

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF111918\
 Data File : VF060788.D
 Acq On : 19 Nov 2018 20:58
 Operator : VA/AP
 Sample : J5950-06
 Misc : 5.00µ/5mL/MSVOA-F/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 SSI-20-TW-1

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\82F111318S.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	1.091	40	47	54	rBV3	15580	47940	3.29%	0.295%
2	2.186	155	158	160	rBV4	7180	14957	1.03%	0.092%
3	2.235	160	163	170	rVB3	10349	27482	1.89%	0.169%
4	3.893	322	331	334	rBV5	3889	14895	1.02%	0.092%
5	4.021	337	344	348	rVV4	9360	33793	2.32%	0.208%
6	4.120	348	354	368	rVB	110167	362438	24.90%	2.232%
7	4.642	402	407	415	rVB2	11701	36911	2.54%	0.227%
8	4.869	422	430	444	rBV2	212322	777039	53.38%	4.785%
9	5.461	484	490	499	rBV4	7575	24688	1.70%	0.152%
10	5.609	499	505	516	rVV	246733	675878	46.43%	4.162%
11	6.280	565	573	578	rBV4	8751	28534	1.96%	0.176%
12	6.457	585	591	595	rBV5	10482	32018	2.20%	0.197%
13	6.832	628	629	631	rBV	29829	19149	1.32%	0.118%
14	7.572	698	704	717	rBV	304771	744098	51.12%	4.582%
15	9.644	908	914	919	rBV6	6431	19269	1.32%	0.119%
16	9.782	922	928	936	rBV	336648	781826	53.71%	4.814%
17	9.910	936	941	946	rVV	352197	727707	49.99%	4.481%
18	10.147	959	965	970	rVB	21984	44650	3.07%	0.275%
19	10.245	970	975	984	rBV5	13908	40110	2.76%	0.247%
20	10.531	1000	1004	1010	rVB4	17084	37199	2.56%	0.229%
21	10.679	1014	1019	1026	rVV6	19991	62171	4.27%	0.383%
22	11.064	1055	1058	1062	rBV3	10216	22701	1.56%	0.140%
23	11.133	1062	1065	1070	rVV	79194	137526	9.45%	0.847%
24	11.281	1076	1080	1083	rVB5	11151	23607	1.62%	0.145%
25	11.419	1090	1094	1100	rBV	322938	539267	37.05%	3.321%
26	11.607	1105	1113	1119	rBV	222936	425681	29.24%	2.621%
27	11.735	1122	1126	1133	rVB3	21843	53014	3.64%	0.326%
28	11.962	1144	1149	1151	rBV4	6514	16664	1.14%	0.103%
29	12.031	1151	1156	1159	rVB	23313	43493	2.99%	0.268%
30	12.139	1164	1167	1172	rVV	55742	90627	6.23%	0.558%
31	12.208	1172	1174	1179	rVB	31667	48412	3.33%	0.298%
32	12.307	1180	1184	1190	rBV	120614	204107	14.02%	1.257%
33	12.553	1206	1209	1213	rVV	496755	829037	56.96%	5.105%
34	12.613	1213	1215	1220	rVV2	71841	116475	8.00%	0.717%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	12.731	1220	1227	1232	rVV2	618307	1237173	84.99%	7.618%
36	12.810	1232	1235	1237	rVV	80116	131459	9.03%	0.809%
37	12.849	1237	1239	1242	rVV	79566	110973	7.62%	0.683%
38	12.938	1245	1248	1252	rVV2	79024	125141	8.60%	0.771%
39	13.007	1252	1255	1260	rVB2	22621	51838	3.56%	0.319%
40	13.086	1260	1263	1265	rBV2	13051	25129	1.73%	0.155%
41	13.165	1267	1271	1276	rVV	842592	1455586	100.00%	8.963%
42	13.234	1276	1278	1282	rVV	330378	505859	34.75%	3.115%
43	13.293	1282	1284	1291	rVV2	108127	193637	13.30%	1.192%
44	13.392	1291	1294	1296	rVB	45381	63148	4.34%	0.389%
45	13.500	1302	1305	1309	rVB	582384	775765	53.30%	4.777%
46	13.570	1310	1312	1317	rVB2	64671	106879	7.34%	0.658%
47	13.648	1317	1320	1323	rBV	82878	106817	7.34%	0.658%
48	13.717	1323	1327	1330	rBV	322075	472835	32.48%	2.911%
49	13.816	1334	1337	1339	rBV	64884	86115	5.92%	0.530%
50	13.856	1339	1341	1343	rBV	538410	1000016	68.70%	6.158%
51	13.925	1347	1348	1351	rVB	73237	71719	4.93%	0.442%
52	13.974	1351	1353	1356	rVB	167872	234922	16.14%	1.447%
53	14.053	1359	1361	1362	rBV	43095	55325	3.80%	0.341%
54	14.092	1362	1365	1367	rVB	121049	185033	12.71%	1.139%
55	14.161	1367	1372	1374	rBV3	99971	263900	18.13%	1.625%
56	14.221	1374	1378	1384	rVB2	166490	337169	23.16%	2.076%
57	14.309	1384	1387	1389	rVB2	11738	14723	1.01%	0.091%
58	14.369	1389	1393	1395	rBV2	26048	38265	2.63%	0.236%
59	14.467	1395	1403	1406	rBV	286587	566102	38.89%	3.486%
60	14.507	1406	1407	1410	rVV	53315	53718	3.69%	0.331%
61	14.605	1414	1417	1419	rVB	87440	124099	8.53%	0.764%
62	14.743	1424	1431	1435	rBV	129433	210590	14.47%	1.297%
63	14.862	1438	1443	1444	rBV	34214	60562	4.16%	0.373%
64	14.891	1444	1446	1449	rVV	51469	73006	5.02%	0.450%
65	14.951	1449	1452	1455	rVB3	9956	17204	1.18%	0.106%
66	15.000	1455	1457	1461	rVB	40962	58175	4.00%	0.358%
67	15.118	1466	1469	1471	rBV3	22450	34058	2.34%	0.210%
68	15.168	1471	1474	1481	rVB7	13038	34313	2.36%	0.211%
69	15.286	1481	1486	1490	rBV	65451	112967	7.76%	0.696%
70	15.414	1496	1499	1503	rVB	67354	101639	6.98%	0.626%
71	15.651	1519	1523	1528	rVB4	11075	25886	1.78%	0.159%

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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

