

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF031919\
 Data File : VF062018.D
 Acq On : 19 Mar 2019 15:14
 Operator : VA/AP
 Sample : K1951-04
 Misc : 5.58g/5mL/MSVOA-F/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 FK-GF-03-006-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\82F031619S.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	1.082	43	46	51	rVB3	5601	13130	1.27%	0.149%
2	2.176	153	157	159	rBV5	8110	14860	1.44%	0.168%
3	2.236	159	163	171	rVV	20088	53542	5.17%	0.607%
4	2.374	173	177	180	rVB3	8334	20514	1.98%	0.233%
5	2.916	227	232	236	rBV4	6161	15470	1.49%	0.175%
6	3.863	322	328	333	rBV4	4788	16251	1.57%	0.184%
7	4.001	333	342	346	rBV2	52593	204602	19.76%	2.319%
8	4.099	346	352	359	rVB	151970	485972	46.95%	5.509%
9	4.405	372	383	399	rVB4	39427	256981	24.82%	2.913%
10	4.612	399	404	407	rBV2	10485	37959	3.67%	0.430%
11	4.849	420	428	439	rBV2	292600	1034440	99.93%	11.726%
12	5.440	484	488	490	rBV3	4275	11543	1.12%	0.131%
13	5.480	490	492	496	rVV4	7508	17840	1.72%	0.202%
14	5.578	496	502	512	rVV	334637	890044	85.98%	10.089%
15	6.791	621	625	628	rBV3	4428	11279	1.09%	0.128%
16	6.998	642	646	650	rVB4	5519	13224	1.28%	0.150%
17	7.531	695	700	706	rVV	435610	1035192	100.00%	11.735%
18	7.600	706	707	713	rVB	19237	28057	2.71%	0.318%
19	8.438	787	792	798	rVB5	9074	27077	2.62%	0.307%
20	8.606	803	809	814	rBV3	7869	25275	2.44%	0.287%
21	8.951	838	844	847	rBV6	6622	16832	1.63%	0.191%
22	9.089	855	858	861	rVV2	5312	11658	1.13%	0.132%
23	9.266	871	876	880	rBV3	7443	21053	2.03%	0.239%
24	9.513	895	901	907	rBV5	4114	17065	1.65%	0.193%
25	9.602	907	910	912	rVV2	4774	10692	1.03%	0.121%
26	9.749	920	925	933	rVV	368360	831395	80.31%	9.425%
27	9.878	936	938	942	rVB4	6715	13325	1.29%	0.151%
28	10.104	958	961	964	rBV3	4950	11322	1.09%	0.128%
29	10.213	967	972	975	rBV4	6726	19182	1.85%	0.217%
30	10.499	997	1001	1004	rVB3	6028	14249	1.38%	0.162%
31	10.657	1013	1017	1018	rBV4	11675	23605	2.28%	0.268%
32	10.696	1018	1021	1024	rVB	19200	38568	3.73%	0.437%
33	10.824	1031	1034	1038	rVB3	10212	23163	2.24%	0.263%
34	11.022	1050	1054	1058	rBV3	6369	13561	1.31%	0.154%

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Integrator: RTE
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35	11.248	1068	1077	1081	rVB5	11869	42032	4.06%	0.476%
36	11.396	1086	1092	1098	rBV	348643	610795	59.00%	6.924%
37	11.544	1103	1107	1110	rBV2	23190	44477	4.30%	0.504%
38	11.643	1115	1117	1120	rVV4	6146	12886	1.24%	0.146%
39	11.702	1120	1123	1127	rVV2	27474	57498	5.55%	0.652%
40	11.810	1130	1134	1137	rVB	41758	67651	6.54%	0.767%
41	12.008	1149	1154	1158	rVB	32231	65497	6.33%	0.742%
42	12.185	1169	1172	1176	rVB	35094	55814	5.39%	0.633%
43	12.274	1176	1181	1187	rBV7	10386	30150	2.91%	0.342%
44	12.412	1193	1195	1197	rBV	15309	24370	2.35%	0.276%
45	12.441	1197	1198	1204	rVB2	16878	27440	2.65%	0.311%
46	12.530	1204	1207	1211	rBV	357812	580669	56.09%	6.582%
47	12.589	1211	1213	1220	rVB	80098	124233	12.00%	1.408%
48	12.708	1220	1225	1228	rBV2	29274	60783	5.87%	0.689%
49	12.757	1228	1230	1232	rVV	36689	49489	4.78%	0.561%
50	12.816	1232	1236	1240	rVB2	46111	97541	9.42%	1.106%
51	12.915	1244	1246	1248	rBV3	8561	13527	1.31%	0.153%
52	12.994	1251	1254	1257	rVB2	26775	40983	3.96%	0.465%
53	13.082	1257	1263	1266	rBV2	29541	72600	7.01%	0.823%
54	13.142	1266	1269	1274	rBV2	38759	87135	8.42%	0.988%
55	13.211	1274	1276	1277	rVV	19418	25345	2.45%	0.287%
56	13.270	1281	1282	1286	rVB3	12920	17989	1.74%	0.204%
57	13.368	1289	1292	1294	rVB3	13519	19445	1.88%	0.220%
58	13.408	1294	1296	1301	rBV2	16230	31974	3.09%	0.362%
59	13.477	1301	1303	1306	rBV	20894	33671	3.25%	0.382%
60	13.526	1306	1308	1310	rBV	42519	54000	5.22%	0.612%
61	13.566	1310	1312	1314	rVB2	15002	20346	1.97%	0.231%
62	13.694	1322	1325	1328	rBV2	54623	86327	8.34%	0.979%
63	13.743	1328	1330	1333	rVB3	14774	22772	2.20%	0.258%
64	13.792	1333	1335	1337	rBV2	15606	23885	2.31%	0.271%
65	13.842	1337	1340	1344	rBV2	70834	155486	15.02%	1.763%
66	13.901	1344	1346	1349	rVB3	20354	33638	3.25%	0.381%
67	13.960	1349	1352	1355	rVB	58332	91307	8.82%	1.035%
68	14.039	1357	1360	1361	rBV2	13507	24046	2.32%	0.273%
69	14.069	1361	1363	1365	rVB	27286	37843	3.66%	0.429%
70	14.138	1365	1370	1373	rBV5	36207	77974	7.53%	0.884%
71	14.207	1375	1377	1383	rVB4	28340	46577	4.50%	0.528%

