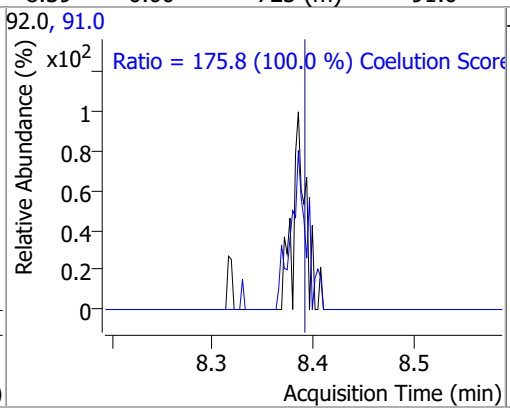


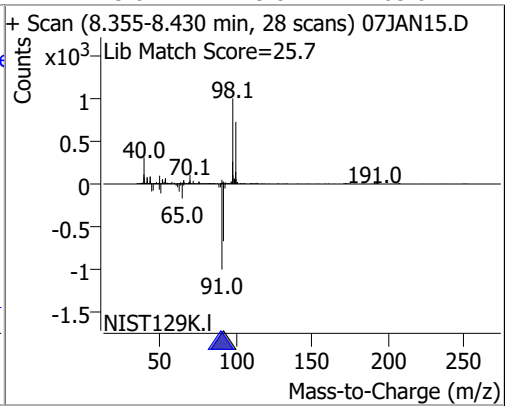




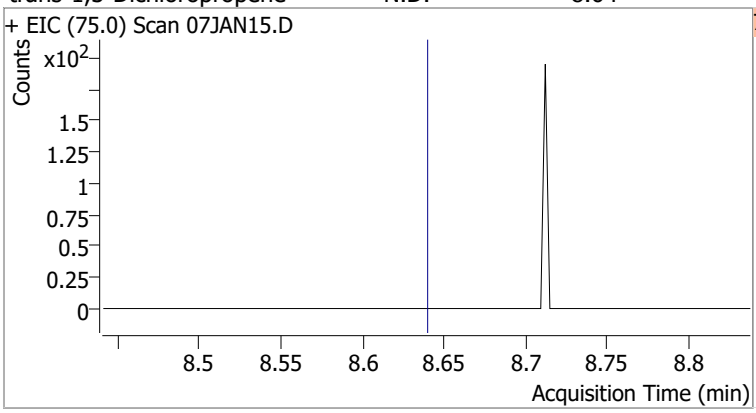
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	0.3502	8.39	0.00	723 (m)	91.0	175.8	145.8	205.8

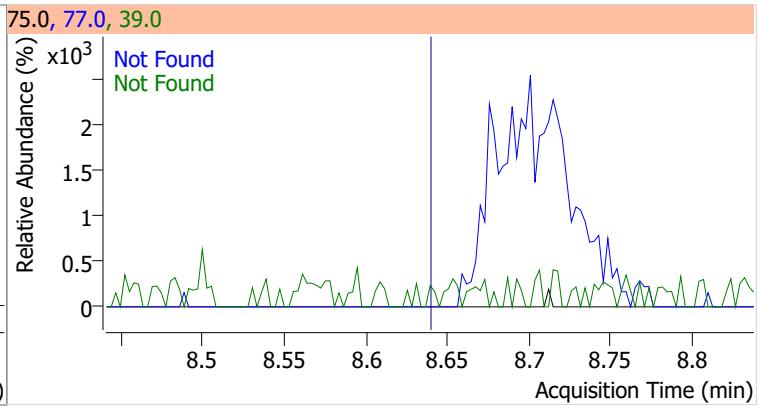




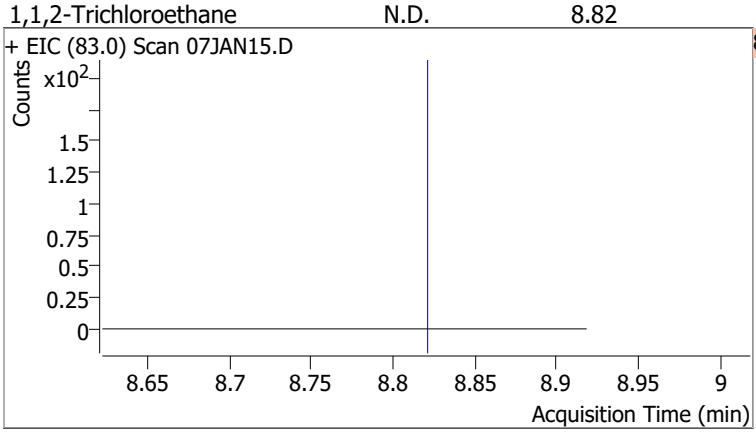


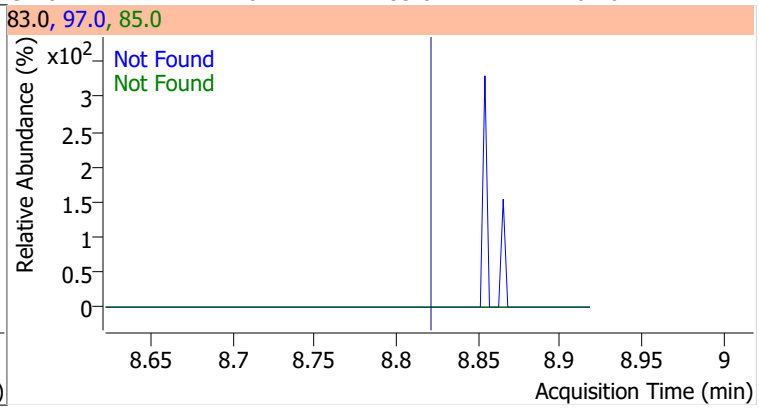
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
trans-1,3-Dichloropropene	N.D.	8.64	39.0	53.4	77.0	32.4



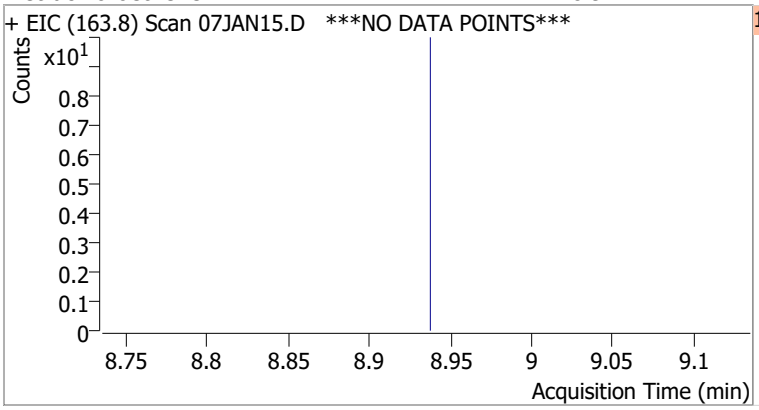


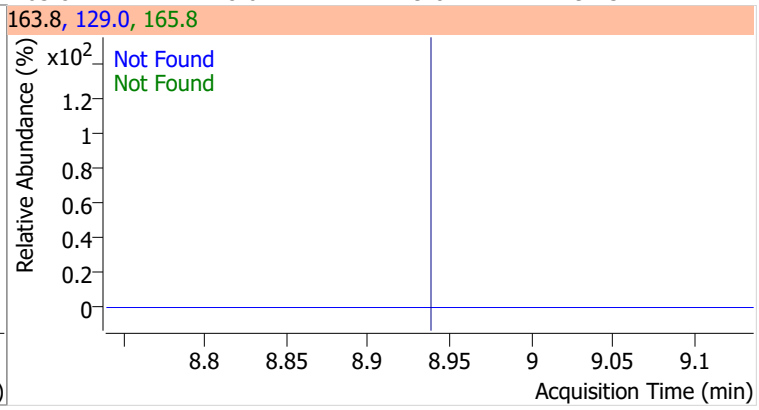
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,1,2-Trichloroethane	N.D.	8.82	97.0	114.6	85.0	67.6



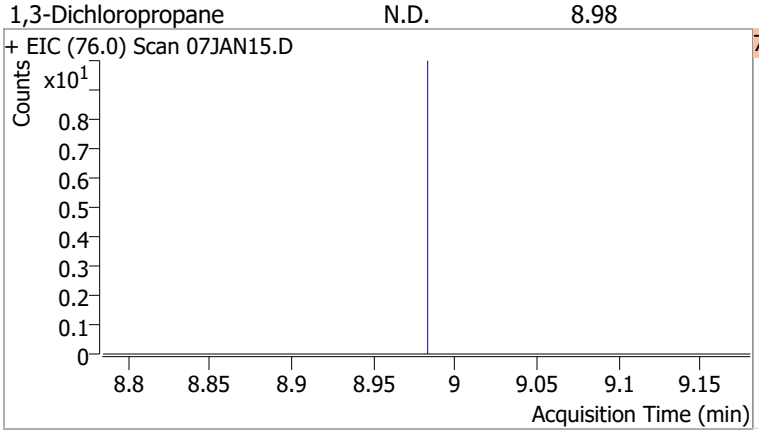


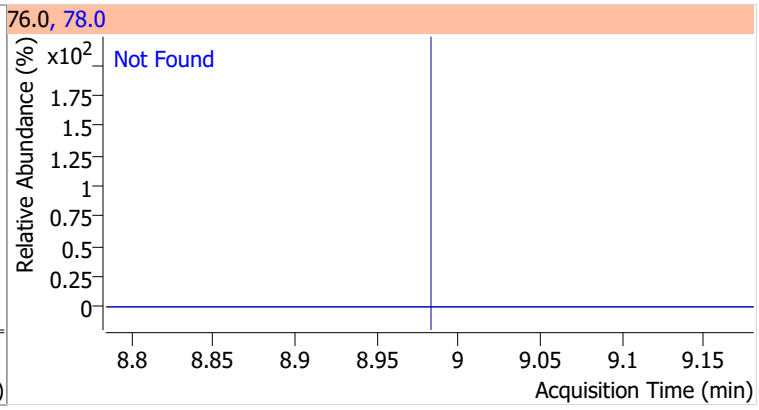
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Tetrachloroethene	N.D.	8.94	165.8	128.6	129.0	91.5



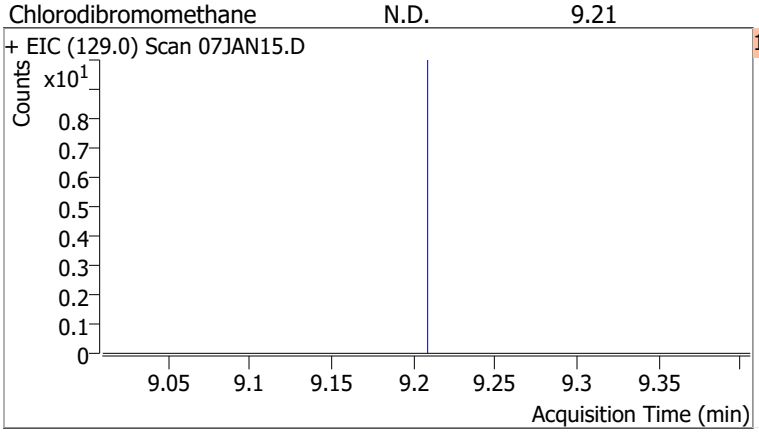


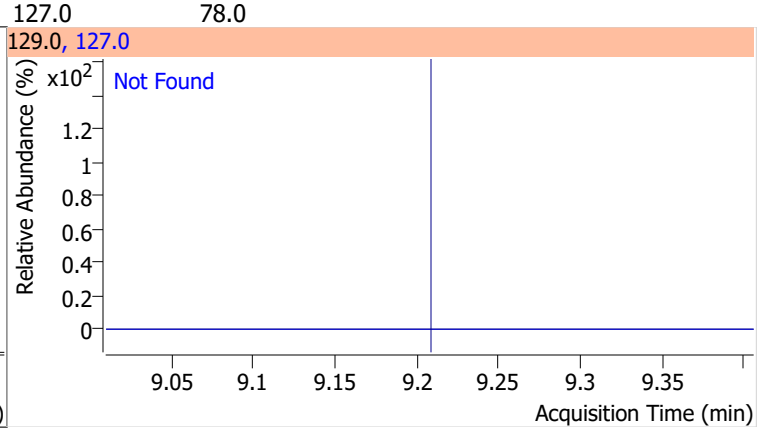
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,3-Dichloropropane	N.D.	8.98	78.0	32.9



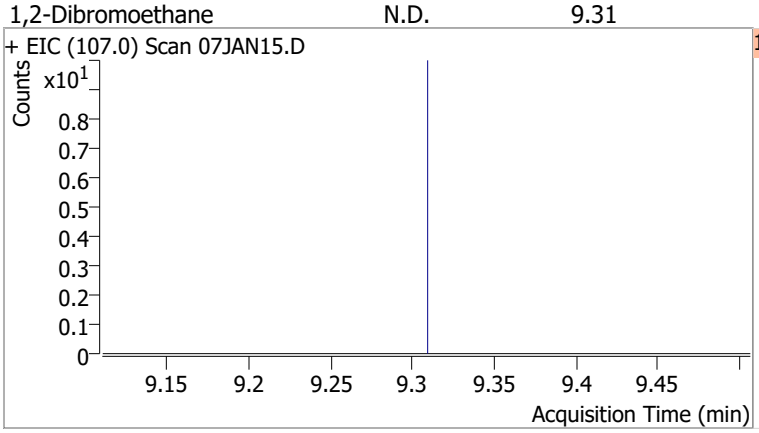


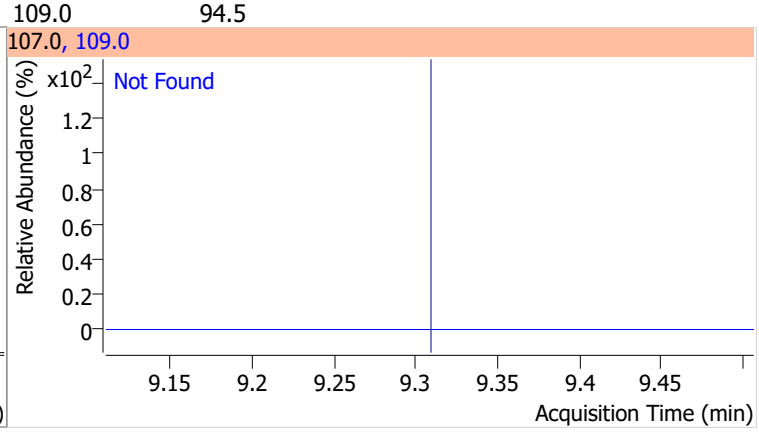
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorodibromomethane	N.D.	9.21	127.0	78.0



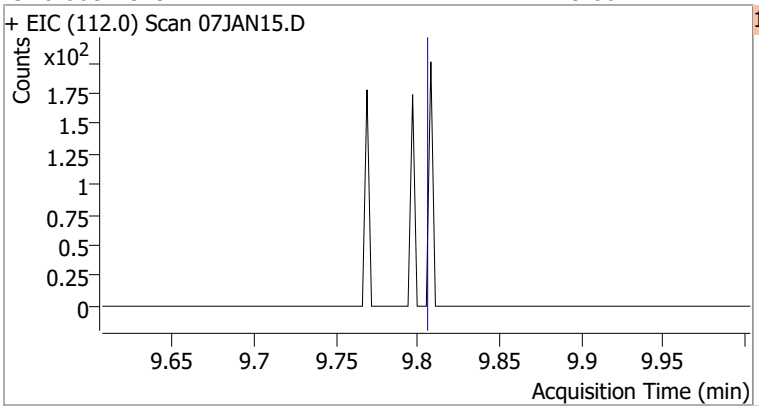


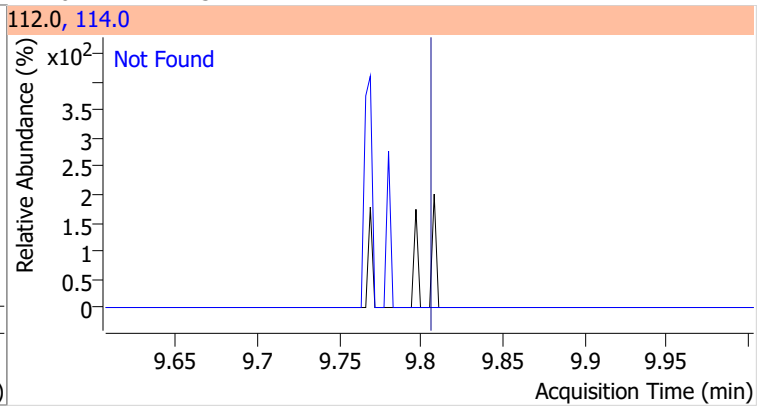
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2-Dibromoethane	N.D.	9.31	109.0	94.5



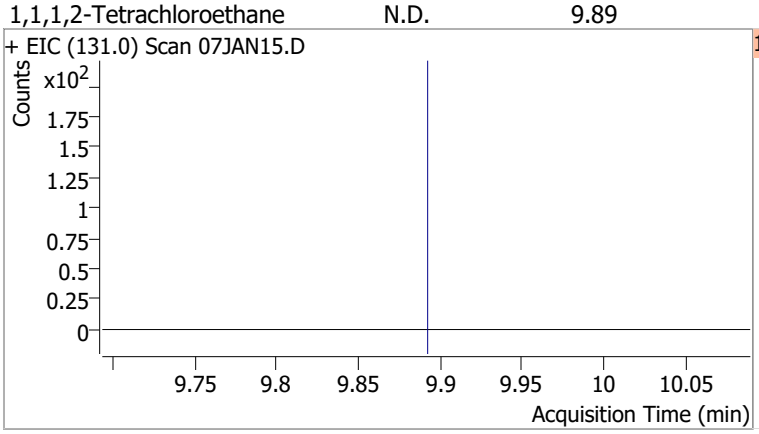


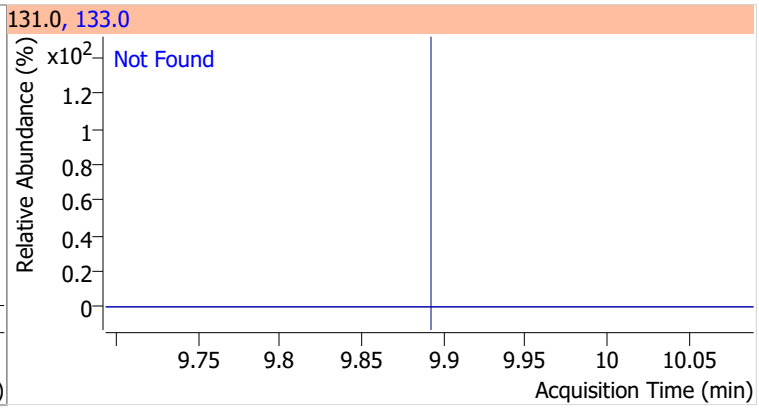
Compound	Conc.	Exp RT	QIon	Exp Ratio
Chlorobenzene	N.D.	9.80	114.0	32.1



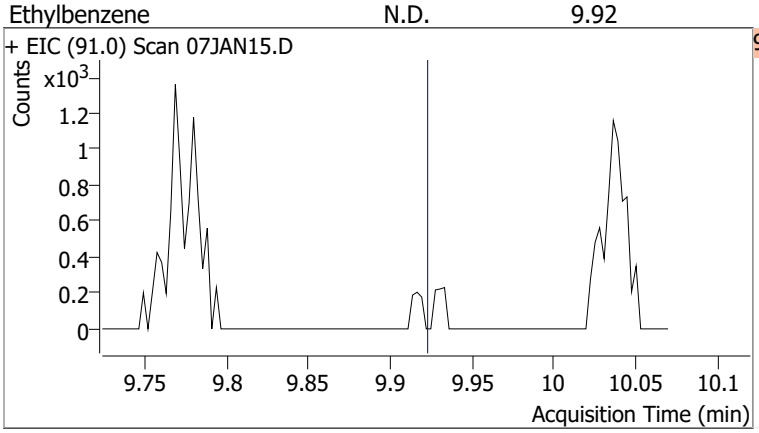


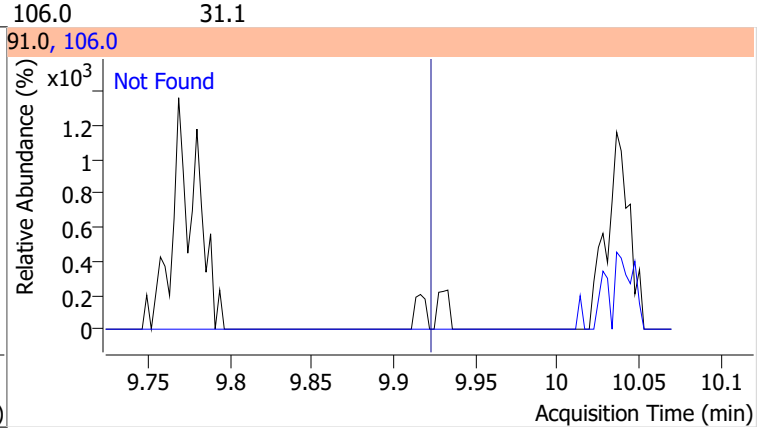
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,1,2-Tetrachloroethane	N.D.	9.89	133.0	98.6



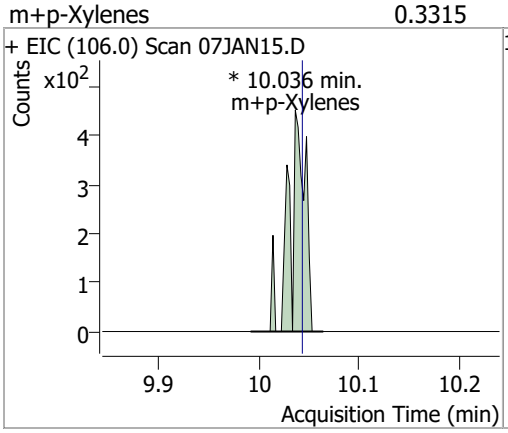


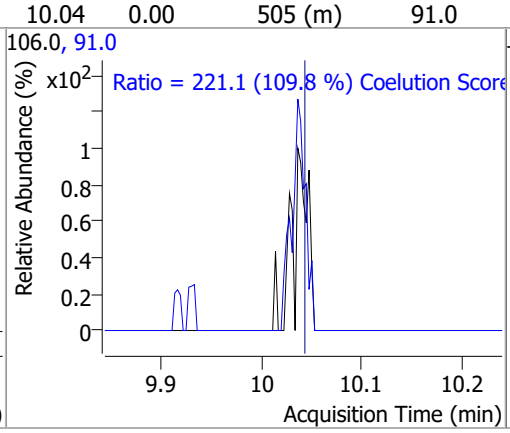
Compound	Conc.	Exp RT	QIon	Exp Ratio
Ethylbenzene	N.D.	9.92	106.0	31.1

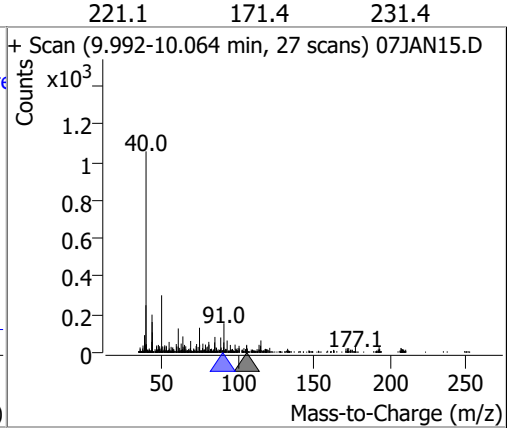




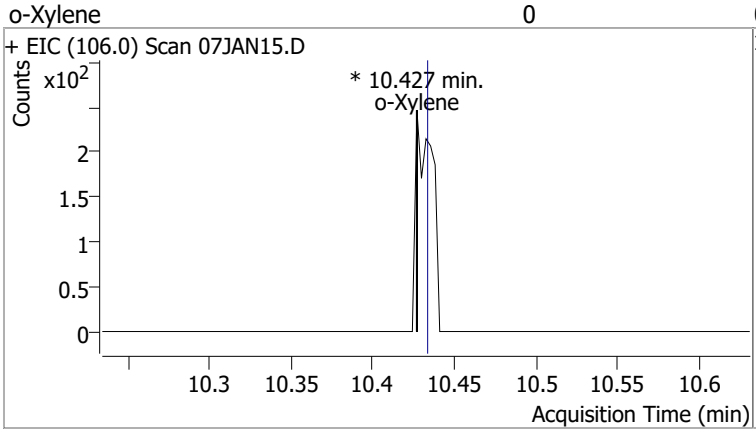
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	0.3315	10.04	0.00	505 (m)	91.0	221.1	171.4	231.4

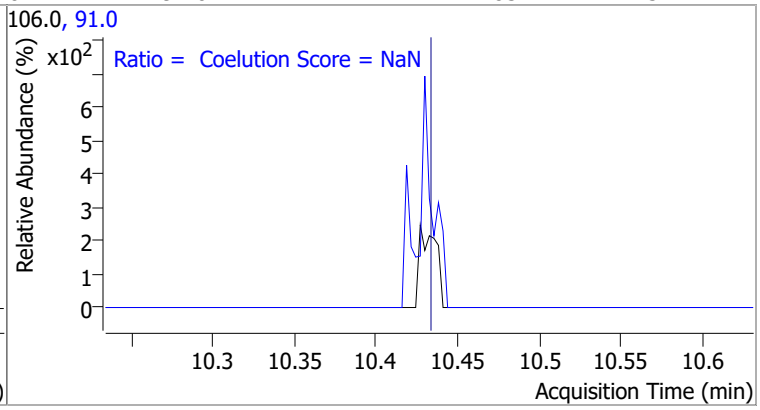




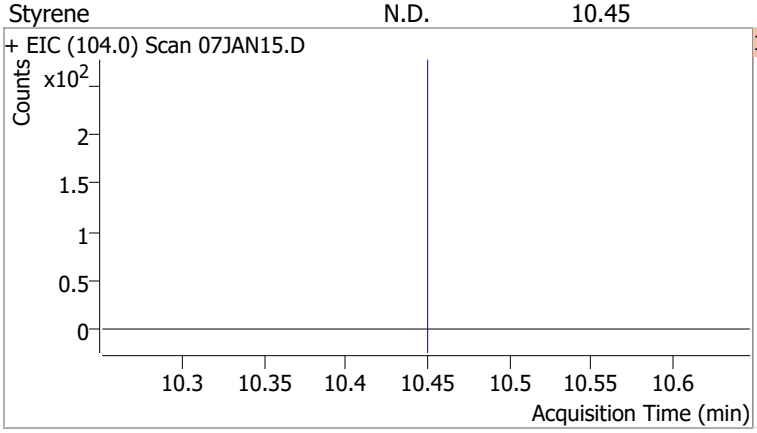


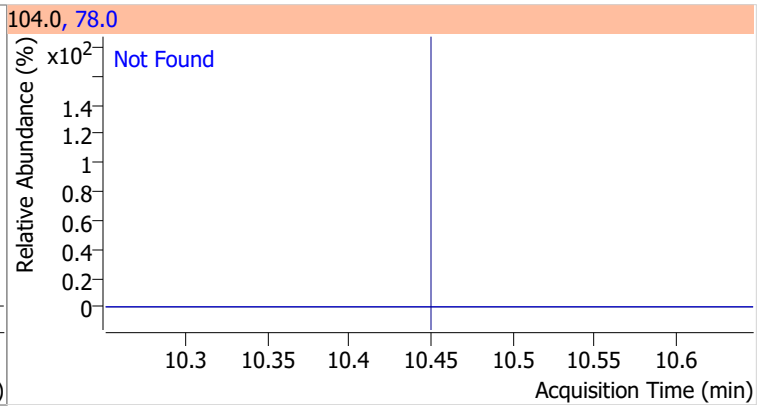
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene		0		0	91.0		183.1	243.1



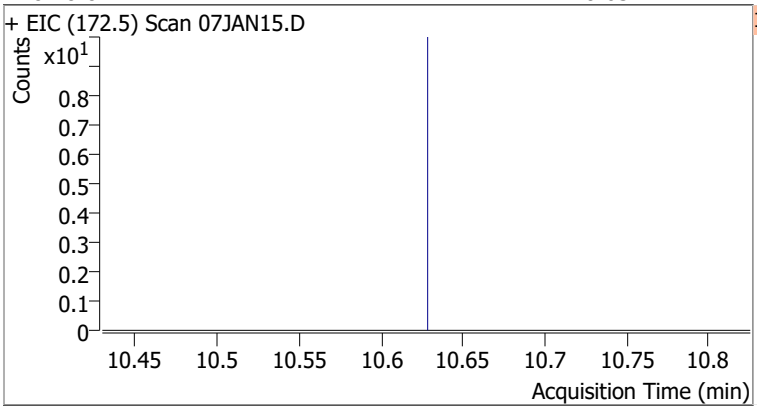


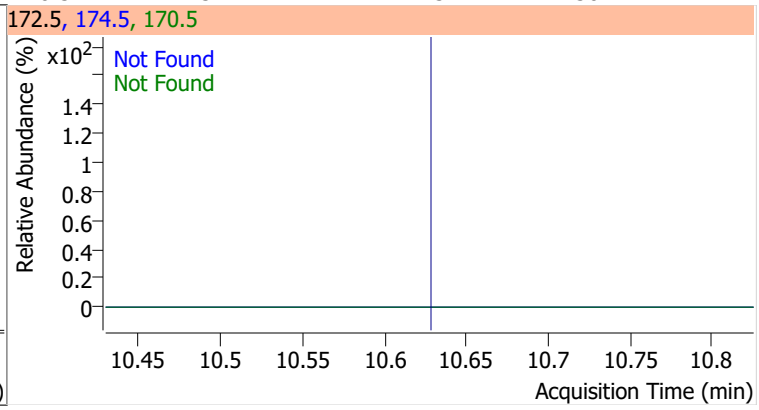
Compound	Conc.	Exp RT	QIon	Exp Ratio
Styrene	N.D.	10.45	78.0	49.6



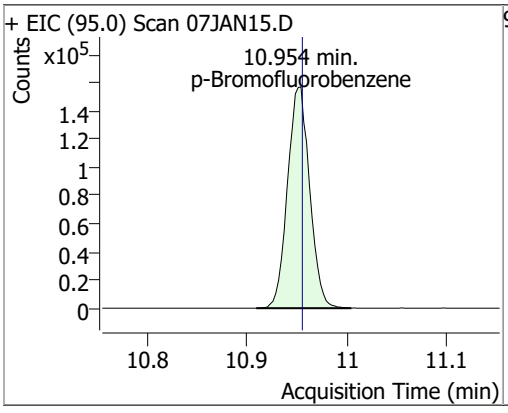


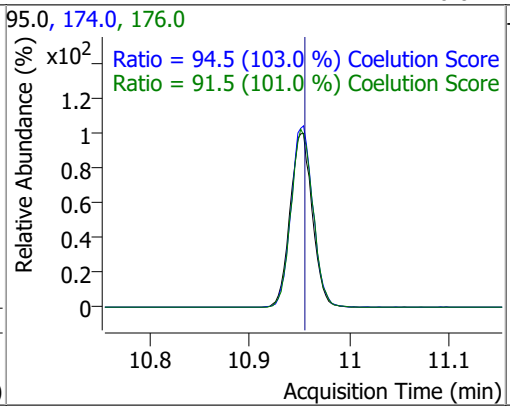
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromoform	N.D.	10.63	170.5	52.1	174.5	50.1

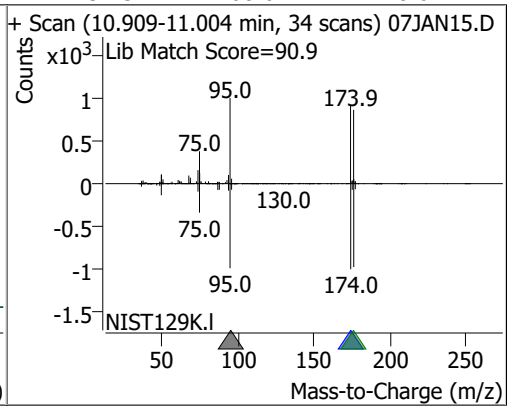




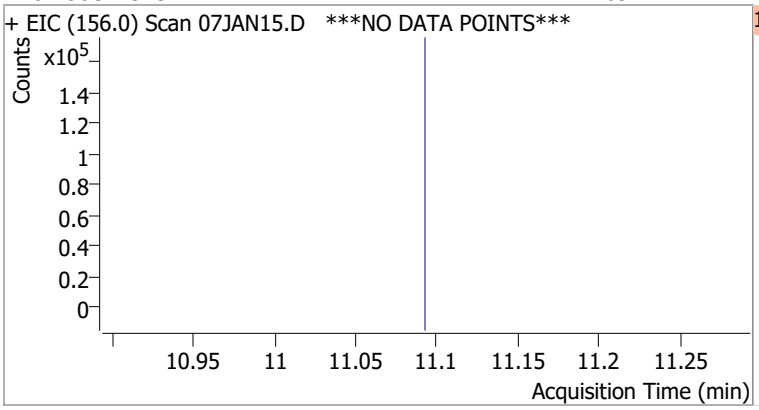
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	269.8409	10.95	0.00	240345	174.0	94.5	61.7	121.7
					176.0	91.5	60.6	120.6

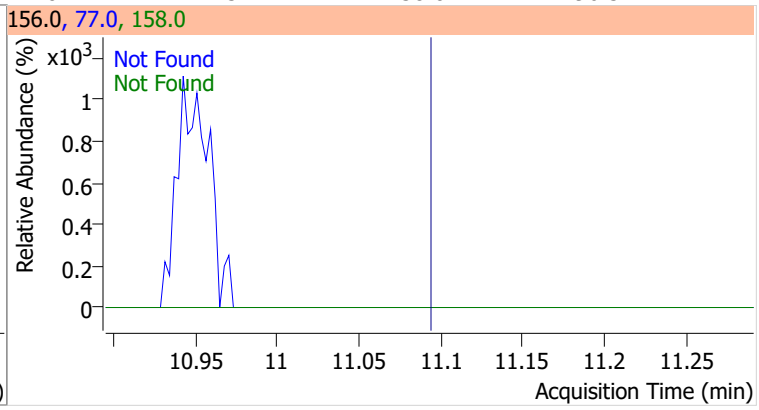




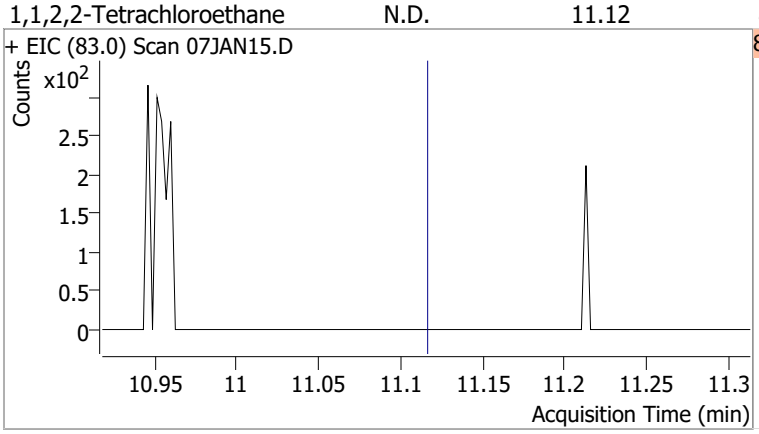


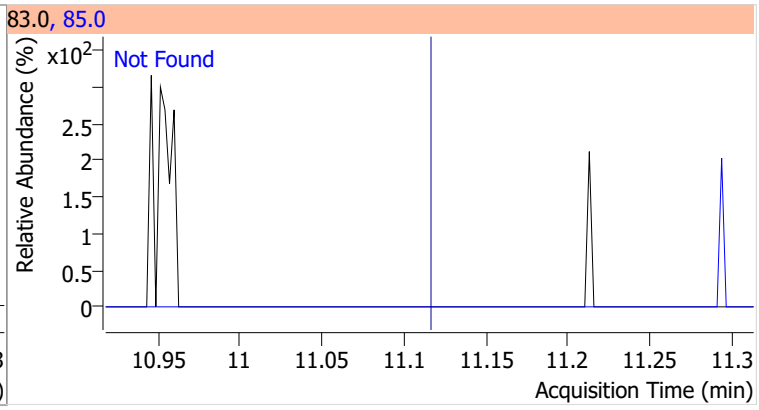
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
Bromobenzene	N.D.	11.09	77.0	145.7	158.0	96.5



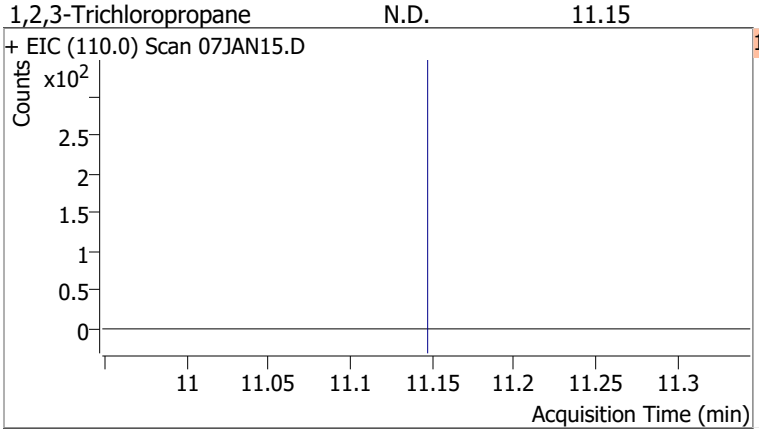


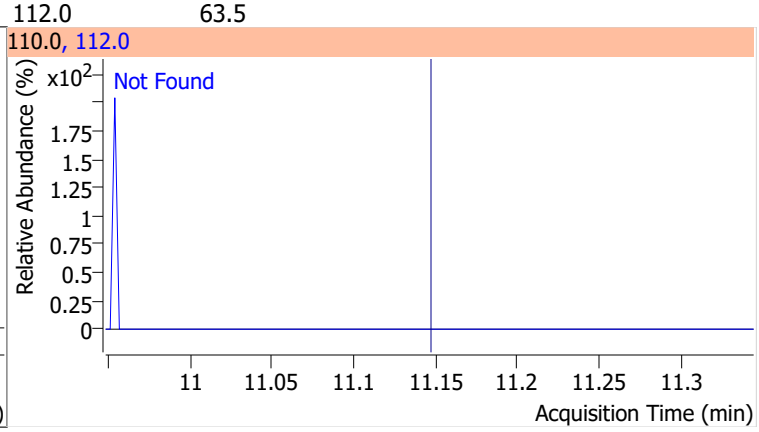
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,1,2,2-Tetrachloroethane	N.D.	11.12	85.0	66.2



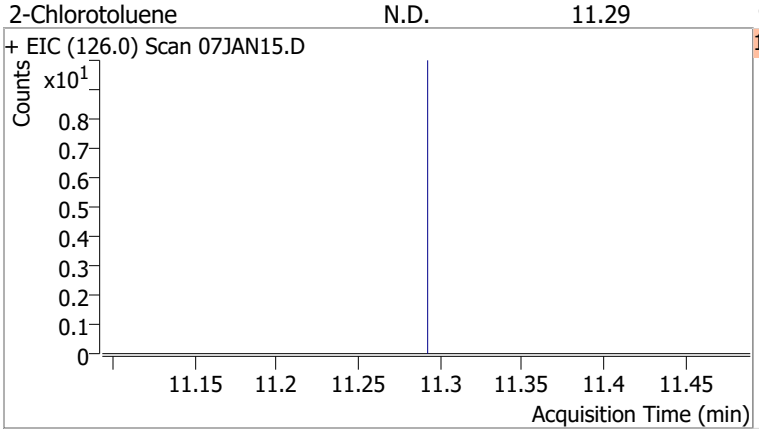


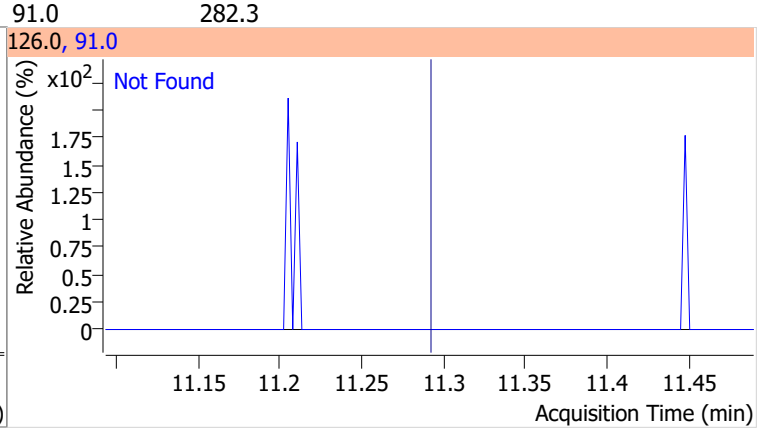
Compound	Conc.	Exp RT	QIon	Exp Ratio
1,2,3-Trichloropropane	N.D.	11.15	112.0	63.5



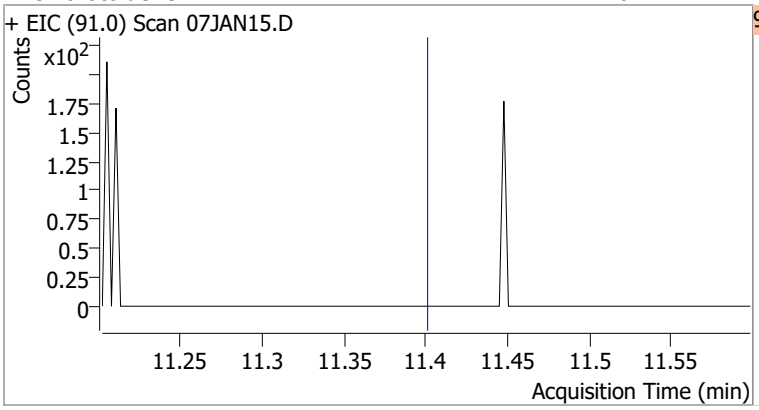


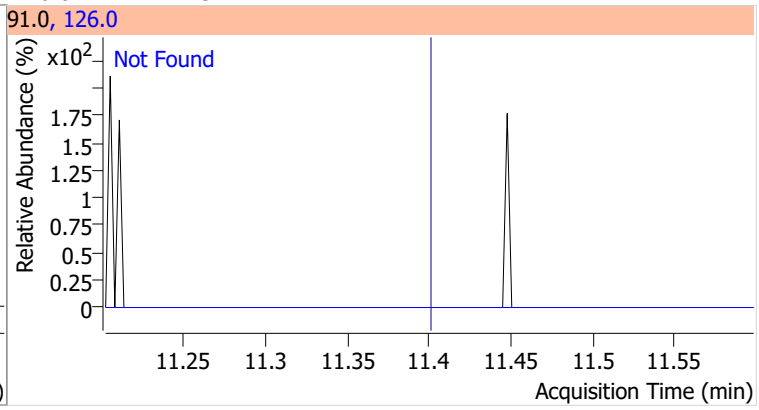
Compound	Conc.	Exp RT	QIon	Exp Ratio
2-Chlorotoluene	N.D.	11.29	91.0	282.3



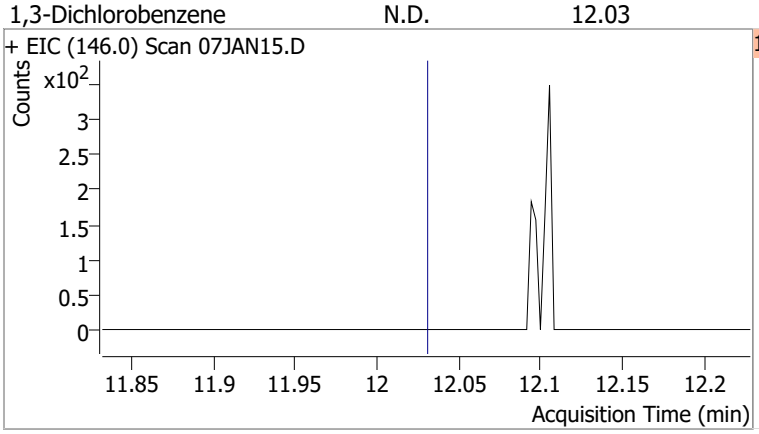


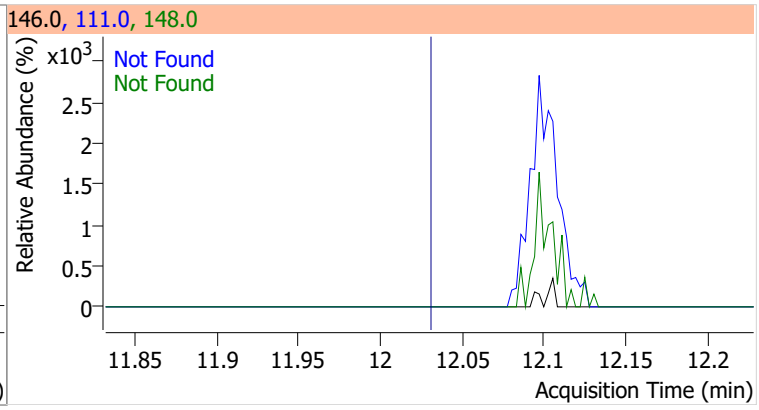
Compound	Conc.	Exp RT	QIon	Exp Ratio
4-Chlorotoluene	N.D.	11.40	126.0	31.7



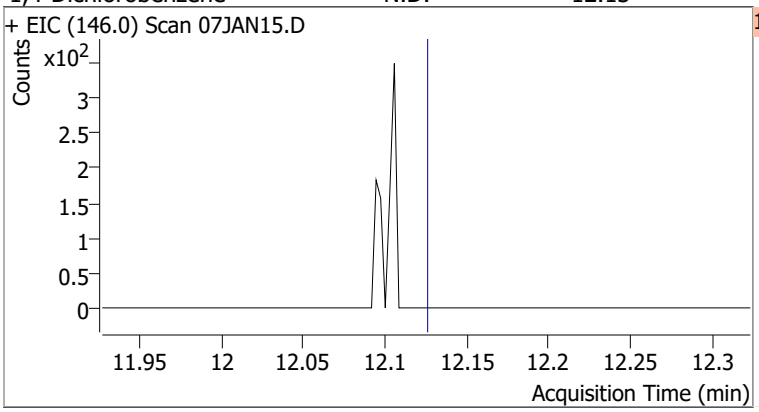


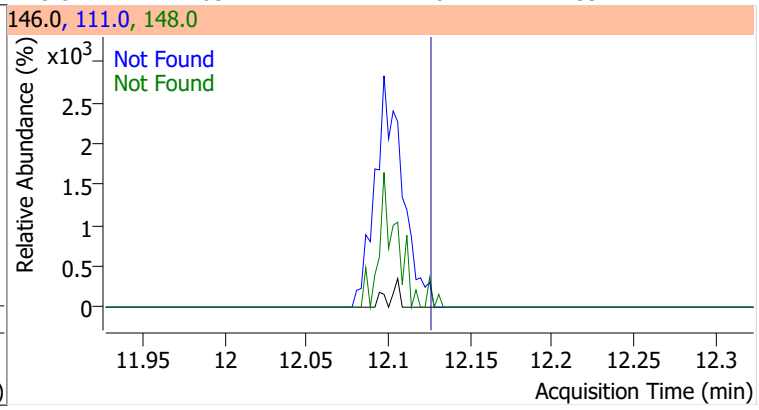
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,3-Dichlorobenzene	N.D.	12.03	148.0	63.6	111.0	39.8



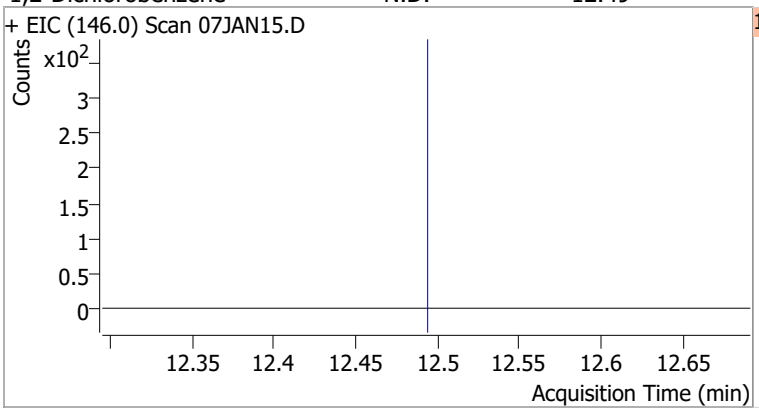


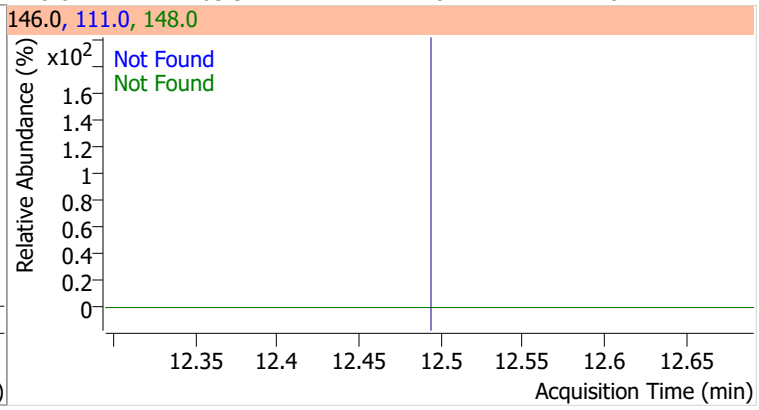
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,4-Dichlorobenzene	N.D.	12.13	148.0	63.1	111.0	39.1





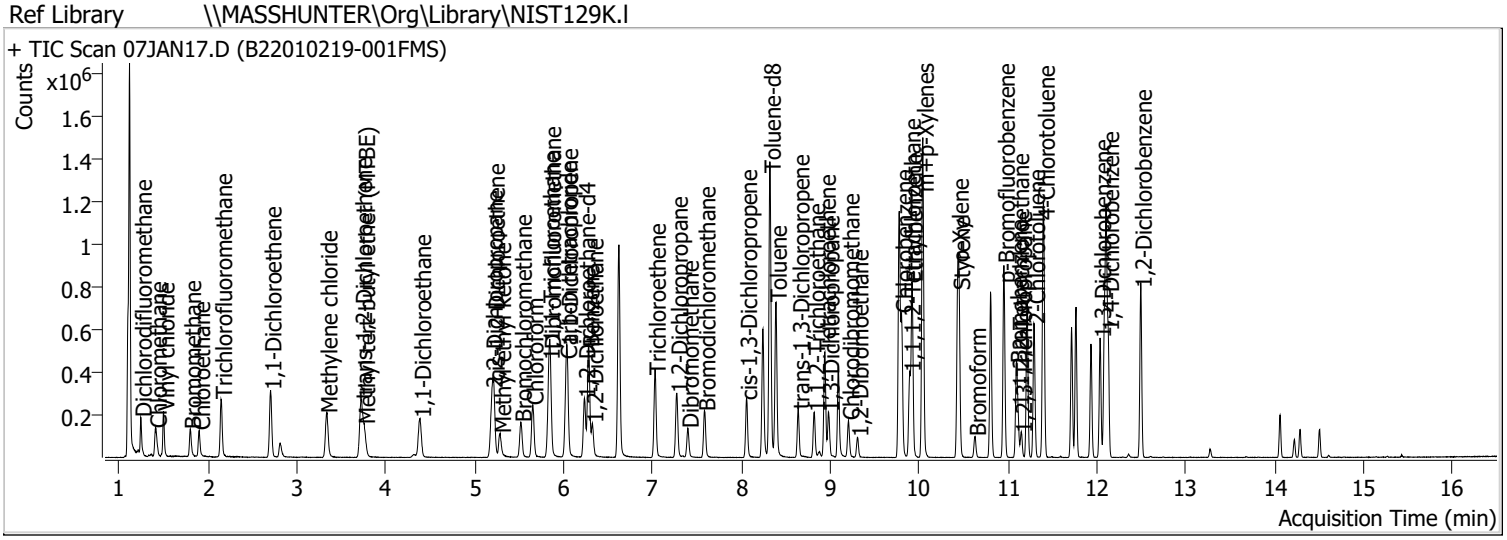
Compound	Conc.	Exp RT	QIon	Exp Ratio	QIon	Exp Ratio
1,2-Dichlorobenzene	N.D.	12.49	148.0	63.9	111.0	41.0





Quantitation Results Report (QT Reviewed)

Data File	07JAN17.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 5:07:38 PM
Sample Name	B22010219-001FMS	Instrument	VOA5975C
Vial	17	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

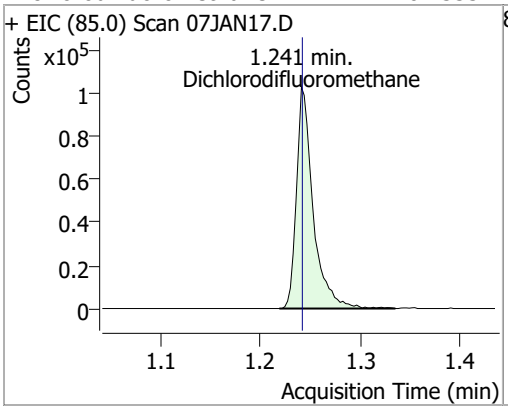


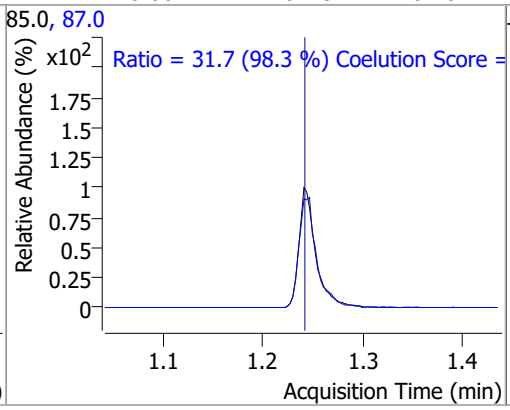
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.620	96.0	855146	250.0000	ng	-0.003
M Chlorobenzene-d5	9.772	82.0	328555	250.0000	ng	0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	264037	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.848	113.0	220360	273.5236	ng	0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 109.41%		
S 1,2-Dichloroethane-d4	6.233	67.0	96662	277.7836	ng	0.000
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 111.11%		
S Toluene-d8	8.319	98.0	860434	271.7626	ng	0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 108.71%		
S p-Bromofluorobenzene	10.951	95.0	265093	274.0538	ng	-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 109.62%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.241	85.0	120279	107.3334	ng	99
T Chloromethane	1.411	50.0	152071	111.8049	ng	99
T Vinyl chloride	1.498	62.0	150050	122.6032	ng	97
T Bromomethane	1.799	96.0	66210	120.9860	ng	98
T Chloroethane	1.899	64.0	85441	141.0182	ng	97
T Trichlorofluoromethane	2.145	101.0	189149	124.5150	ng	98
T 1,1-Dichloroethene	2.702	96.0	112908	131.0796	ng	98
T Methylene chloride	3.333	49.0	154851	121.9493	ng	98
T trans-1,2-Dichloroethene	3.720	96.0	114095	129.8323	ng	98
T Methyl tert-butyl ether (MTBE)	3.757	73.0	159167	140.1252	ng	98
T 1,1-Dichloroethane	4.384	63.0	221763	135.5713	ng	98
T 2,2-Dichloropropane	5.193	77.0	161563	131.8131	ng	99
T cis-1,2-Dichloroethene	5.212	96.0	116966	131.2799	ng	97
T Methyl ethyl ketone	5.285	43.0	151656	1256.6341	ng	100
T Bromochloromethane	5.516	128.0	45934	124.4477	ng	94
T Chloroform	5.650	83.0	197751	121.4742	ng	98

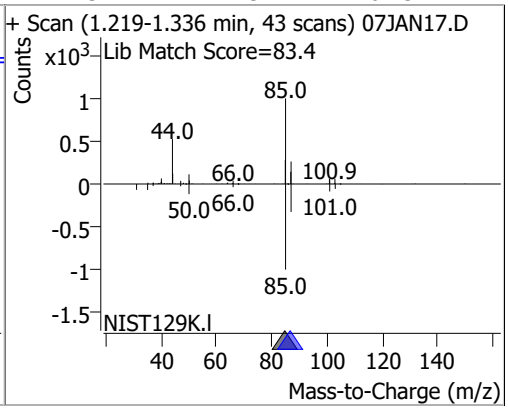
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.831	97.0	190350	124.7685	ng	98
T Carbon tetrachloride	6.026	117.0	182925	121.6947	ng	98
T 1,1-Dichloropropene	6.038	75.0	157497	121.4151	ng	100
T Benzene	6.280	78.0	443412	130.2309	ng	100
T 1,2-Dichloroethane	6.322	62.0	117773	127.8627	ng	98
T Trichloroethene	7.025	95.0	124303	125.4471	ng	98
T 1,2-Dichloropropane	7.273	63.0	109245	125.3364	ng	100
T Dibromomethane	7.396	93.0	46329	125.7798	ng	97
T Bromodichloromethane	7.583	83.0	130361	128.2418	ng	99
T cis-1,3-Dichloropropene	8.054	75.0	139124	121.0494	ng	99
T Toluene	8.386	92.0	276475	129.2716	ng	99
T trans-1,3-Dichloropropene	8.639	75.0	108385	132.4832	ng	97
T 1,1,2-Trichloroethane	8.818	83.0	52652	123.5591	ng	96
T Tetrachloroethene	8.938	163.8	106788	122.3903	ng	99
T 1,3-Dichloropropane	8.980	76.0	106245	126.7568	ng	96
T Chlorodibromomethane	9.203	129.0	83355	125.1595	ng	96
T 1,2-Dibromoethane	9.303	107.0	58167	124.8389	ng	100
T Chlorobenzene	9.799	112.0	300347	128.2720	ng	100
T 1,1,1,2-Tetrachloroethane	9.892	131.0	101471	123.9720	ng	96
T Ethylbenzene	9.917	91.0	515970	127.0574	ng	100
T m+p-Xylenes	10.039	106.0	399658	253.2483	ng	100
T o-Xylene	10.433	106.0	183667	130.7338	ng	98
T Styrene	10.449	104.0	299523	132.4202	ng	100
T Bromoform	10.628	172.5	43824	129.7036	ng	98
T Bromobenzene	11.091	156.0	114173	133.6149	ng	99
T 1,1,2,2-Tetrachloroethane	11.110	83.0	64543	131.2329	ng	97
T 1,2,3-Trichloropropane	11.144	110.0	16295	123.8246	ng	98
T 2-Chlorotoluene	11.291	126.0	111645	131.3133	ng	99
T 4-Chlorotoluene	11.400	91.0	369533	133.3048	ng	99
T 1,3-Dichlorobenzene	12.033	146.0	204772	131.3970	ng	99
T 1,4-Dichlorobenzene	12.125	146.0	206411	129.8964	ng	99
T 1,2-Dichlorobenzene	12.493	146.0	165895	125.9590	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

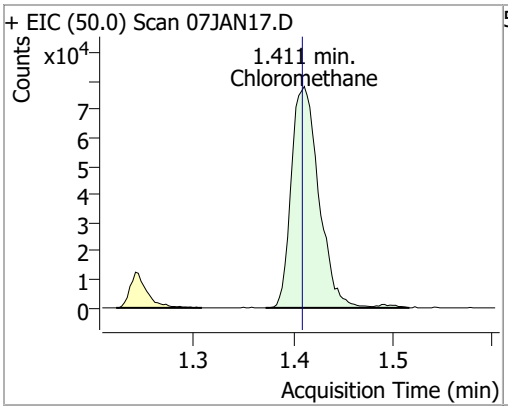
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	107.3334	1.24	0.00	120279	87.0	31.7	2.3	62.3

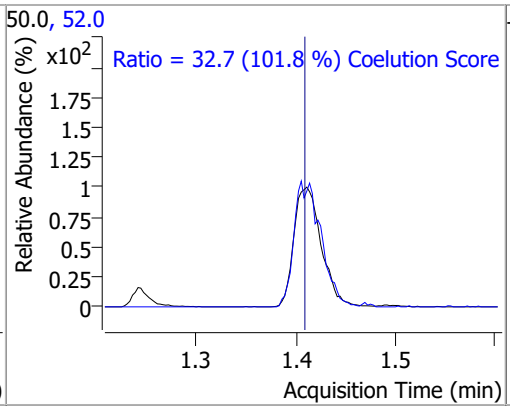


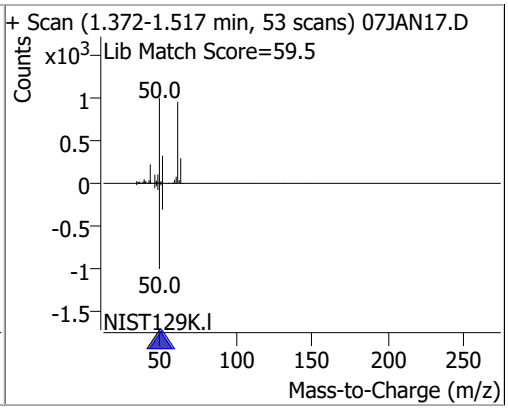




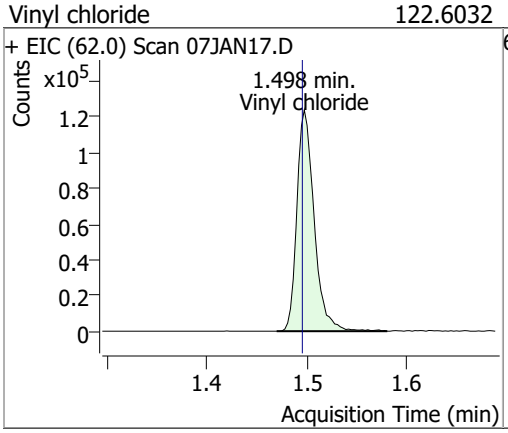
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	111.8049	1.41	0.00	152071	52.0	32.7	2.1	62.1

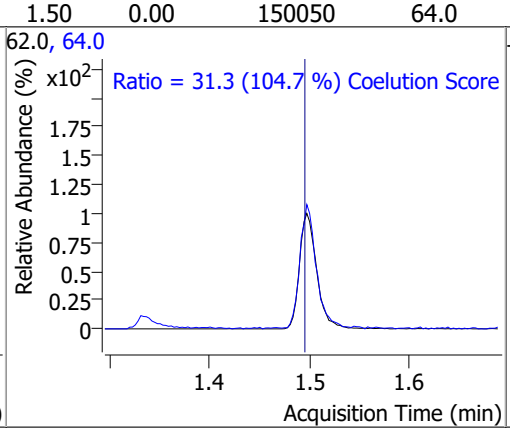


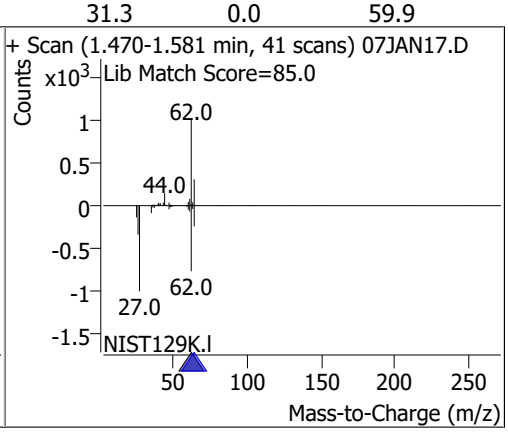




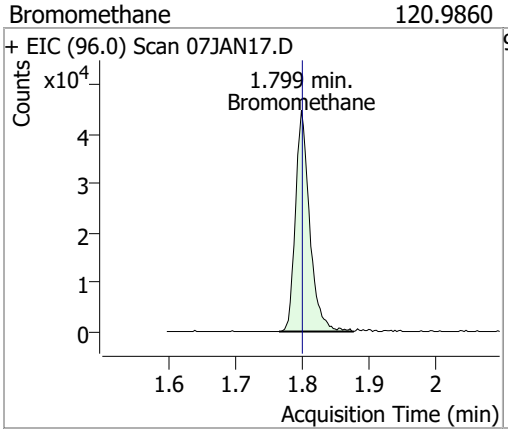
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	122.6032	1.50	0.00	150050	64.0	31.3	0.0	59.9

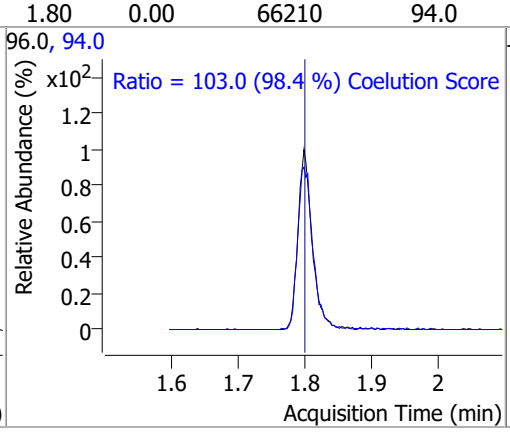


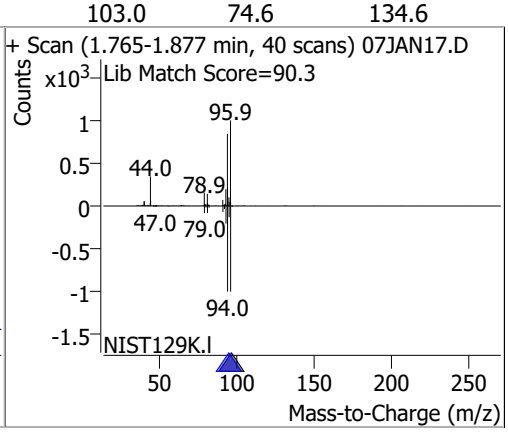




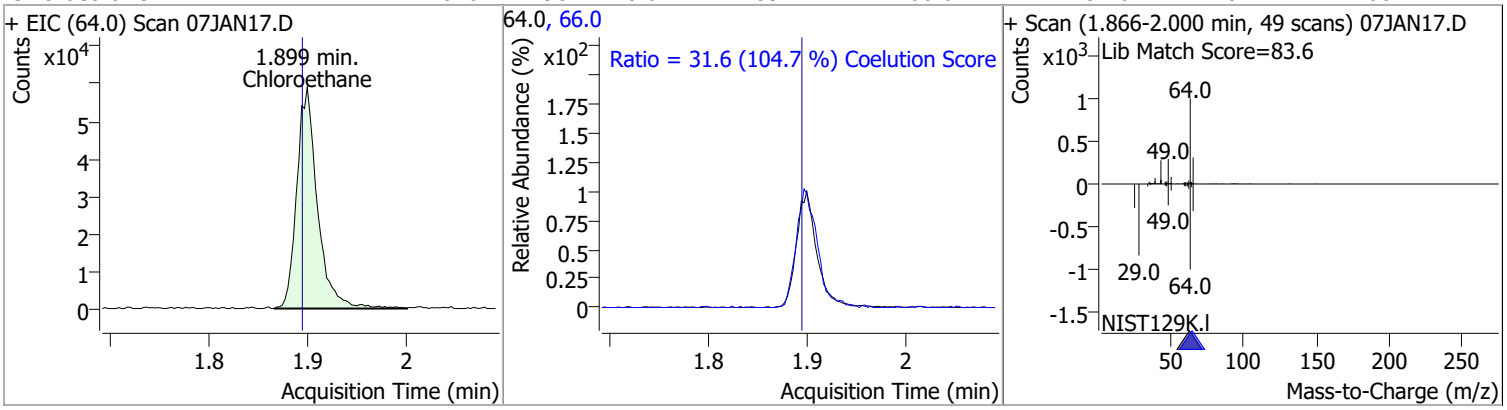
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	120.9860	1.80	0.00	66210	94.0	103.0	74.6	134.6



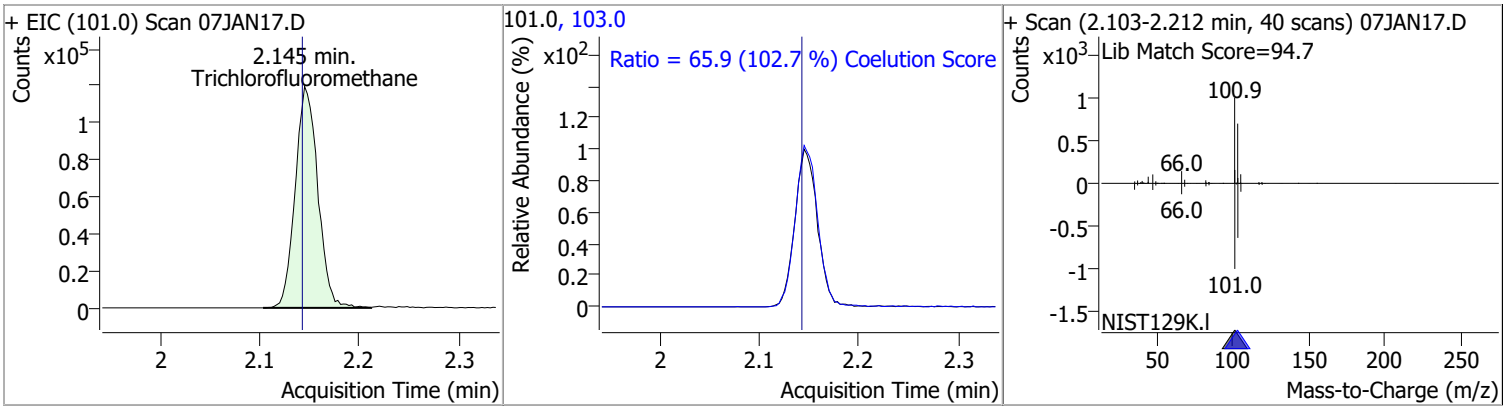




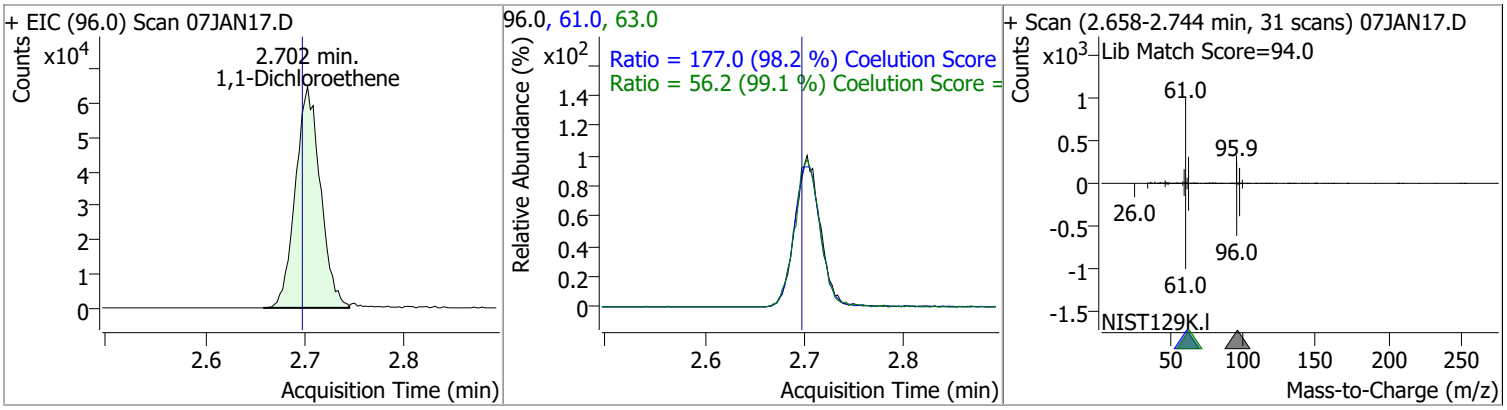
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	141.0182	1.90	0.01	85441	66.0	31.6	0.1	60.1



