

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF050719\
 Data File : VF062357.D
 Acq On : 7 May 2019 15:05
 Operator : FY/SY
 Sample : K2635-01
 Misc : 5.59g/5mL/MSVOA F/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 AU-05-050219-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\82F050619S.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	1.111	56	59	61	rBV2	10922	23068	1.80%	0.144%
2	1.387	81	87	88	rBV5	8938	24728	1.93%	0.155%
3	2.333	180	183	188	rBV7	7598	26744	2.09%	0.168%
4	3.161	265	267	271	rVB5	12199	16412	1.28%	0.103%
5	4.305	376	383	390	rBV	167074	509211	39.82%	3.189%
6	5.044	449	458	470	rBV2	379529	1278939	100.00%	8.010%
7	5.527	506	507	509	rBV	21859	16648	1.30%	0.104%
8	5.773	526	532	545	rBV	396817	1074584	84.02%	6.730%
9	7.705	722	728	739	rBV	357211	850063	66.47%	5.324%
10	8.405	794	799	803	rVB4	10800	29931	2.34%	0.187%
11	8.740	830	833	839	rVB7	5704	15577	1.22%	0.098%
12	9.618	918	922	928	rVB4	10987	26714	2.09%	0.167%
13	9.904	945	951	960	rBV	450234	994174	77.73%	6.227%
14	10.032	960	964	971	rVB5	6658	19750	1.54%	0.124%
15	10.258	981	987	992	rVB5	7984	20854	1.63%	0.131%
16	10.771	1034	1039	1040	rBV4	7802	15650	1.22%	0.098%
17	10.810	1040	1043	1046	rVV2	28436	53966	4.22%	0.338%
18	10.879	1046	1050	1055	rVV6	7074	20714	1.62%	0.130%
19	11.146	1072	1077	1080	rBV7	3992	13871	1.08%	0.087%
20	11.205	1080	1083	1087	rBV5	6064	13273	1.04%	0.083%
21	11.441	1105	1107	1108	rBV2	10155	14677	1.15%	0.092%
22	11.491	1108	1112	1118	rVB	427737	680059	53.17%	4.259%
23	11.629	1123	1126	1128	rVV	18416	32119	2.51%	0.201%
24	11.668	1128	1130	1134	rVV2	11649	21241	1.66%	0.133%
25	11.776	1138	1141	1146	rVV2	67881	138538	10.83%	0.868%
26	11.885	1150	1152	1155	rVB	41918	64596	5.05%	0.405%
27	12.082	1167	1172	1175	rVB	60375	95486	7.47%	0.598%
28	12.200	1178	1184	1185	rBV4	6226	14692	1.15%	0.092%
29	12.259	1187	1190	1195	rVB	42737	66196	5.18%	0.415%
30	12.358	1195	1200	1205	rBV3	9521	24926	1.95%	0.156%
31	12.486	1205	1213	1214	rBV2	28192	56520	4.42%	0.354%
32	12.516	1214	1216	1222	rVB2	25523	39668	3.10%	0.248%
33	12.604	1222	1225	1229	rBV	716056	1063774	83.18%	6.663%
34	12.664	1229	1231	1235	rVB	156935	200692	15.69%	1.257%

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35	12.772	1239	1242	1245	rBV2	76729	141402	11.06%	0.886%
36	12.821	1245	1247	1249	rVV	57318	81513	6.37%	0.511%
37	12.880	1249	1253	1257	rVB2	98410	200636	15.69%	1.257%
38	12.979	1260	1263	1268	rBV5	30942	64045	5.01%	0.401%
39	13.048	1268	1270	1274	rVB	62666	95416	7.46%	0.598%
40	13.137	1274	1279	1282	rBV2	59557	131417	10.28%	0.823%
41	13.196	1282	1285	1287	rBV	134969	218153	17.06%	1.366%
42	13.275	1291	1293	1297	rVB3	66514	138235	10.81%	0.866%
43	13.423	1306	1308	1310	rBV2	21232	26086	2.04%	0.163%
44	13.462	1310	1312	1317	rVB	54349	80887	6.32%	0.507%
45	13.541	1317	1320	1322	rBV	74192	111641	8.73%	0.699%
46	13.580	1322	1324	1326	rVV	95624	136299	10.66%	0.854%
47	13.620	1326	1328	1330	rVV	47037	53533	4.19%	0.335%
48	13.649	1330	1331	1333	rVV	25984	31248	2.44%	0.196%
49	13.679	1333	1334	1337	rVB	41059	33558	2.62%	0.210%
50	13.748	1337	1341	1344	rBV	167190	261533	20.45%	1.638%
51	13.797	1344	1346	1348	rVB	49338	62766	4.91%	0.393%
52	13.846	1348	1351	1353	rBV	67131	91873	7.18%	0.575%
53	13.886	1353	1355	1359	rBV2	334148	597233	46.70%	3.741%
54	13.955	1359	1362	1364	rVB2	88444	144371	11.29%	0.904%
55	14.004	1364	1367	1371	rVB	246882	318506	24.90%	1.995%
56	14.083	1372	1375	1376	rBV	46274	64908	5.08%	0.407%
57	14.113	1376	1378	1381	rVB	93074	120160	9.40%	0.753%
58	14.182	1381	1385	1388	rBV4	161650	368910	28.85%	2.311%
59	14.251	1388	1392	1398	rVB2	194930	399402	31.23%	2.502%
60	14.329	1398	1400	1402	rBV3	14753	16364	1.28%	0.102%
61	14.398	1402	1407	1408	rBV4	29244	73751	5.77%	0.462%
62	14.438	1408	1411	1414	rBV3	160694	278464	21.77%	1.744%
63	14.497	1414	1417	1419	rBV2	134196	187760	14.68%	1.176%
64	14.586	1424	1426	1428	rBV	37420	60677	4.74%	0.380%
65	14.625	1428	1430	1433	rVB2	129297	174704	13.66%	1.094%
66	14.724	1438	1440	1441	rVV	39082	62797	4.91%	0.393%
67	14.763	1441	1444	1448	rVV	403440	634570	49.62%	3.975%
68	14.852	1451	1453	1454	rVV	54219	77616	6.07%	0.486%
69	14.911	1454	1459	1463	rVV3	134764	386449	30.22%	2.420%
70	14.960	1463	1464	1467	rVV2	40500	51537	4.03%	0.323%
71	15.019	1467	1470	1473	rVV	313805	446672	34.93%	2.798%

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72	15.088	1473	1477	1478	rVV2	30797	64762	5.06%	0.406%
73	15.138	1478	1482	1484	rVV	277646	459755	35.95%	2.880%
74	15.167	1484	1485	1493	rVV4	80138	219011	17.12%	1.372%
75	15.266	1493	1495	1496	rVV2	30583	43562	3.41%	0.273%
76	15.305	1496	1499	1501	rVV2	54365	107106	8.37%	0.671%
77	15.374	1501	1506	1509	rVV2	135133	328702	25.70%	2.059%
78	15.433	1509	1512	1518	rVV	47217	104654	8.18%	0.655%
79	15.552	1518	1524	1528	rVB3	28431	79800	6.24%	0.500%
80	15.670	1532	1536	1540	rVV	143929	287743	22.50%	1.802%
81	15.739	1540	1543	1547	rVV5	22569	51357	4.02%	0.322%
82	15.847	1550	1554	1558	rVV6	30557	68809	5.38%	0.431%
83	15.966	1563	1566	1568	rVV3	26927	50125	3.92%	0.314%
84	16.005	1568	1570	1579	rVB3	34320	117355	9.18%	0.735%

