

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF110518\
 Data File : VF060595.D
 Acq On : 6 Nov 2018 5:16
 Operator : VA/AP
 Sample : J5767-03
 Misc : 6.16/5mL/MSVOA-F/SOIL
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 CL-02-110118-B

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_F\METHODS\82F110218S.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	1.093	42	47	54	rBV3	11535	50154	5.31%	0.840%
2	1.389	73	77	79	rBV	10725	24616	2.61%	0.412%
3	2.178	152	157	159	rBV4	18660	38596	4.09%	0.647%
4	2.237	159	163	169	rVB2	26460	79152	8.38%	1.326%
5	3.076	247	248	252	rVB	14103	16632	1.76%	0.279%
6	4.013	338	343	346	rBV6	4061	11949	1.27%	0.200%
7	4.121	348	354	365	rVB2	143464	441240	46.73%	7.392%
8	4.634	401	406	410	rBV8	3731	11872	1.26%	0.199%
9	4.871	422	430	443	rVV2	263914	944283	100.00%	15.819%
10	5.601	498	504	514	rBV	240946	630863	66.81%	10.568%
11	7.564	699	703	706	rVV	14804	39176	4.15%	0.656%
12	7.633	706	710	716	rVV	38153	95582	10.12%	1.601%
13	9.497	896	899	904	rBV5	4066	12270	1.30%	0.206%
14	9.783	921	928	936	rVV	74206	170501	18.06%	2.856%
15	10.148	961	965	968	rVB6	3859	10295	1.09%	0.172%
16	10.247	970	975	979	rBV7	6367	19369	2.05%	0.324%
17	10.533	1001	1004	1009	rVB5	4476	9829	1.04%	0.165%
18	10.721	1020	1023	1026	rVB5	7274	12661	1.34%	0.212%
19	11.273	1076	1079	1083	rVB6	7363	14520	1.54%	0.243%
20	11.421	1090	1094	1100	rBV	173671	292174	30.94%	4.895%
21	11.569	1105	1109	1110	rBV3	5760	9672	1.02%	0.162%
22	11.598	1110	1112	1116	rVB	7043	11211	1.19%	0.188%
23	11.727	1121	1125	1129	rBV3	19202	38997	4.13%	0.653%
24	11.786	1129	1131	1133	rVB3	7223	9470	1.00%	0.159%
25	11.825	1133	1135	1138	rBV3	8122	13561	1.44%	0.227%
26	11.934	1142	1146	1148	rBV5	5011	11544	1.22%	0.193%
27	12.023	1153	1155	1159	rVB2	14283	25943	2.75%	0.435%
28	12.111	1159	1164	1165	rBV4	6299	15040	1.59%	0.252%
29	12.131	1165	1166	1168	rBV	12303	10027	1.06%	0.168%
30	12.171	1168	1170	1172	rBV	16626	20280	2.15%	0.340%
31	12.210	1172	1174	1179	rVB	19727	30735	3.25%	0.515%
32	12.309	1181	1184	1189	rVB	10753	23969	2.54%	0.402%
33	12.437	1189	1197	1199	rBV7	15812	44646	4.73%	0.748%
34	12.476	1199	1201	1205	rVB3	13572	22782	2.41%	0.382%

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35	12.555	1206	1209	1213	rBV	212832	339577	35.96%	5.689%
36	12.614	1213	1215	1219	rVB	54043	76880	8.14%	1.288%
37	12.733	1224	1227	1230	rBV3	12588	21344	2.26%	0.358%
38	12.841	1233	1238	1241	rVB5	23741	63923	6.77%	1.071%
39	12.940	1245	1248	1250	rBV3	11086	21082	2.23%	0.353%
40	13.009	1253	1255	1258	rVB	16236	25260	2.68%	0.423%
41	13.108	1260	1265	1268	rBV3	27493	57327	6.07%	0.960%
42	13.157	1268	1270	1272	rBV	60061	80435	8.52%	1.347%
43	13.206	1272	1275	1278	rVB	31559	39314	4.16%	0.659%
44	13.256	1278	1280	1282	rVB	26375	29608	3.14%	0.496%
45	13.285	1282	1283	1290	rVB4	15183	41023	4.34%	0.687%
46	13.384	1290	1293	1295	rVB2	38954	49597	5.25%	0.831%
47	13.443	1295	1299	1303	rBV3	16921	48665	5.15%	0.815%
48	13.502	1303	1305	1307	rVB	24789	26604	2.82%	0.446%
49	13.542	1307	1309	1311	rBV2	11846	19484	2.06%	0.326%
50	13.581	1311	1313	1316	rBV2	25118	31864	3.37%	0.534%
51	13.640	1316	1319	1322	rVB2	23171	43059	4.56%	0.721%
52	13.709	1323	1326	1329	rBV2	55577	94805	10.04%	1.588%
53	13.759	1329	1331	1334	rVB3	14625	23151	2.45%	0.388%
54	13.818	1334	1337	1338	rBV	20076	30637	3.24%	0.513%
55	13.857	1338	1341	1345	rVB3	65788	109885	11.64%	1.841%
56	13.926	1345	1348	1350	rVB3	48817	71785	7.60%	1.203%
57	13.976	1350	1353	1357	rVB2	43002	65427	6.93%	1.096%
58	14.084	1360	1364	1367	rBV	22067	36655	3.88%	0.614%
59	14.153	1367	1371	1373	rBV3	41793	105155	11.14%	1.762%
60	14.193	1373	1375	1377	rBV2	22622	39590	4.19%	0.663%
61	14.301	1384	1386	1388	rBV3	10710	17421	1.84%	0.292%
62	14.410	1394	1397	1401	rBV2	31854	69799	7.39%	1.169%
63	14.469	1401	1403	1406	rVV	70575	100805	10.68%	1.689%
64	14.568	1410	1413	1414	rBV3	10425	14898	1.58%	0.250%
65	14.597	1414	1416	1419	rVB3	12955	20941	2.22%	0.351%
66	14.647	1419	1421	1423	rBV3	7586	10088	1.07%	0.169%
67	14.745	1425	1431	1434	rBV	89968	146831	15.55%	2.460%
68	14.883	1440	1445	1450	rVB3	33134	85190	9.02%	1.427%
69	15.002	1454	1457	1460	rBV	81451	115945	12.28%	1.942%
70	15.120	1465	1469	1471	rVV	69033	112374	11.90%	1.882%
71	15.169	1471	1474	1479	rVB4	36363	85921	9.10%	1.439%

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Integration Parameters: RTEINT.P

Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	15.248	1479	1482	1483	rBV3	11973	20783	2.20%	0.348%
73	15.288	1483	1486	1488	rVV3	18071	32861	3.48%	0.550%
74	15.357	1488	1493	1496	rVV2	50307	126436	13.39%	2.118%
75	15.406	1496	1498	1502	rVB2	22563	37419	3.96%	0.627%
76	15.653	1519	1523	1528	rBV	45917	91803	9.72%	1.538%
77	15.722	1528	1530	1534	rVV4	8619	20700	2.19%	0.347%
78	15.830	1538	1541	1545	rVB4	13020	25914	2.74%	0.434%
79	15.949	1550	1553	1554	rBV3	7607	12924	1.37%	0.217%
80	15.988	1554	1557	1558	rBV3	9064	14634	1.55%	0.245%

