

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF060619\
 Data File : VF062887.D
 Acq On : 6 Jun 2019 17:23
 Operator : FY/SY
 Sample : K3160-01
 Misc : 6.35g/5mL/MSVOA F/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 EO-02-060619-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_F\METHODS\82F053019S.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	1.124	57	60	68	rBV3	31042	97530	3.67%	0.264%
2	2.356	181	185	192	rVB4	40268	133956	5.04%	0.362%
3	2.504	198	200	204	rBV4	13651	32545	1.23%	0.088%
4	3.086	256	259	263	rBV6	8744	31817	1.20%	0.086%
5	3.135	263	264	271	rVB2	28040	60618	2.28%	0.164%
6	4.239	368	376	378	rBV	16545	45844	1.73%	0.124%
7	4.328	379	385	397	rVB2	346960	1045500	39.37%	2.827%
8	4.840	433	437	443	rVB4	14037	45900	1.73%	0.124%
9	5.067	450	460	473	rBV2	791793	2617063	98.56%	7.076%
10	5.797	528	534	543	rBV	992632	2644318	99.58%	7.149%
11	7.482	699	705	708	rVB7	10583	28401	1.07%	0.077%
12	7.729	724	730	734	rBV	818271	1854228	69.83%	5.013%
13	7.798	734	737	742	rVB	95152	213359	8.04%	0.577%
14	9.927	947	953	962	rBV	1184786	2655338	100.00%	7.179%
15	10.045	963	965	970	rVB5	14133	28402	1.07%	0.077%
16	10.282	985	989	993	rVB4	14193	33141	1.25%	0.090%
17	10.824	1042	1044	1048	rVB3	16398	33559	1.26%	0.091%
18	11.218	1081	1084	1086	rVV4	20147	39342	1.48%	0.106%
19	11.248	1086	1087	1092	rVB5	16605	28081	1.06%	0.076%
20	11.366	1096	1099	1102	rVV3	22291	37496	1.41%	0.101%
21	11.504	1109	1113	1124	rBV	659649	1126841	42.44%	3.047%
22	11.642	1124	1127	1129	rVV3	33252	58247	2.19%	0.157%
23	11.681	1129	1131	1134	rVV2	20785	34096	1.28%	0.092%
24	11.740	1134	1137	1139	rVV	88067	127610	4.81%	0.345%
25	11.800	1139	1143	1147	rVV	55818	124647	4.69%	0.337%
26	11.908	1150	1154	1157	rVB2	48361	80682	3.04%	0.218%
27	12.066	1167	1170	1172	rBV3	64956	101307	3.82%	0.274%
28	12.095	1172	1173	1176	rVB	49366	55473	2.09%	0.150%
29	12.214	1183	1185	1188	rBV3	18535	36940	1.39%	0.100%
30	12.273	1188	1191	1196	rVB	179994	298808	11.25%	0.808%
31	12.371	1197	1201	1205	rBV5	18886	48764	1.84%	0.132%
32	12.440	1205	1208	1211	rBV3	47449	98923	3.73%	0.267%
33	12.490	1211	1213	1215	rVV3	55935	106503	4.01%	0.288%
34	12.519	1215	1216	1218	rVV2	59006	71064	2.68%	0.192%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF060619\
 Data File : VF062887.D
 Acq On : 6 Jun 2019 17:23
 Operator : FY/SY
 Sample : K3160-01
 Misc : 6.35g/5mL/MSVOA F/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 EO-02-060619-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_F\METHODS\82F053019S.M
 Title : SW846 8260

35	12.618	1222	1226	1230	rBV	1212285	1965476	74.02%	5.314%
36	12.677	1230	1232	1235	rVV	203560	268743	10.12%	0.727%
37	12.785	1240	1243	1246	rBV2	152918	285457	10.75%	0.772%
38	12.835	1246	1248	1250	rVV	159856	267882	10.09%	0.724%
39	12.874	1250	1252	1259	rVB3	293081	706277	26.60%	1.910%
40	12.992	1261	1264	1266	rBV2	63712	110720	4.17%	0.299%
41	13.061	1269	1271	1275	rVB	148579	231571	8.72%	0.626%
42	13.150	1275	1280	1284	rBV3	198613	605705	22.81%	1.638%
43	13.209	1284	1286	1288	rVV	213562	294969	11.11%	0.797%
44	13.259	1289	1291	1292	rVV	161494	219582	8.27%	0.594%
45	13.288	1292	1294	1298	rVV3	161013	390272	14.70%	1.055%
46	13.347	1298	1300	1303	rVV2	102071	199421	7.51%	0.539%
47	13.387	1303	1304	1307	rVV2	59980	84004	3.16%	0.227%
48	13.436	1307	1309	1311	rVV	151577	174303	6.56%	0.471%
49	13.475	1311	1313	1314	rVV2	97590	136283	5.13%	0.368%
50	13.544	1317	1320	1322	rVV2	152344	244115	9.19%	0.660%
51	13.594	1322	1325	1327	rVV	256943	401832	15.13%	1.086%
52	13.633	1327	1329	1331	rVV3	107178	165785	6.24%	0.448%
53	13.663	1331	1332	1333	rVV	65657	60080	2.26%	0.162%
54	13.692	1333	1335	1338	rVB2	60901	68272	2.57%	0.185%
55	13.761	1338	1342	1345	rBV2	662049	1114265	41.96%	3.013%
56	13.811	1345	1347	1350	rVV3	152613	278795	10.50%	0.754%
57	13.860	1350	1352	1353	rVV	150440	183744	6.92%	0.497%
58	13.899	1353	1356	1360	rVV3	797777	1502078	56.57%	4.061%
59	13.968	1360	1363	1365	rVV2	213270	359091	13.52%	0.971%
60	14.018	1365	1368	1371	rVV	498757	678585	25.56%	1.835%
61	14.126	1373	1379	1382	rVV	238655	513822	19.35%	1.389%
62	14.195	1382	1386	1389	rVV4	364545	829437	31.24%	2.242%
63	14.264	1391	1393	1399	rVV4	351880	599292	22.57%	1.620%
64	14.343	1399	1401	1403	rVV3	96353	128314	4.83%	0.347%
65	14.402	1404	1407	1408	rVV3	66456	121819	4.59%	0.329%
66	14.451	1409	1412	1415	rVV3	317566	615633	23.18%	1.664%
67	14.501	1415	1417	1425	rVV4	433424	1073748	40.44%	2.903%
68	14.599	1425	1427	1429	rVV3	121120	186179	7.01%	0.503%
69	14.639	1429	1431	1434	rVV2	293083	434002	16.34%	1.173%
70	14.688	1434	1436	1438	rVV3	41373	67530	2.54%	0.183%
71	14.777	1442	1445	1450	rVB	970550	1373649	51.73%	3.714%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF060619\
 Data File : VF062887.D
 Acq On : 6 Jun 2019 17:23
 Operator : FY/SY
 Sample : K3160-01
 Misc : 6.35g/5mL/MSVOA F/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 EO-02-060619-A