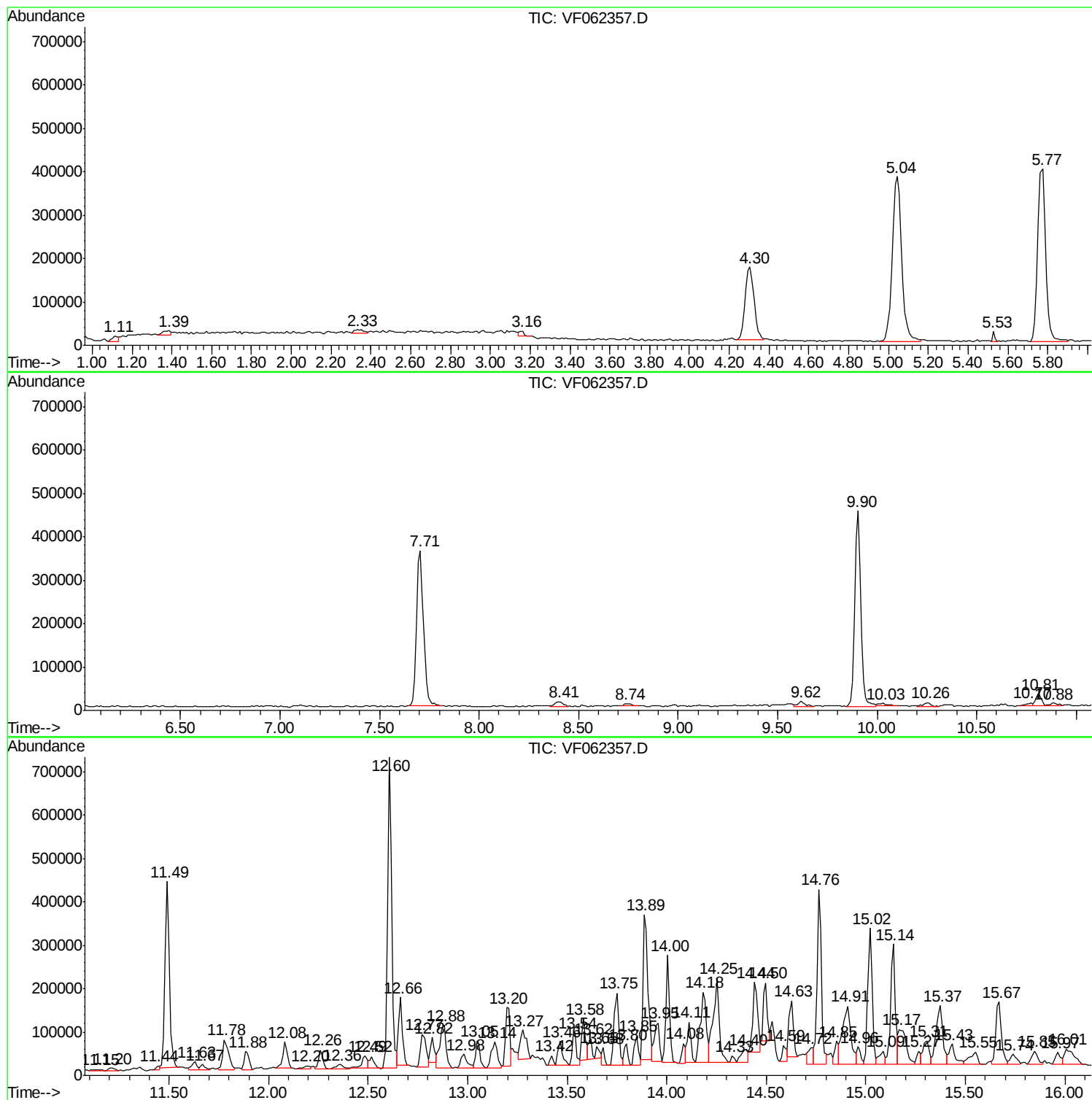
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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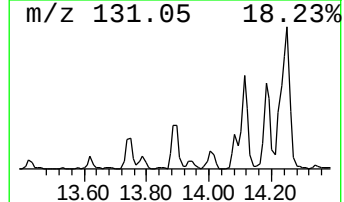
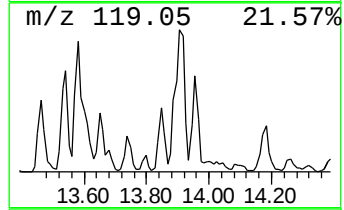
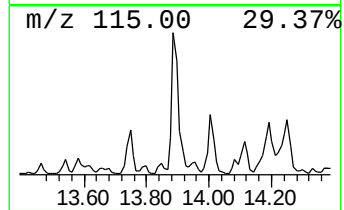
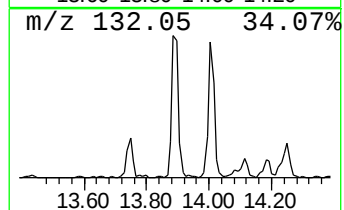
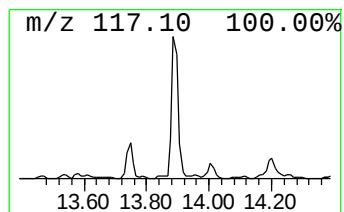
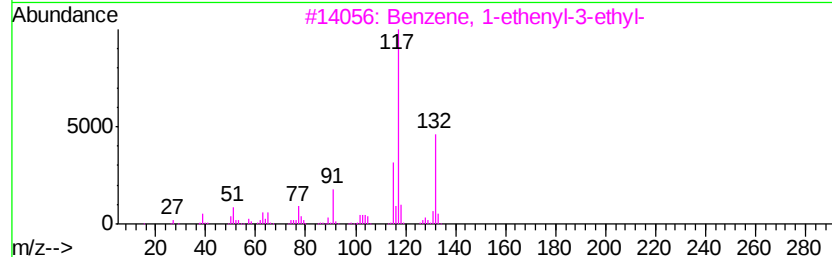
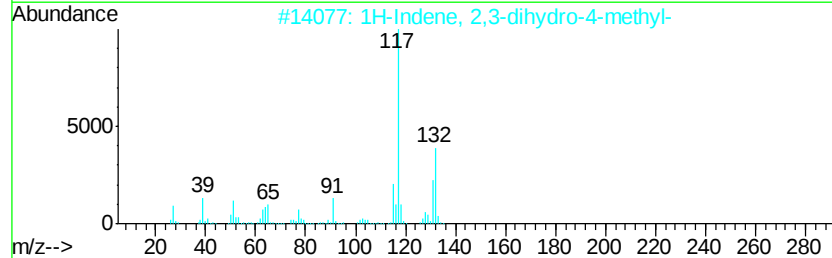
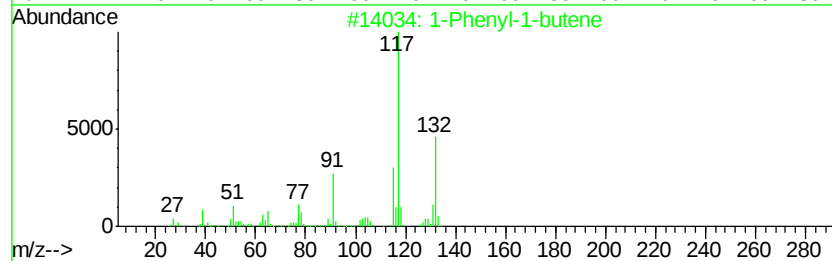
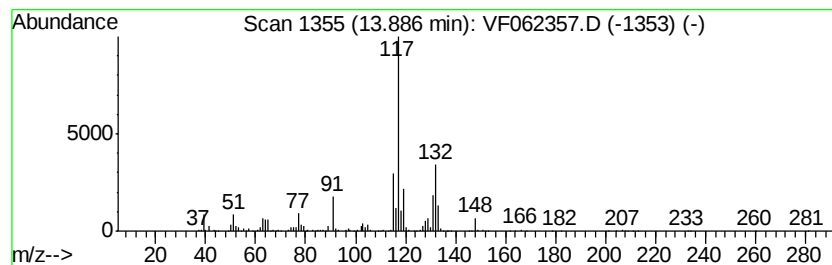
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F050619S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	28.07 ug/l	597233	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



