

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF082418\
 Data File : VF060110.D
 Acq On : 24 Aug 2018 19:58
 Operator : VA/AP
 Sample : J4282-13
 Misc : 5.36g/5mL/MSVOA F/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 SU-02-082318-B

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.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	1.084	53	56	59	rBV2	15668	19954	2.38%	0.202%
2	1.360	78	84	91	rVB5	4605	15997	1.91%	0.162%
3	1.833	129	132	136	rBV	7926	12976	1.55%	0.131%
4	2.277	173	177	179	rBV2	20076	41081	4.89%	0.416%
5	2.326	179	182	189	rVB	23424	51115	6.09%	0.517%
6	2.474	194	197	203	rVB4	5265	11996	1.43%	0.121%
7	4.130	358	365	372	rBV	8254	30087	3.58%	0.304%
8	4.278	372	380	393	rVB	46050	143153	17.05%	1.448%
9	5.027	447	456	468	rBV2	184065	665953	79.31%	6.738%
10	5.628	509	517	524	rBV4	4208	13949	1.66%	0.141%
11	5.757	524	530	544	rBV	192396	516213	61.47%	5.223%
12	7.176	667	674	677	rBV4	8428	22745	2.71%	0.230%
13	7.620	714	719	722	rBV5	4005	10906	1.30%	0.110%
14	7.698	722	727	731	rVV	282281	657730	78.33%	6.655%
15	7.767	731	734	743	rVV	275059	621310	73.99%	6.286%
16	7.905	743	748	753	rVB6	7398	24131	2.87%	0.244%
17	8.290	781	787	790	rBV4	5760	16091	1.92%	0.163%
18	8.339	790	792	798	rVB3	6055	14489	1.73%	0.147%
19	8.526	807	811	813	rBV3	5123	12882	1.53%	0.130%
20	8.595	813	818	825	rVB4	14797	41051	4.89%	0.415%
21	8.763	828	835	839	rBV2	9223	21273	2.53%	0.215%
22	9.098	864	869	874	rBV3	11297	28778	3.43%	0.291%
23	9.187	874	878	880	rVV3	4113	9999	1.19%	0.101%
24	9.246	880	884	890	rVB6	8535	26204	3.12%	0.265%
25	9.423	897	902	906	rVB3	7051	18694	2.23%	0.189%
26	9.680	923	928	933	rVV4	5213	15912	1.89%	0.161%
27	9.749	933	935	939	rVV4	4290	11307	1.35%	0.114%
28	9.837	941	944	945	rVV2	5186	9803	1.17%	0.099%
29	9.906	945	951	959	rVV	284157	639613	76.17%	6.472%
30	10.025	959	963	968	rVV2	49881	125334	14.93%	1.268%
31	10.104	968	971	976	rVB5	3700	10833	1.29%	0.110%
32	10.261	982	987	992	rBV	130720	291472	34.71%	2.949%
33	10.350	992	996	1004	rVB3	14504	44928	5.35%	0.455%
34	10.468	1004	1008	1015	rVB5	3870	11539	1.37%	0.117%

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 ClientSampleId :
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Integration Parameters: RTEINT.P

Integrator: RTE
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35	10.636	1019	1025	1030	rVB5	11355	30031	3.58%	0.304%
36	10.774	1034	1039	1040	rBV2	20295	33943	4.04%	0.343%
37	10.823	1040	1044	1048	rVV	104711	209565	24.96%	2.120%
38	10.882	1048	1050	1058	rVV7	5473	14350	1.71%	0.145%
39	11.050	1062	1067	1070	rBV5	3950	8746	1.04%	0.088%
40	11.148	1073	1077	1080	rBV6	8074	21769	2.59%	0.220%
41	11.217	1080	1084	1091	rVV3	20560	52697	6.28%	0.533%
42	11.355	1095	1098	1102	rVB4	14955	30700	3.66%	0.311%
43	11.503	1109	1113	1119	rBV	246131	425519	50.67%	4.305%
44	11.641	1124	1127	1128	rVV	17071	26246	3.13%	0.266%
45	11.681	1128	1131	1134	rVV	45896	80873	9.63%	0.818%
46	11.740	1134	1137	1139	rVV	49518	81028	9.65%	0.820%
47	11.799	1139	1143	1150	rVB	302620	638111	75.99%	6.456%
48	11.907	1150	1154	1158	rBV	279438	416137	49.56%	4.210%
49	12.095	1168	1173	1177	rBV	112743	198307	23.62%	2.006%
50	12.282	1188	1192	1198	rBV	568382	839729	100.00%	8.496%
51	12.371	1198	1201	1204	rVV3	12543	24956	2.97%	0.253%
52	12.450	1207	1209	1211	rBV2	6038	9311	1.11%	0.094%
53	12.499	1211	1214	1222	rVB	35874	91363	10.88%	0.924%
54	12.627	1223	1227	1230	rBV	263629	413053	49.19%	4.179%
55	12.676	1230	1232	1239	rVB	243009	404953	48.22%	4.097%
56	12.795	1239	1244	1247	rBV2	56826	117237	13.96%	1.186%
57	12.844	1247	1249	1251	rVV	63947	108053	12.87%	1.093%
58	12.893	1251	1254	1259	rVV2	111929	253098	30.14%	2.561%
59	13.002	1262	1265	1267	rBV3	10314	16938	2.02%	0.171%
60	13.071	1269	1272	1276	rVB	29107	55185	6.57%	0.558%
61	13.159	1276	1281	1285	rBV	45395	119715	14.26%	1.211%
62	13.218	1285	1287	1293	rVV	75754	133482	15.90%	1.351%
63	13.317	1295	1297	1299	rVV2	25744	41643	4.96%	0.421%
64	13.396	1303	1305	1308	rVV3	12895	21148	2.52%	0.214%
65	13.445	1308	1310	1312	rVV2	13268	17786	2.12%	0.180%
66	13.485	1312	1314	1318	rVV	19126	29933	3.56%	0.303%
67	13.563	1318	1322	1324	rVV	30950	51812	6.17%	0.524%
68	13.603	1324	1326	1329	rVV	53693	81131	9.66%	0.821%
69	13.642	1329	1330	1332	rVV	12626	11031	1.31%	0.112%
70	13.682	1332	1334	1335	rVV	8285	13139	1.56%	0.133%
71	13.770	1339	1343	1346	rVV2	38958	59918	7.14%	0.606%

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 Stop Thrs : 0 Peak Location: TOP

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72	13.869	1351	1353	1355	rBV2	9234	12098	1.44%	0.122%
73	13.918	1355	1358	1362	rVV2	39457	81063	9.65%	0.820%
74	13.977	1362	1364	1366	rVB2	10252	13664	1.63%	0.138%
75	14.037	1366	1370	1374	rVB	23206	43231	5.15%	0.437%
76	14.106	1374	1377	1379	rBV3	6648	11823	1.41%	0.120%
77	14.135	1379	1380	1383	rVB	11896	13467	1.60%	0.136%
78	14.214	1383	1388	1390	rBV4	14852	35439	4.22%	0.359%
79	14.273	1391	1394	1400	rVB3	16021	33075	3.94%	0.335%
80	14.352	1400	1402	1405	rBV4	4737	8522	1.01%	0.086%
81	14.520	1412	1419	1421	rBV	35020	66280	7.89%	0.671%
82	14.648	1430	1432	1435	rVB	10430	14755	1.76%	0.149%
83	14.796	1444	1447	1451	rVB2	17557	28639	3.41%	0.290%
84	15.052	1470	1473	1478	rBV4	7002	13512	1.61%	0.137%
85	15.170	1482	1485	1487	rVB2	12982	17647	2.10%	0.179%
86	15.219	1487	1490	1492	rBV4	5059	8880	1.06%	0.090%
87	15.338	1498	1502	1505	rBV	28542	46298	5.51%	0.468%
88	15.466	1512	1515	1519	rBV	21142	34057	4.06%	0.345%
89	15.702	1537	1539	1543	rVB3	4498	8783	1.05%	0.089%