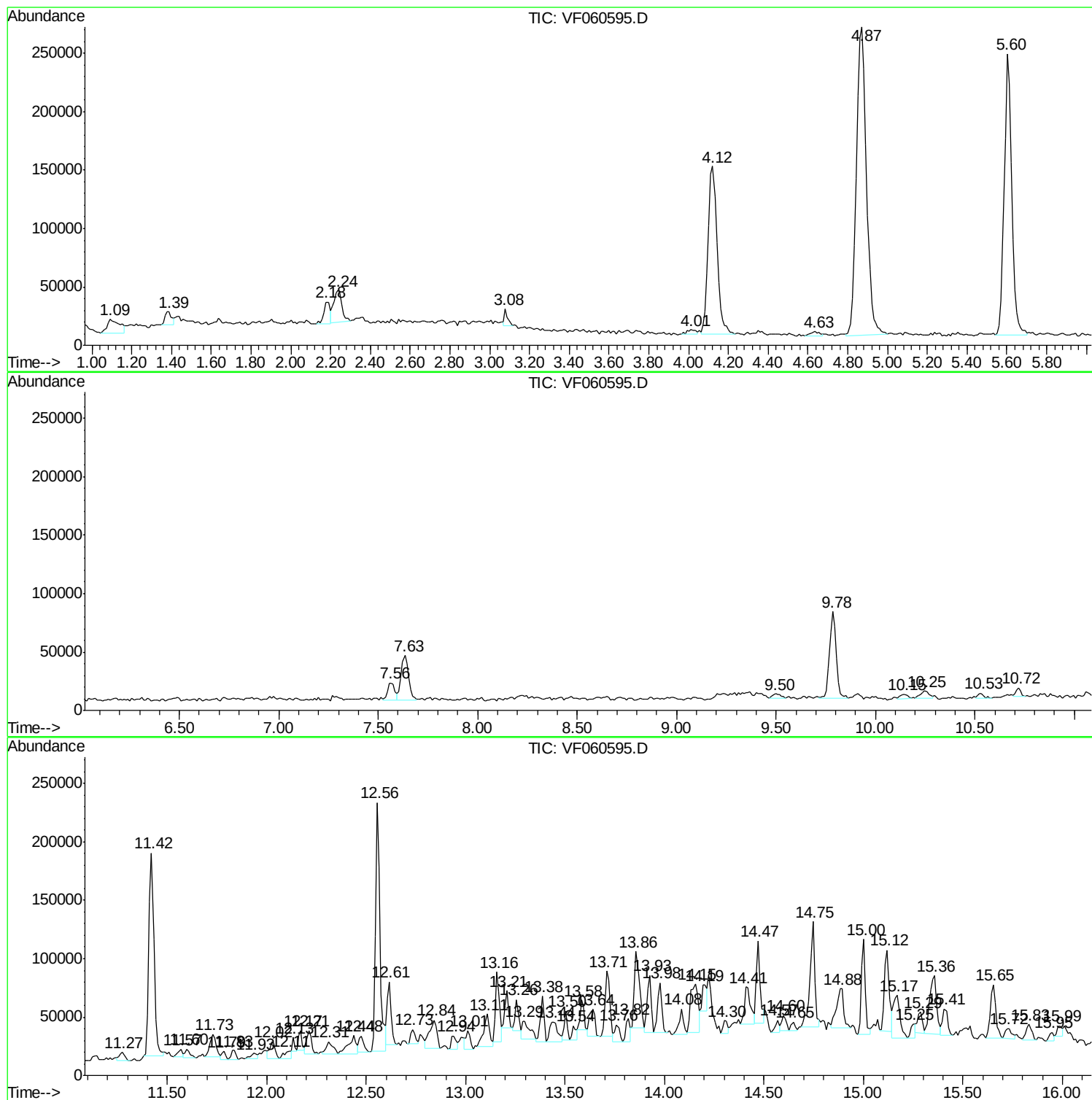
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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



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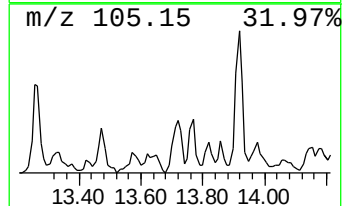
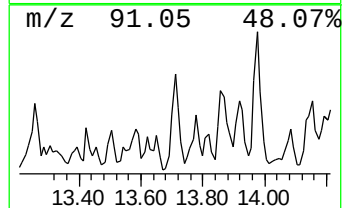
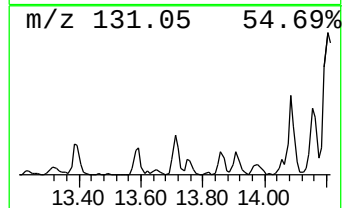
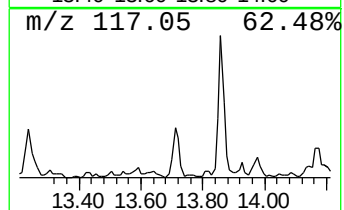
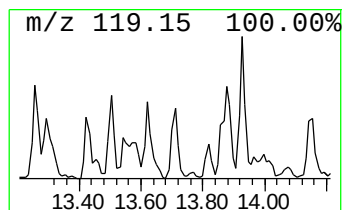
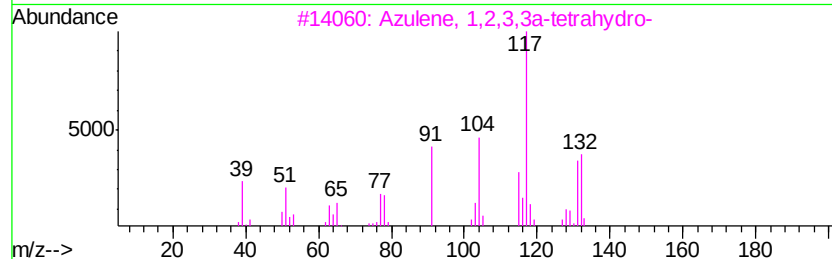
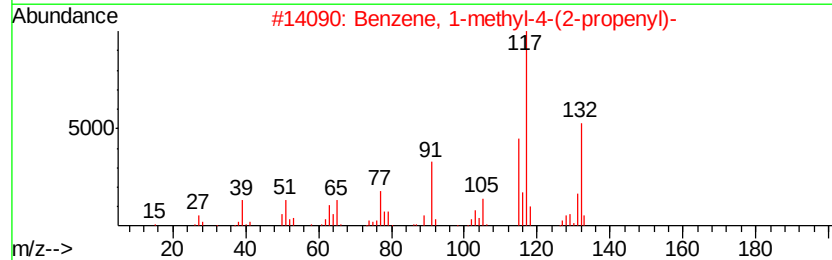
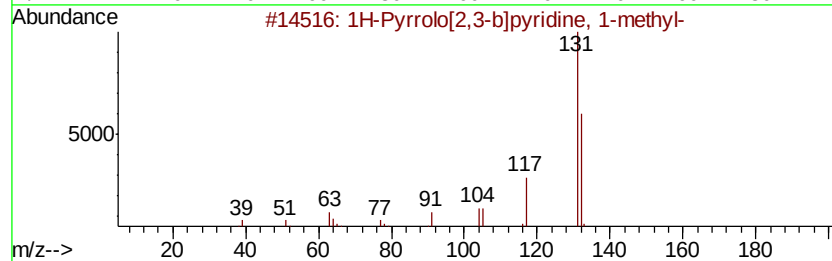
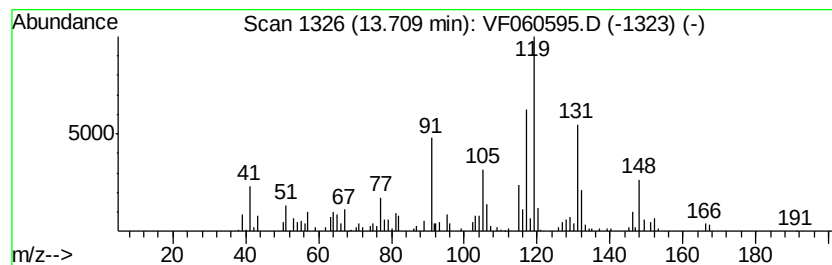
Instrument :
MSVOA_F
ClientSampleId :
CL-02-110118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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