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
 Title : SW846 8260

72	14.865	1452	1454	1455	rBV	84445	98916	3.73%	0.267%
73	14.915	1455	1459	1461	rVV3	162992	382140	14.39%	1.033%
74	14.974	1464	1465	1468	rVB2	82899	113150	4.26%	0.306%
75	15.033	1468	1471	1474	rBV2	518895	811885	30.58%	2.195%
76	15.151	1479	1483	1485	rVV	582339	1022587	38.51%	2.765%
77	15.200	1485	1488	1492	rVV4	243979	589450	22.20%	1.594%
78	15.279	1494	1496	1497	rVV	93686	139868	5.27%	0.378%
79	15.319	1497	1500	1503	rVV2	263369	542324	20.42%	1.466%
80	15.388	1503	1507	1510	rVV3	246546	657102	24.75%	1.777%
81	15.447	1510	1513	1518	rVV	156006	338179	12.74%	0.914%
82	15.536	1520	1522	1523	rVV2	49539	79016	2.98%	0.214%
83	15.565	1523	1525	1527	rVV2	67239	123841	4.66%	0.335%
84	15.634	1530	1532	1533	rVV2	42344	59429	2.24%	0.161%
85	15.683	1533	1537	1542	rVV	221653	499451	18.81%	1.350%
86	15.752	1542	1544	1549	rVV4	53577	113787	4.29%	0.308%
87	15.831	1549	1552	1553	rVV3	28643	51772	1.95%	0.140%
88	15.861	1553	1555	1558	rVB3	45686	79111	2.98%	0.214%
89	15.979	1564	1567	1569	rBV3	45533	98832	3.72%	0.267%
90	16.028	1569	1572	1578	rVB5	56018	169263	6.37%	0.458%