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72	14.394	1393	1396	1399	rBV2	23191	39674	3.83%	0.450%
73	14.443	1399	1401	1404	rBV2	30002	44927	4.34%	0.509%
74	14.552	1410	1412	1413	rBV2	7727	12292	1.19%	0.139%
75	14.581	1413	1415	1418	rVB2	19725	27702	2.68%	0.314%
76	14.719	1426	1429	1433	rBV	75320	109933	10.62%	1.246%
77	14.848	1439	1442	1449	rVV6	12201	41112	3.97%	0.466%
78	14.976	1453	1455	1458	rBV	40871	57559	5.56%	0.652%
79	15.094	1464	1467	1469	rBV	34806	45589	4.40%	0.517%
80	15.143	1471	1472	1478	rVB5	12890	22499	2.17%	0.255%
81	15.262	1482	1484	1488	rVV3	25480	41995	4.06%	0.476%
82	15.331	1488	1491	1494	rVV3	11864	21548	2.08%	0.244%
83	15.390	1494	1497	1504	rVB2	34906	62862	6.07%	0.713%
84	15.617	1518	1520	1526	rVB3	14582	31290	3.02%	0.355%
85	15.794	1535	1538	1545	rVB4	4888	15393	1.49%	0.174%

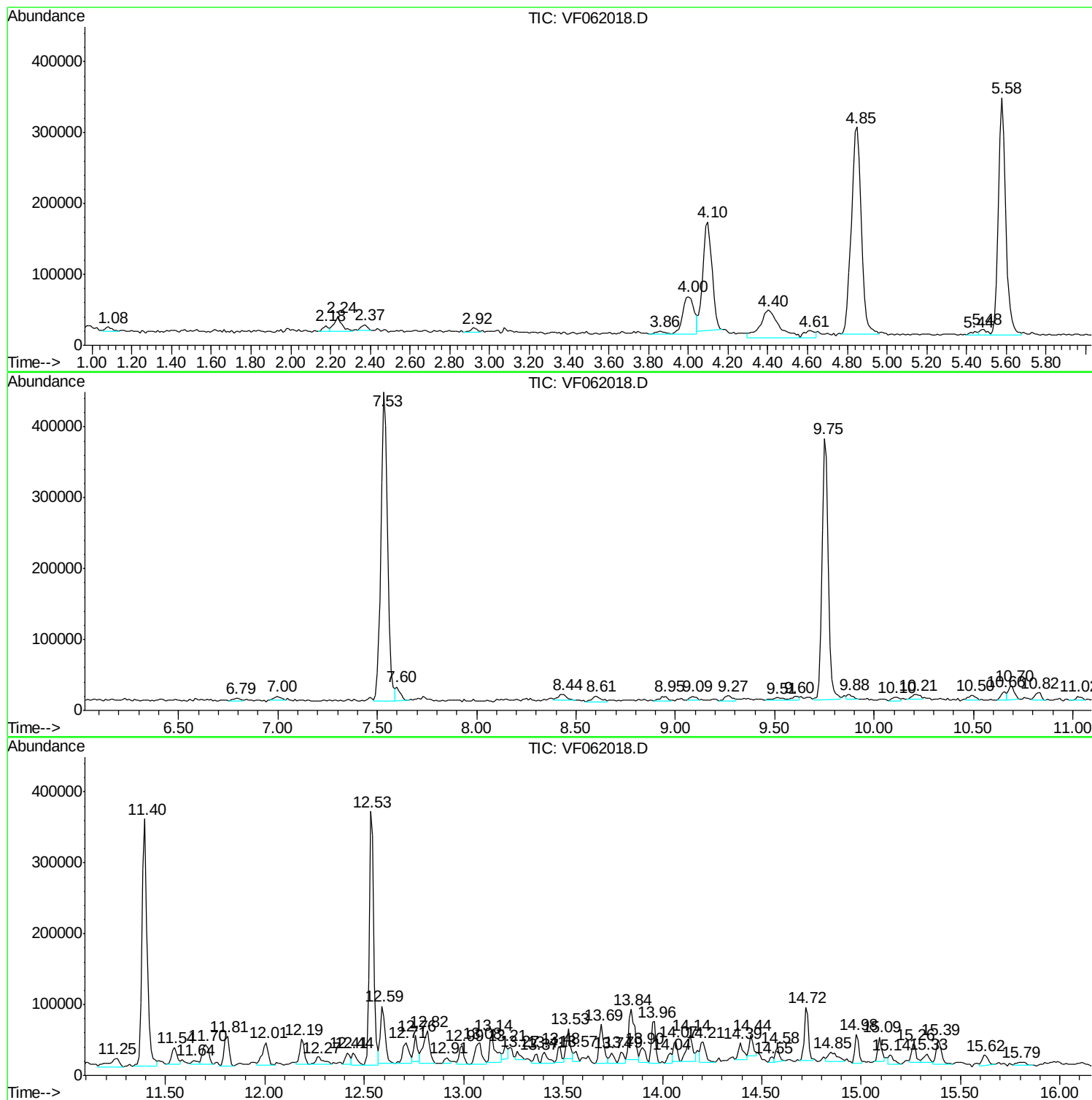
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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



