

Energy Laboratories Inc

ANALYTICAL RUN Summary

09-Mar-22

Run ID GECD.I_220224A

Run Start Date: 2/24/2022
Analyst: Carry L Tran
Ical:
Column ID: RTX-CLP_0.53
Comments:

Std ID	Std Name	Std Amount	Std Units	Samp Amount	Samp Units	SampType	Expiration Date
SeePrepRecord	Standards Tracked in Prep Batch						2/4/2050

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
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15055418	CAL1-163984	PST-8011-W	CAL1	GECD.IG022422\	2/24/2022 1:48:2	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.01137	0.01134158		0.01	0	0	0.0025835	0.01	0	113%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.01045	0.01042388		0.01	0	0	0.0056259	0.02	0	104%	60	140	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
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15055419	CAL7-163984	PST-8011-W	CAL7	GECD.IG022422\	2/24/2022 2:07:5	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.02198	0.02192505		0.02	0	0	0.0025835	0.01	0	110%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.02002	0.01996995		0.02	0	0	0.0056259	0.02	0	100%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist
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15055420	CAL2-163984	PST-8011-W	CAL2	GECD.IG022422\	2/24/2022 2:27:5	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.05472	0.0545832		0.05	0	0	0.0025835	0.01	0	109%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.04879	0.04866803		0.05	0	0	0.0056259	0.02	0	97%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055421	CAL3-163984	PST-8011-W	CAL3	JECD.IMG022422\	2/24/2022 2:47:3	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.10632	0.1060542		0.1	0	0	0.0025835	0.01	0	106%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09788	0.0976353		0.1	0	0	0.0056259	0.02	0	98%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055422	CAL4-163984	PST-8011-W	CAL4	JECD.IMG022422\	2/24/2022 3:07:2	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.1969	0.19640775		0.2	0	0	0.0025835	0.01	0	98%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.1953	0.19481175		0.2	0	0	0.0056259	0.02	0	97%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055423	CAL5-163984	PST-8011-W	CAL5	JECD.IMG022422\	2/24/2022 3:27:0	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.39856	0.3975636		0.4	0	0	0.0025835	0.01	0	99%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.41011	0.40908473		0.4	0	0	0.0056259	0.02	0	102%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055424	CAL6-163984	PST-8011-W	CAL6	JECD.IMG022422\	2/24/2022 3:46:4	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	1.00036	0.9978591		1	0	0	0.0025835	0.01	0	100%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.99739	0.99489653		1	0	0	0.0056259	0.02	0	99%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055425	LCS-163984	PST-8011-W	ICV	JECD.IMG022422\	2/24/2022 4:26:2	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.2423	0.24169425		0.25	0	0	0.0025835	0.01	0	97%	80	120	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10367	0.10341083		0.1	0	0	0.0056259	0.02	0	103%	80	120	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055426	CAL3-163984	PST-8011-W	CCV3	JECD.IG022422\	2/24/2022 4:46:1	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.10246	0.10220385		0.1	0	0	0.0025835	0.01	0	102%	80	120	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09414	0.09390465		0.1	0	0	0.0056259	0.02	0	94%	80	120	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055427	MB-163984	PST-8011-W	MBLK	JECD.IG022422\	2/24/2022 5:06:0	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.005	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.0972	0.096957		0.1	0	0	0.0056259	0.02	0	97%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055428	LCS-163984	PST-8011-W	LCS-DOD	JECD.IG022422\	2/24/2022 5:25:3	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.24299	0.24238253		0.25	0	0	0.0025835	0.01	0	97%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.096	0.09576		0.1	0	0	0.0056259	0.02	0	96%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055429	LCS1-163984	PST-8011-W	LCS1	JECD.IG022422\	2/24/2022 5:45:2	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.10707	0.10680233		0.1	0	0	0.0025835	0.01	0	107%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10224	0.1019844		0.1	0	0	0.0056259	0.02	0	102%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055430	B22010745-008	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 6:24:5	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.01	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.1264	0.126084		0.1	0	0	0.0056259	0.02	0	126%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055431	B22021435-004	PST-8011-W	SAMP	JECD.I\G022422\	2/24/2022 6:44:4	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09381	0.0919338		0.099	0	0	0.0055272	0.02	0	93%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055432	B22021435-006	PST-8011-W	SAMP	JECD.I\G022422\	2/24/2022 7:04:3	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.11585	0.11556038		0.099	0	0	0.0056259	0.02	0	117%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055433	B22021435-010	PST-8011-W	SAMP	JECD.I\G022422\	2/24/2022 7:24:1	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09304	0.0911792		0.098	0	0	0.0055272	0.02	0	93%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055434	B22021435-012	PST-8011-W	SAMP	JECD.I\G022422\	2/24/2022 7:44:0	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09693	0.0949914		0.098	0	0	0.0055272	0.02	0	97%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055435	B22021435-015	PST-8011-W	SAMP	JECD.I\G022422\	2/24/2022 8:03:3	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09912	0.0971376		0.098	0	0	0.0055272	0.02	0	99%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055436	B22021435-017	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 8:23:3	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10481	0.1027138		0.098	0	0	0.0055272	0.02	0	105%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055437	B22021435-020	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 8:43:2	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.0962	0.0959595		0.099	0	0	0.0056259	0.02	0	97%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055438	B22021435-001	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 9:02:5	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10096	0.0989408		0.099	0	0	0.0055272	0.02	0	100%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055439	B22021435-001	PST-8011-W	MS	JECD.IG022422\	2/24/2022 9:22:5	1	163984	2/24/2022 8:	2E+07	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.24168	0.2368464		0.245	0	0	0.0025382	0.01	0	97%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09651	0.0945798		0.098	0	0	0.0055272	0.02	0	97%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055440	B22021435-001	PST-8011-W	MSD	JECD.IG022422\	2/24/2022 9:42:3	1	163984	2/24/2022 8:	2E+07	2E+07						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.2376	0.232848		0.245	0	0.2368464	0.0025382	0.01	0	95%	60	140	2%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09885	0.096873		0.098	0	0	0.0055272	0.02	0	99%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055441	CAL5-163984	PST-8011-W	CCV4	JECD.IG022422\	2/24/2022 10:22:	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.4143	0.41326425		0.4	0	0	0.0025835	0.01	0	103%	80	120	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.42315	0.42209213		0.4	0	0	0.0056259	0.02	0	106%	80	120	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055442	B22021435-022	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 11:01:	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09977	0.0977746		0.098	0	0	0.0055272	0.02	0	100%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055443	B22021435-025	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 11:21:	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09831	0.0963438		0.098	0	0	0.0055272	0.02	0	98%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055444	B22021435-027	PST-8011-W	SAMP	JECD.IG022422\	2/24/2022 11:41:	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.11788	0.1155224		0.098	0	0	0.0055272	0.02	0	118%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055445	B22021435-030	PST-8011-W	SAMP	JECD.IG022422\	2/25/2022 12:00:	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10728	0.1070118		0.099	0	0	0.0056259	0.02	0	108%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055446	B22021528-001	PST-8011-W	SAMP	JECD.ING022422\	2/25/2022 1:00:1	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.15314	0.1500772		0.099	0	0	0.0055272	0.02	0	152%	70	130	0%	S
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055447	B22021528-004	PST-8011-W	SAMP	JECD.ING022422\	2/25/2022 1:59:2	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.11382	0.1115436		0.098	0	0	0.0055272	0.02	0	114%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15055448	CAL3-163984	PST-8011-W	CCV3	JECD.ING022422\	2/25/2022 2:38:5	1	163984	2/24/2022 7:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.11882	0.11852295		0.1	0	0	0.0025835	0.01	0	119%	80	120	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10355	0.10329113		0.1	0	0	0.0056259	0.02	0	103%	80	120	0%	

Write Sequence

Insert Entries(Have the first cell for e

Data File

Sample Name

G:\org\GECD.i\G022422.b\G0224_001	8011Primer ;0.2ug/L\$PST-8011-W,C4
G:\org\GECD.i\G022422.b\G0224_002	8011Primer ;0.2ug/L\$PST-8011-W,C4
G:\org\GECD.i\G022422.b\G0224_003	8011Primer ;0.1ug/L\$PST-8011-W,C3
G:\org\GECD.i\G022422.b\G0224_004	8011Primer ;0.1ug/L\$PST-8011-W,C3
G:\org\GECD.i\G022422.b\G0224_005	8011Primer ;0.2ug/L\$PST-8011-W,C4
G:\org\GECD.i\G022422.b\G0224_006	Hexane ;
G:\org\GECD.i\G022422.b\G0224_007	CAL1-163984 ;
G:\org\GECD.i\G022422.b\G0224_008	CAL7-163984 ;
G:\org\GECD.i\G022422.b\G0224_009	CAL2-163984 ;
G:\org\GECD.i\G022422.b\G0224_010	CAL3-163984 ;
G:\org\GECD.i\G022422.b\G0224_011	CAL4-163984 ;
G:\org\GECD.i\G022422.b\G0224_012	CAL5-163984 ;
G:\org\GECD.i\G022422.b\G0224_013	CAL6-163984 ;
G:\org\GECD.i\G022422.b\G0224_014	Hexane;;
G:\org\GECD.i\G022422.b\G0224_015	LCS-163984 ;
G:\org\GECD.i\G022422.b\G0224_016	CAL3-163984 ;
G:\org\GECD.i\G022422.b\G0224_017	MB-163984 ;
G:\org\GECD.i\G022422.b\G0224_018	LCS-163984 ;
G:\org\GECD.i\G022422.b\G0224_019	LCS1-163984 ;
G:\org\GECD.i\G022422.b\G0224_020	Hexane;;
G:\org\GECD.i\G022422.b\G0224_021	B22010745-008A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_022	B22021435-004A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_023	B22021435-006H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_024	B22021435-010A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_025	B22021435-012H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_026	B22021435-015A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_027	B22021435-017H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_028	B22021435-020A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_029	B22021435-001H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_030	B22021435-001HMS ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_031	B22021435-001HMSD ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_032	Hexane;;
G:\org\GECD.i\G022422.b\G0224_033	CAL5-163984 ;
G:\org\GECD.i\G022422.b\G0224_034	Hexane;;
G:\org\GECD.i\G022422.b\G0224_035	B22021435-022H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_036	B22021435-025A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_037	B22021435-027H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_038	B22021435-030A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_039	B22021435-032H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_040	B22021435-035A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_041	B22021528-001C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_042	B22021528-002C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_043	B22021528-003C ;\$PST-8011-W,

G:\org\GECD.i\G022422.b\G0224_044	B22021528-004C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_045	Hexane;;
G:\org\GECD.i\G022422.b\G0224_046	CAL3-163984 ;
G:\org\GECD.i\G022422.b\G0224_047	8011Primer ;0.2ug/L\$PST-8011-W,C4
G:\org\GECD.i\G022422.b\G0224_048	8011Primer ;0.2ug/L\$PST-8011-W,C4
G:\org\GECD.i\G022422.b\G0224_049	8011Primer ;0.1ug/L\$PST-8011-W,C3
G:\org\GECD.i\G022422.b\G0224_050	8011Primer ;0.1ug/L\$PST-8011-W,C3
G:\org\GECD.i\G022422.b\G0224_051	8011Primer ;0.2ug/L\$PST-8011-W,C4
G:\org\GECD.i\G022422.b\G0224_052	Hexane ;
G:\org\GECD.i\G022422.b\G0224_053	CAL3-163984 ;
G:\org\GECD.i\G022422.b\G0224_054	Hexane ;
G:\org\GECD.i\G022422.b\G0224_055	B22021435-032H ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_056	B22021435-035A ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_057	B22021528-001C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_058	B22021528-002C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_059	B22021528-003C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_060	B22021528-004C ;\$PST-8011-W,
G:\org\GECD.i\G022422.b\G0224_061	Hexane ;
G:\org\GECD.i\G022422.b\G0224_062	CAL5-163984 ;
G:\org\GECD.i\G022422.b\G0224_063	
G:\org\GECD.i\G022422.b\G0224_064	
G:\org\GECD.i\G022422.b\G0224_065	
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G:\org\GECD.i\G022422.b\G0224_087	
G:\org\GECD.i\G022422.b\G0224_088	
G:\org\GECD.i\G022422.b\G0224_089	

Quantitative Analysis Results Summary Report

Batch Path	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin	Analyst Name	BL2000\ctran
Analysis Time	2/25/2022 2:31 PM	Reporter Name	BL2000\srcocx
Report Time	3/9/2022 2:59:37 PM	Batch State	Processed
Last Calib Update	2/25/2022 8:11 AM	Quant Report Version	10.0
Quant Batch Version	10.0		

Sequence Table

Data File	sample Name	Sample Type	Vial Position	Inj Vol	Level	Acq Method File
G0224_007.0007.D	CAL1-163984	CC		0	1	testAcqFileNamePath
G0224_008.0008.D	CAL7-163984	CC		0	7	testAcqFileNamePath
G0224_009.0009.D	CAL2-163984	CC		0	2	testAcqFileNamePath
G0224_010.0010.D	CAL3-163984	CC		0	3	testAcqFileNamePath
G0224_011.0011.D	CAL4-163984	CC		0	4	testAcqFileNamePath
G0224_012.0012.D	CAL5-163984	CC		0	5	testAcqFileNamePath
G0224_013.0013.D	CAL6-163984	CC		0	6	testAcqFileNamePath
G0224_015.0015.D	LCS-163984	QC		0	LCS	testAcqFileNamePath
G0224_017.0017.D	MB-163984	MethodBlank		0		testAcqFileNamePath

Quantitation Results

Compound: 1,2-Dibromoethane

Data File	Sample Type	RT	Resp	Final Conc	Exp. Conc	Accuracy
G0224_007.0007.D	CC	3.453	1929	0.0114	0.0100	113.7
G0224_008.0008.D	CC	3.453	3724	0.0220	0.0200	109.9
G0224_009.0009.D	CC	3.453	9221	0.0547	0.0500	109.4
G0224_010.0010.D	CC	3.453	17768	0.1063	0.1000	106.3
G0224_011.0011.D	CC	3.451	32418	0.1969	0.2000	98.4
G0224_012.0012.D	CC	3.450	63436	0.3986	0.4000	99.6
G0224_013.0013.D	CC	3.449	142853	1.0004	1.0000	100.0
G0224_015.0015.D	QC	3.449	39595	0.2423	0.2500	96.9
G0224_017.0017.D	Blank	3.442	0	ND		

Compound: 1,1,1,2-Tetrachloroethane

Data File	Sample Type	RT	Resp	Final Conc	Exp. Conc	Accuracy
G0224_007.0007.D	CC	3.978	1386	0.0105	0.0100	104.5
G0224_008.0008.D	CC	3.976	4766	0.0200	0.0200	100.1
G0224_009.0009.D	CC	3.975	14968	0.0488	0.0500	97.6
G0224_010.0010.D	CC	3.974	32499	0.0979	0.1000	97.9
G0224_011.0011.D	CC	3.973	67779	0.1953	0.2000	97.6
G0224_012.0012.D	CC	3.973	147845	0.4101	0.4000	102.5
G0224_013.0013.D	CC	3.971	382731	0.9974	1.0000	99.7
G0224_015.0015.D	QC	3.973	34577	0.1037	0.1000	103.7
G0224_017.0017.D	Blank	3.970	32257	0.0972		

Initial Calibration Report - WJB



Method Path \\MASSHUNTER\Org\Data\GECD.I\GECD_methods
 Method File G022422_8011_W_CLT.m
 Batch Name \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT_batch.bin
 Last Calib Update 2/25/2022 8:11:20 AM

Level Name	Calibration Files	Acq. Date-Time	Level Last Update Time
1	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_007.0007.D	2/24/2022 1:48:23 PM	2/25/2022 8:11:20 AM
7	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_008.0008.D	2/24/2022 2:07:59 PM	2/25/2022 8:11:20 AM
2	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_009.0009.D	2/24/2022 2:27:56 PM	2/25/2022 8:11:20 AM
3	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_010.0010.D	2/24/2022 2:47:36 PM	2/25/2022 8:11:20 AM
4	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_011.0011.D	2/24/2022 3:07:26 PM	2/25/2022 8:11:20 AM
5	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_012.0012.D	2/24/2022 3:27:03 PM	2/25/2022 8:11:20 AM
6	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_013.0013.D	2/24/2022 3:46:49 PM	2/25/2022 8:11:20 AM

Compound	Curve Fit	1	7	2	3	4	5	6	Avg RF	%RSD
M 1,2-Dibromoethane	Quadratic	193088	186099	180413	177678	162092	158591	142853	171545	10.298
S 1,1,1,2-Tetrachloroethane	Quadratic	138646	238314	299357	324992	338896	369613	382731	298936	28.536

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

Initial Calibration Report - WJB



Compounds with Curve fitting not using Avg Response Factor:

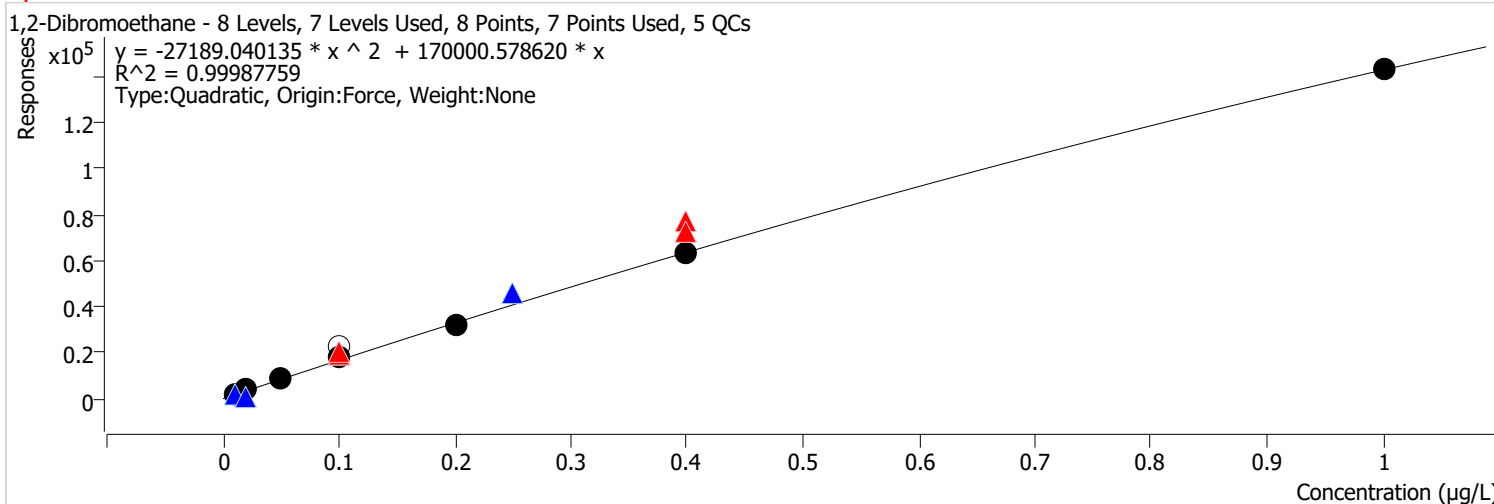
Compound	Curve Fit	Curve Fit Formula	Curve Fit R2
M 1,2-Dibromoethane	Quadratic	$y = -27189.040135 * x^2 + 170000.578620 * x$	0.999878
S 1,1,1,2-Tetrachloroethane	Quadratic	$y = 33932.728826 * x^2 + 352192.061697 * x - 2298.119678$	0.999702

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

Calibration Report

Batch Path	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin	Analyst Name	BL2000\ctran
Analysis Time	2/25/2022 2:31 PM	Reporter Name	BL2000\srcox
Report Time	3/9/2022 3:14:43 PM	Batch State	Processed
Last Calib Update	2/25/2022 8:11 AM	Quant Report Version	10.0
Quant Batch Version	10.0		

1,2-Dibromoethane %RSE =

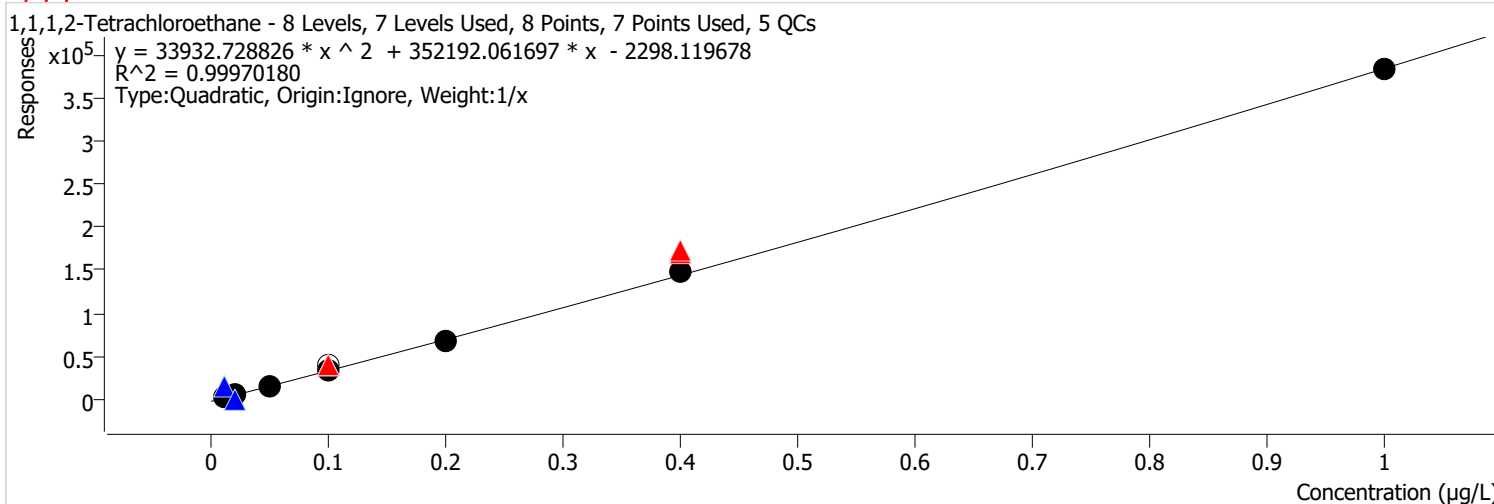


Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_007.0007.D	Calibration	1	x	1931	0.0100	193088.4 351	
D:\Org\Data\GECD.I\G091321\aiexport\G0913_018.0018.D	QC	7	x	1335	0.0200	66739.74 25	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_008.0008.D	Calibration	7	x	3722	0.0200	186099.2 670	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_009.0009.D	Calibration	2	x	9021	0.0500	180413.4 200	
\\MASSHUNTER\Org\Data\GECD.I\G111820\aiexport\G1118_016.0016.D	Calibration	CC3		22970	0.1000		
D:\Org\Data\GECD.I\G081021\aiexport\G0810_016.0016.D	QC	CC3		21004	0.1000	210042.4 247	
D:\Org\Data\GECD.I\G122121\aiexport\G1221_061.0061.D	CC	CC3		19101	0.1000	191007.5 606	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_053.0053.D	CC	3	x	19066	0.1000	190664.4 764	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_030.0030.D	QC	LCS1	x	19195	0.1000	191949.8 990	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_027.0027.D	CC	3	x	20905	0.1000	209045.9 250	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_010.0010.D	Calibration	3	x	17768	0.1000	177677.8 324	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_011.0011.D	Calibration	4	x	32418	0.2000	162091.7 968	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_029.0029.D	QC	LCS	x	46496	0.2500	185984.9 936	
D:\Org\Data\GECD.I\G122121\aiexport\G1221_074.0074.D	CC	CC5	x	77330	0.4000	193324.5 351	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_044.0044.D	CC	5	x	72089	0.4000	180222.6 846	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_012.0012.D	Calibration	5	x	63436	0.4000	158590.6 926	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_013.0013.D	Calibration	6	x	142853	1.0000	142853.4 754	

Calibration Report

Batch Path	\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin	Analyst Name	BL2000\ctran
Analysis Time	2/25/2022 2:31 PM	Reporter Name	BL2000\srcox
Report Time	3/9/2022 3:14:47 PM	Batch State	Processed
Last Calib Update	2/25/2022 8:11 AM	Quant Report Version	10.0
Quant Batch Version	10.0		

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

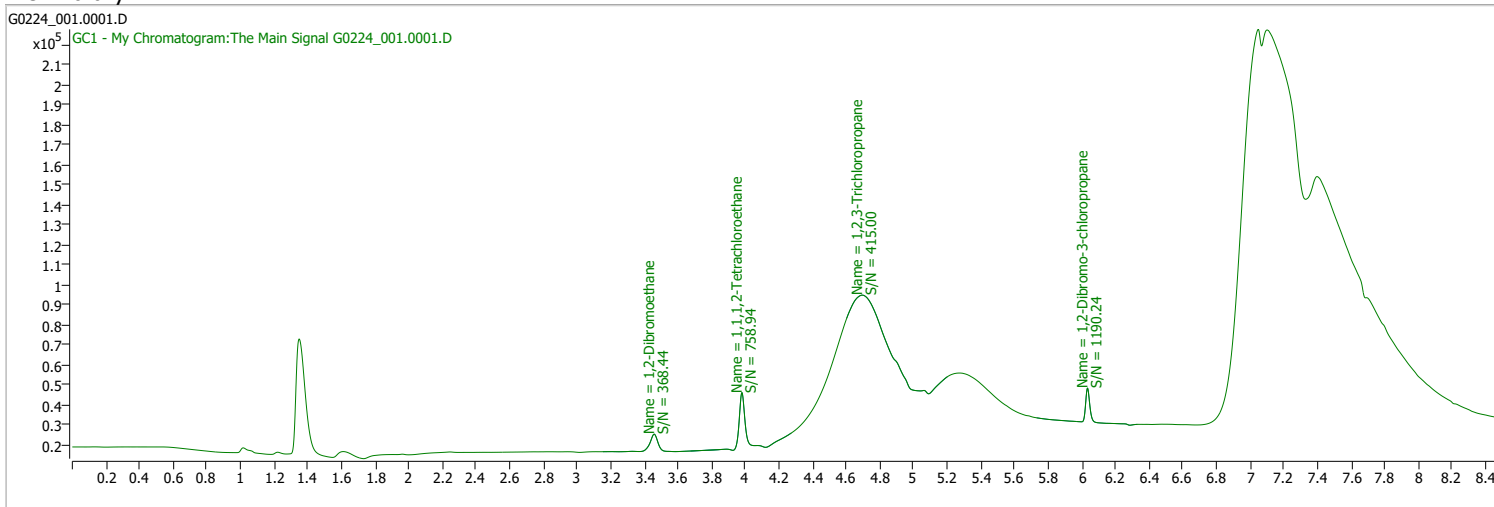


Calibration STD Path	Cal Type	Level	Enabled	Resp.	Exp. Conc	Resp. Factor	Level RSD
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\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_007.0007.D	Calibration	1	x	1386	0.0100	138645.5392	
D:\Org\Data\GECD.I\G091321\aiexport\G0913_018.0018.D	QC	7	x	686	0.0200	34275.7771	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_008.0008.D	Calibration	7	x	4766	0.0200	238314.4042	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_009.0009.D	Calibration	2	x	14968	0.0500	299356.9329	
\\MASSHUNTER\Org\Data\GECD.I\G111820\aiexport\G1118_016.0016.D	Calibration	CC3		41065	0.1000		
D:\Org\Data\GECD.I\G081021\aiexport\G0810_016.0016.D	QC	CC3		42481	0.1000	424813.5788	
D:\Org\Data\GECD.I\G122121\aiexport\G1221_061.0061.D	CC	CC3		29228	0.1000	292276.2189	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_053.0053.D	CC	3	x	38726	0.1000	387264.2087	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_030.0030.D	QC	LCS1	x	41151	0.1000	411505.5261	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_029.0029.D	QC	LCS	x	40669	0.1000	406691.2911	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_027.0027.D	CC	3	x	40786	0.1000	407855.1467	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_010.0010.D	Calibration	3	x	32499	0.1000	324992.1899	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_011.0011.D	Calibration	4	x	67779	0.2000	338895.8219	
D:\Org\Data\GECD.I\G122121\aiexport\G1221_074.0074.D	CC	CC5	x	169695	0.4000	424236.9956	
\\MASSHUNTER\Org\Data\GECD.I\G020922\aiexport\G0209_044.0044.D	CC	5	x	172872	0.4000	432180.2832	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_012.0012.D	Calibration	5	x	147845	0.4000	369612.9718	
\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_013.0013.D	Calibration	6	x	382731	1.0000	382731.2699	

Quantitation Results Report (QT Reviewed)

Data File	G0224_001.0001.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 11:49:39 AM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.979	0.0	59458	0.1725	µg/L	0.005
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 172.48%	*	

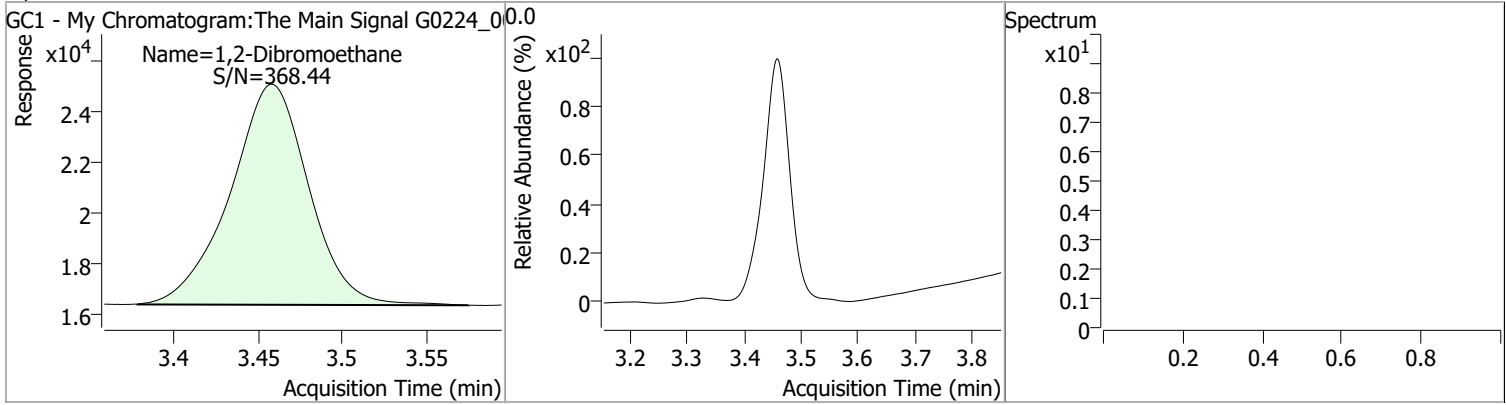
Target Compounds

M 1,2-Dibromoethane	3.458	0.0	28196	0.1705	µg/L	QValue 100
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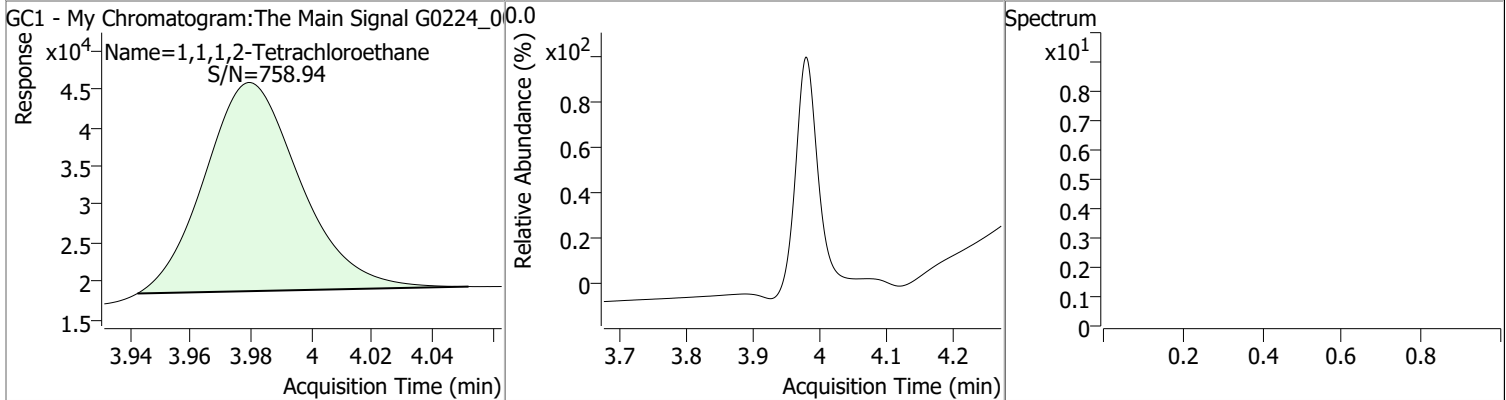
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1705	3.46	0.01	28196				



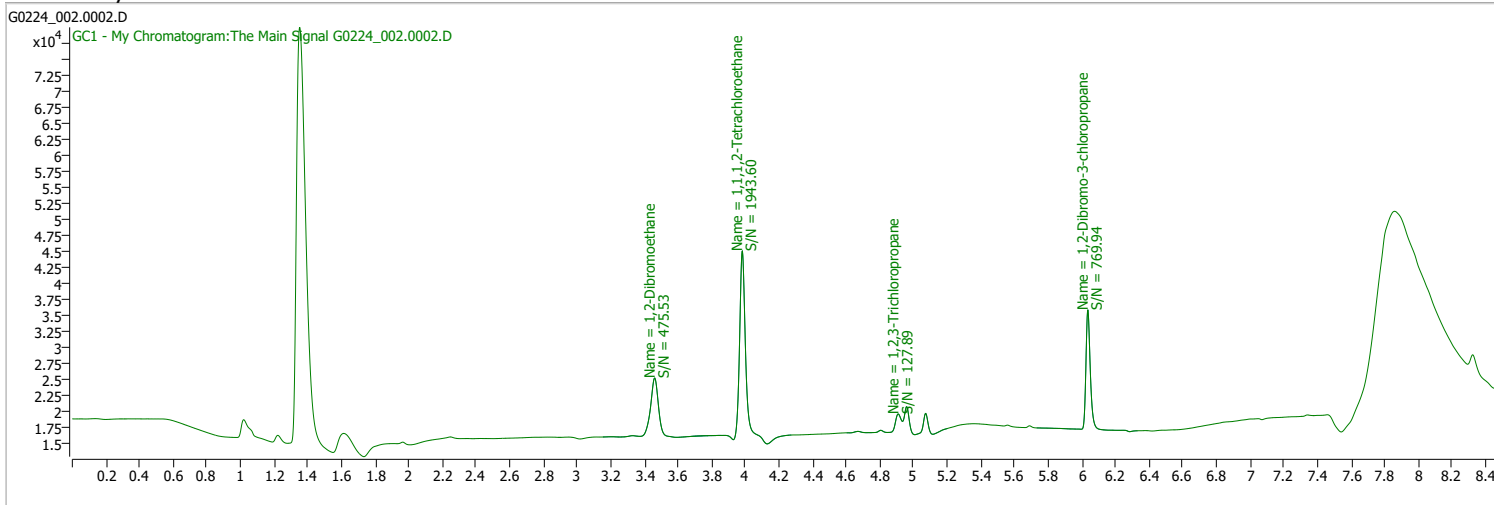
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1725	3.98	0.00	59458				



Quantitation Results Report (QT Reviewed)

Data File	G0224_002.0002.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 12:09:33 PM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.982	0.0	72998	0.2096	µg/L	0.007
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 209.56%	*	

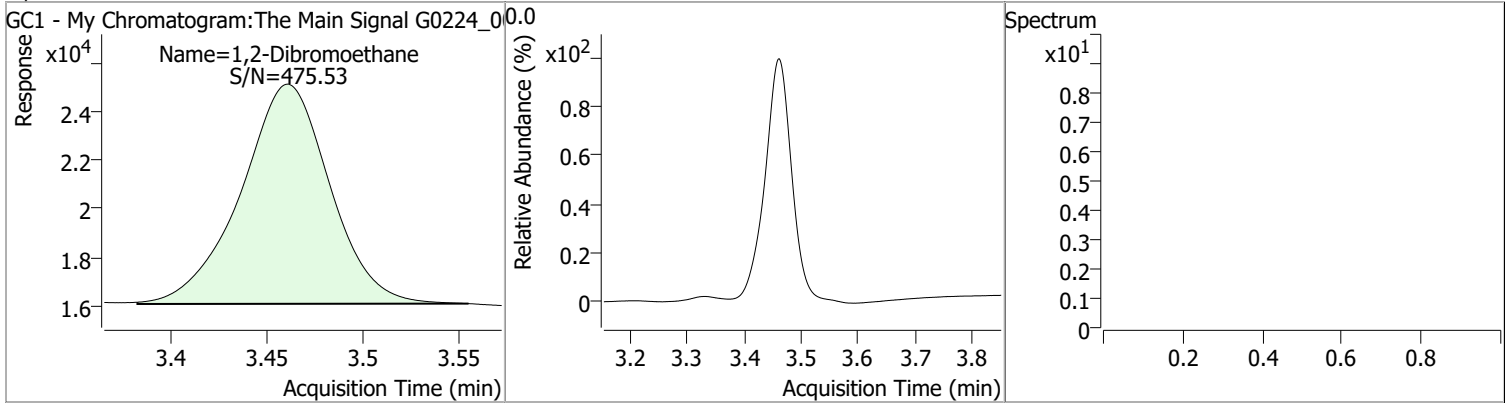
Target Compounds

M 1,2-Dibromoethane	3.462	0.0	29601	0.1793	µg/L	QValue 100
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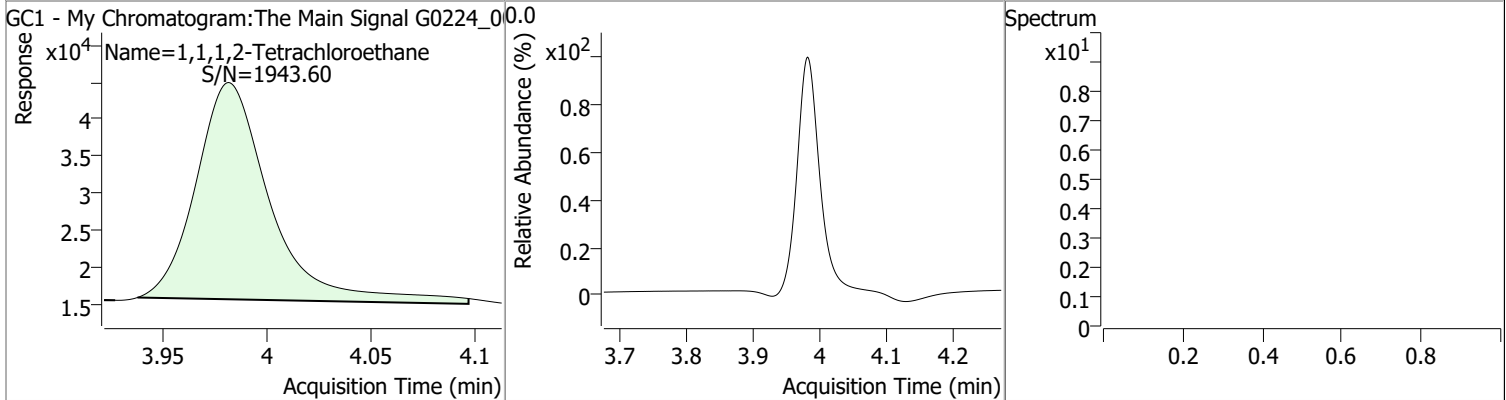
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1793	3.46	0.01	29601				



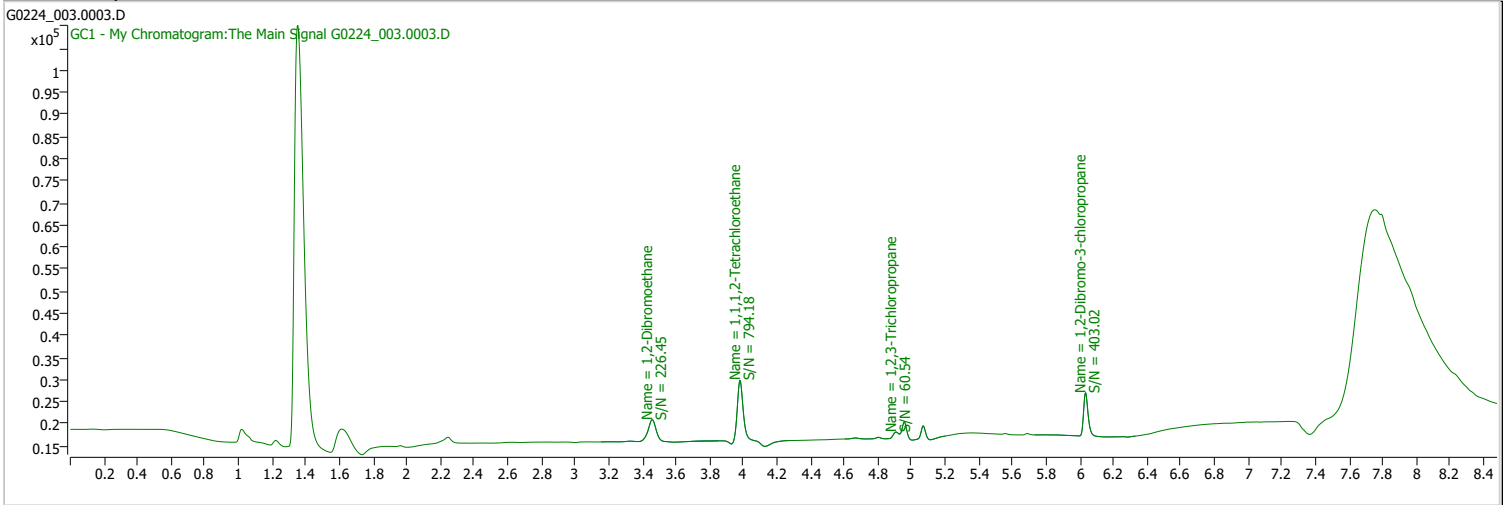
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.2096	3.98	0.01	72998				



Quantitation Results Report (QT Reviewed)

Data File	G0224_003.0003.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 12:29:09 PM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.980	0.0	37683	0.1123	µg/L	0.006
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 112.30%		

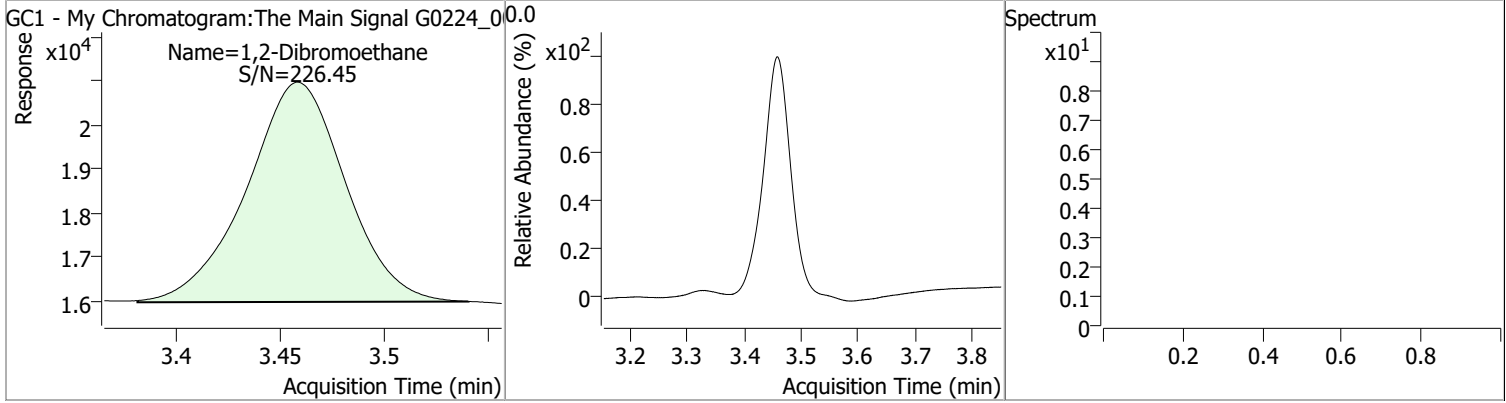
Target Compounds

M 1,2-Dibromoethane	3.458	0.0	16538	0.0988	µg/L	QValue 100
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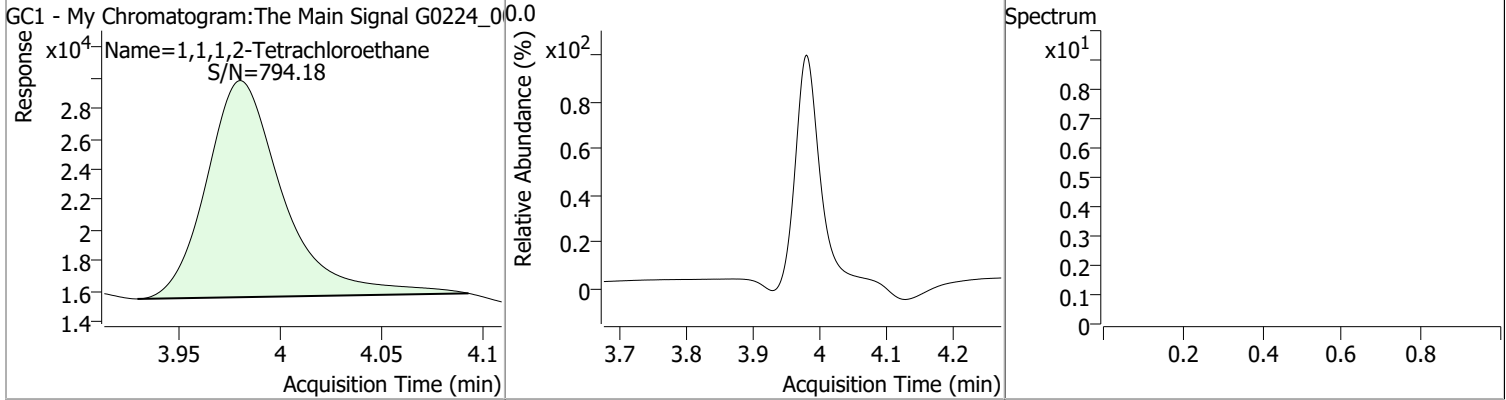
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0988	3.46	0.01	16538				



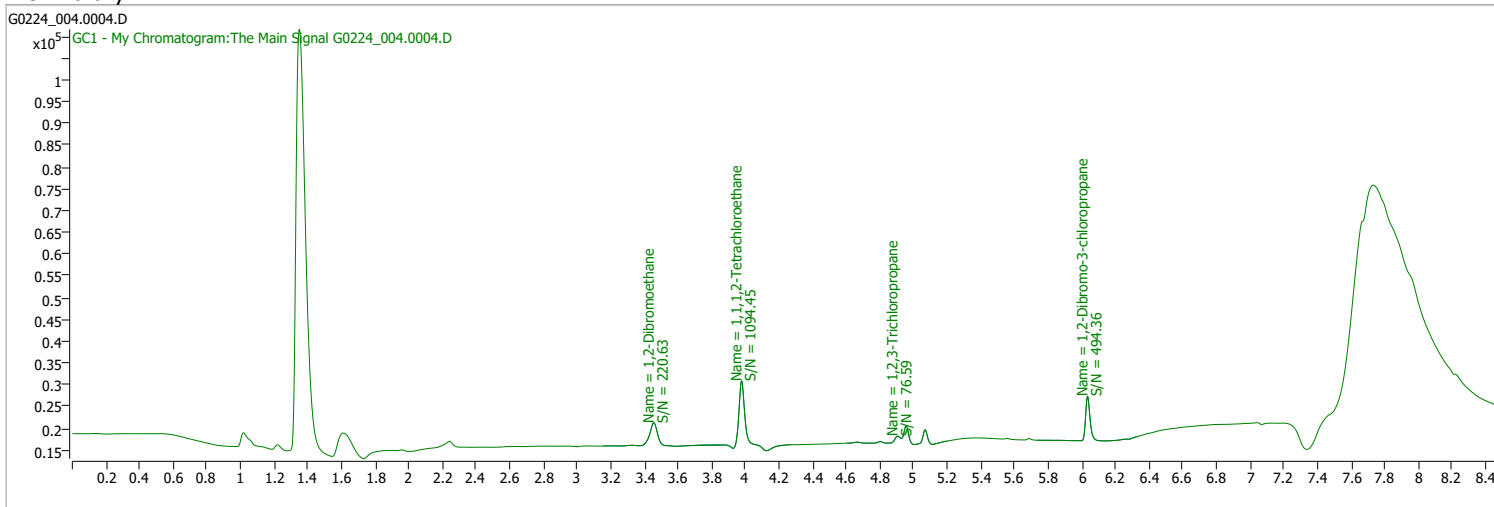
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1123	3.98	0.01	37683				



Quantitation Results Report (QT Reviewed)

Data File	G0224_004.0004.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 12:48:50 PM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

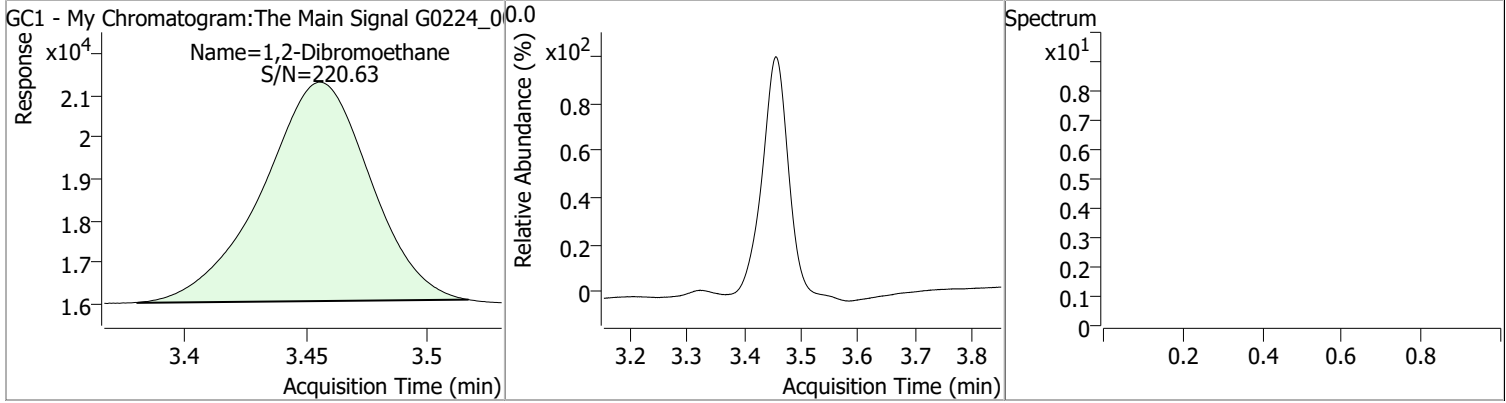


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.978	0.0	42967	0.1270	µg/L	0.003
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 126.97%		
Target Compounds						
M 1,2-Dibromoethane	3.456	0.0	16263	0.0972	µg/L	QValue 100

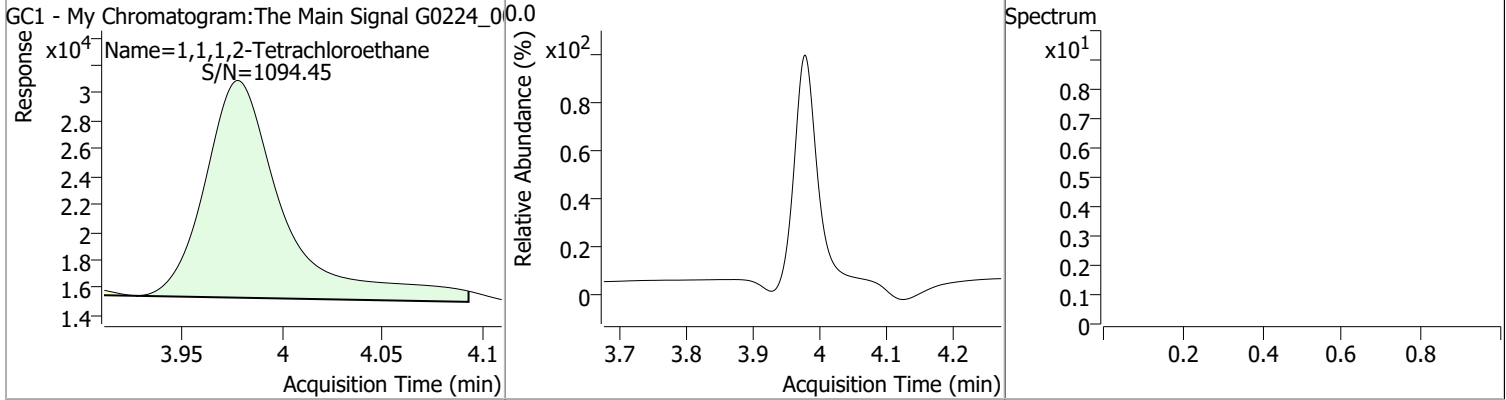
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0972	3.46	0.00	16263				



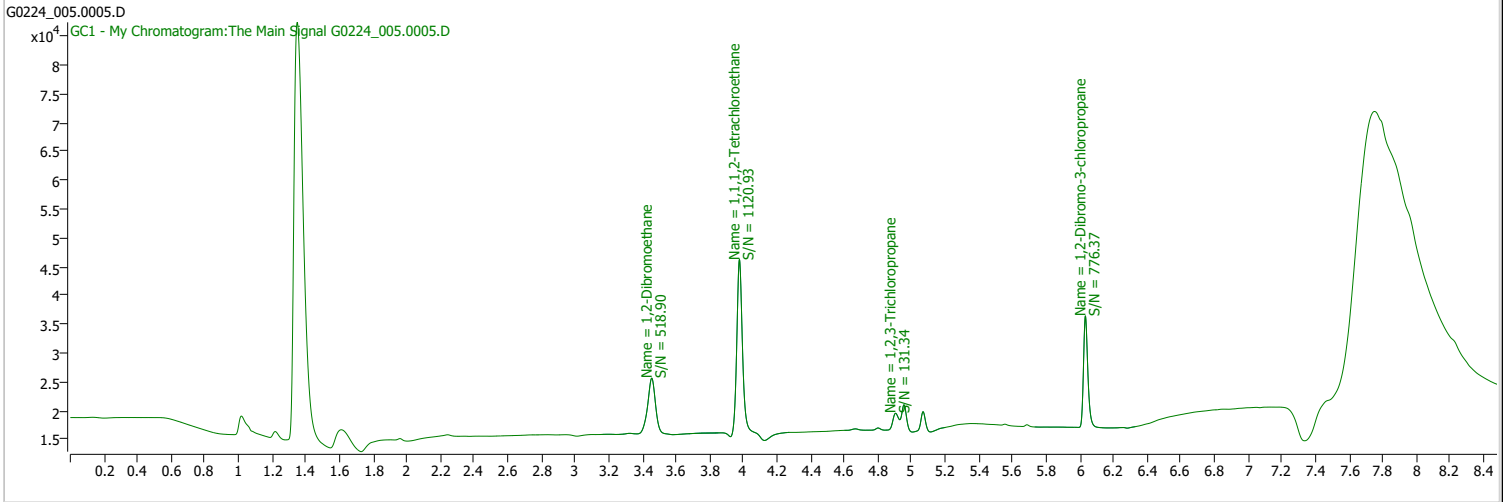
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1270	3.98	0.00	42967				



Quantitation Results Report (QT Reviewed)

Data File	G0224_005.0005.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 1:08:43 PM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

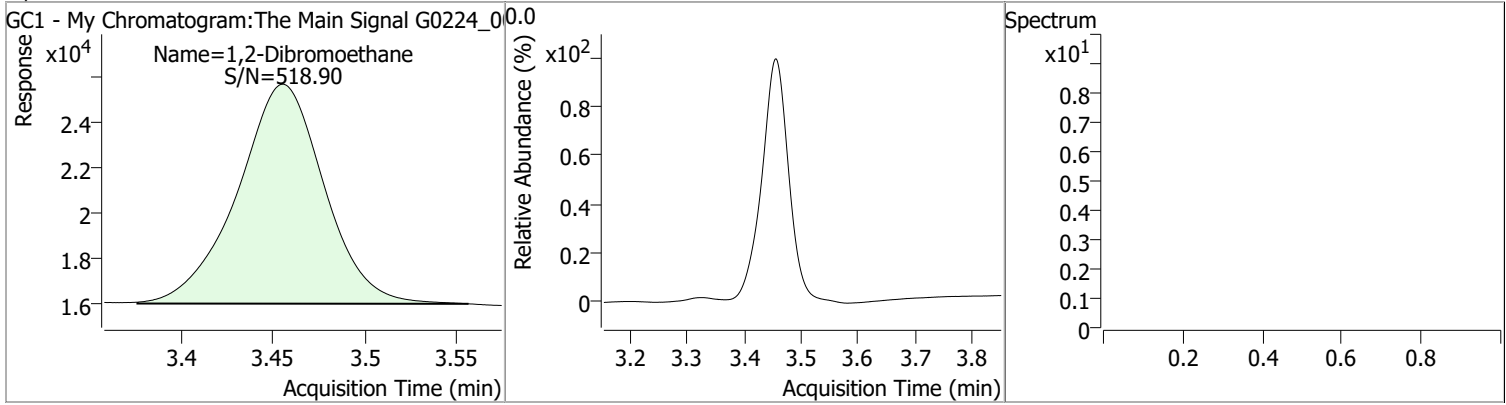


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.977	0.0	73579	0.2111	µg/L	0.002
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 211.15%	*	
Target Compounds						
M 1,2-Dibromoethane	3.456	0.0	31788	0.1929	µg/L	QValue 100

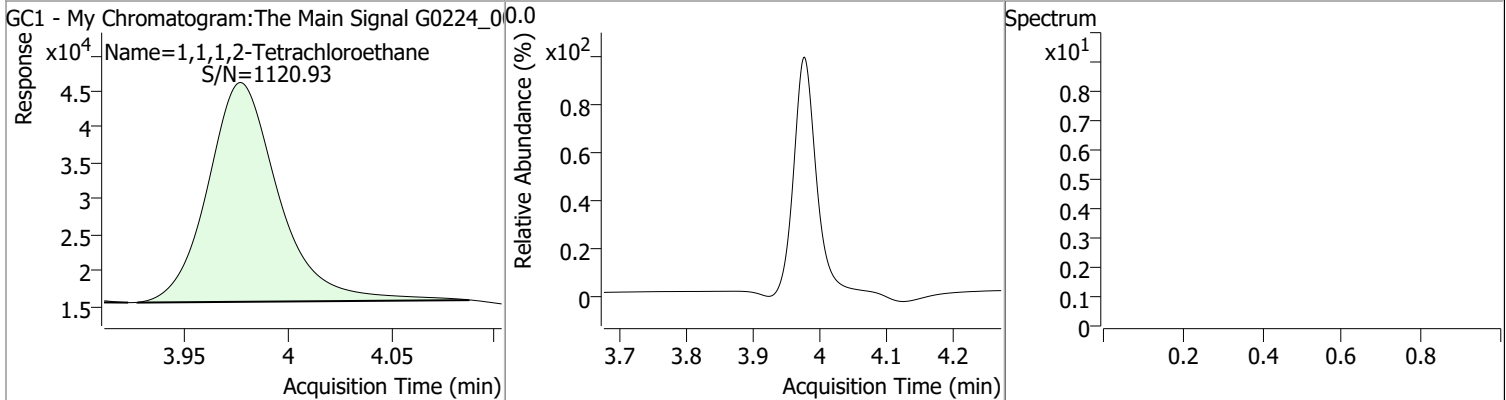
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1929	3.46	0.00	31788				



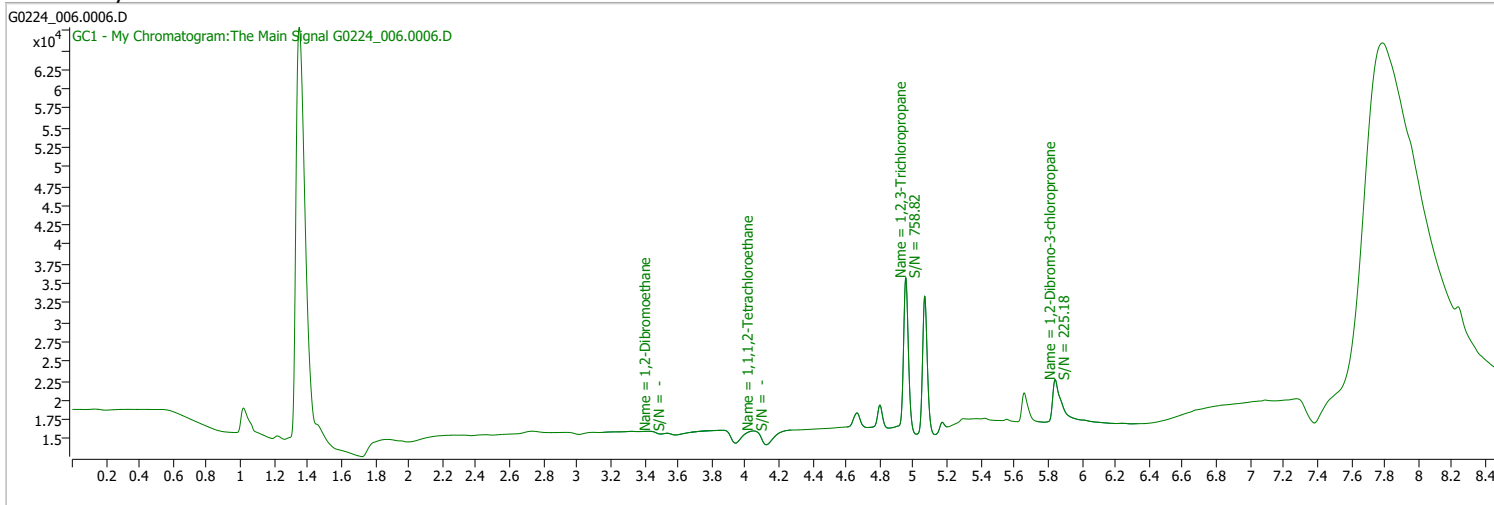
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.2111	3.98	0.00	73579				



Quantitation Results Report (QT Reviewed)

Data File	G0224_006.0006.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 1:28:34 PM
Sample Name	Hexane	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

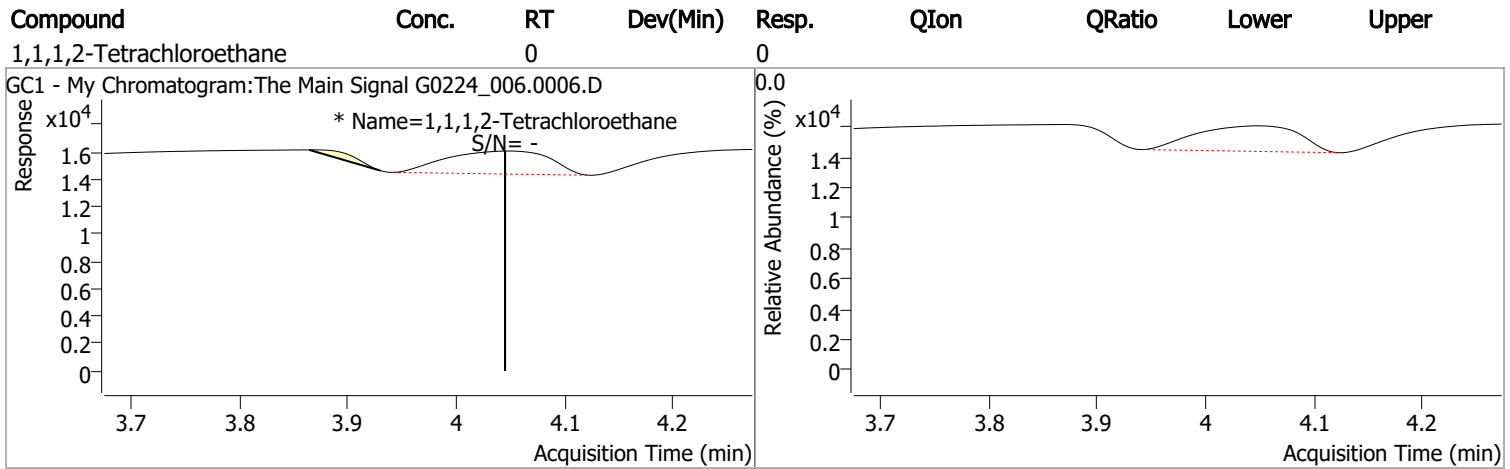
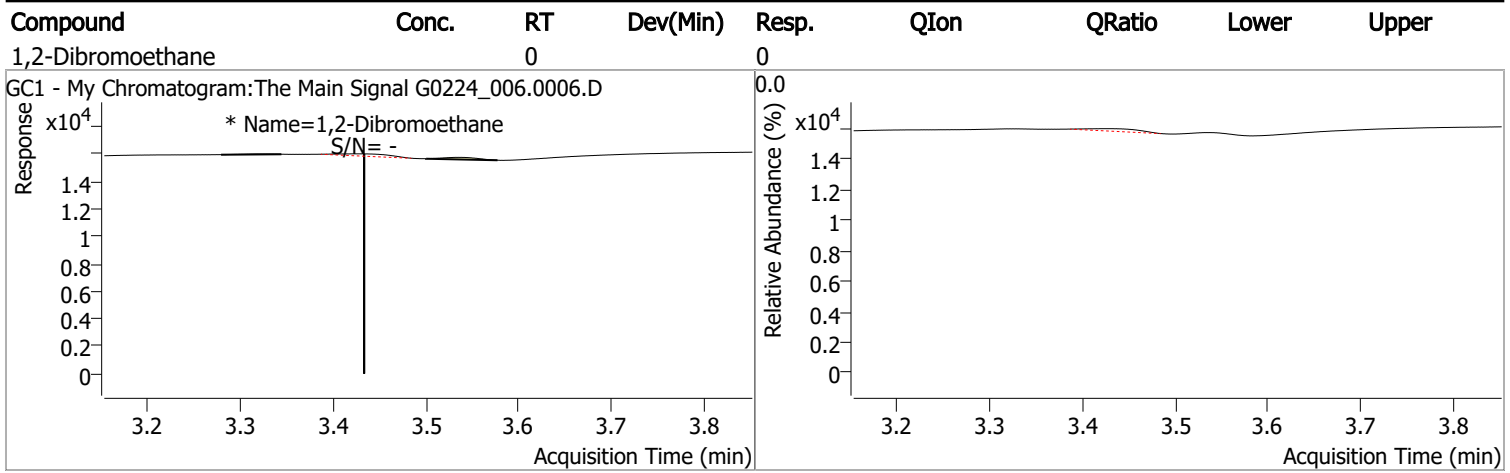
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.045	0.0	0		µg/L	md
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.433	0.0	0		µg/L	md
						QValue
						1

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

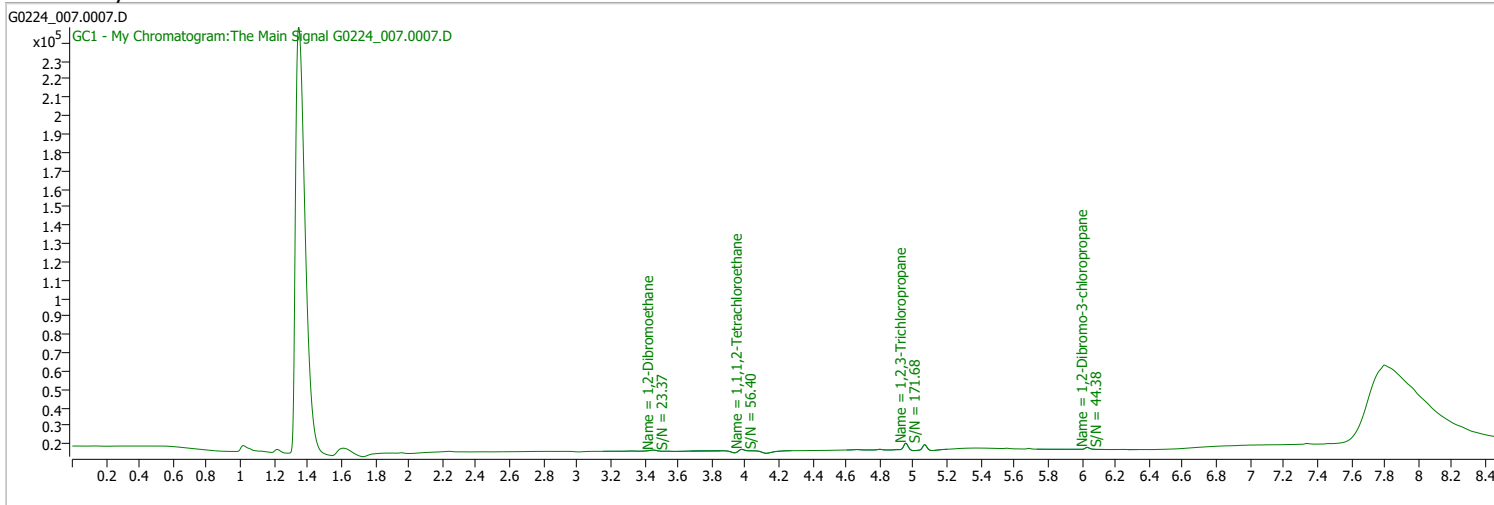
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_007.0007.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 1:48:23 PM
Sample Name	CAL1-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

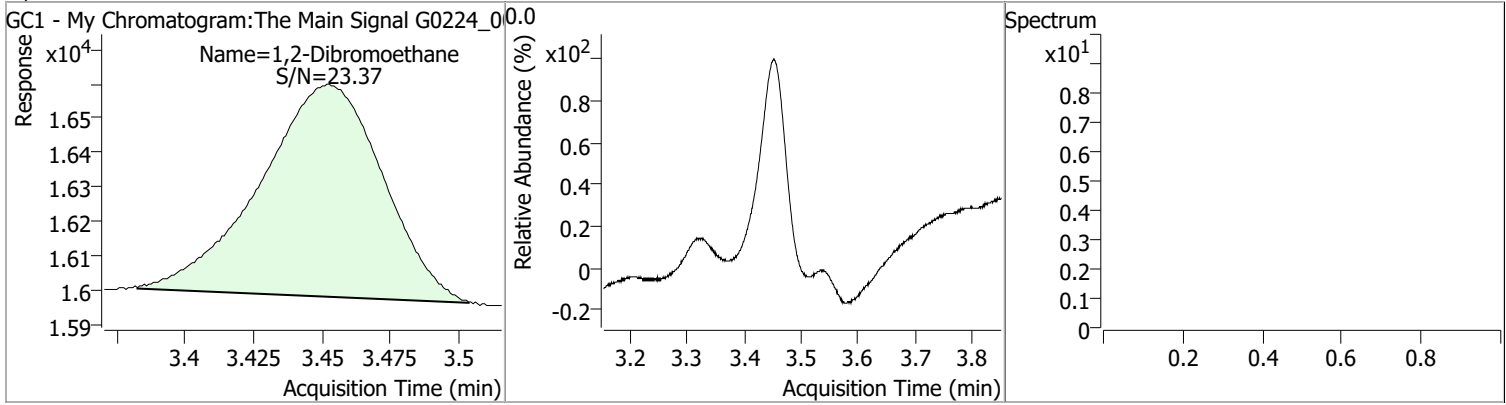


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.978	0.0	1386	0.0105	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 10.45%		*
Target Compounds						
M 1,2-Dibromoethane	3.453	0.0	1929	0.0114	µg/L	QValue 100

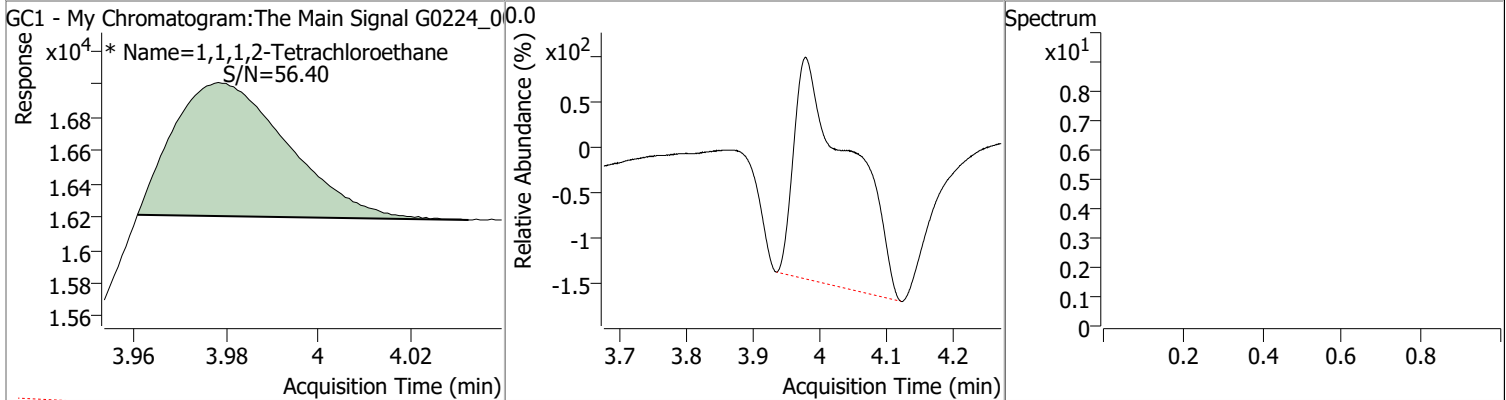
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0114	3.45	0.00	1929				



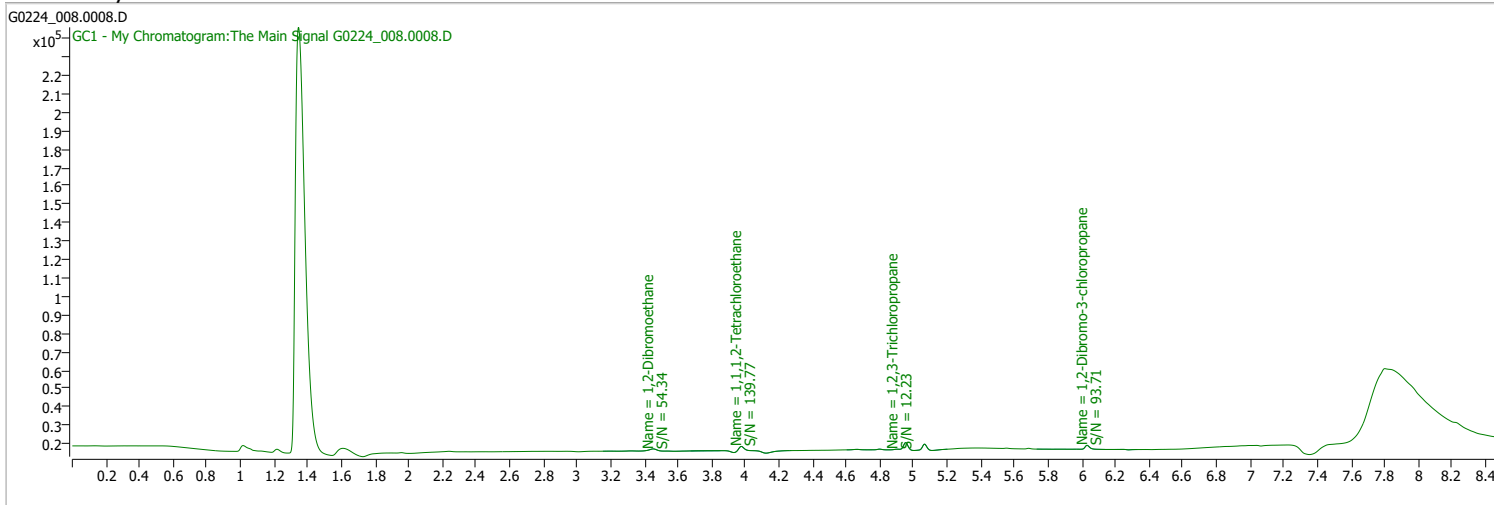
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0105	3.98	0.00	1386 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_008.0008.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/24/2022 2:07:59 PM
Sample Name	CAL7-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

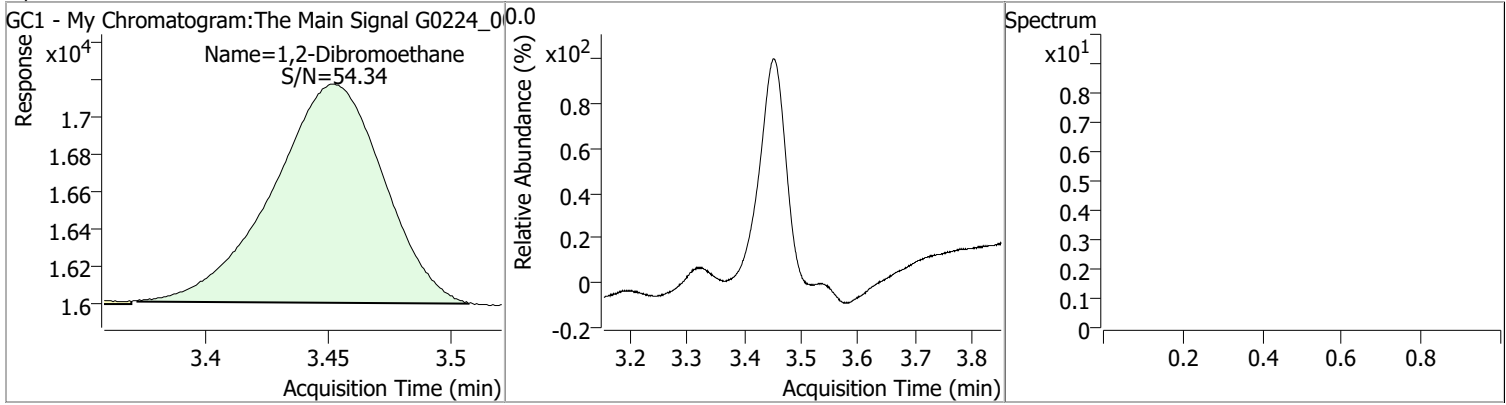


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.976	0.0	4766	0.0200	µg/L	0.002
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 20.02%	*	
Target Compounds						
M 1,2-Dibromoethane	3.453	0.0	3724	0.0220	µg/L	QValue 100

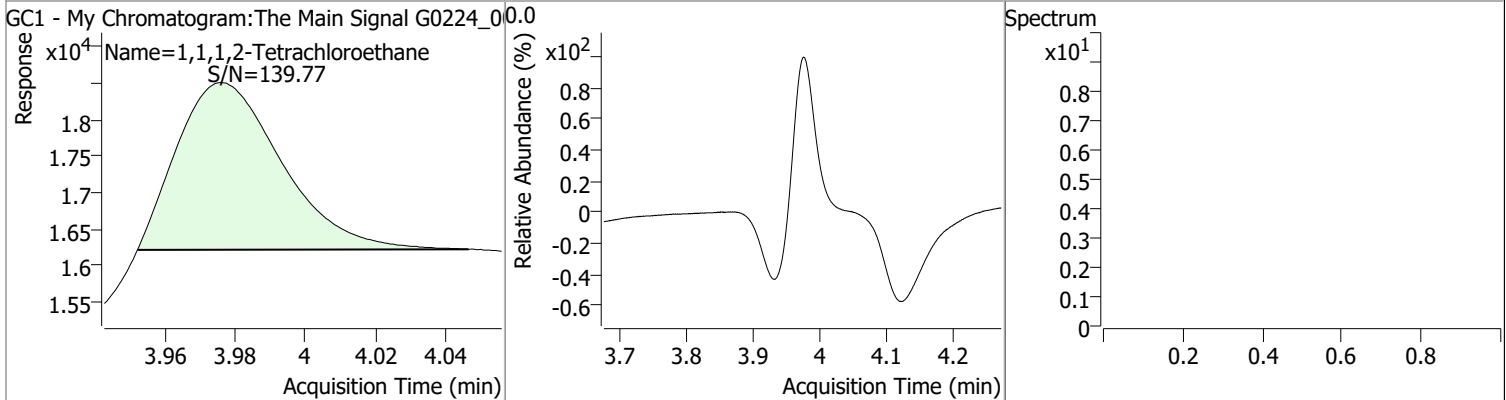
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0220	3.45	0.00	3724				