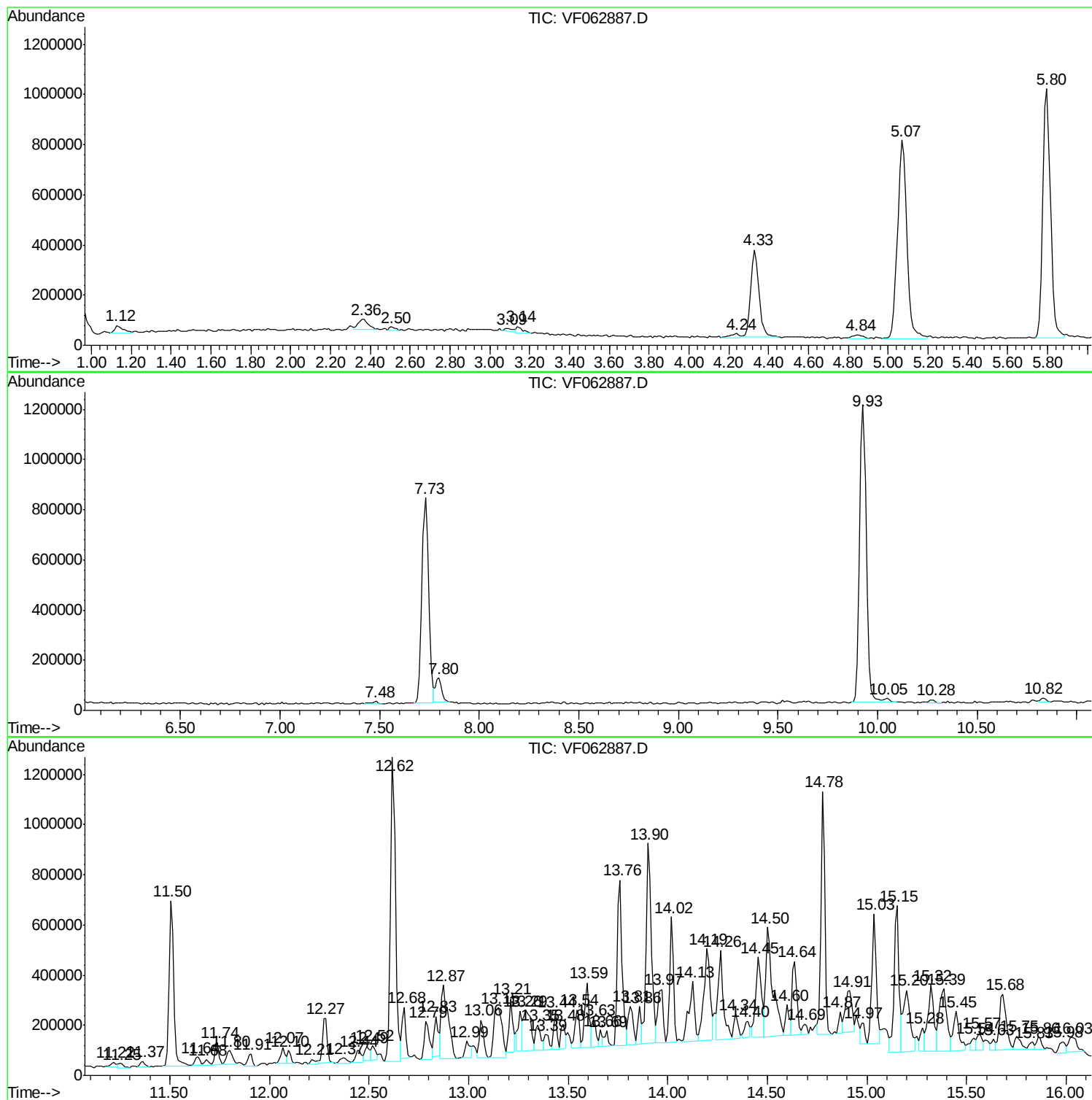
Sum of corrected areas: 36987178

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

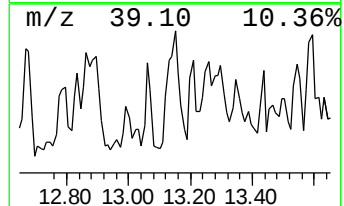
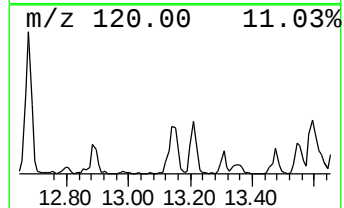
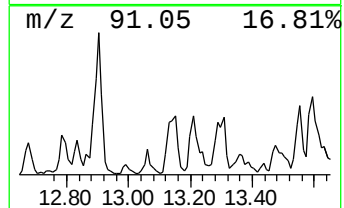
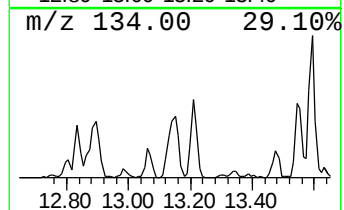
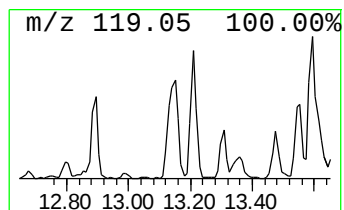
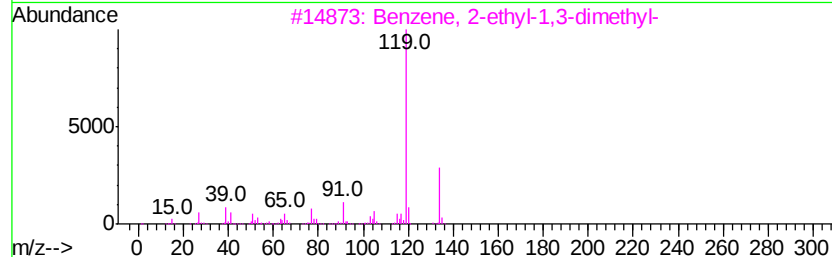
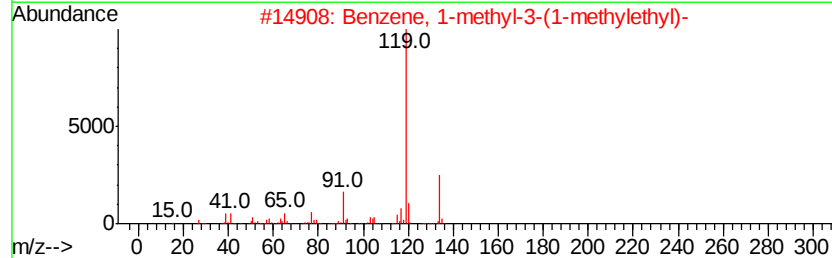
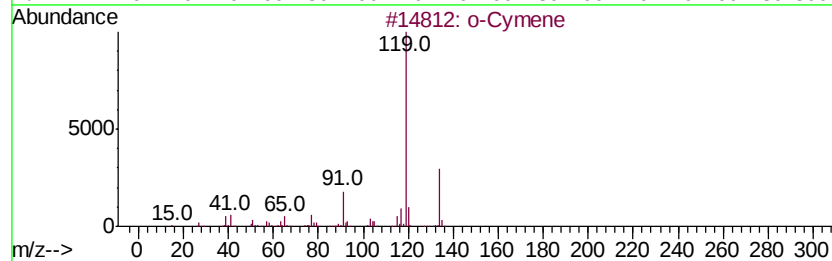
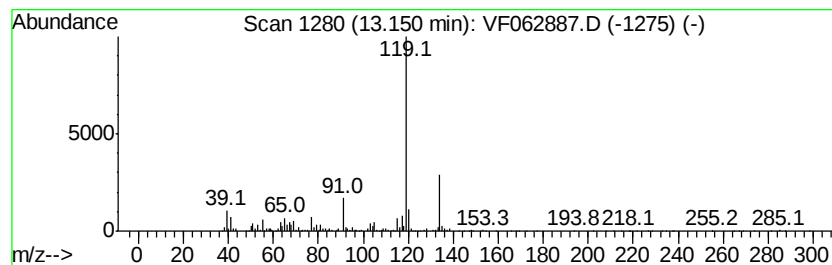
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

Peak Number 1 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	15.41 ug/l	605705	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	95
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4		p-Cymene	134	C10H14	000099-87-6	94
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

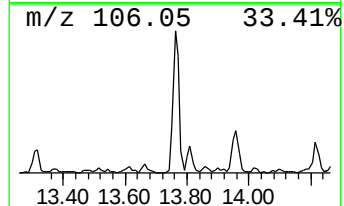
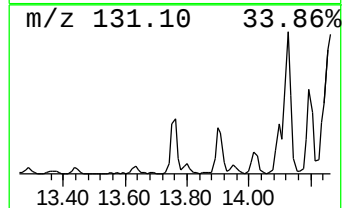
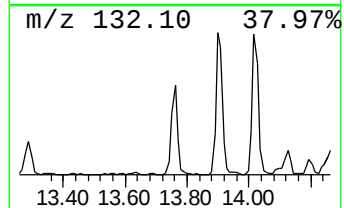
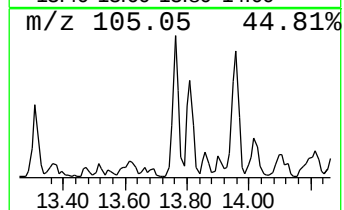
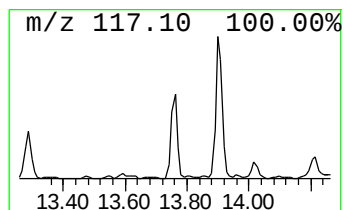
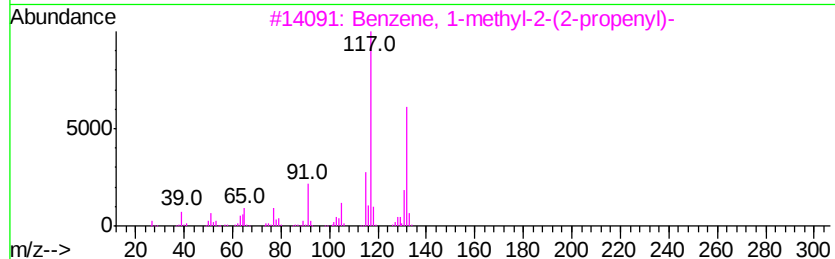
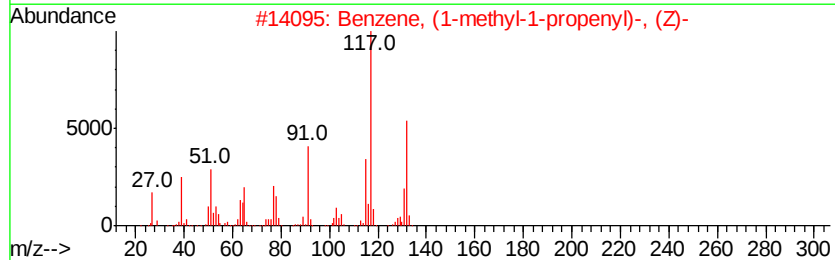
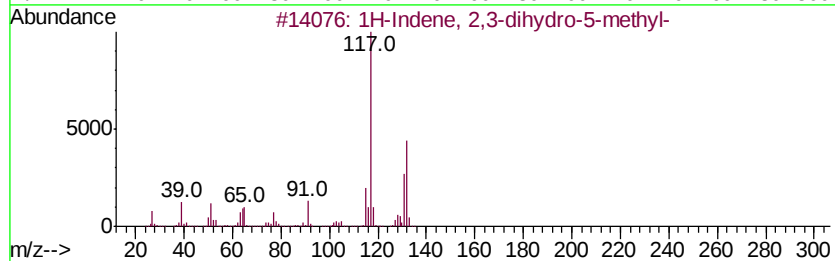
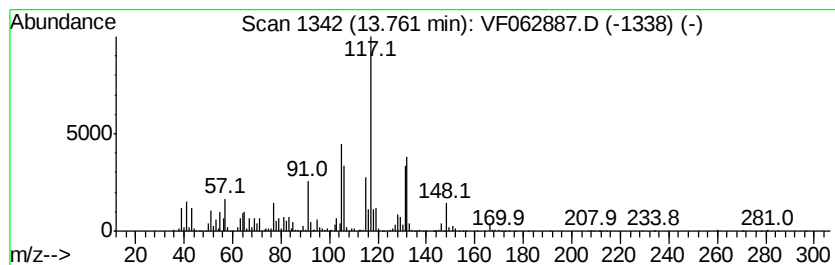
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	28.35 ug/l	1114270	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	89
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5		3a,6-Methano-3aH-indene, 2,3,6,7...	132	C10H12	098640-29-0	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