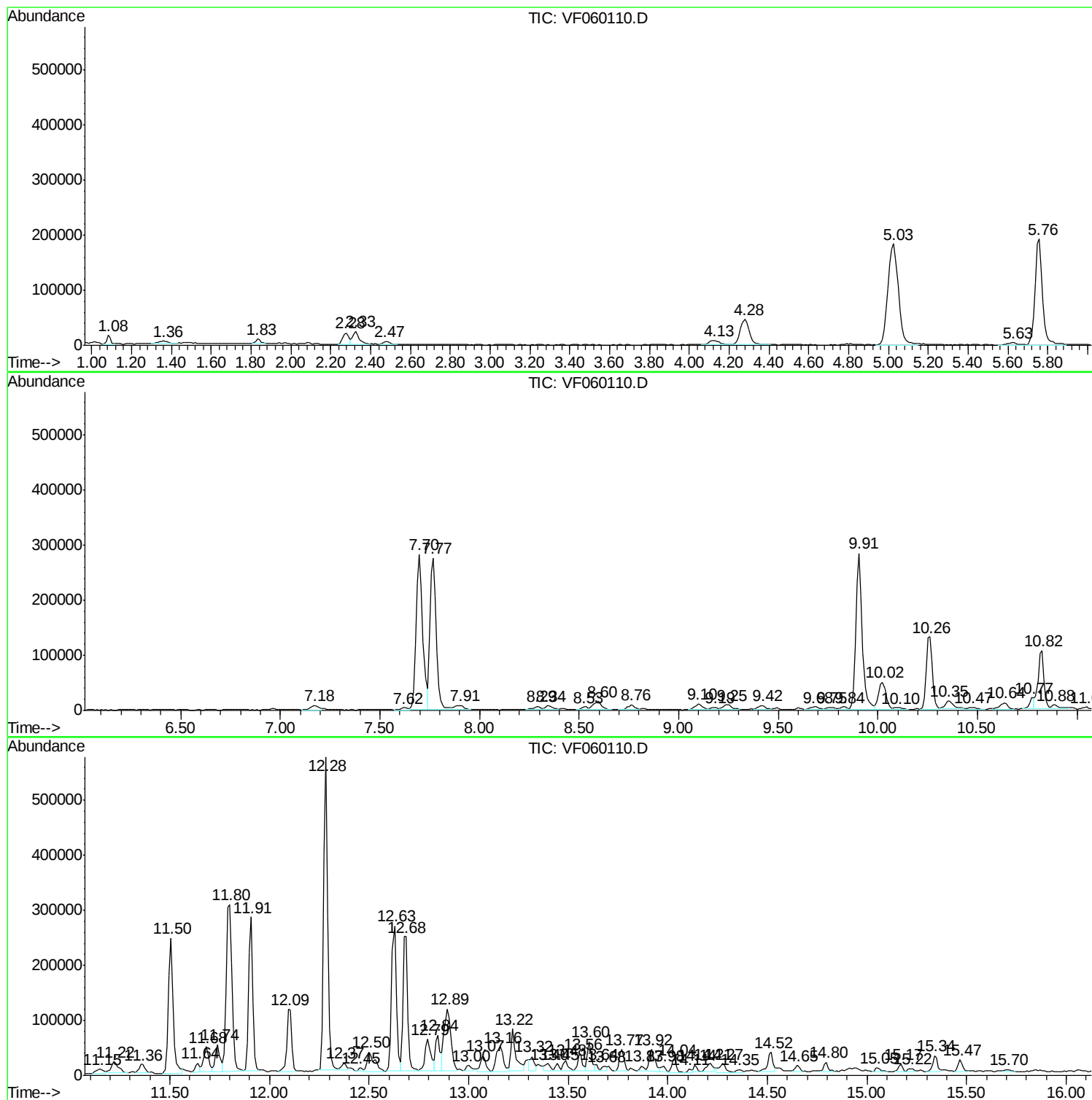
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.M
Quant Title : SW846 8260

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



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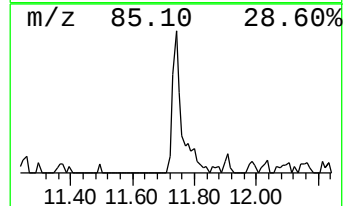
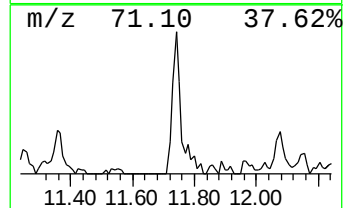
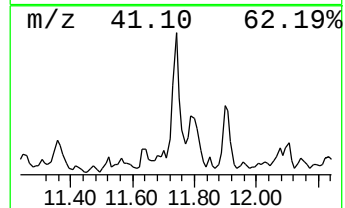
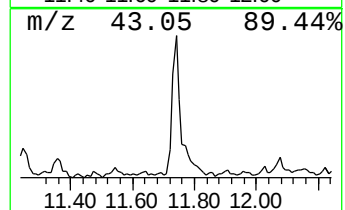
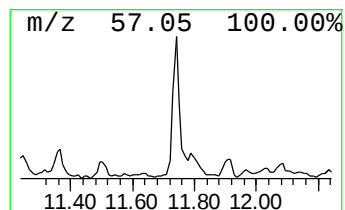
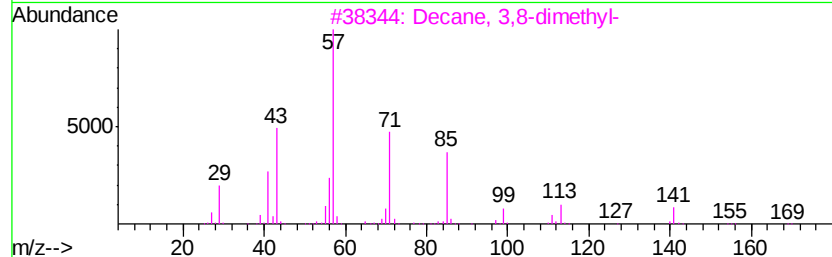
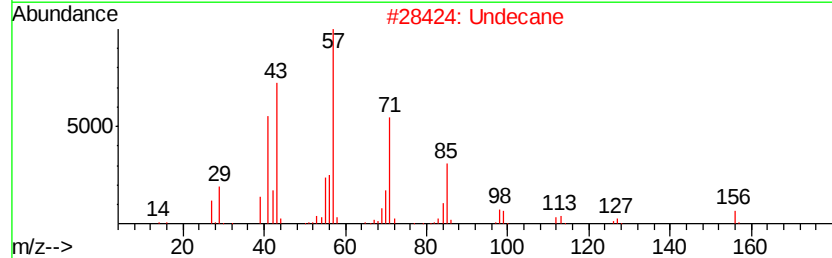
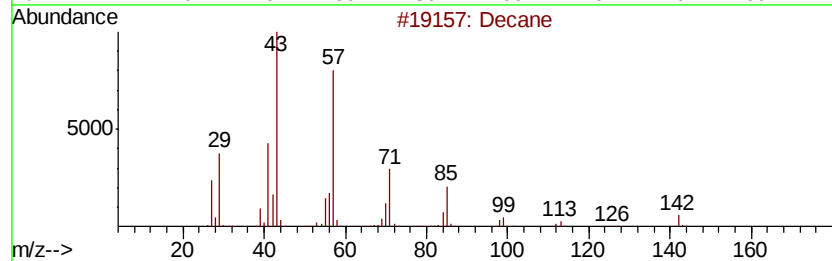
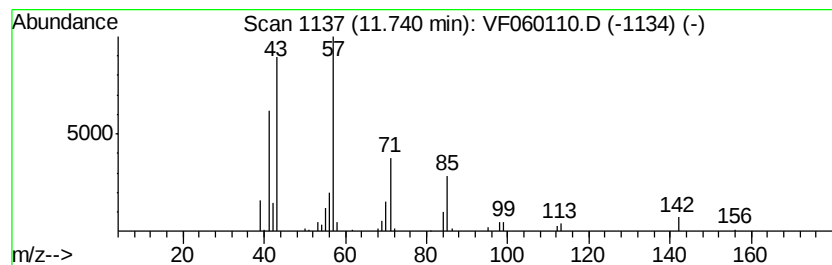
Instrument :
MSVOA_F
ClientSampleId :
SU-02-082318-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.M
Quant Title : SW846 8260

