

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF041819\
 Data File : VF062213.D
 Acq On : 18 Apr 2019 19:06
 Operator : FY/SY
 Sample : K2193-07
 Misc : 6.15g/5mL/MSVOA F/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 HR-05-041719-A

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82F041719S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.229	55	71	72	rBV5	28909	165204	4.63%	0.285%
2	4.215	367	374	377	rBV	23820	83002	2.33%	0.143%
3	4.304	377	383	394	rVB	258848	785039	22.01%	1.352%
4	5.043	449	458	467	rBV2	485795	1577367	44.22%	2.717%
5	5.773	525	532	543	rBV	494930	1300750	36.46%	2.241%
6	7.705	723	728	735	rBV	592328	1397557	39.18%	2.408%
7	8.760	830	835	841	rBV7	12418	39281	1.10%	0.068%
8	9.105	865	870	874	rBV3	14072	37936	1.06%	0.065%
9	9.253	880	885	890	rVB4	15459	39961	1.12%	0.069%
10	9.420	897	902	906	rVB	18025	49688	1.39%	0.086%
11	9.676	919	928	934	rBV	10770	46862	1.31%	0.081%
12	9.913	946	952	959	rVV	541725	1295269	36.31%	2.231%
13	10.012	959	962	969	rVB2	83668	206533	5.79%	0.356%
14	10.258	983	987	991	rVB	45812	85413	2.39%	0.147%
15	10.357	991	997	1001	rBV	55326	138436	3.88%	0.238%
16	10.593	1018	1021	1022	rBV2	22507	41490	1.16%	0.071%
17	10.633	1022	1025	1032	rVB4	46357	114155	3.20%	0.197%
18	10.771	1032	1039	1041	rBV2	146593	309321	8.67%	0.533%
19	10.820	1041	1044	1047	rVB	212141	393413	11.03%	0.678%
20	10.889	1047	1051	1052	rBV2	20620	36664	1.03%	0.063%
21	10.928	1052	1055	1057	rVB	24399	37093	1.04%	0.064%
22	11.047	1063	1067	1070	rVB4	19816	47159	1.32%	0.081%
23	11.135	1070	1076	1078	rBV3	58854	172995	4.85%	0.298%
24	11.204	1080	1083	1085	rBV2	75104	115256	3.23%	0.199%
25	11.244	1085	1087	1090	rVB	135975	189207	5.30%	0.326%
26	11.352	1090	1098	1102	rVB3	233350	562514	15.77%	0.969%
27	11.421	1102	1105	1107	rBV3	18510	44384	1.24%	0.076%
28	11.500	1107	1113	1118	rBV	586576	1137531	31.89%	1.960%
29	11.569	1118	1120	1123	rVB3	44027	66020	1.85%	0.114%
30	11.628	1123	1126	1129	rBV	205368	363902	10.20%	0.627%
31	11.677	1129	1131	1133	rBV2	80774	99447	2.79%	0.171%
32	11.727	1133	1136	1139	rVB	681232	950211	26.64%	1.637%
33	11.796	1139	1143	1146	rBV	344953	732518	20.53%	1.262%
34	11.845	1146	1148	1150	rVB	60623	79401	2.23%	0.137%

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35	11.894	1150	1153	1156	rVB	608166	871439	24.43%	1.501%
36	11.953	1156	1159	1161	rBV	93814	144166	4.04%	0.248%
37	12.022	1161	1166	1167	rBV2	110644	221363	6.21%	0.381%
38	12.062	1167	1170	1171	rBV	327769	464854	13.03%	0.801%
39	12.091	1171	1173	1176	rVB	421450	527827	14.80%	0.909%
40	12.131	1176	1177	1180	rVB2	50951	62396	1.75%	0.107%
41	12.210	1182	1185	1188	rBV2	140241	287755	8.07%	0.496%
42	12.269	1188	1191	1194	rVB	644465	960347	26.92%	1.654%
43	12.358	1194	1200	1202	rBV5	99887	221112	6.20%	0.381%
44	12.436	1205	1208	1210	rBV	281685	449672	12.61%	0.775%
45	12.476	1210	1212	1214	rBV3	216175	356894	10.00%	0.615%
46	12.515	1214	1216	1221	rVB	390819	671332	18.82%	1.156%
47	12.614	1221	1226	1228	rBV2	737318	1464901	41.06%	2.524%
48	12.663	1228	1231	1235	rVB	1114842	1370730	38.42%	2.361%
49	12.722	1235	1237	1240	rVB3	61703	114856	3.22%	0.198%
50	12.781	1240	1243	1245	rBV2	656781	1118821	31.36%	1.927%
51	12.831	1245	1248	1249	rVV	416746	621471	17.42%	1.071%
52	12.860	1249	1251	1258	rVB3	1691114	3567354	100.00%	6.145%
53	12.988	1261	1264	1265	rBV2	210773	368603	10.33%	0.635%
54	13.057	1268	1271	1275	rVB	656877	1036267	29.05%	1.785%
55	13.146	1275	1280	1284	rBV3	475746	1587410	44.50%	2.735%
56	13.205	1284	1286	1289	rBV2	543761	941126	26.38%	1.621%
57	13.274	1292	1293	1297	rVB3	420846	723280	20.27%	1.246%
58	13.333	1297	1299	1302	rBV3	266550	435677	12.21%	0.751%
59	13.383	1302	1304	1306	rVB2	254109	341508	9.57%	0.588%
60	13.432	1306	1309	1310	rBV	325423	385670	10.81%	0.664%
61	13.462	1310	1312	1317	rVB2	528606	681308	19.10%	1.174%
62	13.531	1317	1319	1322	rVB2	440010	743371	20.84%	1.281%
63	13.580	1322	1324	1327	rBV	615220	993706	27.86%	1.712%
64	13.619	1327	1328	1330	rVB2	162284	167578	4.70%	0.289%
65	13.659	1330	1332	1333	rVB	69260	63063	1.77%	0.109%
66	13.678	1333	1334	1336	rVB	82525	91486	2.56%	0.158%
67	13.747	1338	1341	1345	rBV2	2115191	3190561	89.44%	5.496%
68	13.807	1345	1347	1350	rVB4	221846	357915	10.03%	0.617%
69	13.856	1350	1352	1353	rBV2	190692	270240	7.58%	0.466%
70	13.895	1353	1356	1360	rBV3	1599732	3068821	86.03%	5.287%
71	13.954	1360	1362	1365	rVB2	418548	645223	18.09%	1.112%

