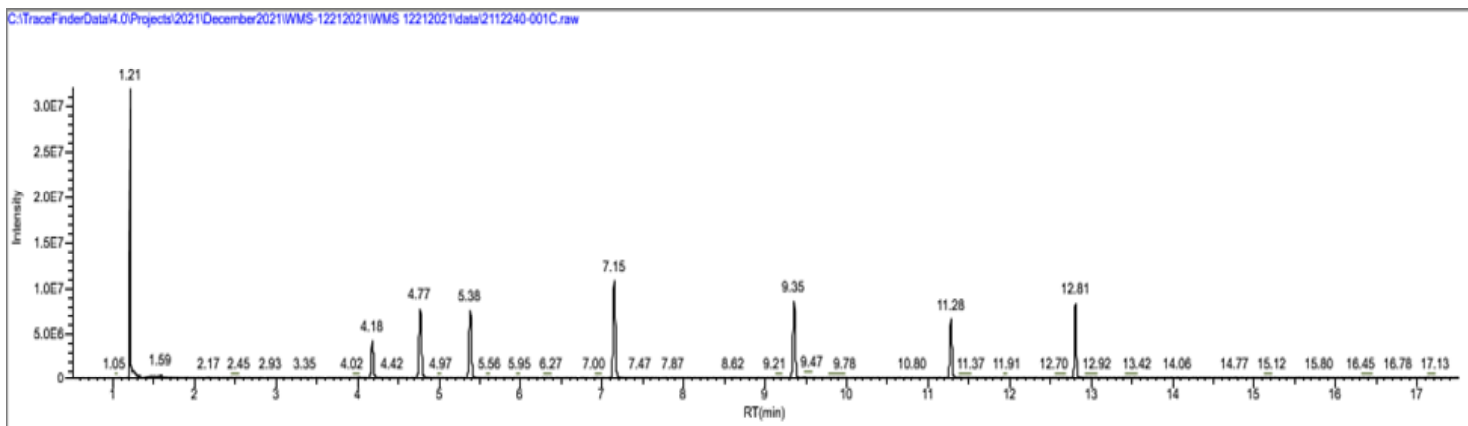


Batch Folder Path: C:\TraceFinderData\4.0\Projects\2021\December2021\WMS-12212021\WMS 12212021
Sample ID: 2112240-001C
User Name: BP
Method Name: WMS 12212021_Torrent Lab Method 8260-W_12142021-ICAL
Instrument Name: GCMS-10
Acquisition Date: 12/21/2021 3:13:24 PM
Cal File: C:\TraceFinderData\4.0\Projects\2021\December2021\WMS-12212021\WMS 12212021\Methods\8260B-ICAL-12142021_WMS 122120



R Time	Name	Type	Conc.	Area	m/z	Conc.Units
4.77	Pentafluorobenzene	internalStandar	11.90	3495838	168	ug/L
5.38	1,4-Difluorobenzene	internalStandar	11.90	6206945	114	ug/L
9.35	Chlorobenzene-d5	internalStandar	11.90	5758629	117	ug/L
12.81	1,4-Dichlorobenzene-d4	internalStandar	11.90	2068523	152	ug/L
R Time	Name	Type	Conc	Total Area	m/z	Units
1.28	Dichlorodifluoromethane	targetCompou	0.00	208	85	ug/L
1.37	Chloromethane	targetCompou	0.03	8194	50	ug/L
1.41	1,3-Butadiene	targetCompou	0.00	652	54	ug/L
1.41	Vinyl Chloride	targetCompou	0.01	1663	62	ug/L
1.57	Bromomethane	targetCompou	0.26	24923	94	ug/L
1.67	Chloroethane	targetCompou	0.00	222	64	ug/L
1.73	Trichlorofluoromethane	targetCompou	0.00	373	101	ug/L
2.03	1,1-Dichloroethene	targetCompou	0.00	497	61	ug/L
2.05	Carbon disulfide	targetCompou	0.02	9186	76	ug/L
2.08	Freon 113	targetCompou	0.00	112	151	ug/L
2.45	Methylene Chloride	targetCompou	0.05	10864	49	ug/L
2.5	Acetone	targetCompou	0.67	17230	43	ug/L
2.58	trans-1,2-Dichloroethene	targetCompou	0.00	363	61	ug/L
2.73	MTBE	targetCompou	0.00	1234	73	ug/L

