

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF082918\
 Data File : VF060183.D
 Acq On : 29 Aug 2018 8:44
 Operator : VA/AP
 Sample : J4277-09
 Misc : 6.48g/5mL/MSVOA F/SOIL
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 OR-01-082718-A

Integration Parameters: RTEINT.P

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

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

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.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	1.085	53	56	59	rBV2	7421	9716	4.25%	0.845%
2	1.223	68	70	72	rVV	1394	2536	1.11%	0.221%
3	1.282	74	76	77	rVV	2647	3568	1.56%	0.310%
4	1.321	78	80	91	rVV4	3235	10035	4.39%	0.873%
5	2.080	153	157	161	rBV3	1626	4343	1.90%	0.378%
6	2.278	175	177	181	rVV3	4471	9986	4.37%	0.868%
7	2.475	195	197	203	rVV4	1860	5122	2.24%	0.445%
8	3.224	271	273	277	rVB	2318	2556	1.12%	0.222%
9	4.279	374	380	387	rBV	22779	66150	28.91%	5.753%
10	5.038	448	457	465	rVV2	68238	228774	100.00%	19.896%
11	5.166	468	470	474	rVB3	1273	2926	1.28%	0.254%
12	5.757	524	530	539	rVV	59026	162628	71.09%	14.144%
13	6.792	634	635	637	rBV	3914	2405	1.05%	0.209%
14	7.709	721	728	734	rBV	34018	84423	36.90%	7.342%
15	7.778	734	735	740	rVV2	2376	3036	1.33%	0.264%
16	8.369	788	795	796	rBV2	1164	2816	1.23%	0.245%
17	8.774	831	836	840	rVV	936	3106	1.36%	0.270%
18	9.907	946	951	957	rBV	38336	85505	37.38%	7.436%
19	10.361	995	997	1000	rBV	1279	2827	1.24%	0.246%
20	10.883	1048	1050	1053	rBV2	1246	3085	1.35%	0.268%
21	10.932	1053	1055	1058	rVV2	1942	2924	1.28%	0.254%
22	11.366	1095	1099	1102	rVB2	3169	6162	2.69%	0.536%
23	11.504	1110	1113	1116	rBV	33553	55291	24.17%	4.809%
24	11.770	1138	1140	1142	rVV2	1734	2720	1.19%	0.237%
25	11.800	1142	1143	1146	rVB2	1929	2613	1.14%	0.227%
26	11.869	1146	1150	1152	rBV4	1706	3088	1.35%	0.269%
27	11.967	1157	1160	1162	rBV	1449	2553	1.12%	0.222%
28	12.086	1169	1172	1176	rVB4	2571	6476	2.83%	0.563%
29	12.155	1176	1179	1181	rBV4	1766	3754	1.64%	0.326%
30	12.243	1185	1188	1189	rVV3	2627	4313	1.89%	0.375%
31	12.273	1189	1191	1193	rVB3	1975	2743	1.20%	0.239%
32	12.362	1198	1200	1203	rVB3	2578	4237	1.85%	0.368%
33	12.411	1203	1205	1208	rBV3	1902	2418	1.06%	0.210%
34	12.539	1212	1218	1219	rBV4	1462	4170	1.82%	0.363%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.M
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35	12.559	1219	1220	1222	rBV2	2031	2638	1.15%	0.229%
36	12.628	1224	1227	1230	rVV	34030	49039	21.44%	4.265%
37	12.687	1230	1233	1236	rVV2	6924	12206	5.34%	1.062%
38	12.756	1238	1240	1242	rVV3	1797	3823	1.67%	0.332%
39	12.835	1245	1248	1251	rBV5	2363	5619	2.46%	0.489%
40	12.923	1252	1257	1259	rVV4	4820	11838	5.17%	1.030%
41	12.963	1259	1261	1264	rVB4	2893	5154	2.25%	0.448%
42	13.012	1264	1266	1267	rBV2	3299	4184	1.83%	0.364%
43	13.061	1269	1271	1273	rVB3	3392	4252	1.86%	0.370%
44	13.180	1278	1283	1285	rBV4	4758	11273	4.93%	0.980%
45	13.219	1285	1287	1288	rVV2	3614	3719	1.63%	0.323%
46	13.268	1288	1292	1297	rVV5	9448	23143	10.12%	2.013%
47	13.406	1301	1306	1308	rBV3	8732	20237	8.85%	1.760%
48	13.446	1308	1310	1314	rVB3	10565	17360	7.59%	1.510%
49	13.594	1321	1325	1328	rBV6	2069	5629	2.46%	0.490%
50	13.732	1334	1339	1341	rBV5	5098	11622	5.08%	1.011%
51	13.771	1341	1343	1347	rVV4	6235	13137	5.74%	1.143%
52	13.860	1351	1352	1355	rVV3	5066	7264	3.18%	0.632%
53	13.929	1355	1359	1361	rVV4	7895	15962	6.98%	1.388%
54	13.968	1361	1363	1366	rVV4	4192	5871	2.57%	0.511%
55	14.008	1366	1367	1371	rVB4	4273	6336	2.77%	0.551%
56	14.087	1371	1375	1376	rBV4	2045	4533	1.98%	0.394%
57	14.165	1380	1383	1388	rVB5	2038	5178	2.26%	0.450%
58	14.294	1390	1396	1397	rVV5	3557	11163	4.88%	0.971%
59	14.363	1401	1403	1404	rVV2	3807	4333	1.89%	0.377%
60	14.412	1407	1408	1413	rVB4	2118	3404	1.49%	0.296%
61	14.491	1414	1416	1418	rVV3	3074	4761	2.08%	0.414%
62	14.560	1421	1423	1425	rVB3	2816	4334	1.89%	0.377%
63	14.599	1425	1427	1429	rVB3	2358	3651	1.60%	0.318%
64	14.648	1430	1432	1434	rBV3	2743	3174	1.39%	0.276%
65	14.688	1434	1436	1440	rVB4	3427	6422	2.81%	0.559%
66	14.747	1440	1442	1443	rBV2	2784	3233	1.41%	0.281%
67	14.925	1459	1460	1463	rBV3	3959	4358	1.90%	0.379%
68	15.141	1480	1482	1483	rBV	5722	4664	2.04%	0.406%
69	15.181	1483	1486	1488	rVV2	10972	12501	5.46%	1.087%
70	15.220	1488	1490	1494	rVB3	3341	7423	3.24%	0.646%
71	15.270	1494	1495	1498	rBV2	2094	3558	1.56%	0.309%