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72	14.014	1365	1368	1371	rBV	1398025	1886976	52.90%	3.251%
73	14.092	1373	1376	1377	rBV2	216837	347748	9.75%	0.599%
74	14.122	1377	1379	1381	rVB	272622	346837	9.72%	0.597%
75	14.191	1381	1386	1389	rBV6	768280	1841814	51.63%	3.173%
76	14.250	1389	1392	1396	rVB3	476159	900932	25.25%	1.552%
77	14.299	1396	1397	1399	rVB	117657	98286	2.76%	0.169%
78	14.339	1399	1401	1403	rVB	268494	334692	9.38%	0.577%
79	14.388	1403	1406	1409	rBV3	130380	209305	5.87%	0.361%
80	14.447	1409	1412	1414	rBV2	389623	600600	16.84%	1.035%
81	14.487	1414	1416	1419	rBV2	634461	886684	24.86%	1.527%
82	14.595	1425	1427	1428	rBV2	169506	204414	5.73%	0.352%
83	14.625	1428	1430	1434	rVB2	452298	666389	18.68%	1.148%
84	14.773	1442	1445	1448	rVB	1331612	1783817	50.00%	3.073%
85	14.861	1452	1454	1455	rVV	128419	185651	5.20%	0.320%
86	14.920	1455	1460	1464	rVV4	314154	966978	27.11%	1.666%
87	14.970	1464	1465	1468	rVB2	125737	158011	4.43%	0.272%
88	15.029	1468	1471	1476	rBV2	754789	1253962	35.15%	2.160%
89	15.137	1479	1482	1485	rBV	566607	911219	25.54%	1.570%
90	15.196	1485	1488	1491	rVV4	195544	371315	10.41%	0.640%
91	15.275	1494	1496	1497	rVV	62274	94956	2.66%	0.164%
92	15.305	1497	1499	1502	rVV3	119647	237351	6.65%	0.409%
93	15.374	1502	1506	1510	rVV2	208073	473460	13.27%	0.816%
94	15.443	1510	1513	1520	rVB2	72985	184759	5.18%	0.318%
95	15.630	1530	1532	1534	rBV3	34501	51433	1.44%	0.089%
96	15.670	1534	1536	1541	rVB	220738	418768	11.74%	0.721%
97	15.748	1541	1544	1547	rVB3	39363	72508	2.03%	0.125%
98	15.817	1548	1551	1553	rBV4	25613	48579	1.36%	0.084%
99	15.965	1563	1566	1569	rBV5	42701	91682	2.57%	0.158%
100	16.015	1569	1571	1576	rVB5	43478	88746	2.49%	0.153%

