

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF100218\
 Data File : VF060414.D
 Acq On : 2 Oct 2018 20:46
 Operator : VA/AP
 Sample : J5165-01
 Misc : 6.23/5mL/MSVOA F/SOIL
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 EP1-MW-G-092718(7-10)

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\82F100118S.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	4.271	364	369	378	rVB	218179	670320	2.19%	0.124%
2	5.010	436	444	455	rBV2	405190	1351149	4.41%	0.249%
3	5.740	511	518	527	rBV	417794	1142417	3.73%	0.210%
4	7.691	706	716	732	rBV2	670257	2505708	8.18%	0.462%
5	7.898	732	737	747	rVB3	97367	427539	1.40%	0.079%
6	8.263	768	774	777	rVV2	159235	427337	1.40%	0.079%
7	8.411	783	789	797	rBV	749398	2296662	7.50%	0.423%
8	8.579	798	806	819	rVV2	522533	2366483	7.73%	0.436%
9	8.766	819	825	841	rVV	1021008	5443254	17.78%	1.003%
10	8.973	841	846	853	rVB2	281243	916394	2.99%	0.169%
11	9.101	853	859	869	rBV2	1886006	7669998	25.05%	1.413%
12	9.229	869	872	879	rVB	1059449	2579911	8.42%	0.475%
13	9.416	879	891	896	rBV	1566922	4698060	15.34%	0.866%
14	9.545	901	904	908	rVB3	185504	391536	1.28%	0.072%
15	9.673	910	917	924	rBV4	695899	2876941	9.39%	0.530%
16	9.821	924	932	936	rBV3	794261	3121108	10.19%	0.575%
17	9.899	936	940	943	rBV2	2651840	6960279	22.73%	1.282%
18	9.959	943	946	957	rVB	2508419	7389762	24.13%	1.362%
19	10.116	957	962	966	rBV2	577565	1665920	5.44%	0.307%
20	10.205	966	971	980	rVB	778209	2943628	9.61%	0.542%
21	10.373	980	988	995	rBV3	3726916	17997962	58.77%	3.316%
22	10.471	995	998	1004	rVB2	2218349	5415461	17.68%	0.998%
23	10.639	1004	1015	1022	rBV6	3538848	17764020	58.01%	3.273%
24	10.787	1022	1030	1038	rBV3	9489170	29652900	96.83%	5.463%
25	10.895	1038	1041	1043	rBV2	1931194	4358302	14.23%	0.803%
26	10.944	1043	1046	1049	rVB	3594010	6664114	21.76%	1.228%
27	11.063	1054	1058	1062	rVB3	2352741	5275131	17.23%	0.972%
28	11.161	1062	1068	1071	rBV3	4635038	14624776	47.76%	2.695%
29	11.230	1071	1075	1080	rVB	6095514	11793447	38.51%	2.173%
30	11.329	1080	1085	1087	rBV3	2881909	7941631	25.93%	1.463%
31	11.378	1087	1090	1094	rVB2	6849030	15768950	51.49%	2.905%
32	11.437	1094	1096	1098	rVB3	280918	375083	1.22%	0.069%
33	11.496	1098	1102	1103	rBV2	2255327	4366831	14.26%	0.805%
34	11.546	1103	1107	1109	rBV3	1829232	3996851	13.05%	0.736%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	11.595	1109	1112	1115	rVB4	3114482	5678394	18.54%	1.046%
36	11.654	1115	1118	1121	rBV	6915389	14795048	48.31%	2.726%
37	11.772	1127	1130	1133	rBV4	1990270	3823276	12.49%	0.704%
38	11.831	1133	1136	1137	rVV2	2249083	3825513	12.49%	0.705%
39	11.861	1137	1139	1142	rVB2	3121777	5201216	16.98%	0.958%
40	11.910	1142	1144	1147	rVB	2661649	4611706	15.06%	0.850%
41	11.979	1147	1151	1153	rBV2	3264899	6847444	22.36%	1.262%
42	12.048	1153	1158	1159	rBV3	3821906	8956352	29.25%	1.650%
43	12.098	1159	1163	1167	rVB2	13528862	27125070	88.58%	4.998%
44	12.167	1167	1170	1174	rBV2	3636058	7440151	24.30%	1.371%
45	12.245	1174	1178	1179	rVV3	3957734	8106940	26.47%	1.494%
46	12.295	1179	1183	1187	rVV	16030729	30622655	100.00%	5.642%
47	12.383	1188	1192	1197	rVV3	6231257	14699422	48.00%	2.708%
48	12.462	1197	1200	1204	rVV2	3094397	6437906	21.02%	1.186%
49	12.571	1208	1211	1219	rBV5	2687932	8073060	26.36%	1.487%
50	12.689	1219	1223	1227	rBV	5602208	10345564	33.78%	1.906%
51	12.817	1232	1236	1238	rVV2	5981700	13657898	44.60%	2.516%
52	12.866	1238	1241	1244	rVV3	5400396	11904870	38.88%	2.193%
53	12.916	1244	1246	1250	rVV2	5889163	11420252	37.29%	2.104%
54	13.004	1250	1255	1260	rVV3	4560284	13824870	45.15%	2.547%
55	13.083	1260	1263	1266	rVV2	2734928	5712020	18.65%	1.052%
56	13.152	1266	1270	1275	rVV3	6016417	19837903	64.78%	3.655%
57	13.221	1275	1277	1283	rVV3	6240775	17295637	56.48%	3.187%
58	13.300	1283	1285	1289	rVV2	3925842	8595558	28.07%	1.584%
59	13.359	1289	1291	1295	rVV2	3746536	7789211	25.44%	1.435%
60	13.448	1298	1300	1303	rVV	3434970	4711040	15.38%	0.868%
61	13.517	1304	1307	1309	rVV4	1352051	2637468	8.61%	0.486%
62	13.566	1309	1312	1314	rVV	4593642	6984579	22.81%	1.287%
63	13.606	1314	1316	1323	rVB2	8085806	15665163	51.16%	2.886%
64	13.704	1323	1326	1330	rBV	2506774	3709200	12.11%	0.683%
65	13.763	1330	1332	1335	rBV	2573219	4091894	13.36%	0.754%
66	13.813	1335	1337	1340	rVB	1165421	1909173	6.23%	0.352%
67	13.872	1340	1343	1345	rBV	2247706	3120263	10.19%	0.575%
68	13.911	1345	1347	1352	rVV3	3833948	6949030	22.69%	1.280%
69	13.980	1352	1354	1356	rVB	1578990	1944846	6.35%	0.358%
70	14.020	1356	1358	1361	rVB3	418630	672929	2.20%	0.124%
71	14.069	1361	1363	1365	rBV	320858	418671	1.37%	0.077%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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72	14.108	1365	1367	1369	rBV2	581720	1069413	3.49%	0.197%
73	14.138	1369	1370	1373	rVB	805777	792302	2.59%	0.146%
74	14.207	1373	1377	1380	rBV3	1760101	4009769	13.09%	0.739%
75	14.276	1382	1384	1390	rVB	1245894	1793254	5.86%	0.330%
76	14.355	1390	1392	1395	rVB2	435776	557627	1.82%	0.103%
77	14.414	1396	1398	1401	rVB	237223	387053	1.26%	0.071%
78	14.463	1401	1403	1407	rBV3	176882	448966	1.47%	0.083%
79	14.798	1434	1437	1439	rVB3	282329	410784	1.34%	0.076%