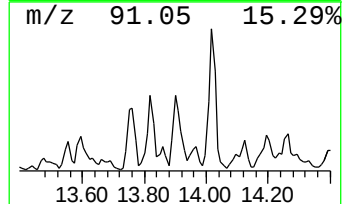
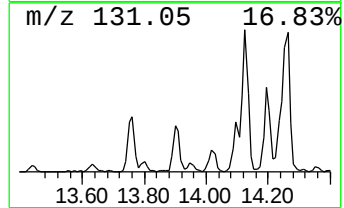
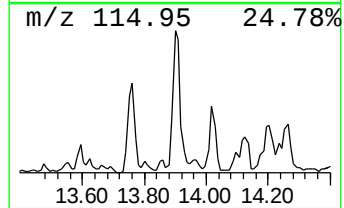
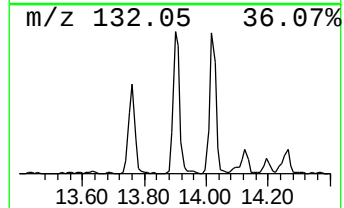
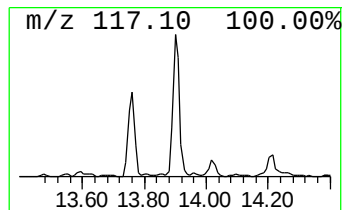
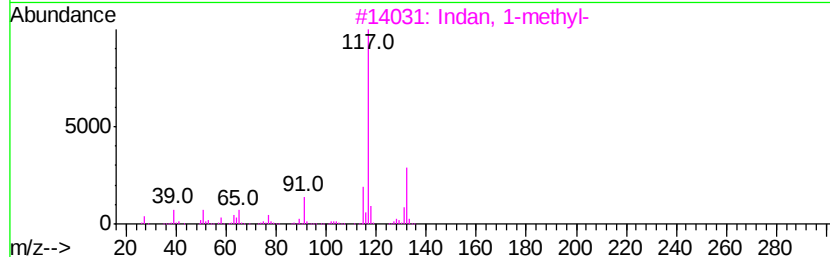
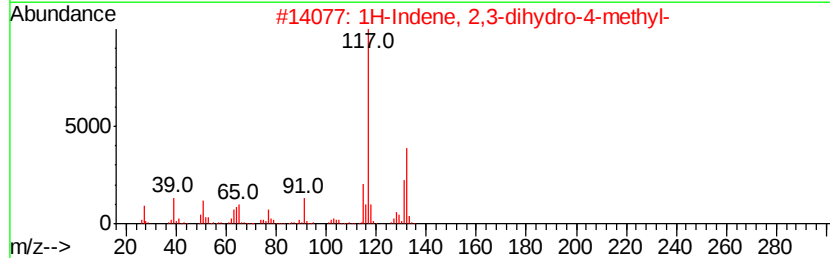
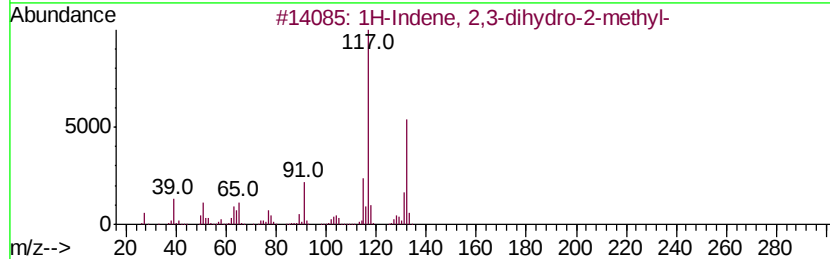
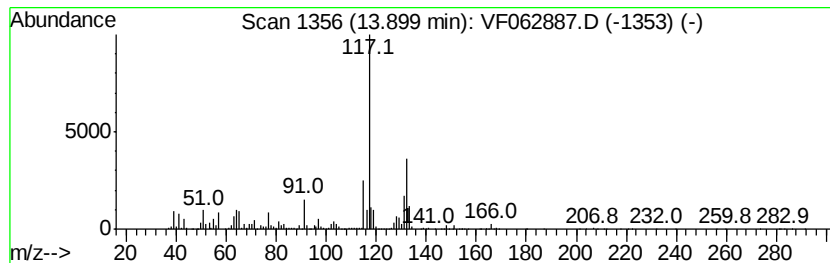
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	38.21 ug/l	1502080	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3		Indan, 1-methyl-	132	C10H12	000767-58-8	81
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