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72	15.309	1498	1499	1501	rBV2	2152	2399	1.05%	0.209%
73	15.486	1516	1517	1520	rBV3	2077	2820	1.23%	0.245%
74	15.615	1528	1530	1534	rVB5	1756	3034	1.33%	0.264%
75	15.684	1534	1537	1540	rBV4	1497	3626	1.58%	0.315%
76	15.723	1540	1541	1544	rVV2	2329	3309	1.45%	0.288%
77	15.792	1547	1548	1552	rBV4	1891	3253	1.42%	0.283%
78	15.891	1555	1558	1562	rVV3	2472	5390	2.36%	0.469%

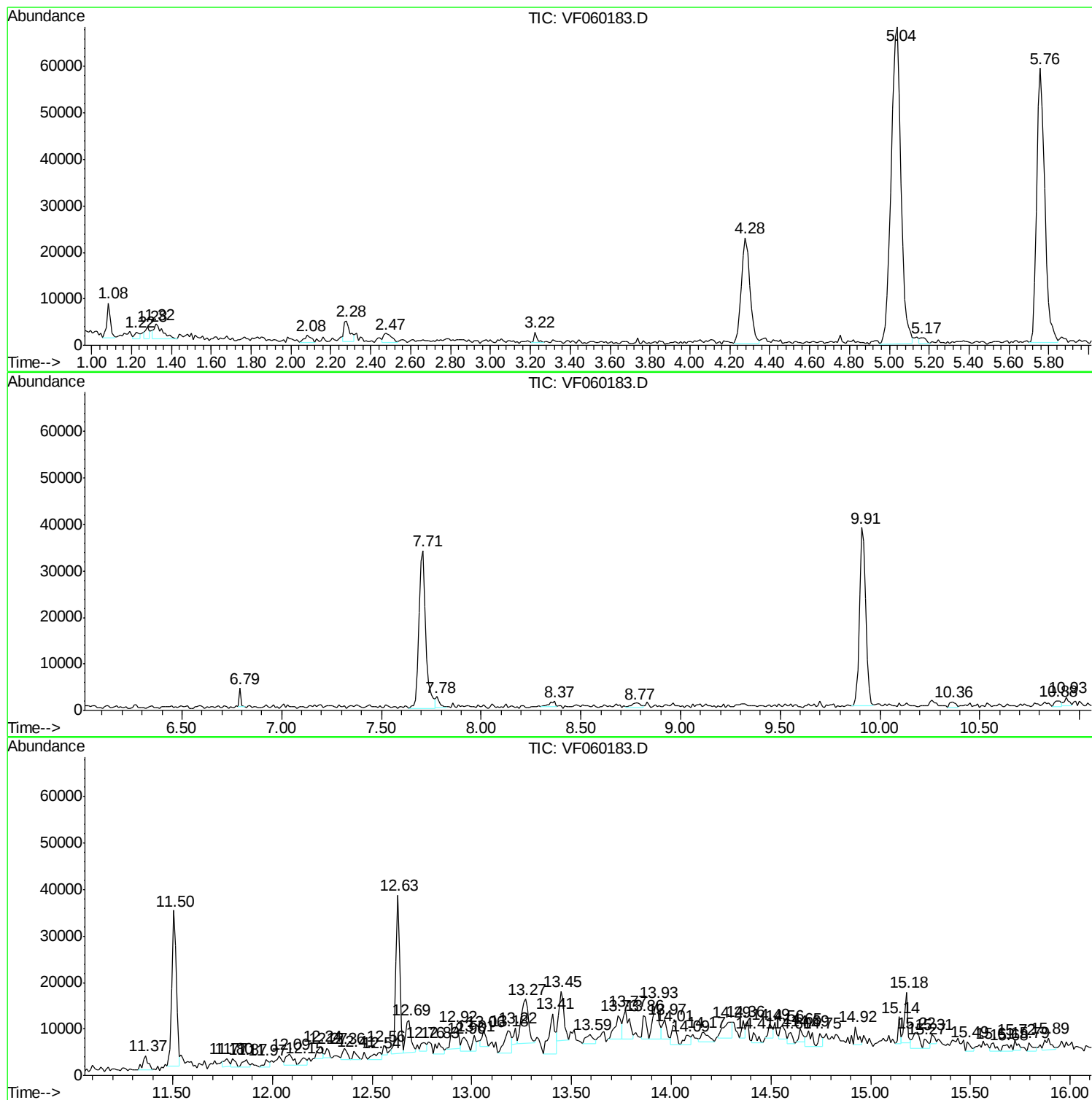
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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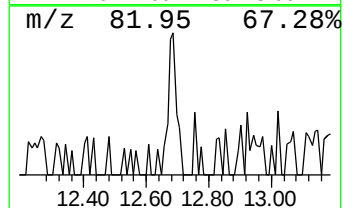
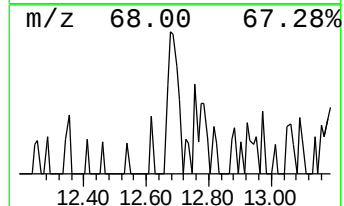
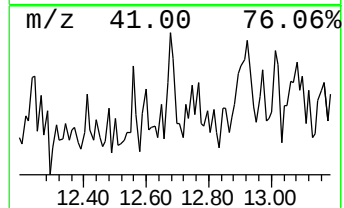
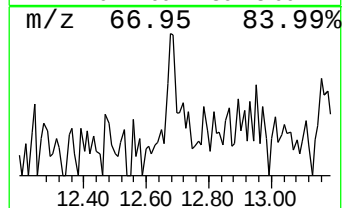
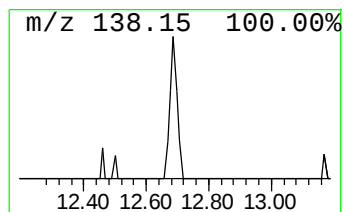
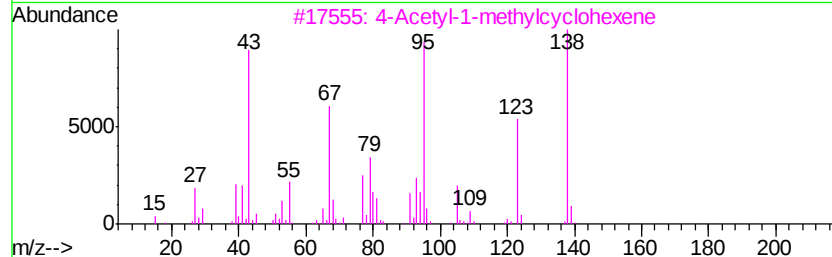
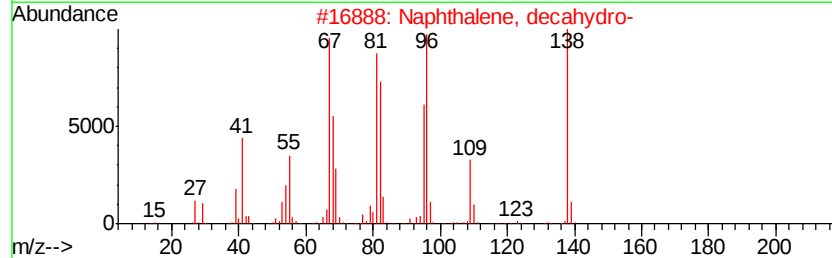
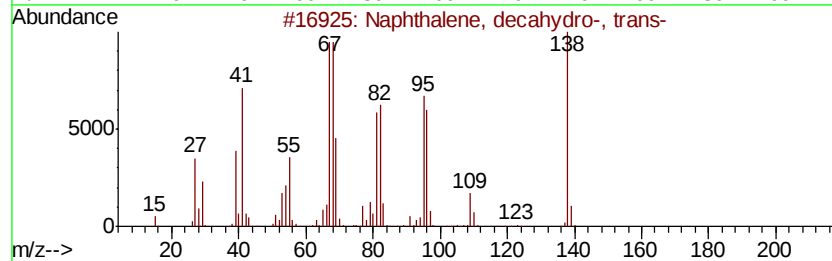
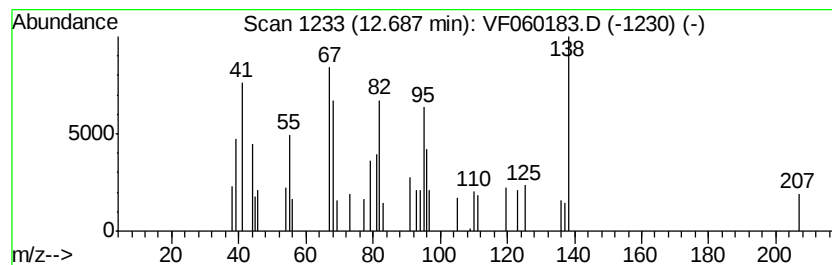
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.M
Quant Title : SW846 8260