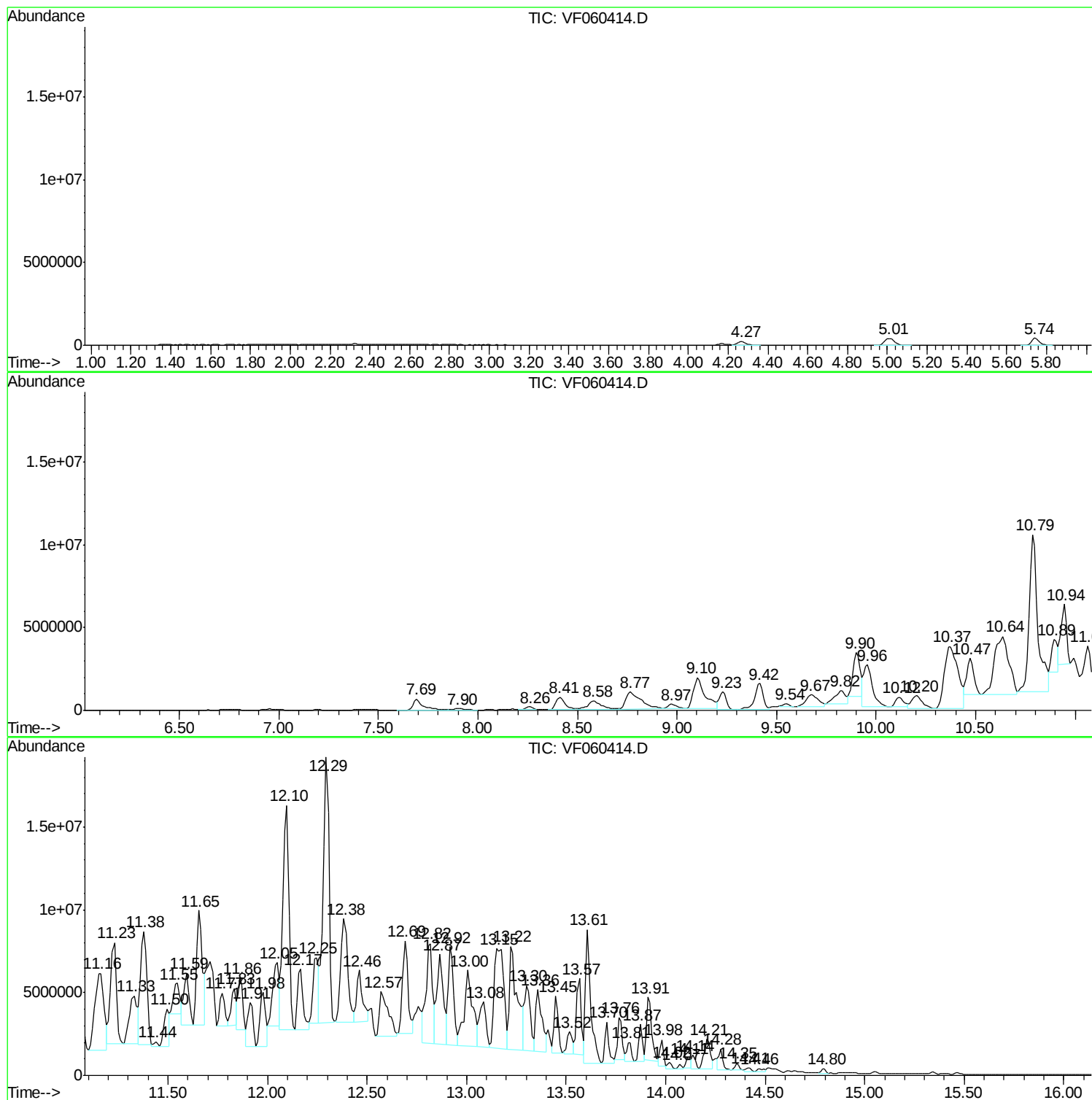
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F100118S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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ALS Vial : 49 Sample Multiplier: 1

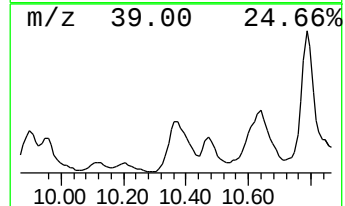
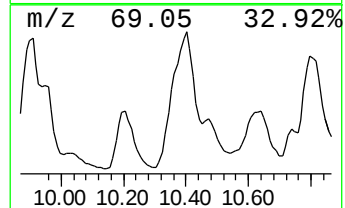
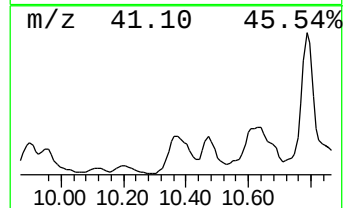
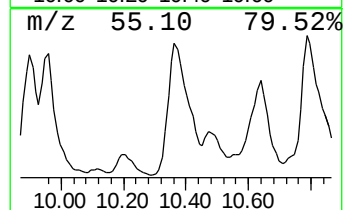
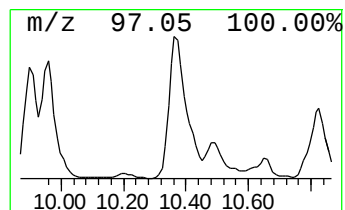
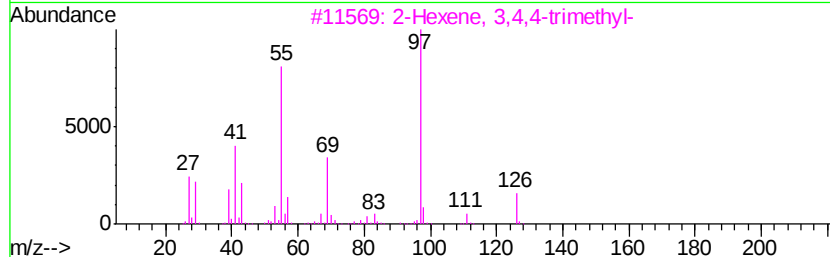
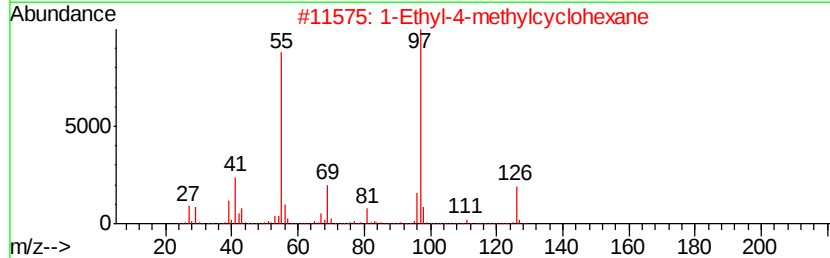
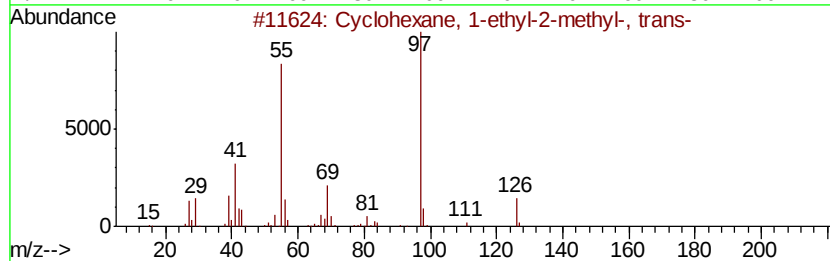
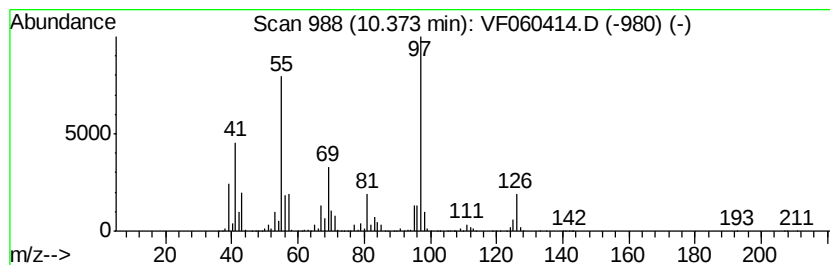
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F100118S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.37	129.29 ug/l	17998000	Chlorobenzene-d5	9.90		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	81
2		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	72
3		2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	72
4		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	64
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	64