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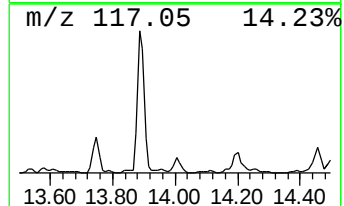
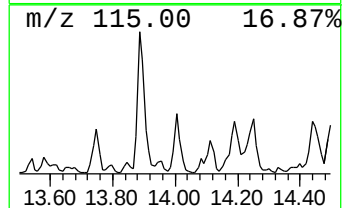
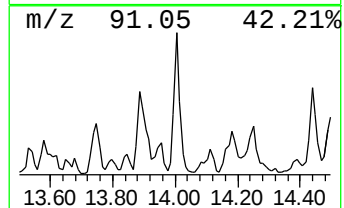
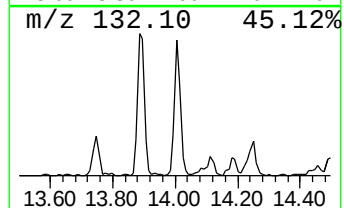
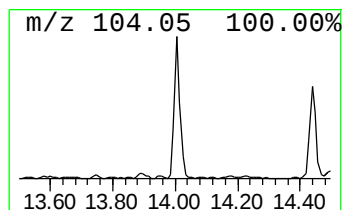
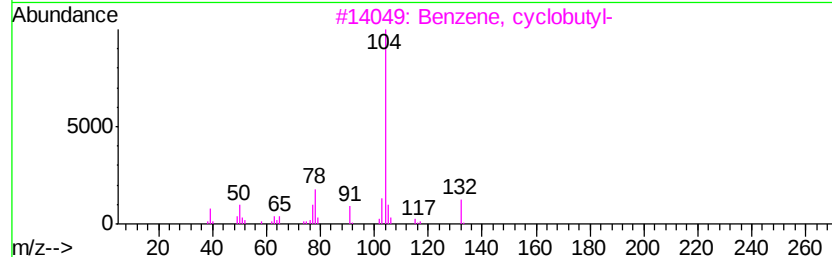
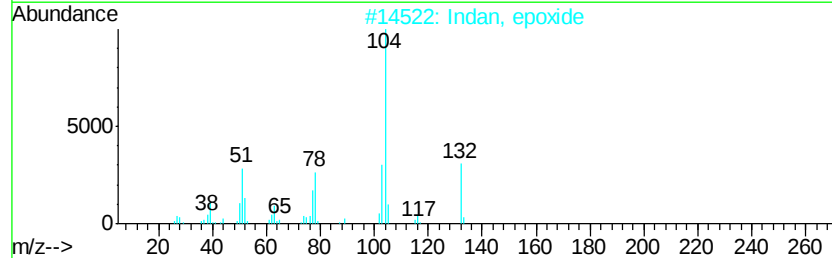
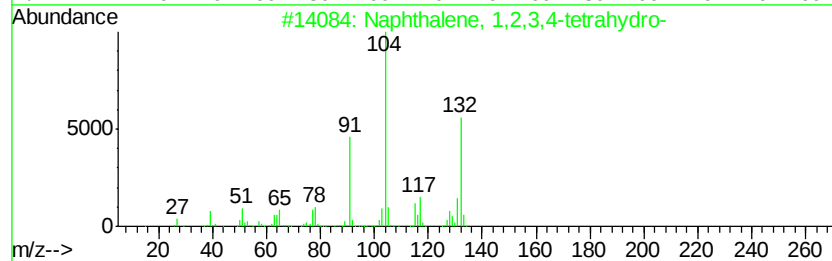
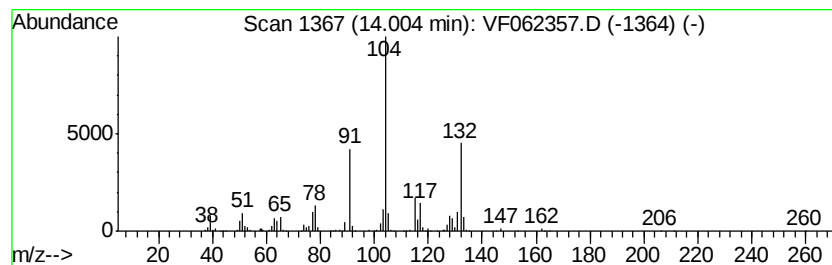
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	14.97 ug/l	318506	1,4-Dichlorobenzene-d4	12.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 96
2		Indan, epoxide	132 C9H8O	000768-22-9 62
3		Benzene, cyclobutyl-	132 C10H12	004392-30-7 59
4		Benzene, 3-cyclohexen-1-yl-	158 C12H14	004994-16-5 47
5		1H-Inden-1-one, 2,3-dihydro-	132 C9H8O	000083-33-0 43