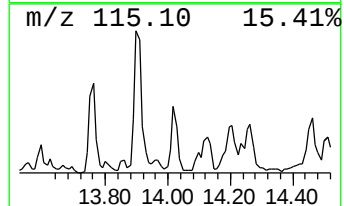
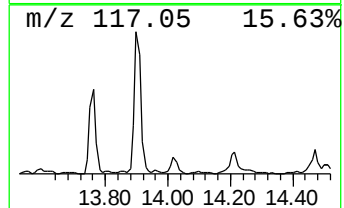
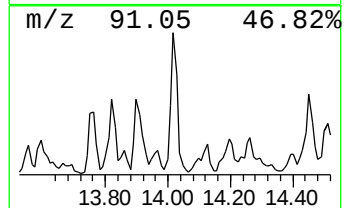
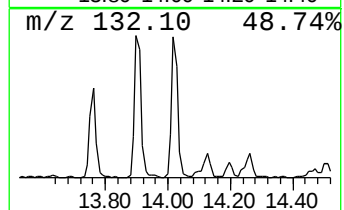
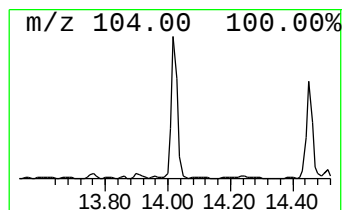
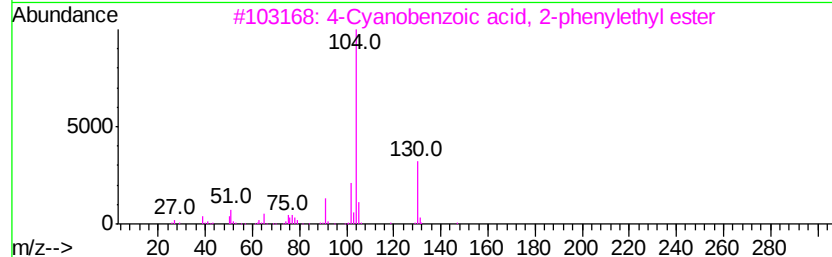
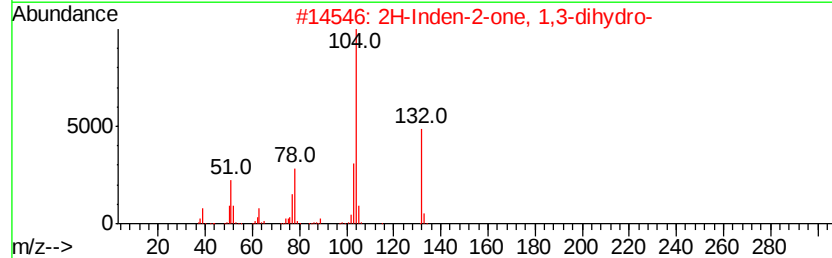
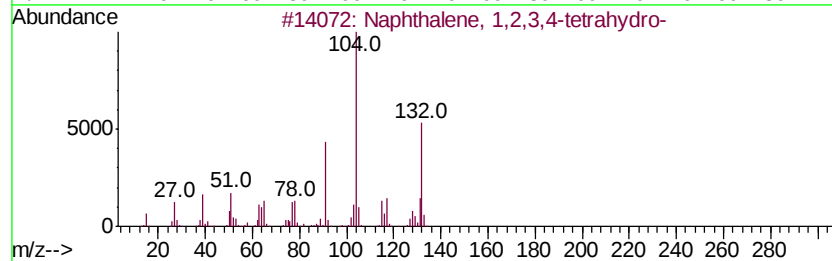
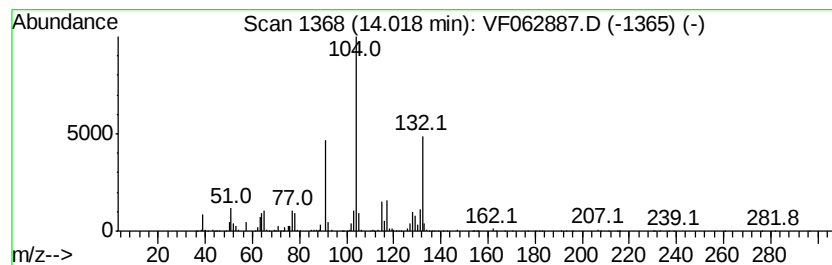
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	17.26 ug/l	678585	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
3		4-Cyanobenzoic acid, 2-phenylethyl...	251	C16H13NO2	131786-46-4	38
4		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	35
5		2-Furanone, 4-phenyltetrahydro	162	C10H10O2	1000197-38-4	32



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

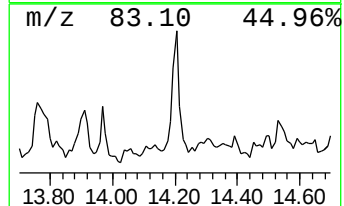
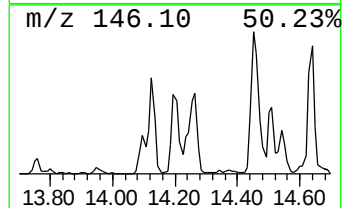
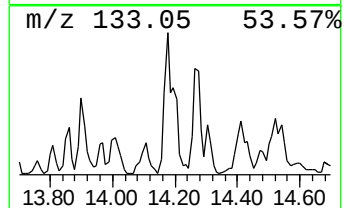
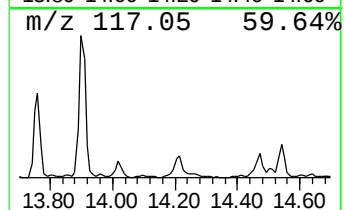
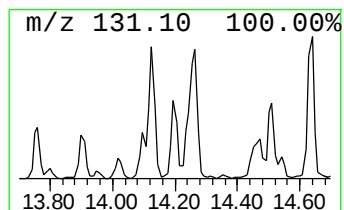
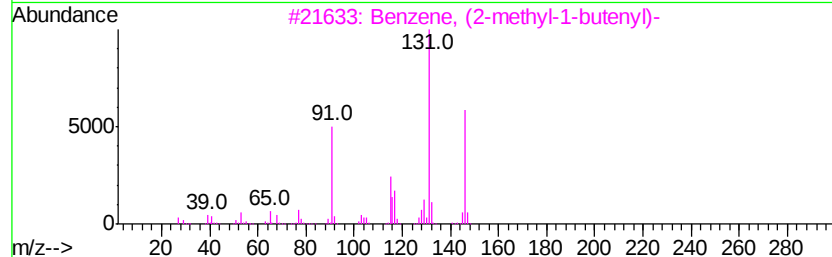
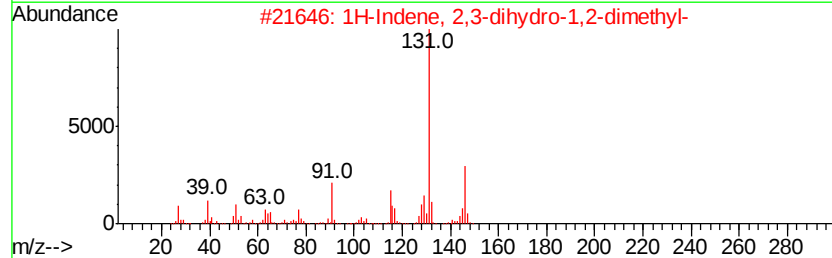
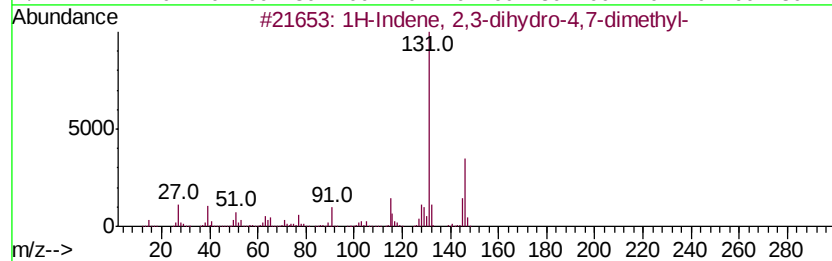
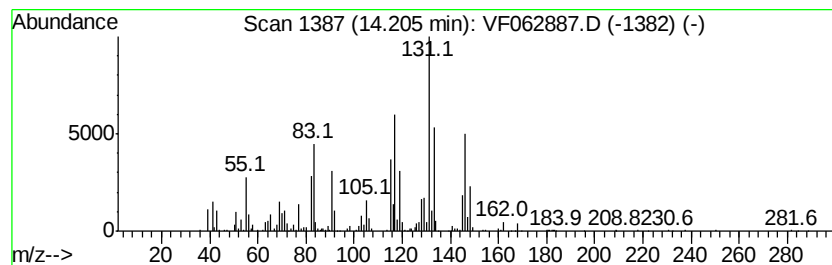
Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	21.10 ug/l	829437	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	90
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	78
3	Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6	78
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	60
5	Benzene, 1-methyl-3-(1-methyl-2-...	146 C11H14	052161-57-6	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