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

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72	16.006	1557	1559	1566	rVB6	5324	15171	1.04%	0.093%
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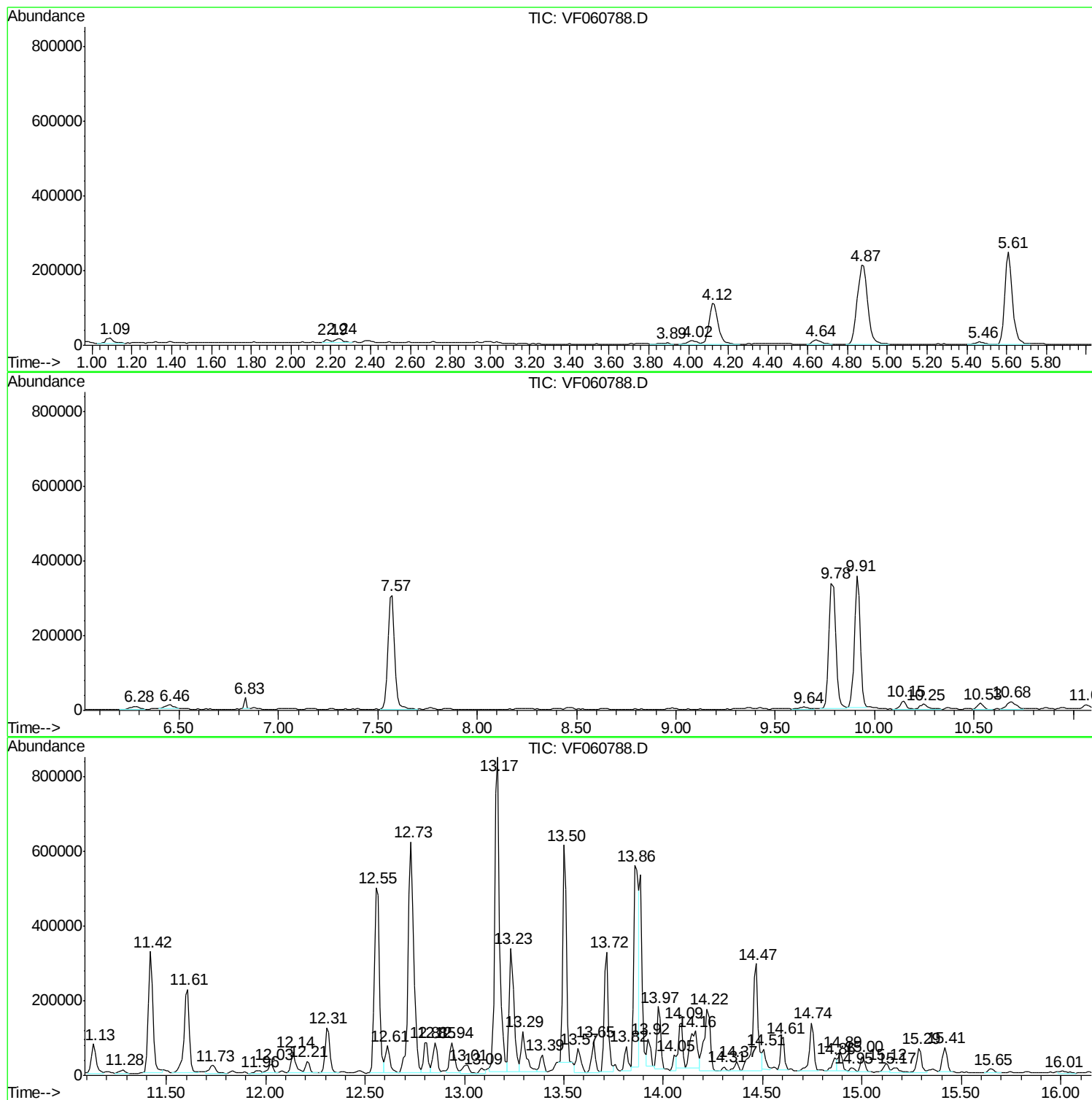
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Data Path : Z:\V0ASRV\HPCHEM1\MSV0A F\DATA\VF111918\
Data File : VF060788.D
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSV0A_F
ClientSampleId :
SSI-20-TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-20-TW-1

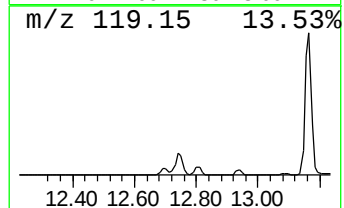
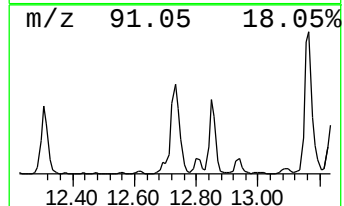
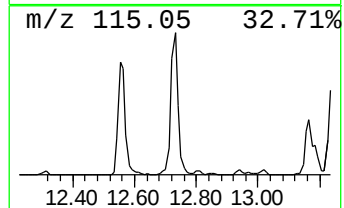
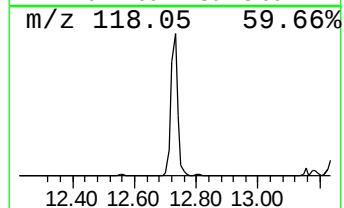
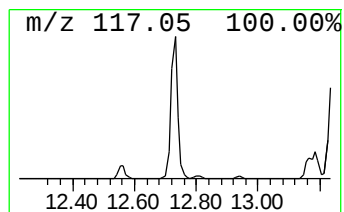
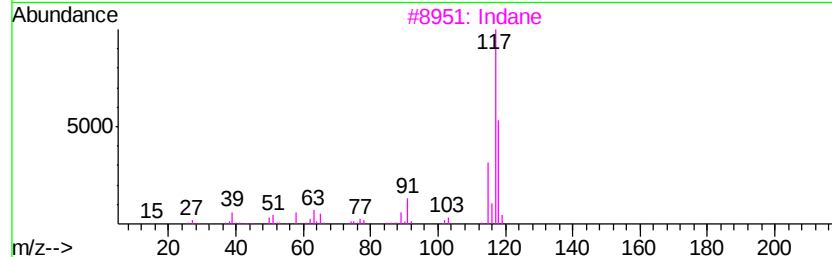
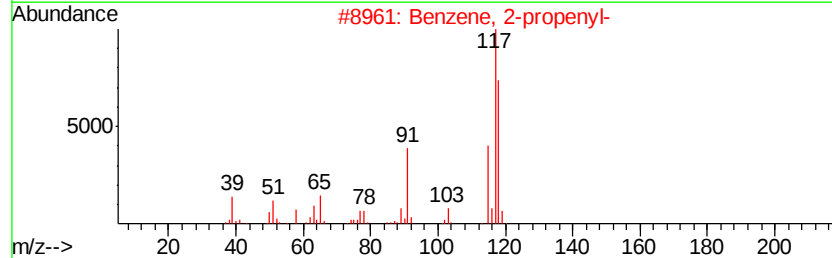
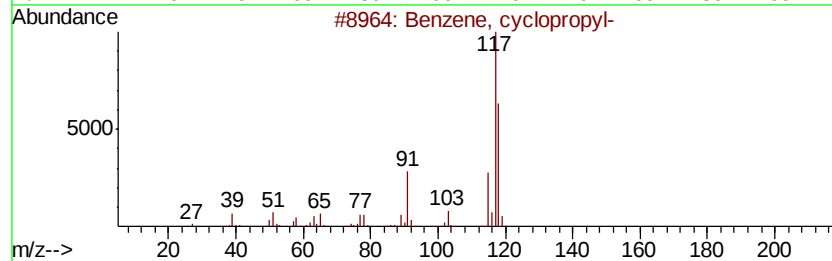
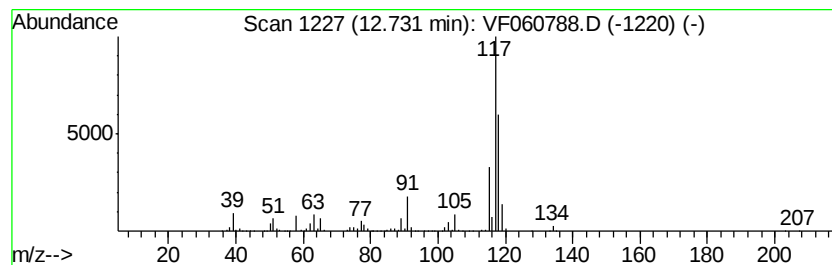
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, cyclopropyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	74.62 ug/l	1237170	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, cvclopropyl-	118	C9H10	000873-49-4	81
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3		Indane	118	C9H10	000496-11-7	81
4		Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	78
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	72