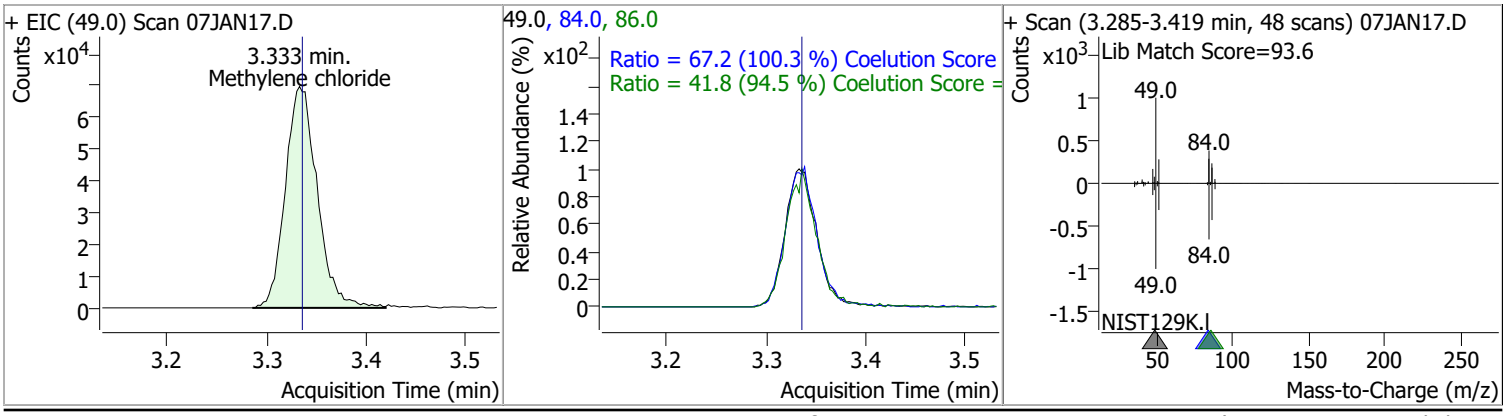
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	124.5150	2.14	0.00	189149	103.0	65.9	34.2	94.2



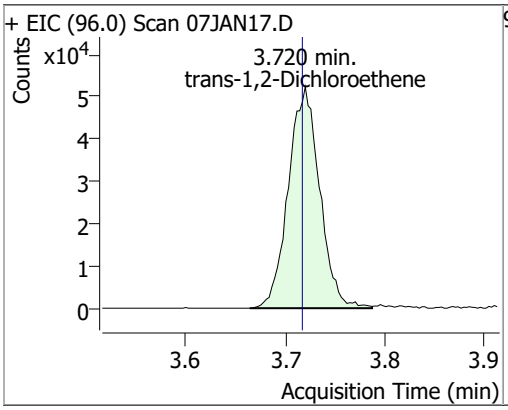
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	131.0796	2.70	0.01	112908	61.0	177.0	150.3	210.3
					63.0	56.2	26.7	86.7

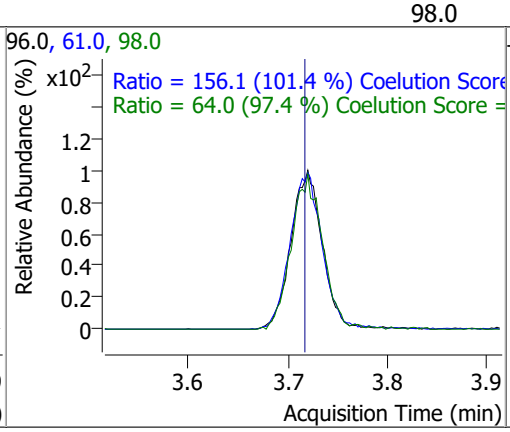


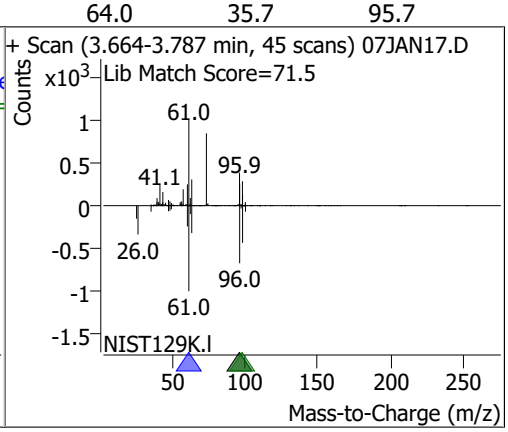
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	121.9493	3.33	0.00	154851	84.0	67.2	36.9	96.9
					86.0	41.8	14.3	74.3



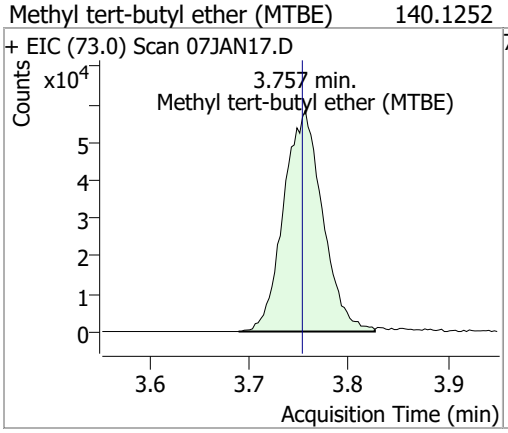
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	129.8323	3.72	0.00	114095	61.0	156.1	123.9	183.9
					98.0	64.0	35.7	95.7

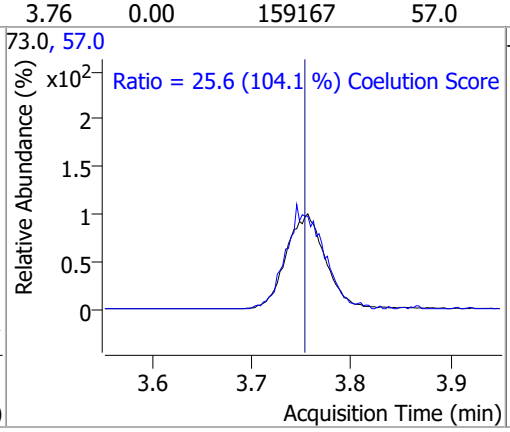


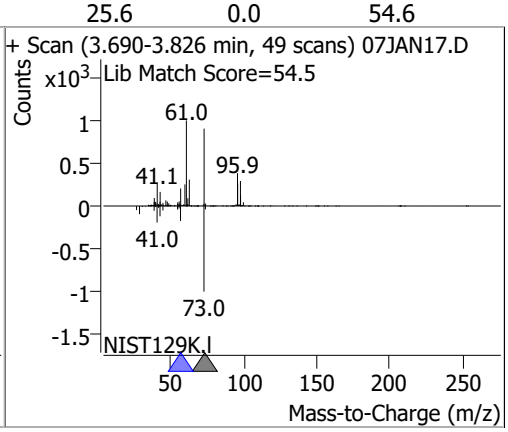




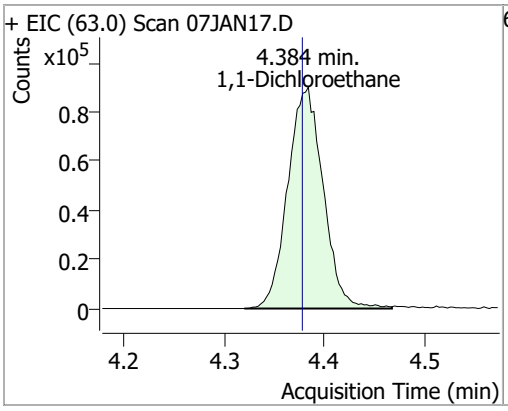
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	140.1252	3.76	0.00	159167	57.0	25.6	0.0	54.6

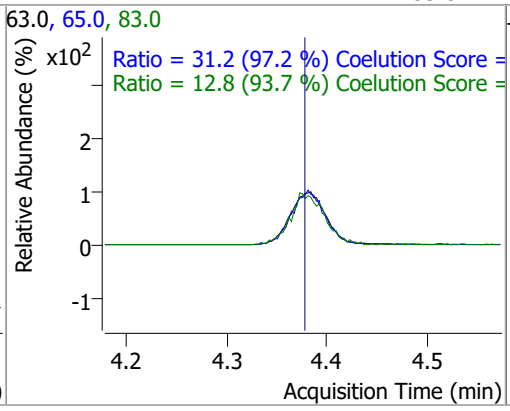


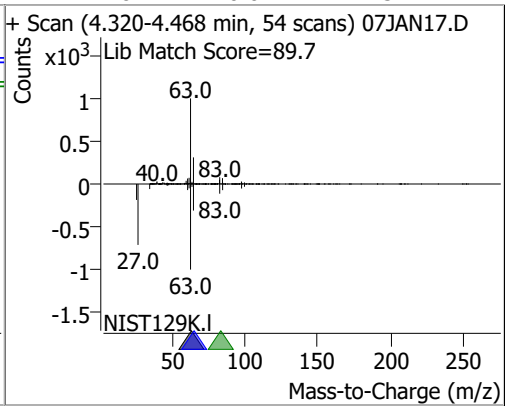




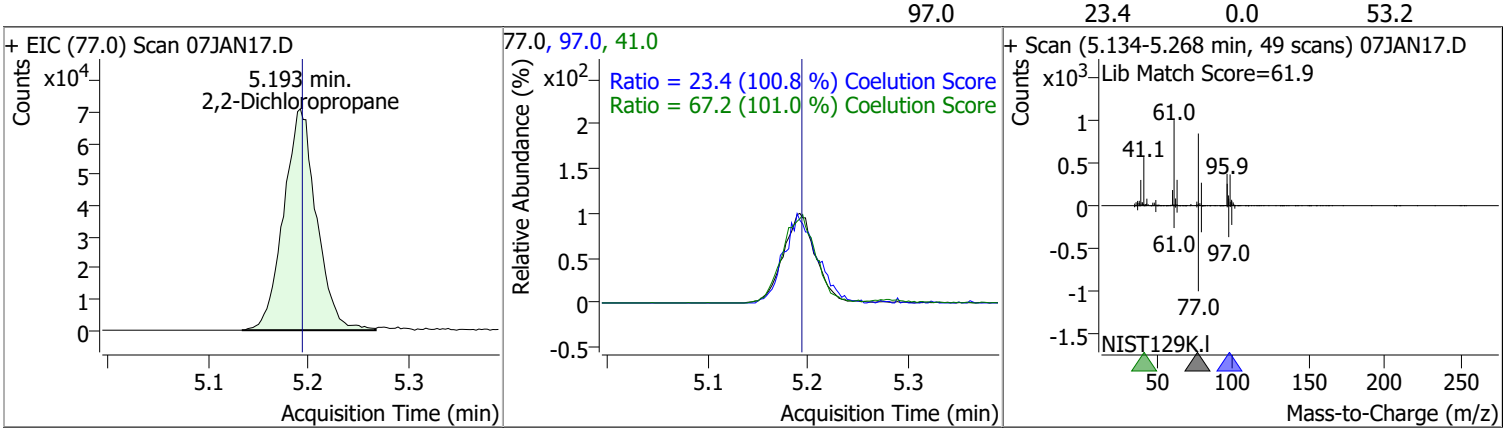
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	135.5713	4.38	0.01	221763	65.0	31.2	2.1	62.1
					83.0	12.8	0.0	43.7



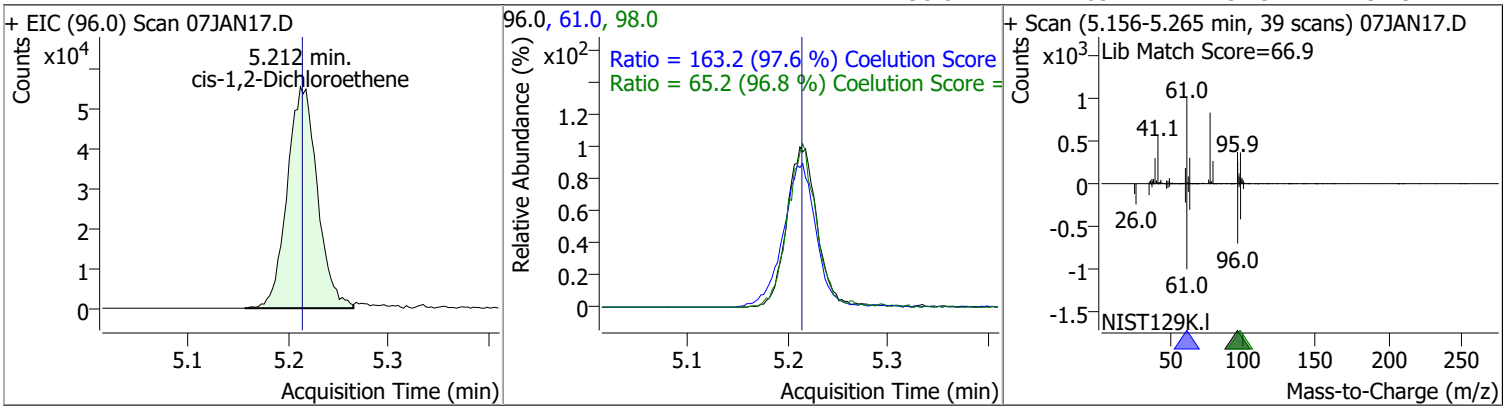




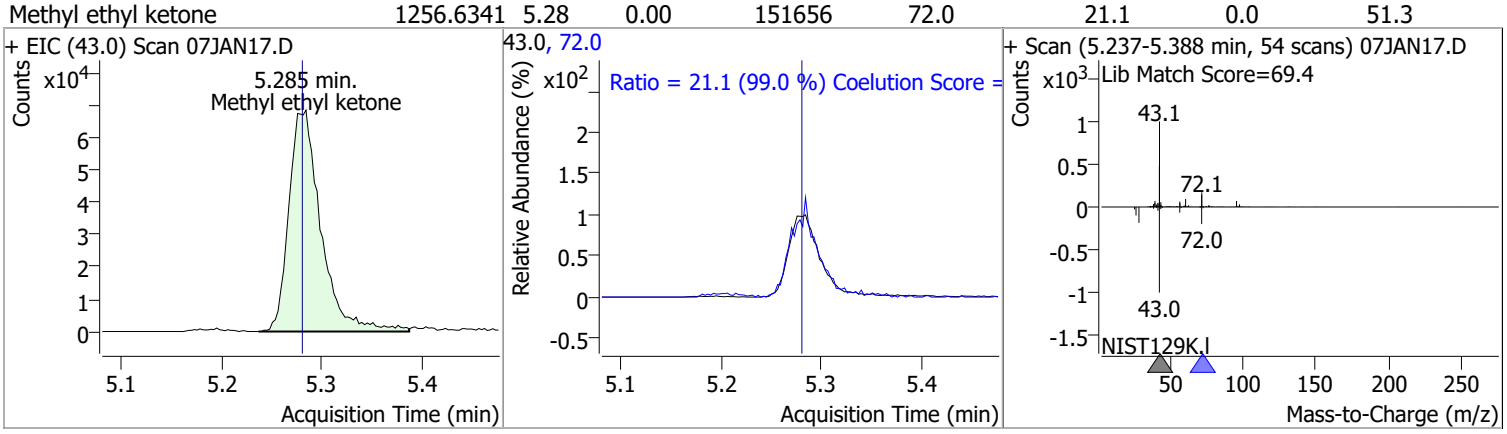
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	131.8131	5.19	0.00	161563	41.0	67.2	36.5	96.5
					97.0	23.4	0.0	53.2



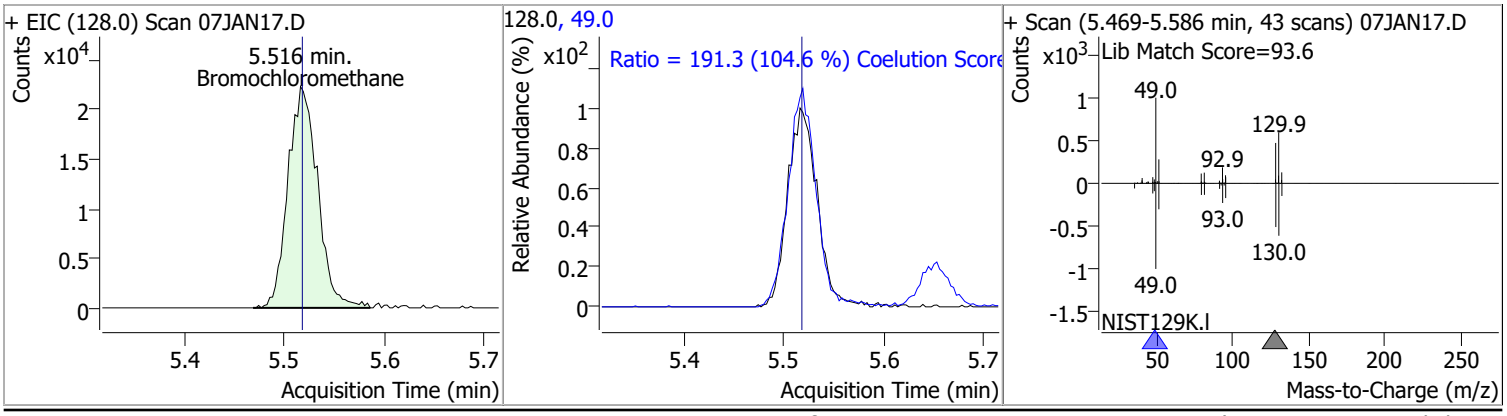
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	131.2799	5.21	0.00	116966	61.0	163.2	137.2	197.2
					98.0	65.2	37.3	97.3



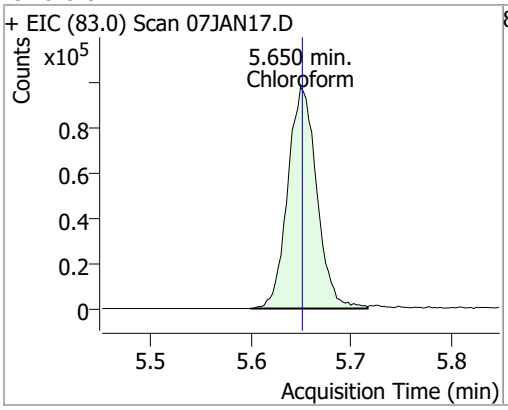
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1256.6341	5.28	0.00	151656	72.0	21.1	0.0	51.3

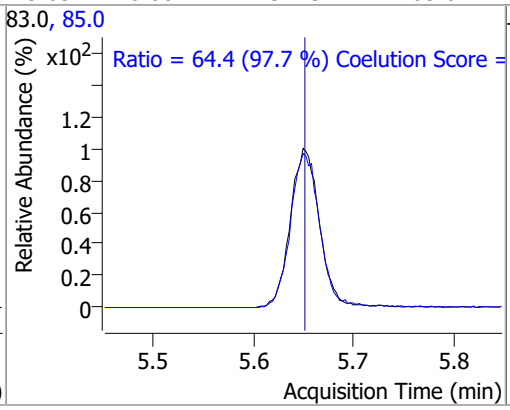


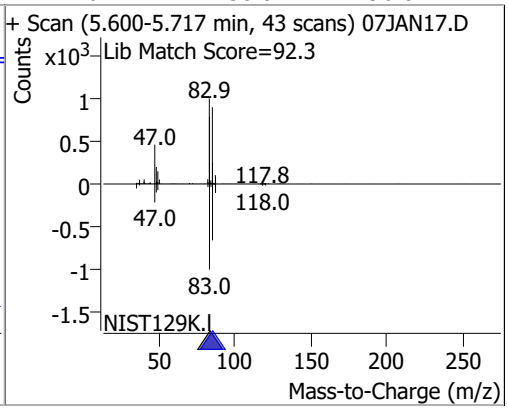
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	124.4477	5.52	0.00	45934	49.0	191.3	152.9	212.9



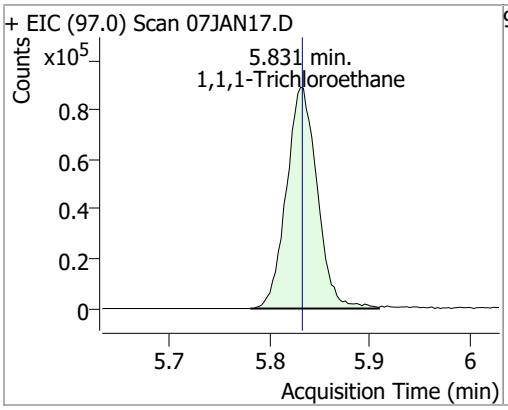
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	121.4742	5.65	0.00	197751	85.0	64.4	36.0	96.0

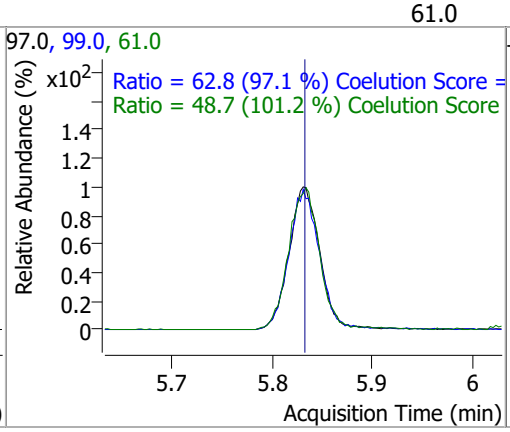


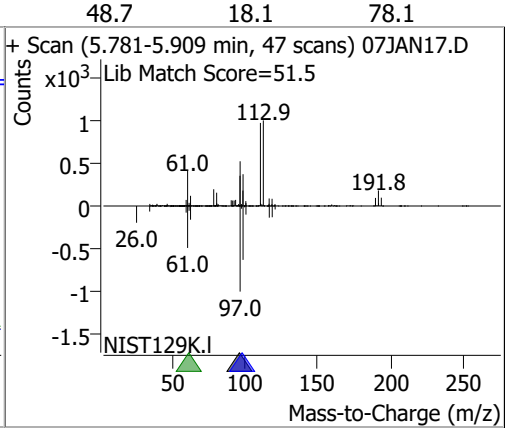




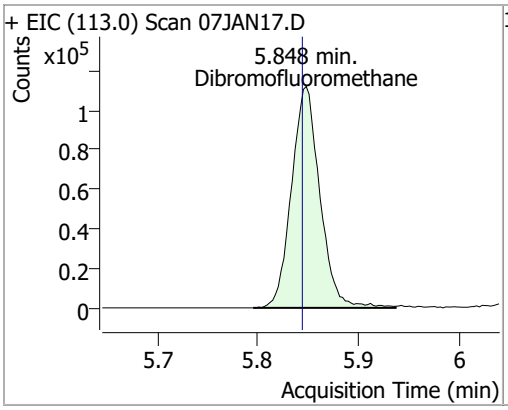
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	124.7685	5.83	0.00	190350	99.0	62.8	34.7	94.7
					61.0	48.7	18.1	78.1

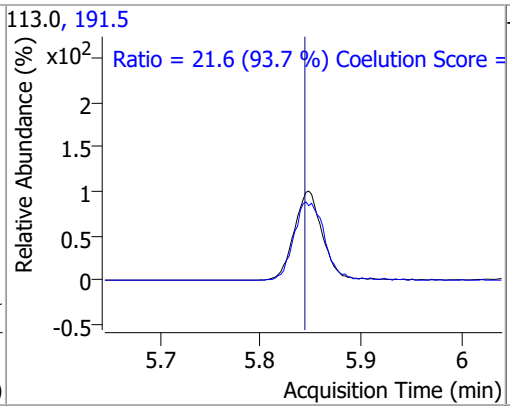


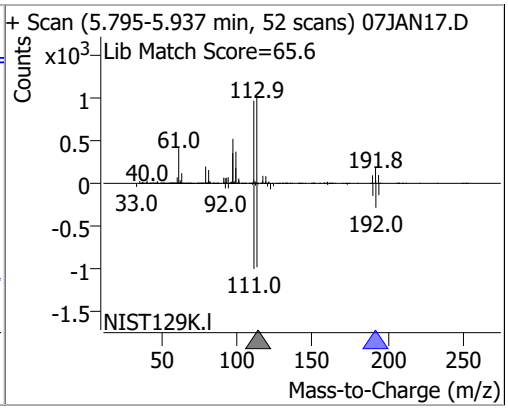




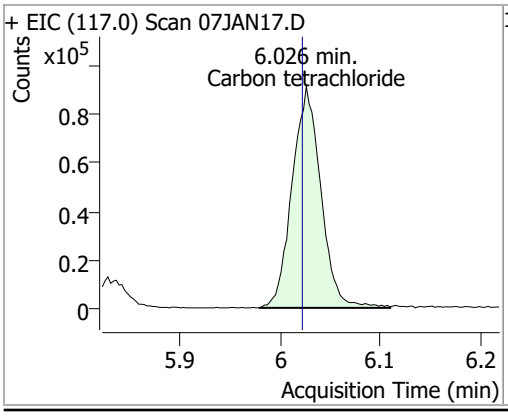
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	273.5236	5.85	0.00	220360	191.5	21.6	0.0	53.1

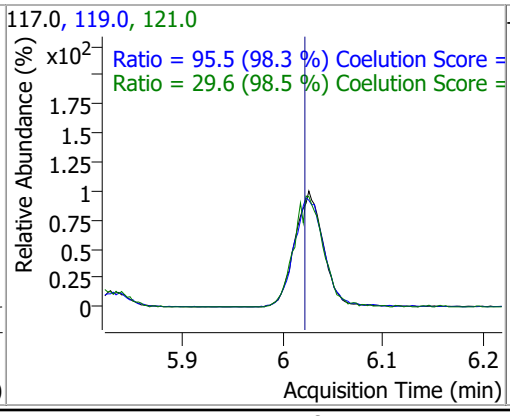


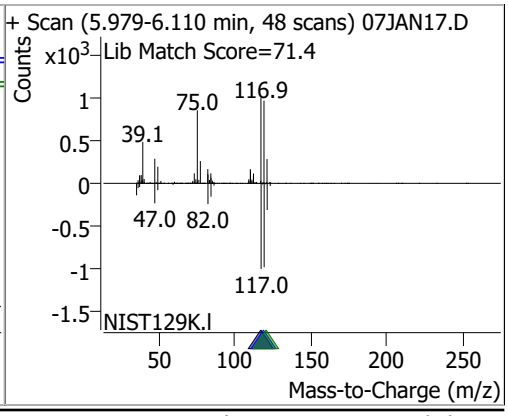




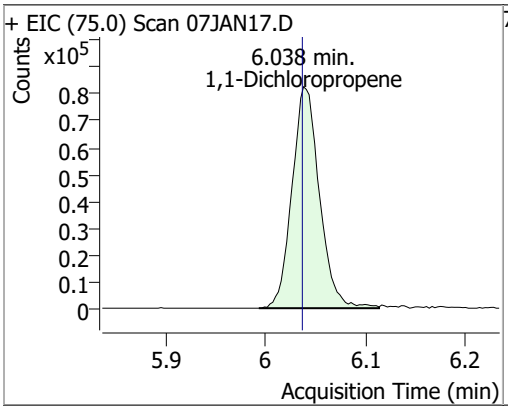
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	121.6947	6.03	0.00	182925	119.0	95.5	67.2	127.2
					121.0	29.6	0.1	60.1

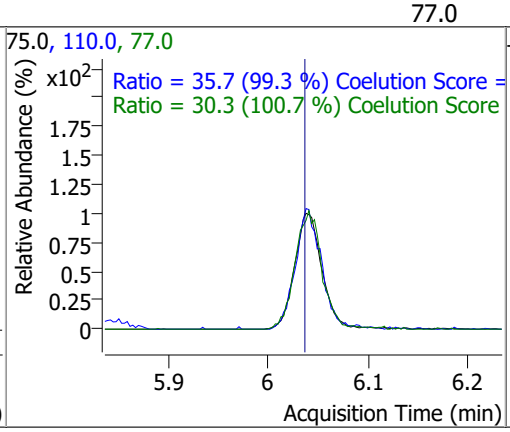


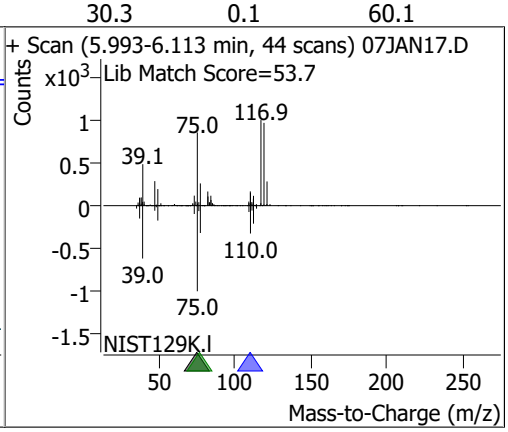




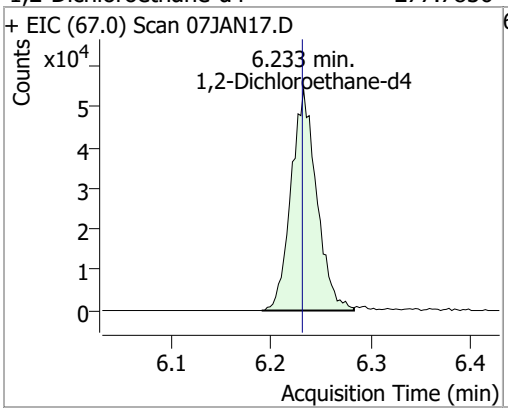
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	121.4151	6.04	0.00	157497	110.0	35.7	5.9	65.9
					77.0	30.3	0.1	60.1

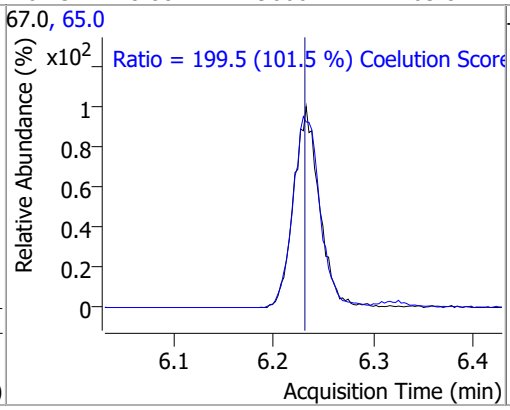


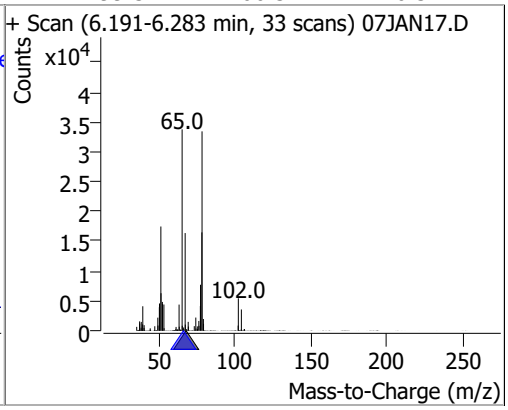




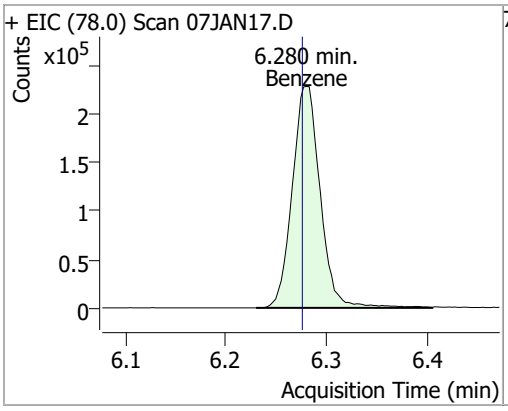
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	277.7836	6.23	0.00	96662	65.0	199.5	166.5	226.5

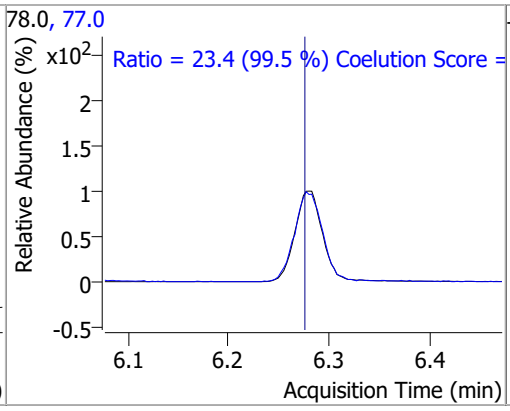


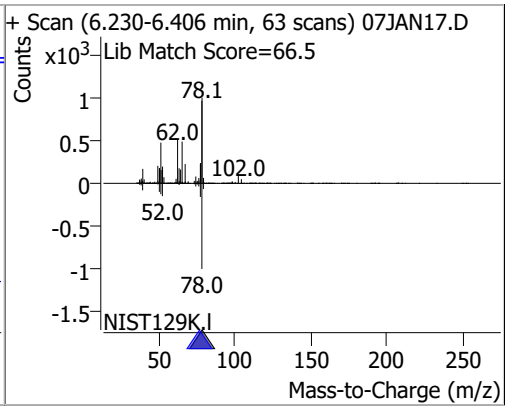




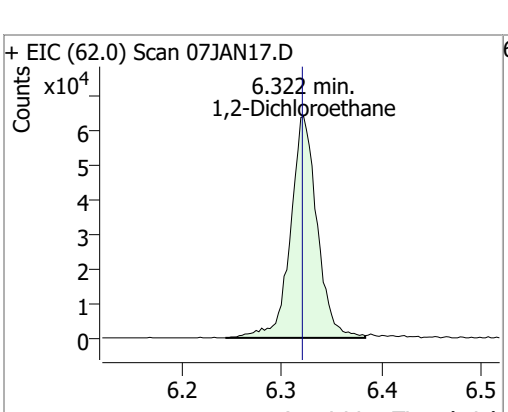
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	130.2309	6.28	0.00	443412	77.0	23.4	0.0	53.5

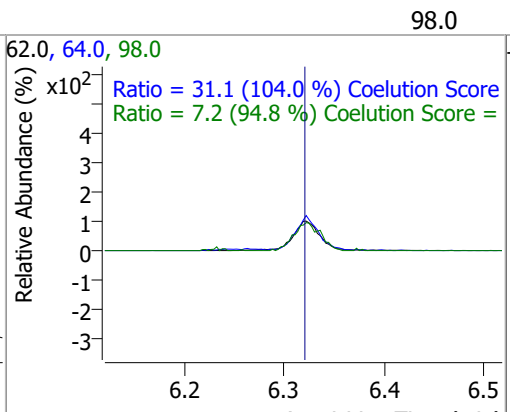


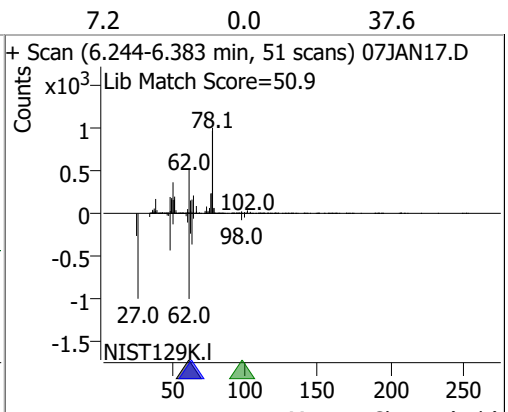




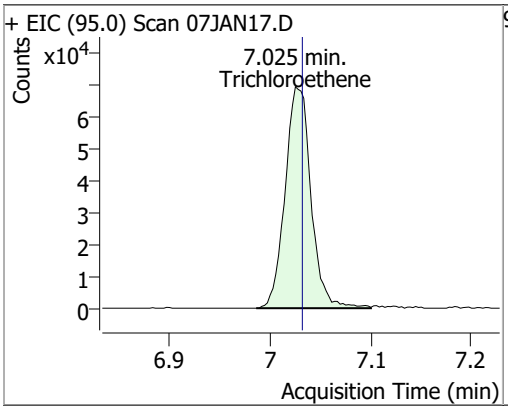
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	127.8627	6.32	0.00	117773	64.0	31.1	0.0	59.9
					98.0	7.2	0.0	37.6

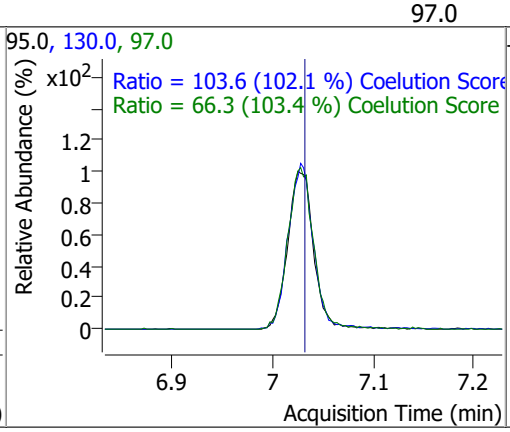


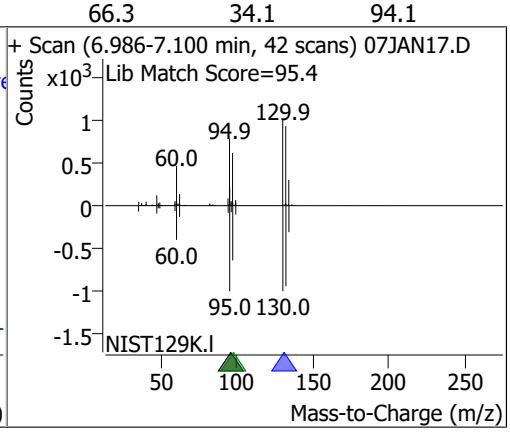




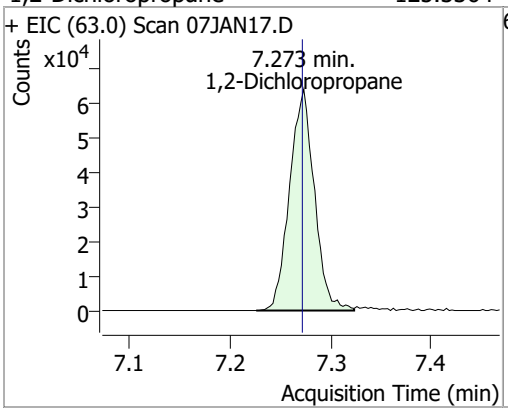
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	125.4471	7.02	-0.01	124303	130.0	103.6	71.5	131.5
					97.0	66.3	34.1	94.1

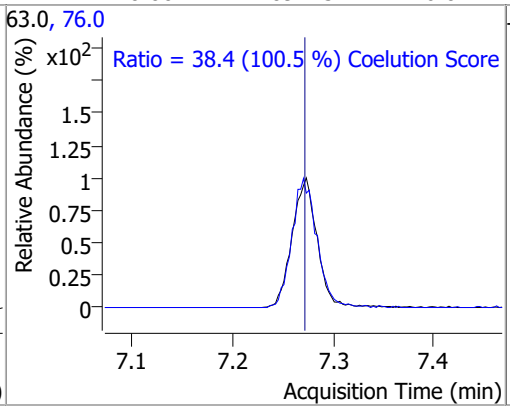


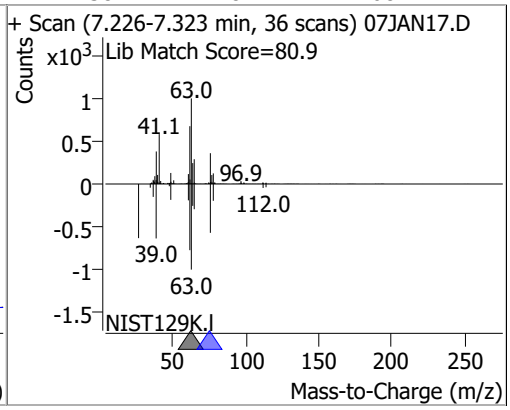




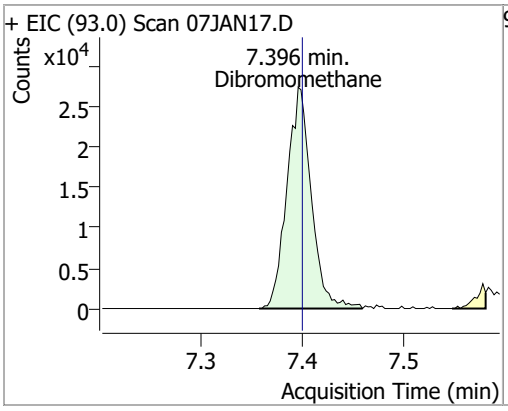
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	125.3364	7.27	0.00	109245	76.0	38.4	8.2	68.2

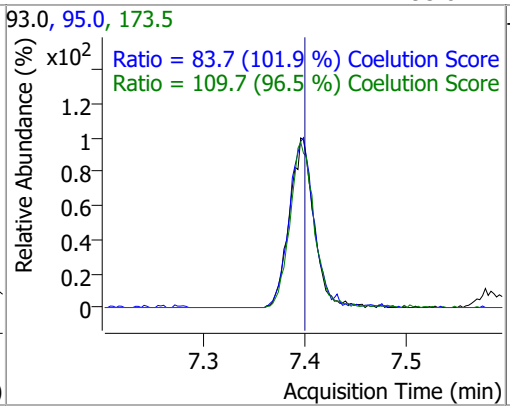


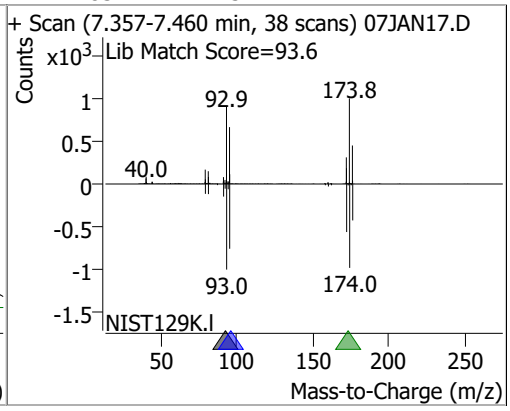




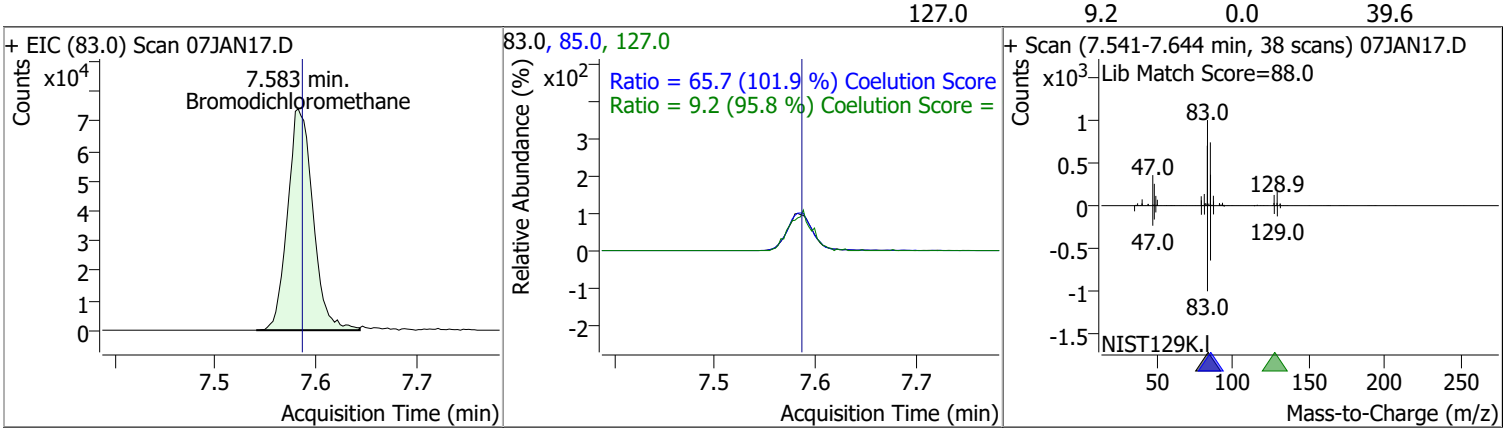
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	125.7798	7.40	0.00	46329	173.5	109.7	83.7	143.7
					95.0	83.7	52.2	112.2



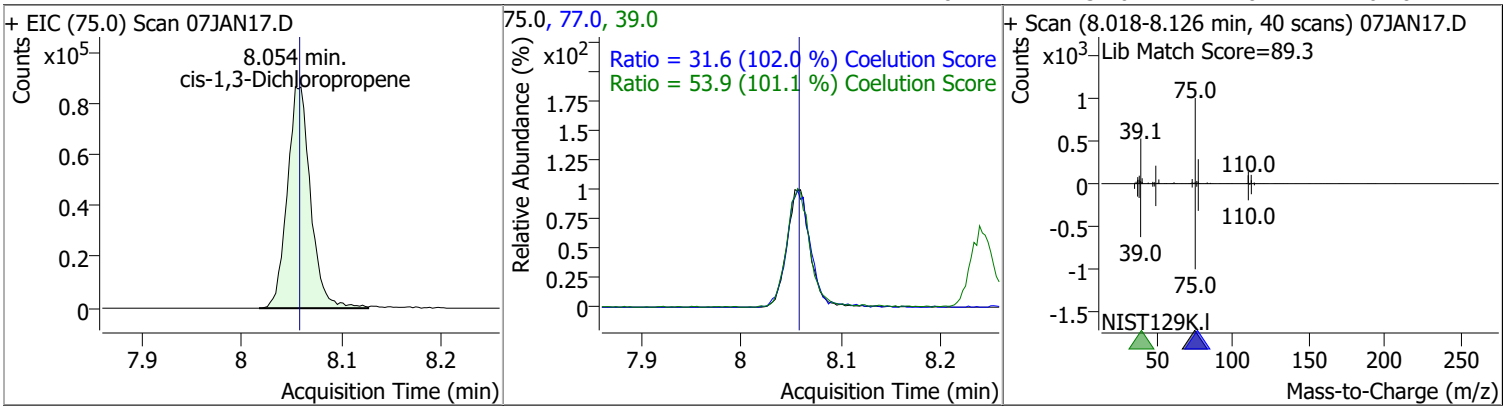




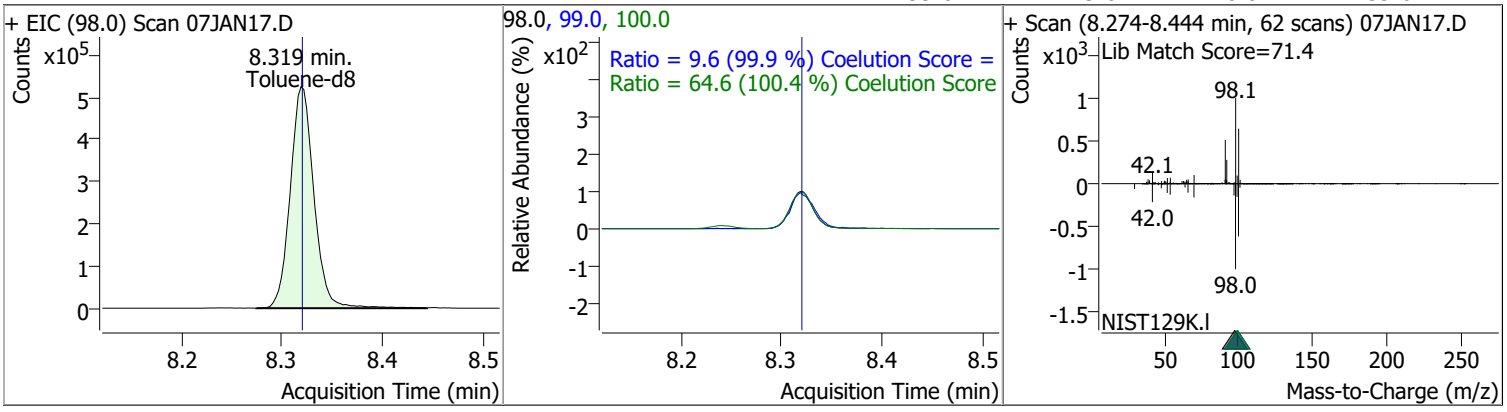
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	128.2418	7.58	0.00	130361	85.0	65.7	34.5	94.5
					127.0	9.2	0.0	39.6



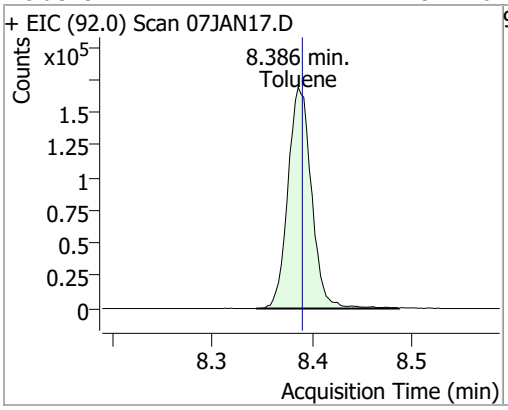
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	121.0494	8.05	-0.01	139124	39.0	53.9	23.3	83.3
					77.0	31.6	1.0	61.0

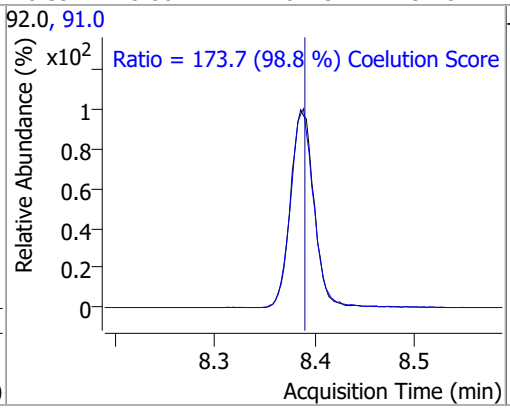


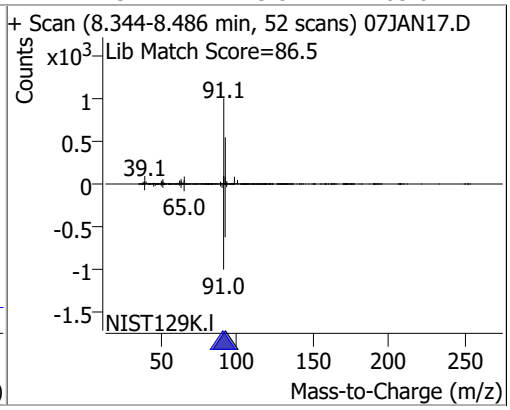
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	271.7626	8.32	0.00	860434	100.0	64.6	34.4	94.4
					99.0	9.6	0.0	39.6



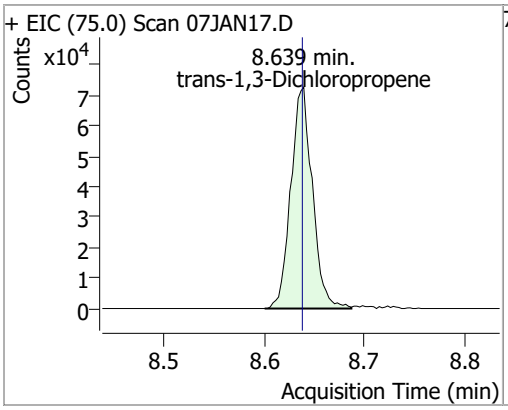
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	129.2716	8.39	0.00	276475	91.0	173.7	145.8	205.8

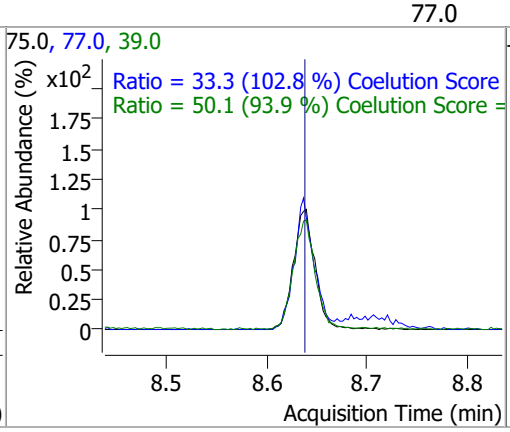


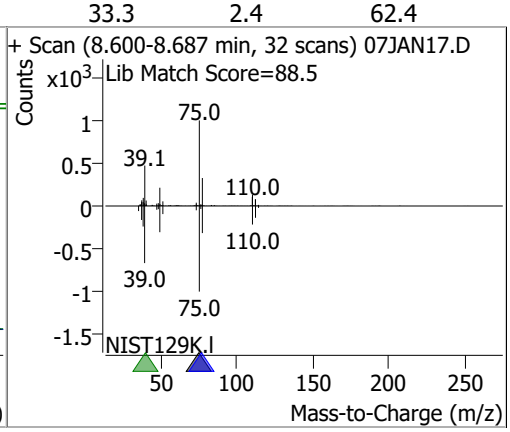




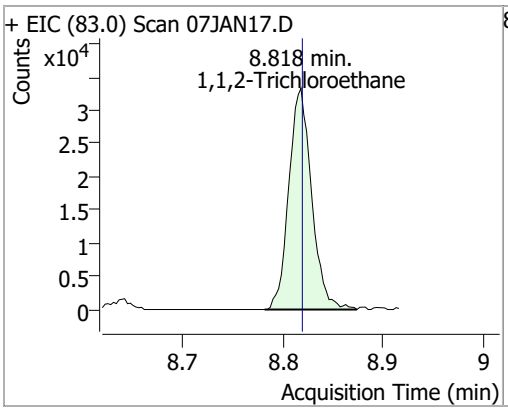
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	132.4832	8.64	0.00	108385	39.0	50.1	23.4	83.4
					77.0	33.3	2.4	62.4

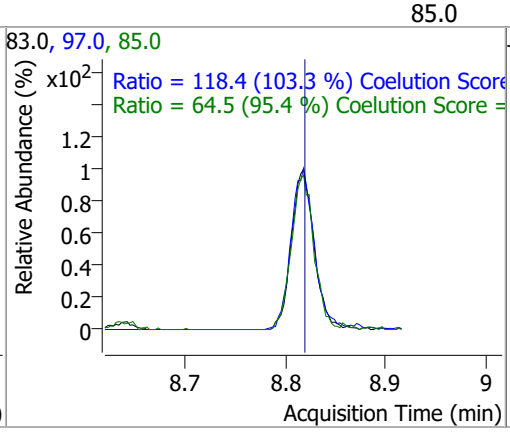


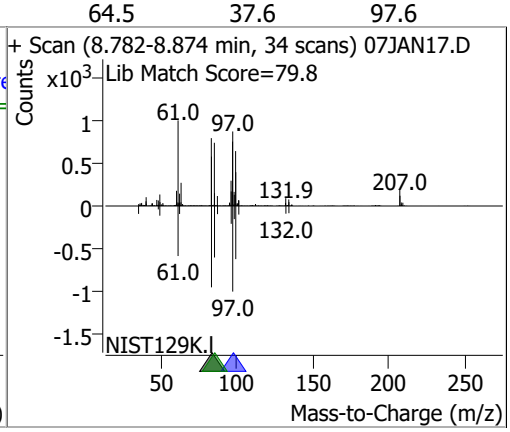




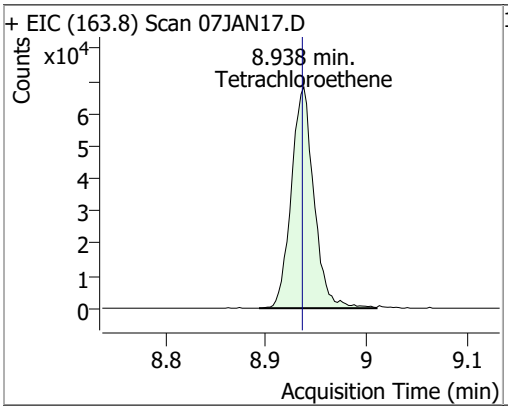
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	123.5591	8.82	0.00	52652	97.0	118.4	84.6	144.6
					85.0	64.5	37.6	97.6

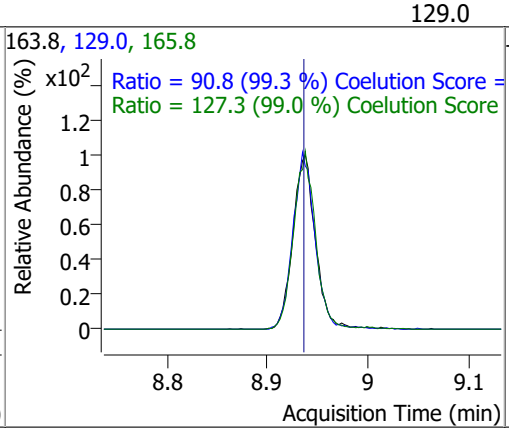


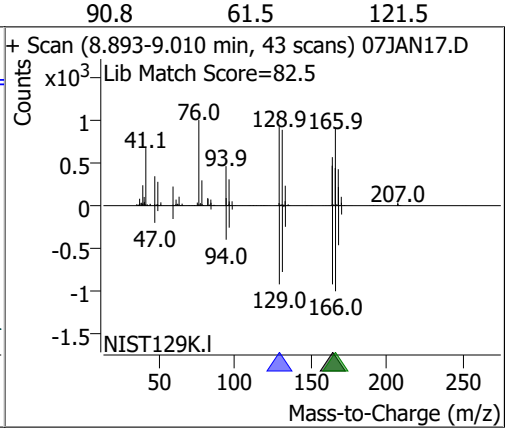




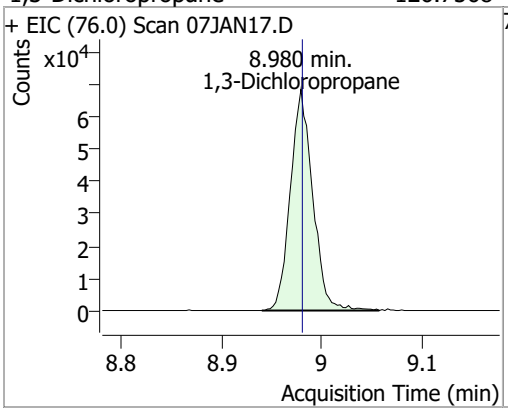
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	122.3903	8.94	0.00	106788	165.8	127.3	98.6	158.6
					129.0	90.8	61.5	121.5

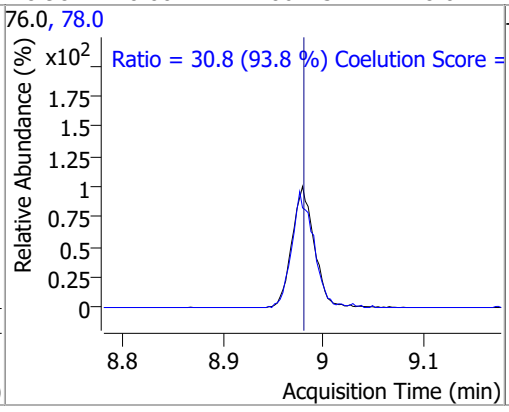


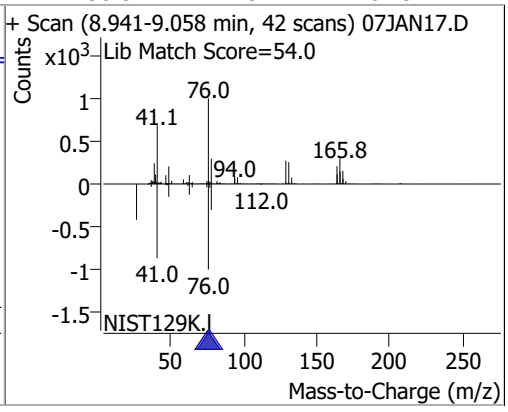




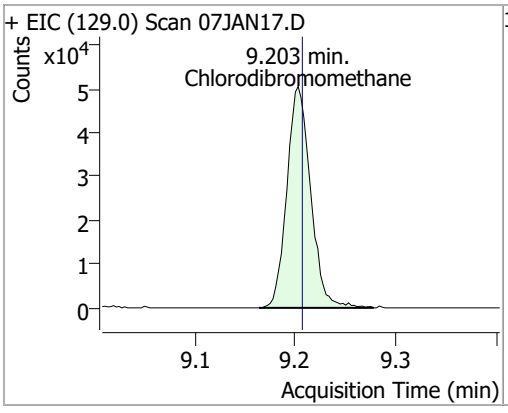
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	126.7568	8.98	0.00	106245	78.0	30.8	2.9	62.9

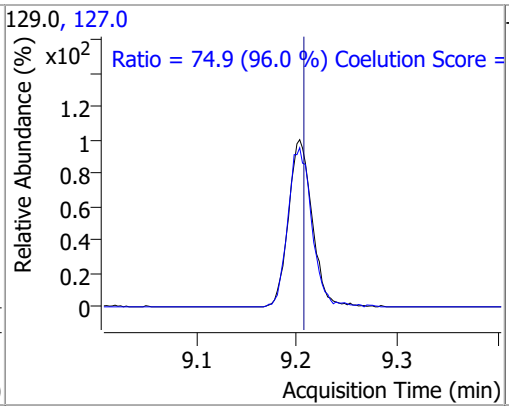


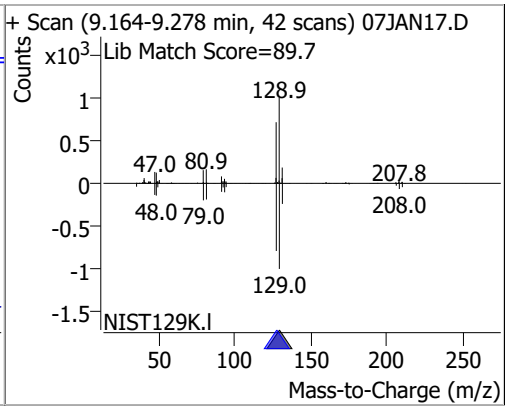




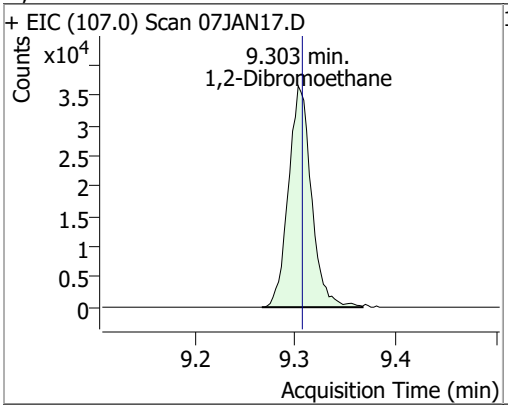
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	125.1595	9.20	0.00	83355	127.0	74.9	48.0	108.0

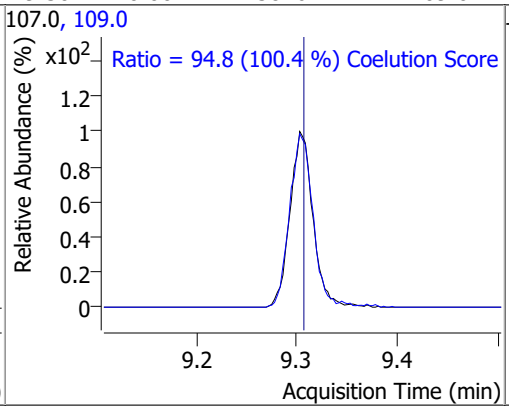


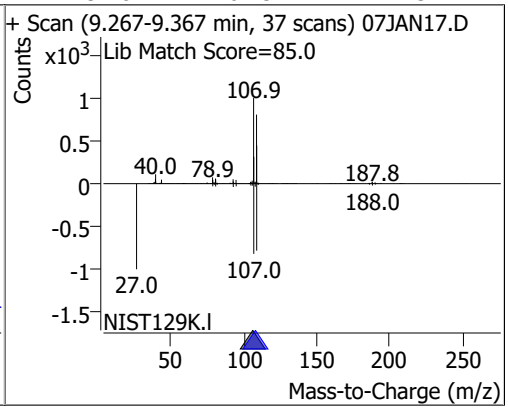




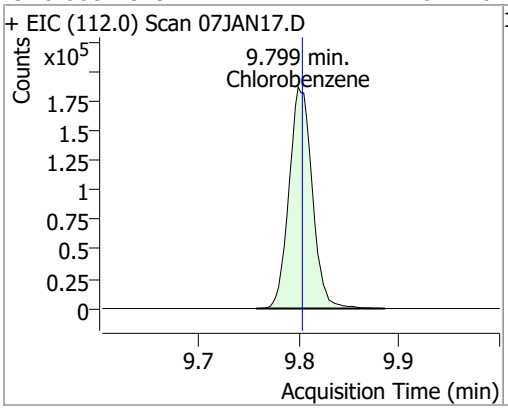
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	124.8389	9.30	0.00	58167	109.0	94.8	64.5	124.5

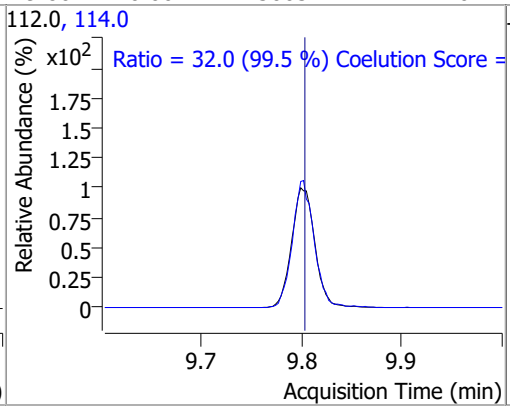


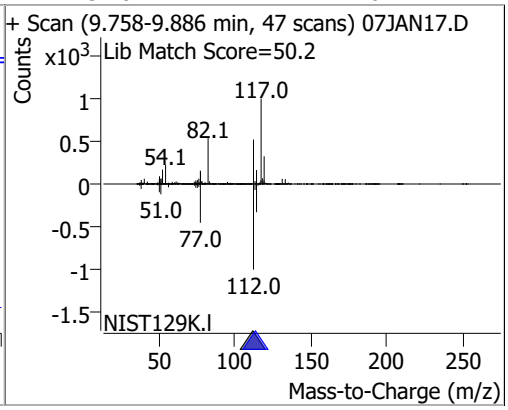




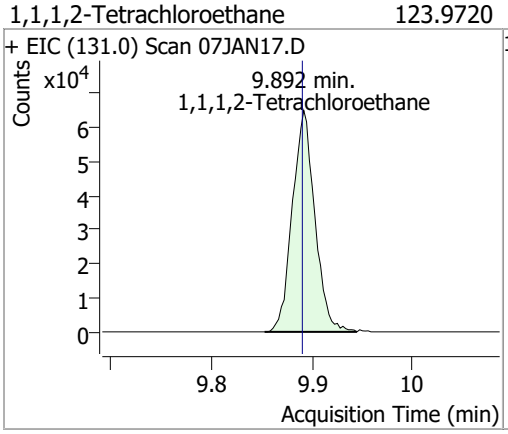
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	128.2720	9.80	0.00	300347	114.0	32.0	2.1	62.1

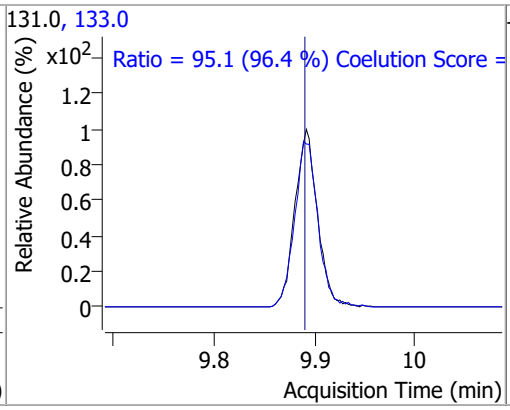


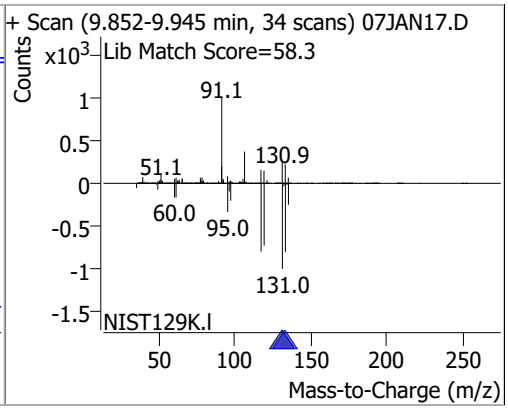




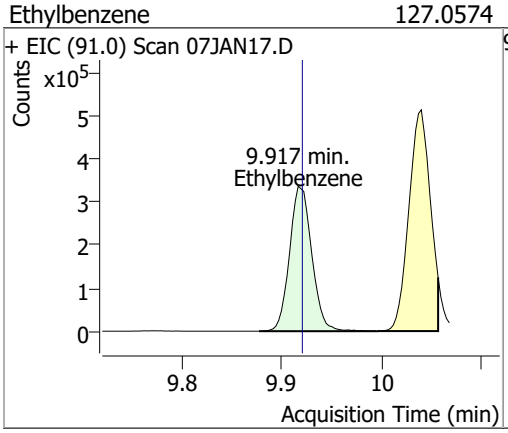
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	123.9720	9.89	0.00	101471	133.0	95.1	68.6	128.6

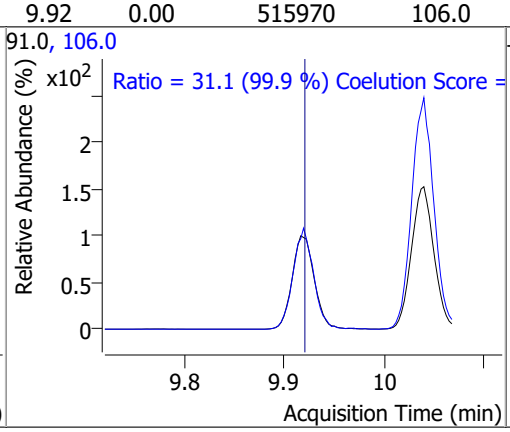


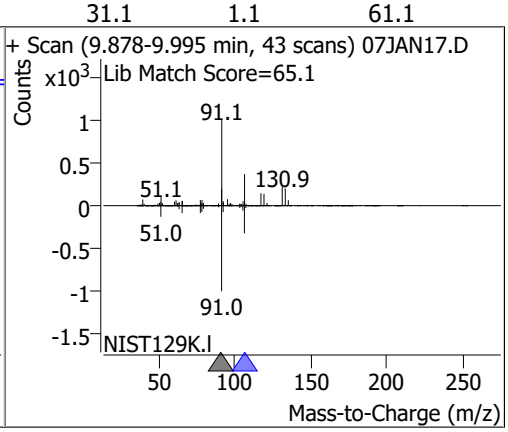




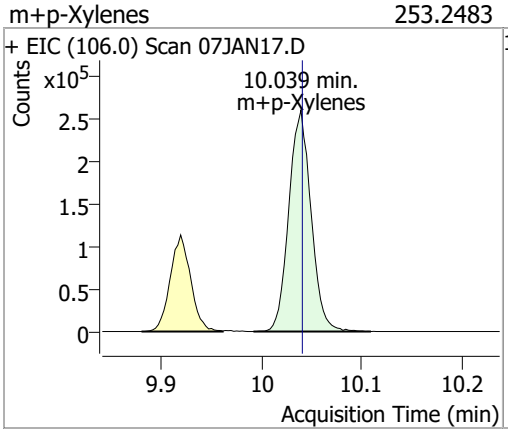
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	127.0574	9.92	0.00	515970	106.0	31.1	1.1	61.1