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

Peak Number 1 Decane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	9.81 ug/l	81028	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane	142 C10H22	000124-18-5	93
2	Undecane	156 C11H24	001120-21-4	78
3	Decane, 3,8-dimethyl-	170 C12H26	017312-55-9	64
4	1-Iodo-2-methylundecane	296 C12H25I	073105-67-6	64
5	Decane, 2,9-dimethyl-	170 C12H26	001002-17-1	59



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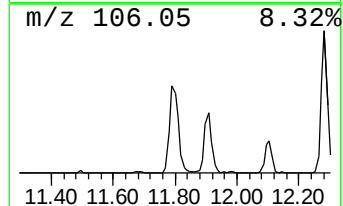
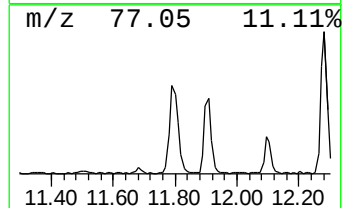
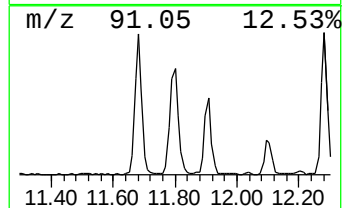
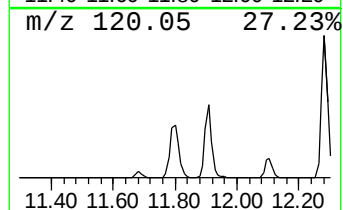
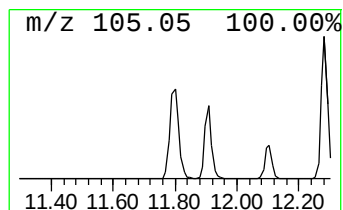
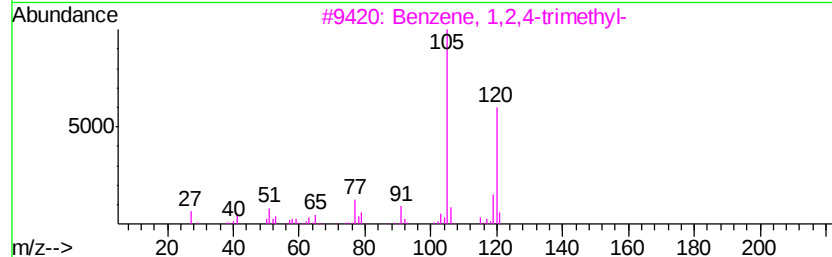
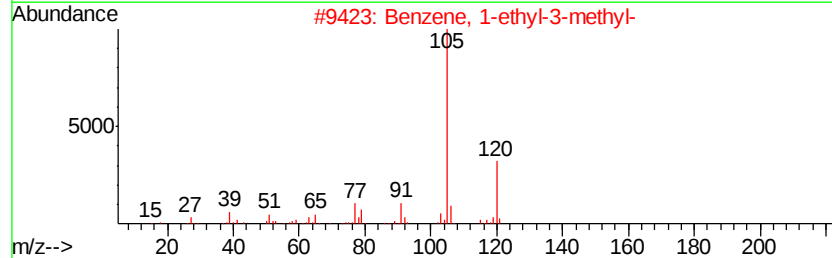
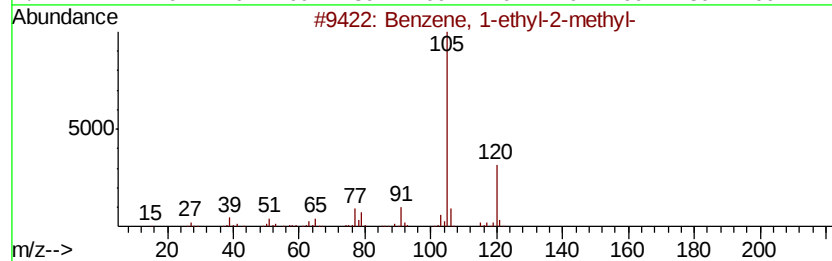
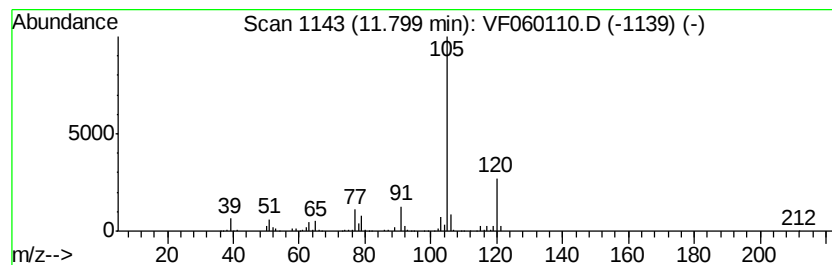
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Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	77.24 ug/l	638111	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