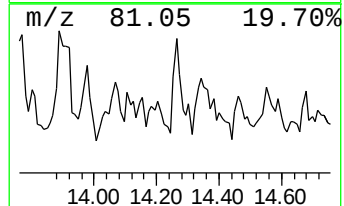
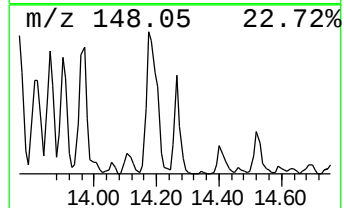
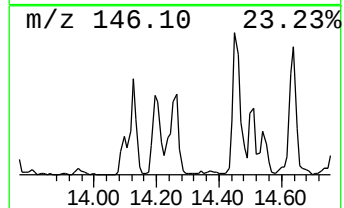
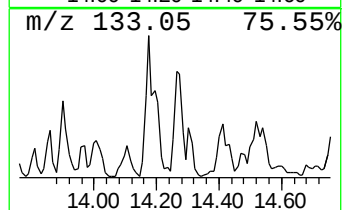
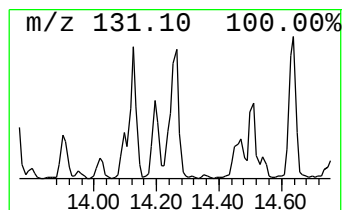
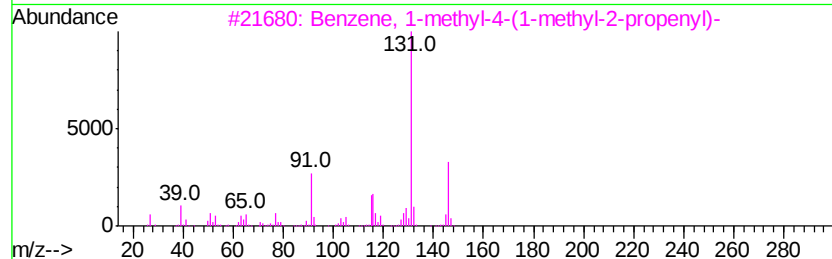
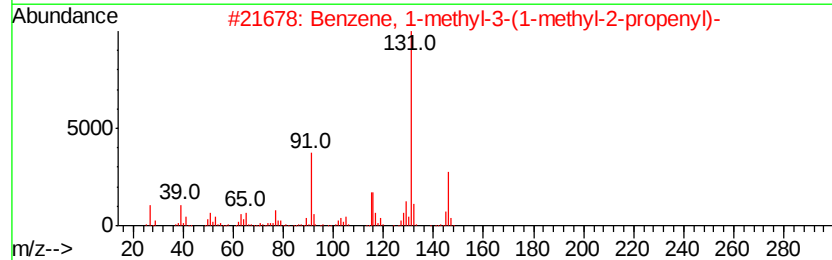
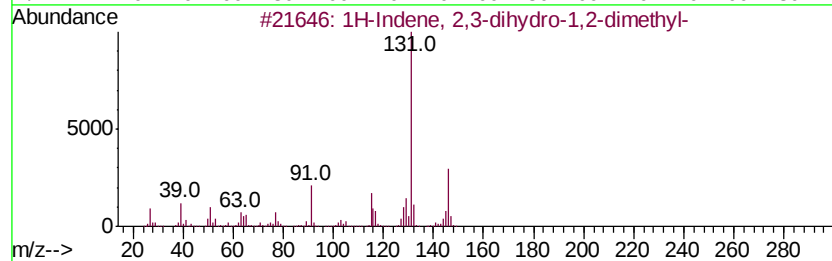
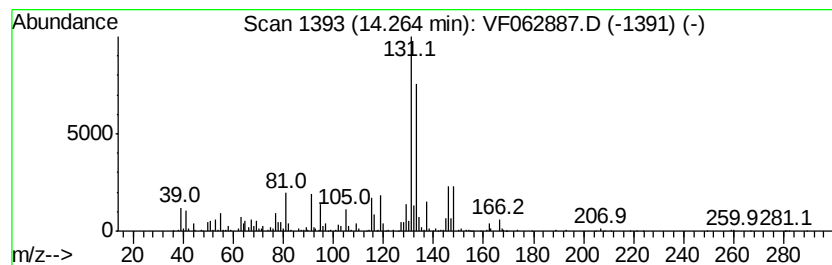
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

Peak Number 6 unknown14.26 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	15.25 ug/l	599292	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	46
2		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	45
3		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	43
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	41
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	41