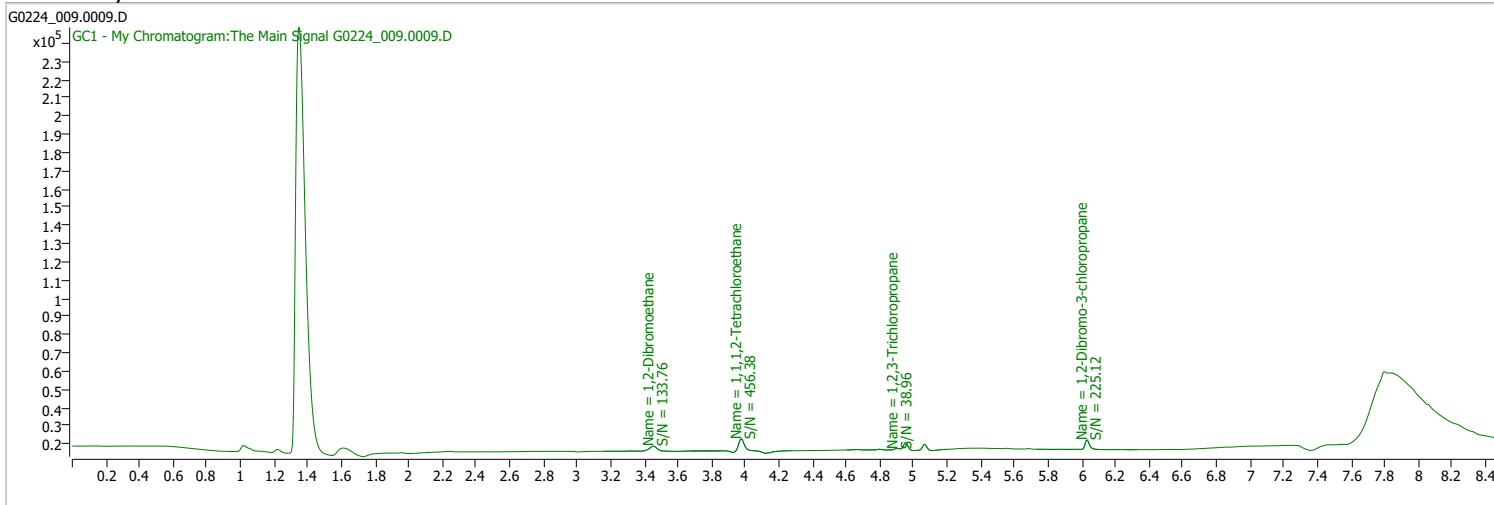
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0200	3.98	0.00	4766				



Quantitation Results Report (QT Reviewed)

Data File	G0224_009.0009.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 2:27:56 PM
Sample Name	CAL2-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

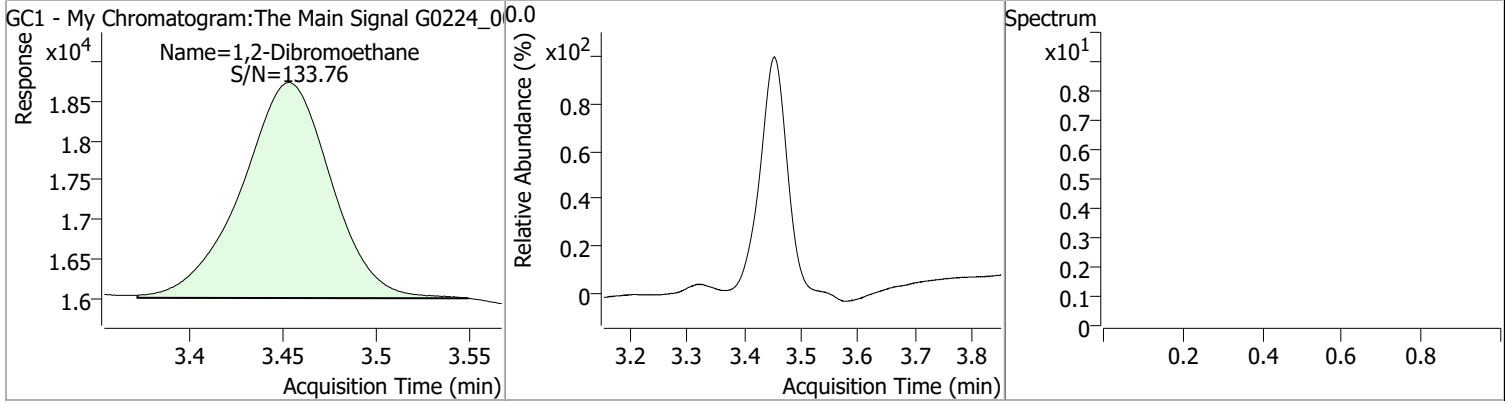


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.975	0.0	14968	0.0488	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 48.79%		*
Target Compounds						
M 1,2-Dibromoethane	3.453	0.0	9221	0.0547	µg/L	QValue 100

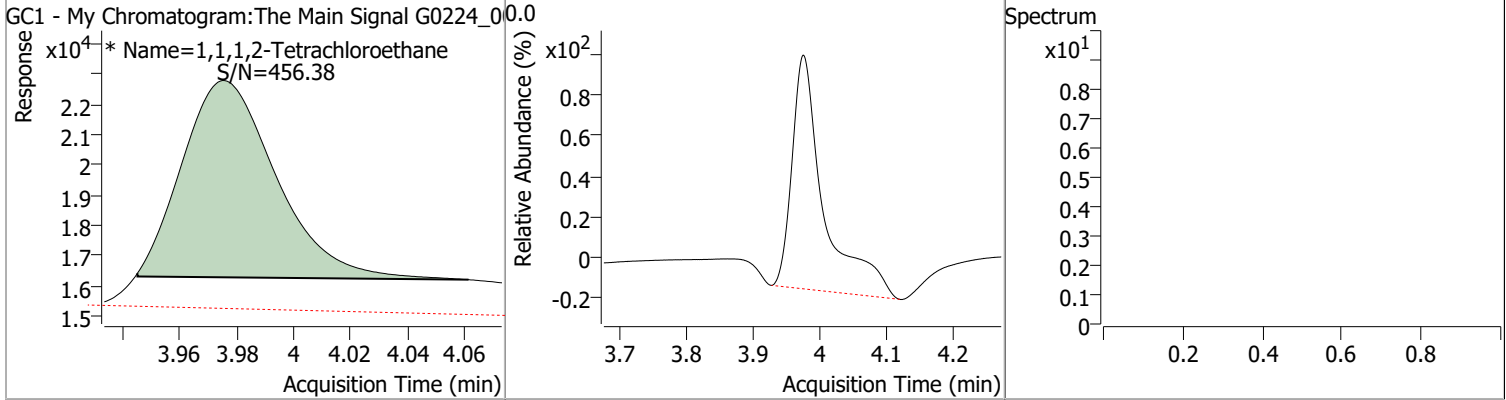
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0547	3.45	0.00	9221				



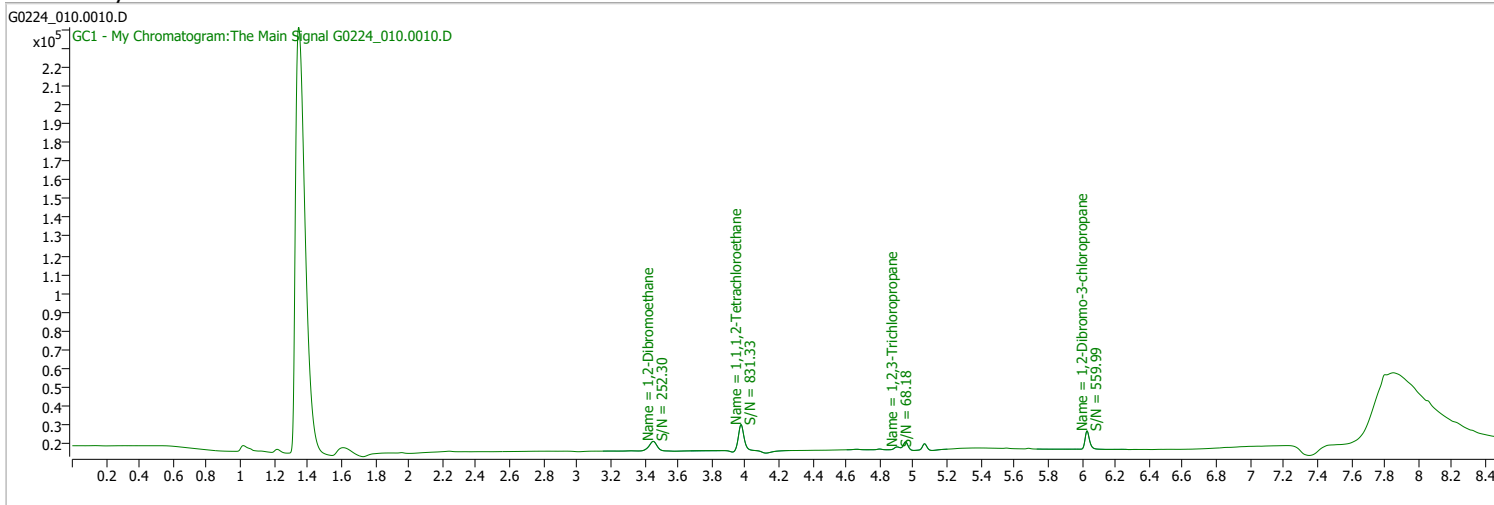
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0488	3.98	0.00	14968 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_010.0010.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/24/2022 2:47:36 PM
Sample Name	CAL3-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

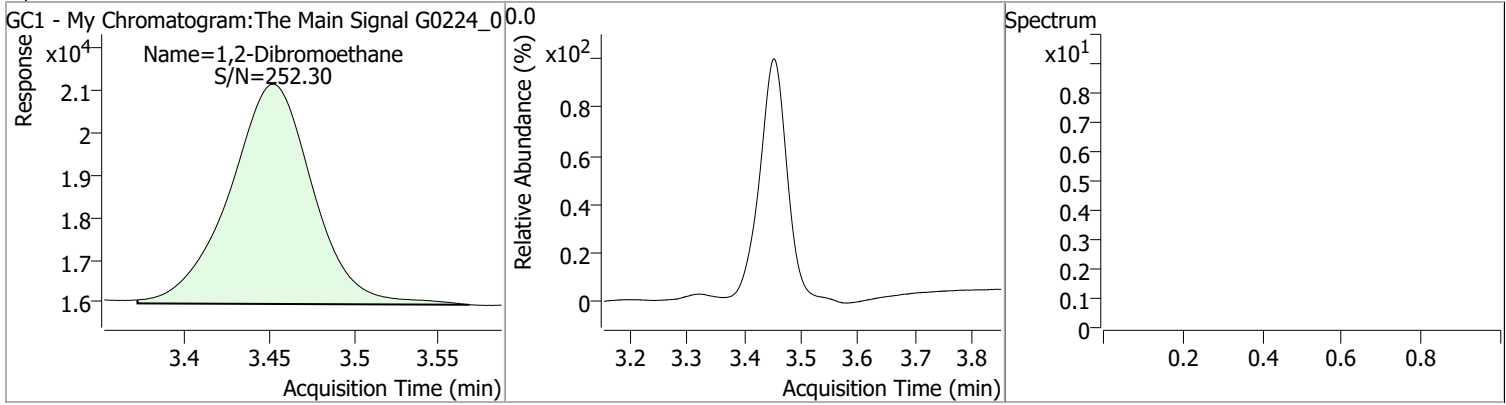


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.974	0.0	32499	0.0979	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%		Recovery = 97.88%			
Target Compounds						
M 1,2-Dibromoethane	3.453	0.0	17768	0.1063	µg/L	QValue 100

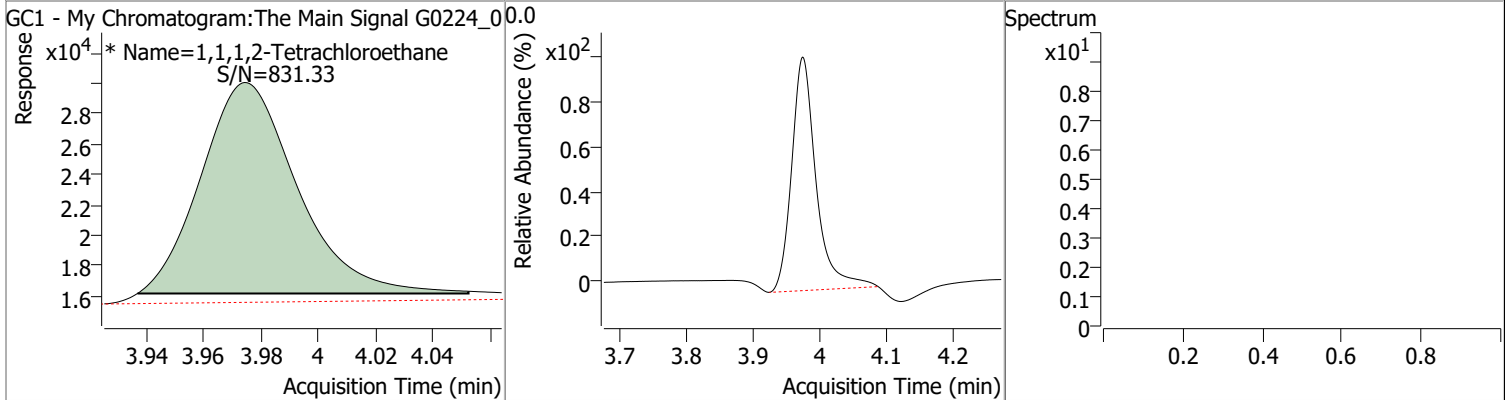
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1063	3.45	0.00	17768				



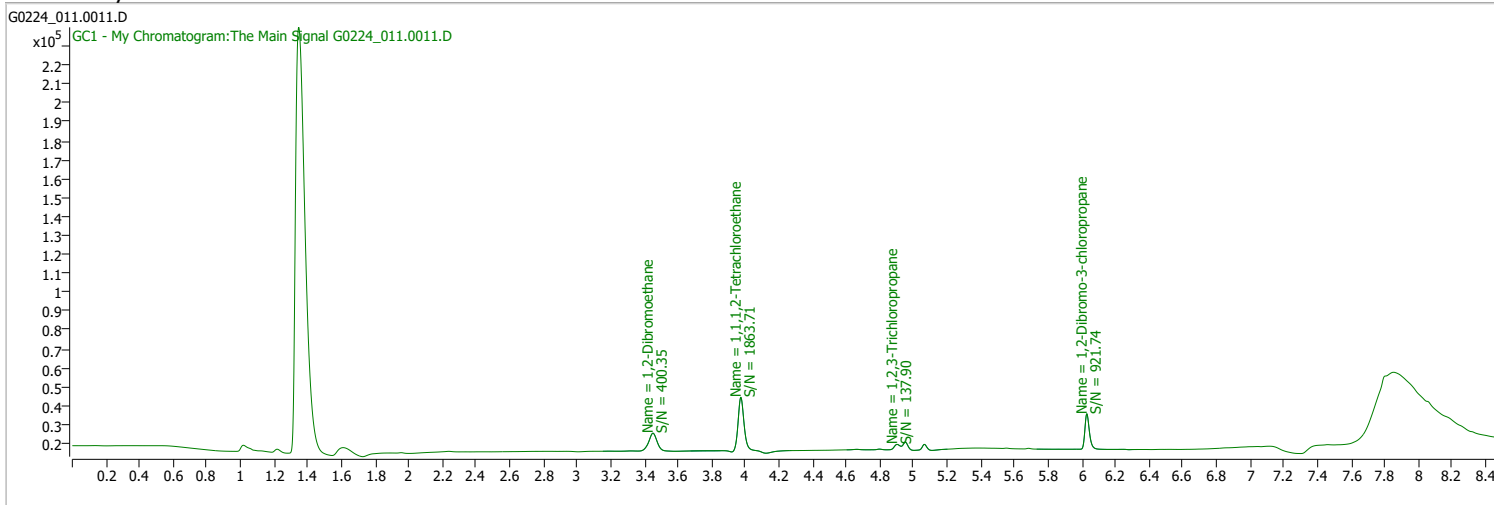
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0979	3.97	0.00	32499 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_011.0011.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/24/2022 3:07:26 PM
Sample Name	CAL4-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.973	0.0	67779	0.1953	µg/L	m	-0.001
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 195.30%		*	

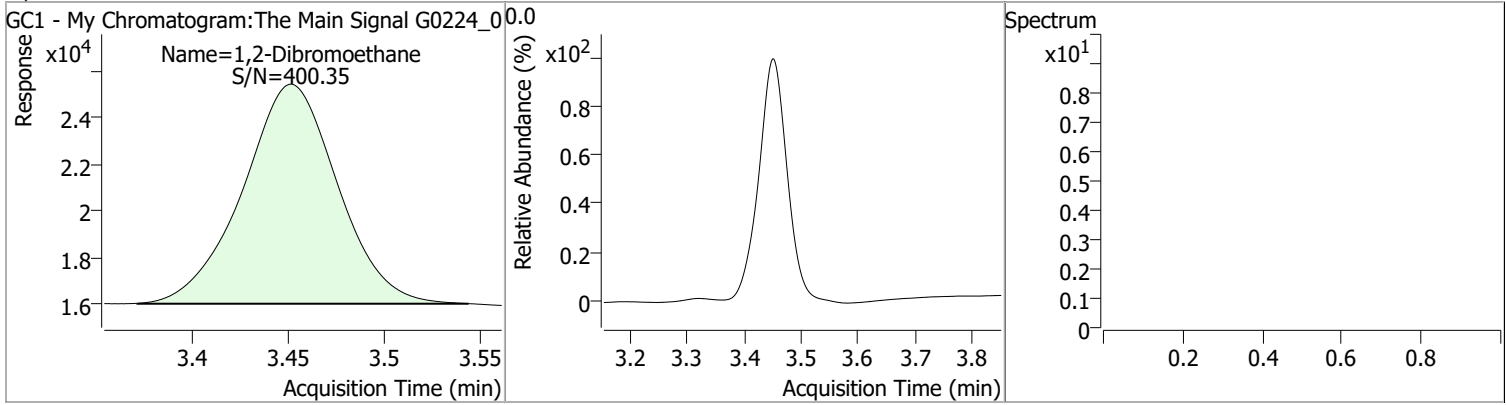
Target Compounds

M 1,2-Dibromoethane	3.451	0.0	32418	0.1969	µg/L		QValue 100
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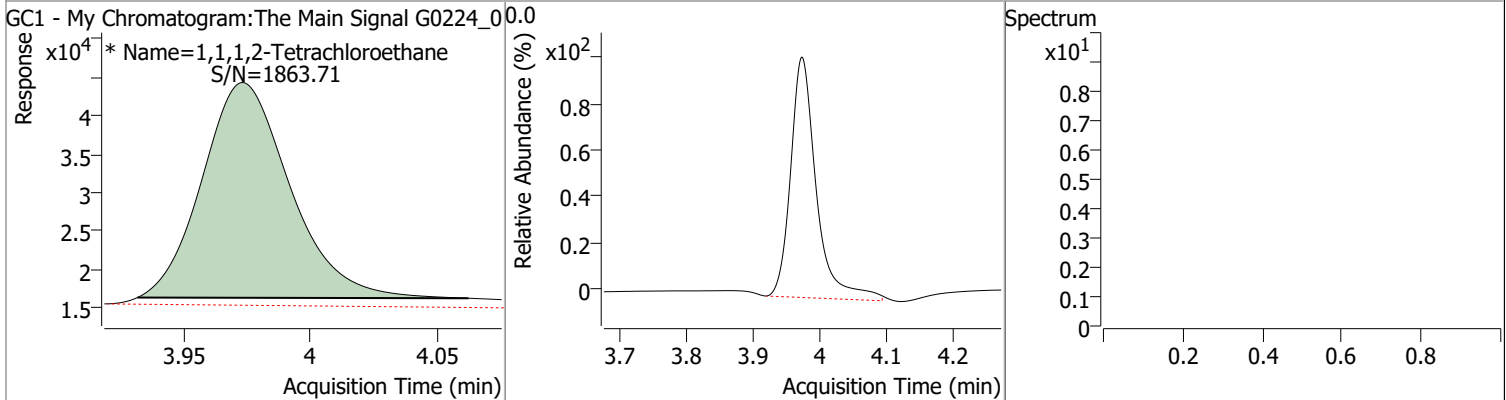
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1969	3.45	0.00	32418				



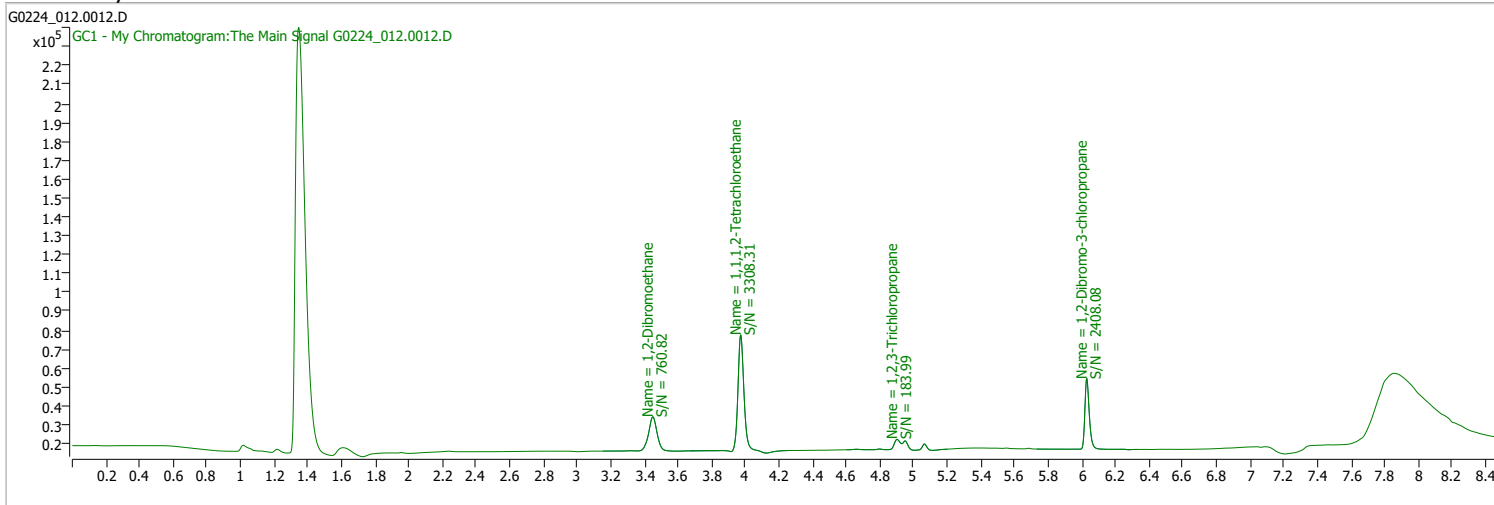
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1953	3.97	0.00	67779 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_012.0012.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 3:27:03 PM
Sample Name	CAL5-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

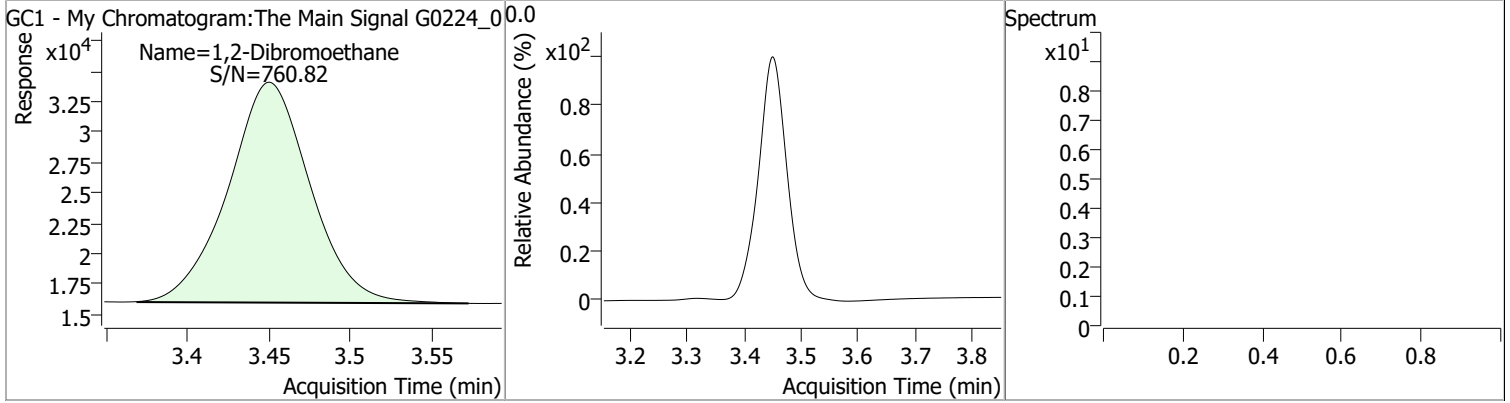


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	147845	0.4101	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 410.11%	*	
Target Compounds						
M 1,2-Dibromoethane	3.450	0.0	63436	0.3986	µg/L	QValue 100

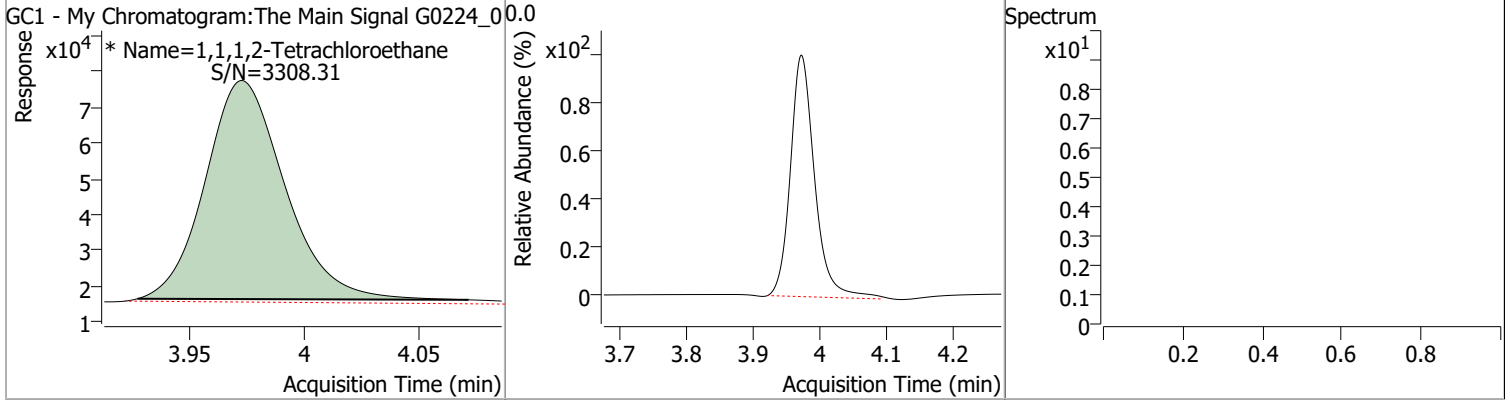
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.3986	3.45	0.00	63436				



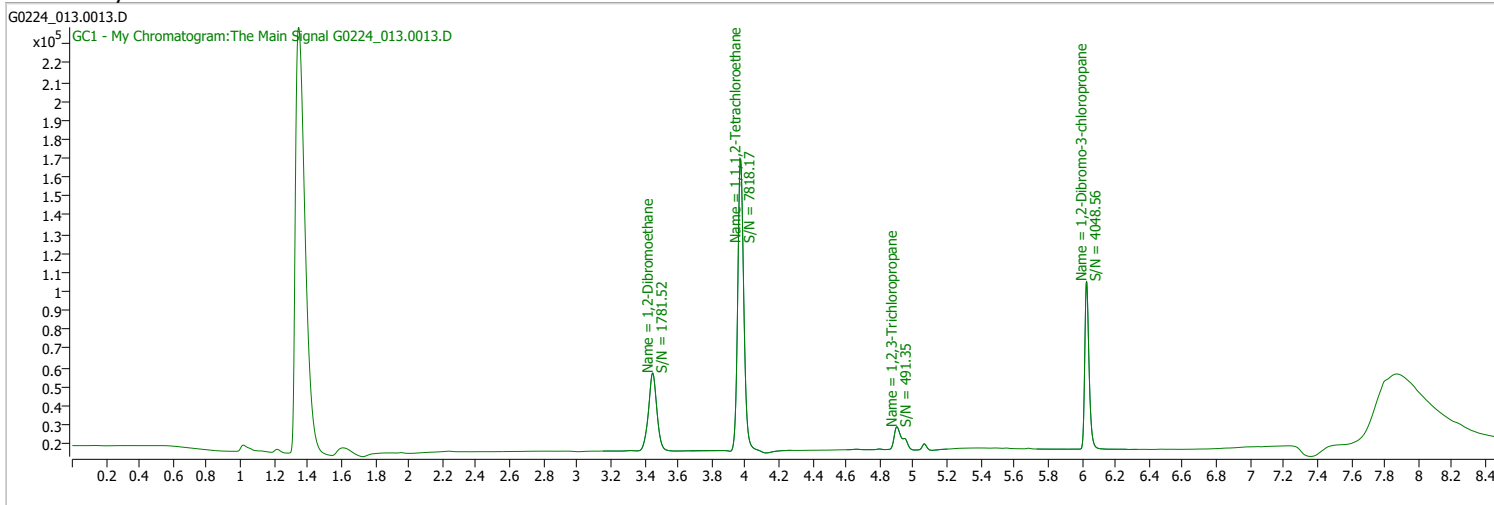
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.4101	3.97	0.00	147845 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_013.0013.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 3:46:49 PM
Sample Name	CAL6-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.971	0.0	382731	0.9974	µg/L	m	-0.003
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 997.39%		*	

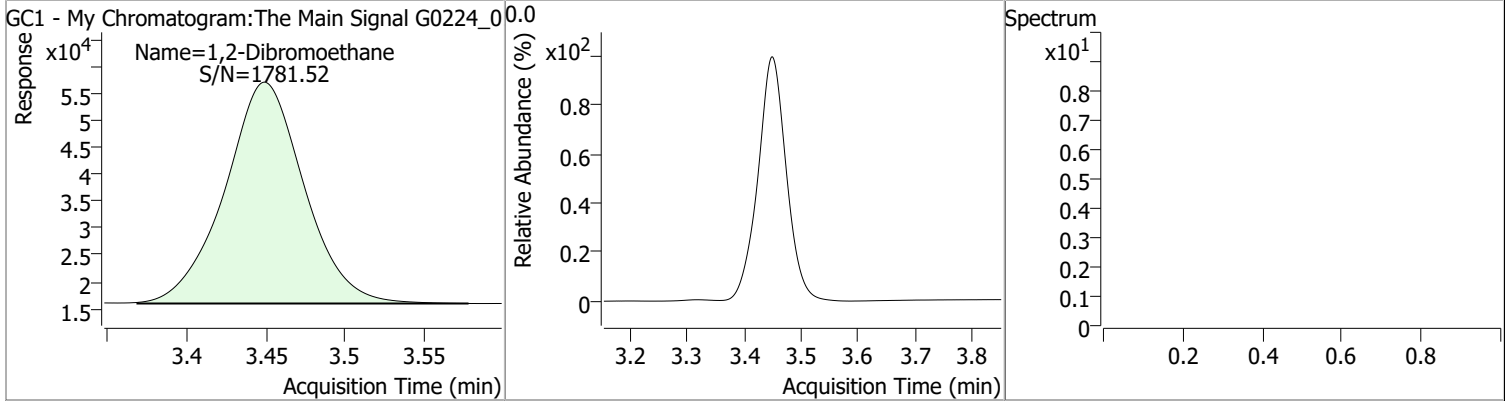
Target Compounds

M 1,2-Dibromoethane	3.449	0.0	142853	1.0004	µg/L		QValue 100
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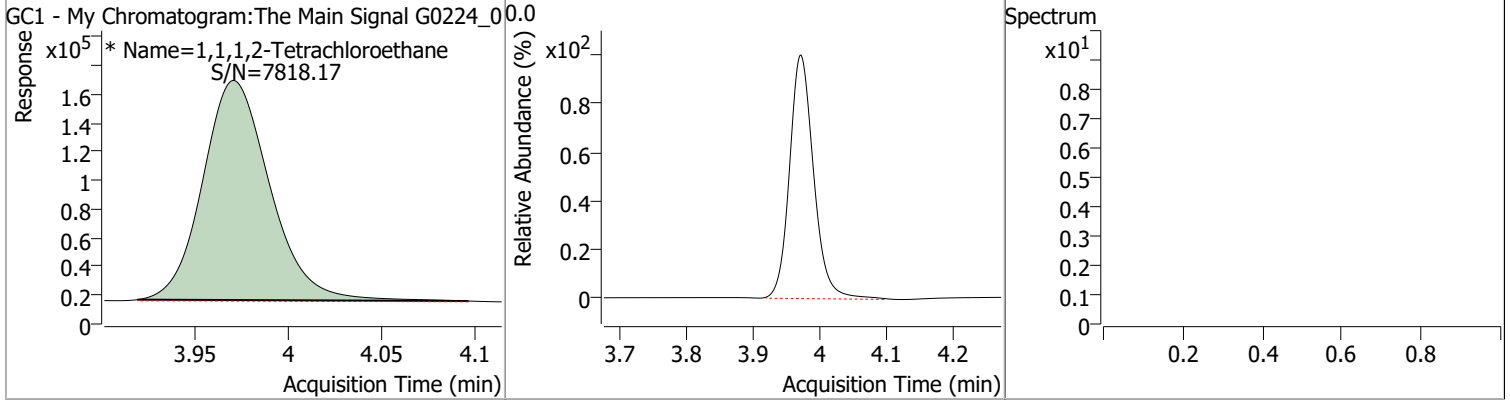
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	1.0004	3.45	0.00	142853				



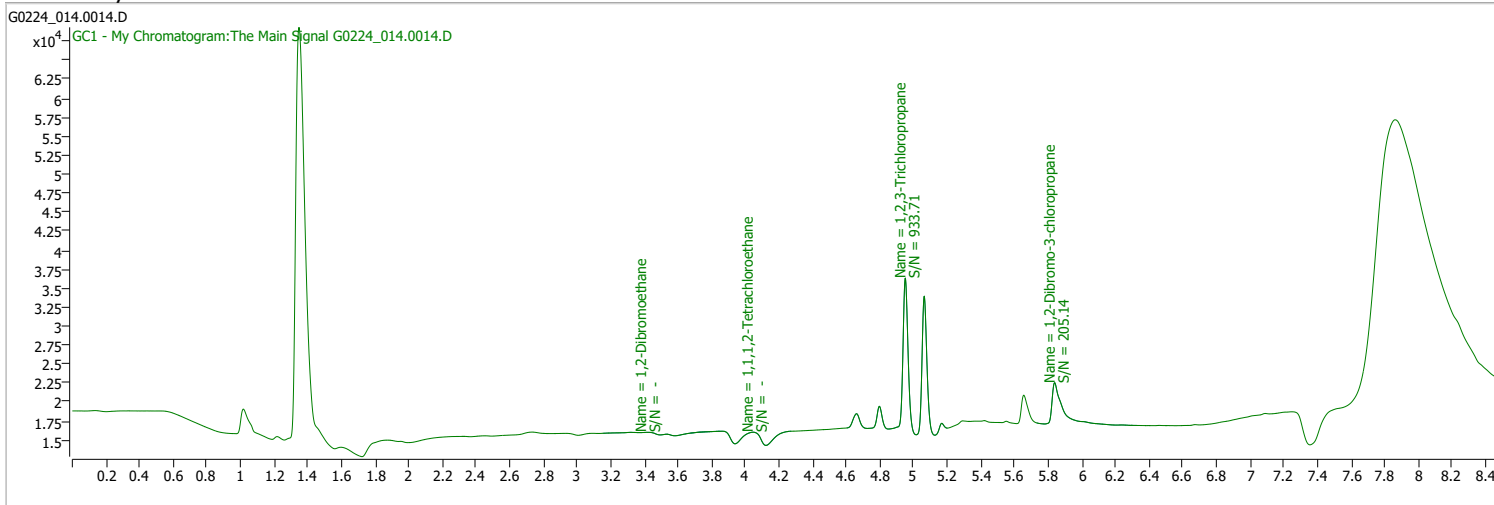
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.9974	3.97	0.00	382731 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_014.0014.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 4:06:39 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

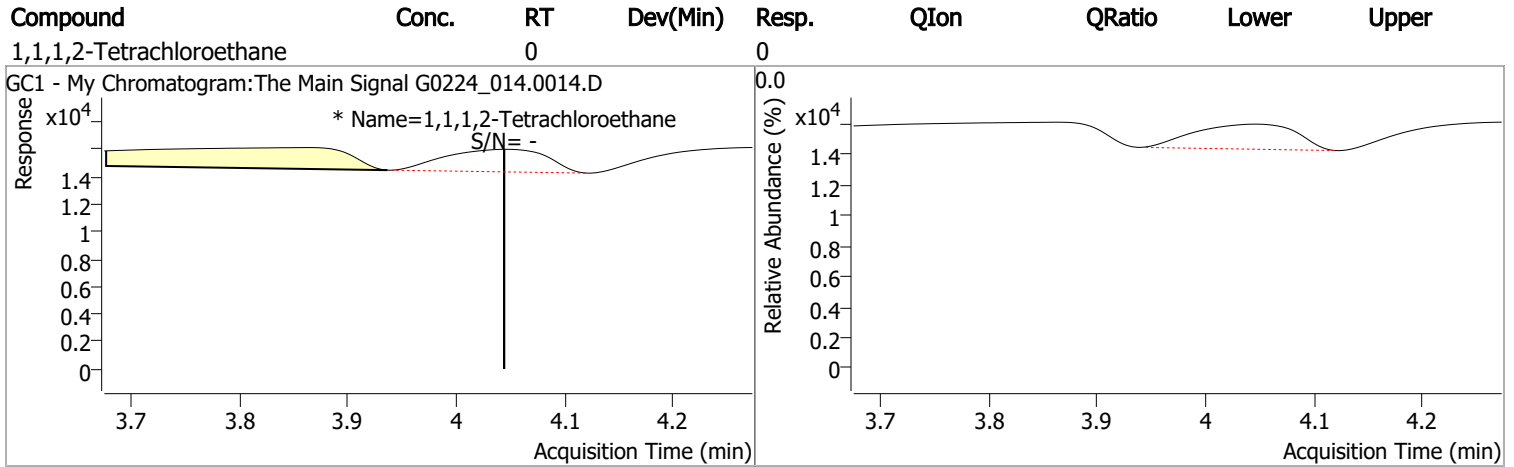
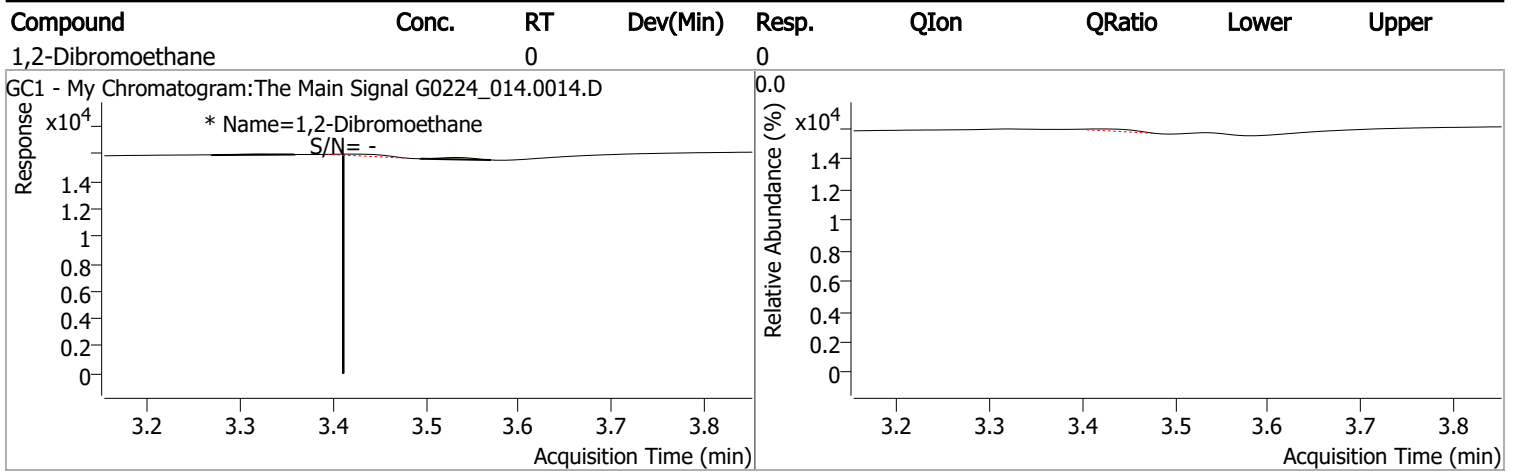
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.044	0.0	0		µg/L	md
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.411	0.0	0		µg/L	md
						QValue
						1

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

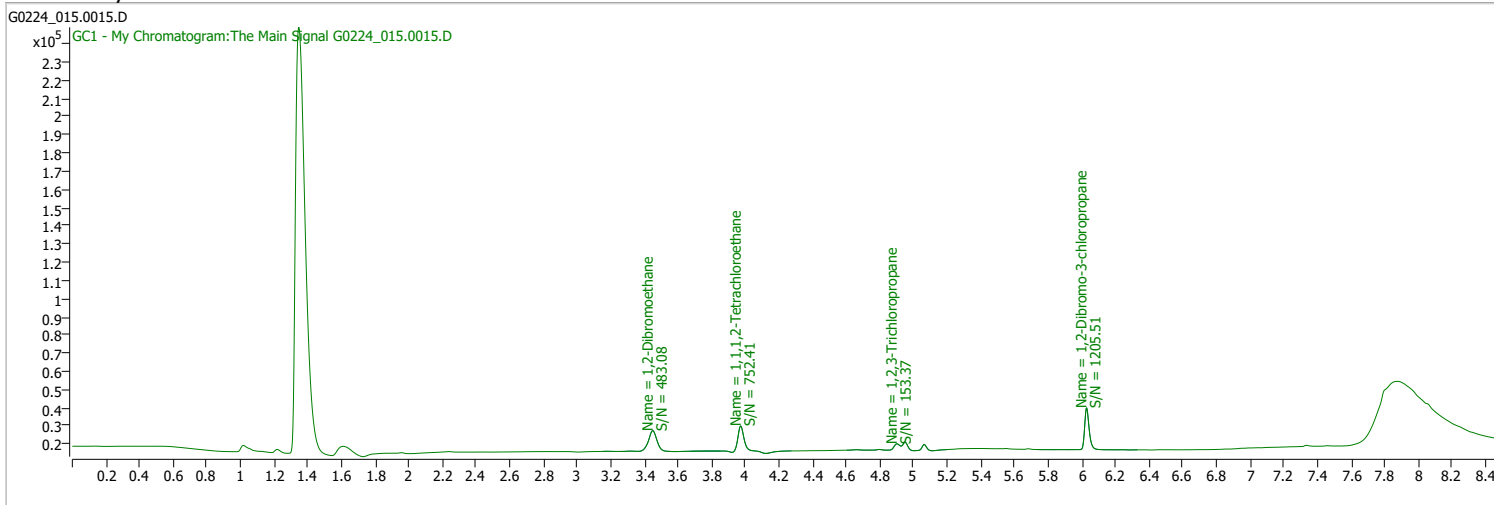
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_015.0015.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/24/2022 4:26:23 PM
Sample Name	LCS-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

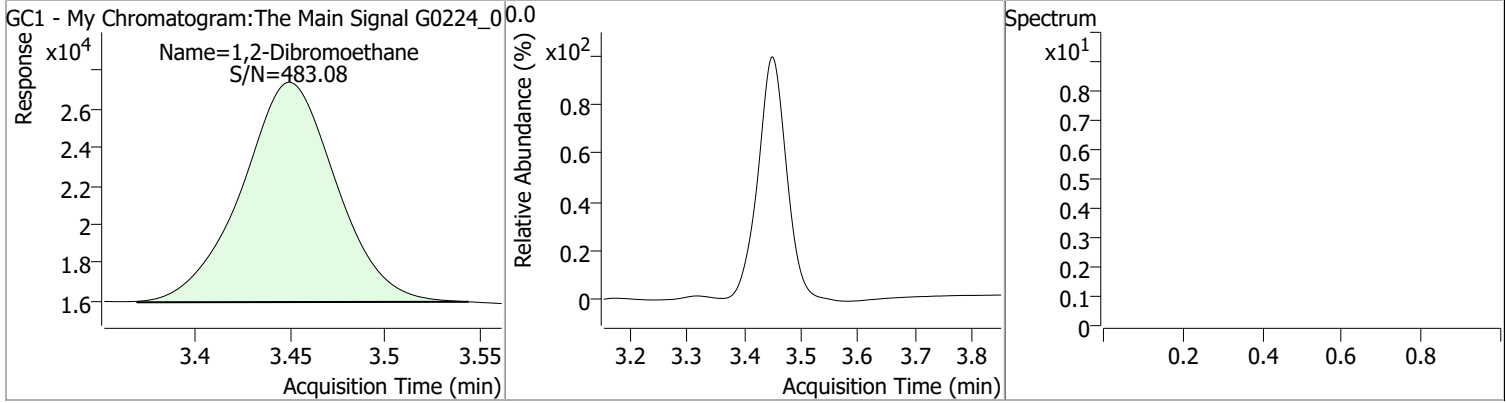


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	34577	0.1037	µg/L	m
Spiked Amount: 0.100				Range: 70.0 - 130.0% Recovery = 103.67%		
Target Compounds						
M 1,2-Dibromoethane	3.449	0.0	39595	0.2423	µg/L	QValue 100

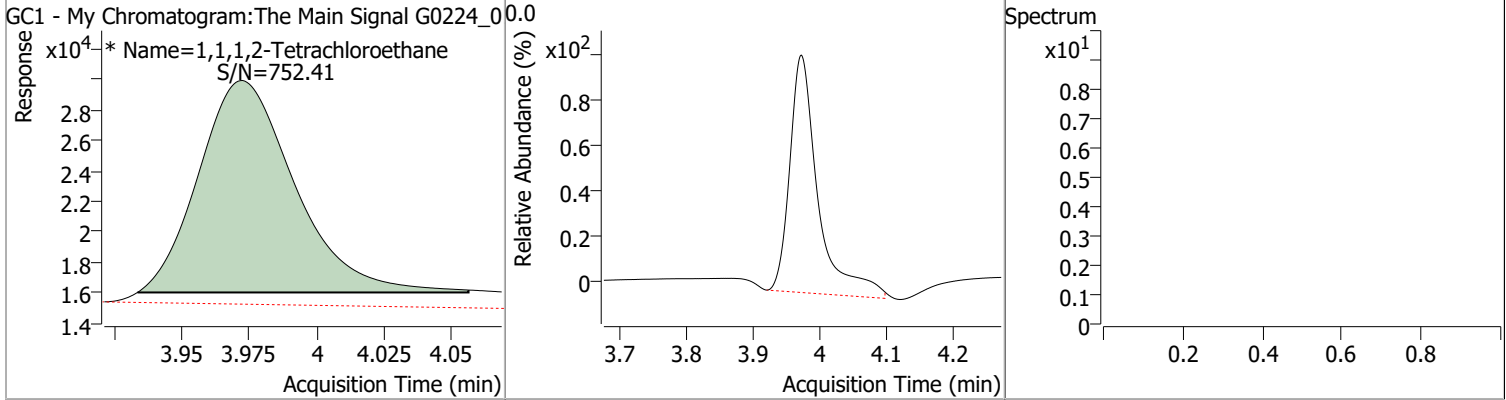
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2423	3.45	0.00	39595				



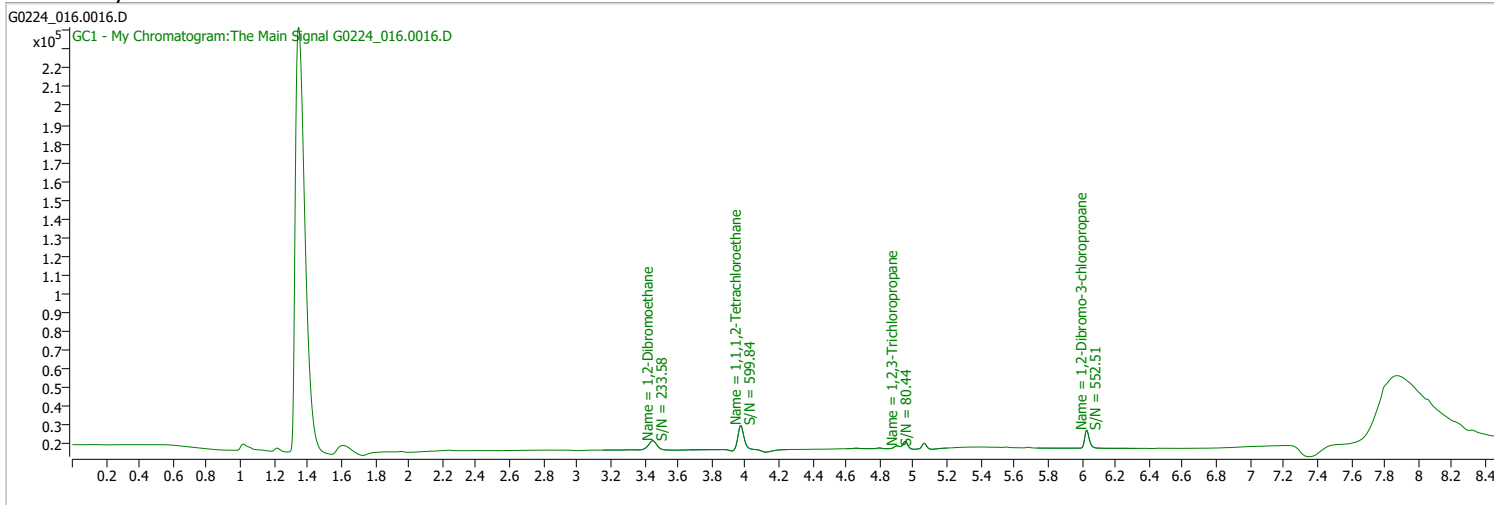
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1037	3.97	0.00	34577 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_016.0016.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 4:46:10 PM
Sample Name	CAL3-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

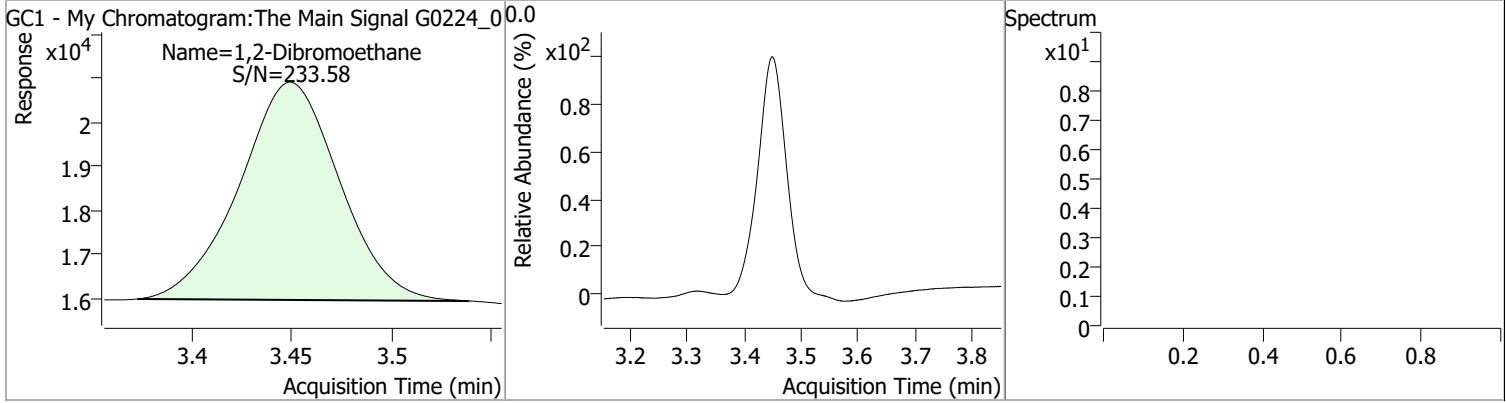


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	31159	0.0941	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 94.14%		
Target Compounds						
M 1,2-Dibromoethane	3.450	0.0	17133	0.1025	µg/L	100

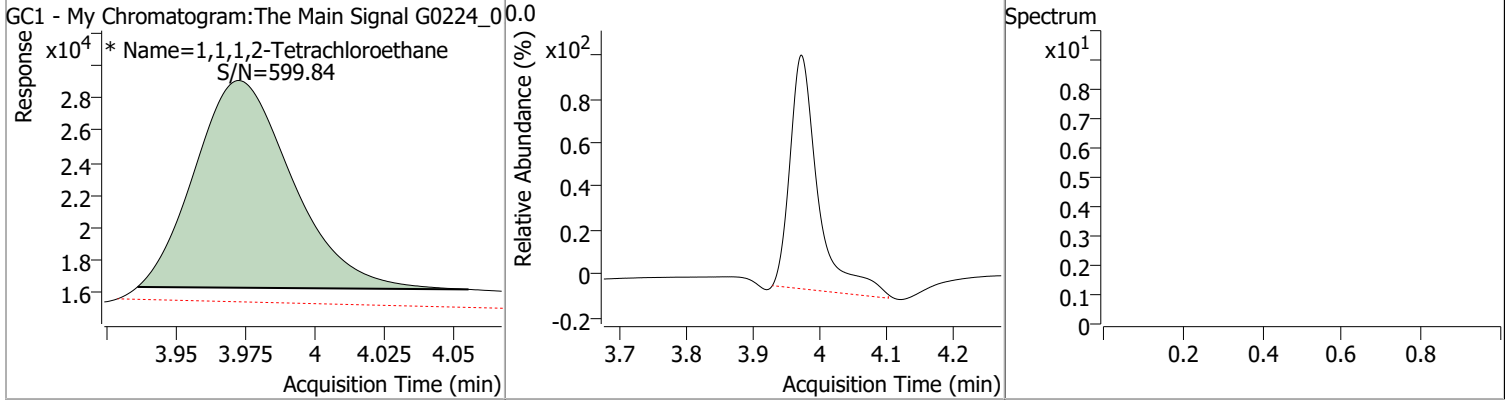
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1025	3.45	0.00	17133				



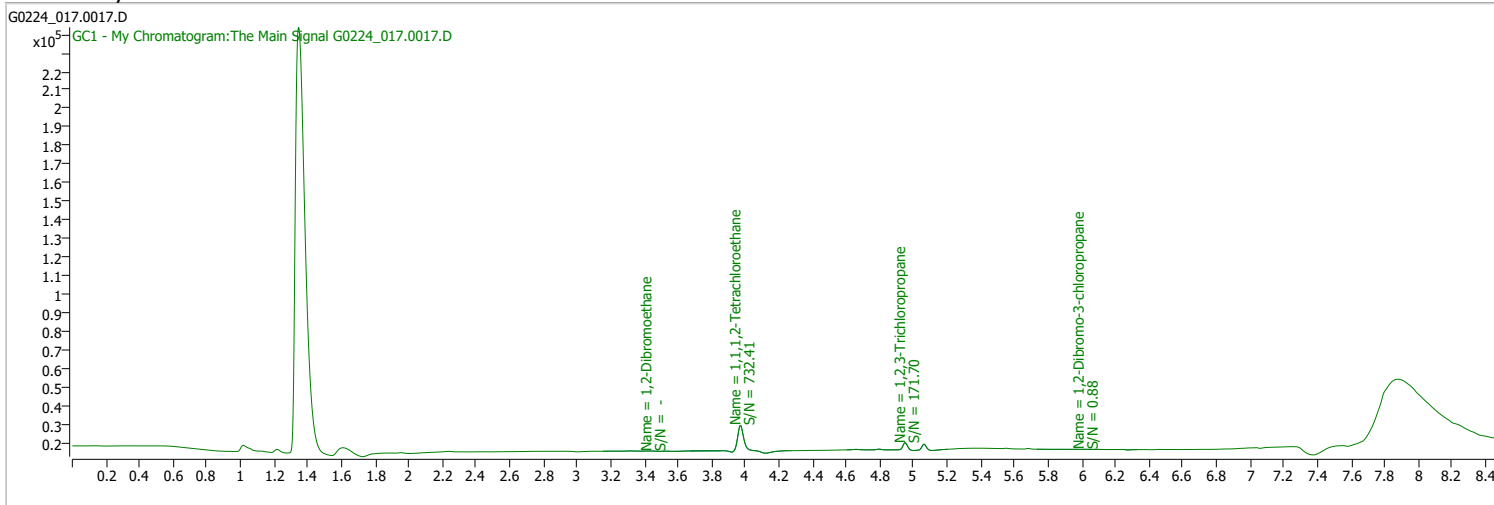
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0941	3.97	0.00	31159 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_017.0017.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 5:06:01 PM
Sample Name	MB-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

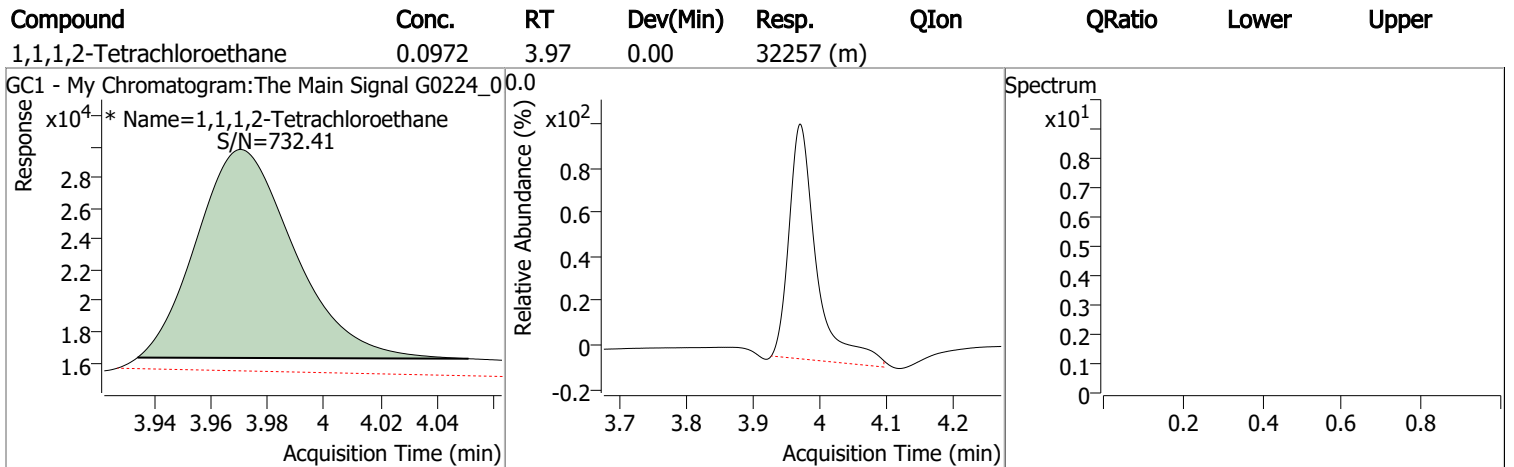
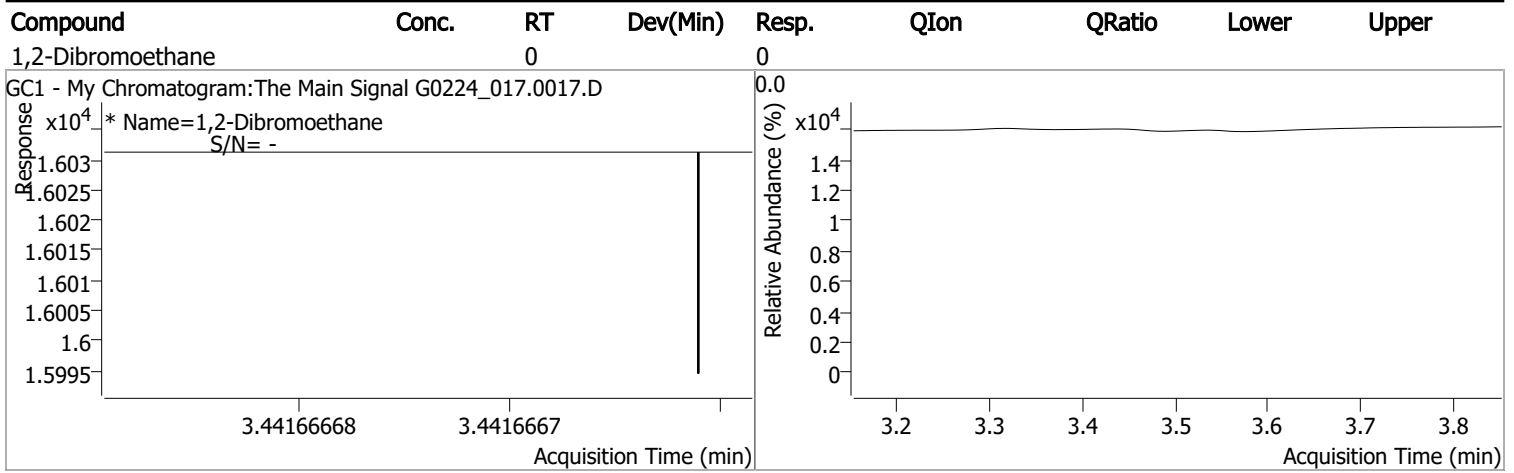
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	32257	0.0972	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 97.20%		
Target Compounds						
M 1,2-Dibromoethane	3.442	0.0	0		µg/L	md
						QValue 1

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

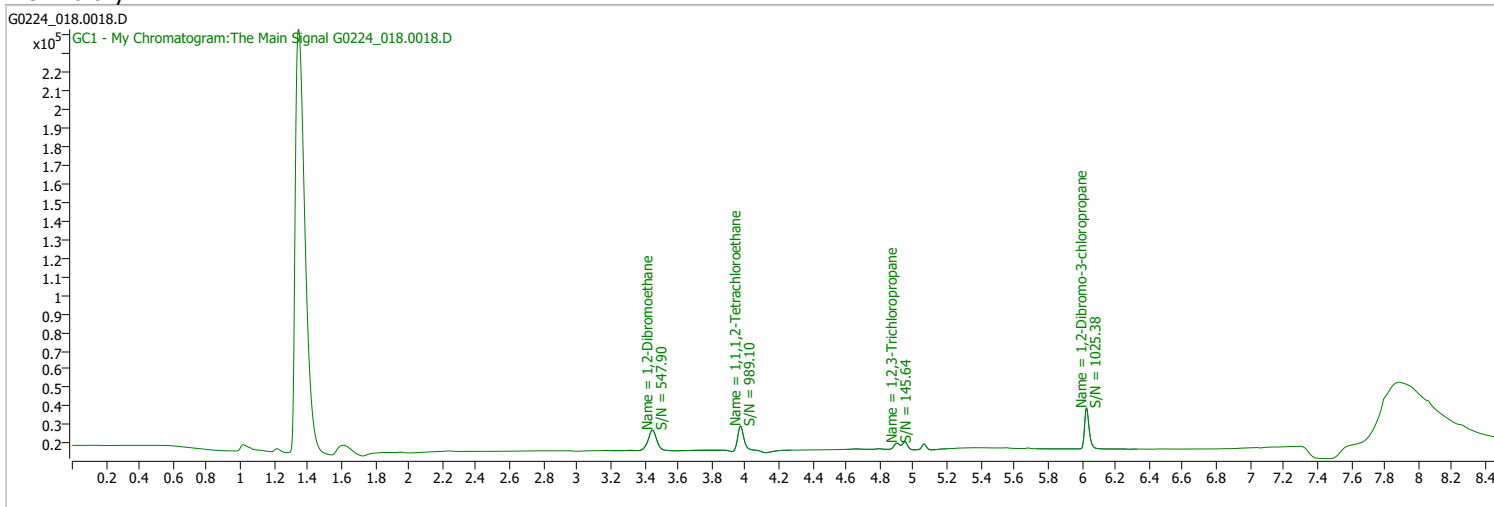
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_018.0018.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/24/2022 5:25:35 PM
Sample Name	LCS-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

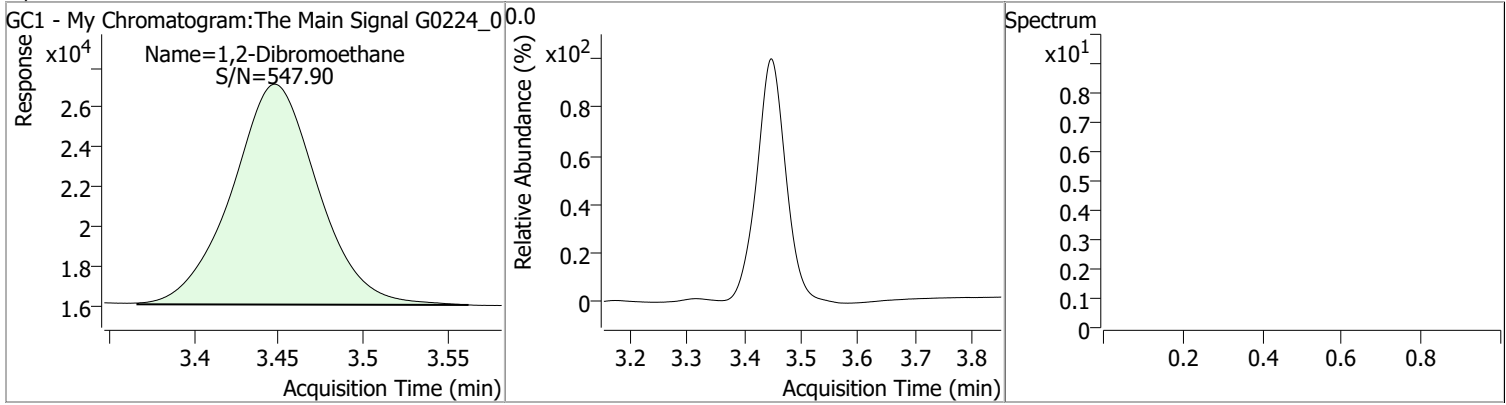


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.971	0.0	31825	0.0960	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%		Recovery = 96.00%			
Target Compounds						
M 1,2-Dibromoethane	3.448	0.0	39703	0.2430	µg/L	QValue 100

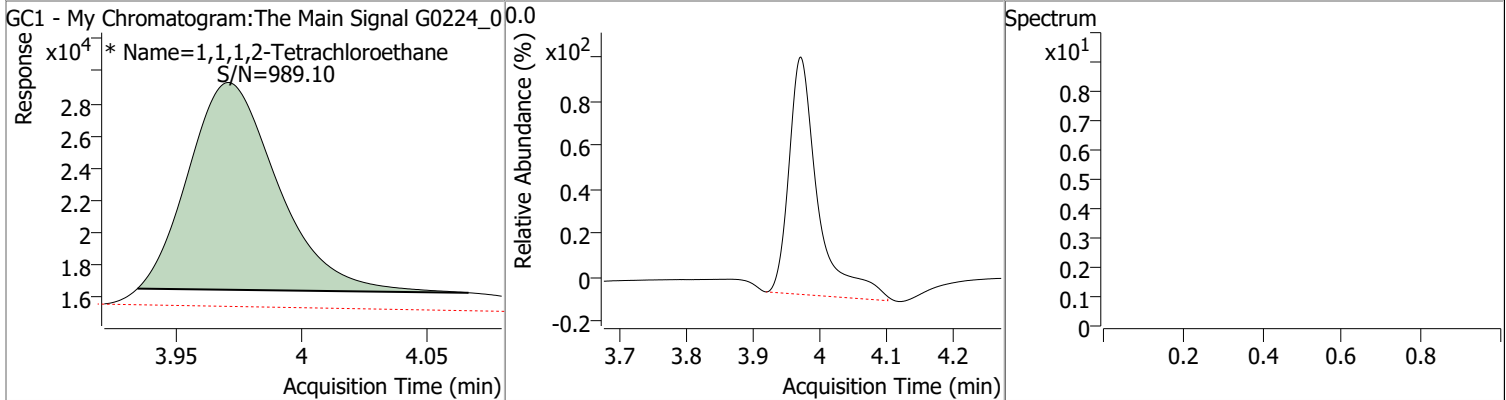
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2430	3.45	-0.01	39703				



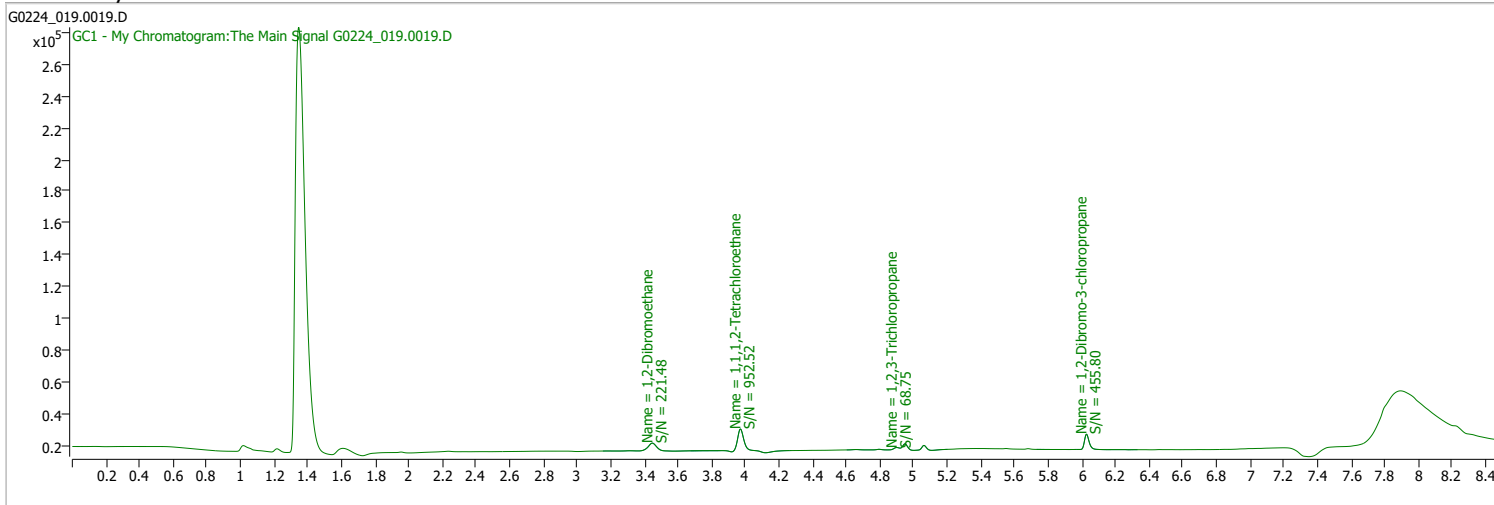
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0960	3.97	0.00	31825 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_019.0019.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/24/2022 5:45:21 PM
Sample Name	LCS1-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

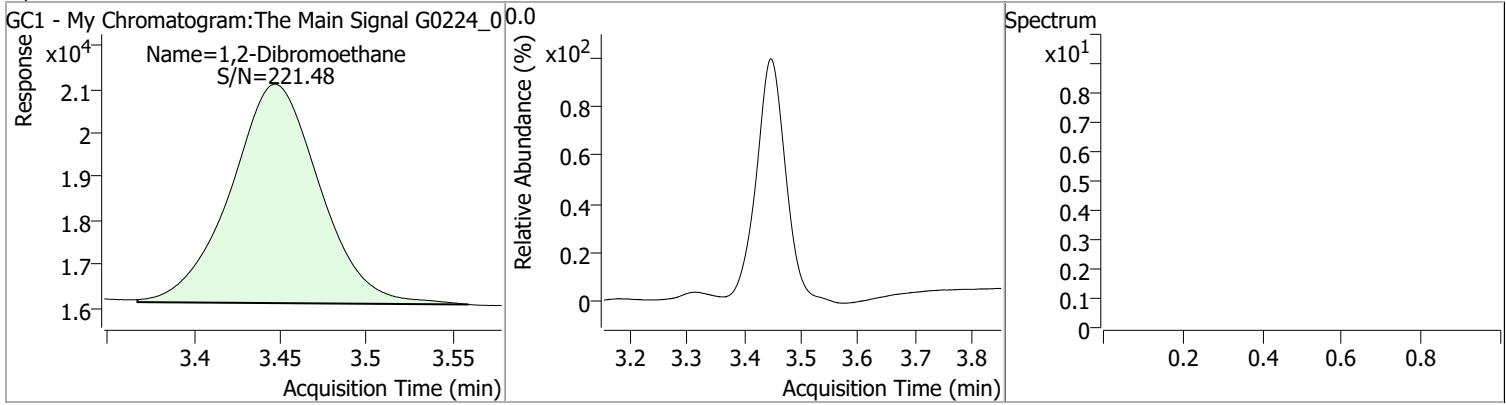


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.971	0.0	34064	0.1022	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 102.24%		
Target Compounds						
M 1,2-Dibromoethane	3.448	0.0	17890	0.1071	µg/L	100

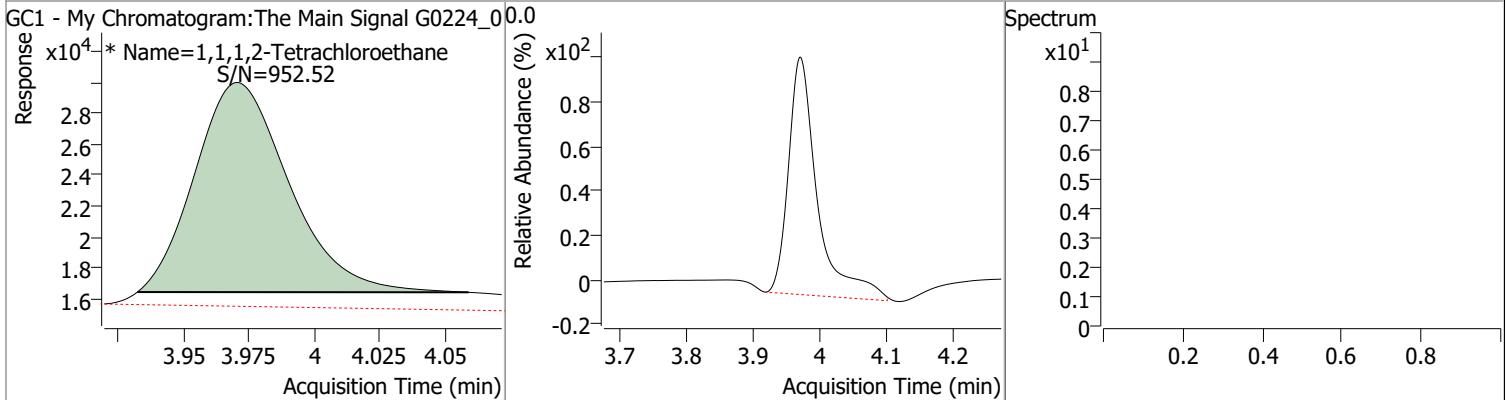
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1071	3.45	-0.01	17890				



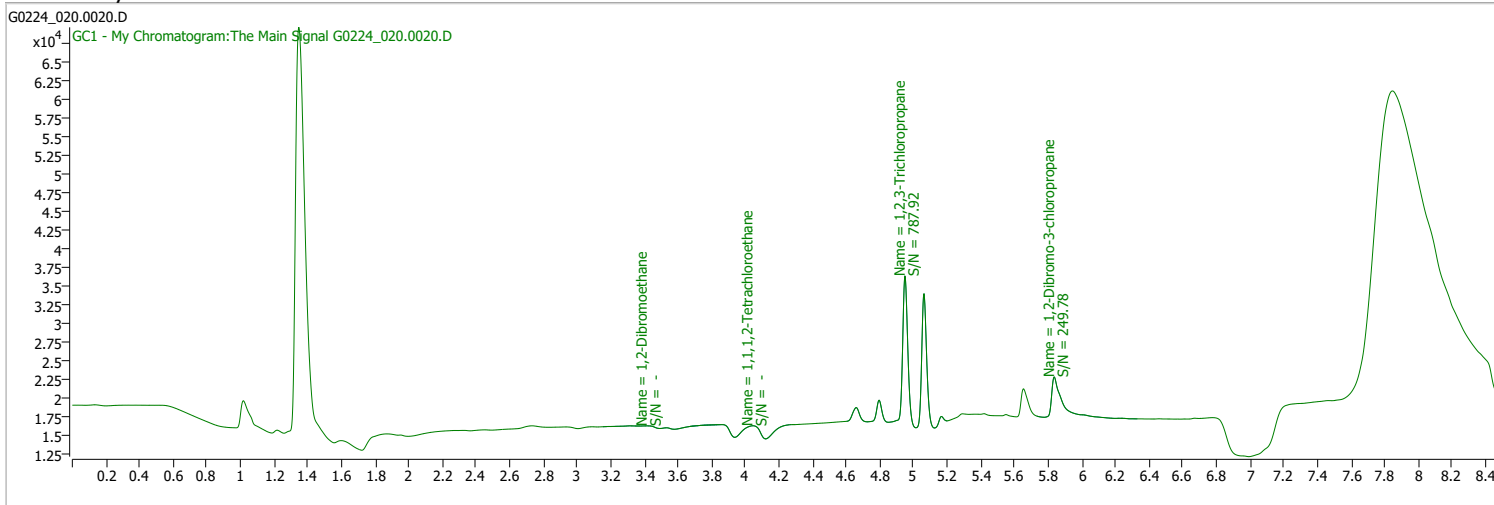
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1022	3.97	0.00	34064 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_020.0020.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 6:05:13 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

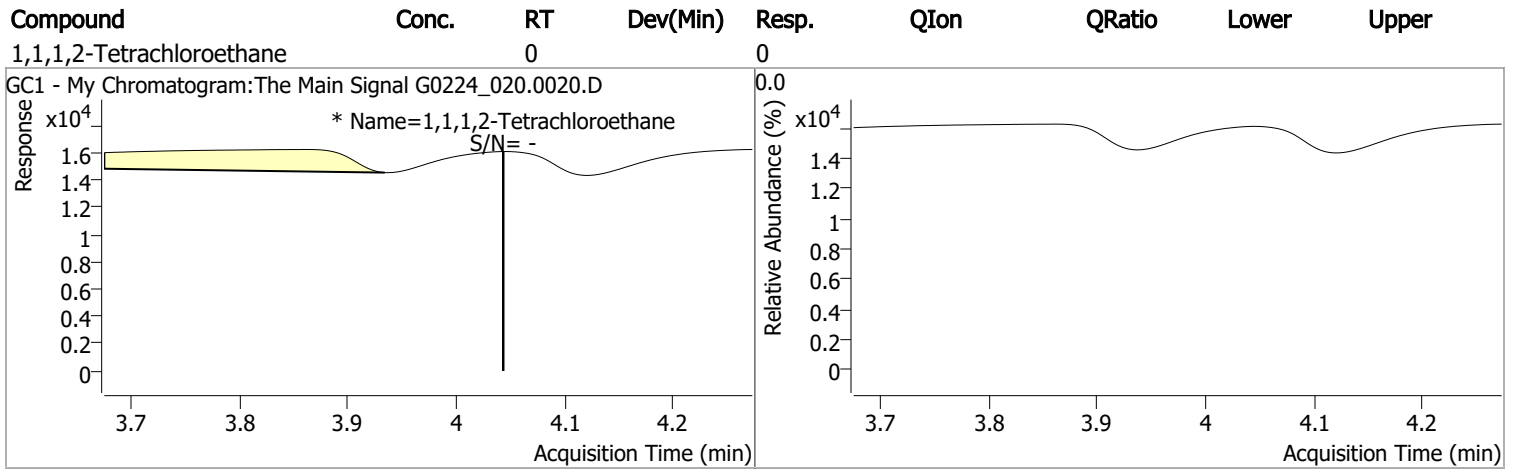
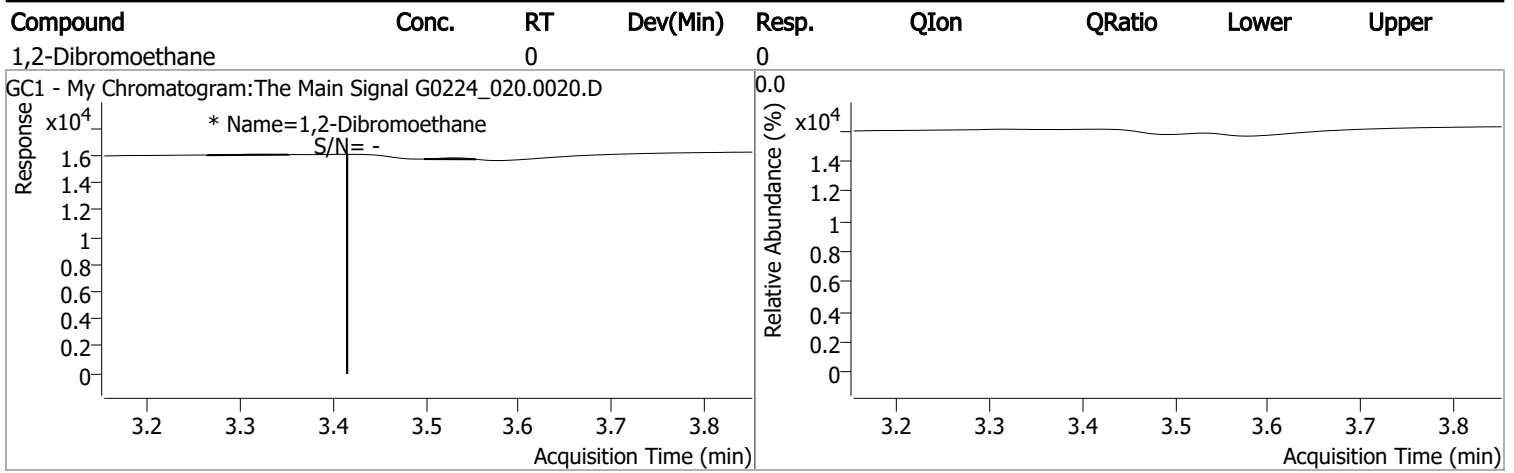
S 1,1,1,2-Tetrachloroethane	4.043	0.0	0		µg/L	md	0.069
Spiked Amount: 0.100	Range: 70.0 - 130.0%				Recovery = NA%		