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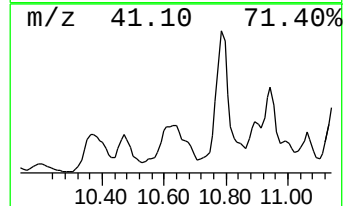
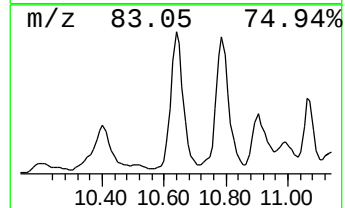
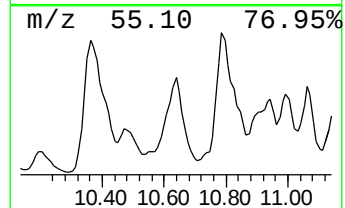
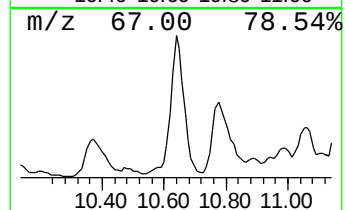
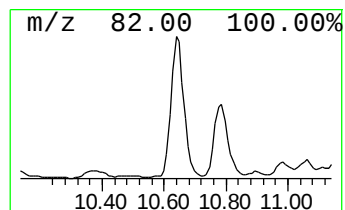
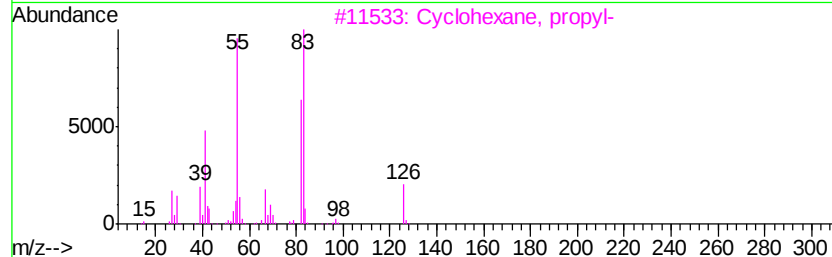
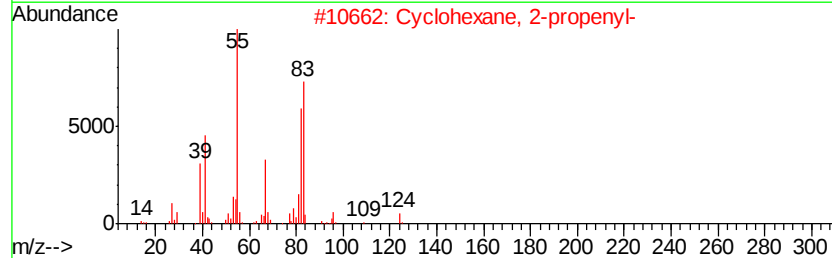
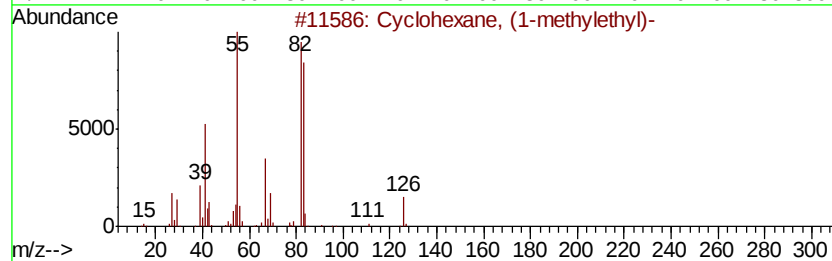
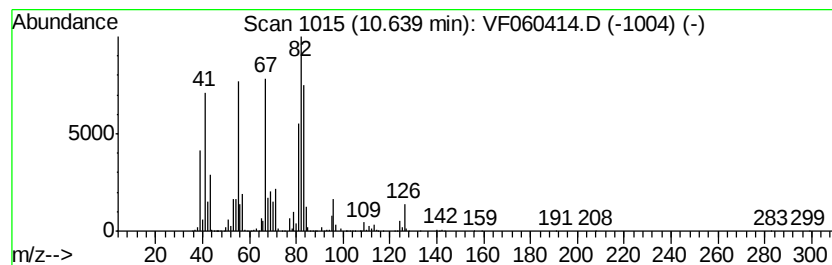
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, (1-methylethyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.64	127.61 ug/l	17764000	Chlorobenzene-d5	9.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	49
2		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	38
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	38
4		1,3-Butadiene, 2,3-dimethyl-	82	C6H10	000513-81-5	38
5		2-Hexyne	82	C6H10	000764-35-2	38