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

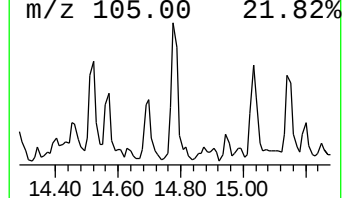
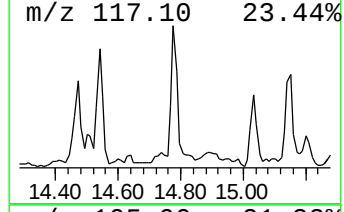
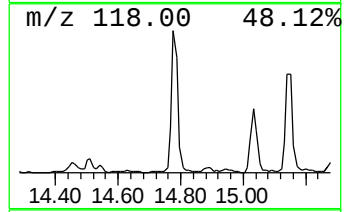
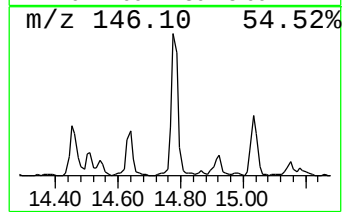
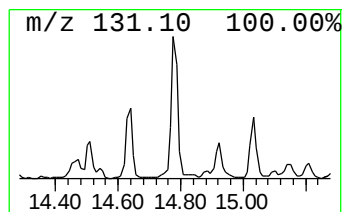
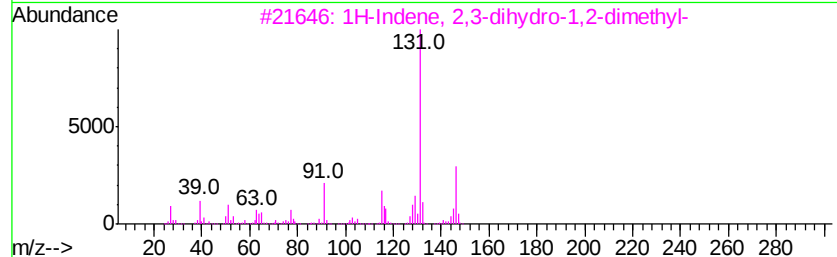
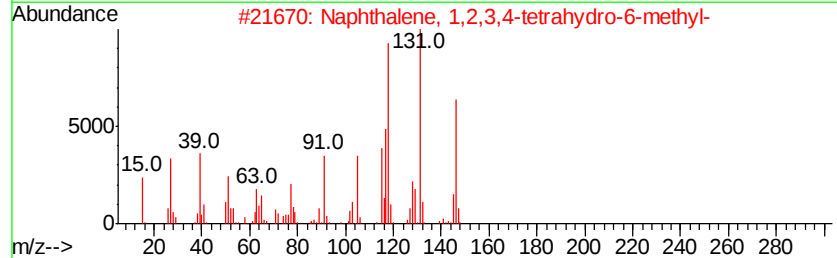
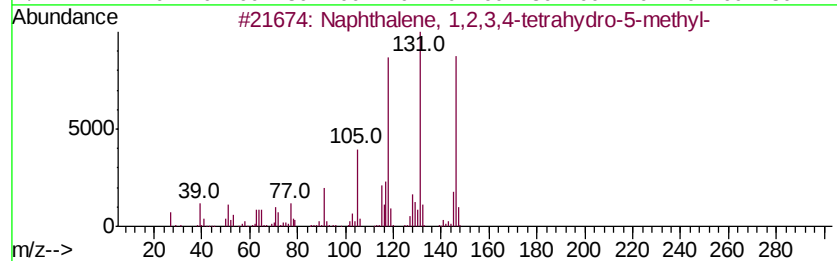
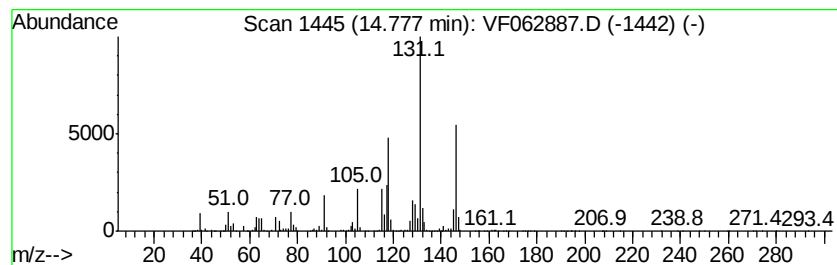
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	34.94 ug/l	1373650	1,4-Dichlorobenzene-d4	12.62

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	94
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	83
4		Benzene, (2-methyl-1-butenvl)-	146	C11H14	056253-64-6	70
5		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

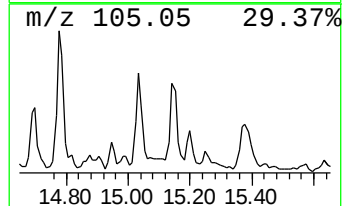
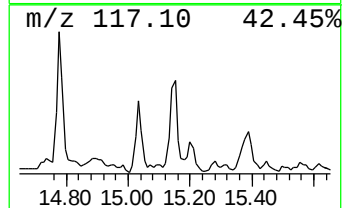
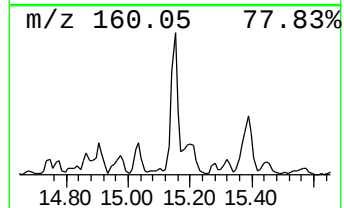
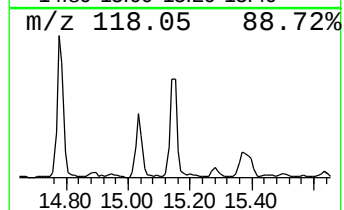
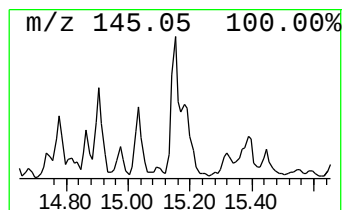
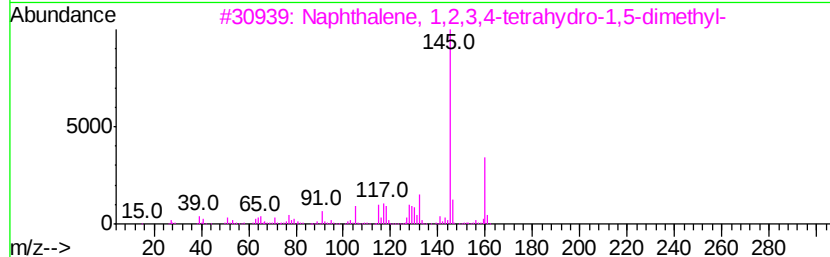
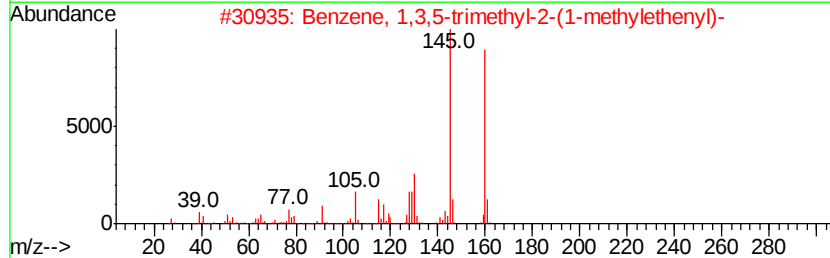
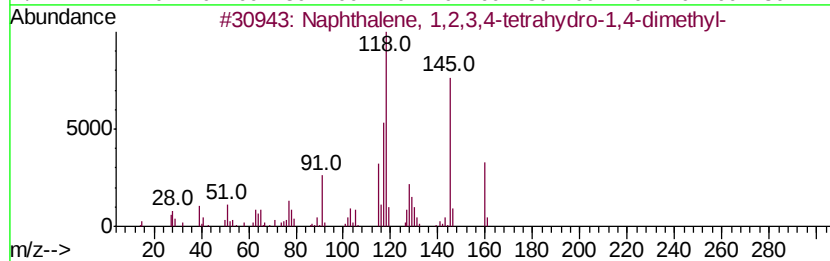
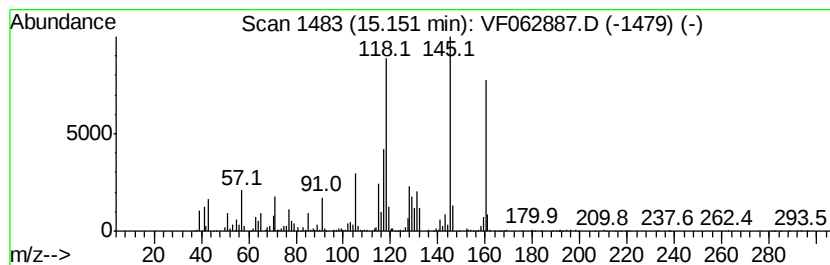
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.15	26.01 ug/l	1022590	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2		Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	78
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	68
4		Benzene, (1,3-dimethyl-2-butenyl)-	160	C12H16	050704-01-3	64
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA F/SOIL
ALS Vial : 15 Sample Multiplier: 1