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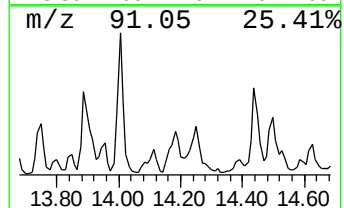
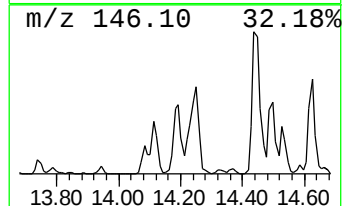
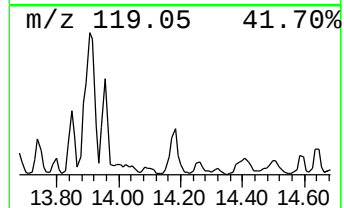
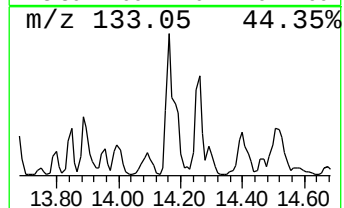
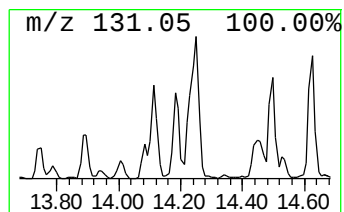
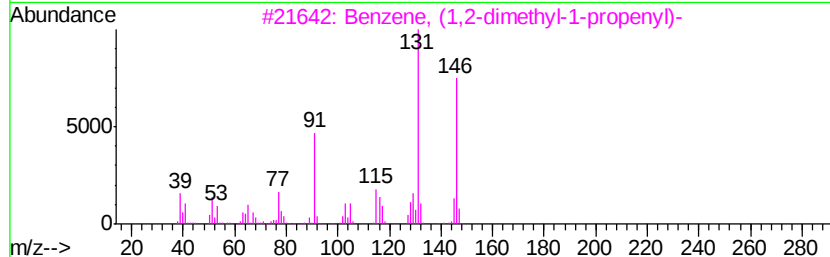
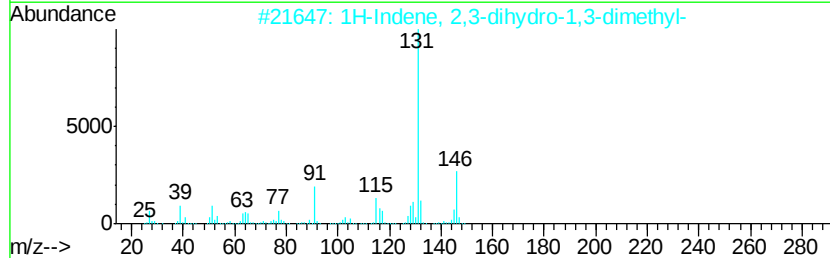
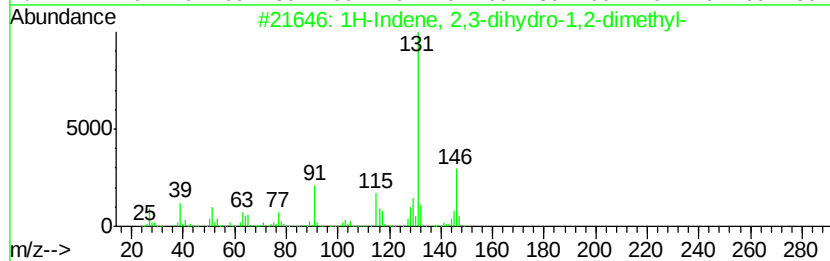
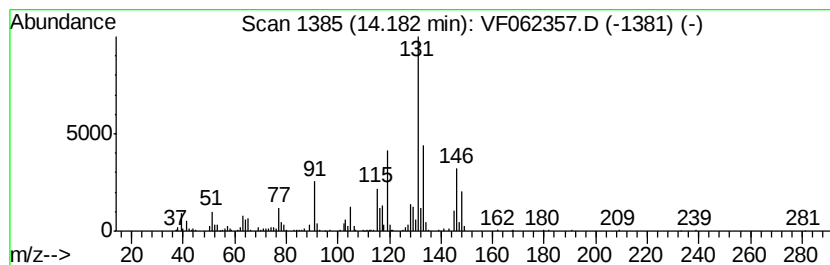
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	17.34 ug/l	368910	1,4-Dichlorobenzene-d4	12.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	92
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	91
3	Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3	89
4	Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2	78
5	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	78



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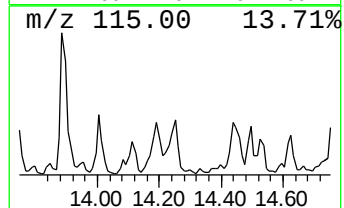
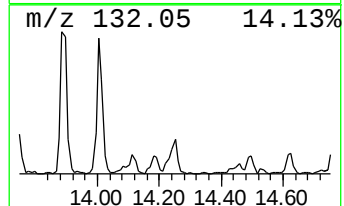
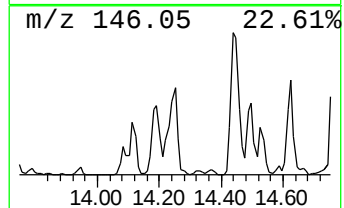
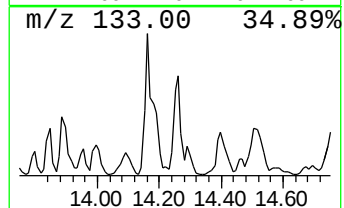
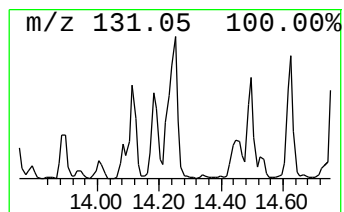
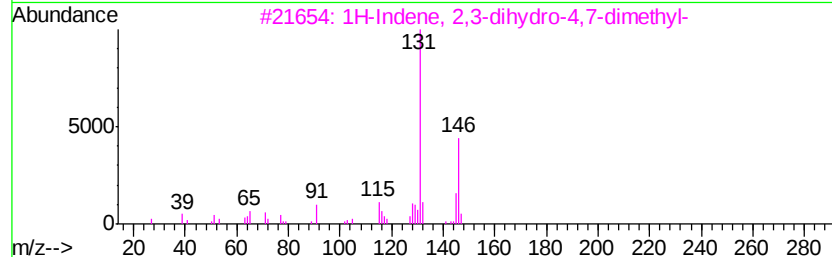
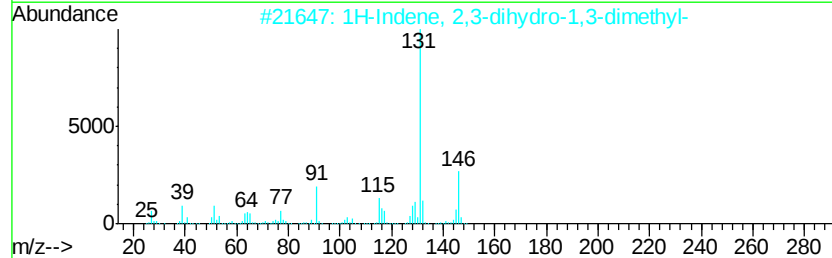
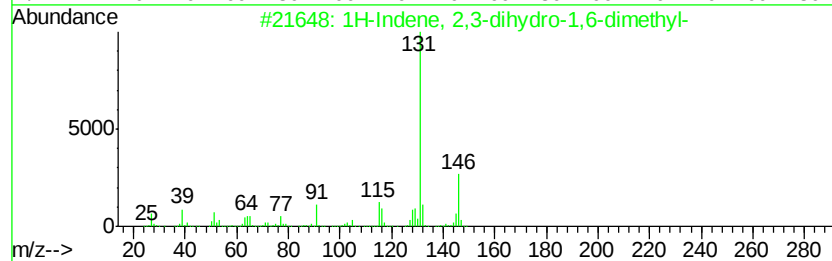
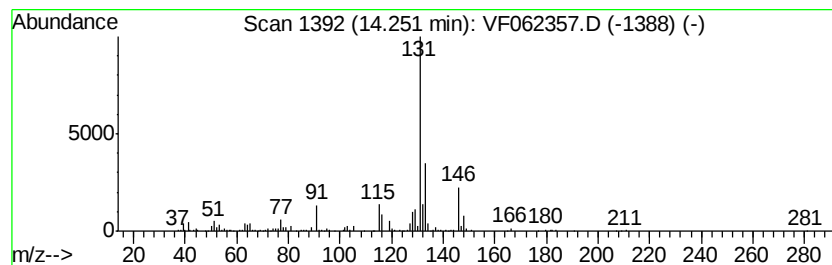
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	18.77 ug/l	399402	1,4-Dichlorobenzene-d4	12.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14		017059-48-2	81
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14		004175-53-5	81
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	81
4	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14		004912-92-9	81
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14		020836-11-7	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA F/SOIL
ALS Vial : 11 Sample Multiplier: 1