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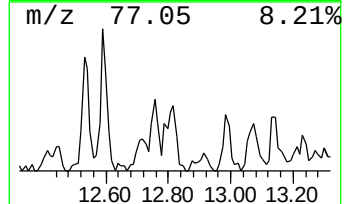
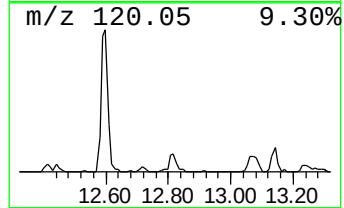
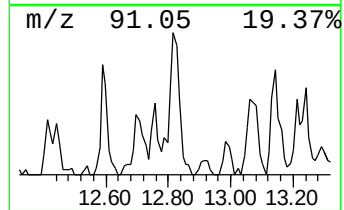
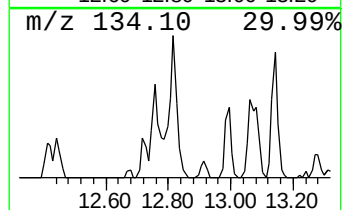
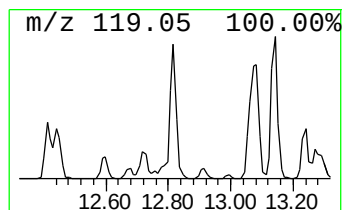
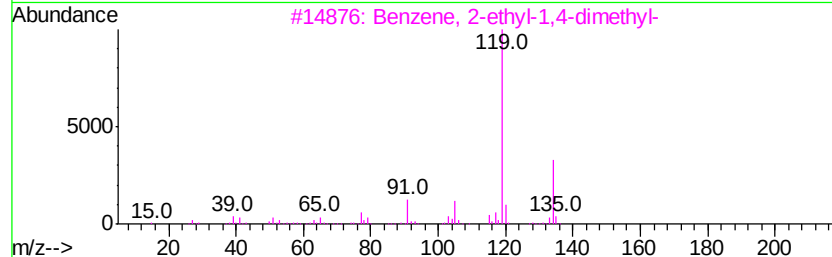
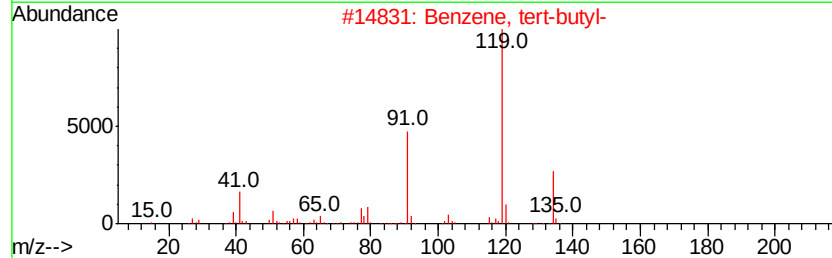
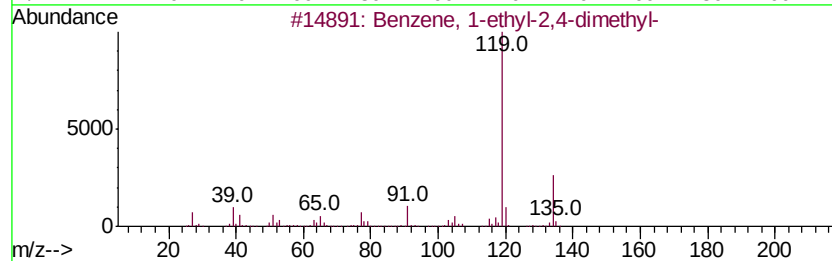
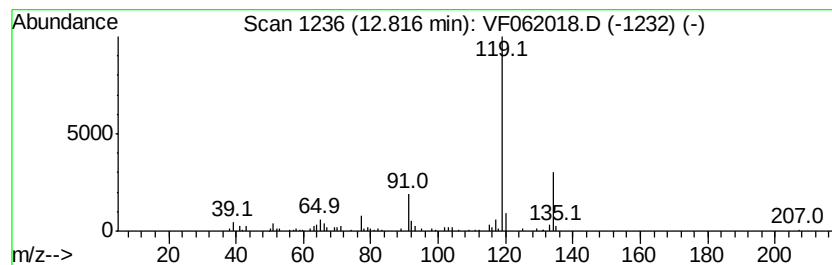
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	8.40 ug/l	97541	1,4-Dichlorobenzene-d4	12.54

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	87
2	Benzene, tert-butyl-	134	C10H14	000098-06-6	83
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	80
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	80
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	80



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Misc : 5.58g/5mL/MSVOA-F/SOIL
ALS Vial : 10 Sample Multiplier: 1