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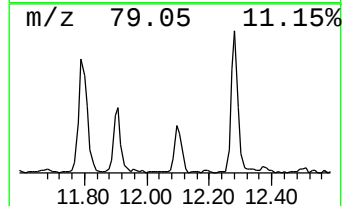
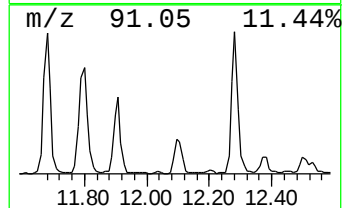
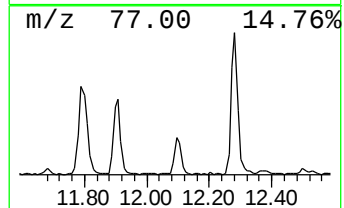
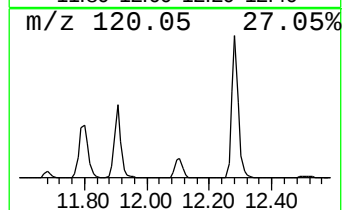
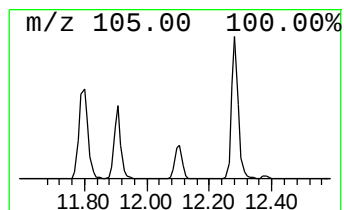
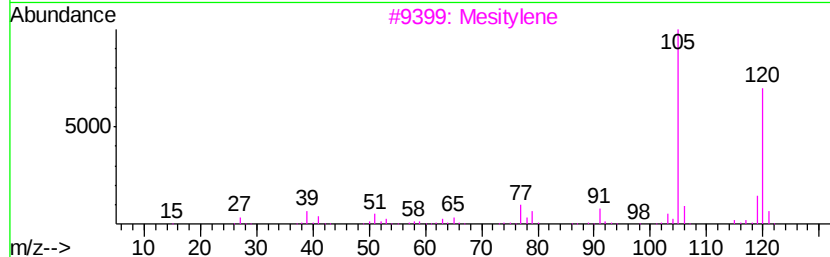
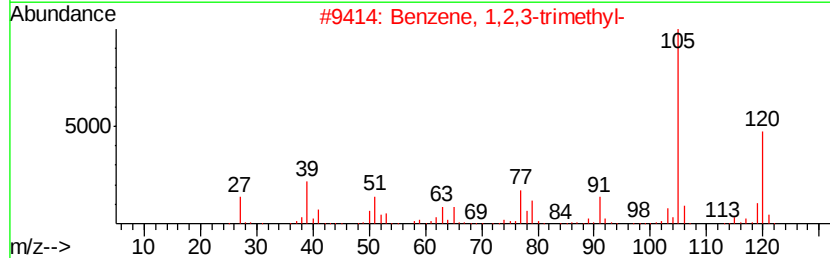
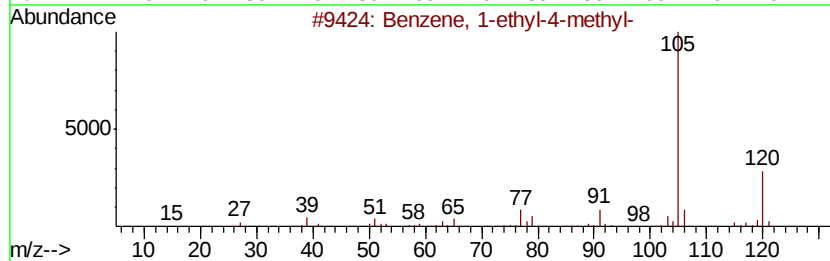
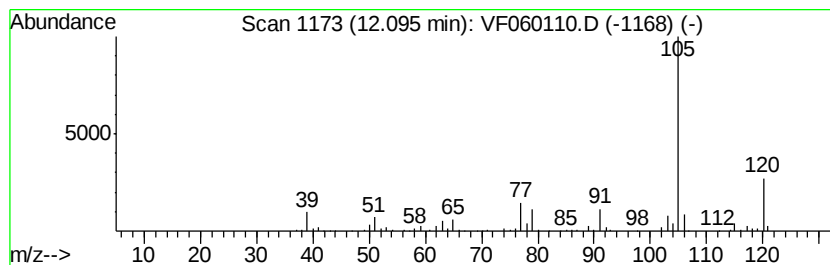
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	24.01 ug/l	198307	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



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ClientSampleID :
SU-02-082318-B

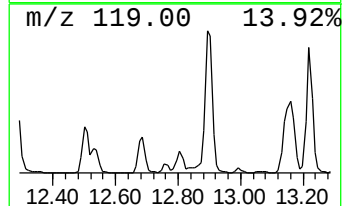
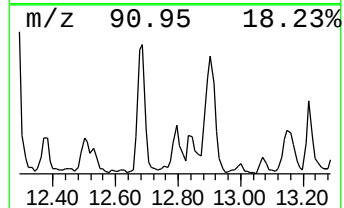
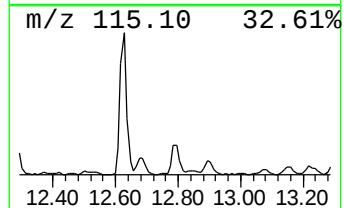
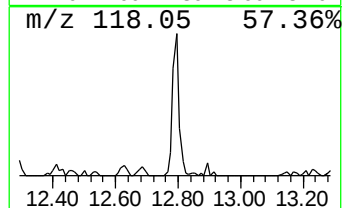
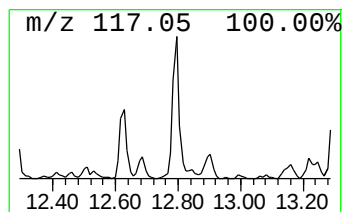
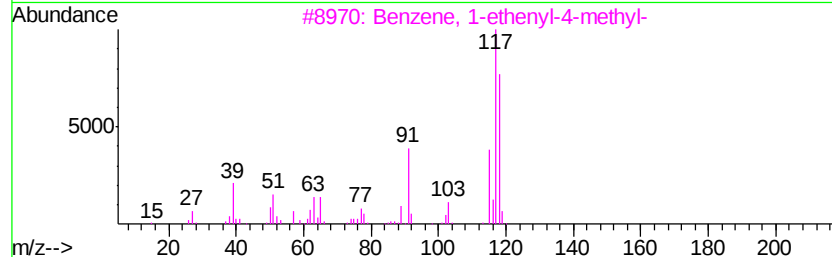
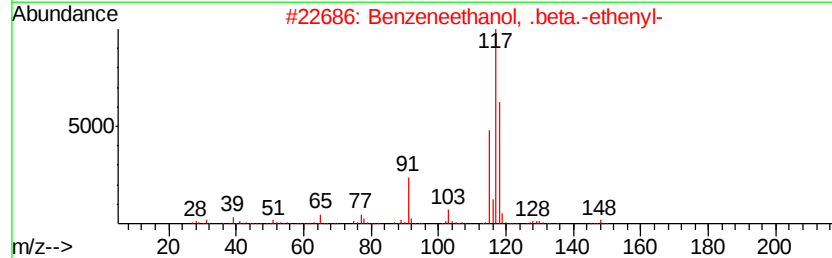
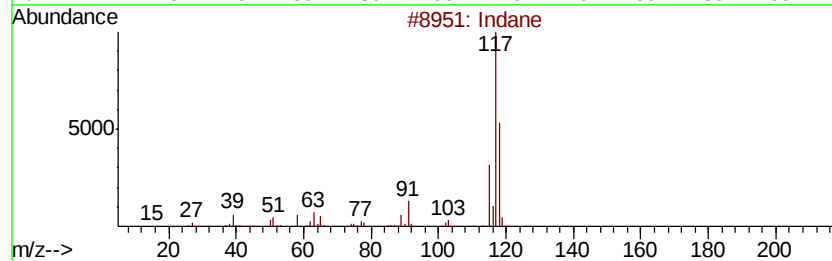
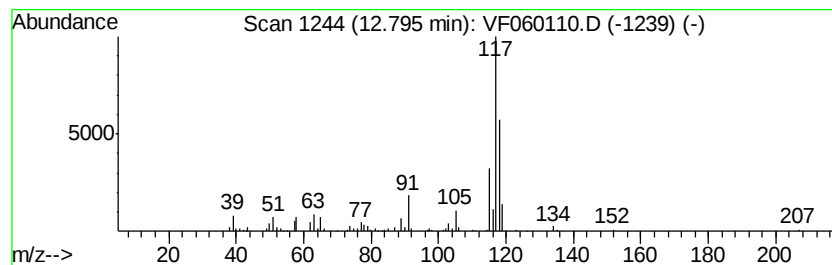
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	14.19 ug/l	117237	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	93
2		Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	80
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	74
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	68
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SU-02-082318-B