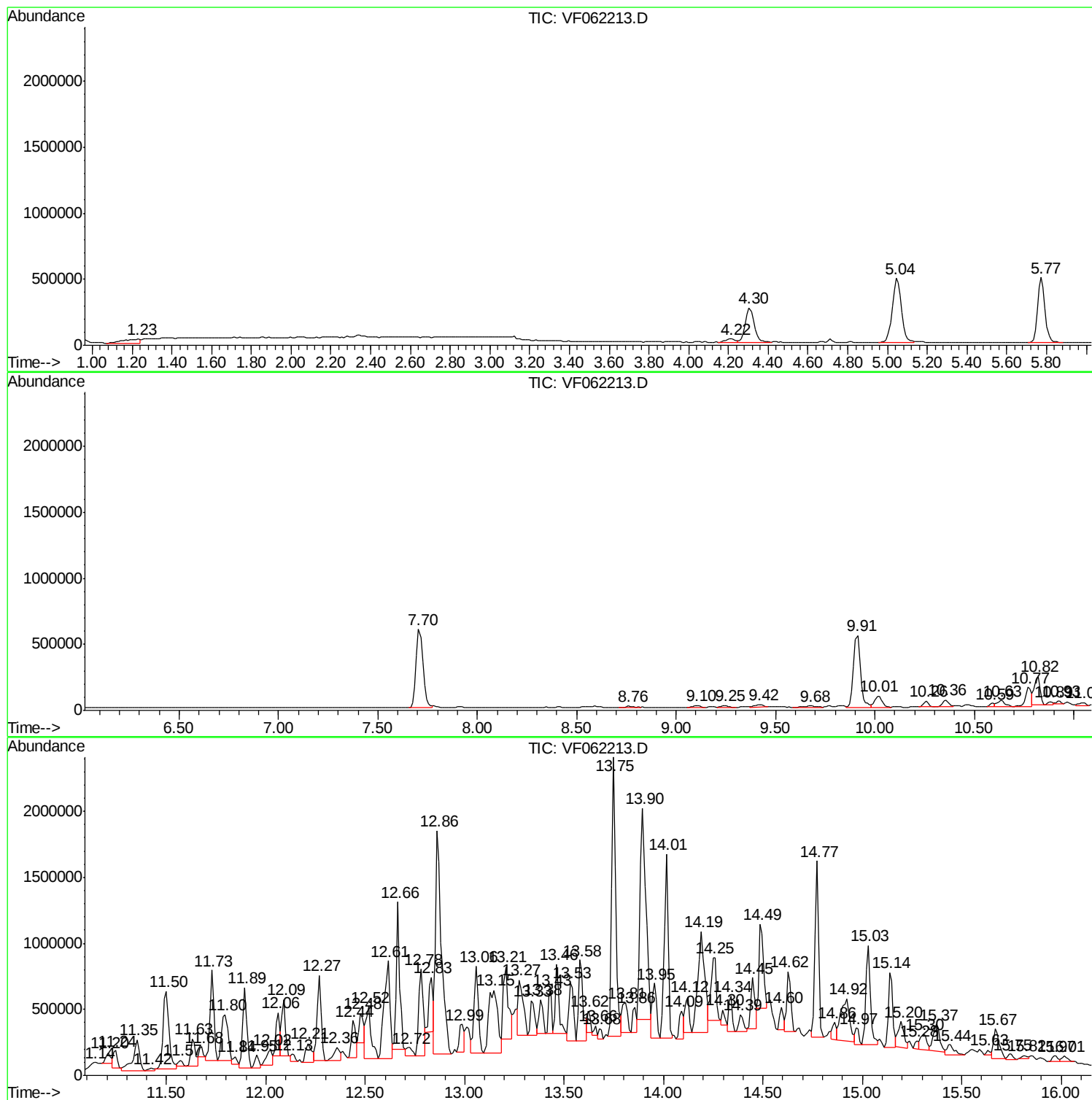
Sum of corrected areas: 58048954

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F041719S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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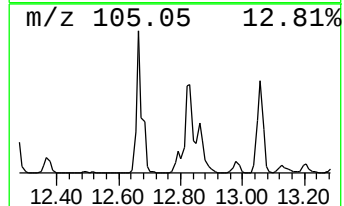
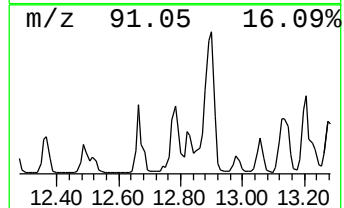
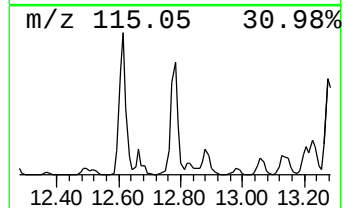
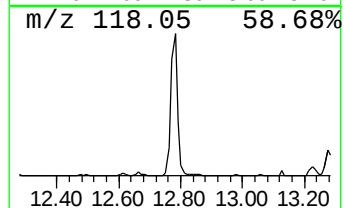
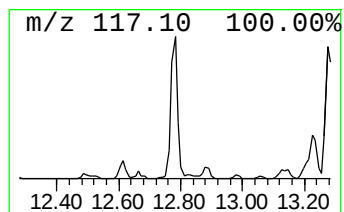
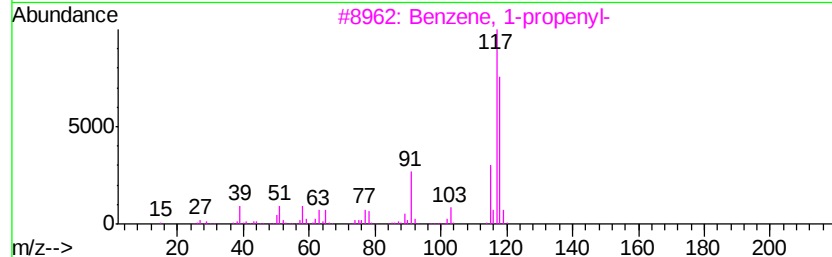
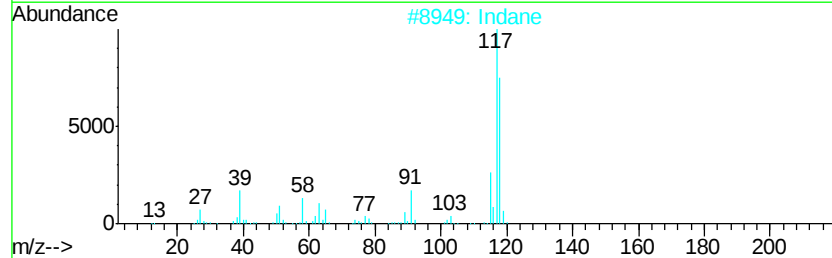
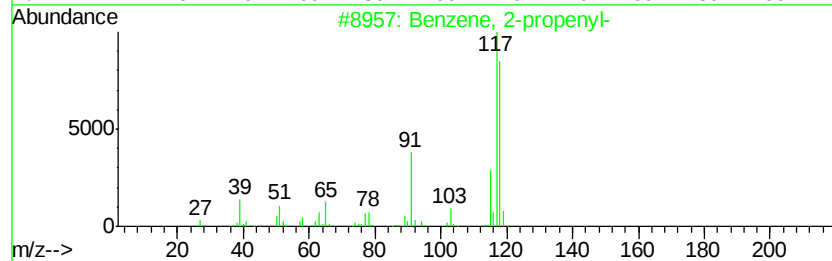
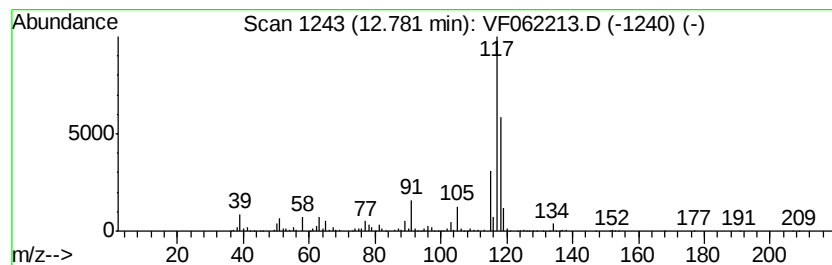
Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F041719S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 2-propenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.78	38.19 ug/l	1118820	1,4-Dichlorobenzene-d4	12.61
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-propenyl-	118 C9H10	000300-57-2 81
2		Indane	118 C9H10	000496-11-7 76
3		Benzene, 1-propenyl-	118 C9H10	000637-50-3 76
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 76
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 64



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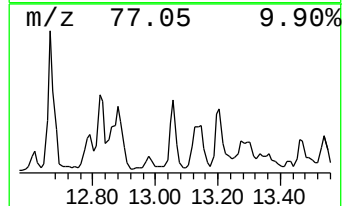
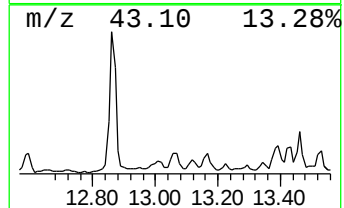
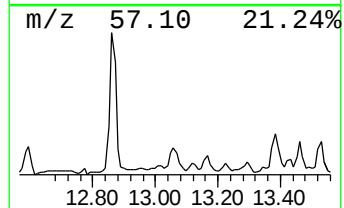
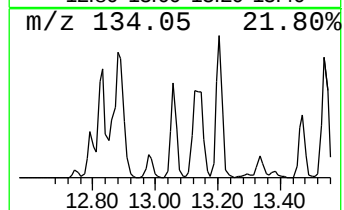
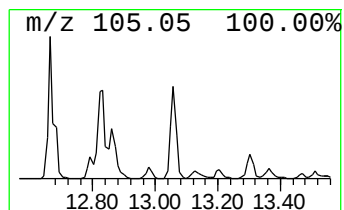
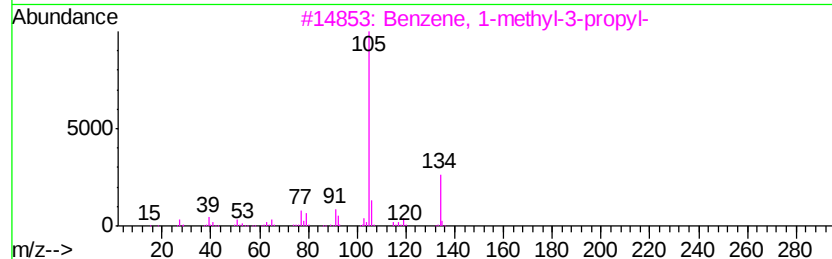
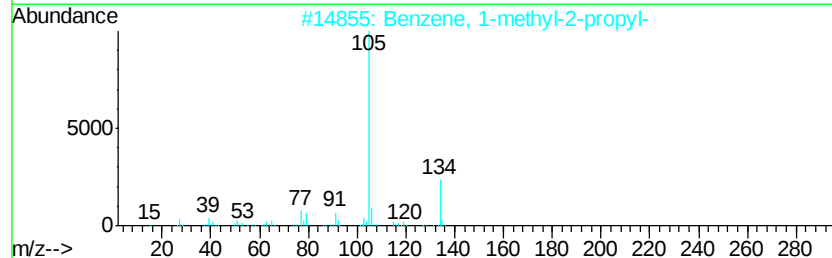
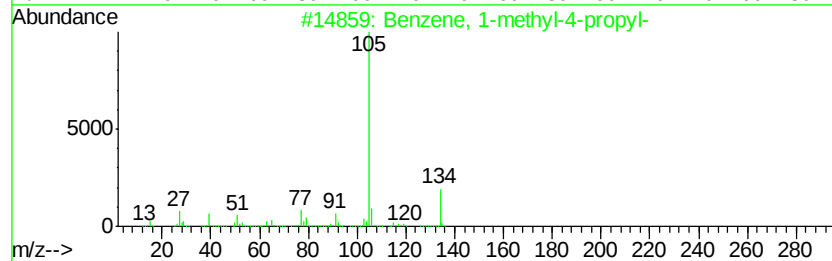
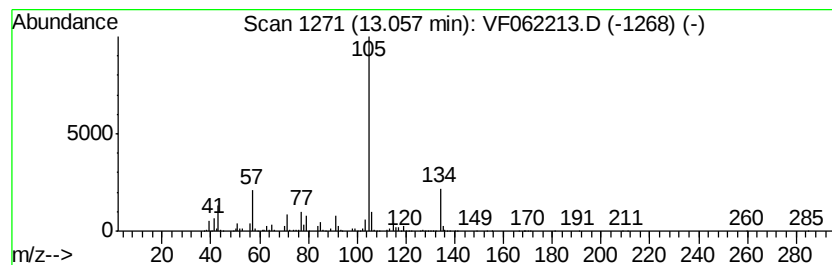
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Peak Number 2 Benzene, 1-methyl-4-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	35.37 ug/l	1036270	1,4-Dichlorobenzene-d4	12.61

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	90
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
5	Ethylenediamine, N,N'-bis(.alpha...	268	C18H24N2	006280-75-7	72