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

Peak Number 1 unknown12.69 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	12.45 ug/l	12206	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-, trans-	138 C10H18	000493-02-7 46
2		Naphthalene, decahydro-	138 C10H18	000091-17-8 35
3		4-Acetyl-1-methylcyclohexene	138 C9H14O	006090-09-1 35
4		2(1H)-Pyridinone	95 C5H5NO	000142-08-5 35
5		Bicyclo[4.1.0]heptane, 3,7,7-tri...	138 C10H18	002778-68-9 35



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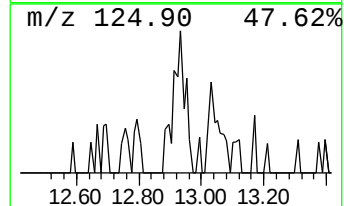
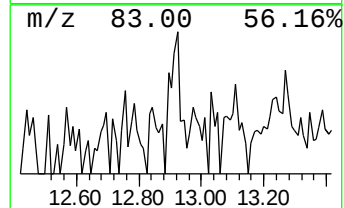
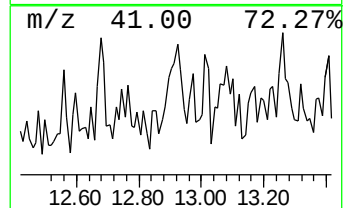
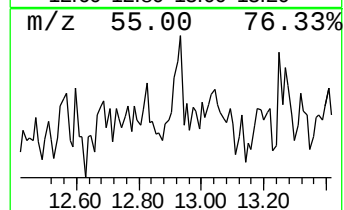
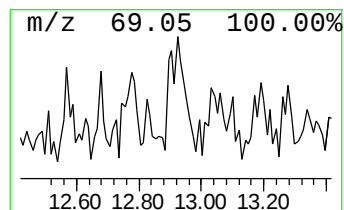
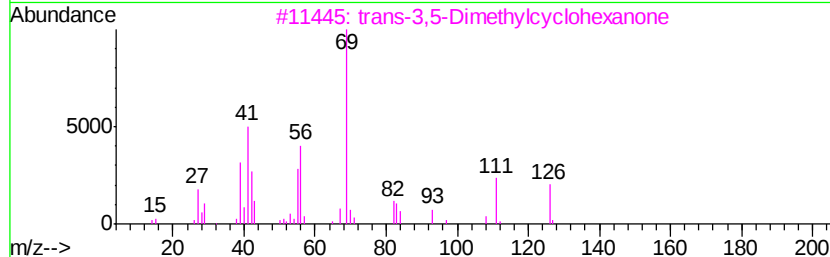
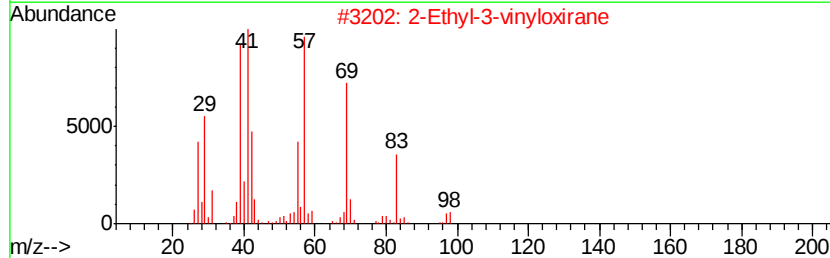
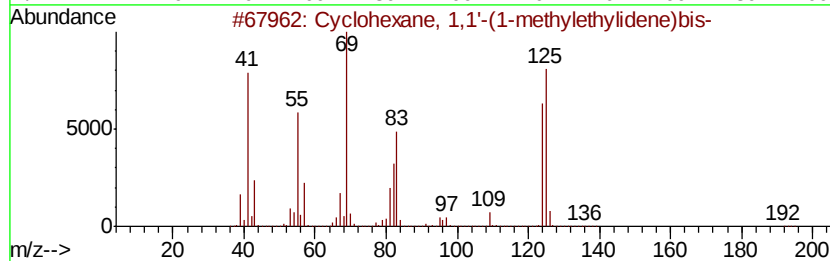
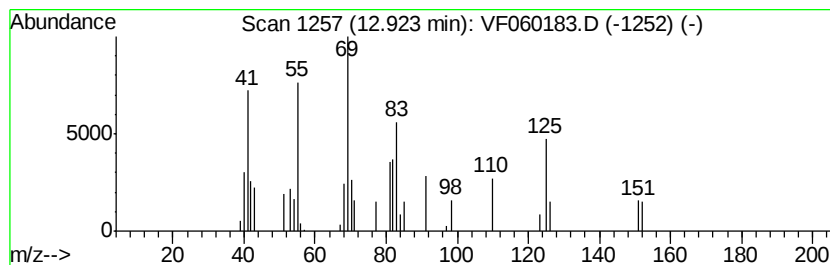