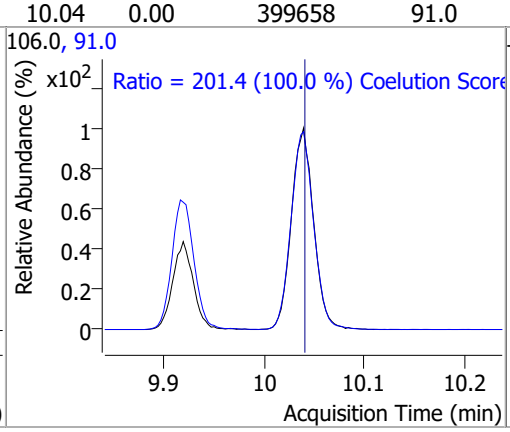


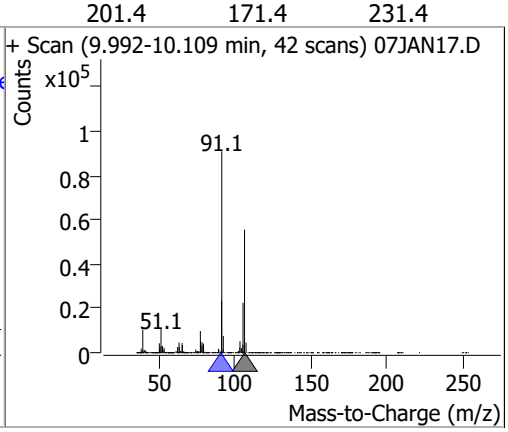




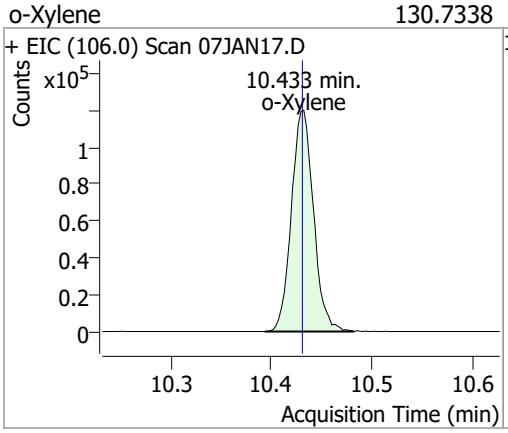
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	253.2483	10.04	0.00	399658	91.0	201.4	171.4	231.4

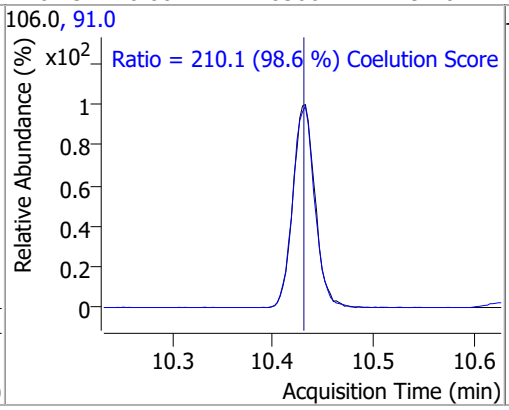


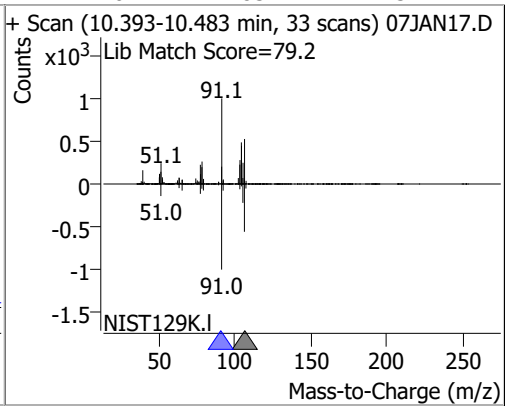




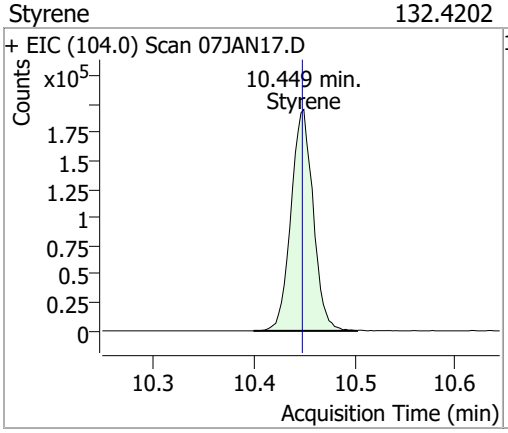
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	130.7338	10.43	0.00	183667	91.0	210.1	183.1	243.1

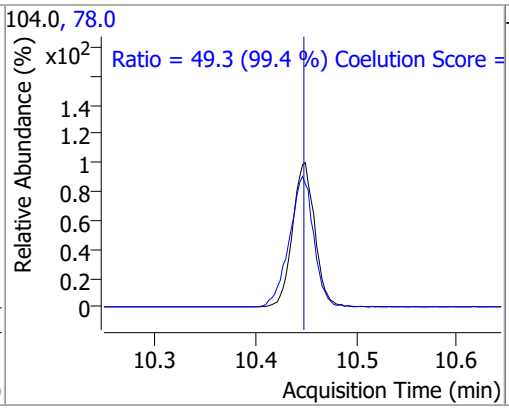


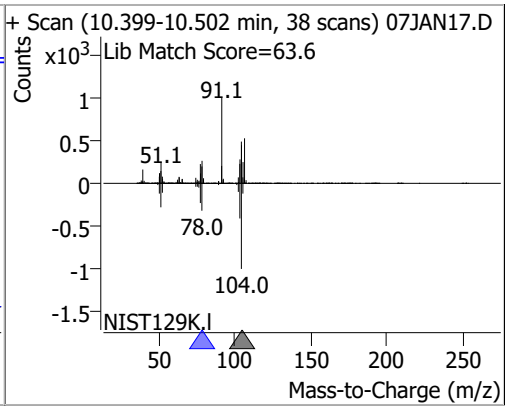




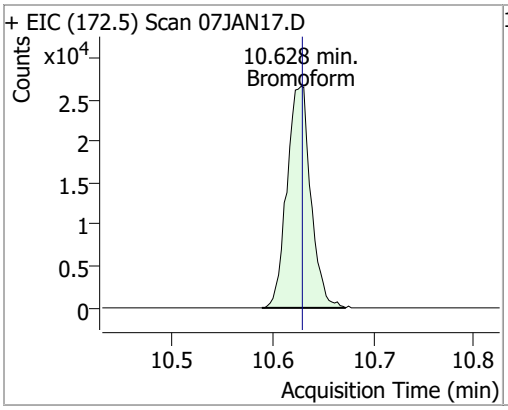
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	132.4202	10.45	0.00	299523	78.0	49.3	19.6	79.6

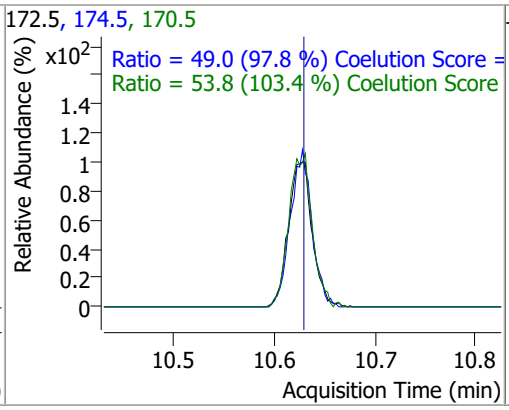


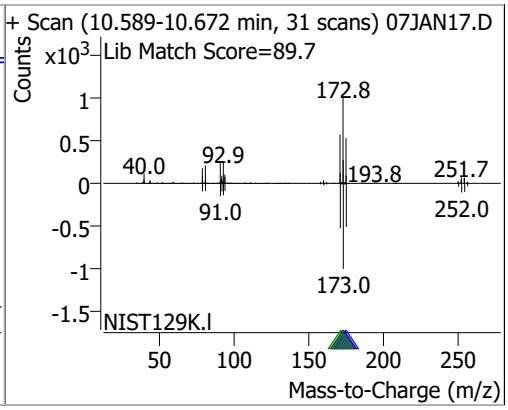




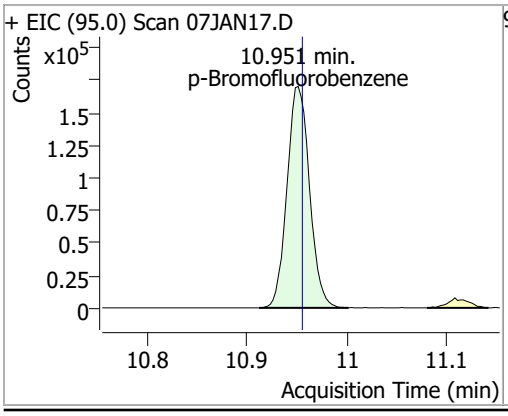
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	129.7036	10.63	0.00	43824	170.5	53.8	22.1	82.1
					174.5	49.0	20.1	80.1

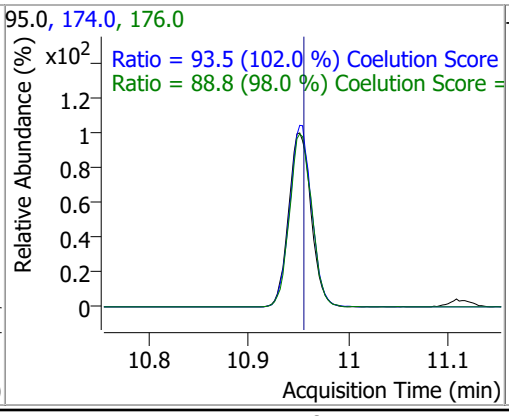


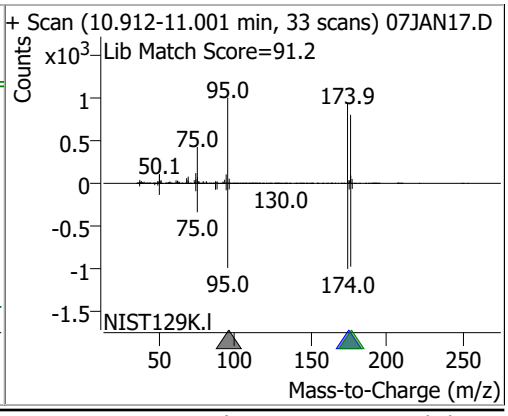




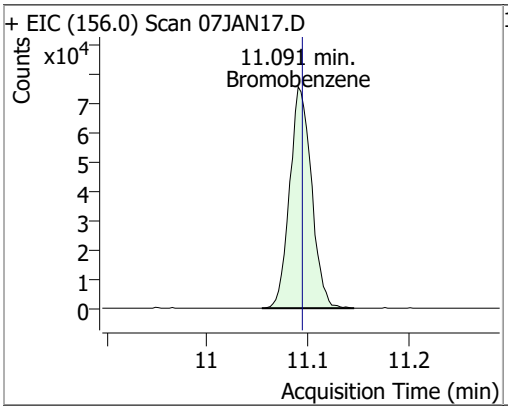
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	274.0538	10.95	0.00	265093	174.0	93.5	61.7	121.7
					176.0	88.8	60.6	120.6

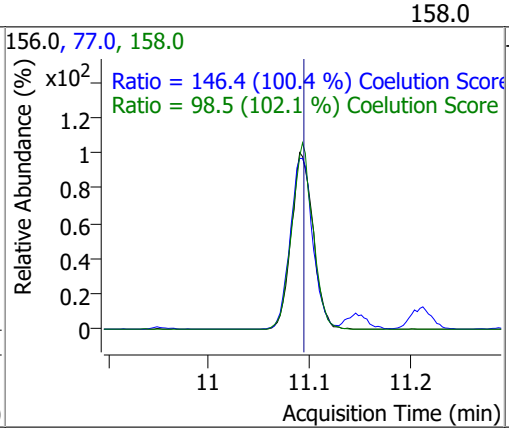


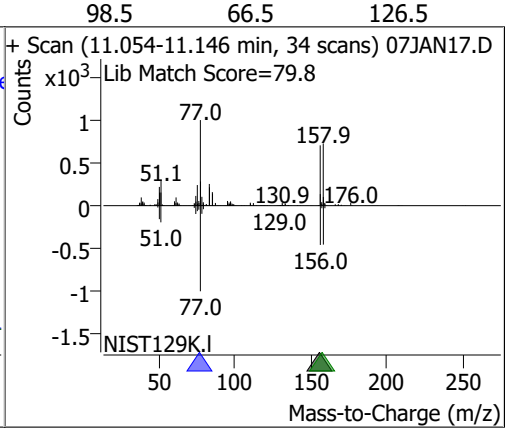




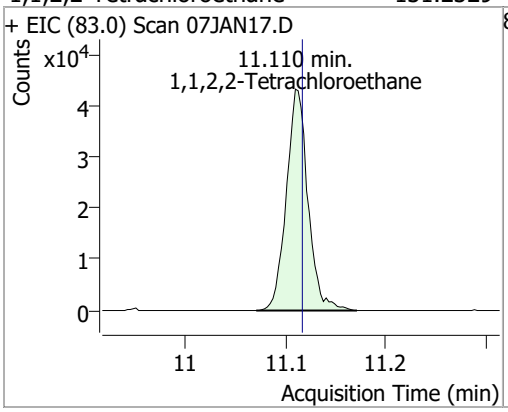
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	133.6149	11.09	0.00	114173	77.0	146.4	115.7	175.7
					158.0	98.5	66.5	126.5

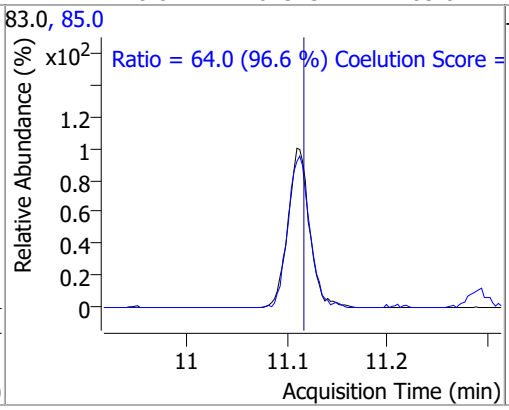


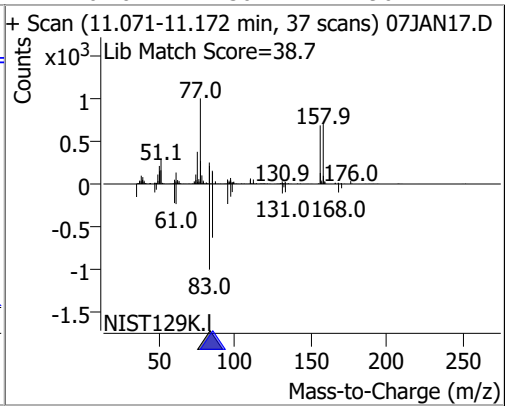




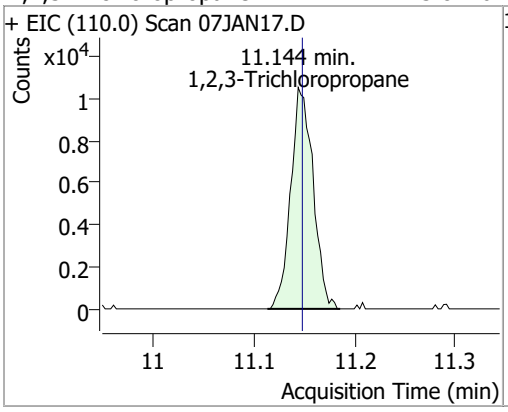
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	131.2329	11.11	-0.01	64543	85.0	64.0	36.2	96.2

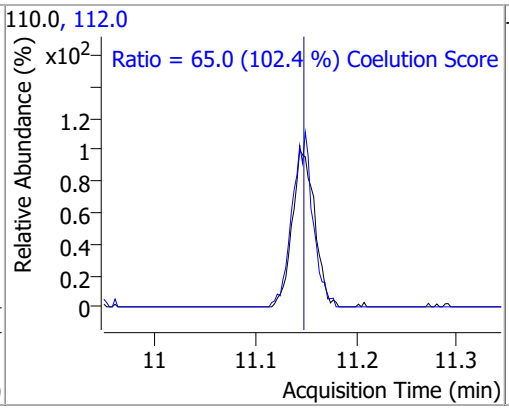


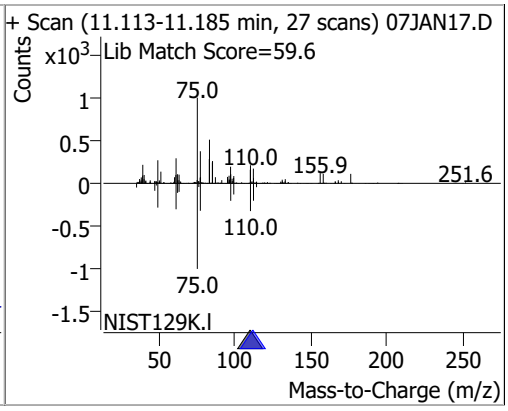




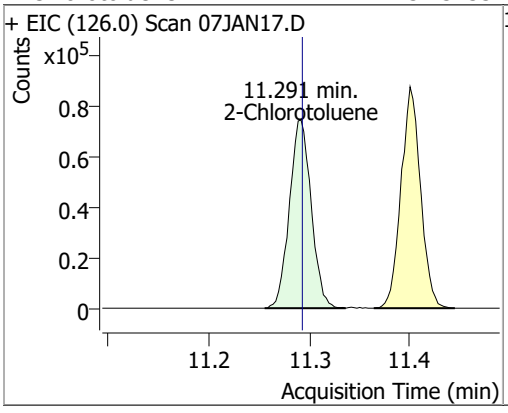
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	123.8246	11.14	0.00	16295	112.0	65.0	33.5	93.5

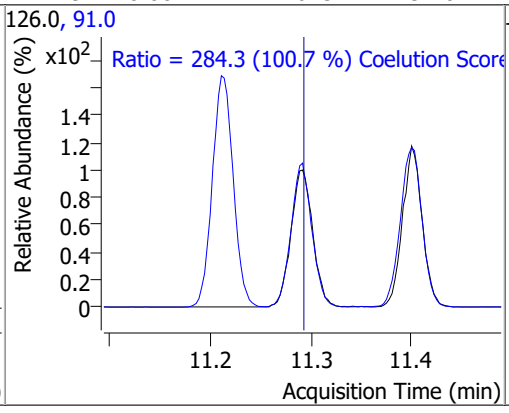


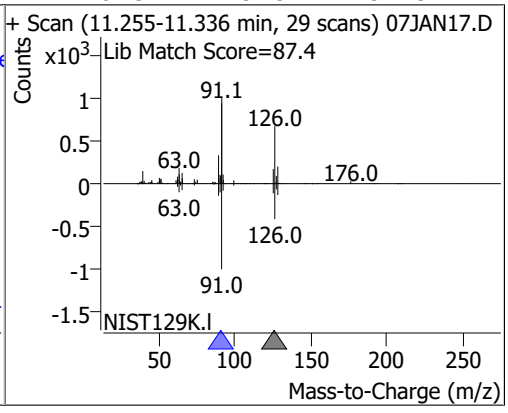




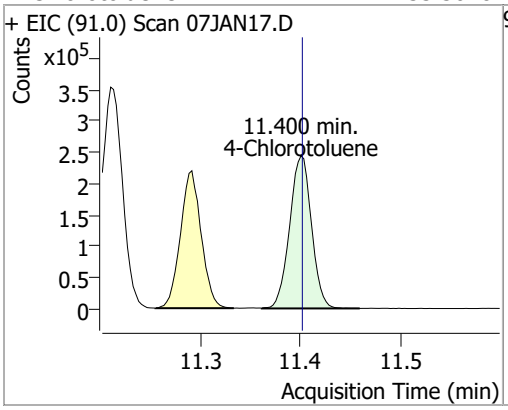
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	131.3133	11.29	0.00	111645	91.0	284.3	252.3	312.3

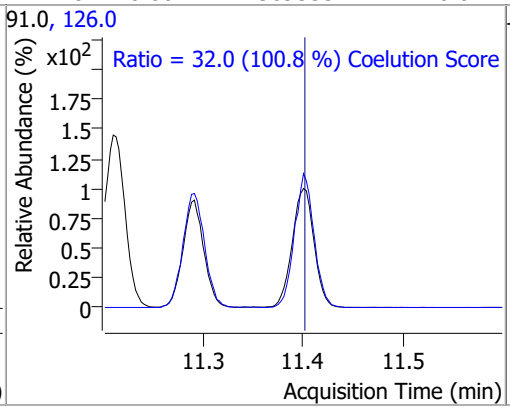


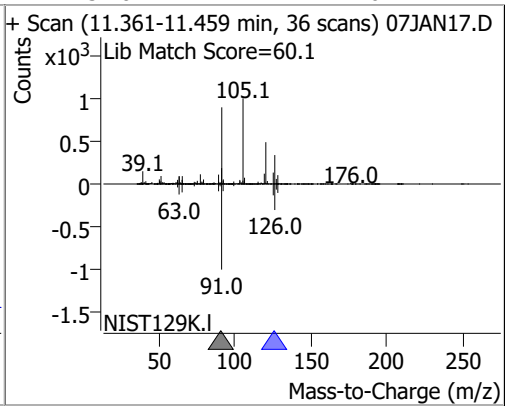




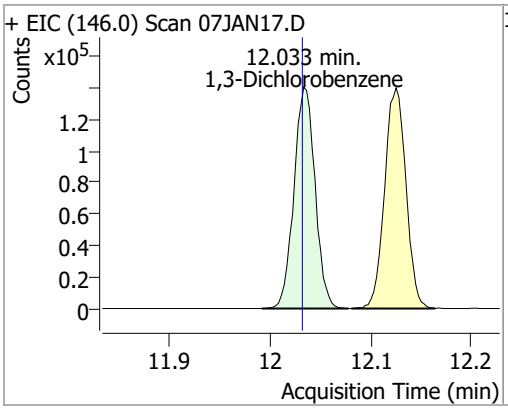
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	133.3048	11.40	0.00	369533	126.0	32.0	1.7	61.7

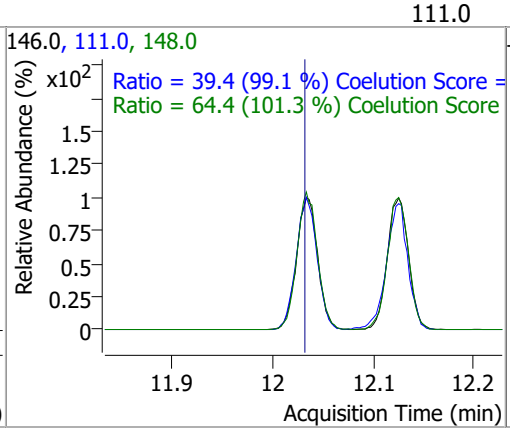


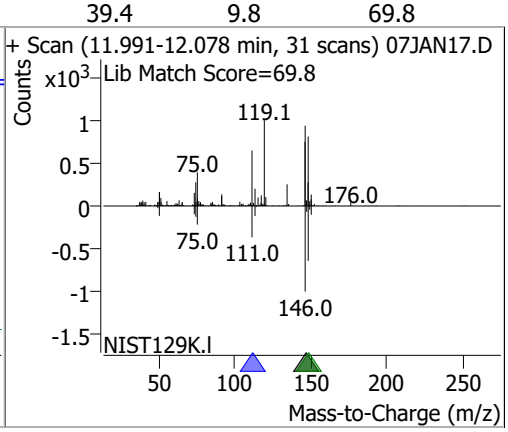




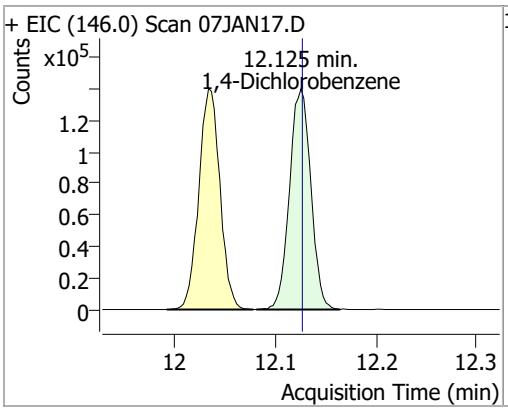
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	131.3970	12.03	0.00	204772	148.0	64.4	33.6	93.6
					111.0	39.4	9.8	69.8

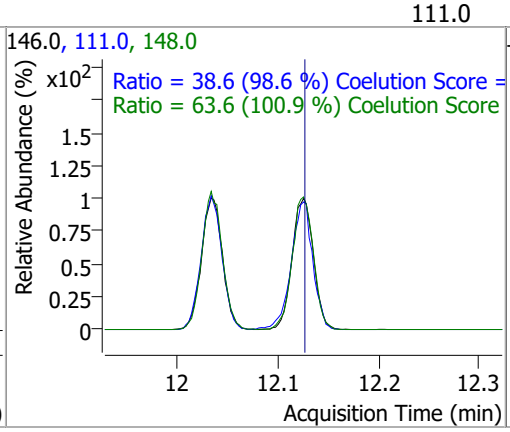


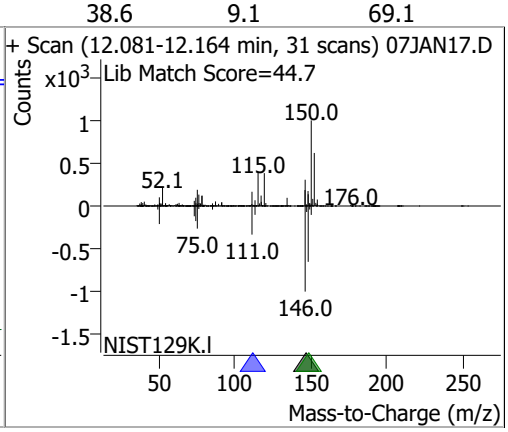




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	129.8964	12.13	0.00	206411	148.0	63.6	33.1	93.1
					111.0	38.6	9.1	69.1

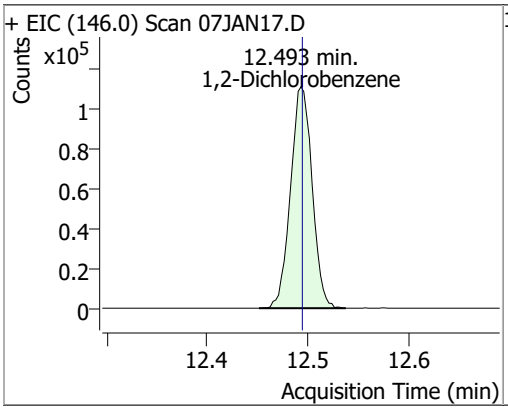


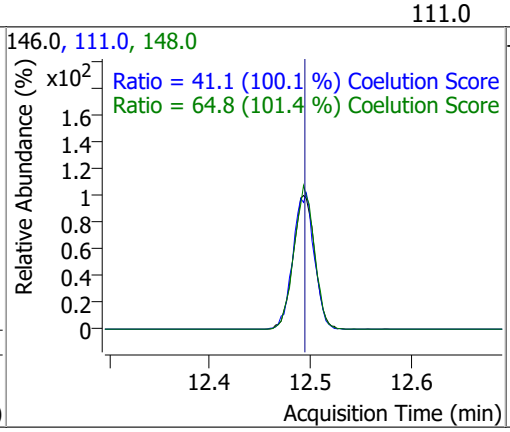


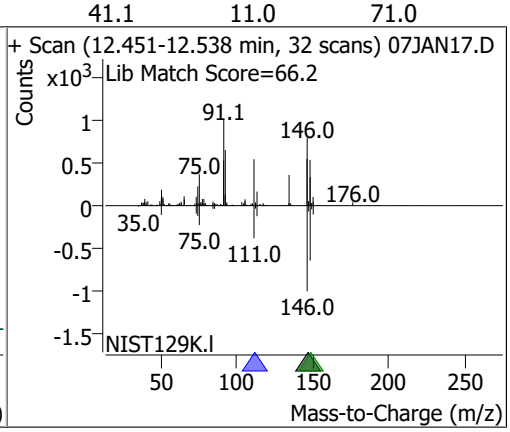


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	125.9590	12.49	0.00	165895	148.0	64.8	33.9	93.9
					111.0	41.1	11.0	71.0

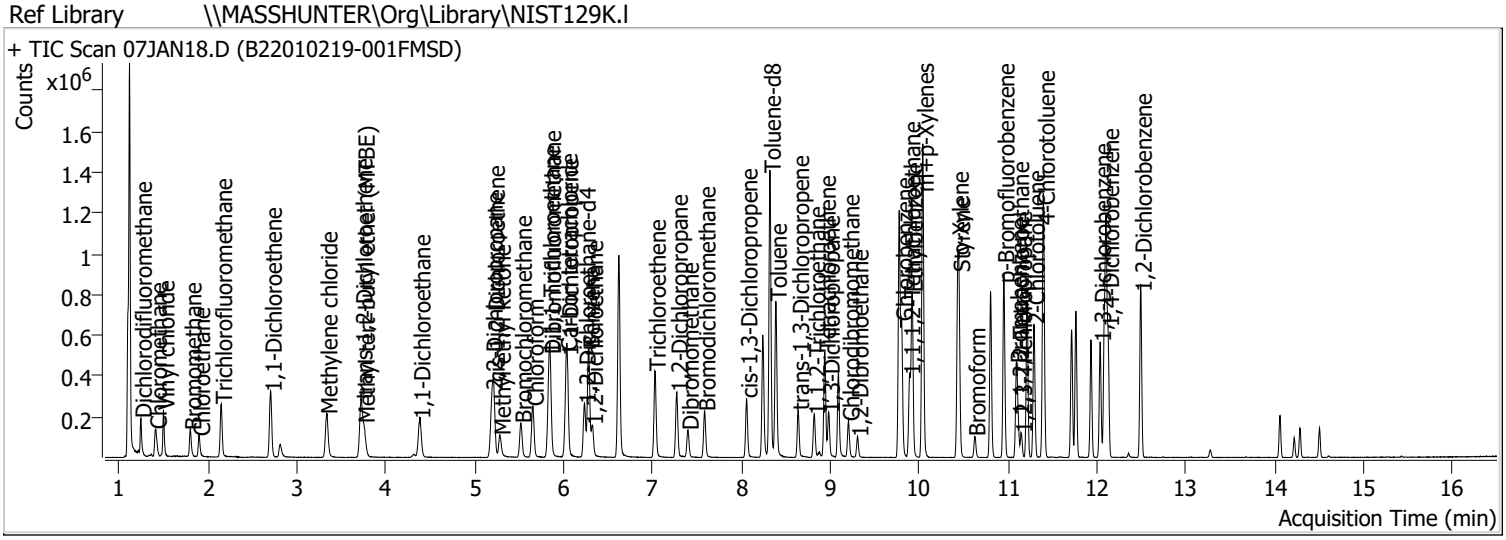






Quantitation Results Report (QT Reviewed)

Data File	07JAN18.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 5:34:50 PM
Sample Name	B22010219-001FMSD	Instrument	VOA5975C
Vial	18	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

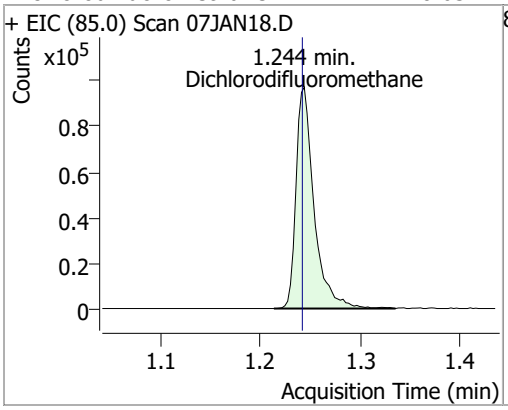


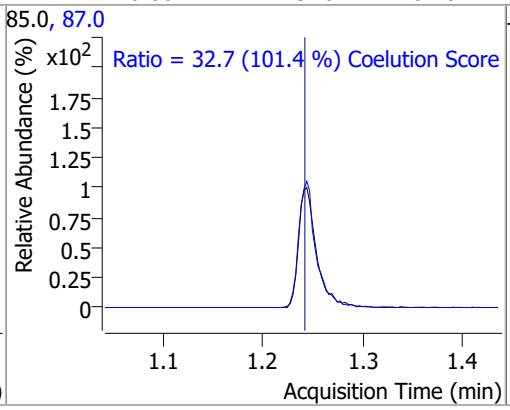
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.621	96.0	847841	250.0000	ng	-0.003
M Chlorobenzene-d5	9.772	82.0	322063	250.0000	ng	0.000
M 1,4-Dichlorobenzene-d4	12.100	152.0	260394	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.848	113.0	214547	268.6027	ng	0.003
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 107.44%		
S 1,2-Dichloroethane-d4	6.236	67.0	94029	272.5452	ng	0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 109.02%		
S Toluene-d8	8.319	98.0	857800	276.3920	ng	0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 110.56%		
S p-Bromofluorobenzene	10.949	95.0	258687	271.1727	ng	-0.006
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 108.47%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.244	85.0	122916	110.6317	ng	99
T Chloromethane	1.411	50.0	158746	117.7180	ng	100
T Vinyl chloride	1.498	62.0	151717	125.0334	ng	98
T Bromomethane	1.802	96.0	65044	119.8794	ng	96
T Chloroethane	1.899	64.0	73037	121.5844	ng	98
T Trichlorofluoromethane	2.148	101.0	181247	120.3412	ng	99
T 1,1-Dichloroethene	2.705	96.0	114899	134.5404	ng	99
T Methylene chloride	3.333	49.0	158297	125.7372	ng	99
T trans-1,2-Dichloroethene	3.723	96.0	120311	138.0852	ng	99
T Methyl tert-butyl ether (MTBE)	3.751	73.0	161012	142.9708	ng	100
T 1,1-Dichloroethane	4.379	63.0	226152	139.4457	ng	99
T 2,2-Dichloropropane	5.193	77.0	171641	141.2419	ng	99
T cis-1,2-Dichloroethene	5.212	96.0	118290	133.9099	ng	99
T Methyl ethyl ketone	5.282	43.0	145750	1218.1021	ng	99
T Bromochloromethane	5.519	128.0	46796	127.8755	ng	97
T Chloroform	5.653	83.0	203762	126.2450	ng	100

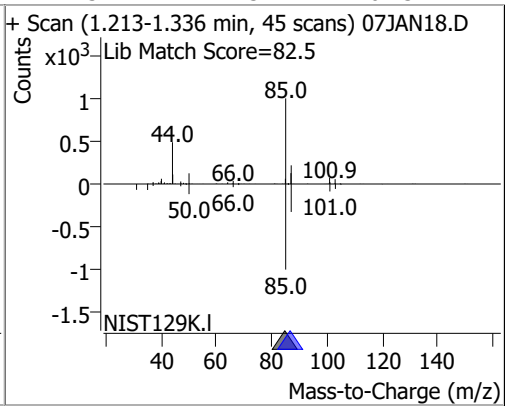
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.834	97.0	196145	129.6747	ng	99
T Carbon tetrachloride	6.029	117.0	185767	124.6502	ng	97
T 1,1-Dichloropropene	6.041	75.0	161545	125.6088	ng	99
T Benzene	6.283	78.0	449224	133.0746	ng	99
T 1,2-Dichloroethane	6.319	62.0	116685	127.7729	ng	96
T Trichloroethene	7.025	95.0	129711	133.5436	ng	99
T 1,2-Dichloropropane	7.270	63.0	113897	133.3077	ng	98
T Dibromomethane	7.396	93.0	46728	129.4203	ng	97
T Bromodichloromethane	7.583	83.0	134905	135.3871	ng	98
T cis-1,3-Dichloropropene	8.057	75.0	143696	127.5477	ng	98
T Toluene	8.386	92.0	289892	138.2772	ng	98
T trans-1,3-Dichloropropene	8.637	75.0	108822	135.6987	ng	98
T 1,1,2-Trichloroethane	8.821	83.0	53613	128.3504	ng	97
T Tetrachloroethene	8.938	163.8	112970	132.0855	ng	98
T 1,3-Dichloropropane	8.983	76.0	109032	132.7040	ng	98
T Chlorodibromomethane	9.203	129.0	83383	127.7253	ng	99
T 1,2-Dibromoethane	9.303	107.0	60123	131.6380	ng	99
T Chlorobenzene	9.802	112.0	312027	135.9465	ng	100
T 1,1,1,2-Tetrachloroethane	9.889	131.0	101872	126.9708	ng	99
T Ethylbenzene	9.917	91.0	539120	135.4342	ng	99
T m+p-Xylenes	10.039	106.0	410298	265.2312	ng	97
T o-Xylene	10.433	106.0	191917	139.3598	ng	98
T Styrene	10.447	104.0	313760	141.5106	ng	99
T Bromoform	10.622	172.5	46577	139.7801	ng	97
T Bromobenzene	11.091	156.0	120119	142.5401	ng	98
T 1,1,2,2-Tetrachloroethane	11.113	83.0	66475	137.0521	ng	95
T 1,2,3-Trichloropropane	11.147	110.0	17465	134.5721	ng	99
T 2-Chlorotoluene	11.289	126.0	118405	141.2126	ng	100
T 4-Chlorotoluene	11.400	91.0	388335	142.0473	ng	100
T 1,3-Dichlorobenzene	12.033	146.0	212018	137.9499	ng	99
T 1,4-Dichlorobenzene	12.123	146.0	215367	137.4287	ng	99
T 1,2-Dichlorobenzene	12.493	146.0	172710	132.9681	ng	99

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

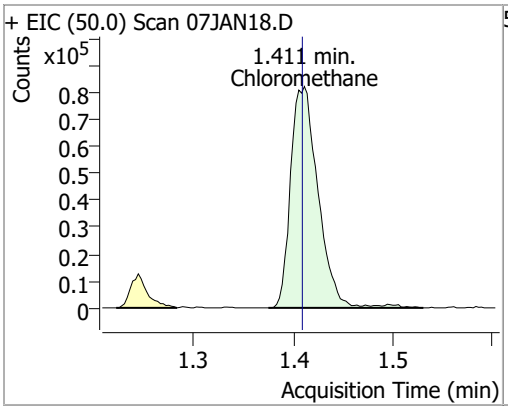
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	110.6317	1.24	0.00	122916	87.0	32.7	2.3	62.3

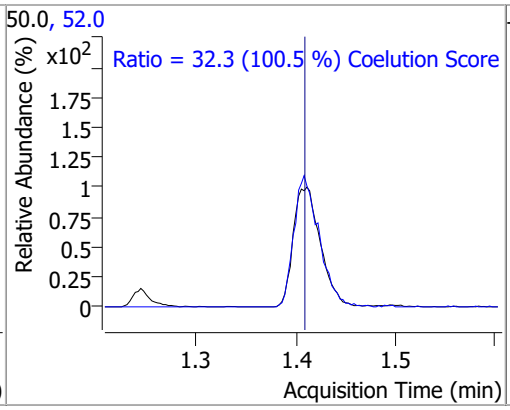


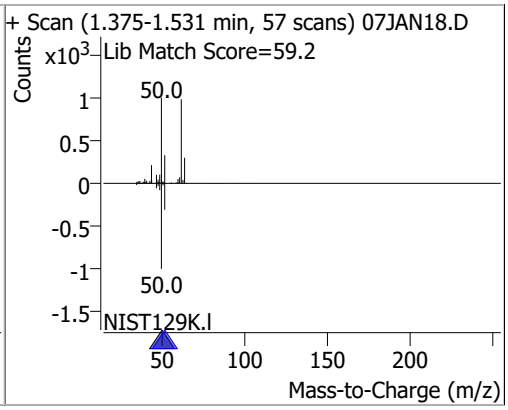




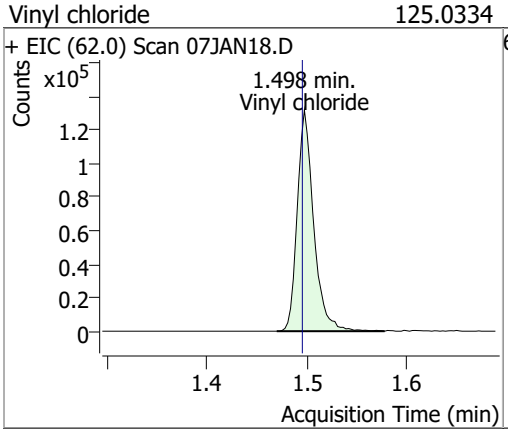
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	117.7180	1.41	0.00	158746	52.0	32.3	2.1	62.1

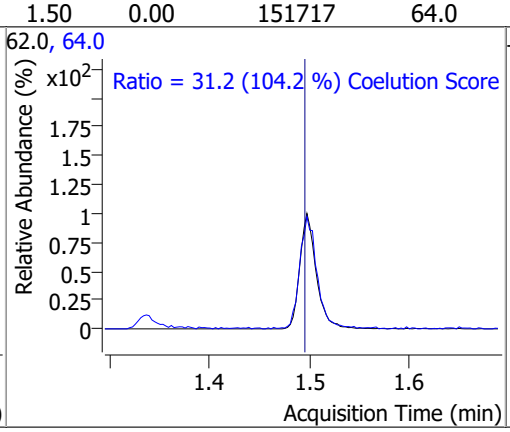


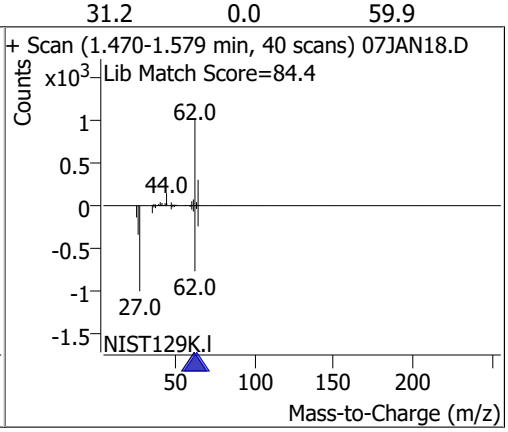




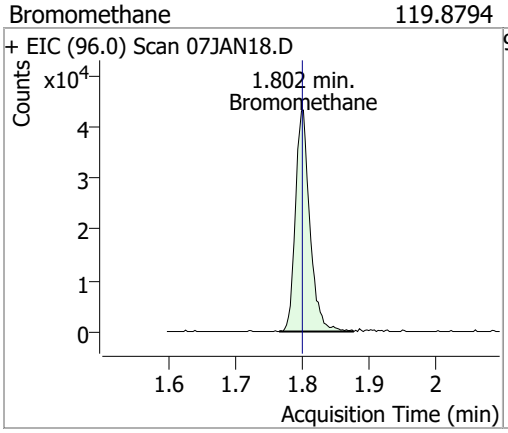
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	125.0334	1.50	0.00	151717	64.0	31.2	0.0	59.9

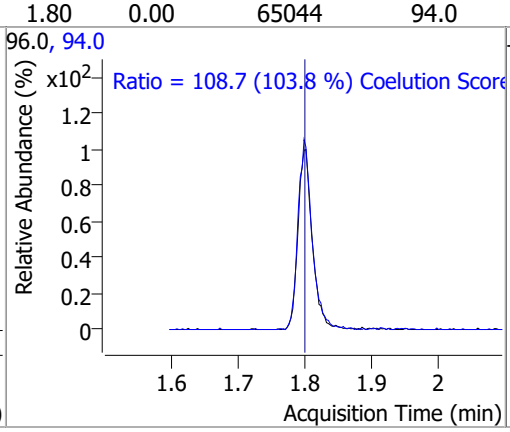


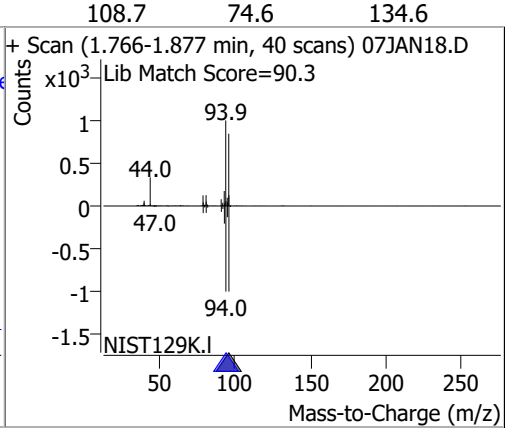




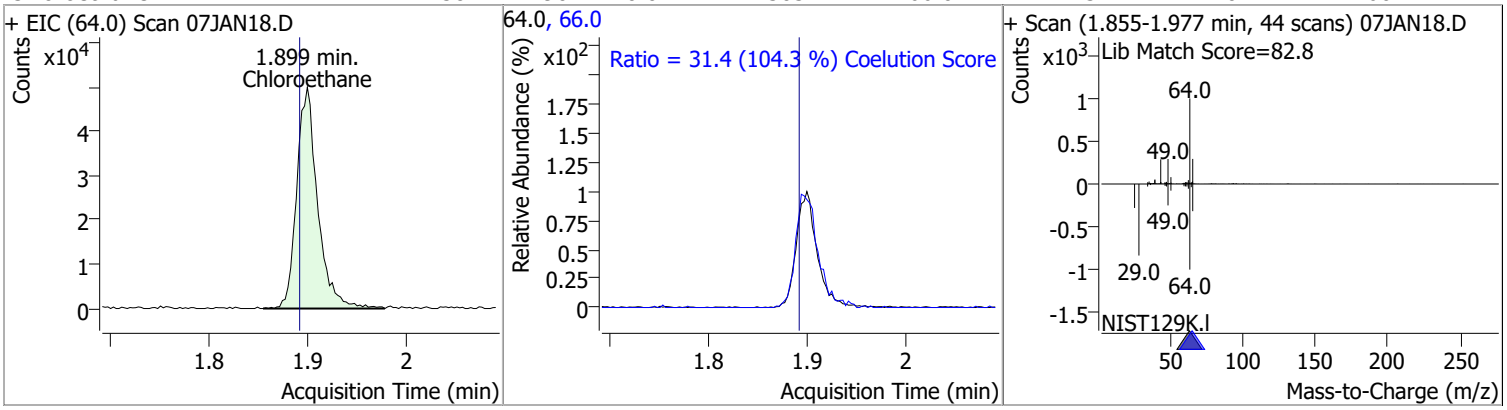
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	119.8794	1.80	0.00	65044	94.0	108.7	74.6	134.6



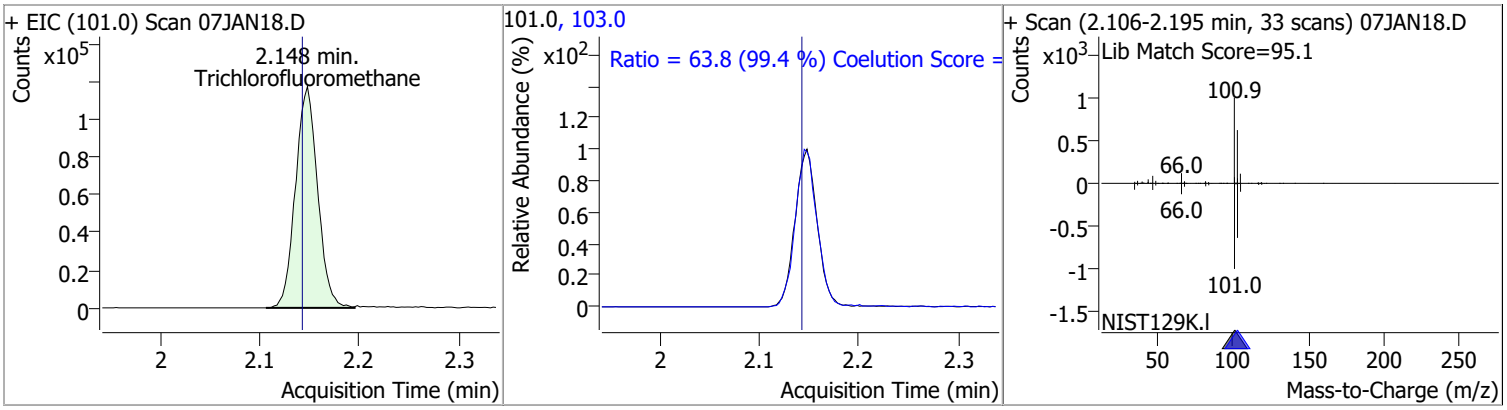




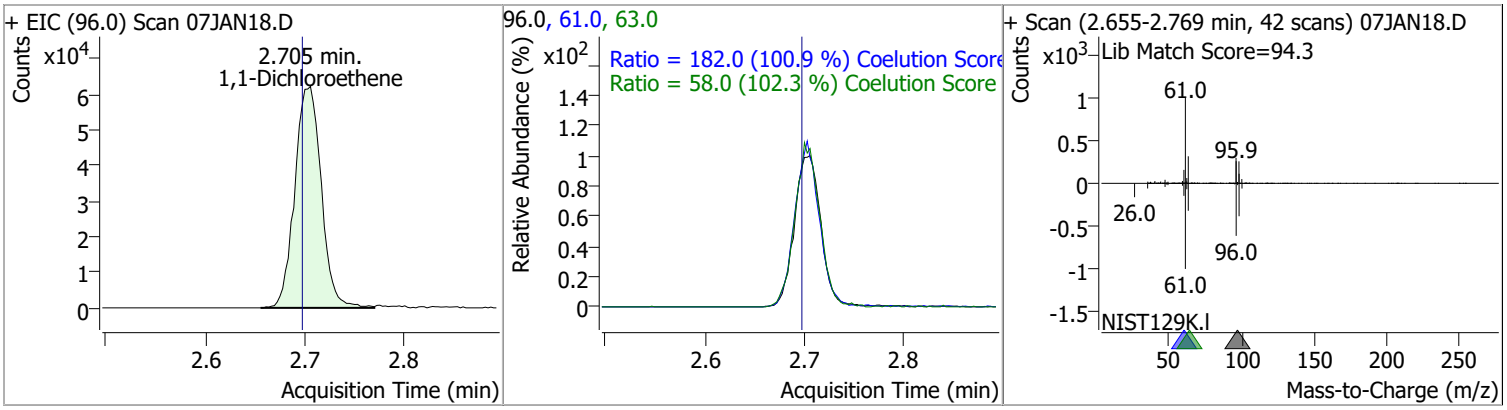
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	121.5844	1.90	0.01	73037	66.0	31.4	0.1	60.1



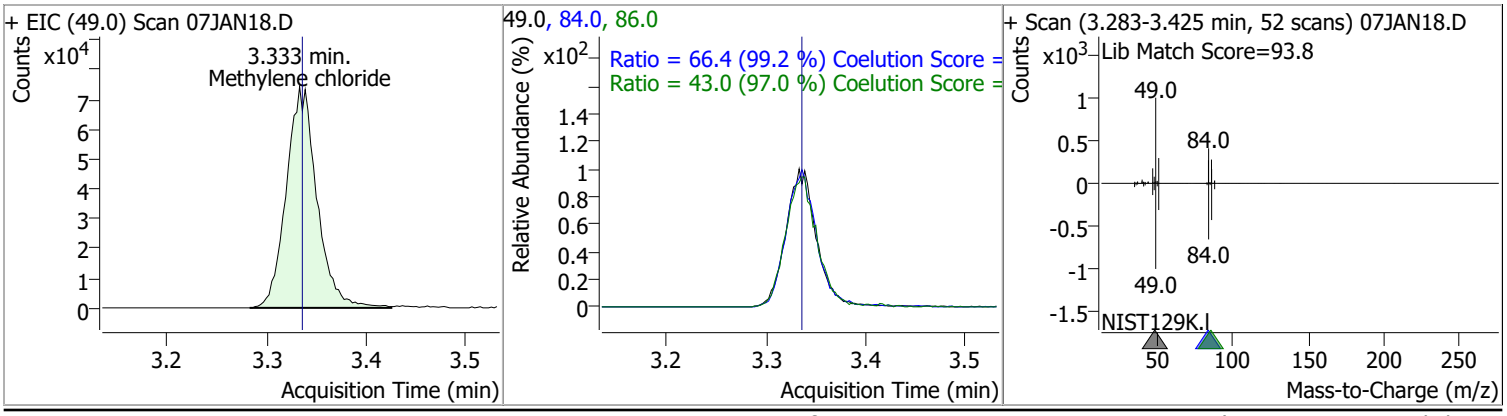
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	120.3412	2.15	0.01	181247	103.0	63.8	34.2	94.2



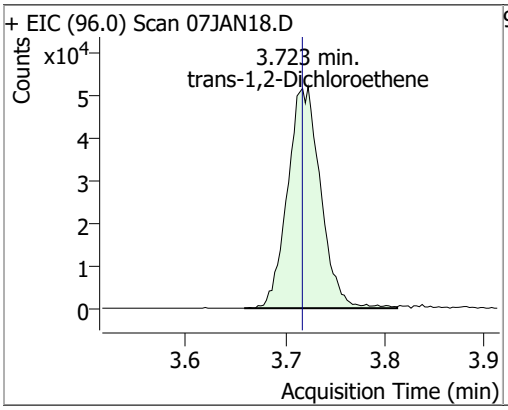
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	134.5404	2.71	0.01	114899	61.0	182.0	150.3	210.3
					63.0	58.0	26.7	86.7

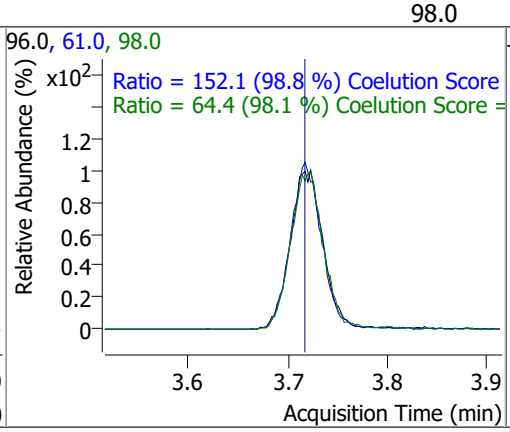


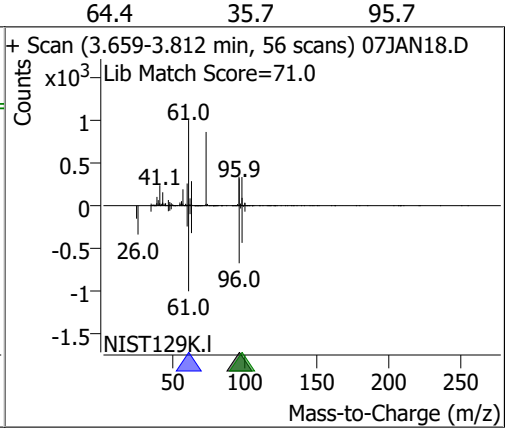
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	125.7372	3.33	0.00	158297	84.0	66.4	36.9	96.9
					86.0	43.0	14.3	74.3



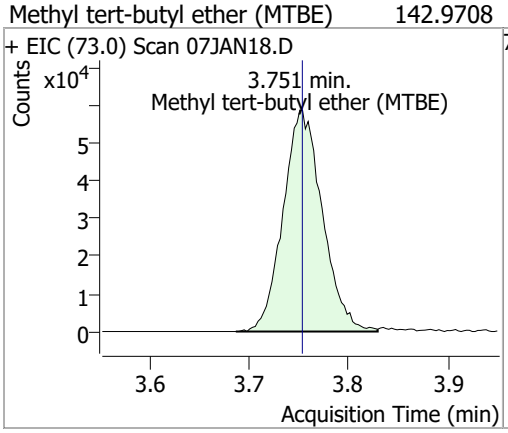
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	138.0852	3.72	0.01	120311	61.0	152.1	123.9	183.9
					98.0	64.4	35.7	95.7

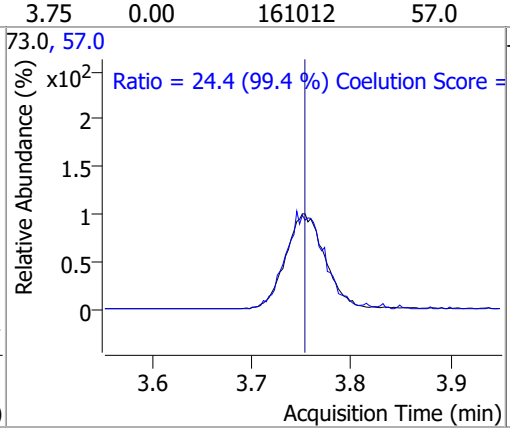


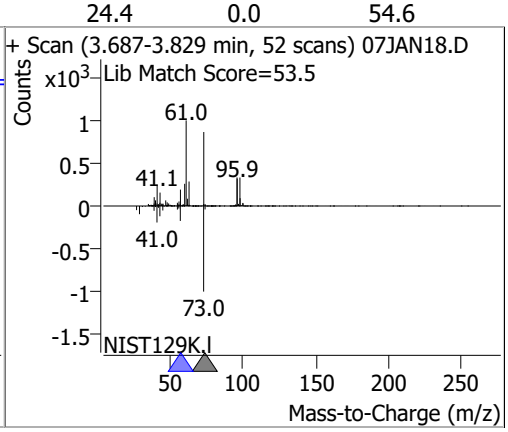




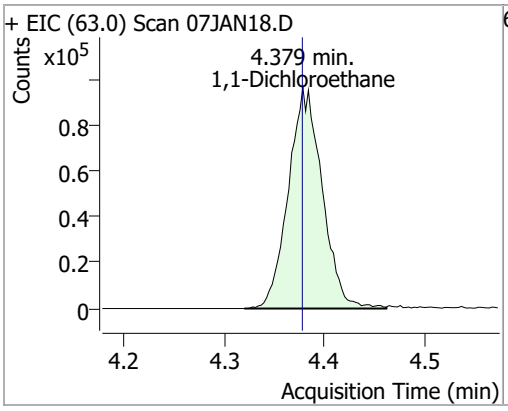
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	142.9708	3.75	0.00	161012	57.0	24.4	0.0	54.6

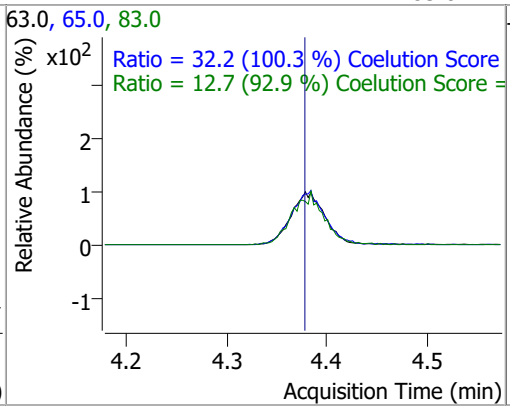


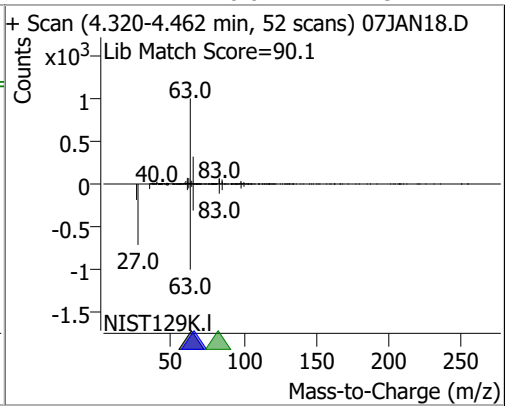




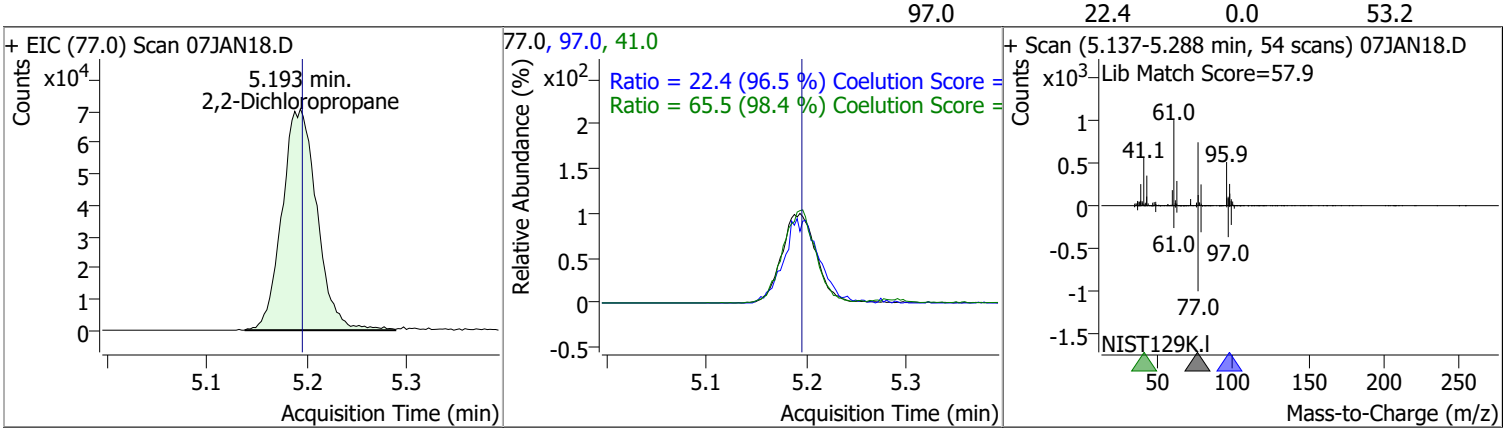
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	139.4457	4.38	0.00	226152	65.0	32.2	2.1	62.1
					83.0	12.7	0.0	43.7



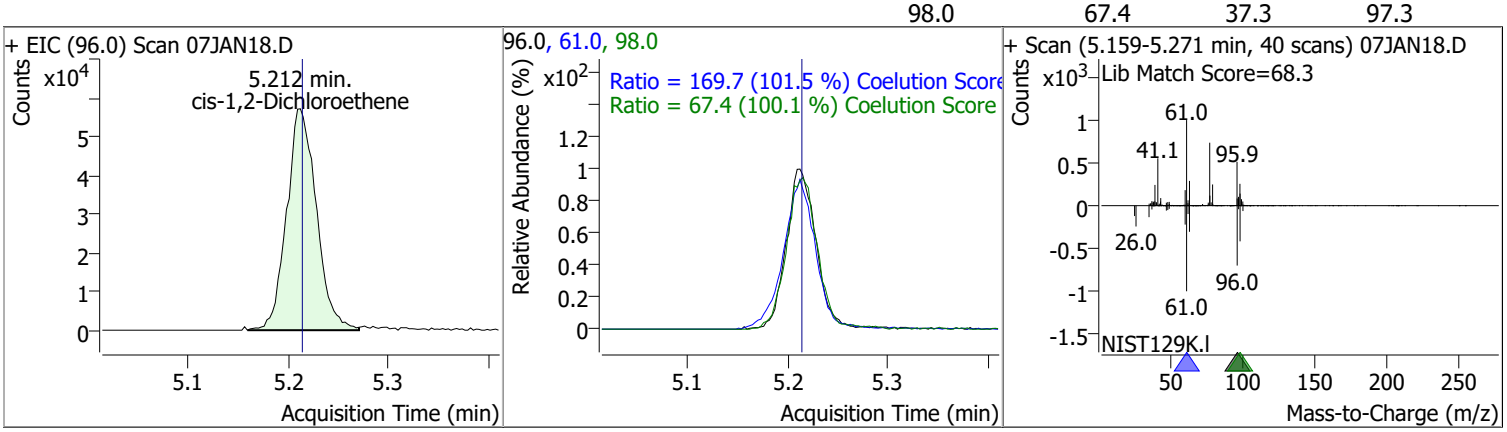




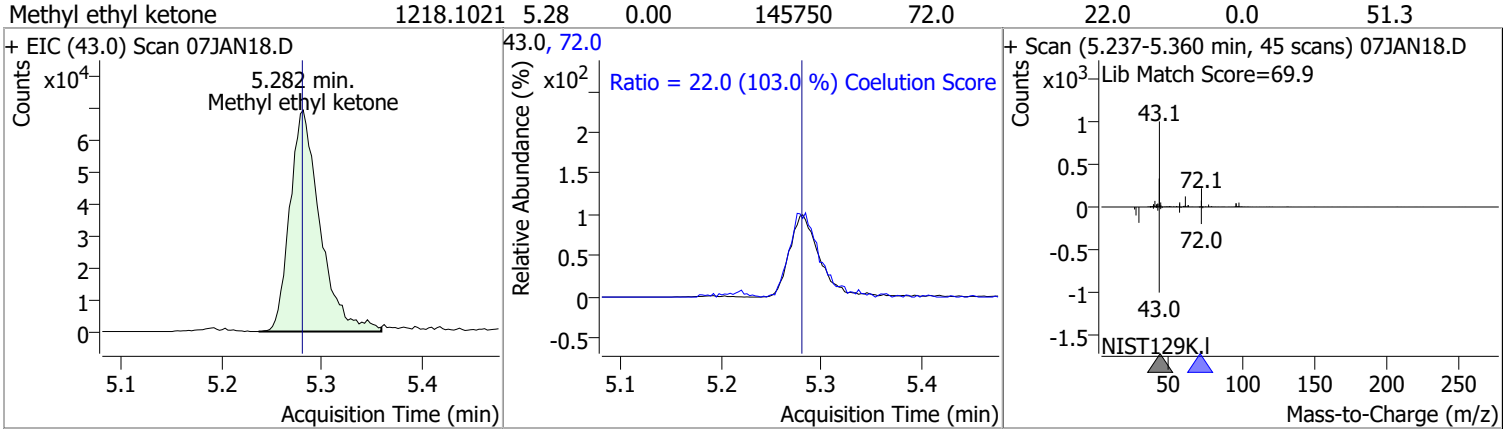
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	141.2419	5.19	0.00	171641	41.0	65.5	36.5	96.5
					97.0	22.4	0.0	53.2



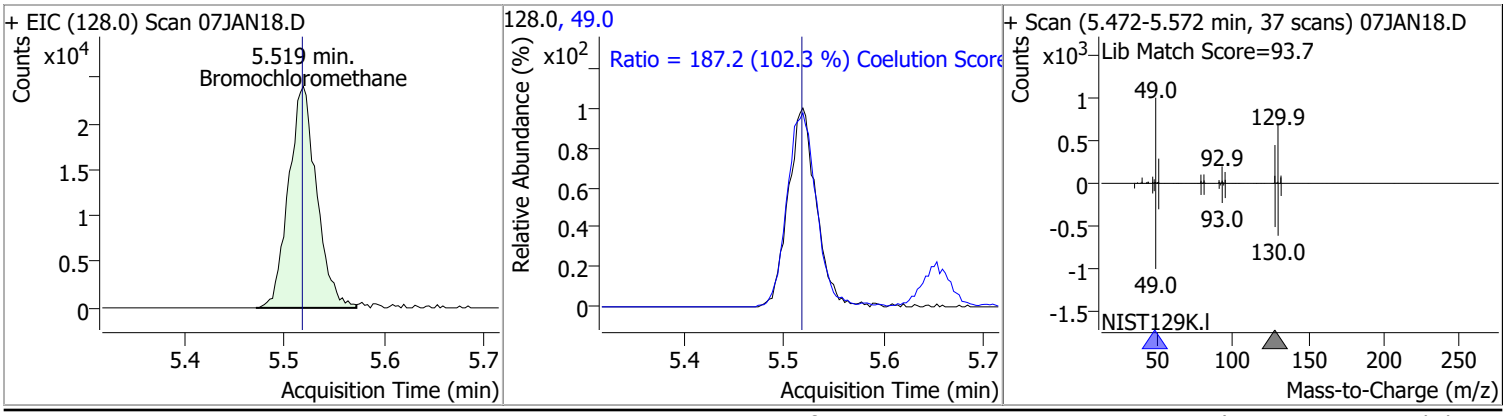
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	133.9099	5.21	0.00	118290	61.0	169.7	137.2	197.2
					98.0	67.4	37.3	97.3



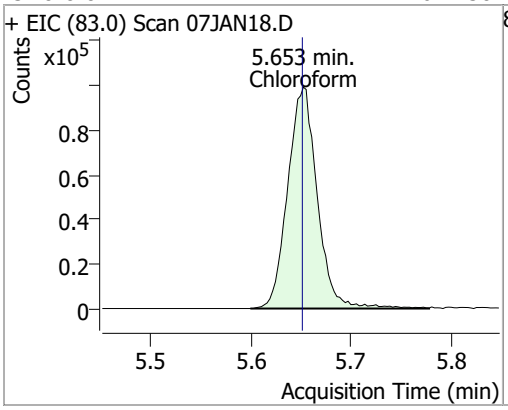
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1218.1021	5.28	0.00	145750	72.0	22.0	0.0	51.3

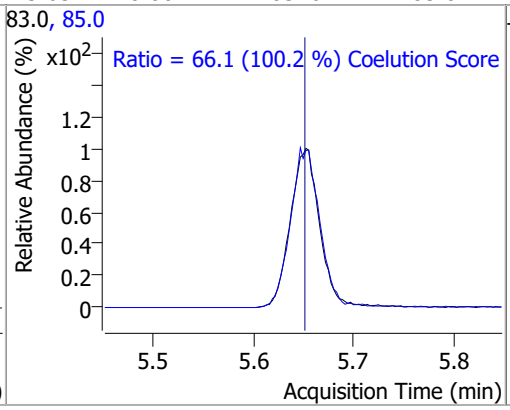


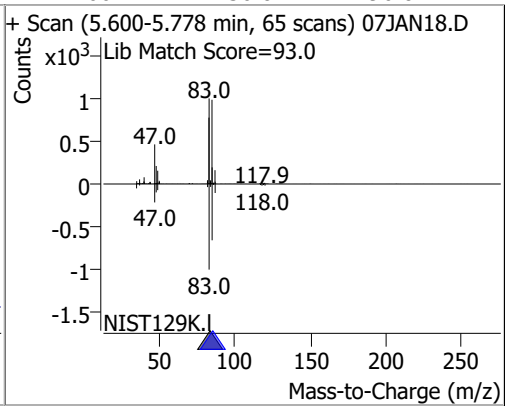
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	127.8755	5.52	0.00	46796	49.0	187.2	152.9	212.9



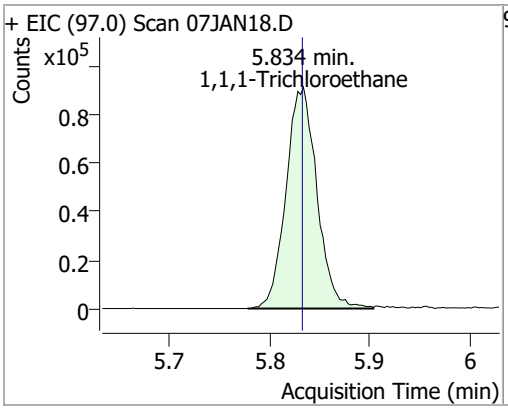
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	126.2450	5.65	0.00	203762	85.0	66.1	36.0	96.0

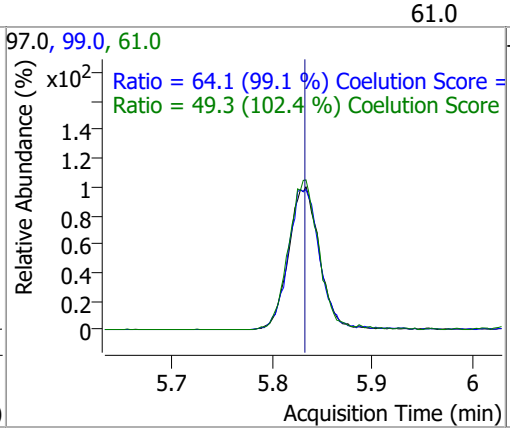


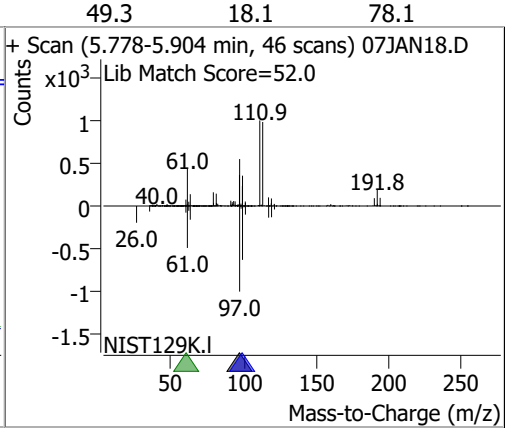




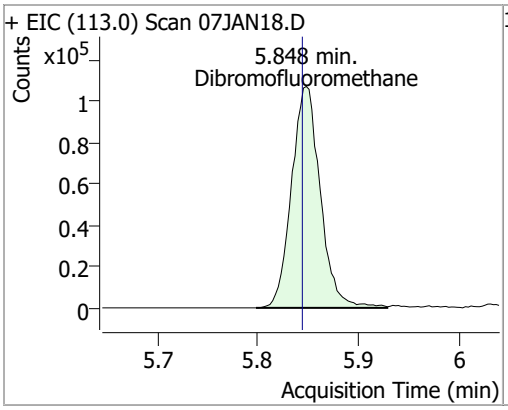
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	129.6747	5.83	0.00	196145	99.0	64.1	34.7	94.7
					61.0	49.3	18.1	78.1