2.84	tert-Butanol	TargetCompou	0.07	522	59	ug/L
3.04	Diisopropyl ether (DIPE)	TargetCompou	0.00	965	45	ug/L
3.18	1,1-Dichloroethane	TargetCompou	0.00	661	63	ug/L
3.44	ETBE	TargetCompou	0.00	944	59	ug/L
3.7	2,2-Dichloropropane	TargetCompou	0.00	799	77	ug/L
3.75	Bromochloromethane	TargetCompou	0.00	503	49	ug/L
3.99	Chloroform	TargetCompou	0.01	1708	83	ug/L
4.18	(S) Dibromofluoromethane	eSurrogate	16.93	2507964	111	ug/L
4.19	1,1,1-Trichloroethane	TargetCompou	0.01	1074	97	ug/L
4.27	1,1-Dichloropropene	TargetCompou	0.00	123	110	ug/L
4.37	2-Butanone	TargetCompou	0.30	4848	43	ug/L
4.61	Benzene	TargetCompou	0.01	3652	78	ug/L
4.76	(S) 1,4-Dichloroethane-d4	eSurrogate	14.82	2322187	65	ug/L
4.76	TAME	TargetCompou	0.02	5844	73	ug/L
4.84	1,2-Dichloroethane	TargetCompou	0.05	8302	62	ug/L
5.32	Trichloroethylene	TargetCompou	0.00	431	132	ug/L
5.83	Dibromomethane	TargetCompou	0.00	366	174	ug/L
5.78	1,2-Dichloropropane	TargetCompou	0.01	2437	63	ug/L
6.06	Bromodichloromethane	TargetCompou	0.00	662	83	ug/L
6.89	cis-1,3-Dichloropropene	TargetCompou	0.00	124	75	ug/L
7.15	(S) Toluene-d8	eSurrogate	11.28	8414081	98	ug/L
7.21	Toluene	TargetCompou	0.00	2409	91	ug/L
7.72	Tetrachloroethylene	TargetCompou	0.00	627	166	ug/L
7.81	4-Methyl-2-Pentanone	TargetCompou	0.02	1870	43	ug/L
7.81	trans-1,3-Dichloropropene	TargetCompou	0.00	148	75	ug/L
8	1,1,2-Trichloroethane	TargetCompou	0.00	324	83	ug/L
8.27	Dibromochloromethane	TargetCompou	0.00	392	129	ug/L
8.63	1,3-Dichloropropane	TargetCompou	0.00	341	76	ug/L
8.55	1,2-Dibromoethane	TargetCompou	0.00	105	107	ug/L
9.04	2-Hexanone	TargetCompou	0.36	9418	43	ug/L
9.38	Chlorobenzene	TargetCompou	0.00	1009	112	ug/L
9.49	Ethyl Benzene	TargetCompou	0.00	1084	91	ug/L
9.49	1,1,1,2-Tetrachloroethane	TargetCompou	0.00	275	133	ug/L
9.73	m,p-Xylene	TargetCompou	0.00	2123	91	ug/L
10.41	o-Xylene	TargetCompou	0.01	3053	91	ug/L
10.39	Bromoform	TargetCompou	0.00	64	173	ug/L
10.49	Styrene	TargetCompou	0.29	3868	104	ug/L
10.92	Isopropyl Benzene	TargetCompou	0.26	1084	105	ug/L
11.28	(S) 4-Bromofluorobenzene	eSurrogate	12.40	2600609	95	ug/L
11.38	Bromobenzene	TargetCompou	0.00	1160	77	ug/L
11.54	n-Propylbenzene	TargetCompou	0.00	1221	91	ug/L
11.66	1,1,2,2-Tetrachloroethane	TargetCompou	0.01	795	83	ug/L
11.68	2-Chlorotoluene	TargetCompou	0.00	473	91	ug/L
11.77	1,2,3-Trichloropropane	TargetCompou	0.00	418	75	ug/L
11.86	1,3,5-Trimethylbenzene	TargetCompou	0.28	766	105	ug/L
11.91	4-Chlorotoluene	TargetCompou	0.00	801	91	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.

12.28	tert-Butylbenzene	TargetCompou	0.22	1260	119	ug/L
12.35	1,2,4-Trimethylbenzene	TargetCompou	0.28	2161	105	ug/L
12.5	sec-Butyl Benzene	TargetCompou	0.00	266	105	ug/L
12.72	1,3-Dichlorobenzene	TargetCompou	0.00	1100	146	ug/L
12.72	p-Isopropyltoluene	TargetCompou	0.28	228	119	ug/L
12.8	1,4-Dichlorobenzene	TargetCompou	0.01	1400	146	ug/L
13.25	n-Butylbenzene	TargetCompou	0.26	747	91	ug/L
13.3	Hexachloroethane	TargetCompou	0.01	501	166	ug/L
13.32	1,2-Dichlorobenzene	TargetCompou	0.00	740	146	ug/L
14.32	1,2-Dibromo-3-Chloropropane	TargetCompou	0.01	200	75	ug/L
15.08	1,2,4-Trichlorobenzene	TargetCompou	0.01	1150	180	ug/L
15.13	Hexachlorobutadiene	TargetCompou	0.01	602	225	ug/L
15.46	Naphthalene	TargetCompou	0.02	2752	128	ug/L
15.66	1,2,3-Trichlorobenzene	TargetCompou	0.02	1547	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
N/F	1,4-Dioxane	TargetCompou	0.00	N/F	88	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.