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown12.92 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	12.07 ug/l	11838	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1'-(1-methylethylidene)bis-	208	C15H28	054934-90-6	38
2		2-Ethyl-3-vinylloxirane	98	C6H10O	034485-78-4	32
3		trans-3,5-Dimethylcyclohexanone	126	C8H14O	007214-49-5	27
4		3,4-Dimethyl cyclohexanone	126	C8H14O	005465-09-8	27
5		Cyclohexanone, 3,5-dimethyl-, cis-	126	C8H14O	007214-52-0	27



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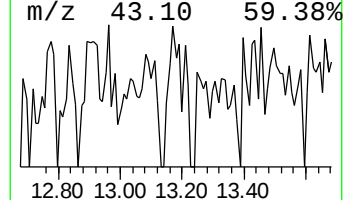
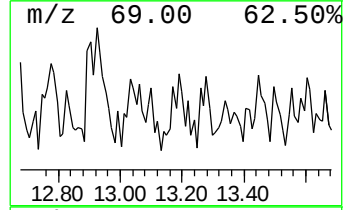
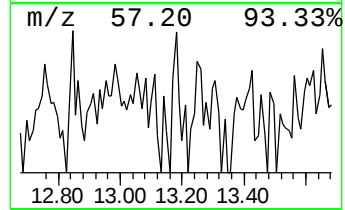
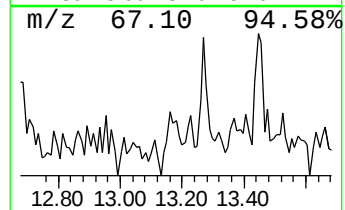
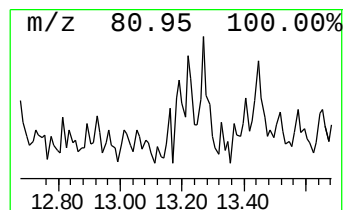
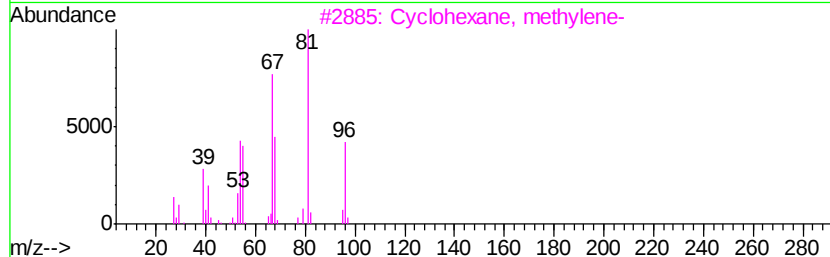
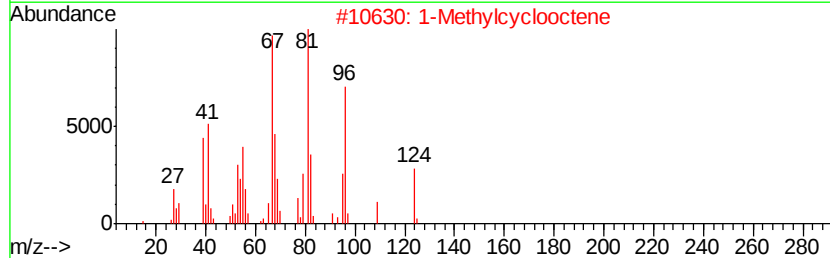
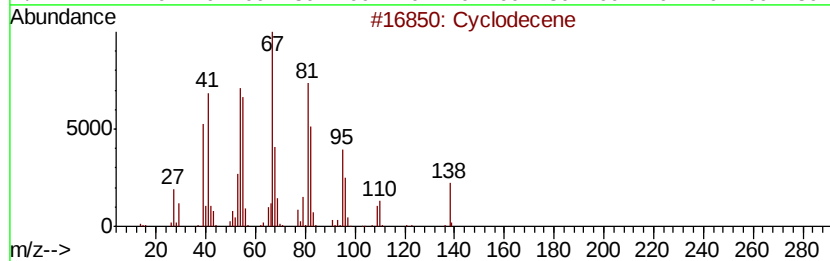
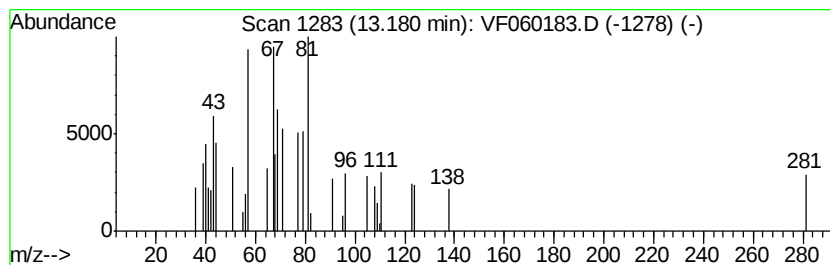
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Quant Title : SW846 8260