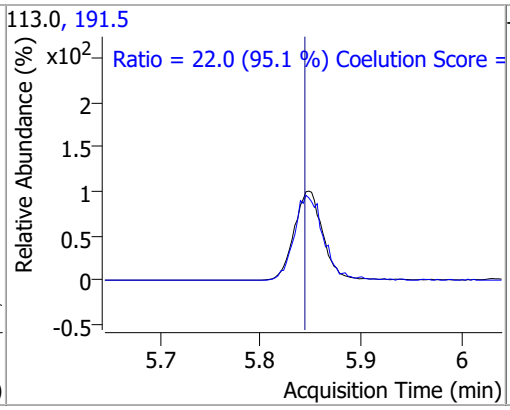


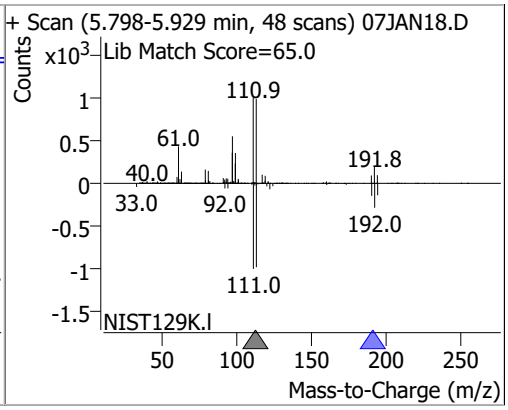




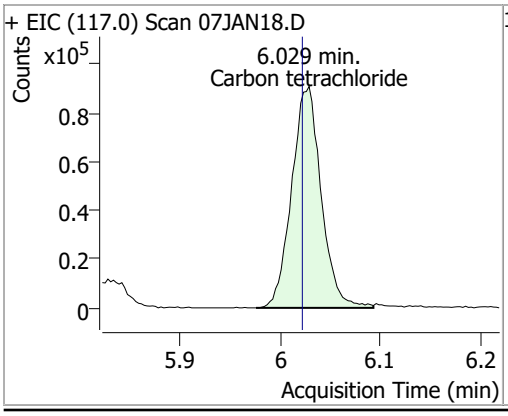
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	268.6027	5.85	0.00	214547	191.5	22.0	0.0	53.1

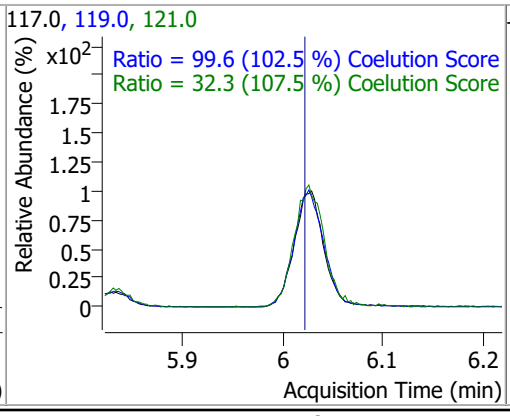


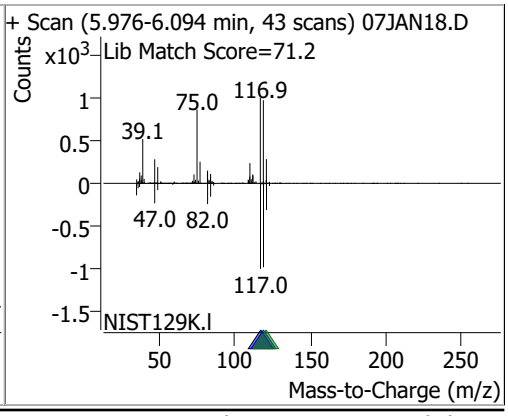




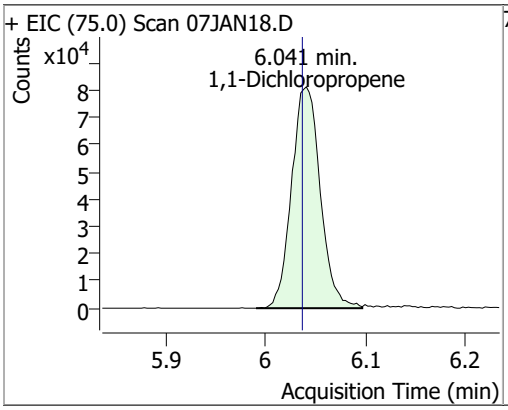
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	124.6502	6.03	0.01	185767	119.0	99.6	67.2	127.2
					121.0	32.3	0.1	60.1

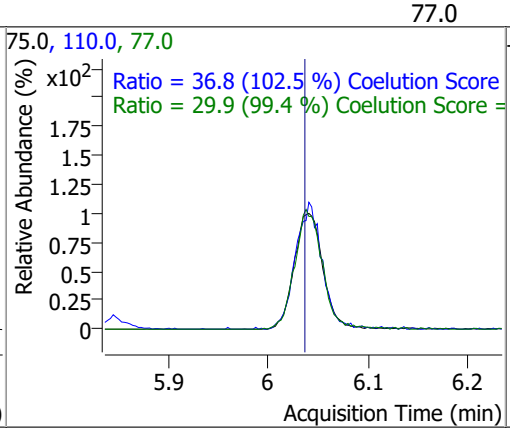


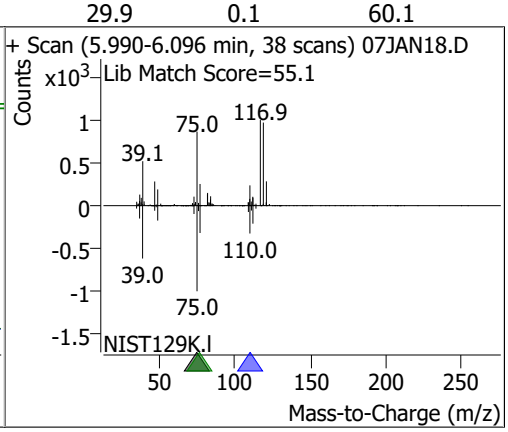




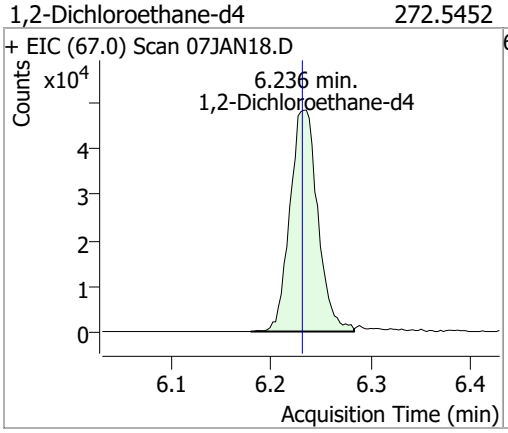
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	125.6088	6.04	0.00	161545	110.0	36.8	5.9	65.9
					77.0	29.9	0.1	60.1

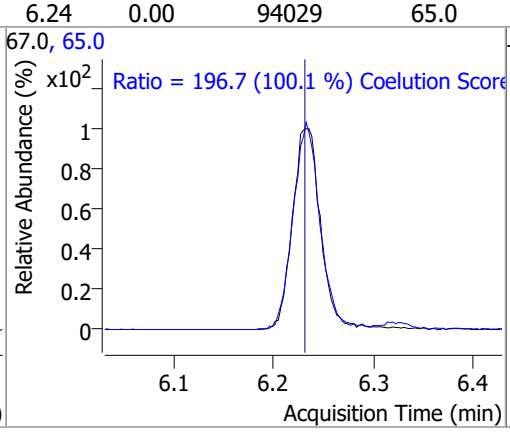


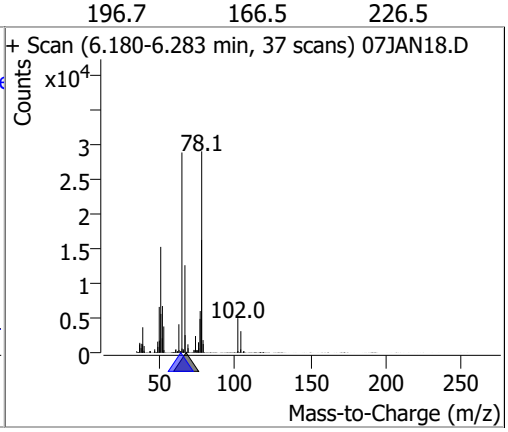




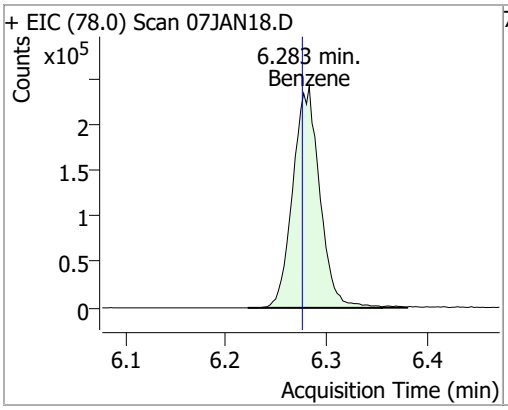
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	272.5452	6.24	0.00	94029	65.0	196.7	166.5	226.5

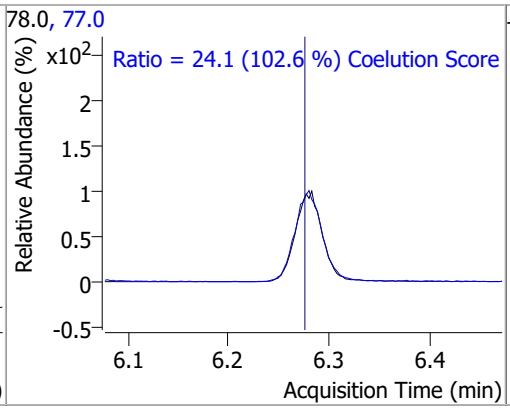


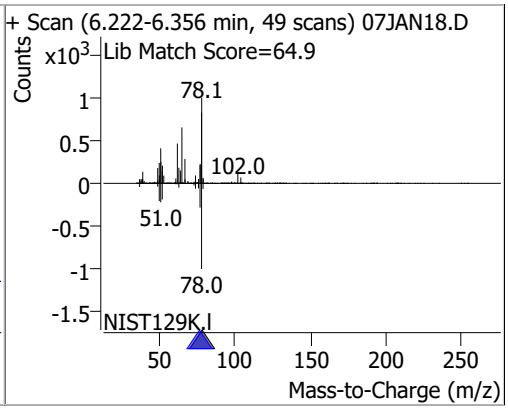




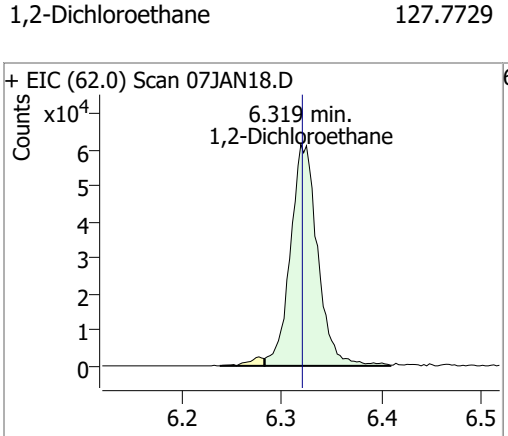
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	133.0746	6.28	0.01	449224	77.0	24.1	0.0	53.5

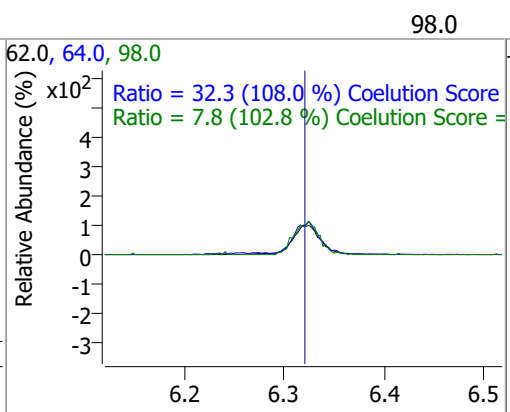


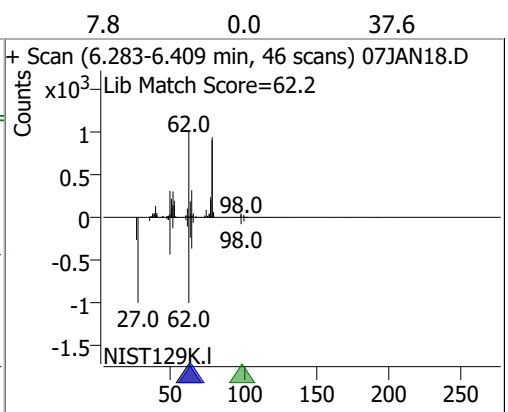




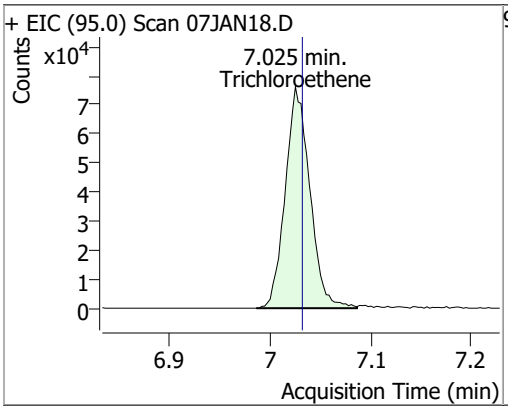
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	127.7729	6.32	0.00	116685	64.0	32.3	0.0	59.9
					98.0	7.8	0.0	37.6

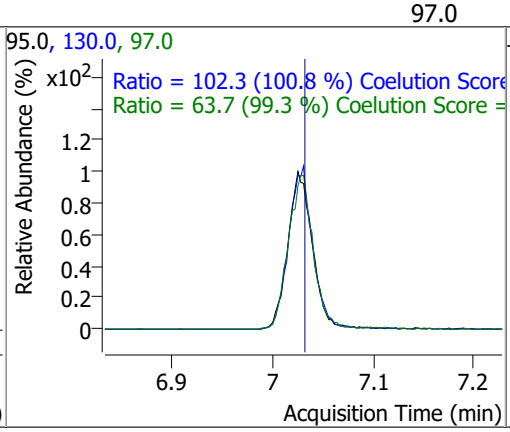


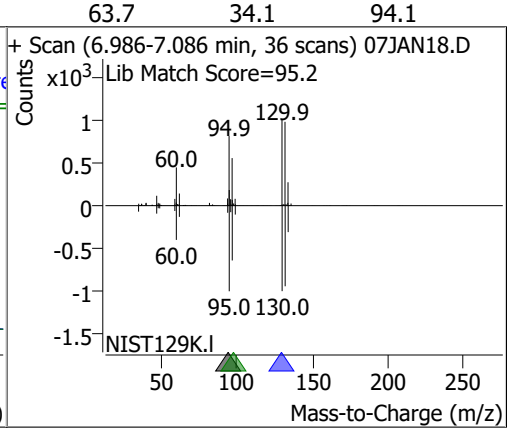




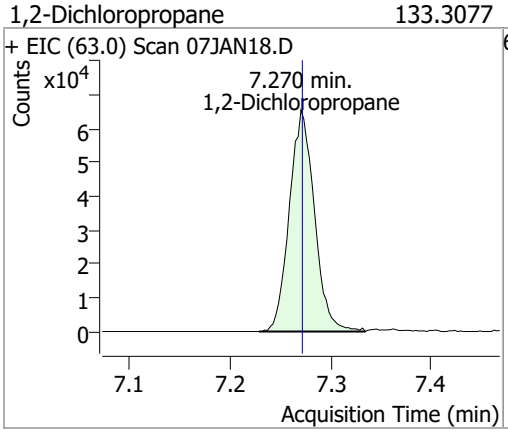
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	133.5436	7.02	-0.01	129711	130.0	102.3	71.5	131.5
					97.0	63.7	34.1	94.1

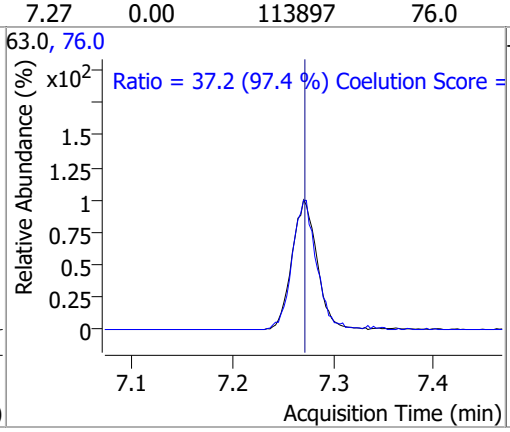


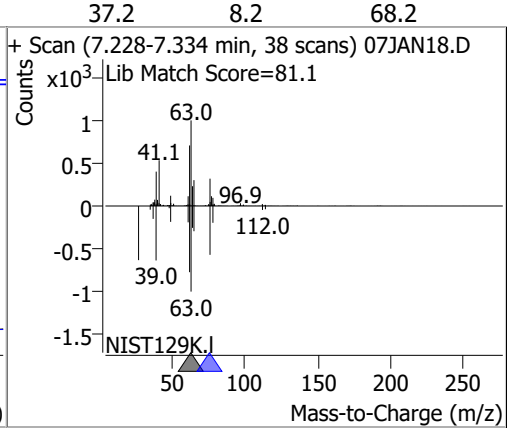




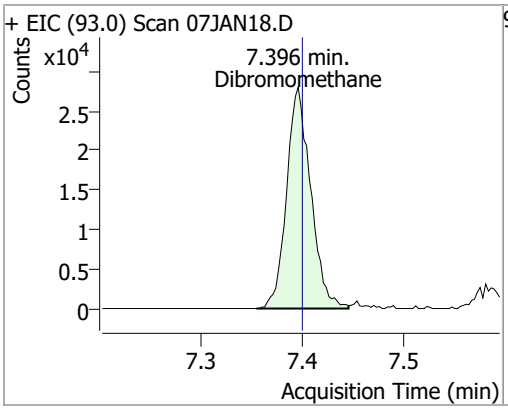
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	133.3077	7.27	0.00	113897	76.0	37.2	8.2	68.2

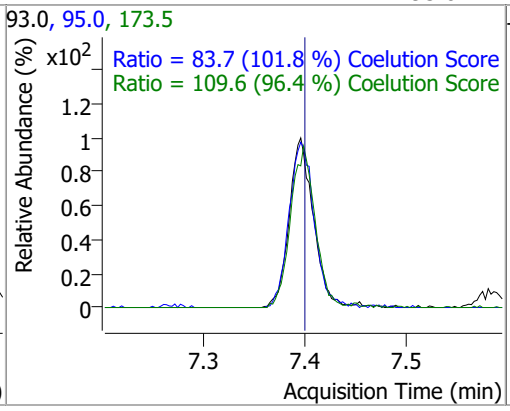


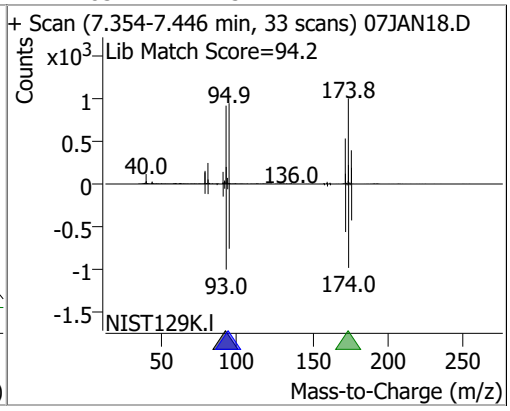




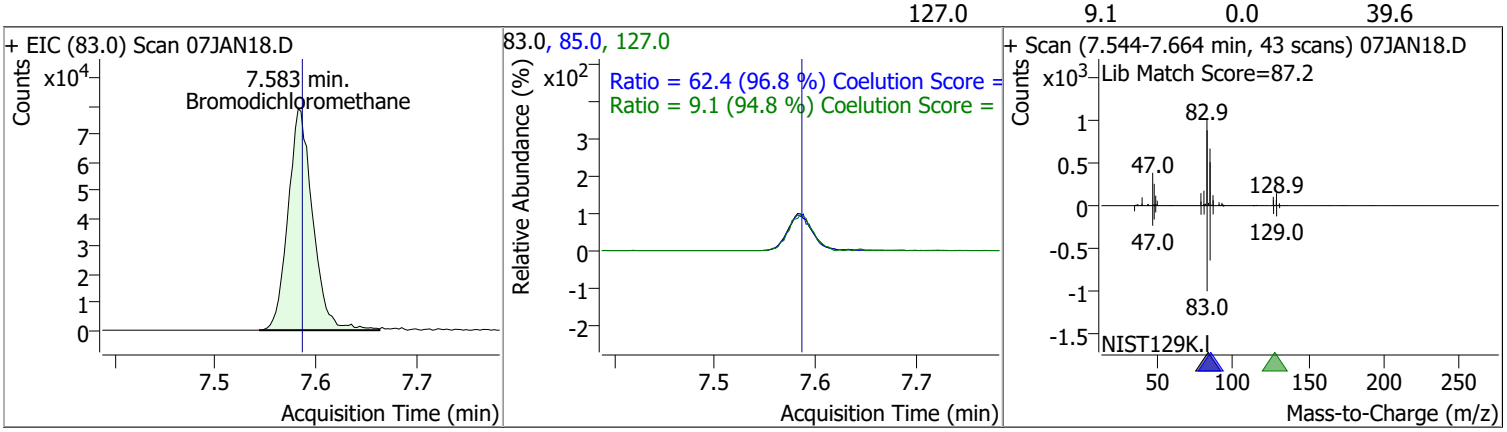
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	129.4203	7.40	0.00	46728	173.5	109.6	83.7	143.7
					95.0	83.7	52.2	112.2



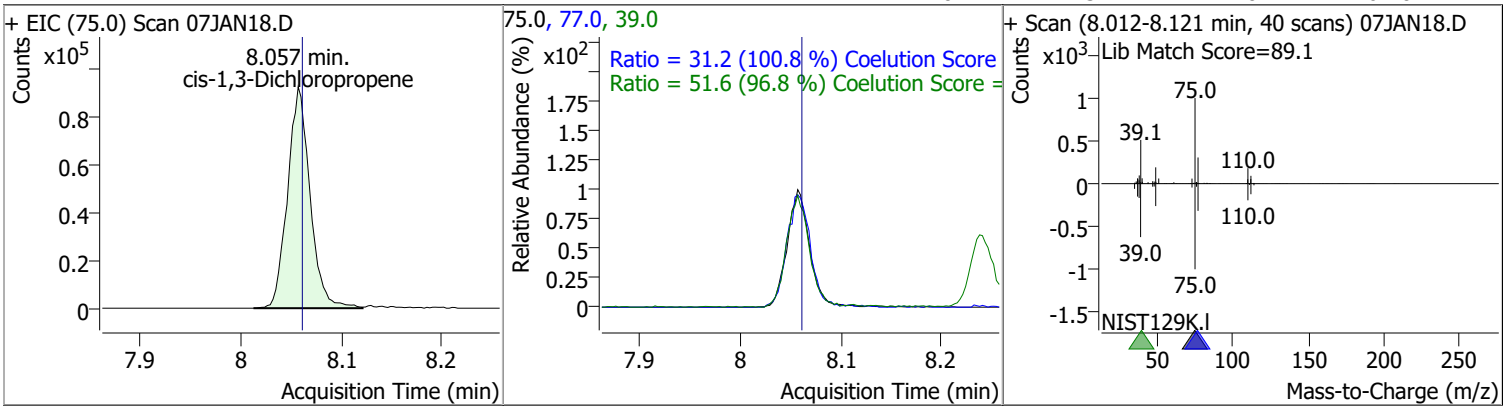




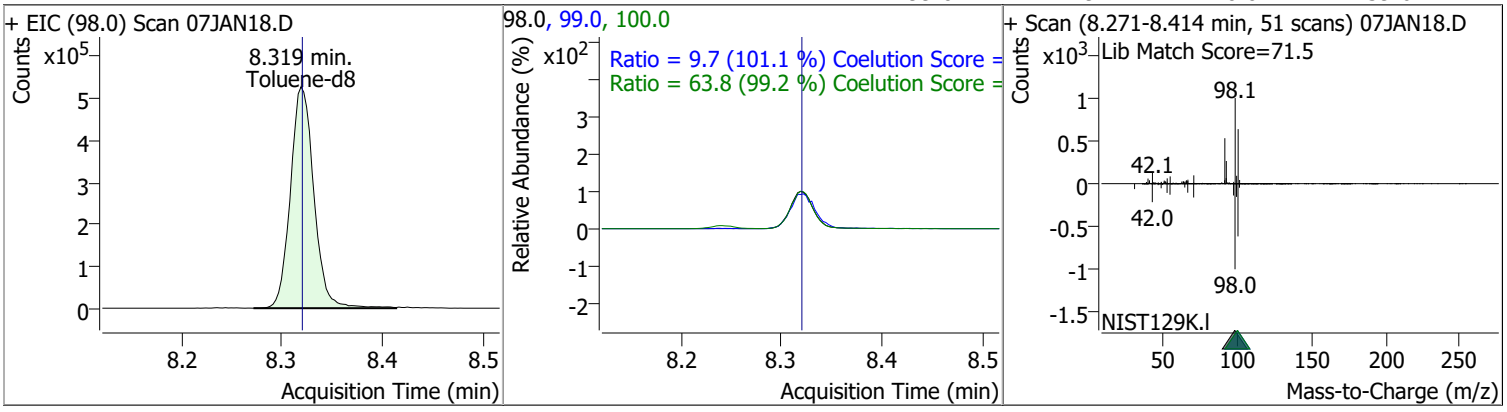
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	135.3871	7.58	0.00	134905	85.0	62.4	34.5	94.5
					127.0	9.1	0.0	39.6



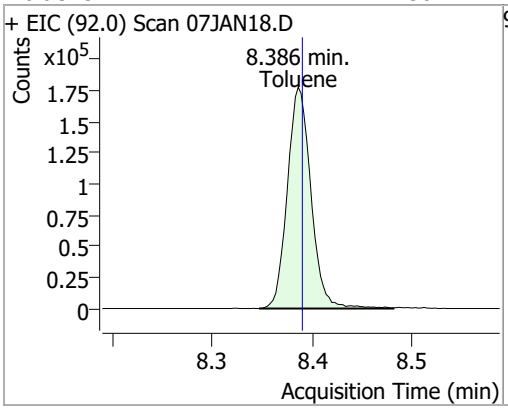
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	127.5477	8.06	0.00	143696	39.0	51.6	23.3	83.3
					77.0	31.2	1.0	61.0

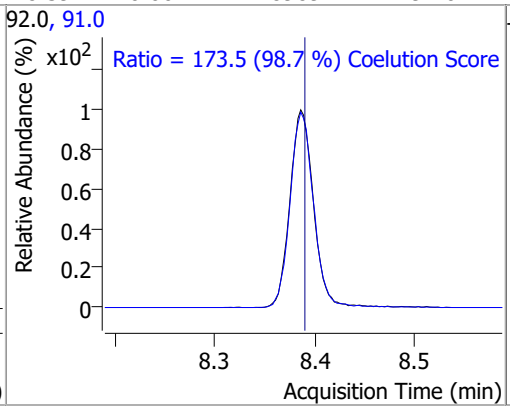


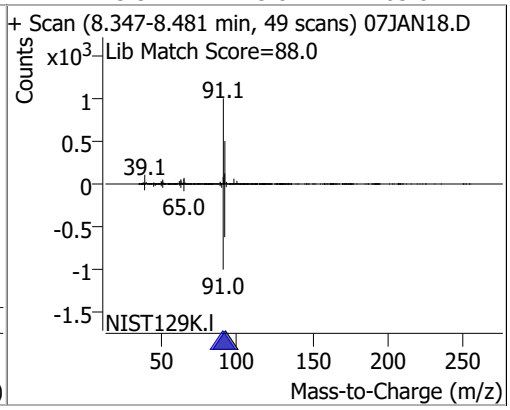
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	276.3920	8.32	0.00	857800	100.0	63.8	34.4	94.4
					99.0	9.7	0.0	39.6



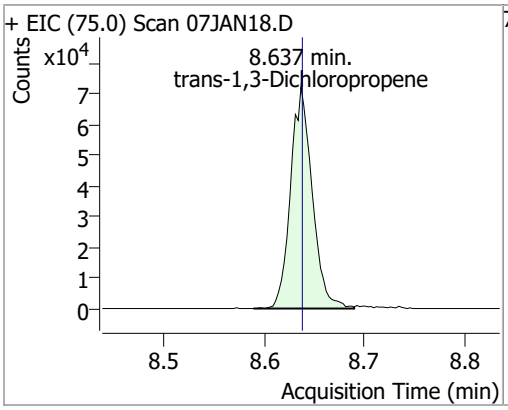
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	138.2772	8.39	0.00	289892	91.0	173.5	145.8	205.8

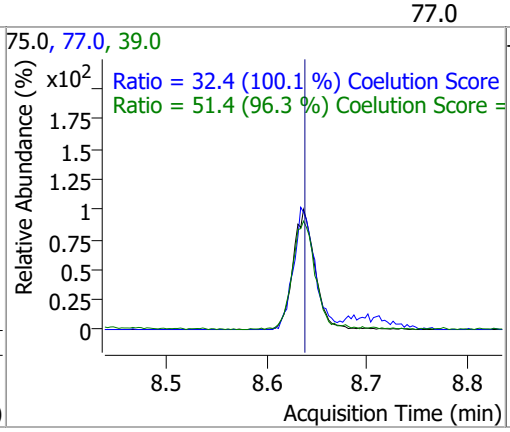


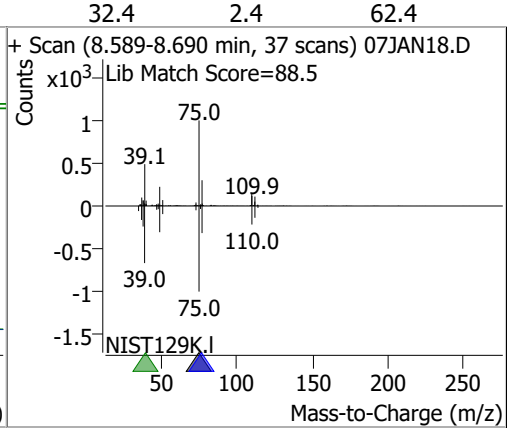




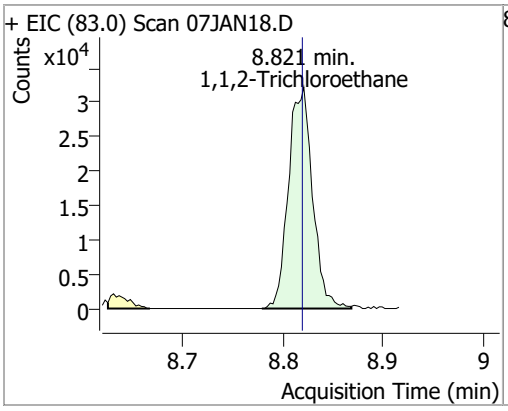
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	135.6987	8.64	0.00	108822	39.0	51.4	23.4	83.4
					77.0	32.4	2.4	62.4

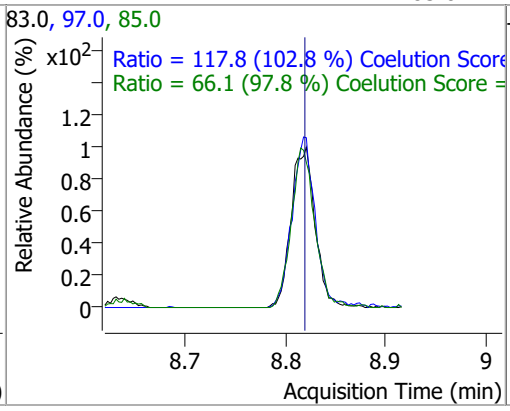


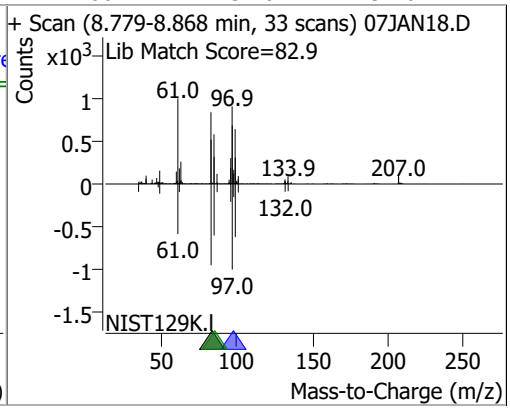




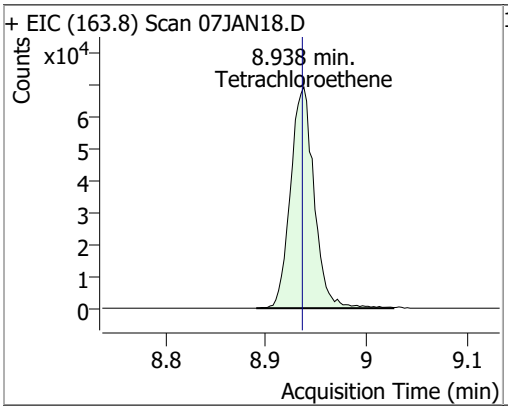
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	128.3504	8.82	0.00	53613	97.0	117.8	84.6	144.6
					85.0	66.1	37.6	97.6

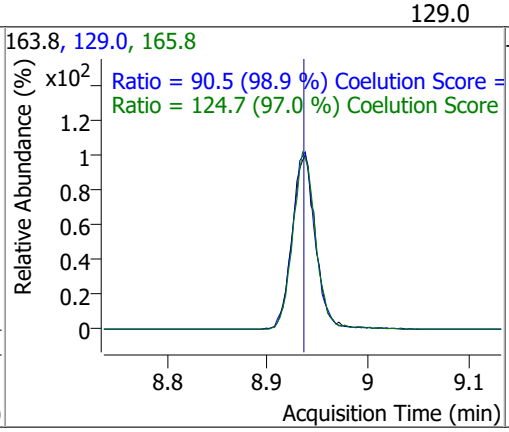


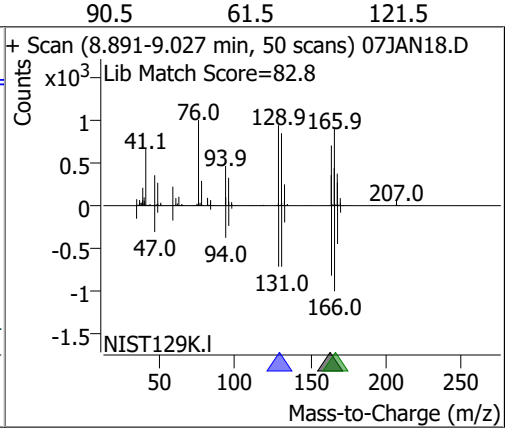




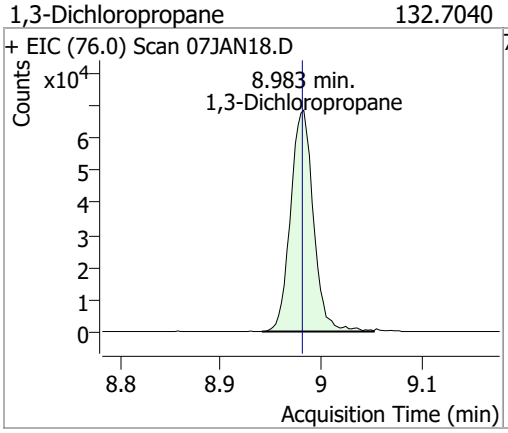
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	132.0855	8.94	0.00	112970	165.8	124.7	98.6	158.6
					129.0	90.5	61.5	121.5

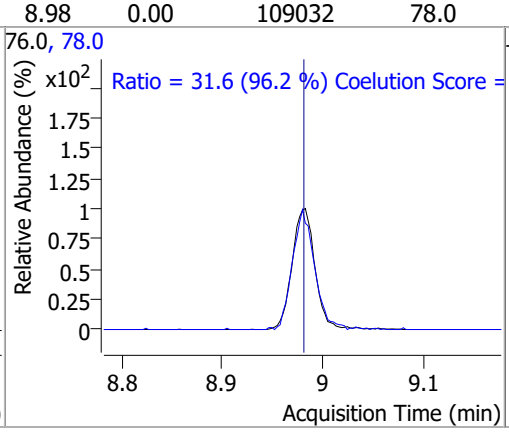


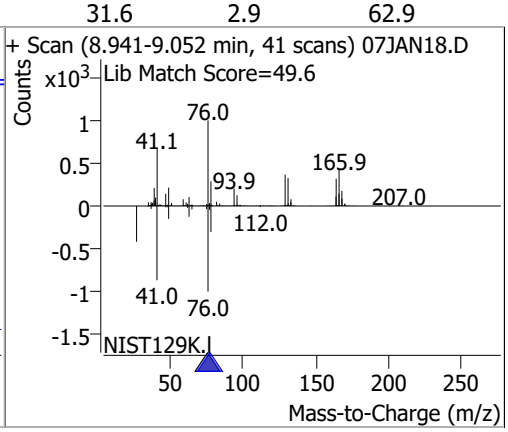




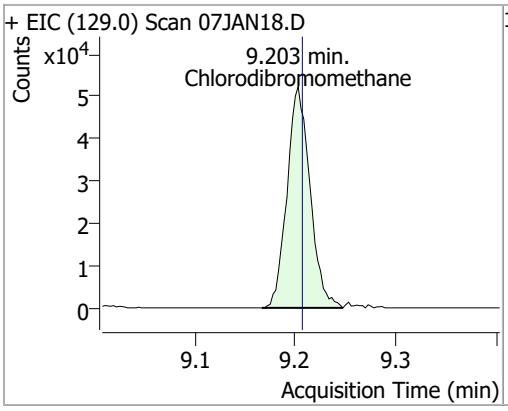
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	132.7040	8.98	0.00	109032	78.0	31.6	2.9	62.9

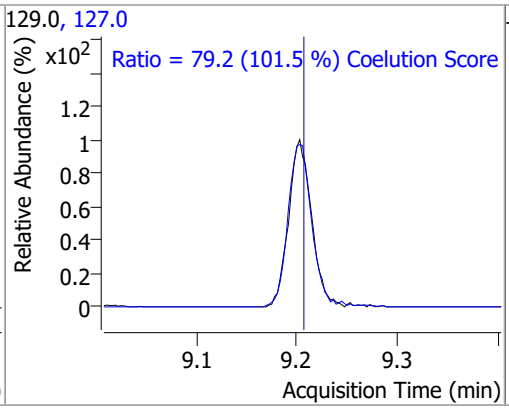


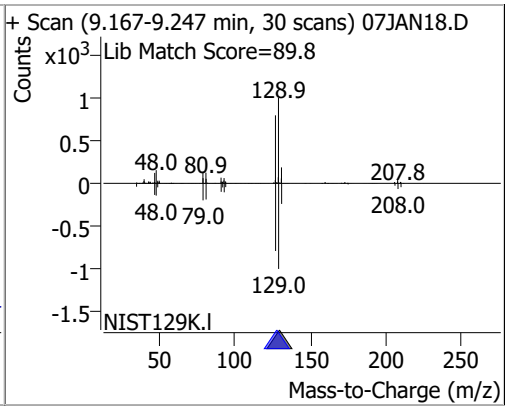




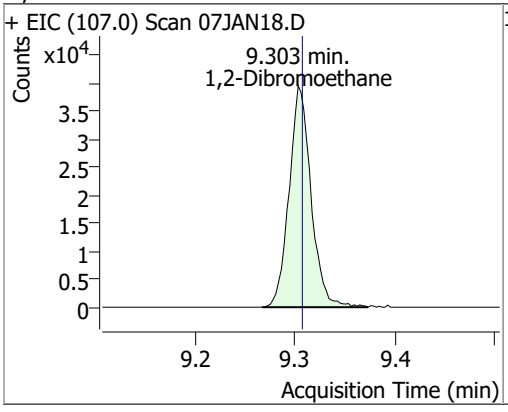
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	127.7253	9.20	0.00	83383	127.0	79.2	48.0	108.0

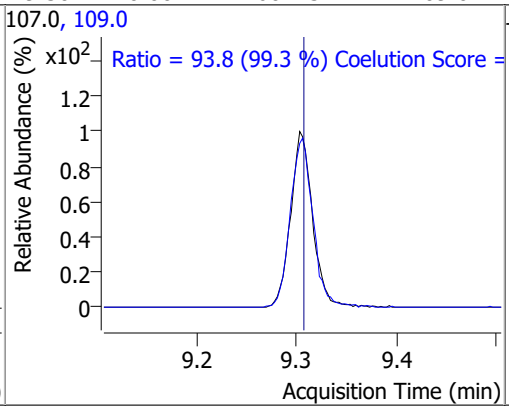


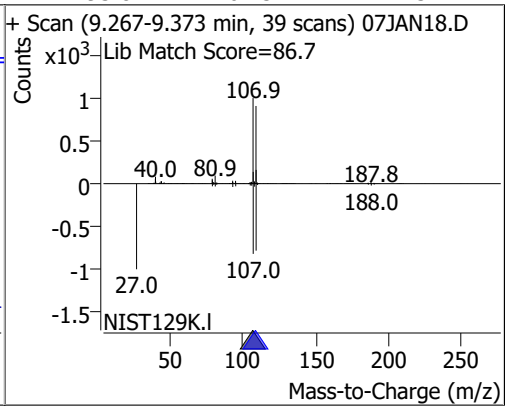




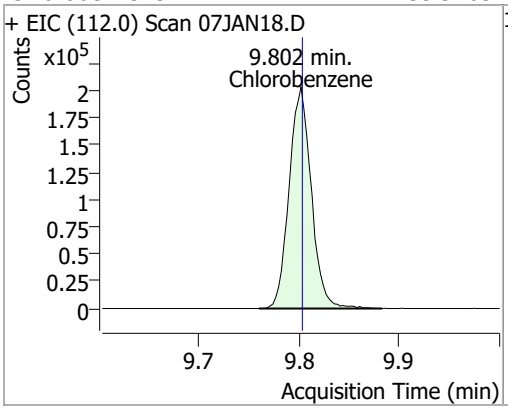
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	131.6380	9.30	0.00	60123	109.0	93.8	64.5	124.5

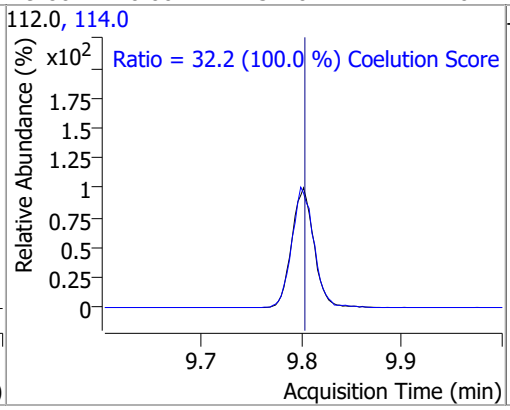


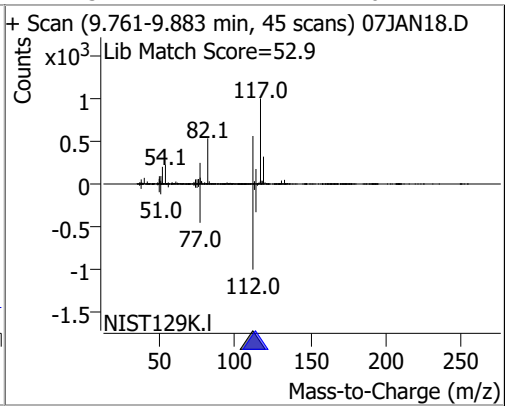




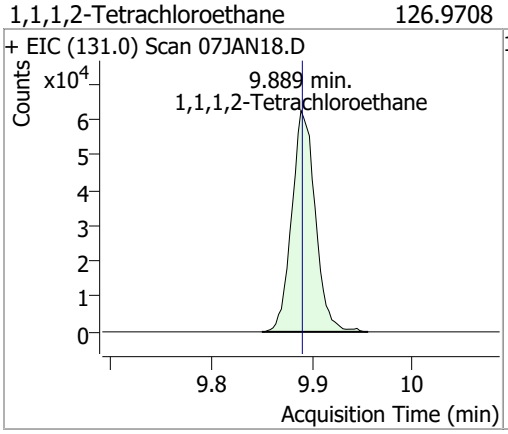
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	135.9465	9.80	0.00	312027	114.0	32.2	2.1	62.1

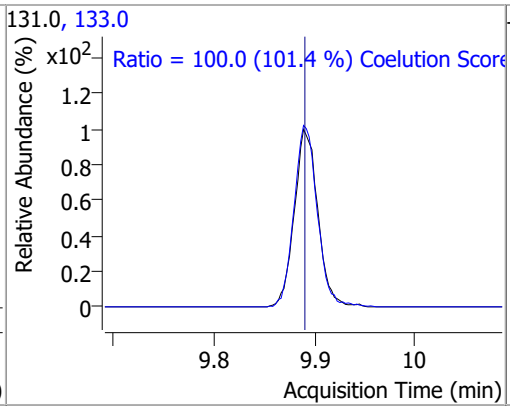


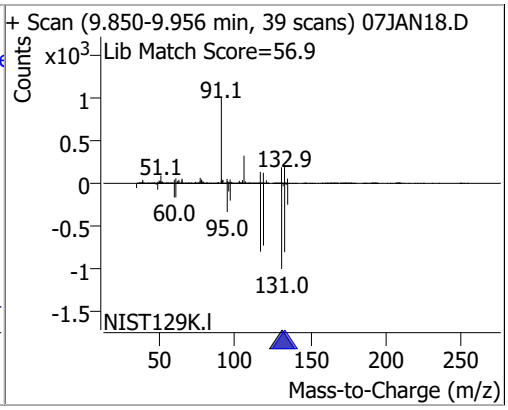




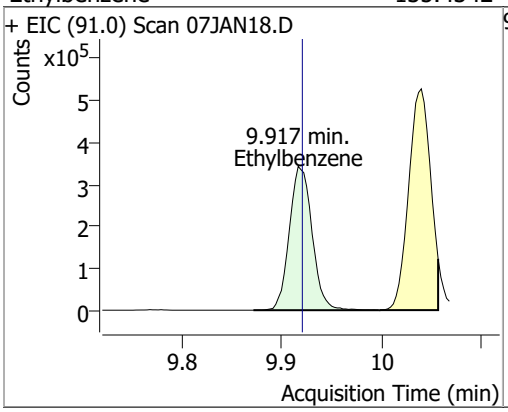
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	126.9708	9.89	0.00	101872	133.0	100.0	68.6	128.6

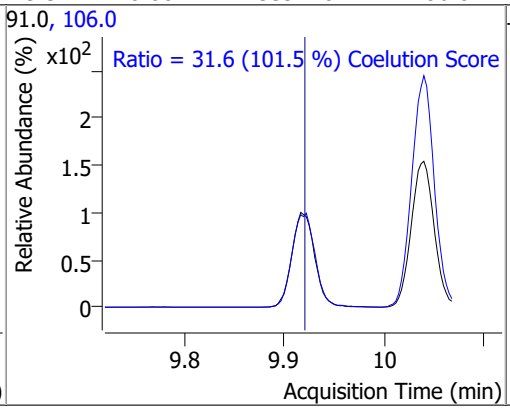


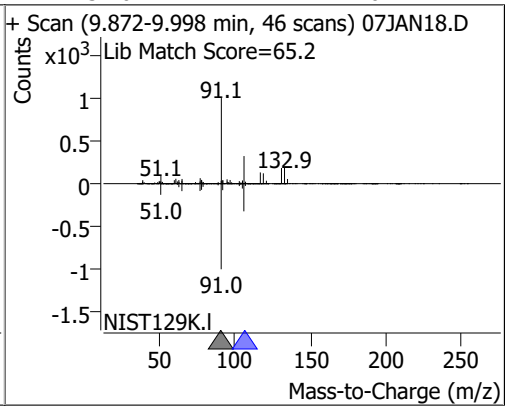




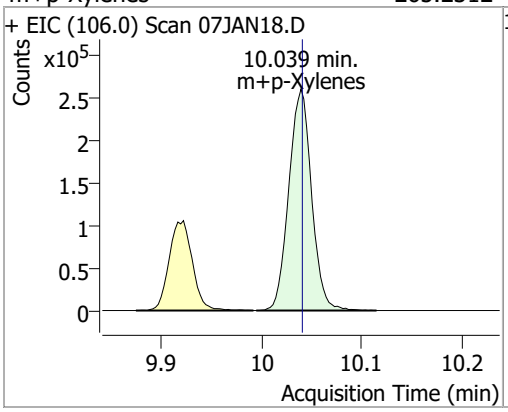
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	135.4342	9.92	0.00	539120	106.0	31.6	1.1	61.1

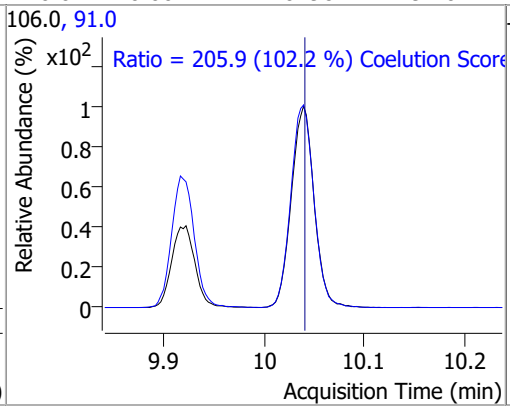


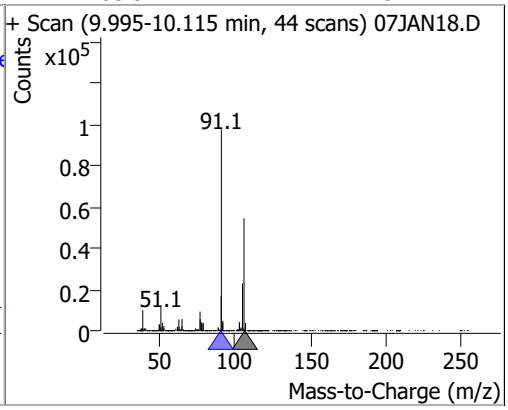




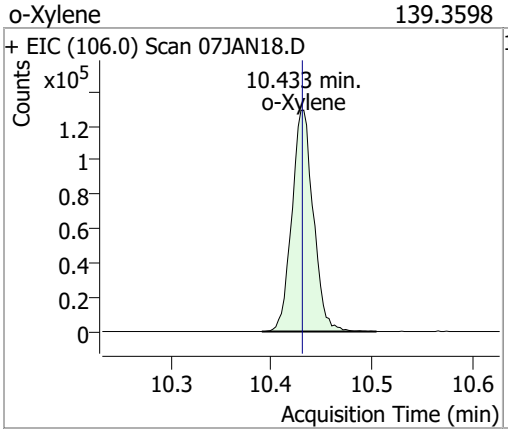
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	265.2312	10.04	0.00	410298	91.0	205.9	171.4	231.4

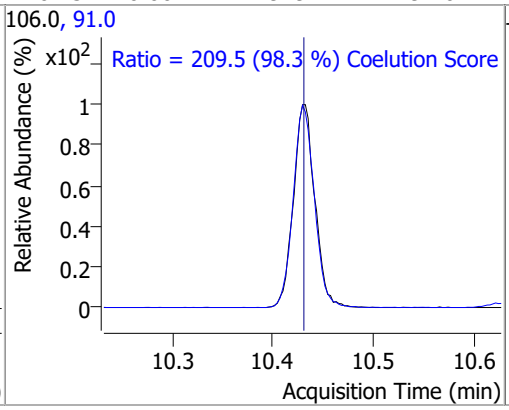


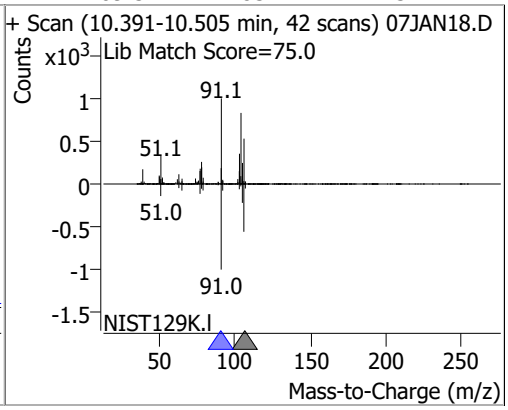




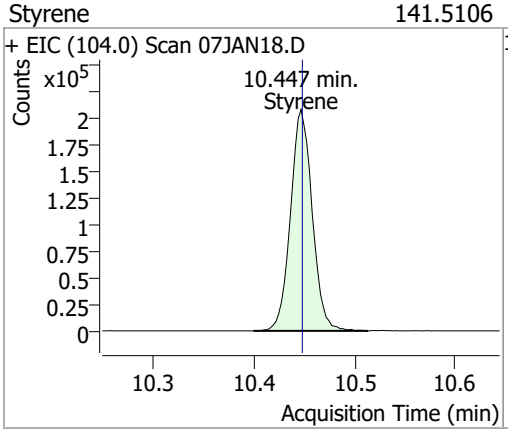
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	139.3598	10.43	0.00	191917	91.0	209.5	183.1	243.1

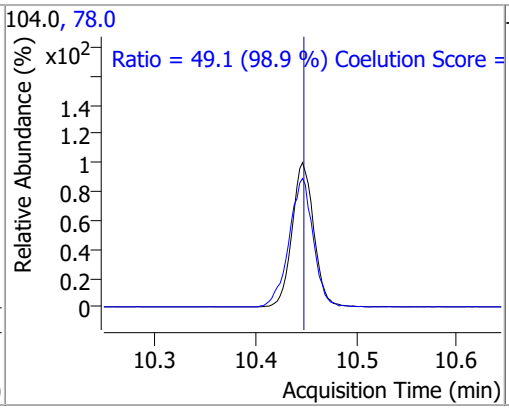


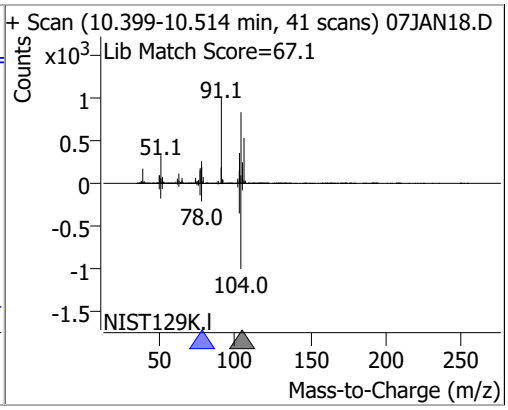




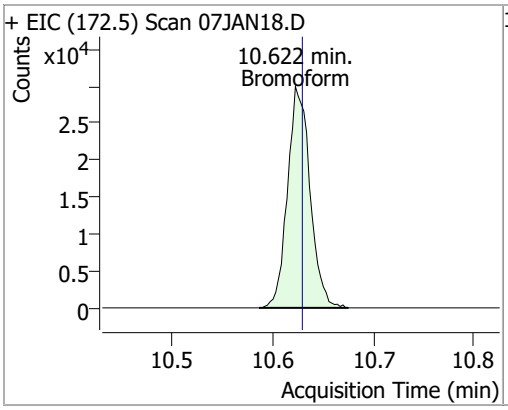
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	141.5106	10.45	0.00	313760	78.0	49.1	19.6	79.6

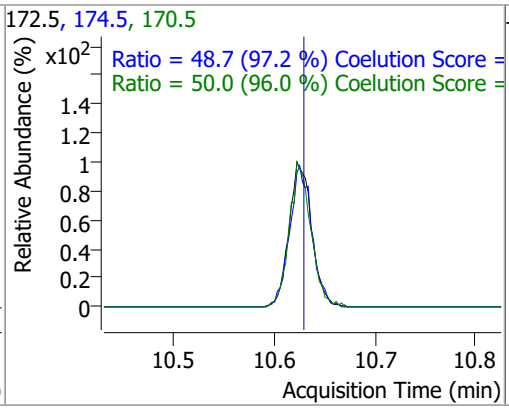


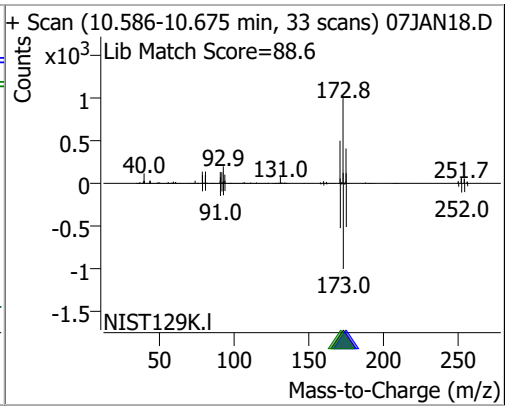




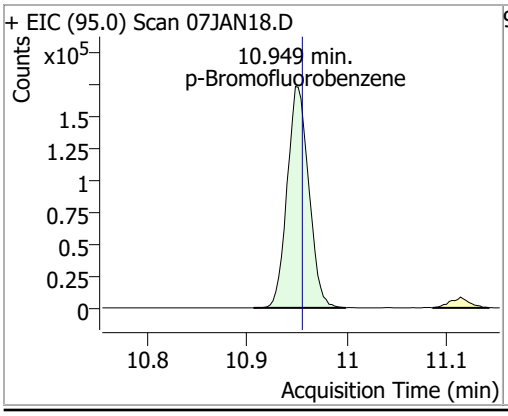
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	139.7801	10.62	-0.01	46577	170.5	50.0	22.1	82.1
					174.5	48.7	20.1	80.1

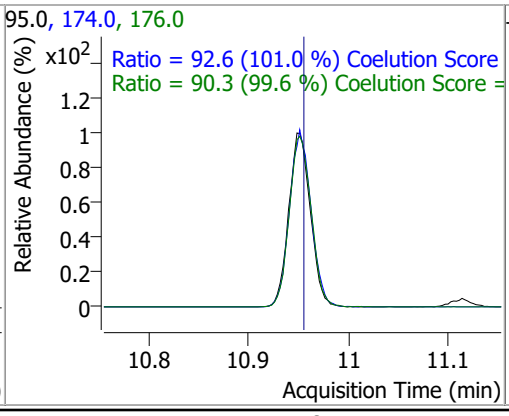


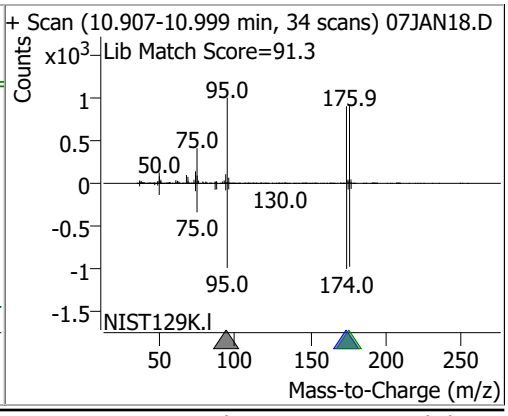




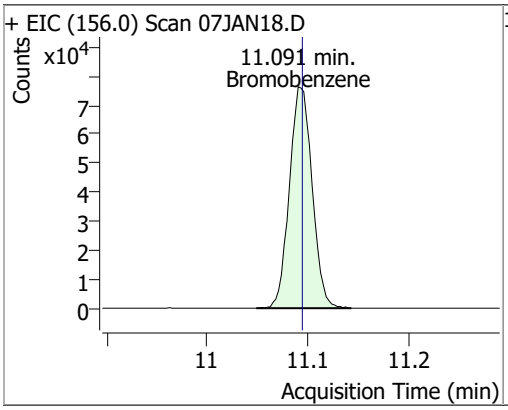
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	271.1727	10.95	-0.01	258687	174.0	92.6	61.7	121.7
					176.0	90.3	60.6	120.6

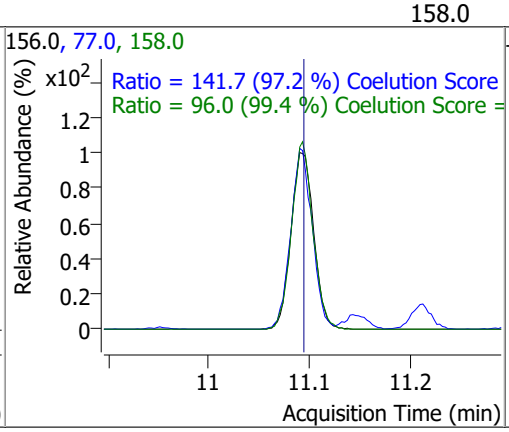


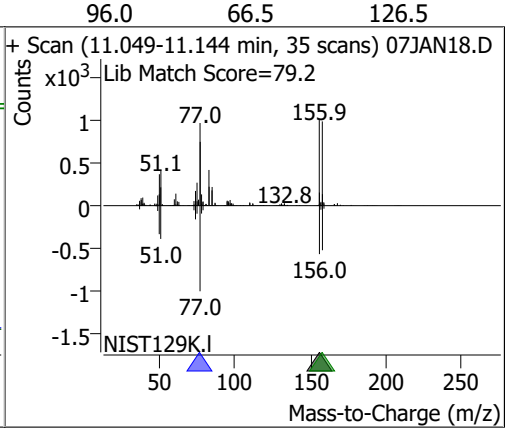




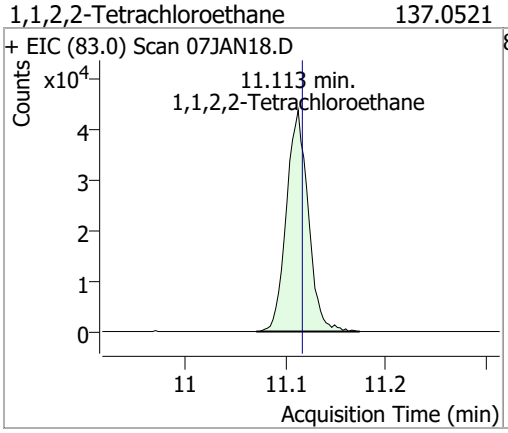
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	142.5401	11.09	0.00	120119	77.0	141.7	115.7	175.7
					158.0	96.0	66.5	126.5

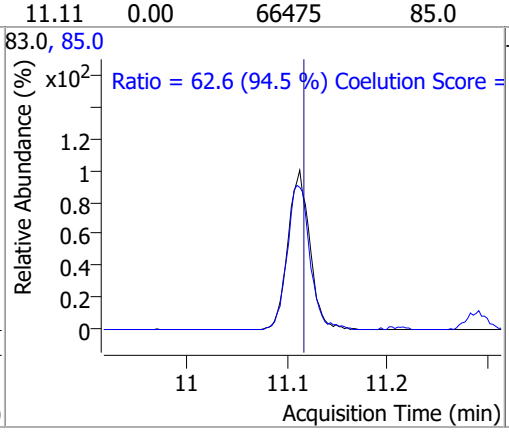


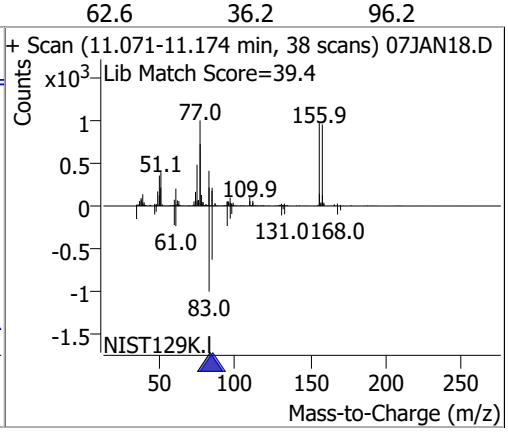




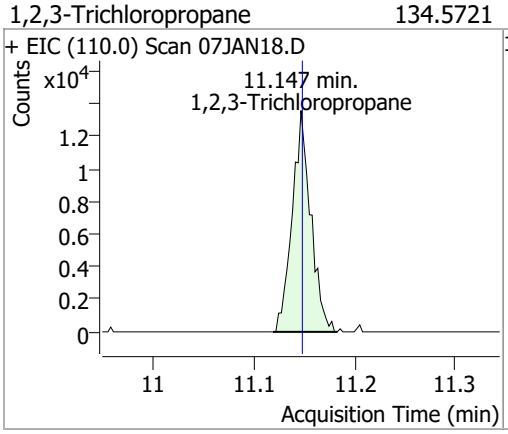
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	137.0521	11.11	0.00	66475	85.0	62.6	36.2	96.2

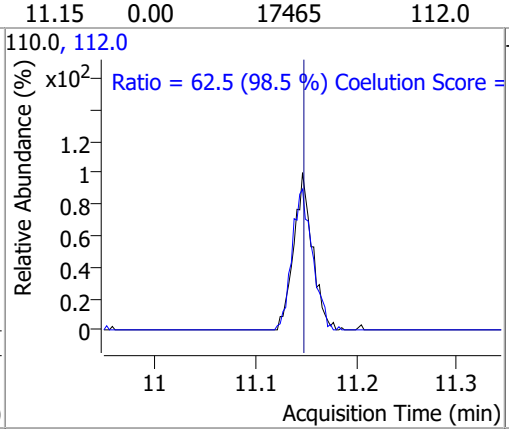


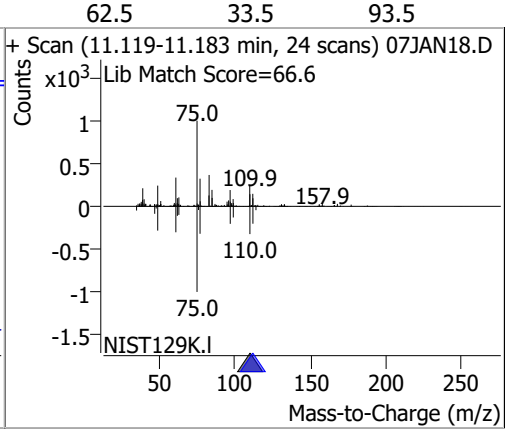




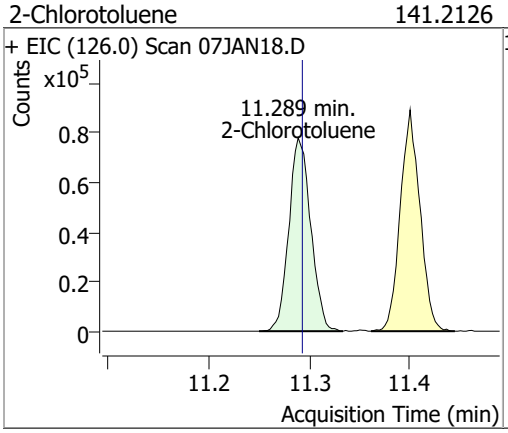
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	134.5721	11.15	0.00	17465	112.0	62.5	33.5	93.5

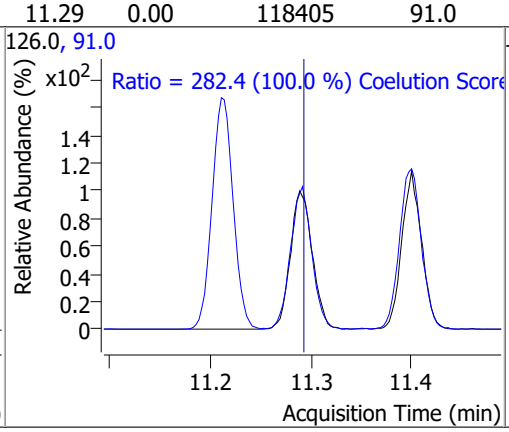


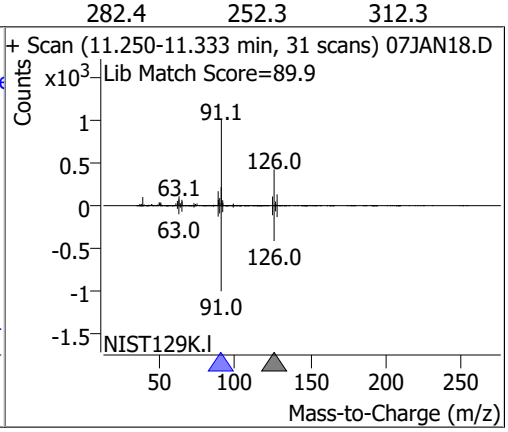




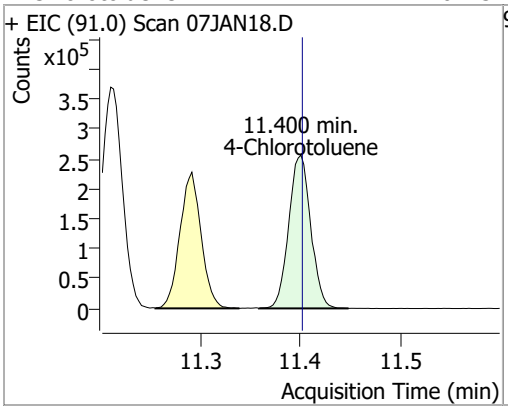
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	141.2126	11.29	0.00	118405	91.0	282.4	252.3	312.3

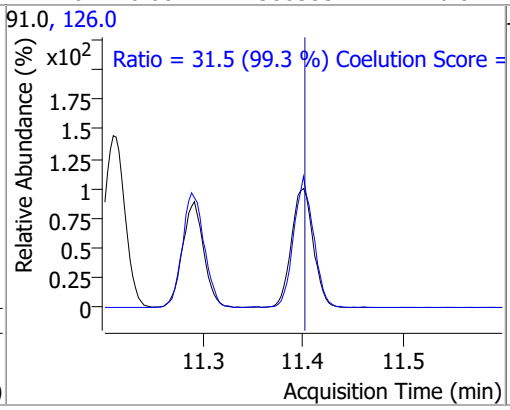


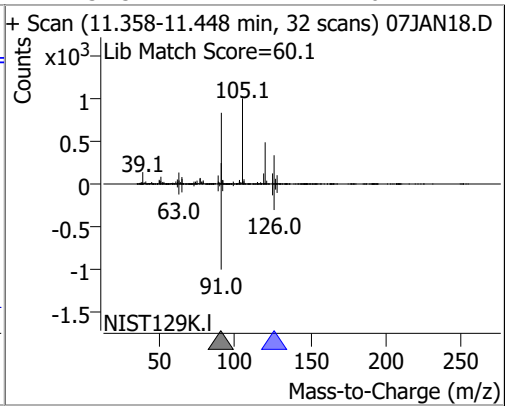




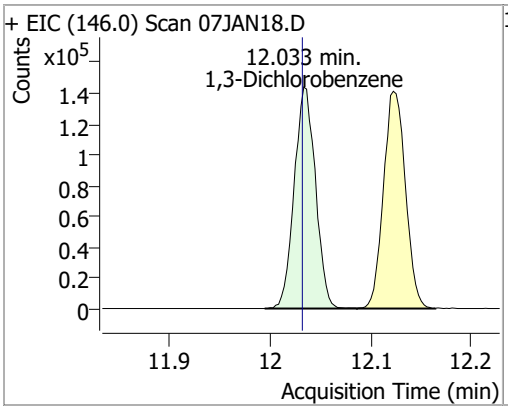
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	142.0473	11.40	0.00	388335	126.0	31.5	1.7	61.7