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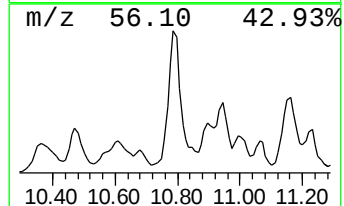
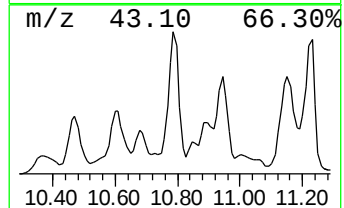
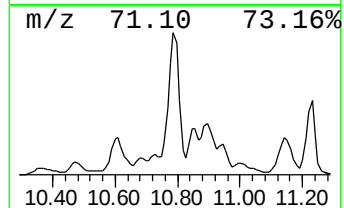
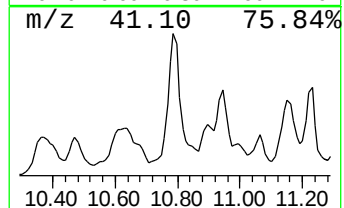
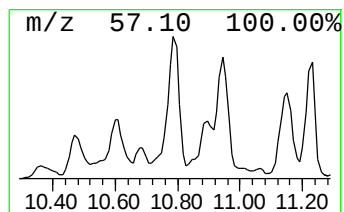
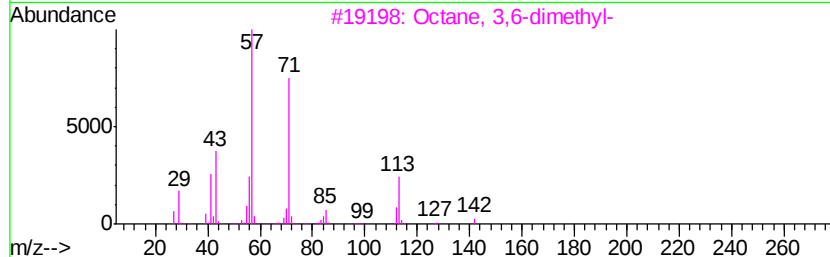
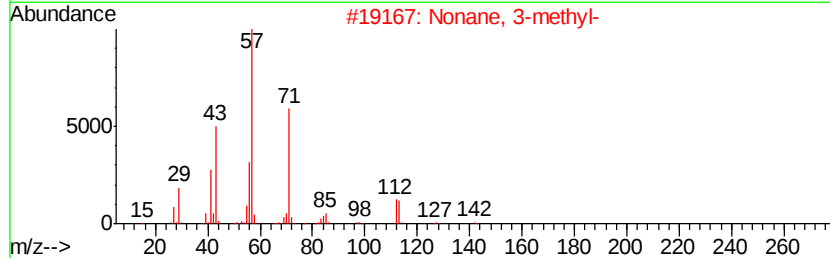
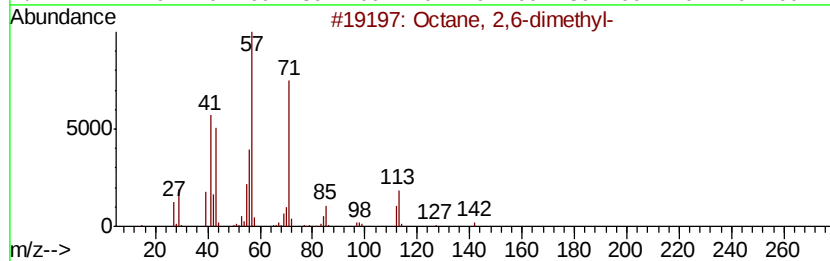
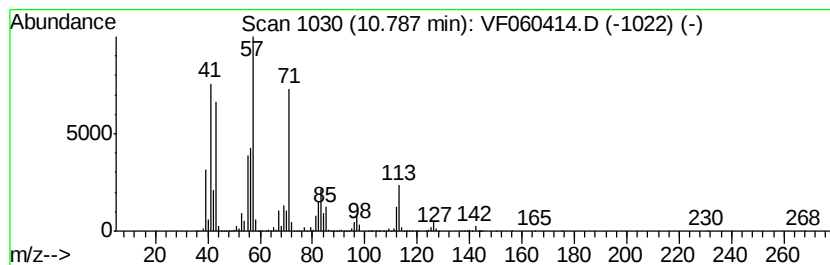
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.79	213.02 ug/l	29652900	Chlorobenzene-d5	9.90		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	95
2	Nonane, 3-methyl-		142	C10H22	005911-04-6	76
3	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	70
4	Sulfurous acid, 2-ethylhexyl oct...		446	C26H54O3S	1000309-20-1	58
5	Octane, 3,5-dimethyl-		142	C10H22	015869-93-9	58



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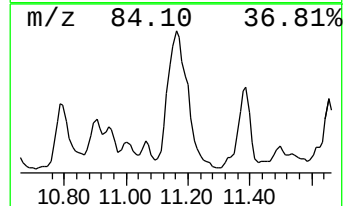
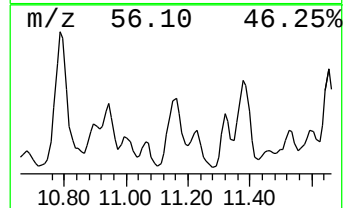
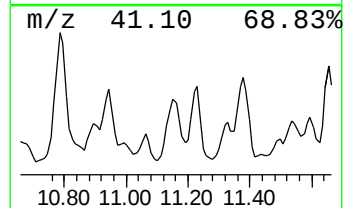
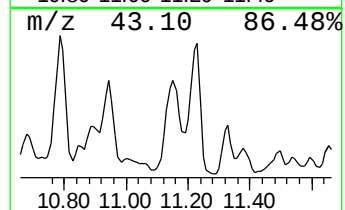
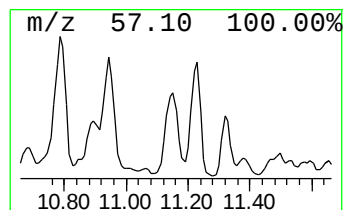
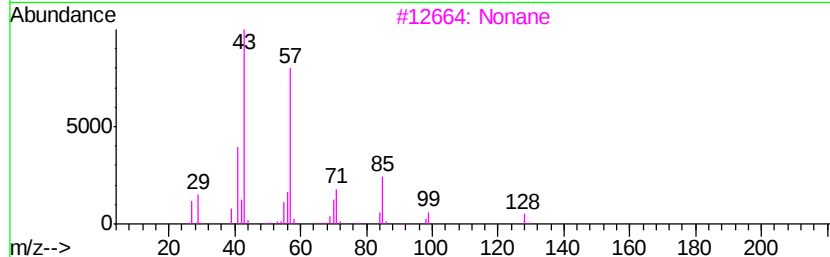
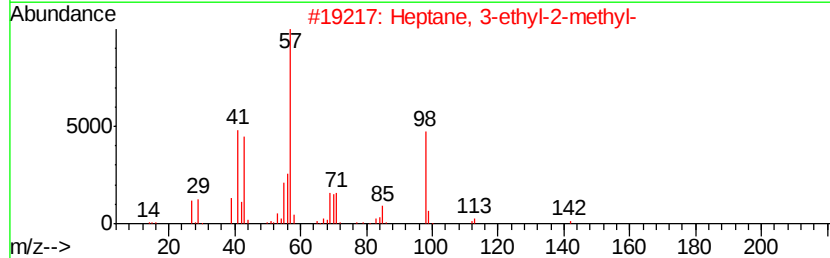
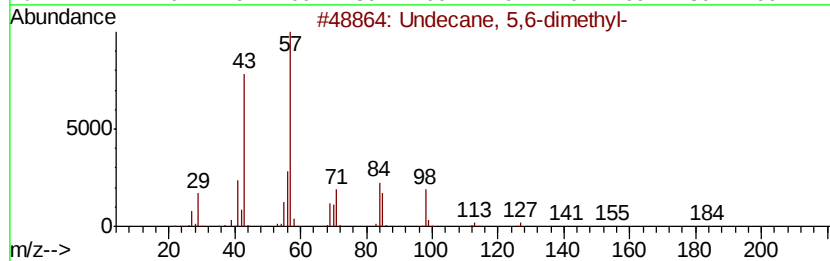
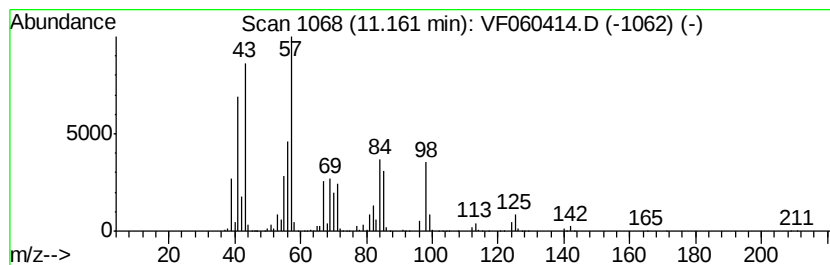
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Quant Title : SW846 8260