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

Peak Number 3 unknown13.18 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	11.49 ug/l	11273	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclodecene	138	C10H18	003618-12-0	47
2		1-Methylcyclooctene	124	C9H16	000933-11-9	46
3		Cyclohexane, methylene-	96	C7H12	001192-37-6	43
4		1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	40
5		9-Oxabicyclo[3.3.1]non-6-en-2-on...	153	C8H11NO2	063827-16-7	40



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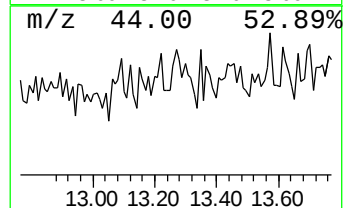
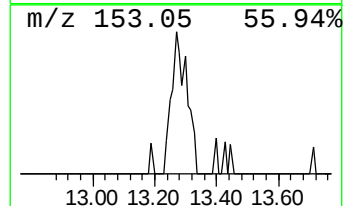
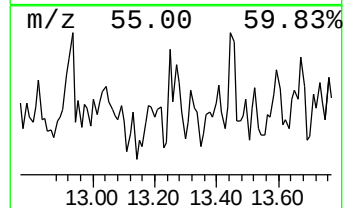
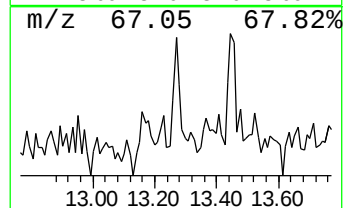
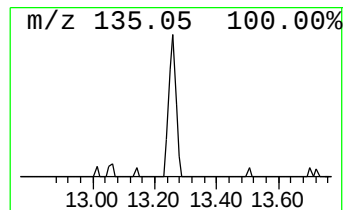
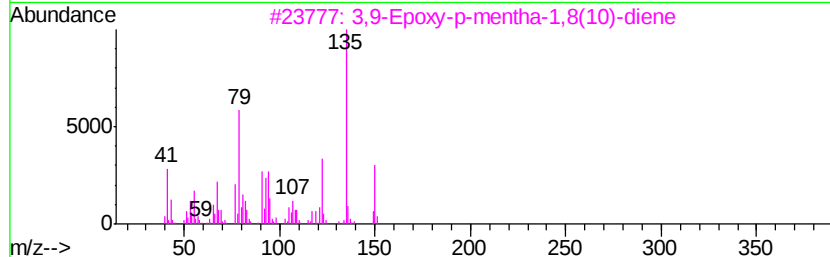
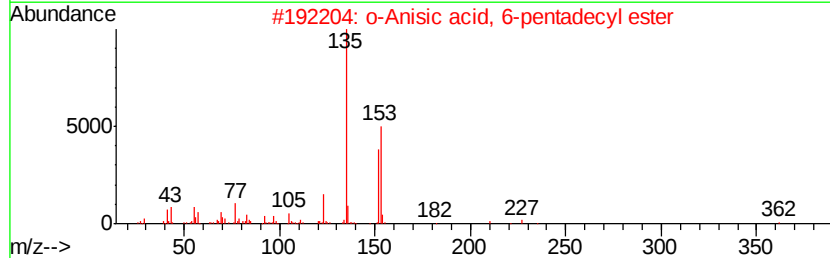
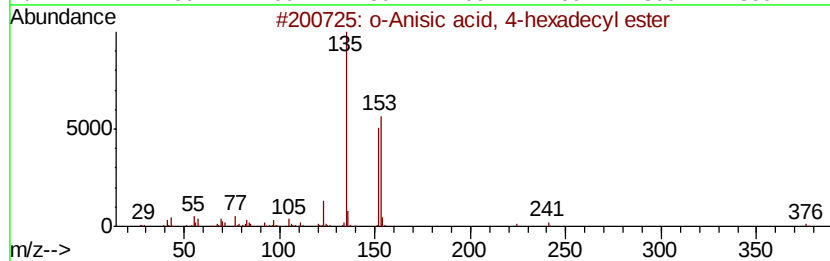
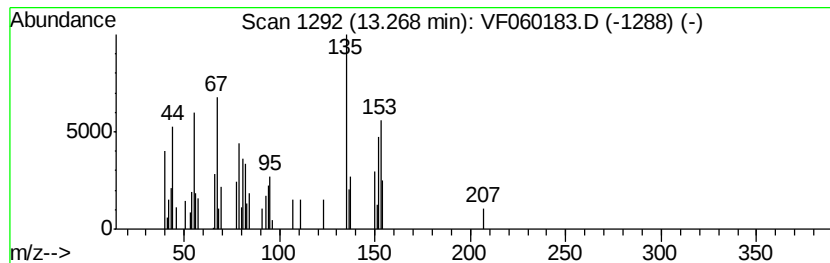
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Quant Title : SW846 8260