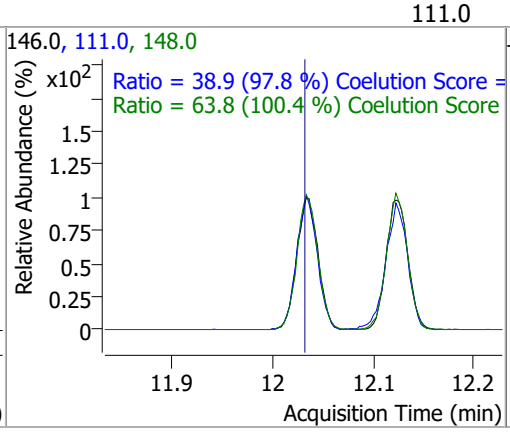


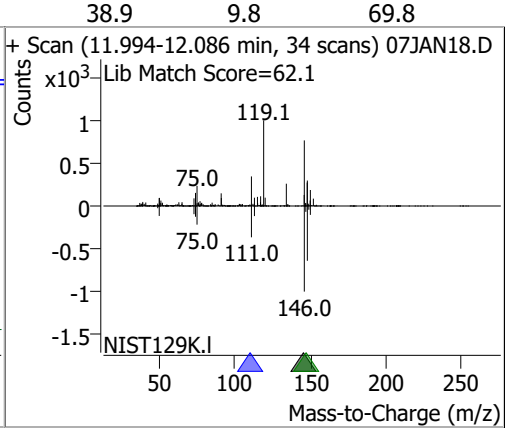




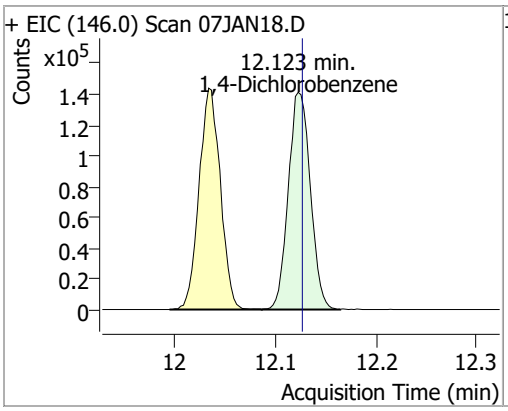
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	137.9499	12.03	0.00	212018	148.0	63.8	33.6	93.6
					111.0	38.9	9.8	69.8

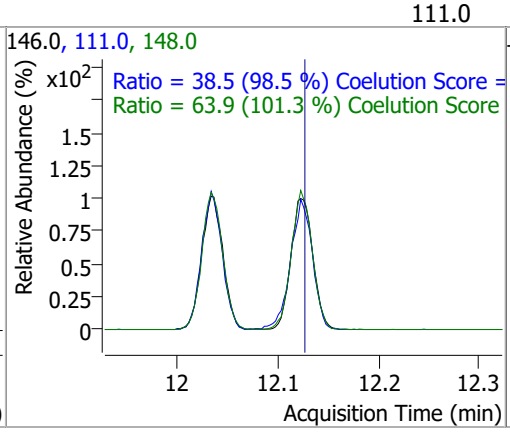


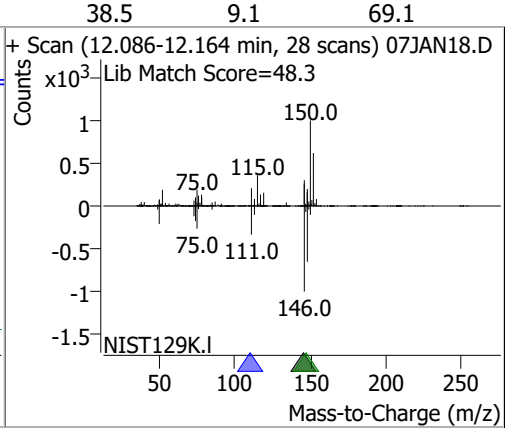




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	137.4287	12.12	0.00	215367	148.0	63.9	33.1	93.1
					111.0	38.5	9.1	69.1

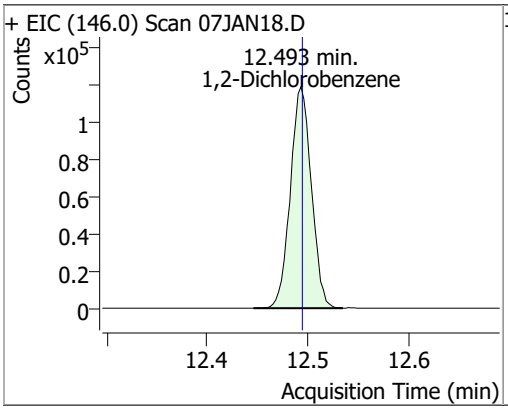


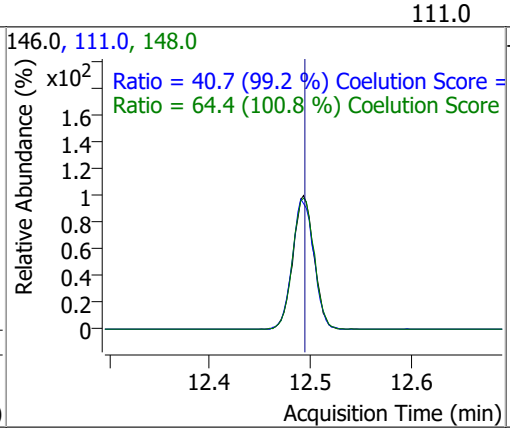


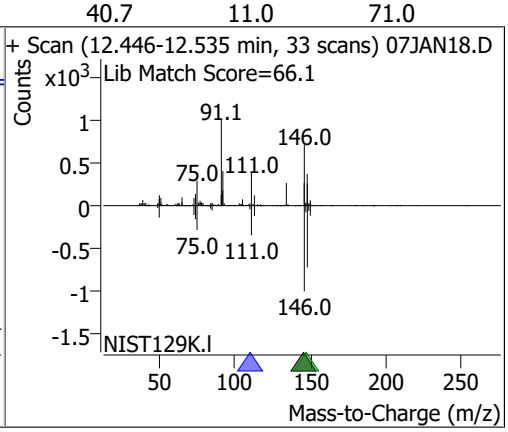


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	132.9681	12.49	0.00	172710	148.0	64.4	33.9	93.9
					111.0	40.7	11.0	71.0







Quantitation Results Report (QT Reviewed)

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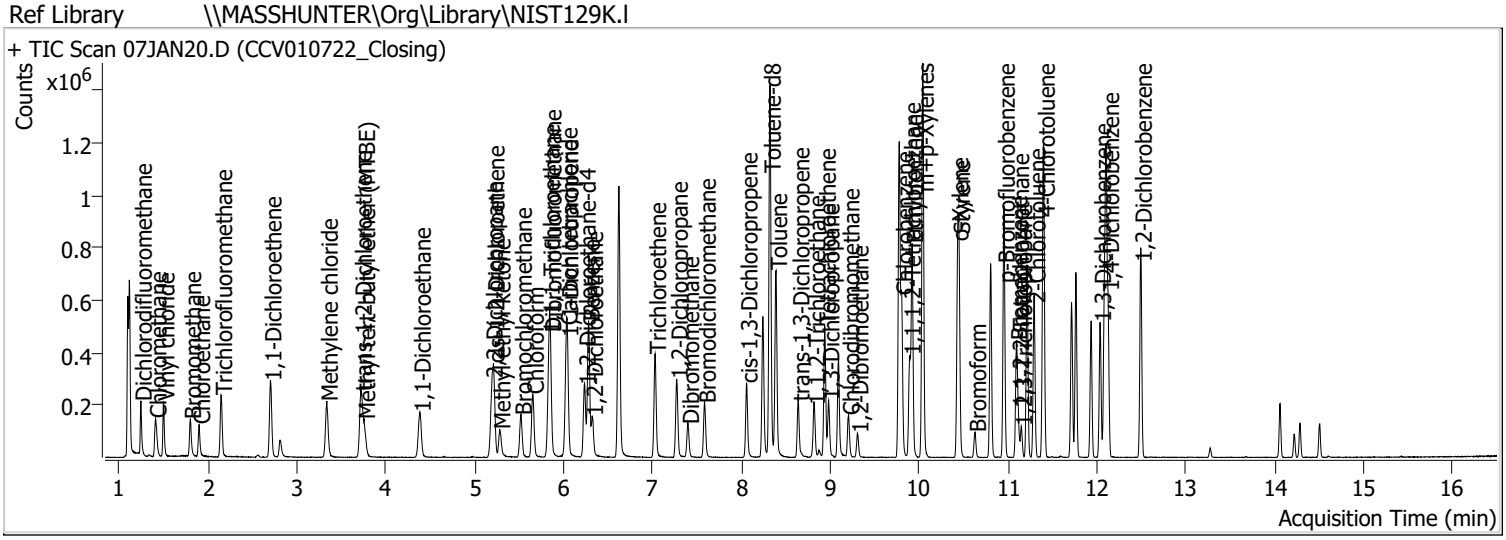
Data File	07JAN20.D	Operator	MSC
Acq. Method	5975CACQF.M	Acq. Date-Time	1/7/2022 6:29:09 PM
Sample Name	CCV010722_Closing	Instrument	VOA5975C
Vial	20	Multiplier	1.00
DA Method File	VOA5975C_8260B_SHT_DoD_L4_010422.m	Comment	
Tune File	BFB_Atune3.u	Tune Date	10/11/2021 4:02:00 PM
Batch Name	VG010722_8260B.batch.bin	Last Calib Update	3/2/2022 10:41:44 AM
Ref Library	\\MASSHUNTER\Org\Library\NIST129K.l		

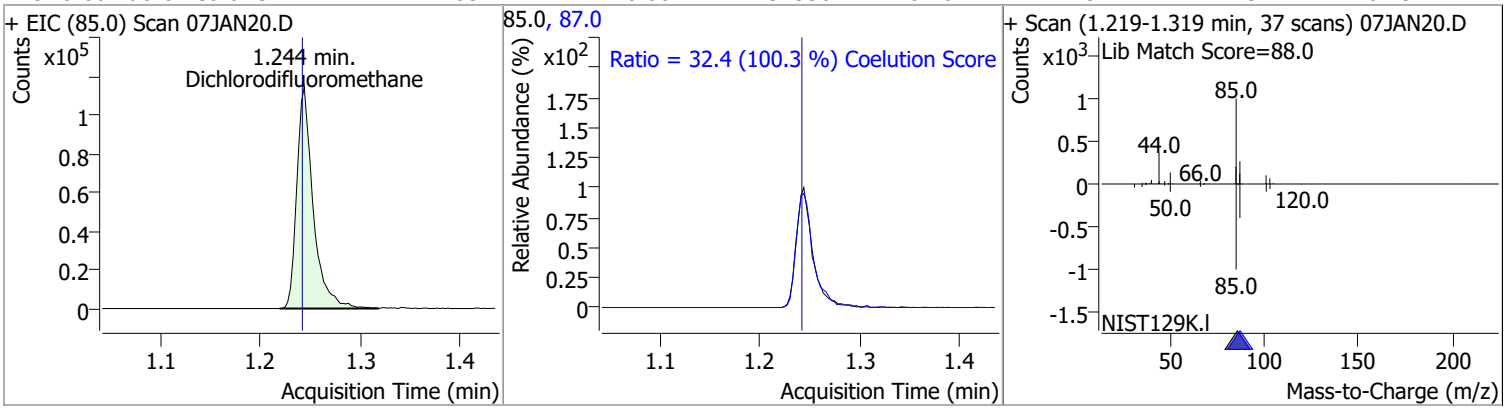


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
M Fluorobenzene	6.621	96.0	875212	250.0000	ng	-0.003
M Chlorobenzene-d5	9.775	82.0	332712	250.0000	ng	0.003
M 1,4-Dichlorobenzene-d4	12.100	152.0	273694	250.0000	ng	0.000
System Monitoring Compounds						
S Dibromofluoromethane	5.851	113.0	220533	267.4624	ng	0.006
Spiked Amount: 250.000	Range: 80.0 - 119.0%			Recovery = 106.98%		
S 1,2-Dichloroethane-d4	6.236	67.0	97554	273.9195	ng	0.003
Spiked Amount: 250.000	Range: 81.0 - 118.0%			Recovery = 109.57%		
S Toluene-d8	8.319	98.0	873247	272.3635	ng	0.000
Spiked Amount: 250.000	Range: 89.0 - 112.0%			Recovery = 108.95%		
S p-Bromofluorobenzene	10.951	95.0	269134	268.4143	ng	-0.003
Spiked Amount: 250.000	Range: 85.0 - 114.0%			Recovery = 107.37%		
Target Compounds						QValue
T Dichlorodifluoromethane	1.244	85.0	131556	114.7051	ng	100
T Chloromethane	1.409	50.0	159080	114.2765	ng	99
T Vinyl chloride	1.498	62.0	147497	117.7541	ng	98
T Bromomethane	1.799	96.0	69603	124.2700	ng	98
T Chloroethane	1.899	64.0	75250	121.3507	ng	98
T Trichlorofluoromethane	2.148	101.0	163895	105.4169	ng	98
T 1,1-Dichloroethene	2.700	96.0	104720	118.7865	ng	100
T Methylene chloride	3.336	49.0	149219	114.8197	ng	98
T trans-1,2-Dichloroethene	3.718	96.0	108710	120.8683	ng	97
T Methyl tert-butyl ether (MTBE)	3.754	73.0	138605	119.2255	ng	99
T 1,1-Dichloroethane	4.381	63.0	209364	125.0569	ng	98
T 2,2-Dichloropropane	5.193	77.0	152733	121.7521	ng	99
T cis-1,2-Dichloroethene	5.210	96.0	111394	122.1596	ng	99
T Methyl ethyl ketone	5.279	43.0	144340	1168.5921	ng	100
T Bromochloromethane	5.516	128.0	43567	115.3287	ng	95
T Chloroform	5.653	83.0	197264	118.3968	ng	96

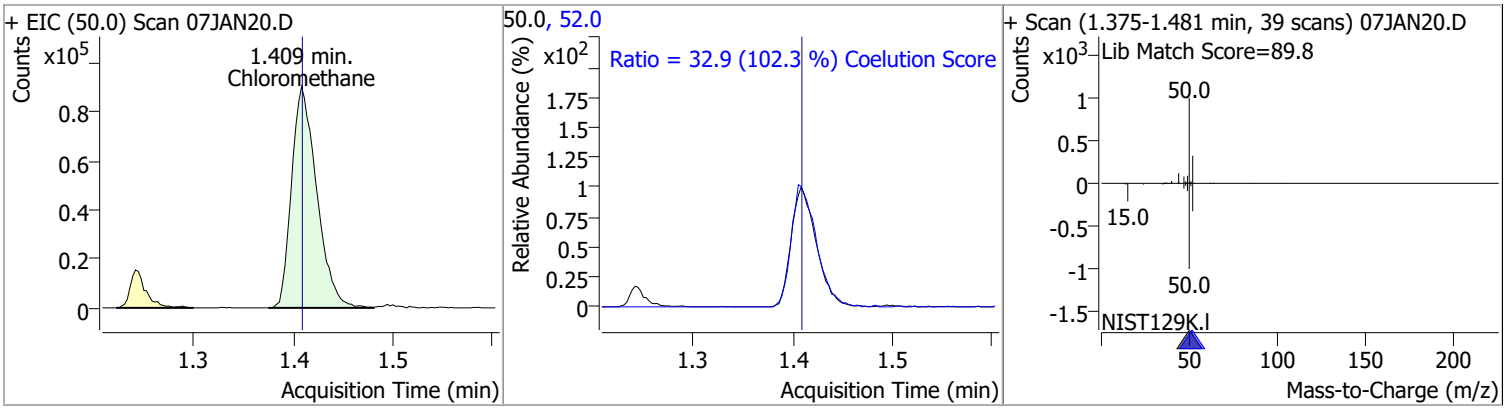
Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
T 1,1,1-Trichloroethane	5.834	97.0	185675	118.9139	ng	99
T Carbon tetrachloride	6.029	117.0	180946	117.6182	ng	98
T 1,1-Dichloropropene	6.043	75.0	159136	119.8660	ng	99
T Benzene	6.280	78.0	426971	122.5270	ng	99
T 1,2-Dichloroethane	6.319	62.0	108503	115.0977	ng	96
T Trichloroethene	7.028	95.0	122705	122.2872	ng	99
T 1,2-Dichloropropane	7.270	63.0	109262	123.7897	ng	98
T Dibromomethane	7.399	93.0	45080	120.8597	ng	96
T Bromodichloromethane	7.585	83.0	121732	118.2568	ng	96
T cis-1,3-Dichloropropene	8.057	75.0	138637	119.1185	ng	99
T Toluene	8.389	92.0	268266	123.8661	ng	98
T trans-1,3-Dichloropropene	8.640	75.0	103086	124.4317	ng	99
T 1,1,2-Trichloroethane	8.818	83.0	50481	116.9843	ng	99
T Tetrachloroethene	8.935	163.8	105808	119.7520	ng	98
T 1,3-Dichloropropane	8.980	76.0	103169	121.5491	ng	99
T Chlorodibromomethane	9.203	129.0	79716	118.1999	ng	98
T 1,2-Dibromoethane	9.303	107.0	56624	120.0089	ng	94
T Chlorobenzene	9.802	112.0	287084	121.0758	ng	99
T 1,1,1,2-Tetrachloroethane	9.892	131.0	96577	116.5186	ng	97
T Ethylbenzene	9.920	91.0	500217	121.6392	ng	99
T m+p-Xylenes	10.039	106.0	396787	248.2876	ng	99
T o-Xylene	10.430	106.0	175221	123.1636	ng	100
T Styrene	10.447	104.0	289229	126.2715	ng	100
T Bromoform	10.628	172.5	42320	120.8329	ng	99
T Bromobenzene	11.094	156.0	110486	124.7379	ng	98
T 1,1,2,2-Tetrachloroethane	11.113	83.0	61840	121.3005	ng	96
T 1,2,3-Trichloropropane	11.147	110.0	15684	114.9765	ng	96
T 2-Chlorotoluene	11.294	126.0	109275	123.9909	ng	98
T 4-Chlorotoluene	11.400	91.0	363039	126.3413	ng	99
T 1,3-Dichlorobenzene	12.033	146.0	196110	121.3987	ng	100
T 1,4-Dichlorobenzene	12.123	146.0	198916	120.7629	ng	98
T 1,2-Dichlorobenzene	12.493	146.0	162097	118.7328	ng	100

(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

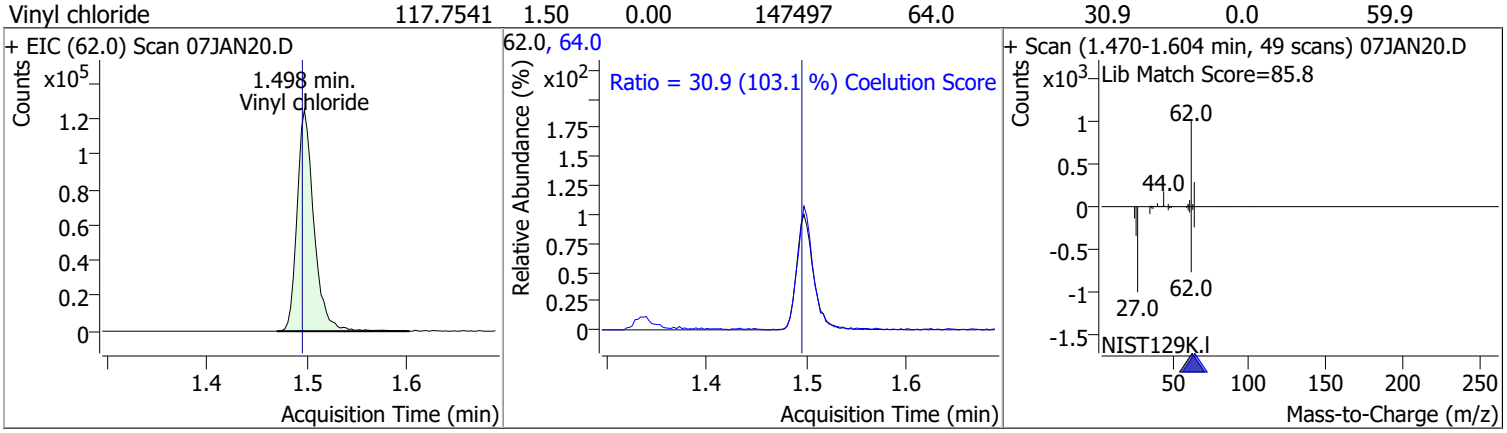
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dichlorodifluoromethane	114.7051	1.24	0.00	131556	87.0	32.4	2.3	62.3



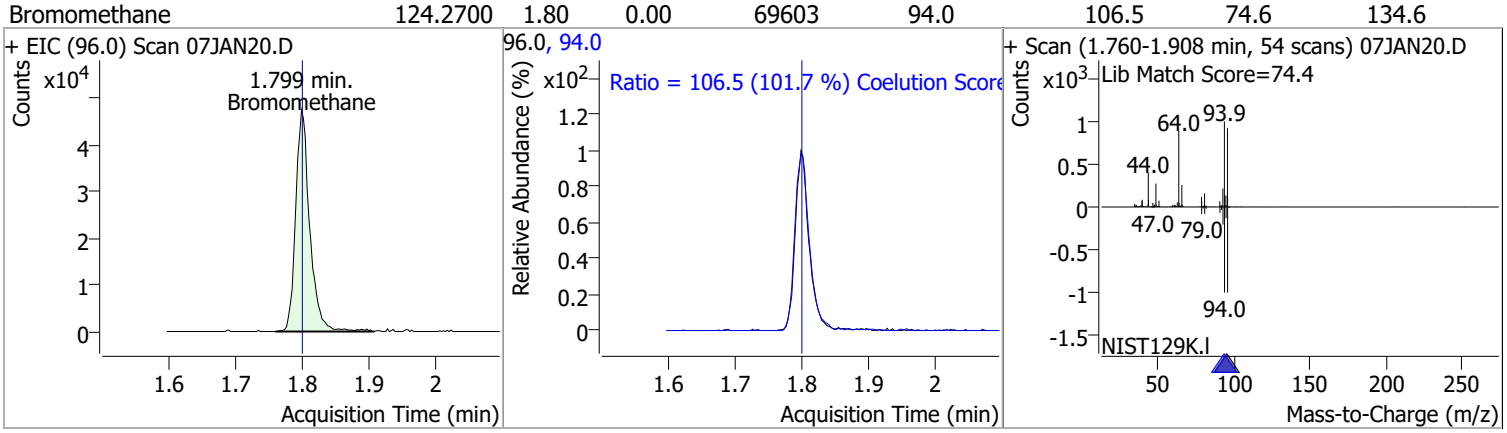
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloromethane	114.2765	1.41	0.00	159080	52.0	32.9	2.1	62.1



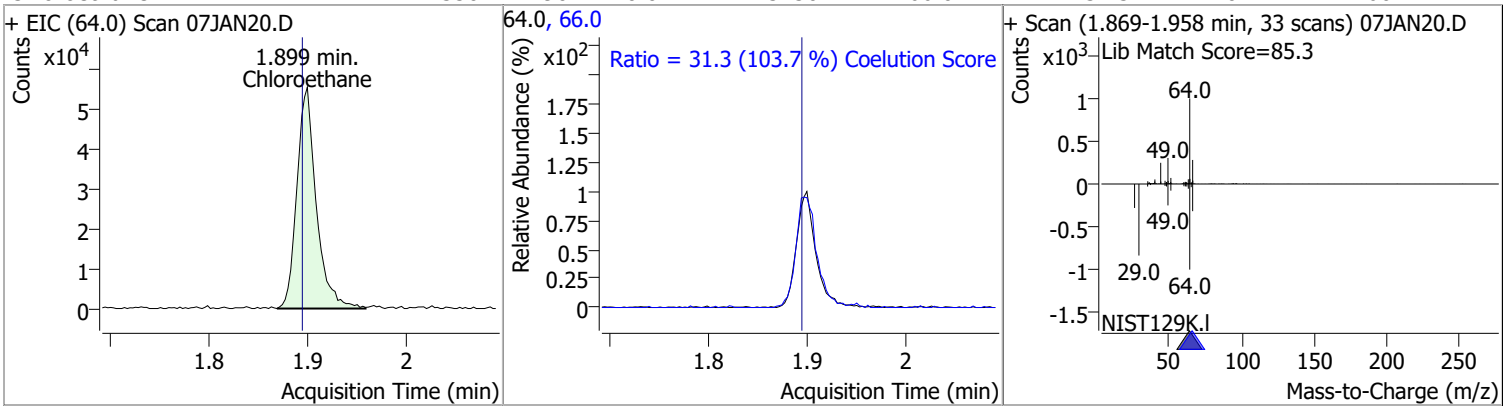
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Vinyl chloride	117.7541	1.50	0.00	147497	64.0	30.9	0.0	59.9



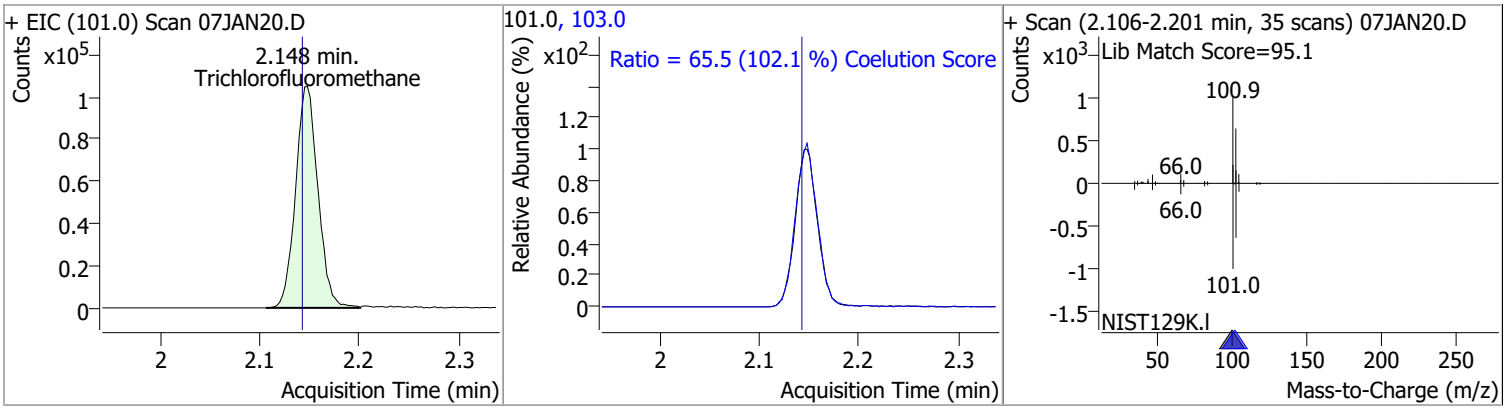
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromomethane	124.2700	1.80	0.00	69603	94.0	106.5	74.6	134.6



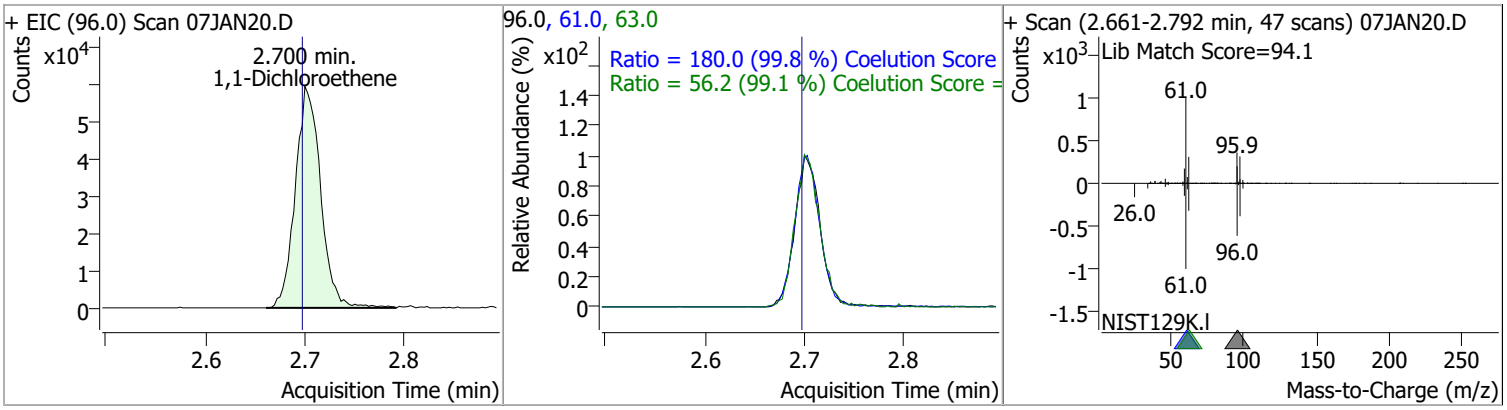
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroethane	121.3507	1.90	0.01	75250	66.0	31.3	0.1	60.1



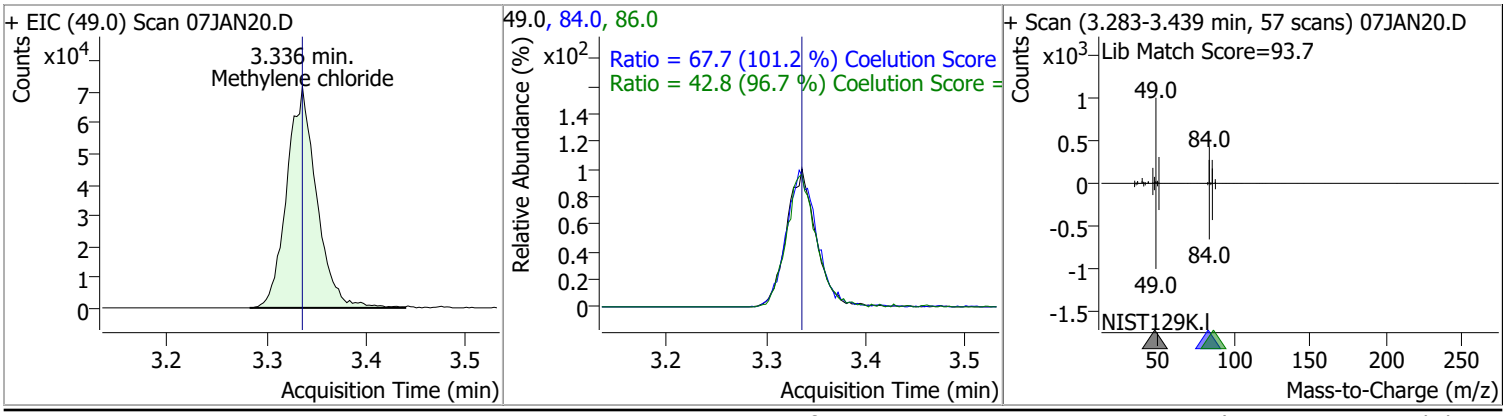
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichlorofluoromethane	105.4169	2.15	0.01	163895	103.0	65.5	34.2	94.2



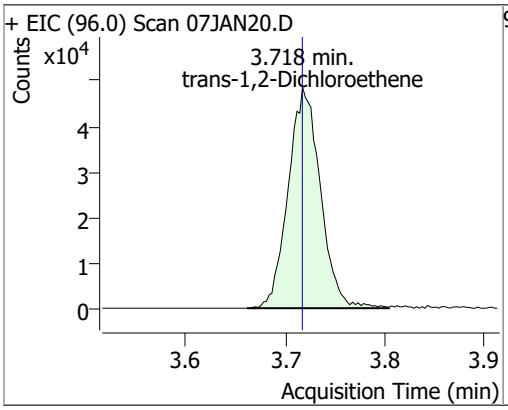
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethene	118.7865	2.70	0.00	104720	61.0	180.0	150.3	210.3
					63.0	56.2	26.7	86.7

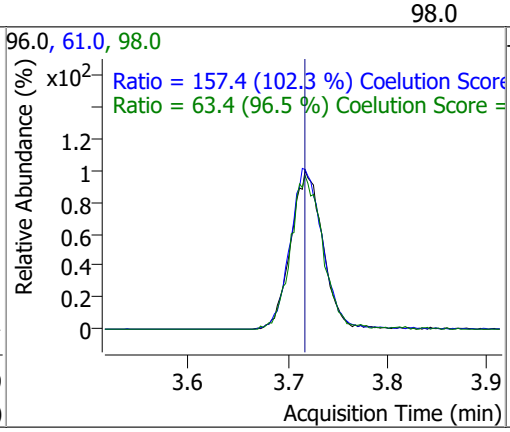


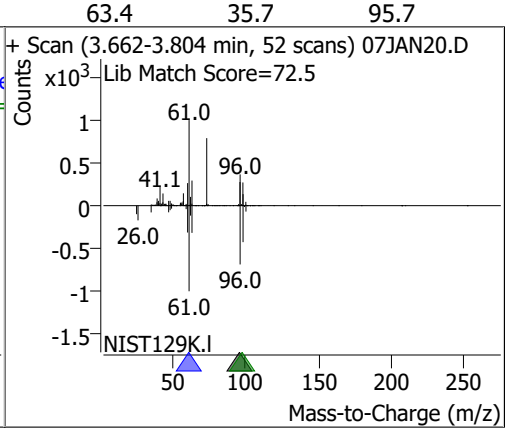
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methylene chloride	114.8197	3.34	0.00	149219	84.0	67.7	36.9	96.9
					86.0	42.8	14.3	74.3



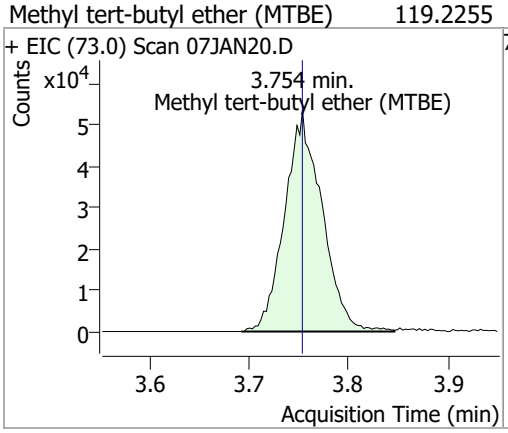
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,2-Dichloroethene	120.8683	3.72	0.00	108710	61.0	157.4	123.9	183.9
					98.0	63.4	35.7	95.7

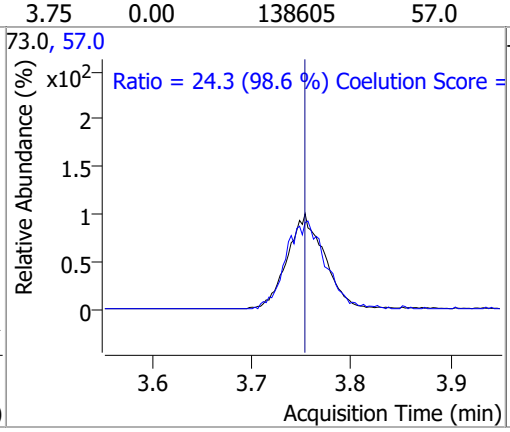


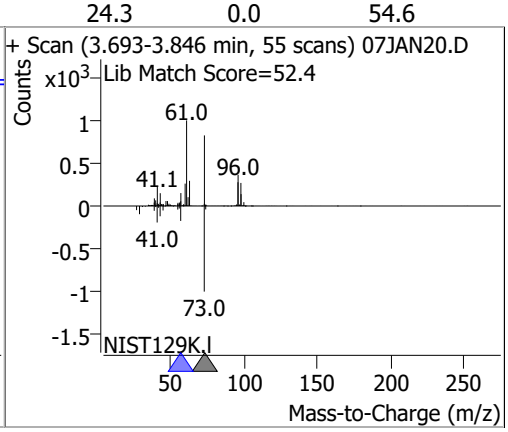




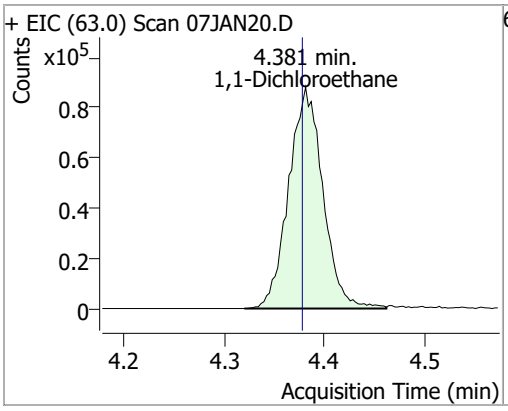
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl tert-butyl ether (MTBE)	119.2255	3.75	0.00	138605	57.0	24.3	0.0	54.6

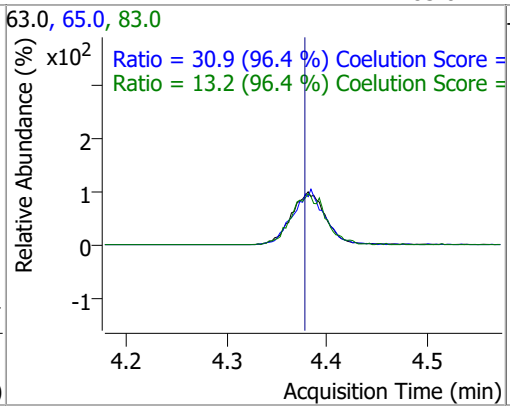


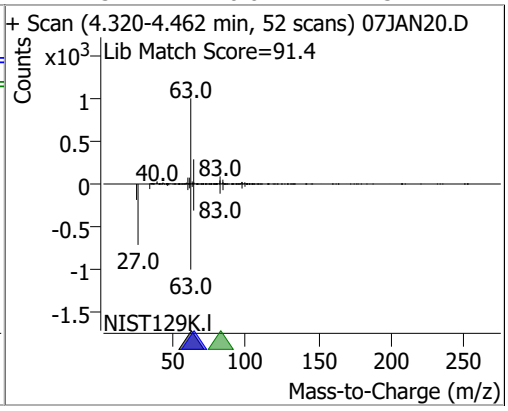




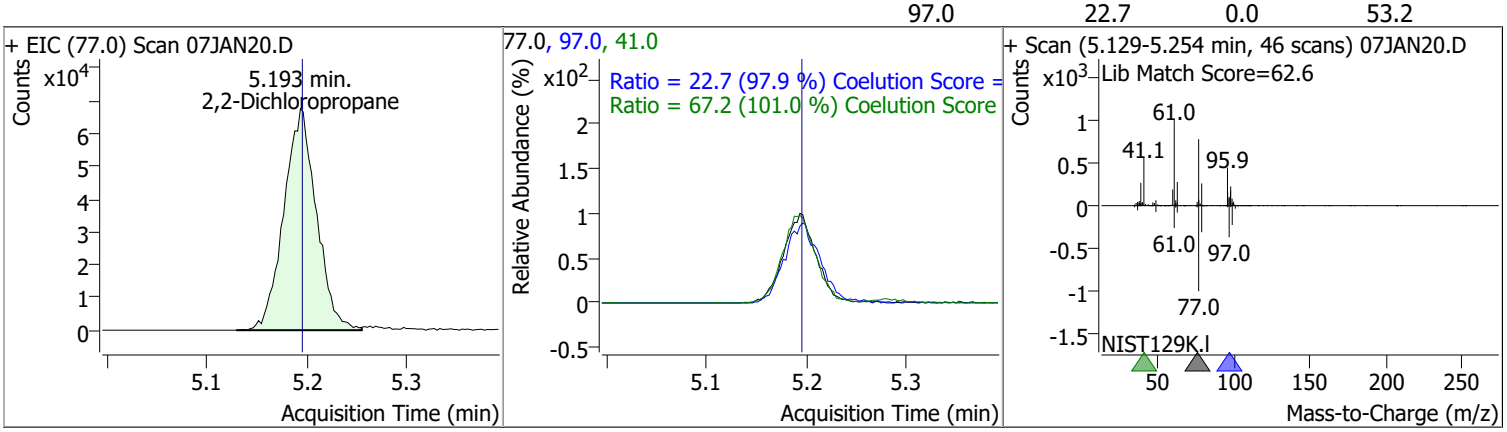
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloroethane	125.0569	4.38	0.00	209364	65.0	30.9	2.1	62.1
					83.0	13.2	0.0	43.7



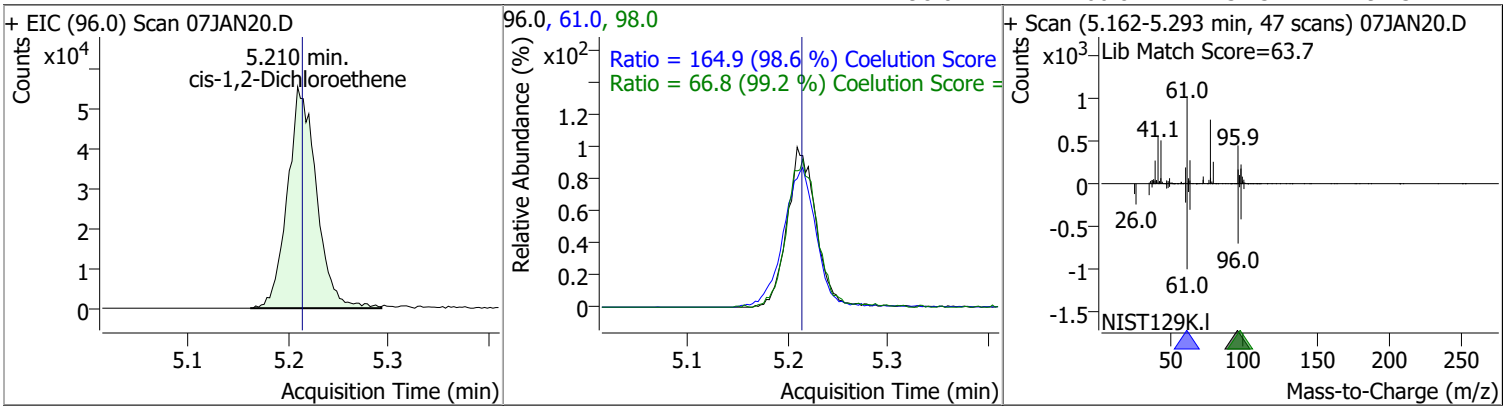




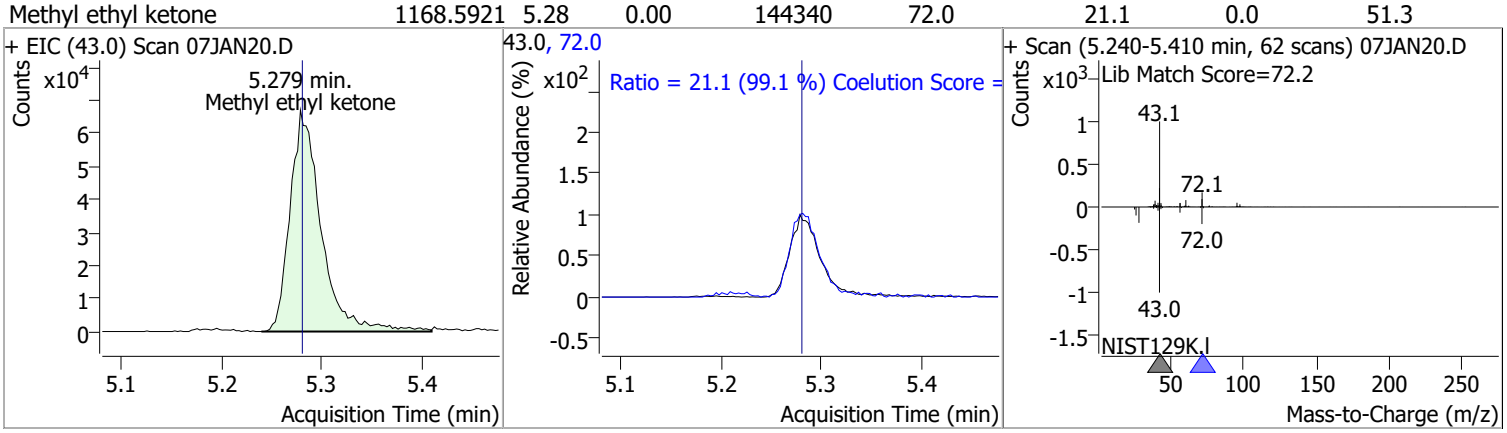
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2,2-Dichloropropane	121.7521	5.19	0.00	152733	41.0	67.2	36.5	96.5
					97.0	22.7	0.0	53.2



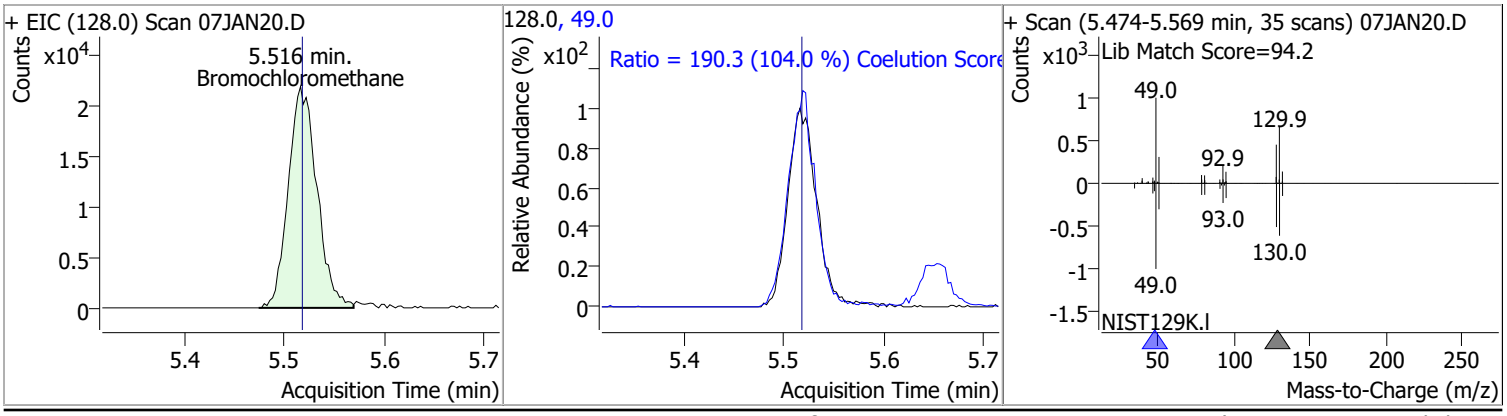
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,2-Dichloroethene	122.1596	5.21	-0.01	111394	61.0	164.9	137.2	197.2
					98.0	66.8	37.3	97.3



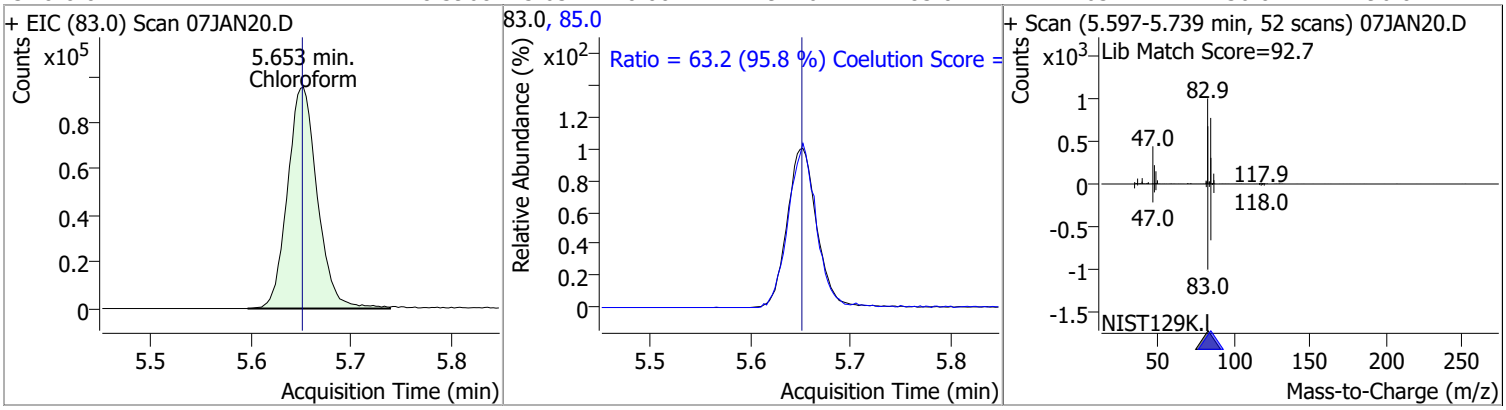
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Methyl ethyl ketone	1168.5921	5.28	0.00	144340	72.0	21.1	0.0	51.3



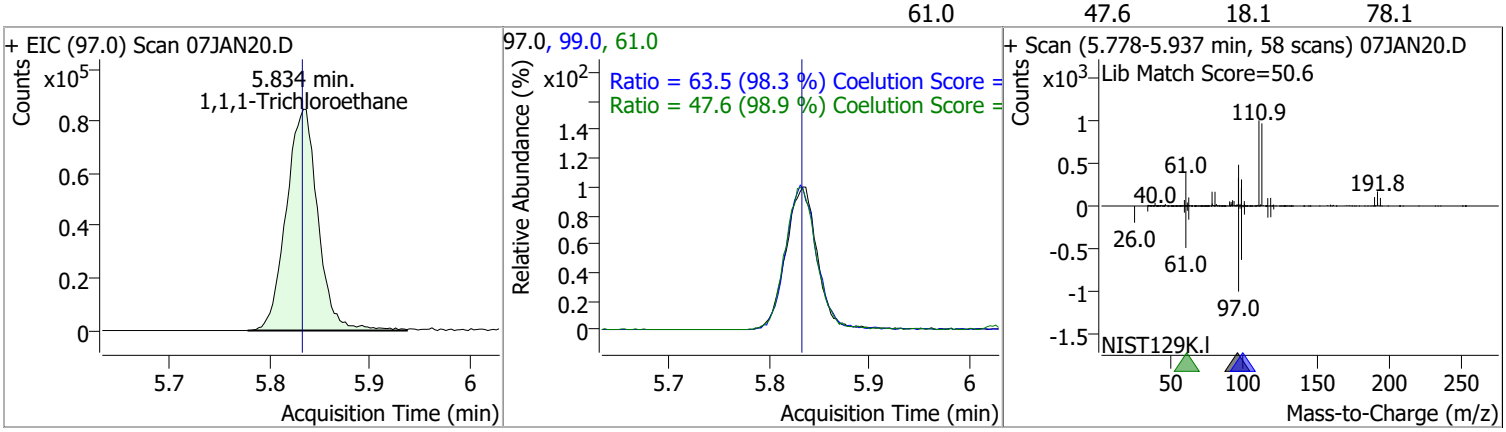
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromochloromethane	115.3287	5.52	0.00	43567	49.0	190.3	152.9	212.9



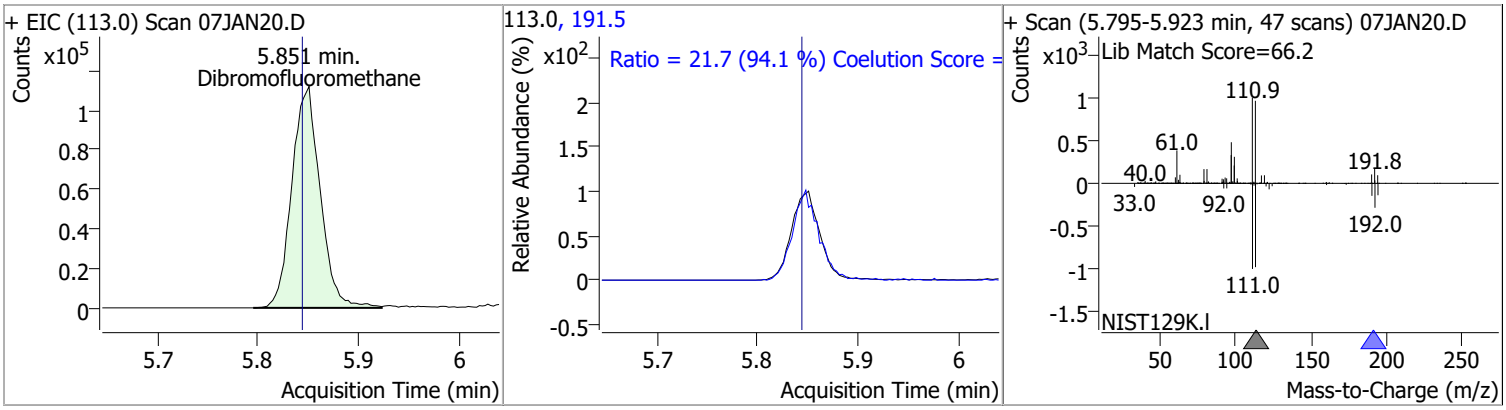
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chloroform	118.3968	5.65	0.00	197264	85.0	63.2	36.0	96.0



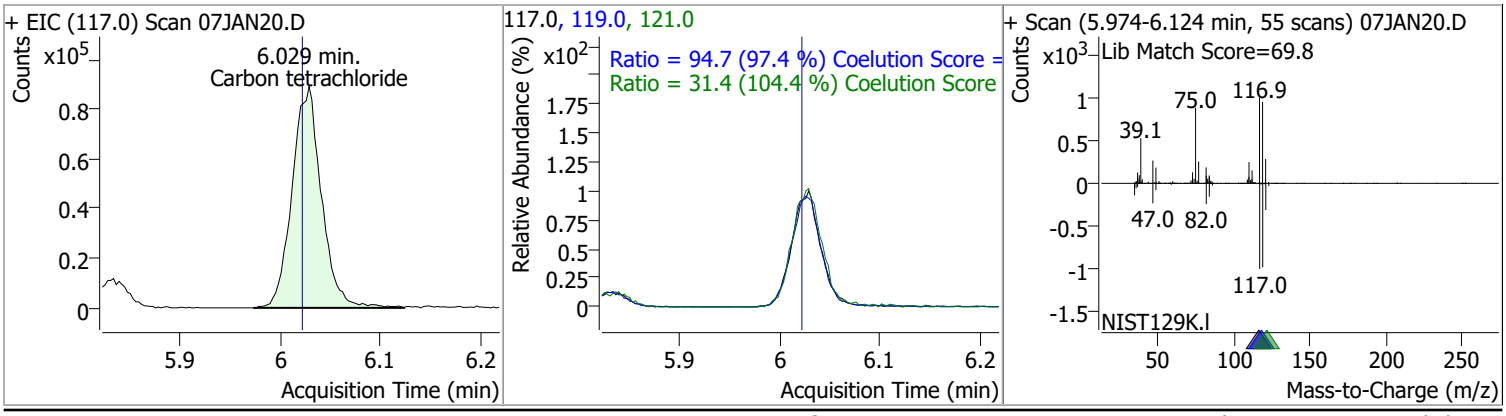
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1-Trichloroethane	118.9139	5.83	0.00	185675	99.0	63.5	34.7	94.7
					61.0	47.6	18.1	78.1



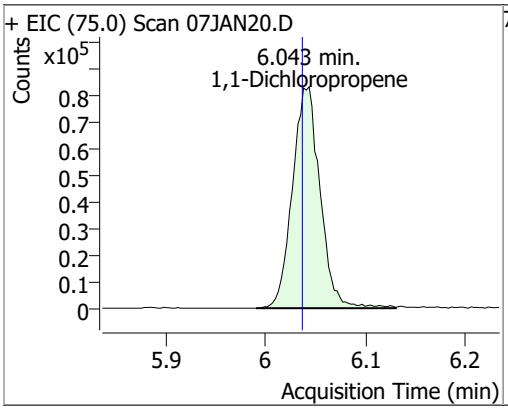
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromofluoromethane	267.4624	5.85	0.01	220533	191.5	21.7	0.0	53.1

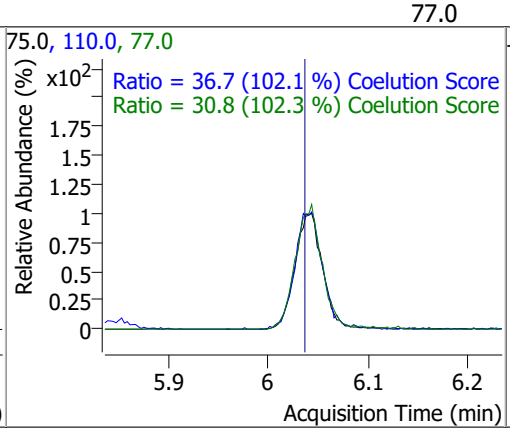


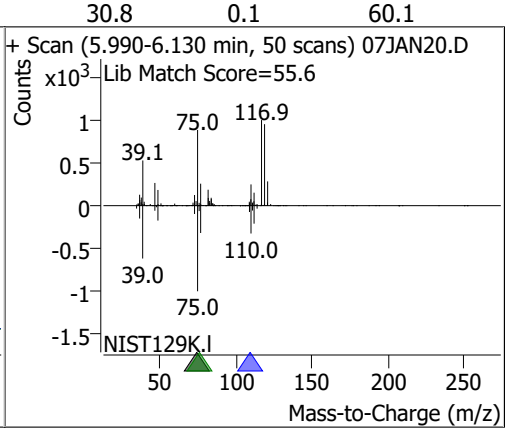
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Carbon tetrachloride	117.6182	6.03	0.01	180946	119.0	94.7	67.2	127.2
					121.0	31.4	0.1	60.1



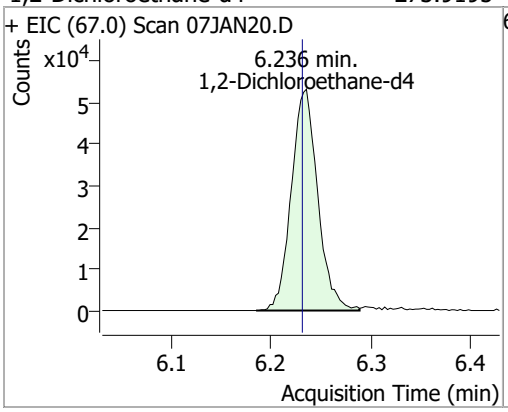
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1-Dichloropropene	119.8660	6.04	0.01	159136	110.0	36.7	5.9	65.9
					77.0	30.8	0.1	60.1

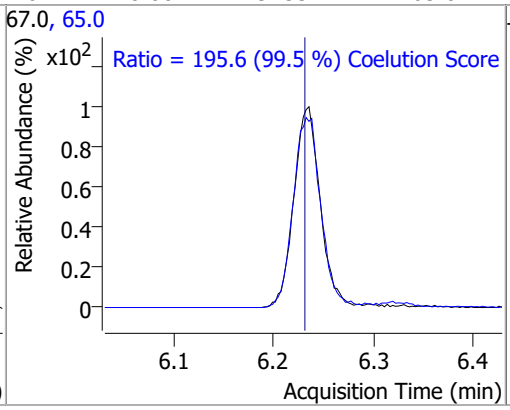


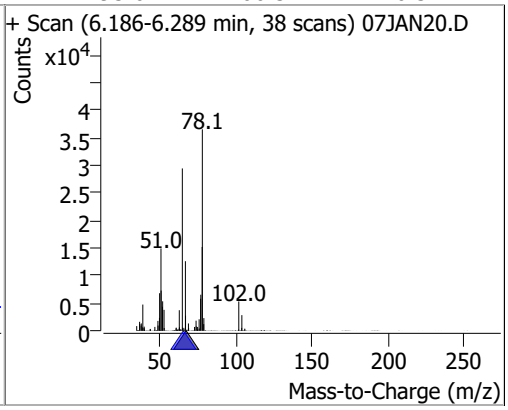




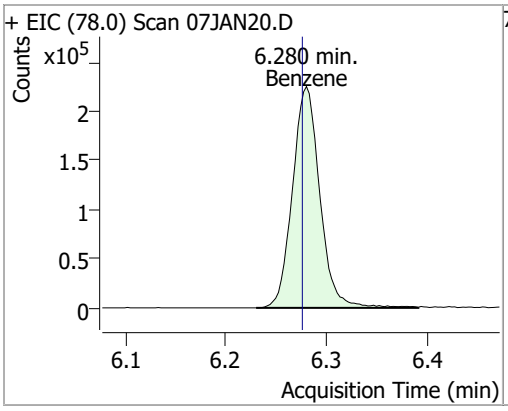
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane-d4	273.9195	6.24	0.00	97554	65.0	195.6	166.5	226.5

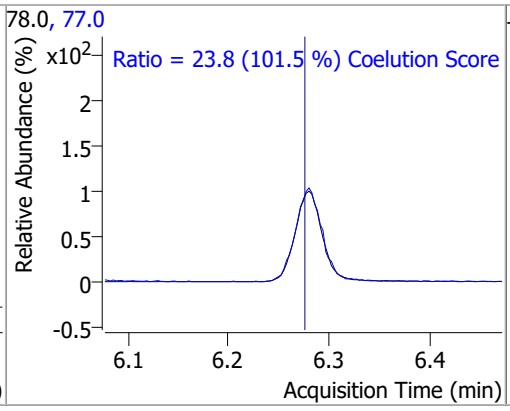


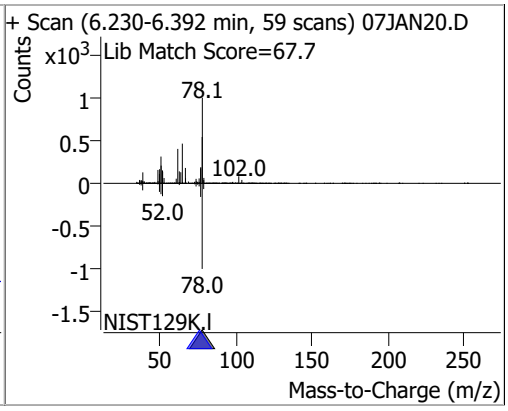




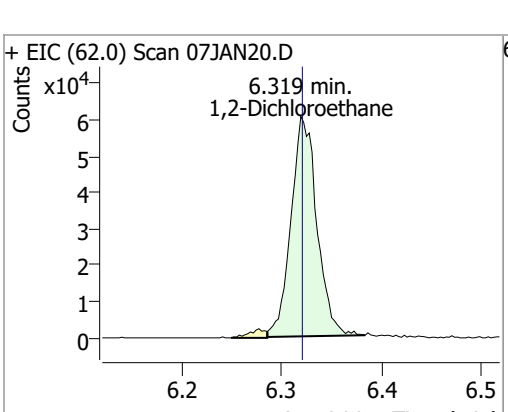
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Benzene	122.5270	6.28	0.00	426971	77.0	23.8	0.0	53.5

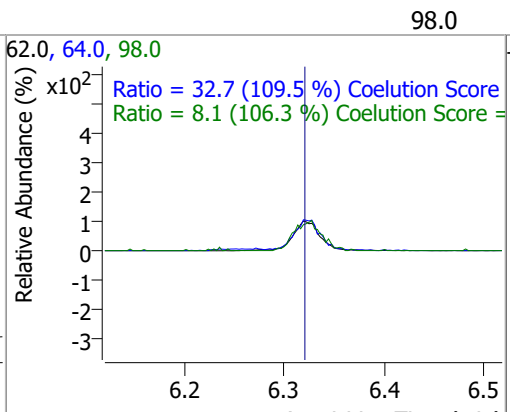


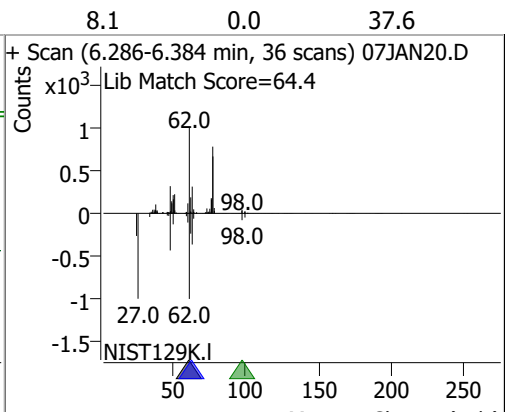




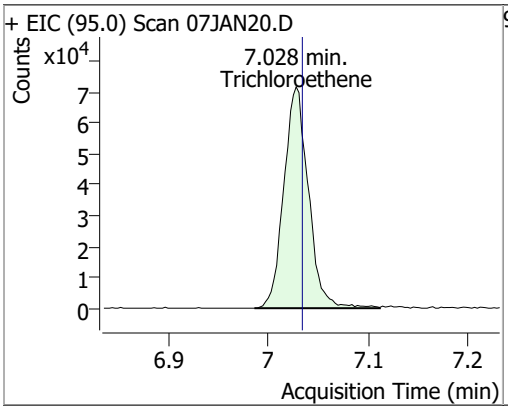
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloroethane	115.0977	6.32	0.00	108503	64.0	32.7	0.0	59.9
					98.0	8.1	0.0	37.6

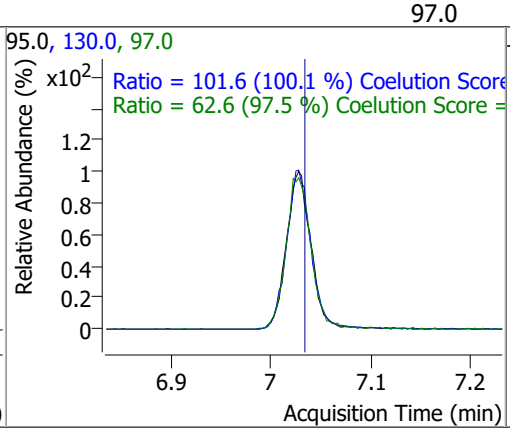


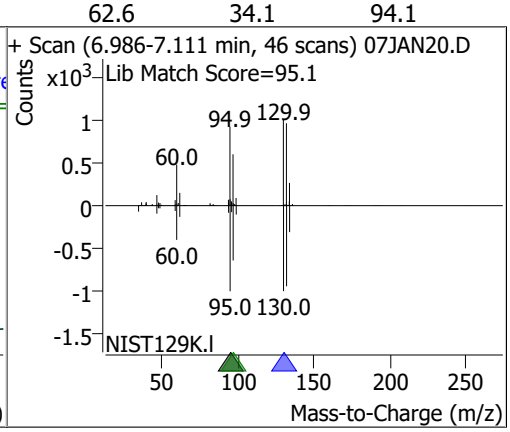




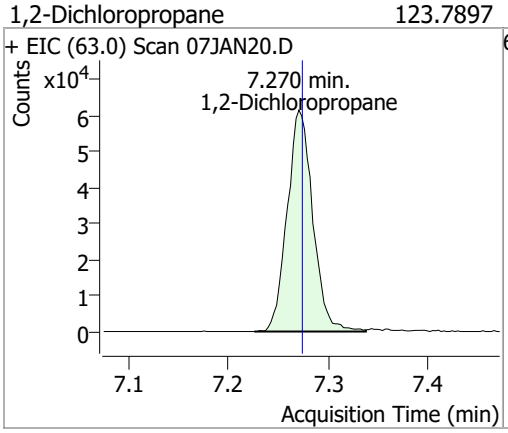
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Trichloroethene	122.2872	7.03	0.00	122705	130.0	101.6	71.5	131.5
					97.0	62.6	34.1	94.1

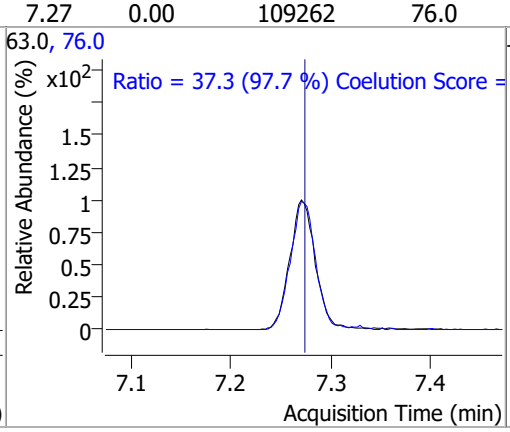


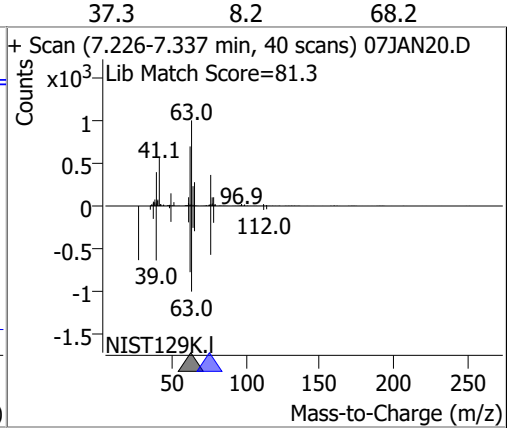




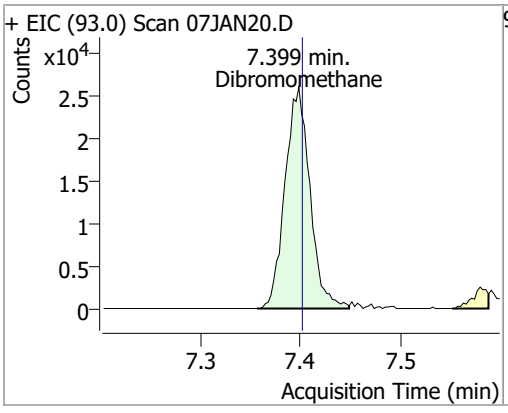
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichloropropane	123.7897	7.27	0.00	109262	76.0	37.3	8.2	68.2

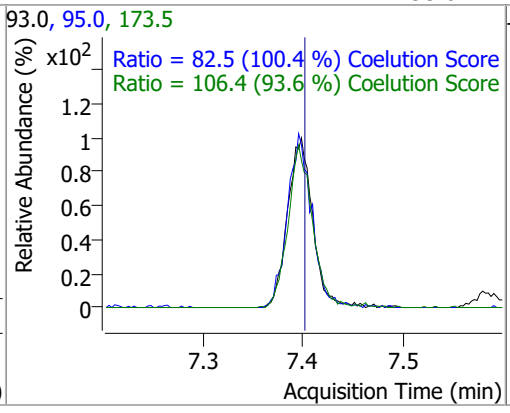


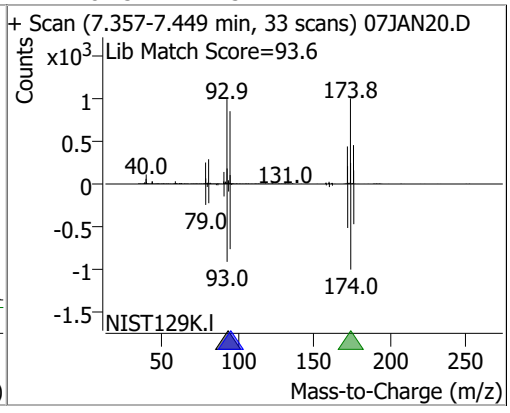




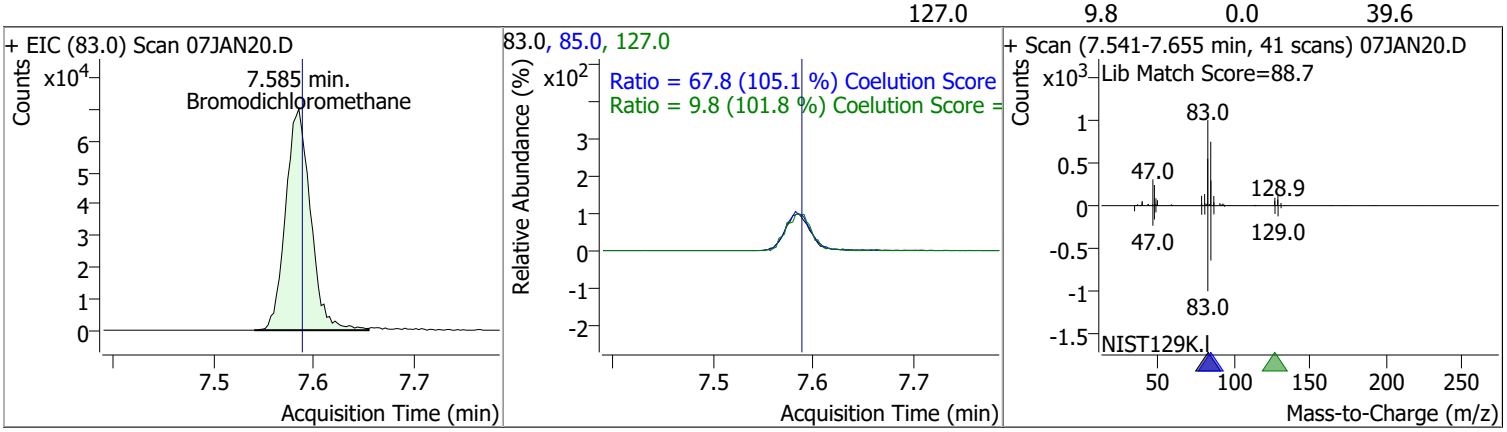
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Dibromomethane	120.8597	7.40	0.00	45080	173.5	106.4	83.7	143.7
					95.0	82.5	52.2	112.2