Target Compounds

M 1,2-Dibromoethane	3.415	0.0	0		µg/L	md	QValue 1
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(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

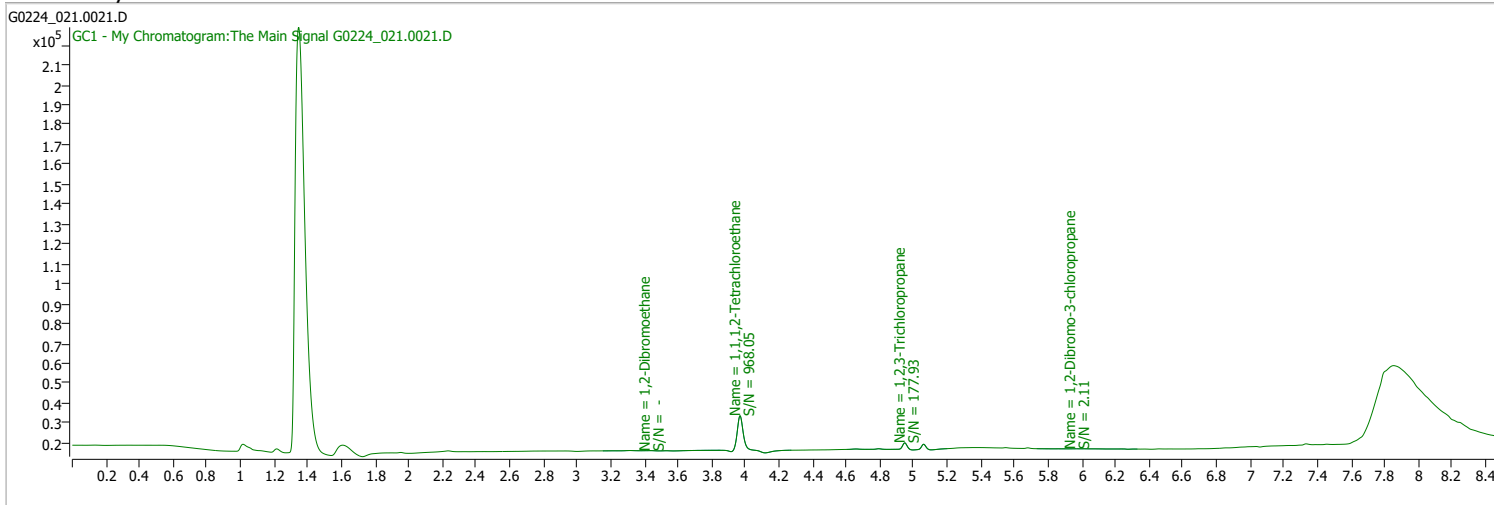
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_021.0021.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 6:24:52 PM
Sample Name	B22010745-008A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

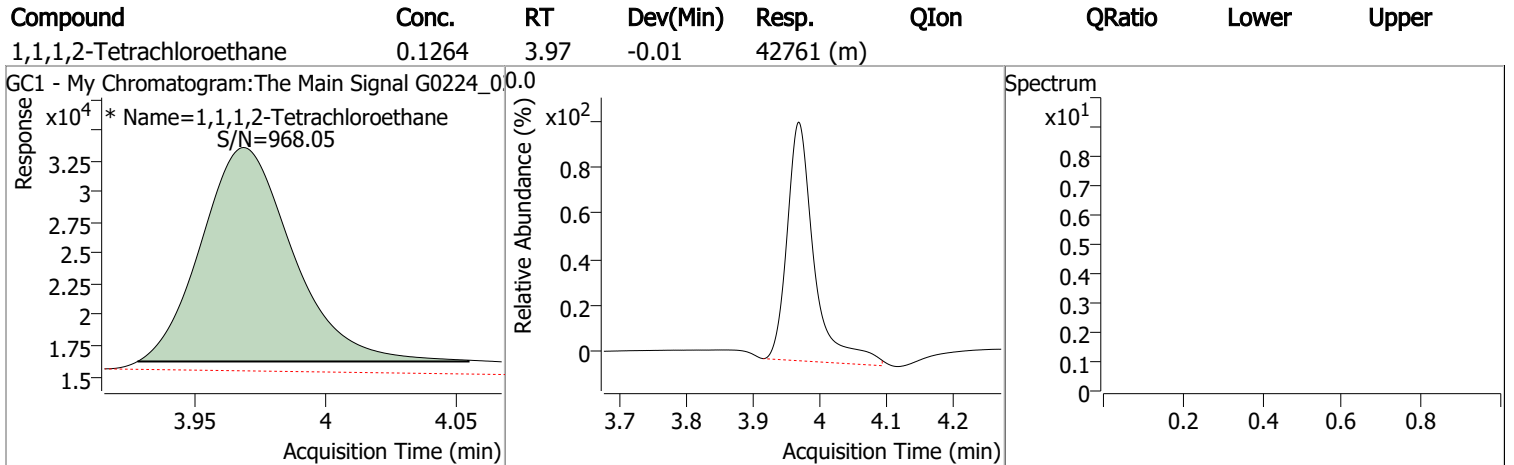
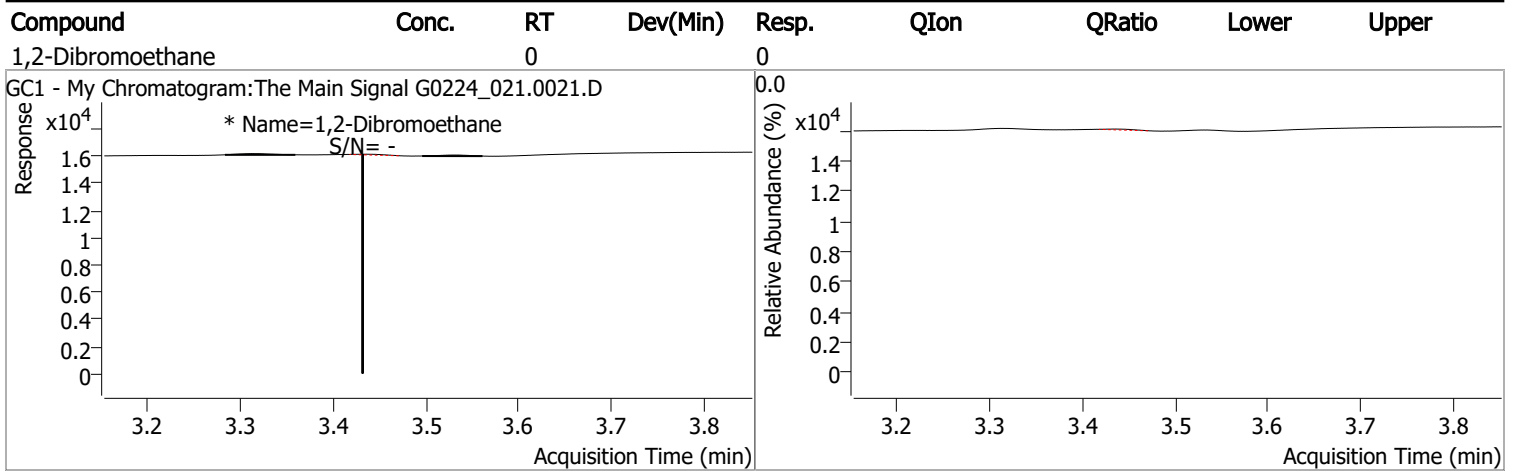
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	42761	0.1264	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 126.40%		
Target Compounds						
M 1,2-Dibromoethane	3.432	0.0	0		µg/L	md
						QValue
						1

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

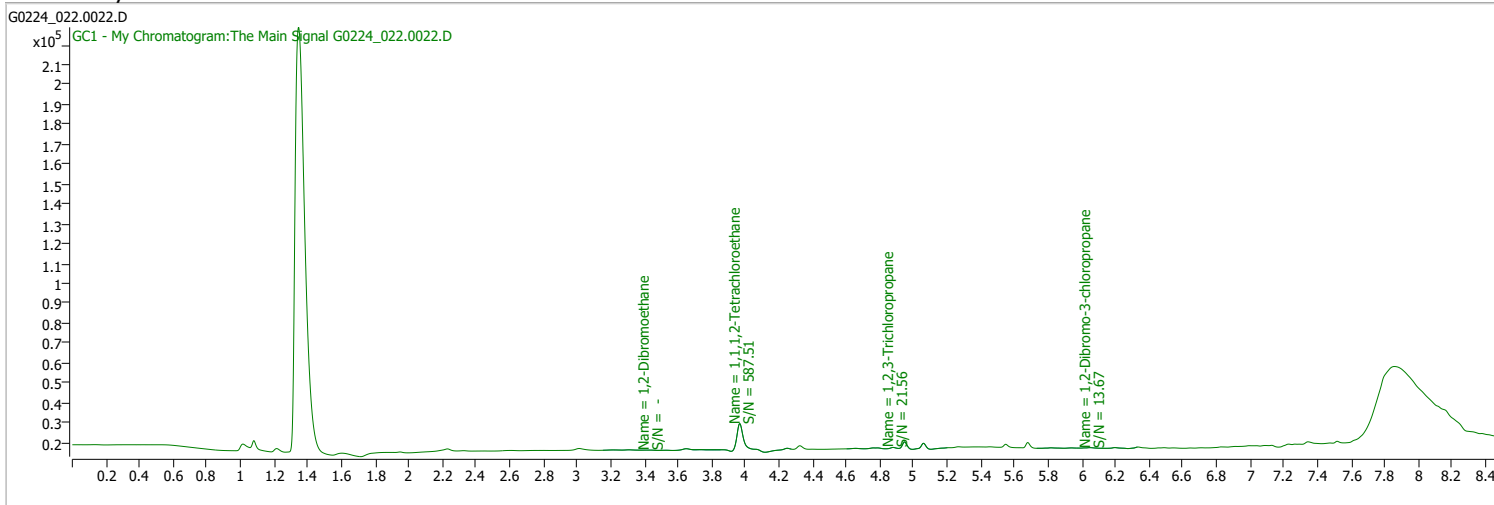
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_022.0022.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 6:44:41 PM
Sample Name	B22021435-004A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

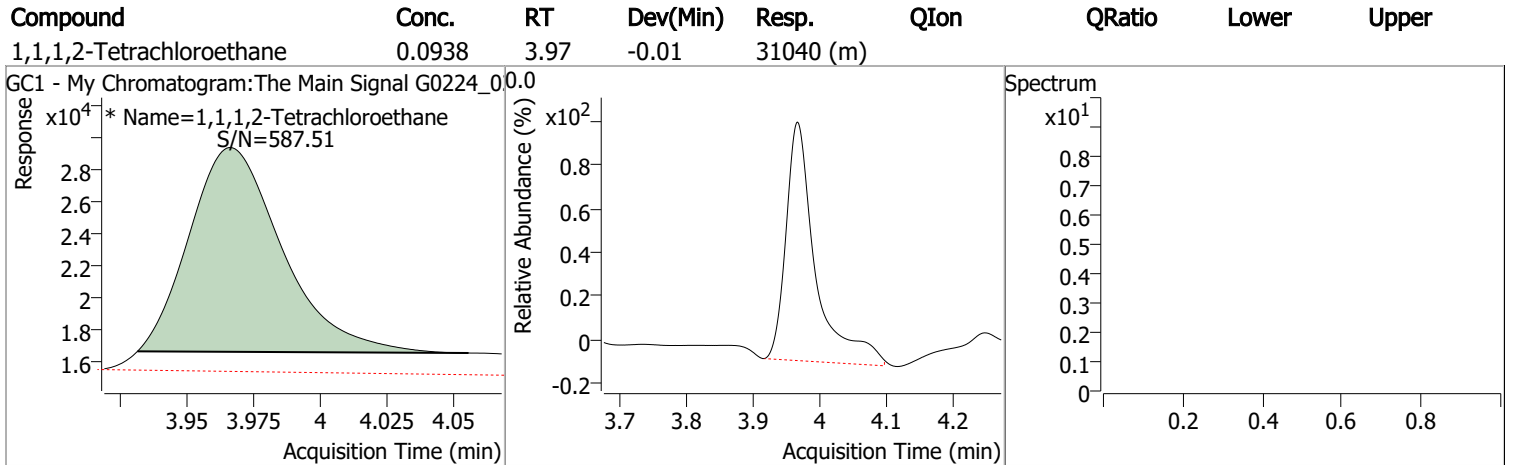
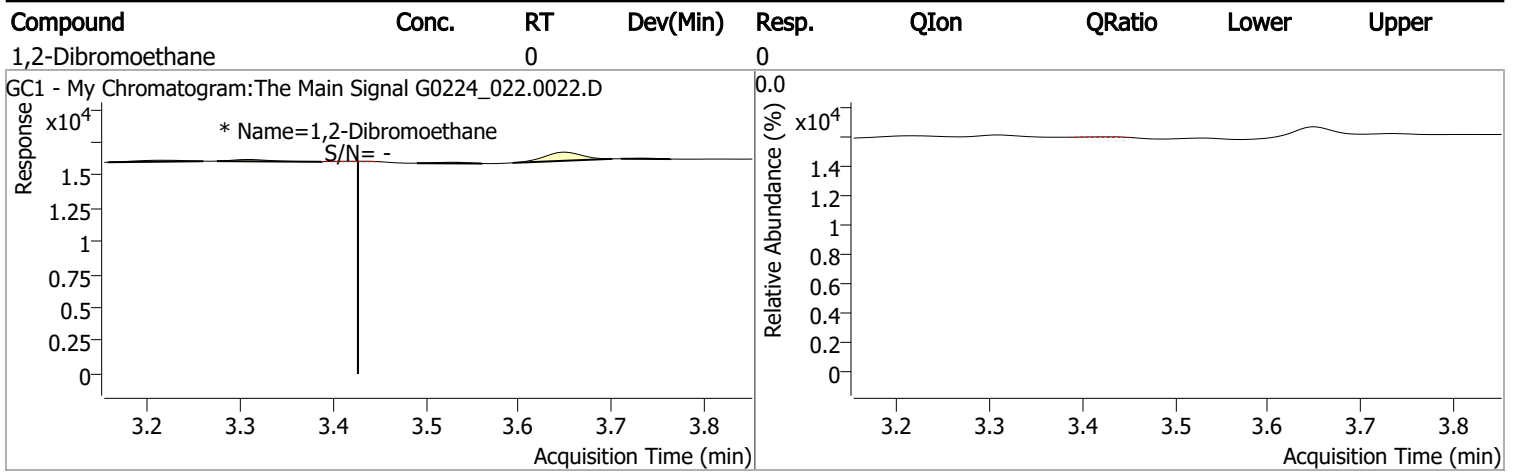
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.967	0.0	31040	0.0938	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 93.81%		
Target Compounds						
M 1,2-Dibromoethane	3.427	0.0	0		µg/L	md
						QValue
						1

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

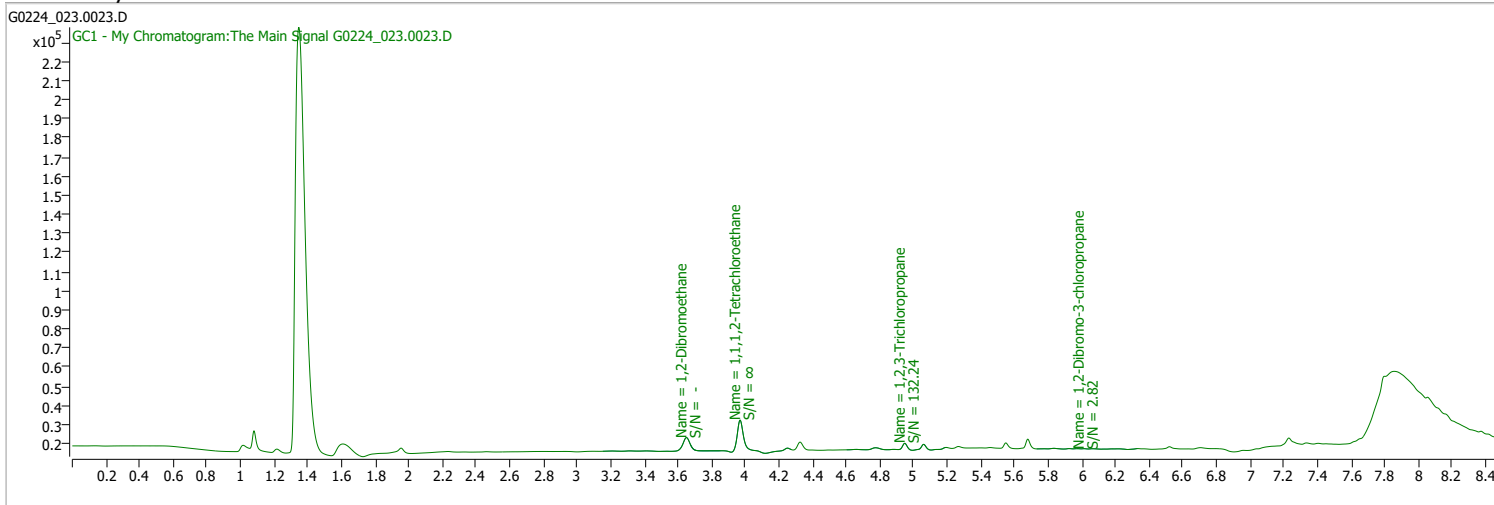
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_023.0023.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 7:04:37 PM
Sample Name	B22021435-006H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

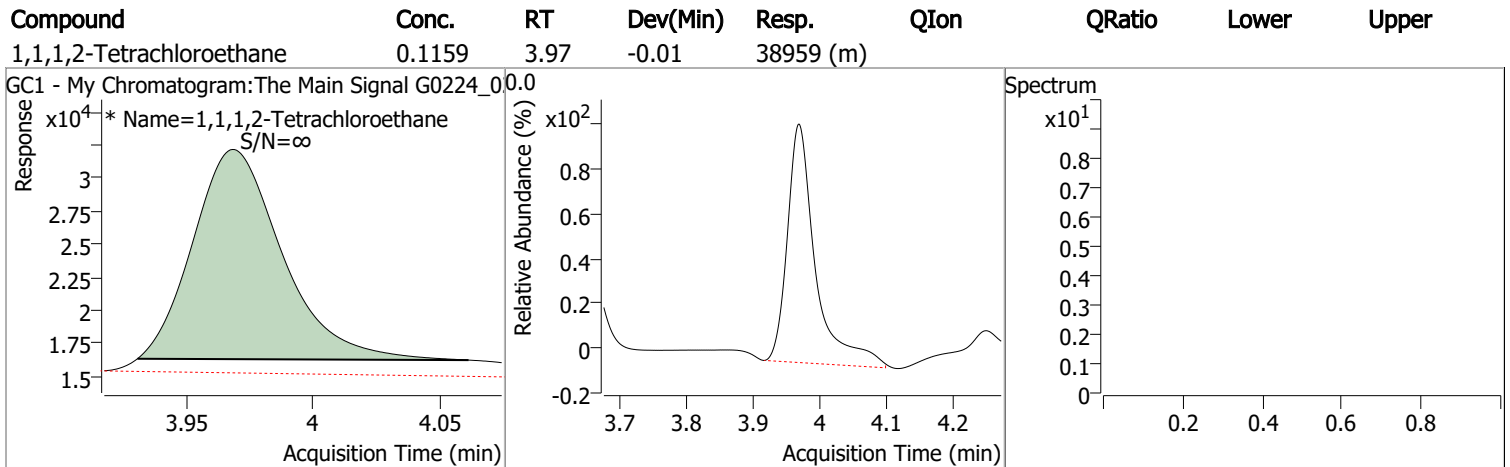
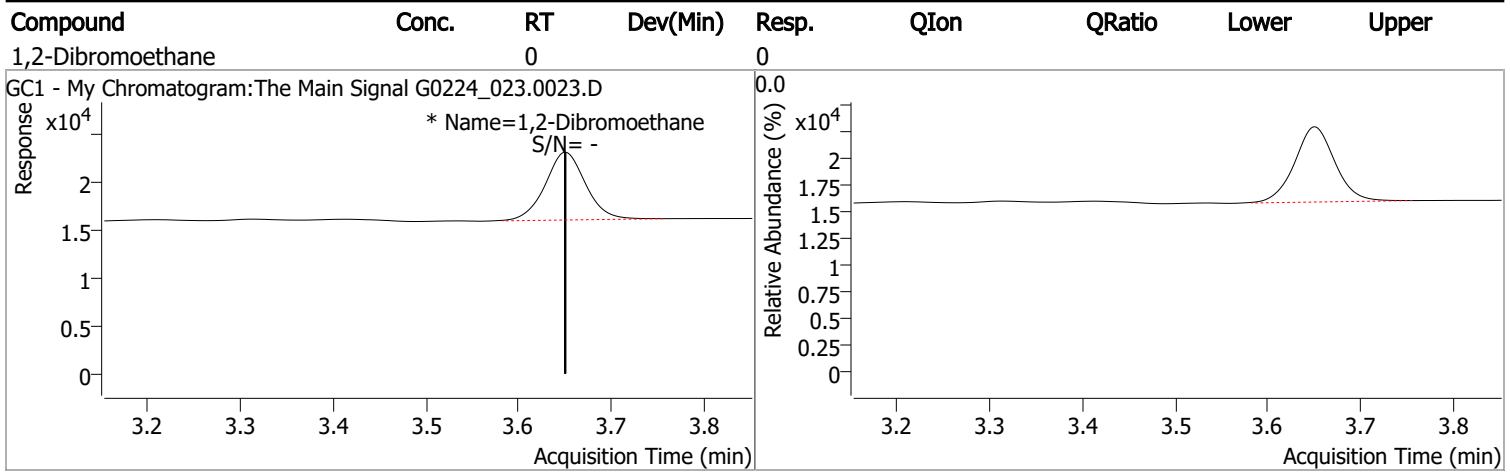
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	38959	0.1159	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 115.85%		
Target Compounds						
M 1,2-Dibromoethane	3.650	0.0	0		µg/L	md
						QValue 1

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

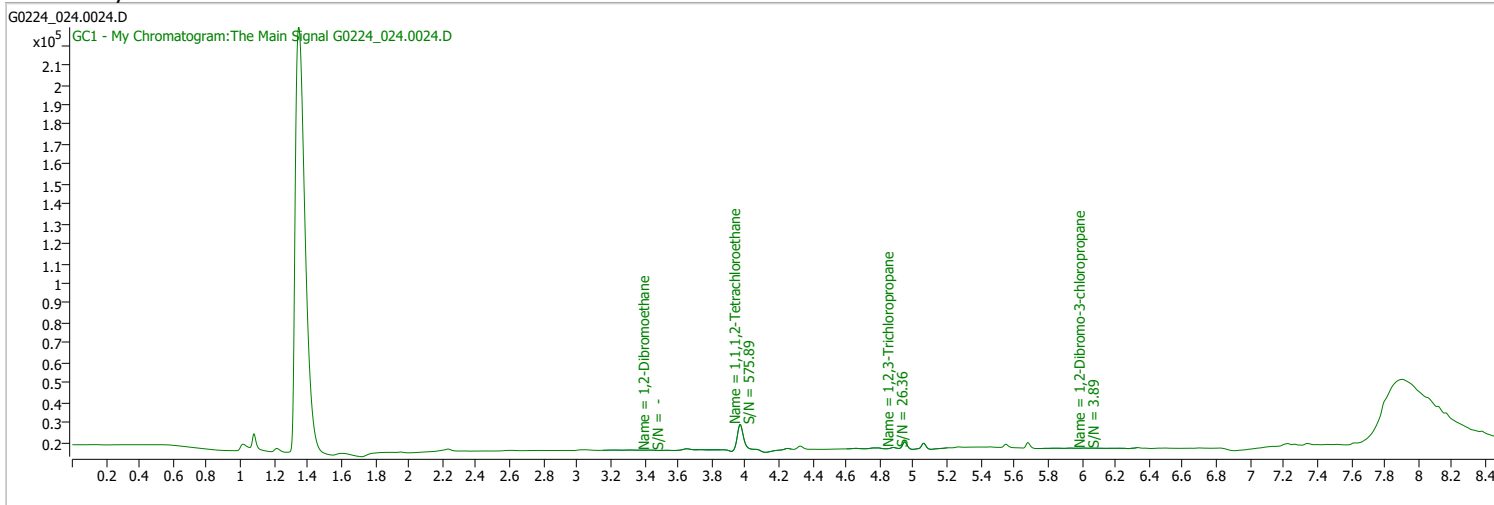
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_024.0024.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 7:24:14 PM
Sample Name	B22021435-010A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

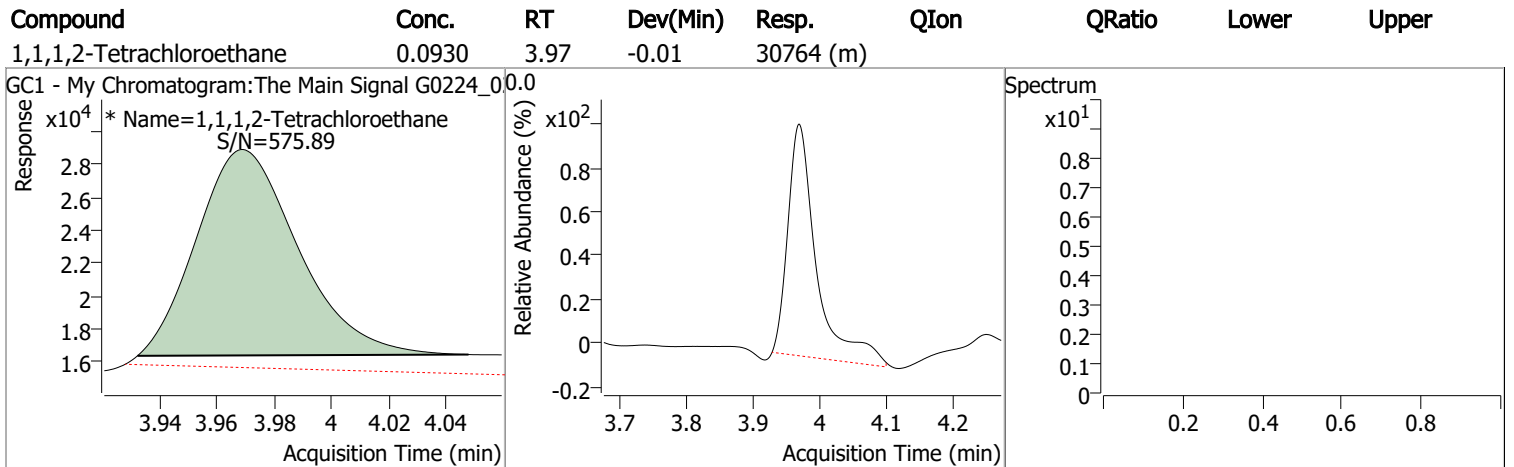
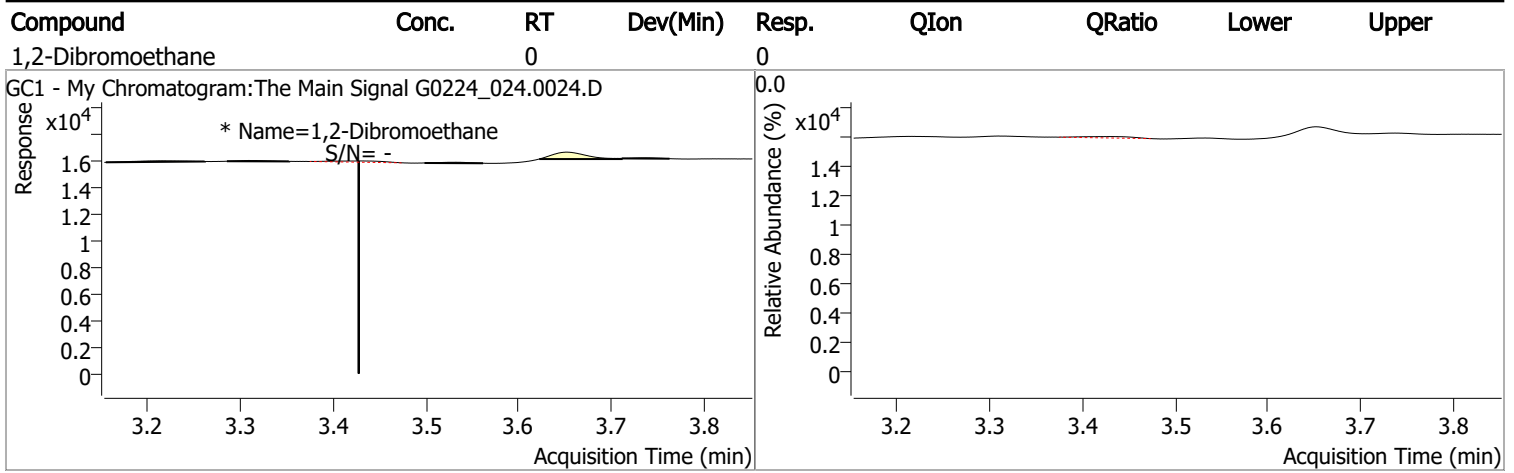
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.969	0.0	30764	0.0930	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 93.04%		
Target Compounds						
M 1,2-Dibromoethane	3.428	0.0	0		µg/L	md
						QValue 1

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

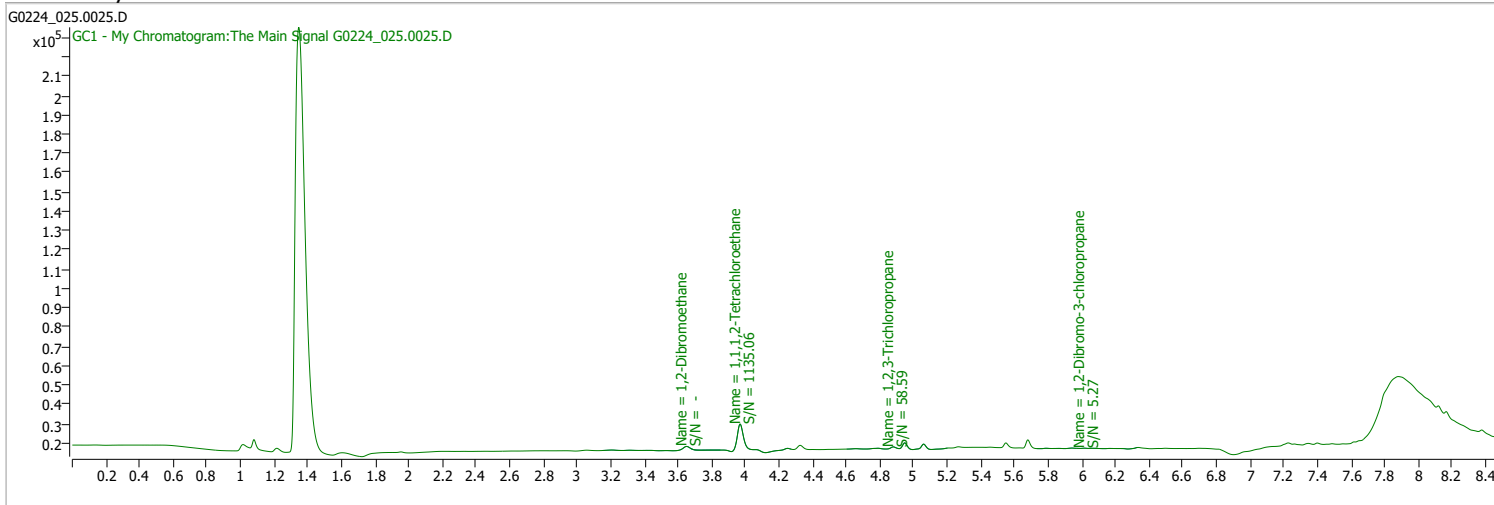
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_025.0025.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 7:44:02 PM
Sample Name	B22021435-012H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

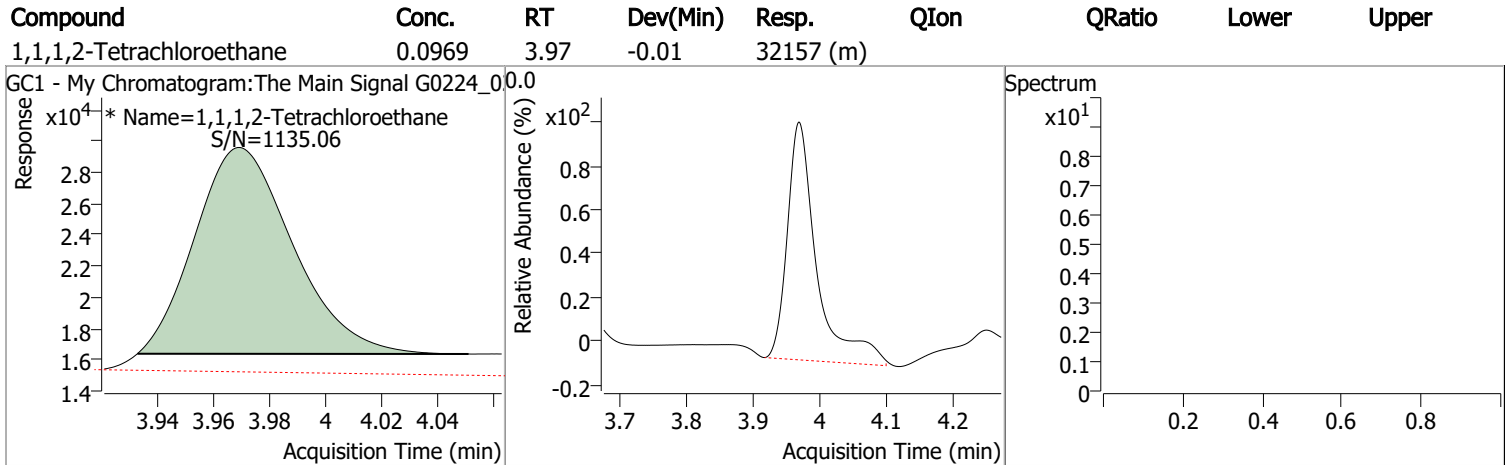
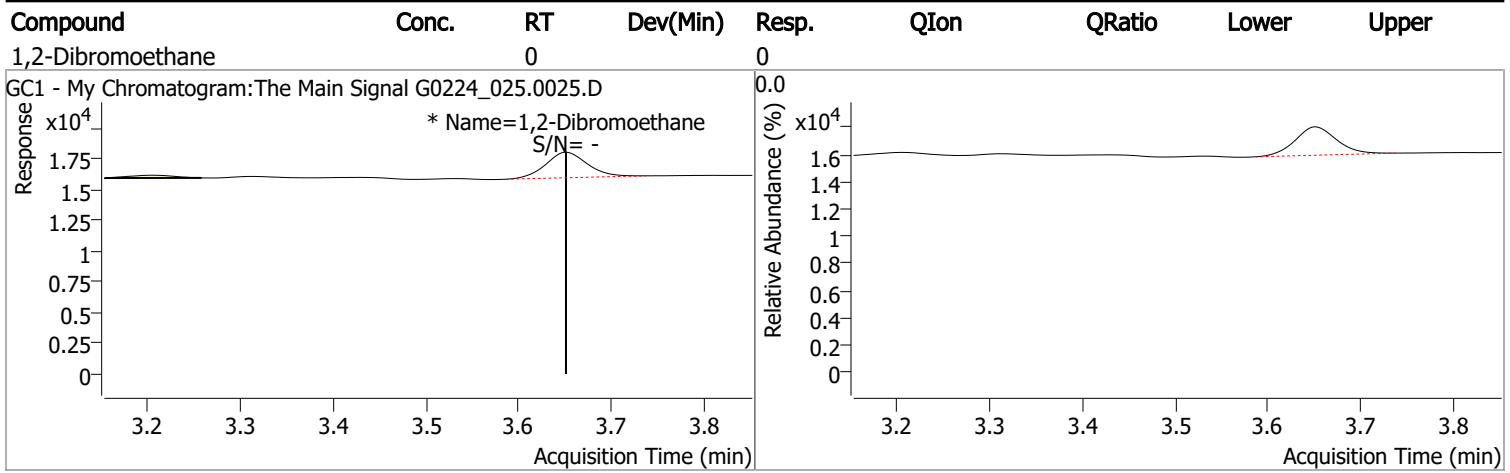
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	32157	0.0969	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 96.93%		
Target Compounds						
M 1,2-Dibromoethane	3.651	0.0	0		µg/L	md
						QValue 1

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

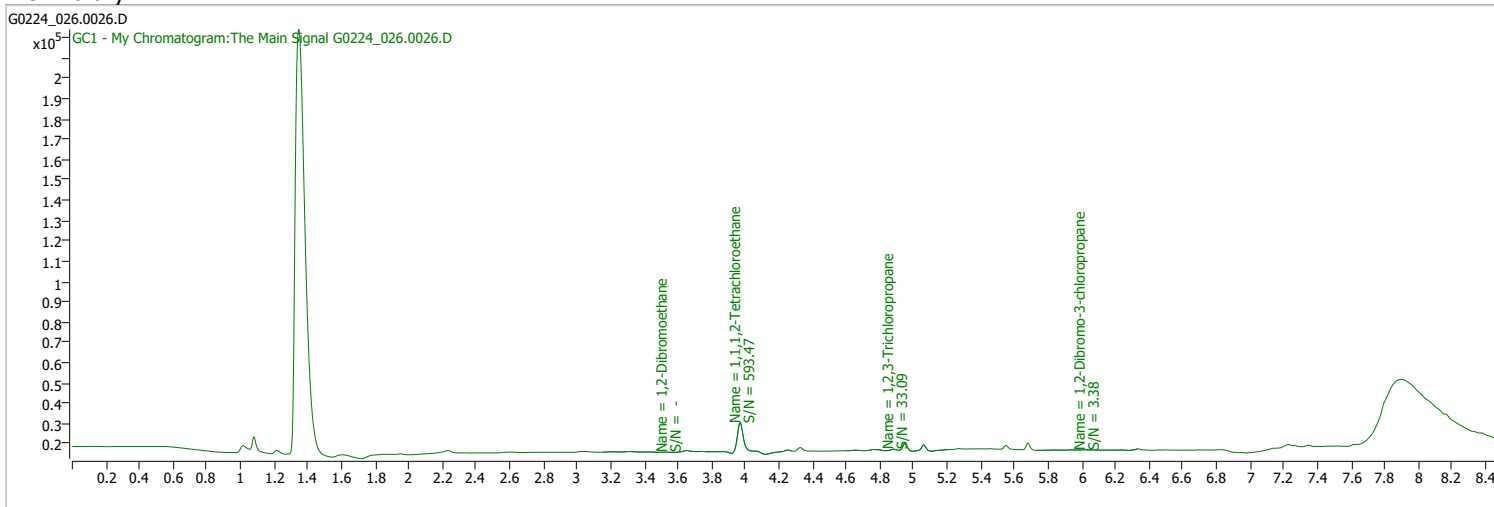
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_026.0026.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 8:03:38 PM
Sample Name	B22021435-015A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

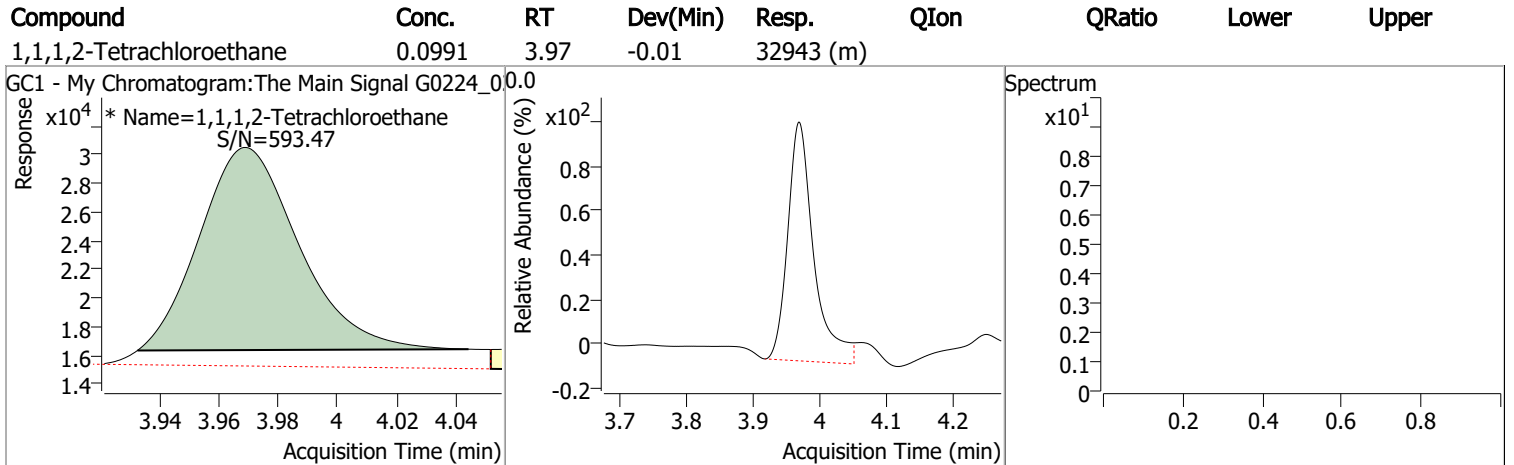
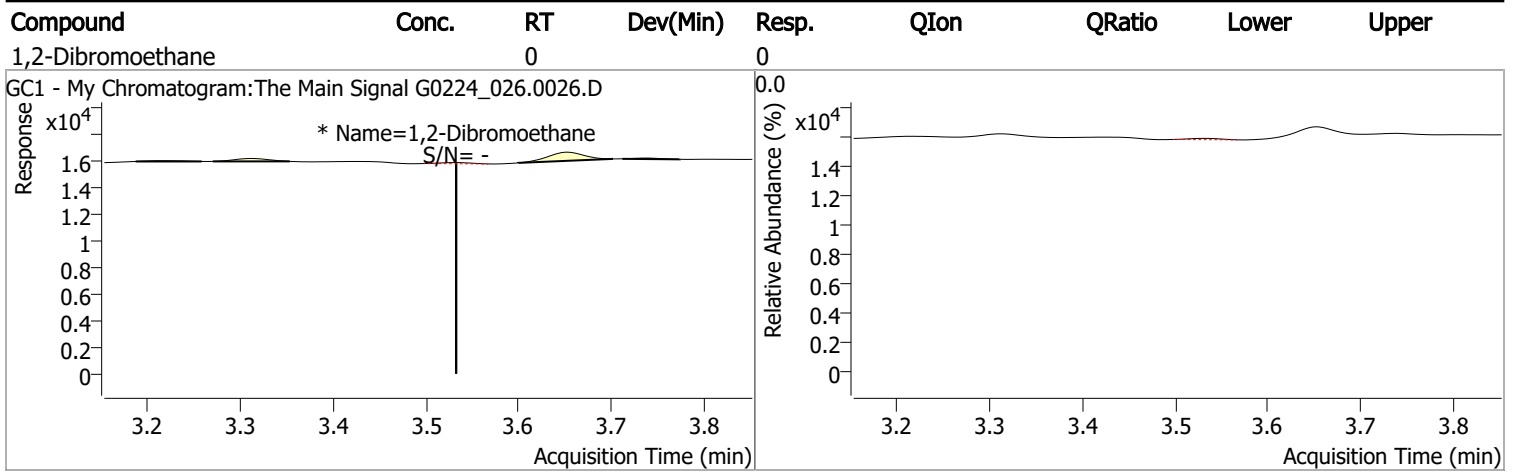
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	32943	0.0991	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%		Recovery = 99.12%			
Target Compounds						
M 1,2-Dibromoethane	3.533	0.0	0		µg/L	md
						QValue
						1

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

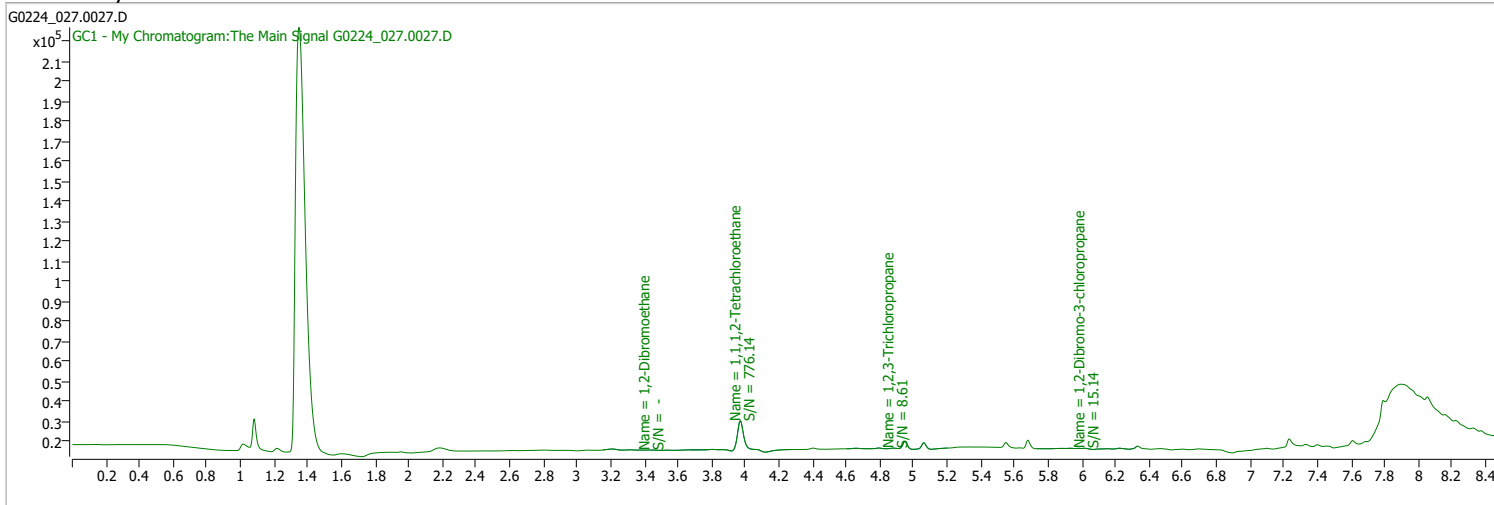
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_027.0027.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 8:23:35 PM
Sample Name	B22021435-017H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

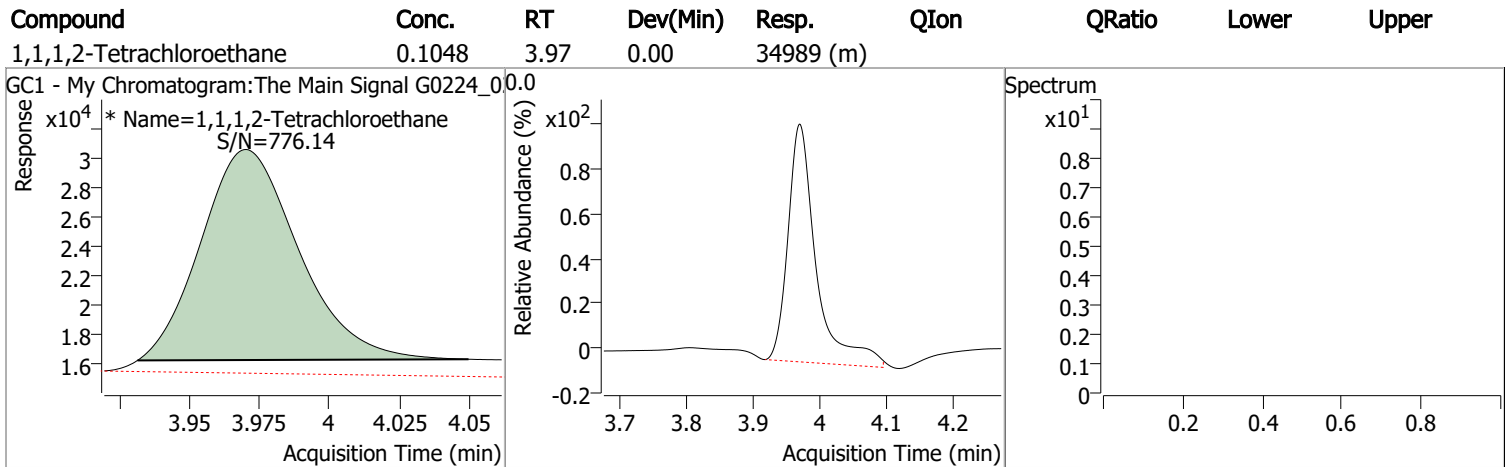
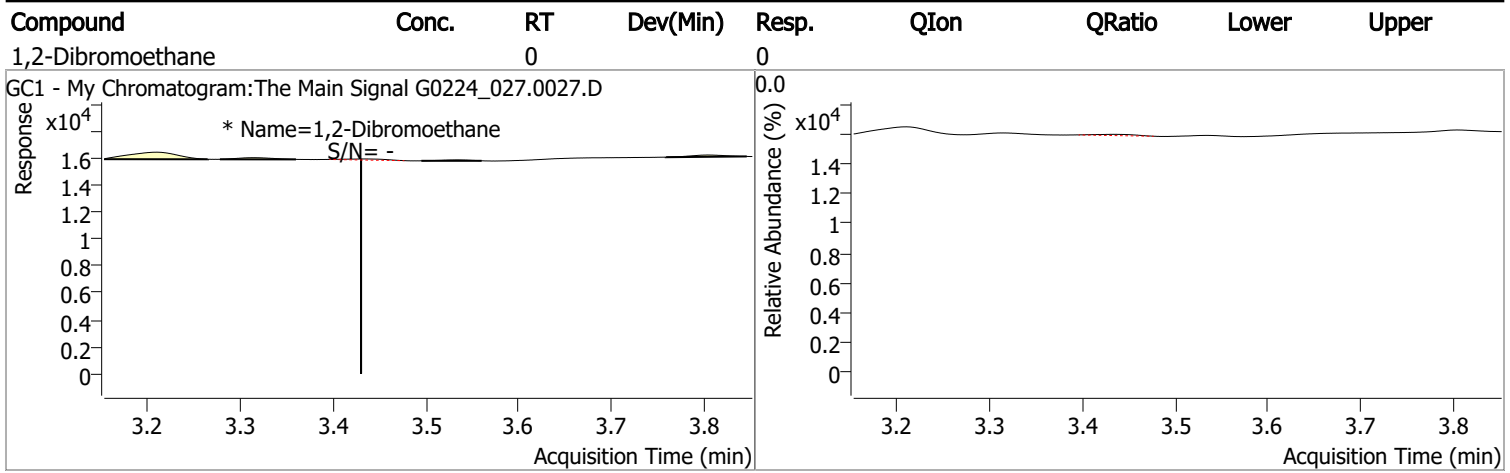
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	34989	0.1048	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 104.81%		
Target Compounds						
M 1,2-Dibromoethane	3.430	0.0	0		µg/L	md
						QValue 1

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

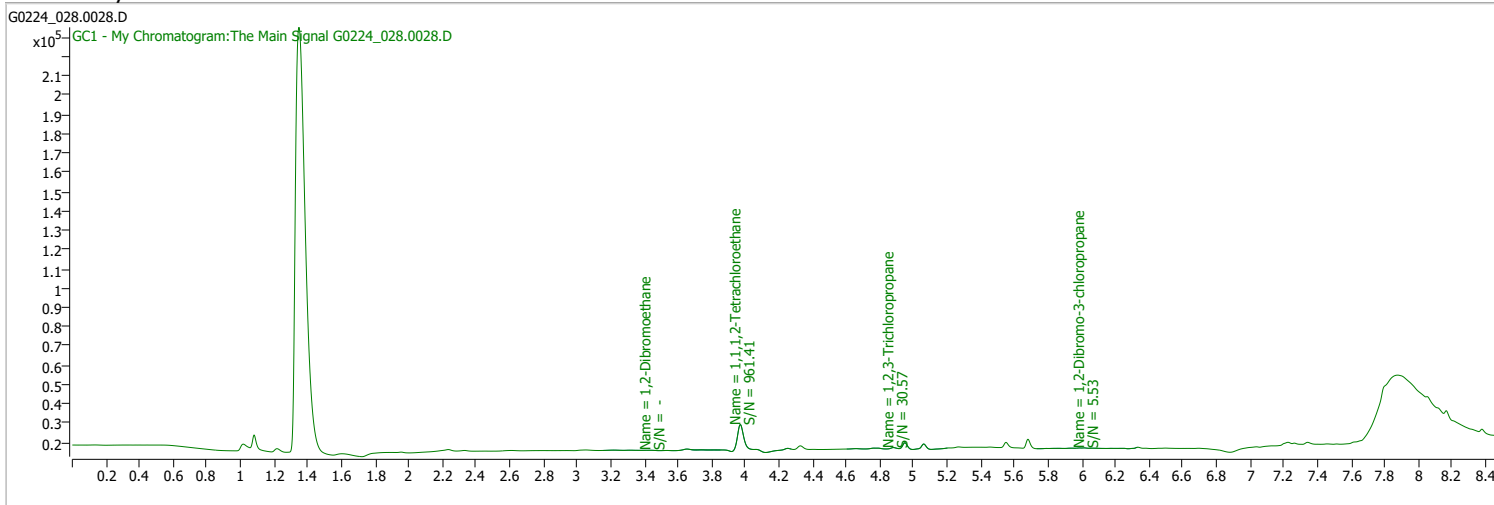
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_028.0028.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 8:43:24 PM
Sample Name	B22021435-020A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

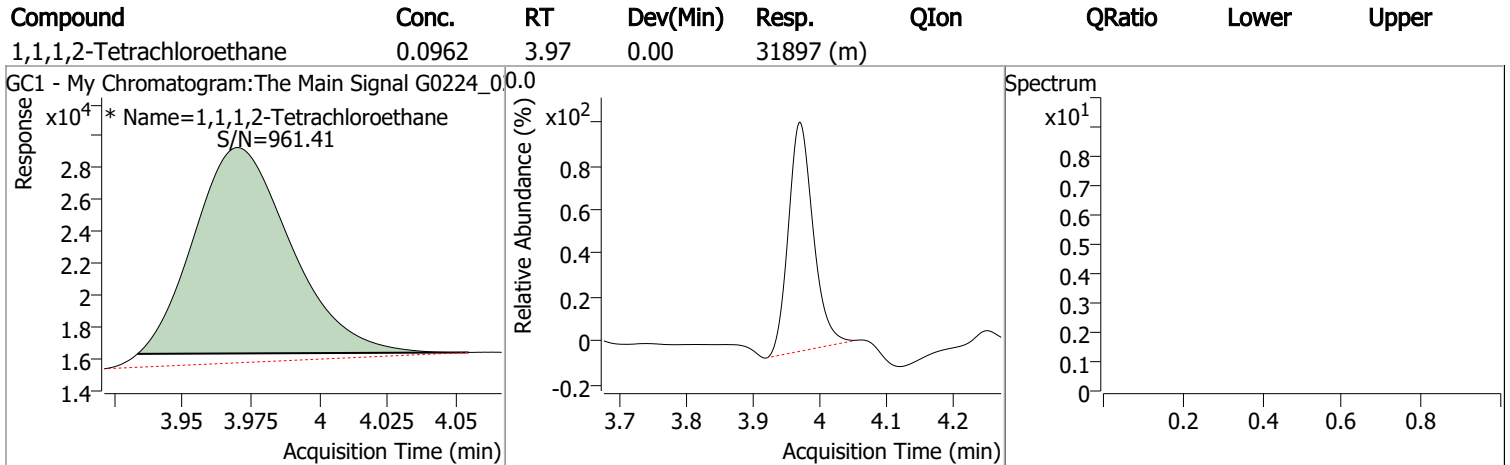
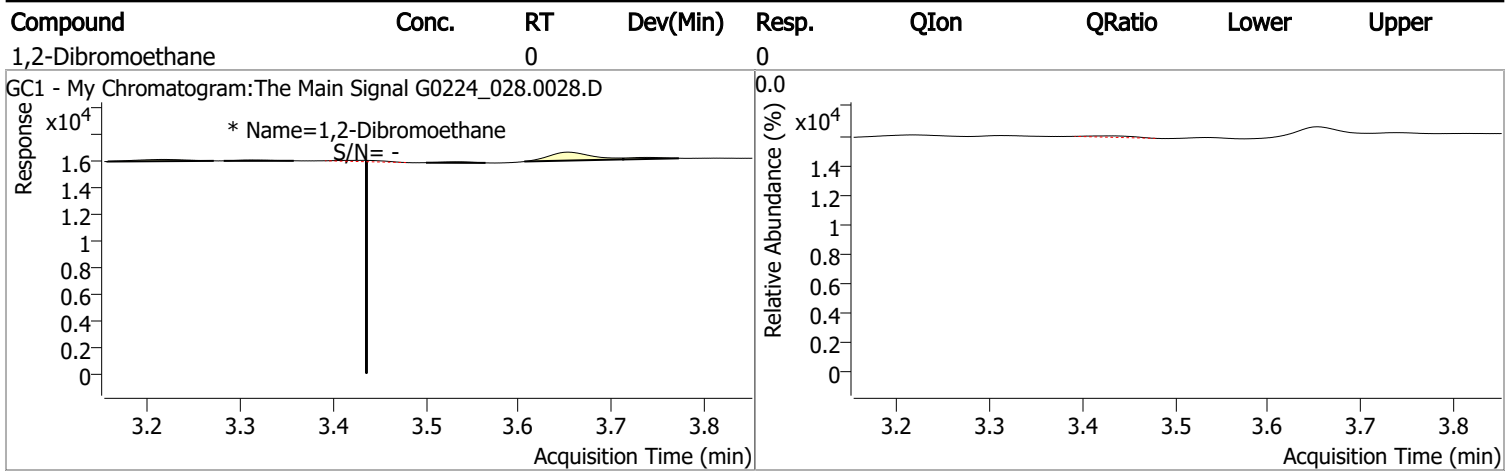
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	31897	0.0962	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 96.20%		
Target Compounds						
M 1,2-Dibromoethane	3.436	0.0	0		µg/L	md
						QValue
						1

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

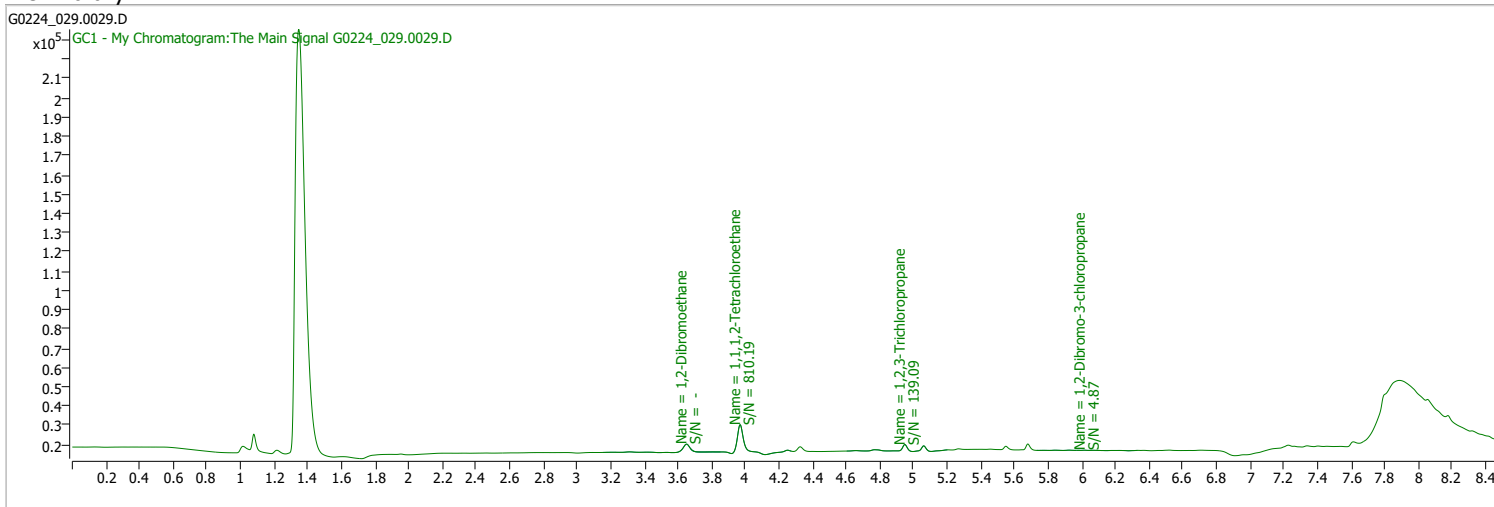
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_029.0029.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 9:02:58 PM
Sample Name	B22021435-001H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

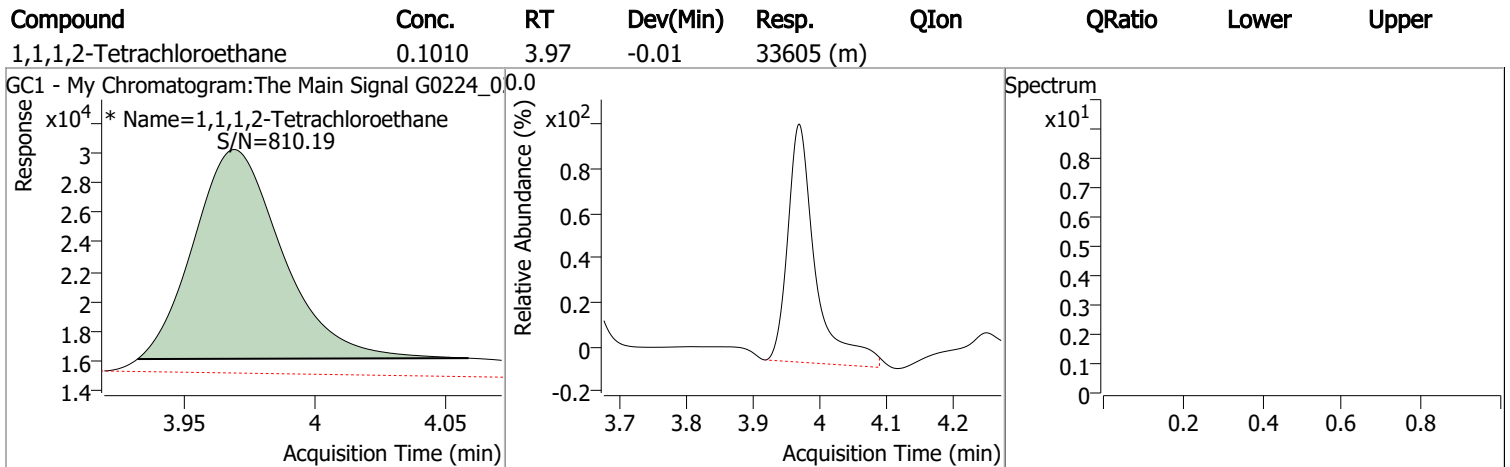
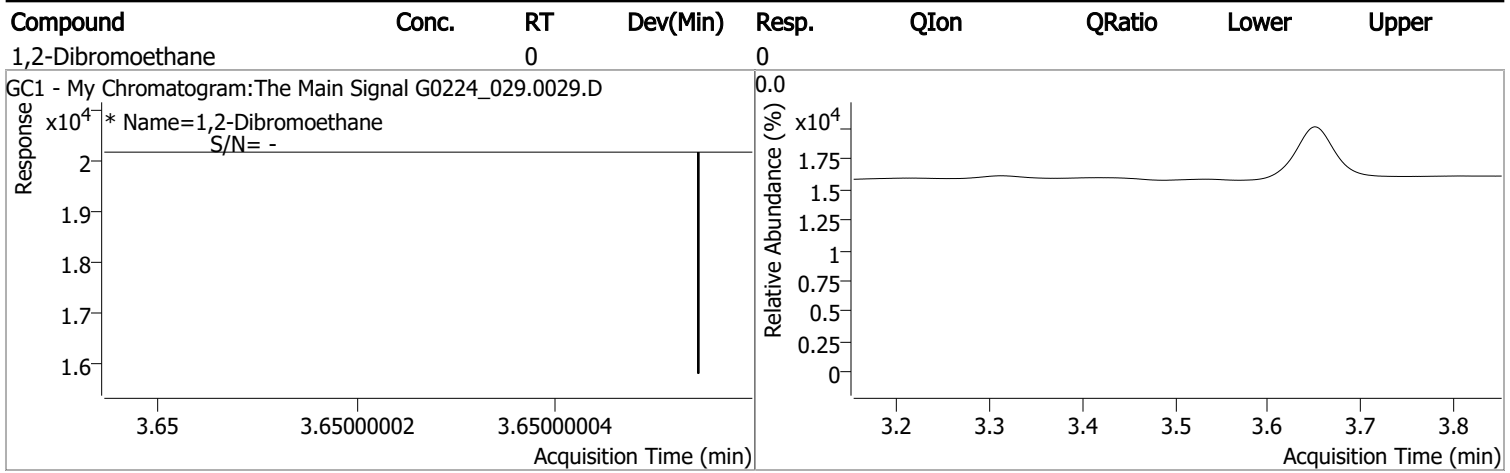
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.969	0.0	33605	0.1010	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 100.96%		
Target Compounds						
M 1,2-Dibromoethane	3.650	0.0	0		µg/L	md
						QValue
						1

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

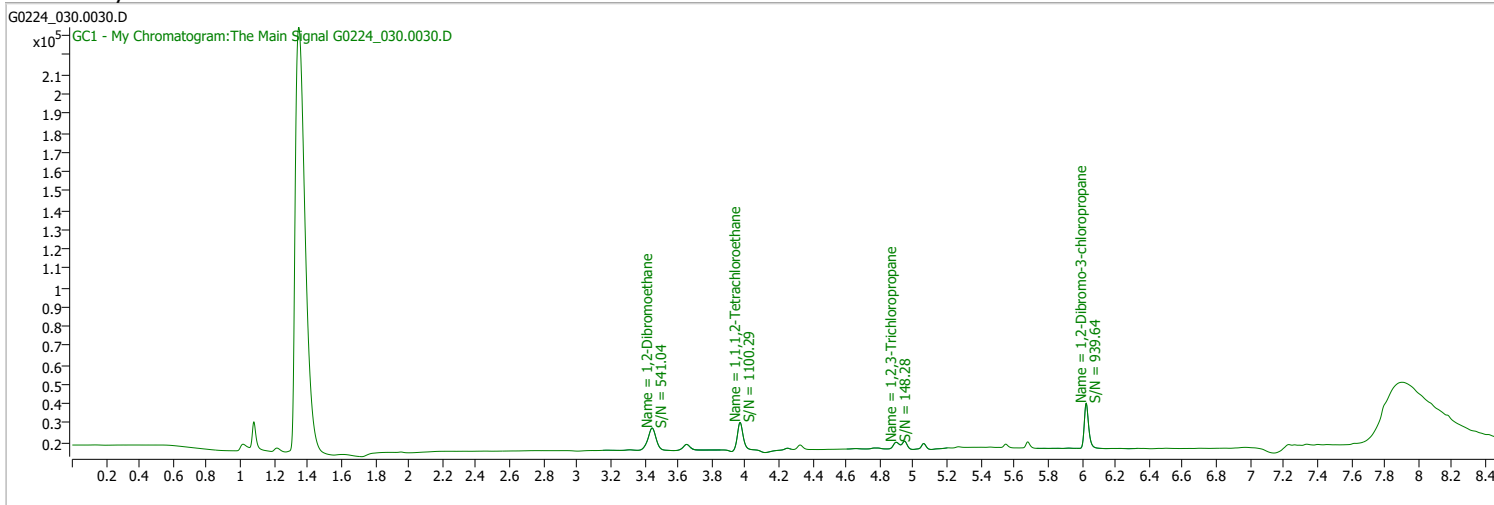
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_030.0030.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 9:22:50 PM
Sample Name	B22021435-001HMS	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

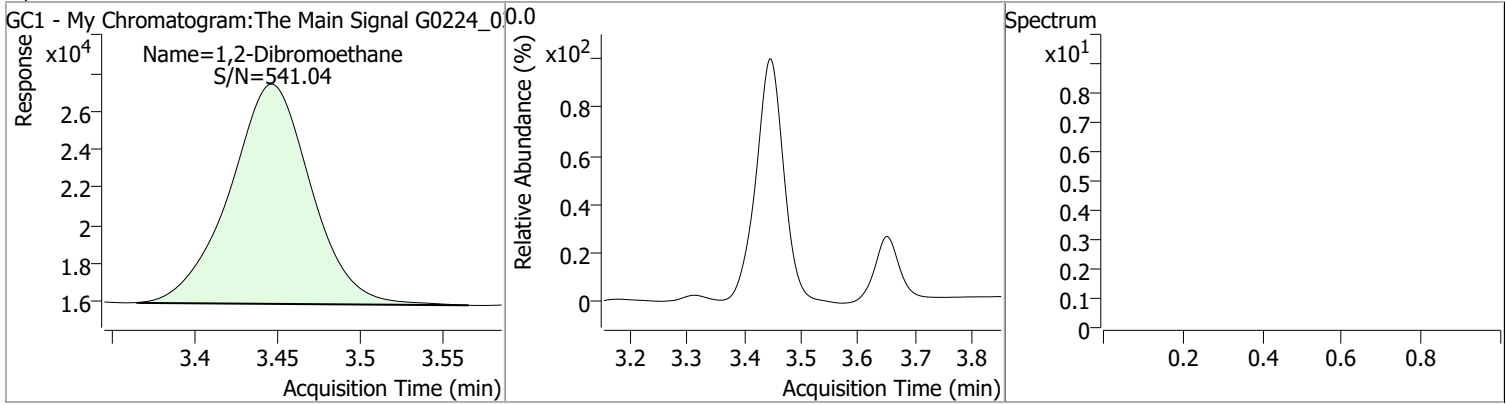


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	32008	0.0965	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 96.51%		
Target Compounds						
M 1,2-Dibromoethane	3.447	0.0	39497	0.2417	µg/L	100

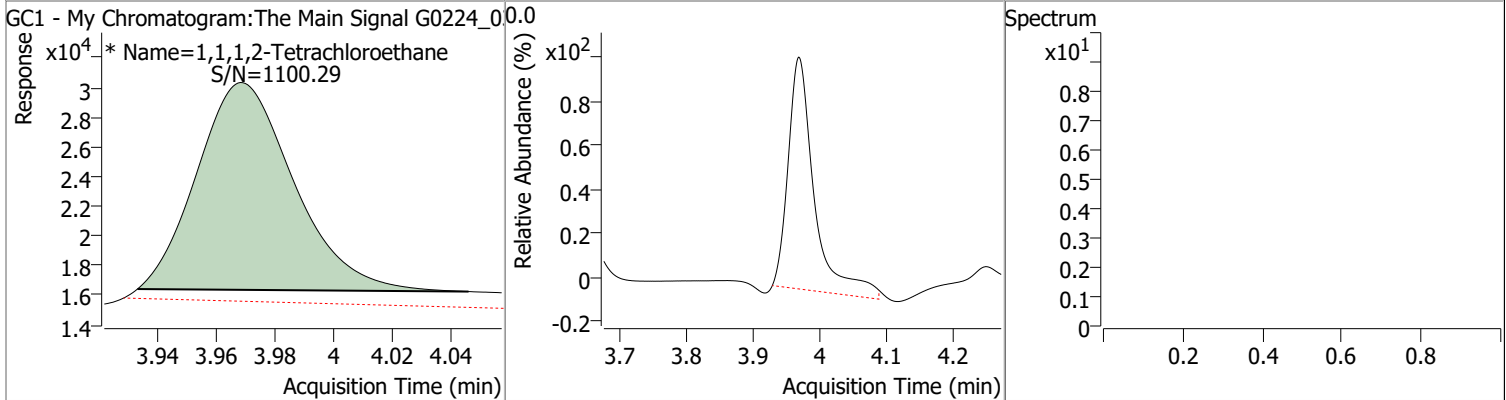
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2417	3.45	-0.01	39497				



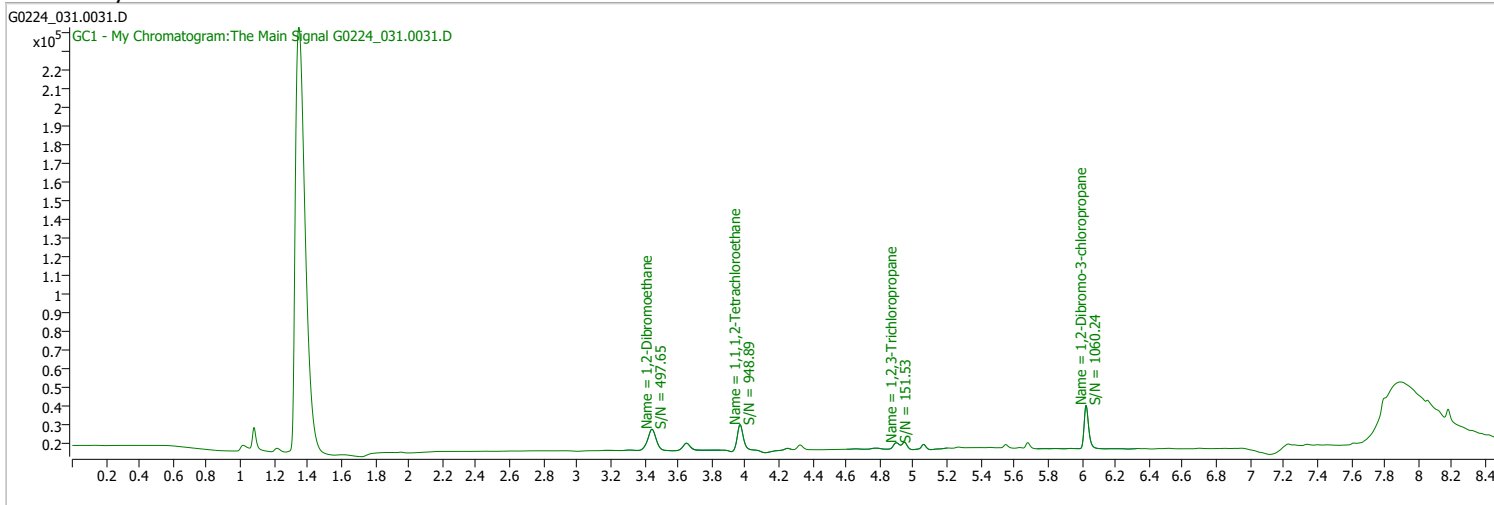
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0965	3.97	-0.01	32008 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_031.0031.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 9:42:32 PM
Sample Name	B22021435-001HMSD	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

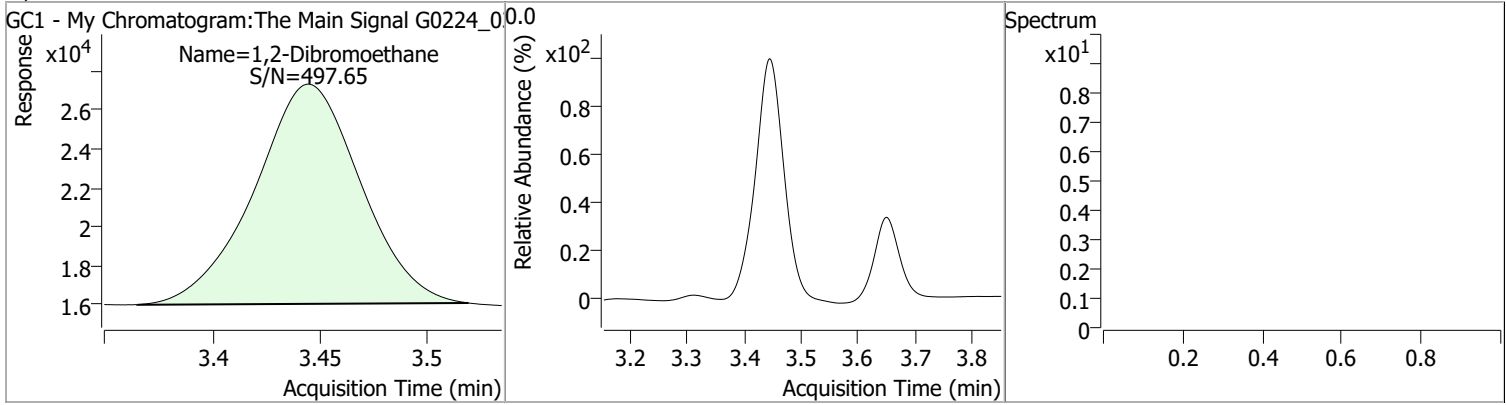


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	32849	0.0989	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 98.85%		
Target Compounds						
M 1,2-Dibromoethane	3.445	0.0	38858	0.2376	µg/L	100

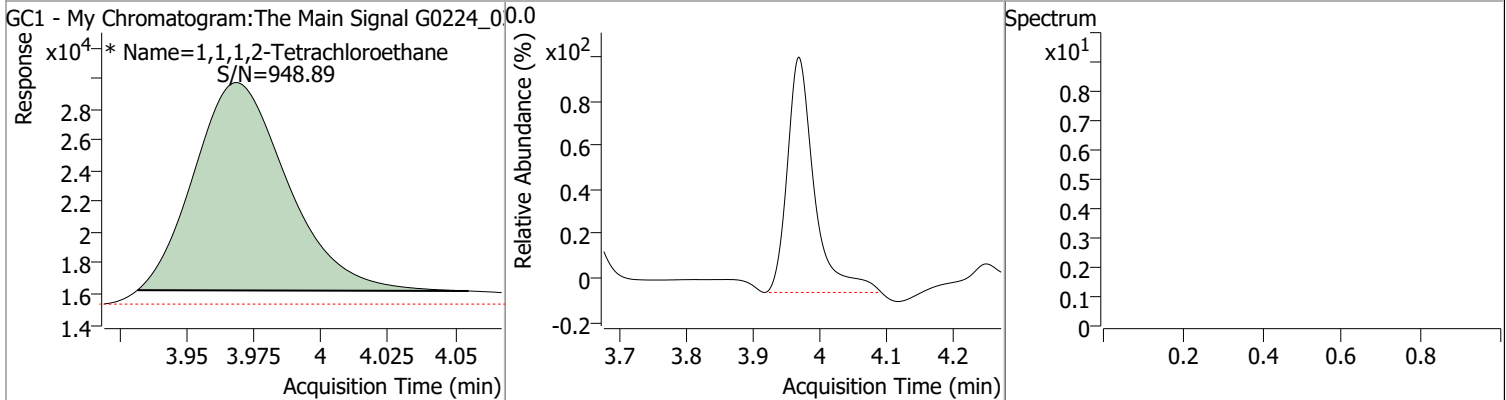
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2376	3.45	-0.01	38858				



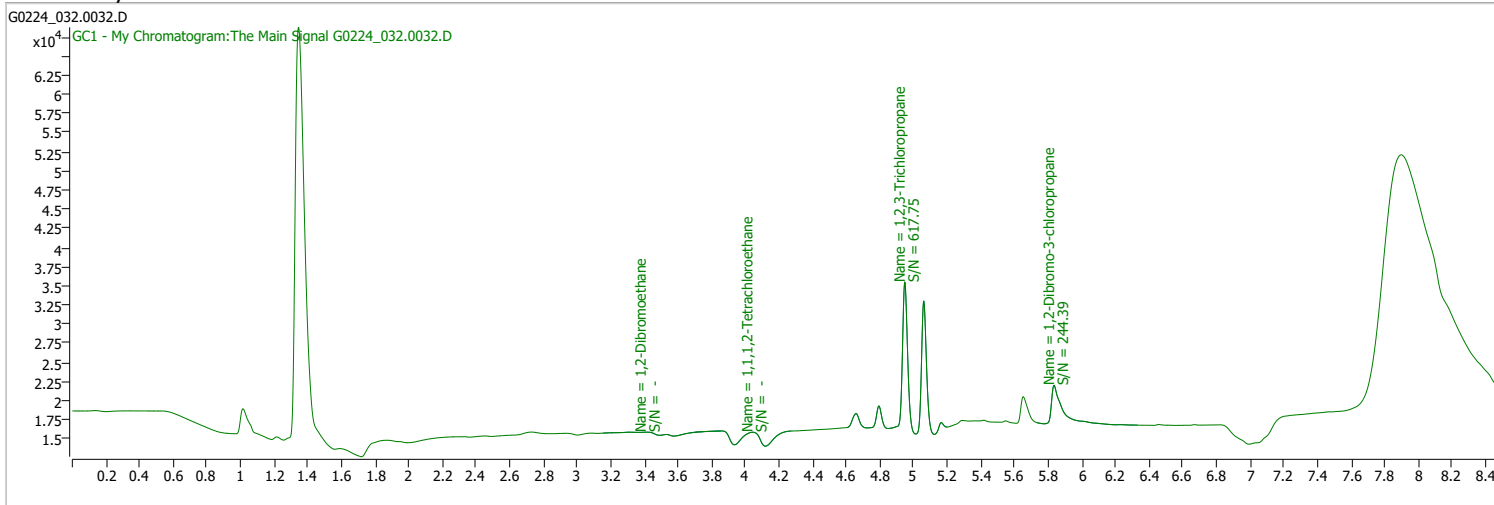
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0989	3.97	-0.01	32849 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_032.0032.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 10:02:17 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