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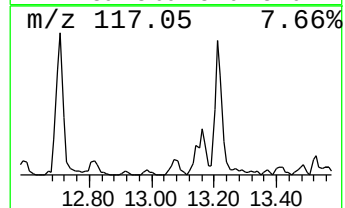
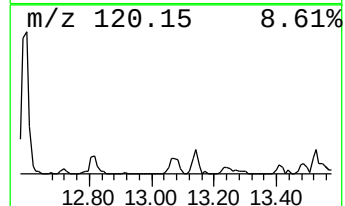
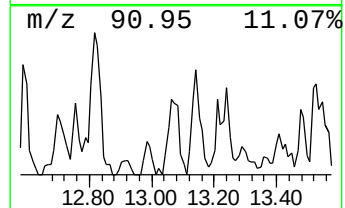
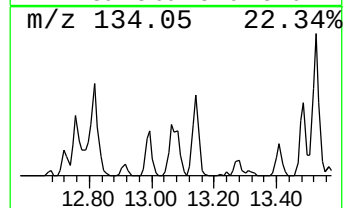
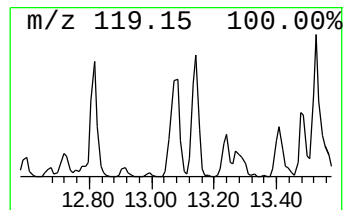
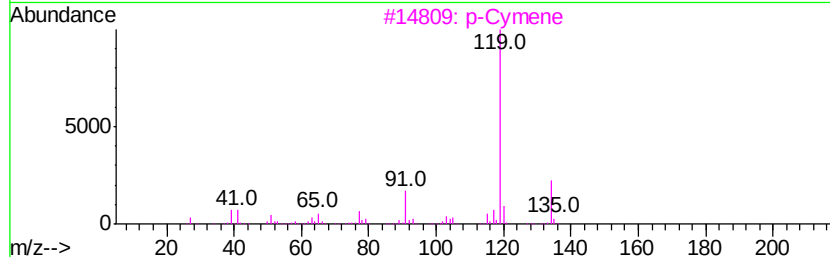
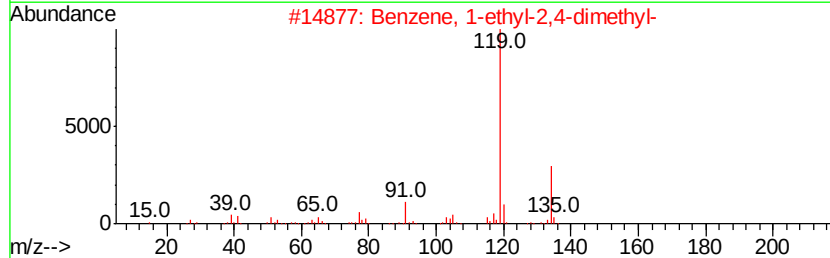
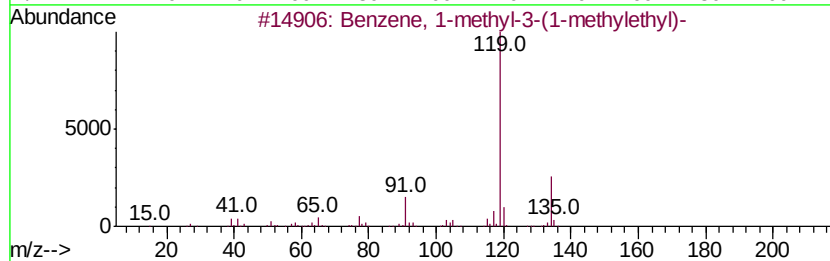
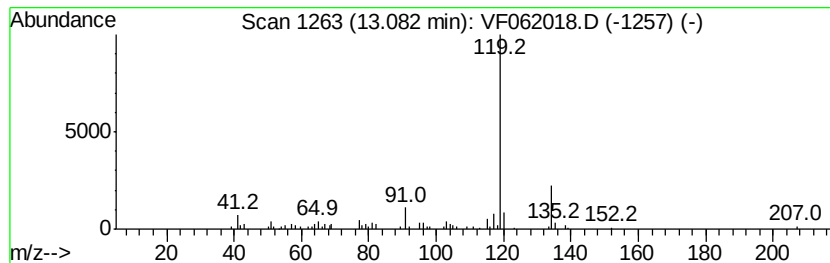
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-methyl-3-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	6.25 ug/l	72600	1,4-Dichlorobenzene-d4	12.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	1-methyl-3-(1-methyleth...	134	C10H14		000535-77-3	95
2	Benzene,	1-ethyl-2,4-dimethyl-	134	C10H14		000874-41-9	94
3	p-Cymene		134	C10H14		000099-87-6	91
4	Benzene,	1,2,3,4-tetramethyl-	134	C10H14		000488-23-3	91
5	Benzene,	2-ethyl-1,3-dimethyl-	134	C10H14		002870-04-4	91