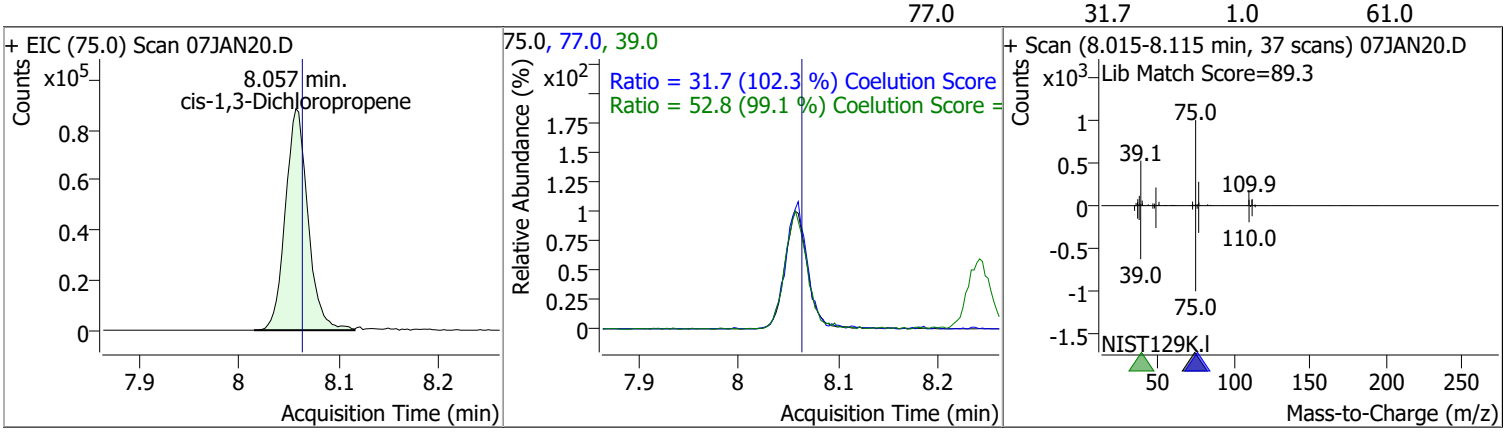




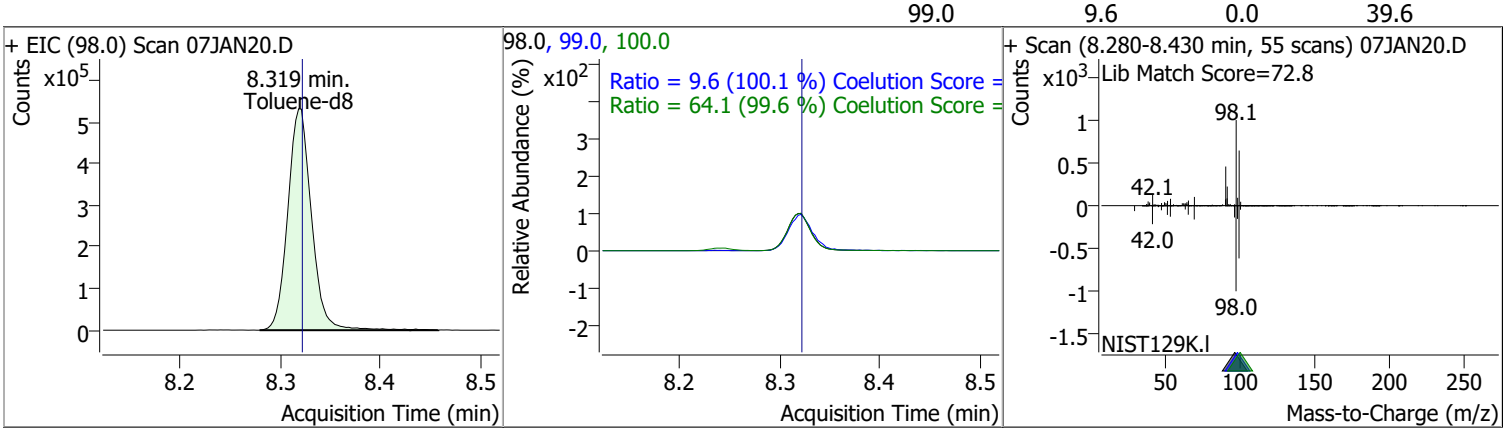
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromodichloromethane	118.2568	7.59	0.00	121732	85.0	67.8	34.5	94.5
					127.0	9.8	0.0	39.6



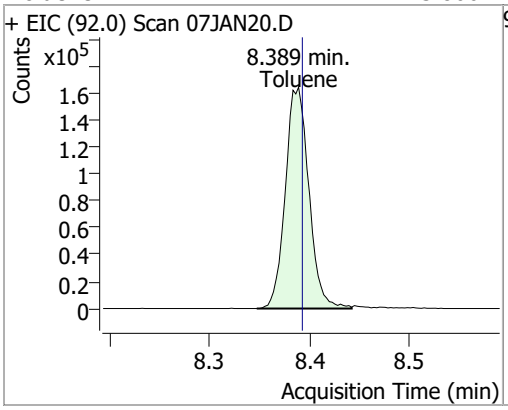
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
cis-1,3-Dichloropropene	119.1185	8.06	0.00	138637	39.0	52.8	23.3	83.3
					77.0	31.7	1.0	61.0

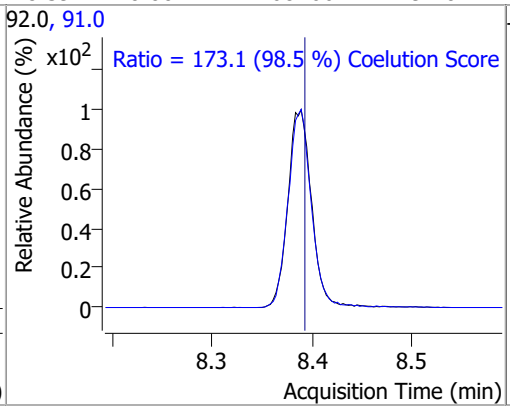


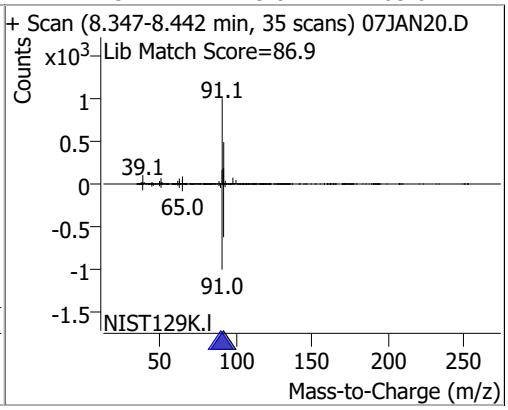
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene-d8	272.3635	8.32	0.00	873247	100.0	64.1	34.4	94.4
					99.0	9.6	0.0	39.6



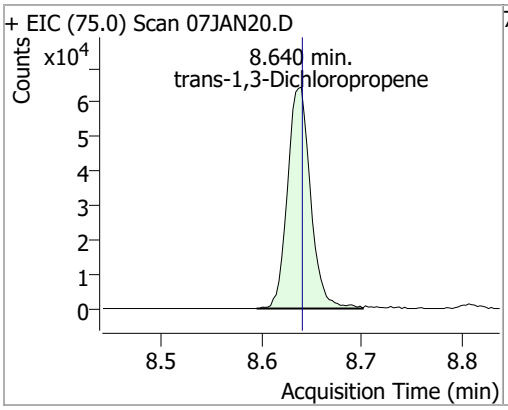
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Toluene	123.8661	8.39	0.00	268266	91.0	173.1	145.8	205.8

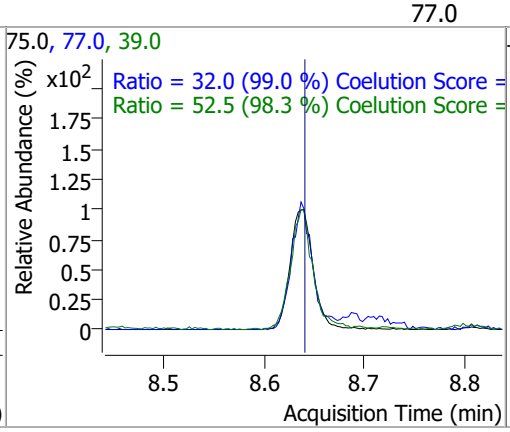


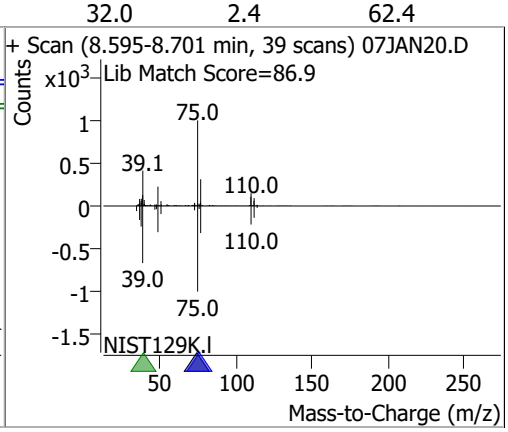




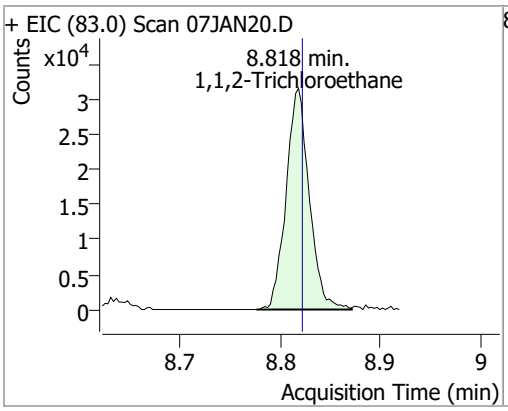
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
trans-1,3-Dichloropropene	124.4317	8.64	0.00	103086	39.0	52.5	23.4	83.4
					77.0	32.0	2.4	62.4

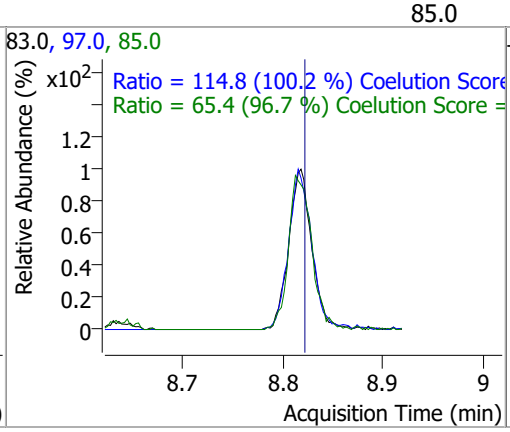


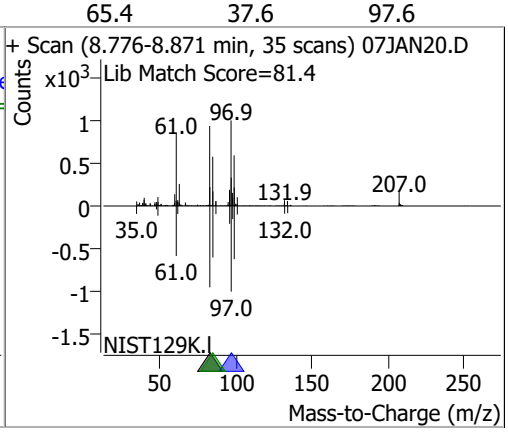




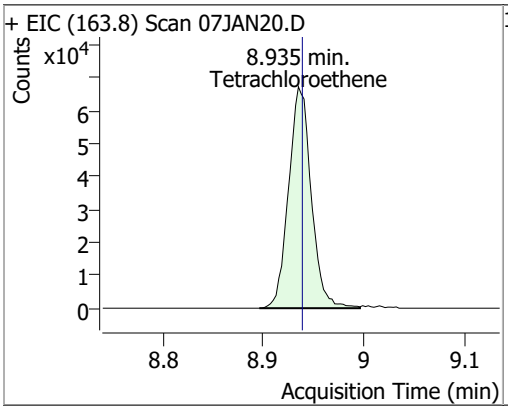
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2-Trichloroethane	116.9843	8.82	0.00	50481	97.0	114.8	84.6	144.6
					85.0	65.4	37.6	97.6

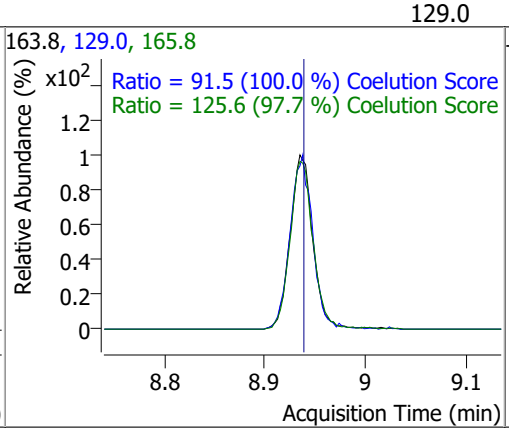


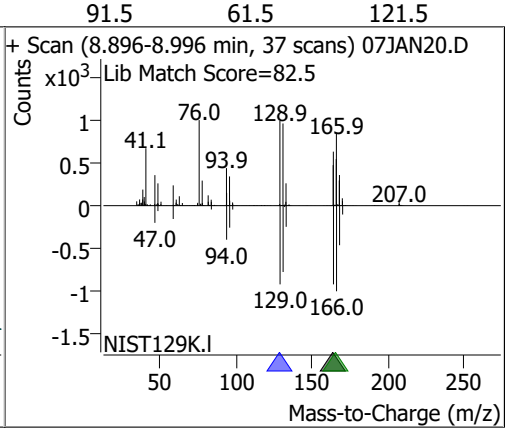




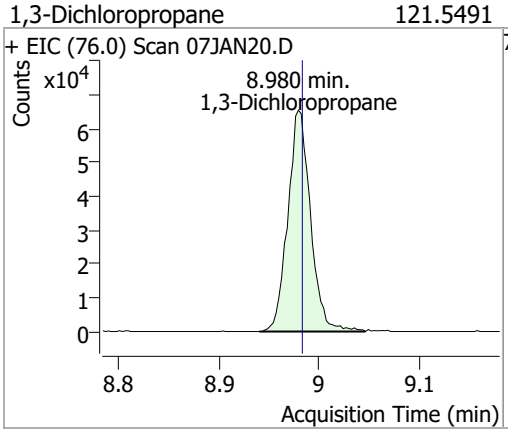
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Tetrachloroethene	119.7520	8.94	0.00	105808	165.8	125.6	98.6	158.6
					129.0	91.5	61.5	121.5

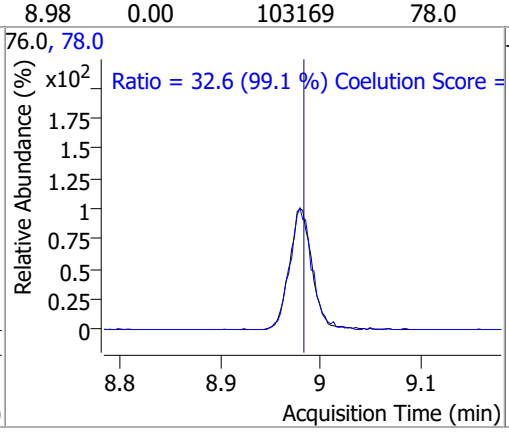


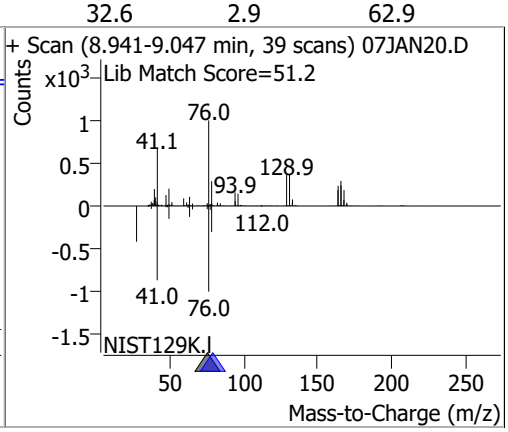




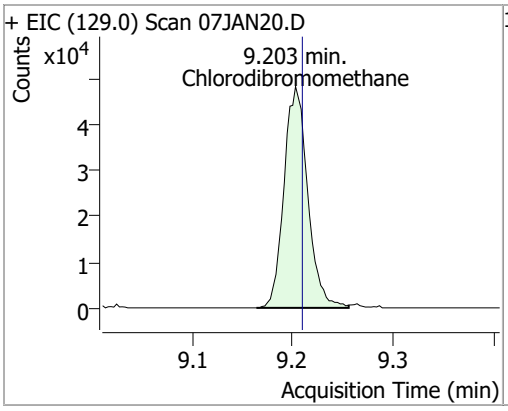
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichloropropane	121.5491	8.98	0.00	103169	78.0	32.6	2.9	62.9

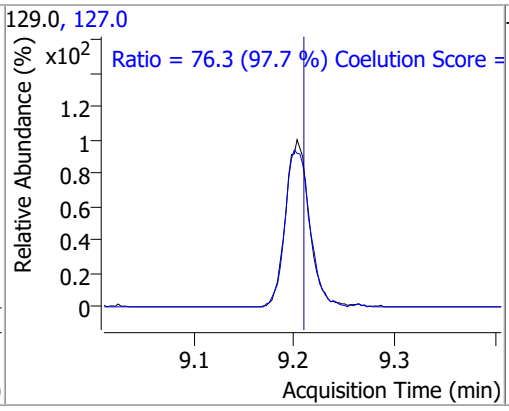


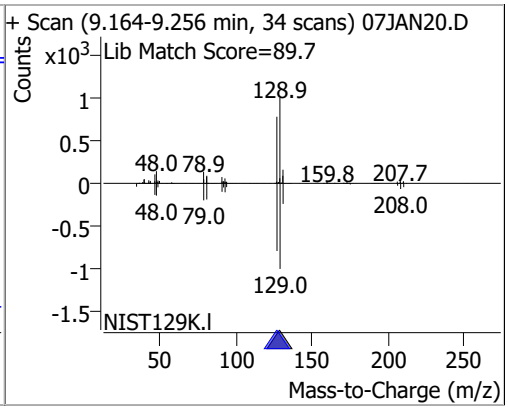




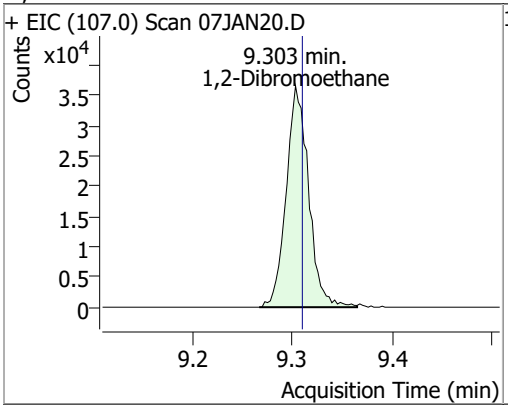
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorodibromomethane	118.1999	9.20	0.00	79716	127.0	76.3	48.0	108.0

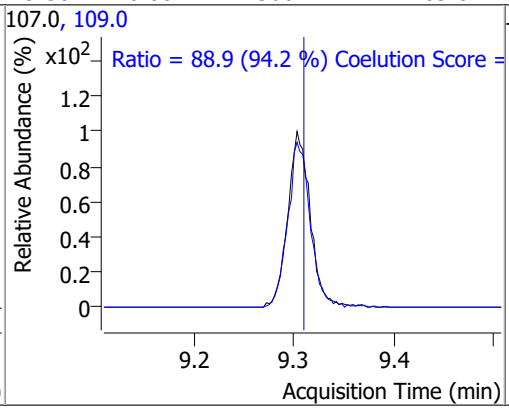


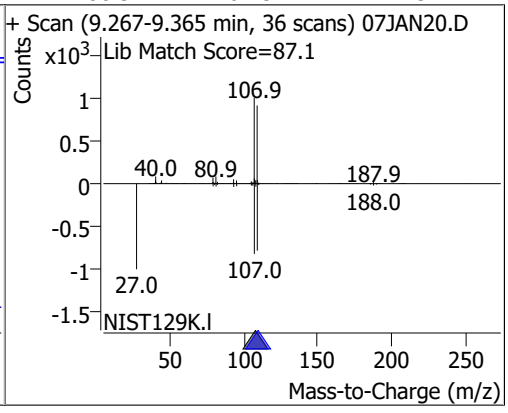




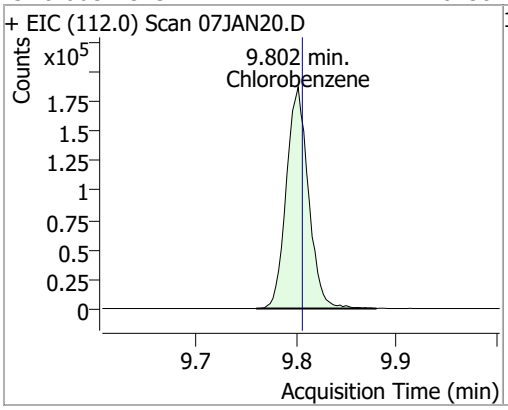
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	120.0089	9.30	0.00	56624	109.0	88.9	64.5	124.5

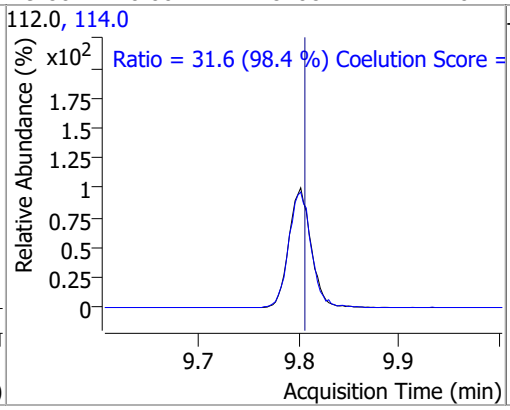


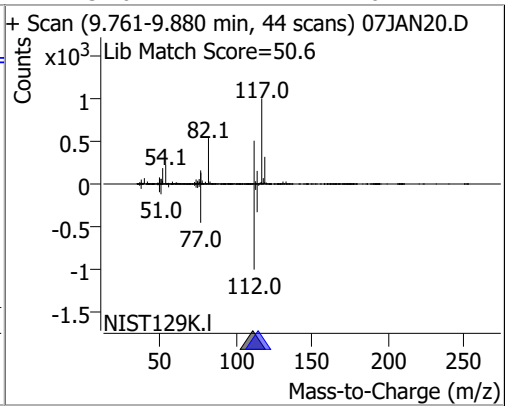




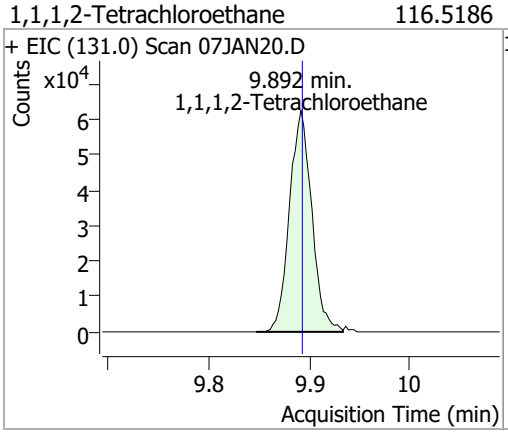
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Chlorobenzene	121.0758	9.80	0.00	287084	114.0	31.6	2.1	62.1

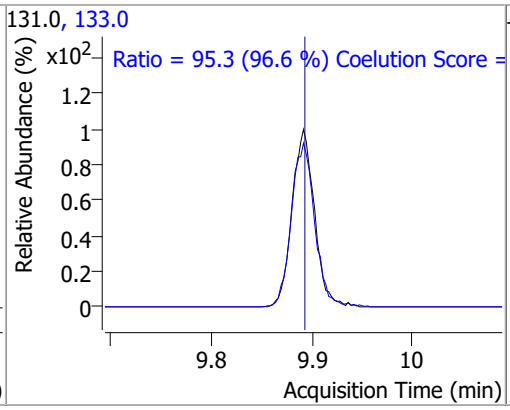


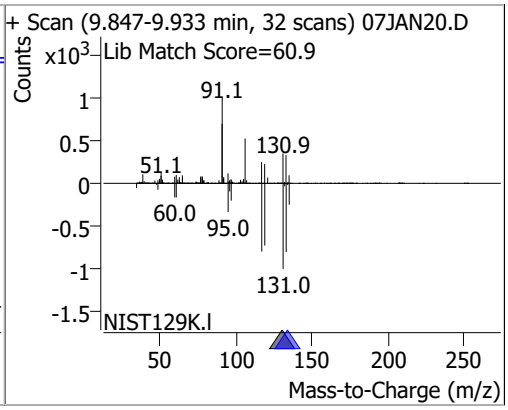




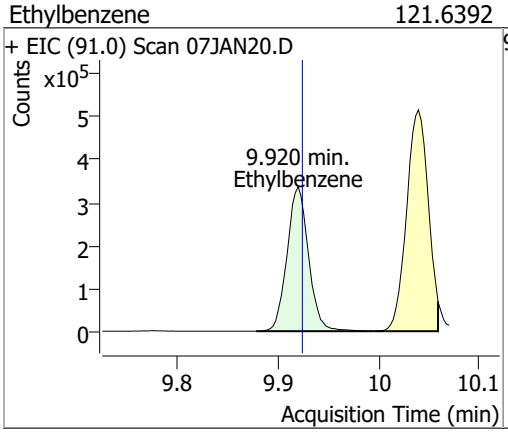
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	116.5186	9.89	0.00	96577	133.0	95.3	68.6	128.6

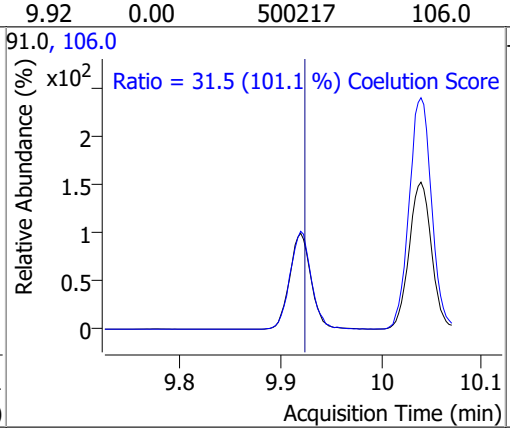


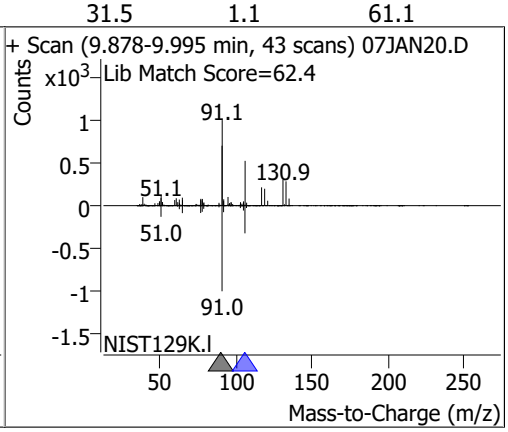




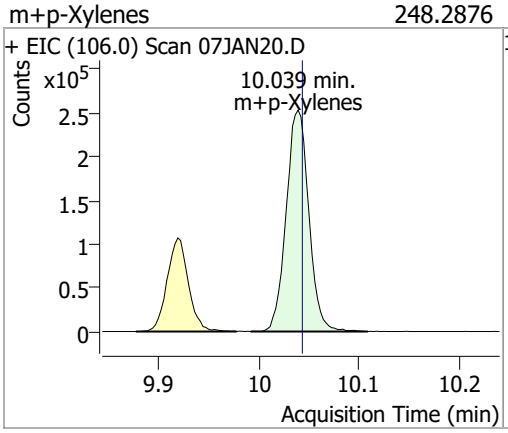
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Ethylbenzene	121.6392	9.92	0.00	500217	106.0	31.5	1.1	61.1

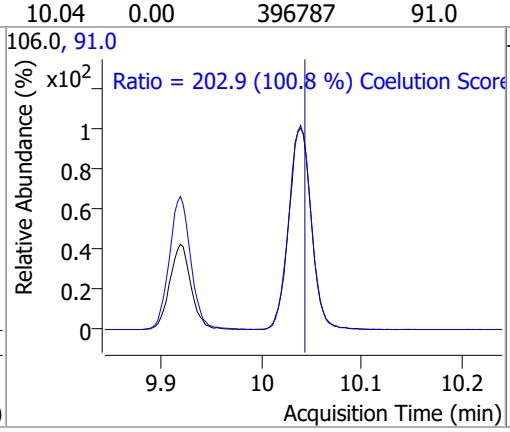


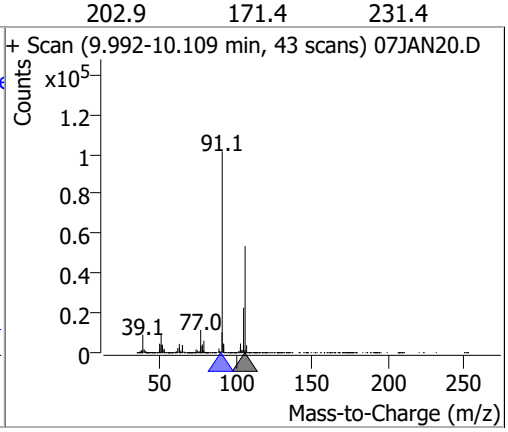




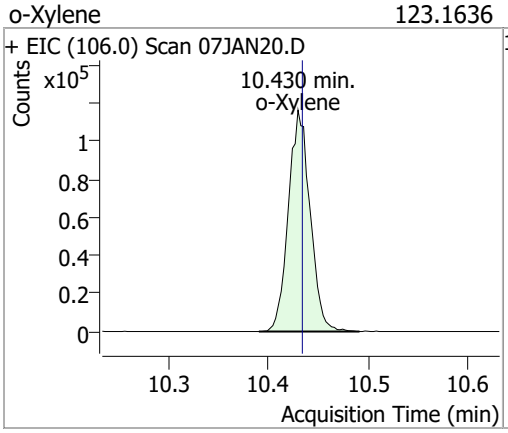
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
m+p-Xylenes	248.2876	10.04	0.00	396787	91.0	202.9	171.4	231.4

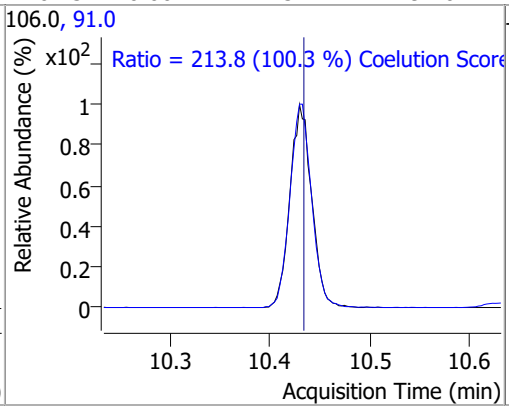


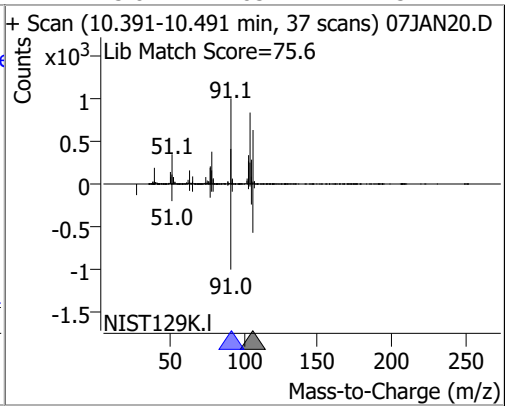




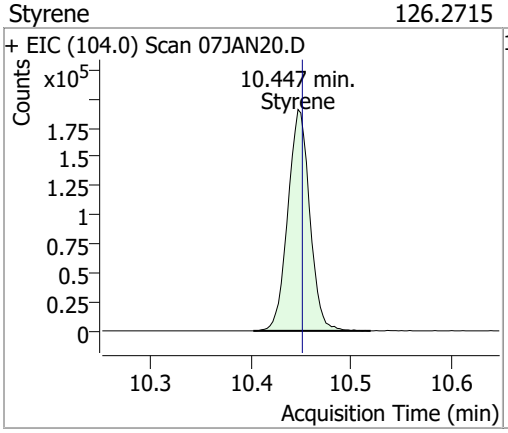
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
o-Xylene	123.1636	10.43	0.00	175221	91.0	213.8	183.1	243.1

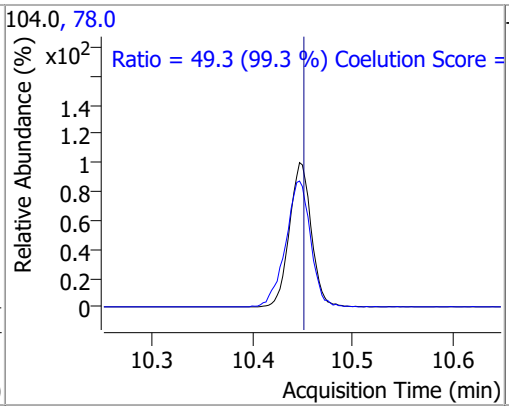


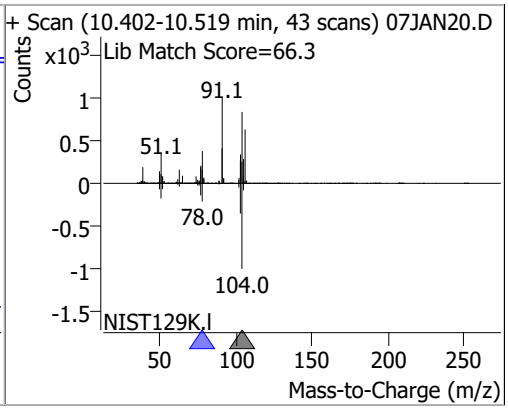




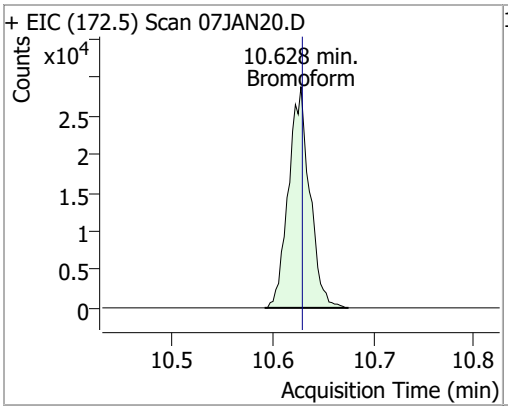
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Styrene	126.2715	10.45	0.00	289229	78.0	49.3	19.6	79.6

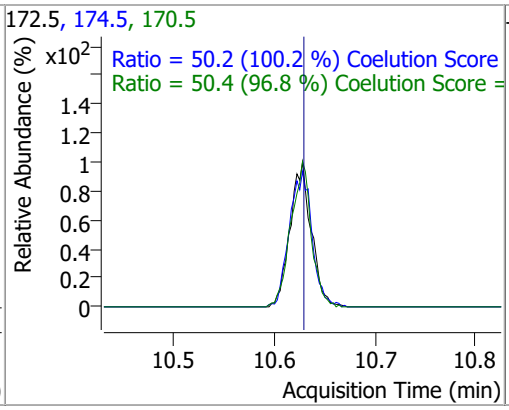


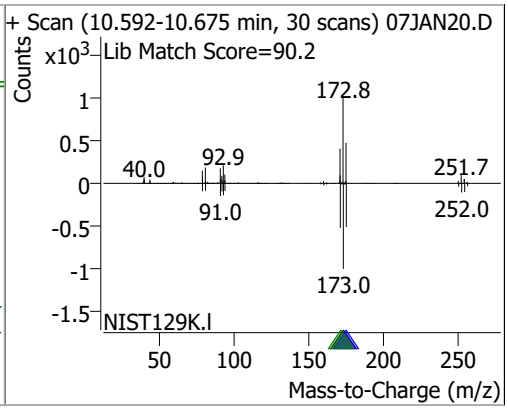




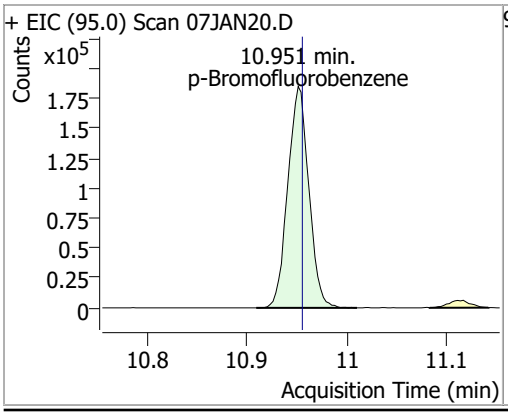
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromoform	120.8329	10.63	0.00	42320	170.5	50.4	22.1	82.1
					174.5	50.2	20.1	80.1

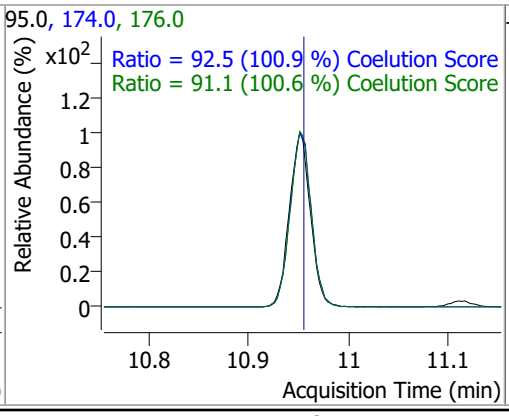


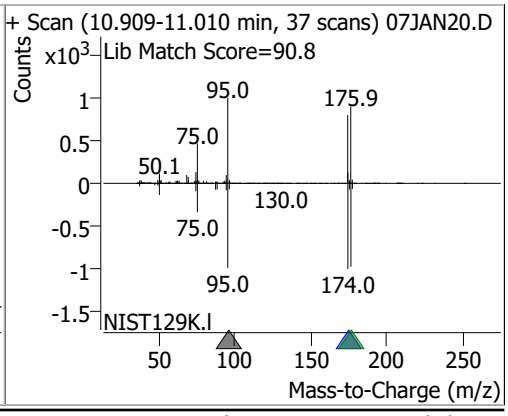




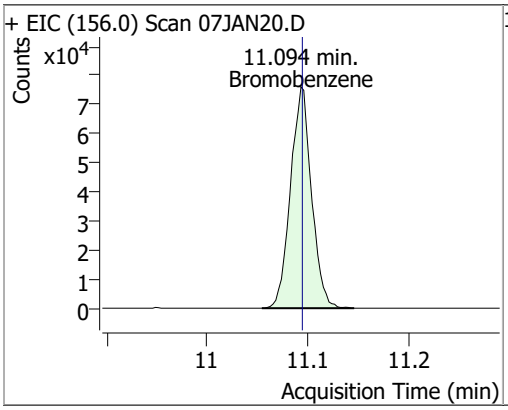
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
p-Bromofluorobenzene	268.4143	10.95	0.00	269134	174.0	92.5	61.7	121.7
					176.0	91.1	60.6	120.6

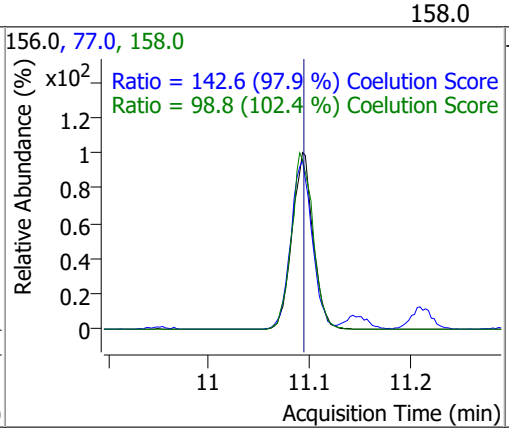


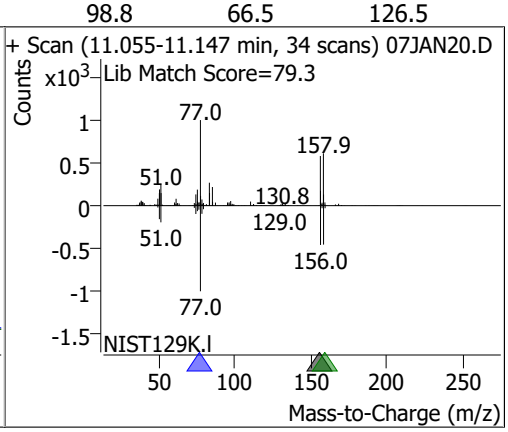




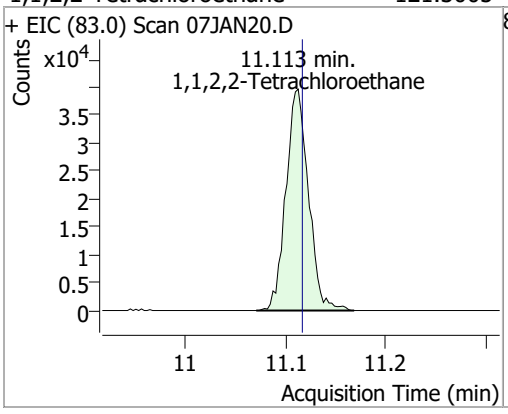
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
Bromobenzene	124.7379	11.09	0.00	110486	77.0	142.6	115.7	175.7
					158.0	98.8	66.5	126.5

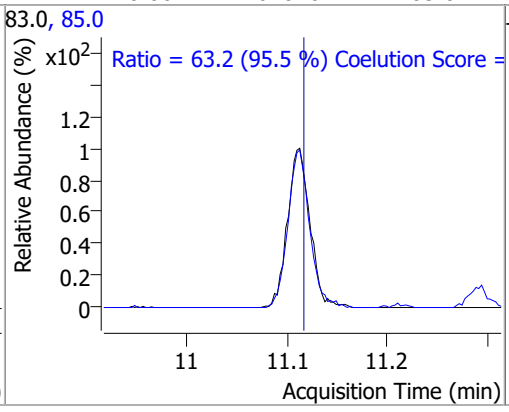


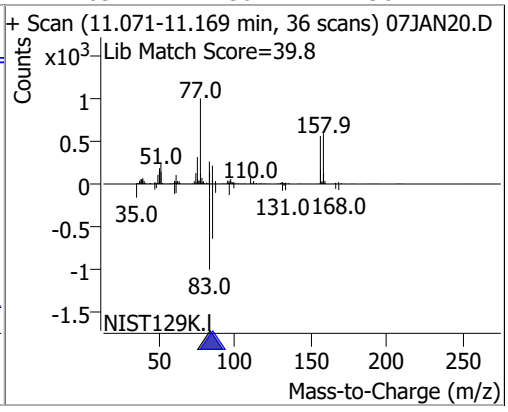




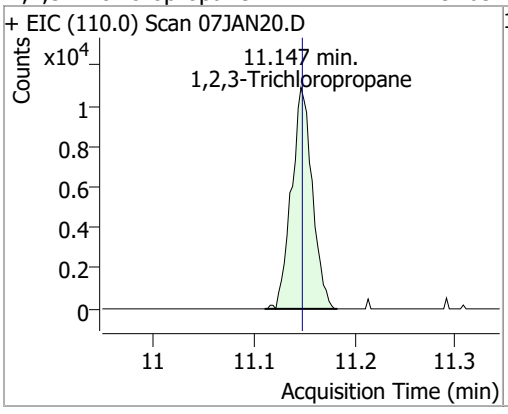
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,2,2-Tetrachloroethane	121.3005	11.11	0.00	61840	85.0	63.2	36.2	96.2

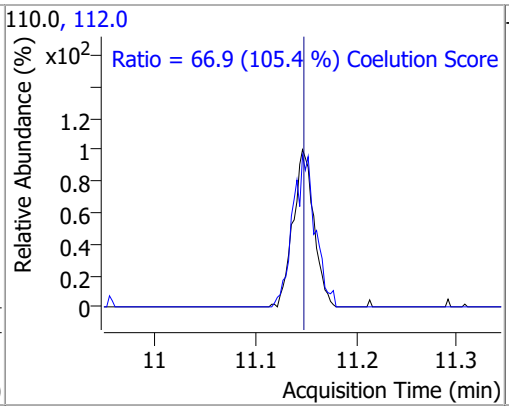


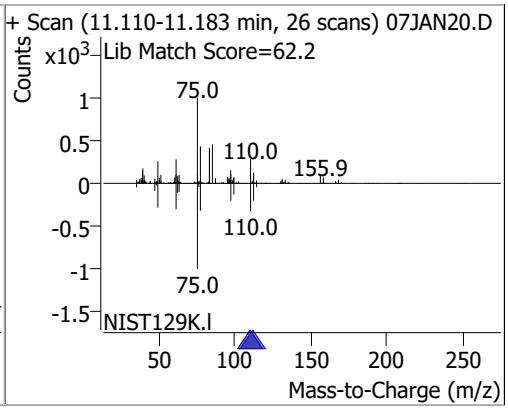




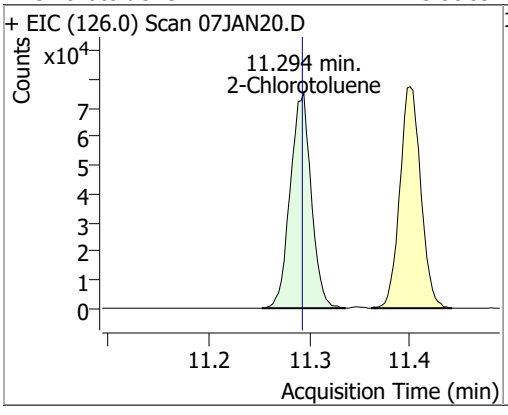
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2,3-Trichloropropane	114.9765	11.15	0.00	15684	112.0	66.9	33.5	93.5

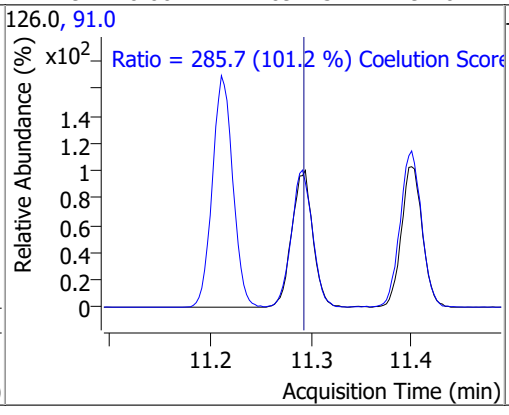


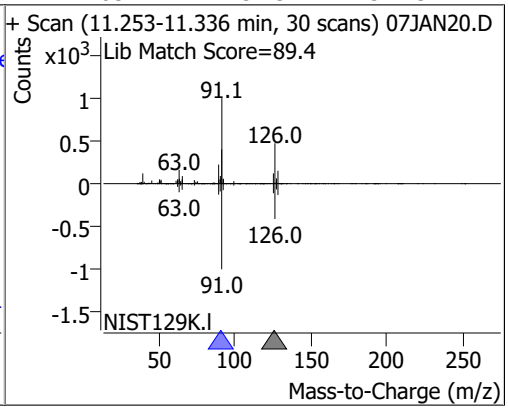




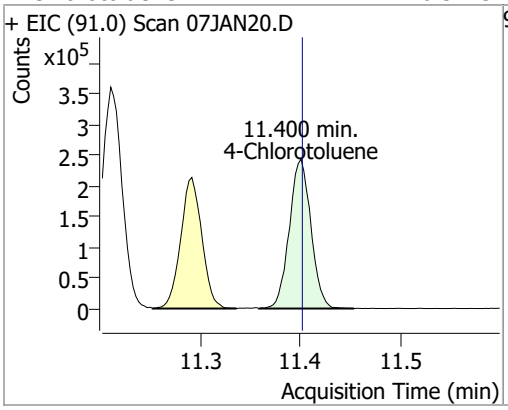
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
2-Chlorotoluene	123.9909	11.29	0.00	109275	91.0	285.7	252.3	312.3

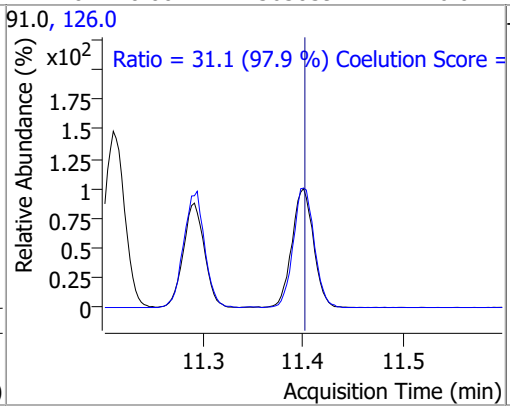


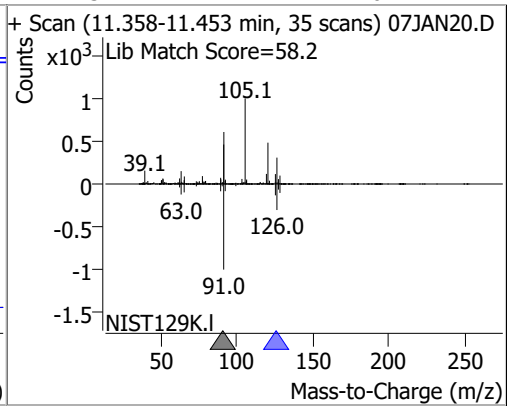




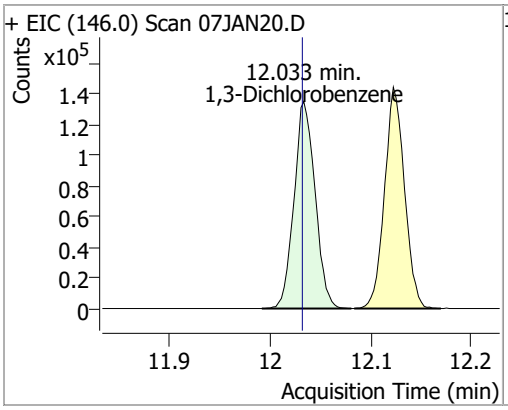
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
4-Chlorotoluene	126.3413	11.40	0.00	363039	126.0	31.1	1.7	61.7

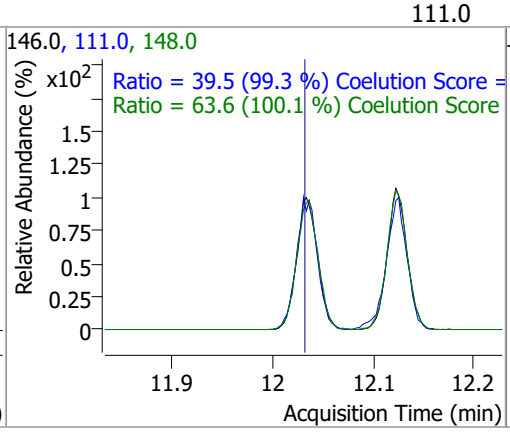


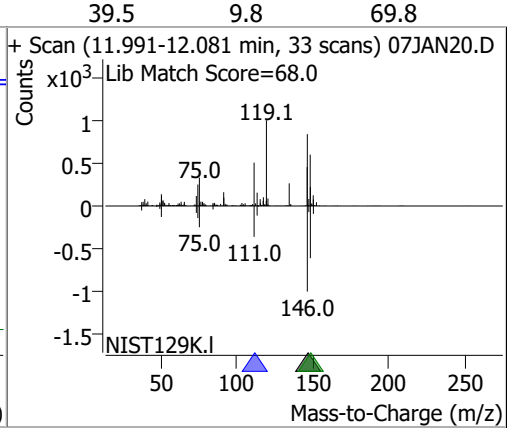




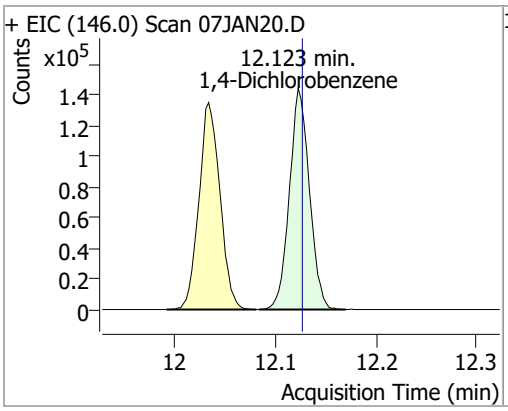
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,3-Dichlorobenzene	121.3987	12.03	0.00	196110	148.0	63.6	33.6	93.6
					111.0	39.5	9.8	69.8

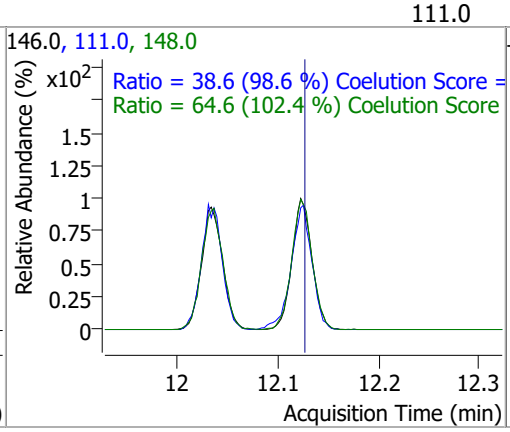


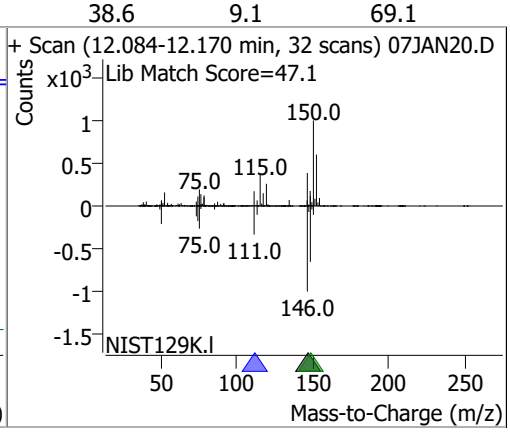




Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,4-Dichlorobenzene	120.7629	12.12	0.00	198916	148.0	64.6	33.1	93.1
					111.0	38.6	9.1	69.1

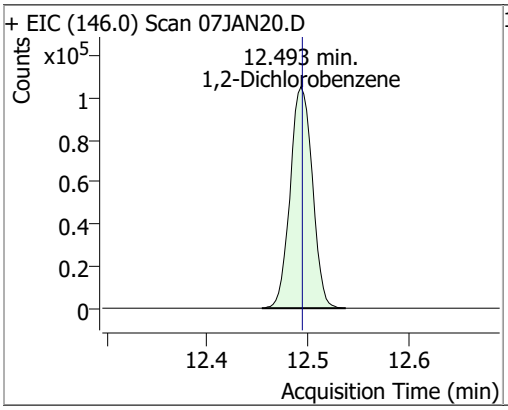


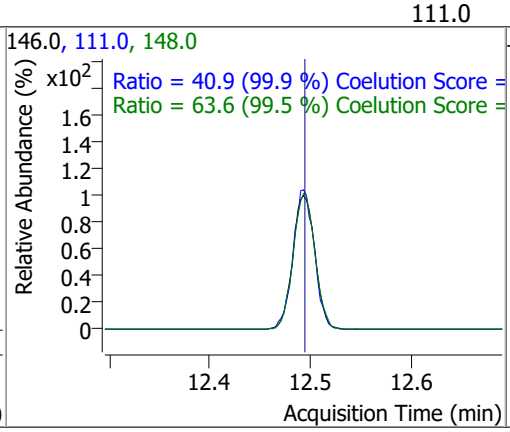


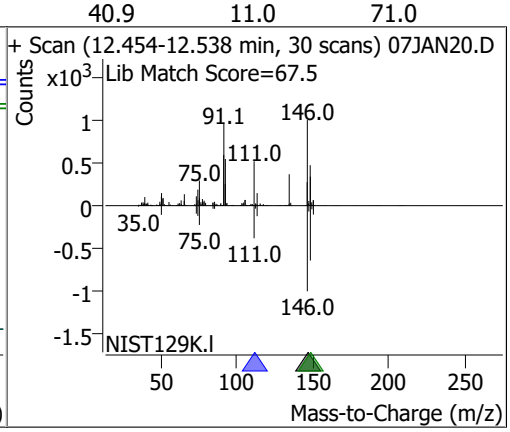


Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dichlorobenzene	118.7328	12.49	0.00	162097	148.0	63.6	33.9	93.9
					111.0	40.9	11.0	71.0







Batch name and path: D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin
Quant batch version: 10.0
Quant reporting version: 10.0

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdNewBatchTable	BL2000\mchavez	1/7/2022 9:13:24 AM	Create new batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 9:13:32 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN01.D			✓	
CmdStartMethodEditing	BL2000\mchavez	1/7/2022 9:14:04 AM	Start method editing			✓	
CmdImportMethodFromBatch	BL2000\mchavez	1/7/2022 9:14:05 AM	Import method from batch D:\Org\Data\VOA5975C\VG010622\VG010622_8260B.batch.bin			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/7/2022 9:14:10 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/7/2022 9:14:10 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/7/2022 9:14:11 AM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 9:14:14 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 9:24:01 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 9:25:03 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN02.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 9:25:07 AM	Set SampleType = TuneCheck for sample 07JAN02.D; previous value = Sample			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 9:26:45 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 10:31:59 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 10:32:17 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN03.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 10:32:20 AM	Set SampleType = CC for sample 07JAN03.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 10:32:26 AM	Set LevelName = CC for sample 07JAN03.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 10:32:29 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 10:33:53 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 11:19:56 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 11:20:13 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN04.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 11:20:18 AM	Set SampleType = QC for sample 07JAN04.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 11:20:22 AM	Set LevelName = QC for sample 07JAN04.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 11:20:25 AM	Set SampleInformation = LCSA for sample 07JAN04.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 11:20:29 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 11:20:55 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 11:55:20 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 11:56:57 AM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN05.D			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 11:57:04 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 11:57:25 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 12:03:31 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 12:03:46 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN06.D			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 12:03:51 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/7/2022 12:04:34 PM	Set SampleType = Blank for sample 07JAN06.D; previous value = Sample			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 12:04:38 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 12:11:24 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 1:16:43 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN08.D, D:\Org\Data\VOA5975C\VG010722\07JAN07.D			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 1:16:50 PM	Quantitate all compounds in all samples			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 1:33:40 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 2:44:00 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 2:44:40 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN11.D, D:\Org\Data\VOA5975C\VG010722\07JAN10.D, D:\Org\Data\VOA5975C\VG010722\07JAN09.D			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 2:44:51 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 2:51:46 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 3:59:33 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 4:00:03 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN14.D, D:\Org\Data\VOA5975C\VG010722\07JAN13.D, D:\Org\Data\VOA5975C\VG010722\07JAN12.D			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 4:00:13 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 4:15:06 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/7/2022 4:21:12 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\mchavez	1/7/2022 4:21:51 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN15.D			✓	
CmdQuantitate	BL2000\mchavez	1/7/2022 4:22:04 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/7/2022 4:34:37 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/9/2022 12:17:18 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\mchavez	1/9/2022 12:18:19 PM	Add samples from worklist: D:\Org\Data\VOA5975C\VG010722\07JAN21.D, D:\Org\Data\VOA5975C\VG010722\07JAN20.D, D:\Org\Data\VOA5975C\VG010722\07JAN19.D, D:\Org\Data\VOA5975C\VG010722\07JAN18.D, D:\Org\Data\VOA5975C\VG010722\07JAN17.D, D:\Org\Data\VOA5975C\VG010722\07JAN16.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:19:03 PM	Set SampleType = CC for sample 07JAN20.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:19:10 PM	Set LevelName = CC for sample 07JAN20.D; previous value =			✓	
CmdQuantitate	BL2000\mchavez	1/9/2022 12:19:23 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:19:53 PM	Set SampleType = Matrix for sample 07JAN17.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:19:58 PM	Set SampleType = MatrixDup for sample 07JAN18.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:20:04 PM	Set SampleInformation = MatrixA for sample 07JAN17.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:20:08 PM	Set SampleInformation = MatrixA for sample 07JAN18.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:20:12 PM	Set MatrixSpikeGroup = 219 for sample 07JAN17.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:20:14 PM	Set MatrixSpikeGroup = 219 for sample 07JAN18.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/9/2022 12:24:41 PM	Set MatrixSpikeGroup = 219 for sample 07JAN11.D; previous value =			✓	
CmdSaveBatchTable	BL2000\mchavez	1/9/2022 12:26:45 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/9/2022 9:08:38 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdStartMethodEditing	BL2000\mchavez	1/9/2022 9:08:59 PM	Start method editing			✓	
CmdImportMethodFromFile	BL2000\mchavez	1/9/2022 9:09:00 PM	Import method from file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/9/2022 9:09:10 PM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/9/2022 9:09:10 PM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/9/2022 9:09:11 PM	End method editing			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdQuantitate	BL2000\mchavez	1/9/2022 9:09:23 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/9/2022 9:14:28 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/11/2022 10:34:48 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdStartMethodEditing	BL2000\mchavez	1/11/2022 10:35:03 AM	Start method editing			✓	
CmdImportMethodFromFile	BL2000\mchavez	1/11/2022 10:35:04 AM	Import method from file \\MASSHUNTER\Org\Data\Methods\Quant\VOA5975C\VOA5975C_010422_CAL\VOA5975C_8260B_SHT_DoD_L4_010422.m			✓	
CmdApplyMethodToAllSamples	BL2000\mchavez	1/11/2022 10:35:12 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\mchavez	1/11/2022 10:35:12 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\mchavez	1/11/2022 10:35:12 AM	End method editing			✓	
CmdQuantitate	BL2000\mchavez	1/11/2022 10:35:24 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	1/11/2022 10:35:53 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/11/2022 3:58:37 PM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/11/2022 4:00:44 PM	Set SampleApproved = True for sample 07JAN03.D; previous value = False			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:01:01 PM	Manually integrate qualifier 57.0 of compound Methyl tert-butyl ether (MTBE) in sample 07JAN04.D, from x, y = 3.681, 0 to 3.809, 338, result = 38793; previous integration is from x, y = 3.745, 144 to 3.809, 338 and previous response = 25564.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\mchavez	1/11/2022 4:01:06 PM	Drop baseline for qualifier 57.0 of compound Methyl tert-butyl ether (MTBE) in sample 07JAN04.D to y = 0, new integration is from x, y = 3.681, 0 to 3.809, 0 and new response = 40094; previous integration is from x, y = 3.681, 0 to 3.809, 338 and previous response = 38793.			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/11/2022 4:01:44 PM	Set SampleApproved = True for sample 07JAN04.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/11/2022 4:03:47 PM	Set SampleApproved = True for sample 07JAN06.D; previous value = False			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:04:04 PM	Manually integrate compound Methylene chloride in sample 07JAN06.D from x, y = 3.310, 0 to 3.377, 0; result = 1664			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/11/2022 4:04:06 PM	Set UserAnnotation = NI for compound Methylene chloride in sample 07JAN06.D; previous value =			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:04:08 PM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN06.D from x, y = 3.299, 0 to 3.394, 0; result = 1299			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:04:10 PM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN06.D from x, y = 3.285, 0 to 3.400, 0; result = 522			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:05:13 PM	Manually integrate compound Chloromethane in sample 07JAN07.D from x, y = 1.364, 0 to 1.450, 0; result = 1730			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:05:20 PM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN07.D from x, y = 1.364, 0 to 1.450, 0; result = 574			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/11/2022 4:06:09 PM	Set UserAnnotation = NI for compound Chloromethane in sample 07JAN07.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:06:18 PM	Manually integrate compound Methylene chloride in sample 07JAN07.D from x, y = 3.282, 0 to 3.394, 0; result = 2222			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:06:20 PM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN07.D from x, y = 3.282, 0 to 3.419, 0; result = 1662			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:06:23 PM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN07.D from x, y = 3.294, 0 to 3.388, 0; result = 1096			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/11/2022 4:06:25 PM	Set UserAnnotation = NI for compound Methylene chloride in sample 07JAN07.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:07:09 PM	Manually integrate compound Chloroform in sample 07JAN07.D from x, y = 5.617, 0 to 5.697, 0; result = 209			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:07:11 PM	Manually integrate qualifier85.0 of compound Chloroform in sample 07JAN07.D from x, y = 5.619, 0 to 5.675, 0; result = 105			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/11/2022 4:07:15 PM	Set UserAnnotation = NI for compound Chloroform in sample 07JAN07.D; previous value =			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:07:54 PM	Manually integrate compound Benzene in sample 07JAN07.D from x, y = 6.263, 0 to 6.305, 0; result = 144			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:07:56 PM	Manually integrate qualifier77.0 of compound Benzene in sample 07JAN07.D from x, y = 6.252, 0 to 6.328, 0; result = 26			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/11/2022 4:08:01 PM	Set UserAnnotation = NI for compound Benzene in sample 07JAN07.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:08:54 PM	Manually integrate compound Ethylbenzene in sample 07JAN07.D from x, y = 9.891, 0 to 9.953, 0; result = 389			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:08:56 PM	Manually integrate qualifier106.0 of compound Ethylbenzene in sample 07JAN07.D from x, y = 9.917, 0 to 9.958, 0; result = 37			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/11/2022 4:08:58 PM	Set UserAnnotation = NI for compound Ethylbenzene in sample 07JAN07.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:09:04 PM	Manually integrate compound m+p-Xylenes in sample 07JAN07.D from x, y = 10.009, 0 to 10.059, 0; result = 88			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/11/2022 4:09:06 PM	Manually integrate qualifier91.0 of compound m+p-Xylenes in sample 07JAN07.D from x, y = 10.006, 0 to 10.078, 0; result = 443			✓	
CmdZeroOutPeak	BL2000\mchavez	1/11/2022 4:09:08 PM	Zero out primary peak of compound Ethylbenzene in sample 07JAN07.D			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/11/2022 4:09:13 PM	Manually integrate compound Ethylbenzene in sample 07JAN07.D, from x, y = 9.903, 0 to 9.942, 0, result = 389; previous integration is from x, y = 9.917, 0 to 9.917, 0 and previous response = 0.			✓	
CmdZeroOutPeak	BL2000\mchavez	1/11/2022 4:09:17 PM	Zero out primary peak of compound m+p-Xylenes in sample 07JAN07.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/11/2022 4:10:30 PM	Set SampleApproved = True for sample 07JAN07.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/11/2022 4:10:42 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/11/2022 4:11:11 PM	Set UserDefined = Qualifier ratio did not meet method criteria for m+p Xylenes for sample 07JAN07.D; previous value =			✓	
CmdSaveBatchTable	BL2000\mchavez	1/11/2022 4:54:29 PM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdOpenBatchTable	BL2000\mchavez	1/12/2022 9:04:15 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:04:32 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN08.D from x, y = 1.369, 0 to 1.445, 0; result = 737			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:04:52 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN08.D from x, y = 3.294, 0 to 3.391, 0; result = 1567			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:04:56 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN08.D from x, y = 3.294, 0 to 3.361, 0; result = 978			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:05:17 AM	Manually integrate compound Benzene in sample 07JAN08.D from x, y = 6.264, 0 to 6.305, 0; result = 212			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:05:19 AM	Set UserAnnotation = NI for compound Benzene in sample 07JAN08.D; previous value =			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:05:21 AM	Manually integrate qualifier77.0 of compound Benzene in sample 07JAN08.D from x, y = 6.264, 0 to 6.314, 0; result = 84			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:05:53 AM	Manually integrate compound Ethylbenzene in sample 07JAN08.D from x, y = 9.903, 0 to 9.939, 0; result = 483			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:05:55 AM	Manually integrate qualifier106.0 of compound Ethylbenzene in sample 07JAN08.D from x, y = 9.900, 0 to 9.936, 0; result = 26			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:05:58 AM	Set UserAnnotation = NI for compound Ethylbenzene in sample 07JAN08.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:06:35 AM	Manually integrate compound m+p-Xylenes in sample 07JAN08.D from x, y = 10.014, 0 to 10.042, 0; result = 200			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:06:37 AM	Manually integrate qualifier91.0 of compound m+p-Xylenes in sample 07JAN08.D from x, y = 10.003, 0 to 10.064, 0; result = 712			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:06:41 AM	Set UserAnnotation = NI for compound m+p-Xylenes in sample 07JAN08.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:06:44 AM	Set UserAnnotation = for compound m+p-Xylenes in sample 07JAN08.D; previous value = NI			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:06:47 AM	Zero out primary peak of compound m+p-Xylenes in sample 07JAN08.D			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:07:09 AM	Set UserDefined = Qualifier ratio did not meet method criteria for m+p Xylenes for sample 07JAN08.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:07:32 AM	Set SampleApproved = True for sample 07JAN08.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/12/2022 9:07:50 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:09:43 AM	Manually integrate compound Chloroform in sample 07JAN09.D from x, y = 5.617, 0 to 5.689, 0; result = 381			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:09:46 AM	Manually integrate qualifier85.0 of compound Chloroform in sample 07JAN09.D from x, y = 5.608, 0 to 5.675, 0; result = 250			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:10:07 AM	Manually integrate compound Methylene chloride in sample 07JAN09.D, from x, y = 3.274, 0 to 3.375, 0, result = 2658; previous integration is from x, y = 3.316, 0 to 3.375, 0 and previous response = 2339.			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:10:09 AM	Set UserAnnotation = LT for compound Methylene chloride in sample 07JAN09.D; previous value =			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:10:12 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN09.D from x, y = 3.305, 0 to 3.386, 0; result = 1818			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:10:14 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN09.D from x, y = 3.305, 0 to 3.377, 0; result = 1137			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:10:26 AM	Manually integrate compound Chloromethane in sample 07JAN09.D from x, y = 1.372, 0 to 1.456, 0; result = 1041			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:10:28 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN09.D from x, y = 1.367, 0 to 1.462, 0; result = 421			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:10:35 AM	Set UserAnnotation = NI for compound Chloromethane in sample 07JAN09.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:10:41 AM	Set UserAnnotation = NI for compound Chloroform in sample 07JAN09.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:10:44 AM	Set SampleApproved = True for sample 07JAN09.D; previous value = False			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:11:07 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN10.D from x, y = 1.381, 0 to 1.445, 0; result = 734			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:11:21 AM	Manually integrate compound Methylene chloride in sample 07JAN10.D from x, y = 3.291, 0 to 3.400, 0; result = 1975			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:11:23 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN10.D from x, y = 3.291, 0 to 3.400, 0; result = 1289			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:11:25 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN10.D from x, y = 3.299, 0 to 3.414, 0; result = 796			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:11:28 AM	Set UserAnnotation = NI for compound Methylene chloride in sample 07JAN10.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:11:40 AM	Manually integrate compound Chloroform in sample 07JAN10.D from x, y = 5.617, 0 to 5.689, 0; result = 53			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:11:42 AM	Manually integrate qualifier85.0 of compound Chloroform in sample 07JAN10.D from x, y = 5.625, 0 to 5.709, 0; result = 110			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:11:47 AM	Zero out primary peak of compound Chloroform in sample 07JAN10.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:11:58 AM	Set UserDefined = Qualifier ratio did not meet method criteria for m+p Xylenes for sample 07JAN10.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:12:09 AM	Set UserDefined = Qualifier ratio did not meet method criteria for Chloroform for sample 07JAN10.D; previous value = Qualifier ratio did not meet method criteria for m+p Xylenes			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:12:31 AM	Manually integrate compound Benzene in sample 07JAN10.D from x, y = 6.250, 0 to 6.328, 0; result = 229			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:12:32 AM	Manually integrate qualifier77.0 of compound Benzene in sample 07JAN10.D from x, y = 6.258, 0 to 6.319, 0; result = 33			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:12:43 AM	Set UserAnnotation = NI for compound Benzene in sample 07JAN10.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:12:53 AM	Manually integrate compound Bromodichloromethane in sample 07JAN10.D from x, y = 7.541, 0 to 7.633, 0; result = 903			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:12:55 AM	Manually integrate qualifier85.0 of compound Bromodichloromethane in sample 07JAN10.D from x, y = 7.544, 0 to 7.636, 0; result = 309			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:12:57 AM	Manually integrate qualifier127.0 of compound Bromodichloromethane in sample 07JAN10.D from x, y = 7.558, 0 to 7.630, 0; result = 35			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:12:58 AM	Manually integrate qualifier 127.0 of compound Bromodichloromethane in sample 07JAN10.D, from x, y = 7.613, 0 to 7.630, 0, result = 0; previous integration is from x, y = 7.558, 0 to 7.630, 0 and previous response = 35.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:13:00 AM	Manually integrate qualifier 127.0 of compound Bromodichloromethane in sample 07JAN10.D, from x, y = 7.563, 0 to 7.599, 0, result = 35; previous integration is from x, y = 7.613, 0 to 7.630, 0 and previous response = 0.			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:13:05 AM	Zero out primary peak of compound Bromodichloromethane in sample 07JAN10.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:13:26 AM	Set UserDefined = Qualifier ratio did not meet method criteria for Chloroform, Bromodichloromethane for sample 07JAN10.D; previous value = Qualifier ratio did not meet method criteria for Chloroform			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/12/2022 9:13:44 AM	Manually integrate compound Chlorodibromomethane in sample 07JAN10.D, from x, y = 9.172, 0 to 9.253, 0, result = 5690; previous integration is from x, y = 9.172, 0 to 9.203, 0 and previous response = 2830.			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/12/2022 9:14:23 AM	Manually integrate compound m+p-Xylenes in sample 07JAN10.D from x, y = 10.020, 0 to 10.062, 0; result = 93			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:14:26 AM	Manually integrate qualifier91.0 of compound m+p-Xylenes in sample 07JAN10.D from x, y = 10.009, 0 to 10.078, 0; result = 397			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:14:30 AM	Zero out primary peak of compound m+p-Xylenes in sample 07JAN10.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:15:43 AM	Set UserDefined = Qualifier ratio did not meet method criteria for Chloroform, Bromodichloromethane, m+p Xylenes for sample 07JAN10.D; previous value = Qualifier ratio did not meet method criteria for Chloroform, Bromodichloromethane			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSaveBatchTable	BL2000\mchavez	1/12/2022 9:15:57 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	1/12/2022 9:29:47 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:30:26 AM	Manually integrate qualifier174.5 of compound Bromoform in sample 07JAN10.D from x, y = 10.594, 0 to 10.661, 0; result = 1221			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:30:28 AM	Manually integrate qualifier170.5 of compound Bromoform in sample 07JAN10.D from x, y = 10.580, 0 to 10.692, 0; result = 1208			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:30:43 AM	Set SampleApproved = True for sample 07JAN10.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:31:16 AM	Manually integrate compound o-Xylene in sample 07JAN11.D from x, y = 10.405, 0 to 10.466, 0; result = 64			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:31:17 AM	Manually integrate qualifier91.0 of compound o-Xylene in sample 07JAN11.D from x, y = 10.416, 0 to 10.458, 0; result = 44			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:31:20 AM	Zero out primary peak of compound o-Xylene in sample 07JAN11.D			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:31:30 AM	Manually integrate compound m+p-Xylenes in sample 07JAN11.D from x, y = 10.028, 0 to 10.056, 0; result = 316			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:31:34 AM	Manually integrate qualifier91.0 of compound m+p-Xylenes in sample 07JAN11.D from x, y = 10.031, 0 to 10.062, 0; result = 279			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:31:37 AM	Zero out primary peak of compound m+p-Xylenes in sample 07JAN11.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:31:47 AM	Set UserDefined = Qualifier ratio did not meet method criteria for m+p Xylenes for sample 07JAN11.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:32:00 AM	Set UserDefined = Qualifier ratio did not meet method criteria for m+p Xylenes, o-Xylene for sample 07JAN11.D; previous value = Qualifier ratio did not meet method criteria for m+p Xylenes			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:32:23 AM	Manually integrate compound 1,3-Dichloropropane in sample 07JAN11.D from x, y = 8.946, 0 to 9.022, 0; result = 319			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:32:25 AM	Manually integrate qualifier78.0 of compound 1,3-Dichloropropane in sample 07JAN11.D from x, y = 8.946, 0 to 9.005, 0; result = 72			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:32:28 AM	Set UserAnnotation = NI for compound 1,3-Dichloropropane in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:32:37 AM	Manually integrate compound 1,1,2-Trichloroethane in sample 07JAN11.D from x, y = 8.776, 0 to 8.854, 0; result = 383			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:32:39 AM	Manually integrate qualifier97.0 of compound 1,1,2-Trichloroethane in sample 07JAN11.D from x, y = 8.782, 0 to 8.865, 0; result = 527			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:32:43 AM	Manually integrate qualifier85.0 of compound 1,1,2-Trichloroethane in sample 07JAN11.D from x, y = 8.787, 0 to 8.854, 0; result = 86			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:32:47 AM	Zero out primary peak of compound 1,1,2-Trichloroethane in sample 07JAN11.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:33:16 AM	Set UserDefined = Qualifier ratio did not meet method criteria for m+p Xylenes, o-Xylene, 1,1,2-Trichloroethane for sample 07JAN11.D; previous value = Qualifier ratio did not meet method criteria for m+p Xylenes, o-Xylene			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:33:28 AM	Manually integrate compound Toluene in sample 07JAN11.D from x, y = 8.361, 0 to 8.430, 0; result = 699			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:33:30 AM	Manually integrate qualifier91.0 of compound Toluene in sample 07JAN11.D from x, y = 8.349, 0 to 8.433, 0; result = 1384			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:33:36 AM	Set UserAnnotation = NI for compound Toluene in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:35:15 AM	Manually integrate compound Benzene in sample 07JAN11.D from x, y = 6.205, 0 to 6.333, 0; result = 1404			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:35:16 AM	Manually integrate qualifier77.0 of compound Benzene in sample 07JAN11.D from x, y = 6.250, 0 to 6.314, 0; result = 156			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:35:20 AM	Set UserAnnotation = NI for compound Benzene in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:35:40 AM	Manually integrate compound Chloroform in sample 07JAN11.D from x, y = 5.614, 0 to 5.711, -70; result = 983			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:35:45 AM	Manually integrate compound Chloroform in sample 07JAN11.D, from x, y = 5.614, 0 to 5.689, 0, result = 777; previous integration is from x, y = 5.614, 0 to 5.711, -70 and previous response = 983.			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:35:47 AM	Manually integrate qualifier85.0 of compound Chloroform in sample 07JAN11.D from x, y = 5.628, 0 to 5.700, 0; result = 368			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:35:49 AM	Set UserAnnotation = NI for compound Chloroform in sample 07JAN11.D; previous value =			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:36:02 AM	Manually integrate compound 1,1-Dichloroethane in sample 07JAN11.D from x, y = 4.328, 0 to 4.459, 0; result = 0				Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for compound 1,1-Dichloroethane in sample B22010219-001F. ---> Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for compound 1,1-Dichloroethane in sample B22010219-001F. ---> System.IndexOutOfRangeException: Index was outside the bounds of the array. at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.a(Double[] A_0, Single[] A_1, Int32 A_2, Int32 A_3, Int32 A_4, Double A_5, Double A_6, Double& A_7, Double& A_8, Int32& A_9, Int32& A_10, Int32& A_11, Int32& A_12) at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.ComputeChromatographicMetrics(Double[] xArray, Single[] yArray, Int32 startIndex, Int32 apexIndex, Int32 endIndex, Double baselineSlope, Double yIntercept, Double& fullWidthHalfMaximum, Double& symmetry) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.a(IChromatogram A_0, IPeakList A_1) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.ManualIntegrate(IChromatogram chromatogram, Double xStart, Double yStart, Double xEnd, Double yEnd, IChromPeakList& peaklist) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd) --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.QuantifierIon.SetManualIntegrationFailureMessage(Exception e) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd)