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Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampled :
SSI-20-TW-1

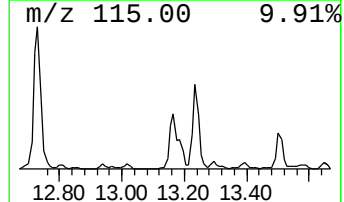
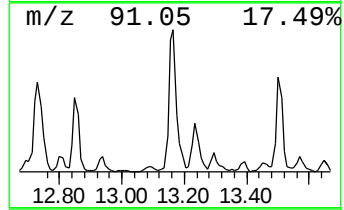
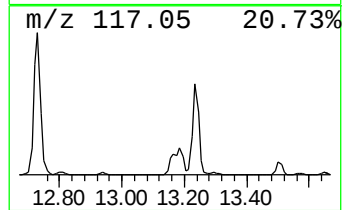
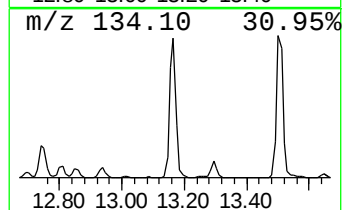
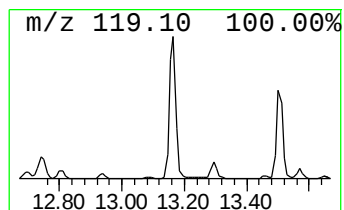
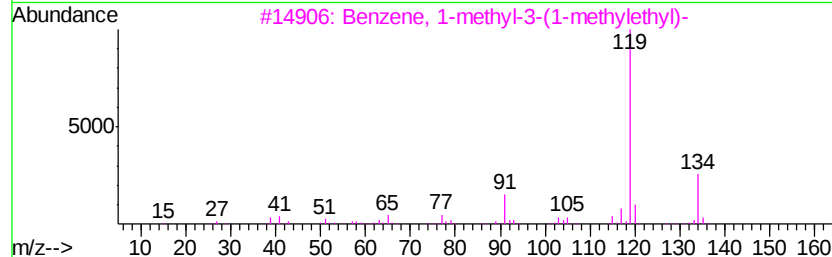
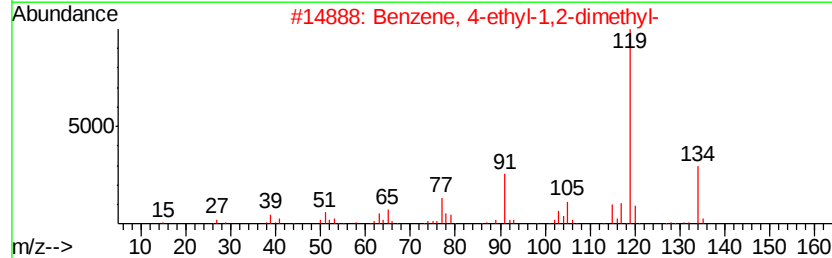
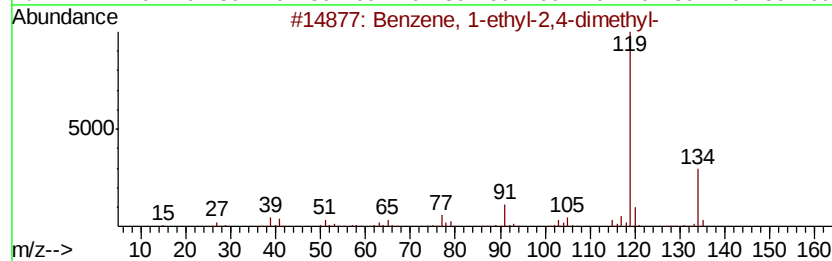
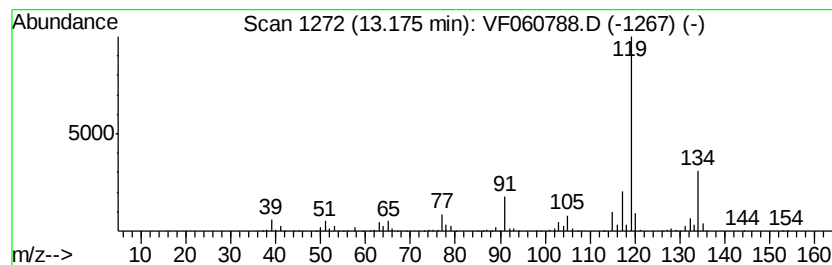
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.17	87.79 ug/l	1455590	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5		o-Cymene	134	C10H14	000527-84-4	94