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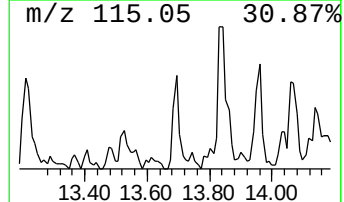
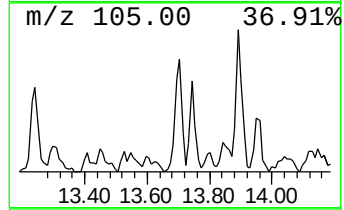
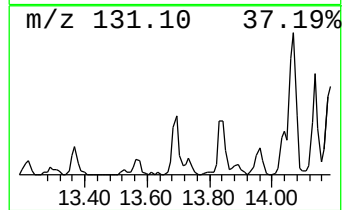
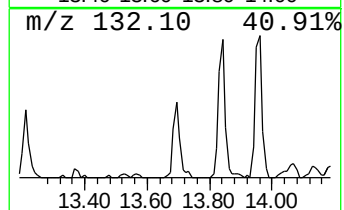
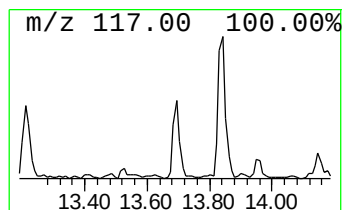
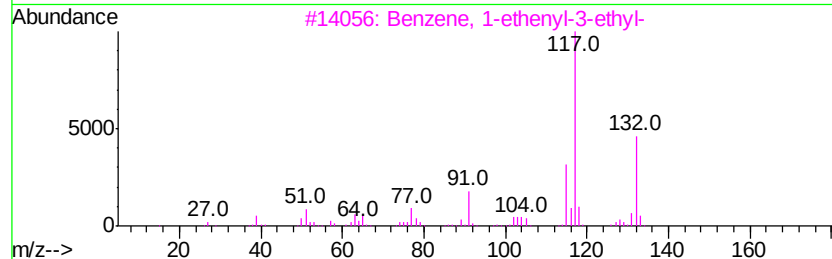
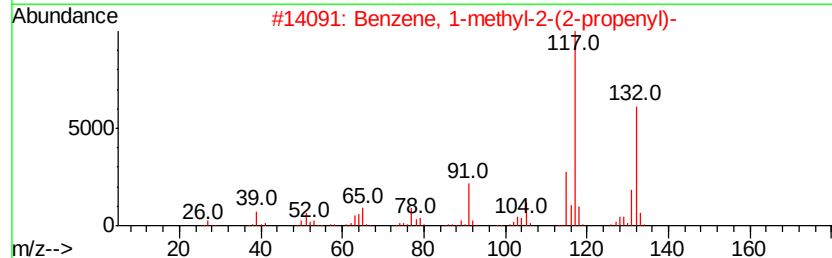
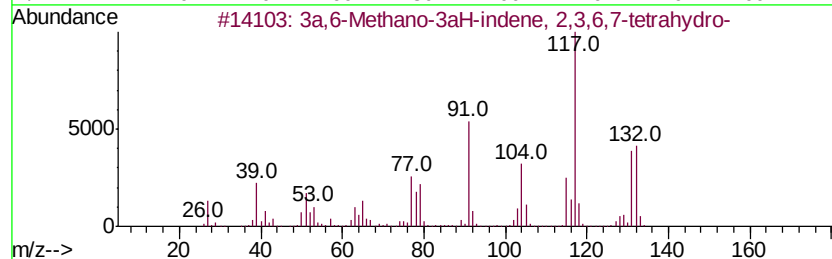
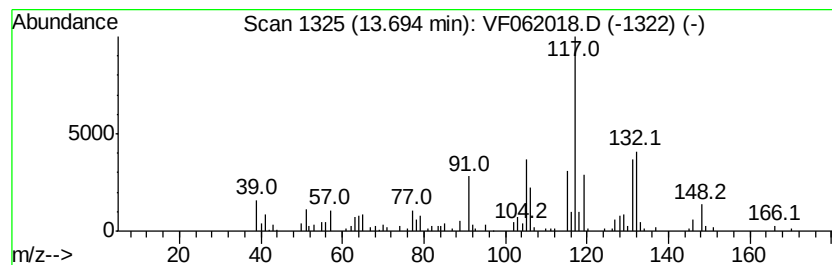
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Quant Title : SW846 8260

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

Peak Number 6 3a,6-Methano-3aH-indene, 2,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	7.43 ug/l	86327	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3a,6-Methano-3aH-indene, 2,3,6,7...	132	C10H12	098640-29-0	70
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	64



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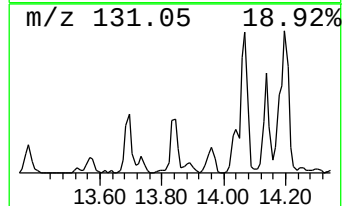
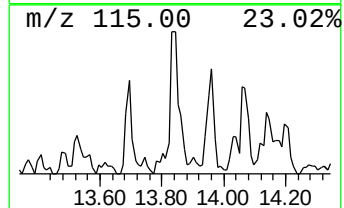
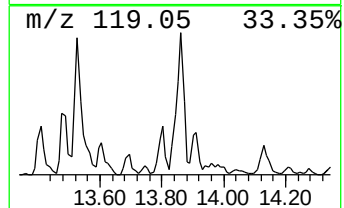
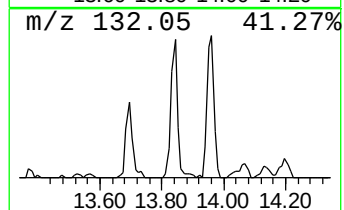
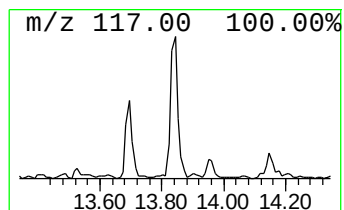
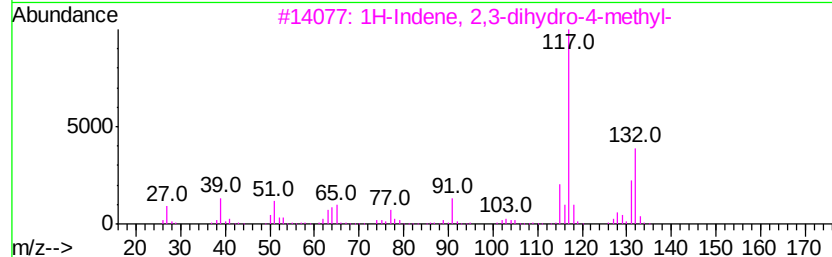
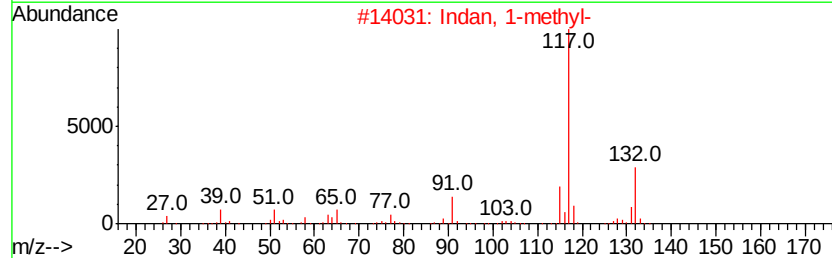
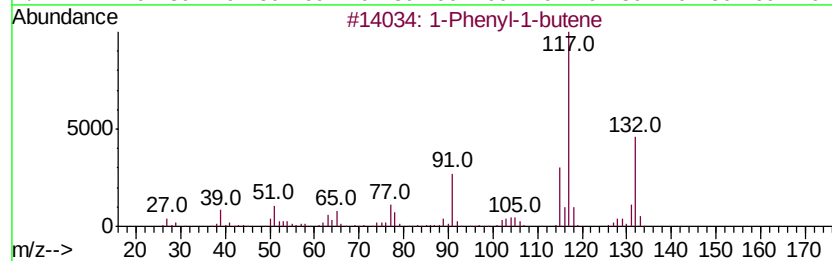
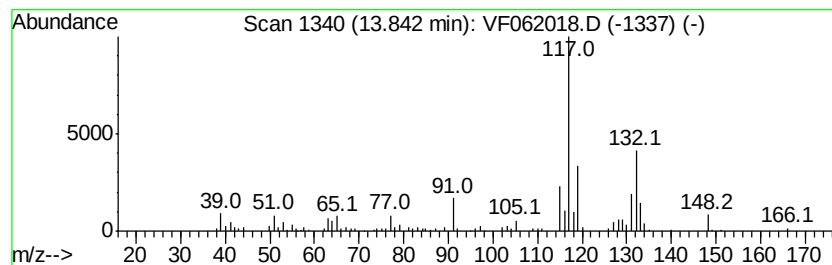
Instrument :
MSVOA_F
ClientSampleId :
FK-GF-03-006-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	13.39 ug/l	155486	1,4-Dichlorobenzene-d4	12.54
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132 C10H12	000824-90-8	81
2	Indan, 1-methyl-	132 C10H12	000767-58-8	81
3	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	74
4	Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7	74
5	1H-Indene, 2,3-dihydro-2-methyl-	132 C10H12	000824-63-5	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
Data File : VF062018.D
Acq On : 19 Mar 2019 15:14
Operator : VA/AP
Sample : K1951-04
Misc : 5.58g/5mL/MSVOA-F/SOIL
ALS Vial : 10 Sample Multiplier: 1