Name	User	Time	Action	Reason	Comment	Succeed	Exception
							at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegratePeak.Do() --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegratePeak.Do() at Agilent.MassSpectrometry.CommandModel.CommandHistory.Invoke(ICommand cmd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.AppCommandContext._Invoke(ICommand cmd)

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:36:06 AM	Manually integrate qualifier65.0 of compound 1,1-Dichloroethane in sample 07JAN11.D from x, y = 4.328, 0 to 4.401, 0; result = 0				Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for qualifier 65.0 of compound 1,1-Dichloroethane in sample B22010219-001F. ---> Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for qualifier 65.0 of compound 1,1-Dichloroethane in sample B22010219-001F. ---> System.IndexOutOfRangeException: Index was outside the bounds of the array. at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.a(Double[] A_0, Single[] A_1, Int32 A_2, Int32 A_3, Int32 A_4, Double A_5, Double A_6, Double& A_7, Double& A_8, Int32& A_9, Int32& A_10, Int32& A_11, Int32& A_12) at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.ComputeChromatographicMetrics(Double[] xArray, Single[] yArray, Int32 startIndex, Int32 apexIndex, Int32 endIndex, Double baselineSlope, Double yIntercept, Double& fullWidthHalfMaximum, Double& symmetry) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.a(IChromatogram A_0, IPeakList A_1) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.ManualIntegrate(IChromatogram chromatogram, Double xStart, Double yStart, Double xEnd, Double yEnd, IChromPeakList& peaklist) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd) --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.QualifierIon.SetManualIntegrationFailureMessage(Exception e) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.M

Name	User	Time	Action	Reason	Comment	Succeed	Exception
							anualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegrateQualifierPeak.Do() --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegrateQualifierPeak.Do() at Agilent.MassSpectrometry.CommandModel.CommandHistory.Invoke(ICommand cmd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.AppCommandContext._Invoke(ICommand cmd)
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:36:10 AM	Manually integrate qualifier83.0 of compound 1,1-Dichloroethane in sample 07JAN11.D from x, y = 4.348, 0 to 4.445, 0; result = 0			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:36:13 AM	Manually integrate compound 1,1-Dichloroethane in sample 07JAN11.D from x, y = 4.339, 0 to 4.448, 0; result = 953			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:36:16 AM	Manually integrate qualifier65.0 of compound 1,1-Dichloroethane in sample 07JAN11.D from x, y = 4.320, 0 to 4.412, 0; result = 177			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:36:19 AM	Set UserAnnotation = NI for compound 1,1-Dichloroethane in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:37:13 AM	Manually integrate compound Methylene chloride in sample 07JAN11.D from x, y = 3.291, 0 to 3.394, 0; result = 1285			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:37:15 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN11.D from x, y = 3.296, 0 to 3.377, 0; result = 1048			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:37:18 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN11.D from x, y = 3.299, 0 to 3.394, 0; result = 702			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:37:20 AM	Set UserAnnotation = NI for compound Methylene chloride in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:37:32 AM	Manually integrate compound Chloroethane in sample 07JAN11.D from x, y = 1.869, 0 to 1.922, 0; result = 1604			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:37:35 AM	Manually integrate qualifier66.0 of compound Chloroethane in sample 07JAN11.D from x, y = 1.869, 0 to 1.911, -2; result = 433			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:37:37 AM	Set UserAnnotation = NI for compound Chloroethane in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:37:48 AM	Manually integrate compound Chloromethane in sample 07JAN11.D from x, y = 1.381, 0 to 1.456, 0; result = 1302			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:37:51 AM	Set UserAnnotation = NI for compound Chloromethane in sample 07JAN11.D; previous value =			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:37:53 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN11.D from x, y = 1.378, 0 to 1.459, 0; result = 301			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:37:58 AM	Set SampleApproved = True for sample 07JAN11.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/12/2022 9:38:16 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:38:53 AM	Manually integrate compound Chloromethane in sample 07JAN12.D from x, y = 1.372, 0 to 1.442, 0; result = 854			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:38:56 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN12.D from x, y = 1.403, 0 to 1.450, 0; result = 97			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:38:59 AM	Set UserAnnotation = NI for compound Chloromethane in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:39:27 AM	Manually integrate compound Methyl ethyl ketone in sample 07JAN12.D from x, y = 5.265, 0 to 5.352, 0; result = 565			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:39:30 AM	Manually integrate qualifier72.0 of compound Methyl ethyl ketone in sample 07JAN12.D from x, y = 5.254, 0 to 5.352, 0; result = 130			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:39:43 AM	Manually integrate compound Chloroform in sample 07JAN12.D from x, y = 5.608, 0 to 5.706, 0; result = 1655			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:39:46 AM	Manually integrate qualifier85.0 of compound Chloroform in sample 07JAN12.D from x, y = 5.608, 0 to 5.714, 0; result = 947			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:40:08 AM	Manually integrate compound Benzene in sample 07JAN12.D from x, y = 6.252, 0 to 6.314, 0; result = 409			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:40:11 AM	Manually integrate qualifier 65.0 of compound 1,2-Dichloroethane-d4 in sample 07JAN12.D, from x, y = 6.261, 0 to 6.275, 2250, result = 3576; previous integration is from x, y = 6.188, 0 to 6.316, 0 and previous response = 182216.			✓	
CmdClearManualIntegra tion	BL2000\mchavez	1/12/2022 9:40:16 AM	Clear manual integration of qualifier 65.0 for compound 1,2-Dichloroethane-d4 in sample 07JAN12.D			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:40:23 AM	Manually integrate qualifier77.0 of compound Benzene in sample 07JAN12.D from x, y = 6.244, 0 to 6.305, 0; result = 43			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/12/2022 9:40:34 AM	Manually integrate compound Dibromomethane in sample 07JAN12.D from x, y = 7.373, 0 to 7.426, 0; result = 111			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:40:36 AM	Manually integrate qualifier95.0 of compound Dibromomethane in sample 07JAN12.D from x, y = 7.376, 0 to 7.429, 0; result = 67			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:40:38 AM	Manually integrate qualifier173.5 of compound Dibromomethane in sample 07JAN12.D from x, y = 7.373, 0 to 7.437, 0; result = 135			✓	
CmdSetTargetCompoun dAttribute	BL2000\mchavez	1/12/2022 9:40:40 AM	Set UserAnnotation = NI for compound Dibromomethane in sample 07JAN12.D; previous value =			✓	
CmdSetTargetCompoun dAttribute	BL2000\mchavez	1/12/2022 9:40:44 AM	Set UserAnnotation = NI for compound Benzene in sample 07JAN12.D; previous value =			✓	
CmdSetTargetCompoun dAttribute	BL2000\mchavez	1/12/2022 9:40:47 AM	Set UserAnnotation = NI for compound Methyl ethyl ketone in sample 07JAN12.D; previous value =			✓	
CmdSetTargetCompoun dAttribute	BL2000\mchavez	1/12/2022 9:40:51 AM	Set UserAnnotation = NI for compound Chloroform in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegrateP eak	BL2000\mchavez	1/12/2022 9:40:57 AM	Manually integrate compound Bromodichloromethane in sample 07JAN12.D from x, y = 7.557, 0 to 7.624, 0; result = 654			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:40:59 AM	Manually integrate qualifier85.0 of compound Bromodichloromethane in sample 07JAN12.D from x, y = 7.543, 0 to 7.633, 0; result = 391			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:41:01 AM	Manually integrate qualifier127.0 of compound Bromodichloromethane in sample 07JAN12.D from x, y = 7.552, 0 to 7.608, 0; result = 34			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:41:10 AM	Set UserAnnotation = NI for compound Bromodichloromethane in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:41:22 AM	Manually integrate compound Toluene in sample 07JAN12.D from x, y = 8.369, 0 to 8.422, 0; result = 493			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:41:25 AM	Manually integrate qualifier91.0 of compound Toluene in sample 07JAN12.D from x, y = 8.338, 0 to 8.436, 0; result = 0				Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for qualifier 91.0 of compound Toluene in sample B22010213-002C. ---> Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for qualifier 91.0 of compound Toluene in sample B22010213-002C. ---> System.IndexOutOfRangeException: Index was outside the bounds of the array. at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.a(Double[] A_0, Single[] A_1, Int32 A_2, Int32 A_3, Int32 A_4, Double A_5, Double A_6, Double& A_7, Double& A_8, Int32& A_9, Int32& A_10, Int32& A_11, Int32& A_12) at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.ComputeChromatographicMetrics(Double[] xArray, Single[] yArray, Int32 startIndex, Int32 apexIndex, Int32 endIndex, Double baselineSlope, Double yIntercept, Double& fullWidthHalfMaximum, Double& symmetry) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.a(IChromatogram A_0, IPeakList A_1) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.ManualIntegrate(IChromatogram chromatogram, Double xStart, Double yStart, Double xEnd, Double yEnd, IChromPeakList& peaklist) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd) --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.QualifierIon.SetManualIntegrationFailureMessage(Exception e) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd)

Name	User	Time	Action	Reason	Comment	Succeed	Exception
							at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegrateQualifierPeak.Do() --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegrateQualifierPeak.Do() at Agilent.MassSpectrometry.CommandModel.CommandHistory.Invoke(ICommand cmd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.AppCommandContext._Invoke(ICommand cmd)
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:41:29 AM	Manually integrate qualifier91.0 of compound Toluene in sample 07JAN12.D from x, y = 8.352, 0 to 8.427, 0; result = 888			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:41:32 AM	Set UserAnnotation = NI for compound Toluene in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:41:48 AM	Manually integrate compound Chlorodibromomethane in sample 07JAN12.D from x, y = 9.169, 0 to 9.239, 0; result = 1602			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:41:50 AM	Set UserAnnotation = NI for compound Chlorodibromomethane in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:41:52 AM	Manually integrate qualifier127.0 of compound Chlorodibromomethane in sample 07JAN12.D from x, y = 9.175, 0 to 9.244, 0; result = 973			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:42:09 AM	Manually integrate compound Ethylbenzene in sample 07JAN12.D, from x, y = 9.875, 0 to 9.944, 0, result = 1441; previous integration is from x, y = 10.009, 0 to 10.059, 0 and previous response = 3201.			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:42:12 AM	Manually integrate qualifier106.0 of compound Ethylbenzene in sample 07JAN12.D from x, y = 9.908, 0 to 9.958, 0; result = 144			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:42:16 AM	Set UserAnnotation = NI for compound Ethylbenzene in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:42:21 AM	Manually integrate compound m+p-Xylenes in sample 07JAN12.D from x, y = 10.006, 0 to 10.070, 0; result = 1346			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:42:32 AM	Manually integrate qualifier 91.0 of compound m+p-Xylenes in sample 07JAN12.D, from x, y = 10.009, 0 to 10.059, 26, result = 3162; previous integration is from x, y = 10.009, 0 to 10.073, 0 and previous response = 3521.			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/12/2022 9:42:35 AM	Drop baseline for qualifier 91.0 of compound m+p-Xylenes in sample 07JAN12.D to y = 0, new integration is from x, y = 10.009, 0 to 10.059, 0 and new response = 3201; previous integration is from x, y = 10.009, 0 to 10.059, 26 and previous response = 3162.			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:42:39 AM	Manually integrate qualifier 91.0 of compound m+p-Xylenes in sample 07JAN12.D, from x, y = 10.017, 19 to 10.059, 0, result = 3051; previous integration is from x, y = 10.009, 0 to 10.059, 0 and previous response = 3201.			✓	
CmdManuallyIntegrate DropBaseline	BL2000\mchavez	1/12/2022 9:42:41 AM	Drop baseline for qualifier 91.0 of compound m+p-Xylenes in sample 07JAN12.D to y = 0, new integration is from x, y = 10.017, 0 to 10.059, 0 and new response = 3074; previous integration is from x, y = 10.017, 19 to 10.059, 0 and previous response = 3051.			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:42:48 AM	Set UserAnnotation = NI for compound m+p-Xylenes in sample 07JAN12.D; previous value =			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:43:02 AM	Manually integrate qualifier174.5 of compound Bromoform in sample 07JAN12.D from x, y = 10.586, 0 to 10.681, 0; result = 1218			✓	
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:43:04 AM	Manually integrate qualifier170.5 of compound Bromoform in sample 07JAN12.D from x, y = 10.575, 0 to 10.678, 0; result = 1221			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:43:09 AM	Zero out primary peak of compound 1,1,2,2-Tetrachloroethane in sample 07JAN12.D			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:43:18 AM	Zero out primary peak of compound 1,2,3-Trichloropropane in sample 07JAN12.D			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:43:27 AM	Zero out primary peak of compound 4-Chlorotoluene in sample 07JAN12.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:43:35 AM	Set SampleApproved = True for sample 07JAN12.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/12/2022 9:43:56 AM	Quantitate all compounds in all samples			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate QualifierPeak	BL2000\mchavez	1/12/2022 9:45:41 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN14.D from x, y = 1.372, 0 to 1.436, 0; result = 725			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:46:20 AM	Manually integrate compound Toluene in sample 07JAN14.D from x, y = 8.338, 0 to 8.444, 0; result = 0				Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for compound Toluene in sample B22010214-001F. ---> Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Manual integration failed for compound Toluene in sample B22010214-001F. ---> System.IndexOutOfRangeException: Index was outside the bounds of the array. at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.a(Double[] A_0, Single[] A_1, Int32 A_2, Int32 A_3, Int32 A_4, Double A_5, Double A_6, Double& A_7, Double& A_8, Int32& A_9, Int32& A_10, Int32& A_11, Int32& A_12) at Agilent.MassSpectrometry.DataAnalysis.AgileIntegrator.Peak.ComputeChromatographicMetrics(Double[] xArray, Single[] yArray, Int32 startIndex, Int32 apexIndex, Int32 endIndex, Double baselineSlope, Double yIntercept, Double& fullWidthHalfMaximum, Double& symmetry) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.a(IChromatogram A_0, IPeakList A_1) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Integrator.ManualIntegrate(IChromatogram chromatogram, Double xStart, Double yStart, Double xEnd, Double yEnd, IChromPeakList& peaklist) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd) --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.QuantifierIon.SetManualIntegrationFailureMessage(Exception e) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ManualIntegrate(Double xStart, Double yStart, Double xEnd, Double yEnd)

Name	User	Time	Action	Reason	Comment	Succeed	Exception
							at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegratePeak.Do() --- End of inner exception stack trace --- at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdManuallyIntegratePeak.Do() at Agilent.MassSpectrometry.CommandModel.CommandHistory.Invoke(ICommand cmd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.AppCommandContext._Invoke(ICommand cmd)
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:46:25 AM	Manually integrate qualifier91.0 of compound Toluene in sample 07JAN14.D from x, y = 8.341, 0 to 8.419, 0; result = 0			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:46:28 AM	Manually integrate compound Toluene in sample 07JAN14.D from x, y = 8.344, 0 to 8.433, 0; result = 667			✓	
CmdClearManualIntegration	BL2000\mchavez	1/12/2022 9:46:49 AM	Clear manual integration of target signal for compound Toluene in sample 07JAN14.D				Agilent.MassSpectrometry.DataAnalysis.Quantitative.ApplicationCommandException: Integrator did not find any peaks at Agilent.MassSpectrometry.DataAnalysis.Quantitative.Analysis.MeasuredIon.ClearManualIntegration() at Agilent.MassSpectrometry.DataAnalysis.Quantitative.CmdClearManualIntegration.Do() at Agilent.MassSpectrometry.CommandModel.CommandHistory.Invoke(ICommand cmd) at Agilent.MassSpectrometry.DataAnalysis.Quantitative.AppCommandContext._Invoke(ICommand cmd)
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:47:15 AM	Manually integrate compound Toluene in sample 07JAN14.D from x, y = 8.355, 0 to 8.436, 0; result = 667			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:47:17 AM	Zero out primary peak of compound Toluene in sample 07JAN14.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:47:40 AM	Set UserDefined = Qualifier ratio did not meet method criteria for Toluene for sample 07JAN14.D; previous value =			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:48:16 AM	Manually integrate compound Methylene chloride in sample 07JAN14.D from x, y = 3.302, 0 to 3.400, 0; result = 885			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:48:18 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN14.D from x, y = 3.302, 0 to 3.375, 0; result = 448			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:48:20 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN14.D from x, y = 3.302, 0 to 3.383, 0; result = 313			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:48:24 AM	Set UserAnnotation = NI for compound Methylene chloride in sample 07JAN14.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:48:49 AM	Set SampleApproved = True for sample 07JAN14.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:49:15 AM	Manually integrate compound Dichlorodifluoromethane in sample 07JAN15.D from x, y = 1.219, 0 to 1.300, 0; result = 1185			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:49:17 AM	Manually integrate qualifier87.0 of compound Dichlorodifluoromethane in sample 07JAN15.D from x, y = 1.202, 0 to 1.280, 0; result = 350			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:49:20 AM	Set UserAnnotation = NI for compound Dichlorodifluoromethane in sample 07JAN15.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:49:24 AM	Manually integrate compound Chloromethane in sample 07JAN15.D from x, y = 1.386, 0 to 1.422, 0; result = 626			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:49:27 AM	Manually integrate qualifier52.0 of compound Chloromethane in sample 07JAN15.D from x, y = 1.397, 0 to 1.434, 0; result = 61			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:49:32 AM	Set UserAnnotation = NI for compound Chloromethane in sample 07JAN15.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:49:51 AM	Manually integrate compound Methylene chloride in sample 07JAN15.D from x, y = 3.299, 0 to 3.372, 0; result = 648			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:49:54 AM	Manually integrate qualifier84.0 of compound Methylene chloride in sample 07JAN15.D from x, y = 3.280, 0 to 3.405, 0; result = 450			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:49:56 AM	Manually integrate qualifier86.0 of compound Methylene chloride in sample 07JAN15.D from x, y = 3.291, 0 to 3.383, 0; result = 284			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:49:58 AM	Set UserAnnotation = NI for compound Methylene chloride in sample 07JAN15.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:50:12 AM	Manually integrate compound Chloroform in sample 07JAN15.D from x, y = 5.605, 0 to 5.697, 0; result = 600			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:50:14 AM	Manually integrate qualifier85.0 of compound Chloroform in sample 07JAN15.D from x, y = 5.636, 0 to 5.695, 0; result = 88			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:50:22 AM	Manually integrate compound Benzene in sample 07JAN15.D from x, y = 6.230, 0 to 6.328, 0; result = 748			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:50:24 AM	Manually integrate qualifier77.0 of compound Benzene in sample 07JAN15.D from x, y = 6.238, 0 to 6.314, 0; result = 108			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:50:39 AM	Manually integrate compound Toluene in sample 07JAN15.D from x, y = 8.355, 0 to 8.430, 0; result = 723			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:50:42 AM	Manually integrate qualifier91.0 of compound Toluene in sample 07JAN15.D from x, y = 8.352, 0 to 8.439, 0; result = 1271			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:50:59 AM	Manually integrate compound m+p-Xylenes in sample 07JAN15.D from x, y = 9.992, 0 to 10.064, 0; result = 505			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:51:01 AM	Manually integrate qualifier91.0 of compound m+p-Xylenes in sample 07JAN15.D from x, y = 10.009, 0 to 10.064, 0; result = 1116			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	1/12/2022 9:51:06 AM	Manually integrate compound o-Xylene in sample 07JAN15.D from x, y = 10.419, 0 to 10.463, 0; result = 172			✓	
CmdManuallyIntegrateQualifierPeak	BL2000\mchavez	1/12/2022 9:51:08 AM	Manually integrate qualifier91.0 of compound o-Xylene in sample 07JAN15.D from x, y = 10.396, 0 to 10.477, 0; result = 450			✓	
CmdZeroOutPeak	BL2000\mchavez	1/12/2022 9:51:13 AM	Zero out primary peak of compound o-Xylene in sample 07JAN15.D			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:51:25 AM	Set UserDefined = Qualifier ratio did not meet method criteria for Toluene for sample 07JAN15.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:51:35 AM	Set UserDefined = Qualifier ratio did not meet method criteria for o-Xylene for sample 07JAN15.D; previous value = Qualifier ratio did not meet method criteria for Toluene			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:51:44 AM	Set SampleName = B22010262-002A for sample 07JAN09.D; previous value = B22010226-002A			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:52:03 AM	Set UserDefined = Corrected sample ID from vial run on autosampler. for sample 07JAN09.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	1/12/2022 9:52:16 AM	Set UserAnnotation = NI for compound m+p-Xylenes in sample 07JAN15.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:52:52 AM	Set SampleApproved = True for sample 07JAN15.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/12/2022 9:53:07 AM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:55:03 AM	Set SampleApproved = True for sample 07JAN17.D; previous value = False			✓	
CmdQuantitate	BL2000\mchavez	1/12/2022 9:55:20 AM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 9:59:39 AM	Set SampleApproved = True for sample 07JAN18.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\mchavez	1/12/2022 10:01:17 AM	Set SampleApproved = True for sample 07JAN20.D; previous value = False			✓	
CmdSaveBatchTable	BL2000\mchavez	1/12/2022 10:01:26 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
CmdOpenBatchTable	BL2000\mchavez	3/2/2022 10:30:43 AM	Open batch D:\Org\Data\VOA5975C\VG010722\VG010722_8260B.batch.bin			✓	
CmdSetSampleAttribute	BL2000\mchavez	3/2/2022 10:31:09 AM	Set SampleType = MatrixBlank for sample 07JAN11.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\mchavez	3/2/2022 10:31:15 AM	Set SampleApproved = True for sample 07JAN02.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	3/2/2022 10:31:59 AM	Set UserAnnotation = LT for compound Chlorodibromomethane in sample 07JAN10.D; previous value =			✓	
CmdZeroOutPeak	BL2000\mchavez	3/2/2022 10:33:29 AM	Zero out primary peak of compound Chloroform in sample 07JAN15.D			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	3/2/2022 10:33:57 AM	Manually integrate compound Chloroform in sample 07JAN15.D, from x, y = 5.653, 0 to 5.689, 0, result = 129; previous integration is from x, y = 5.653, 0 to 5.653, 0 and previous response = 0.			✓	
CmdManuallyIntegratePeak	BL2000\mchavez	3/2/2022 10:34:00 AM	Manually integrate compound Chloroform in sample 07JAN15.D, from x, y = 5.600, 0 to 5.689, 0, result = 600; previous integration is from x, y = 5.653, 0 to 5.689, 0 and previous response = 129.			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\mchavez	3/2/2022 10:34:14 AM	Set UserAnnotation = NI for compound Chloroform in sample 07JAN15.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	3/2/2022 10:34:20 AM	Set UserAnnotation = NI for compound Benzene in sample 07JAN15.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\mchavez	3/2/2022 10:34:23 AM	Set UserAnnotation = NI for compound Toluene in sample 07JAN15.D; previous value =			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdCalibrate	BL2000\mchavez	3/2/2022 10:36:54 AM	Replace level QC with QC sample 07JAN04.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Toluene}; Replace level CC with CC sample 07JAN03.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Toluene};				
CmdQuantitate	BL2000\mchavez	3/2/2022 10:37:16 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	3/2/2022 10:40:15 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
GenerateReport	BL2000\mchavez	3/2/2022 10:41:04 AM	Generates report - Method: \\MASSHUNTER\Org\reports\LevelIV_Reports\SampleSequence\CC_mid_rpt.m, Output Path: D:\Org\Data\VOA5975C\VG010722\QuantReports\VG010722_8260B			✓	
CmdCalibrate	BL2000\mchavez	3/2/2022 10:41:44 AM	Replace level CC with CC sample 07JAN20.D for compounds {1,2-Dichlorobenzene, 1,4-Dichlorobenzene, 1,3-Dichlorobenzene, 4-Chlorotoluene, 2-Chlorotoluene, 1,2,3-Trichloropropane, 1,1,2,2-Tetrachloroethane, Bromobenzene, p-Bromofluorobenzene, Bromoform, Styrene, o-Xylene, m+p-Xylenes, Ethylbenzene, 1,1,1,2-Tetrachloroethane, Chlorobenzene, 1,2-Dibromoethane, Chlorodibromomethane, 1,3-Dichloropropane, Tetrachloroethene, 1,1,2-Trichloroethane, trans-1,3-Dichloropropene, Toluene-d8, cis-1,3-Dichloropropene, Bromodichloromethane, Dibromomethane, 1,2-Dichloropropane, Trichloroethene, 1,2-Dichloroethane, Benzene, 1,2-Dichloroethane-d4, 1,1-Dichloropropene, Carbon tetrachloride, Dibromofluoromethane, 1,1,1-Trichloroethane, Chloroform, Bromochloromethane, Methyl ethyl ketone, cis-1,2-Dichloroethene, 2,2-Dichloropropane, 1,1-Dichloroethane, Methyl tert-butyl ether (MTBE), trans-1,2-Dichloroethene, Methylene chloride, 1,1-Dichloroethene, Trichlorofluoromethane, Chloroethane, Bromomethane, Vinyl chloride, Chloromethane, Dichlorodifluoromethane, Toluene};			✓	

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdQuantitate	BL2000\mchavez	3/2/2022 10:43:26 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\mchavez	3/2/2022 10:44:13 AM	Save batch D:\Org\Data\VOA5975C\VG010722\QuantResults\VG010722_8260B.batch.bin			✓	
GenerateReport	BL2000\mchavez	3/2/2022 10:44:51 AM	Generates report - Method: \\MASSHUNTER\Org\reports\LevelIV_Reports\SampleSequence\CC_mid_rpt.m, Output Path: D:\Org\Data\VOA5975C\VG010722\QuantReports\VG010722_8260B-1			✓	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF0313
Standard Name: Liquids
Type: Primary
Prep Date: 6/23/2020
Prep By: Alethea M. Shaules
Exp Date: 4/13/2023
Status: New
Department: gcmsvoa
Vendor: AccuStd
Final Volume: 10 mL
Lot Number: 220041126
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL. Catalog # M502A-R-10X. Corrected lot number to match CI. MSC 01/14/2022

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Volatile Organic Compounds - Liquids	12797	1	mL	4/13/2023

Stock Source	Base Units	Amount Added
VOCF0313	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF0352
Spike Name: 2nd Source Liquids
Prep Date: 11/23/2020
Exp Date: 12/31/2023
Department: gcmsvoa
Vendor: Agilent
Lot Number: 0006570990
Balance ID:
Comments: Date Prepared is same as Date Received. 2000 ug/mL in MeOH. Catalog # DWM-589N-1.

Type: Primary
Prep By: Steve Dilts
Status: New
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
VOC Standard	13292	1	mL	12/31/2023

Stock Source	Base Units	Amount Added
VOCF0352	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF0364
Spike Name: Surrogates 2.0 mg/mL
Type: Primary
Prep Date: 1/6/2021
Prep By: Jerran D. Brenden
Exp Date: 4/18/2029
Status: New
Department: gcmsvoa
Vendor: AccuStandard
Final Volume: 10 mL
Lot Number: 219041458
Balance ID:
Comments: Date Received 01/04/2021. 2.0 mg/mL. Catalog # M-8260A-B-SS-10X

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Surrogate Standard Mix	13385	1	mL	4/18/2029

Stock Source	Base Units	Amount Added
VOCF0364	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF0373
Standard Name: MtBE (Methy tert-Butyl Ether)
Type: Primary
Prep Date: 2/26/2021
Prep By: Steve Dilts
Exp Date: 8/31/2022
Status: New
Department: gcmsvoa
Vendor: Agilent
Final Volume: 5 mL
Lot Number: 0006555762
Balance ID:
Comments: Date Prepared is same as Date Received. 2000 ug/mL in MeOH. Catalog # STS-440

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methyl tert-Butyl Ether Standard	13578	1	mL	8/31/2022

Stock Source	Base Units	Amount Added
VOCF0373	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF0401
Spike Name: 2nd Source MtBE
Type: Primary
Prep Date: 6/7/2021
Prep By: Alethea M. Shaules
Exp Date: 12/11/2029
Status: New
Department: gcmsvoa
Vendor: AccuStandard
Final Volume: 5 mL
Lot Number: 220051182
Balance ID:
Comments: Date Prepared is same as Date Receive. 2,000 ug/mL in MeOH. Catalog # S-078-10X.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
MTBE	13920	1	mL	12/11/2029

Stock Source	Base Units	Amount Added
VOCF0401	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF0417
Spike Name: Chem Service Gases
Prep Date: 8/3/2021
Exp Date: 2/28/2022
Department: gcmsvoa
Vendor: Chemservice
Lot Number: 11882100
Balance ID:
Type: Primary
Prep By: Steve Dilts
Status: New
Final Volume: 5 mL
Comments: Date Prepared is same as Date Received. 2000 ug/mL in MeOH. Catalog # M-VOHC6M5-1ML

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Volatile Organics High Concentration Mixture #6	14142	5	mL	2/28/2022

Stock Source	Base Units	Amount Added
VOCF0417	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF0425
Standard Name: Internals
Type: Primary
Prep Date: 9/8/2021
Prep By: Jerran D. Brenden
Exp Date: 12/31/2022
Status: New
Department: gcmsvoa
Vendor: Agilent
Final Volume: 10 mL
Lot Number: 0006582580
Balance ID:
Comments: Date Prepared is same as Date Received. 2,500 ug/mL in MeOH. Catalog # STM-520-1.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Internal Standard	14251	1	mL	12/31/2022

Stock Source	Base Units	Amount Added
VOCF0425	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF0426
Spike Name: Surrogates 2.0 mg/mL
Prep Date: 9/14/2021
Exp Date: 4/18/2029
Department: gcmsvoa
Vendor: AccuStandard
Lot Number: 219041458
Balance ID:
Comments: Date Received 01/04/2021. 2.0 mg/mL. Catalog # M-8260A-B-SS-10X

Type: Primary
Prep By: Jerran D. Brenden
Status: New
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Surrogate Standard Mix	14269	1	mL	4/18/2029

Stock Source	Base Units	Amount Added
VOCF0426	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF0427
Standard Name: Gases
Type: Primary
Prep Date: 9/17/2021
Prep By: Alethea M. Shaules
Exp Date: 8/3/2024
Status: New
Department: gcmsvoa
Vendor: Absolute
Final Volume: 10 mL
Lot Number: 080321
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL in MeOH. Catalog # 30058.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
EPA Method 502-524 - Volatile Gases Mix #1	14285	1	mL	8/3/2024

Stock Source	Base Units	Amount Added
VOCF0427	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF0434
Standard Name: Ketones
Type: Primary
Prep Date: 10/26/2021
Prep By: Steve Dilts
Exp Date: 6/30/2023
Status: New
Department: gcmsvoa
Vendor: Chem Service
Final Volume: 5 mL
Lot Number: 10251200
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL in 90:10 MeOH:H2O. Catalog # M-TCL-1AN5-5ML.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketone Mix	14443	1	mL	6/30/2023

Stock Source	Base Units	Amount Added
VOCF0434	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF0439
Standard Name: 2nd Source Ketones
Type: Primary
Prep Date: 11/30/2021
Prep By: Melissa Chavez
Exp Date: 11/26/2022
Status: New
Department: gcmsvoa
Vendor: AccuStandard
Final Volume: 1 mL
Lot Number: 221101480
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL in Methanol. Catalog # CLP-022K-10X.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketones Mixture	14567	2	mL	11/26/2022

Stock Source	Base Units	Amount Added
VOCF0439	ug/mL	

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF3473
Standard Name: Calibration Surrogates
Type: Secondary
Prep Date: 9/14/2021
Prep By: Jerran D. Brenden
Exp Date: 3/14/2022
Status: New
Department: gcmsvoa
Vendor:
Final Volume: 5 mL
Lot Number:
Balance ID:
Comments: Final Concentration 0.2 ug/uL in MeOH

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EA226	13754	4.5	mL	3/14/2022

Stock Source	Base Units	Amount Added
VOCF0364	ug/mL	0.5 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF3517
Spike Name: Internal Standard / Surrogates (INT/SURR)
Type: Secondary
Prep Date: 11/10/2021
Prep By: Alethea M. Shaules
Exp Date: 12/31/2022
Status: New
Department: gcmsvoa
Vendor:
Final Volume: 100 mL
Lot Number:
Balance ID:
Comments: Final Concentration 0.05 ug/uL in MeOH.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	95.5	mL	12/31/2022

Stock Source	Base Units	Amount Added
VOCF0425	ug/mL	2 mL
VOCF0426	ug/mL	2.5 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF3529B
Spike Name: 2nd Source MtBE
Prep Date: 11/29/2021
Exp Date: 1/29/2022
Department: gcmsvoa
Vendor:
Lot Number:
Balance ID:
Comments: Final Concentration 0.2 ug/uL

Type: Secondary
Prep By: Alethea M. Shaules
Status:
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	9	mL	1/29/2022

Stock Source	Base Units	Amount Added
VOCF0401	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF3546A
Standard Name: Liquids
Type: Secondary
Prep Date: 12/13/2021
Prep By: Alethea M. Shaules
Exp Date: 1/13/2022
Status:
Department: GCMSVOA
Vendor:
Final Volume: 10 mL
Lot Number:
Balance ID:
Comments: 1.0 ml/10 ml final volume. Final Concentration 0.2 ug/uL. Corrected comment and analyte list 11/9/2021 sbd

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EA899	13926	9	mL	1/13/2022

Stock Source	Base Units	Amount Added
VOCF0313	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF3549
Spike Name: 2nd Source Ketones
Prep Date: 12/15/2021
Exp Date: 1/15/2022
Department: gcmsvoa
Vendor: AccuStandard
Lot Number: 221101480
Balance ID:
Comments: Vial opened for use. 2.0 µg/µL

Type: Primary
Prep By: Melissa Chavez
Status:
Final Volume: 1 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketones Mixture	14567	1	mL	1/15/2022

Stock Source	Base Units	Amount Added
VOCF0439	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF3550
Standard Name: Ketones
Type: Primary
Prep Date: 12/16/2021
Prep By: Melissa Chavez
Exp Date: 1/16/2022
Status:
Department: gcmsvoa
Vendor: Chem Service
Final Volume: 1 mL
Lot Number: 10251200
Balance ID:
Comments: Vial Opened For Use . 2.0 ug/uL in 90:10 MeOH:H2O.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketone Mix	14443	1	mL	1/16/2022

Stock Source	Base Units	Amount Added
VOCF0434	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF3558B
Spike Name: 2nd Source Liquids
Prep Date: 12/27/2021
Exp Date: 2/27/2022
Department: gcmsvoa
Vendor:
Lot Number:
Balance ID:
Comments: Final Concentration 0.2ug/uL.

Type: Secondary
Prep By: Steve Dilts
Status: Open
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	9	mL	2/27/2022

Stock Source	Base Units	Amount Added
VOCF0352	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF3559A
Standard Name: MtBE
Type: Secondary
Prep Date: 12/27/2021
Prep By: Melissa Chavez
Exp Date: 1/27/2022
Status:
Department: gcmsvoa
Vendor:
Final Volume: 10 mL
Lot Number:
Balance ID:
Comments: Final Concentration 0.2 ug/uL.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	9	mL	1/27/2022

Stock Source	Base Units	Amount Added
VOCF0373	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF3562A
Standard Name: Gases
Type: Secondary
Prep Date: 1/3/2022
Prep By: Alethea M. Shaules
Exp Date: 1/10/2022
Status:
Department: GCMSVOA
Vendor:
Final Volume: 10 mL
Lot Number:
Balance ID:
Comments: 1.0 ml/10 ml final volume. Final Concentration 0.2 ug/uL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EB373	14519	9	mL	1/10/2022

Stock Source	Base Units	Amount Added
VOCF0427	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Standard ID: VOCF3563
Standard Name: Internals
Type: Secondary
Prep Date: 1/3/2022
Prep By: Alethea M. Shaules
Exp Date: 7/3/2022
Status: New
Department: gcmsvoa
Vendor:
Final Volume: 50 mL
Lot Number:
Balance ID:
Comments: Final Concentration 0.05 ug/uL.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EB373	14519	49	mL	7/3/2022

Stock Source	Base Units	Amount Added
VOCF0425	ug/mL	1 mL

Analytical RunID VOA5975C.I_220104A Standards Traceability Report

Spike ID: VOCF3566A
Spike Name: 2nd Source Gases
Prep Date: 1/4/2022
Exp Date: 1/11/2022
Department: gcmsvoa
Vendor:
Lot Number:
Balance ID:
Comments: 1.0 ml/10 ml final volume. Final Concentration 0.2 ug/uL. Corrected final volume column to match comments and added final concentrations of analytes. MSC 01/14/2021

Type: Secondary
Prep By: Melissa Chavez
Status:
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EB373	14519	9	mL	1/11/2022

Stock Source	Base Units	Amount Added
VOCF0417	ug/mL	1 mL

125 Market Street
New Haven, CT 06513
USA

AccuStandard®

Tel (203)786-5290
Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: M-502A-R-10X
Description: Volatile Organic Compounds - Liquids
Lot: 220041126
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Apr 13, 2020
Expiration: Apr 13, 2023
Sample Size: 1 mL
Components: 54
Storage Condition: Refrig (0-5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
Benzene	71-43-2	100.0	2002	2002
Bromobenzene	108-86-1	100.0	2003	2003
Bromochloromethane	74-97-5	99.1	2001	1983
Bromodichloromethane	75-27-4	99.0	2002	1982
Bromoform	75-25-2	99.2	2001	1985
n-Butylbenzene	104-51-8	100.0	2002	2002
sec-Butylbenzene	135-98-8	100.0	2001	2001
tert-Butylbenzene	98-06-6	99.0	2003	1983
Carbon tetrachloride	56-23-5	100.0	2003	2003
Chlorobenzene	108-90-7	99.6	2001	1993
Chloroform	67-66-3	99.2	2004	1988
2-Chlorotoluene	95-49-8	99.0	2003	1983
4-Chlorotoluene	106-43-4	99.8	2002	1998
Dibromochloromethane	124-48-1	97.8	2049*	2004
1,2-Dibromo-3-chloropropane	96-12-8	99.2	2001	1985
1,2-Dibromoethane	106-93-4	100.0	2006	2006
Dibromomethane	74-95-3	99.0	2002	1982
1,2-Dichlorobenzene	95-50-1	98.2	2003	1967
1,3-Dichlorobenzene	541-73-1	100.0	2000	2000
1,4-Dichlorobenzene	106-46-7	100.0	2002	2002
1,1-Dichloroethane	75-34-3	98.6	2001	1973
1,2-Dichloroethane	107-06-2	99.8	2010	2006
1,1-Dichloroethene	75-35-4	99.0	2000	1980
cis-1,2-Dichloroethene	156-59-2	99.0	2002	1982
trans-1,2-Dichloroethene	156-60-5	99.5	2001	1991
1,2-Dichloropropane	78-87-5	99.5	2003	1993
1,3-Dichloropropane	142-28-9	96.7	2073*	2005
2,2-Dichloropropane	594-20-7	99.9	2012	2010
1,1-Dichloropropene	563-58-6	98.9	2001	1979
cis-1,3-Dichloropropene **	10061-01-5	93.9	2041*	1916
trans-1,3-Dichloropropene **	10061-02-6	93.9	1968*	1848
Ethylbenzene	100-41-4	99.7	2000	1994
Hexachlorobutadiene	87-68-3	98.0	2003	1963
Isopropylbenzene	98-82-8	100.0	2002	2002
p-Isopropyltoluene	99-87-6	99.4	2000	1988
Methylene chloride	75-09-2	99.9	2001	1999
Naphthalene	91-20-3	100.0	2002	2002
n-Propylbenzene	103-65-1	100.0	2001	2001
Styrene	100-42-5	100.0	2003	2003
1,1,1,2-Tetrachloroethane	630-20-6	98.9	2005	1983
1,1,2,2-Tetrachloroethane	79-34-5	96.0	2087*	2004
Tetrachloroethene	127-18-4	99.4	2017	2005
Toluene	108-88-3	100.0	2001	2001
1,2,3-Trichlorobenzene	87-61-6	100.0	2002	2002

For use in routine laboratory analysis.

AccuStandard is accredited to ISO 17034, ISO/IEC 17025 and certified to ISO 9001:2015

125 Market Street
New Haven, CT 06513
USA

AccuStandard®

Tel (203)786-5290
Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: M-502A-R-10X
Description: Volatile Organic Compounds - Liquids
Lot: 220041126
Solvent: Methanol

Date Certified: Apr 13, 2020
Expiration: Apr 13, 2023
Sample Size: 1 mL
Components: 54

Component - *continued*	CAS #	Purity % (GC/MS)	Prepared Concentration[2] (µg/mL)	Certified Analyte Concentration[1] (µg/mL)
1,2,4-Trichlorobenzene	120-82-1	99.6	2001	1993
1,1,1-Trichloroethane	71-55-6	100.0	2002	2002
1,1,2-Trichloroethane	79-00-5	98.6	2000	1972
Trichloroethene	79-01-6	100.0	2003	2003
1,2,3-Trichloropropane	96-18-4	97.5	2055*	2004
1,2,4-Trimethylbenzene	95-63-6	98.2	2001	1965
1,3,5-Trimethylbenzene	108-67-8	98.8	2001	1977
o-Xylene	95-47-6	99.0	2000	1980
m-Xylene	108-38-3	99.2	2002	1986
p-Xylene	106-42-3	95.4	2097*	2001

* Weight compensated to 100% purity.

** 47.8% cis isomer, 46.1% trans isomer

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

[2] All weights are traceable through NIST, Test No. 684/289871-17

[1] Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Certified By: Larry Decker, Organic QC Manager

ID #: 12797
Opened:_____
Volatile Organic Compounds - Liquids
Expires: 4/13/2023
Rec'd: 6/23/2020
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

For use in routine laboratory analysis.

AccuStandard is accredited to ISO 17034, ISO/IEC 17025 and certified to ISO 9001:2015

QR-ORG/INO-001
Rev. 5/18

Agilent

ID #: 13292
Opened:
VOC Standard
Expires: 12/31/2023
Rec'd: 11/23/2020
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Certificate of Analysis

Product Name: VOC Standard

Product Number: DWM-589N-1

Lot Number: 0006570990

Lot Issue Date: 17-Nov-2020

Expiration Date: 31-Dec-2023

Description:
This analytical reference material (RM) was manufactured and verified in accordance with an ISO 9001 registered quality system and analyte concentrations were verified by an ISO 17025 accredited laboratory. The concentration and uncertainty value at the 95% confidence level for each analyte, determined gravimetrically, is listed below.

Analyte	CAS#	Analyte Lot	Concentration ± Uncertainty
bromochloromethane	000074-97-5	RM00009	2010 ± 10 µg/mL
bromodichloromethane	000075-27-4	RM12585	2009 ± 10 µg/mL
bromoform	000075-25-2	RM13987	2010 ± 10 µg/mL
carbon tetrachloride	000056-23-5	RM07576	2010 ± 10 µg/mL
chloroform	000067-66-3	RM13988	2009 ± 10 µg/mL
dibromochloromethane	000124-48-1	RM14843	2009 ± 10 µg/mL
dibromomethane	000074-95-3	RM12878	2009 ± 10 µg/mL
methylene chloride	000075-09-2	RM11650	2009 ± 10 µg/mL
1,2-dibromoethane	000106-93-4	RM00018	2010 ± 10 µg/mL
1,1-dichloroethane	000075-34-3	RM16217	2006 ± 10 µg/mL
1,2-dichloroethane	000107-06-2	RM04655	2005 ± 10 µg/mL
1,1-dichloroethene	000075-35-4	RM14486	2010 ± 10 µg/mL
cis-1,2-dichloroethene	000156-59-2	RM15008	2007 ± 10 µg/mL
trans-1,2-dichloroethene	000156-60-5	RM07565	2008 ± 10 µg/mL
1,1,1,2-tetrachloroethane	000630-20-6	RM12632	2005 ± 10 µg/mL
1,1,2,2-tetrachloroethane	000079-34-5	RM02540	2009 ± 10 µg/mL
tetrachloroethene	000127-18-4	RM06491	2008 ± 10 µg/mL

ANAB ACCREDITED ISO17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
Page: 1 of 4
www.agilent.com/quality/
CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert No. AT-1937

Certificate of Analysis

Product Number: DWM-589N-1

Lot Number: 0006570990

1,1,1-trichloroethane	000071-55-6	RM16539	2004 ± 10 µg/mL
1,1,2-trichloroethane	000079-00-5	RM01175	2009 ± 10 µg/mL
trichloroethene	000079-01-6	RM14232	2009 ± 10 µg/mL
1,2-dibromo-3-chloropropane	000096-12-8	RM13666	2009 ± 10 µg/mL
1,2-dichloropropane	000078-87-5	RM12821	2008 ± 10 µg/mL
1,3-dichloropropane	000142-28-9	RM02080	2008 ± 10 µg/mL
2,2-dichloropropane	000594-20-7	RM12927	2005 ± 10 µg/mL
1,1-dichloropropene	000563-58-6	RM16190	2010 ± 10 µg/mL
cis-1,3-dichloropropene	010061-01-5	RM12891	2007 ± 10 µg/mL
trans-1,3-dichloropropene	010061-02-6	RM12254	2006 ± 10 µg/mL
hexachlorobutadiene	000087-68-3	RM09157	2005 ± 10 µg/mL
1,2,3-trichloropropane	000096-18-4	RM13082	2004 ± 10 µg/mL
benzene	000071-43-2	RM12931	2009 ± 10 µg/mL
n-butylbenzene	000104-51-8	RM03651	2008 ± 10 µg/mL
sec-butylbenzene	000135-98-8	RM10905	2005 ± 10 µg/mL
tert-butylbenzene	000098-06-6	RM14040	2007 ± 10 µg/mL
ethylbenzene	000100-41-4	RM12195	2006 ± 10 µg/mL
isopropylbenzene	000098-82-8	RM00835	2009 ± 10 µg/mL
4-isopropyltoluene	000099-87-6	RM09747	2009 ± 10 µg/mL
naphthalene	000091-20-3	NT00970	2006 ± 10 µg/mL
n-propylbenzene	000103-65-1	RM12785	2010 ± 10 µg/mL
styrene	000100-42-5	RM13393	2010 ± 10 µg/mL

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER

ISO 17034 Cert
No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026

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CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY

ISO 17025 Cert
No. AT-1937

Agilent

Certificate of Analysis

Product Number: DWM-589N-1

Lot Number: 0006570990

toluene	000108-88-3	RM06650	2008 ± 10 µg/mL
1,2,4-trimethylbenzene	000095-63-6	RM06731	2002 ± 10 µg/mL
1,3,5-trimethylbenzene	000108-67-8	RM12905	2009 ± 10 µg/mL
o-xylene	000095-47-6	RM15639	2005 ± 10 µg/mL
m-xylene	000108-38-3	RM15919	2006 ± 10 µg/mL
p-xylene	000106-42-3	RM02647	2009 ± 10 µg/mL
bromobenzene	000108-86-1	RM10227	2008 ± 10 µg/mL
chlorobenzene	000108-90-7	RM01874	2008 ± 10 µg/mL
2-chlorotoluene	000095-49-8	RM13774	2007 ± 10 µg/mL
4-chlorotoluene	000106-43-4	RM11750	2009 ± 10 µg/mL
1,2-dichlorobenzene	000095-50-1	RM13636	2005 ± 10 µg/mL
1,3-dichlorobenzene	000541-73-1	NT00356	2009 ± 10 µg/mL
1,4-dichlorobenzene	000106-46-7	RM12826	2009 ± 10 µg/mL
1,2,3-trichlorobenzene	000087-61-6	RM10193	2007 ± 10 µg/mL
1,2,4-trichlorobenzene	000120-82-1	RM09454	2009 ± 10 µg/mL

Matrix: methanol (methyl alcohol)

Storage Conditions: Store Frozen (-25° to -10°C).

Traceability:
The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCSL Z540.3, ISO 9001, ISO 17025, and ISO 17034. Calibrated Class A glassware is used for volumetric measurements. Thermometers are calibrated against a NIST traceable thermometer in accordance with NIST Special Publication 1088.

Homogeneity:
This RM was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

ANAB ACCREDITED ISO17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026

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CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert No. AT-1937

Certificate of Analysis

Product Number: DWM-589N-1

Lot Number: 0006570990

Intended Use:
This RM is intended for the preparation of working reference samples for use in routine laboratory analyses, calibration of instruments, validation of analytical methods, assessments of measurement methods, and continuing calibration verification.

Instructions for Use:
Sample aliquots for analysis should be withdrawn at 20°C to 25°C immediately after opening the container and should be processed without delay for the certified values to be valid within the stated uncertainties.

Hazards:
Refer to the Safety Data Sheet on www.agilent.com for information regarding this RM.

Expiration of Certification:
The certification of this RM is valid until the expiration date specified above, provided the RM is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the RM is damaged, contaminated, or otherwise modified.

Maintenance of Certification:
If substantive changes are noted that affect the certification before the expiration of this certificate, Agilent will notify the purchaser.

Sample lot approver:

Monica Bourgeois
QMS Representative

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert
No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
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CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert
No. AT-1937

125 Market Street
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USA

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CERTIFICATE OF ANALYSIS

Catalog No: M-8260A-B-SS-10X
Description: Surrogate Standard Mix
Lot: 219041458
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Apr 18, 2019
Expiration: Apr 18, 2029
Sample Size: 1 mL
Components: 4
Storage Condition: Ambient (>5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
p-Bromofluorobenzene	460-00-4	99.9	2004	2002
Dibromofluoromethane	1868-53-7	99.8	2005	2001
1,2-Dichloroethane-d4	17060-07-0	100.0	2001	2001
Toluene-d8	2037-26-5	100.0	2000	2000

ID #: 13385
Opened:
Surrogate Standard Mix
Expires: 4/18/2029
Rec'd: 1/4/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

² All weights are traceable through NIST, Test No. 684/289871-17

¹ Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

 See reverse side for additional information

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: Larry Decker, Organic QC Manager

For use in routine laboratory analysis.

AccuStandard is accredited to ISO 17034, ISO/IEC 17025 and certified to ISO 9001:2015

QR-ORG/INO-001
Rev. 7/20

Agilent

ID #: 13578
Opened:
Methyl tert-Butyl Ether Standard
Expires: 8/31/2022
Rec'd: 2/26/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Certificate of Analysis

Product Name: Methyl tert-Butyl Ether Standard

Product Number: STS-440-1

Lot Number: 0006555762

Lot Issue Date: 19-Aug-2020

Expiration Date: 31-Aug-2022

Description:
This analytical reference material (RM) was manufactured and verified in accordance with an ISO 9001 registered quality system, and analyte concentrations were verified by an ISO 17025 accredited laboratory. The concentration and uncertainty value at the 95% confidence level for each analyte, determined gravimetrically, is listed below.

Analyte	CAS#	Analyte Lot	Concentration ± Uncertainty
tert-butylmethyl ether	001634-04-4	RM06568	2006 ± 10 µg/mL

Matrix: methanol (methyl alcohol)

Storage Conditions: Store Frozen (-25° to -10°C).

Traceability:
The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCSL Z540.3, ISO 9001, ISO 17025, and ISO 17034. Calibrated Class A glassware is used for volumetric measurements. Thermometers are calibrated against a NIST traceable thermometer in accordance with NIST Special Publication 1088.

Homogeneity:
This RM was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

Intended Use:
This RM is intended for the preparation of working reference samples for use in routine laboratory analyses, calibration of instruments, validation of analytical methods, assessments of measurement methods, and continuing calibration verification.

Instructions for Use:
Sample aliquots for analysis should be withdrawn at 20°C to 25°C immediately after opening the container and should be processed without delay for the certified values to be valid within the stated uncertainties.

Hazards:
Refer to the Safety Data Sheet on www.agilent.com for information regarding this RM.

Expiration of Certification:
The certification of this RM is valid until the expiration date specified above, provided the RM is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the RM is damaged, contaminated, or otherwise modified.

Maintenance of Certification:
If substantive changes are noted that affect the certification before the expiration of this certificate, Agilent will notify the purchaser.

Sample lot approver:

Monica Bourgeois
QMS Representative

ANAB ACCREDITED ISO17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert
No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
Page: 1 of 1
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CSD-QA-015.1

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ISO 17025 Cert
No. AT-1937

125 Market Street
New Haven, CT 06513
USA

AccuStandard®

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Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: S-078-10X
Description: MtBE
Lot: 220051182
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: May 18, 2020
Expiration: May 18, 2030
Sample Size: 1 mL
Components: 1
Storage Condition: Ambient (>5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration[2] (µg/mL)	Certified Analyte Concentration[1] (µg/mL)
MtBE	1634-04-4	100.0	2002	2002

ID #: 13920
Opened:_____
MTBE
Expires: 5/18/2030
Rec'd: 6/7/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

[2] All weights are traceable through NIST, Test No. 684/289871-17

[1] Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: _____
Larry Decker, Organic QC Manager

For use in routine laboratory analysis.

660 Tower Lane • P.O. Box 599 • West Chester, PA 19381-0599
1-800-452-9994 • 1-610-692-3026 • Fax 1-610-692-8729
info@chemservice.com • www.chemservice.com

CERTIFICATE OF ANALYSIS

Volatile Organics High Concentration Mixture #6

CONCENTRATION	2000ug/ml in Methanol
CATALOG NUMBER	M-VOHC6M5-1ML
LOT NUMBER	11882100
DATE CERTIFIED	05/25/21
EXPIRATION DATE	02/28/22
STORAGE	Store at room temperature (20 - 25 °C).
HANDLING	See Safety Data Sheet
INTENDED USE	For laboratoy use only.
ISO 17034:2016 CERTIFIED	[ ]

ID #: 14142
Opened:
Volatile Organics High Concentration Mixture
Expires: 2/28/2022
Rec'd: 8/3/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

ID	Analyte	CAS	Weight Analyte (mg)	Lot	Purity	Certified Concentration (ug/mL)
N-11446	Chloroethane	75-00-3	96.300	00001728	100.0	2006.3
N-11665	Dichlorodifluoromethane	75-71-8	96.610	00001729	100.0	2012.7
N-12417	Methyl bromide	74-83-9	96.910	00024694	100.0	2019.0
N-12421	Methyl chloride	74-87-3	96.150	00001731	100.0	2003.1
N-13655	Trichlorofluoromethane	75-69-4	96.300	00027239	99.4	1994.2
N-13748	Vinyl chloride	75-01-4	96.150	00019298	100.0	2003.1

Analytical Test	Value
CONCENTRATION (GC/MSD)	VERIFIED

COA Form
Revision 3 (3/2015)

PJLA Testing
Accreditation #63529

Print Date: 07/28/21

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info@chemservice.com • www.chemservice.com

Instructions for Use:
Shake mixture prior to use. If particles are present, sonicate for homogeneity. If sample is diluted to lower concentrations, Class A volumetric glassware must be used.
Minimum Sample Size- 0.2 uL for Direct Injection.
Chem Service Inc. guarantees the expanded uncertainty of the above analytes to be +/- 2.0% of the certified concentrations based on gravimetric preparation. The test results published in this report were obtained using equipment capable of producing results that are traceable to NIST and through NIST to the International System of Units (SI). The reported expanded uncertainty of measurement is stated as the combined standard uncertainty of measurement multiplied by the coverage factor k (k=2) such that the coverage probability corresponds to approximately 95%. For certified reference materials, homogeneity and thermal stability testing are available upon request.

Certified By:

Mary Beth O'Donnell

Mary Beth O'Donnell
CSM/TC

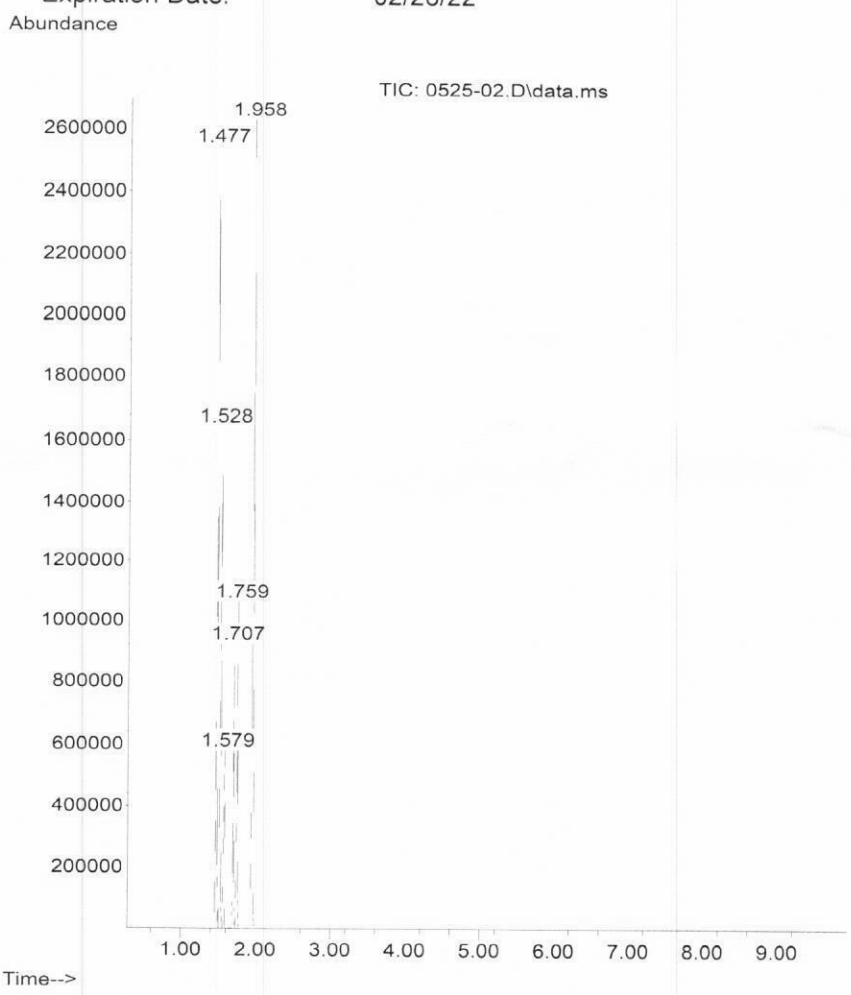
PJLA
Testing
Accreditation #63520

660 Tower Lane • P.O. Box 599 • West Chester, PA 19381-0599
1-800-452-9994 • 1-610-692-3026 • Fax 1-610-692-8729
info@chemservice.com • www.chemservice.com

CERTIFICATE OF ANALYSIS

Analysis Method:

Catalog Number:	M-VOHC6M5-1ML
Description:	Volatile Organics High Concentration Mixture #6
Lot Number:	11882100
Expiration Date:	02/28/22



PJLA Testing

ID #: 14251
Opened:
Internal Standard
Expires: 12/31/2022
Rec'd: 9/8/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Agilent

Certificate of Analysis

Product Name: Internal Standard

Product Number: STM-520-1

Lot Number: 0006582580

Lot Issue Date: 05-Jan-2021

Expiration Date: 31-Dec-2022

Description:
This analytical reference material (RM) was manufactured and verified in accordance with an ISO 9001 registered quality system and analyte concentrations were verified by an ISO 17025 accredited laboratory. The concentration and uncertainty value at the 95% confidence level for each analyte, determined gravimetrically, is listed below.

Analyte	CAS#	Analyte Lot	Concentration ± Uncertainty
chlorobenzene-d5	003114-55-4	RM12274	2501 ± 13 µg/mL
1,4-dichlorobenzene-d4	003855-82-1	RM12517	2501 ± 13 µg/mL
fluorobenzene	000462-06-6	RM13378	2512 ± 13 µg/mL

Matrix: methanol (methyl alcohol)

Storage Conditions: Store Frozen (-25° to -10°C).

Traceability:
The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCSL Z540.3, ISO 9001, ISO 17025, and ISO 17034. Calibrated Class A glassware is used for volumetric measurements. Thermometers are calibrated against a NIST traceable thermometer in accordance with NIST Special Publication 1088.

Homogeneity:
This RM was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

Intended Use:
This RM is intended for the preparation of working reference samples for use in routine laboratory analyses, calibration of instruments, validation of analytical methods, assessments of measurement methods, and continuing calibration verification.

Instructions for Use:
Sample aliquots for analysis should be withdrawn at 20°C to 25°C immediately after opening the container and should be processed without delay for the certified values to be valid within the stated uncertainties.

Hazards:
Refer to the Safety Data Sheet on www.agilent.com for information regarding this RM.

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
Page: 1 of 2

www.agilent.com/quality/
CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert No. AT-1937

Agilent

Certificate of Analysis

Product Number: STM-520-1 **Lot Number:** 0006582580

Expiration of Certification:
The certification of this RM is valid until the expiration date specified above, provided the RM is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the RM is damaged, contaminated, or otherwise modified.

Maintenance of Certification:
If substantive changes are noted that affect the certification before the expiration of this certificate, Agilent will notify the purchaser.

Sample lot approver:

Monica Bourgeois
QMS Representative

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert
No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
Page: 2 of 2
www.agilent.com/quality/
CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert
No. AT-1937

125 Market Street
New Haven, CT 06513
USA

AccuStandard®

Tel (203)786-5290
Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: M-8260A-B-SS-10X
Description: Surrogate Standard Mix
Lot: 219041458
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Apr 18, 2019
Expiration: Apr 18, 2029
Sample Size: 1 mL
Components: 4
Storage Condition: Ambient (>5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
p-Bromofluorobenzene	460-00-4	99.9	2004	2002
Dibromofluoromethane	1868-53-7	99.8	2005	2001
1,2-Dichloroethane-d4	17060-07-0	100.0	2001	2001
Toluene-d8	2037-26-5	100.0	2000	2000

ID #: 14269
Opened:_____
Surrogate Standard Mix
Expires: 4/18/2029
Rec'd: 9/14/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

² All weights are traceable through NIST, Test No. 684/289871-17

¹ Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: _____
Larry Decker, Organic QC Manager

For use in routine laboratory analysis.