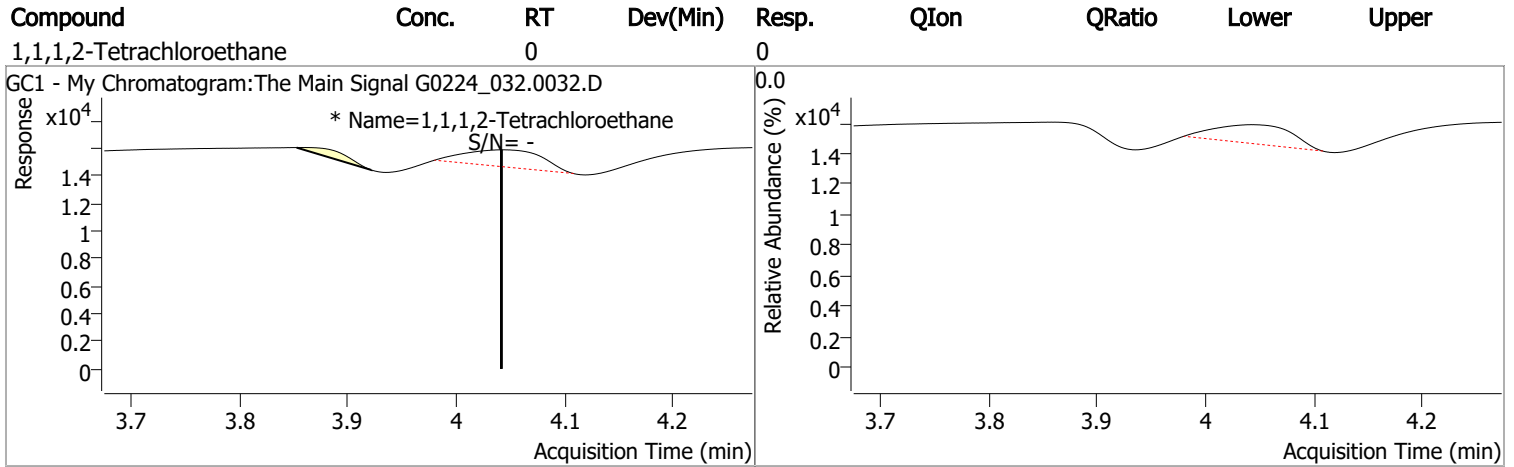
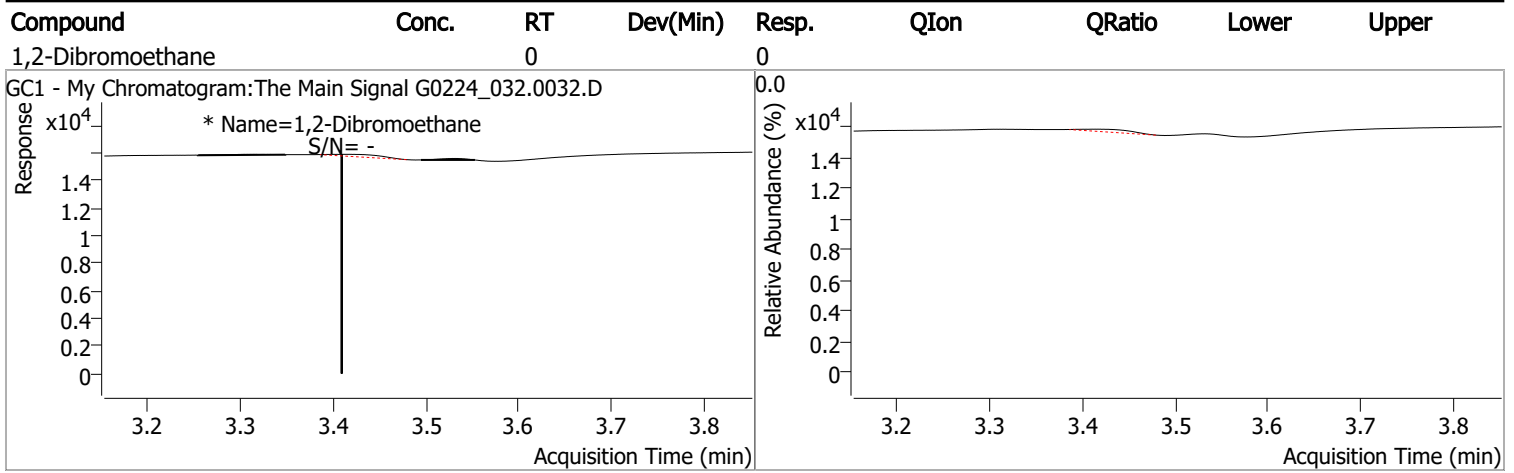
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.042	0.0	0		µg/L	md
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.409	0.0	0		µg/L	md
						QValue
						1

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

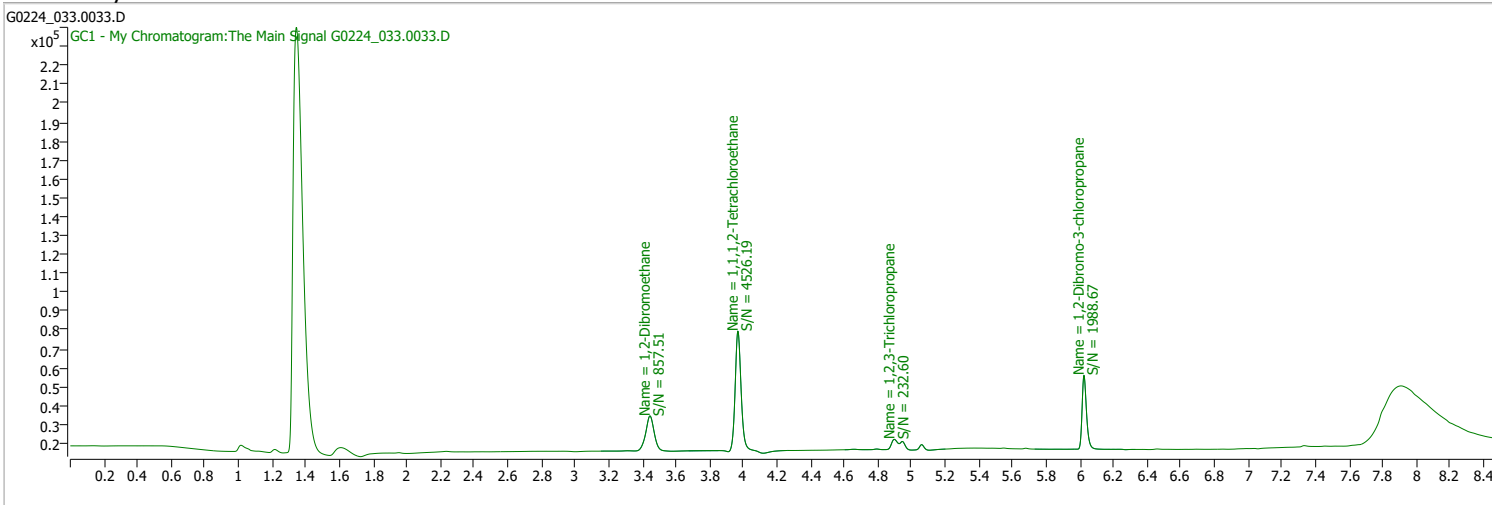
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_033.0033.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 10:22:01 PM
Sample Name	CAL5-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

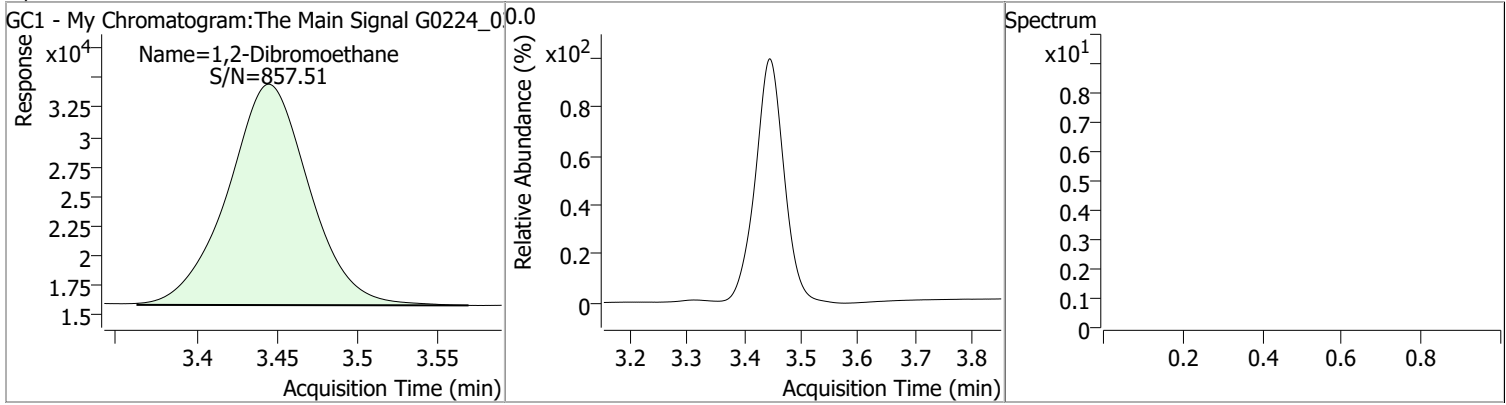


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	152809	0.4232	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 423.15%		*
Target Compounds						
M 1,2-Dibromoethane	3.445	0.0	65765	0.4143	µg/L	QValue 100

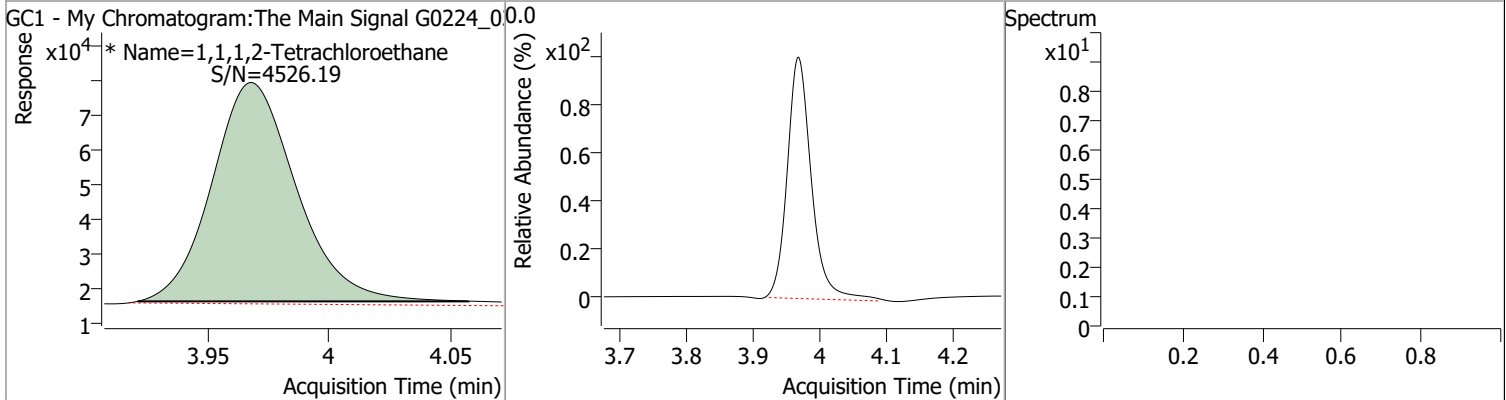
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.4143	3.45	-0.01	65765				



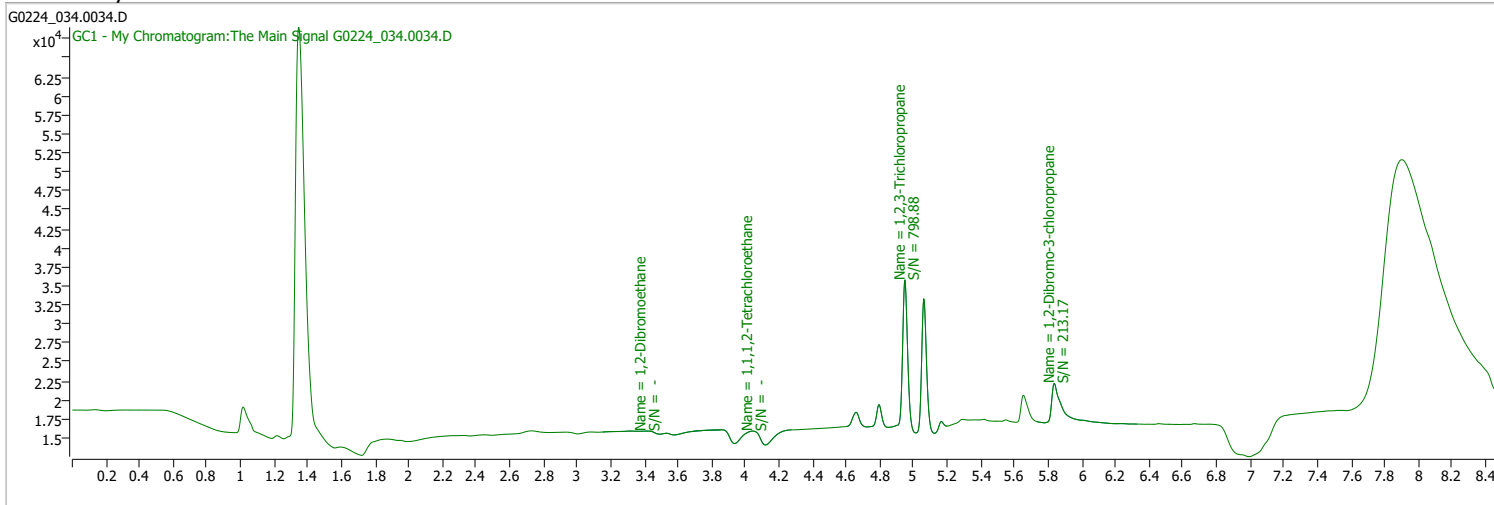
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.4232	3.97	-0.01	152809 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_034.0034.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 10:41:51 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

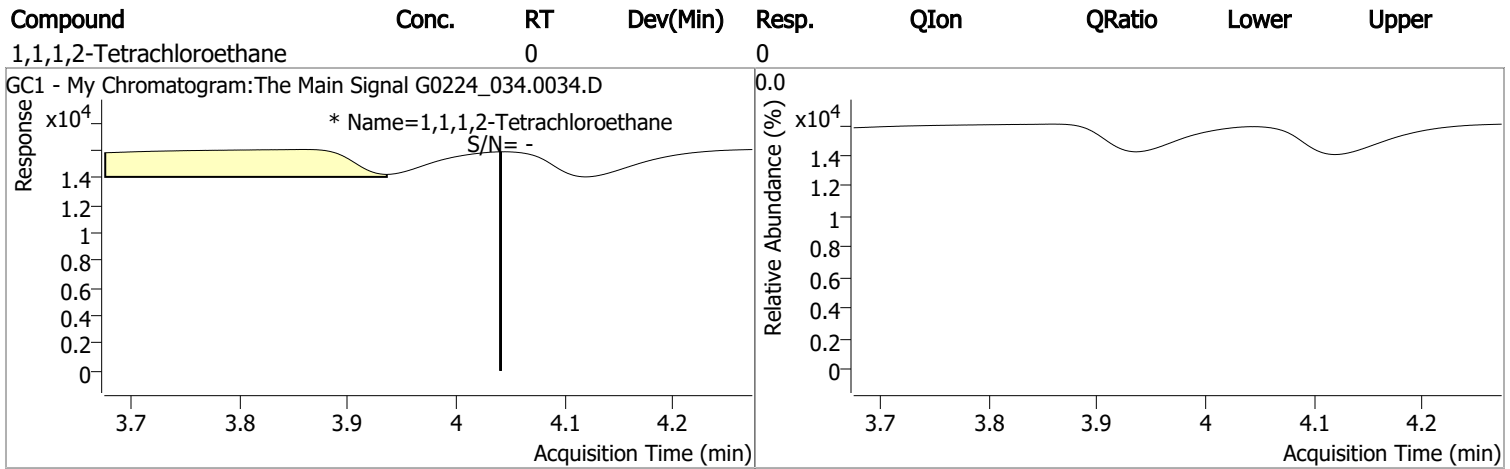
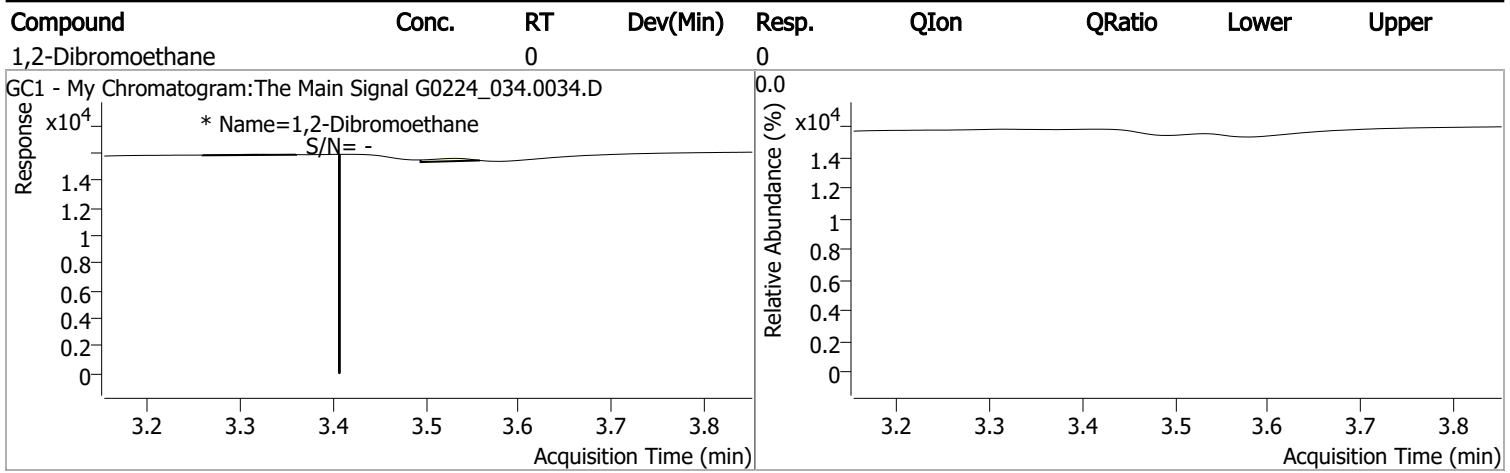
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.041	0.0	0		µg/L	md
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.407	0.0	0		µg/L	md
						QValue
						1

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

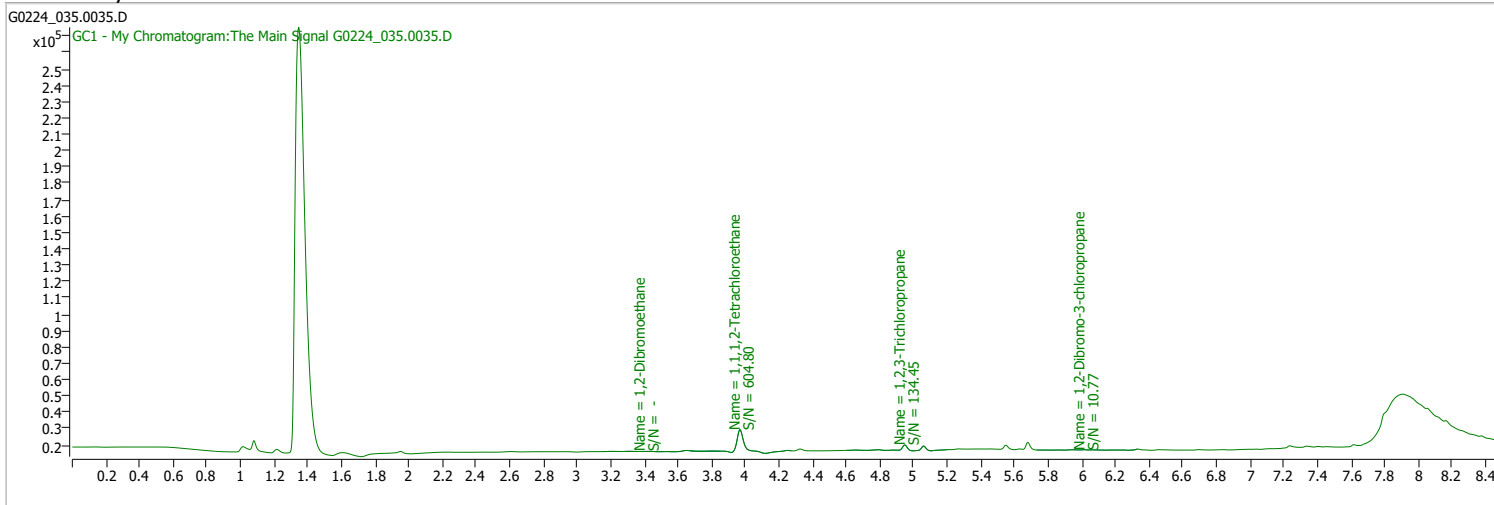
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_035.0035.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 11:01:45 PM
Sample Name	B22021435-022H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

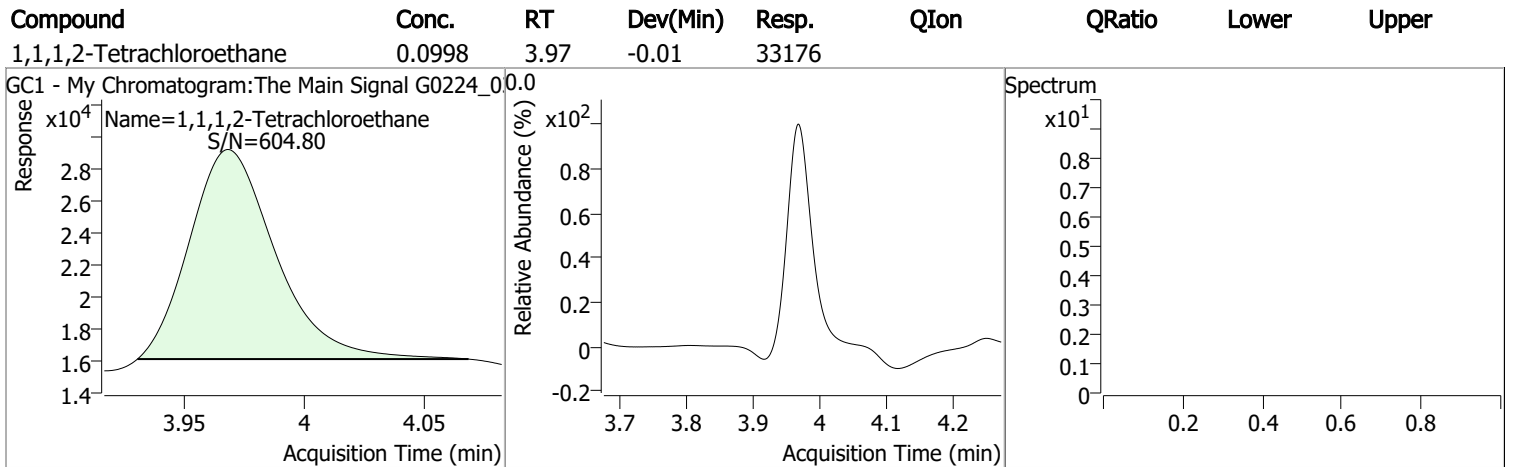
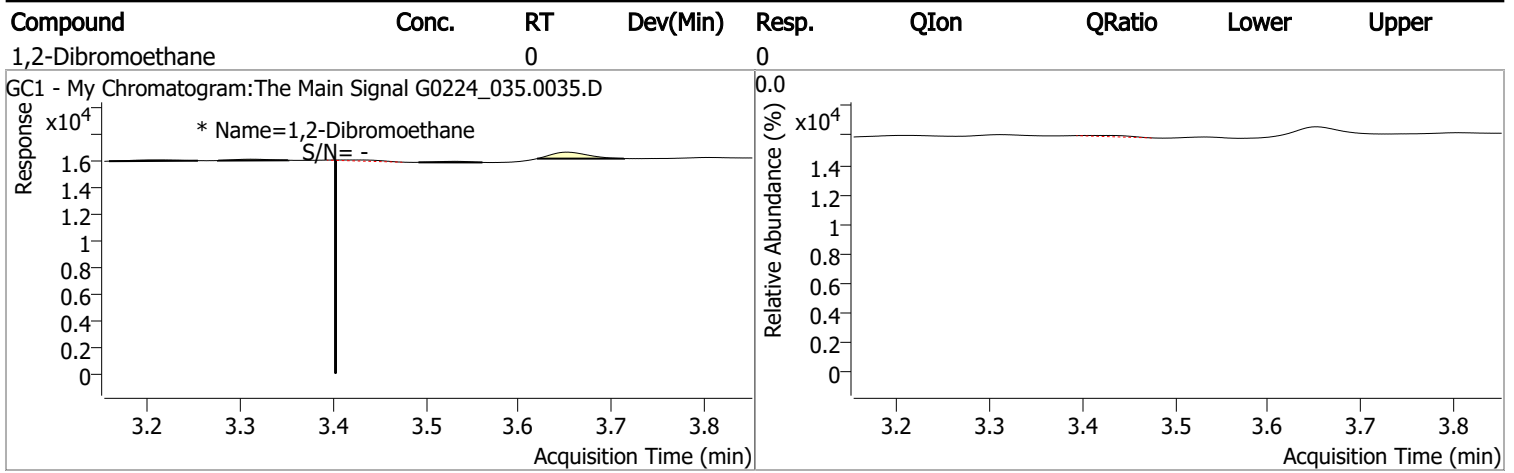
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	33176	0.0998	µg/L	-0.007
Spiked Amount: 0.100	Range: 70.0 - 130.0%		Recovery = 99.77%			
Target Compounds						
M 1,2-Dibromoethane	3.403	0.0	0		µg/L md	QValue 1

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

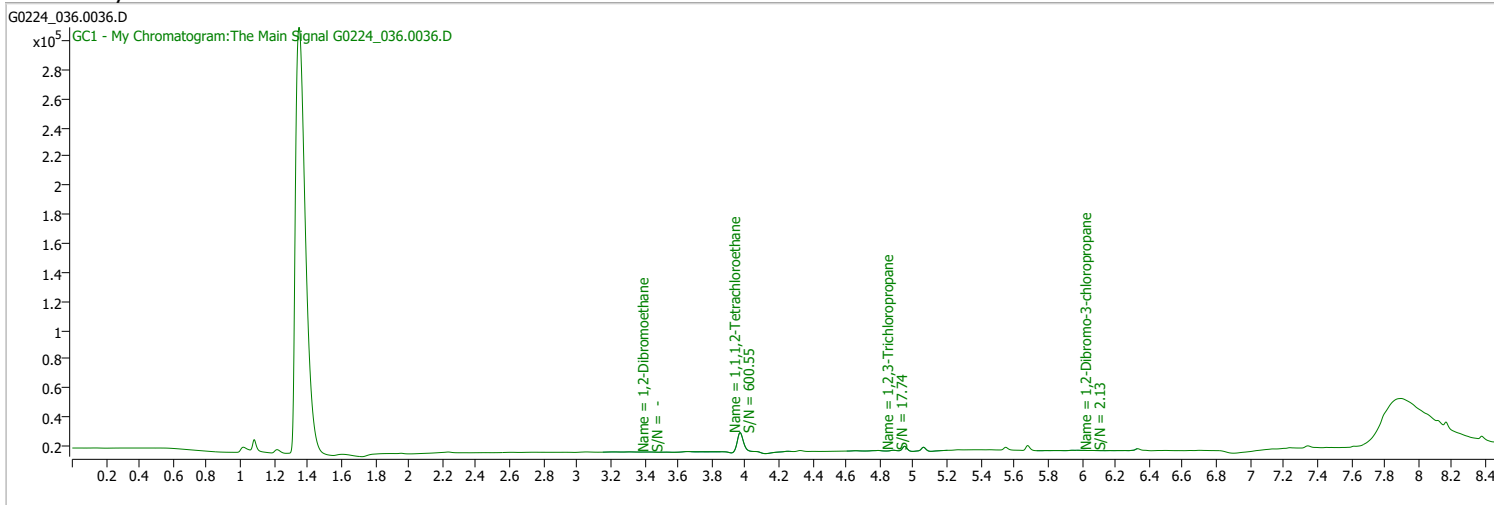
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_036.0036.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 11:21:30 PM
Sample Name	B22021435-025A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

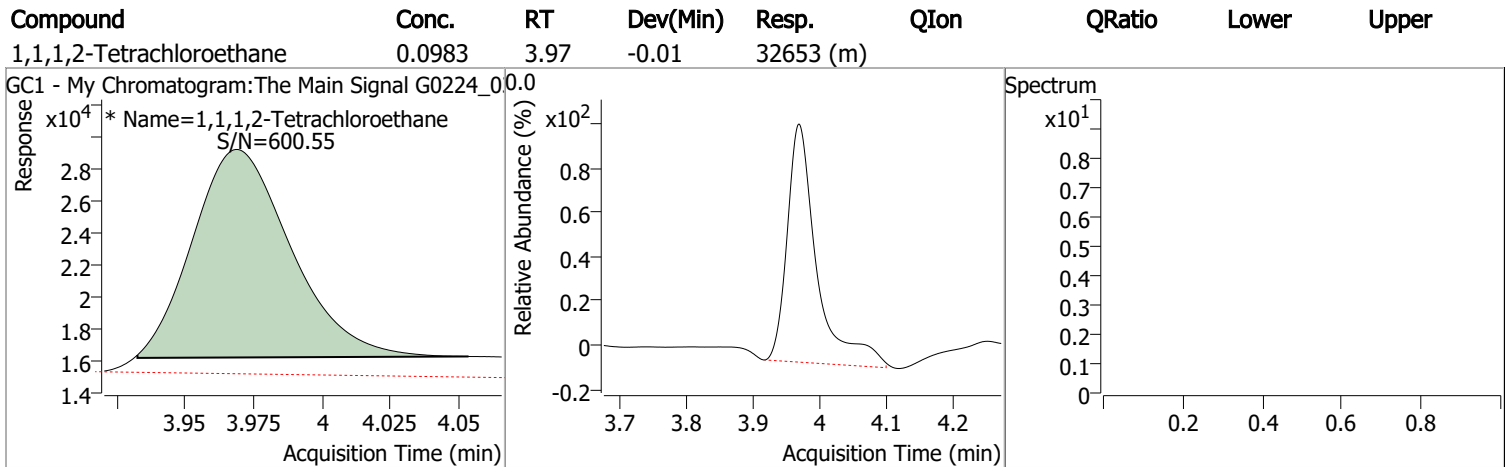
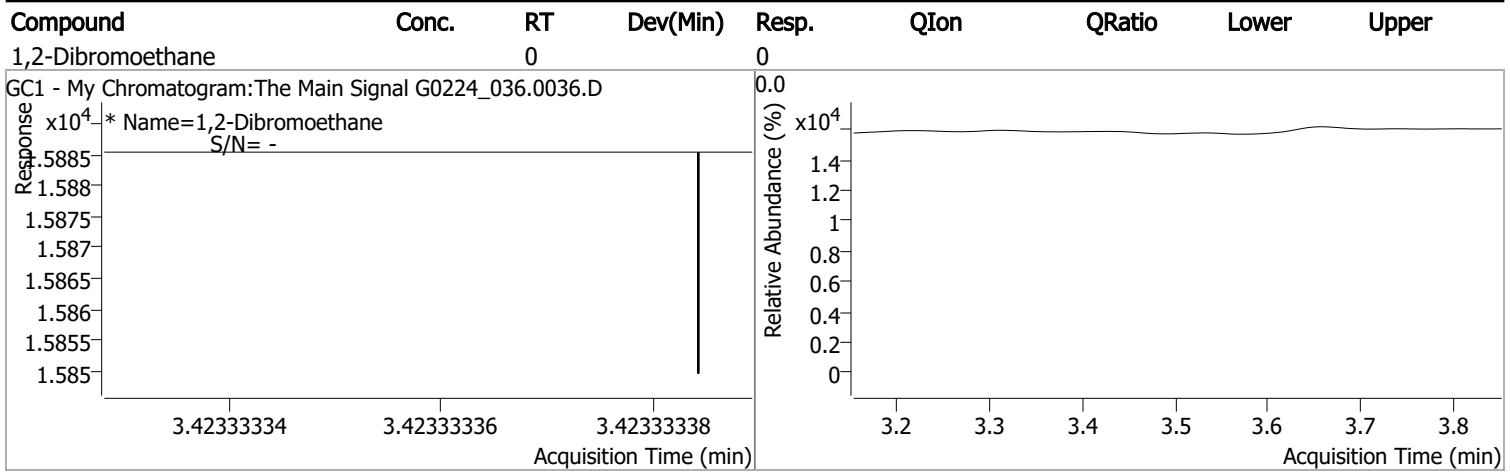
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	32653	0.0983	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 98.31%		
Target Compounds						
M 1,2-Dibromoethane	3.423	0.0	0		µg/L	md
						QValue
						1

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

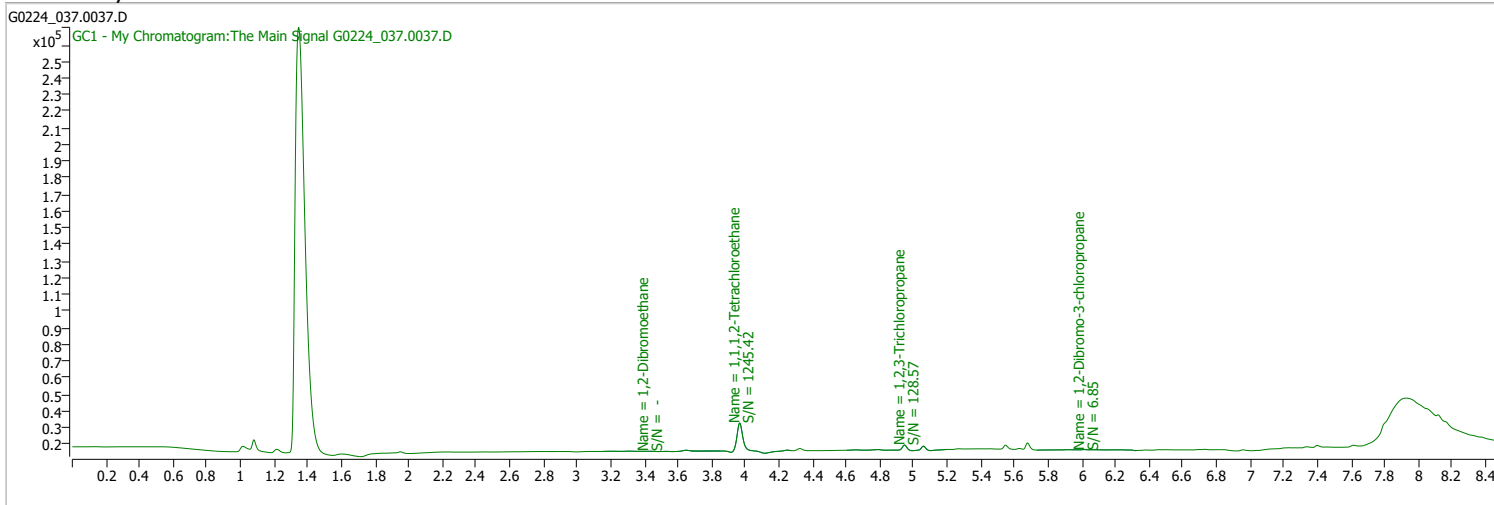
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_037.0037.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/24/2022 11:41:09 PM
Sample Name	B22021435-027H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

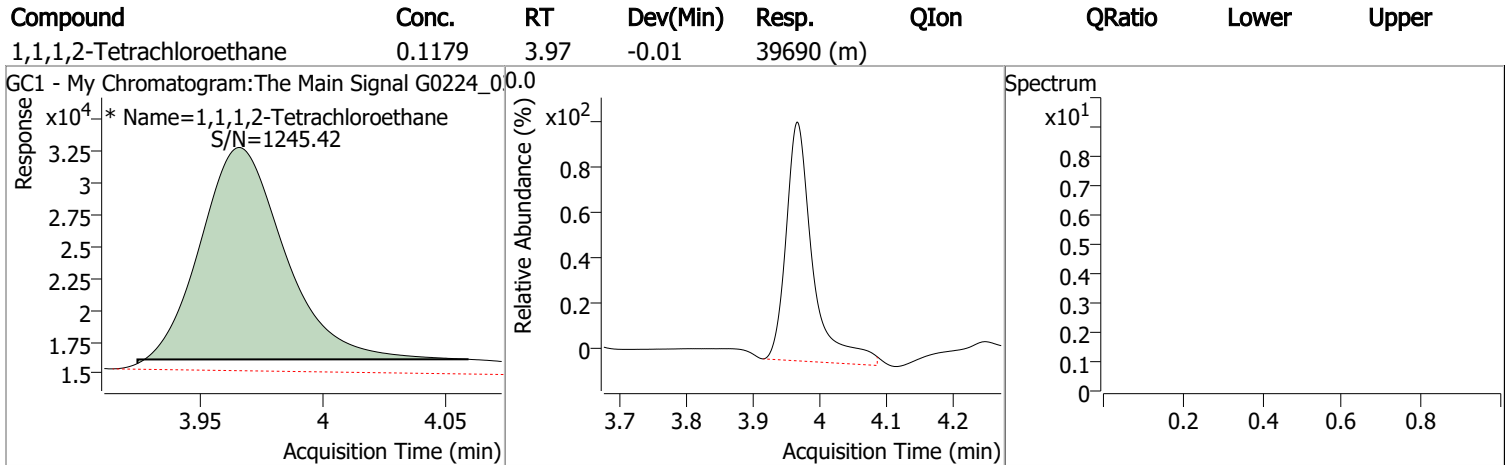
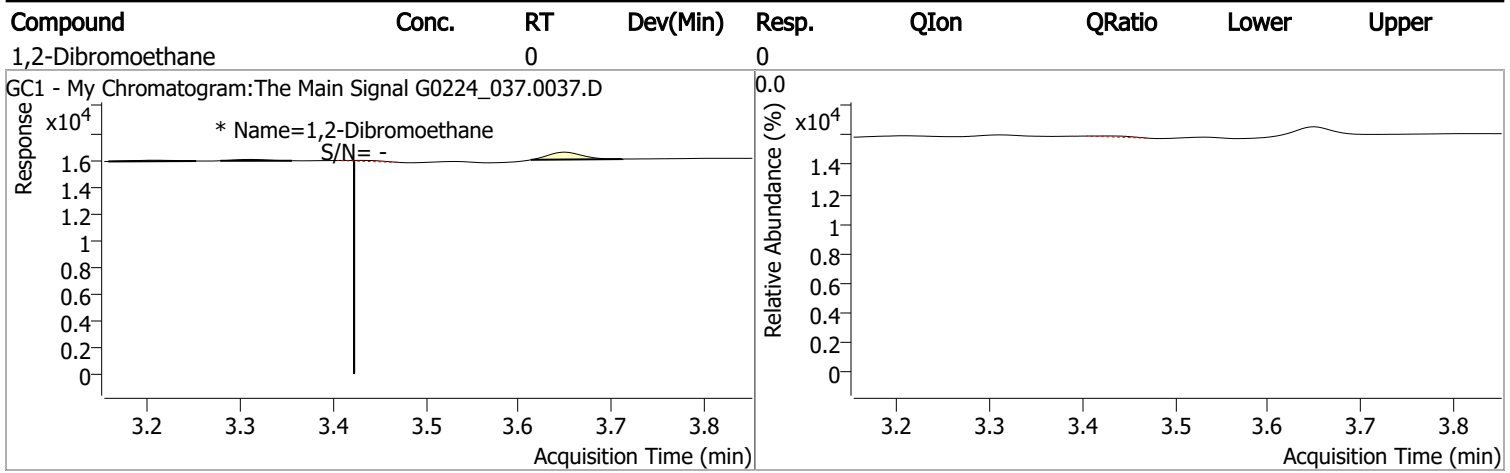
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.966	0.0	39690	0.1179	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 117.88%		
Target Compounds						
M 1,2-Dibromoethane	3.423	0.0	0		µg/L	md
						QValue 1

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

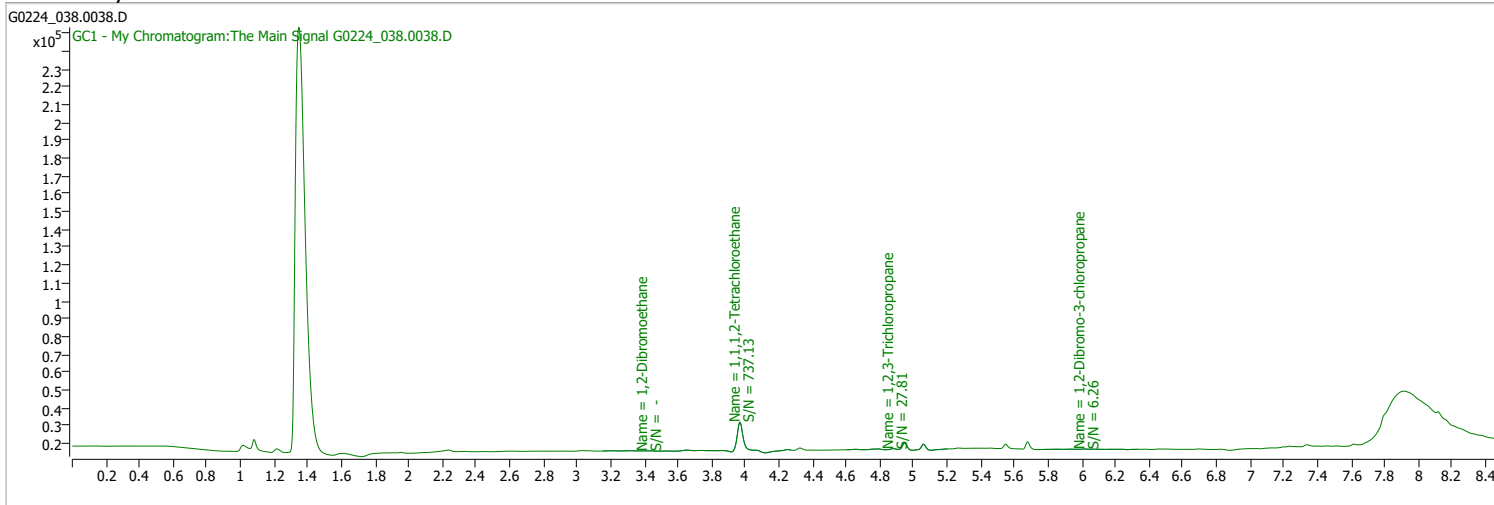
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_038.0038.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 12:00:50 AM
Sample Name	B22021435-030A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

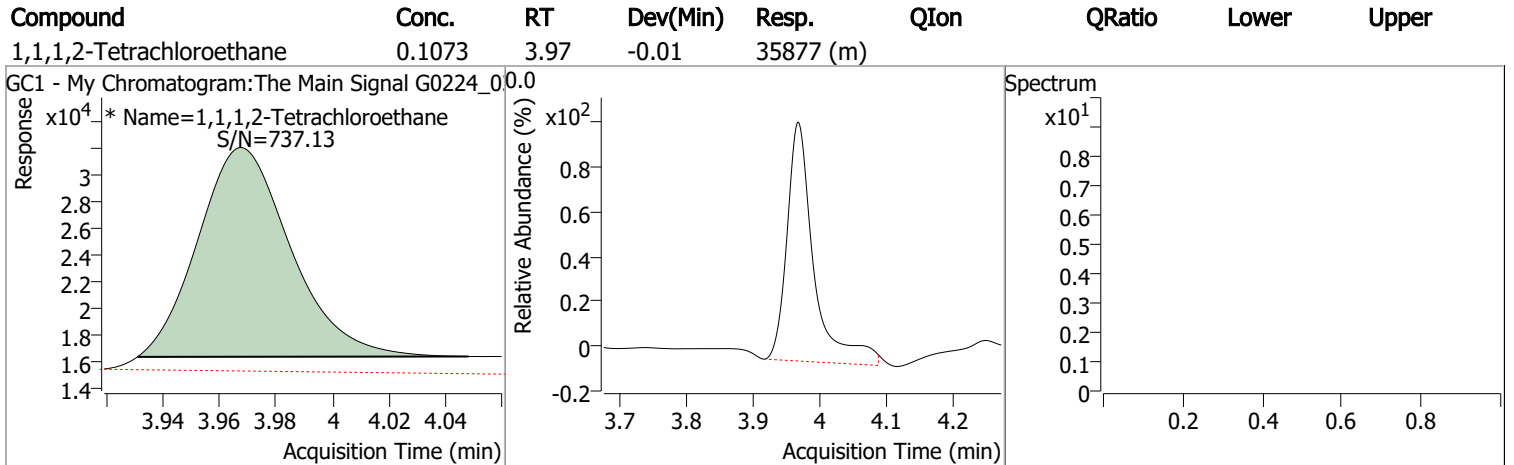
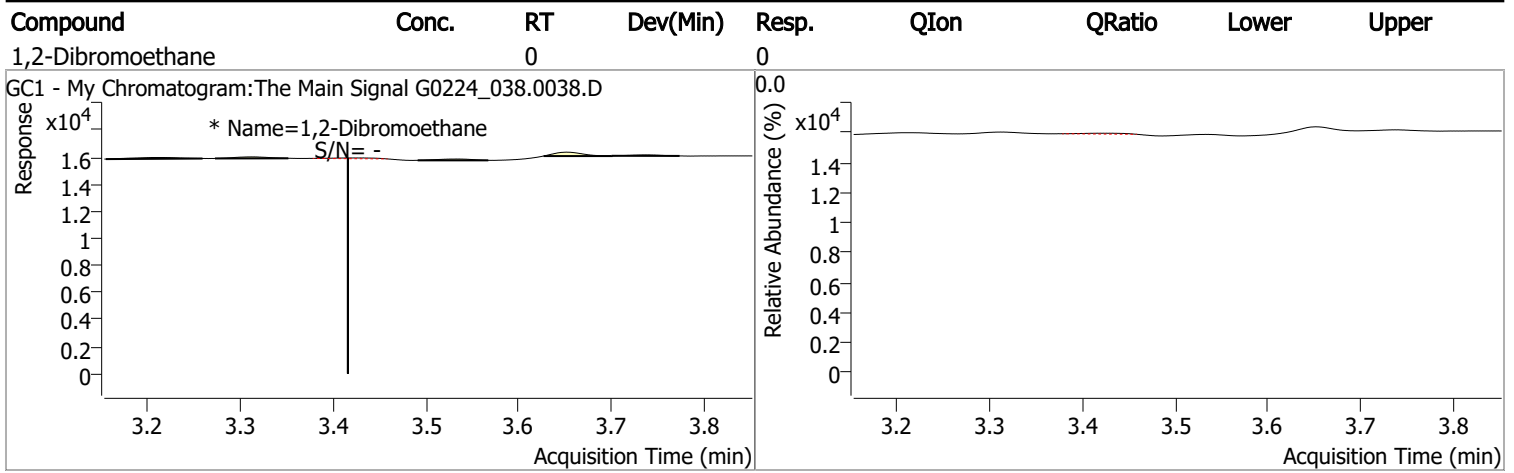
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	35877	0.1073	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 107.28%		
Target Compounds						
M 1,2-Dibromoethane	3.416	0.0	0		µg/L	md
						QValue 1

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

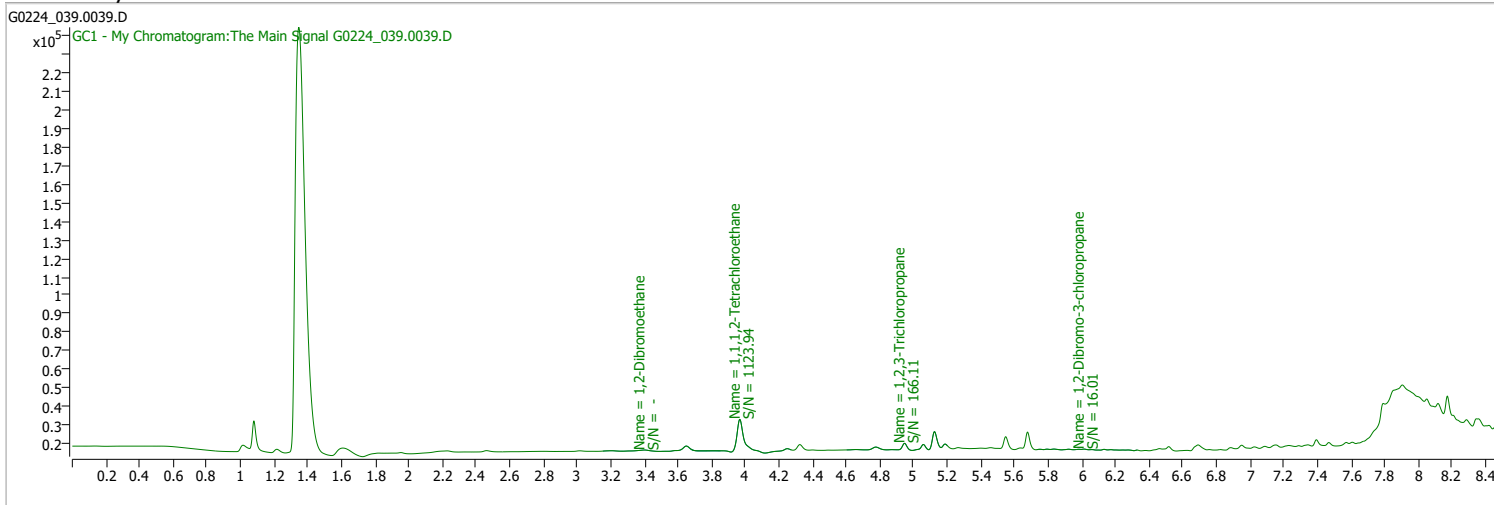
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_039.0039.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 12:20:35 AM
Sample Name	B22021435-032H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	re-extract
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

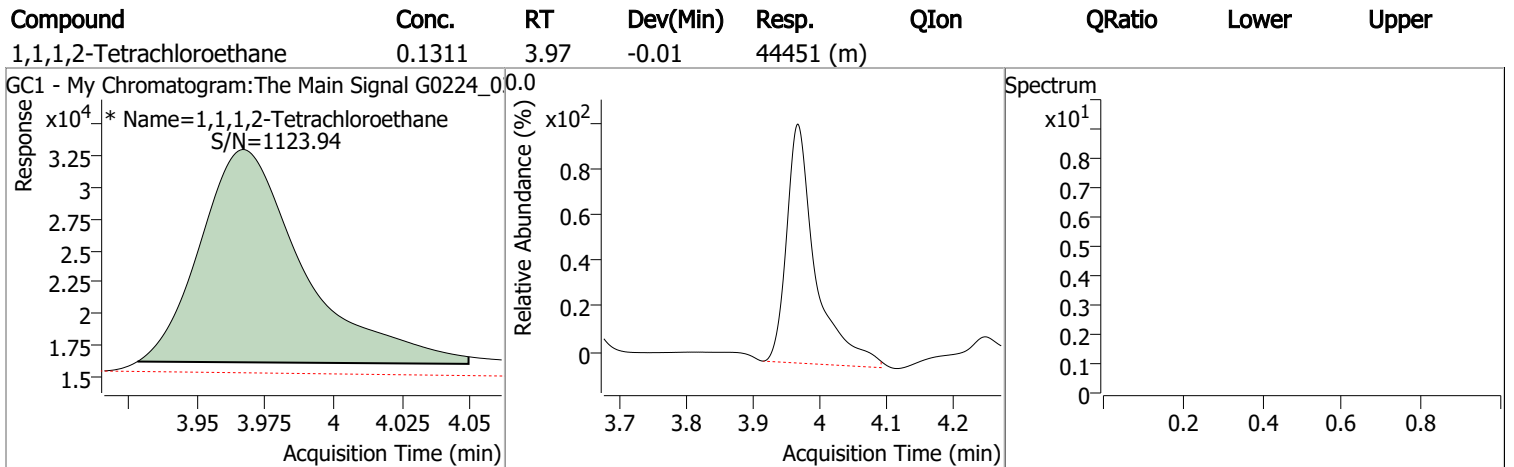
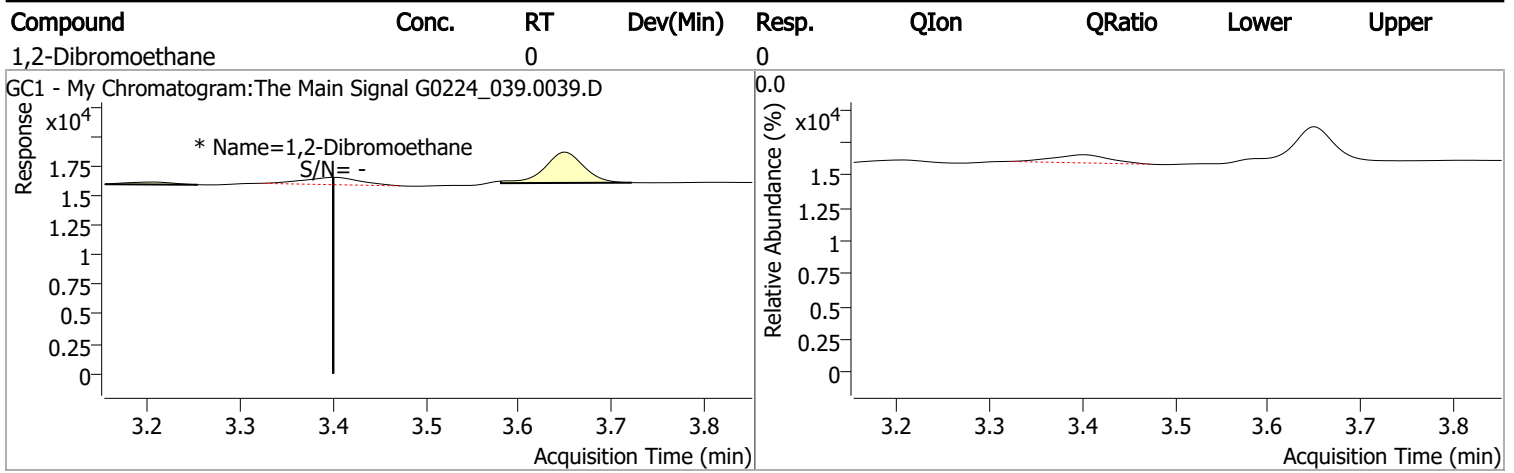
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.967	0.0	44451	0.1311	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 131.08%		*
Target Compounds						
M 1,2-Dibromoethane	3.400	0.0	0		µg/L	md
						QValue
						1

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

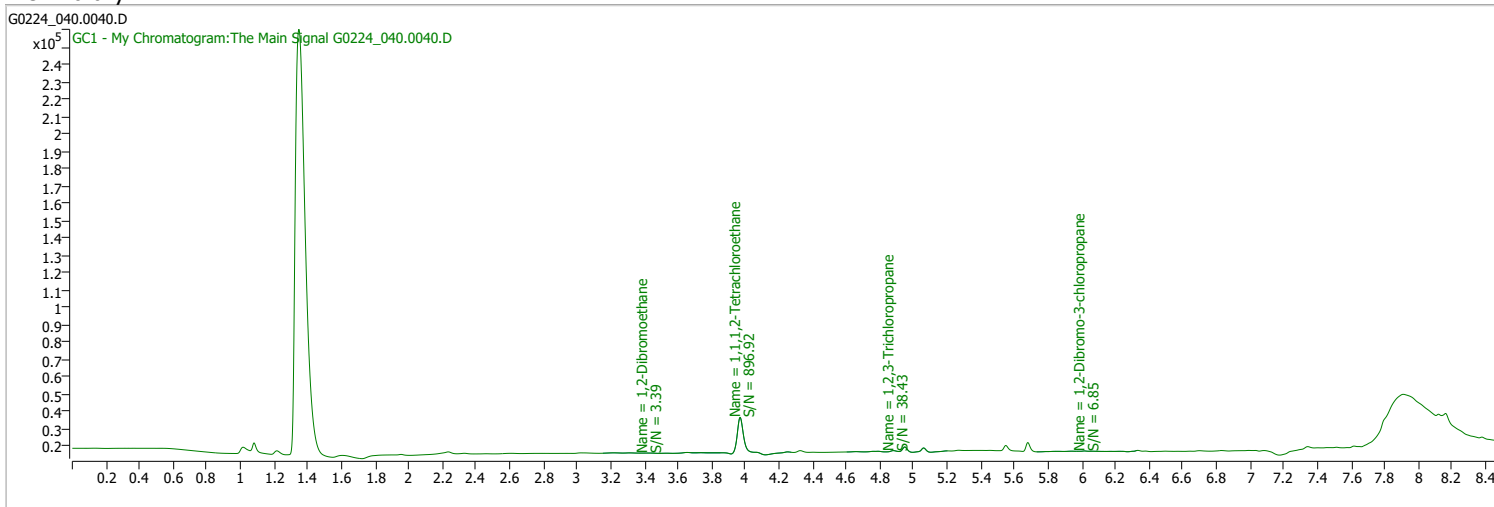
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_040.0040.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 12:40:21 AM
Sample Name	B22021435-035A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	re-extract
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

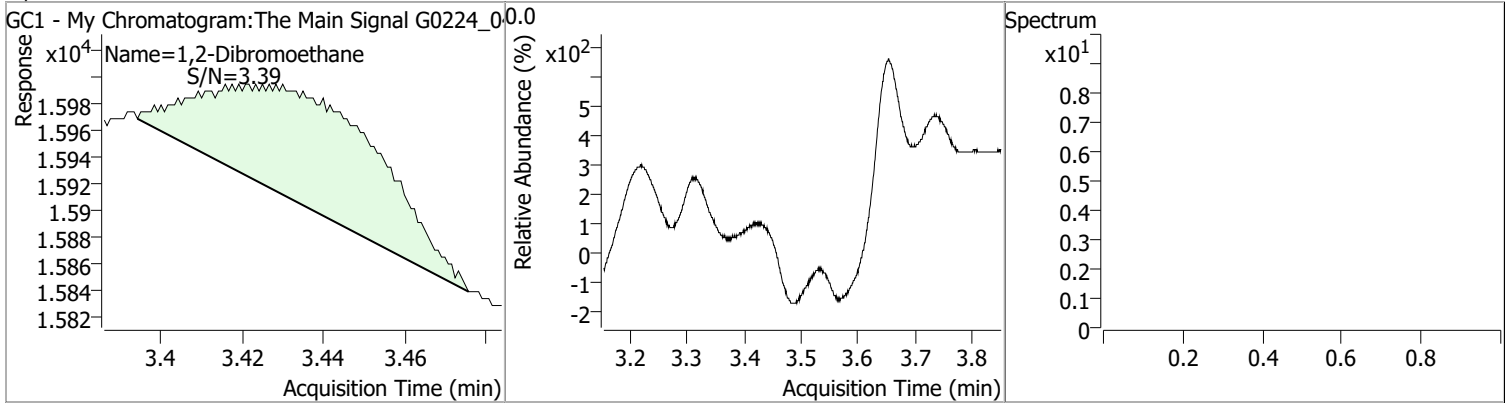


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	49542	0.1452	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 145.16%		*
Target Compounds						
M 1,2-Dibromoethane	3.416	0.0	257	0.0015	µg/L	QValue 100

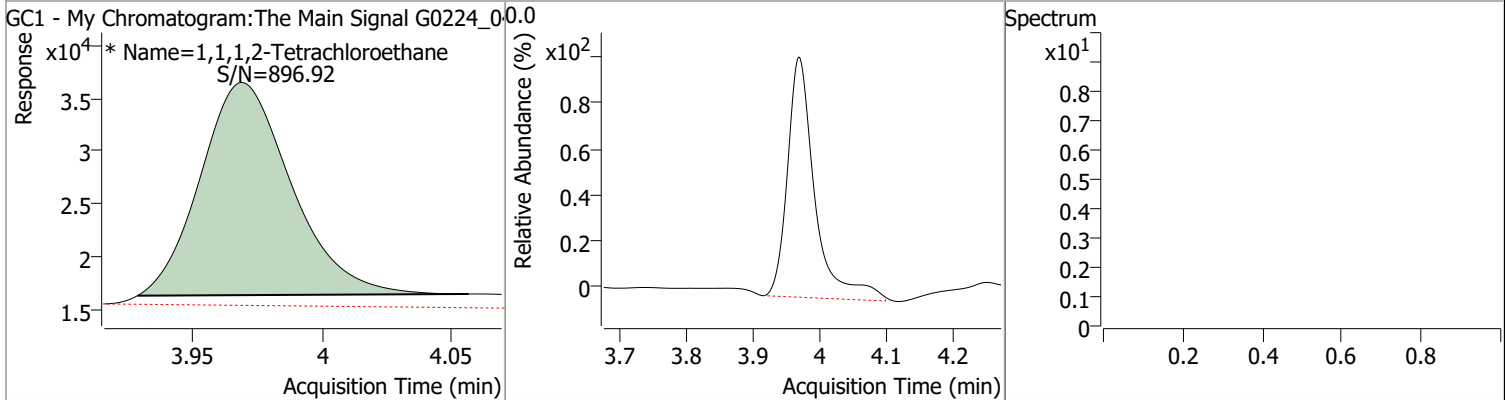
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0015	3.42	-0.04	257				



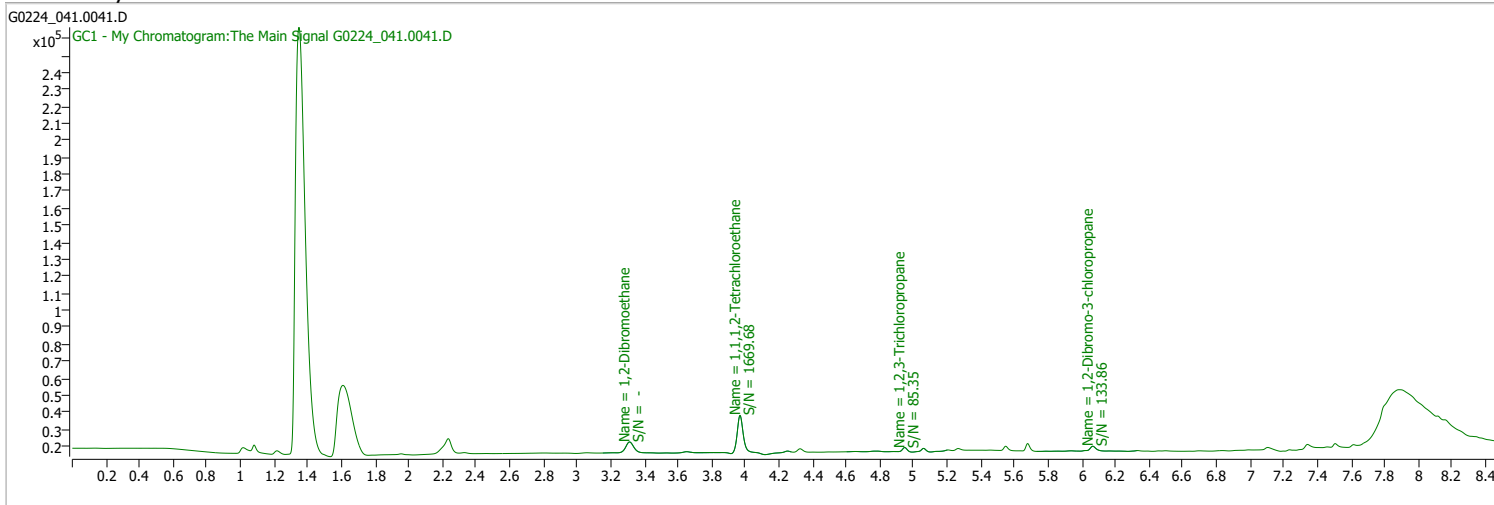
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1452	3.97	-0.01	49542 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_041.0041.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:00:10 AM
Sample Name	B22021528-001C	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	cannot re-extract due to limited sample volume
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

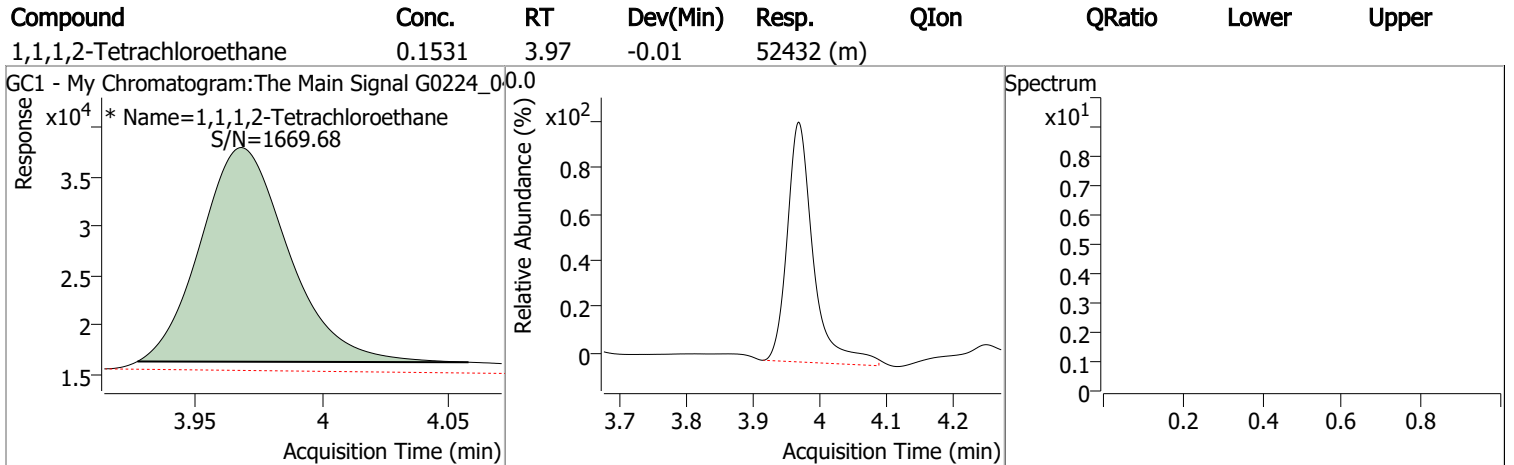
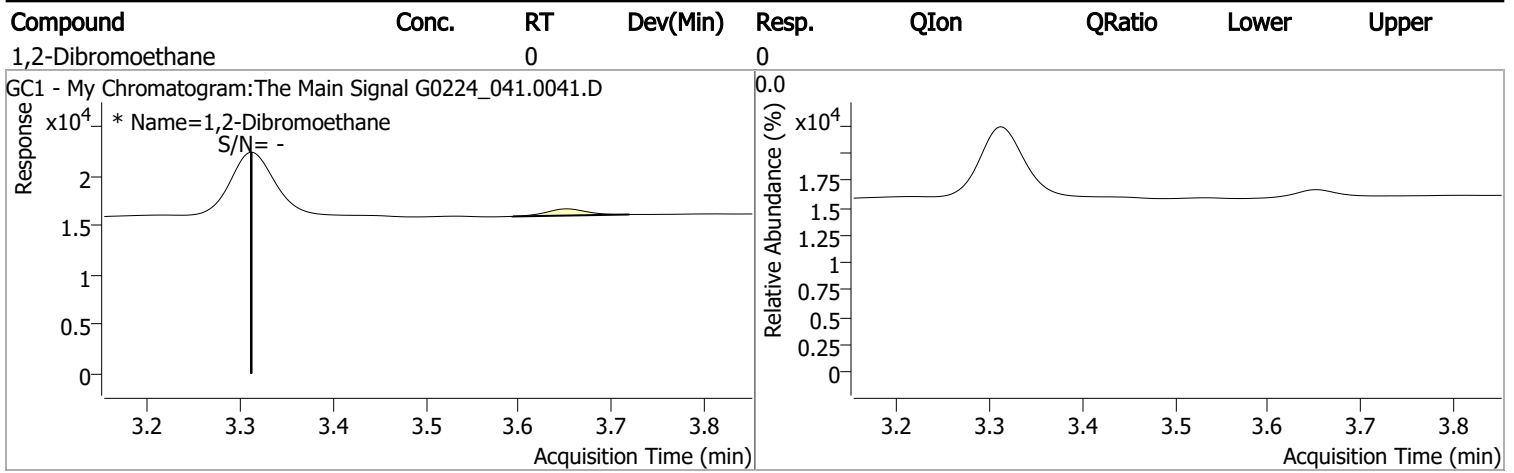
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	52432	0.1531	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 153.14%		*
Target Compounds						
M 1,2-Dibromoethane	3.312	0.0	0		µg/L	md
						QValue
						1

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

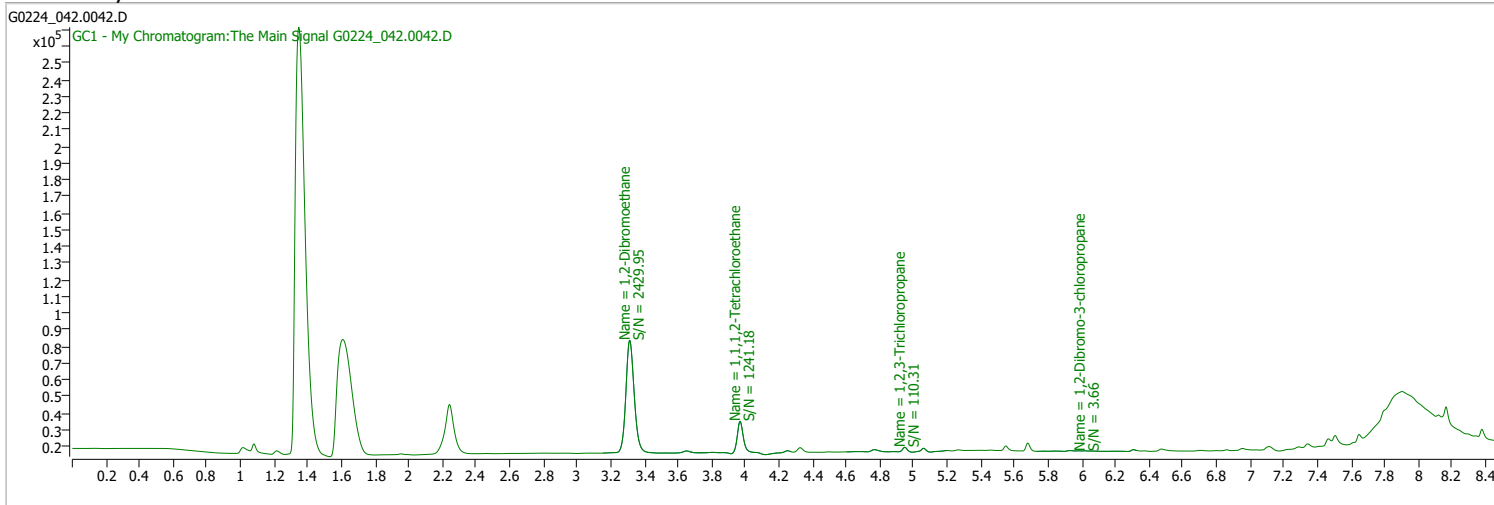
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_042.0042.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:19:57 AM
Sample Name	B22021528-002C	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	re-extract
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.968	0.0	45163	0.1331	µg/L	m	-0.006
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 133.05%		*	

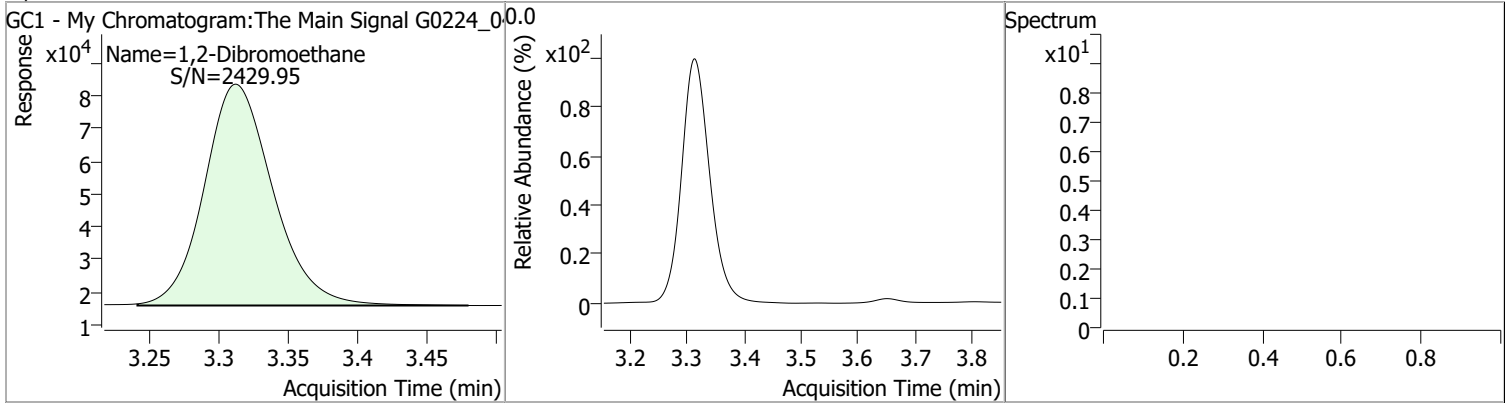
Target Compounds

M 1,2-Dibromoethane	3.313	0.0	233854	2.0434	µg/L		QValue 100
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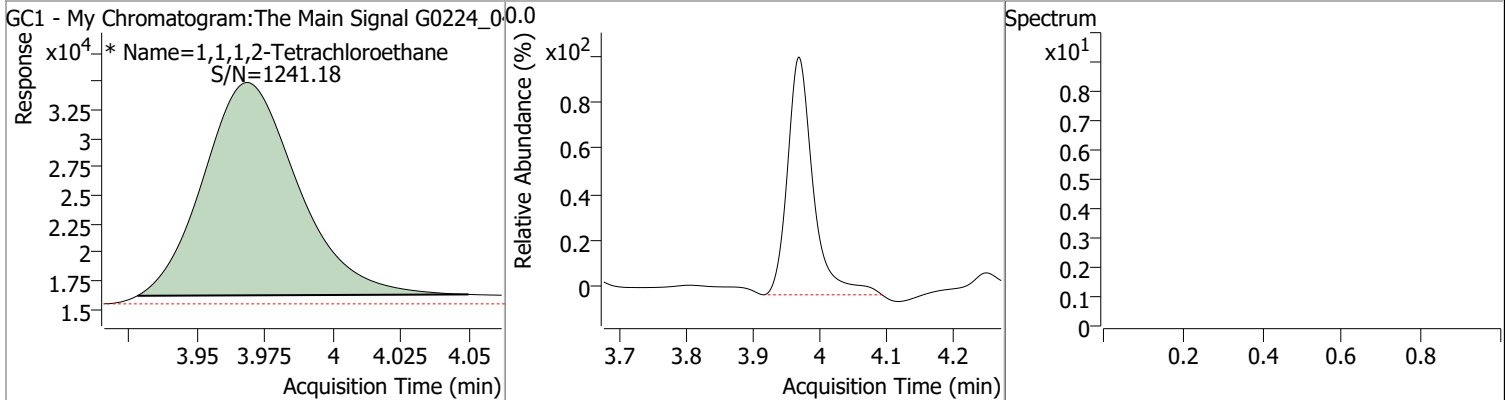
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	2.0434	3.31	-0.14	233854				



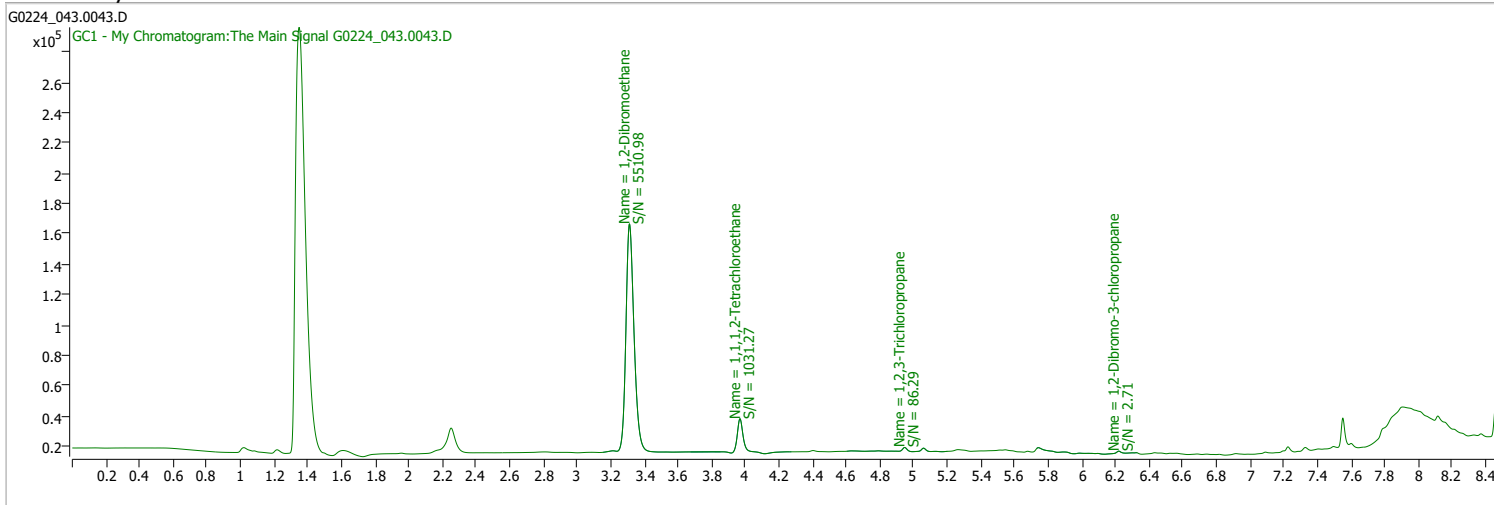
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1331	3.97	-0.01	45163 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_043.0043.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:39:37 AM
Sample Name	B22021528-003C	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	re-extract
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.968	0.0	50827	0.1487	µg/L	-0.006
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 148.71%		*

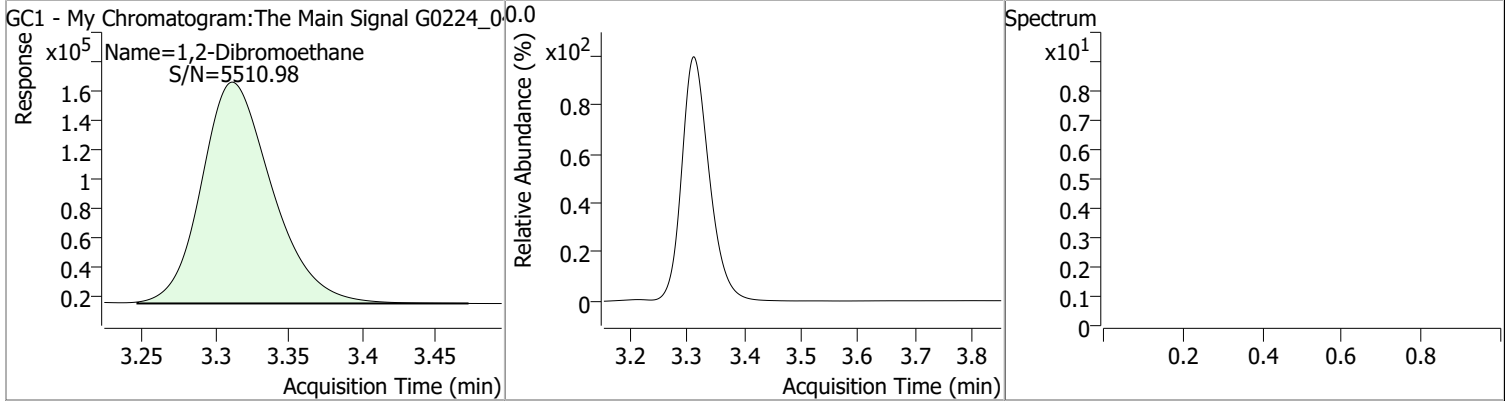
Target Compounds

M 1,2-Dibromoethane	3.311	0.0	505159		µg/L	QValue 100
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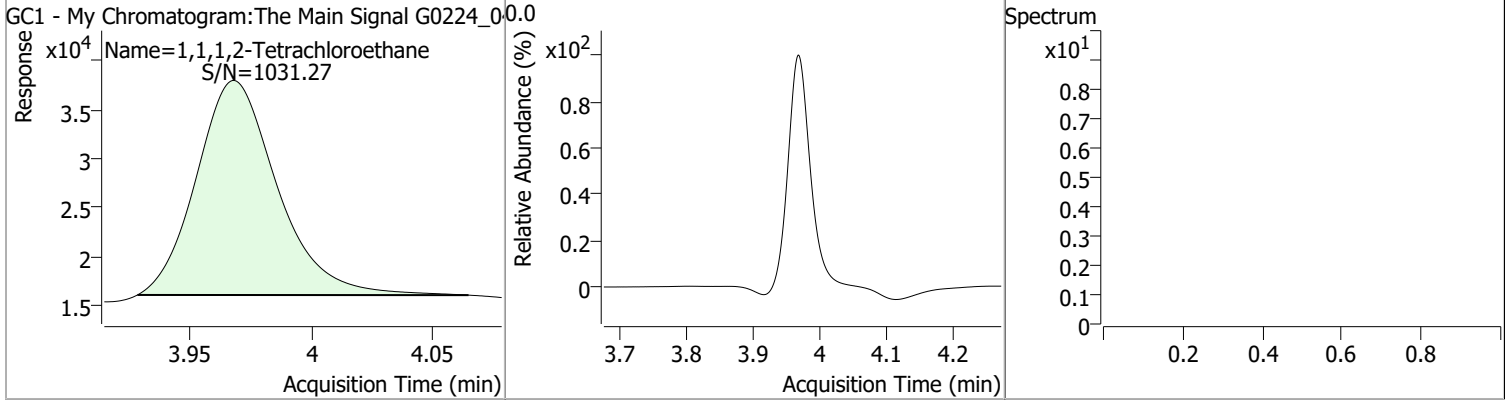
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane		3.31	-0.14	505159				