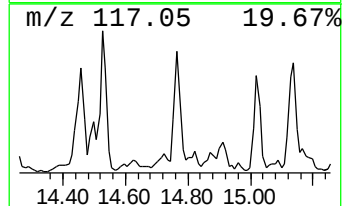
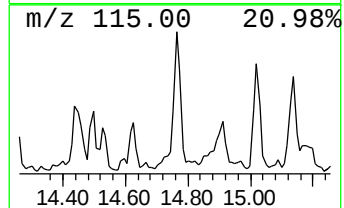
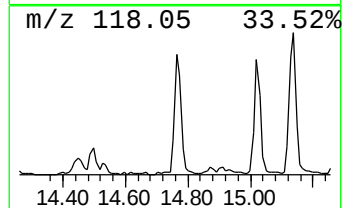
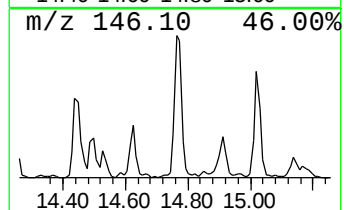
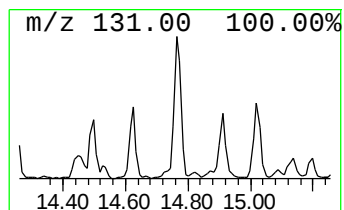
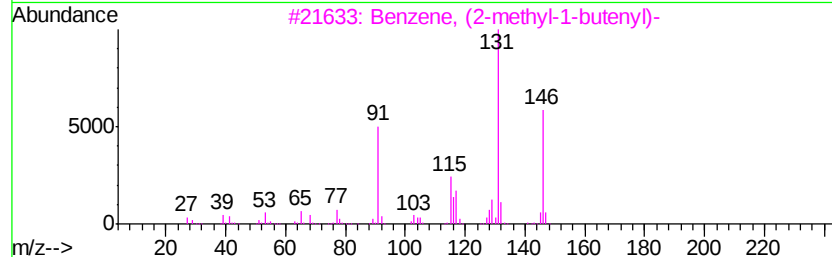
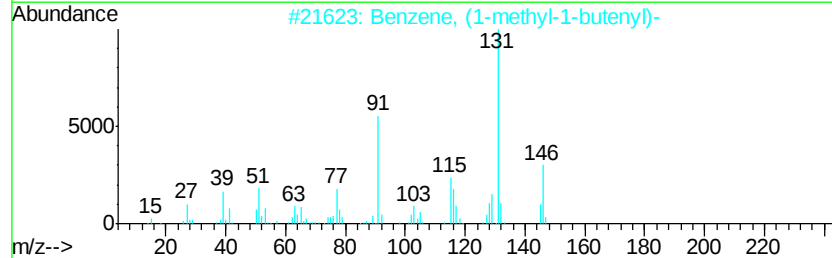
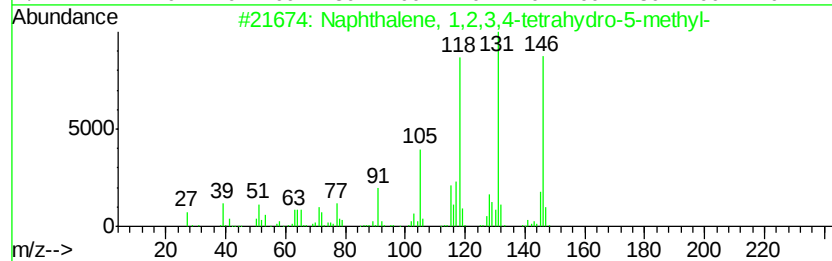
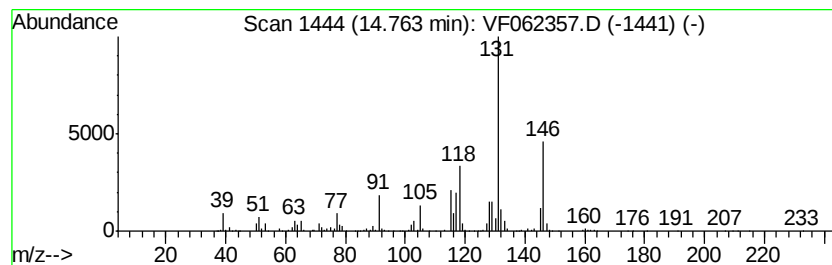
Instrument :
MSVOA_F
ClientSampleId :
AU-05-050219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F050619S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	29.83 ug/l	634570	1,4-Dichlorobenzene-d4	12.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 91
2		Benzene, (1-methyl-1-butenvl)-	146 C11H14	053172-84-2 90
3		Benzene, (2-methyl-1-butenvl)-	146 C11H14	056253-64-6 87
4		1,3-Cyclopentadiene, 5-(trans-2-...	146 C11H14	079209-36-2 87
5		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
AU-05-050219-A