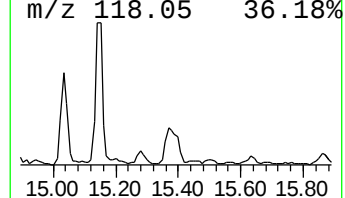
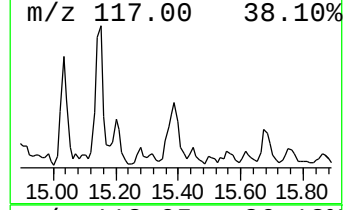
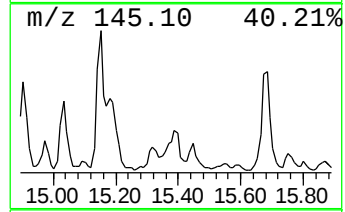
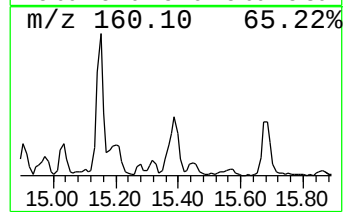
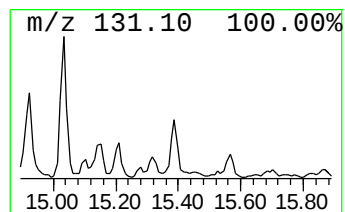
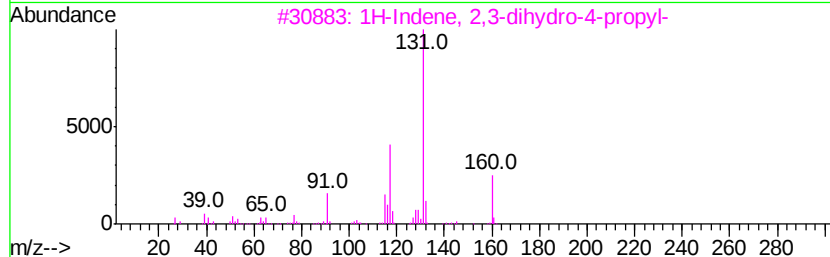
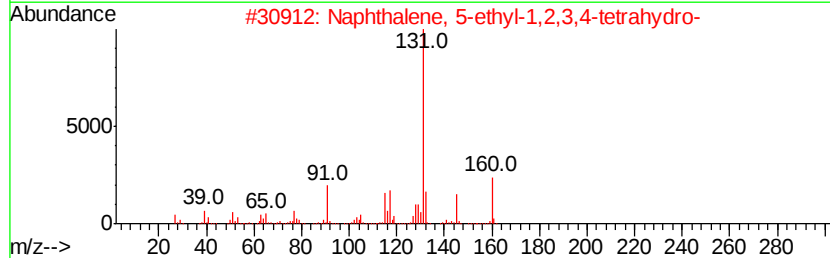
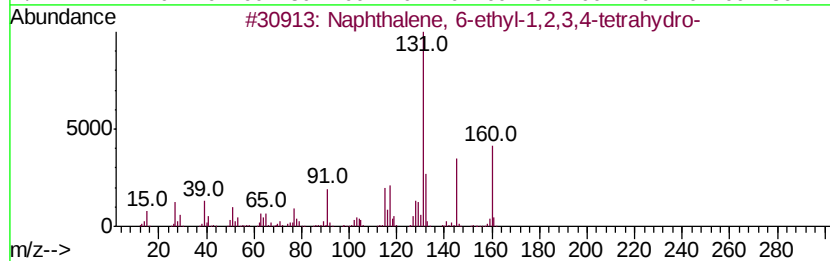
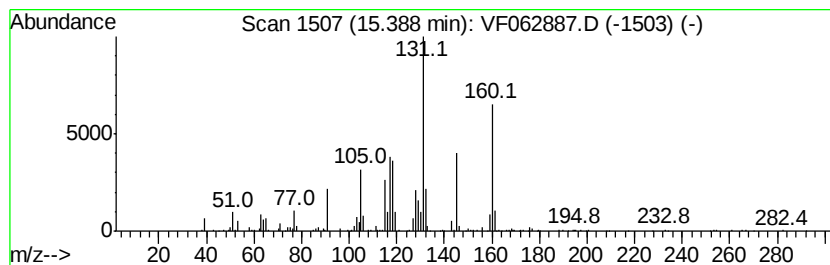
Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 6-ethyl-1,2,3,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.39	16.72 ug/l	657102	1,4-Dichlorobenzene-d4	12.62
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		Naphthalene, 5-ethyl-1,2,3,4-tet...	160 C12H16	042775-75-7 64
3		1H-Indene, 2,3-dihydro-4-propyl-	160 C12H16	092013-16-6 52
4		1H-Pyrrolo[2,3-b]pyridine, 2-n-p...	160 C10H12N2	1000212-02-1 49
5		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF060619\
Data File : VF062887.D
Acq On : 6 Jun 2019 17:23
Operator : FY/SY
Sample : K3160-01
Misc : 6.35g/5mL/MSVOA_F/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.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
o-Cymene	13.15	15.4	ug/l	605705	4	12.62	1965480	50.0
1H-Indene, 2,3-di...	13.76	28.4	ug/l	1114270	4	12.62	1965480	50.0
1H-Indene, 2,3-di...	13.90	38.2	ug/l	1502080	4	12.62	1965480	50.0
Naphthalene, 1,2,...	14.02	17.3	ug/l	678585	4	12.62	1965480	50.0
1H-Indene, 2,3-di...	14.20	21.1	ug/l	829437	4	12.62	1965480	50.0
unknown14.26	14.26	15.3	ug/l	599292	4	12.62	1965480	50.0
Naphthalene, 1,2,...	14.78	34.9	ug/l	1373650	4	12.62	1965480	50.0
Naphthalene, 1,2,...	15.15	26.0	ug/l	1022590	4	12.62	1965480	50.0
Naphthalene, 6-et...	15.39	16.7	ug/l	657102	4	12.62	1965480	50.0