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

Peak Number 4 unknown11.16 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.16	105.06 ug/l	14624800	Chlorobenzene-d5	9.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 5,6-dimethyl-	184	C13H28	017615-91-7	47
2		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	43
3		Nonane	128	C9H20	000111-84-2	38
4		1-Octene, 3,4-dimethyl-	140	C10H20	056728-11-1	38
5		Carbonic acid, butyl heptyl ester	216	C12H24O3	1000314-63-4	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA F/SOIL
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EP1-MW-G-092718(7-10)

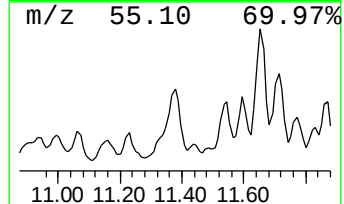
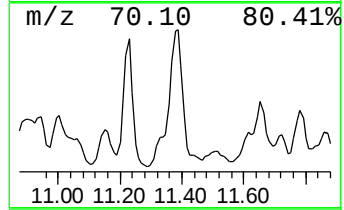
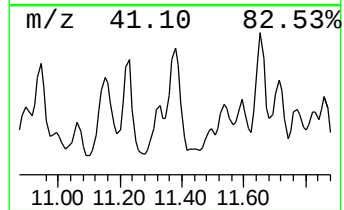
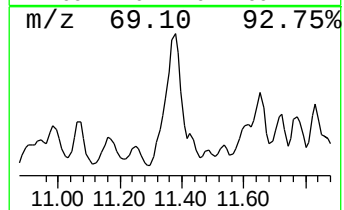
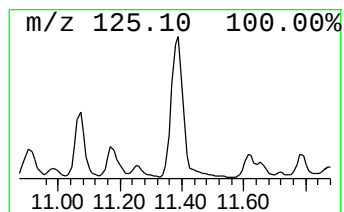
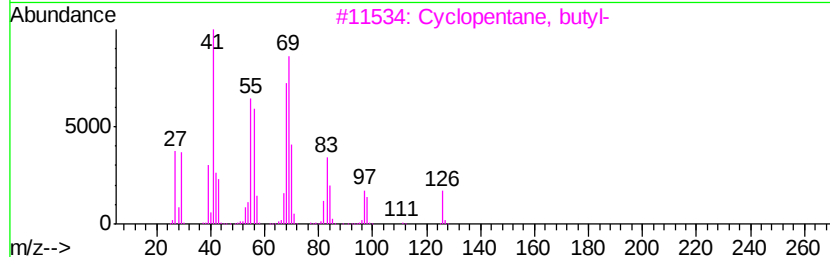
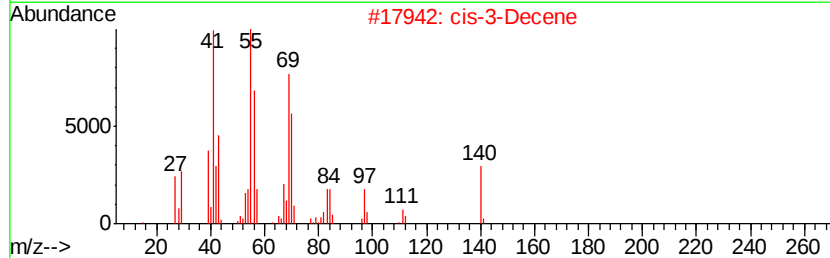
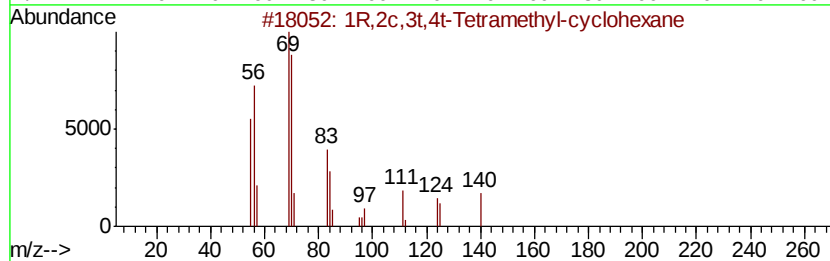
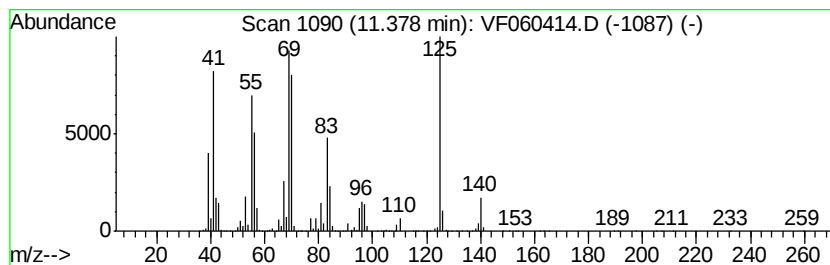
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Quant Title : SW846 8260

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

Peak Number 5 1R,2c,3t,4t-Tetramethyl-cyc... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.38	76.21 ug/l	15769000	1,4-Dichlorobenzene-d4	12.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1R,2c,3t,4t-Tetramethyl-cyclohexane	140	C10H20	1000144-07-3	58		
2	cis-3-Decene	140	C10H20	019398-86-8	52		
3	Cyclopentane, butyl-	126	C9H18	002040-95-1	45		
4	1-Octene, 2,6-dimethyl-	140	C10H20	006874-29-9	43		
5	2-Dodecenal, (E)-	182	C12H22O	020407-84-5	35		



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA F/SOIL
ALS Vial : 49 Sample Multiplier: 1