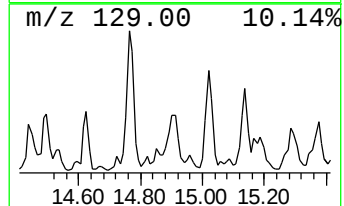
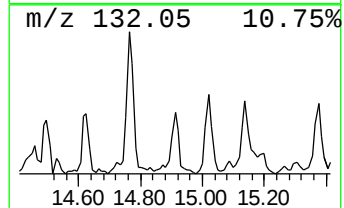
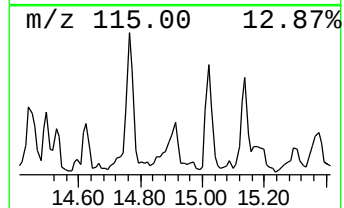
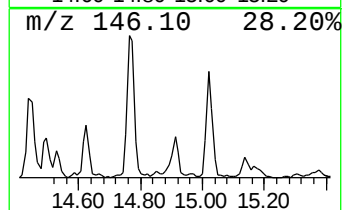
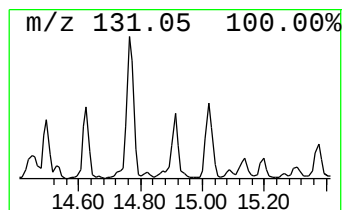
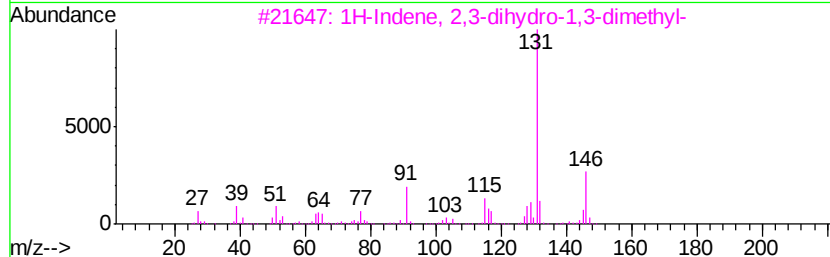
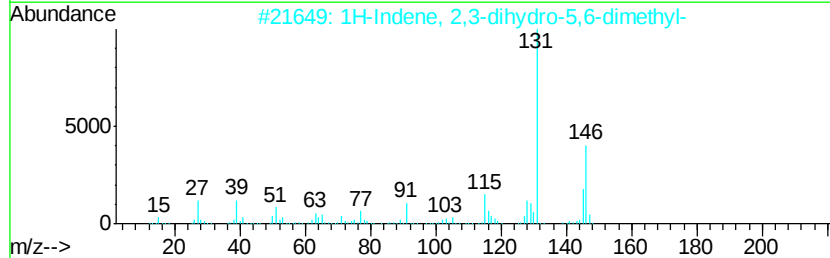
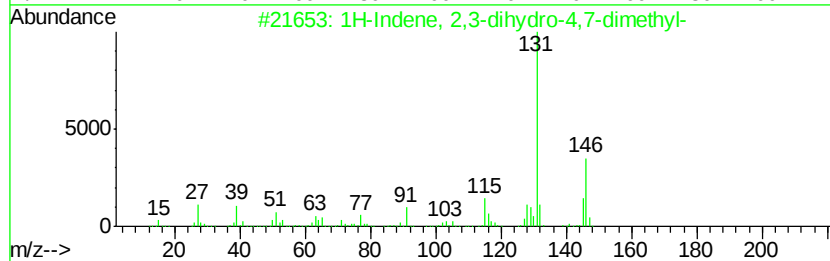
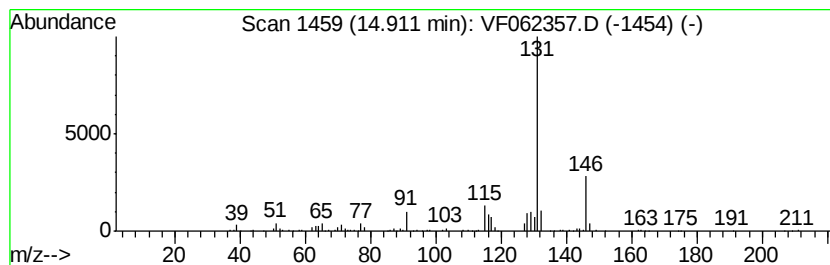
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	18.16 ug/l	386449	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA F/SOIL
ALS Vial : 11 Sample Multiplier: 1

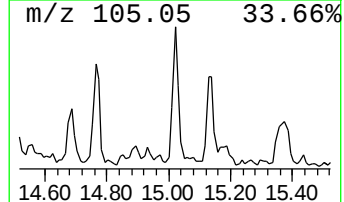
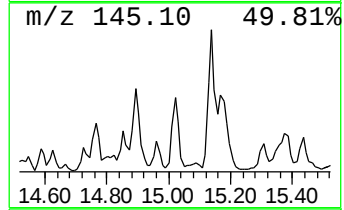
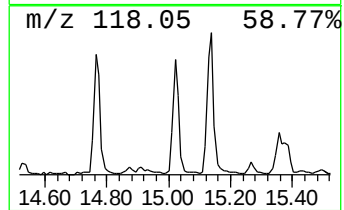
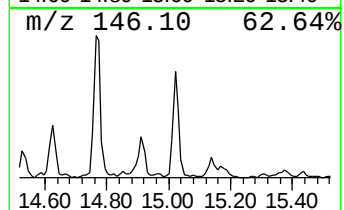
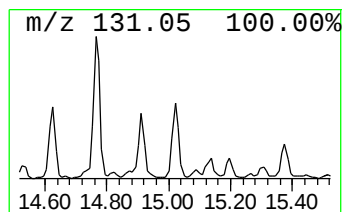
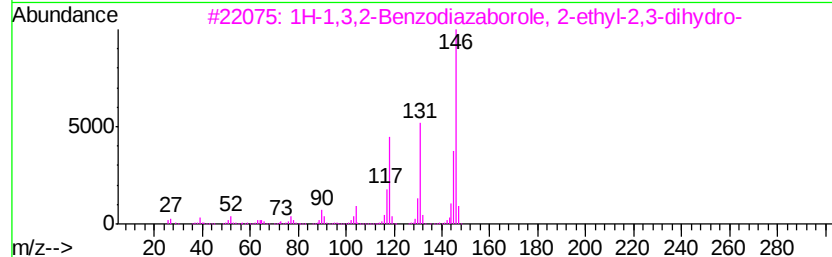
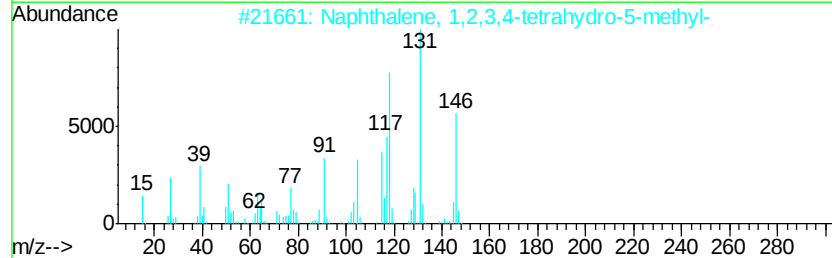
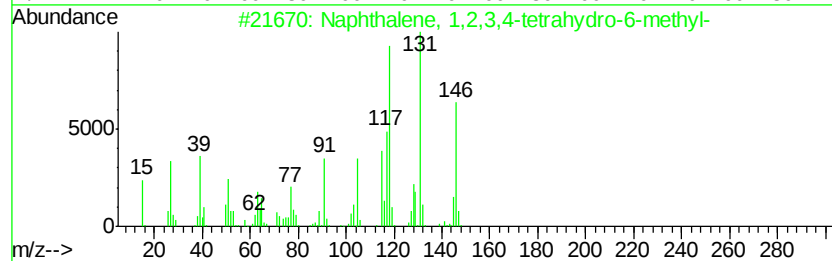
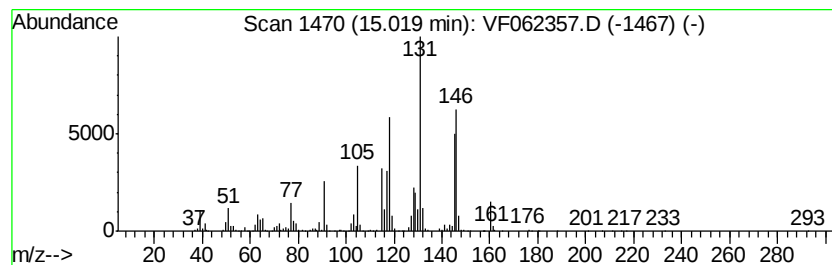
Instrument :
MSVOA_F
ClientSampled :
AU-05-050219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F050619S.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.
15.02	20.99 ug/l	446672	1,4-Dichlorobenzene-d4	12.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 76
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 70
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 60
5		1H-Benzimidazole, 5,6-dimethyl-	146 C9H10N2	000582-60-5 60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
AU-05-050219-A