Peak Number 1 unknown13.71 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	13.96 ug/l	94805	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Pyrrolo[2,3-b]pyridine, 1-met...	132 C8H8N2	027257-15-4 42
2		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 38
3		Azulene, 1,2,3,3a-tetrahydro-	132 C10H12	033877-87-1 38
4		Benzene, 2-butenyl-	132 C10H12	001560-06-1 38
5		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 38



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ALS Vial : 37 Sample Multiplier: 1

Instrument :
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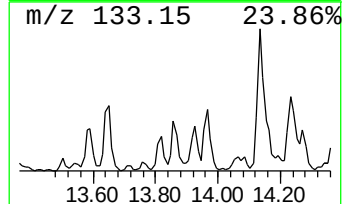
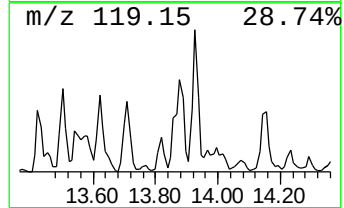
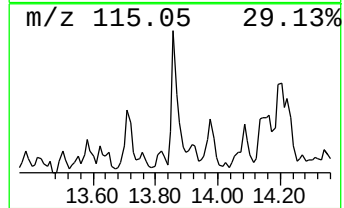
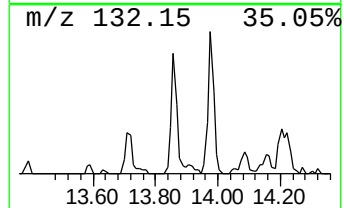
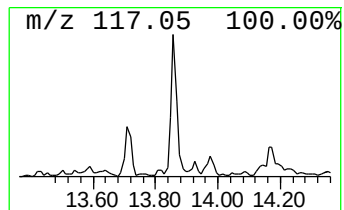
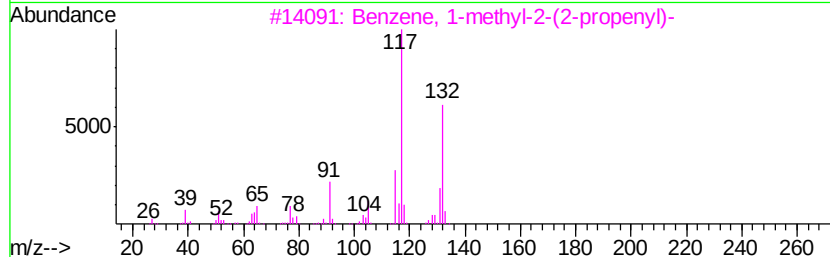
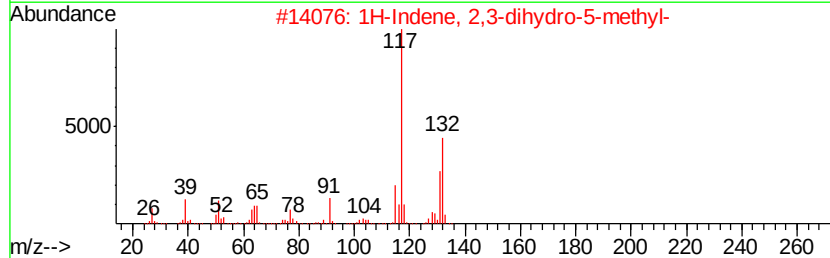
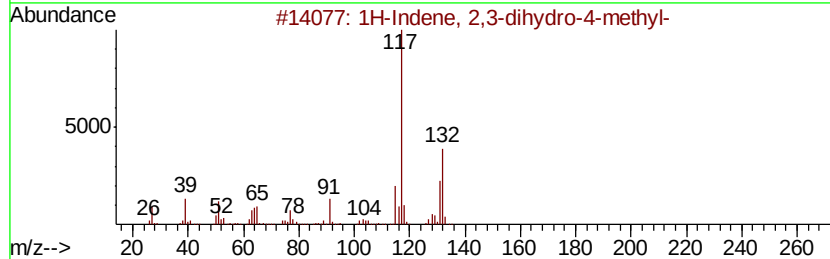
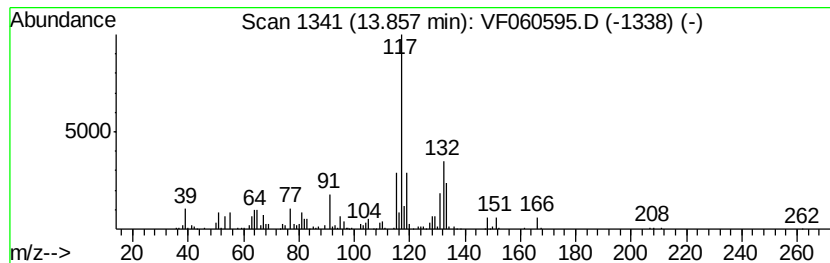
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Quant Title : SW846 8260

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

Peak Number 2 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	16.18 ug/l	109885	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	62
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	62



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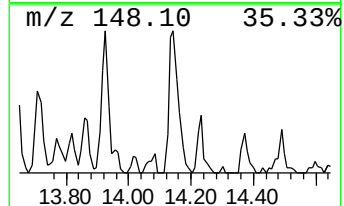
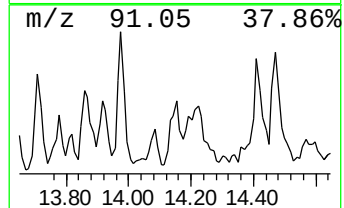
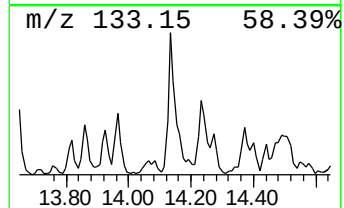
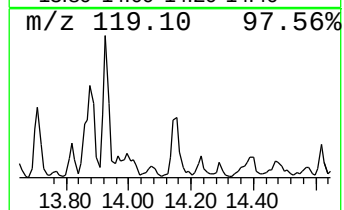
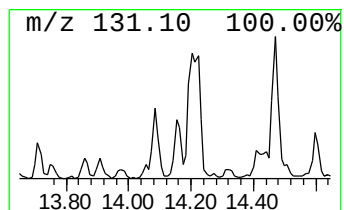
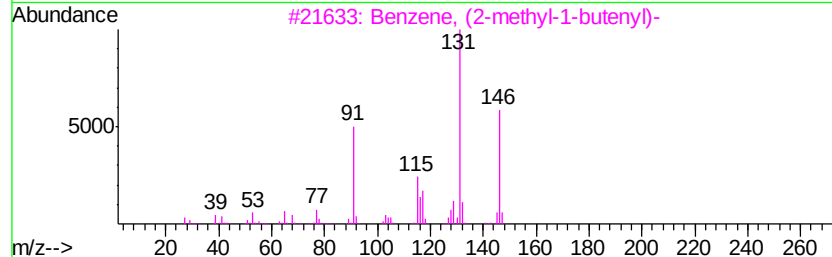
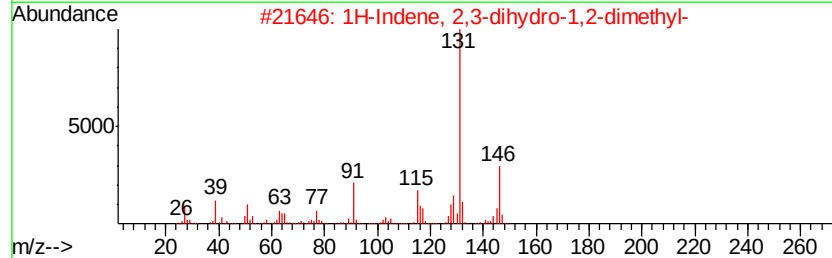
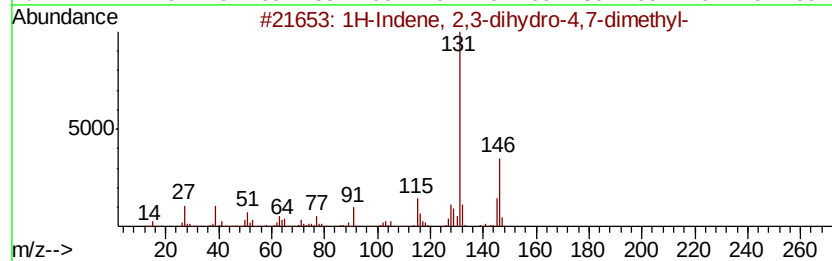
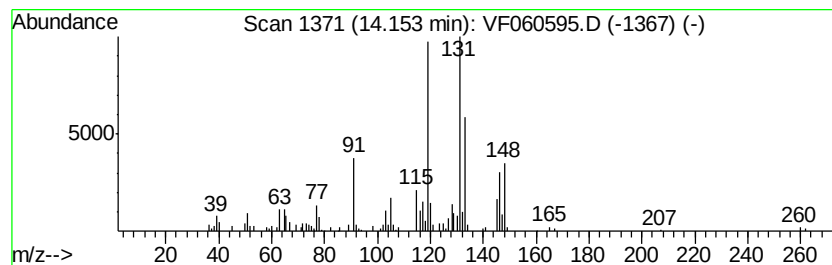
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Quant Title : SW846 8260