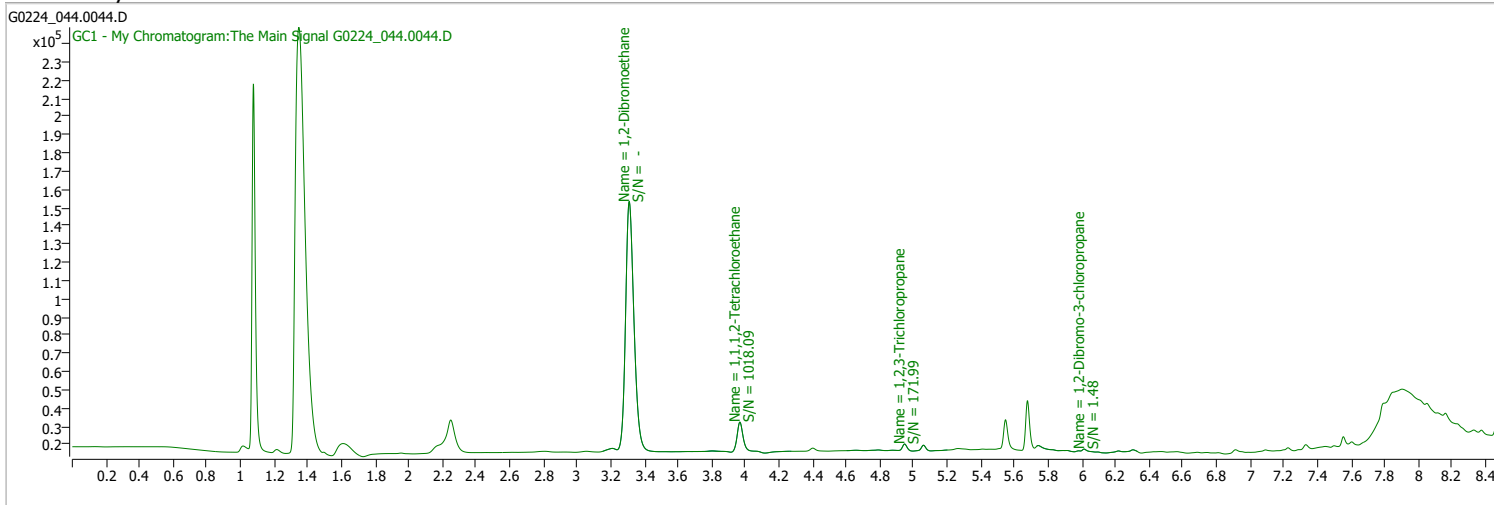
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1487	3.97	-0.01	50827				



Quantitation Results Report (QT Reviewed)

Data File	G0224_044.0044.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:59:23 AM
Sample Name	B22021528-004C	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

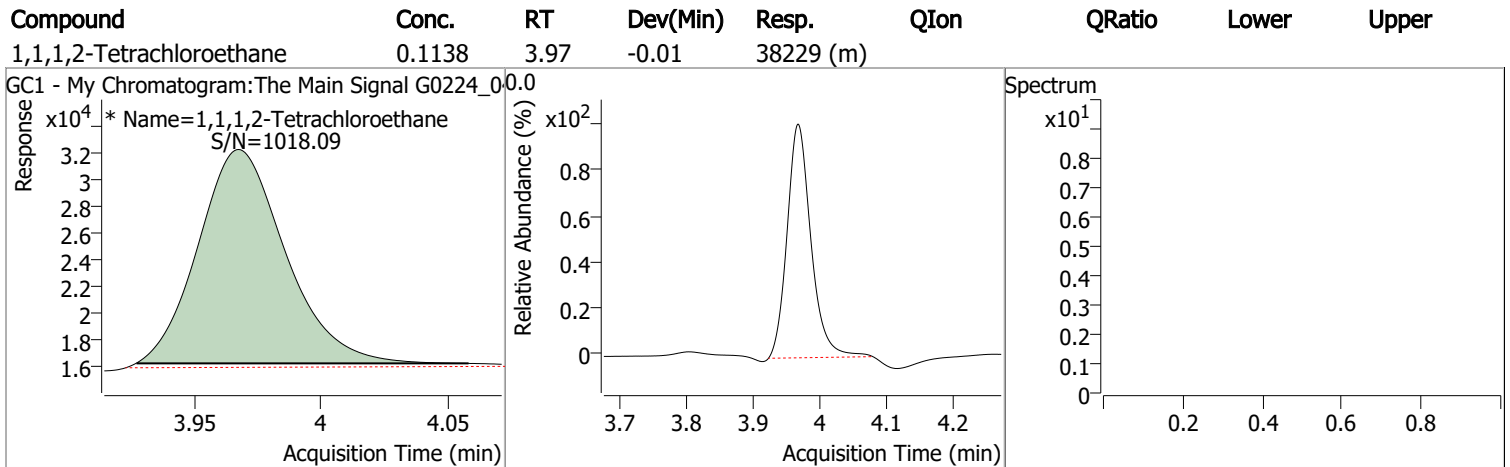
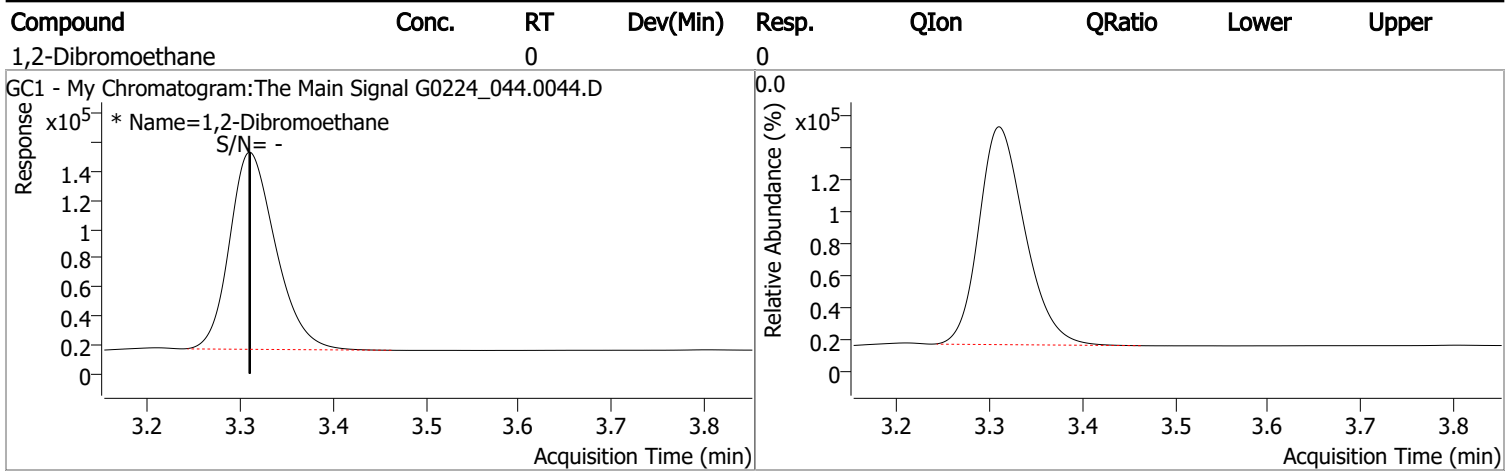
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	38229	0.1138	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 113.82%		
Target Compounds						
M 1,2-Dibromoethane	3.310	0.0	0		µg/L	md
						QValue 1

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

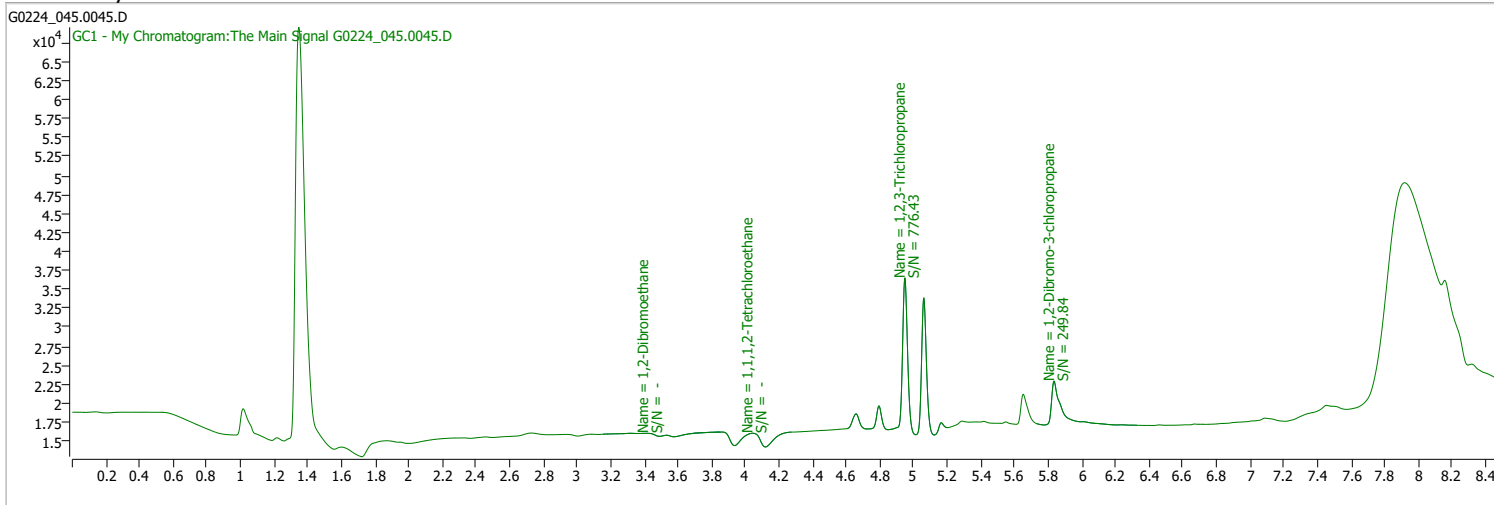
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_045.0045.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 2:18:58 AM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

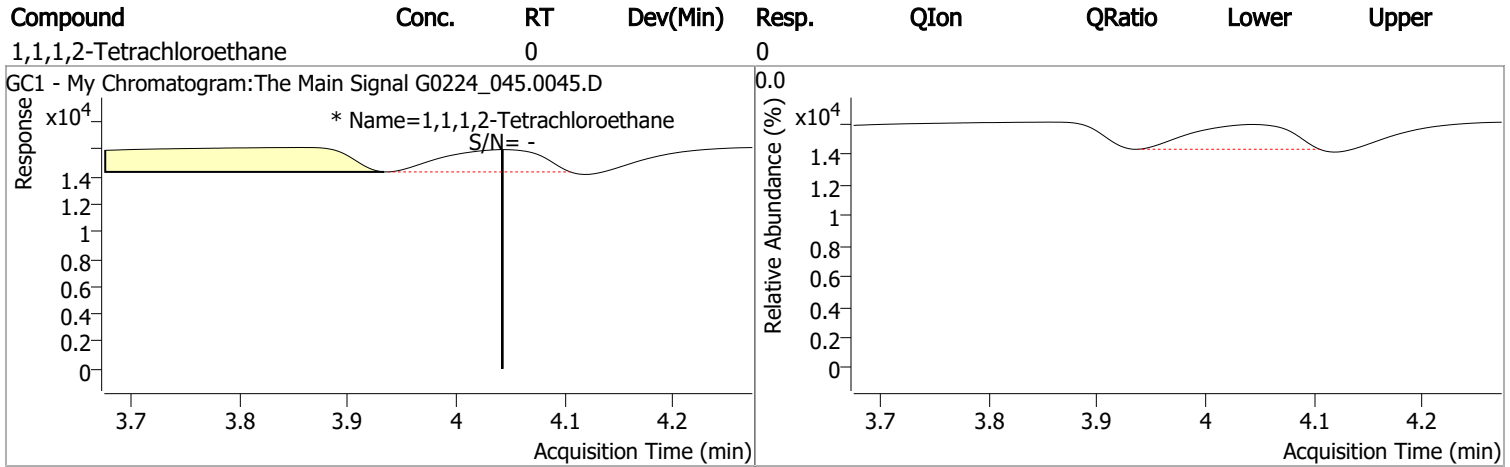
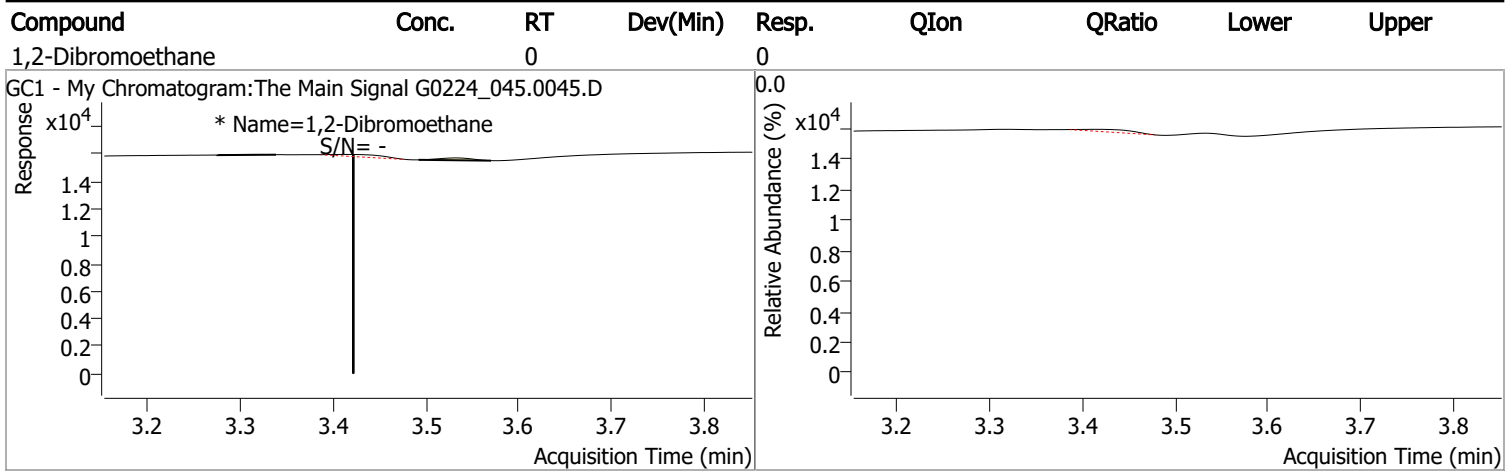
S 1,1,1,2-Tetrachloroethane	4.043	0.0	0		µg/L	md	0.068
Spiked Amount: 0.100	Range: 70.0 - 130.0%				Recovery = NA%		

Target Compounds

M 1,2-Dibromoethane	3.422	0.0	0		µg/L	md	QValue 1
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(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

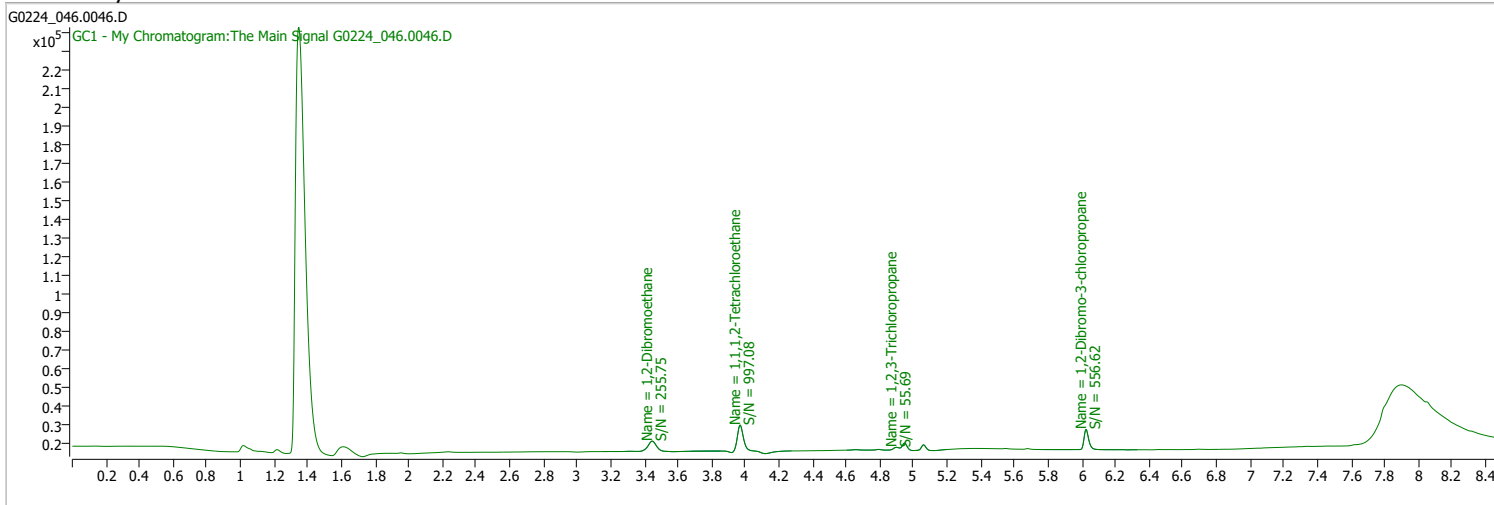
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_046.0046.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 2:38:50 AM
Sample Name	CAL3-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

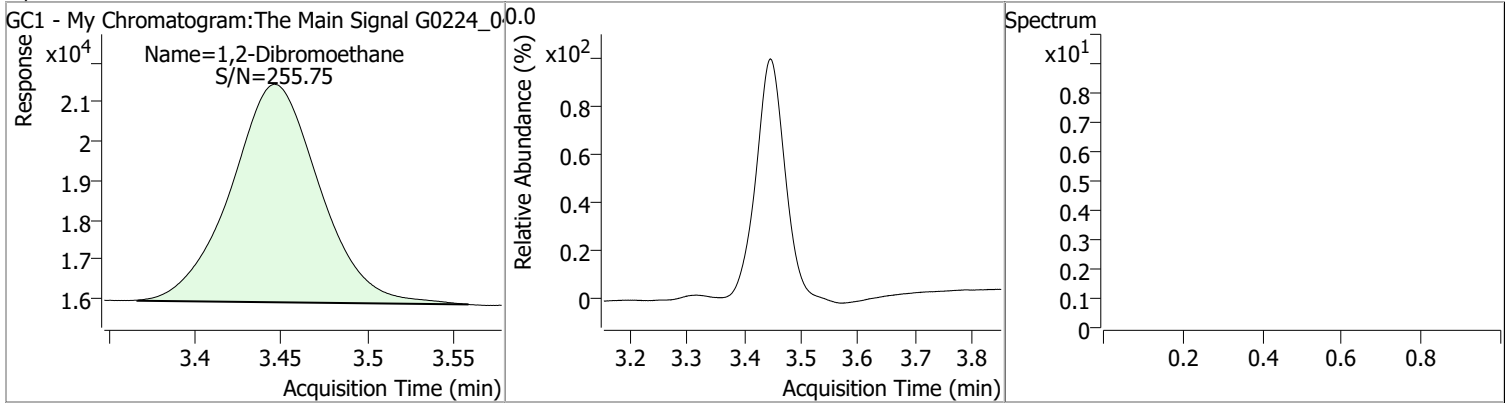


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	34534	0.1035	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 103.55%		
Target Compounds						
M 1,2-Dibromoethane	3.446	0.0	19816	0.1188	µg/L	QValue 100

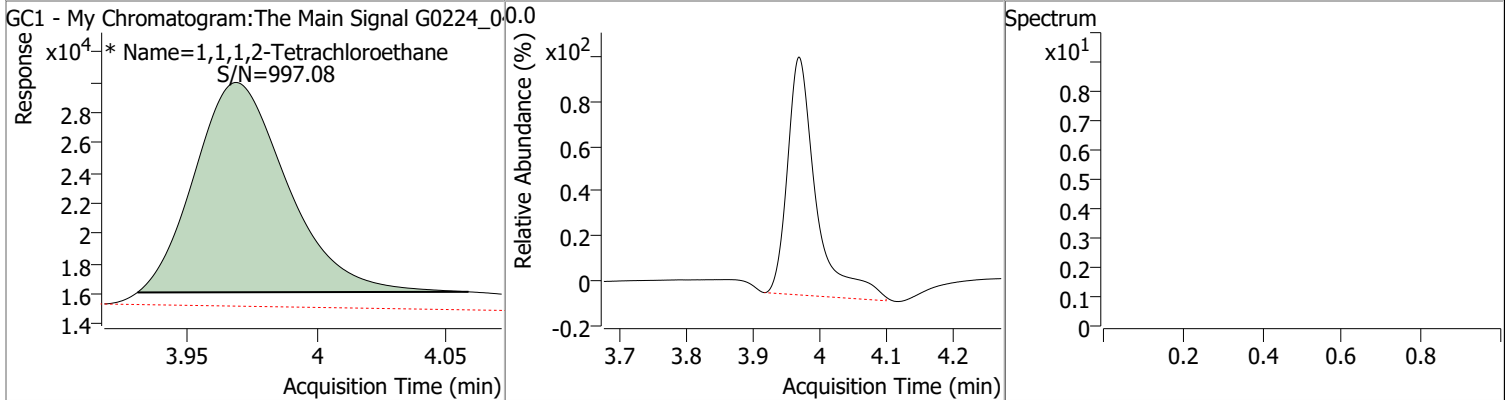
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1188	3.45	-0.01	19816				



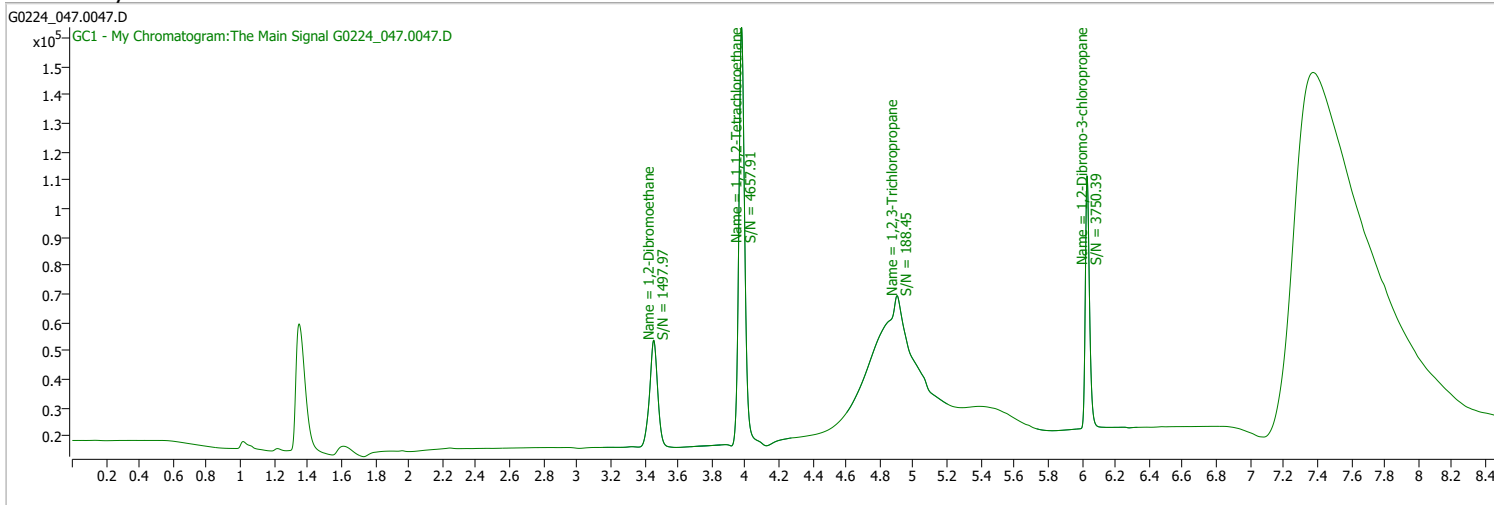
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1035	3.97	-0.01	34534 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_047.0047.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 8:57:34 AM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

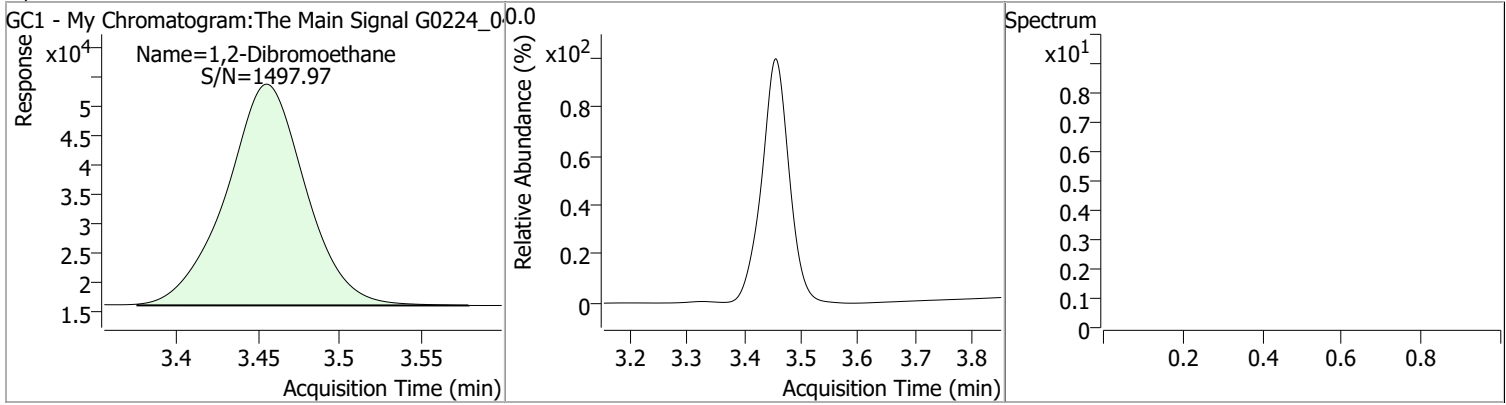


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.978	0.0	362681	0.9495	µg/L	0.003
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 949.45%		*
Target Compounds						
M 1,2-Dibromoethane	3.455	0.0	126642	0.8645	µg/L	100

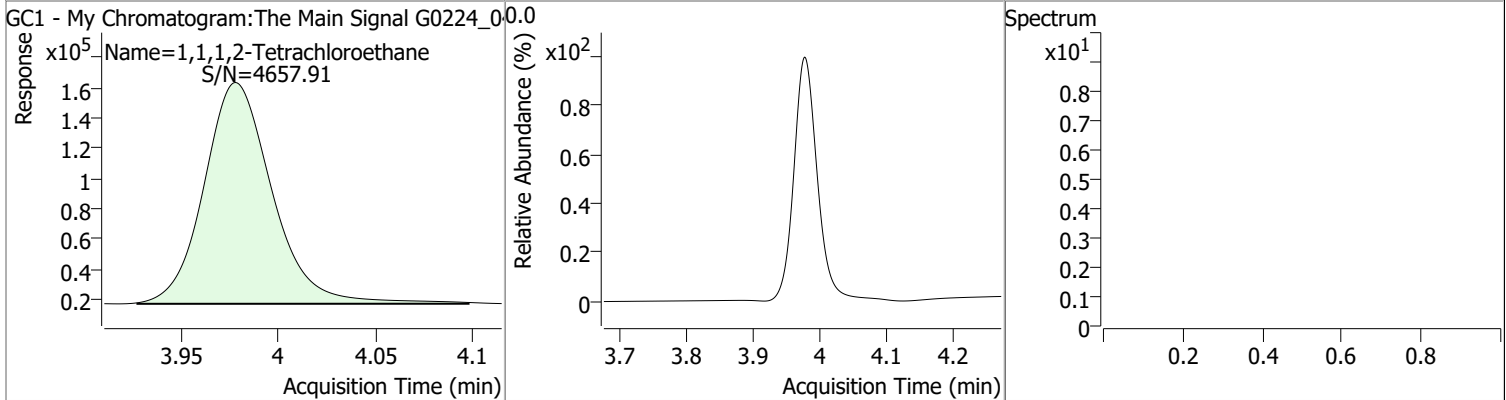
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.8645	3.46	0.00	126642				



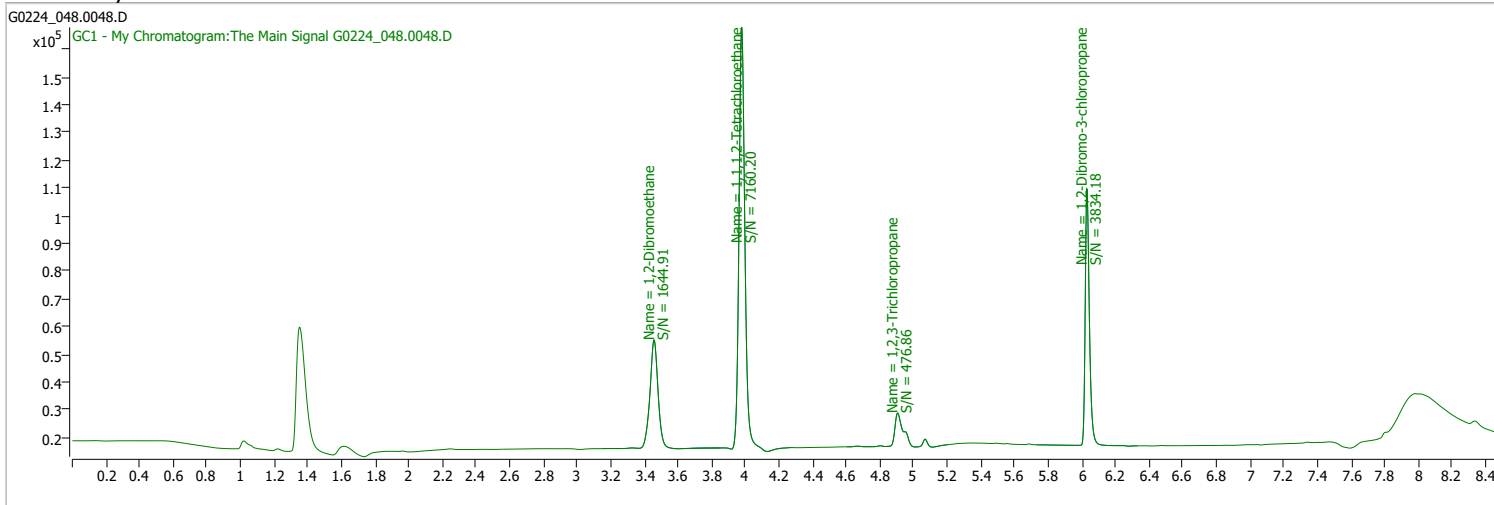
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.9495	3.98	0.00	362681				



Quantitation Results Report (QT Reviewed)

Data File	G0224_048.0048.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 9:16:31 AM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

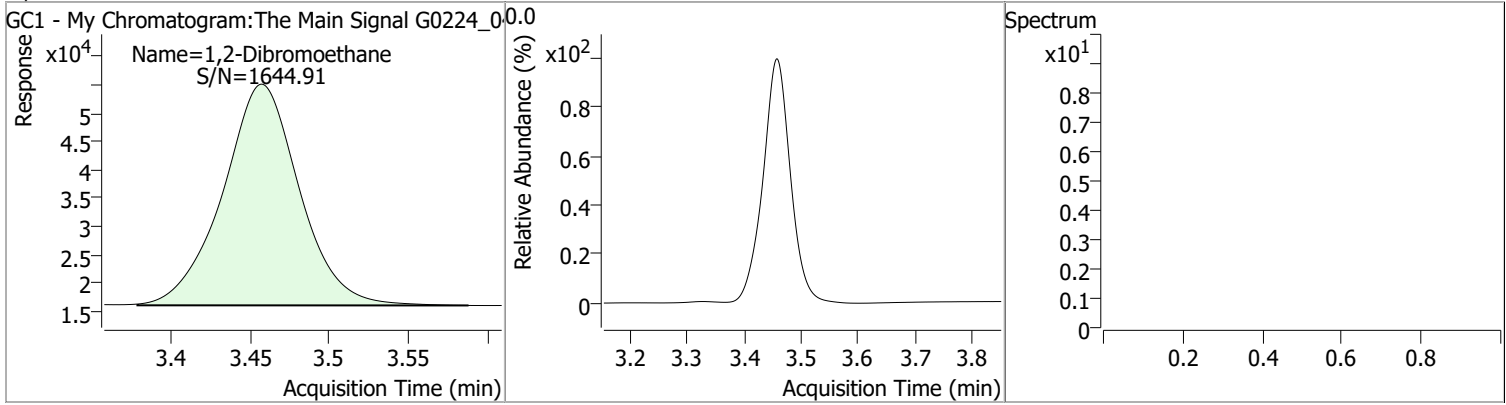


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.978	0.0	374159	0.9769	µg/L	0.004
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 976.94%		*
Target Compounds						
M 1,2-Dibromoethane	3.458	0.0	132080	0.9091	µg/L	QValue 100

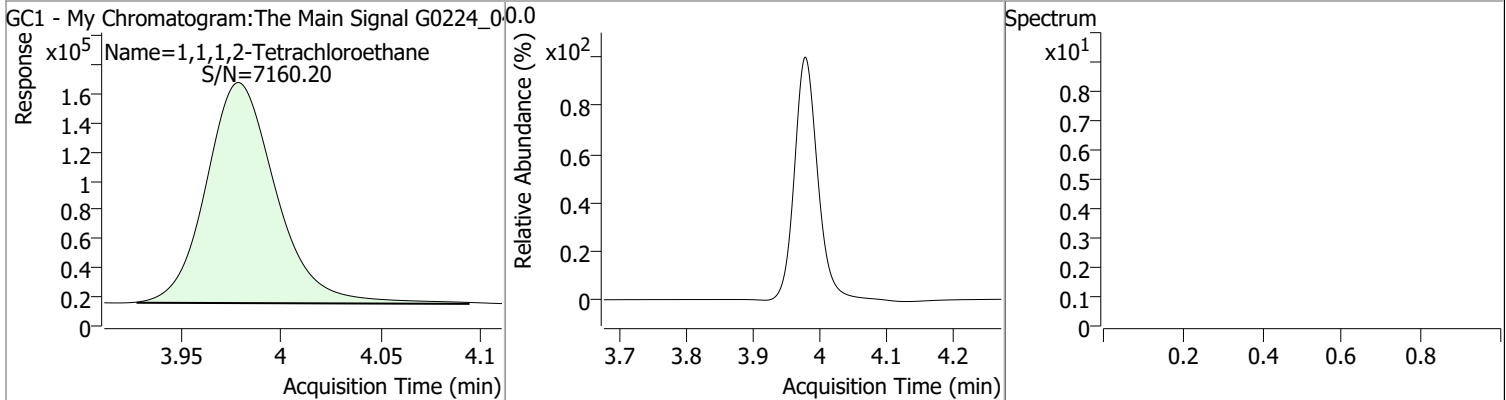
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.9091	3.46	0.00	132080				



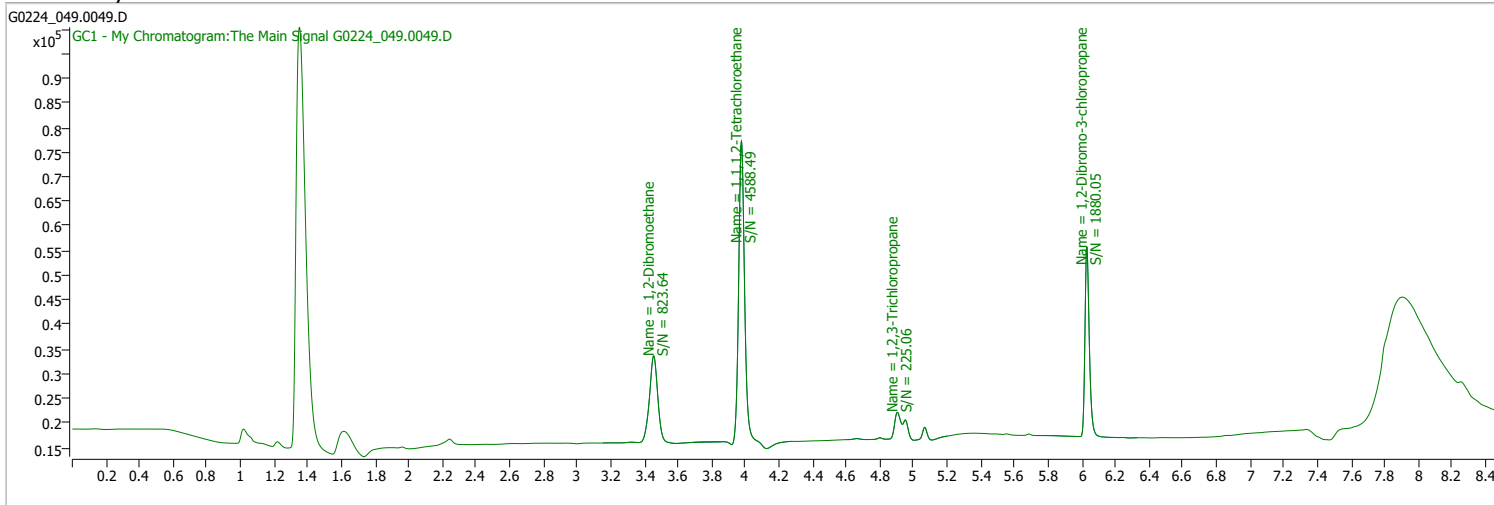
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.9769	3.98	0.00	374159				



Quantitation Results Report (QT Reviewed)

Data File	G0224_049.0049.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 9:36:14 AM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

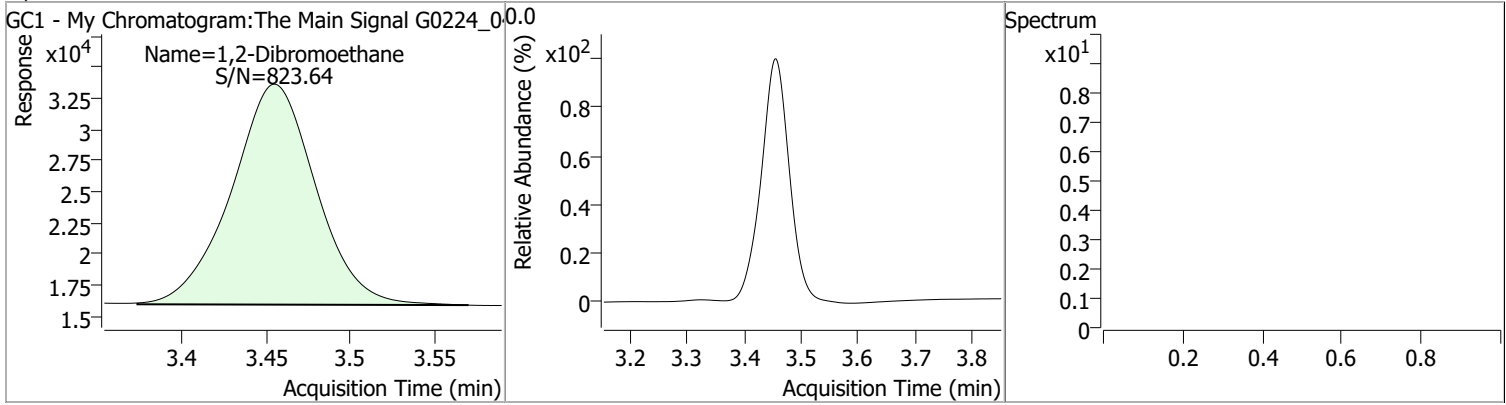


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.977	0.0	157427	0.4353	µg/L	0.002
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 435.26%	*	
Target Compounds						
M 1,2-Dibromoethane	3.455	0.0	61367	0.3846	µg/L	QValue 100

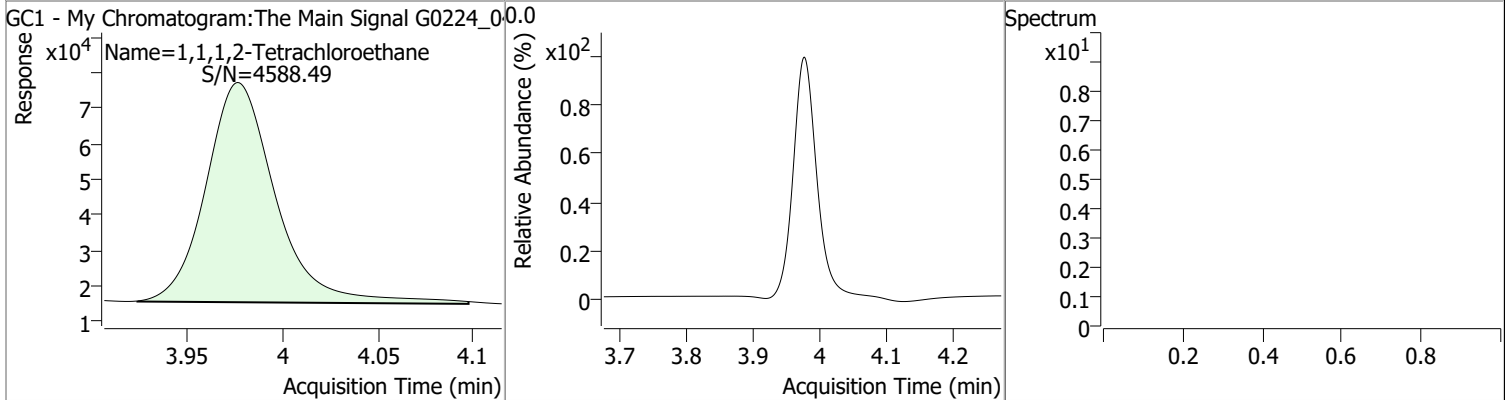
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.3846	3.46	0.00	61367				



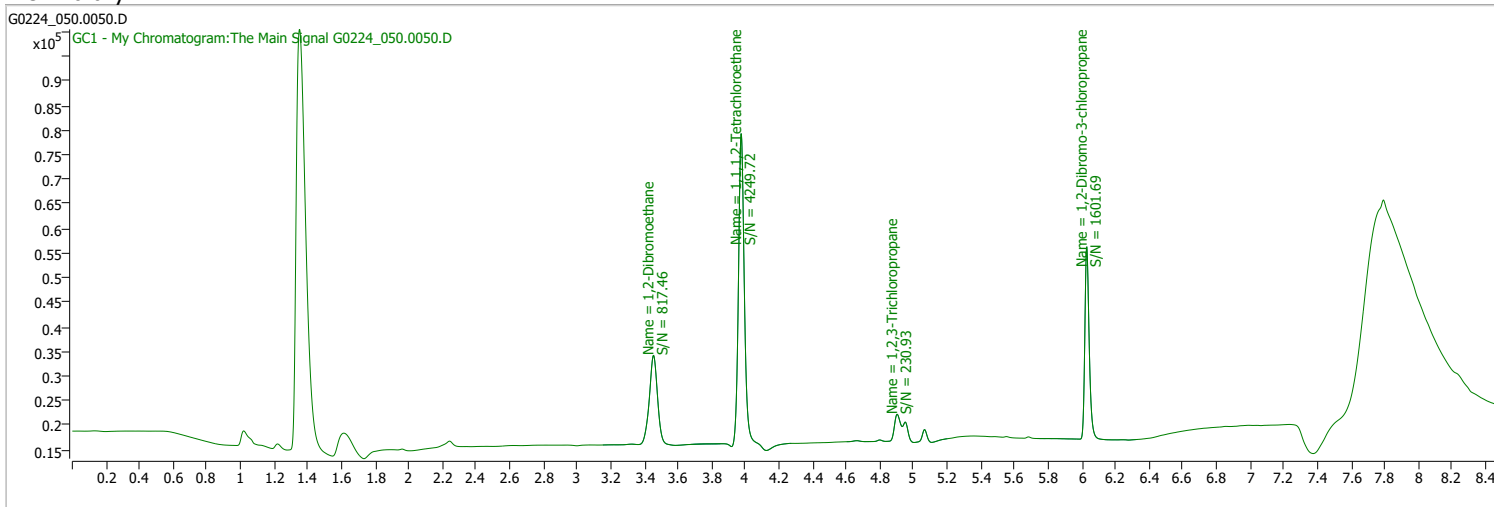
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.4353	3.98	0.00	157427				



Quantitation Results Report (QT Reviewed)

Data File	G0224_050.0050.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 9:56:08 AM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

S 1,1,1,2-Tetrachloroethane	3.975	0.0	159739	0.4413	µg/L	0.001
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 441.32%	*	

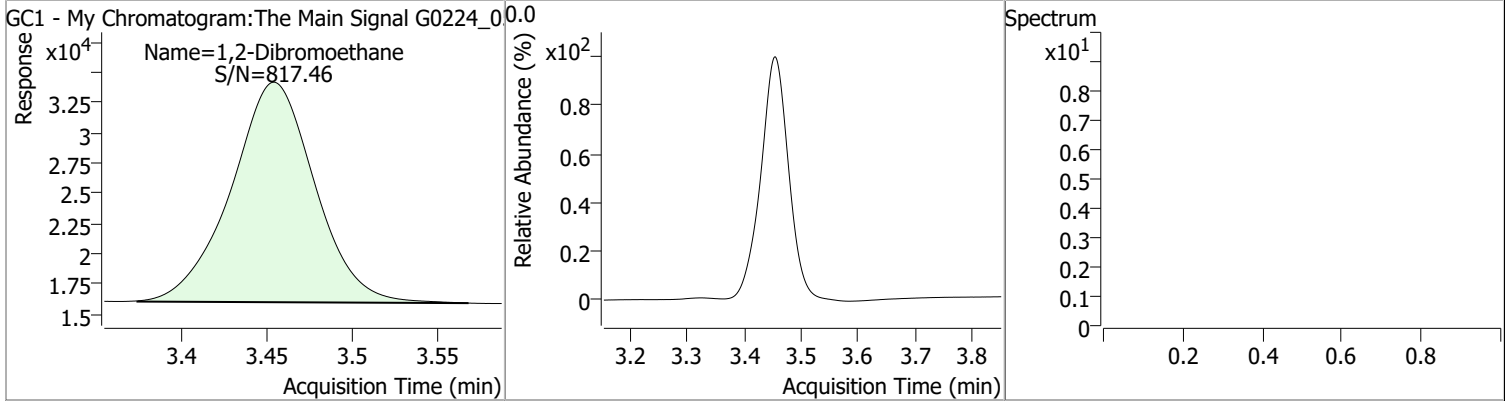
Target Compounds

M 1,2-Dibromoethane	3.454	0.0	62210	0.3903	µg/L	QValue 100
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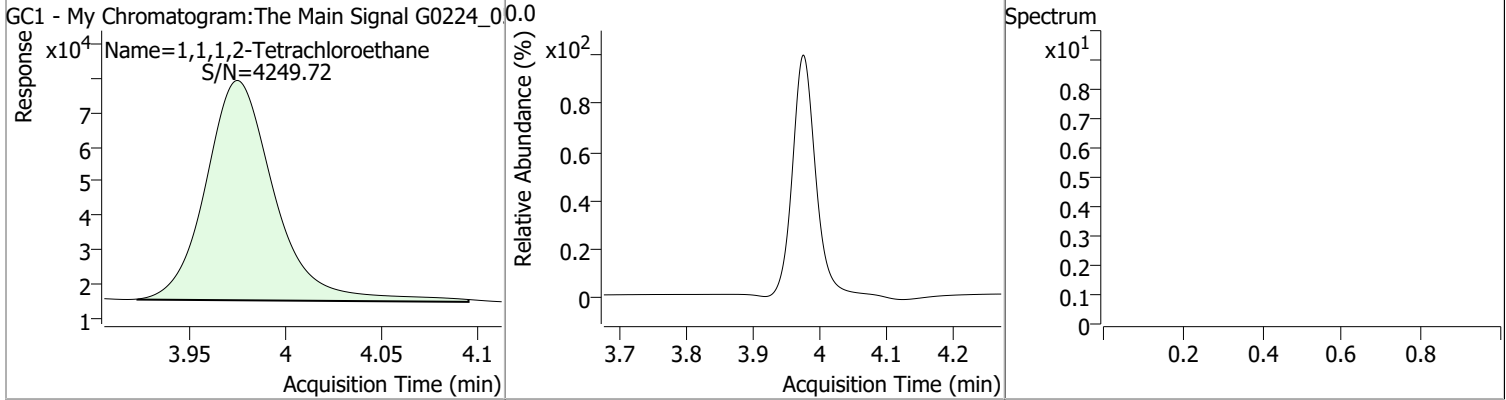
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.3903	3.45	0.00	62210				



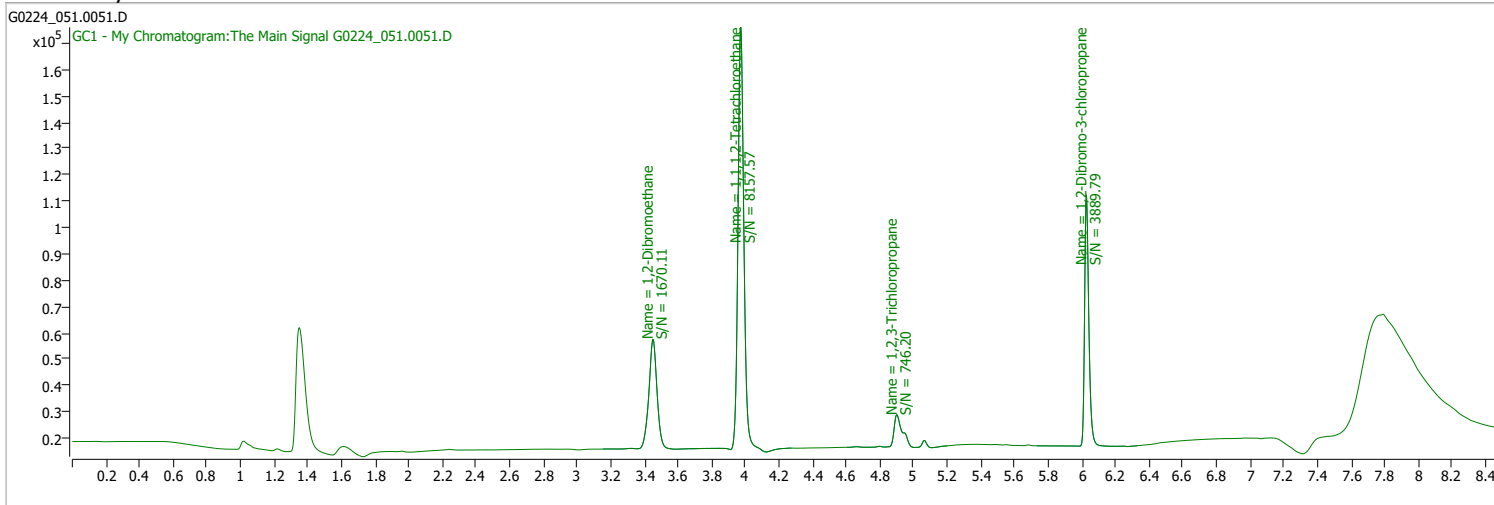
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.4413	3.98	0.00	159739				



Quantitation Results Report (QT Reviewed)

Data File	G0224_051.0051.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 10:15:47 AM
Sample Name	8011Primer	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

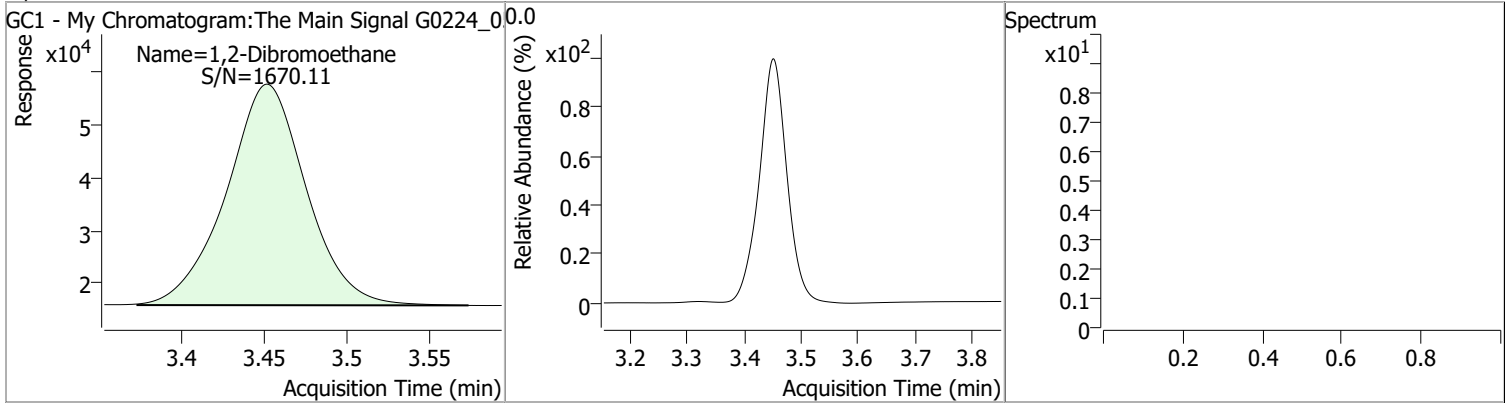


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	395059	1.0267	µg/L	-0.002
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 1026.68% *		
Target Compounds						
M 1,2-Dibromoethane	3.451	0.0	140701	0.9818	µg/L	QValue 100

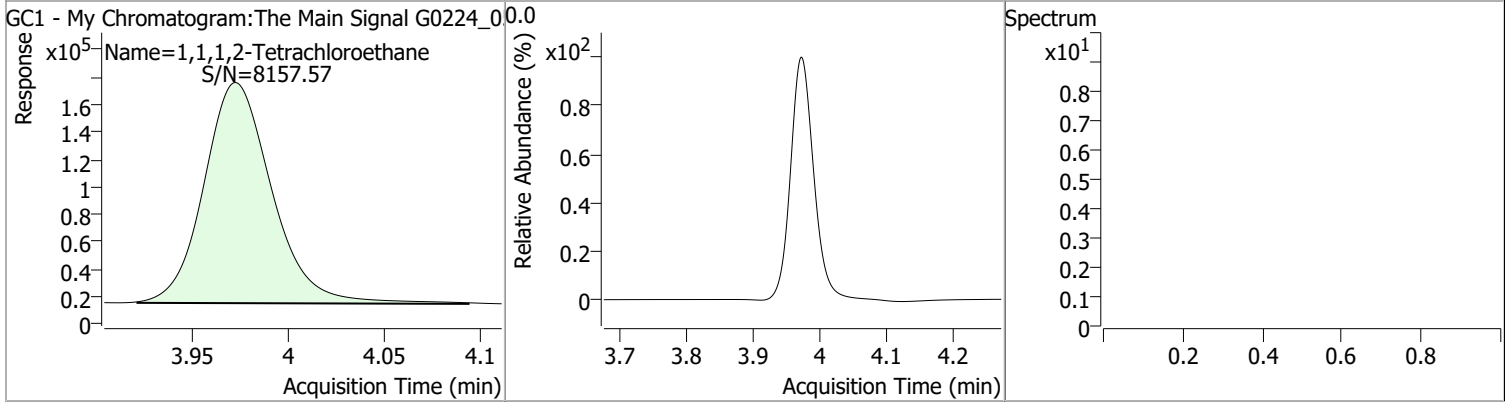
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.9818	3.45	0.00	140701				



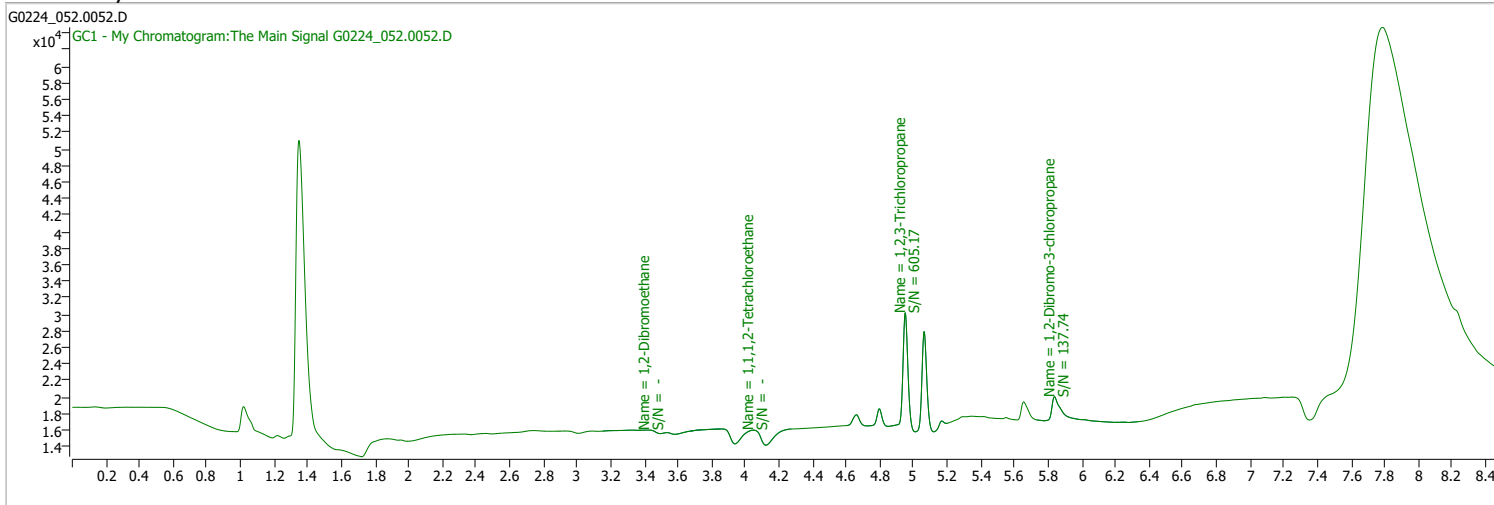
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	1.0267	3.97	0.00	395059				



Quantitation Results Report (QT Reviewed)

Data File	G0224_052.0052.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/25/2022 10:35:30 AM
Sample Name	Hexane	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

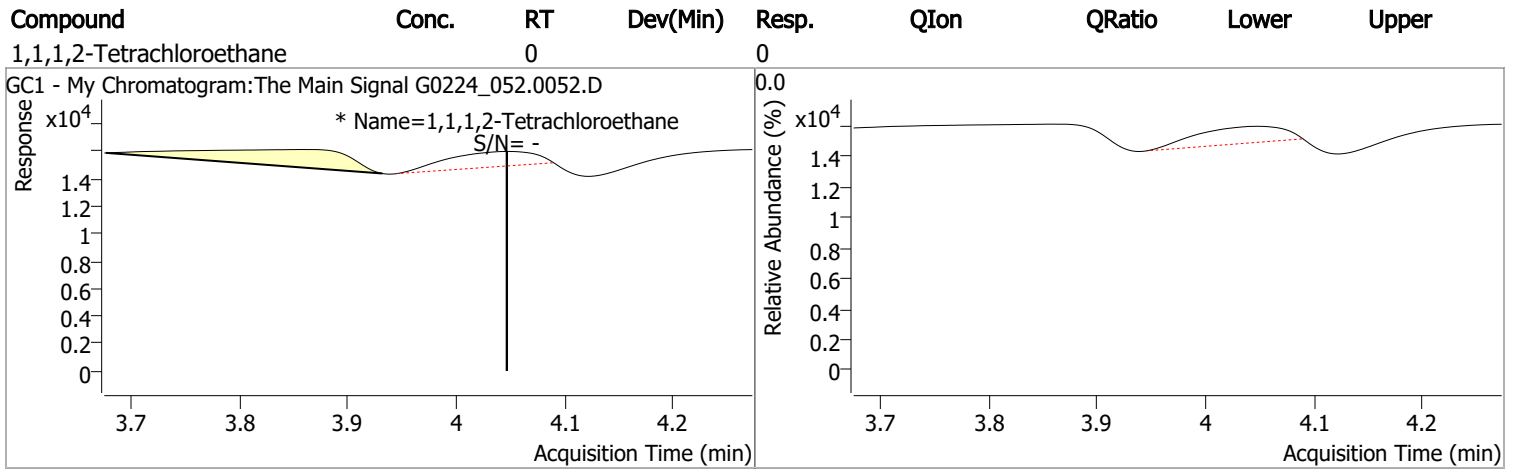
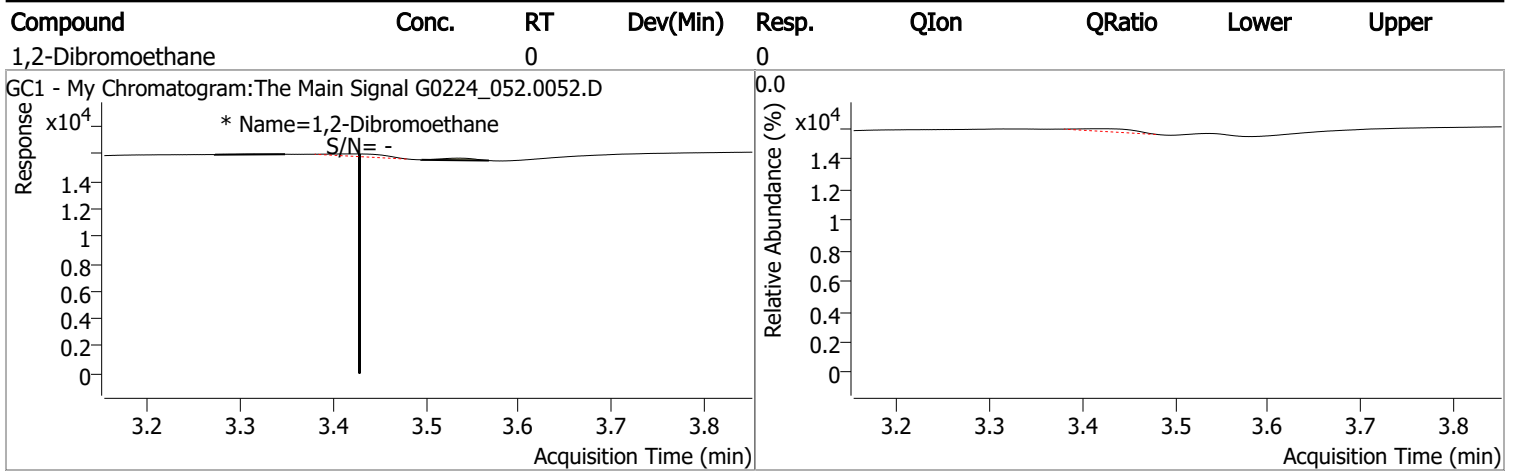
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.047	0.0	0		µg/L	md
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.428	0.0	0		µg/L	md
						QValue
						1

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

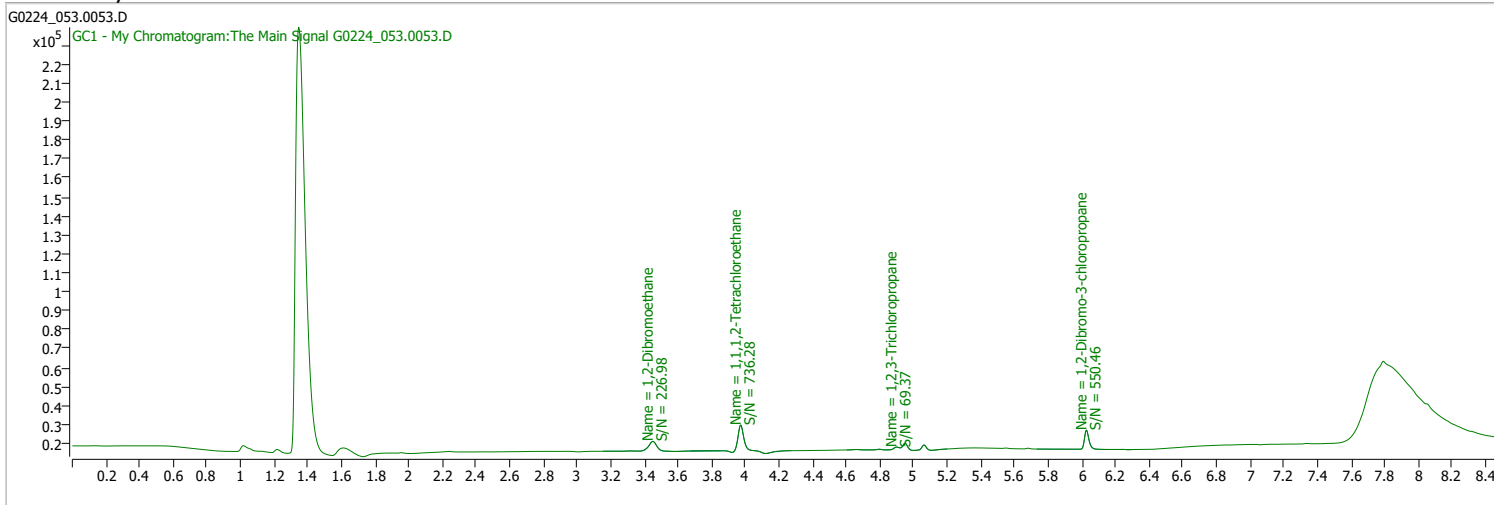
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_053.0053.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/25/2022 10:55:11 AM
Sample Name	CAL3-163984	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

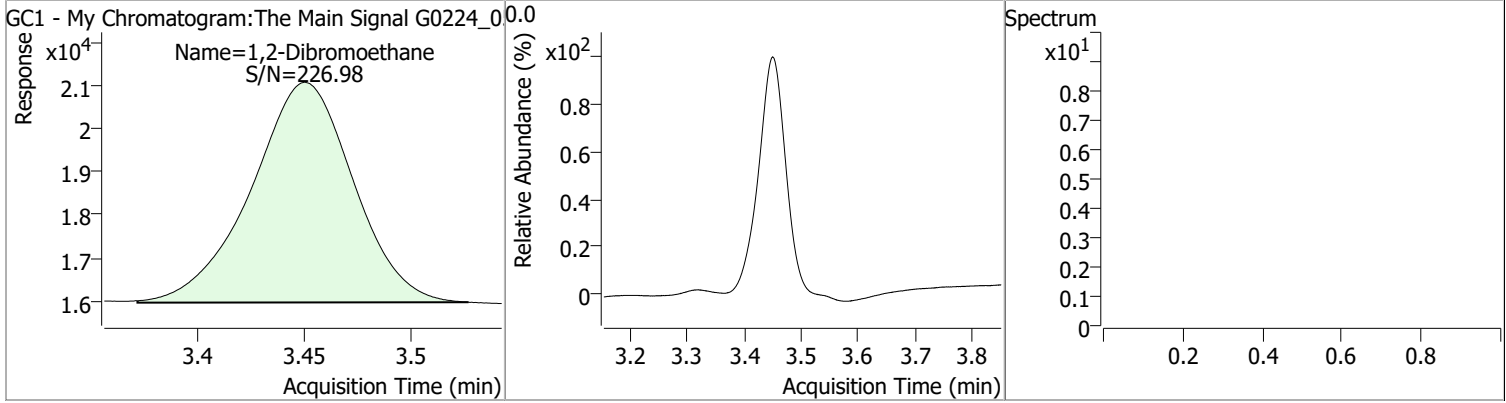


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	32143	0.0969	µg/L	-0.002
Spiked Amount: 0.100				Range: 70.0 - 130.0% Recovery = 96.89%		
Target Compounds						
M 1,2-Dibromoethane	3.450	0.0	17084	0.1022	µg/L	QValue 100

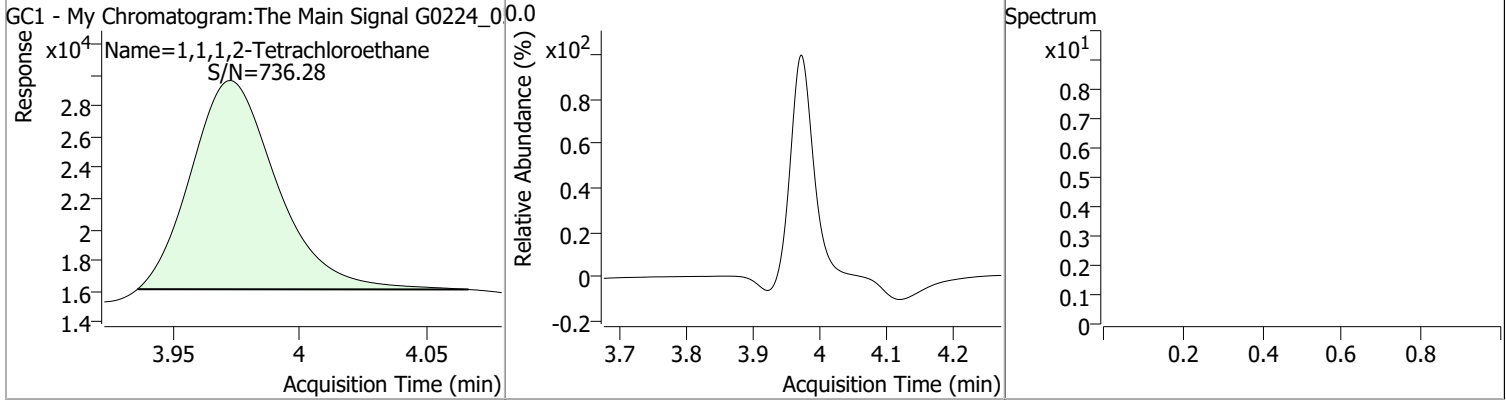
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1022	3.45	0.00	17084				



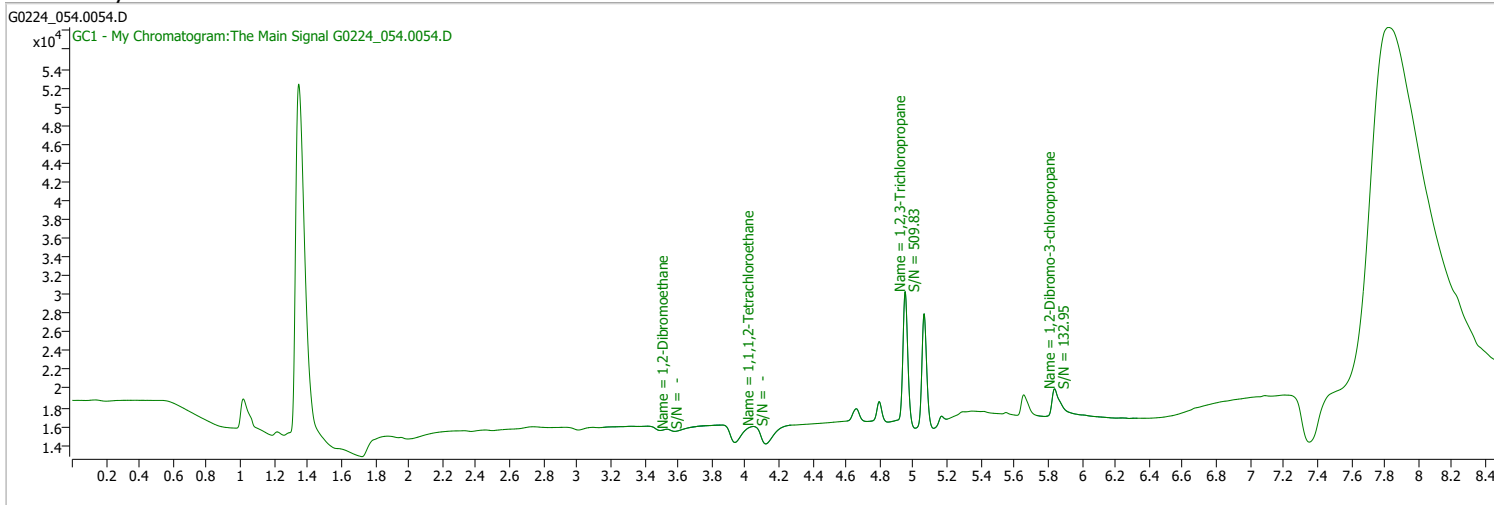
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0969	3.97	0.00	32143				



Quantitation Results Report (QT Reviewed)

Data File	G0224_054.0054.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 11:14:57 AM
Sample Name	Hexane	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

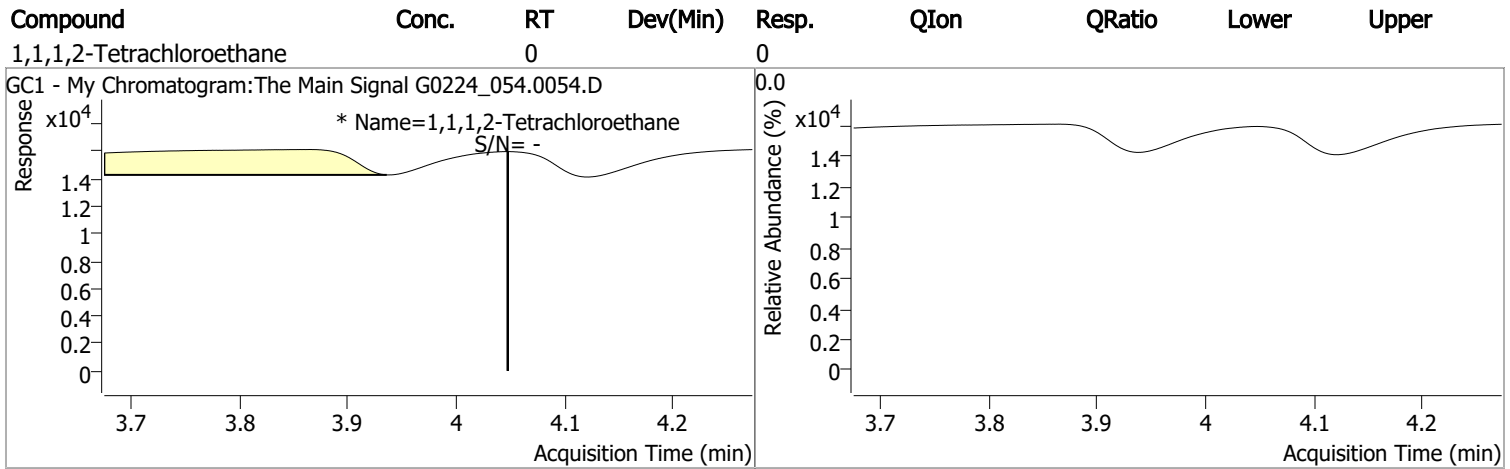
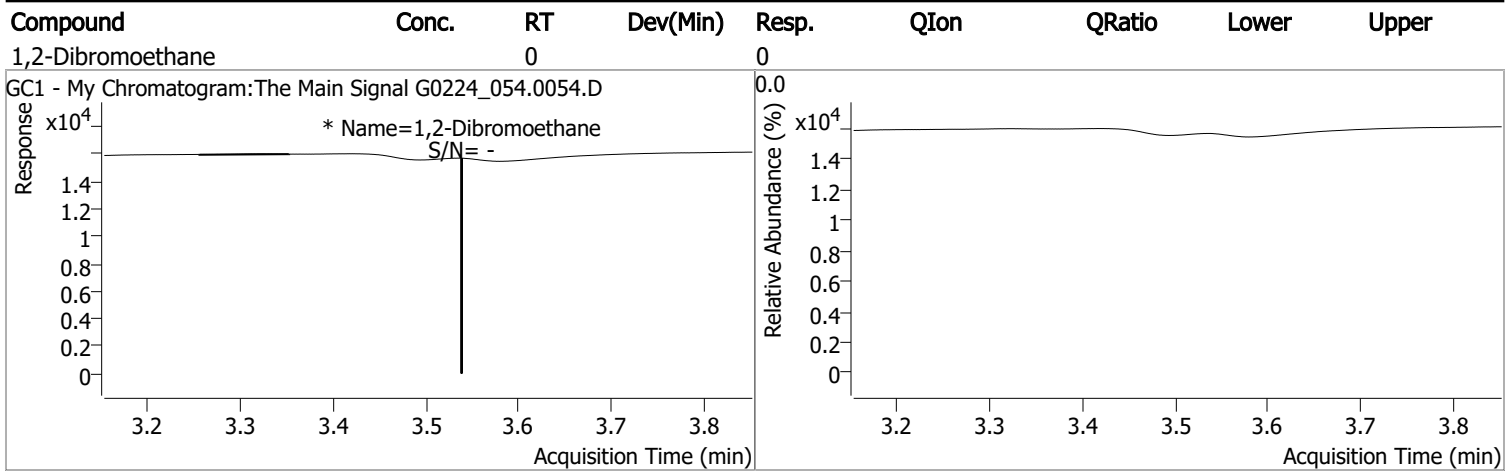
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.048	0.0	0		µg/L	md
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.538	0.0	0		µg/L	md
						QValue
						1

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

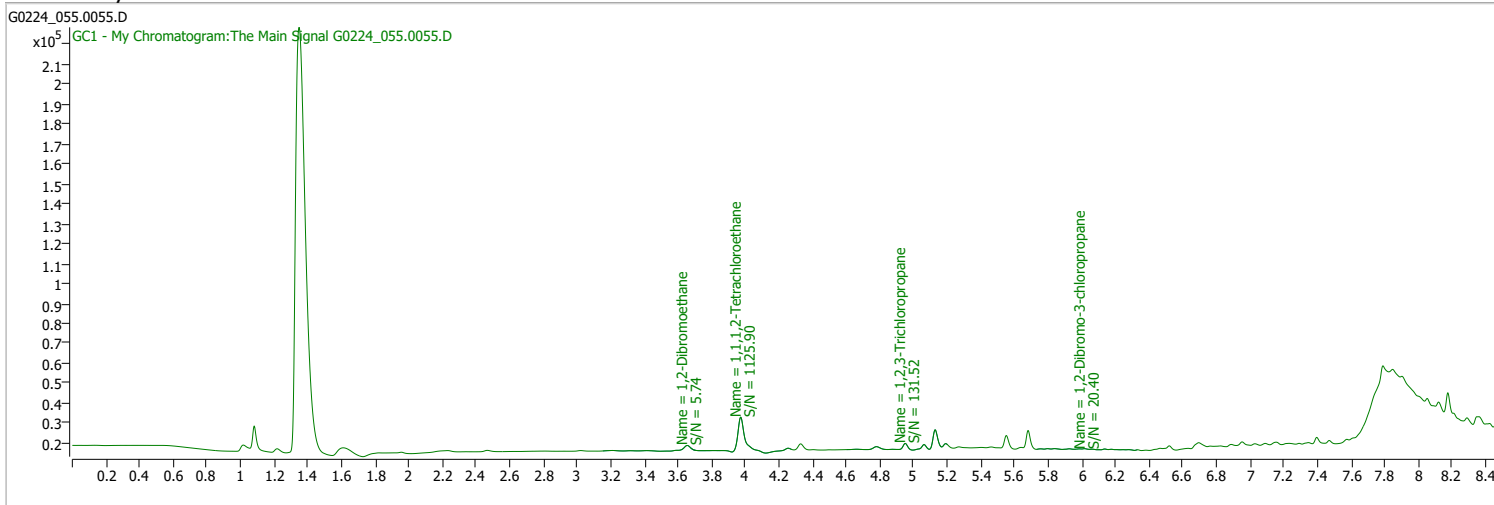
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0224_055.0055.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 11:34:46 AM
Sample Name	B22021435-032H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	re-extract
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

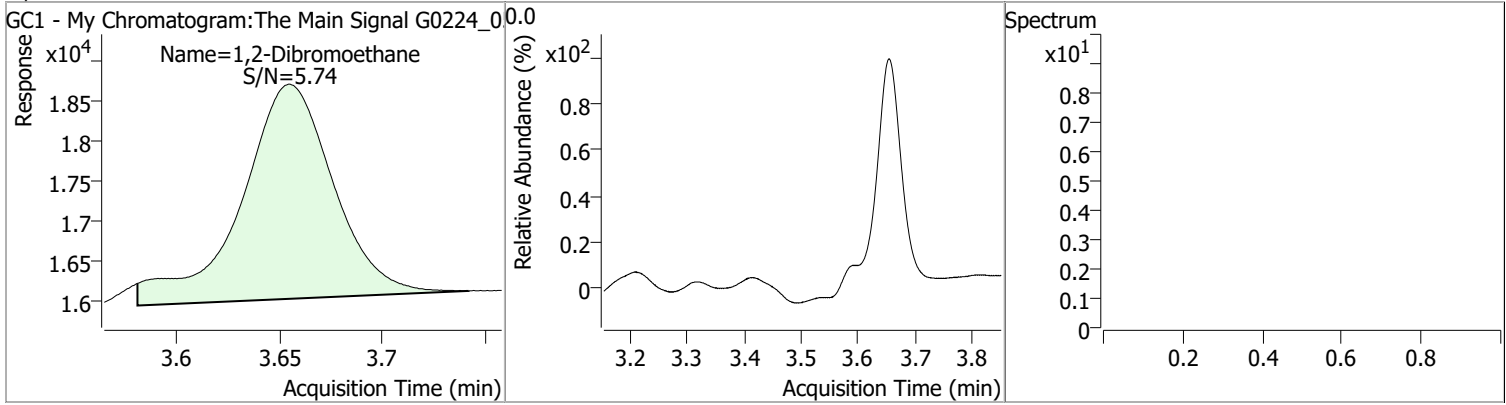


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	45726	0.1346	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 134.61%		*
Target Compounds						
M 1,2-Dibromoethane	3.655	0.0	8425	0.0500	µg/L	QValue 100