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

Peak Number 4 unknown13.27 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	23.60 ug/l	23143	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Anisic acid, 4-hexadecyl ester	376	C24H40O3	1000299-95-3	35
2		o-Anisic acid, 6-pentadecyl ester	362	C23H38O3	1000299-76-8	35
3		3,9-Epoxy-p-mentha-1,8(10)-diene	150	C10H14O	1000111-14-8	25
4		3-Aminosalicylic acid	153	C7H7NO3	000570-23-0	22
5		2-Ethoxy-5-nitropyridine	168	C7H8N2O3	031594-45-3	22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA F/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-01-082718-A

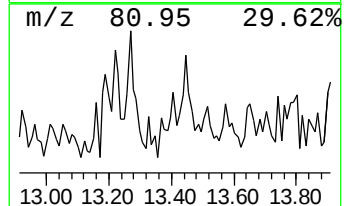
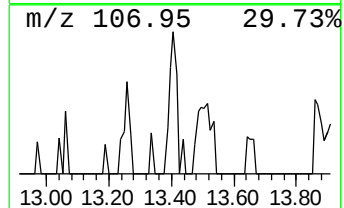
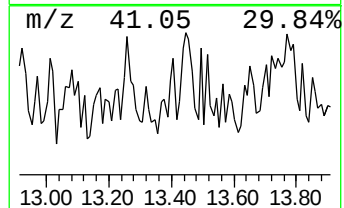
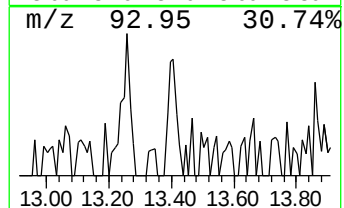
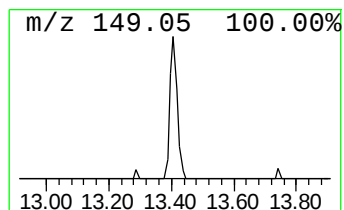
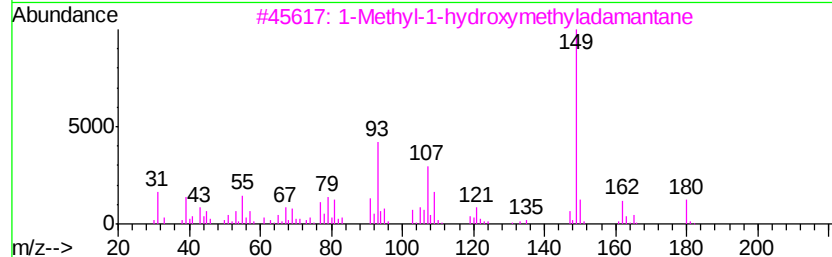
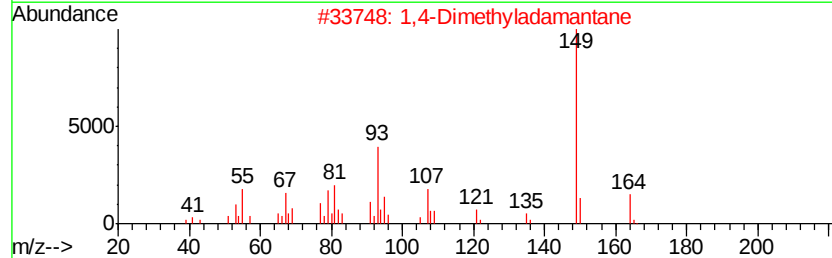
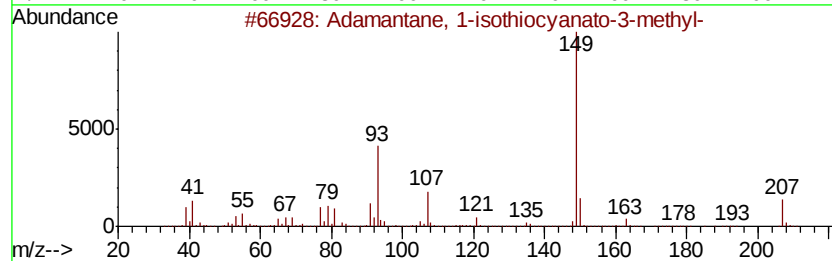
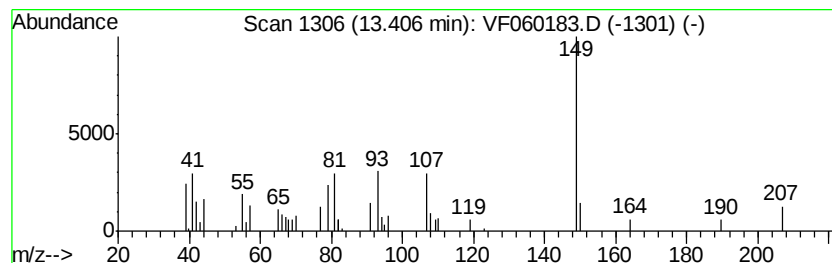
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Quant Title : SW846 8260

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

Peak Number 5 Adamantane, 1-isothiocyanat... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	20.63 ug/l	20237	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Adamantane, 1-isothiocyanato-3-m...	207	C12H17NS	136860-48-5	64
2		1,4-Dimethyladamantane	164	C12H20	024145-88-8	59
3		1-Methyl-1-hydroxymethyladamantane	180	C12H20O	001200-78-8	59
4		Adamantane, 1,3-dimethyl-	164	C12H20	000702-79-4	53
5		Pyridinium, 1-(acetylamino)-2,6-...	164	C9H12N2O	031020-35-6	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA F/SOIL
ALS Vial : 46 Sample Multiplier: 1