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

Peak Number 3 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	15.48 ug/l	105155	1,4-Dichlorobenzene-d4	12.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	50	
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	46	
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	46	
4	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	46	
5	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	41	



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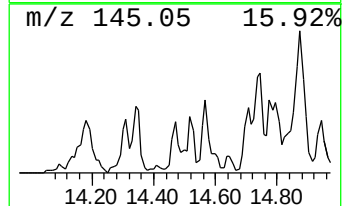
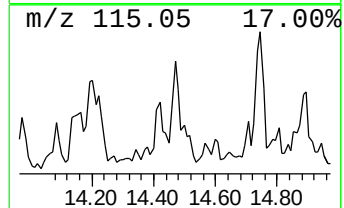
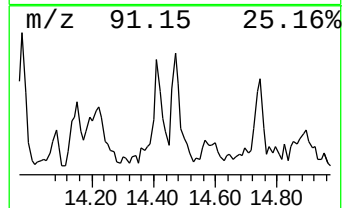
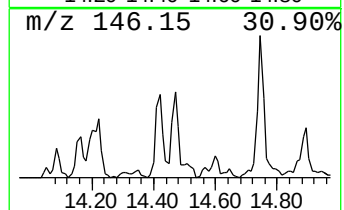
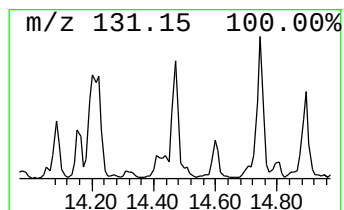
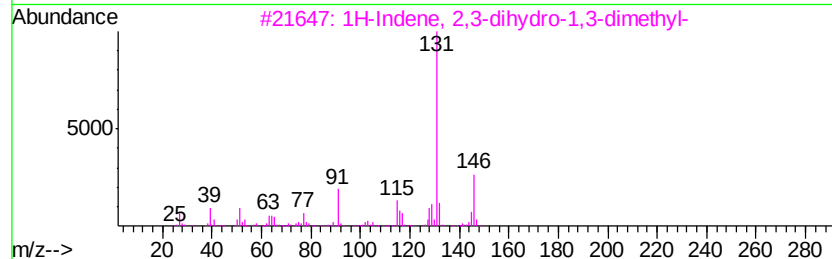
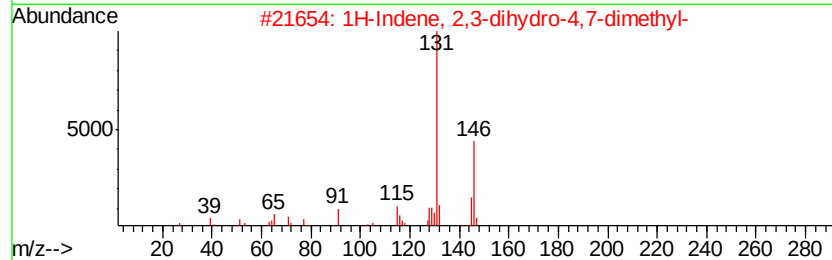
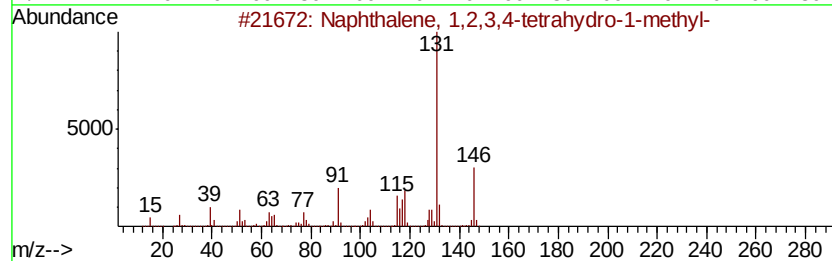
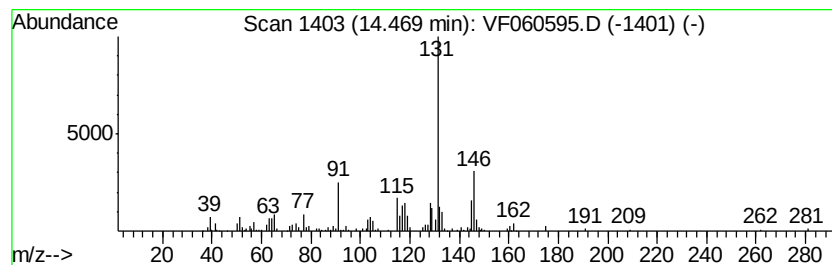
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	14.84 ug/l	100805	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	96
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF110518\
Data File : VF060595.D
Acq On : 6 Nov 2018 5:16
Operator : VA/AP
Sample : J5767-03
Misc : 6.16/5mL/MSVOA-F/SOIL
ALS Vial : 37 Sample Multiplier: 1