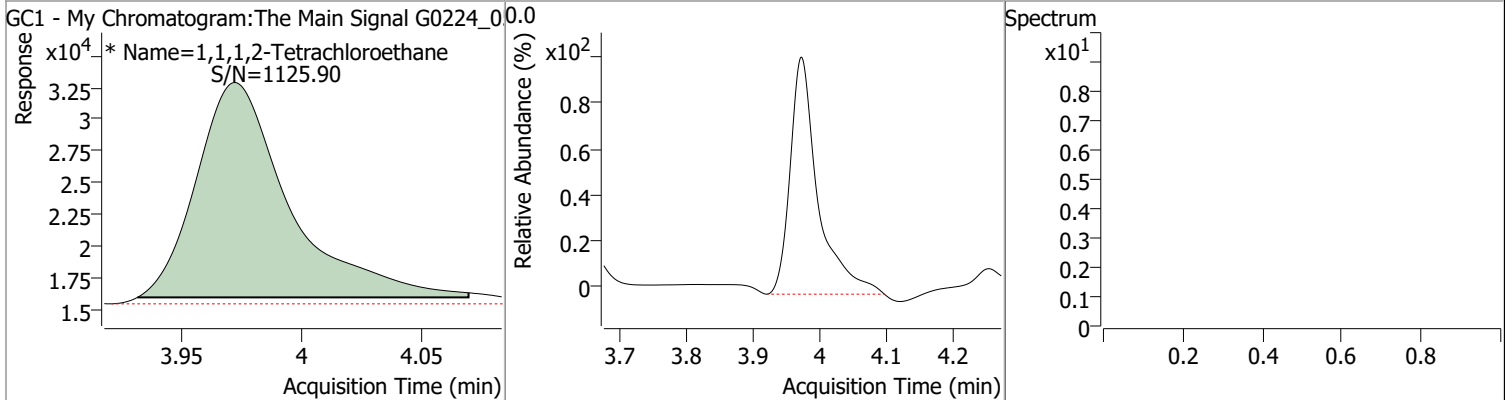
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0500	3.66	0.20	8425				



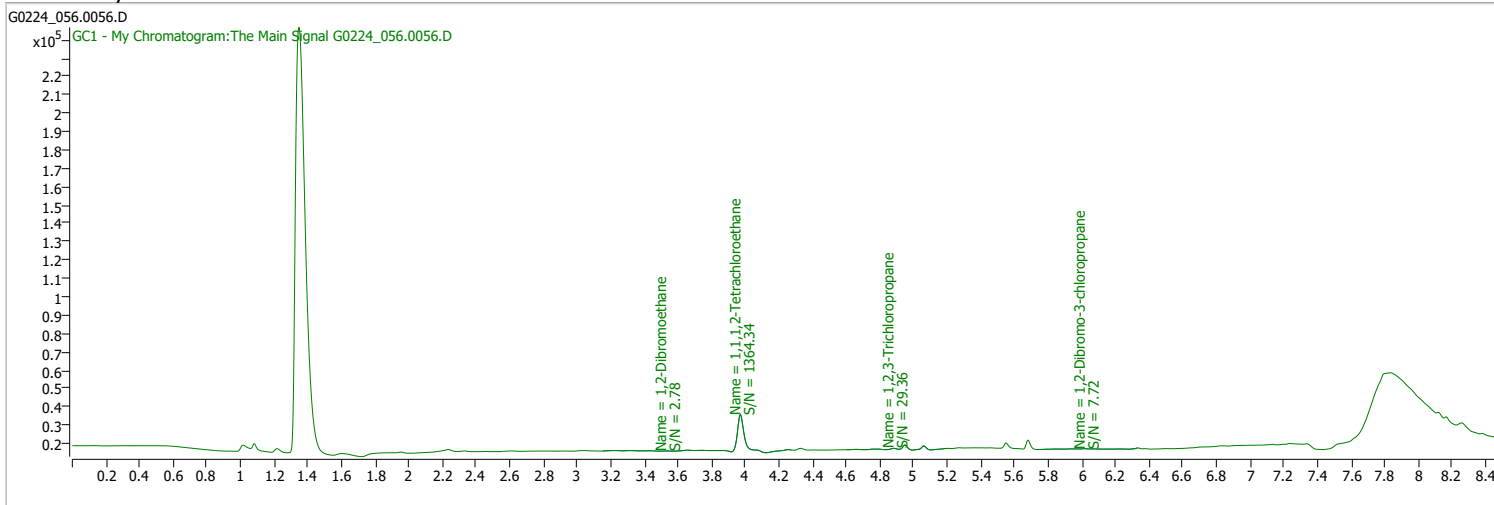
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1346	3.97	0.00	45726 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0224_056.0056.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 11:54:37 AM
Sample Name	B22021435-035A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	re-extract
Tune File		Tune Date	
Batch Name	G022422_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

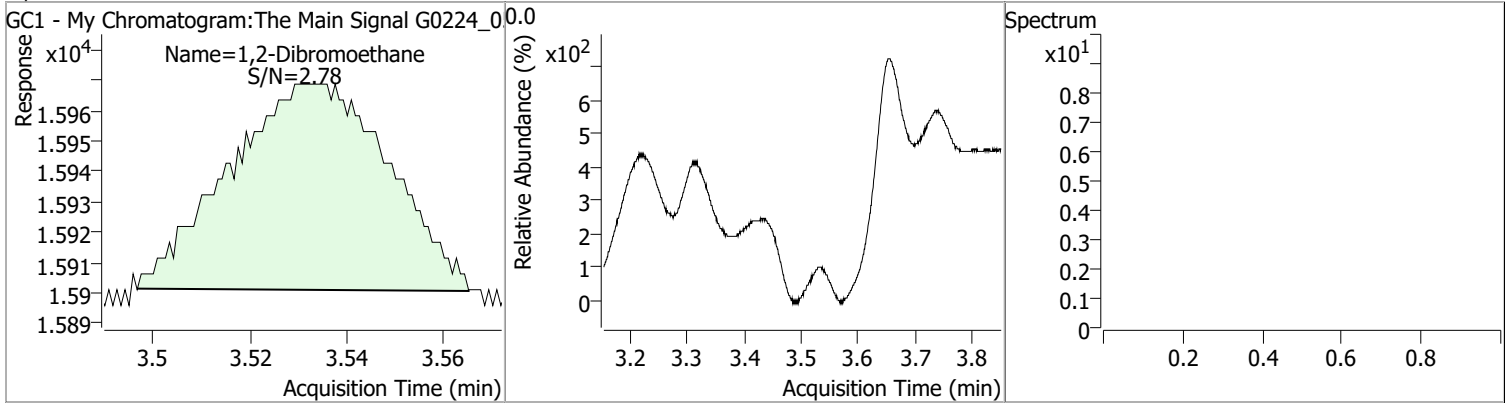


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	48775	0.1430	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 143.04%		*
Target Compounds						
M 1,2-Dibromoethane	3.529	0.0	157	0.0009	µg/L	QValue 100

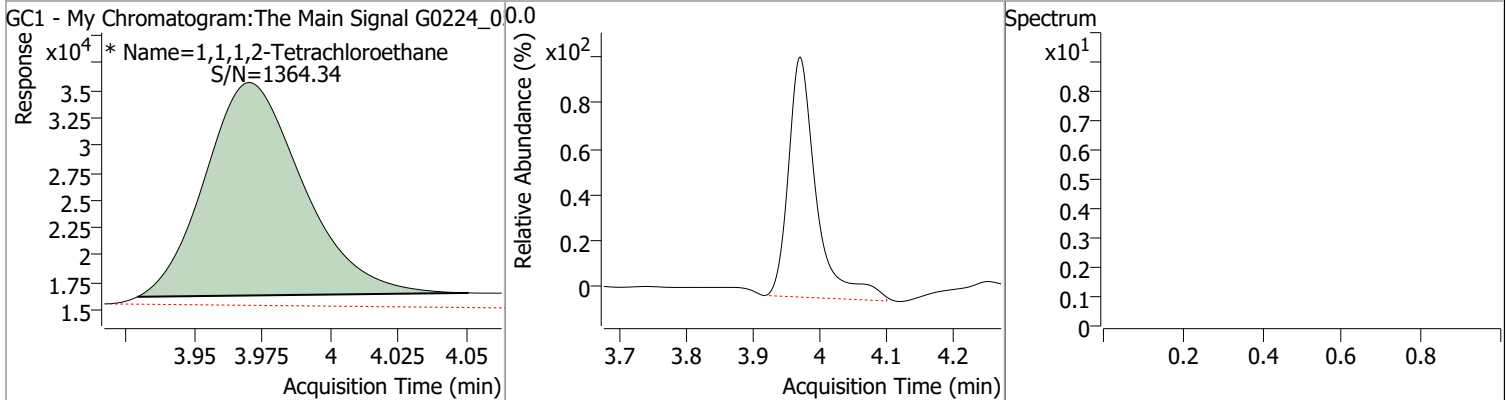
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.0009	3.53	0.08	157				



Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1430	3.97	0.00	48775 (m)				



Audit Trail report

Batch name and path: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin
Quant batch version: 10.0
Quant reporting version: 10.0

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdNewBatchTable	BL2000\ctran	2/24/2022 12:02:23 PM	Create new batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/24/2022 12:02:26 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_001.0001.D			✓	
CmdStartMethodEditing	BL2000\ctran	2/24/2022 12:04:12 PM	Start method editing			✓	
CmdImportMethodFromFile	BL2000\ctran	2/24/2022 12:04:13 PM	Import method from file \\MASSHUNTER\Org\Data\GECD.I\GECD_methods\G022322_8011_W_CLT.m			✓	
CmdApplyMethodToAllSamples	BL2000\ctran	2/24/2022 12:04:22 PM	Apply method to all samples			✓	
CmdMethodClear	BL2000\ctran	2/24/2022 12:04:22 PM	Clear method			✓	
CmdEndMethodEditing	BL2000\ctran	2/24/2022 12:04:23 PM	End method editing			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 12:04:23 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/24/2022 12:07:17 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/24/2022 12:21:18 PM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/24/2022 12:22:01 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_002.0002.D			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 12:22:03 PM	Quantitate all compounds in all samples			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 12:23:09 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/24/2022 12:25:10 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/24/2022 2:15:31 PM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\ctran	2/24/2022 2:15:46 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_007.0007.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_006.0006.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_005.0005.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_004.0004.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_003.0003.D			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:15:51 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\ctran	2/24/2022 2:17:22 PM	Set SampleType = Calibration for sample G0224_007.0007.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/24/2022 2:17:24 PM	Set LevelName = 1 for sample G0224_007.0007.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:17:26 PM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/24/2022 2:18:14 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_007.0007.D, from x, y = 3.961, 16214 to 4.033, 16182, result = 1386; previous integration is from x, y = 3.935, 15109 to 4.122, 14851 and previous response = 11629.			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:23:22 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/24/2022 2:23:23 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/24/2022 2:54:49 PM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/24/2022 2:55:00 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_009.0009.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_008.0008.D			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:55:02 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\ctran	2/24/2022 2:55:06 PM	Set SampleType = Calibration for sample G0224_008.0008.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/24/2022 2:55:09 PM	Set LevelName = 7 for sample G0224_008.0008.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/24/2022 2:55:11 PM	Set SampleType = Calibration for sample G0224_009.0009.D; previous value = Sample			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/24/2022 2:55:13 PM	Set LevelName = 2 for sample G0224_009.0009.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:55:15 PM	Quantitate all compounds in all samples			✓	
CmdUpdateRetentionTimes	BL2000\ctran	2/24/2022 2:56:02 PM	Update retention time for compound 1,1,1,2-Tetrachloroethane;			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:56:05 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/24/2022 2:56:07 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/24/2022 2:56:20 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_009.0009.D, from x, y = 3.945, 16303 to 4.062, 16198, result = 14968; previous integration is from x, y = 3.928, 15359 to 4.121, 14908 and previous response = 25109.			✓	
CmdQuantitate	BL2000\ctran	2/24/2022 2:56:27 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/24/2022 2:56:28 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdSaveBatchTable	BL2000\ctran	2/24/2022 3:31:58 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/25/2022 8:07:50 AM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\ctran	2/25/2022 8:08:11 AM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_046.0046.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_045.0045.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_044.0044.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_043.0043.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_042.0042.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_041.0041.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_040.0040.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_039.0039.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_038.0038.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_037.0037.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_036.0036.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_035.0035.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_034.0034.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_033.0033.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_032.0032.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_031.0031.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_030.0030.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_029.0029.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_028.0028.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_027.0027.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_026.0026.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_025.0025.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_024.0024.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_023.0023.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_022.0022.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_021.0021.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_020.0020.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_019.0019.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_018.0018.D,			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
			\\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_017.0017.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_016.0016.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_015.0015.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_014.0014.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_013.0013.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_012.0012.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_011.0011.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_010.0010.D				
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:23 AM	Set SampleType = Calibration for sample G0224_010.0010.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:25 AM	Set LevelName = 3 for sample G0224_010.0010.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:27 AM	Set SampleType = Calibration for sample G0224_011.0011.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:30 AM	Set LevelName = 4 for sample G0224_011.0011.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:32 AM	Set SampleType = Calibration for sample G0224_012.0012.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:35 AM	Set LevelName = 5 for sample G0224_012.0012.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:37 AM	Set SampleType = Calibration for sample G0224_013.0013.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:39 AM	Set LevelName = 6 for sample G0224_013.0013.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:42 AM	Set SampleType = DoubleBlank for sample G0224_014.0014.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:45 AM	Set SampleType = QC for sample G0224_015.0015.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:48 AM	Set LevelName = LCS for sample G0224_015.0015.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:51 AM	Set SampleType = CC for sample G0224_016.0016.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:53 AM	Set LevelName = 3 for sample G0224_016.0016.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:55 AM	Set SampleType = Blank for sample G0224_017.0017.D; previous value = Sample			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:57 AM	Set SampleType = QC for sample G0224_018.0018.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:08:59 AM	Set LevelName = LCS for sample G0224_018.0018.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:04 AM	Set SampleType = QC for sample G0224_019.0019.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:06 AM	Set LevelName = LCS1 for sample G0224_019.0019.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:08 AM	Set SampleType = DoubleBlank for sample G0224_020.0020.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:23 AM	Set SampleType = MatrixBlank for sample G0224_029.0029.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:26 AM	Set SampleType = Matrix for sample G0224_030.0030.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:28 AM	Set SampleType = MatrixDup for sample G0224_031.0031.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:32 AM	Set SampleType = DoubleBlank for sample G0224_032.0032.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:33 AM	Set SampleType = DoubleBlank for sample G0224_034.0034.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:35 AM	Set SampleType = CC for sample G0224_033.0033.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:38 AM	Set LevelName = 5 for sample G0224_033.0033.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:43 AM	Set SampleType = DoubleBlank for sample G0224_045.0045.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:45 AM	Set SampleType = CC for sample G0224_046.0046.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:09:47 AM	Set LevelName = 3 for sample G0224_046.0046.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 8:09:50 AM	Quantitate all compounds in all samples			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:10:15 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_009.0009.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:10:23 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_010.0010.D, from x, y = 3.937, 16177 to 4.053, 16323, result = 31992; previous integration is from x, y = 3.924, 15484 to 4.088, 15842 and previous response = 37124.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/25/2022 8:10:25 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_010.0010.D to y = 16177, new integration is from x, y = 3.937, 16177 to 4.053, 16177 and new response = 32499; previous integration is from x, y = 3.937, 16177 to 4.053, 16323 and previous response = 31992.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:10:26 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_010.0010.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:10:43 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_011.0011.D, from x, y = 3.932, 16370 to 4.063, 16276, result = 67779; previous integration is from x, y = 3.920, 15543 to 4.094, 14970 and previous response = 77954.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:10:44 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_011.0011.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:10:52 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_012.0012.D, from x, y = 3.927, 16583 to 4.072, 16286, result = 147845; previous integration is from x, y = 3.921, 15934 to 4.093, 15068 and previous response = 156878.			✓	
CmdManuallyIntegrateSnapBaseline	BL2000\ctran	2/25/2022 8:11:03 AM	Snap baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_013.0013.D, from x = 3.920 to x = 4.097, new integration is from x, y = 3.920, 16599 to 4.097, 15641 and new response = 382731; previous integration is from x, y = 3.920, 15819 to 4.097, 15081 and previous response = 389829.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:11:07 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_013.0013.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:11:10 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_012.0012.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdCalibrate	BL2000\ctran	2/25/2022 8:11:20 AM	Replace level 6 with Calibration sample G0224_013.0013.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane}; Replace level 5 with Calibration sample G0224_012.0012.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane}; Replace level 4 with Calibration sample G0224_011.0011.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane}; Replace level 3 with Calibration sample G0224_010.0010.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane}; Replace level 2 with Calibration sample G0224_009.0009.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane}; Replace level 7 with Calibration sample G0224_008.0008.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane}; Replace level 1 with Calibration sample G0224_007.0007.D for compounds {1,2-Dibromo-3-chloropropane, 1,2,3-Trichloropropane, 1,2-Dibromoethane, 1,1,1,2-Tetrachloroethane};			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 8:11:24 AM	Quantitate all compounds in all samples			✓	
CmdStartMethodEditing	BL2000\ctran	2/25/2022 8:11:52 AM	Start method editing			✓	
CmdImportMethodFrom Sample	BL2000\ctran	2/25/2022 8:11:52 AM	Import method from sample G0224_012.0012.D			✓	
CmdSaveMethodAs	BL2000\ctran	2/25/2022 8:12:18 AM	Save method to file \\MASSHUNTER\Org\Data\GEC.D.I\GEC D_methods\G022422_8011_W_CLT.m			✓	
CmdApplyMethodToAll Samples	BL2000\ctran	2/25/2022 8:12:22 AM	Apply method to all samples			✓	
CmdMethodClear	BL2000\ctran	2/25/2022 8:12:22 AM	Clear method			✓	
CmdEndMethodEditing	BL2000\ctran	2/25/2022 8:12:23 AM	End method editing			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 8:12:25 AM	Quantitate all compounds in all samples			✓	
CmdUpdateRetentionTimes	BL2000\ctran	2/25/2022 8:12:35 AM	Update retention time for compound 1,2,3-Trichloropropane; 1,2-Dibromo-3-chloropropane; 1,2-Dibromoethane; 1,1,1,2-Tetrachloroethane;			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdQuantitate	BL2000\ctran	2/25/2022 8:12:39 AM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:00 AM	Set SampleApproved = True for sample G0224_013.0013.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:01 AM	Set SampleApproved = True for sample G0224_012.0012.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:01 AM	Set SampleApproved = True for sample G0224_011.0011.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:02 AM	Set SampleApproved = True for sample G0224_010.0010.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:02 AM	Set SampleApproved = True for sample G0224_009.0009.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:02 AM	Set SampleApproved = True for sample G0224_008.0008.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:14:03 AM	Set SampleApproved = True for sample G0224_007.0007.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:14:16 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_015.0015.D, from x, y = 3.933, 16047 to 4.057, 16203, result = 33999; previous integration is from x, y = 3.920, 15432 to 4.098, 14921 and previous response = 43119.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/25/2022 8:14:17 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_015.0015.D to y = 16047, new integration is from x, y = 3.933, 16047 to 4.057, 16047 and new response = 34577; previous integration is from x, y = 3.933, 16047 to 4.057, 16203 and previous response = 33999.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:14:39 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_016.0016.D, from x, y = 3.936, 16370 to 4.055, 16219, result = 31159; previous integration is from x, y = 3.929, 15641 to 4.103, 14908 and previous response = 40314.			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:14:58 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_017.0017.D			✓	
CmdClearManualIntegration	BL2000\ctran	2/25/2022 8:15:06 AM	Clear manual integration of target signal for compound 1,2-Dibromoethane in sample G0224_017.0017.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:15:09 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_017.0017.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:15:18 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_018.0018.D, from x, y = 3.935, 16495 to 4.067, 16032, result = 32624; previous integration is from x, y = 3.919, 15536 to 4.103, 15020 and previous response = 42117.			✓	
CmdManuallyIntegrateSnapBaseline	BL2000\ctran	2/25/2022 8:15:19 AM	Snap baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_018.0018.D, from x = 3.935 to x = 4.067, new integration is from x, y = 3.935, 16495 to 4.067, 16234 and new response = 31825; previous integration is from x, y = 3.935, 16495 to 4.067, 16032 and previous response = 32624.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:15:22 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_018.0018.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:15:26 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_019.0019.D, from x, y = 3.933, 16411 to 4.059, 16396, result = 34064; previous integration is from x, y = 3.918, 15641 to 4.102, 15115 and previous response = 43834.			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:15:30 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_020.0020.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:15:31 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_020.0020.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:15:42 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_021.0021.D, from x, y = 3.928, 16313 to 4.058, 16292, result = 42127; previous integration is from x, y = 3.915, 15620 to 4.094, 15081 and previous response = 51299.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:15:53 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_022.0022.D, from x, y = 3.932, 16589 to 4.056, 16484, result = 31040; previous integration is from x, y = 3.917, 15448 to 4.098, 15024 and previous response = 43454.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:01 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_023.0023.D, from x, y = 3.931, 16323 to 4.062, 16219, result = 38959; previous integration is from x, y = 3.916, 15406 to 4.099, 14881 and previous response = 49761.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:06 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_024.0024.D, from x, y = 3.933, 16344 to 4.048, 16422, result = 30764; previous integration is from x, y = 3.928, 15828 to 4.100, 14991 and previous response = 39797.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:12 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_025.0025.D, from x, y = 3.933, 16375 to 4.051, 16359, result = 32157; previous integration is from x, y = 3.917, 15357 to 4.100, 14867 and previous response = 44286.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:17 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_026.0026.D, from x, y = 3.933, 16323 to 4.044, 16411, result = 32943; previous integration is from x, y = 3.918, 15365 to 4.052, 15039 and previous response = 41697.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:21 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_027.0027.D, from x, y = 3.931, 16229 to 4.050, 16307, result = 34989; previous integration is from x, y = 3.918, 15511 to 4.096, 15011 and previous response = 44769.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:28 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_028.0028.D, from x, y = 3.933, 16313 to 4.055, 16403, result = 31897; previous integration is from x, y = 3.921, 15375 to 4.055, 16403 and previous response = 35174.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:35 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_029.0029.D, from x, y = 3.932, 16109 to 4.059, 16172, result = 33605; previous integration is from x, y = 3.918, 15307 to 4.090, 14837 and previous response = 43741.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:40 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_030.0030.D, from x, y = 3.933, 16411 to 4.046, 16240, result = 32008; previous integration is from x, y = 3.929, 15805 to 4.089, 14939 and previous response = 40281.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:45 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_031.0031.D, from x, y = 3.932, 16250 to 4.055, 16198, result = 32849; previous integration is from x, y = 3.918, 15344 to 4.093, 15344 and previous response = 40891.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:16:53 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_033.0033.D, from x, y = 3.921, 16276 to 4.058, 16359, result = 152467; previous integration is from x, y = 3.917, 15846 to 4.088, 14923 and previous response = 161426.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/25/2022 8:16:53 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_033.0033.D to y = 16276, new integration is from x, y = 3.921, 16276 to 4.058, 16276 and new response = 152809; previous integration is from x, y = 3.921, 16276 to 4.058, 16359 and previous response = 152467.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:16:55 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_033.0033.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:17:19 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_034.0034.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:17:21 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_034.0034.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:17:28 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_036.0036.D, from x, y = 3.933, 16148 to 4.053, 16240, result = 32653; previous integration is from x, y = 3.917, 15281 to 4.100, 14823 and previous response = 43774.			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:17:32 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_036.0036.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:17:40 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_037.0037.D, from x, y = 3.925, 16098 to 4.059, 16109, result = 39690; previous integration is from x, y = 3.915, 15354 to 4.087, 14868 and previous response = 49184.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:17:47 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_038.0038.D, from x, y = 3.931, 16292 to 4.048, 16323, result = 35877; previous integration is from x, y = 3.918, 15367 to 4.088, 14928 and previous response = 46963.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:17:54 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_039.0039.D, from x, y = 3.928, 16198 to 4.079, 15948, result = 45043; previous integration is from x, y = 3.915, 15447 to 4.093, 14964 and previous response = 53670.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:17:59 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_039.0039.D, from x, y = 3.928, 16198 to 4.049, 16012, result = 44451; previous integration is from x, y = 3.928, 16198 to 4.079, 15948 and previous response = 45043.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:18:16 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_040.0040.D, from x, y = 3.928, 16276 to 4.057, 16438, result = 49542; previous integration is from x, y = 3.915, 15495 to 4.099, 15023 and previous response = 60525.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:18:25 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_041.0041.D, from x, y = 3.928, 16281 to 4.058, 16203, result = 52432; previous integration is from x, y = 3.915, 15531 to 4.089, 15014 and previous response = 61699.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:18:32 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_042.0042.D, from x, y = 3.928, 16193 to 4.050, 16313, result = 45163; previous integration is from x, y = 3.916, 15484 to 4.094, 15484 and previous response = 52453.			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:19:18 AM	Set LevelName = LCS for sample G0224_030.0030.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 8:19:25 AM	Set LevelName = LCS for sample G0224_031.0031.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 8:19:27 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:22:43 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_046.0046.D, from x, y = 3.931, 16094 to 4.058, 16135, result = 34534; previous integration is from x, y = 3.917, 15333 to 4.100, 14826 and previous response = 44450.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 8:25:29 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_044.0044.D, from x, y = 3.928, 16370 to 4.058, 16307, result = 37984; previous integration is from x, y = 3.923, 15978 to 4.080, 16092 and previous response = 40659.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/25/2022 8:25:31 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_044.0044.D to y = 16307, new integration is from x, y = 3.928, 16307 to 4.058, 16307 and new response = 38229; previous integration is from x, y = 3.928, 16370 to 4.058, 16307 and previous response = 37984.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 8:25:33 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_044.0044.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 8:58:46 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_029.0029.D			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 8:58:49 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 9:00:01 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_017.0017.D, from x, y = 3.934, 16333 to 4.051, 16260, result = 32257; previous integration is from x, y = 3.927, 15647 to 4.097, 14974 and previous response = 41183.			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 9:03:52 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_041.0041.D			✓	
CmdRemoveSamples	BL2000\ctran	2/25/2022 9:46:55 AM	Remove 1 sample(s): Remove Sample sample B22010745-007A, data file G0224_021.0021.D ;			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/25/2022 9:47:07 AM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_021.0021.D			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/25/2022 9:47:22 AM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_048.0048.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_047.0047.D			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 9:56:13 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/25/2022 9:56:44 AM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\ctran	2/25/2022 12:26:45 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_056.0056.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_055.0055.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_054.0054.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_053.0053.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_052.0052.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_051.0051.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_050.0050.D, \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G0224_049.0049.D			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 12:26:50 PM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 12:27:01 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_056.0056.D, from x, y = 3.929, 16167 to 4.051, 16521, result = 48775; previous integration is from x, y = 3.918, 15521 to 4.101, 15058 and previous response = 59527.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 12:27:06 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_055.0055.D, from x, y = 3.933, 16047 to 4.067, 16370, result = 43633; previous integration is from x, y = 3.921, 15451 to 4.097, 15451 and previous response = 50885.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 12:27:09 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_055.0055.D, from x, y = 3.933, 16047 to 4.058, 15980, result = 44961; previous integration is from x, y = 3.933, 16047 to 4.067, 16370 and previous response = 43633.			✓	
CmdClearManualIntegration	BL2000\ctran	2/25/2022 12:27:12 PM	Clear manual integration of target signal for compound 1,1,1,2-Tetrachloroethane in sample G0224_055.0055.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 12:27:20 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_055.0055.D, from x, y = 3.932, 15958 to 4.070, 16323, result = 44213; previous integration is from x, y = 3.921, 15451 to 4.097, 15451 and previous response = 50885.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrate DropBaseline	BL2000\ctran	2/25/2022 12:27:21 PM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_055.0055.D to y = 15958, new integration is from x, y = 3.932, 15958 to 4.070, 15958 and new response = 45726; previous integration is from x, y = 3.932, 15958 to 4.070, 16323 and previous response = 44213.			✓	
CmdSaveBatchTable	BL2000\ctran	2/25/2022 1:27:26 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/25/2022 1:32:01 PM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:32:38 PM	Set SampleApproved = True for sample G0224_001.0001.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:32:39 PM	Set SampleApproved = True for sample G0224_002.0002.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:32:40 PM	Set SampleApproved = True for sample G0224_003.0003.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:32:40 PM	Set SampleApproved = True for sample G0224_004.0004.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:32:40 PM	Set SampleApproved = True for sample G0224_005.0005.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:32:44 PM	Set SampleType = DoubleBlank for sample G0224_006.0006.D; previous value = Sample			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:32:46 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_006.0006.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:32:48 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_006.0006.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:32:53 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_014.0014.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:33:04 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_014.0014.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:33:06 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_014.0014.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:33:08 PM	Set SampleApproved = True for sample G0224_014.0014.D; previous value = False			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:38:21 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_015.0015.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:38:25 PM	Set SampleApproved = True for sample G0224_015.0015.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:38:30 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_016.0016.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:38:32 PM	Set SampleApproved = True for sample G0224_016.0016.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:38:36 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_017.0017.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:38:39 PM	Set SampleApproved = True for sample G0224_017.0017.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:38:48 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_019.0019.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:38:52 PM	Set SampleApproved = True for sample G0224_018.0018.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:38:53 PM	Set SampleApproved = True for sample G0224_019.0019.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:38:55 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_020.0020.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:38:57 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_020.0020.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:38:58 PM	Set SampleApproved = True for sample G0224_020.0020.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:41:37 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_022.0022.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:41:40 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_022.0022.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:41:41 PM	Set SampleApproved = True for sample G0224_022.0022.D; previous value = False			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:41:44 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_023.0023.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:41:47 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_023.0023.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:41:48 PM	Set SampleApproved = True for sample G0224_023.0023.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:41:52 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_024.0024.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:41:56 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_024.0024.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:41:57 PM	Set SampleApproved = True for sample G0224_024.0024.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:41:59 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_025.0025.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:42:02 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_025.0025.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:03 PM	Set SampleApproved = True for sample G0224_025.0025.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:42:07 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_026.0026.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:42:10 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_026.0026.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:11 PM	Set SampleApproved = True for sample G0224_026.0026.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:42:14 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_027.0027.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:16 PM	Set SampleApproved = True for sample G0224_027.0027.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:42:17 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_027.0027.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:42:20 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_028.0028.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:42:23 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_028.0028.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:24 PM	Set SampleApproved = True for sample G0224_028.0028.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:42:26 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_029.0029.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:30 PM	Set SampleApproved = True for sample G0224_029.0029.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:42:37 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_030.0030.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:37 PM	Set SampleApproved = True for sample G0224_030.0030.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:42:40 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_031.0031.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:46 PM	Set SampleApproved = True for sample G0224_031.0031.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:42:49 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_032.0032.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:42:51 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_032.0032.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:42:58 PM	Set SampleApproved = True for sample G0224_033.0033.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:43:00 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_034.0034.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:43:02 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_034.0034.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:43:05 PM	Set SampleApproved = True for sample G0224_034.0034.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:43:10 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_035.0035.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:43:14 PM	Set SampleApproved = True for sample G0224_035.0035.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:43:17 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_036.0036.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:43:20 PM	Set SampleApproved = True for sample G0224_036.0036.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:43:22 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_037.0037.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:43:28 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_037.0037.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:43:29 PM	Set SampleApproved = True for sample G0224_037.0037.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:44:03 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_038.0038.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:44:06 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_038.0038.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:07 PM	Set SampleApproved = True for sample G0224_038.0038.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:44:09 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_039.0039.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:44:14 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_039.0039.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:20 PM	Set Comment = re-extract for sample G0224_039.0039.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:25 PM	Set Comment = re-extract for sample G0224_040.0040.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:25 PM	Set SampleApproved = True for sample G0224_040.0040.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:26 PM	Set SampleApproved = True for sample G0224_039.0039.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:32 PM	Set Comment = re-extract for sample G0224_041.0041.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:32 PM	Set SampleApproved = True for sample G0224_041.0041.D; previous value = False			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:53 PM	Set Comment = cannot re-extract due to limited sample volume for sample G0224_041.0041.D; previous value = re-extract			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:53 PM	Set SampleApproved = True for sample G0224_042.0042.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:58 PM	Set Comment = re-extract for sample G0224_042.0042.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:44:59 PM	Set SampleApproved = True for sample G0224_043.0043.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:45:03 PM	Set Comment = re-extract for sample G0224_043.0043.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:45:11 PM	Set SampleApproved = True for sample G0224_044.0044.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:45:12 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_044.0044.D; previous value = GT			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:45:16 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_044.0044.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:45:36 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_045.0045.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:45:38 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_045.0045.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:45:39 PM	Set SampleApproved = True for sample G0224_045.0045.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:47:42 PM	Set SampleApproved = True for sample G0224_056.0056.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:47:43 PM	Set SampleApproved = True for sample G0224_055.0055.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:47:44 PM	Set SampleApproved = True for sample G0224_054.0054.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:47:52 PM	Set Comment = re-extract for sample G0224_055.0055.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:47:56 PM	Set Comment = re-extract for sample G0224_056.0056.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:47:59 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_054.0054.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:48:00 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_054.0054.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:07 PM	Set SampleApproved = True for sample G0224_053.0053.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:09 PM	Set SampleType = DoubleBlank for sample G0224_052.0052.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:11 PM	Set SampleType = DoubleBlank for sample G0224_054.0054.D; previous value = Sample			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:48:13 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0224_052.0052.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:48:19 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_054.0054.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:48:21 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_052.0052.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:25 PM	Set SampleApproved = True for sample G0224_052.0052.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:26 PM	Set SampleApproved = True for sample G0224_051.0051.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:27 PM	Set SampleApproved = True for sample G0224_050.0050.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:27 PM	Set SampleApproved = True for sample G0224_049.0049.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:28 PM	Set SampleApproved = True for sample G0224_048.0048.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:29 PM	Set SampleApproved = True for sample G0224_047.0047.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 1:48:39 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0224_021.0021.D, from x, y = 3.928, 16219 to 4.055, 16323, result = 42363; previous integration is from x, y = 3.915, 15620 to 4.094, 15081 and previous response = 51299.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/25/2022 1:48:40 PM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0224_021.0021.D to y = 16219, new integration is from x, y = 3.928, 16219 to 4.055, 16219 and new response = 42761; previous integration is from x, y = 3.928, 16219 to 4.055, 16323 and previous response = 42363.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:48:41 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_021.0021.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 1:48:44 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0224_021.0021.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:45 PM	Set SampleApproved = True for sample G0224_021.0021.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:48:52 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_046.0046.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:48:54 PM	Set SampleApproved = True for sample G0224_046.0046.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:49:01 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_042.0042.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:49:03 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_041.0041.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 1:49:07 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0224_040.0040.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:49:27 PM	Set SampleApproved = True for sample G0224_032.0032.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:49:55 PM	Set SampleApproved = True for sample G0224_006.0006.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:27 PM	Set SampleType = CC for sample G0224_007.0007.D; previous value = Calibration			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:28 PM	Set SampleType = CC for sample G0224_008.0008.D; previous value = Calibration			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:31 PM	Set SampleType = CC for sample G0224_009.0009.D; previous value = Calibration			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:32 PM	Set SampleType = CC for sample G0224_010.0010.D; previous value = Calibration			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:34 PM	Set SampleType = CC for sample G0224_011.0011.D; previous value = Calibration			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:36 PM	Set SampleType = CC for sample G0224_012.0012.D; previous value = Calibration			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 2:30:38 PM	Set SampleType = CC for sample G0224_013.0013.D; previous value = Calibration			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 2:30:42 PM	Quantitate all compounds in all samples			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 2:31:38 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/25/2022 3:20:51 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022422\aiexport\QuantResults\G022422_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\srcox	3/9/2022 2:55:32 PM	Open batch D:\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	
GenerateReport	BL2000\srcox	3/9/2022 2:59:41 PM	Generates report - Method: \\MASSHUNTER\Org\reports\Gen_ResultsSummary.m, Output Path: D:\Org\Data\GECD.I\G022422\aiexport\QuantReports\G022422_8011_W_CLT			✓	
GenerateReport	BL2000\srcox	3/9/2022 3:02:22 PM	Generates report - Method: \\MASSHUNTER\Org\reports\init_cal_report.m, Output Path: D:\Org\Data\GECD.I\G022422\aiexport\QuantReports\G022422_8011_W_CLT-1			✓	
CmdOpenBatchTable	BL2000\srcox	3/9/2022 3:13:56 PM	Open batch D:\Org\Data\GECD.I\G022422\aiexport\G022422_8011_W_CLT.batch.bin			✓	
GenerateReport	BL2000\srcox	3/9/2022 3:14:49 PM	Generates report - Method: \\MASSHUNTER\Org\reports\Gen_Calibration.m, Output Path: D:\Org\Data\GECD.I\G022422\aiexport\QuantReports\G022422_8011_W_CLT-2			✓	
GenerateReport	BL2000\srcox	3/9/2022 3:18:07 PM	Generates report - Method: \\MASSHUNTER\Org\reports\Env_QuantResults_wGraphics+Chromatogram.m, Output Path: D:\Org\Data\GECD.I\G022422\aiexport\QuantReports\G022422_8011_W_CLT-3			✓	

PREP BATCH REPORT

Prep Code: **PRP-8011-W**
 Prep Batch **164037** Prep Temp **NA °C**

Technician: **Carry L Tran**
 Batch Units: **ML**

Prep Start Date: **2/25/2022 9:49:45 AM**
 Prep End Date: **2/25/2022 12:07:00 P**

Sample ID	Matrix	pH	Initial Samp Amt	Sol Added	Sol Recovered	Final Vol (mL)	Factor	Balance	Prep Start Date	Prep End Date
MB-164037		6	35	0	0	2.0	0.057		2/25/2022	2/25/2022
	CLT spiked and surrogated. ORR witnessed. NDM assisted.									
LCS-164037		6	35	0	0	2.0	0.057		2/25/2022	2/25/2022
	Samples went on solvent at 11:30am									
LCS1-164037		6	35	0	0	2.0	0.057	Bal #25	2/25/2022	2/25/2022
	5mL_19K50667 calibrated/passed on 02/25/2022 prior to the extraction.									
CK3-164037		6	35	0	0	2.0	0.057	Bal #25	2/25/2022	2/25/2022
	All samples poured to 35mL using a gravimetrically determined standard made by CLT on 02/25/2022									
CK5-164037		6	35	0	0	2.0	0.057	Bal #25	2/25/2022	2/25/2022
	Unlocked to add comments concerning re-extracts- CLT 2/25/22. Unlocked to add comments and final masses - CLT 2/25/22									
B22021627-001H	Ground Water	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 1/3. Combined vial and sample weight of 61.38g with cap on. Empty vial weight with cap on 25.75g=35.63g. Entire sample consumed in extraction.									
B22021627-001HMS	Ground Water	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 2/3. Combined vial and sample weight of 61.29g with cap on. Empty vial weight with cap on 25.52g=35.77g. Entire sample consumed in extraction.									
B22021627-001HMSD	Ground Water	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 3/3. Combined vial and sample weight of 61.69g with cap on. Empty vial weight with cap on 25.95g=36.01g. Entire sample consumed in extraction.									
B22021627-004A	Trip Blank	1	35	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 1/2. Combined vial and sample weight of 61.27g with cap on. Empty vial weight with cap on 25.82g=35.45g.									
B22021627-006H	Ground Water	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 1/3. Combined vial and sample weight of 61.02g with cap on. Empty vial weight with cap on 25.52g=35.50g. Dark brown sediment present in sample.									
B22021627-009A	Trip Blank	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 1/2. Combined vial and sample weight of 61.19g with cap on. Empty vial weight with cap on 25.63g=35.56g.									
B22021627-011H	Ground Water	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 1/3. Combined vial and sample weight of 61.18g with cap on. Empty vial weight with cap on 25.33g=35.85g.									
B22021627-014A	Trip Blank	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
	Vial 1/2. Combined vial and sample weight of 61.21 g with cap on. Empty vial weight with cap on 25.52g=35.69g.									

Number	Reagent Name	Exp Date
11	Carbon Filter Water	1/1/2023
14206	pH-indicator Strips 0-14 HC160347	8/26/2026
14543	Hexane EB754	6/4/2023
14729	Laboratory Fortified Blank Sample Concentrate	2/6/2023
14776	4ML, Amber Vial, 20220118	1/18/2023
14853	40 mL Clear VOA Lot 00081943	2/8/2023

Spk ID	Spike Name	SampType	AmtAdd	Exp Date
NaCl 01/22/22(13	Baked Sodium Chloride	ALL	7g	6/15/2026
PH021622504Su	504.1 Surrogate (0.1ug/mL)MeOH	ALL Except CKs	35uL	3/20/2023
PH011122504C2	504.1 Cal Stock 2(0.07ug/mL) MeOH	CK3	50uL	2/12/2023
PH011122504C3	504.1 Cal Stock 3(0.7ug/mL) MeOH	CK5	20uL	2/12/2023

PREP BATCH REPORT

Prep Code: **PRP-8011-W**
 Prep Batch **164037** Prep Temp **NA °C**

Technician: **Carry L Tran**
 Batch Units: **ML**

Prep Start Date: **2/25/2022 9:49:45 AM**
 Prep End Date: **2/25/2022 12:07:00 P**

Sample ID	Matrix	pH	Initial Samp Amt	Sol Added	Sol Recovered	Final Vol (mL)	Factor	Balance	Prep Start Date	Prep End Date
B22021435-032H	Ground Water	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
This is a re-extract from batch #163984 due to the hexane dispensette not accurately delivering the required solvent due to a low amount of hexane in the bottle. Vial 2/3. Combined vial and sample weight of 61.41g with cap on. Empty vial weight with cap on 25.65g=35.76g.										
B22021435-035A	Trip Blank	1	35	0	0	2.0	0.057	Bal #25	2/25/2022	2/25/2022
This is a re-extract from batch #163984 due to the hexane dispensette not accurately delivering the required solvent due to a low amount of hexane in the bottle. Vial 2/2. Combined vial and sample weight of 61.13g with cap on. Empty vial weight with cap on 25.79g=35.34g. Entire sample consumed in extraction.										
B22021528-002C	Groundwater	1	36	0	0	2.0	0.056	Bal #25	2/25/2022	2/25/2022
This is a re-extract from batch #163984 due to the hexane dispensette not accurately delivering the required solvent due to a low amount of hexane in the bottle. Vial 2/2. Combined vial and sample weight of 61.28g with cap on. Empty vial weight with cap on 25.72g=35.56g. Entire sample consumed in extraction. Sediment present in sample.										
B22021528-003C	Groundwater	1	35	0	0	2.0	0.057	Bal #25	2/25/2022	2/25/2022
This is a re-extract from batch #163984 due to the hexane dispensette not accurately delivering the required solvent due to a low amount of hexane in the bottle. Vial 3/3. Combined vial and sample weight of 61.20g with cap on. Empty vial weight with cap on 25.94g=35.26g. Sediment is present in sample. Vial 2/3 is still in storage cooler.										

Number	Reagent Name	Exp Date
11	Carbon Filter Water	1/1/2023
14206	pH-indicator Strips 0-14 HC160347	8/26/2026
14543	Hexane EB754	6/4/2023
14729	Laboratory Fortified Blank Sample Concentrate	2/6/2023
14776	4ML, Amber Vial, 20220118	1/18/2023
14853	40 mL Clear VOA Lot 00081943	2/8/2023

Spk ID	Spike Name	SampType	AmtAdd	Exp Date
NaCl 01/22/22(13	Baked Sodium Chloride	ALL	7g	6/15/2026
PH021622504Su	504.1 Surrogate (0.1ug/mL) MeOH	ALL Except CKs	35uL	3/20/2023
PH011122504C2	504.1 Cal Stock 2(0.07ug/mL) MeOH	CK3	50uL	2/12/2023
PH011122504C3	504.1 Cal Stock 3(0.7ug/mL) MeOH	CK5	20uL	2/12/2023

Energy Laboratories Inc

ANALYTICAL RUN Summary

09-Mar-22

Run ID GECD.I_220225A

Run Start Date: 2/25/2022
Analyst: Carry L Tran
Ical:
Column ID: RTX-CLP_0.53
Comments:

Std ID	Std Name	Std Amount	Std Units	Samp Amount	Samp Units	SampType	Expiration Date
SeePrepRecord	Standards Tracked in Prep Batch						2/4/2050

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist				
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15056846	CK3-164037	PST-8011-W	CCV3	ECD.I\G022522\ai	2/25/2022 12:56:	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.11139	0.11111153		0.1	0	0	0.0025835	0.01	0	111%	80	120	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10089	0.10063778		0.1	0	0	0.0056259	0.02	0	101%	80	120	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist				
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15056847	MB-164037	PST-8011-W	MBLK	ECD.I\G022522\	2/25/2022 1:16:5	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.005	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10034	0.10008915		0.1	0	0	0.0056259	0.02	0	100%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist				
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15056848	LCS-164037	PST-8011-W	LCS-DOD	ECD.I\G022522\	2/25/2022 1:36:4	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.24587	0.24525533		0.25	0	0	0.0025835	0.01	0	98%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09866	0.09841335		0.1	0	0	0.0056259	0.02	0	98%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056849	LCS1-164037	PST-8011-W	LCS1	JECD.IG022522\	2/25/2022 1:56:4	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.1024	0.102144		0.1	0	0	0.0025835	0.01	0	102%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09806	0.09781485		0.1	0	0	0.0056259	0.02	0	98%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056850	B22021627-004	PST-8011-W	SAMP	JECD.IG022522\	2/25/2022 2:36:1	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.09656	0.0946288		0.099	0	0	0.0055272	0.02	0	96%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056851	B22021627-006	PST-8011-W	SAMP	JECD.IG022522\	2/25/2022 2:55:5	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10647	0.1043406		0.099	0	0	0.0055272	0.02	0	105%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056852	B22021627-009	PST-8011-W	SAMP	JECD.IG022522\	2/25/2022 3:15:3	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10571	0.1035958		0.098	0	0	0.0055272	0.02	0	106%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056853	B22021627-011	PST-8011-W	SAMP	JECD.IG022522\	2/25/2022 3:35:2	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10112	0.0990976		0.098	0	0	0.0055272	0.02	0	101%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056854	B22021627-014	PST-8011-W	SAMP	JECD.I\G022522\	2/25/2022 3:55:1	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10216	0.1001168		0.098	0	0	0.0055272	0.02	0	102%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056855	B22021435-032	PST-8011-W	SAMP	JECD.I\G022522\	2/25/2022 4:15:0	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10745	0.105301		0.098	0	0	0.0055272	0.02	0	107%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056856	B22021435-035	PST-8011-W	SAMP	JECD.I\G022522\	2/25/2022 4:34:4	1	163984	2/24/2022 8:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10088	0.0988624		0.099	0	0	0.0055272	0.02	0	100%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056857	B22021627-001	PST-8011-W	SAMP	JECD.I\G022522\	2/25/2022 4:54:2	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	U
1,1,1,2-Tetrachloroethane	S	ug/L	0.10764	0.1054872		0.098	0	0	0.0055272	0.02	0	108%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056858	B22021627-001	PST-8011-W	MS-DOD	JECD.I\G022522\	2/25/2022 5:14:0	1	164037	2/25/2022 9:	2E+07	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.25382	0.2487436		0.245	0	0	0.0025382	0.01	0	102%	60	140	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10177	0.0997346		0.098	0	0	0.0055272	0.02	0	102%	70	130	0%	

Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056859	B22021627-001	PST-8011-W	MSD-DOD	JECD.IG022522\	2/25/2022 5:34:0	1	164037	2/25/2022 9:	2E+07	2E+07						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.25931	0.2541238		0.2425	0	0.2487436	0.0025382	0.01	0	105%	60	140	2%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10629	0.1041642		0.097	0	0	0.0055272	0.02	0	107%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056860	CK5-164037	PST-8011-W	CCV4	JECD.IG022522\	2/25/2022 6:13:4	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.4496	0.448476		0.4	0	0	0.0025835	0.01	0	112%	80	120	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.45915	0.45800213		0.4	0	0	0.0056259	0.02	0	115%	80	120	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056861	B22021528-002	PST-8011-W	SAMP	JECD.IG022522\	2/25/2022 6:53:0	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025382	0.01	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.09888	0.0969024		0.099	0	0	0.0055272	0.02	0	98%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056862	B22021528-003	PST-8011-W	SAMP	JECD.IG022522\	2/25/2022 7:12:5	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0	0		0	0	0	0.0025835	0.01	0	0%	0	0	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.10501	0.10474748		0.099	0	0	0.0056259	0.02	0	106%	70	130	0%	
Seq No	Lab ID	Test Code	Sample Typ	File ID	Analysis Date	DF	Batch ID	Prep Date	SPKref	RPDref	pmoist					
15056863	CK3-164037	PST-8011-W	CCV1	JECD.IG022522\	2/25/2022 7:52:2	1	164037	2/25/2022 9:	0	0						
Analyte	T	Units	RAW	Final	Text	Spike	SPKref	RPDref	MDL	PQL	UQL	%REC	LOW	HIGH	%RPD	Q
1,2-Dibromoethane	A	ug/L	0.12359	0.12328103		0.1	0	0	0.0025835	0.01	0	123%	70	130	0%	
1,1,1,2-Tetrachloroethane	S	ug/L	0.1062	0.1059345		0.1	0	0	0.0056259	0.02	0	106%	70	130	0%	

Write Sequence

Insert Entries(Have the first cell for e

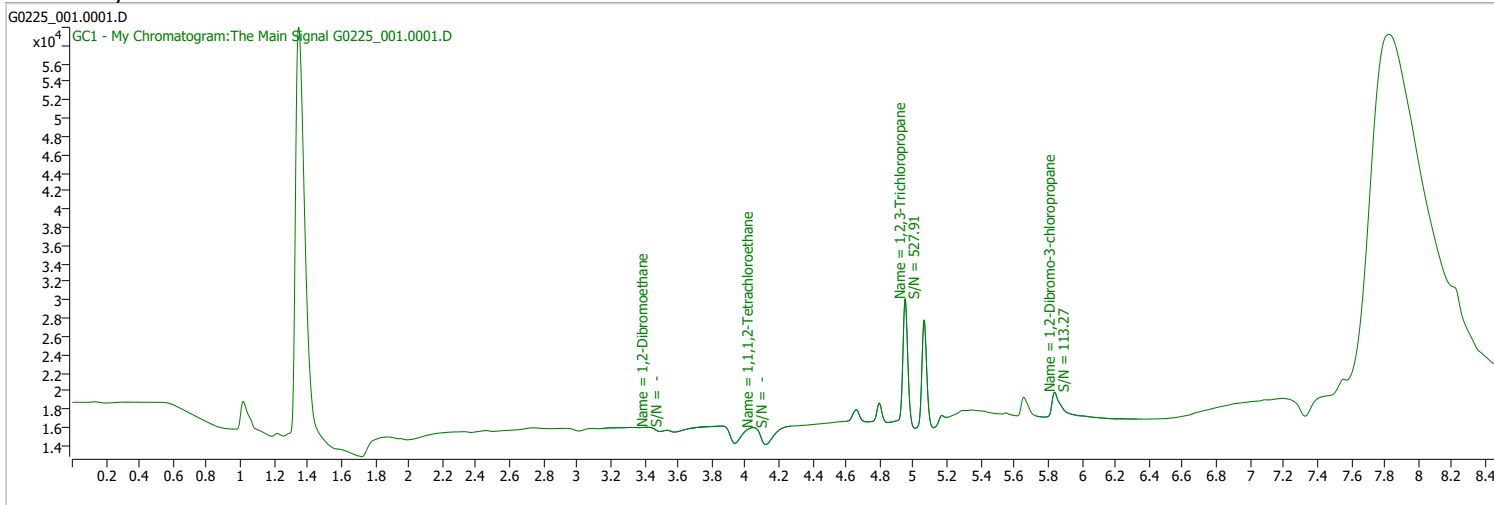
Data File**Sample Name**

Data File	Sample Name
G:\org\GECD.i\G022522.b\G0225_001	Hexane ;
G:\org\GECD.i\G022522.b\G0225_002	CK3-164037 ;
G:\org\GECD.i\G022522.b\G0225_003	MB-164037 ;
G:\org\GECD.i\G022522.b\G0225_004	LCS-164037 ;
G:\org\GECD.i\G022522.b\G0225_005	LCS1-164037 ;
G:\org\GECD.i\G022522.b\G0225_006	Hexane;;
G:\org\GECD.i\G022522.b\G0225_007	B22021627-004A ;\$PST-8011-W,
G:\org\GECD.i\G022522.b\G0225_008	B22021627-006H ;\$PST-8011-W,
G:\org\GECD.i\G022522.b\G0225_009	B22021627-009A ;\$PST-8011-W,
G:\org\GECD.i\G022522.b\G0225_010	B22021627-011H ;\$PST-8011-W,
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G:\org\GECD.i\G022522.b\G0225_012	B22021435-032H ;\$PST-8011-W,
G:\org\GECD.i\G022522.b\G0225_013	B22021435-035A ;\$PST-8011-W,
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G:\org\GECD.i\G022522.b\G0225_017	Hexane;;
G:\org\GECD.i\G022522.b\G0225_018	CK5-164037 ;
G:\org\GECD.i\G022522.b\G0225_019	Hexane;;
G:\org\GECD.i\G022522.b\G0225_020	B22021528-002C ;\$PST-8011-W,
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Quantitation Results Report (QT Reviewed)

Data File	G0225_001.0001.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 12:37:08 PM
Sample Name	Hexane	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

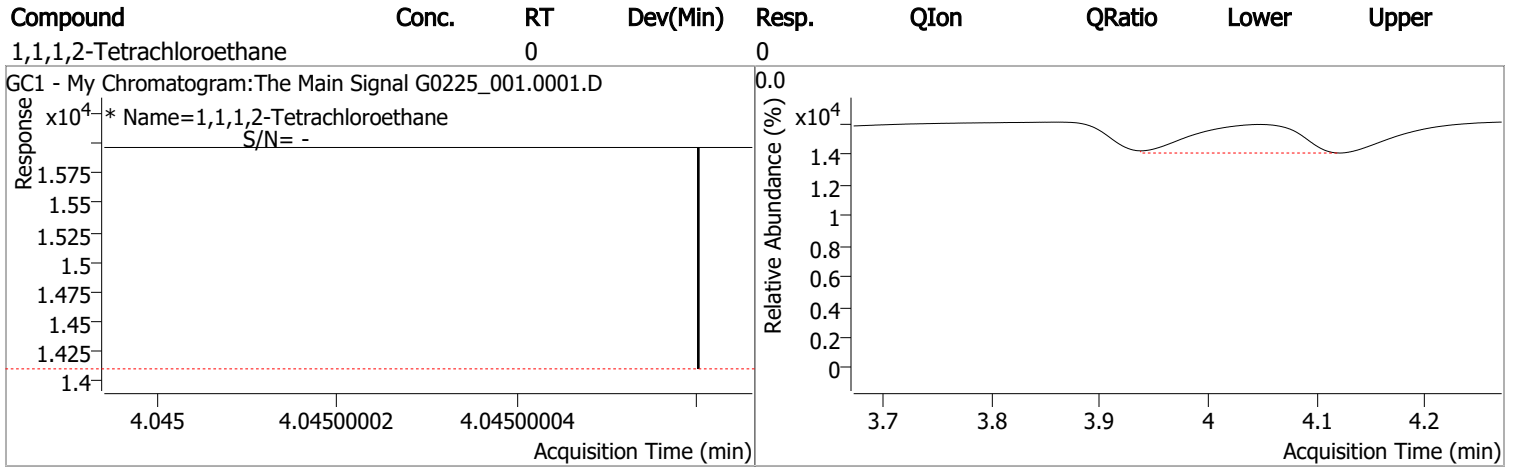
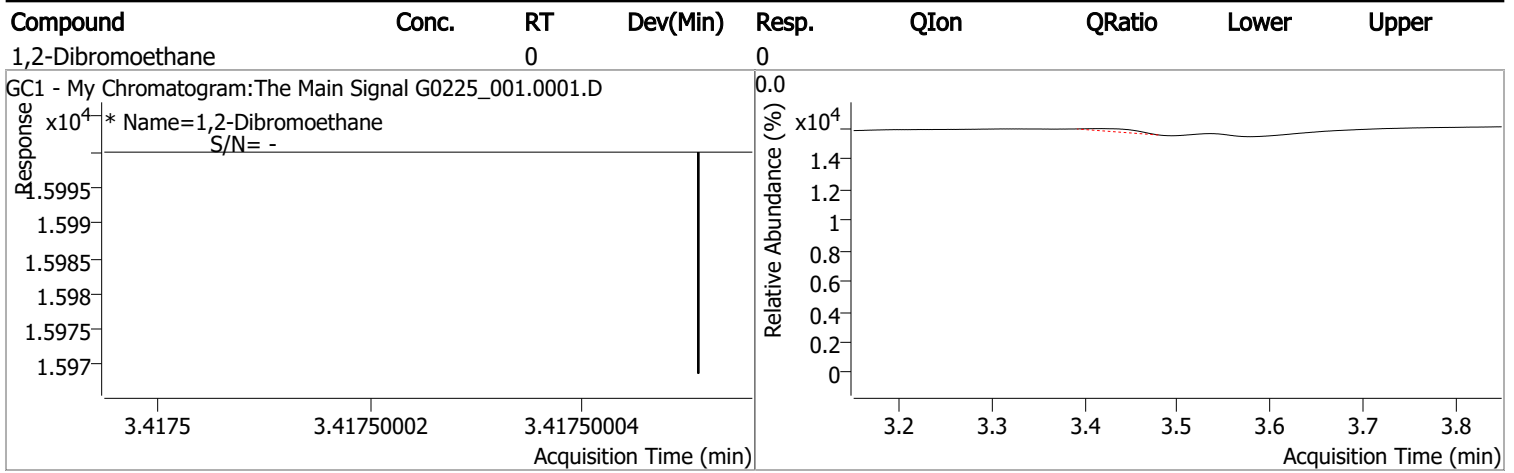
S 1,1,1,2-Tetrachloroethane	4.045	0.0	0		µg/L	md	0.073
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = NA%			

Target Compounds

M 1,2-Dibromoethane	3.418	0.0	0		µg/L	md	QValue 1
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(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

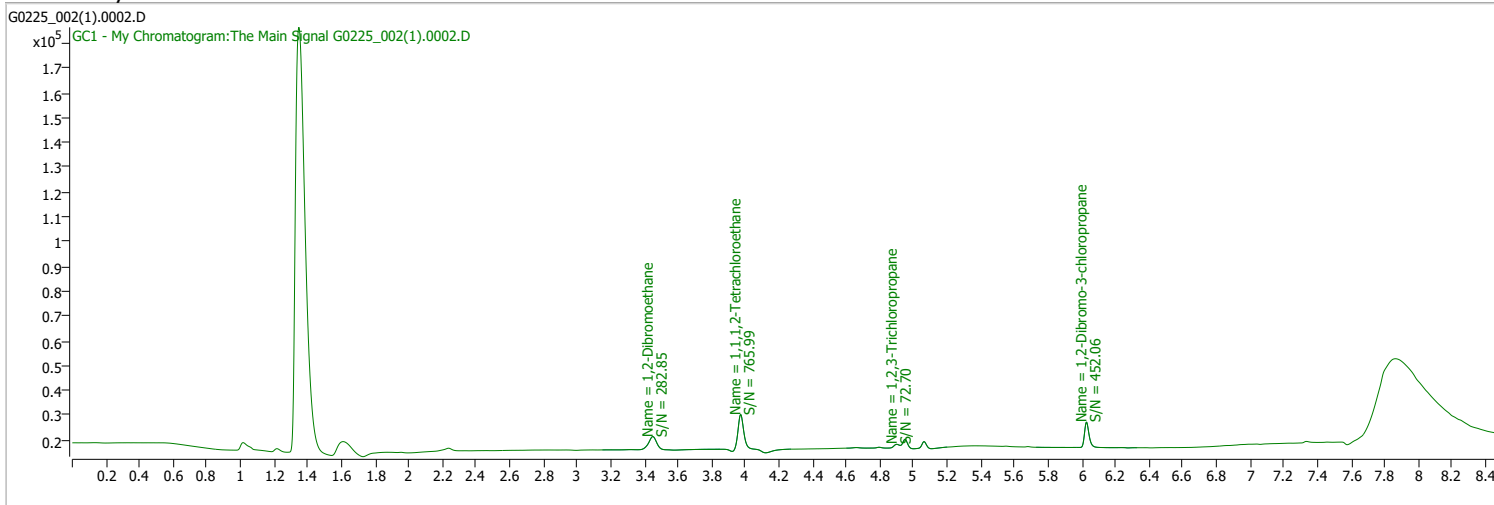
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_002(1).0002.D	Operator	
Acq. Method	testAcqFilePath	Acq. Date-Time	2/25/2022 12:56:57 PM
Sample Name	CK3-164037	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

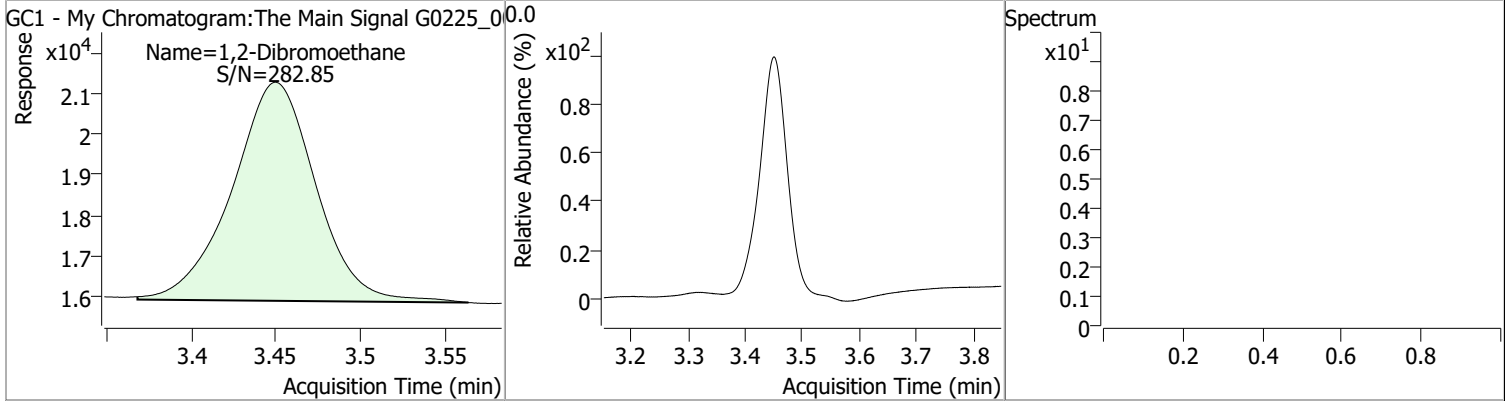


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.972	0.0	33580	0.1009	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 100.89%		
Target Compounds						
M 1,2-Dibromoethane	3.450	0.0	18599	0.1114	µg/L	QValue 100

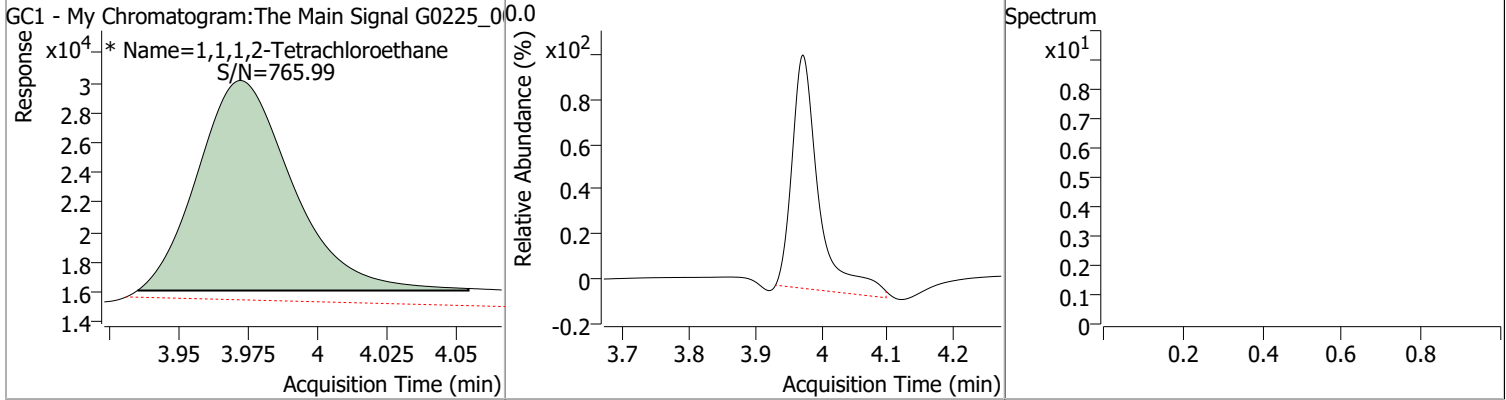
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1114	3.45	0.00	18599				



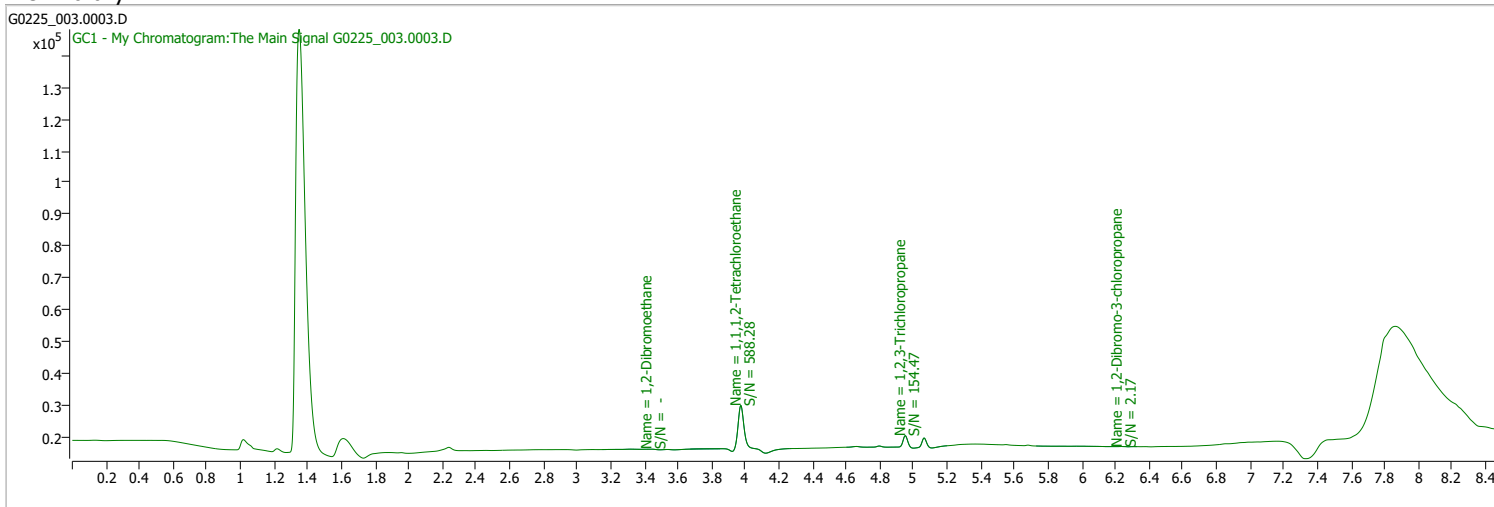
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1009	3.97	0.00	33580 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0225_003.0003.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:16:58 PM
Sample Name	MB-164037	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

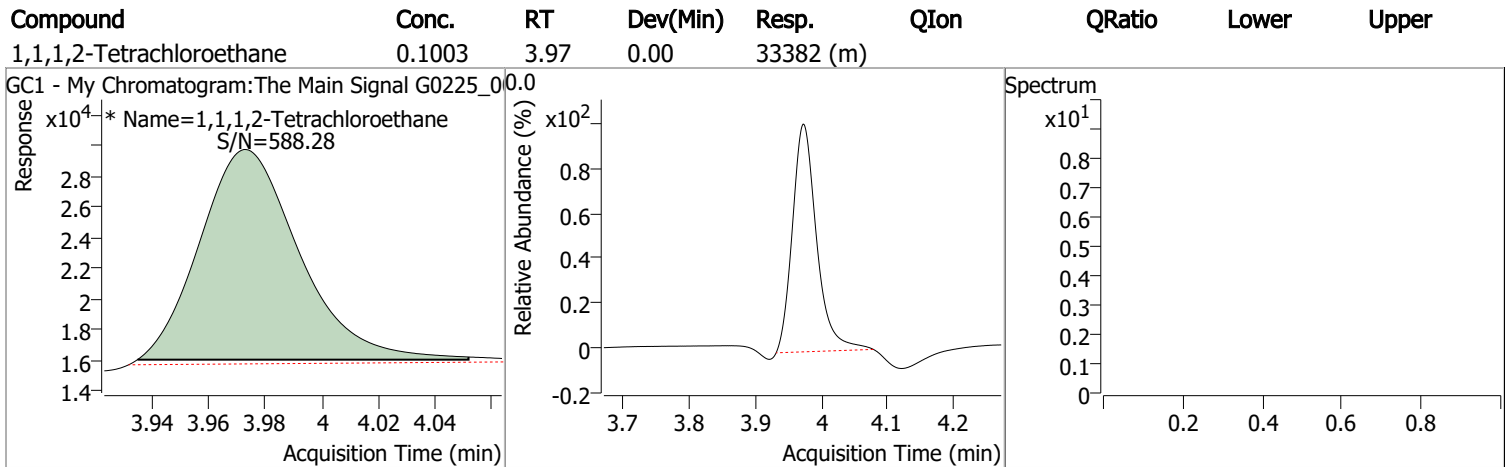
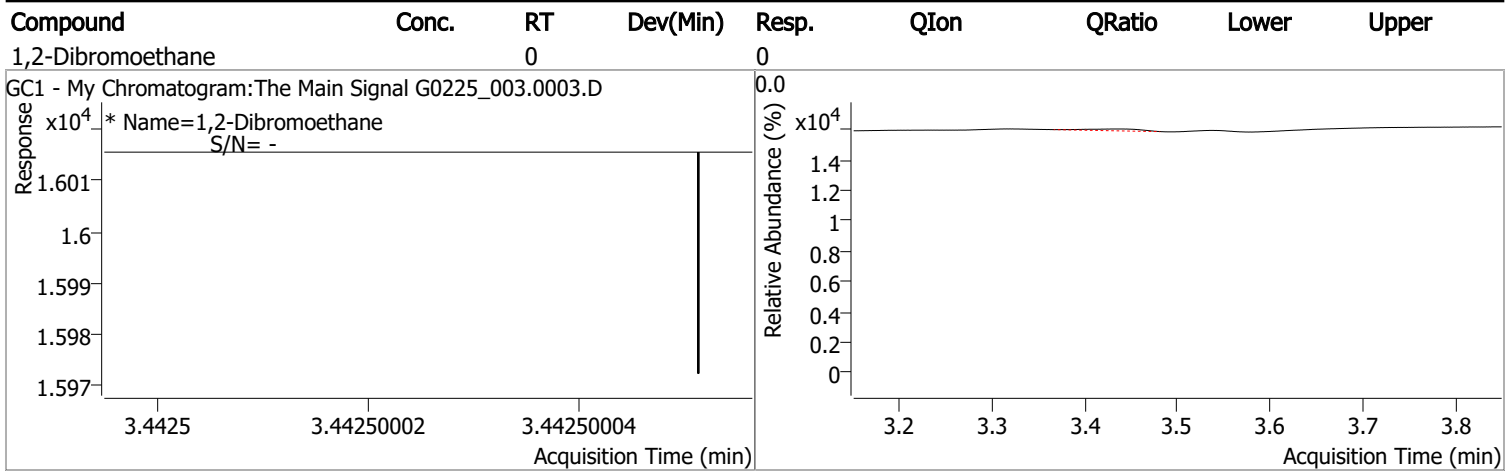
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.973	0.0	33382	0.1003	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 100.34%		
Target Compounds						
M 1,2-Dibromoethane	3.443	0.0	0		µg/L	md
						QValue
						1

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

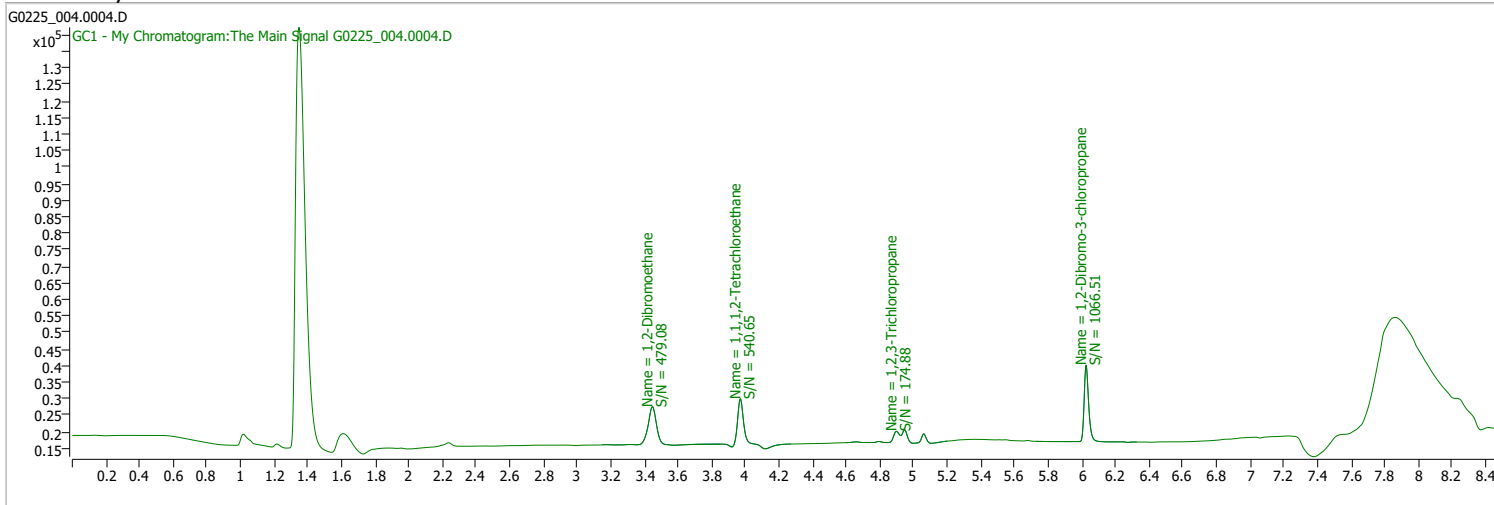
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_004.0004.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:36:47 PM
Sample Name	LCS-164037	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

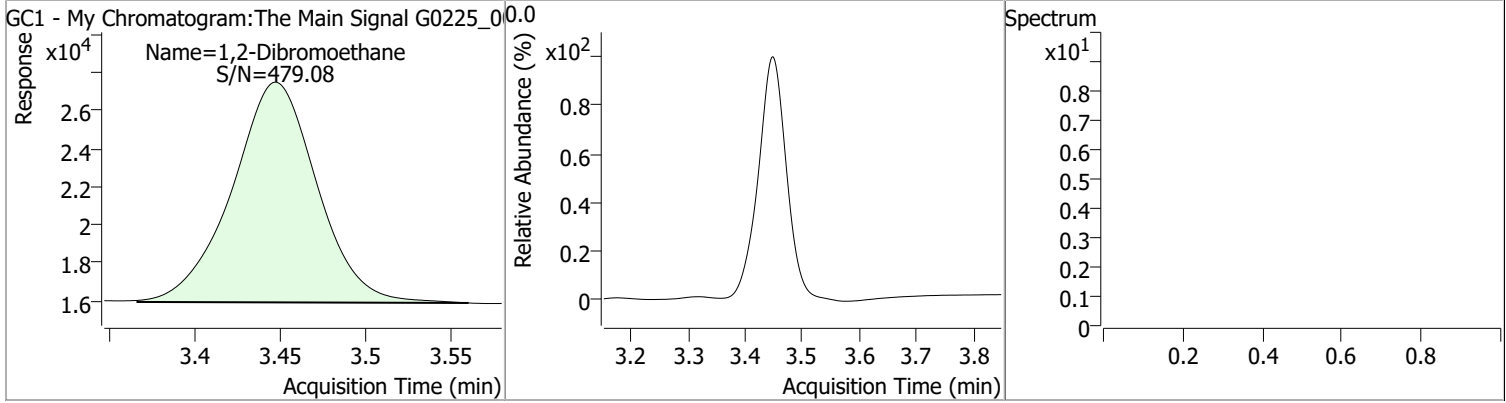


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	32779	0.0987	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 98.66%		
Target Compounds						
M 1,2-Dibromoethane	3.448	0.0	40154	0.2459	µg/L	QValue 100

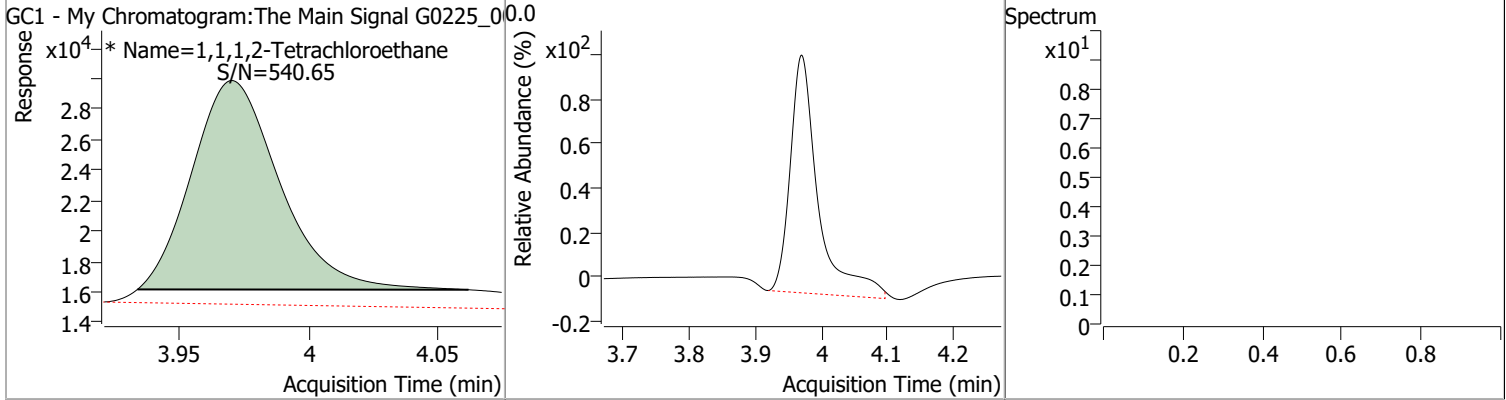
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2459	3.45	0.00	40154				



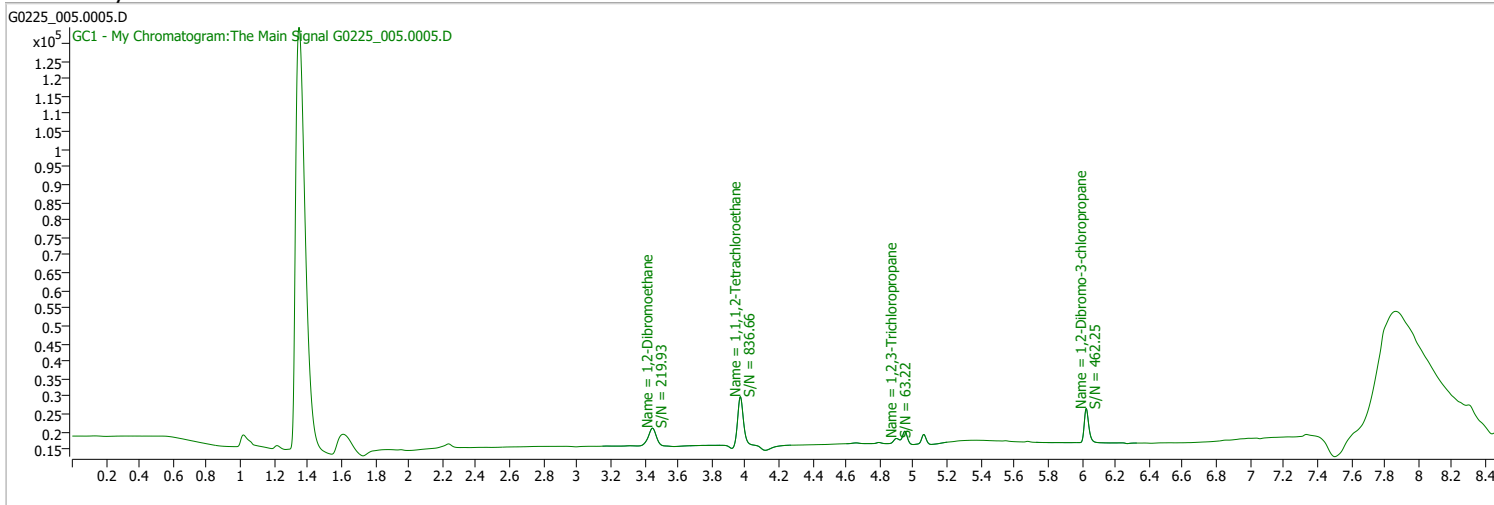
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0987	3.97	0.00	32779 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0225_005.0005.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 1:56:44 PM
Sample Name	LCS1-164037	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