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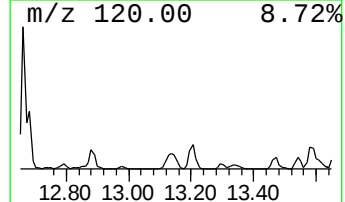
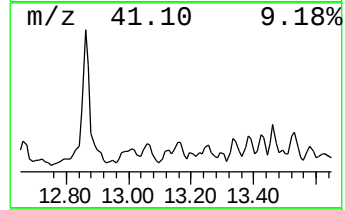
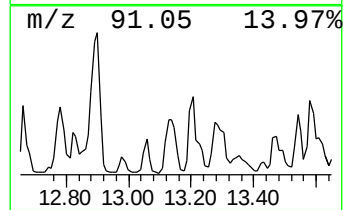
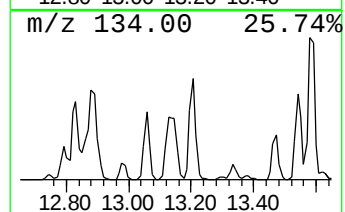
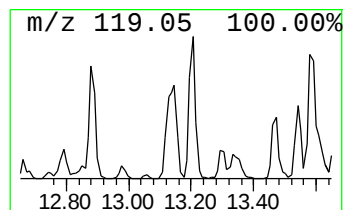
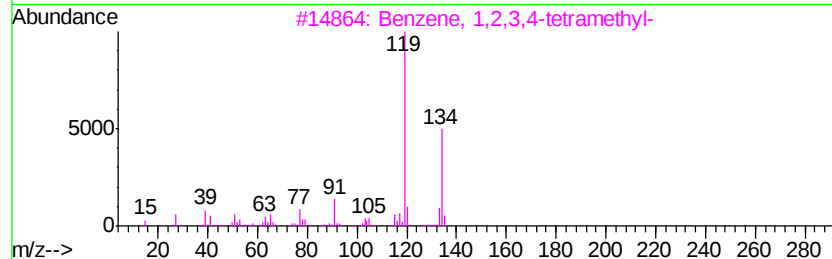
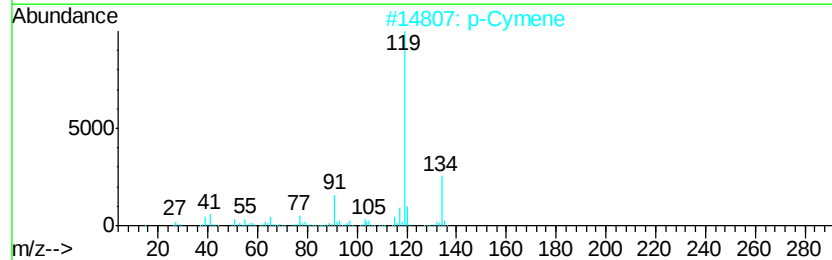
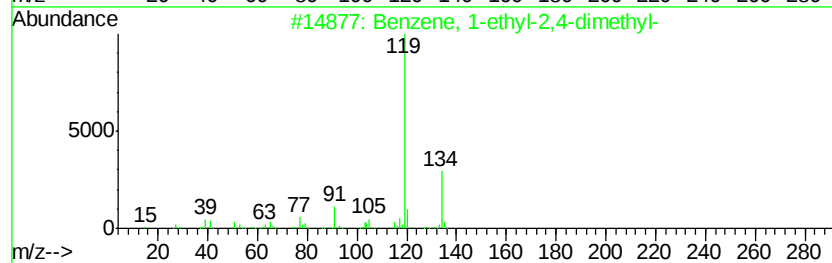
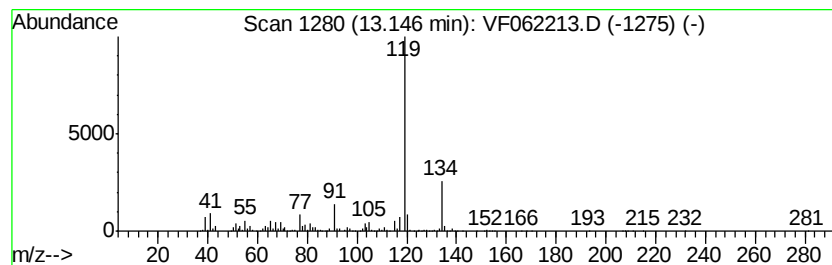
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	54.18 ug/l	1587410	1,4-Dichlorobenzene-d4	12.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
2		p-Cymene	134	C10H14	000099-87-6	91
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
4		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15a/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
HR-05-041719-A

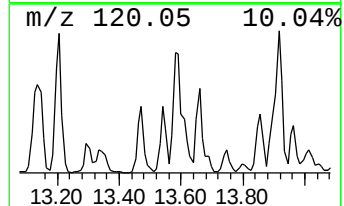
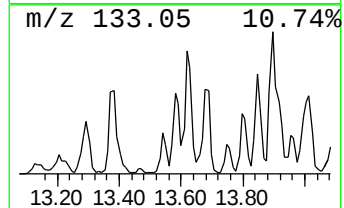
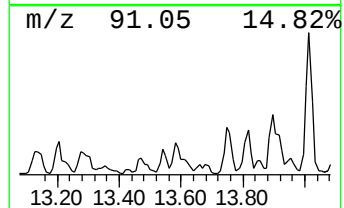
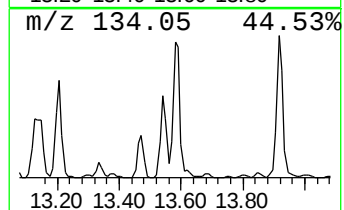
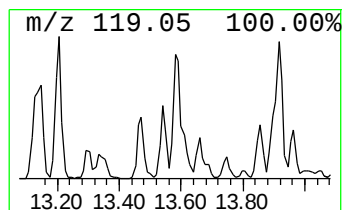
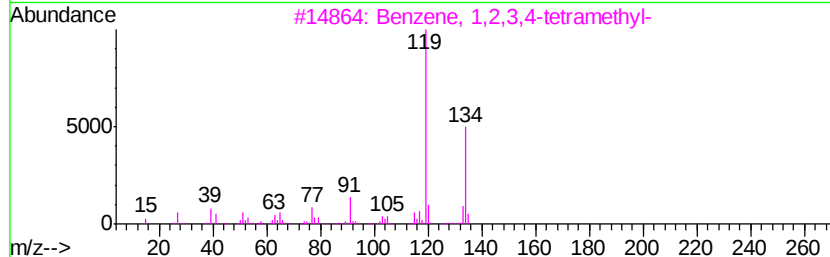
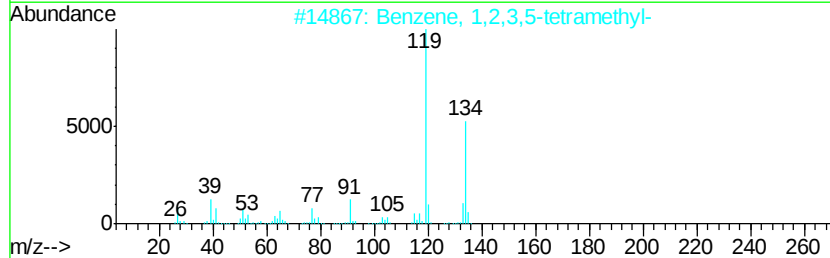
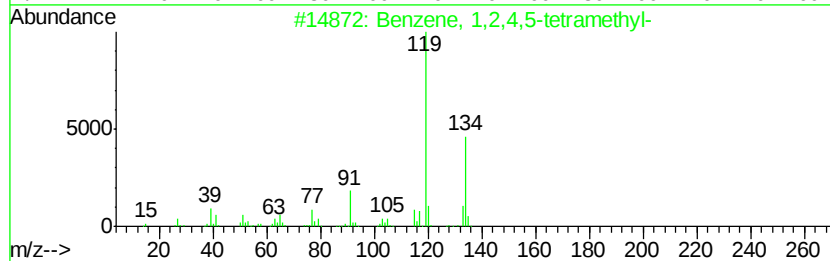
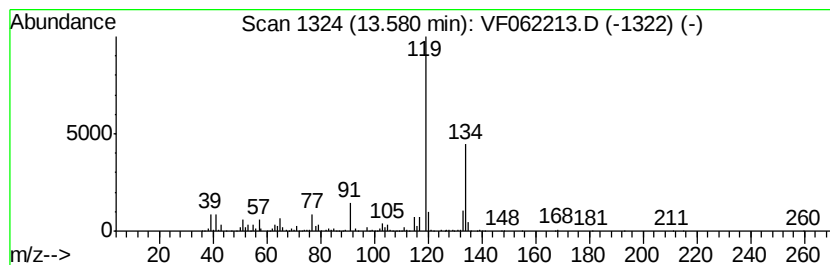
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.58	33.92 ug/l	993706	1,4-Dichlorobenzene-d4	12.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5		p-Cymene	134	C10H14	000099-87-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15a/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
HR-05-041719-A