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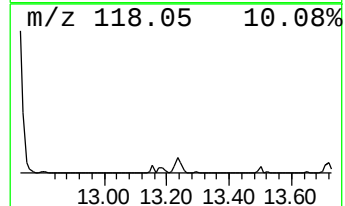
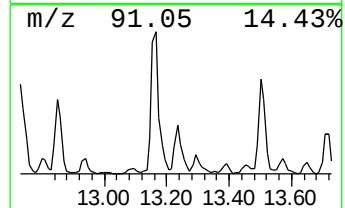
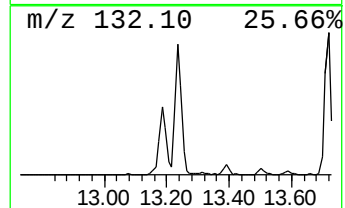
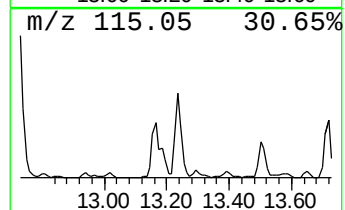
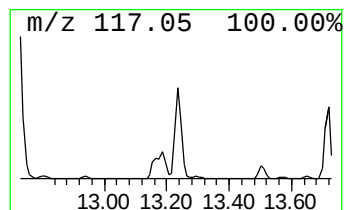
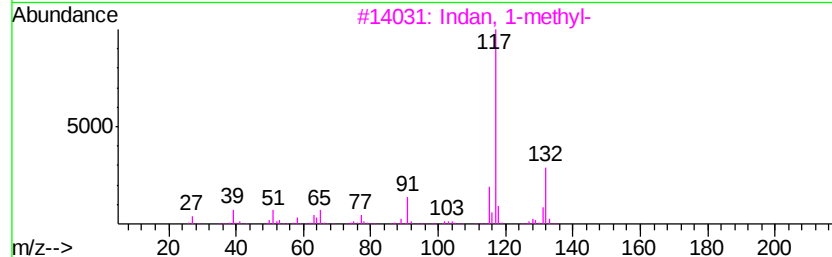
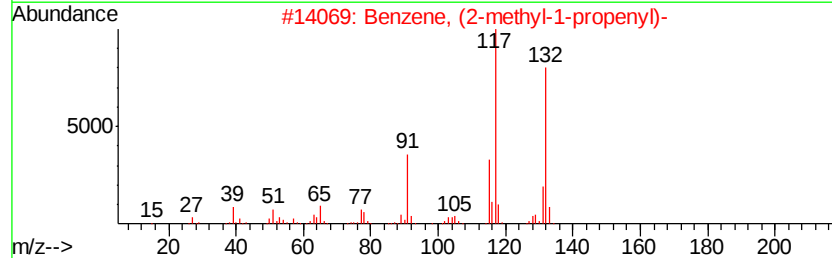
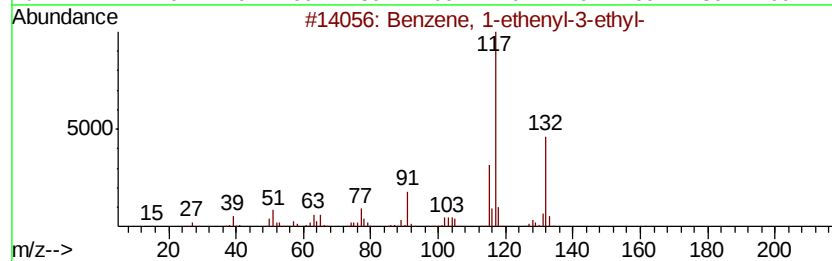
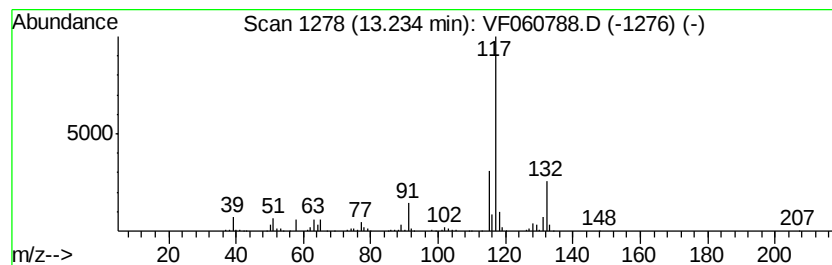
Instrument :
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.23	30.51 ug/l	505859	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 91
2		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 90
3		Indan, 1-methyl-	132 C10H12	000767-58-8 90
4		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 90
5		Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7 87



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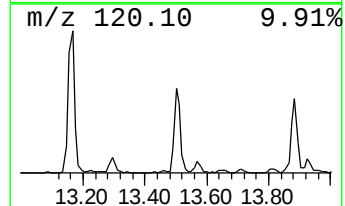
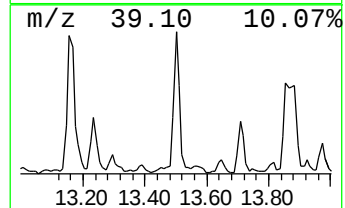
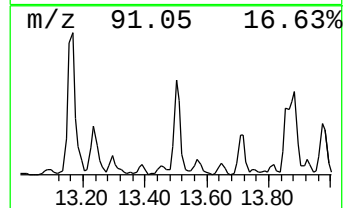
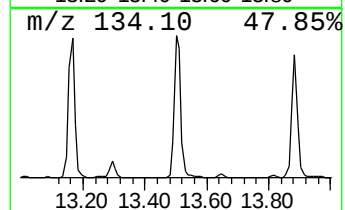
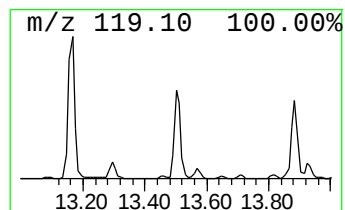
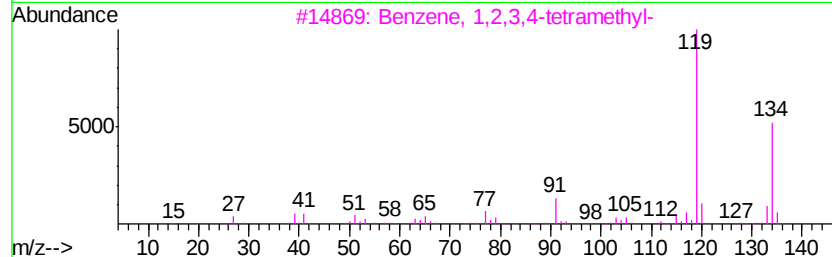
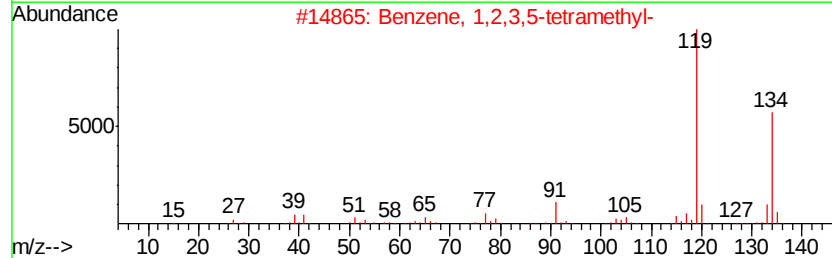
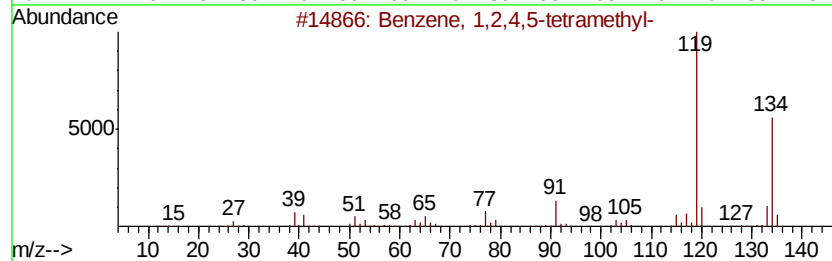
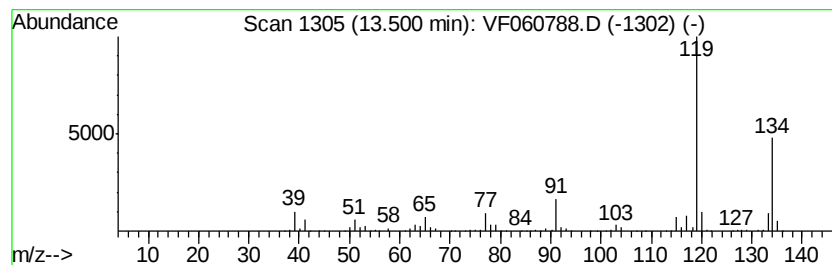
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.50	46.79 ug/l	775765	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5		o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampled :
SSI-20-TW-1