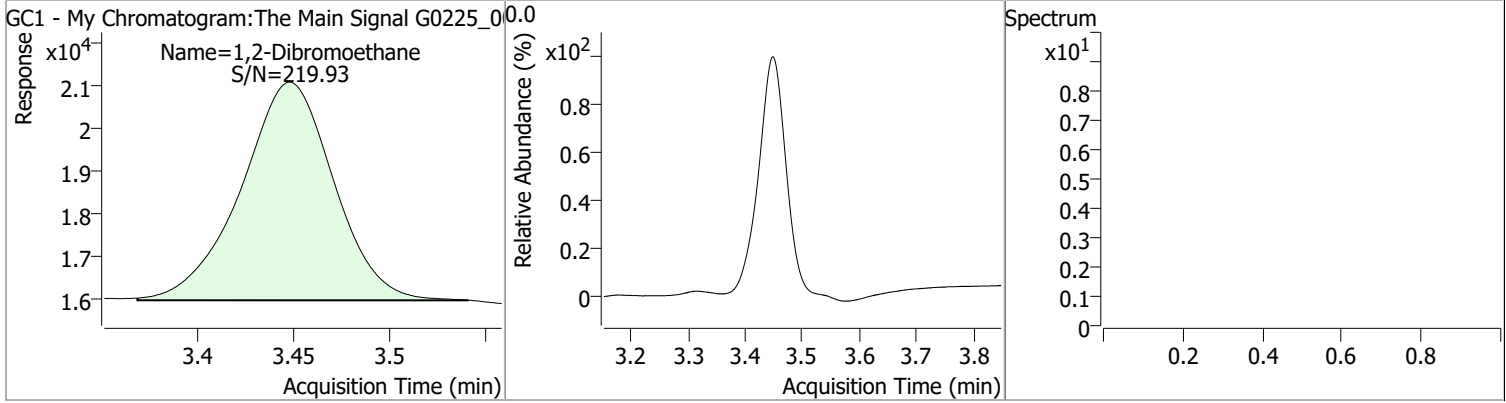


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	32562	0.0981	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 98.06%		
Target Compounds						
M 1,2-Dibromoethane	3.448	0.0	17122	0.1024	µg/L	QValue 100

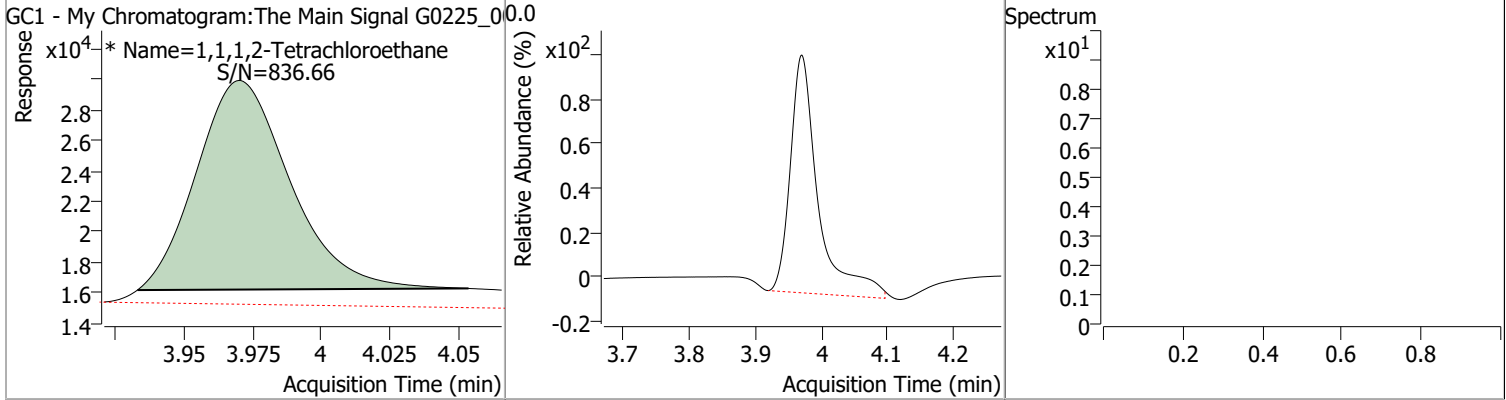
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1024	3.45	0.00	17122				



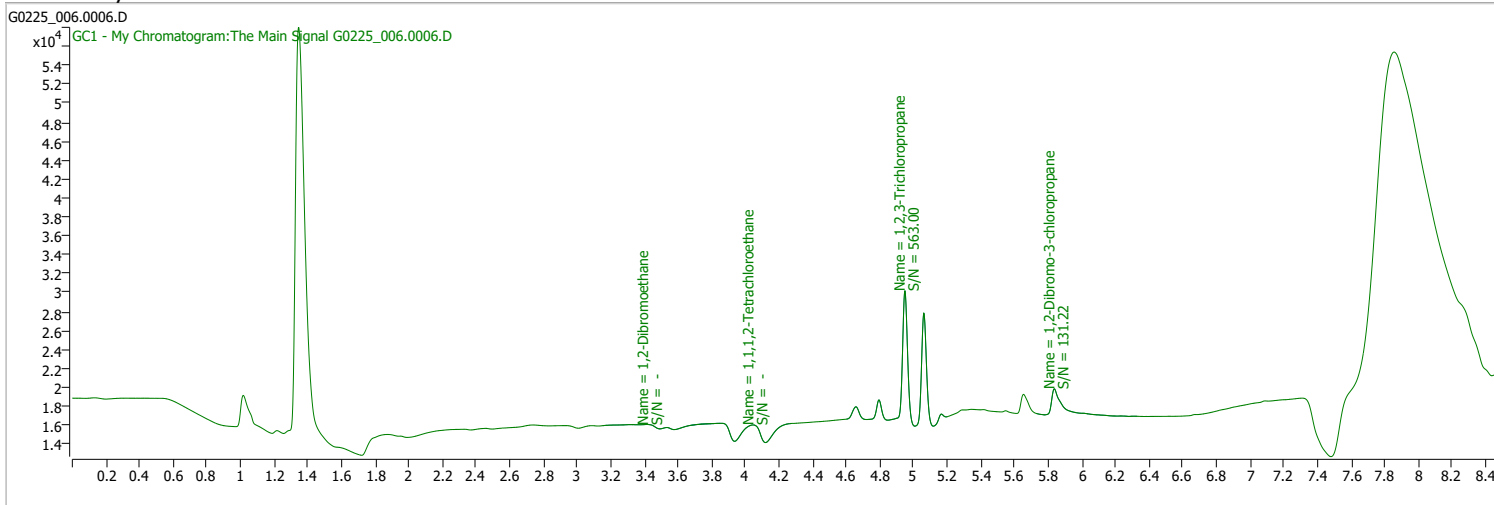
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.0981	3.97	0.00	32562 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0225_006.0006.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 2:16:20 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

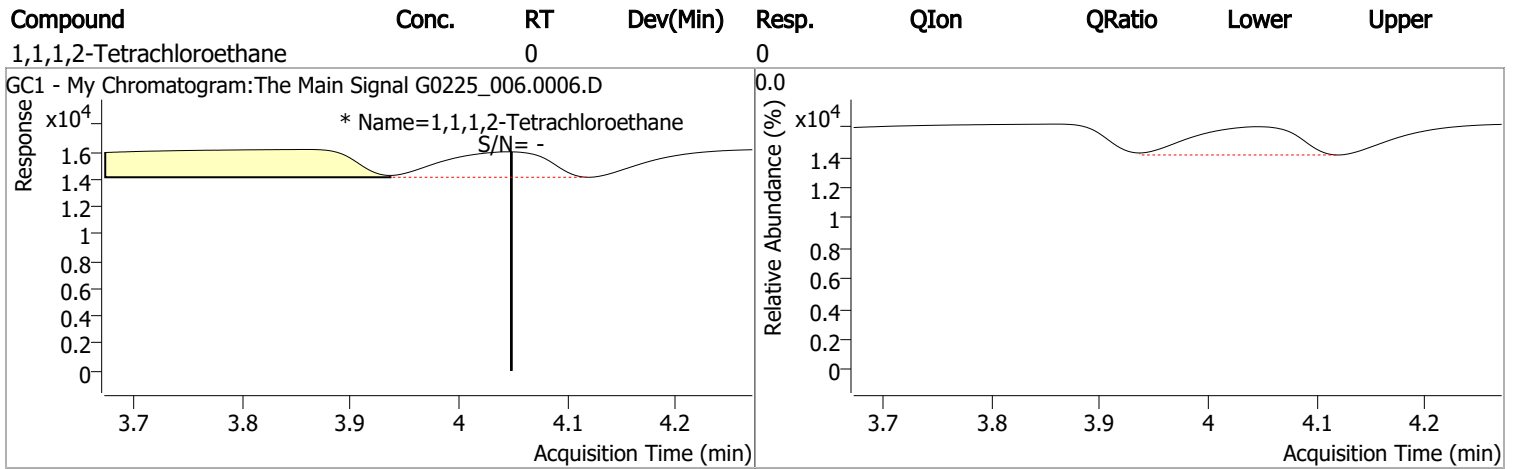
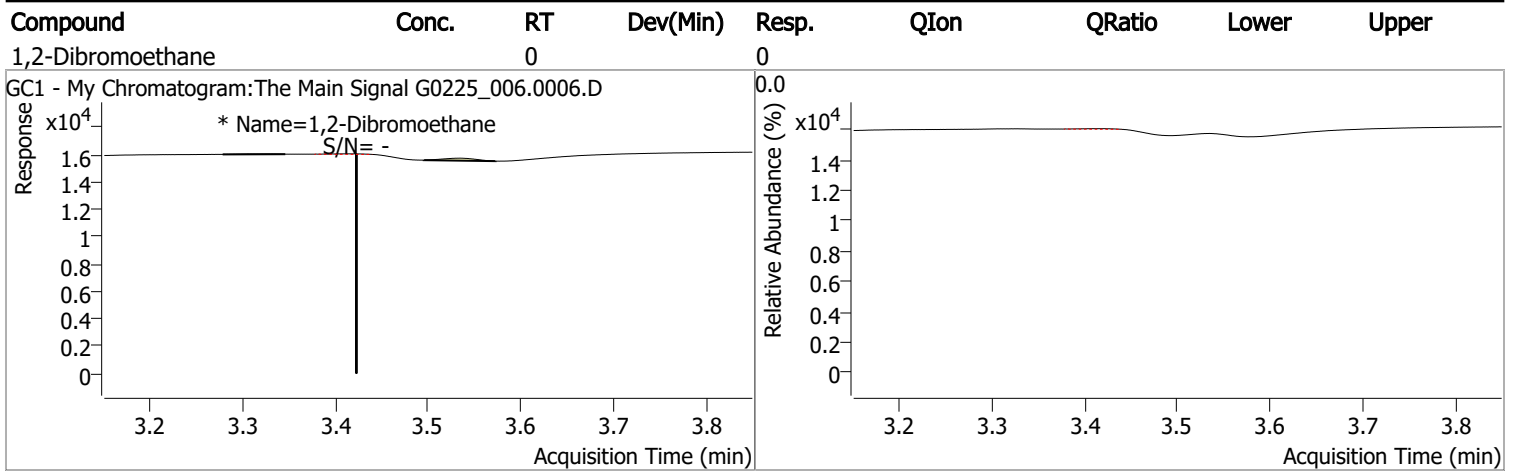
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.048	0.0	0		µg/L	md
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.423	0.0	0		µg/L	md
						QValue
						1

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

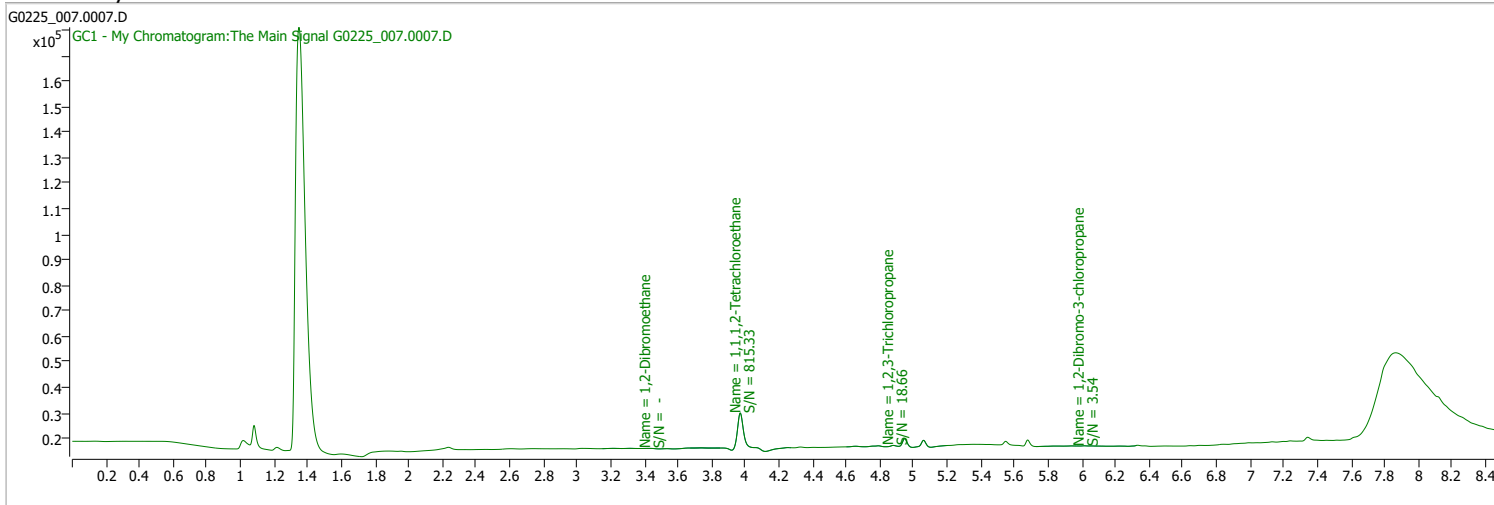
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_007.0007.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 2:36:12 PM
Sample Name	B22021627-004A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

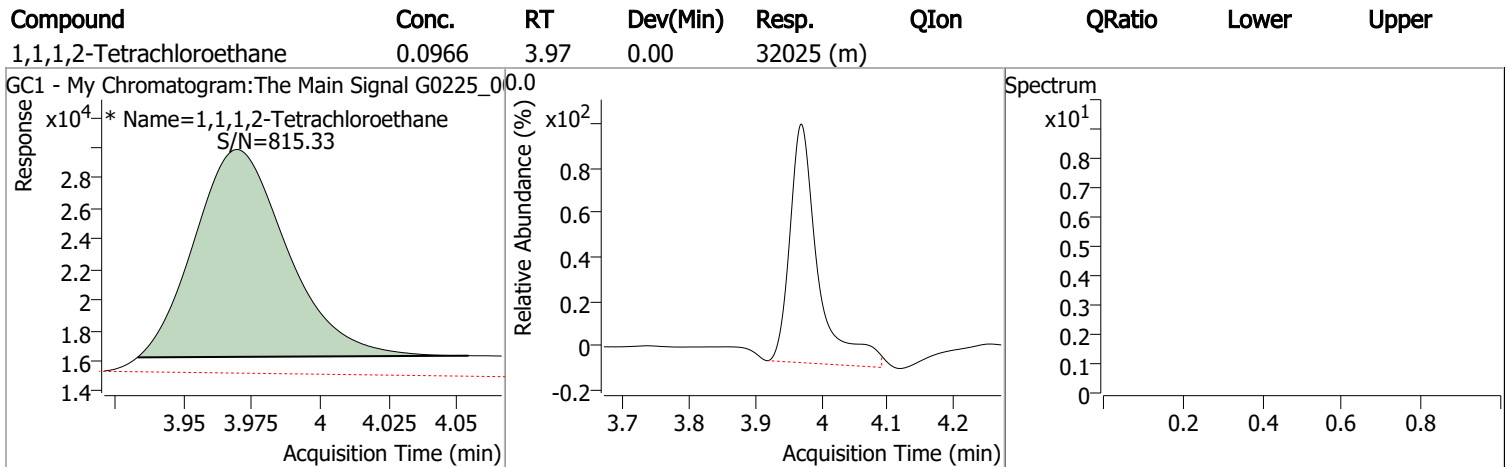
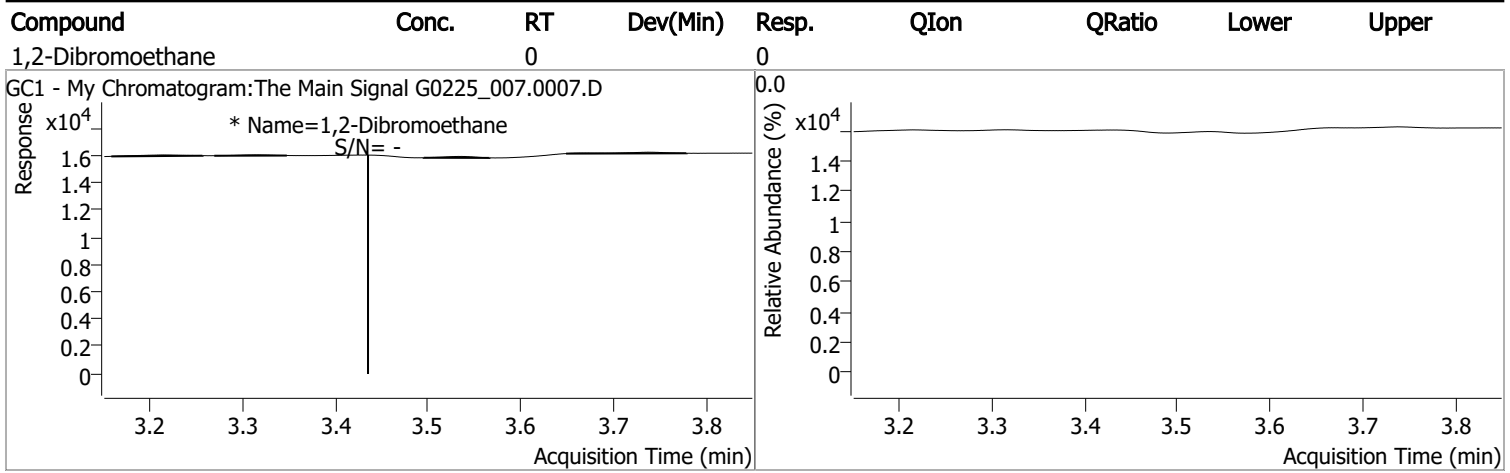
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	32025	0.0966	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 96.56%		
Target Compounds						
M 1,2-Dibromoethane	3.435	0.0	0		µg/L	md
						QValue 1

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

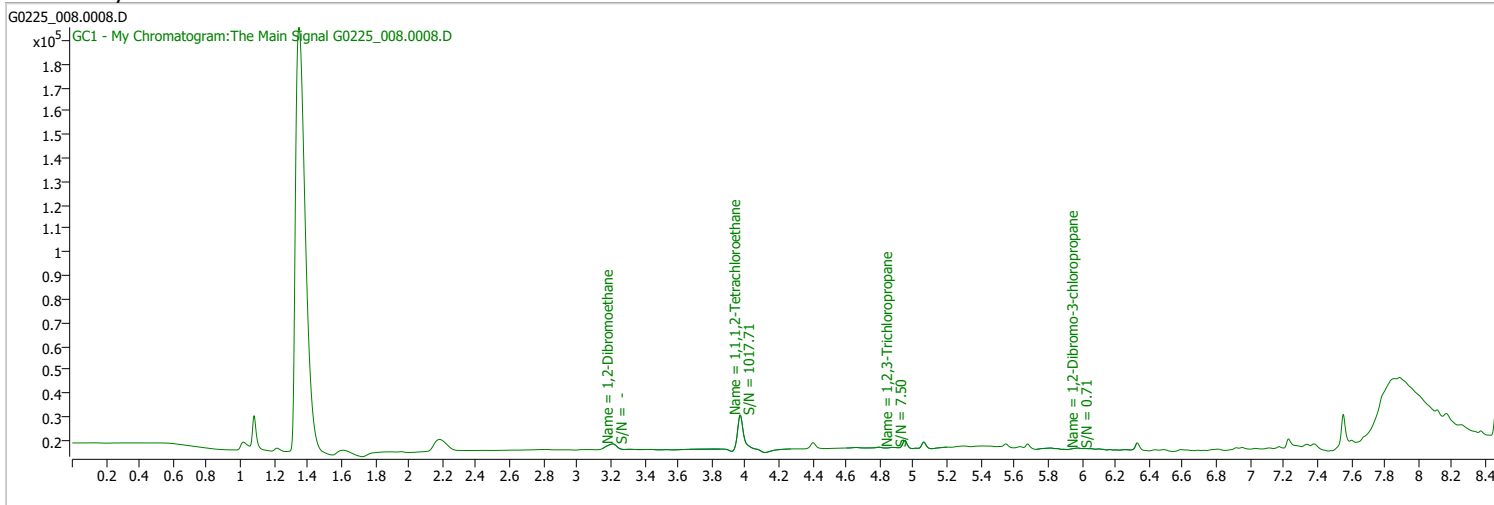
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_008.0008.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 2:55:59 PM
Sample Name	B22021627-006H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

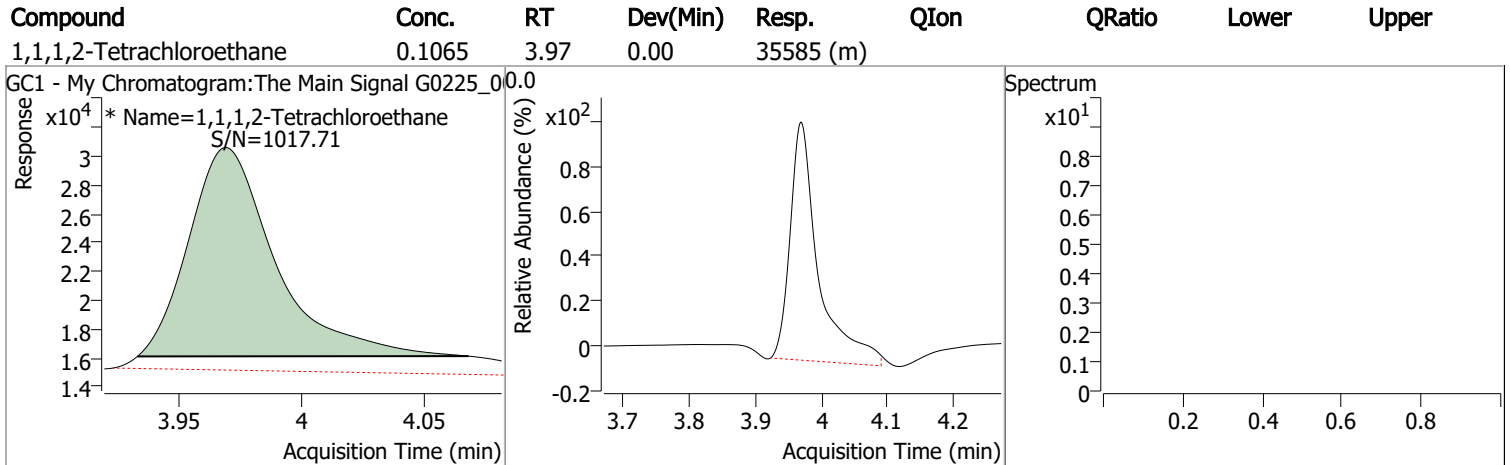
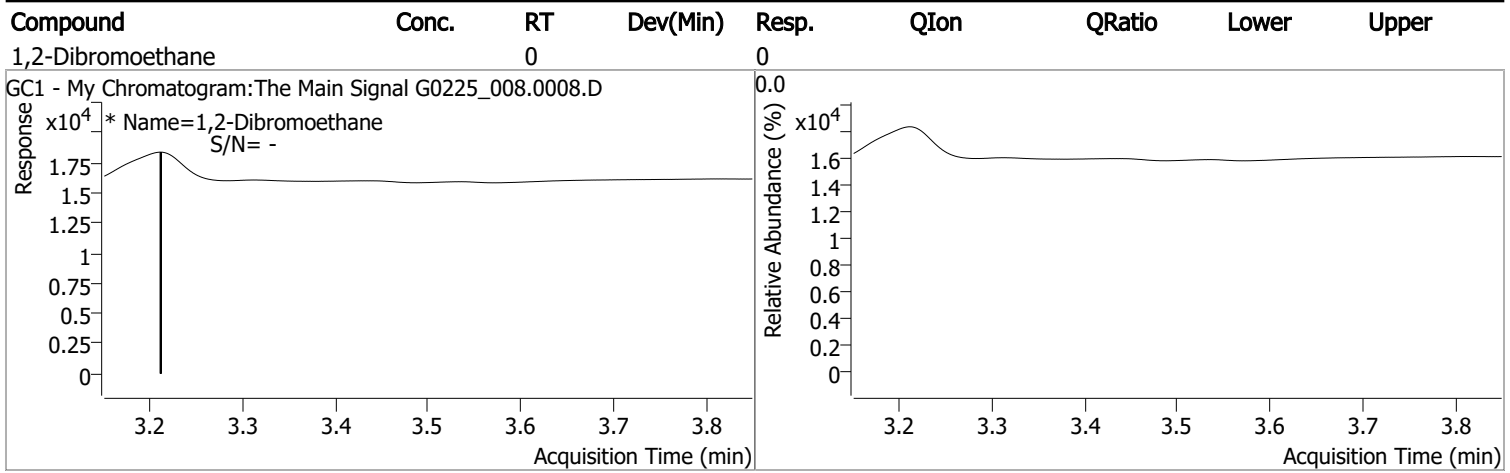
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.969	0.0	35585	0.1065	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 106.47%		
Target Compounds						
M 1,2-Dibromoethane	3.212	0.0	0		µg/L	md
						QValue 1

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

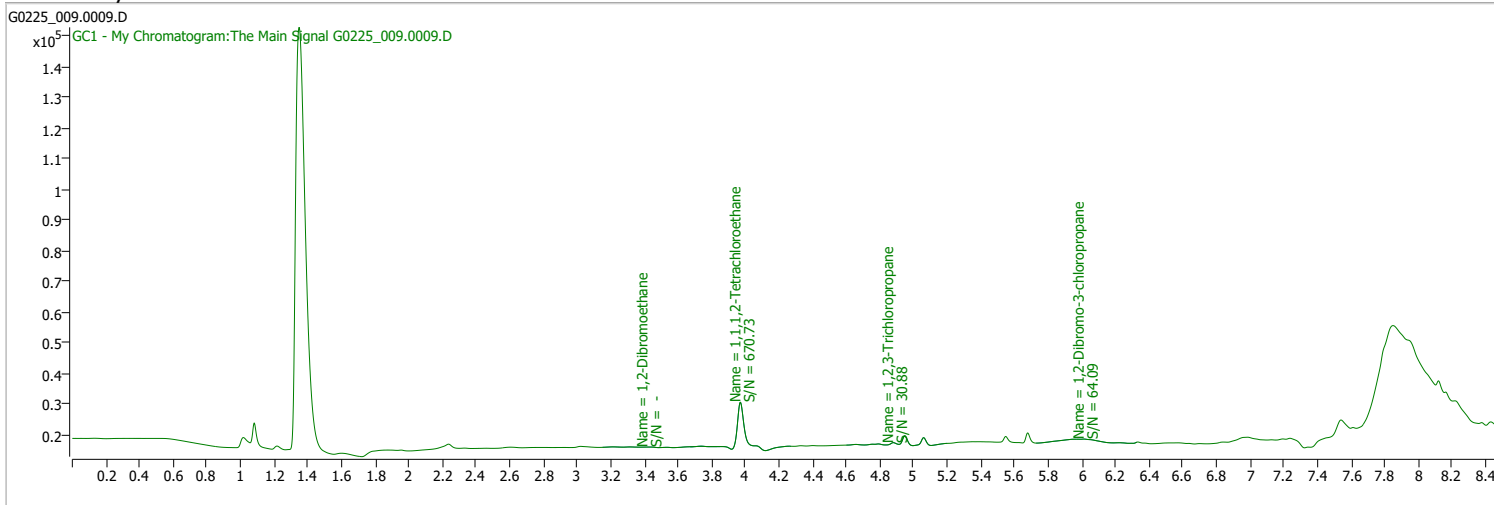
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_009.0009.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 3:15:39 PM
Sample Name	B22021627-009A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

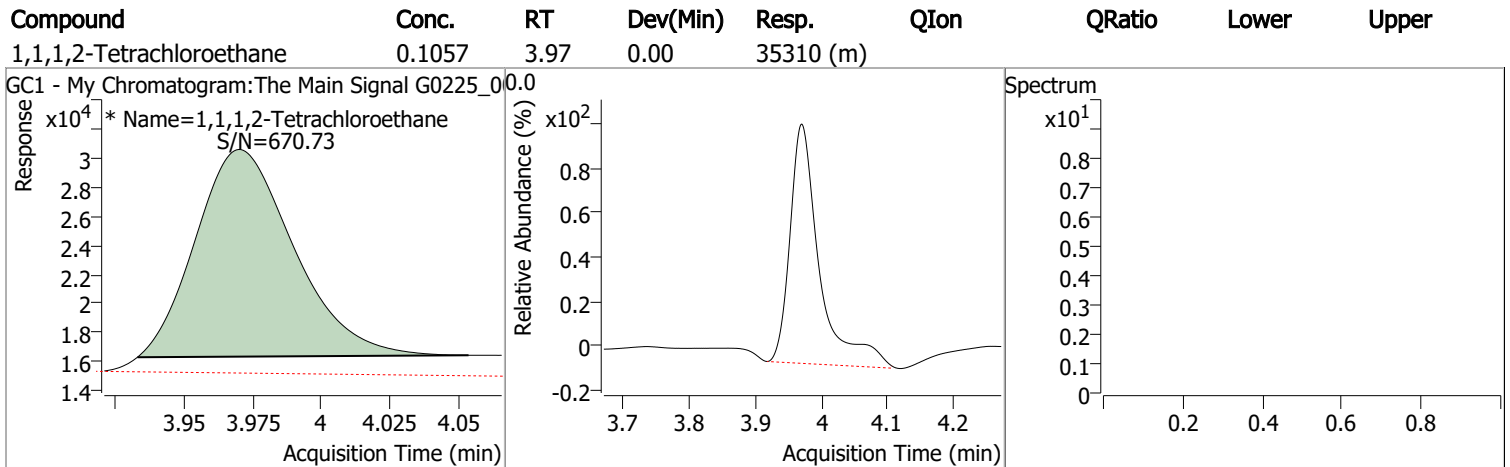
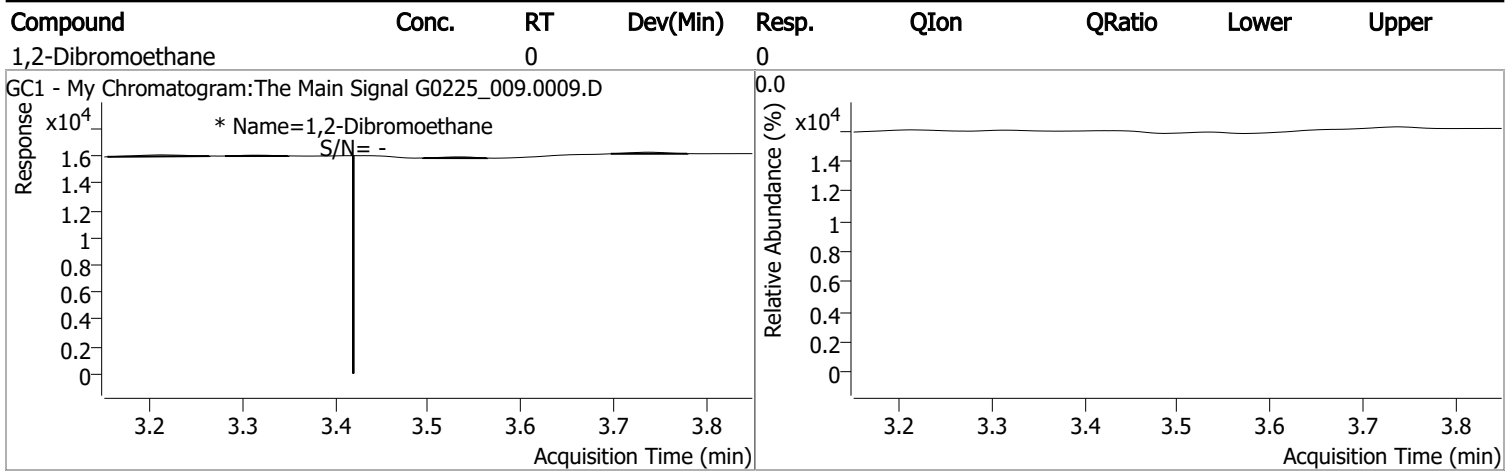
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.970	0.0	35310	0.1057	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%		Recovery = 105.71%			
Target Compounds						
M 1,2-Dibromoethane	3.419	0.0	0		µg/L	md
						QValue
						1

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

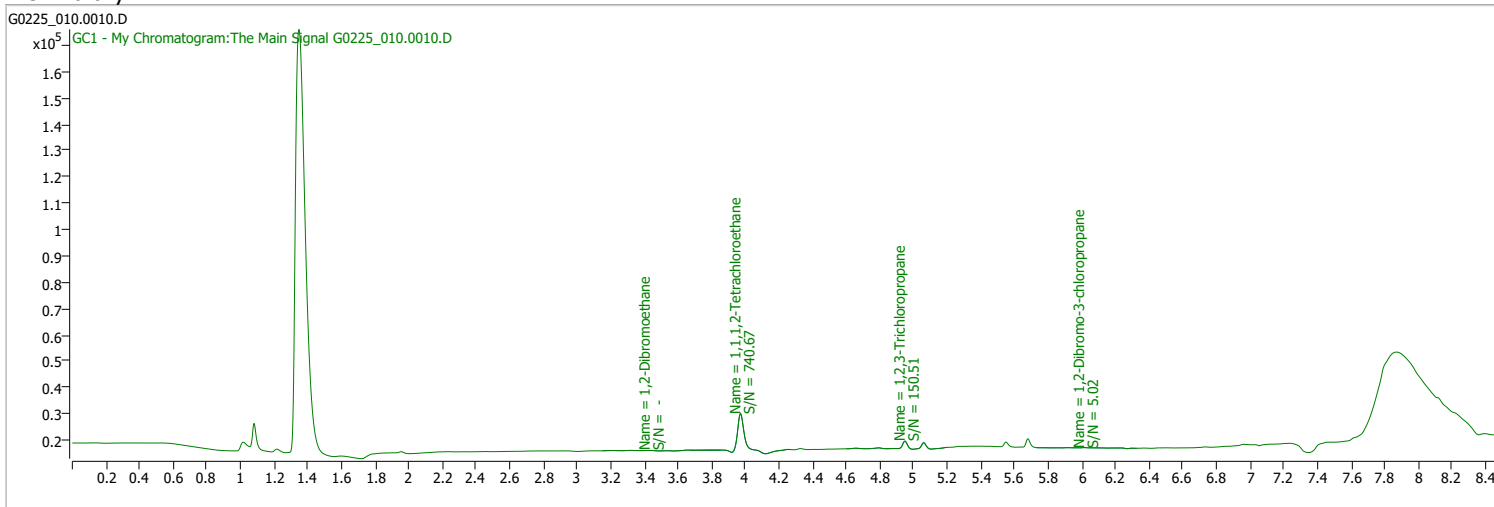
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_010.0010.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 3:35:21 PM
Sample Name	B22021627-011H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

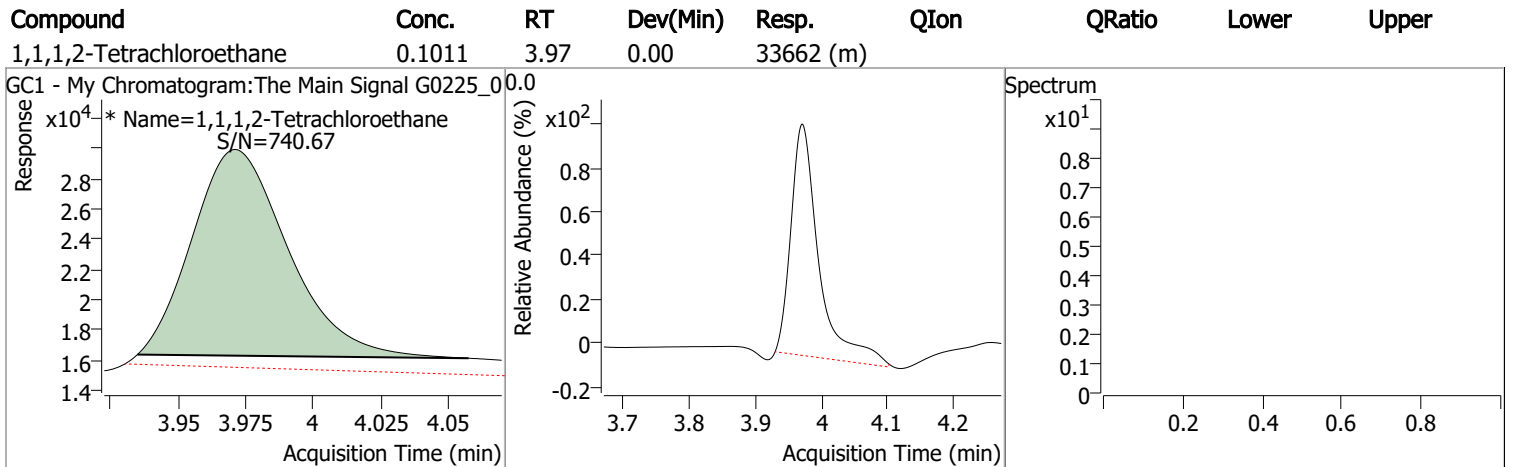
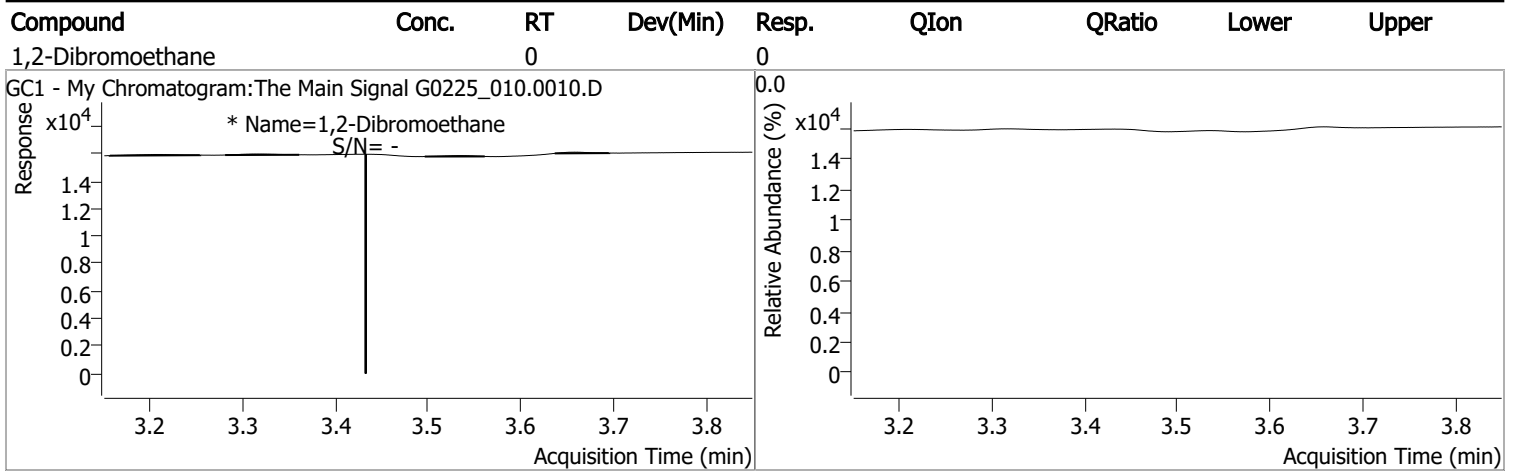
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.971	0.0	33662	0.1011	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 101.12%		
Target Compounds						
M 1,2-Dibromoethane	3.433	0.0	0		µg/L	md
						QValue 1

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

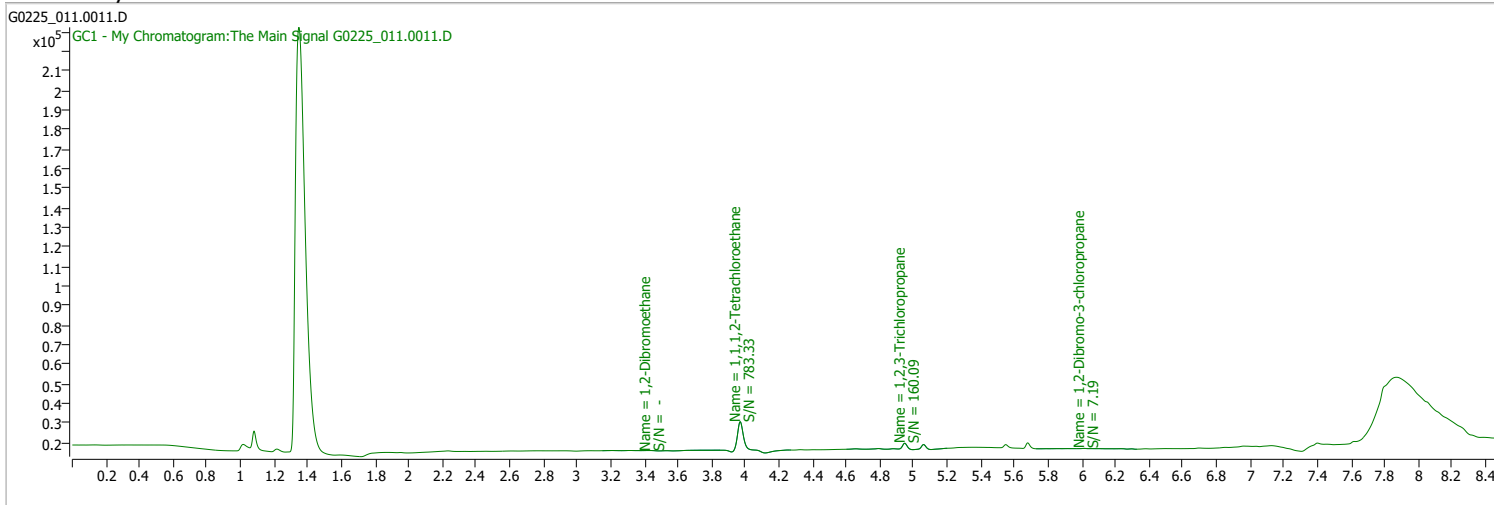
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_011.0011.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 3:55:15 PM
Sample Name	B22021627-014A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

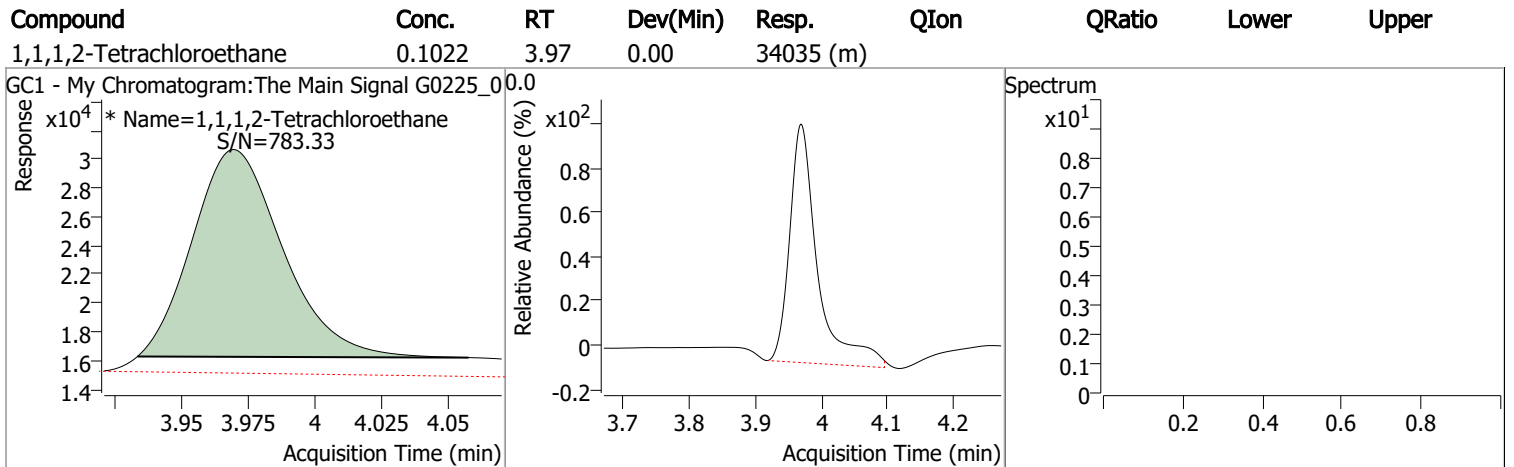
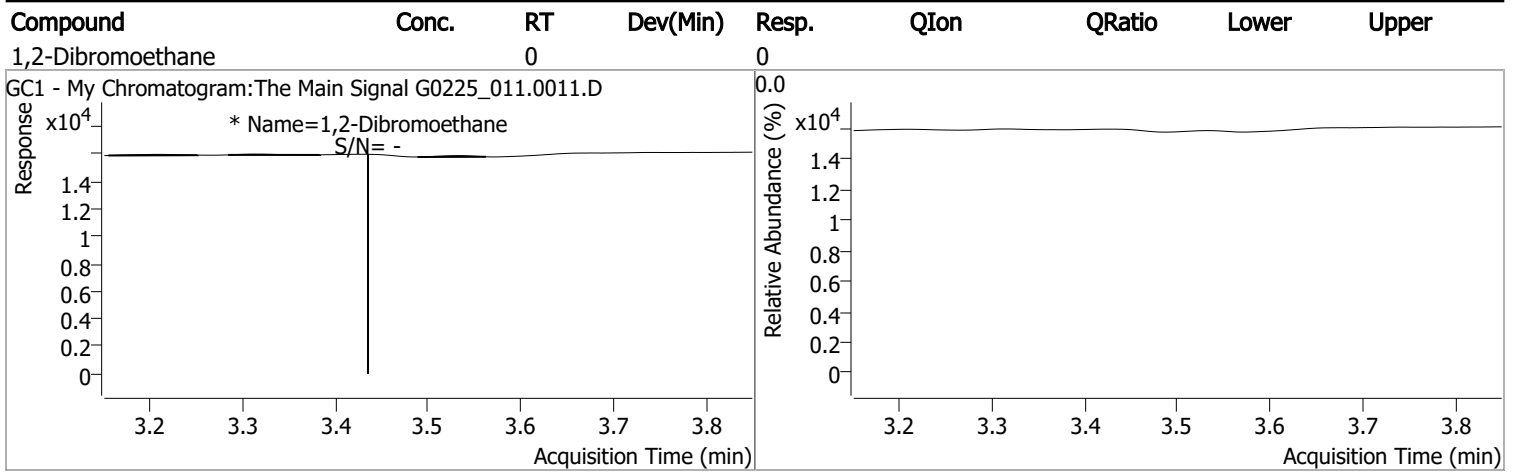
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.969	0.0	34035	0.1022	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 102.16%		
Target Compounds						
M 1,2-Dibromoethane	3.435	0.0	0		µg/L	md
						QValue
						1

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

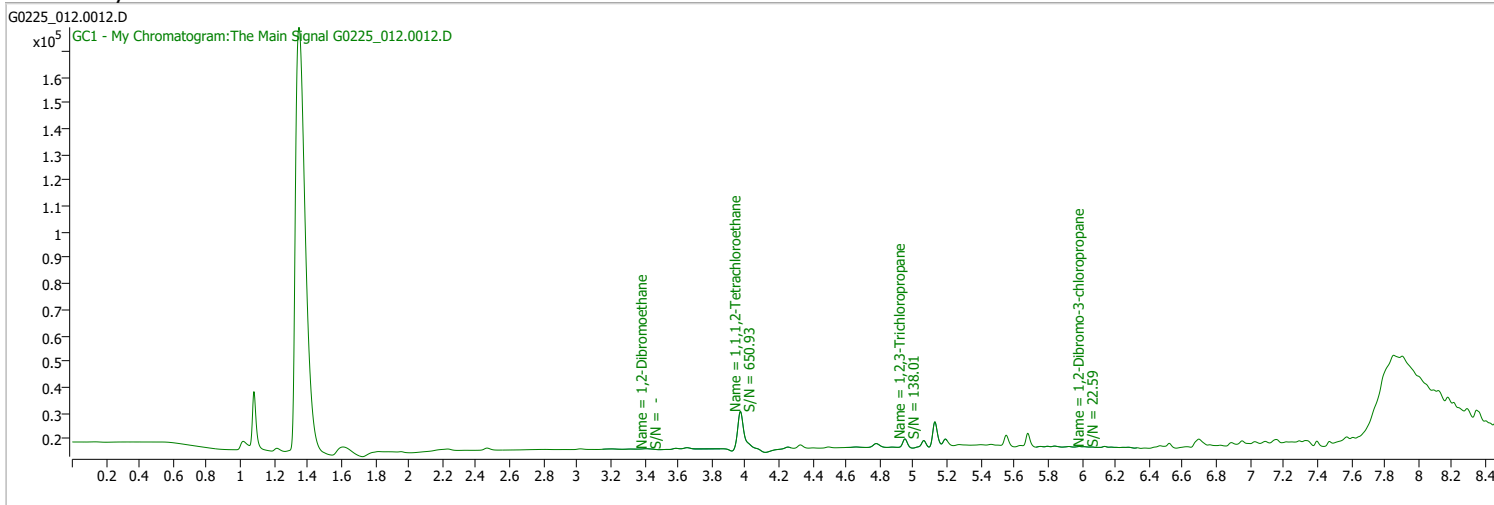
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_012.0012.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 4:15:00 PM
Sample Name	B22021435-032H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

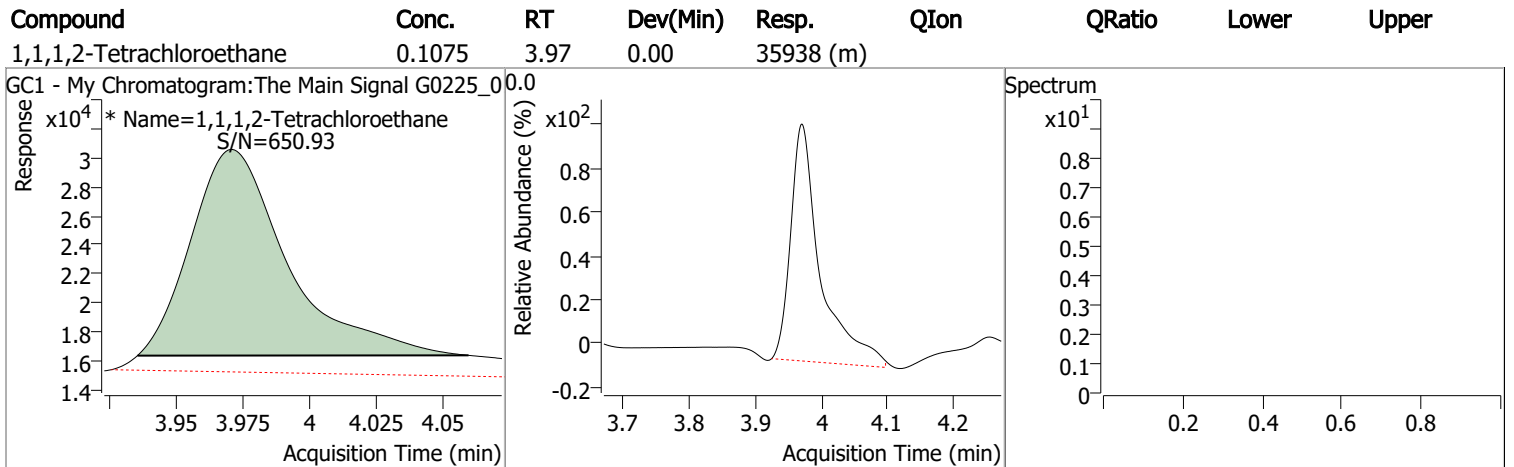
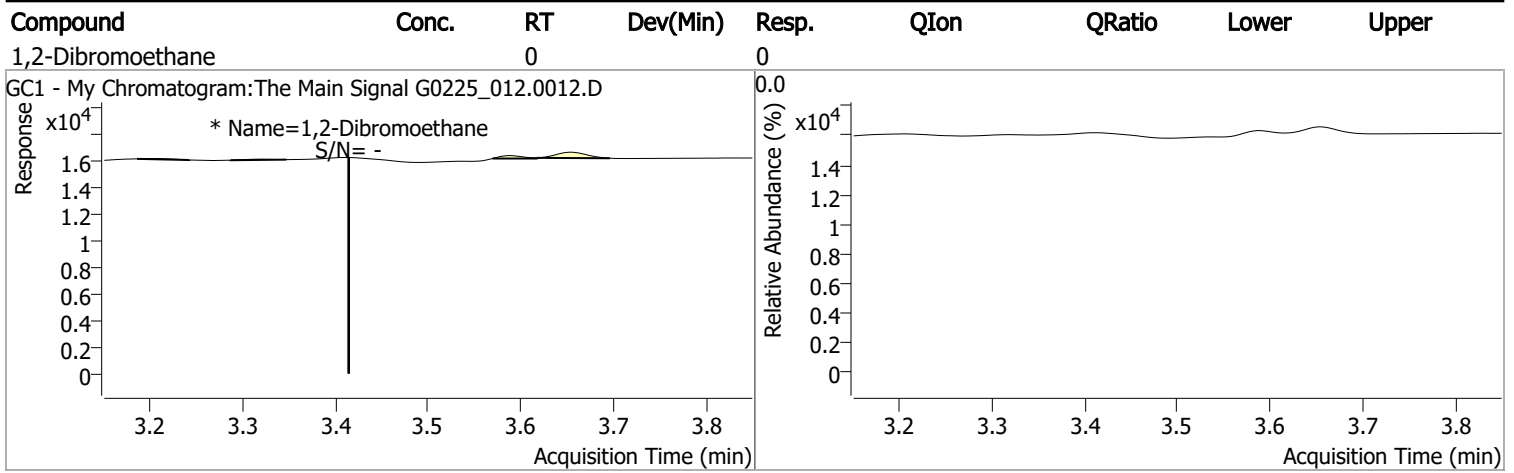
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.971	0.0	35938	0.1075	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 107.45%		
Target Compounds						
M 1,2-Dibromoethane	3.414	0.0	0		µg/L	md
						QValue 1

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

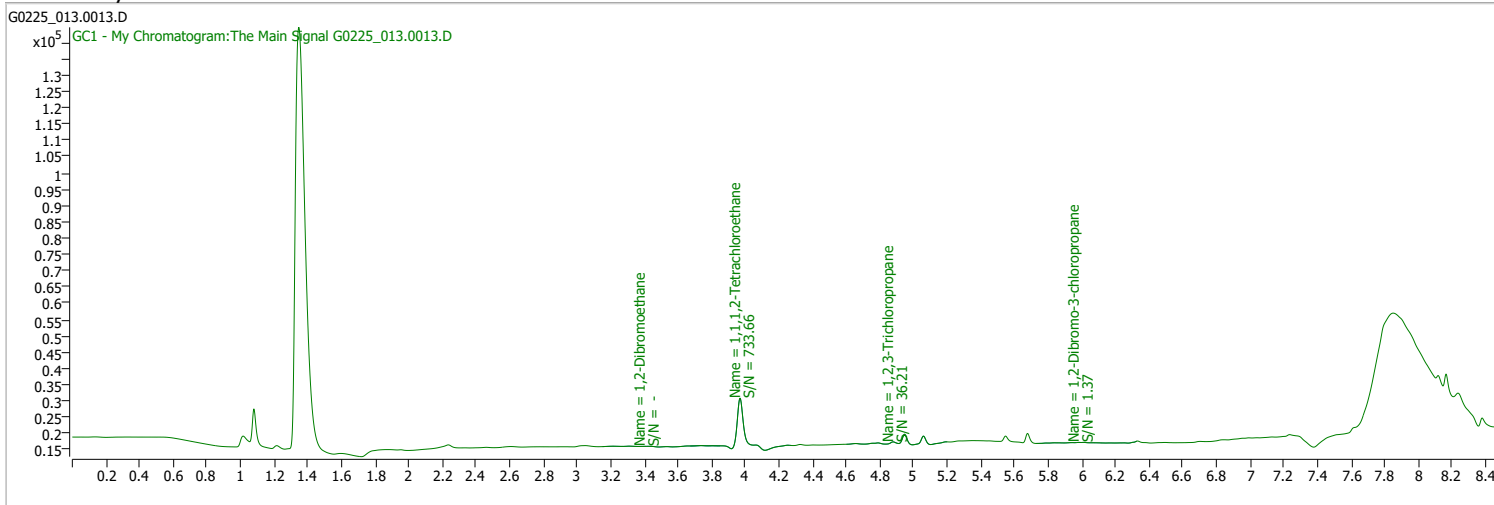
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_013.0013.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 4:34:40 PM
Sample Name	B22021435-035A	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

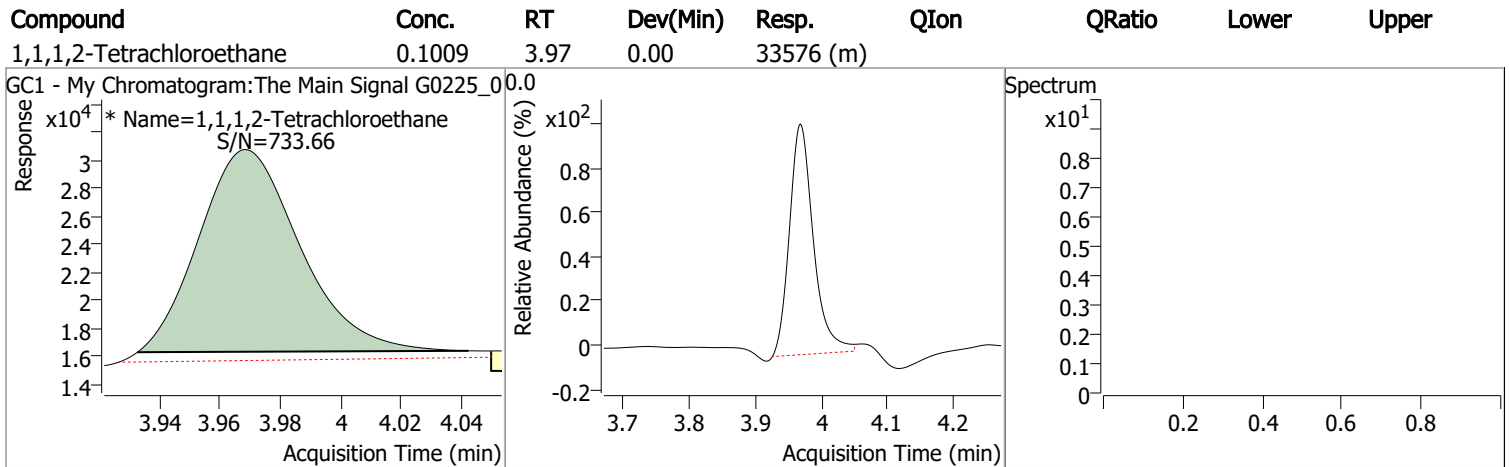
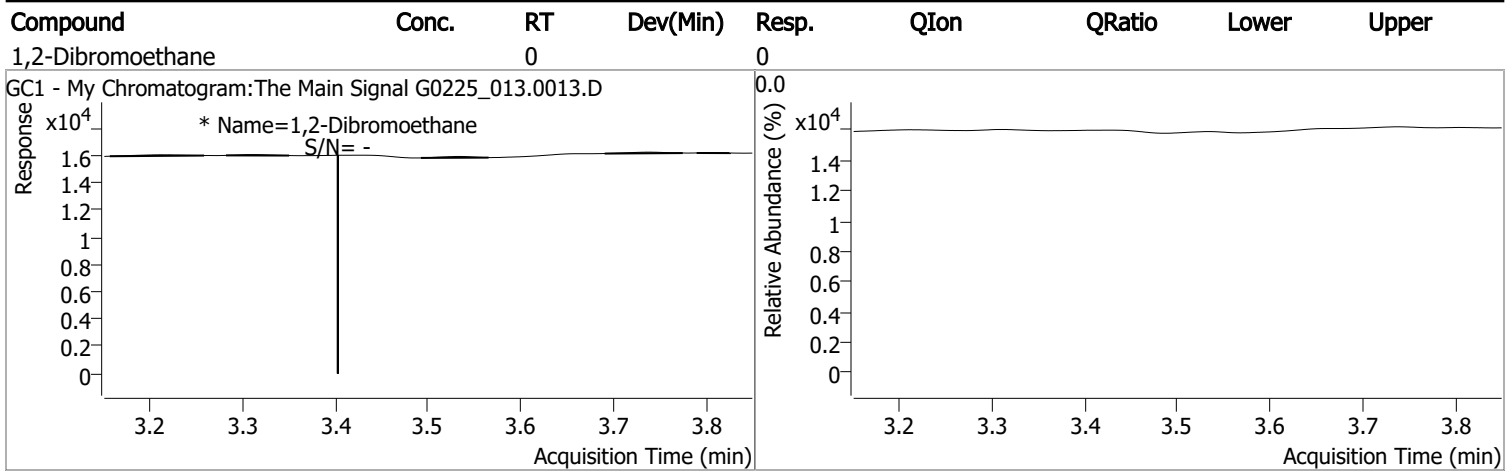
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.968	0.0	33576	0.1009	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 100.88%		
Target Compounds						
M 1,2-Dibromoethane	3.403	0.0	0		µg/L	md
						QValue 1

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

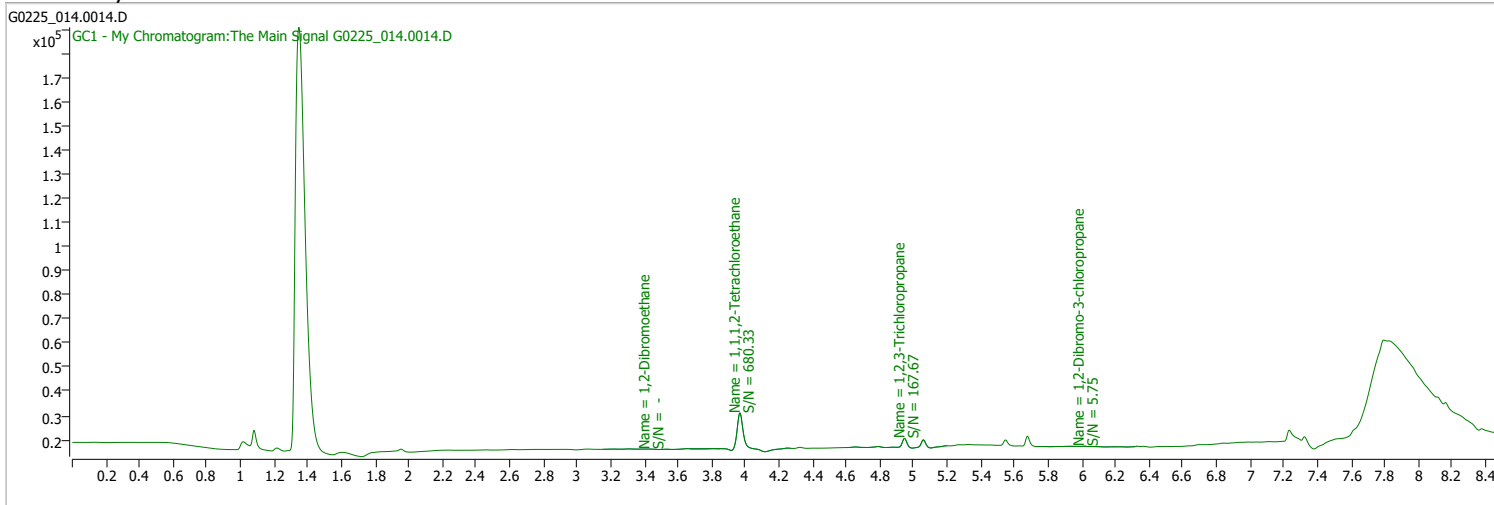
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_014.0014.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 4:54:24 PM
Sample Name	B22021627-001H	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

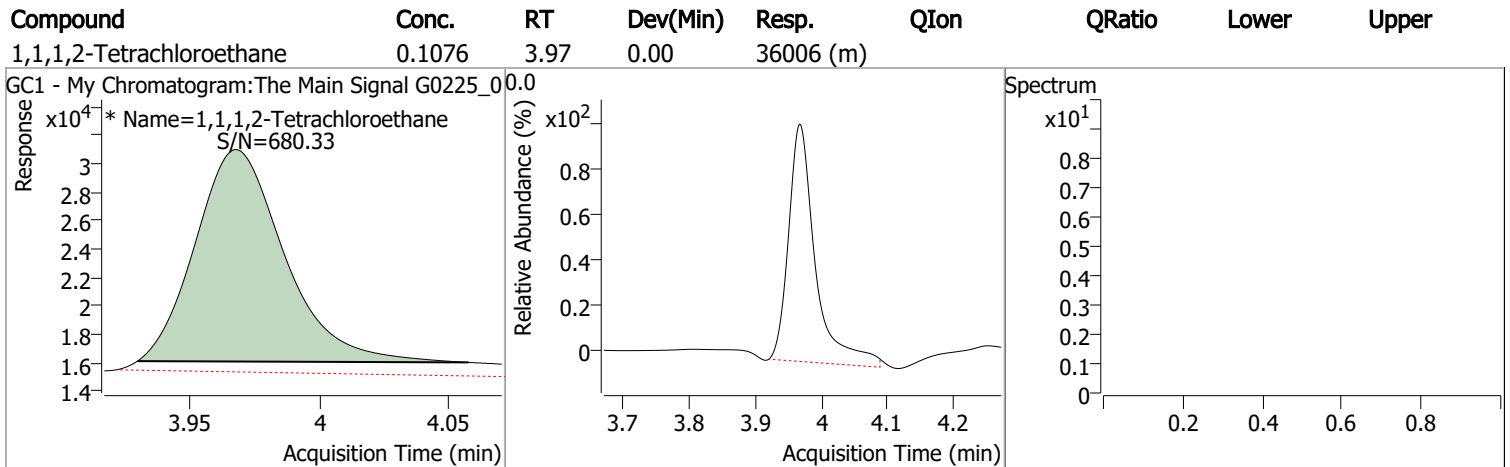
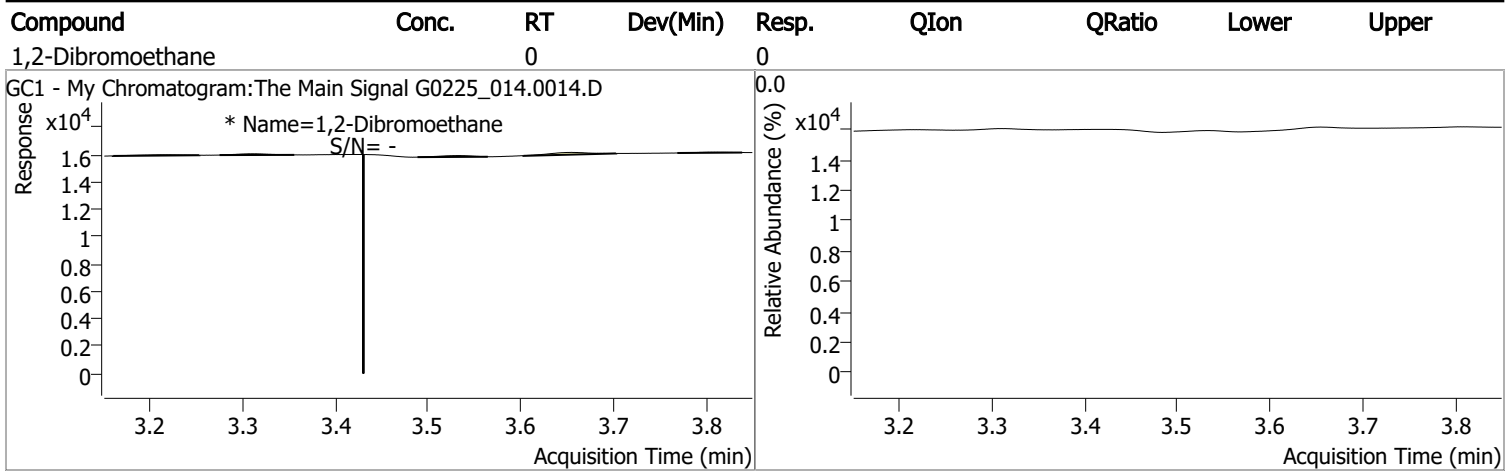
S 1,1,1,2-Tetrachloroethane	3.968	0.0	36006	0.1076	µg/L	m	-0.004
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 107.64%			

Target Compounds

M 1,2-Dibromoethane	3.430	0.0	0	µg/L	md	QValue	1
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(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

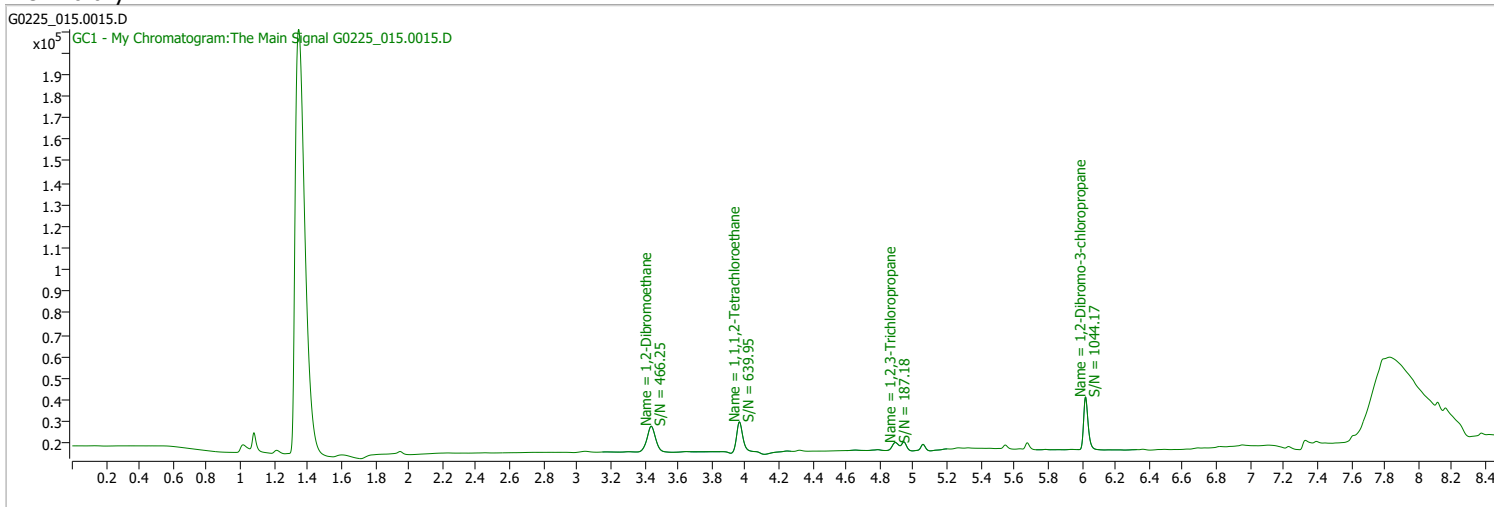
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_015.0015.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 5:14:06 PM
Sample Name	B22021627-001HMS	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

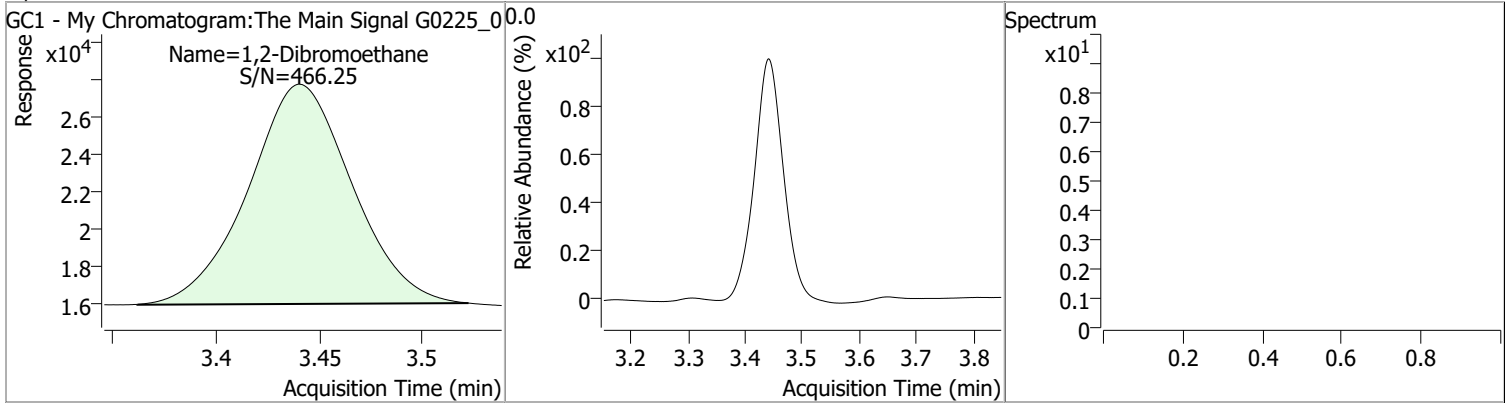


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.965	0.0	33895	0.1018	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 101.77%		
Target Compounds						
M 1,2-Dibromoethane	3.440	0.0	41398	0.2538	µg/L	QValue 100

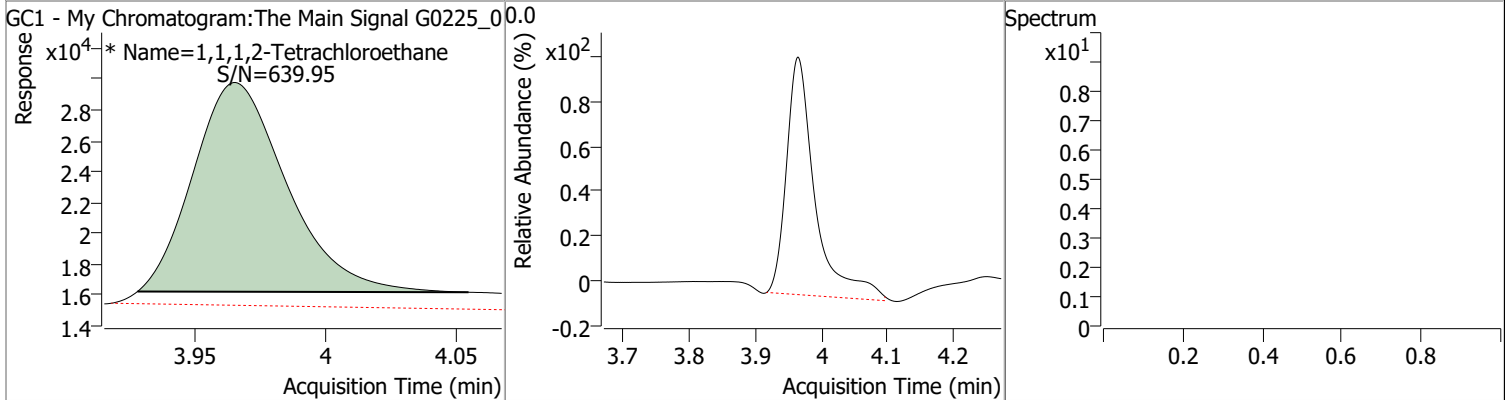
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2538	3.44	-0.01	41398				



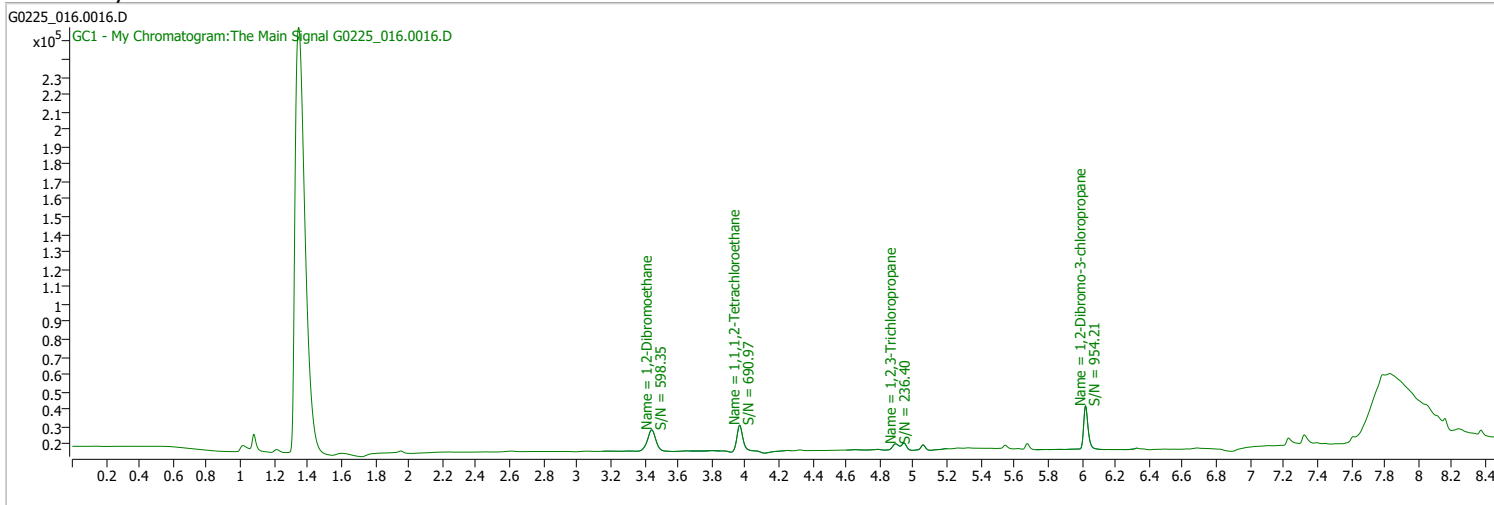
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1018	3.97	-0.01	33895 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0225_016.0016.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 5:34:08 PM
Sample Name	B22021627-001HMSD	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

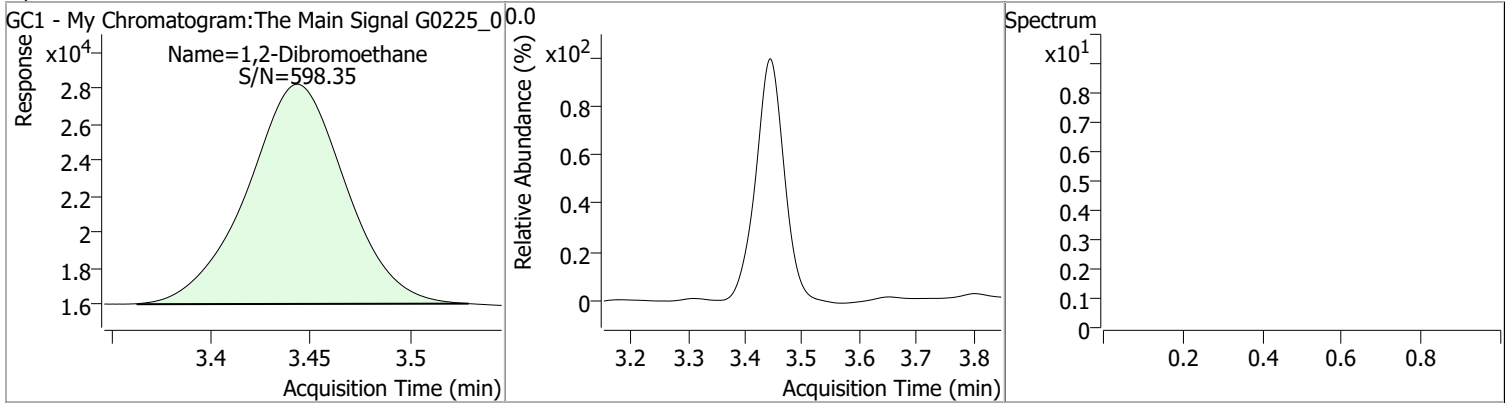


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.966	0.0	35520	0.1063	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 106.29%		
Target Compounds						
M 1,2-Dibromoethane	3.443	0.0	42255	0.2593	µg/L	100