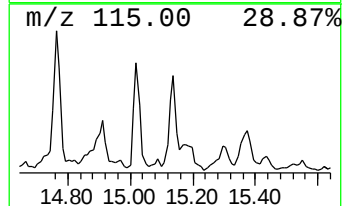
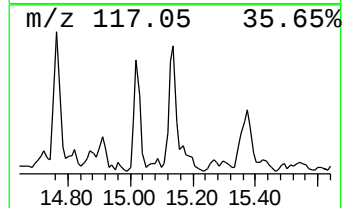
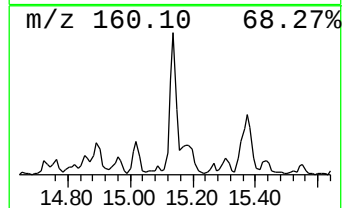
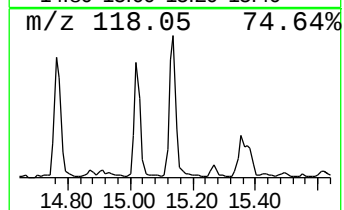
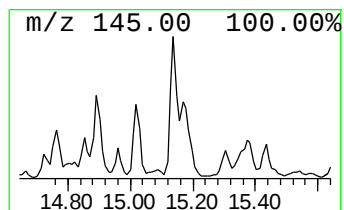
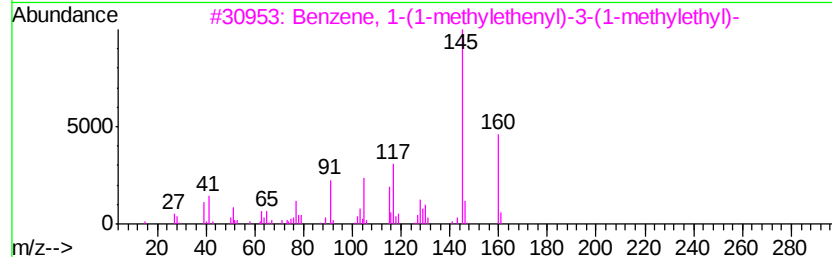
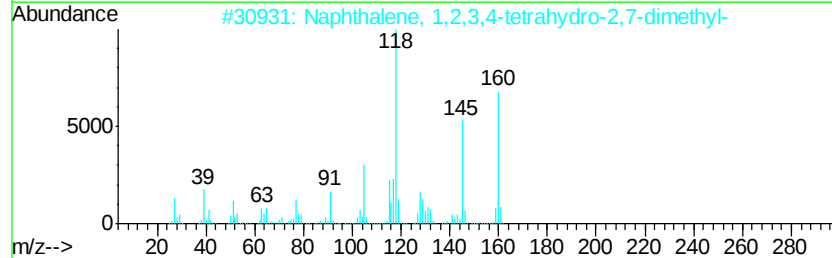
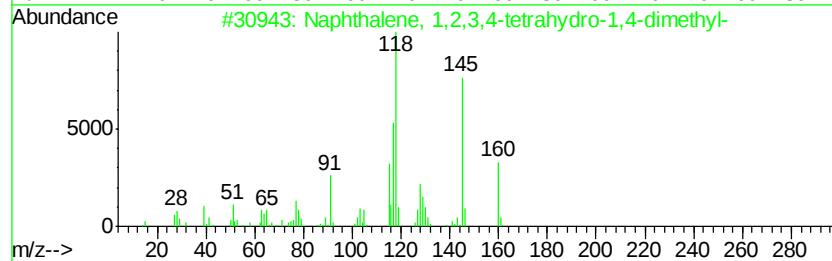
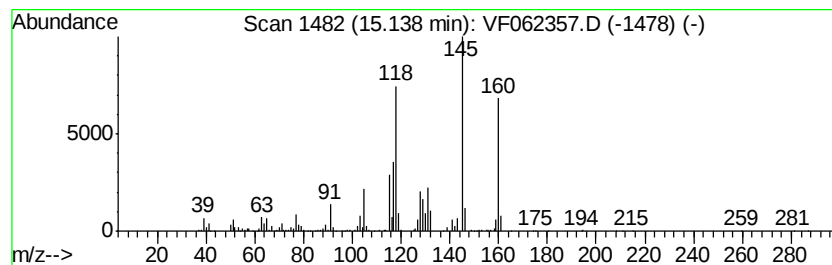
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F050619S.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	21.61 ug/l	459755	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	70
3		Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	60
4		1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	60
5		Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
AU-05-050219-A

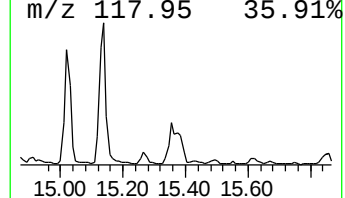
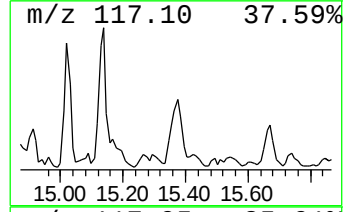
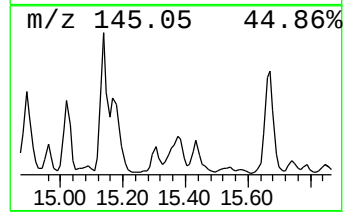
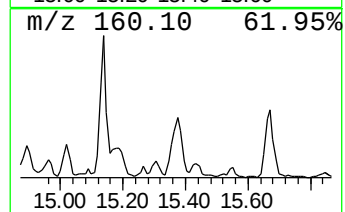
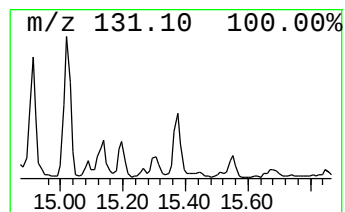
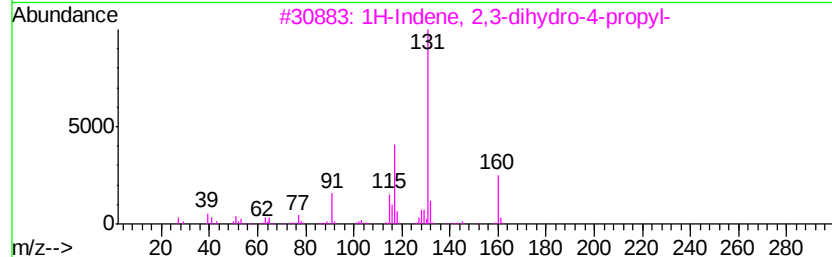
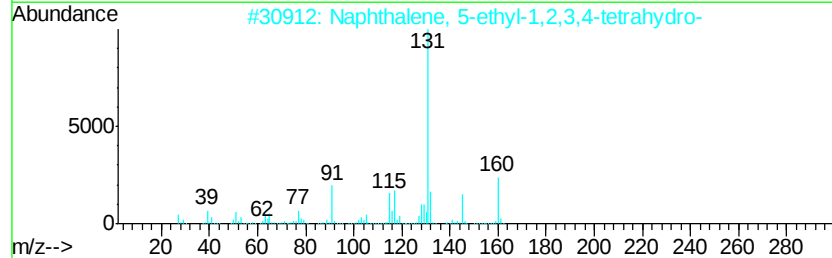
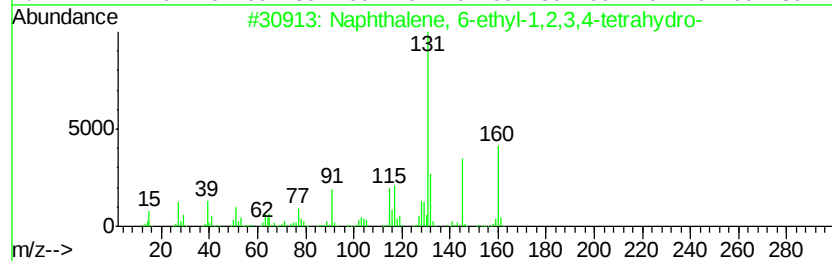
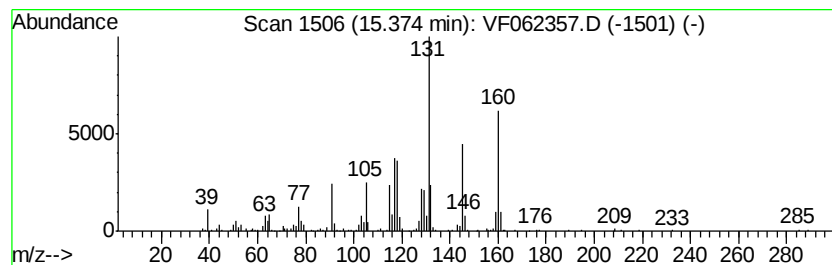
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 6-ethyl-1,2,3,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	15.45 ug/l	328702	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	87
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	49
4		1H-Pyrrolo[2,3-b]pyridine, 2-n-p...	160	C10H12N2	1000212-02-1	46
5		Hex-3-ene-1,5-diyne, 3,4-diisopr...	160	C12H16	1000211-22-8	41



