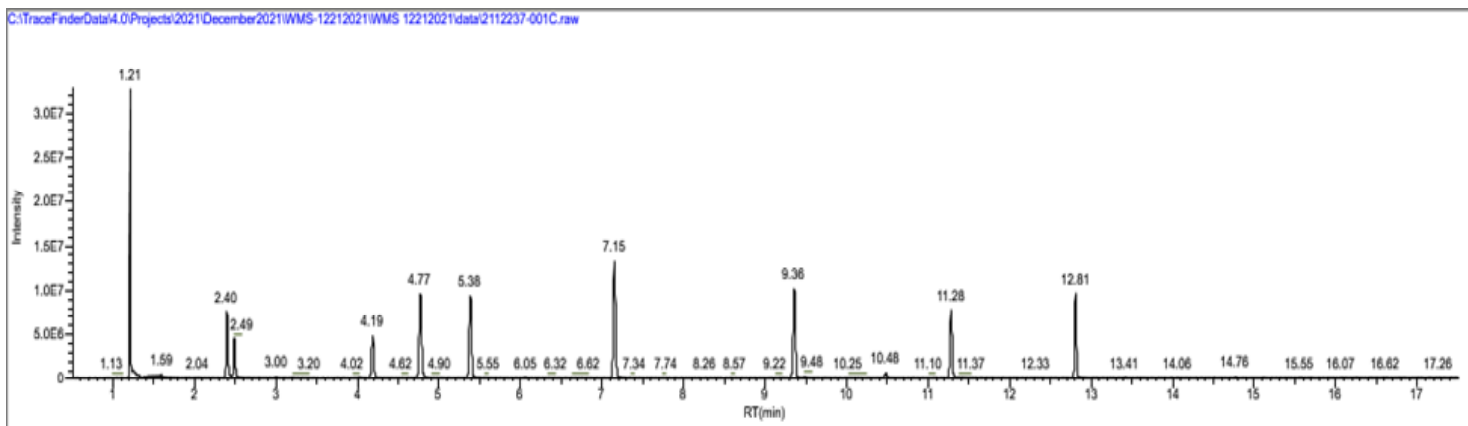


Batch Folder Path: C:\TraceFinderData\4.0\Projects\2021\December2021\WMS-12212021\WMS 12212021
Sample ID: 2112237-001C
User Name: BP
Method Name: WMS 12212021_Torrent Lab Method 8260-W_12142021-ICAL
Instrument Name: GCMS-10
Acquisition Date: 12/21/2021 2:43:26 PM
Cal File: C:\TraceFinderData\4.0\Projects\2021\December2021\WMS-12212021\WMS 12212021\Methods\8260B-ICAL-12142021_WMS 12212C



R Time	Name	Type	Conc.	Area	m/z	Conc.Units
4.78	Pentafluorobenzene	internalStandar	11.90	4518154	168	ug/L
5.38	1,4-Difluorobenzene	internalStandar	11.90	7906376	114	ug/L
9.36	Chlorobenzene-d5	internalStandar	11.90	6786287	117	ug/L
12.81	1,4-Dichlorobenzene-d4	internalStandar	11.90	2403909	152	ug/L
R Time	Name	Type	Conc	Total Area	m/z	Units
1.28	Dichlorodifluoromethane	targetCompou	0.00	187	85	ug/L
1.37	Chloromethane	targetCompou	0.03	8754	50	ug/L
1.42	1,3-Butadiene	targetCompou	0.01	1525	54	ug/L
1.41	Vinyl Chloride	targetCompou	0.01	1604	62	ug/L
1.57	Bromomethane	targetCompou	0.22	27411	94	ug/L
1.64	Chloroethane	targetCompou	0.01	853	64	ug/L
1.68	Trichlorofluoromethane	targetCompou	0.00	247	101	ug/L
2.05	1,1-Dichloroethene	targetCompou	0.00	313	61	ug/L
2.05	Carbon disulfide	targetCompou	0.02	16830	76	ug/L
1.97	Freon 113	targetCompou	0.00	153	151	ug/L
2.45	Methylene Chloride	targetCompou	0.05	14435	49	ug/L
2.49	Acetone	targetCompou	119.89	3996256	43	ug/L
2.58	trans-1,2-Dichloroethene	targetCompou	0.00	673	61	ug/L
2.74	MTBE	targetCompou	0.00	959	73	ug/L