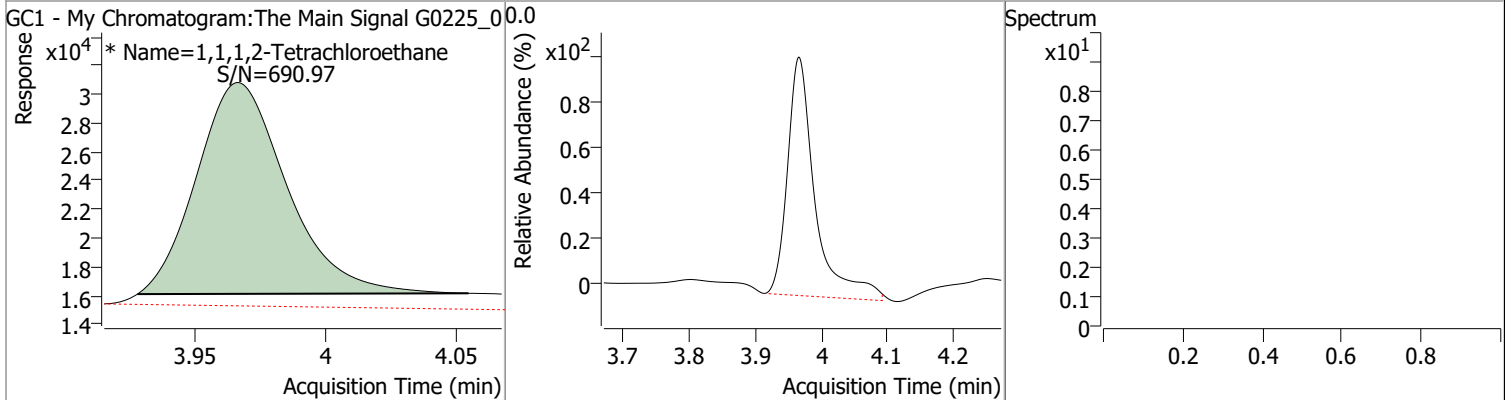
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.2593	3.44	-0.01	42255				



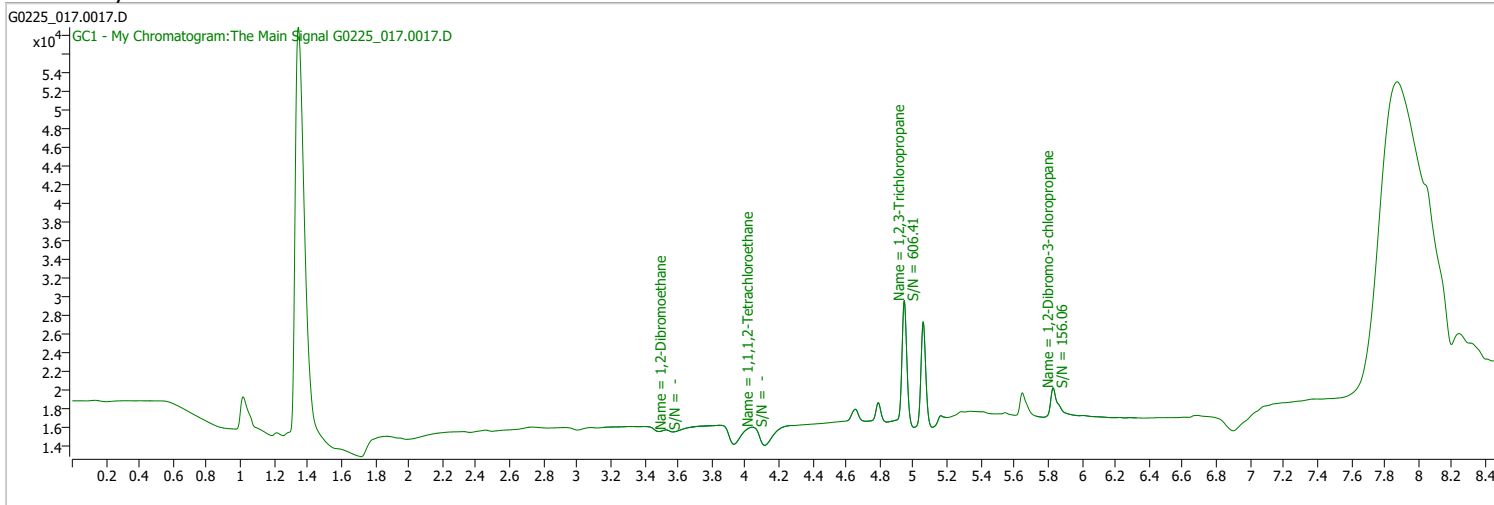
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1063	3.97	-0.01	35520 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0225_017.0017.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 5:53:54 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

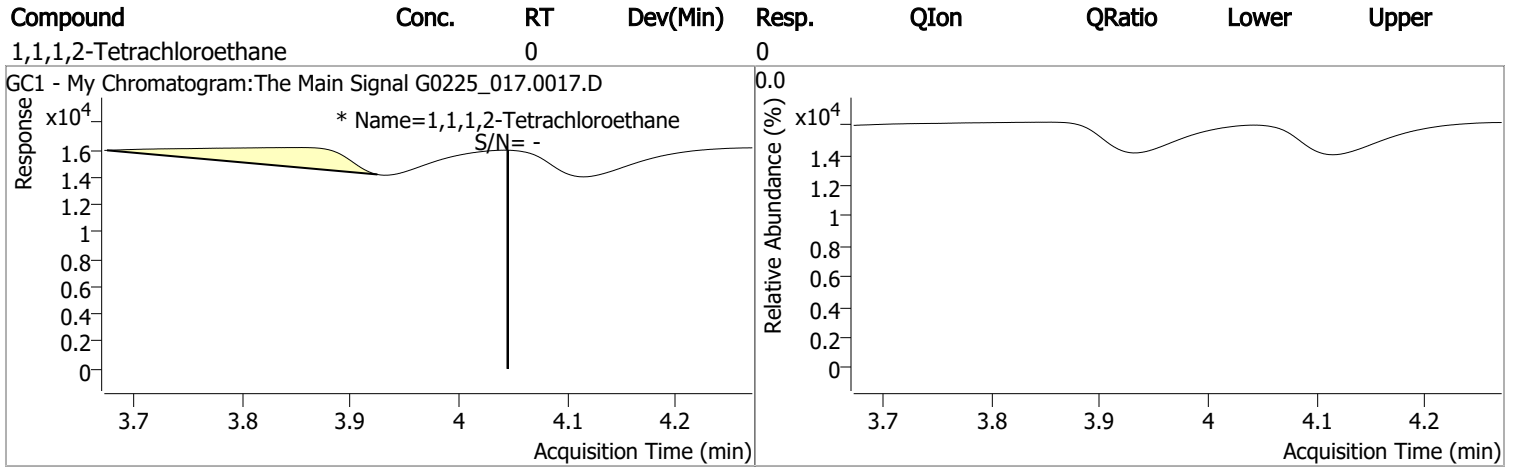
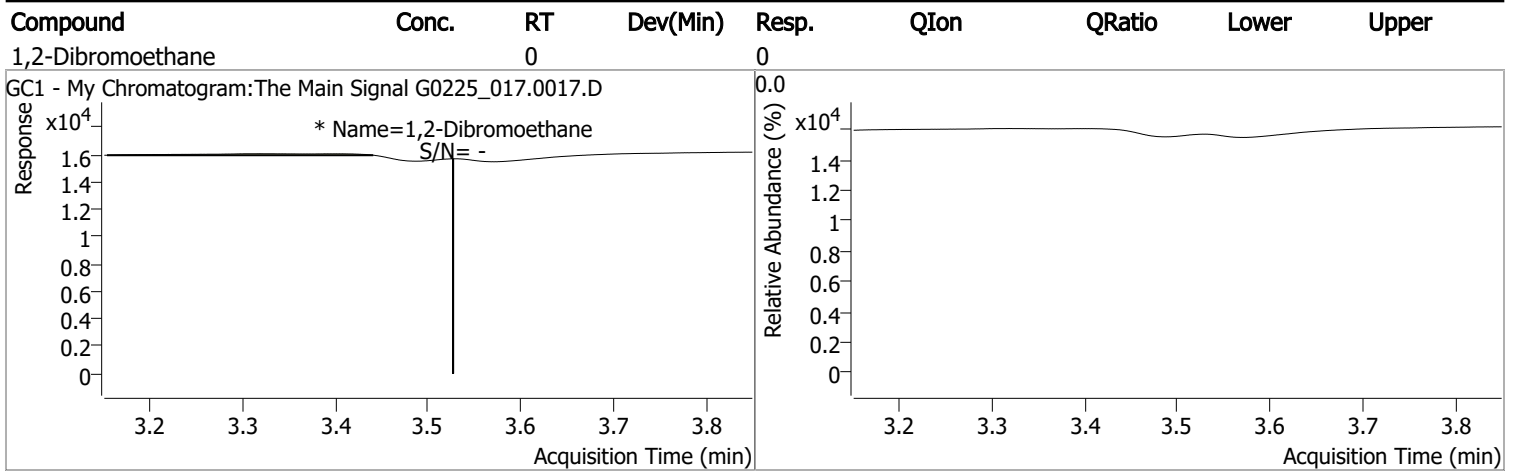
S 1,1,1,2-Tetrachloroethane	4.045	0.0	0		µg/L	md	0.073
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = NA%			

Target Compounds

M 1,2-Dibromoethane	3.527	0.0	0		µg/L	md	QValue 1
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(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

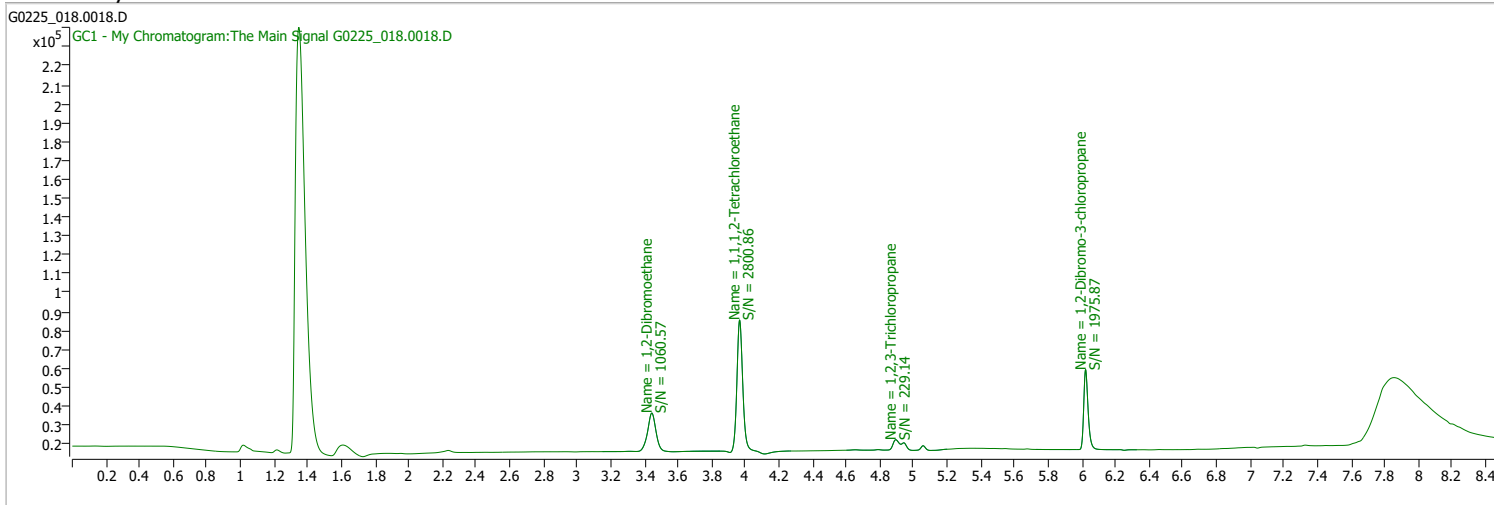
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_018.0018.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 6:13:41 PM
Sample Name	CK5-164037	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

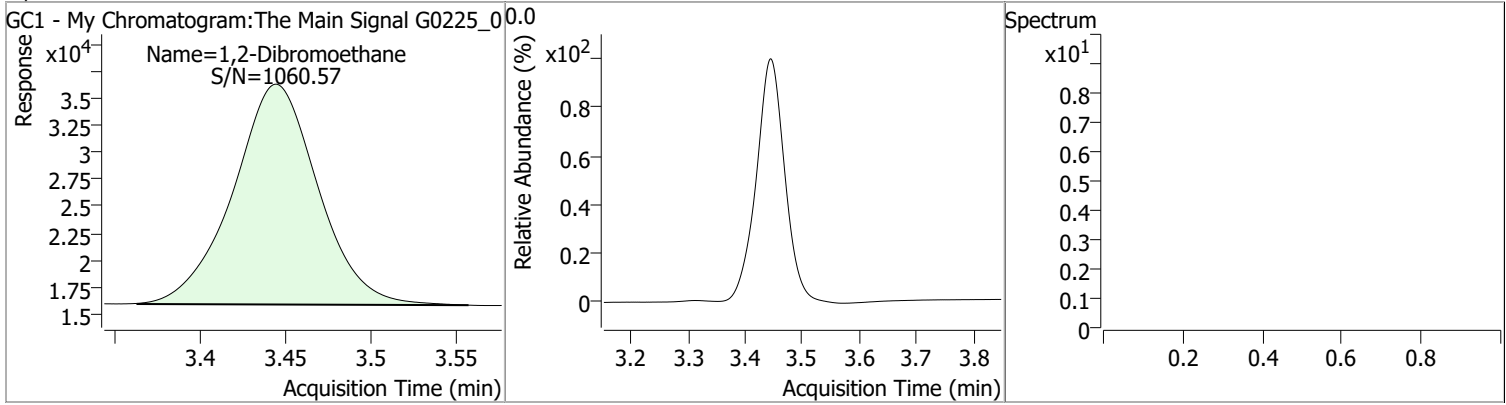


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.966	0.0	166563	0.4591	µg/L	m
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 459.15%		*
Target Compounds						
M 1,2-Dibromoethane	3.444	0.0	70936	0.4496	µg/L	QValue 100

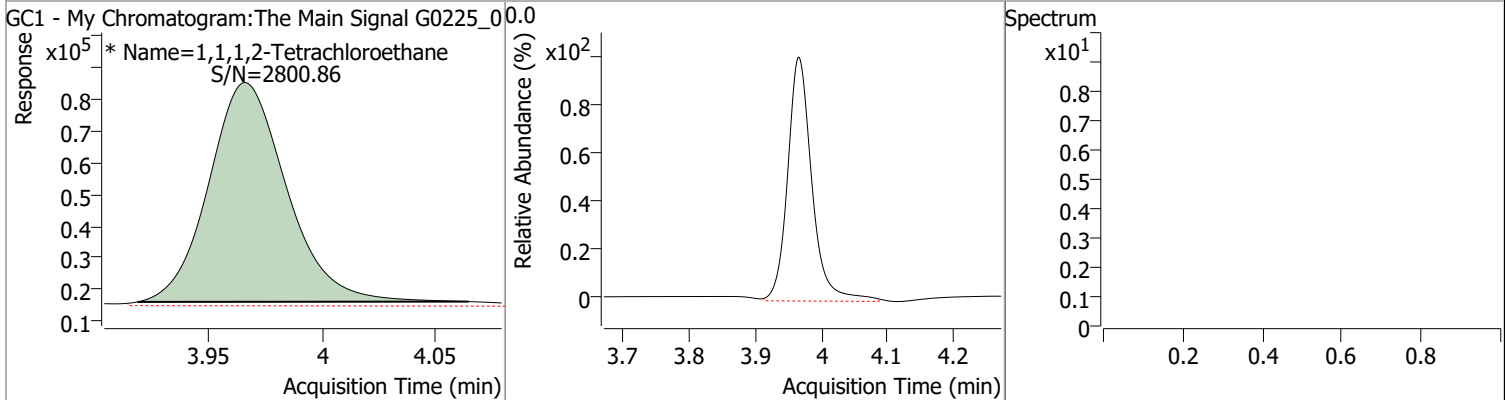
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.4496	3.44	-0.01	70936				



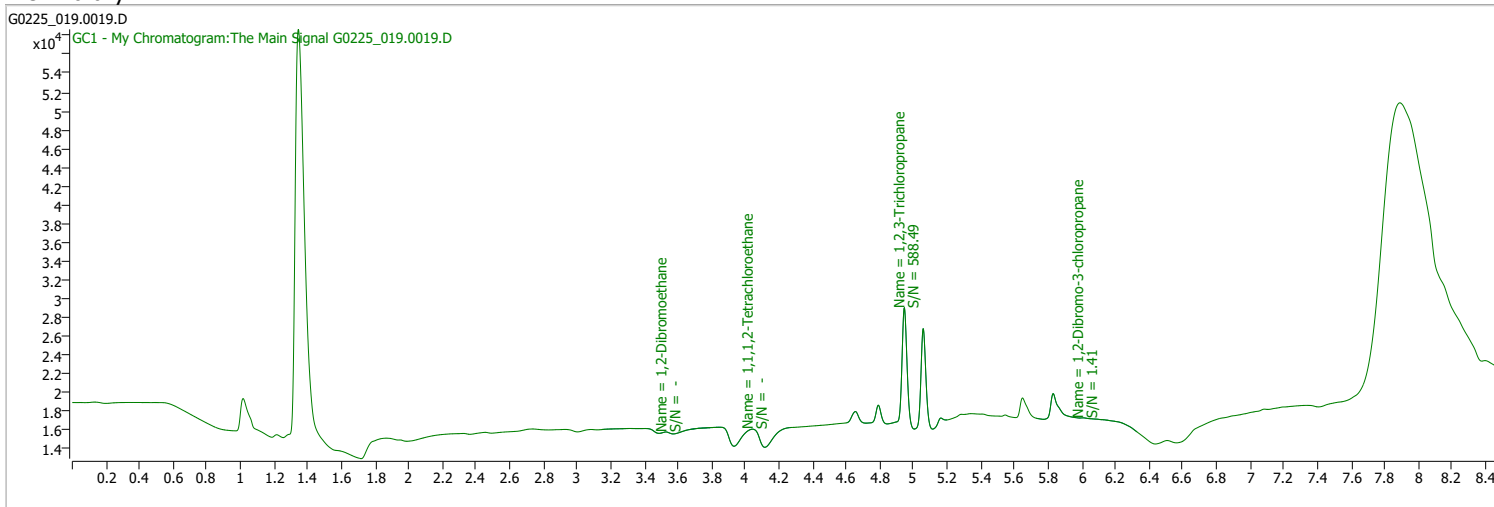
Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.4591	3.97	-0.01	166563 (m)				



Quantitation Results Report (QT Reviewed)

Data File	G0225_019.0019.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 6:33:22 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

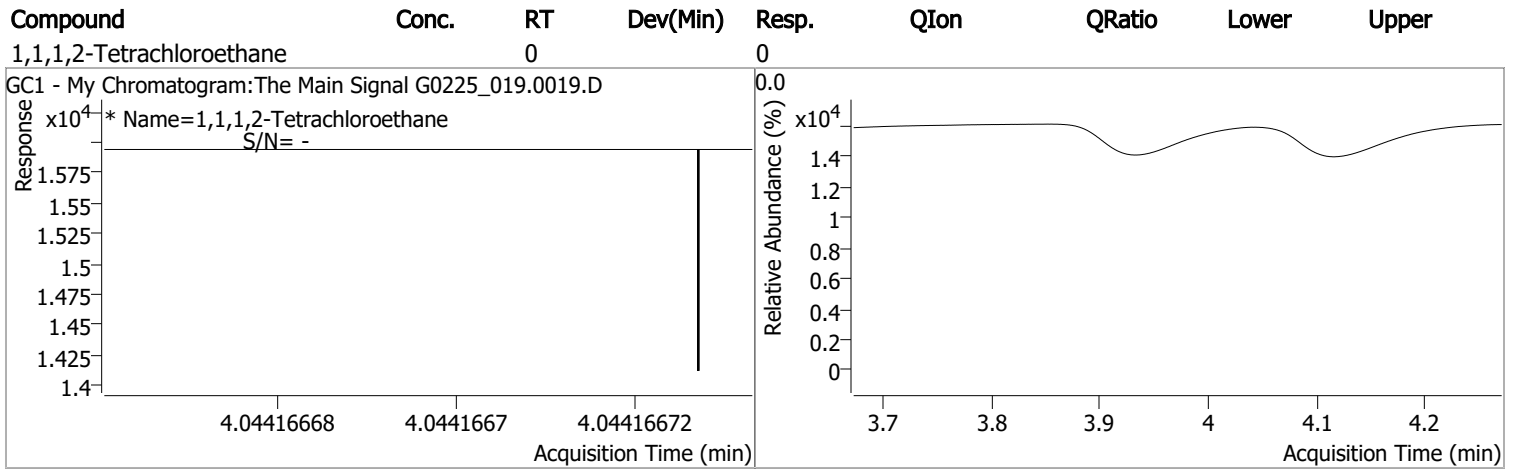
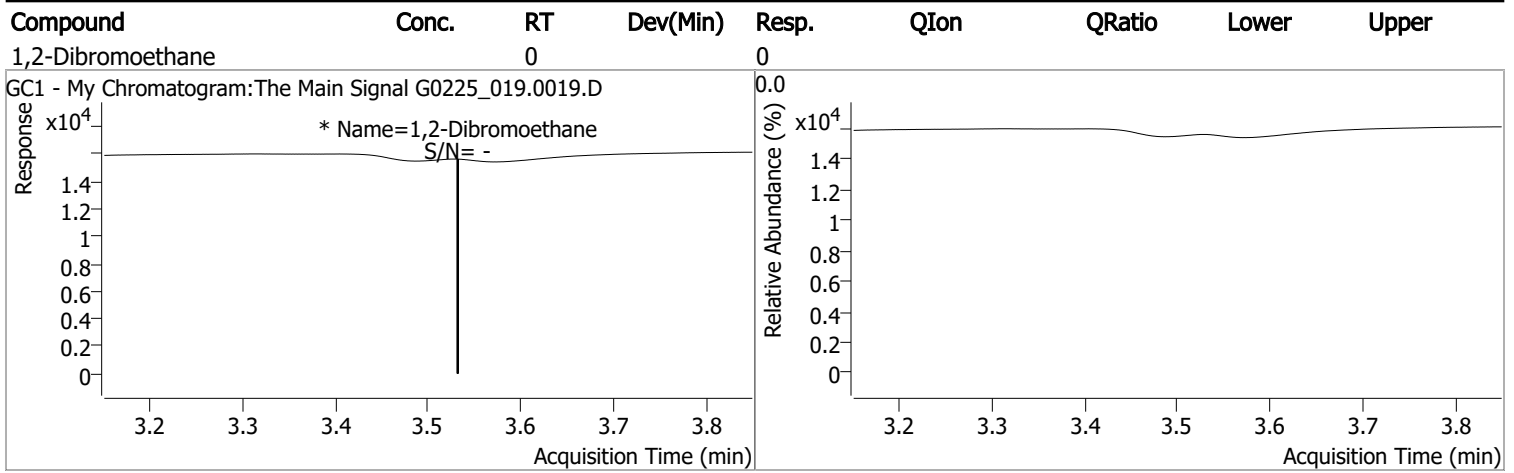
S 1,1,1,2-Tetrachloroethane	4.044	0.0	0		µg/L	md	0.072
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = NA%			

Target Compounds

M 1,2-Dibromoethane	3.532	0.0	0		µg/L	md	QValue
							1

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

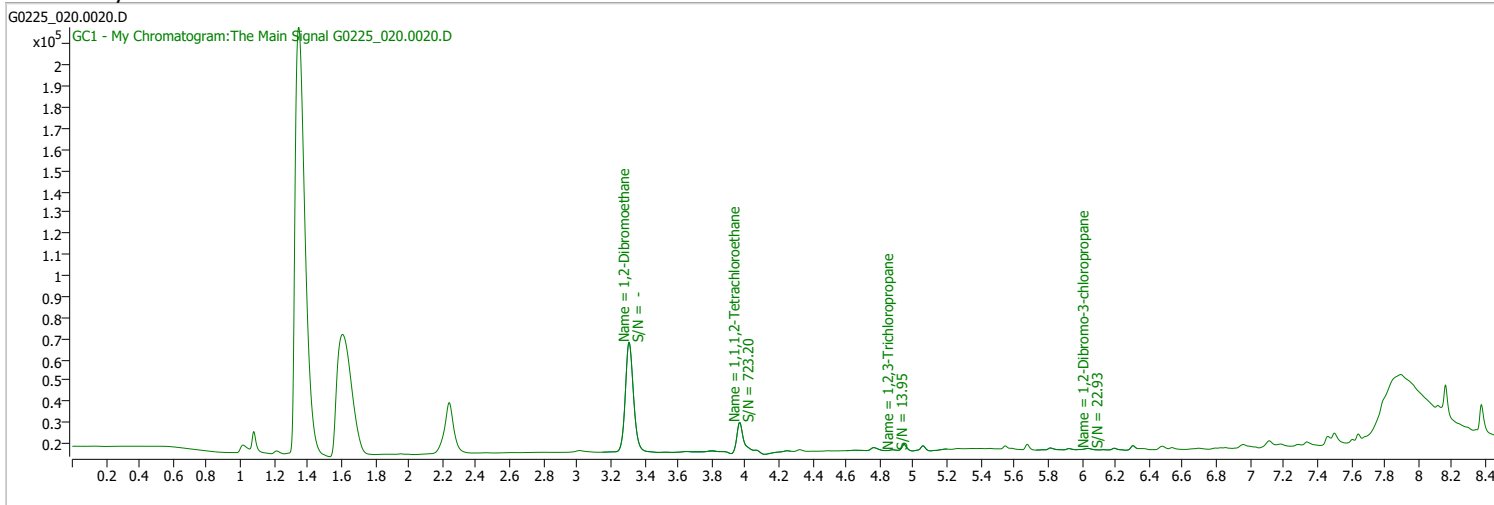
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_020.0020.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 6:53:05 PM
Sample Name	B22021528-002C	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
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Internal Standards

System Monitoring Compounds

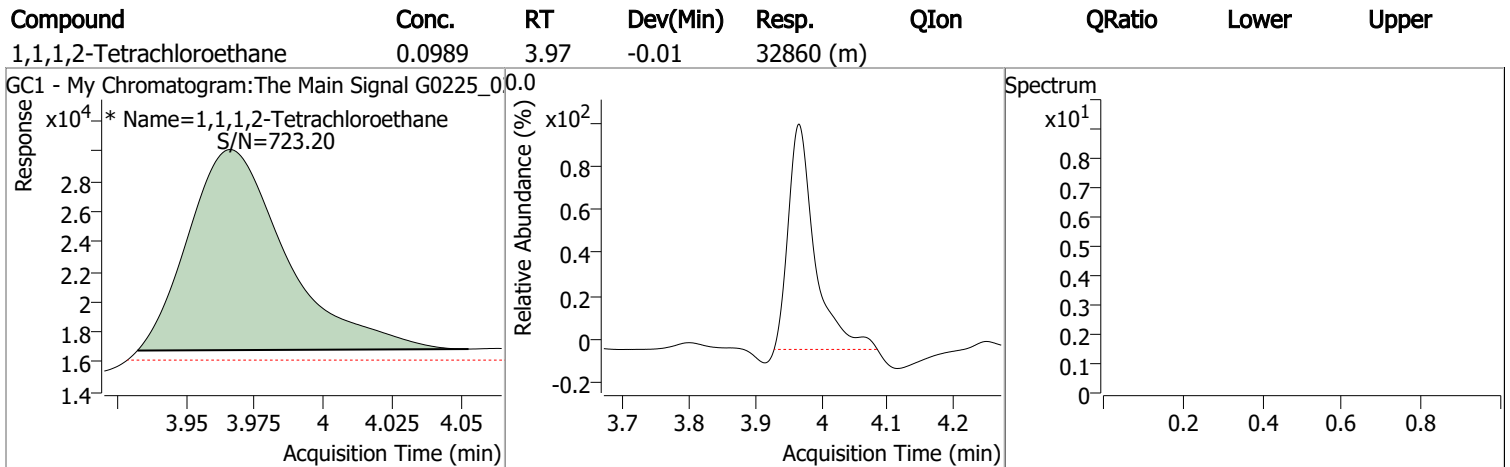
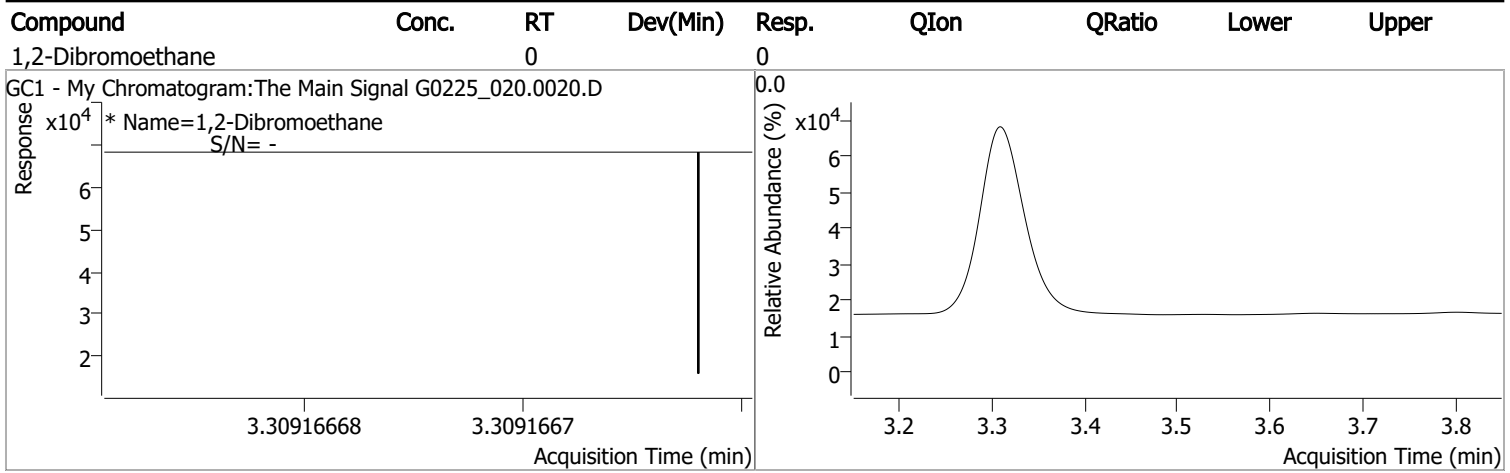
S 1,1,1,2-Tetrachloroethane	3.966	0.0	32860	0.0989	µg/L	m	-0.006
Spiked Amount: 0.100	Range: 70.0 - 130.0%			Recovery = 98.88%			

Target Compounds

M 1,2-Dibromoethane	3.309	0.0	0	µg/L	md	QValue	1
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(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

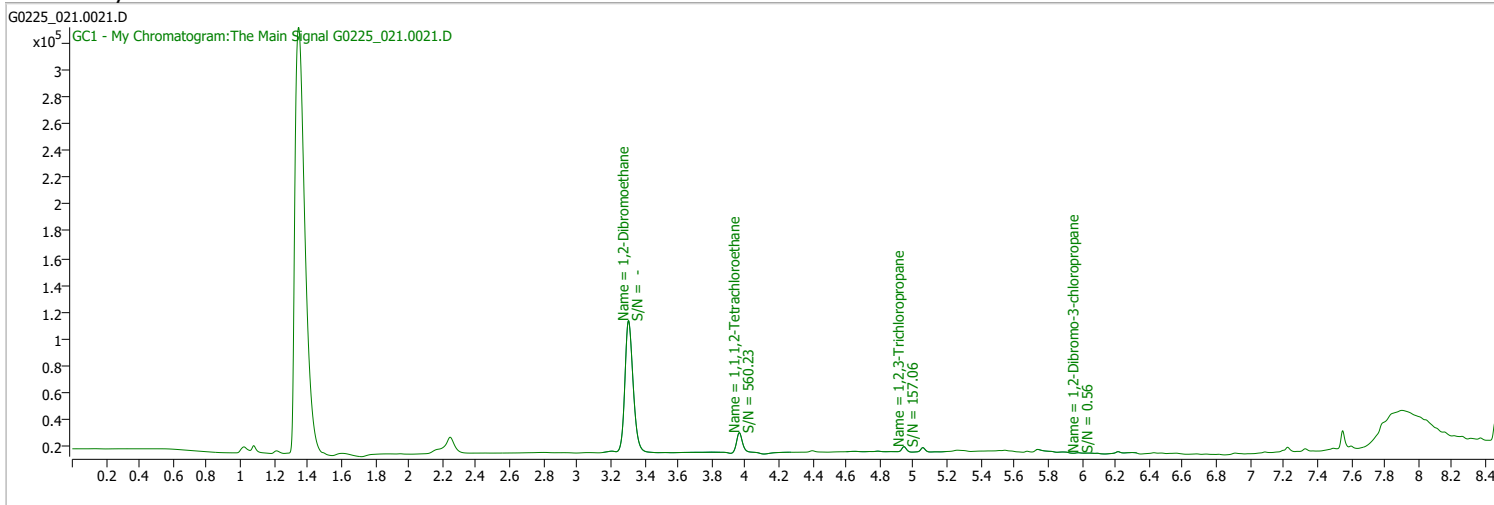
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_021.0021.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 7:12:52 PM
Sample Name	B22021528-003C	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

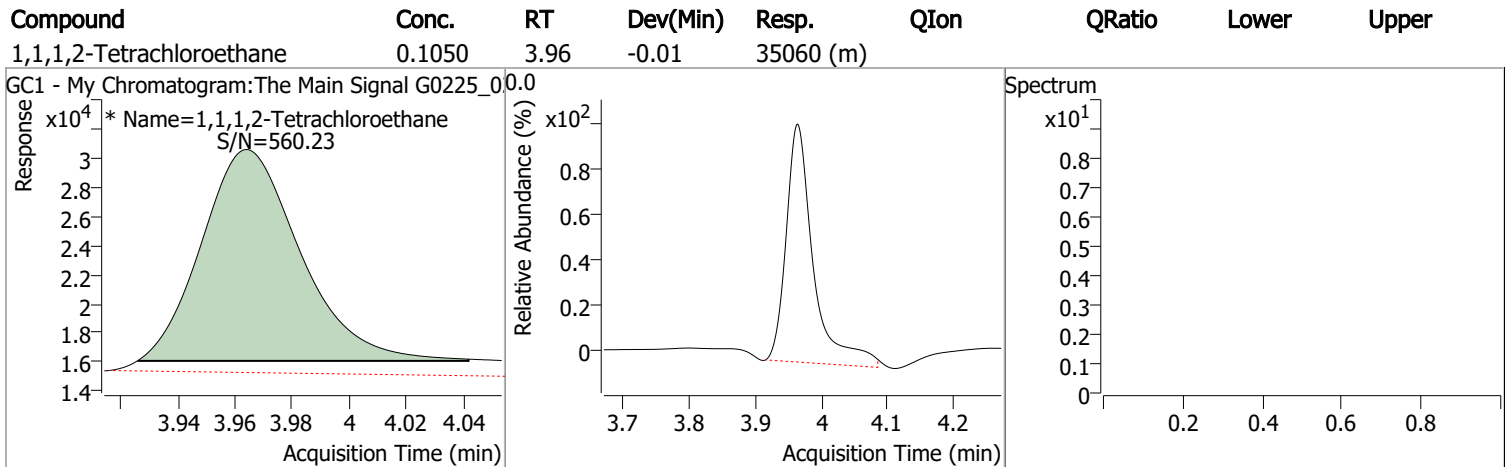
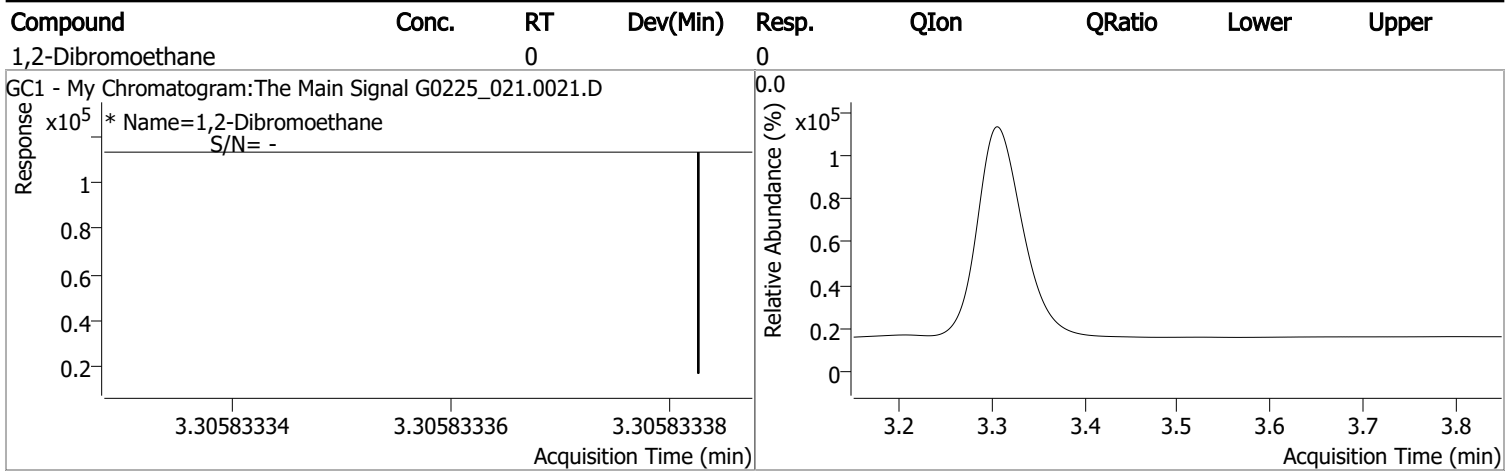
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.964	0.0	35060	0.1050	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 105.01%		
Target Compounds						
M 1,2-Dibromoethane	3.306	0.0	0		µg/L	md
						QValue
						1

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

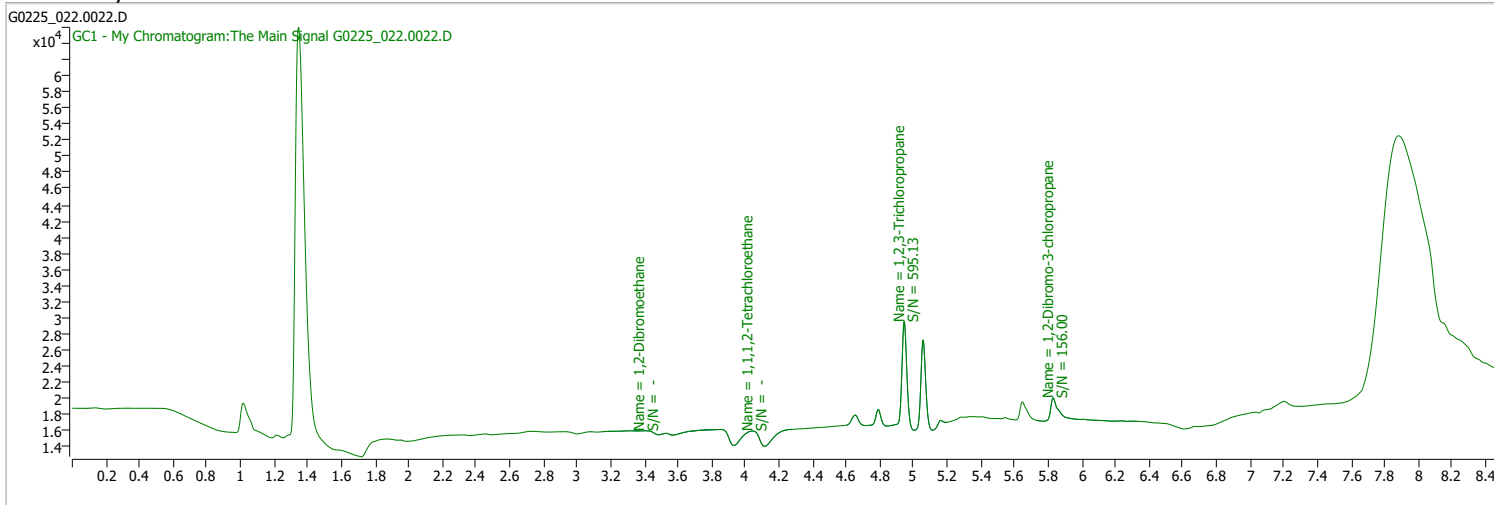
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_022.0022.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 7:32:33 PM
Sample Name	Hexan	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

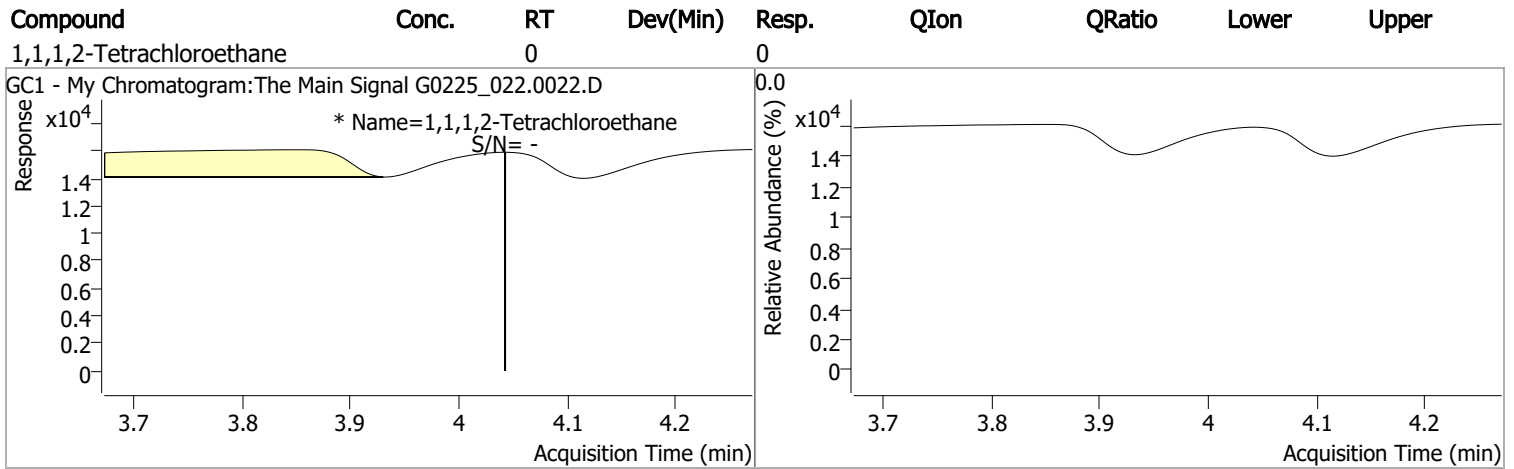
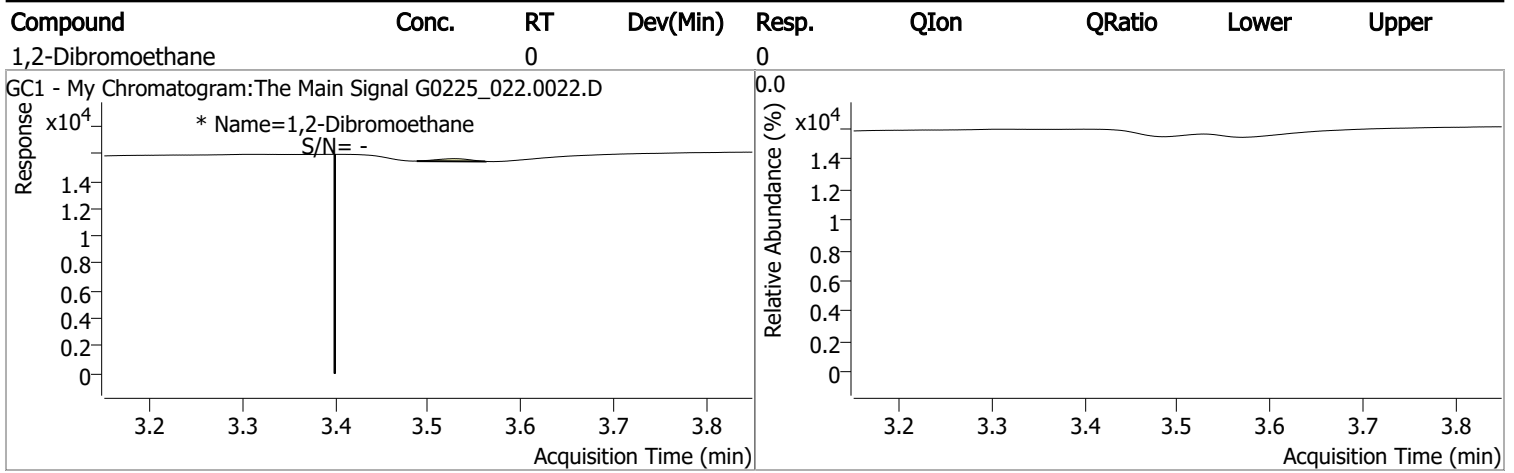
Ref Library



Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	4.043	0.0	0		µg/L	md
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = NA%		
Target Compounds						
M 1,2-Dibromoethane	3.399	0.0	0		µg/L	md
						QValue
						1

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

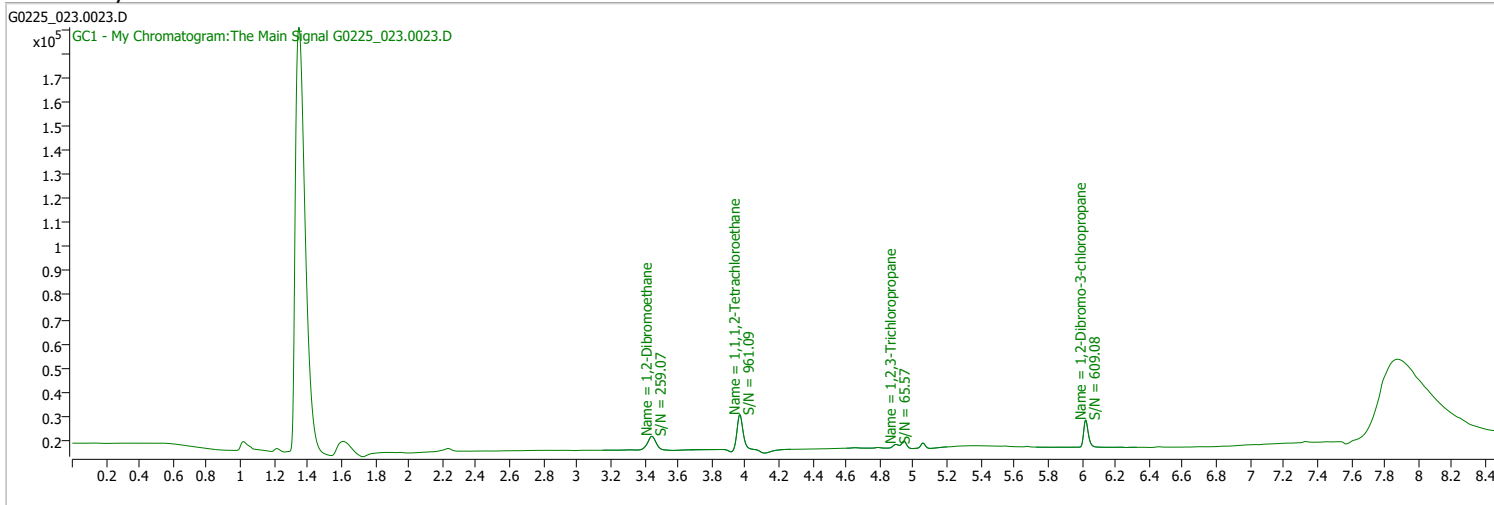
Quantitation Results Report (QT Reviewed)



Quantitation Results Report (QT Reviewed)

Data File	G0225_023.0023.D	Operator	
Acq. Method	testAcqFileNamePath	Acq. Date-Time	2/25/2022 7:52:24 PM
Sample Name	CK3-164037	Instrument	WJB
Vial		Multiplier	1.00
DA Method File	G022422_8011_W_CLT.m	Comment	
Tune File		Tune Date	
Batch Name	G022522_8011_W_CLT.batch.bin	Last Calib Update	2/25/2022 8:11:20 AM

Ref Library

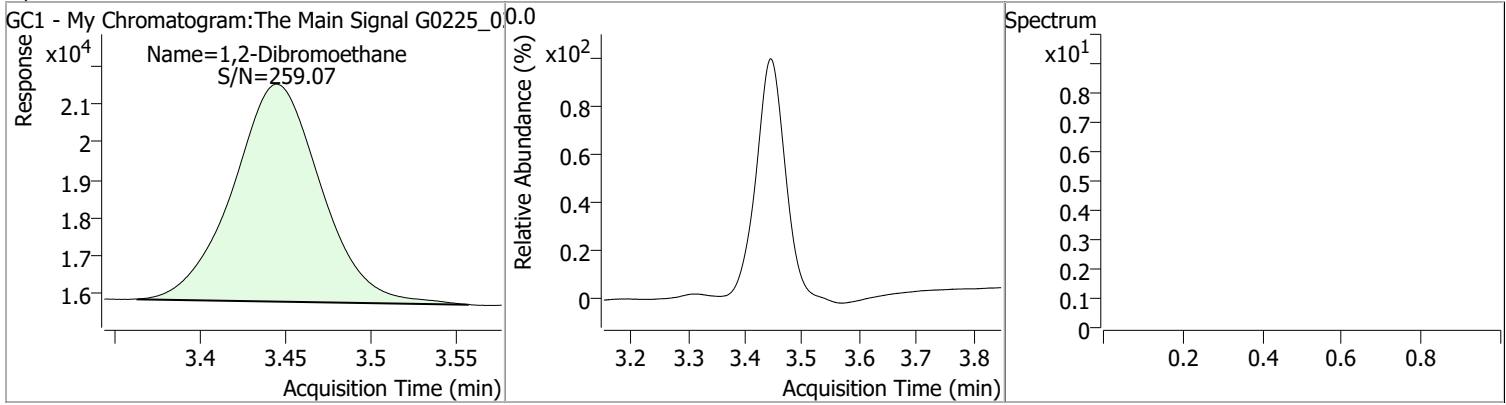


Compound	RT	QIon	Resp.	Conc.	Units	Dev(Min)
Internal Standards						
System Monitoring Compounds						
S 1,1,1,2-Tetrachloroethane	3.967	0.0	35489	0.1062	µg/L	m
Spiked Amount: 0.100		Range: 70.0 - 130.0%		Recovery = 106.20%		
Target Compounds						
M 1,2-Dibromoethane	3.444	0.0	20595	0.1236	µg/L	QValue 100

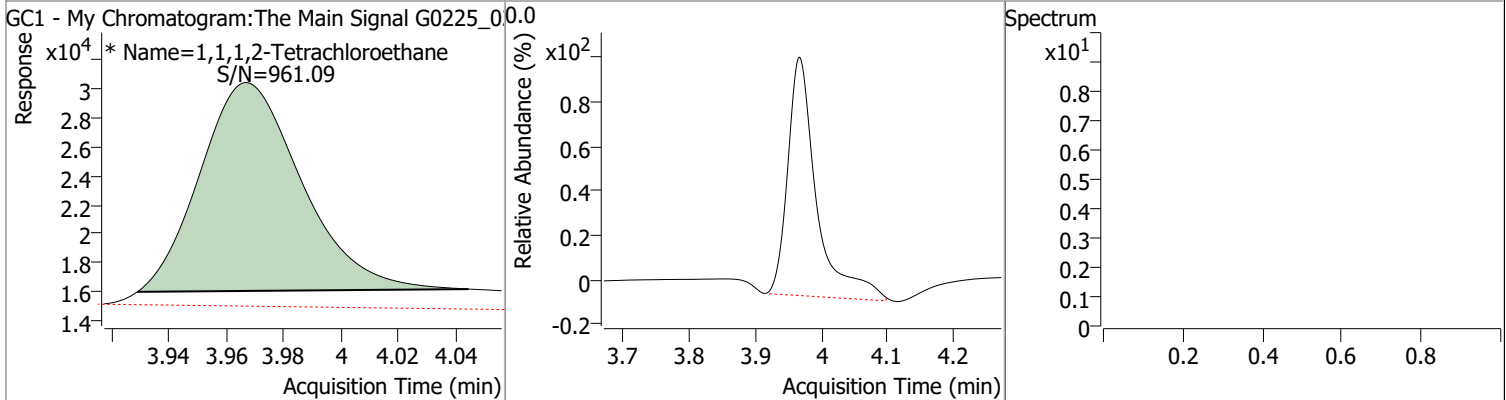
(#) = Qualifier Out of Range; (m) = Manual Integration; (+) = Area Summed; (*) = Surrogate Percent Recovery Out of Range; (d): Zeroed Peak

Quantitation Results Report (QT Reviewed)

Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,2-Dibromoethane	0.1236	3.44	-0.01	20595				



Compound	Conc.	RT	Dev(Min)	Resp.	QIon	QRatio	Lower	Upper
1,1,1,2-Tetrachloroethane	0.1062	3.97	-0.01	35489 (m)				



Audit Trail report

Batch name and path: \\MASSHUNTER\Org\Data\GEC.D.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin
Quant batch version: 10.0
Quant reporting version: 10.0

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdNewBatchTable	BL2000\ctran	2/25/2022 1:28:42 PM	Create new batch \\MASSHUNTER\Org\Data\GEC.D.I\G022522\aiexport\G022522_8011_W_CLT.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/25/2022 1:28:53 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GEC.D.I\G022522\aiexport\G0225_003.0003.D, \\MASSHUNTER\Org\Data\GEC.D.I\G022522\aiexport\G0225_002(1).0002.D, \\MASSHUNTER\Org\Data\GEC.D.I\G022522\aiexport\G0225_001.0001.D			✓	
CmdStartMethodEditing	BL2000\ctran	2/25/2022 1:29:06 PM	Start method editing			✓	
CmdImportMethodFromFile	BL2000\ctran	2/25/2022 1:29:06 PM	Import method from file \\MASSHUNTER\Org\Data\GEC.D.I\GEC.D_methods\G022422_8011_W_CLT.m			✓	
CmdApplyMethodToAllSamples	BL2000\ctran	2/25/2022 1:30:21 PM	Apply method to all samples			✓	
CmdMethodClear	BL2000\ctran	2/25/2022 1:30:21 PM	Clear method			✓	
CmdEndMethodEditing	BL2000\ctran	2/25/2022 1:30:22 PM	End method editing			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 1:30:22 PM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:30:31 PM	Set SampleType = CC for sample G0225_002(1).0002.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 1:30:33 PM	Set LevelName = 3 for sample G0225_002(1).0002.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 1:30:35 PM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 1:30:45 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_002(1).0002.D, from x, y = 3.935, 16073 to 4.054, 16214, result = 33077; previous integration is from x, y = 3.931, 15646 to 4.098, 14852 and previous response = 41286.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/25/2022 1:30:47 PM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_002(1).0002.D to y = 16073, new integration is from x, y = 3.935, 16073 to 4.054, 16073 and new response = 33580; previous integration is from x, y = 3.935, 16073 to 4.054, 16214 and previous response = 33077.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSaveBatchTable	BL2000\ctran	2/25/2022 1:31:26 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/25/2022 3:21:09 PM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G022522_8011_W_CLT.batch.bin			✓	
CmdImportSamplesFromWorklist	BL2000\ctran	2/25/2022 3:22:36 PM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_008.0008.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_007.0007.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_006.0006.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_005.0005.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_004.0004.D			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 3:22:51 PM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 3:23:08 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_008.0008.D, from x, y = 3.933, 16094 to 4.068, 16115, result = 35585; previous integration is from x, y = 3.923, 15286 to 4.090, 14776 and previous response = 45658.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 3:23:18 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_007.0007.D, from x, y = 3.933, 16240 to 4.054, 16349, result = 32025; previous integration is from x, y = 3.919, 15329 to 4.091, 14915 and previous response = 43217.			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:23 PM	Set SampleType = DoubleBlank for sample G0225_006.0006.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:26 PM	Set SampleType = Blank for sample G0225_003.0003.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:28 PM	Set SampleType = QC for sample G0225_004.0004.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:31 PM	Set LevelName = LCS for sample G0225_004.0004.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:34 PM	Set SampleType = QC for sample G0225_005.0005.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:37 PM	Set LevelName = LCS1 for sample G0225_005.0005.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 3:23:47 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_004.0004.D, from x, y = 3.933, 16161 to 4.063, 16141, result = 32779; previous integration is from x, y = 3.920, 15333 to 4.096, 14837 and previous response = 42913.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/25/2022 3:23:52 PM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_005.0005.D, from x, y = 3.933, 16151 to 4.053, 16260, result = 32562; previous integration is from x, y = 3.920, 15359 to 4.096, 14877 and previous response = 42864.			✓	
CmdSetSampleAttribute	BL2000\ctran	2/25/2022 3:23:57 PM	Set SampleType = DoubleBlank for sample G0225_001.0001.D; previous value = Sample			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 3:23:58 PM	Quantitate all compounds in all samples			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:25:32 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_006.0006.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:25:34 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_007.0007.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:25:35 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_008.0008.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:25:37 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_006.0006.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:26:36 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_003.0003.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:26:54 PM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_001.0001.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/25/2022 3:26:55 PM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_001.0001.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 3:27:10 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_002(1).0002.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 3:27:17 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_004.0004.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 3:27:19 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_005.0005.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 3:27:24 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_007.0007.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/25/2022 3:27:25 PM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_008.0008.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 3:27:36 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/25/2022 3:27:38 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 3:28:22 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/25/2022 3:28:23 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	
CmdQuantitate	BL2000\ctran	2/25/2022 3:30:22 PM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/25/2022 3:30:24 PM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\ctran	2/28/2022 7:31:48 AM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G022522_8011_W_CLT.batch.bin			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdImportSamplesFromWorklist	BL2000\ctran	2/28/2022 7:34:29 AM	Add samples from worklist: \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_023.0023.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_022.0022.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_021.0021.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_020.0020.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_019.0019.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_018.0018.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_017.0017.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_016.0016.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_015.0015.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_014.0014.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_013.0013.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_012.0012.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_011.0011.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_010.0010.D, \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G0225_009.0009.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:01 AM	Set SampleType = MatrixBlank for sample G0225_014.0014.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:04 AM	Set SampleType = Matrix for sample G0225_015.0015.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:06 AM	Set SampleType = MatrixDup for sample G0225_016.0016.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:13 AM	Set SampleType = Blank for sample G0225_017.0017.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:14 AM	Set SampleType = DoubleBlank for sample G0225_017.0017.D; previous value = Blank			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:16 AM	Set SampleType = CC for sample G0225_018.0018.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:19 AM	Set LevelName = 5 for sample G0225_018.0018.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:21 AM	Set SampleType = DoubleBlank for sample G0225_019.0019.D; previous value = Sample			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:23 AM	Set SampleType = DoubleBlank for sample G0225_022.0022.D; previous value = Sample			✓	
CmdQuantitate	BL2000\ctran	2/28/2022 7:35:28 AM	Quantitate all compounds in all samples			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:35:40 AM	Set SampleApproved = True for sample G0225_002(1).0002.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:36:06 AM	Set SampleApproved = True for sample G0225_003.0003.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:36:13 AM	Set UserAnnotation = LT for compound 1,1,1,2-Tetrachloroethane in sample G0225_004.0004.D; previous value = GT			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:36:14 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_004.0004.D; previous value = LT			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:36:16 AM	Set SampleApproved = True for sample G0225_004.0004.D; previous value = False			✓	
CmdManuallyIntegrateSnapBaseline	BL2000\ctran	2/28/2022 7:36:28 AM	Snap baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_003.0003.D, from x = 3.932 to x = 4.078, new integration is from x, y = 3.932, 15708 to 4.078, 15938 and new response = 35537; previous integration is from x, y = 3.932, 15727 to 4.078, 15937 and previous response = 35455.			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:36:29 AM	Clear manual integration of target signal for compound 1,1,1,2-Tetrachloroethane in sample G0225_003.0003.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:36:33 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_003.0003.D, from x, y = 3.935, 16063 to 4.052, 16240, result = 32762; previous integration is from x, y = 3.932, 15727 to 4.078, 15937 and previous response = 35455.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/28/2022 7:36:34 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_003.0003.D to y = 16063, new integration is from x, y = 3.935, 16063 to 4.052, 16063 and new response = 33382; previous integration is from x, y = 3.935, 16063 to 4.052, 16240 and previous response = 32762.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:36:36 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_003.0003.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:36:43 AM	Set SampleApproved = True for sample G0225_005.0005.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:36:46 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_006.0006.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:36:48 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_006.0006.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:36:49 AM	Set SampleApproved = True for sample G0225_006.0006.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:37:00 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_007.0007.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:37:02 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_008.0008.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:37:04 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_009.0009.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:37:11 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_012.0012.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:37:23 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_014.0014.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:37:30 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_017.0017.D			✓	
CmdManuallyIntegrateSnapBaseline	BL2000\ctran	2/28/2022 7:37:37 AM	Snap baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_018.0018.D, from x = 3.914 to x = 4.087, new integration is from x, y = 3.914, 15635 to 4.087, 15422 and new response = 172249; previous integration is from x, y = 3.914, 14890 to 4.087, 14711 and previous response = 179787.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:37:51 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_018.0018.D, from x, y = 3.918, 16068 to 4.065, 16172, result = 166563; previous integration is from x, y = 3.914, 15635 to 4.087, 15422 and previous response = 172249.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:37:53 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_018.0018.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdUpdateRetentionTimes	BL2000\ctran	2/28/2022 7:38:18 AM	Update retention time for compound 1,2,3-Trichloropropane; 1,2-Dibromo-3-chloropropane; 1,2-Dibromoethane; 1,1,1,2-Tetrachloroethane;			✓	
CmdQuantitate	BL2000\ctran	2/28/2022 7:38:42 AM	Quantitate all compounds in all samples			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:38:44 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_019.0019.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:38:46 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_019.0019.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:38:48 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_017.0017.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:38:51 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_017.0017.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:39:05 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_020.0020.D, from x, y = 3.933, 16766 to 4.053, 16865, result = 32860; previous integration is from x, y = 3.929, 16130 to 4.084, 16130 and previous response = 38994.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:39:08 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_020.0020.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:39:11 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_020.0020.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:39:13 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_021.0021.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:39:19 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_021.0021.D, from x, y = 3.926, 16057 to 4.042, 16177, result = 34644; previous integration is from x, y = 3.916, 15393 to 4.085, 14909 and previous response = 43641.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/28/2022 7:39:20 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_021.0021.D to y = 16057, new integration is from x, y = 3.926, 16057 to 4.042, 16057 and new response = 35060; previous integration is from x, y = 3.926, 16057 to 4.042, 16177 and previous response = 34644.			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:39:28 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_023.0023.D, from x, y = 3.929, 15974 to 4.044, 16161, result = 35489; previous integration is from x, y = 3.915, 15120 to 4.098, 14639 and previous response = 46715.			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:39:36 AM	Set SampleType = CC for sample G0225_023.0023.D; previous value = Sample			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:39:38 AM	Set LevelName = 3 for sample G0225_023.0023.D; previous value =			✓	
CmdQuantitate	BL2000\ctran	2/28/2022 7:39:40 AM	Quantitate all compounds in all samples			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:39:45 AM	Manually integrate compound 1,2-Dibromoethane in sample G0225_023.0023.D, from x, y = 3.363, 15856 to 3.537, 15818, result = 20077; previous integration is from x, y = 3.363, 15856 to 3.557, 15714 and previous response = 20595.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:39:51 AM	Manually integrate compound 1,2-Dibromoethane in sample G0225_023.0023.D, from x, y = 3.363, 15856 to 3.528, 15854, result = 19896; previous integration is from x, y = 3.363, 15856 to 3.537, 15818 and previous response = 20077.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:39:55 AM	Set UserAnnotation = GT for compound 1,2-Dibromoethane in sample G0225_023.0023.D; previous value =			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:40:28 AM	Manually integrate compound 1,2-Dibromoethane in sample G0225_002(1).0002.D, from x, y = 3.368, 15921 to 3.537, 15953, result = 18024; previous integration is from x, y = 3.368, 15921 to 3.564, 15847 and previous response = 18599.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:40:30 AM	Set UserAnnotation = GT for compound 1,2-Dibromoethane in sample G0225_002(1).0002.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:40:57 AM	Set SampleApproved = True for sample G0225_023.0023.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:41:02 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_023.0023.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:41:05 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_022.0022.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:41:07 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_022.0022.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:41:09 AM	Set SampleApproved = True for sample G0225_022.0022.D; previous value = False			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:41:14 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_021.0021.D; previous value =			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:41:20 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_020.0020.D; previous value = GT			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:41:24 AM	Set SampleApproved = True for sample G0225_021.0021.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:41:25 AM	Set SampleApproved = True for sample G0225_020.0020.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:45:27 AM	Set SampleApproved = True for sample G0225_001.0001.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:45:32 AM	Set SampleApproved = True for sample G0225_007.0007.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:45:36 AM	Set SampleApproved = True for sample G0225_008.0008.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:45:44 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D, from x, y = 3.932, 16068 to 4.050, 16443, result = 36091; previous integration is from x, y = 3.918, 15333 to 4.106, 14907 and previous response = 47512.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/28/2022 7:45:45 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D to y = 16068, new integration is from x, y = 3.932, 16068 to 4.050, 16068 and new response = 37422; previous integration is from x, y = 3.932, 16068 to 4.050, 16443 and previous response = 36091.			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:45:47 AM	Clear manual integration of target signal for compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:45:51 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D, from x, y = 3.935, 16167 to 4.053, 16438, result = 35774; previous integration is from x, y = 3.918, 15333 to 4.106, 14907 and previous response = 47512.			✓	
CmdManuallyIntegrateDropBaseline	BL2000\ctran	2/28/2022 7:45:53 AM	Drop baseline for compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D to y = 16167, new integration is from x, y = 3.935, 16167 to 4.053, 16167 and new response = 36726; previous integration is from x, y = 3.935, 16167 to 4.053, 16438 and previous response = 35774.			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:45:57 AM	Clear manual integration of target signal for compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:00 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D, from x, y = 3.985, 16221 to 4.053, 16167, result = 9529; previous integration is from x, y = 3.918, 15333 to 4.106, 14907 and previous response = 47512.			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:46:00 AM	Clear manual integration of target signal for compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:05 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D, from x, y = 3.933, 16302 to 4.053, 16438, result = 35310; previous integration is from x, y = 3.918, 15333 to 4.106, 14907 and previous response = 47512.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:46:06 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_009.0009.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:46:08 AM	Set SampleApproved = True for sample G0225_009.0009.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:12 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_010.0010.D, from x, y = 3.983, 16383 to 4.057, 15898, result = 10657; previous integration is from x, y = 3.930, 15777 to 4.102, 14825 and previous response = 42130.			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:46:13 AM	Clear manual integration of target signal for compound 1,1,1,2-Tetrachloroethane in sample G0225_010.0010.D			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:16 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_010.0010.D, from x, y = 3.935, 16396 to 4.058, 16151, result = 33662; previous integration is from x, y = 3.930, 15777 to 4.102, 14825 and previous response = 42130.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:46:17 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_010.0010.D; previous value =			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:46:20 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_010.0010.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:46:22 AM	Set SampleApproved = True for sample G0225_010.0010.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:46:25 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_011.0011.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:35 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_011.0011.D, from x, y = 3.933, 16281 to 4.058, 16203, result = 34035; previous integration is from x, y = 3.919, 15259 to 4.095, 14798 and previous response = 45490.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:46:36 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_011.0011.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:46:38 AM	Set SampleApproved = True for sample G0225_011.0011.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:51 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_012.0012.D, from x, y = 3.935, 16385 to 4.054, 16521, result = 35505; previous integration is from x, y = 3.925, 15417 to 4.097, 14850 and previous response = 47607.			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:46:56 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_012.0012.D, from x, y = 3.935, 16385 to 4.060, 16406, result = 35938; previous integration is from x, y = 3.935, 16385 to 4.054, 16521 and previous response = 35505.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:46:58 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_012.0012.D; previous value =			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:47:03 AM	Set SampleApproved = True for sample G0225_012.0012.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:47:06 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_013.0013.D			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:47:14 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_013.0013.D, from x, y = 3.933, 16281 to 4.043, 16380, result = 33576; previous integration is from x, y = 3.926, 15563 to 4.050, 15921 and previous response = 37796.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:47:17 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_013.0013.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:47:21 AM	Set SampleApproved = True for sample G0225_013.0013.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:47:32 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_014.0014.D, from x, y = 3.929, 16125 to 4.058, 16036, result = 36006; previous integration is from x, y = 3.921, 15517 to 4.088, 14990 and previous response = 43740.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:47:34 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_014.0014.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:47:43 AM	Set SampleApproved = True for sample G0225_014.0014.D; previous value = False			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:47:49 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_015.0015.D, from x, y = 3.928, 16193 to 4.054, 16146, result = 33895; previous integration is from x, y = 3.917, 15432 to 4.096, 14929 and previous response = 43342.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:47:50 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_015.0015.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:47:54 AM	Set SampleApproved = True for sample G0225_015.0015.D; previous value = False			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:48:04 AM	Manually integrate compound 1,1,1,2-Tetrachloroethane in sample G0225_016.0016.D, from x, y = 3.928, 16109 to 4.055, 16167, result = 35520; previous integration is from x, y = 3.915, 15430 to 4.092, 14963 and previous response = 44700.			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:48:06 AM	Set UserAnnotation = GT for compound 1,1,1,2-Tetrachloroethane in sample G0225_016.0016.D; previous value =			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:48:11 AM	Set SampleApproved = True for sample G0225_016.0016.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:48:12 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_017.0017.D			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:48:14 AM	Zero out primary peak of compound 1,1,1,2-Tetrachloroethane in sample G0225_017.0017.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:48:16 AM	Set SampleApproved = True for sample G0225_017.0017.D; previous value = False			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:48:37 AM	Set SampleApproved = True for sample G0225_018.0018.D; previous value = False			✓	
CmdZeroOutPeak	BL2000\ctran	2/28/2022 7:48:42 AM	Zero out primary peak of compound 1,2-Dibromoethane in sample G0225_019.0019.D			✓	
CmdSetSampleAttribute	BL2000\ctran	2/28/2022 7:48:44 AM	Set SampleApproved = True for sample G0225_019.0019.D; previous value = False			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:50:51 AM	Clear manual integration of target signal for compound 1,2-Dibromoethane in sample G0225_023.0023.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:50:51 AM	Set UserAnnotation = for compound 1,2-Dibromoethane in sample G0225_023.0023.D; previous value = GT			✓	
CmdManuallyIntegratePeak	BL2000\ctran	2/28/2022 7:50:57 AM	Manually integrate compound 1,2-Dibromoethane in sample G0225_023.0023.D, from x, y = 3.363, 15856 to 3.528, 15854, result = 19896; previous integration is from x, y = 3.363, 15856 to 3.557, 15714 and previous response = 20595.			✓	
CmdSaveBatchTable	BL2000\ctran	2/28/2022 7:51:20 AM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	

Audit Trail report

Name	User	Time	Action	Reason	Comment	Succeed	Exception
CmdOpenBatchTable	BL2000\ctran	2/28/2022 7:51:55 AM	Open batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\G022522_8011_W_CLT.batch.bin			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:52:02 AM	Clear manual integration of target signal for compound 1,2-Dibromoethane in sample G0225_023.0023.D			✓	
CmdClearManualIntegration	BL2000\ctran	2/28/2022 7:53:31 AM	Clear manual integration of target signal for compound 1,2-Dibromoethane in sample G0225_002(1).0002.D			✓	
CmdSetTargetCompoundAttribute	BL2000\ctran	2/28/2022 7:53:31 AM	Set UserAnnotation = for compound 1,2-Dibromoethane in sample G0225_002(1).0002.D; previous value = GT			✓	
CmdSaveBatchTable	BL2000\ctran	2/28/2022 8:34:39 AM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	
CmdQuantitate	BL2000\ctran	2/28/2022 8:34:47 AM	Quantitate all compounds in all samples			✓	
CmdSaveBatchTable	BL2000\ctran	2/28/2022 8:34:49 AM	Save batch \\MASSHUNTER\Org\Data\GECD.I\G022522\aiexport\QuantResults\G022522_8011_W_CLT.batch.bin			✓	
CmdOpenBatchTable	BL2000\srcox	3/9/2022 3:37:49 PM	Open batch D:\Org\Data\GECD.I\G022522\aiexport\G022522_8011_W_CLT.batch.bin			✓	
GenerateReport	BL2000\srcox	3/9/2022 3:46:37 PM	Generates report - Method: \\MASSHUNTER\Org\reports\Env_QuantResults_wGraphics+Chromatogram.m, Output Path: D:\Org\Data\GECD.I\G022522\aiexport\QuantReports\G022522_8011_W_CLT			✓	



ID #: 13327

Opened:

Calibration Standard

Expires: 12/31/2023

Rec'd: 12/11/2020

Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Certificate of Analysis

Product Name: Calibration Standard

Product Number: DWM-514-1

Lot Issue Date: 08-Dec-2020

Lot Number: 0006573696

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
1,2-dibromo-3-chloropropane	000096-12-8	RM12895	200.7 ± 1.0 µg/mL
1,2-dibromoethane	000106-93-4	RM00018	200.2 ± 1.0 µg/mL
1,2,3-trichloropropane	000096-18-4	RM13082	200.4 ± 1.0 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.



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



ISO 17025 Cert
No. AT-1937

Certificate of Analysis

Product Number: DWM-514-1

Lot Number: 0006573696

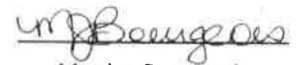
Expiration of Certification:

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



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



ISO 17025 Cert
No. AT-1937

Energy Laboratories Inc

Spike LOG

Standard ID: PH121120504P
Standard Name: 504.1 Mix (200ug/mL) MeOH
Date Prepared: 12/11/2019
Date Expires: 12/31/2023
Department: PST/HRBPR
Vendor: Agilent
Lot Number: 0006573696
Balance ID:

Type: Primary
BY: Selina R. Cox
Status: New

Comments: Date prepped is same as date received. [200ug/mL] MeOH. Recieved x4 1mL vials.

Chemical / Solvent Used	BottleNo	Amt	Units	Exp
Calibration Standard	13327	4	mL	12/31

Final Volume: 1 mL

Stock Source

Base Units

Amount Added

Analvtes

CAS

Conc: **ug/mL**

Energy Laboratories Inc

Standard LOG

Standard ID: PH011122504C3
Standard Name: 504.1 Cal Stock 3(0.7ug/mL) MeOH
Date Prepared: 1/11/2022
Date Expires: 2/12/2023
Department: PST/HRBPR
Vendor:
Lot Number:
Balance ID:

Type: Secondary
BY: Carry L Tran
Status: New

Comments: Final concentration = 0.7ug/mL Vol Flask# - EX-0117. Concentration represents both calmix and surrogate. 4/27/21 SRC.

Chemical / Solvent Used	BottleNo	Amt	Units	Exp
Methanol, Purge and Trap EA899	13926	9.895	mL	2/12/
1, 1, 1, 2-Tetrachloroethane Standard	14248	0.07	mL	11/30

Final Volume: 10 mL

Stock Source

PH121120504P 504.1 Mix (200ug/mL) MeOH

Base Units

ug/mL

Amount Added

0.035 mL

Analvtes

CAS

Conc: ug/mL

Energy Laboratories Inc

Standard LOG

Standard ID: PH011122504C2
Standard Name: 504.1 Cal Stock 2(0.07ug/mL) MeOH
Date Prepared: 1/11/2022
Date Expires: 2/12/2023
Department: PST/HRBPR
Vendor:
Lot Number:
Balance ID:
Type: Tertiary
BY: Carry L Tran
Status: New
Comments: Final concentration = 0.07ug/mL Vol Flask# - EX-0117

Chemical / Solvent Used	BottleNo	Amt	Units	Exp
Methanol, Purge and Trap EA899	13926	9	mL	2/12/

Final Volume: 10 mL

Stock Source
PH011122504C3 504.1 Cal Stock 3(0.7ug/mL) MeOH

Base Units
ug/mL

Amount Added
1 mL

Analvtes

CAS

Conc: **ug/mL**

Energy Laboratories Inc

Standard LOG

Standard ID: PH011122504C1
Standard Name: 504.1 Cal Stock 1(0.007ug/mL) MeOH
Date Prepared: 1/11/2022
Date Expires: 2/12/2023
Department: PST/HRBPR
Vendor:
Lot Number:
Balance ID:

Type: Tertiary
BY: Carry L Tran

Status: New

Comments: Final concentration = 0.007ug/mL Vol Flask# - EX-0117

Chemical / Solvent Used	BottleNo	Amt	Units	Exp
Methanol, Purge and Trap EA899	13926	9	mL	2/12/

Final Volume: 10 mL

Stock Source
PH011122504C2 504.1 Cal Stock 2(0.07ug/mL) MeOH

Base Units
ug/mL

Amount Added
1 mL

Analvtes

CAS

Conc: **ug/mL**



Certificate of Analysis

ID #: 14248

Opened: _____

1, 1, 1, 2-Tetrachloroethane Standard

Expires: 11/30/2024

Rec'd: 9/7/2021

Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107

Product Name: 1,1,1,2-Tetrachloroethane Standard

Product Number: HC-410-1

Lot Issue Date: 27-Oct-2020

Lot Number: 0006567948

Expiration Date: 30-Nov-2024

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
1,1,1,2-tetrachloroethane	000630-20-6	RM12632	99.9 ± 0.5 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/NCCL 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



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

www.agilent.com/quality/
CSD-QA-015.1



ISO 17025 Cert
No. AT-1937

CERTIFICATE OF ANALYSIS

Catalog No: M-504.1-LFB
Description: Laboratory Fortified Blank Sample Concentrate
Lot: 220021015
Solvent: Methanol
Hazards: Refer to SDS for complete safety information

Date Certified: Feb 6, 2020
Expiration: Feb 6, 2023
Sample Size: 1 mL
Components: 3
Storage Condition: Ambient (>5 °C)



Signal Word: Danger

Certified Reference Material



Component	CAS #	Purity % (GC/FID)	Prepared Concentration ² (µg/mL)	Certified Analyte Concentration ¹ (µg/mL)
1,2-Dibromoethane	106-93-4	99.9	0.2503	0.2500
1,2-Dibromo-3-chloropropane	96-12-8	100.0	0.2505	0.2505
1,2,3-Trichloropropane	96-18-4	99.0	0.2503	0.2478

ID #: 14729

Opened:

Laboratory Fortified Blank Sample Concentrate
Expires: 2/6/2023

Rec'd: 1/6/2022

Energy Laboratories Inc 1120 Sp. 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. Matrix blank to be used for background correction.

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.

The information on this certificate may not be reproduced without the express permission of the manufacturer. 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

ISO 17034

ID #: 14328

Opened:

1,1,1,2-Tetrachloroethane Solution

Expires: 10/31/2025

Rec'd: 9/29/2021

Energy Laboratories Inc 1120 So. 27th Street
Billings MT 59107



Agilent

Trusted Answers

Reference Material Certificate

Product Name: 1,1,1,2-Tetrachloroethane Standard

Lot Number: 0006633034

Product Number: HC-410-1

Lot Issue Date: 27-Sep-2021

Storage Conditions: Store Frozen (-25° to -10°C).

Expiration Date: 31-Oct-2025

Component Name	CERTIFIED VALUES		CAS#	Analyte Lot
	Concentration	Expanded Uncertainty		
1,1,1,2-tetrachloroethane	100.4	± 0.5 µg/mL	000630-20-6	RM12632

Matrix: methanol (methyl alcohol)

Description:

This document is prepared in accordance with ISO 17034 and Guide 31. This analytical reference material standard 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 above.

Traceability:

The balances used for these measurements are calibrated with weights traceable to NIST in compliance with ANSI/NCCL 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 analytical reference standard was unitized according to an in-house procedure and is guaranteed to be homogeneous. There is no minimum sub-sample size required.

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.

Safety:

Refer to the Safety Data Sheet on www.agilent.com for information regarding this analytical reference material.

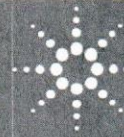
Intended Use:

This analytical reference standard 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.

Expiration of Certification:

The certification of this analytical reference standard is valid until the expiration date specified above, provided the material is handled and stored in accordance with the instructions given in this certificate. This certification is nullified if the material is damaged, contaminated, or otherwise modified.

ISO 17034



Agilent

Trusted Answers

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



ISO 17034 Cert
No. AR-1936

RM was produced in accordance with the TUV/SUD registered ISO 9001:2015
Quality Management System. Cert# 951215321

Page: 2 of 2

www.agilent.com/quality/
CSD-QA-015.1



ISO 17025 Cert
No. AT-1937

Energy Laboratories Inc

Standard LOG

Standard ID: PH021622504SU
Standard Name: 504.1 Surrogate (0.1ug/mL)MeOH
Date Prepared: 2/16/2022
Date Expires: 3/20/2023
Department: PST/HRBPR
Vendor:
Lot Number:
Balance ID:
Type: Secondary
BY: Carry L Tran
Status: New
Comments: Final Concentration = (0.1ug/mL) Vol Flask: EX-0117

Chemical / Solvent Used	BottleNo	Amt	Units	Exp
Methanol, Purge and Trap - EB679	14746	9.99	mL	5/28/
1,1,1,2-Tetrachloroethane Solution	14328	0.01	mL	10/31

Final Volume: 10 mL

Stock Source

Base Units

Amount Added

Analtes

CAS

Conc: **ug/mL**