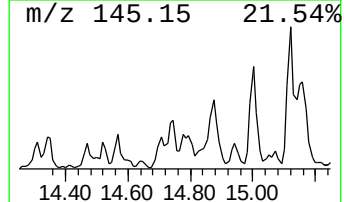
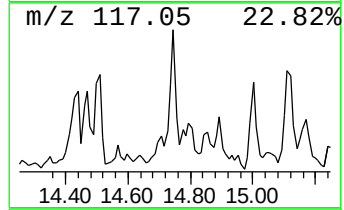
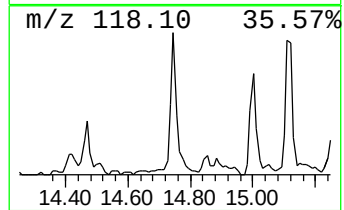
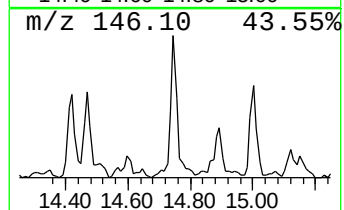
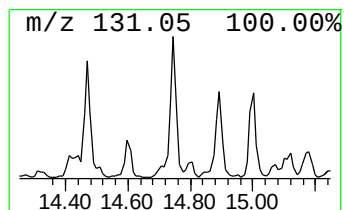
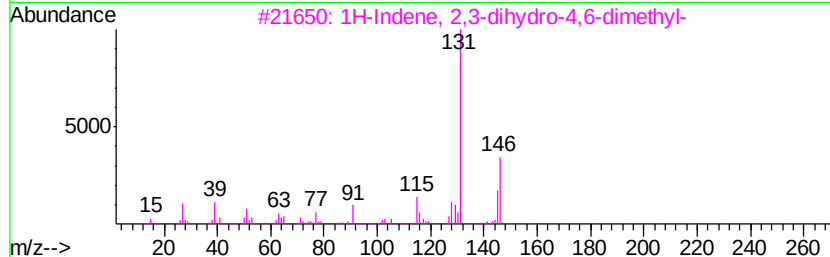
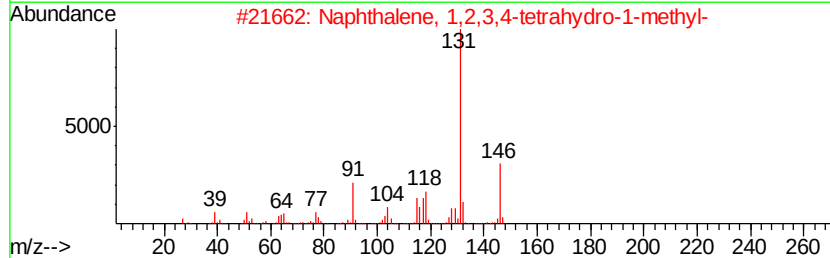
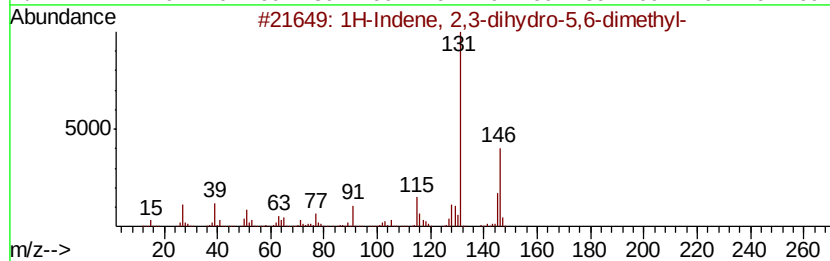
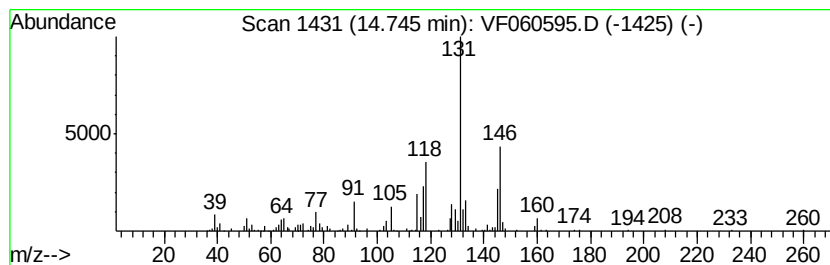
Instrument :
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ClientSampleId :
CL-02-110118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-5,6-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	21.62 ug/l	146831	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-5,6-dimet...	146 C11H14	001075-22-5	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	76
3	1H-Indene, 2,3-dihydro-4,6-dimet...	146 C11H14	001685-82-1	70
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146 C10H10O	017496-14-9	70
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF110518\
Data File : VF060595.D
Acq On : 6 Nov 2018 5:16
Operator : VA/AP
Sample : J5767-03
Misc : 6.16/5mL/MSVOA-F/SOIL
ALS Vial : 37 Sample Multiplier: 1

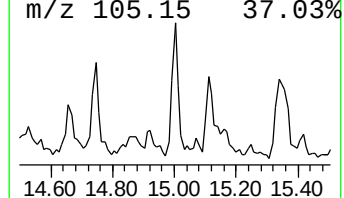
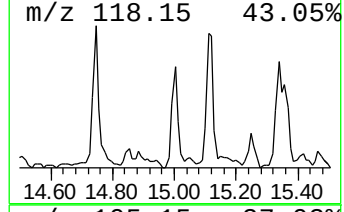
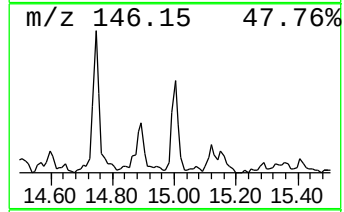
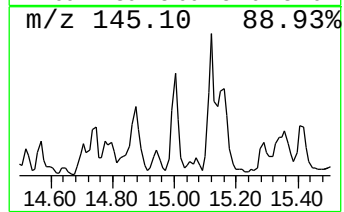
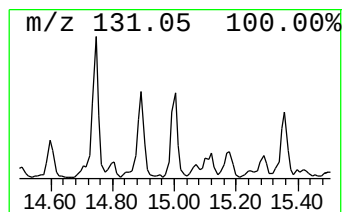
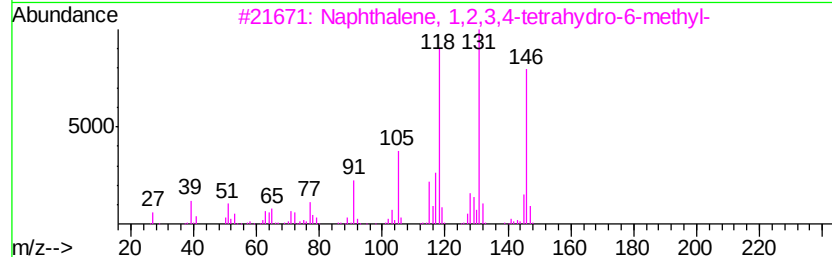
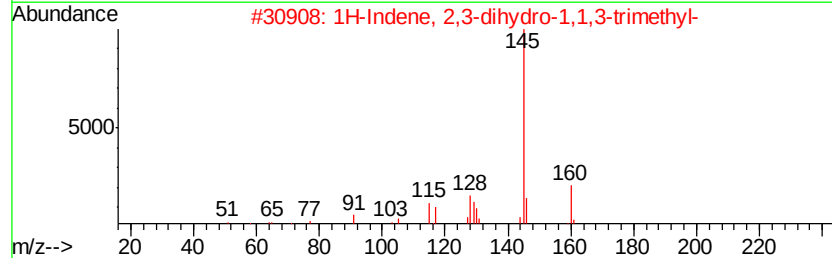
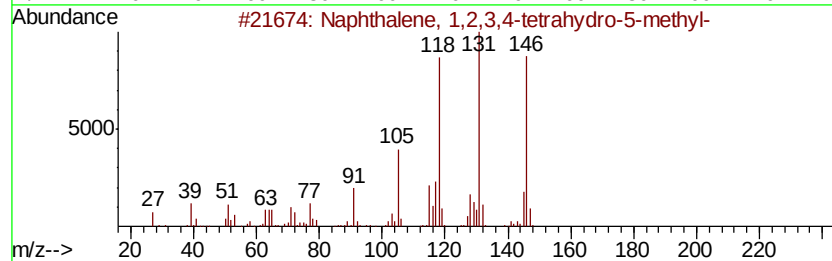
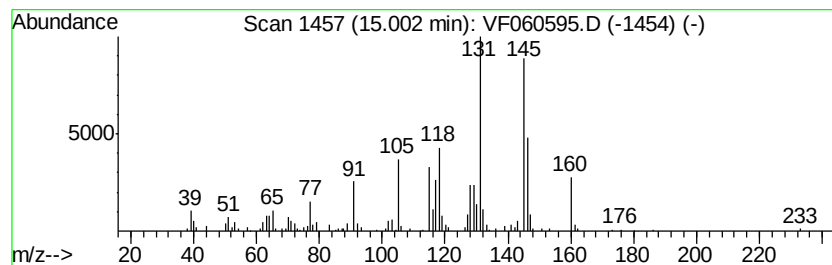
Instrument :
MSVOA_F
ClientSampleId :
CL-02-110118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	17.07 ug/l	115945	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 58
2		1H-Indene, 2,3-dihydro-1,1,3-tri...	160 C12H16	002613-76-5 55
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 49
4		2-Propenal, 2-methyl-3-phenyl-	146 C10H10O	000101-39-3 46
5		Benzene, 1-(1-methylethenyl)-2-(...	160 C12H16	005557-93-7 46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF110518\
Data File : VF060595.D
Acq On : 6 Nov 2018 5:16
Operator : VA/AP
Sample : J5767-03
Misc : 6.16/5mL/MSVOA-F/SOIL
ALS Vial : 37 Sample Multiplier: 1