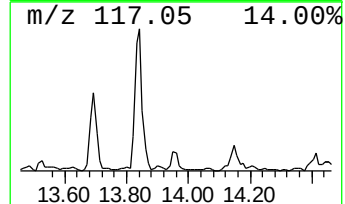
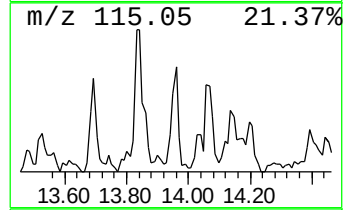
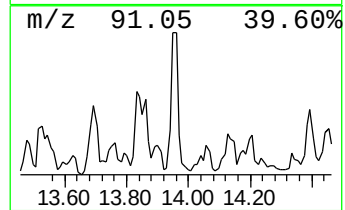
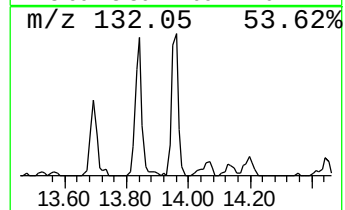
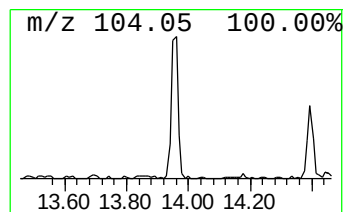
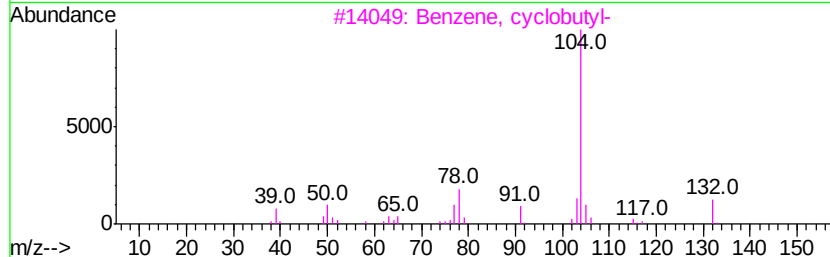
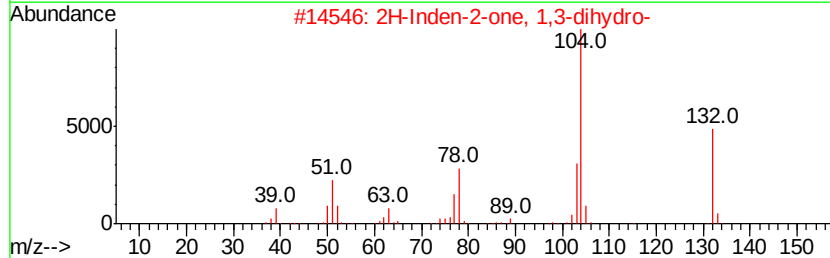
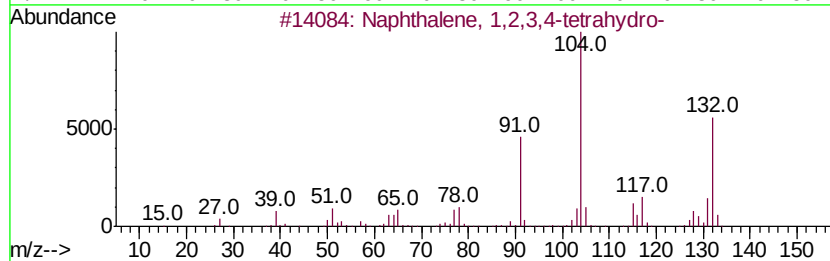
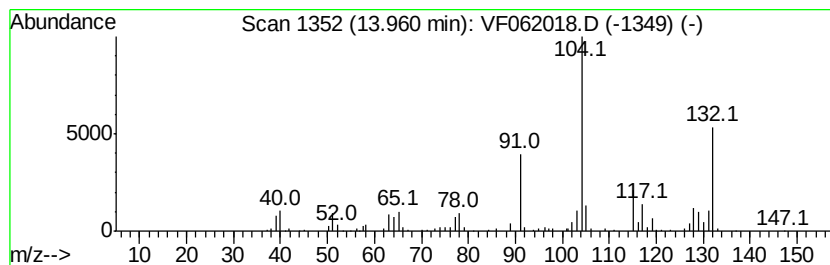
Instrument :
MSVOA_F
ClientSampled :
FK-GF-03-006-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.96	7.86 ug/l	91307	1,4-Dichlorobenzene-d4	12.54	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	47
3	Benzene, cyclobutyl-	132	C10H12	004392-30-7	47
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	35
5	Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
 Data File : VF062018.D
 Acq On : 19 Mar 2019 15:14
 Operator : VA/AP
 Sample : K1951-04
 Misc : 5.58g/5mL/MSVOA-F/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 FK-GF-03-006-B