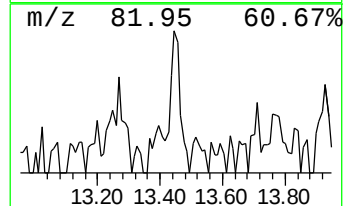
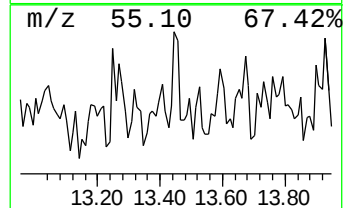
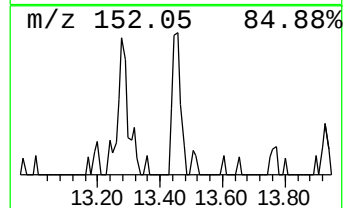
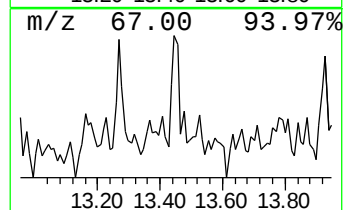
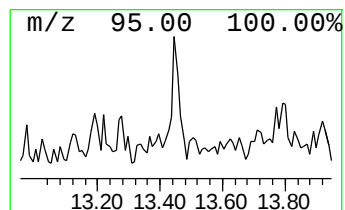
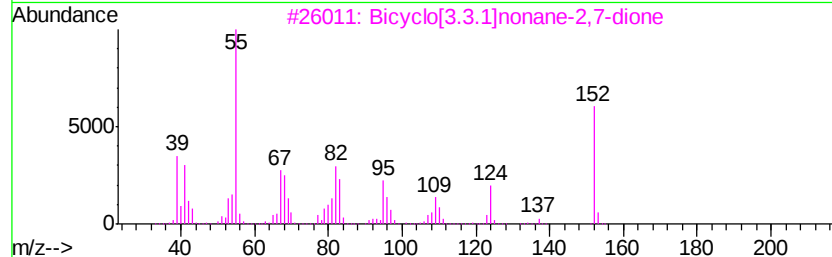
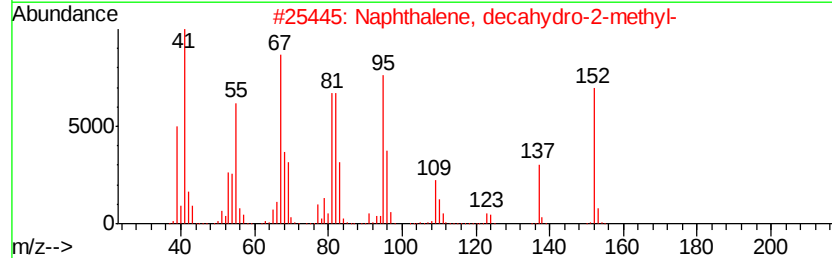
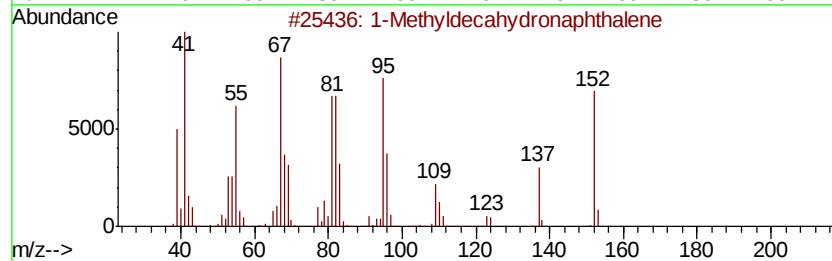
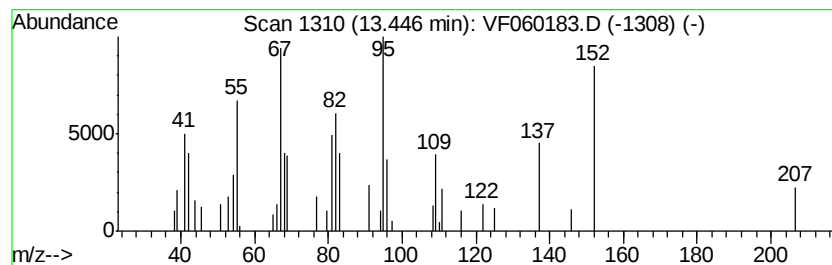
Instrument :
MSVOA_F
ClientSampleId :
OR-01-082718-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.M
Quant Title : SW846 8260

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

Peak Number 6 1-Methyldecahydronaphthalene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	17.70 ug/l	17360	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Methyldecahydronaphthalene	152 C11H20	002958-75-0	91
2	Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1	91
3	Bicyclo[3.3.1]nonane-2,7-dione	152 C9H12O2	1000210-30-3	60
4	Decalin, syn-1-methyl-, cis-	152 C11H20	1000158-89-1	49
5	1-Adamantanol	152 C10H16O	000768-95-6	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA F/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-01-082718-A