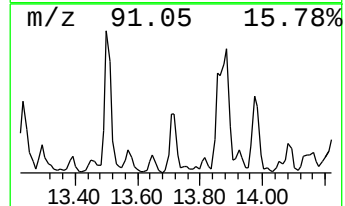
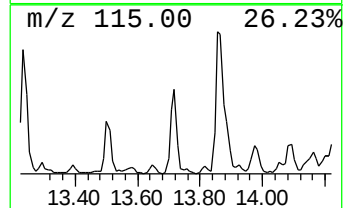
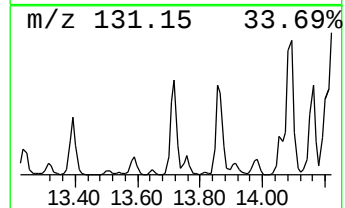
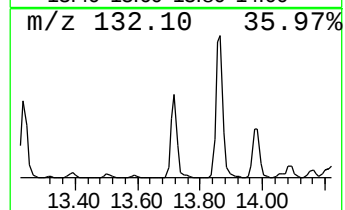
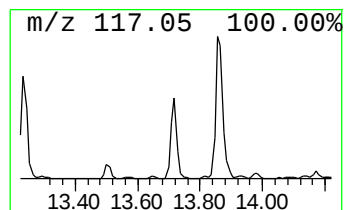
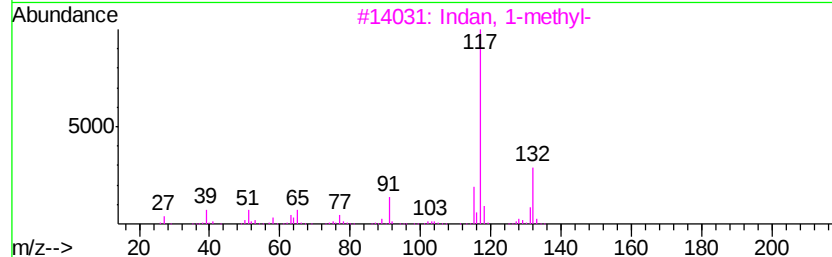
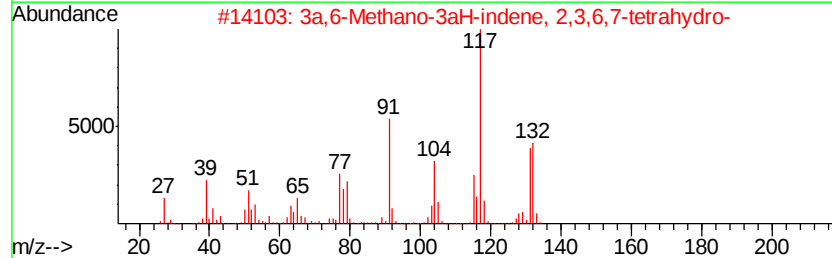
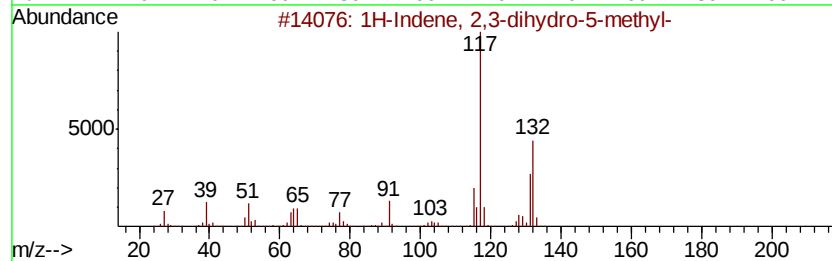
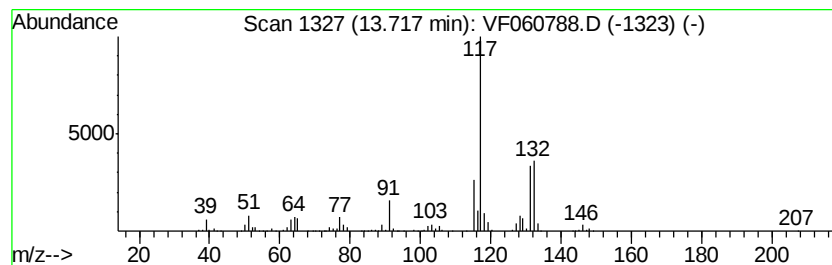
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	28.52 ug/l	472835	1,4-Dichlorobenzene-d4	12.56

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	3a,6-Methano-3aH-indene, 2,3,6,7...	132	C10H12	098640-29-0	91
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

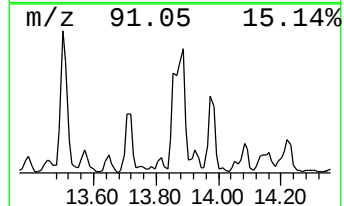
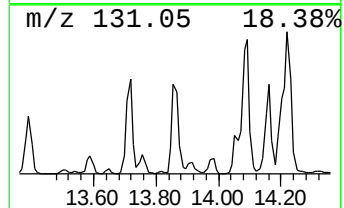
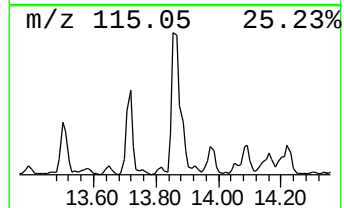
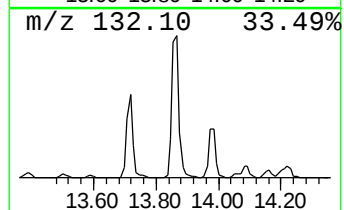
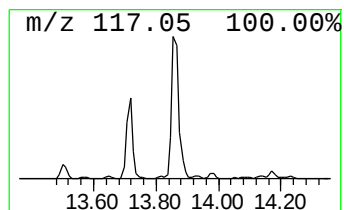
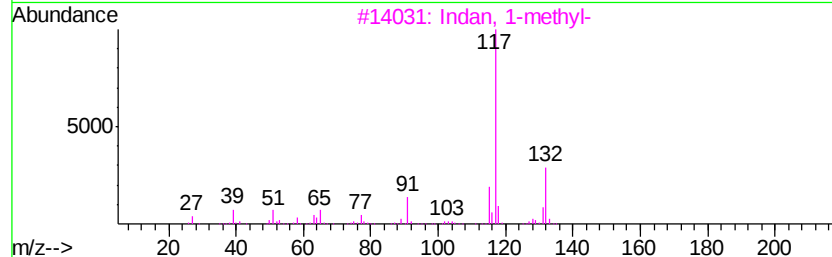
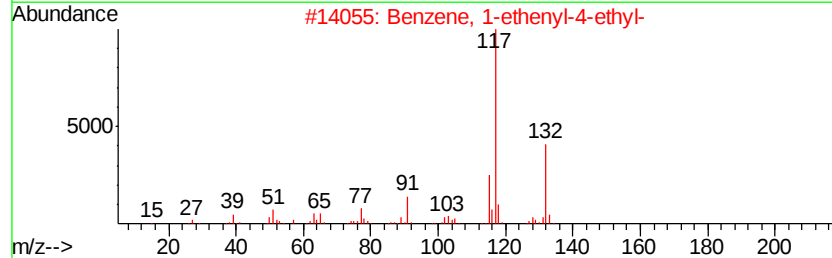
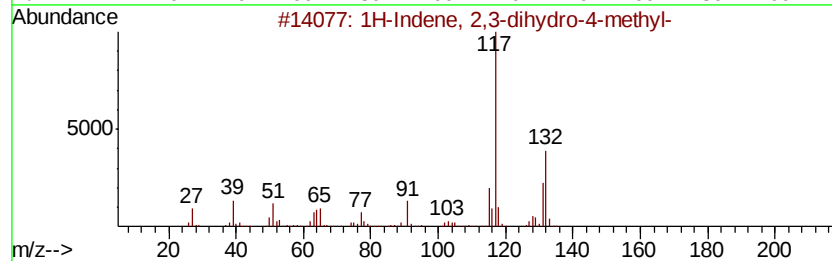
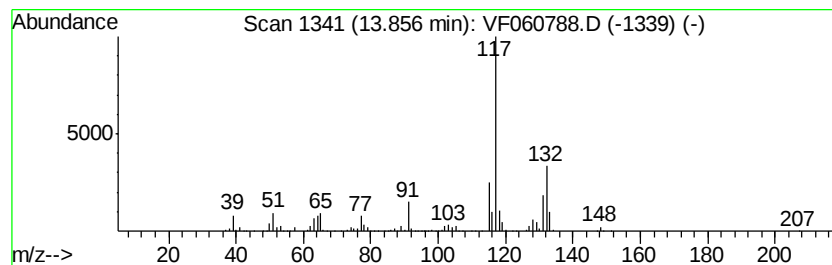
Instrument :
MSVOA_F
ClientSampleId :
SSI-20-TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	60.31 ug/l	1000020	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 93
2		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 93
3		Indan, 1-methyl-	132 C10H12	000767-58-8 87
4		1H-Indene, 2,3-dihydro-2-methyl-	132 C10H12	000824-63-5 86
5		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-20-TW-1