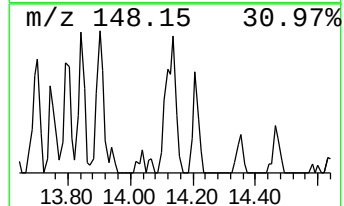
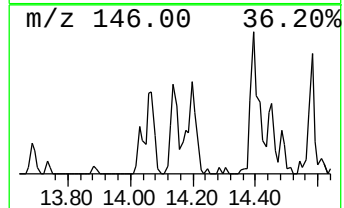
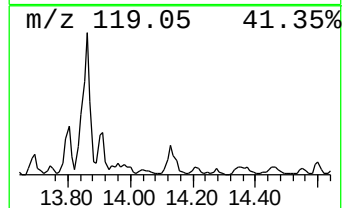
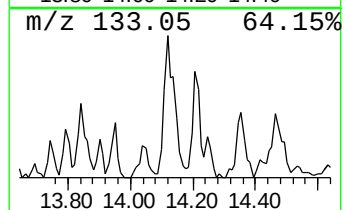
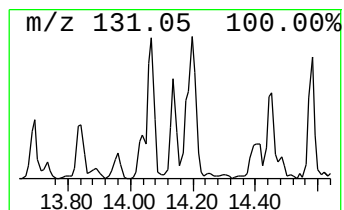
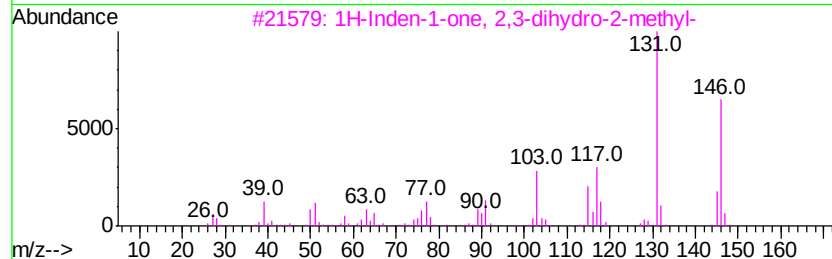
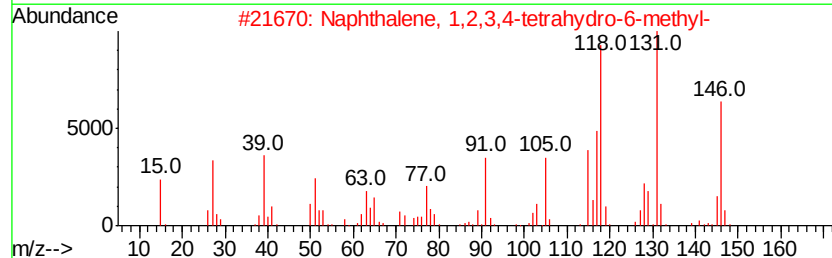
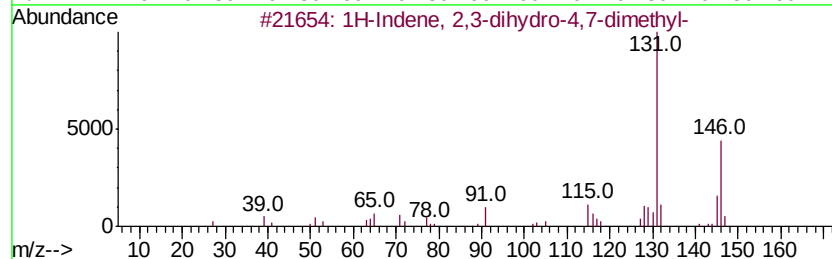
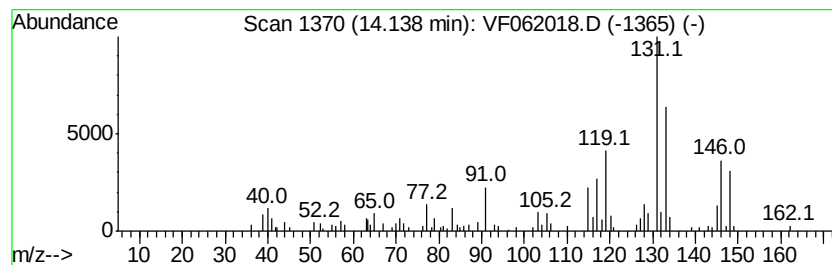
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
 Quant Title : SW846 8260