2.77	tert-Butanol	TargetCompou	0.87	7945	59	ug/L
3.11	Diisopropyl ether (DIPE)	TargetCompou	0.00	2311	45	ug/L
3.05	1,1-Dichloroethane	TargetCompou	0.00	300	63	ug/L
3.43	ETBE	TargetCompou	0.00	302	59	ug/L
3.75	2,2-Dichloropropane	TargetCompou	0.00	615	77	ug/L
4.04	Bromochloromethane	TargetCompou	0.00	299	49	ug/L
3.99	Chloroform	TargetCompou	0.00	744	83	ug/L
4.18	(S) Dibromofluoromethane	eSurrogate	15.52	2971853	111	ug/L
4.18	1,1,1-Trichloroethane	TargetCompou	0.00	238	97	ug/L
4.35	1,1-Dichloropropene	TargetCompou	0.00	348	110	ug/L
4.32	2-Butanone	TargetCompou	0.25	2819	43	ug/L
4.61	Benzene	TargetCompou	0.01	8566	78	ug/L
4.76	(S) 1,4-Dichloroethane-d4	eSurrogate	13.94	2782235	65	ug/L
4.8	TAME	TargetCompou	0.01	4630	73	ug/L
4.85	1,2-Dichloroethane	TargetCompou	0.03	5592	62	ug/L
5.33	Trichloroethylene	TargetCompou	0.00	93	132	ug/L
5.82	Dibromomethane	TargetCompou	0.00	246	174	ug/L
5.94	1,2-Dichloropropane	TargetCompou	0.00	536	63	ug/L
6.07	Bromodichloromethane	TargetCompou	0.01	2023	83	ug/L
6.35	1,4-Dioxane	TargetCompou	0.93	888	88	ug/L
6.91	cis-1,3-Dichloropropene	TargetCompou	0.00	400	75	ug/L
7.15	(S) Toluene-d8	eSurrogate	11.77	10350365	98	ug/L
7.22	Toluene	TargetCompou	0.05	52403	91	ug/L
7.74	Tetrachloroethylene	TargetCompou	0.00	497	166	ug/L
7.8	4-Methyl-2-Pentanone	TargetCompou	0.02	2982	43	ug/L
7.86	trans-1,3-Dichloropropene	TargetCompou	0.00	522	75	ug/L
8.03	1,1,2-Trichloroethane	TargetCompou	0.01	1079	83	ug/L
8.28	Dibromochloromethane	TargetCompou	0.00	203	129	ug/L
8.45	1,3-Dichloropropane	TargetCompou	0.00	412	76	ug/L
8.58	1,2-Dibromoethane	TargetCompou	0.00	156	107	ug/L
9.06	2-Hexanone	TargetCompou	0.35	10358	43	ug/L
9.38	Chlorobenzene	TargetCompou	0.00	990	112	ug/L
9.5	Ethyl Benzene	TargetCompou	0.03	22536	91	ug/L
9.52	1,1,1,2-Tetrachloroethane	TargetCompou	0.01	971	133	ug/L
9.73	m,p-Xylene	TargetCompou	0.01	7647	91	ug/L
10.38	o-Xylene	TargetCompou	0.01	4480	91	ug/L
10.4	Bromoform	TargetCompou	0.00	194	173	ug/L
10.48	Styrene	TargetCompou	0.79	321548	104	ug/L
10.92	Isopropyl Benzene	TargetCompou	0.26	4439	105	ug/L
11.28	(S) 4-Bromofluorobenzene	eSurrogate	12.84	3129431	95	ug/L
11.38	Bromobenzene	TargetCompou	0.00	1181	77	ug/L
11.55	n-Propylbenzene	TargetCompou	0.01	4094	91	ug/L
11.7	1,1,2,2-Tetrachloroethane	TargetCompou	0.00	556	83	ug/L
11.7	2-Chlorotoluene	TargetCompou	0.00	1452	91	ug/L
11.78	1,2,3-Trichloropropane	TargetCompou	0.01	536	75	ug/L
11.87	1,3,5-Trimethylbenzene	TargetCompou	0.28	2235	105	ug/L

Flag legend: LOD<J<LOQ; I=Ion ratio failure; C=Carryover; ?=Linearity limit; D=Detection limit; Q=Quan limit; POS=Rpt limit; b=Blank; s=Solvent blank; est=Estimated; Bordered Cell=Manually integrated.

11.94	4-Chlorotoluene	TargetCompou	0.00	862	91	ug/L
12.21	tert-Butylbenzene	TargetCompou	0.22	369	119	ug/L
12.37	1,2,4-Trimethylbenzene	TargetCompou	0.27	939	105	ug/L
12.51	sec-Butyl Benzene	TargetCompou	0.00	1595	105	ug/L
12.72	1,3-Dichlorobenzene	TargetCompou	0.00	1348	146	ug/L
12.73	p-Isopropyltoluene	TargetCompou	0.28	596	119	ug/L
12.82	1,4-Dichlorobenzene	TargetCompou	0.01	1843	146	ug/L
13.25	n-Butylbenzene	TargetCompou	0.26	1958	91	ug/L
13.3	Hexachloroethane	TargetCompou	0.00	195	166	ug/L
13.36	1,2-Dichlorobenzene	TargetCompou	0.00	1261	146	ug/L
14.29	1,2-Dibromo-3-Chloropropane	TargetCompou	0.04	828	75	ug/L
15.09	1,2,4-Trichlorobenzene	TargetCompou	0.01	1762	180	ug/L
15.11	Hexachlorobutadiene	TargetCompou	0.02	1686	225	ug/L
15.45	Naphthalene	TargetCompou	0.03	6471	128	ug/L
15.66	1,2,3-Trichlorobenzene	TargetCompou	0.02	1827	180	ug/L
N/F	cis-1,2-Dichloroethene	TargetCompou	0.00	N/F	61	ug/L
N/F	Carbon Tetrachloride	TargetCompou	0.00	N/F	119	ug/L

Flag legend: LOD<J<LOQ; I=Ion ratio failure; C=Carryover; ?=Linearity limit; D=Detection limit; Q=Quan limit; POS=Rpt limit; b=Blank; s=Solvent blank; est=Estimated; Bordered Cell=Manually integrated.