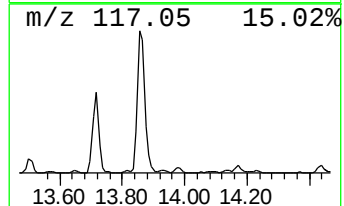
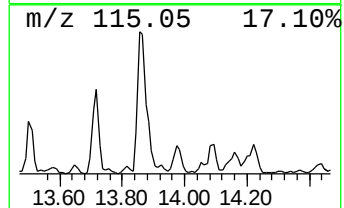
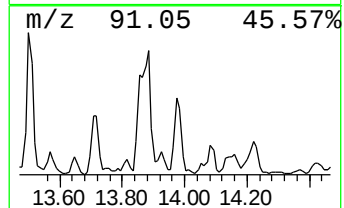
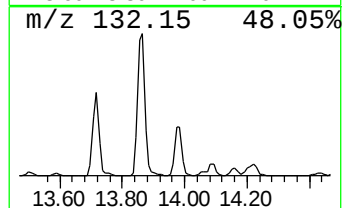
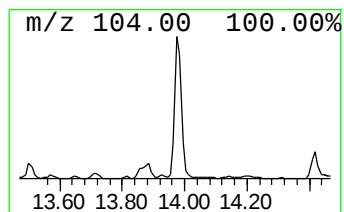
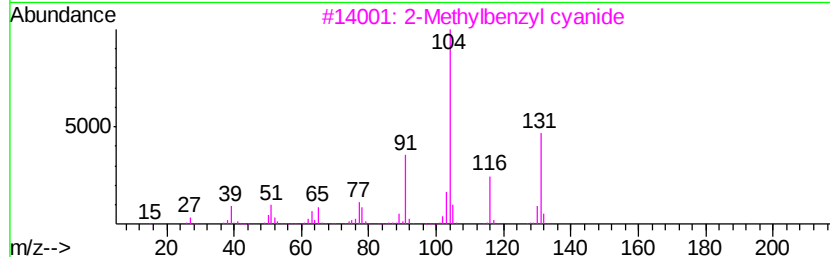
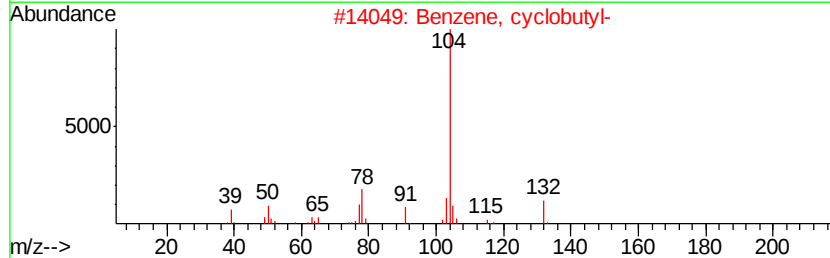
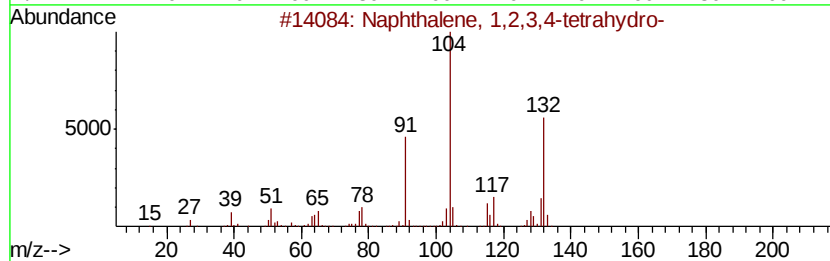
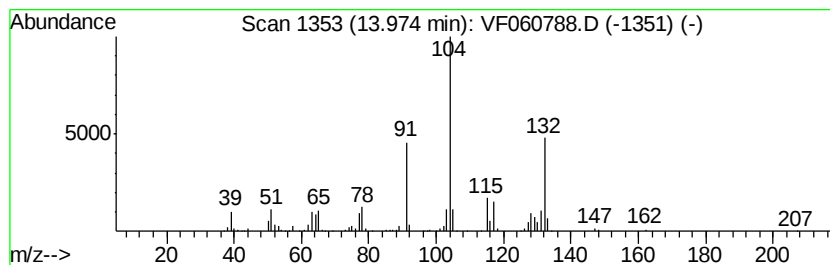
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	14.17 ug/l	234922	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
3		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	47
4		Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47
5		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