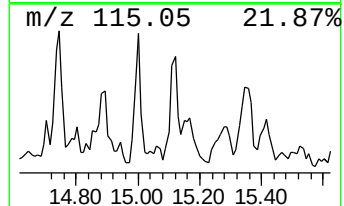
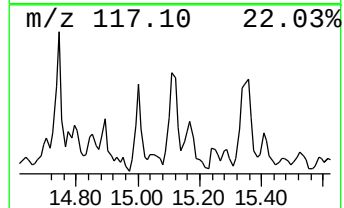
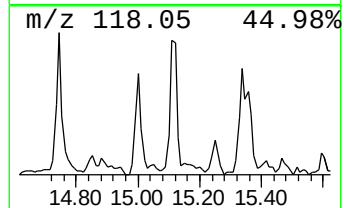
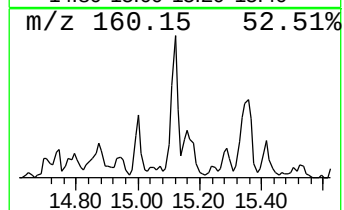
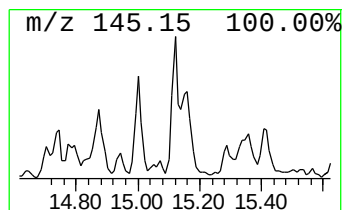
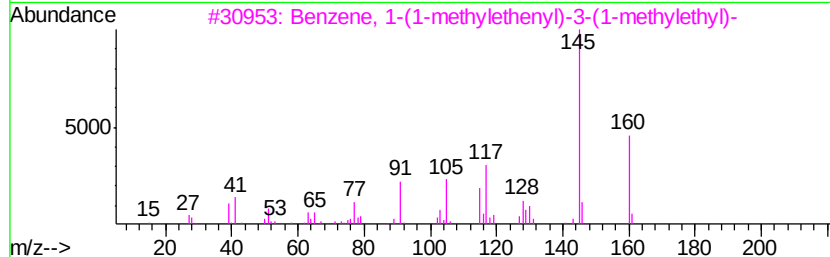
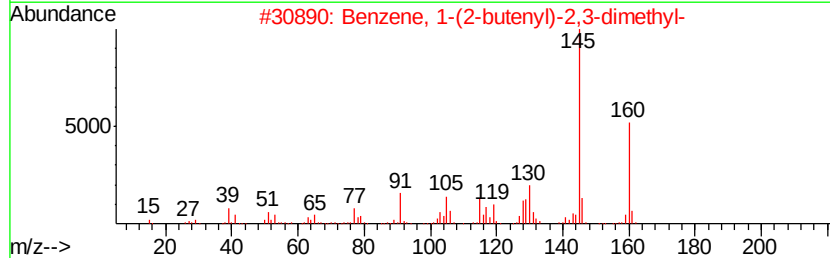
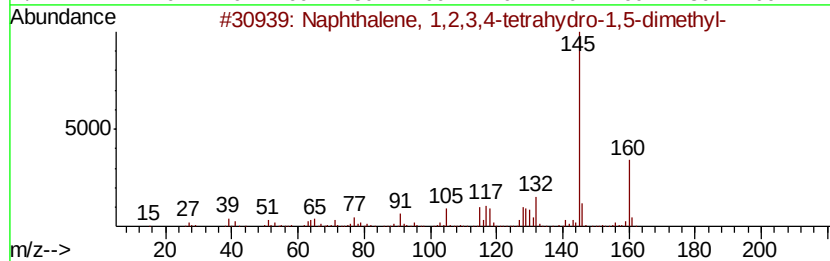
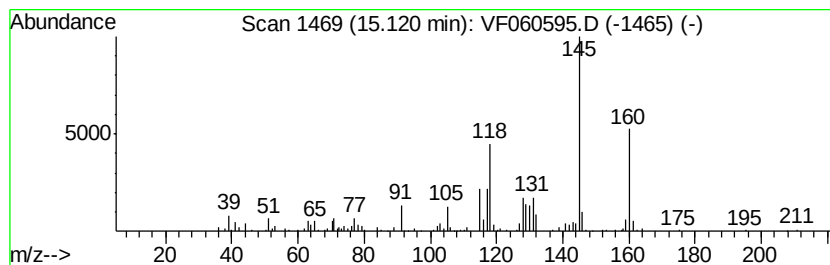
Instrument :
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ClientSampleId :
CL-02-110118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	16.55 ug/l	112374	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	021564-91-0 87
2		Benzene, 1-(2-butenyl)-2,3-dimet...	160 C12H16	054340-85-1 83
3		Benzene, 1-(1-methylethenyl)-3-(...	160 C12H16	001129-29-9 76
4		1H-Indene, 2,3-dihydro-1,1,3-tri...	160 C12H16	002613-76-5 76
5		Benzene, 4-(2-butenyl)-1,2-dimet...	160 C12H16	054340-86-2 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF110518\
Data File : VF060595.D
Acq On : 6 Nov 2018 5:16
Operator : VA/AP
Sample : J5767-03
Misc : 6.16/5mL/MSVOA-F/SOIL
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
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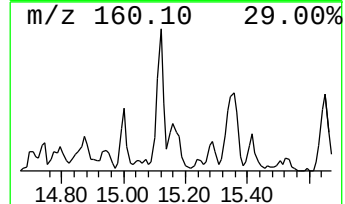
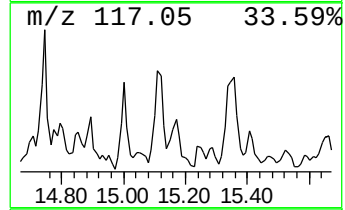
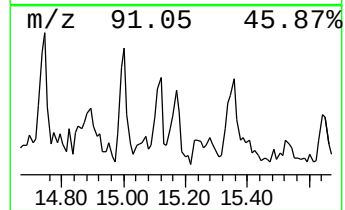
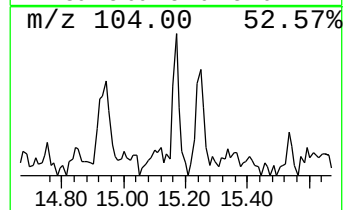
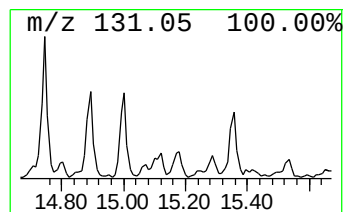
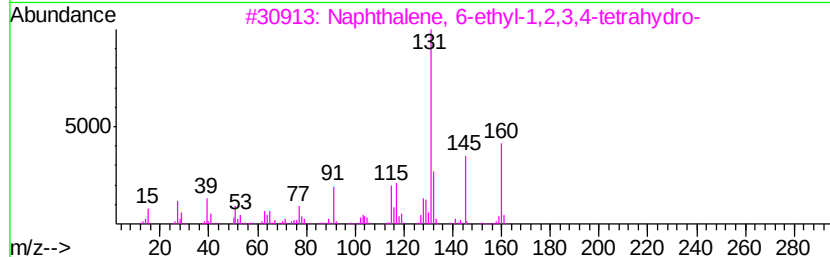
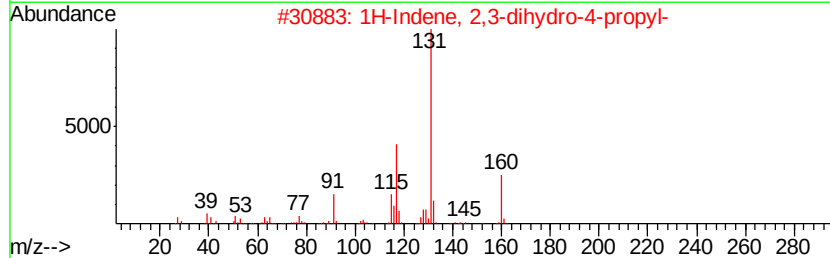
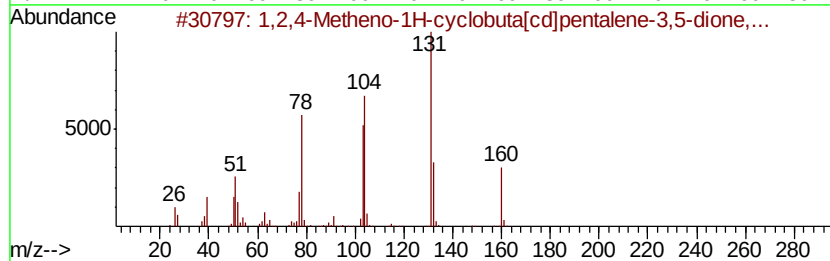
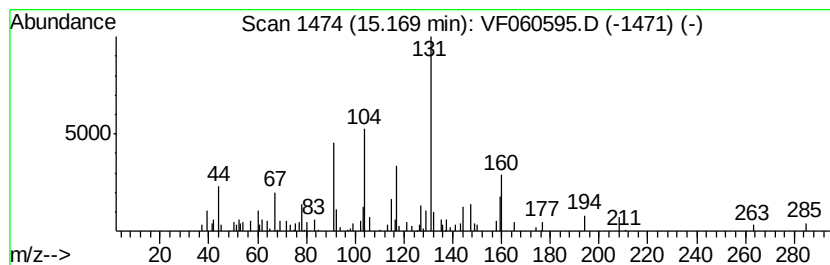
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Quant Title : SW846 8260