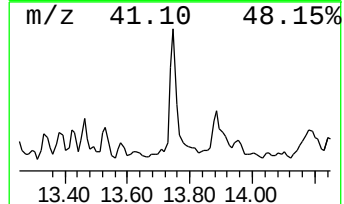
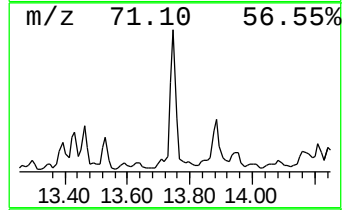
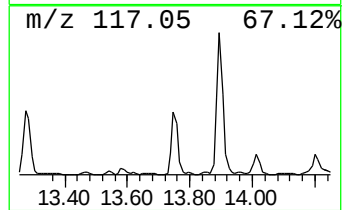
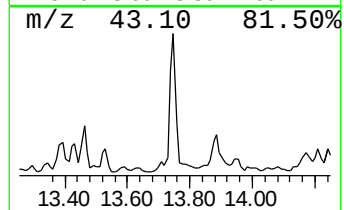
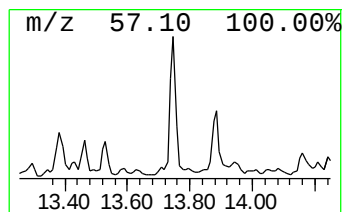
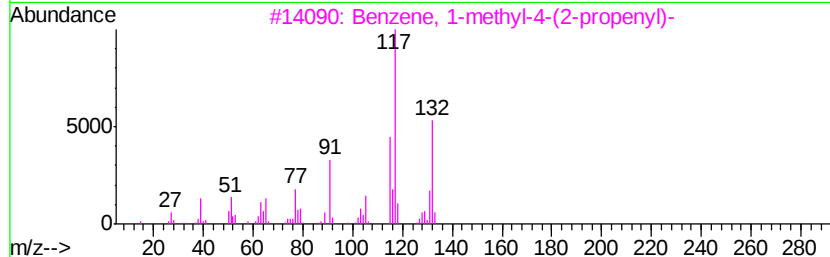
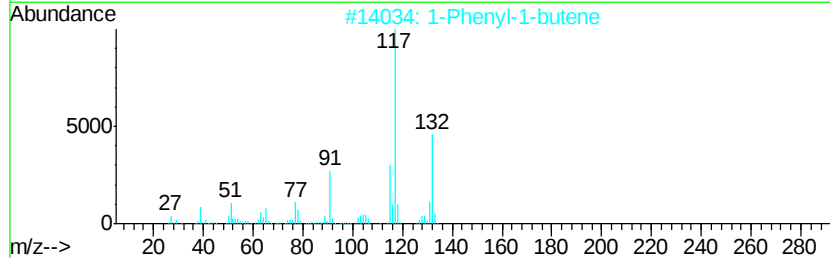
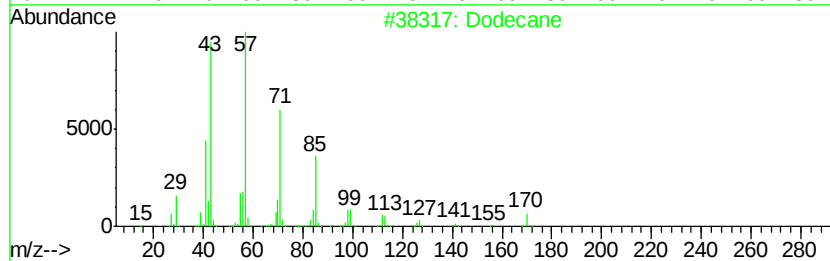
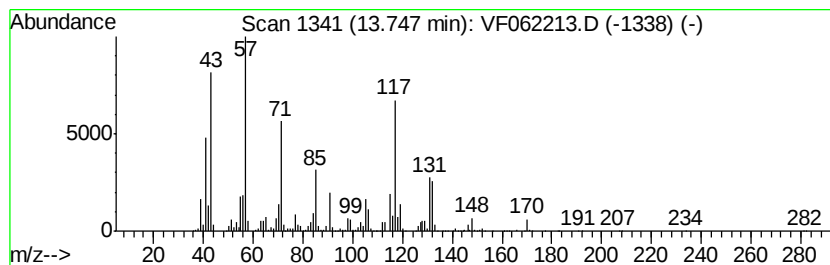
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Dodecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	108.90 ug/l	3190560	1,4-Dichlorobenzene-d4	12.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane		170	C12H26	000112-40-3	93
2	1-Phenyl-1-butene		132	C10H12	000824-90-8	80
3	Benzene, 1-methyl-4-(2-propenyl)-		132	C10H12	003333-13-9	42
4	Benzene, 1-methyl-2-(2-propenyl)-		132	C10H12	001587-04-8	42
5	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15a/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-05-041719-A