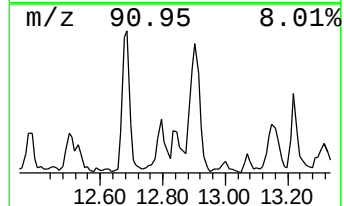
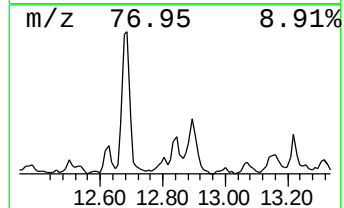
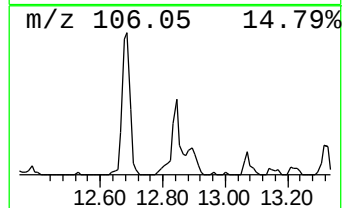
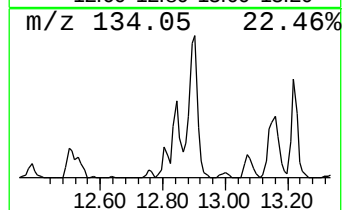
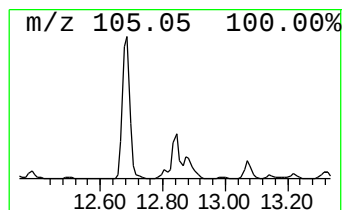
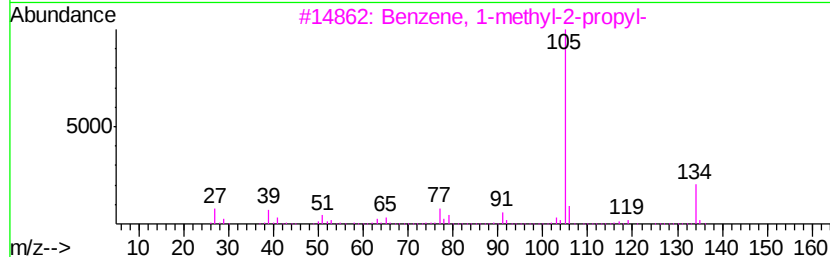
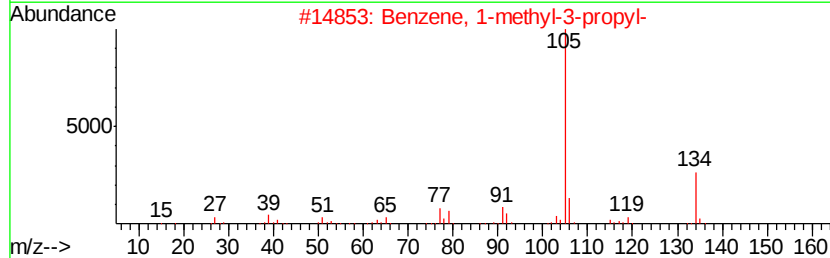
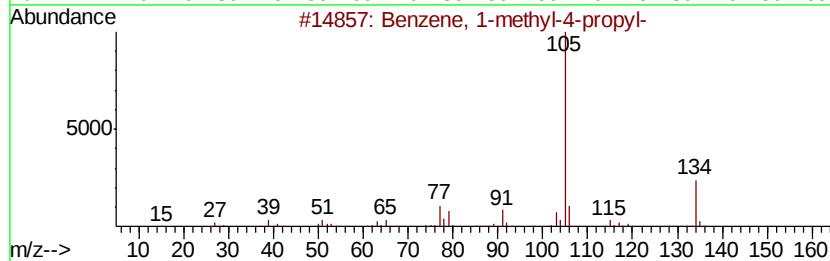
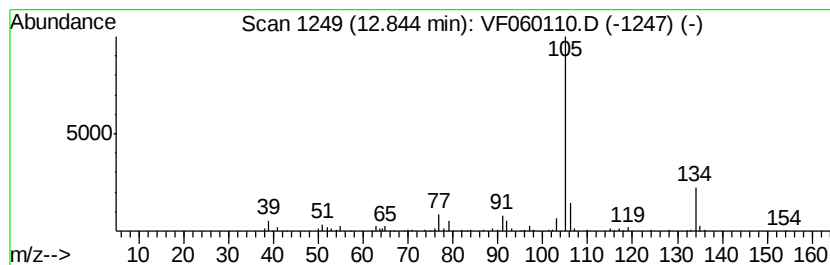
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-methyl-4-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.84	13.08 ug/l	108053	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	83
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	80
4		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	64
5		2-Tolylloxirane	134	C9H10O	002783-26-8	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