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

Peak Number 8 1,2,4-Metheno-1H-cyclobuta[...] Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	12.65 ug/l	85921	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,4-Metheno-1H-cyclobuta[cd]pe...	160	C10H8O2	014725-77-0	50
2		1H-Indene, 2,3-dihydro-4-propyl-	160	C12H16	092013-16-6	46
3		Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	43
4		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	43
5		Quinoxalin-6-yl-methanol	160	C9H8N2O	1000318-21-1	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF110518\
Data File : VF060595.D
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Operator : VA/AP
Sample : J5767-03
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ALS Vial : 37 Sample Multiplier: 1

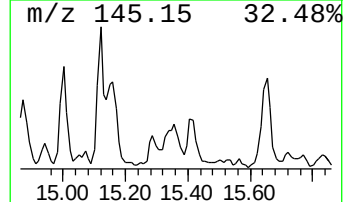
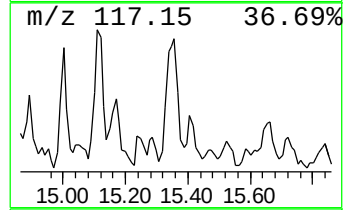
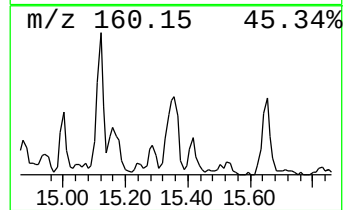
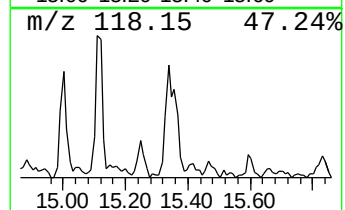
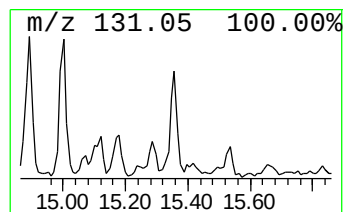
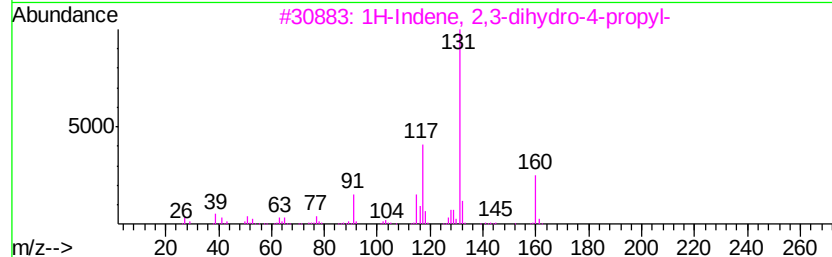
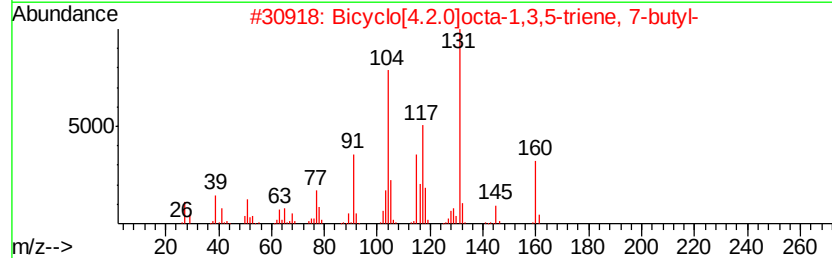
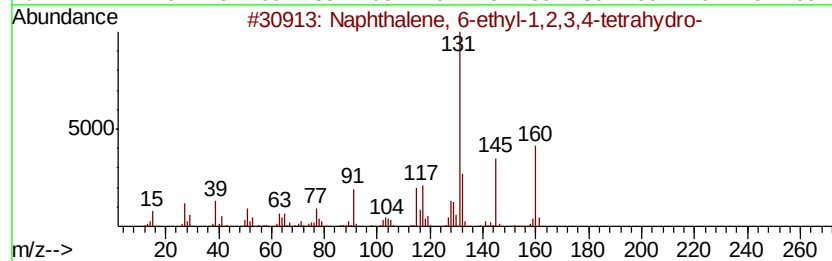
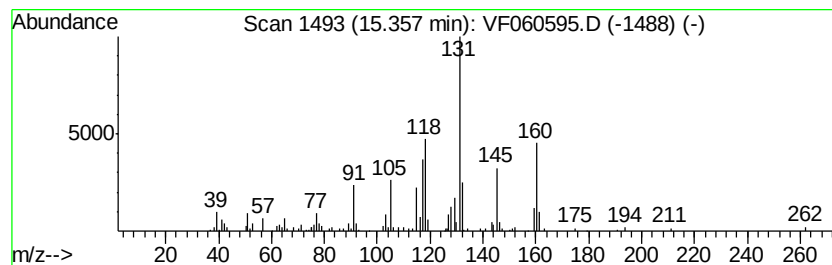
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ClientSampleId :
CL-02-110118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 6-ethyl-1,2,3,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	18.62 ug/l	126436	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 6-ethyl-1,2,3,4-tet...	160 C12H16	022531-20-0 76
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	160 C12H16	078920-29-3 59
3		1H-Indene, 2,3-dihydro-4-propyl-	160 C12H16	092013-16-6 53
4		1,4-Methanonaphthalen-9-ol, 1,2,...	160 C11H12O	055255-94-2 49
5		Quinoxalin-6-yl-methanol	160 C9H8N2O	1000318-21-1 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF110518\
Data File : VF060595.D
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Sample : J5767-03
Misc : 6.16/5mL/MSVOA-F/SOIL
ALS Vial : 37 Sample Multiplier: 1