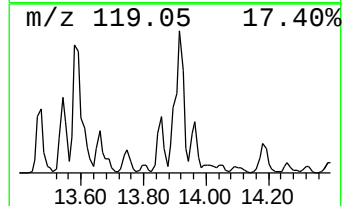
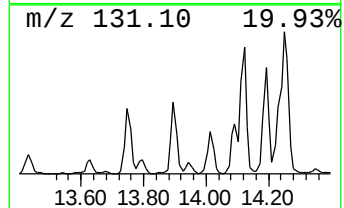
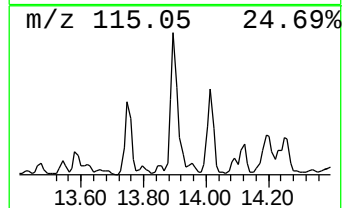
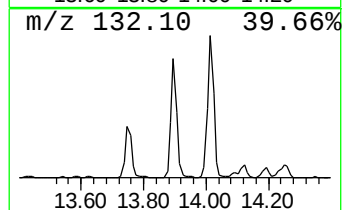
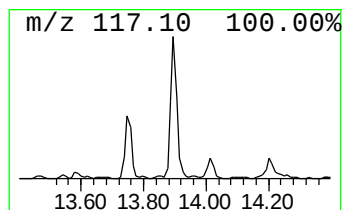
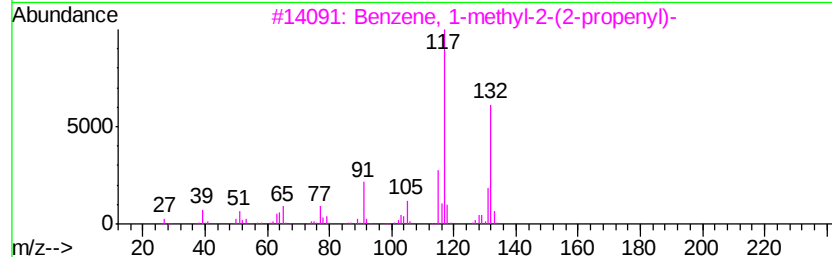
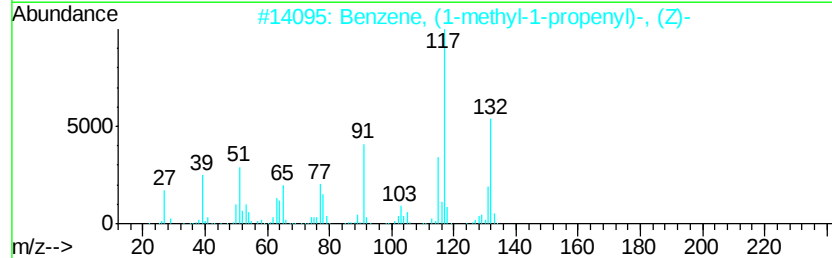
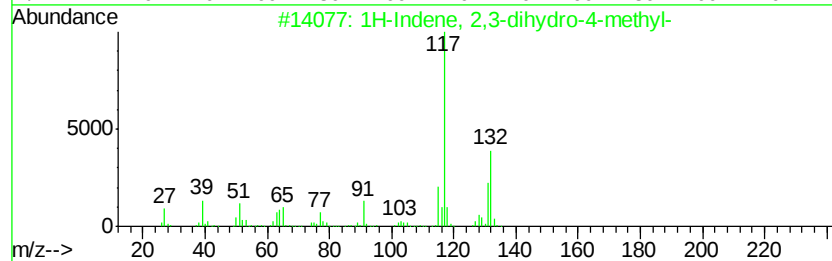
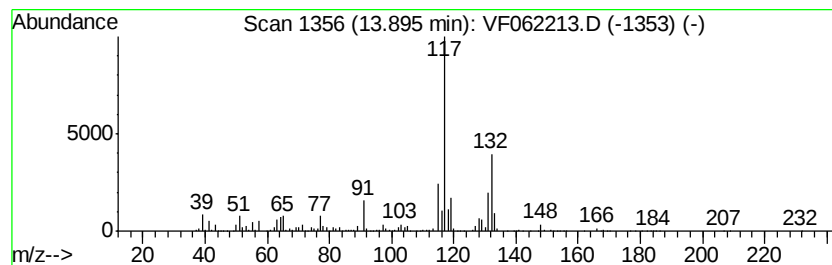
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	104.75 ug/l	3068820	1,4-Dichlorobenzene-d4	12.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	86
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15a/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-05-041719-A

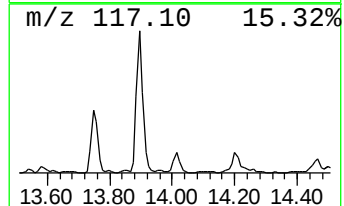
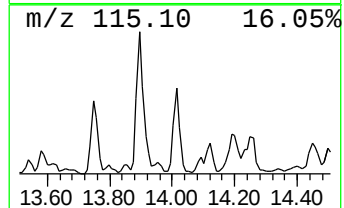
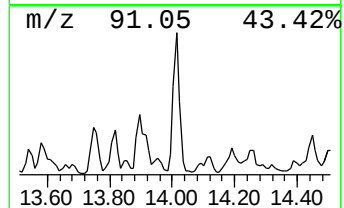
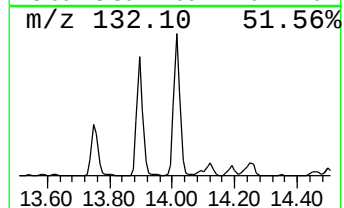
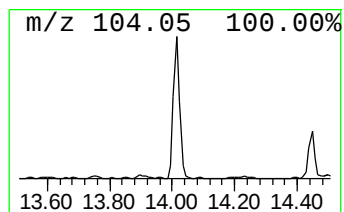
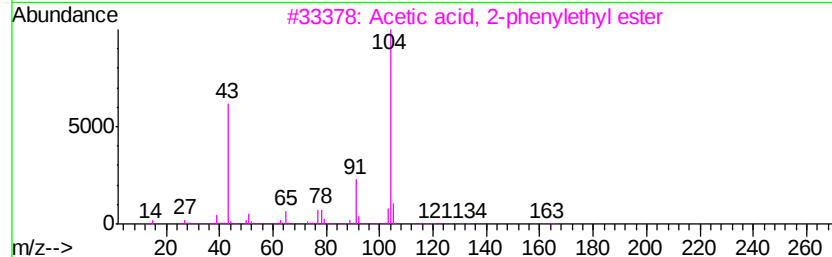
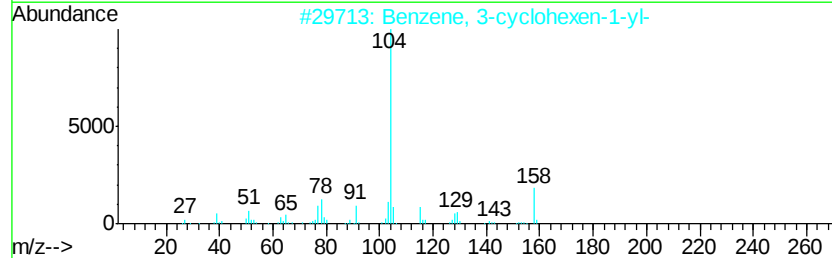
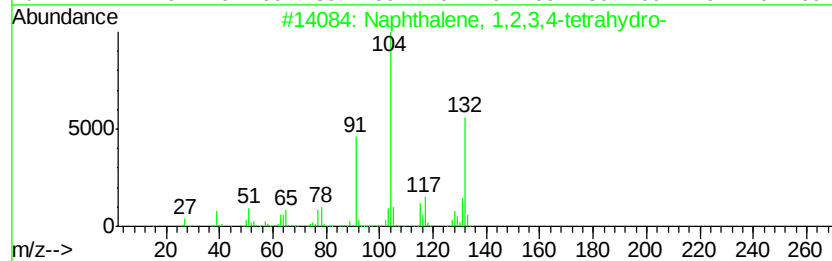
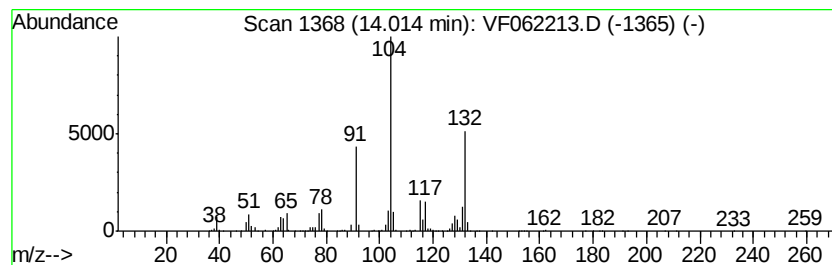
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	64.41 ug/l	1886980	1,4-Dichlorobenzene-d4	12.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47
3		Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	38
4		2-Furanone, 4-phenyltetrahydro	162	C10H10O2	1000197-38-4	38
5		2-Propenoic acid, 2-phenylethyl ...	176	C11H12O2	003530-36-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15a/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-05-041719-A