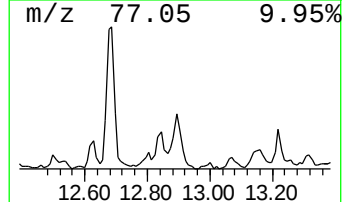
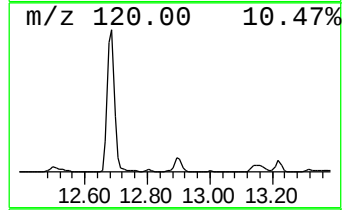
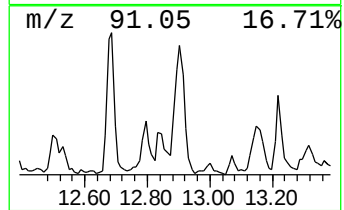
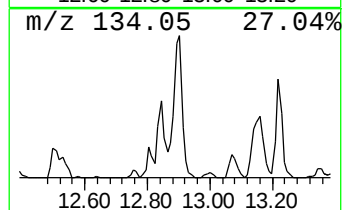
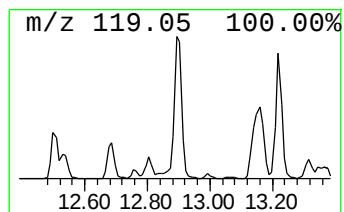
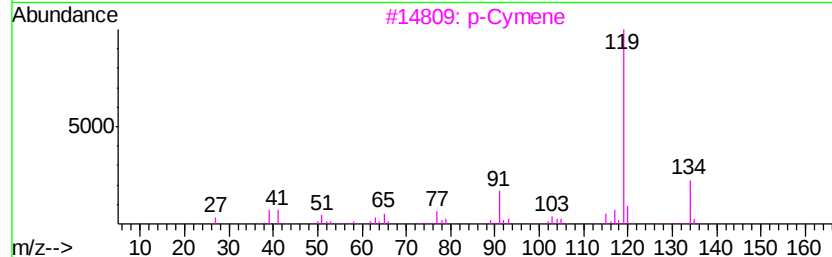
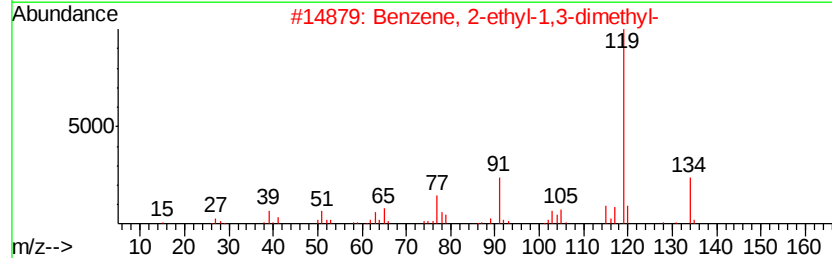
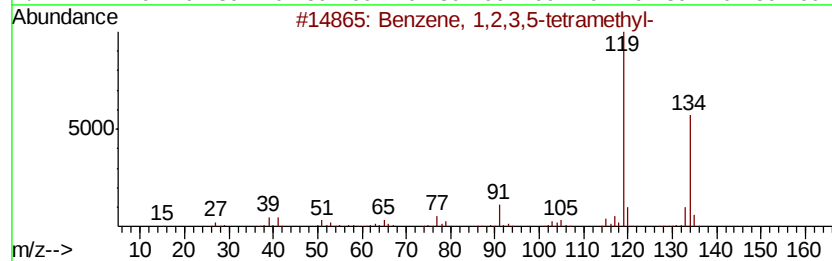
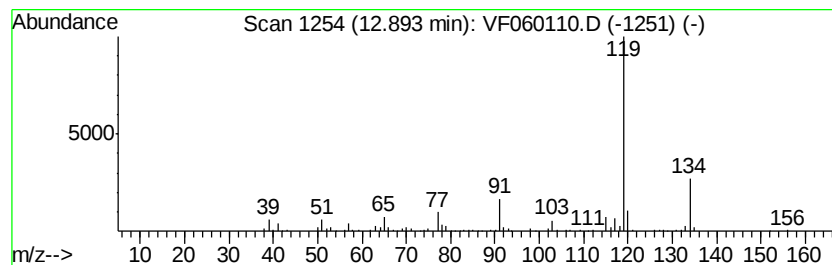
Instrument :
MSVOA_F
ClientSampled :
SU-02-082318-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	30.64 ug/l	253098	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7 91
2		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 91
3		p-Cymene	134 C10H14	000099-87-6 91
4		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 91
5		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SU-02-082318-B

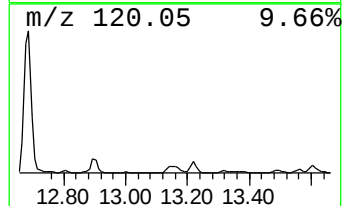
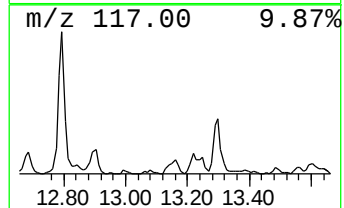
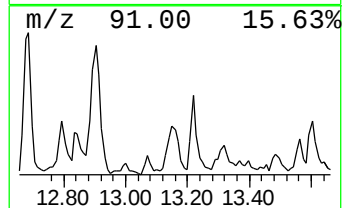
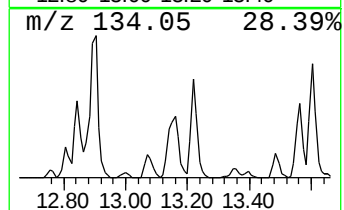
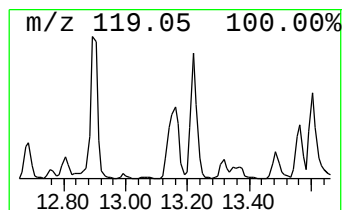
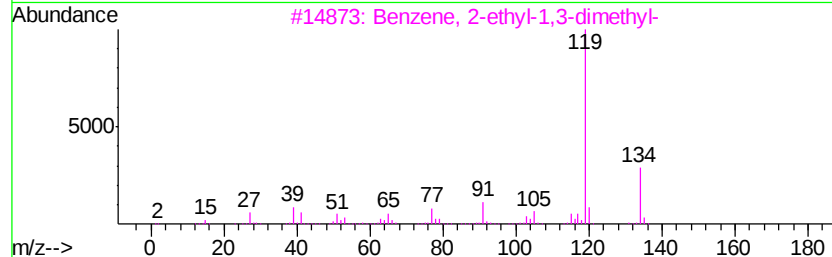
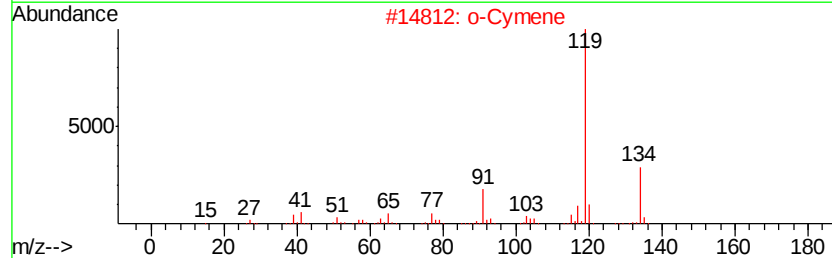
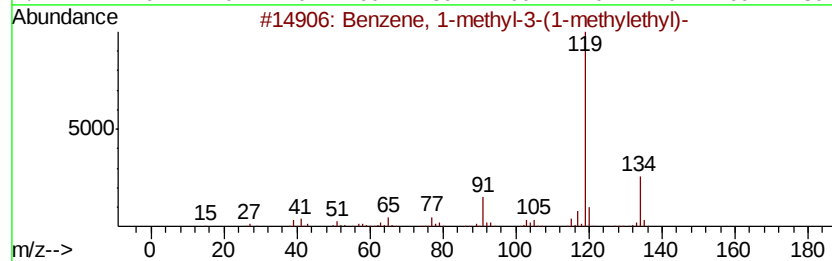
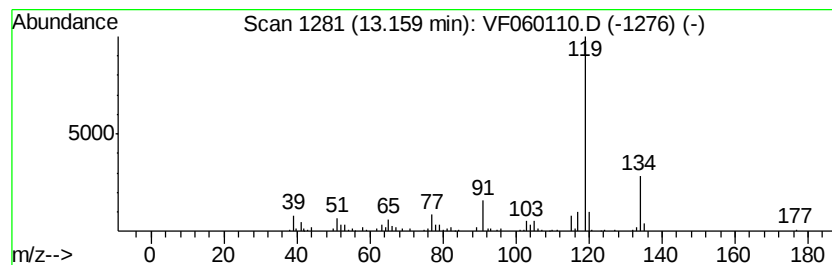
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-3-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	14.49 ug/l	119715	1,4-Dichlorobenzene-d4	12.63