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
AU-05-050219-A

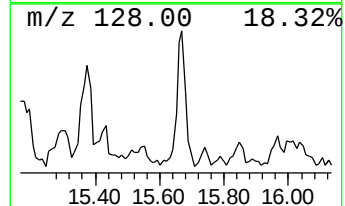
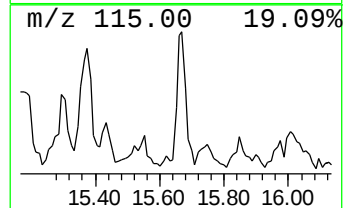
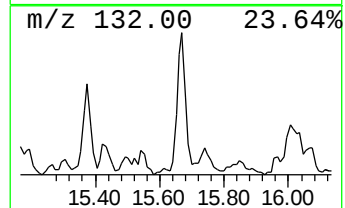
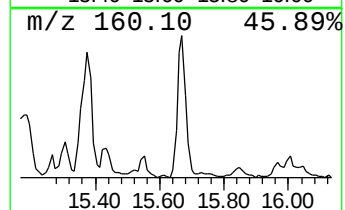
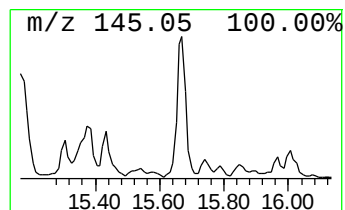
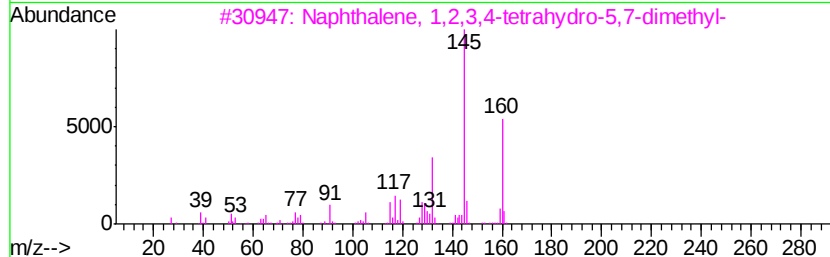
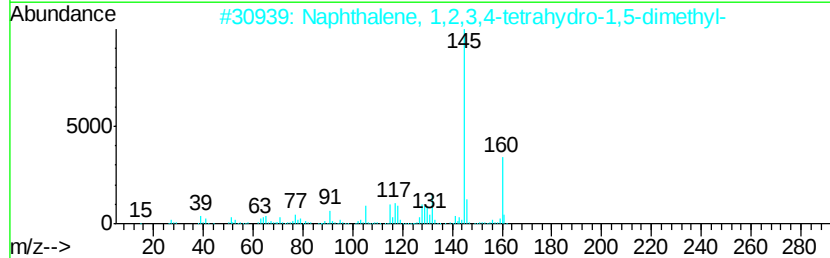
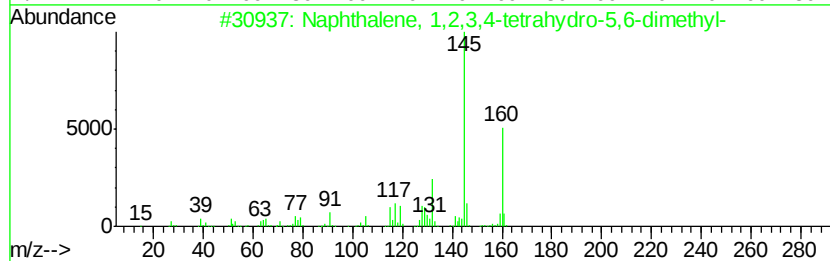
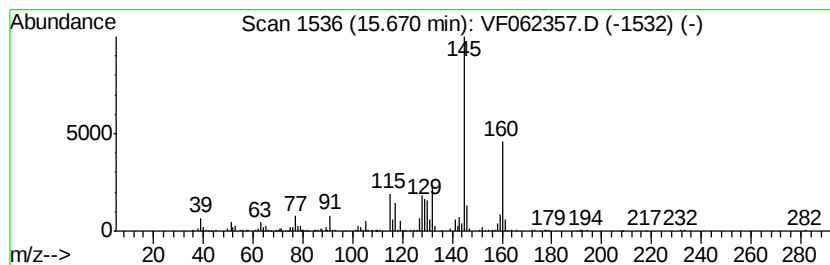
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F050619S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	13.52 ug/l	287743	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	020027-77-4	91
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	91
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	90
4		Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	90
5		1H-Inden-1-one, 2,3-dihydro-3,3-...	160	C11H12O	026465-81-6	89



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF050719\
Data File : VF062357.D
Acq On : 7 May 2019 15:05
Operator : FY/SY
Sample : K2635-01
Misc : 5.59g/5mL/MSVOA_F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
AU-05-050219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82F050619S.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
1-Phenyl-1-butene	13.89	28.1	ug/l	597233	4	12.60	1063770	50.0
Naphthalene, 1,2,...	14.00	15.0	ug/l	318506	4	12.60	1063770	50.0
1H-Indene, 2,3-di...	14.18	17.3	ug/l	368910	4	12.60	1063770	50.0
1H-Indene, 2,3-di...	14.25	18.8	ug/l	399402	4	12.60	1063770	50.0
Naphthalene, 1,2,...	14.76	29.8	ug/l	634570	4	12.60	1063770	50.0
1H-Indene, 2,3-di...	14.91	18.2	ug/l	386449	4	12.60	1063770	50.0
Naphthalene, 1,2,...	15.02	21.0	ug/l	446672	4	12.60	1063770	50.0
Naphthalene, 1,2,...	15.14	21.6	ug/l	459755	4	12.60	1063770	50.0
Naphthalene, 6-et...	15.37	15.4	ug/l	328702	4	12.60	1063770	50.0
Naphthalene, 1,2,...	15.67	13.5	ug/l	287743	4	12.60	1063770	50.0