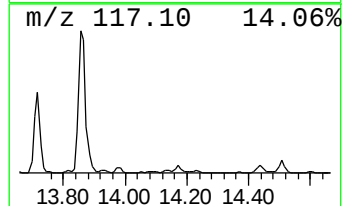
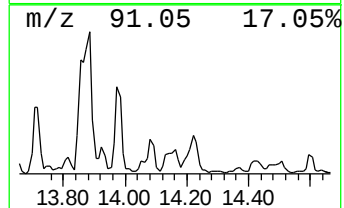
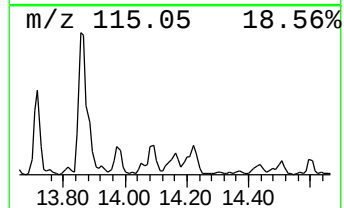
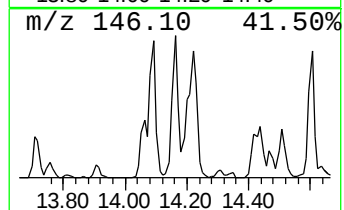
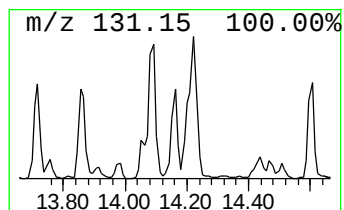
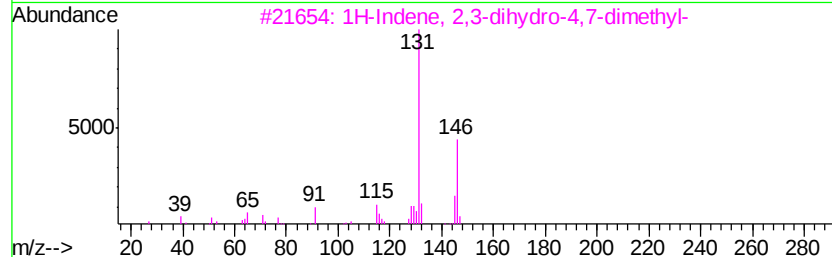
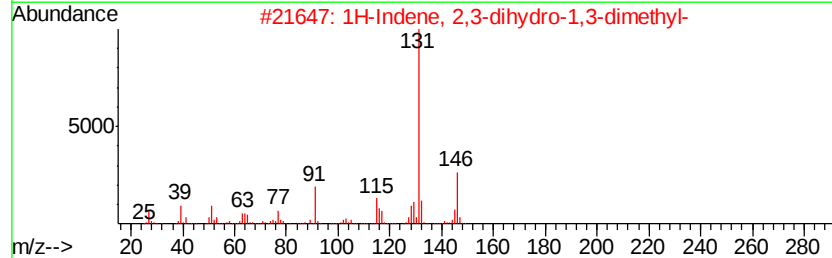
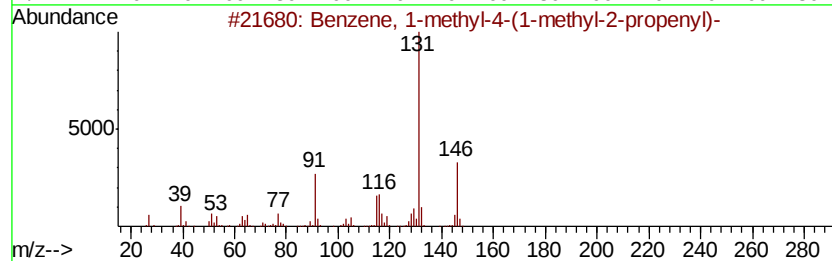
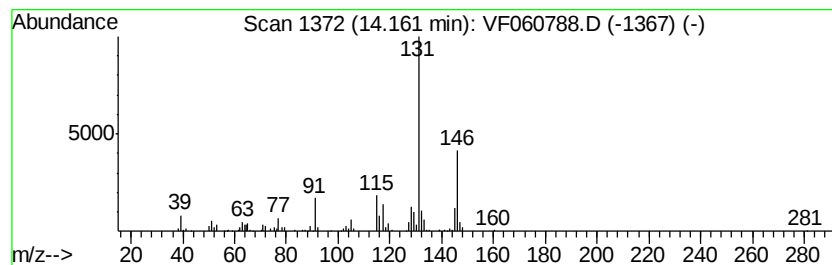
Instrument :
MSVOA_F
ClientSampled :
SSI-20-TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.16	15.92 ug/l	263900	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	91
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	91
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	90
4	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	90
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-20-TW-1

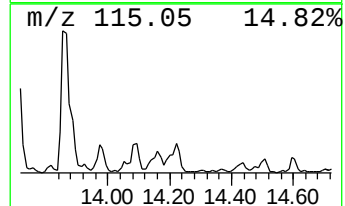
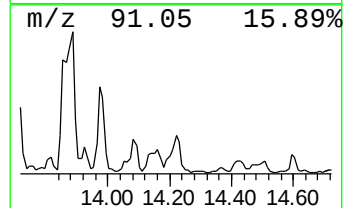
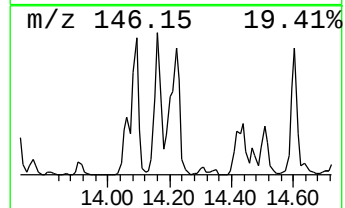
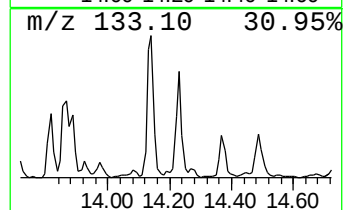
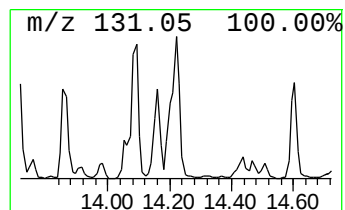
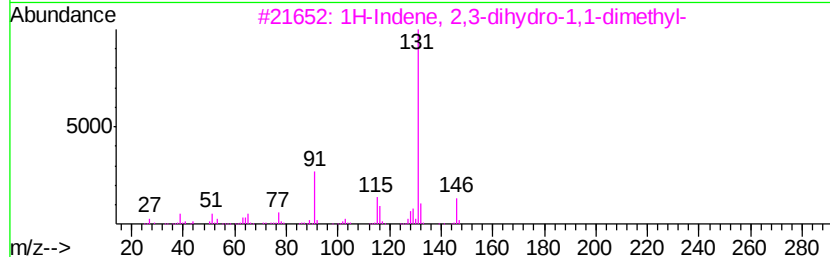
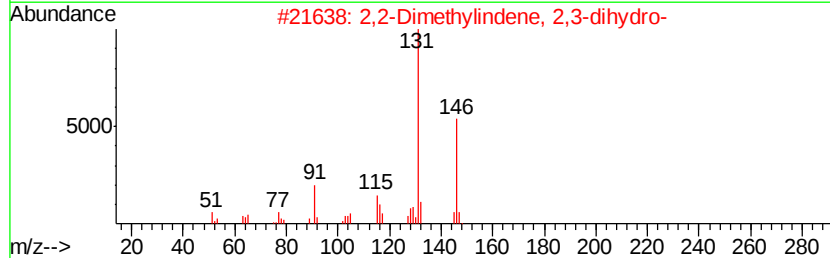
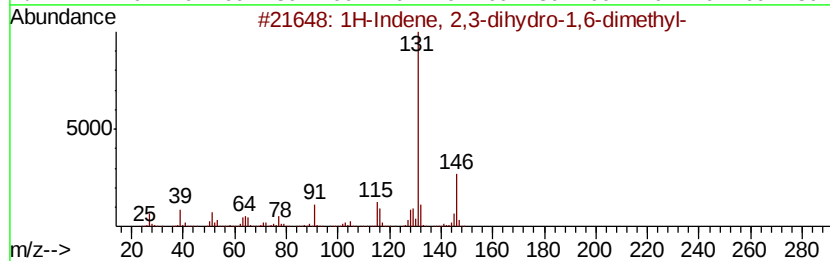
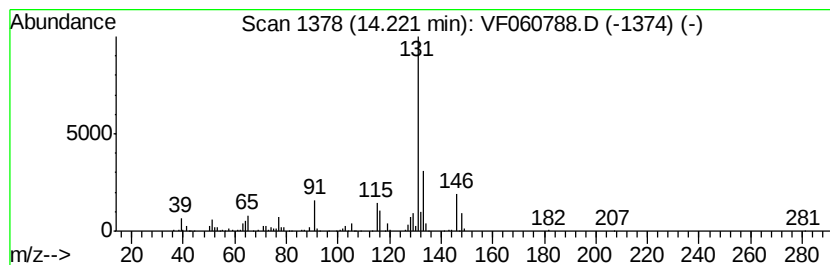
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	20.34 ug/l	337169	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76
2		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	76
3		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	76
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	76
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00µ/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