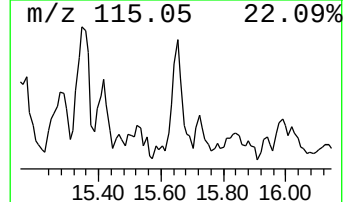
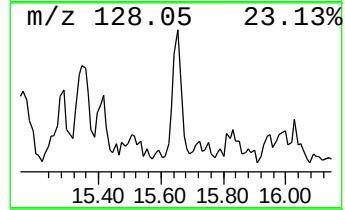
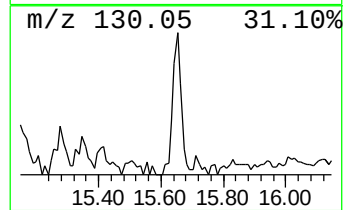
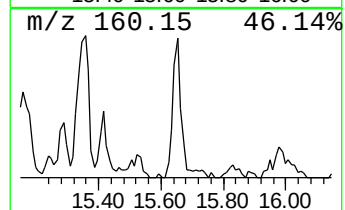
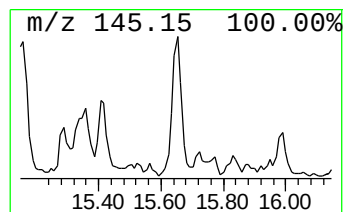
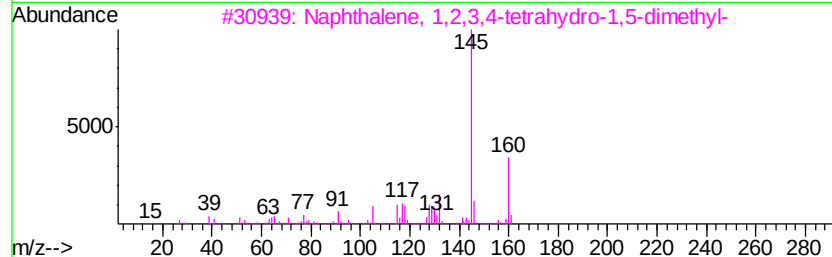
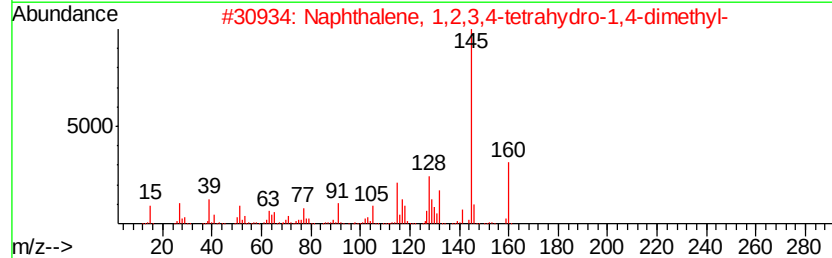
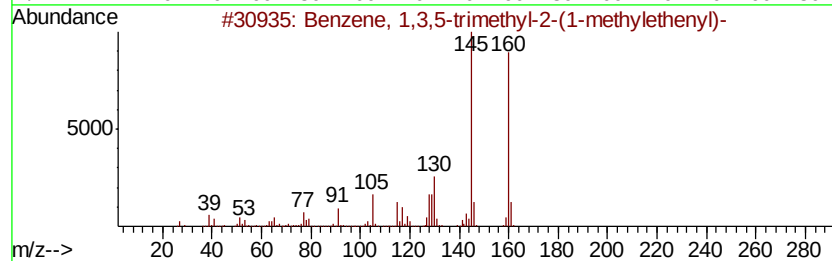
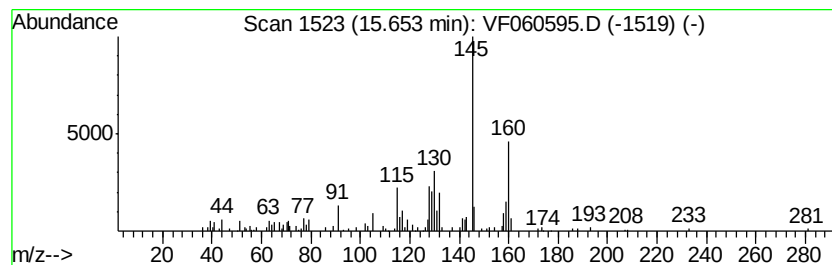
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	13.52 ug/l	91803	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3,5-trimethyl-2-(1-me...	160 C12H16	014679-13-1 94
2		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6 90
3		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	021564-91-0 81
4		Benzene, 1-(2-butenyl)-2,3-dimet...	160 C12H16	054340-85-1 81
5		1H-Inden-1-one, 2,3-dihydro-3,3-...	160 C11H12O	026465-81-6 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF110518\
Data File : VF060595.D
Acq On : 6 Nov 2018 5:16
Operator : VA/AP
Sample : J5767-03
Misc : 6.16/5mL/MSVOA-F/SOIL
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
CL-02-110118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F110218S.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
unknown	13.71	14.0	ug/l	94805	4	12.56	339577	50.0
1H-Indene, 2,3-di...	13.86	16.2	ug/l	109885	4	12.56	339577	50.0
1H-Indene, 2,3-di...	14.15	15.5	ug/l	105155	4	12.56	339577	50.0
Naphthalene, 1,2,...	14.47	14.8	ug/l	100805	4	12.56	339577	50.0
1H-Indene, 2,3-di...	14.75	21.6	ug/l	146831	4	12.56	339577	50.0
Naphthalene, 1,2,...	15.00	17.1	ug/l	115945	4	12.56	339577	50.0
Naphthalene, 1,2,...	15.12	16.6	ug/l	112374	4	12.56	339577	50.0
1,2,4-Metheno-1H-...	15.17	12.7	ug/l	85921	4	12.56	339577	50.0
Naphthalene, 6-et...	15.36	18.6	ug/l	126436	4	12.56	339577	50.0
Benzene, 1,3,5-tr...	15.65	13.5	ug/l	91803	4	12.56	339577	50.0