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

 Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	6.71 ug/l	77974	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	55
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	46
3		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	45
4		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	43
5		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
Data File : VF062018.D
Acq On : 19 Mar 2019 15:14
Operator : VA/AP
Sample : K1951-04
Misc : 5.58g/5mL/MSVOA-F/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
FK-GF-03-006-B

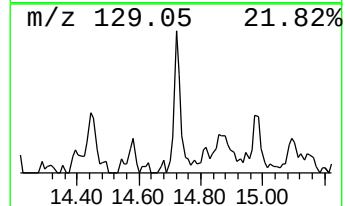
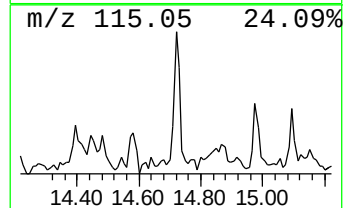
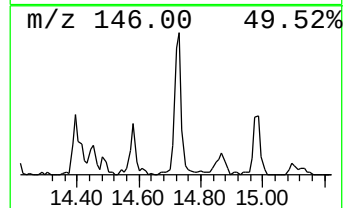
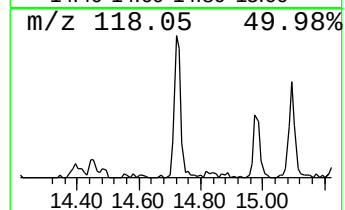
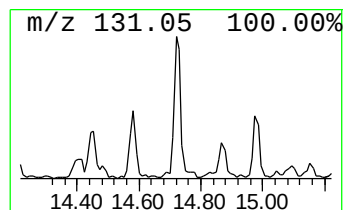
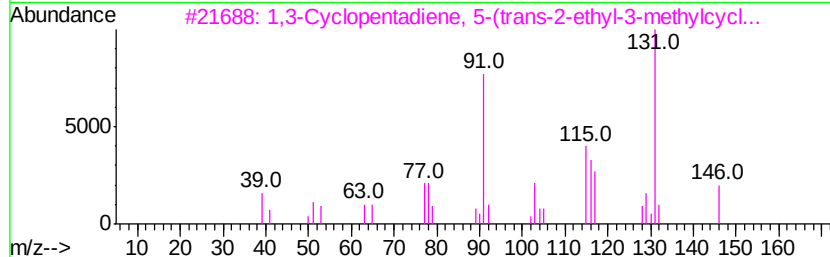
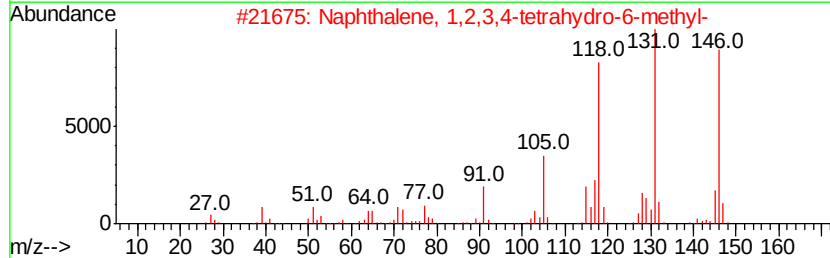
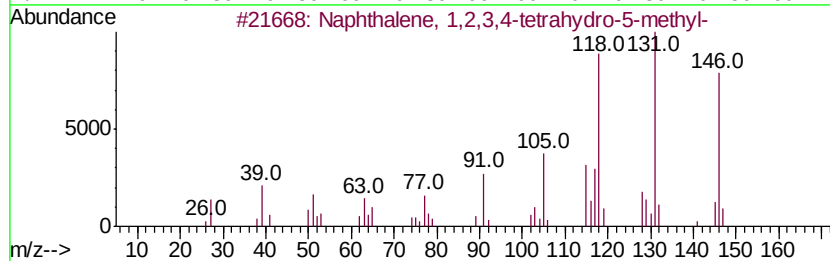
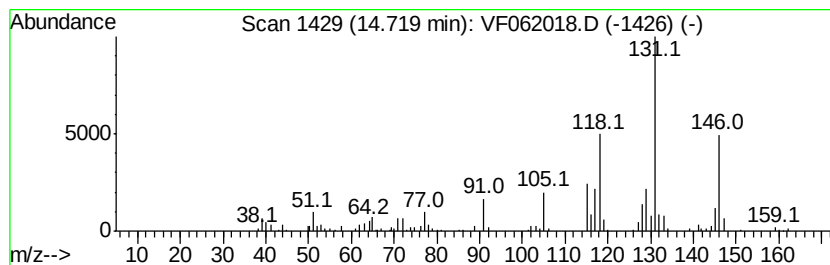
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	9.47 ug/l	109933	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	92
3		1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	62
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60
5		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF031919\
Data File : VF062018.D
Acq On : 19 Mar 2019 15:14
Operator : VA/AP
Sample : K1951-04
Misc : 5.58g/5mL/MSVOA-F/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
FK-GF-03-006-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.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, 1-ethyl-...	12.82	8.4	ug/l	97541	4	12.54	580669	50.0
Benzene, 1-methyl...	13.08	6.3	ug/l	72600	4	12.54	580669	50.0
3a,6-Methano-3aH-...	13.69	7.4	ug/l	86327	4	12.54	580669	50.0
1-Phenyl-1-butene	13.84	13.4	ug/l	155486	4	12.54	580669	50.0
Naphthalene, 1,2,...	13.96	7.9	ug/l	91307	4	12.54	580669	50.0
1H-Indene, 2,3-di...	14.14	6.7	ug/l	77974	4	12.54	580669	50.0
Naphthalene, 1,2,...	14.72	9.5	ug/l	109933	4	12.54	580669	50.0