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

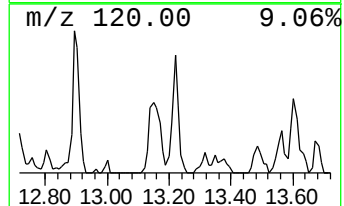
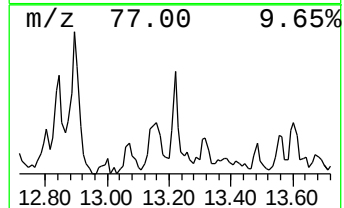
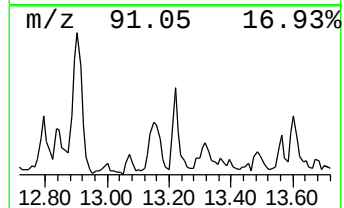
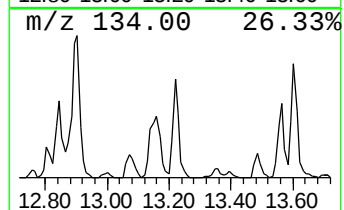
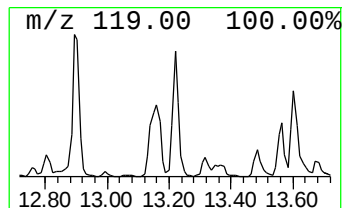
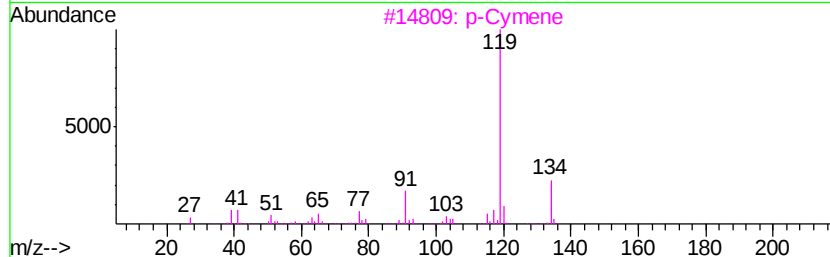
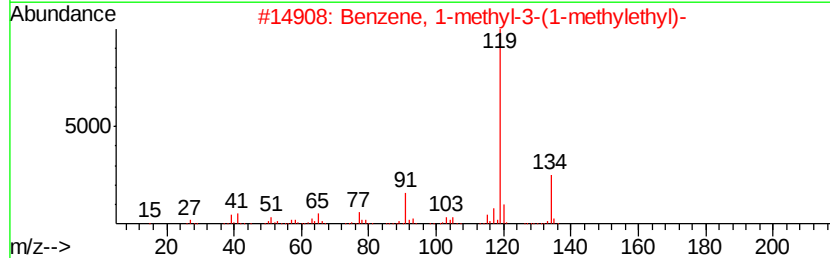
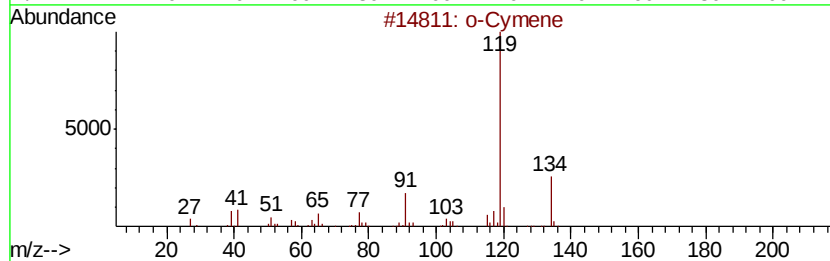
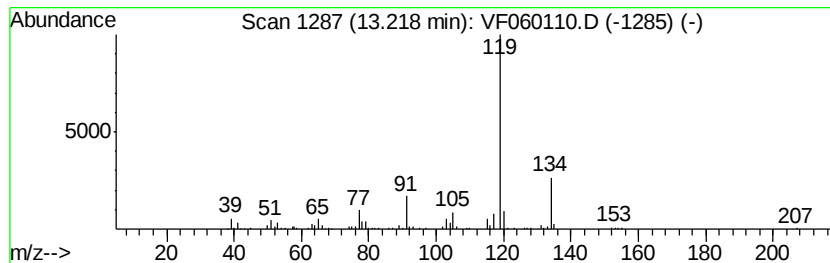
Instrument :
MSVOA_F
ClientSampleId :
SU-02-082318-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.M
Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	16.16 ug/l	133482	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 95
2		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 94
3		p-Cymene	134 C10H14	000099-87-6 94
4		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 91
5		Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampled :
SU-02-082318-B

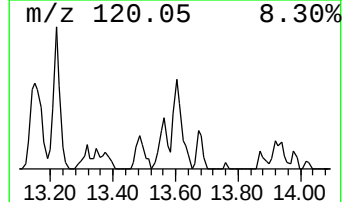
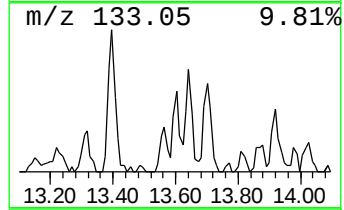
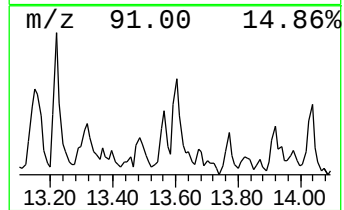
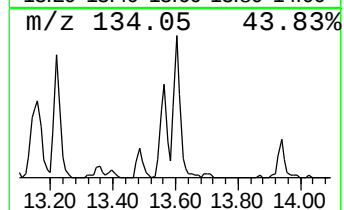
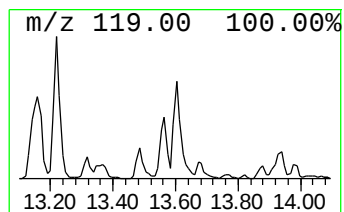
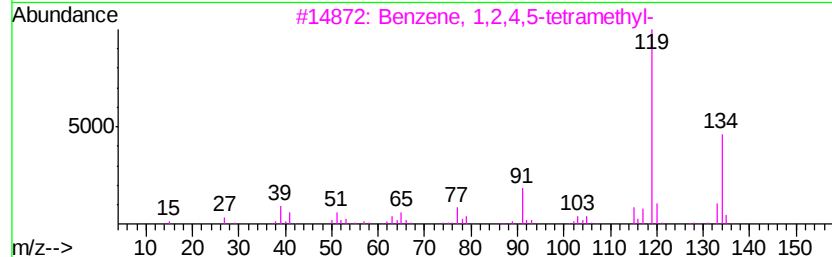
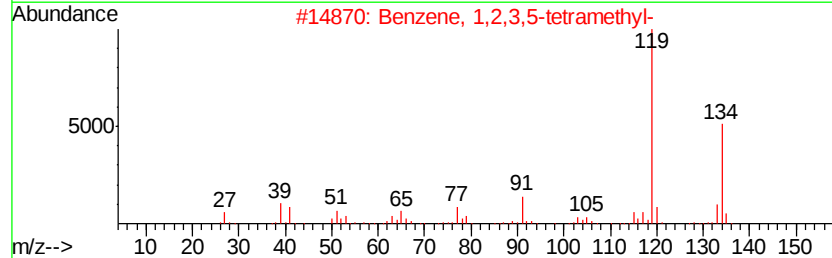
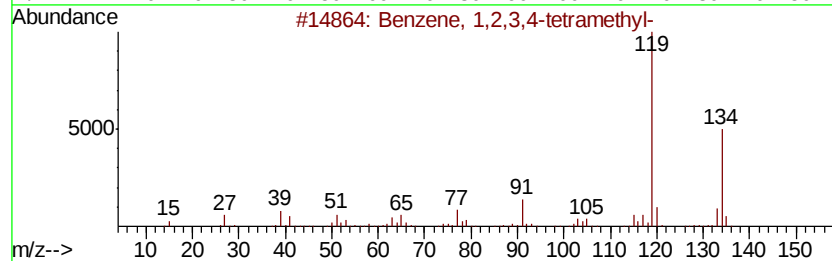
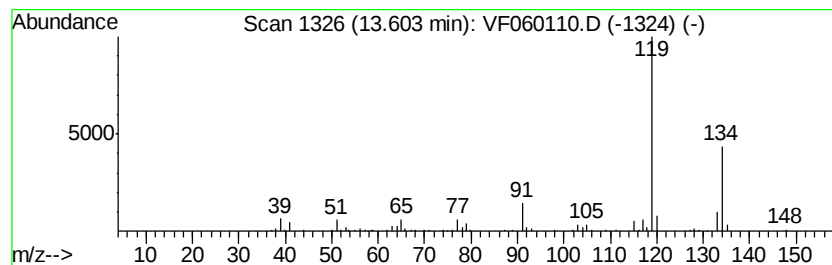
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	9.82 ug/l	81131	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SU-02-082318-B