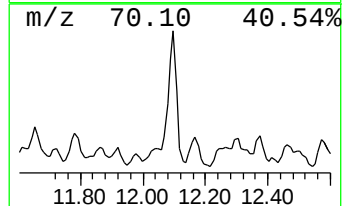
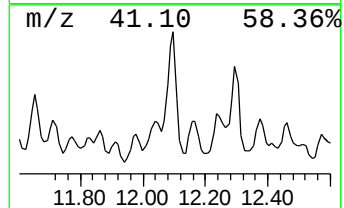
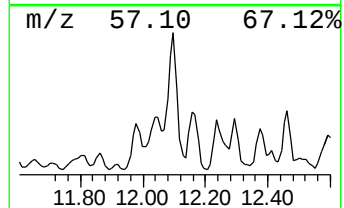
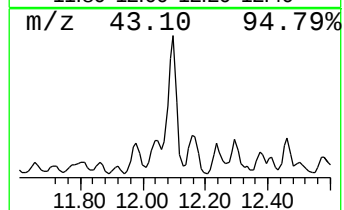
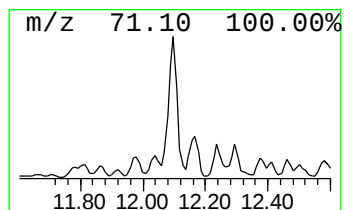
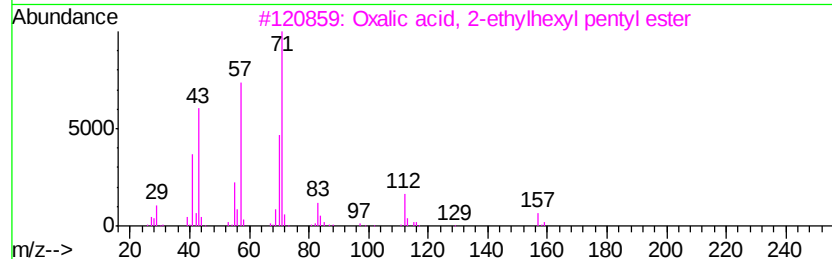
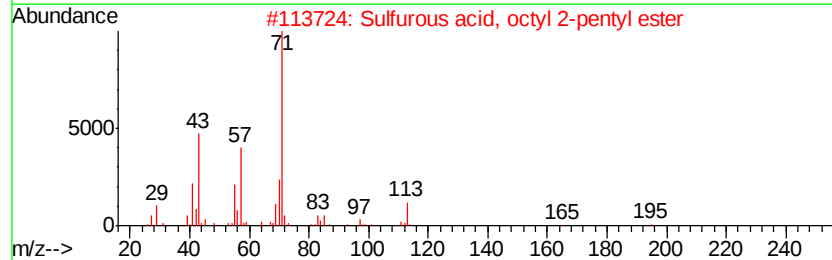
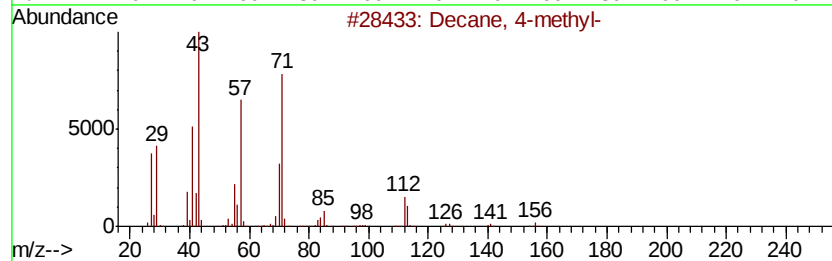
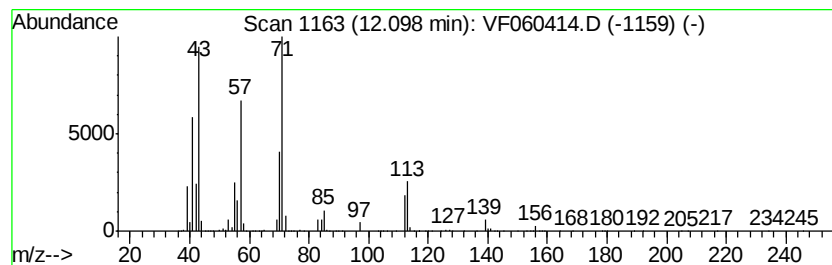
Instrument :
MSVOA_F
ClientSampleId :
EP1-MW-G-092718(7-10)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F100118S.M
Quant Title : SW846 8260

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

Peak Number 6 Decane, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	131.10 ug/l	27125100	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 4-methyl-	156 C11H24	002847-72-5	87
2	Sulfurous acid, octyl 2-pentyl e...	264 C13H28O3S	1000309-15-7	72
3	Oxalic acid, 2-ethylhexyl pentyl...	272 C15H28O4	1000309-38-7	59
4	Heptane, 3,3,5-trimethyl-	142 C10H22	007154-80-5	59
5	Octane, 3,3-dimethyl-	142 C10H22	004110-44-5	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA F/SOIL
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EP1-MW-G-092718(7-10)

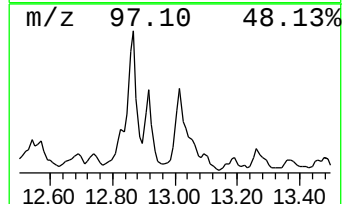
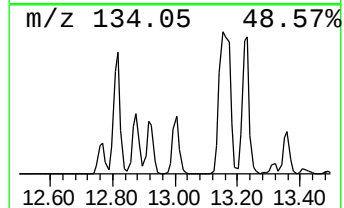
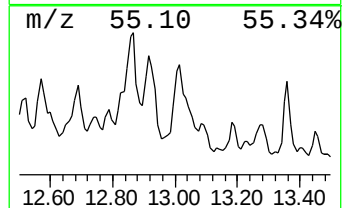
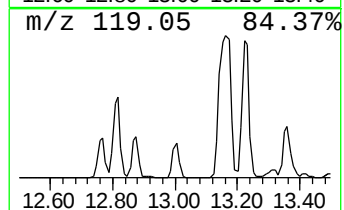
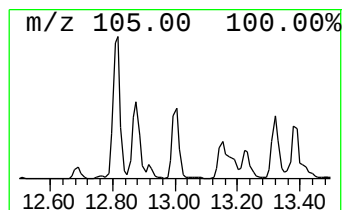
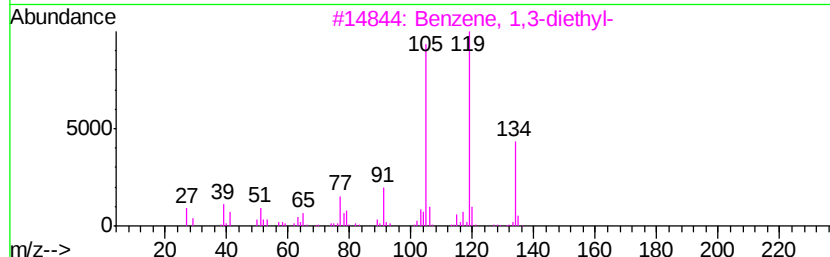
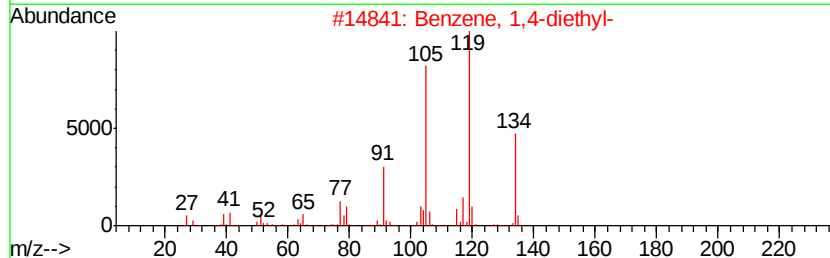
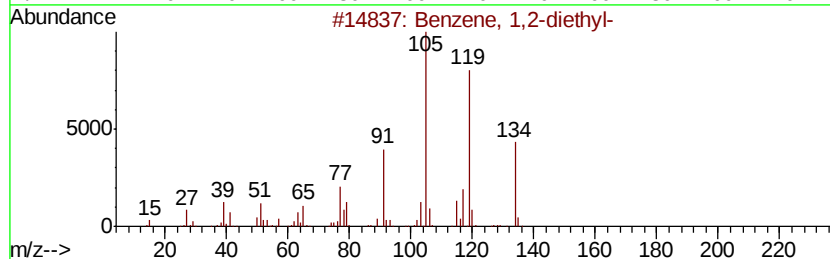
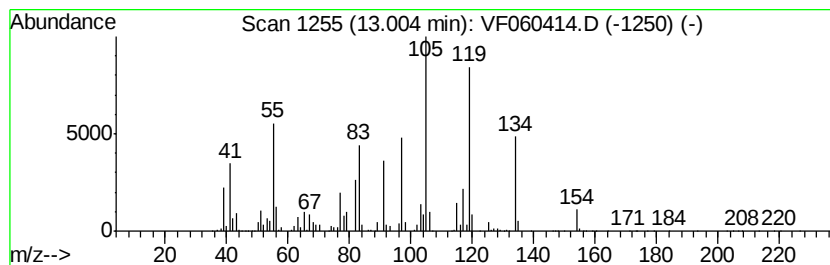
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	66.82 ug/l	13824900	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	87
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	43
5		Ethanone, 1-(4-methylphenyl)-	134	C9H10O	000122-00-9	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA F/SOIL
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EP1-MW-G-092718(7-10)