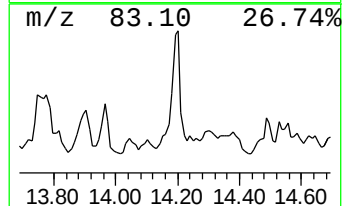
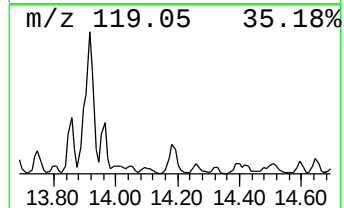
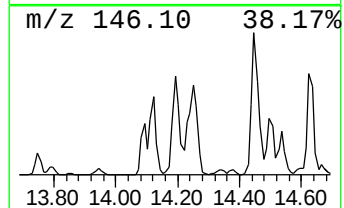
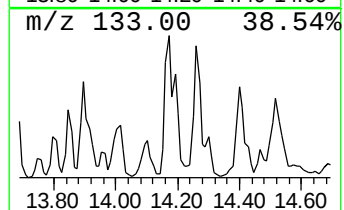
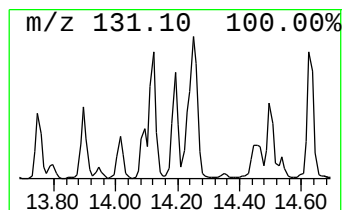
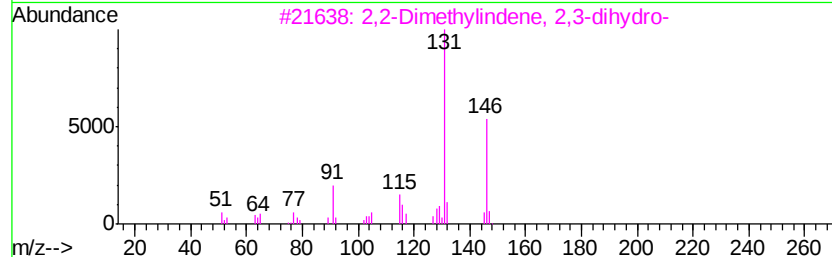
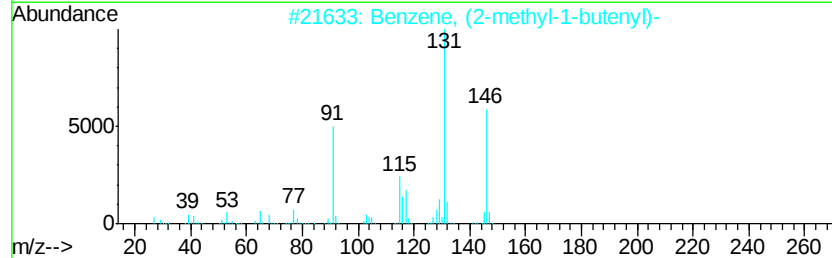
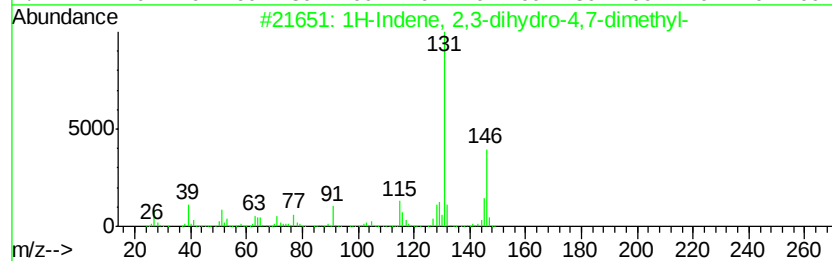
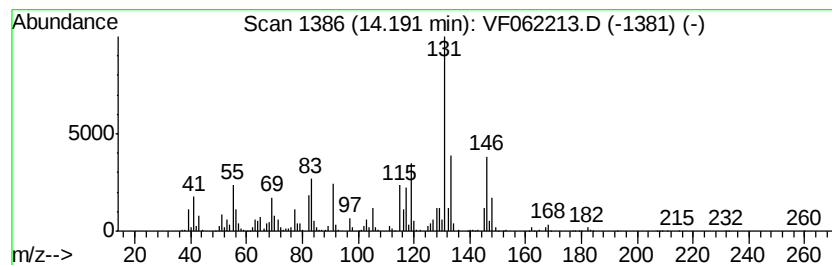
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.19	62.86 ug/l	1841810	1,4-Dichlorobenzene-d4	12.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91	
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86	
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	78	
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64	
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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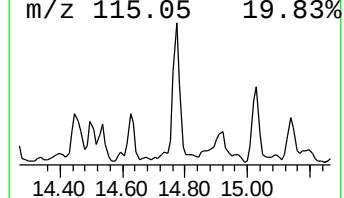
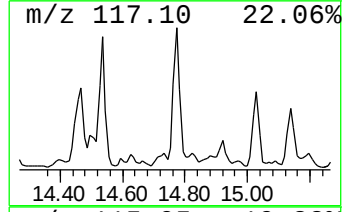
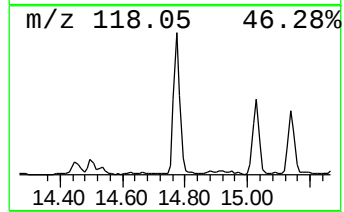
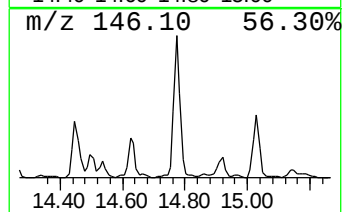
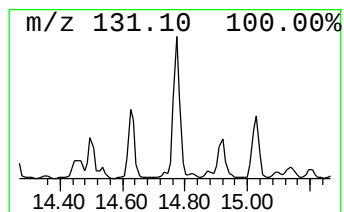
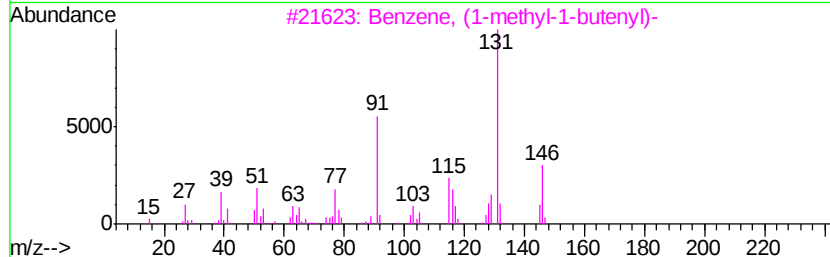
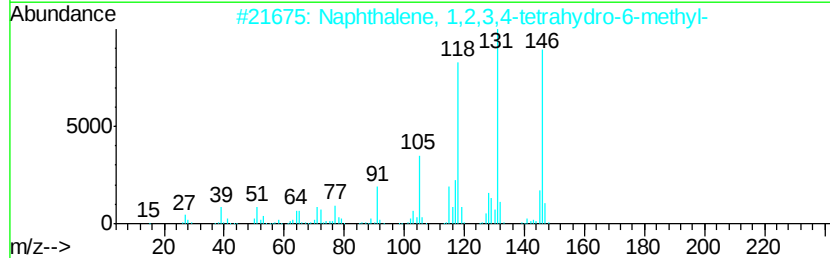
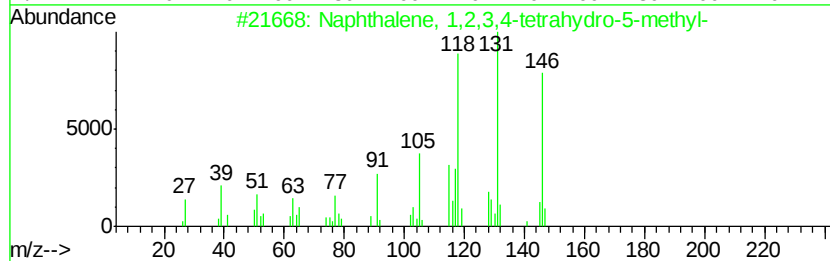
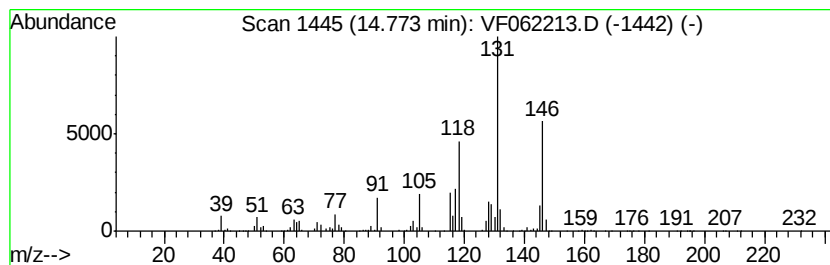
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	60.89 ug/l	1783820	1,4-Dichlorobenzene-d4	12.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
3		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
HR-05-041719-A