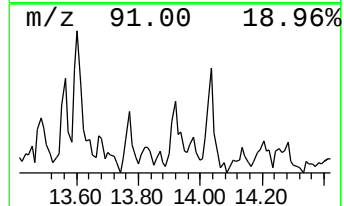
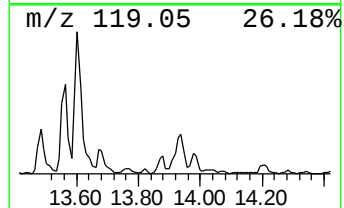
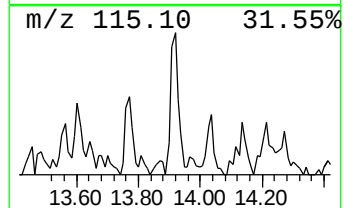
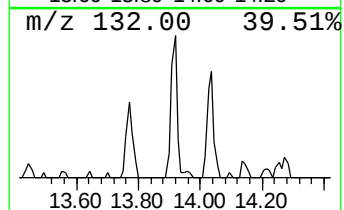
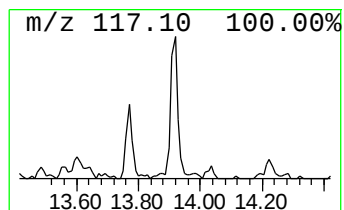
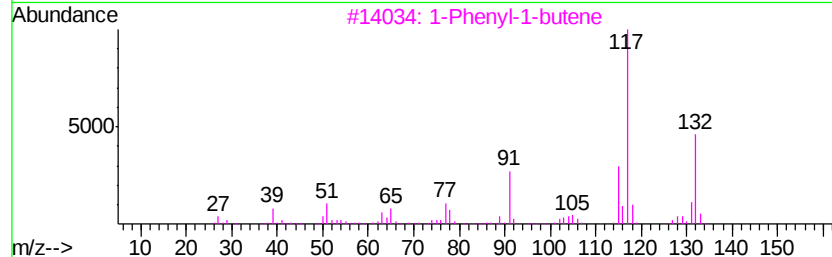
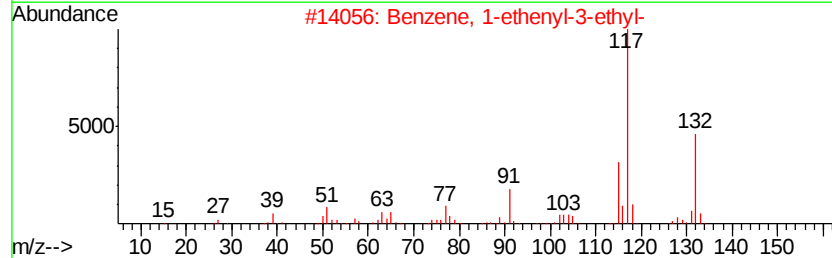
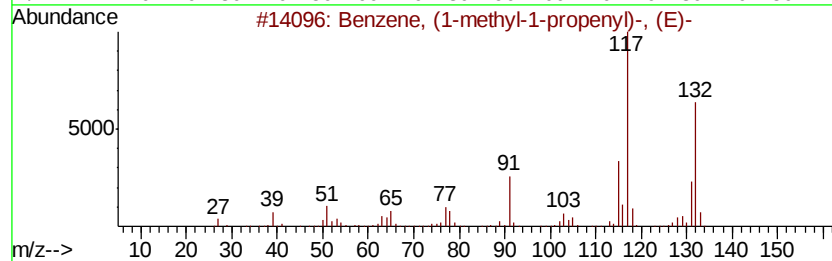
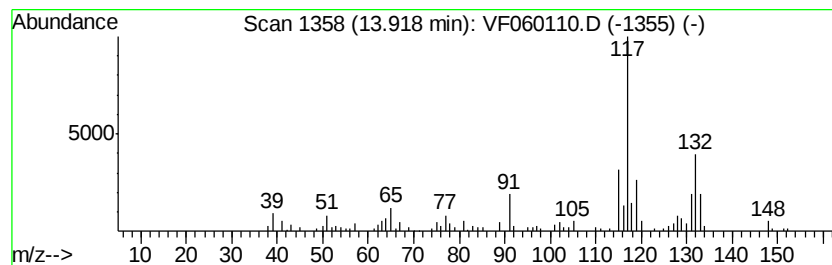
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	9.81 ug/l	81063	1,4-Dichlorobenzene-d4	12.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	80
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
5			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF082418\
Data File : VF060110.D
Acq On : 24 Aug 2018 19:58
Operator : VA/AP
Sample : J4282-13
Misc : 5.36g/5mL/MSVOA_F/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SU-02-082318-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082118S.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
Decane	11.74	9.8	ug/l	81028	4	12.63	413053	50.0
Benzene, 1-ethyl-...	11.80	77.2	ug/l	638111	4	12.63	413053	50.0
Benzene, 1-ethyl-...	12.09	24.0	ug/l	198307	4	12.63	413053	50.0
Indane	12.79	14.2	ug/l	117237	4	12.63	413053	50.0
Benzene, 1-methyl...	12.84	13.1	ug/l	108053	4	12.63	413053	50.0
Benzene, 1,2,3,5-...	12.89	30.6	ug/l	253098	4	12.63	413053	50.0
Benzene, 1-methyl...	13.16	14.5	ug/l	119715	4	12.63	413053	50.0
o-Cymene	13.22	16.2	ug/l	133482	4	12.63	413053	50.0
Benzene, 1,2,3,4-...	13.60	9.8	ug/l	81131	4	12.63	413053	50.0
Benzene, (1-methy...	13.92	9.8	ug/l	81063	4	12.63	413053	50.0