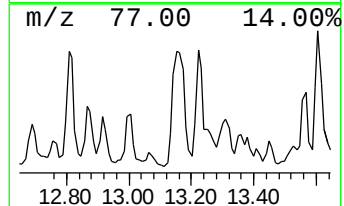
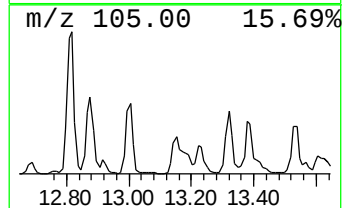
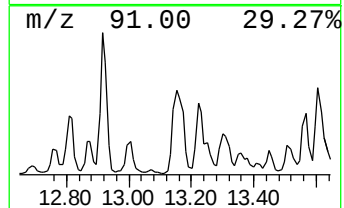
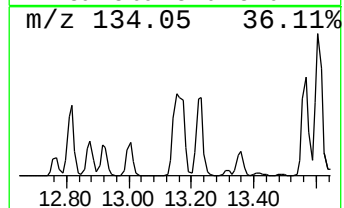
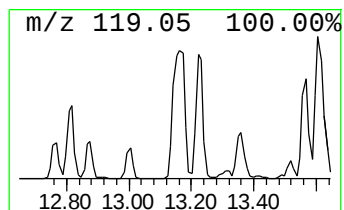
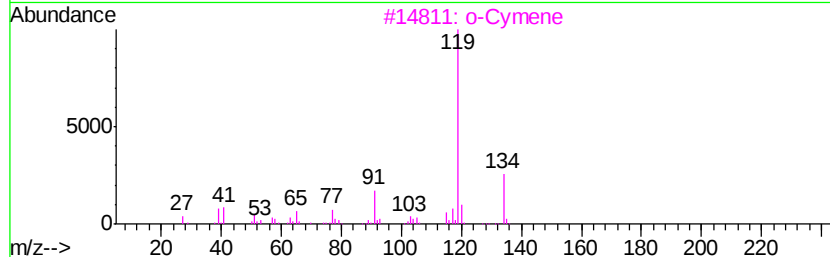
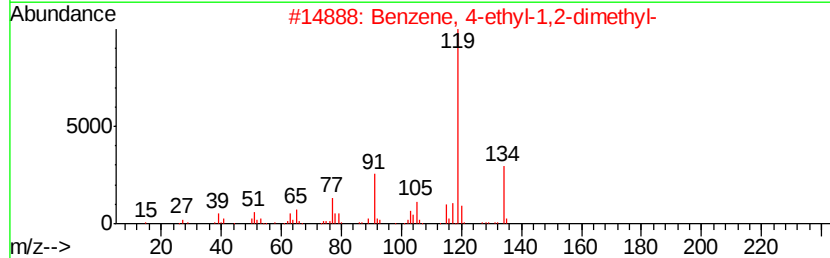
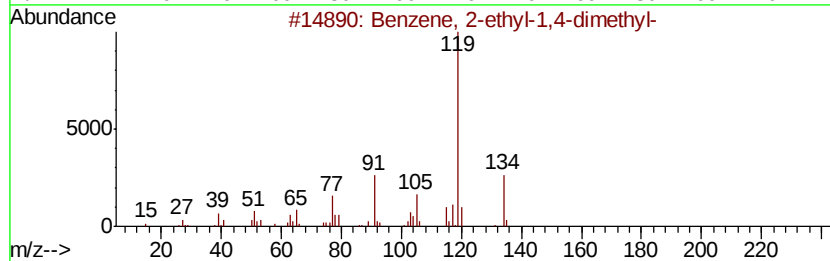
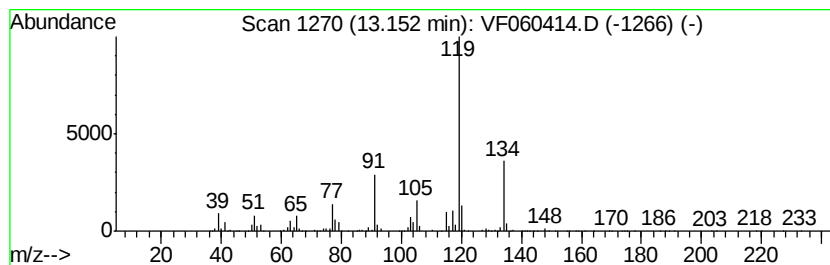
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	95.88 ug/l	19837900	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA F/SOIL
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EP1-MW-G-092718(7-10)

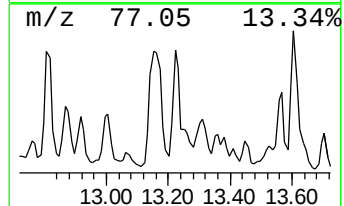
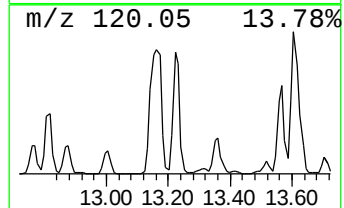
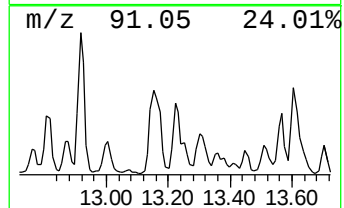
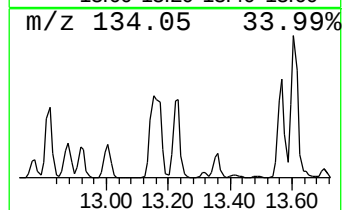
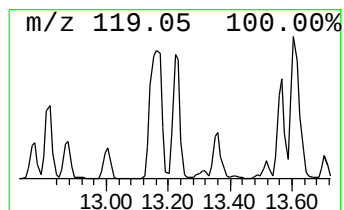
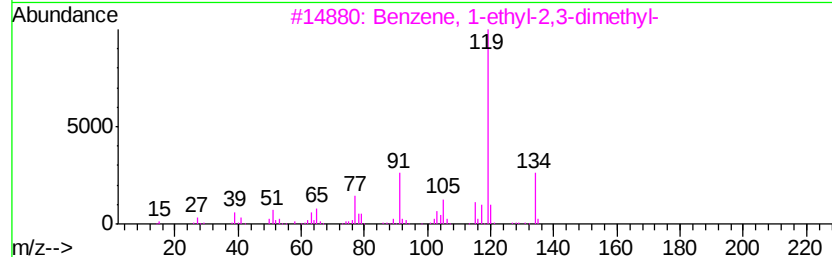
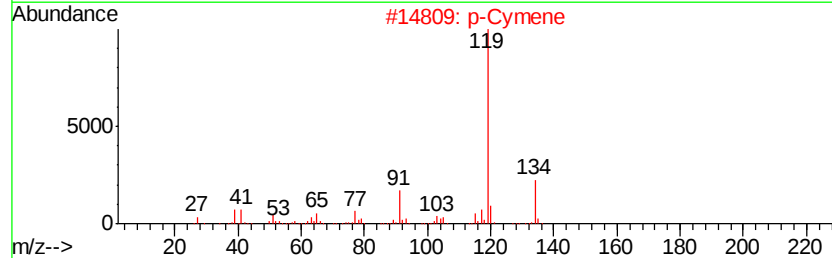
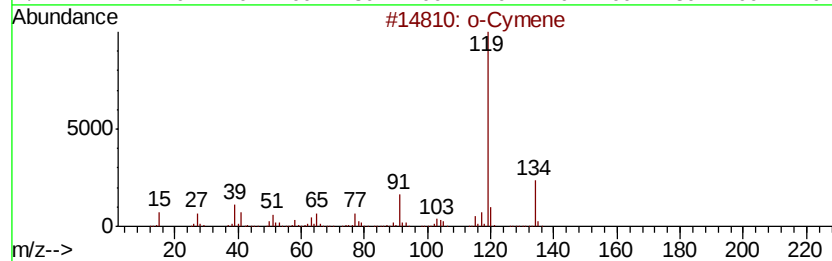
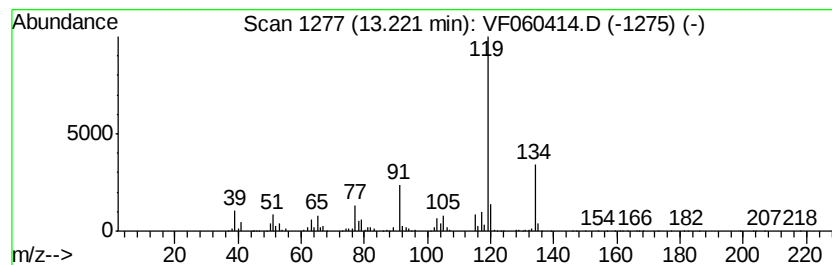
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Quant Title : SW846 8260