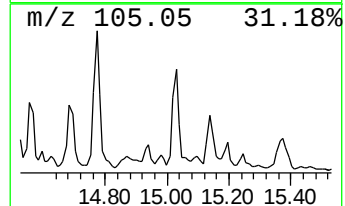
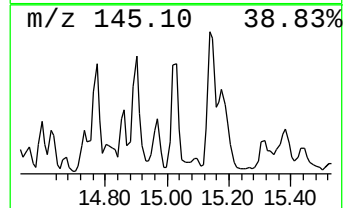
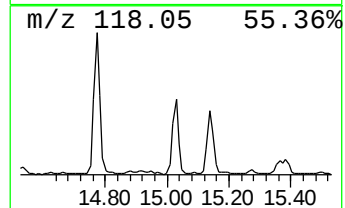
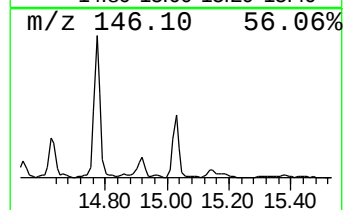
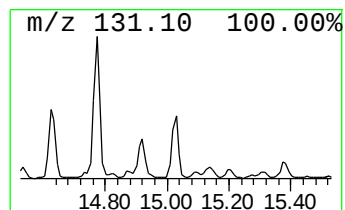
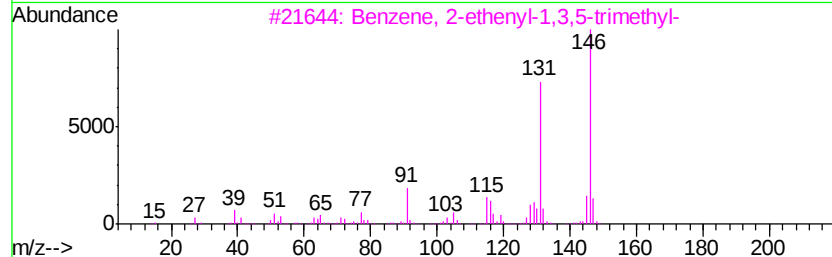
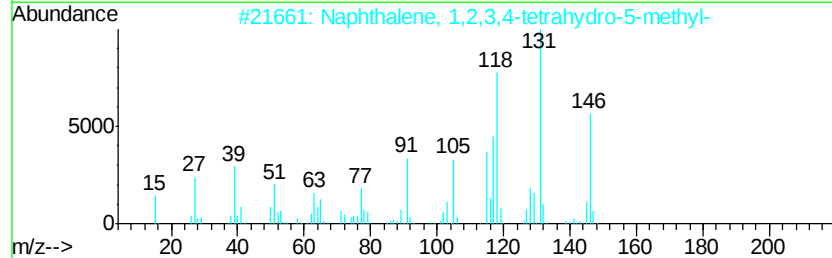
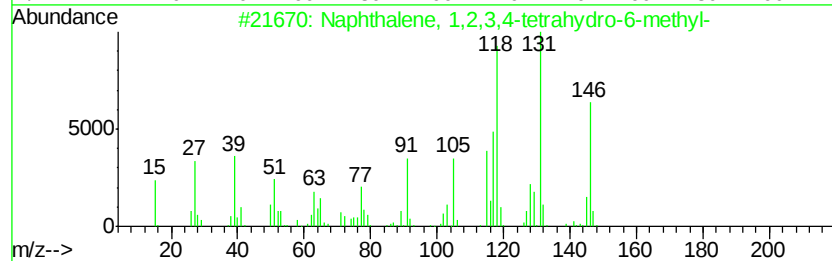
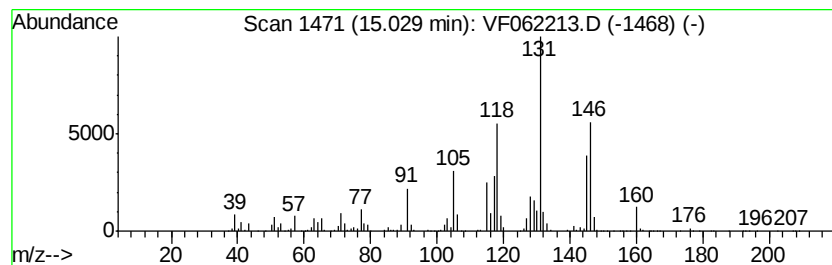
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.03	42.80 ug/l	1253960	1,4-Dichlorobenzene-d4	12.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4		2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	64
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF041819\
Data File : VF062213.D
Acq On : 18 Apr 2019 19:06
Operator : FY/SY
Sample : K2193-07
Misc : 6.15g/5mL/MSVOA_F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
HR-05-041719-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F041719S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 2-propenyl-	12.78	38.2	ug/l	1118820	4	12.61	1464900	50.0
Benzene, 1-methyl...	13.06	35.4	ug/l	1036270	4	12.61	1464900	50.0
Benzene, 1-ethyl...	13.15	54.2	ug/l	1587410	4	12.61	1464900	50.0
Benzene, 1,2,4,5-...	13.58	33.9	ug/l	993706	4	12.61	1464900	50.0
Dodecane	13.75	108.9	ug/l	3190560	4	12.61	1464900	50.0
1H-Indene, 2,3-di...	13.90	104.8	ug/l	3068820	4	12.61	1464900	50.0
Naphthalene, 1,2,...	14.01	64.4	ug/l	1886980	4	12.61	1464900	50.0
1H-Indene, 2,3-di...	14.19	62.9	ug/l	1841810	4	12.61	1464900	50.0
Naphthalene, 1,2,...	14.77	60.9	ug/l	1783820	4	12.61	1464900	50.0
Naphthalene, 1,2,...	15.03	42.8	ug/l	1253960	4	12.61	1464900	50.0