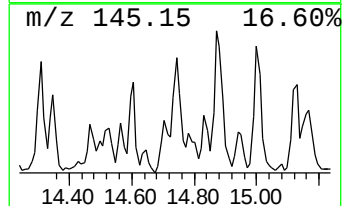
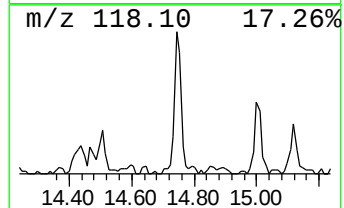
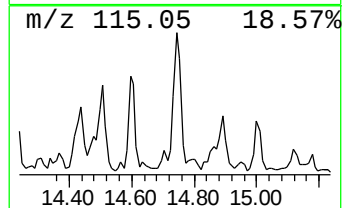
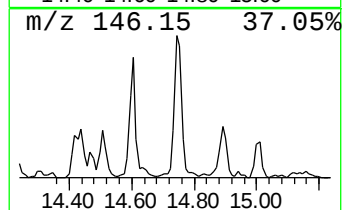
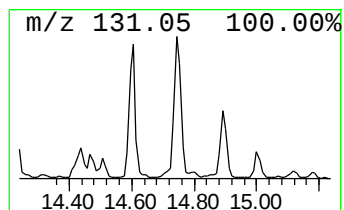
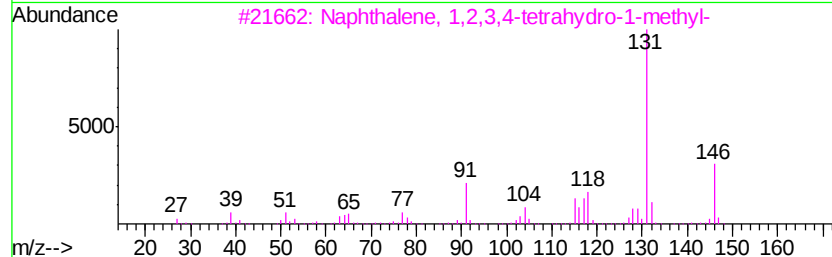
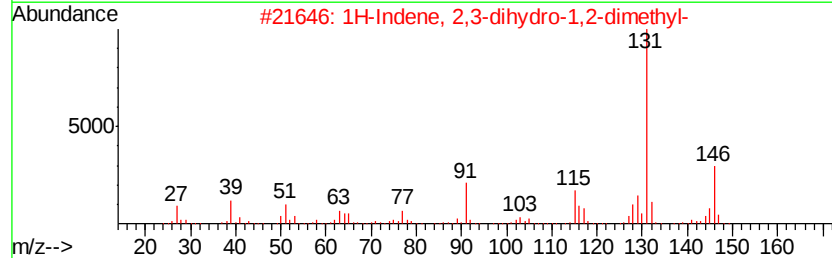
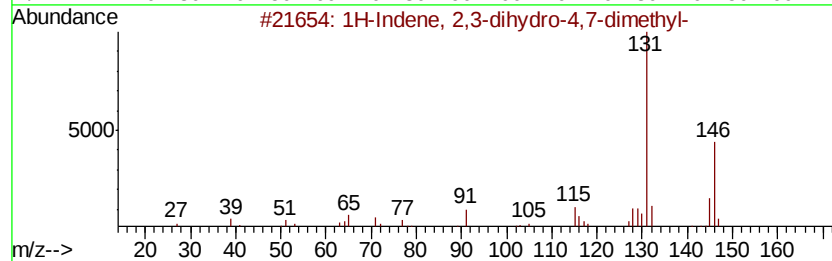
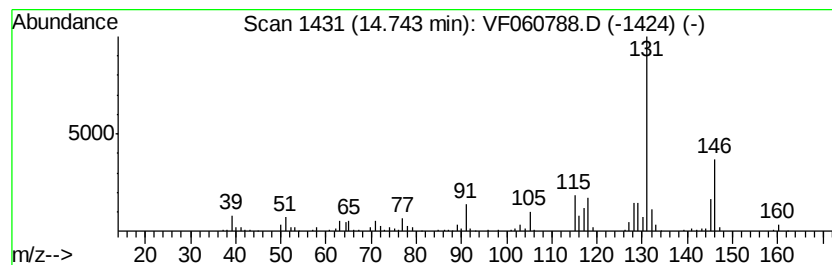
Instrument :
MSVOA_F
ClientSampleId :
SSI-20-TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	12.70 ug/l	210590	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	91
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	90
3	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	90
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3	87
5	1,3-Cyclopentadiene, 5-(trans-2-...	146 C11H14	079209-36-2	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF111918\
Data File : VF060788.D
Acq On : 19 Nov 2018 20:58
Operator : VA/AP
Sample : J5950-06
Misc : 5.00g/5mL/MSVOA-F/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-20-TW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.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, cyclopro...	12.73	74.6	ug/l	1237170	4	12.56	829037	50.0
Benzene, 1-ethyl-...	13.17	87.8	ug/l	1455590	4	12.56	829037	50.0
Benzene, 1-etheny...	13.23	30.5	ug/l	505859	4	12.56	829037	50.0
Benzene, 1,2,4,5-...	13.50	46.8	ug/l	775765	4	12.56	829037	50.0
1H-Indene, 2,3-di...	13.72	28.5	ug/l	472835	4	12.56	829037	50.0
1H-Indene, 2,3-di...	13.86	60.3	ug/l	1000020	4	12.56	829037	50.0
Naphthalene, 1,2,...	13.97	14.2	ug/l	234922	4	12.56	829037	50.0
Benzene, 1-methyl...	14.16	15.9	ug/l	263900	4	12.56	829037	50.0
1H-Indene, 2,3-di...	14.22	20.3	ug/l	337169	4	12.56	829037	50.0
1H-Indene, 2,3-di...	14.74	12.7	ug/l	210590	4	12.56	829037	50.0