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

Peak Number 9 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	83.59 ug/l	17295600	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	97
2		p-Cymene	134	C10H14	000099-87-6	97
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	96
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA F/SOIL
ALS Vial : 49 Sample Multiplier: 1

Instrument :
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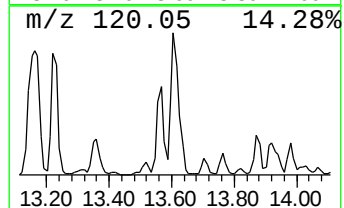
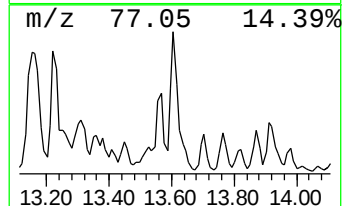
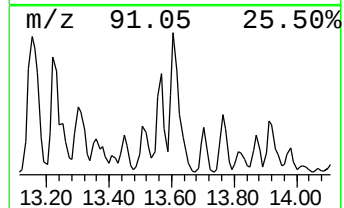
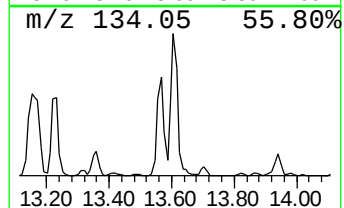
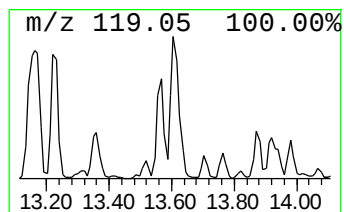
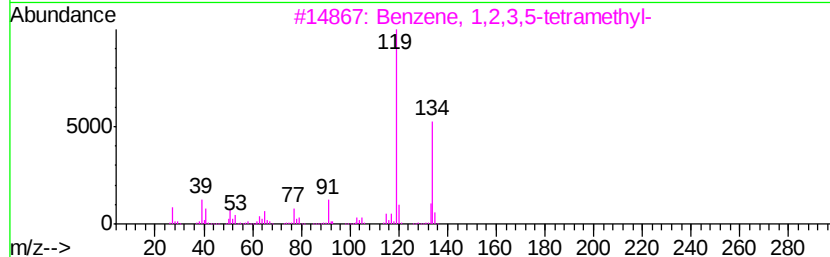
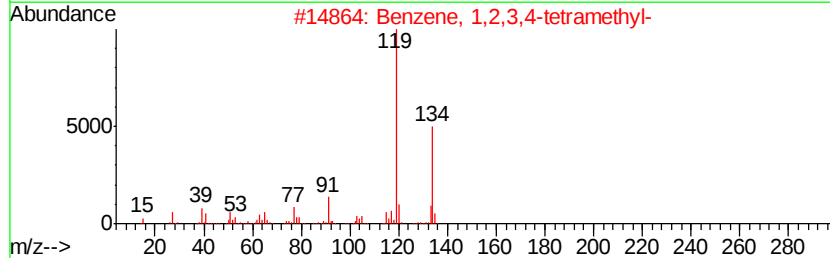
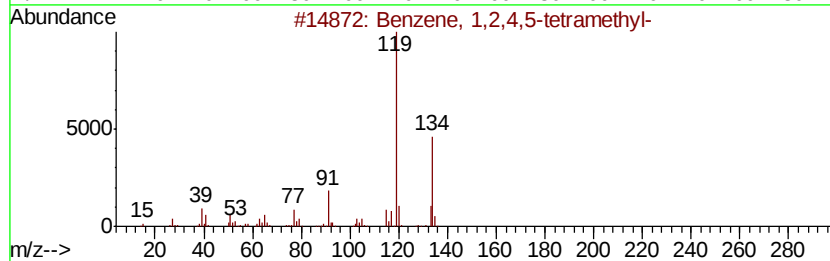
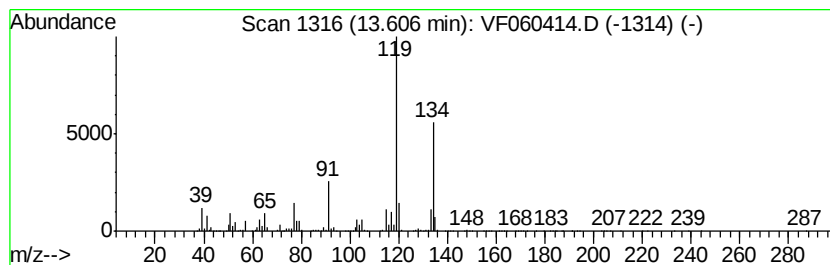
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.61	75.71 ug/l	15665200	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF100218\
Data File : VF060414.D
Acq On : 2 Oct 2018 20:46
Operator : VA/AP
Sample : J5165-01
Misc : 6.23/5mL/MSVOA_F/SOIL
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EP1-MW-G-092718(7-10)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F100118S.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
Cyclohexane, 1-et...	10.37	129.3	ug/l	17998000	3	9.90	6960280	50.0
Cyclohexane, (1-m...	10.64	127.6	ug/l	17764000	3	9.90	6960280	50.0
Octane, 2,6-dimet...	10.79	213.0	ug/l	29652900	3	9.90	6960280	50.0
unknown11.16	11.16	105.1	ug/l	14624800	3	9.90	6960280	50.0
1R,2c,3t,4t-Tetra...	11.38	76.2	ug/l	15769000	4	12.63	10345600	50.0
Decane, 4-methyl-	12.10	131.1	ug/l	27125100	4	12.63	10345600	50.0
Benzene, 1,2-diet...	13.00	66.8	ug/l	13824900	4	12.63	10345600	50.0
Benzene, 2-ethyl-...	13.15	95.9	ug/l	19837900	4	12.63	10345600	50.0
o-Cymene	13.22	83.6	ug/l	17295600	4	12.63	10345600	50.0
Benzene, 1,2,4,5-...	13.61	75.7	ug/l	15665200	4	12.63	10345600	50.0