Absolute Standards, Inc.
800-368-1131
www.absolutestandards.com

Certified Reference Material CRM

ANAB ISO 17034 Accredited
AR-1539 Certificate Number
https://Absolutestandards.com

CERTIFIED WEIGHT REPORT

Part Number: 30058
Lot Number: 080321
Description: EPA Method 502/524 - Volatile Gases Mix #1
6 components
Expiration Date: 080324
Recommended Storage: Freezer (0 °C)
Nominal Concentration (µg/mL): 2000
NIST Test ID#: 6UTB

Solvent: Methanol **Lot#** EA783-US

5E-05 Balance Uncertainty
Weight(s) shown below were combined and diluted to (mL): 500.0 0.058 Flask Uncertainty

Formulated By:	Mario Luis	080321 DATE
Reviewed By:	Pedro L. Rentas	080321 DATE

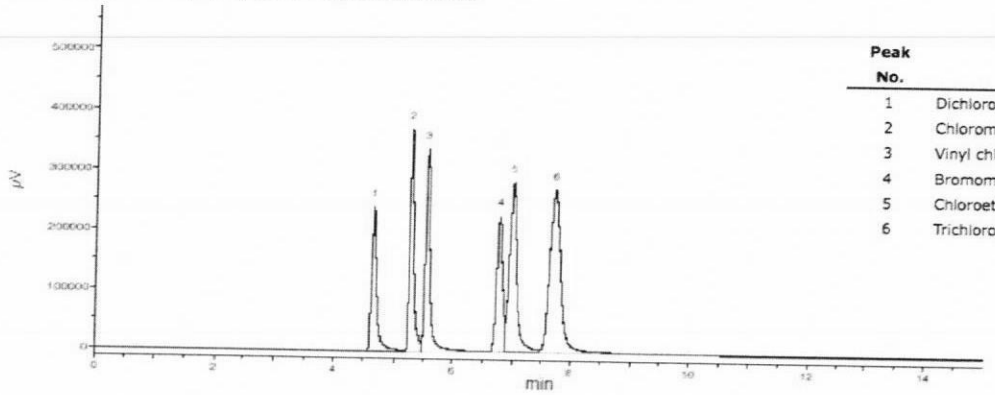
Compound	RM#	Lot Number	Nominal Conc (µg/mL)	Purity (%)	Uncertainty Purity (%)	Target Weight (g)	Actual Weight (g)	Actual Conc(µg/mL)	Expanded Uncertainty (+/-) (µg/mL)	SDS Information (Solvent Safety Info. On Attached pg.) CAS#	OSHA PEL (TWA)	LD50
1. Bromomethane	50	01611JX	2000	99.5	0.2	1.00508	1.0098	2009.4	8.1	74-83-9	5 ppm (20mg/m3/8H) (skin)	orl-rat 214mg/kg
2. Chloroethane	72	062617	2000	99	0.2	1.01016	1.0146	2008.8	8.1	75-00-3	1000 ppm (2600mg/m3/8H)	N/A
3. Chloromethane	79	06908MS	2000	99.5	0.2	1.00508	1.0154	2020.5	8.1	74-87-3	100 ppm	orl-rat 1800mg/kg
4. Dichlorodifluoromethane	134	92-0487	2000	99	0.2	1.01016	1.0224	2024.2	8.2	75-71-8	1000 ppm (4950mg/m3/8H)	N/A
5. Trichlorofluoromethane	294	01823MW	2000	99	0.2	1.01016	1.0110	2001.7	8.1	75-69-4	1000 ppm (5600mg/m3/8H)	ipr-mus 1743mg/kg
6. Vinyl chloride	305	04854EA	2000	99.5	0.2	1.00508	1.0071	2004.0	8.1	75-01-4	N/A	N/A

- The certified value is the concentration calculated from gravimetric and volumetric measurements unless otherwise stated.
- Standards are prepared gravimetrically using balances that are calibrated with weights traceable to NIST (see above).
- Standards are certifed (+/-) 0.5% of the stated value, unless otherwise stated.
- All Standards, after opening ampule, should be stored with caps tight and under appropriate laboratory conditions.
- Uncertainty Reference: Taylor, B.N. and Kuyat, C.E., "Guidelines for Evaluating and Expressing the Uncertainty of NIST Measurement Result," NIST Technical Note 1297, U.S. Government Printing Office, Washington, DC, (1994).

Comments

GC15-M9 Analysis by Melissa Stonier
Column ID SPB-Vocol 105 meter X 0.53mm X 3.0µm film thickness
Flow rates: Total flow=150mL/min., Helium (carrier)=10mL/min.,
Helium(make-up)=40mL/min., Hydogen(make-up)=100mL/min.
Oven Profile: Temp. 1=35°C (Time 1=9 min.), Temp 2=200°C (Time 2=1 min.),
Rate = 33°C/min., Total run time=15 min. Injector temp.=200°C, FID Temp.=200°C.
ELCD Signal = Edaq Channel 1 PID Signal = Edaq Channel 2
Standard injection = 0.5µL, Range=3 Purge Valve = 0 min.

ID #: 14285
Opened: _____
EPA Method 502-524 - Volatile Gases Mix #1
Expires: 8/3/2024
Rec'd: 9/17/2021
Enerav Laboratories Inc 1120 So. 27th Street
Billings MT 59107



Peak No.	Analyte	ELCD RT (min.)
1	Dichlorodifluoromethane	4.67
2	Chloromethane	5.28
3	Vinyl chloride	5.56
4	Bromomethane	6.75
5	Chloroethane	6.99
6	Trichlorofluoromethane	7.72

660 Tower Lane • P.O. Box 599 • West Chester, PA 19381-0599
1-800-452-9994 • 1-610-692-3026 • Fax 1-610-692-8729
info@chemservice.com • www.chemservice.com

ID #: 14443
Opened:
TCL Ketone Mix
Expires: 6/30/2023
Rec'd: 10/26/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

CERTIFICATE OF ANALYSIS

TCL Ketones Mixture

CONCENTRATION	2000ug/ml in Methanol:Water (90:10)
CATALOG NUMBER	M-TCL1AN5-1ML
LOT NUMBER	10251200
DATE CERTIFIED	06/16/20
EXPIRATION DATE	06/30/23
STORAGE	Freezer storage (-20 - -25 °C)
HANDLING	See Safety Data Sheet
INTENDED USE	For laboratory use only.
ISO 17034:2016 CERTIFIED	[ ]

ID	Analyte	CAS	Weight Analyte (mg)	Lot	Purity	Certified Concentration (ug/mL)
N-11014	Acetone	67-64-1	203.300	00026182	98.7	2006.6
N-10297	2-Butanone	78-93-3	202.800	00027454	99.5	2017.9
N-10369	2-Hexanone	591-78-6	202.600	00025720	99.5	2015.9
N-10844	4-Methyl-2-pentanone	108-10-1	204.700	6403300	99.5	2036.8

Analytical Test	Value
CONCENTRATION (GC/FID)	VERIFIED

Chem Service, Inc. is accredited to ISO 17034:2016, ISO/IEC 17025:2017 and certified to ISO 9001:2015.

PJLA Testing

PJLA Reference Material Producer

COA Form
Revision 3 (3/2015)

Print Date: 10/22/21

660 Tower Lane • P.O. Box 599 • West Chester, PA 19381-0599
1-800-452-9994 • 1-610-692-3026 • Fax 1-610-692-8729
info@chemservice.com • www.chemservice.com

Instructions for Use:
Shake mixture prior to use. If particles are present, sonicate for homogeneity. If sample is diluted to lower concentrations, Class A volumetric glassware must be used.
Minimum Sample Size- 0.2 uL for Direct Injection.
Chem Service Inc. guarantees the expanded uncertainty of the above analytes to be +/- 2.0% of the certified concentrations based on gravimetric preparation. The test results published in this report were obtained using equipment capable of producing results that are traceable to NIST and through NIST to the International System of Units (SI). The reported expanded uncertainty of measurement is stated as the combined standard uncertainty of measurement multiplied by the coverage factor k (k=2) such that the coverage probability corresponds to approximately 95%. For certified reference materials, homogeneity and thermal stability testing are available upon request.

Certified By:

Mary Beth O'Donnell

Mary Beth O'Donnell
CSM/TC

Chem Service, Inc. is accredited to ISO 17034:2016, ISO/IEC 17025:2017 and certified to ISO 9001:2015.

PJLA Testing

PJLA Reference Material Producer

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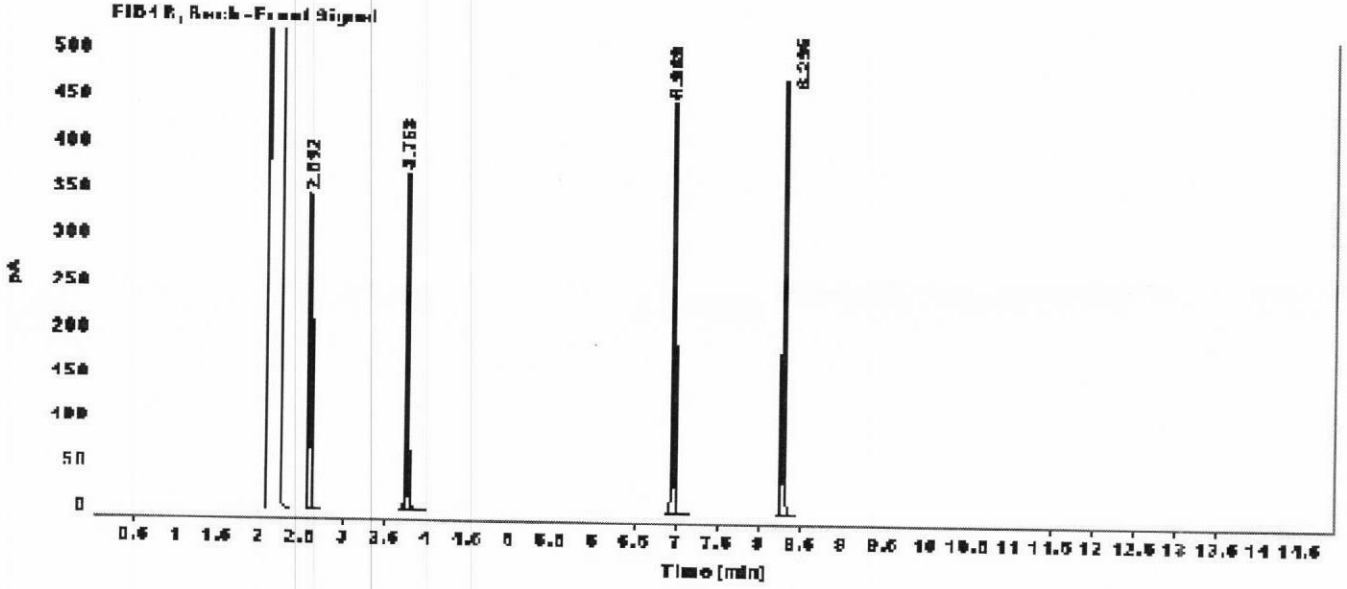
info@chemservice.com • www.chemservice.com

CERTIFICATE OF ANALYSIS

Gas Chromatography / Flame Ionization Detector (GC/FID)

Data file:	C:\CHEM32\1\DATA\2020 DATA\0620\M-TCL1AN5.D		
Sample name:	M-TCL1AN5		
Acq. method:	N-14276.M		
Instrument:	GC3	**Location:**	202
Injection date:	6/16/2020 2:52:35 PM	**Injection Vol:**	1.000
Column name:	RTx-5MS (30m x 0.25mm x 0.5µm)	**# Of Injections:**	1



Signal: FID1 B, Back - Front Signal

RT [min]	Type	Width [min]	Area	Height	Area%
2.592	BB	0.0277	580.2505	343.4966	18.4655
3.763	BB	0.0323	735.4804	367.8491	23.4054
6.989	BB	0.0326	904.3389	447.8770	28.7791
8.295	BB	0.0307	922.2798	474.3798	29.3500
		Sum	3142.3497		

Chem Service, Inc. is accredited to ISO 17034:2016, ISO/IEC 17025:2017 and certified to ISO 9001:2015.

PJLA Testing

PJLA Reference Material Producer

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New Haven, CT 06513
USA

AccuStandard®

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Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: CLP-022K-10X
Description: TCL Ketone Mix
Lot: 221101480
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Oct 26, 2021
Expiration: Nov 26, 2022
Sample Size: 1 mL
Components: 4
Storage Condition: Freeze (<-10 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463

ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
Acetone	67-64-1	100.0	2004	2004
2-Butanone	78-93-3	100.0	2004	2004
2-Hexanone	591-78-6	98.7	2004	1978
4-Methyl-2-pentanone	108-10-1	100.0	2004	2004

ID #: 14567
Opened:_____
TCL Ketones Mixture
Expires: 11/26/2022
Rec'd: 11/30/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

This Certified Reference Material was verified in accordance with ISO/IEC 17025

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

² All weights are traceable through NIST, Test No. 684/289871-17

¹ Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: _____
Larry Decker, Organic QC Manager

For use in routine laboratory analysis.

125 Market Street
New Haven, CT 06513
USA

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Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: CLP-022K-10X
Description: TCL Ketone Mix
Lot: 221101480
Solvent: Methanol

Date Certified: Oct 26, 2021
Expiration: Nov 26, 2022
Sample Size: 1 mL
Components: 4

T-Test

AccuStandard, Inc.
Statistical Report for CLP (SOW 1991)
26-Oct-2021

QR-QCO-003 rev. 1/16

RT	Component	CLP-022K-10X 221101480 Run #1	Run #2	Run #3	Run #4	Mean	Std Dev	% RSD	CLP-022K-10X 221041075 Run #1	Run #2	Run #3	Run #4	Mean	Std Dev	% RSD	NOTES: t.025 test	CI 221101480	Component	CI 221041075	# of Runs	Conc.	10 % error check of means
3.74	Acetone (67-64-1)	1925	1881	1854	1803	1866	51.05	2.74%	1751	1712	1750	1764	1744	22.43	1.29%	4.36	119.2	Acetone (67-64-1)	55	4	2000	7 %
5.77	2-Butanone (78-93-3)	2275	2223	2237	2149	2221	52.79	2.38%	2157	2103	2145	2177	2146	31.26	1.46%	2.46	58.5	2-Butanone (78-93-3)	35.9	4	2000	3 %
8.34	4-Methyl-2-pentanone (108-10-1)	3373	3302	3408	3225	3327	81.05	2.44%	3349	3240	3296	3415	3325	74.70	2.25%	0.04	0.9	4-Methyl-2-pentanone (108	0.8	4	2000	0 %
9.13	2-Hexanone (591-78-6)	3260	3199	3332	3118	3227	90.88	2.82%	3186	3072	3120	3239	3154	73.32	2.32%	1.25	35.2	2-Hexanone (591-78-6)	29.1	4	2000	2 %

For use in routine laboratory analysis.

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New Haven, CT 06513
USA

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CERTIFICATE OF ANALYSIS

Catalog No: CLP-022K-10X
Description: TCL Ketone Mix
Lot: 221101480
Solvent: Methanol

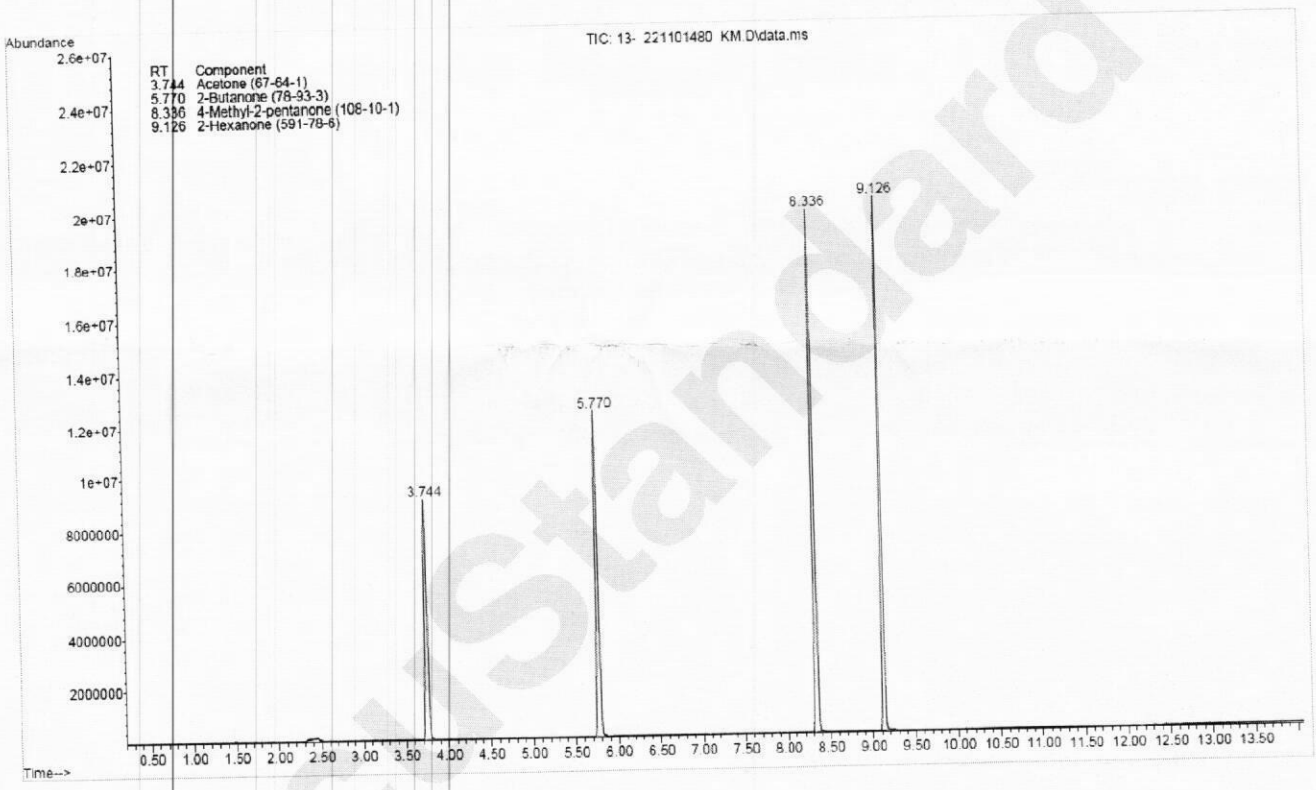
Date Certified: Oct 26, 2021
Expiration: Nov 26, 2022
Sample Size: 1 mL
Components: 4

Chromatogram

```
File       :Q:\GCMS-06 Minimal\DATA\102521\13-  221101480  KM.D
Operator   : Organic QC Lab
Acquired   : 25 Oct 2021  21:00     using AcqMethod VOC-Split100.M
Instrument :   GCMS 6
Sample Name: CLP-022K-10X  (221101480)
Misc Info  : TCL Ketone Mix @ 2000 ug/mL in MeOH
Vial Number: 34
```

AccuStandard®

Column: DB-624 UI, 30m x 0.25mm ID x 1.4µm
Oven Program: 35°C (hold 5min), 11°C/min to 60°C, 22°C/min to 230°C (hold 4min)
GC Parameters: Split 100:1, 1µl inj.; GC/MS; injector temp 240°C



For use in routine laboratory analysis.

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF0313
Standard Name: Liquids
Type: Primary
Prep Date: 6/23/2020
Prep By: Alethea M. Shaules
Exp Date: 4/13/2023
Status: New
Department: gcmsvoa
Vendor: AccuStd
Final Volume: 10 mL
Lot Number: 220041126
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL. Catalog # M502A-R-10X. Corrected lot number to match CI. MSC 01/14/2022

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Volatile Organic Compounds - Liquids	12797	1	mL	4/13/2023

Stock Source	Base Units	Amount Added
VOCF0313	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF0352
Spike Name: 2nd Source Liquids
Prep Date: 11/23/2020
Exp Date: 12/31/2023
Department: gcmsvoa
Vendor: Agilent
Lot Number: 0006570990
Balance ID:
Comments: Date Prepared is same as Date Received. 2000 ug/mL in MeOH. Catalog # DWM-589N-1.

Type: Primary
Prep By: Steve Dilts
Status: New
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
VOC Standard	13292	1	mL	12/31/2023

Stock Source	Base Units	Amount Added
VOCF0352	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF0373
Standard Name: MtBE (Methy tert-Butyl Ether)
Type: Primary
Prep Date: 2/26/2021
Prep By: Steve Dilts
Exp Date: 8/31/2022
Status: New
Department: gcmsvoa
Vendor: Agilent
Final Volume: 5 mL
Lot Number: 0006555762
Balance ID:
Comments: Date Prepared is same as Date Received. 2000 ug/mL in MeOH. Catalog # STS-440

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methyl tert-Butyl Ether Standard	13578	1	mL	8/31/2022

Stock Source	Base Units	Amount Added
VOCF0373	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF0401
Spike Name: 2nd Source MtBE
Type: Primary
Prep Date: 6/7/2021
Prep By: Alethea M. Shaules
Exp Date: 12/11/2029
Status: New
Department: gcmsvoa
Vendor: AccuStandard
Final Volume: 5 mL
Lot Number: 220051182
Balance ID:
Comments: Date Prepared is same as Date Receive. 2,000 ug/mL in MeOH. Catalog # S-078-10X.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
MTBE	13920	1	mL	12/11/2029

Stock Source	Base Units	Amount Added
VOCF0401	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF0417
Spike Name: Chem Service Gases
Prep Date: 8/3/2021
Exp Date: 2/28/2022
Department: gcmsvoa
Vendor: Chemservice
Lot Number: 11882100
Balance ID:
Type: Primary
Prep By: Steve Dilts
Status:
Final Volume: 5 mL
Comments: Date Prepared is same as Date Received. 2000 ug/mL in MeOH. Catalog # M-VOHC6M5-1ML

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Volatile Organics High Concentration Mixture #6	14142	5	mL	2/28/2022

Stock Source	Base Units	Amount Added
VOCF0417	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF0425
Standard Name: Internals
Type: Primary
Prep Date: 9/8/2021
Prep By: Jerran D. Brenden
Exp Date: 12/31/2022
Status: New
Department: gcmsvoa
Vendor: Agilent
Final Volume: 10 mL
Lot Number: 0006582580
Balance ID:
Comments: Date Prepared is same as Date Received. 2,500 ug/mL in MeOH. Catalog # STM-520-1.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Internal Standard	14251	1	mL	12/31/2022

Stock Source	Base Units	Amount Added
VOCF0425	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF0426
Spike Name: Surrogates 2.0 mg/mL
Prep Date: 9/14/2021
Exp Date: 4/18/2029
Department: gcmsvoa
Vendor: AccuStandard
Lot Number: 219041458
Balance ID:
Comments: Date Received 01/04/2021. 2.0 mg/mL. Catalog # M-8260A-B-SS-10X

Type: Primary
Prep By: Jerran D. Brenden
Status: New
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Surrogate Standard Mix	14269	1	mL	4/18/2029

Stock Source	Base Units	Amount Added
VOCF0426	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF0427
Standard Name: Gases
Type: Primary
Prep Date: 9/17/2021
Prep By: Alethea M. Shaules
Exp Date: 8/3/2024
Status: New
Department: gcmsvoa
Vendor: Absolute
Final Volume: 10 mL
Lot Number: 080321
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL in MeOH. Catalog # 30058.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
EPA Method 502-524 - Volatile Gases Mix #1	14285	1	mL	8/3/2024

Stock Source	Base Units	Amount Added
VOCF0427	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF0434
Standard Name: Ketones
Type: Primary
Prep Date: 10/26/2021
Prep By: Steve Dilts
Exp Date: 6/30/2023
Status: New
Department: gcmsvoa
Vendor: Chem Service
Final Volume: 5 mL
Lot Number: 10251200
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL in 90:10 MeOH:H2O. Catalog # M-TCL-1AN5-5ML.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketone Mix	14443	1	mL	6/30/2023

Stock Source	Base Units	Amount Added
VOCF0434	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF0439
Standard Name: 2nd Source Ketones
Type: Primary
Prep Date: 11/30/2021
Prep By: Melissa Chavez
Exp Date: 11/26/2022
Status: New
Department: gcmsvoa
Vendor: AccuStandard
Final Volume: 1 mL
Lot Number: 221101480
Balance ID:
Comments: Date Prepared is same as Date Received. 2,000 ug/mL in Methanol. Catalog # CLP-022K-10X.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketones Mixture	14567	2	mL	11/26/2022

Stock Source	Base Units	Amount Added
VOCF0439	ug/mL	

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF3517
Spike Name: Internal Standard / Surrogates (INT/SURR)
Type: Secondary
Prep Date: 11/10/2021
Prep By: Alethea M. Shaules
Exp Date: 12/31/2022
Status: New
Department: gcmsvoa
Vendor:
Final Volume: 100 mL
Lot Number:
Balance ID:
Comments: Final Concentration 0.05 ug/uL in MeOH.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	95.5	mL	12/31/2022

Stock Source	Base Units	Amount Added
VOCF0425	ug/mL	2 mL
VOCF0426	ug/mL	2.5 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF3529B
Spike Name: 2nd Source MtBE
Prep Date: 11/29/2021
Exp Date: 1/29/2022
Department: gcmsvoa
Vendor:
Lot Number:
Balance ID:
Comments: Final Concentration 0.2 ug/uL

Type: Secondary
Prep By: Alethea M. Shaules
Status:
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	9	mL	1/29/2022

Stock Source	Base Units	Amount Added
VOCF0401	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF3546A
Standard Name: Liquids
Type: Secondary
Prep Date: 12/13/2021
Prep By: Alethea M. Shaules
Exp Date: 1/13/2022
Status:
Department: GCMSVOA
Vendor:
Final Volume: 10 mL
Lot Number:
Balance ID:
Comments: 1.0 ml/10 ml final volume. Final Concentration 0.2 ug/uL. Corrected comment and analyte list 11/9/2021 sbd

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EA899	13926	9	mL	1/13/2022

Stock Source	Base Units	Amount Added
VOCF0313	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF3549
Spike Name: 2nd Source Ketones
Prep Date: 12/15/2021
Exp Date: 1/15/2022
Department: gcmsvoa
Vendor: AccuStandard
Lot Number: 221101480
Balance ID:
Comments: Vial opened for use. 2.0 µg/µL

Type: Primary
Prep By: Melissa Chavez
Status:
Final Volume: 1 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketones Mixture	14567	1	mL	1/15/2022

Stock Source	Base Units	Amount Added
VOCF0439	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF3550
Standard Name: Ketones
Type: Primary
Prep Date: 12/16/2021
Prep By: Melissa Chavez
Exp Date: 1/16/2022
Status:
Department: gcmsvoa
Vendor: Chem Service
Final Volume: 1 mL
Lot Number: 10251200
Balance ID:
Comments: Vial Opened For Use . 2.0 ug/uL in 90:10 MeOH:H2O.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
TCL Ketone Mix	14443	1	mL	1/16/2022

Stock Source	Base Units	Amount Added
VOCF0434	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF3558B
Spike Name: 2nd Source Liquids
Prep Date: 12/27/2021
Exp Date: 2/27/2022
Department: gcmsvoa
Vendor:
Lot Number:
Balance ID:
Comments: Final Concentration 0.2ug/uL.

Type: Secondary
Prep By: Steve Dilts
Status: Open
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	9	mL	2/27/2022

Stock Source	Base Units	Amount Added
VOCF0352	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF3559A
Standard Name: MtBE
Type: Secondary
Prep Date: 12/27/2021
Prep By: Melissa Chavez
Exp Date: 1/27/2022
Status:
Department: gcmsvoa
Vendor:
Final Volume: 10 mL
Lot Number:
Balance ID:
Comments: Final Concentration 0.2 ug/uL.

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap - EB199-US	14334	9	mL	1/27/2022

Stock Source	Base Units	Amount Added
VOCF0373	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Standard ID: VOCF3562A
Standard Name: Gases
Type: Secondary
Prep Date: 1/3/2022
Prep By: Alethea M. Shaules
Exp Date: 1/10/2022
Status:
Department: GCMSVOA
Vendor:
Final Volume: 10 mL
Lot Number:
Balance ID:
Comments: 1.0 ml/10 ml final volume. Final Concentration 0.2 ug/uL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EB373	14519	9	mL	1/10/2022

Stock Source	Base Units	Amount Added
VOCF0427	ug/mL	1 mL

Analytical RunID VOA5975C.I_220107A Standards Traceability Report

Spike ID: VOCF3566A
Spike Name: 2nd Source Gases
Prep Date: 1/4/2022
Exp Date: 1/11/2022
Department: gcmsvoa
Vendor:
Lot Number:
Balance ID:
Comments: 1.0 ml/10 ml final volume. Final Concentration 0.2 ug/uL. Corrected final volume column to match comments and added final concentrations of analytes. MSC 01/14/2021

Type: Secondary
Prep By: Melissa Chavez
Status:
Final Volume: 10 mL

Chemical/Solvent Used	Bottle No	Amt	Units	Expires
Methanol, Purge and Trap EB373	14519	9	mL	1/11/2022

Stock Source	Base Units	Amount Added
VOCF0417	ug/mL	1 mL

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USA

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CERTIFICATE OF ANALYSIS

Catalog No: M-502A-R-10X
Description: Volatile Organic Compounds - Liquids
Lot: 220041126
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Apr 13, 2020
Expiration: Apr 13, 2023
Sample Size: 1 mL
Components: 54
Storage Condition: Refrig (0-5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
Benzene	71-43-2	100.0	2002	2002
Bromobenzene	108-86-1	100.0	2003	2003
Bromochloromethane	74-97-5	99.1	2001	1983
Bromodichloromethane	75-27-4	99.0	2002	1982
Bromoform	75-25-2	99.2	2001	1985
n-Butylbenzene	104-51-8	100.0	2002	2002
sec-Butylbenzene	135-98-8	100.0	2001	2001
tert-Butylbenzene	98-06-6	99.0	2003	1983
Carbon tetrachloride	56-23-5	100.0	2003	2003
Chlorobenzene	108-90-7	99.6	2001	1993
Chloroform	67-66-3	99.2	2004	1988
2-Chlorotoluene	95-49-8	99.0	2003	1983
4-Chlorotoluene	106-43-4	99.8	2002	1998
Dibromochloromethane	124-48-1	97.8	2049*	2004
1,2-Dibromo-3-chloropropane	96-12-8	99.2	2001	1985
1,2-Dibromoethane	106-93-4	100.0	2006	2006
Dibromomethane	74-95-3	99.0	2002	1982
1,2-Dichlorobenzene	95-50-1	98.2	2003	1967
1,3-Dichlorobenzene	541-73-1	100.0	2000	2000
1,4-Dichlorobenzene	106-46-7	100.0	2002	2002
1,1-Dichloroethane	75-34-3	98.6	2001	1973
1,2-Dichloroethane	107-06-2	99.8	2010	2006
1,1-Dichloroethene	75-35-4	99.0	2000	1980
cis-1,2-Dichloroethene	156-59-2	99.0	2002	1982
trans-1,2-Dichloroethene	156-60-5	99.5	2001	1991
1,2-Dichloropropane	78-87-5	99.5	2003	1993
1,3-Dichloropropane	142-28-9	96.7	2073*	2005
2,2-Dichloropropane	594-20-7	99.9	2012	2010
1,1-Dichloropropene	563-58-6	98.9	2001	1979
cis-1,3-Dichloropropene **	10061-01-5	93.9	2041*	1916
trans-1,3-Dichloropropene **	10061-02-6	93.9	1968*	1848
Ethylbenzene	100-41-4	99.7	2000	1994
Hexachlorobutadiene	87-68-3	98.0	2003	1963
Isopropylbenzene	98-82-8	100.0	2002	2002
p-Isopropyltoluene	99-87-6	99.4	2000	1988
Methylene chloride	75-09-2	99.9	2001	1999
Naphthalene	91-20-3	100.0	2002	2002
n-Propylbenzene	103-65-1	100.0	2001	2001
Styrene	100-42-5	100.0	2003	2003
1,1,1,2-Tetrachloroethane	630-20-6	98.9	2005	1983
1,1,2,2-Tetrachloroethane	79-34-5	96.0	2087*	2004
Tetrachloroethene	127-18-4	99.4	2017	2005
Toluene	108-88-3	100.0	2001	2001
1,2,3-Trichlorobenzene	87-61-6	100.0	2002	2002

For use in routine laboratory analysis.

AccuStandard is accredited to ISO 17034, ISO/IEC 17025 and certified to ISO 9001:2015

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www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: M-502A-R-10X
Description: Volatile Organic Compounds - Liquids
Lot: 220041126
Solvent: Methanol

Date Certified: Apr 13, 2020
Expiration: Apr 13, 2023
Sample Size: 1 mL
Components: 54

Component - *continued*	CAS #	Purity % (GC/MS)	Prepared Concentration[2] (µg/mL)	Certified Analyte Concentration[1] (µg/mL)
1,2,4-Trichlorobenzene	120-82-1	99.6	2001	1993
1,1,1-Trichloroethane	71-55-6	100.0	2002	2002
1,1,2-Trichloroethane	79-00-5	98.6	2000	1972
Trichloroethene	79-01-6	100.0	2003	2003
1,2,3-Trichloropropane	96-18-4	97.5	2055*	2004
1,2,4-Trimethylbenzene	95-63-6	98.2	2001	1965
1,3,5-Trimethylbenzene	108-67-8	98.8	2001	1977
o-Xylene	95-47-6	99.0	2000	1980
m-Xylene	108-38-3	99.2	2002	1986
p-Xylene	106-42-3	95.4	2097*	2001

* Weight compensated to 100% purity.

** 47.8% cis isomer, 46.1% trans isomer

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

[2] All weights are traceable through NIST, Test No. 684/289871-17

[1] Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Certified By: Larry Decker, Organic QC Manager

ID #: 12797
Opened:_____
Volatile Organic Compounds - Liquids
Expires: 4/13/2023
Rec'd: 6/23/2020
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

For use in routine laboratory analysis.

AccuStandard is accredited to ISO 17034, ISO/IEC 17025 and certified to ISO 9001:2015

QR-ORG/INO-001
Rev. 5/18

Agilent

ID #: 13292
Opened:
VOC Standard
Expires: 12/31/2023
Rec'd: 11/23/2020
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Certificate of Analysis

Product Name: VOC Standard

Product Number: DWM-589N-1

Lot Number: 0006570990

Lot Issue Date: 17-Nov-2020

Expiration Date: 31-Dec-2023

Description:
This analytical reference material (RM) was manufactured and verified in accordance with an ISO 9001 registered quality system and analyte concentrations were verified by an ISO 17025 accredited laboratory. The concentration and uncertainty value at the 95% confidence level for each analyte, determined gravimetrically, is listed below.

Analyte	CAS#	Analyte Lot	Concentration ± Uncertainty
bromochloromethane	000074-97-5	RM00009	2010 ± 10 µg/mL
bromodichloromethane	000075-27-4	RM12585	2009 ± 10 µg/mL
bromoform	000075-25-2	RM13987	2010 ± 10 µg/mL
carbon tetrachloride	000056-23-5	RM07576	2010 ± 10 µg/mL
chloroform	000067-66-3	RM13988	2009 ± 10 µg/mL
dibromochloromethane	000124-48-1	RM14843	2009 ± 10 µg/mL
dibromomethane	000074-95-3	RM12878	2009 ± 10 µg/mL
methylene chloride	000075-09-2	RM11650	2009 ± 10 µg/mL
1,2-dibromoethane	000106-93-4	RM00018	2010 ± 10 µg/mL
1,1-dichloroethane	000075-34-3	RM16217	2006 ± 10 µg/mL
1,2-dichloroethane	000107-06-2	RM04655	2005 ± 10 µg/mL
1,1-dichloroethene	000075-35-4	RM14486	2010 ± 10 µg/mL
cis-1,2-dichloroethene	000156-59-2	RM15008	2007 ± 10 µg/mL
trans-1,2-dichloroethene	000156-60-5	RM07565	2008 ± 10 µg/mL
1,1,1,2-tetrachloroethane	000630-20-6	RM12632	2005 ± 10 µg/mL
1,1,2,2-tetrachloroethane	000079-34-5	RM02540	2009 ± 10 µg/mL
tetrachloroethene	000127-18-4	RM06491	2008 ± 10 µg/mL

ANAB ACCREDITED ISO17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
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www.agilent.com/quality/
CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert No. AT-1937

Certificate of Analysis

Product Number: DWM-589N-1

Lot Number: 0006570990

1,1,1-trichloroethane	000071-55-6	RM16539	2004 ± 10 µg/mL
1,1,2-trichloroethane	000079-00-5	RM01175	2009 ± 10 µg/mL
trichloroethene	000079-01-6	RM14232	2009 ± 10 µg/mL
1,2-dibromo-3-chloropropane	000096-12-8	RM13666	2009 ± 10 µg/mL
1,2-dichloropropane	000078-87-5	RM12821	2008 ± 10 µg/mL
1,3-dichloropropane	000142-28-9	RM02080	2008 ± 10 µg/mL
2,2-dichloropropane	000594-20-7	RM12927	2005 ± 10 µg/mL
1,1-dichloropropene	000563-58-6	RM16190	2010 ± 10 µg/mL
cis-1,3-dichloropropene	010061-01-5	RM12891	2007 ± 10 µg/mL
trans-1,3-dichloropropene	010061-02-6	RM12254	2006 ± 10 µg/mL
hexachlorobutadiene	000087-68-3	RM09157	2005 ± 10 µg/mL
1,2,3-trichloropropane	000096-18-4	RM13082	2004 ± 10 µg/mL
benzene	000071-43-2	RM12931	2009 ± 10 µg/mL
n-butylbenzene	000104-51-8	RM03651	2008 ± 10 µg/mL
sec-butylbenzene	000135-98-8	RM10905	2005 ± 10 µg/mL
tert-butylbenzene	000098-06-6	RM14040	2007 ± 10 µg/mL
ethylbenzene	000100-41-4	RM12195	2006 ± 10 µg/mL
isopropylbenzene	000098-82-8	RM00835	2009 ± 10 µg/mL
4-isopropyltoluene	000099-87-6	RM09747	2009 ± 10 µg/mL
naphthalene	000091-20-3	NT00970	2006 ± 10 µg/mL
n-propylbenzene	000103-65-1	RM12785	2010 ± 10 µg/mL
styrene	000100-42-5	RM13393	2010 ± 10 µg/mL

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER

ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026

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CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY

ISO 17025 Cert No. AT-1937

Agilent

Certificate of Analysis

Product Number: DWM-589N-1

Lot Number: 0006570990

toluene	000108-88-3	RM06650	2008 ± 10 µg/mL
1,2,4-trimethylbenzene	000095-63-6	RM06731	2002 ± 10 µg/mL
1,3,5-trimethylbenzene	000108-67-8	RM12905	2009 ± 10 µg/mL
o-xylene	000095-47-6	RM15639	2005 ± 10 µg/mL
m-xylene	000108-38-3	RM15919	2006 ± 10 µg/mL
p-xylene	000106-42-3	RM02647	2009 ± 10 µg/mL
bromobenzene	000108-86-1	RM10227	2008 ± 10 µg/mL
chlorobenzene	000108-90-7	RM01874	2008 ± 10 µg/mL
2-chlorotoluene	000095-49-8	RM13774	2007 ± 10 µg/mL
4-chlorotoluene	000106-43-4	RM11750	2009 ± 10 µg/mL
1,2-dichlorobenzene	000095-50-1	RM13636	2005 ± 10 µg/mL
1,3-dichlorobenzene	000541-73-1	NT00356	2009 ± 10 µg/mL
1,4-dichlorobenzene	000106-46-7	RM12826	2009 ± 10 µg/mL
1,2,3-trichlorobenzene	000087-61-6	RM10193	2007 ± 10 µg/mL
1,2,4-trichlorobenzene	000120-82-1	RM09454	2009 ± 10 µg/mL

Matrix: methanol (methyl alcohol)

Storage Conditions: Store Frozen (-25° to -10°C).

Traceability:
The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCSL Z540.3, ISO 9001, ISO 17025, and ISO 17034. Calibrated Class A glassware is used for volumetric measurements. Thermometers are calibrated against a NIST traceable thermometer in accordance with NIST Special Publication 1088.

Homogeneity:
This RM was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026

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CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert No. AT-1937

Certificate of Analysis

Product Number: DWM-589N-1

Lot Number: 0006570990

Intended Use:
This RM is intended for the preparation of working reference samples for use in routine laboratory analyses, calibration of instruments, validation of analytical methods, assessments of measurement methods, and continuing calibration verification.

Instructions for Use:
Sample aliquots for analysis should be withdrawn at 20°C to 25°C immediately after opening the container and should be processed without delay for the certified values to be valid within the stated uncertainties.

Hazards:
Refer to the Safety Data Sheet on www.agilent.com for information regarding this RM.

Expiration of Certification:
The certification of this RM is valid until the expiration date specified above, provided the RM is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the RM is damaged, contaminated, or otherwise modified.

Maintenance of Certification:
If substantive changes are noted that affect the certification before the expiration of this certificate, Agilent will notify the purchaser.

Sample lot approver:

Monica Bourgeois
QMS Representative

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert
No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
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CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert
No. AT-1937

Agilent

ID #: 13578
Opened:
Methyl tert-Butyl Ether Standard
Expires: 8/31/2022
Rec'd: 2/26/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Certificate of Analysis

Product Name: Methyl tert-Butyl Ether Standard

Product Number: STS-440-1

Lot Number: 0006555762

Lot Issue Date: 19-Aug-2020

Expiration Date: 31-Aug-2022

Description:
This analytical reference material (RM) was manufactured and verified in accordance with an ISO 9001 registered quality system, and analyte concentrations were verified by an ISO 17025 accredited laboratory. The concentration and uncertainty value at the 95% confidence level for each analyte, determined gravimetrically, is listed below.

Analyte	CAS#	Analyte Lot	Concentration ± Uncertainty
tert-butylmethyl ether	001634-04-4	RM06568	2006 ± 10 µg/mL

Matrix: methanol (methyl alcohol)

Storage Conditions: Store Frozen (-25° to -10°C).

Traceability:
The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCSL Z540.3, ISO 9001, ISO 17025, and ISO 17034. Calibrated Class A glassware is used for volumetric measurements. Thermometers are calibrated against a NIST traceable thermometer in accordance with NIST Special Publication 1088.

Homogeneity:
This RM was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

Intended Use:
This RM is intended for the preparation of working reference samples for use in routine laboratory analyses, calibration of instruments, validation of analytical methods, assessments of measurement methods, and continuing calibration verification.

Instructions for Use:
Sample aliquots for analysis should be withdrawn at 20°C to 25°C immediately after opening the container and should be processed without delay for the certified values to be valid within the stated uncertainties.

Hazards:
Refer to the Safety Data Sheet on www.agilent.com for information regarding this RM.

Expiration of Certification:
The certification of this RM is valid until the expiration date specified above, provided the RM is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the RM is damaged, contaminated, or otherwise modified.

Maintenance of Certification:
If substantive changes are noted that affect the certification before the expiration of this certificate, Agilent will notify the purchaser.

Sample lot approver:

Monica Bourgeois
QMS Representative

ANAB ACCREDITED ISO17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
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CSD-QA-015.1

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ISO 17025 Cert No. AT-1937

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CERTIFICATE OF ANALYSIS

Catalog No: S-078-10X
Description: MtBE
Lot: 220051182
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: May 18, 2020
Expiration: May 18, 2030
Sample Size: 1 mL
Components: 1
Storage Condition: Ambient (>5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration[2] (µg/mL)	Certified Analyte Concentration[1] (µg/mL)
MtBE	1634-04-4	100.0	2002	2002

ID #: 13920
Opened:_____
MTBE
Expires: 5/18/2030
Rec'd: 6/7/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

[2] All weights are traceable through NIST, Test No. 684/289871-17

[1] Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: _____
Larry Decker, Organic QC Manager

Page 1 of 1

For use in routine laboratory analysis.

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info@chemservice.com • www.chemservice.com

CERTIFICATE OF ANALYSIS

Volatile Organics High Concentration Mixture #6

CONCENTRATION	2000ug/ml in Methanol
CATALOG NUMBER	M-VOHC6M5-1ML
LOT NUMBER	11882100
DATE CERTIFIED	05/25/21
EXPIRATION DATE	02/28/22
STORAGE	Store at room temperature (20 - 25 °C).
HANDLING	See Safety Data Sheet
INTENDED USE	For laboratoy use only.
ISO 17034:2016 CERTIFIED	[ ]

ID #: 14142
Opened:
Volatile Organics High Concentration Mixture
Expires: 2/28/2022
Rec'd: 8/3/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

ID	Analyte	CAS	Weight Analyte (mg)	Lot	Purity	Certified Concentration (ug/mL)
N-11446	Chloroethane	75-00-3	96.300	00001728	100.0	2006.3
N-11665	Dichlorodifluoromethane	75-71-8	96.610	00001729	100.0	2012.7
N-12417	Methyl bromide	74-83-9	96.910	00024694	100.0	2019.0
N-12421	Methyl chloride	74-87-3	96.150	00001731	100.0	2003.1
N-13655	Trichlorofluoromethane	75-69-4	96.300	00027239	99.4	1994.2
N-13748	Vinyl chloride	75-01-4	96.150	00019298	100.0	2003.1

Analytical Test	Value
CONCENTRATION (GC/MSD)	VERIFIED

COA Form
Revision 3 (3/2015)

PJLA Testing
Accreditation #63529

Print Date: 07/28/21

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Instructions for Use:
Shake mixture prior to use. If particles are present, sonicate for homogeneity. If sample is diluted to lower concentrations, Class A volumetric glassware must be used.
Minimum Sample Size- 0.2 uL for Direct Injection.
Chem Service Inc. guarantees the expanded uncertainty of the above analytes to be +/- 2.0% of the certified concentrations based on gravimetric preparation. The test results published in this report were obtained using equipment capable of producing results that are traceable to NIST and through NIST to the International System of Units (SI). The reported expanded uncertainty of measurement is stated as the combined standard uncertainty of measurement multiplied by the coverage factor k (k=2) such that the coverage probability corresponds to approximately 95%. For certified reference materials, homogeneity and thermal stability testing are available upon request.

Certified By:

Mary Beth O'Donnell

Mary Beth O'Donnell
CSM/TC

PJLA
Testing
Accreditation #63520

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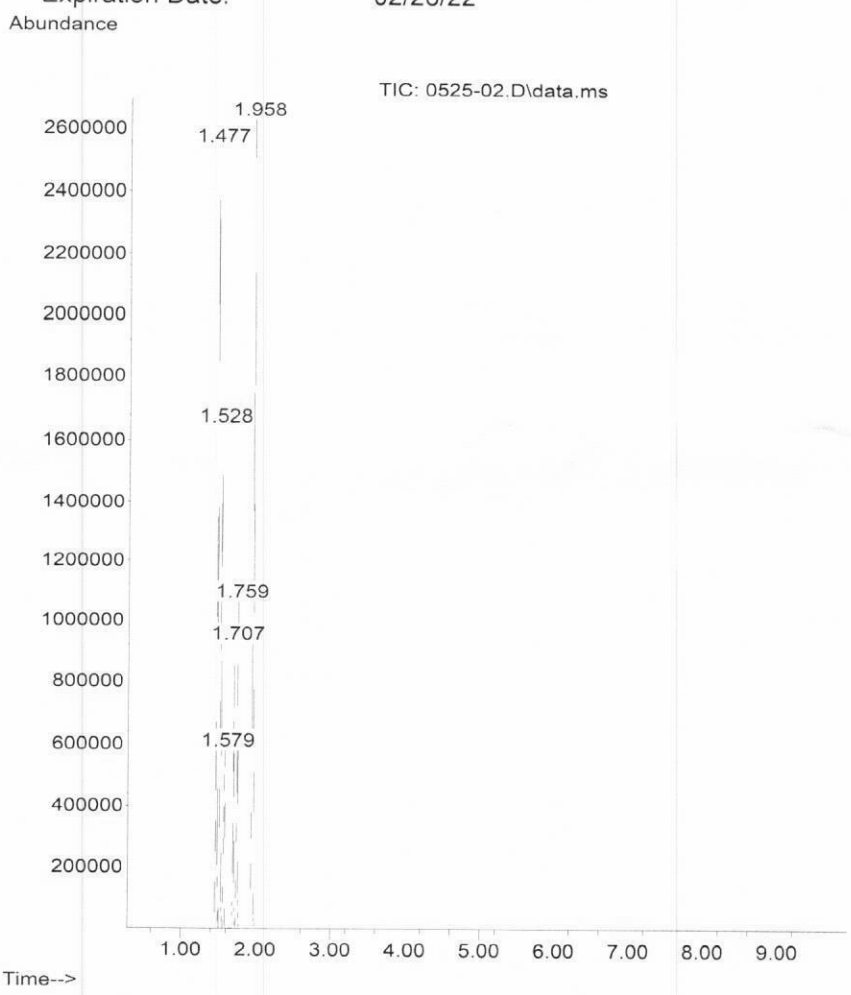
CERTIFICATE OF ANALYSIS

Analysis Method:

Catalog Number:	M-VOHC6M5-1ML
Description:	Volatile Organics High Concentration Mixture #6
Lot Number:	11882100
Expiration Date:	02/28/22



PJLA Testing

ID #: 14251
Opened:
Internal Standard
Expires: 12/31/2022
Rec'd: 9/8/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Agilent

Certificate of Analysis

Product Name: Internal Standard

Product Number: STM-520-1

Lot Number: 0006582580

Lot Issue Date: 05-Jan-2021

Expiration Date: 31-Dec-2022

Description:
This analytical reference material (RM) was manufactured and verified in accordance with an ISO 9001 registered quality system and analyte concentrations were verified by an ISO 17025 accredited laboratory. The concentration and uncertainty value at the 95% confidence level for each analyte, determined gravimetrically, is listed below.

Analyte	CAS#	Analyte Lot	Concentration ± Uncertainty
chlorobenzene-d5	003114-55-4	RM12274	2501 ± 13 µg/mL
1,4-dichlorobenzene-d4	003855-82-1	RM12517	2501 ± 13 µg/mL
fluorobenzene	000462-06-6	RM13378	2512 ± 13 µg/mL

Matrix: methanol (methyl alcohol)

Storage Conditions: Store Frozen (-25° to -10°C).

Traceability:
The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCSL Z540.3, ISO 9001, ISO 17025, and ISO 17034. Calibrated Class A glassware is used for volumetric measurements. Thermometers are calibrated against a NIST traceable thermometer in accordance with NIST Special Publication 1088.

Homogeneity:
This RM was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

Intended Use:
This RM is intended for the preparation of working reference samples for use in routine laboratory analyses, calibration of instruments, validation of analytical methods, assessments of measurement methods, and continuing calibration verification.

Instructions for Use:
Sample aliquots for analysis should be withdrawn at 20°C to 25°C immediately after opening the container and should be processed without delay for the certified values to be valid within the stated uncertainties.

Hazards:
Refer to the Safety Data Sheet on www.agilent.com for information regarding this RM.

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
Page: 1 of 2

www.agilent.com/quality/
CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert No. AT-1937

Agilent

Certificate of Analysis

Product Number: STM-520-1 **Lot Number:** 0006582580

Expiration of Certification:
The certification of this RM is valid until the expiration date specified above, provided the RM is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the RM is damaged, contaminated, or otherwise modified.

Maintenance of Certification:
If substantive changes are noted that affect the certification before the expiration of this certificate, Agilent will notify the purchaser.

Sample lot approver:

Monica Bourgeois
QMS Representative

ANAB ACCREDITED ISO 17034 REFERENCE MATERIAL PRODUCER
ISO 17034 Cert
No. AR-1936

RM was produced in accordance with TUV USA Inc registered ISO 9001 Quality Management System. Cert # 56 100 18560026
Page: 2 of 2
www.agilent.com/quality/
CSD-QA-015.1

ANAB ACCREDITED ISO/IEC 17025 TESTING LABORATORY
ISO 17025 Cert
No. AT-1937

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CERTIFICATE OF ANALYSIS

Catalog No: M-8260A-B-SS-10X
Description: Surrogate Standard Mix
Lot: 219041458
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Apr 18, 2019
Expiration: Apr 18, 2029
Sample Size: 1 mL
Components: 4
Storage Condition: Ambient (>5 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463
ANAB ACCREDITED TESTING LABORATORY AT-1339
ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
p-Bromofluorobenzene	460-00-4	99.9	2004	2002
Dibromofluoromethane	1868-53-7	99.8	2005	2001
1,2-Dichloroethane-d4	17060-07-0	100.0	2001	2001
Toluene-d8	2037-26-5	100.0	2000	2000

ID #: 14269
Opened:_____
Surrogate Standard Mix
Expires: 4/18/2029
Rec'd: 9/14/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

² All weights are traceable through NIST, Test No. 684/289871-17

¹ Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: _____
Larry Decker, Organic QC Manager

Page 1 of 1

For use in routine laboratory analysis.

Absolute Standards, Inc.
800-368-1131
www.absolutestandards.com

Certified Reference Material CRM

ANAB ISO 17034 Accredited
AR-1539 Certificate Number
https://Absolutestandards.com

CERTIFIED WEIGHT REPORT

Part Number: 30058
Lot Number: 080321
Description: EPA Method 502/524 - Volatile Gases Mix #1
6 components
Expiration Date: 080324
Recommended Storage: Freezer (0 °C)
Nominal Concentration (µg/mL): 2000
NIST Test ID#: 6UTB

Solvent: Methanol **Lot#** EA783-US

5E-05 Balance Uncertainty
Weight(s) shown below were combined and diluted to (mL): 500.0 0.058 Flask Uncertainty

Formulated By:	Mario Luis	080321 DATE
Reviewed By:	Pedro L. Rentas	080321 DATE

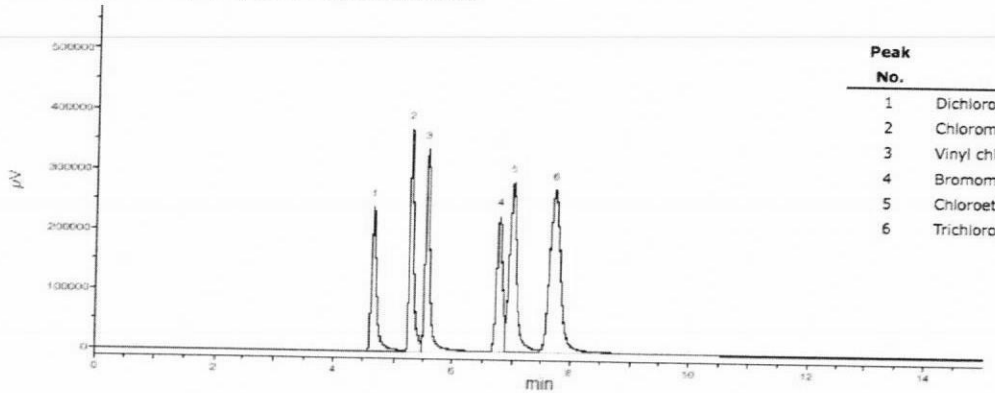
Compound	RM#	Lot Number	Nominal Conc (µg/mL)	Purity (%)	Uncertainty Purity (%)	Target Weight (g)	Actual Weight (g)	Actual Conc(µg/mL)	Expanded Uncertainty (+/-) (µg/mL)	SDS Information (Solvent Safety Info. On Attached pg.) CAS#	OSHA PEL (TWA)	LD50
1. Bromomethane	50	01611JX	2000	99.5	0.2	1.00508	1.0098	2009.4	8.1	74-83-9	5 ppm (20mg/m3/8H) (skin)	orl-rat 214mg/kg
2. Chloroethane	72	062617	2000	99	0.2	1.01016	1.0146	2008.8	8.1	75-00-3	1000 ppm (2600mg/m3/8H)	N/A
3. Chloromethane	79	06908MS	2000	99.5	0.2	1.00508	1.0154	2020.5	8.1	74-87-3	100 ppm	orl-rat 1800mg/kg
4. Dichlorodifluoromethane	134	92-0487	2000	99	0.2	1.01016	1.0224	2024.2	8.2	75-71-8	1000 ppm (4950mg/m3/8H)	N/A
5. Trichlorofluoromethane	294	01823MW	2000	99	0.2	1.01016	1.0110	2001.7	8.1	75-69-4	1000 ppm (5600mg/m3/8H)	ipr-mus 1743mg/kg
6. Vinyl chloride	305	04854EA	2000	99.5	0.2	1.00508	1.0071	2004.0	8.1	75-01-4	N/A	N/A

- The certified value is the concentration calculated from gravimetric and volumetric measurements unless otherwise stated.
- Standards are prepared gravimetrically using balances that are calibrated with weights traceable to NIST (see above).
- Standards are certified (+/-) 0.5% of the stated value, unless otherwise stated.
- All Standards, after opening ampule, should be stored with caps tight and under appropriate laboratory conditions.
- Uncertainty Reference: Taylor, B.N. and Kuyat, C.E., "Guidelines for Evaluating and Expressing the Uncertainty of NIST Measurement Result," NIST Technical Note 1297, U.S. Government Printing Office, Washington, DC, (1994).

Comments

GC15-M9 Analysis by Melissa Stonier
Column ID SPB-Vocol 105 meter X 0.53mm X 3.0µm film thickness
Flow rates: Total flow=150mL/min., Helium (carrier)=10mL/min.,
Helium(make-up)=40mL/min., Hydogen(make-up)=100mL/min.
Oven Profile: Temp. 1=35°C (Time 1=9 min.), Temp 2=200°C (Time 2=1 min.),
Rate = 33°C/min., Total run time=15 min. Injector temp.=200°C, FID Temp.=200°C.
ELCD Signal = Edaq Channel 1 PID Signal = Edaq Channel 2
Standard injection = 0.5µL, Range=3 Purge Valve = 0 min.

ID #: 14285
Opened: _____
EPA Method 502-524 - Volatile Gases Mix #1
Expires: 8/3/2024
Rec'd: 9/17/2021
Eneray Laboratories Inc 1120 So. 27th Street
Billings MT 59107



Peak No.	Analyte	ELCD RT (min.)
1	Dichlorodifluoromethane	4.67
2	Chloromethane	5.28
3	Vinyl chloride	5.56
4	Bromomethane	6.75
5	Chloroethane	6.99
6	Trichlorofluoromethane	7.72

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ID #: 14443
Opened:
TCL Ketone Mix
Expires: 6/30/2023
Rec'd: 10/26/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

CERTIFICATE OF ANALYSIS

TCL Ketones Mixture

CONCENTRATION	2000ug/ml in Methanol:Water (90:10)
CATALOG NUMBER	M-TCL1AN5-1ML
LOT NUMBER	10251200
DATE CERTIFIED	06/16/20
EXPIRATION DATE	06/30/23
STORAGE	Freezer storage (-20 - -25 °C)
HANDLING	See Safety Data Sheet
INTENDED USE	For laboratory use only.
ISO 17034:2016 CERTIFIED	[ ]

ID	Analyte	CAS	Weight Analyte (mg)	Lot	Purity	Certified Concentration (ug/mL)
N-11014	Acetone	67-64-1	203.300	00026182	98.7	2006.6
N-10297	2-Butanone	78-93-3	202.800	00027454	99.5	2017.9
N-10369	2-Hexanone	591-78-6	202.600	00025720	99.5	2015.9
N-10844	4-Methyl-2-pentanone	108-10-1	204.700	6403300	99.5	2036.8

Analytical Test	Value
CONCENTRATION (GC/FID)	VERIFIED

Chem Service, Inc. is accredited to ISO 17034:2016, ISO/IEC 17025:2017 and certified to ISO 9001:2015.

PJLA Testing

PJLA Reference Material Producer

COA Form
Revision 3 (3/2015)

Print Date: 10/22/21

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Instructions for Use:
Shake mixture prior to use. If particles are present, sonicate for homogeneity. If sample is diluted to lower concentrations, Class A volumetric glassware must be used.
Minimum Sample Size- 0.2 uL for Direct Injection.
Chem Service Inc. guarantees the expanded uncertainty of the above analytes to be +/- 2.0% of the certified concentrations based on gravimetric preparation. The test results published in this report were obtained using equipment capable of producing results that are traceable to NIST and through NIST to the International System of Units (SI). The reported expanded uncertainty of measurement is stated as the combined standard uncertainty of measurement multiplied by the coverage factor k (k=2) such that the coverage probability corresponds to approximately 95%. For certified reference materials, homogeneity and thermal stability testing are available upon request.

Certified By:

Mary Beth O'Donnell

Mary Beth O'Donnell
CSM/TC

Chem Service, Inc. is accredited to ISO 17034:2016, ISO/IEC 17025:2017 and certified to ISO 9001:2015.

PJLA Testing

PJLA Reference Material Producer

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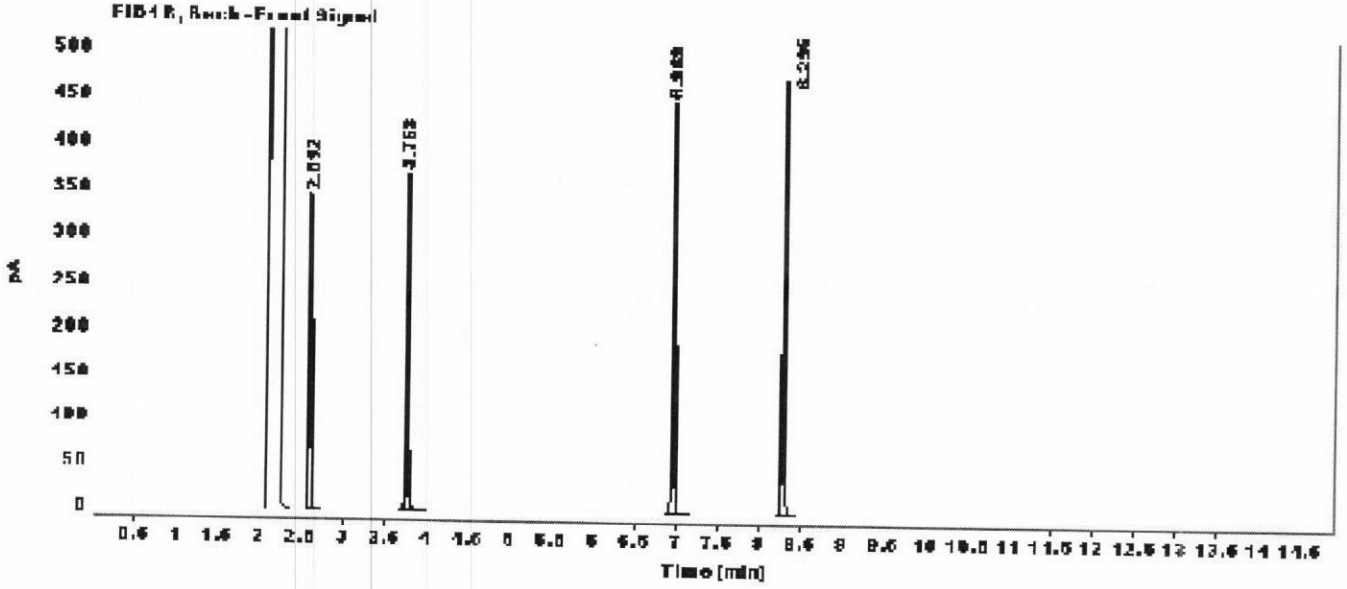
info@chemservice.com • www.chemservice.com

CERTIFICATE OF ANALYSIS

Gas Chromatography / Flame Ionization Detector (GC/FID)

Data file:	C:\CHEM32\1\DATA\2020 DATA\0620\M-TCL1AN5.D		
Sample name:	M-TCL1AN5		
Acq. method:	N-14276.M		
Instrument:	GC3	**Location:**	202
Injection date:	6/16/2020 2:52:35 PM	**Injection Vol:**	1.000
Column name:	RTx-5MS (30m x 0.25mm x 0.5µm)	**# Of Injections:**	1



Signal: FID1 B, Back - Front Signal

RT [min]	Type	Width [min]	Area	Height	Area%
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3.763	BB	0.0323	735.4804	367.8491	23.4054
6.989	BB	0.0326	904.3389	447.8770	28.7791
8.295	BB	0.0307	922.2798	474.3798	29.3500
		Sum	3142.3497		

Chem Service, Inc. is accredited to ISO 17034:2016, ISO/IEC 17025:2017 and certified to ISO 9001:2015.

PJLA Testing

PJLA Reference Material Producer

125 Market Street
New Haven, CT 06513
USA

AccuStandard®

Tel (203)786-5290
Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: CLP-022K-10X
Description: TCL Ketone Mix
Lot: 221101480
Solvent: Methanol
Hazards: *Refer to SDS for complete safety information*

Signal Word: Danger

Date Certified: Oct 26, 2021
Expiration: Nov 26, 2022
Sample Size: 1 mL
Components: 4
Storage Condition: Freeze (<-10 °C)

Certified Reference Material

ANAB ACCREDITED REFERENCE MATERIAL PRODUCER AR-1463

ilac-MRA

Component	CAS #	Purity % (GC/MS)	Prepared Concentration² (µg/mL)	Certified Analyte Concentration¹ (µg/mL)
Acetone	67-64-1	100.0	2004	2004
2-Butanone	78-93-3	100.0	2004	2004
2-Hexanone	591-78-6	98.7	2004	1978
4-Methyl-2-pentanone	108-10-1	100.0	2004	2004

ID #: 14567
Opened:_____
TCL Ketones Mixture
Expires: 11/26/2022
Rec'd: 11/30/2021
Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

This Certified Reference Material was verified in accordance with ISO/IEC 17025

A product with a suffix (-1A, -2B, etc. or -01, -02, etc.) on its lot number has had its expiration date extended and is identical to the same lot number without the suffix.

² All weights are traceable through NIST, Test No. 684/289871-17

¹ Certified Analyte Concentration = Purity x Prepared Concentration.

The Uncertainty associated with the certified concentration reported on this certificate is ±2.4%. This value is the combined expanded uncertainty and represents an estimated standard deviation equal to the positive square root of the total variation of the uncertainty of components. A normal distribution is assumed and a coverage factor of K=2 is chosen using approximately a 95% confidence level.

Labels and certificates follow U.S. Conventions in reporting numerical values: A comma (,) is used to separate units of one-thousand or greater. A period (.) is used as a decimal place marker.

Hazard Information: Please refer to the SDS for information regarding the hazards associated with using this material.

This product was prepared according to in-house procedures and is guaranteed to be homogeneous.

Certified By: _____
Larry Decker, Organic QC Manager

For use in routine laboratory analysis.

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New Haven, CT 06513
USA

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CERTIFICATE OF ANALYSIS

Catalog No: CLP-022K-10X
Description: TCL Ketone Mix
Lot: 221101480
Solvent: Methanol

Date Certified: Oct 26, 2021
Expiration: Nov 26, 2022
Sample Size: 1 mL
Components: 4

T-Test

AccuStandard, Inc.
Statistical Report for CLP (SOW 1991)
26-Oct-2021

QR-QCO-003 rev. 1/16

		CLP-022K-10X 221101480							CLP-022K-10X 221041075							NOTES:						
RT	Component	Run #1	Run #2	Run #3	Run #4	Mean	Std Dev	% RSD	Run #1	Run #2	Run #3	Run #4	Mean	Std Dev	% RSD	t.025 test	CI 221101480	Component	CI 221041075	# of Runs	Conc.	10 % error check of means
3.74	Acetone (67-64-1)	1925	1881	1854	1803	1866	51.05	2.74%	1751	1712	1750	1764	1744	22.43	1.29%	4.36	119.2	Acetone (67-64-1)	55	4	2000	7 %
5.77	2-Butanone (78-93-3)	2275	2223	2237	2149	2221	52.79	2.38%	2157	2103	2145	2177	2146	31.26	1.46%	2.46	58.5	2-Butanone (78-93-3)	35.9	4	2000	3 %
8.34	4-Methyl-2-pentanone (108-10-1)	3373	3302	3408	3225	3327	81.05	2.44%	3349	3240	3296	3415	3325	74.70	2.25%	0.04	0.9	4-Methyl-2-pentanone (108	0.8	4	2000	0 %
9.13	2-Hexanone (591-78-6)	3260	3199	3332	3118	3227	90.88	2.82%	3186	3072	3120	3239	3154	73.32	2.32%	1.25	35.2	2-Hexanone (591-78-6)	29.1	4	2000	2 %

For use in routine laboratory analysis.

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USA

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Fax (203)786-5287
www.AccuStandard.com

CERTIFICATE OF ANALYSIS

Catalog No: CLP-022K-10X
Description: TCL Ketone Mix
Lot: 221101480
Solvent: Methanol

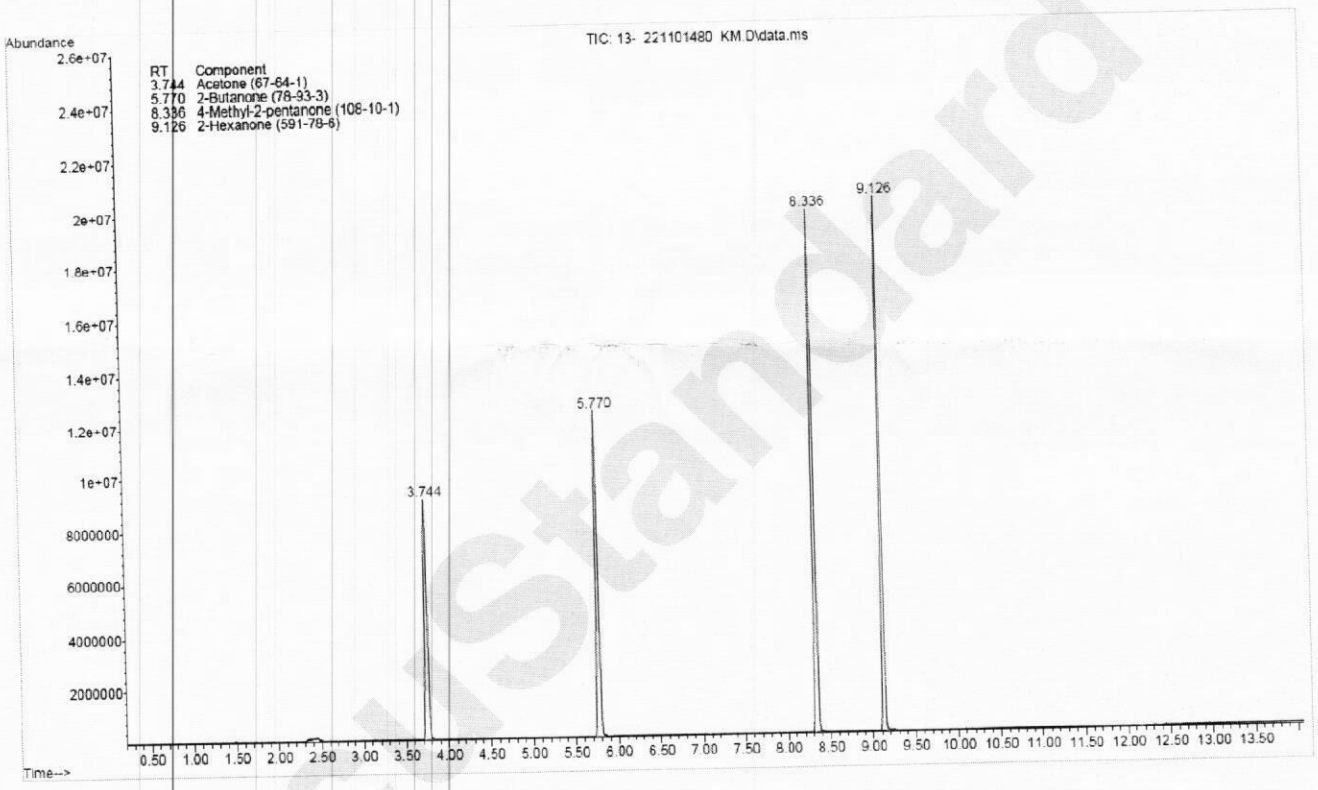
Date Certified: Oct 26, 2021
Expiration: Nov 26, 2022
Sample Size: 1 mL
Components: 4

Chromatogram

```
File       :Q:\GCMS-06 Minimal\DATA\102521\13-  221101480  KM.D
Operator   : Organic QC Lab
Acquired   : 25 Oct 2021  21:00     using AcqMethod VOC-Split100.M
Instrument :   GCMS 6
Sample Name: CLP-022K-10X  (221101480)
Misc Info  : TCL Ketone Mix @ 2000 ug/mL in MeOH
Vial Number: 34
```

AccuStandard®

Column: DB-624 UI, 30m x 0.25mm ID x 1.4µm
Oven Program: 35°C (hold 5min), 11°C/min to 60°C, 22°C/min to 230°C (hold 4min)
GC Parameters: Split 100:1, 1µl inj.; GC/MS; injector temp 240°C



For use in routine laboratory analysis.