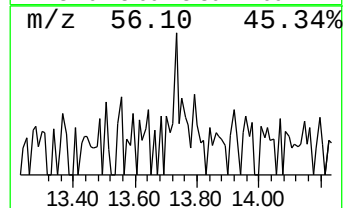
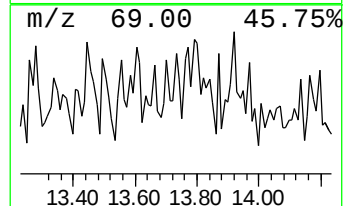
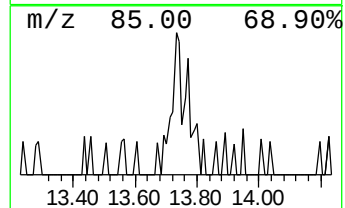
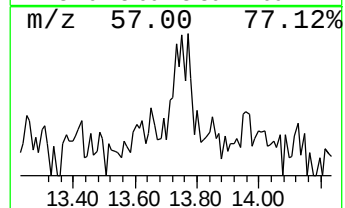
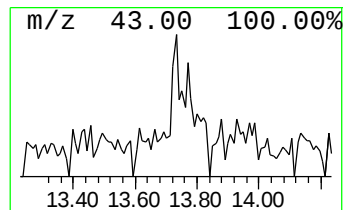
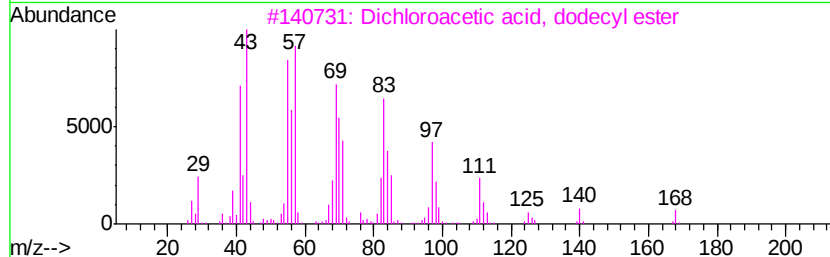
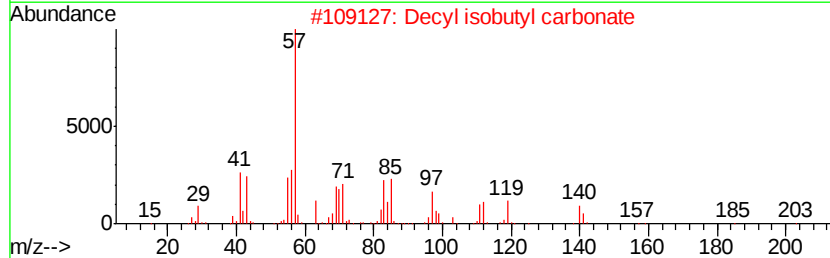
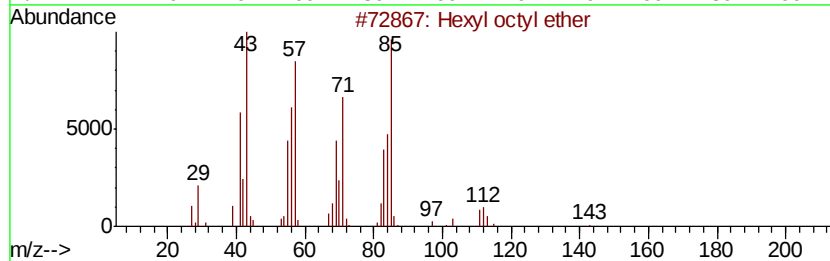
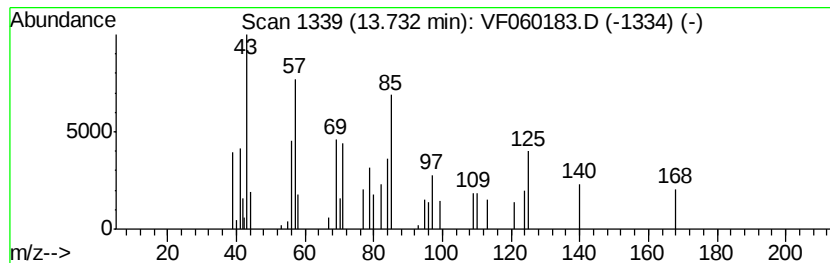
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Quant Title : SW846 8260

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

Peak Number 7 unknown13.73 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	11.85 ug/l	11622	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexyl octyl ether	214	C14H30O	017071-54-4	38
2		Decyl isobutyl carbonate	258	C15H30O3	959226-07-4	38
3		Dichloroacetic acid, dodecyl ester	296	C14H26Cl2O2	083005-01-0	35
4		Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	27
5		Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	27



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA F/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-01-082718-A

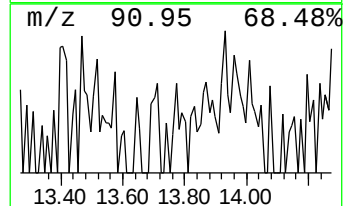
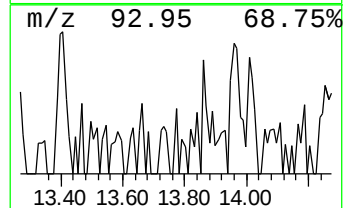
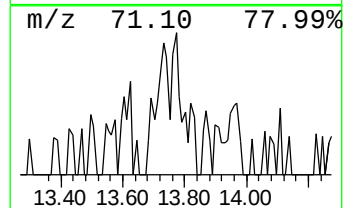
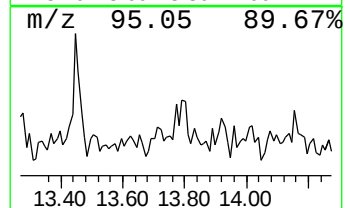
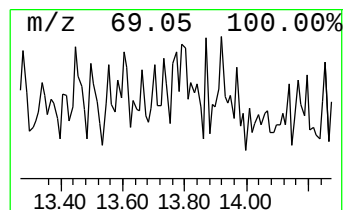
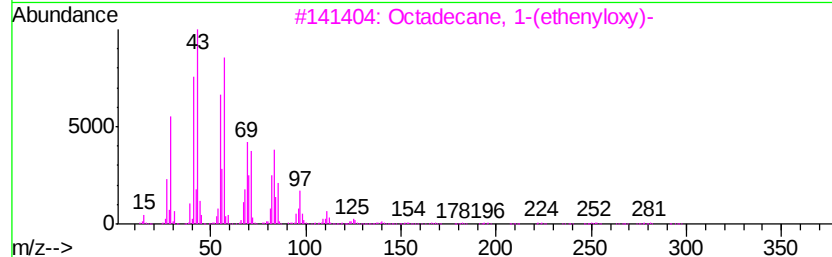
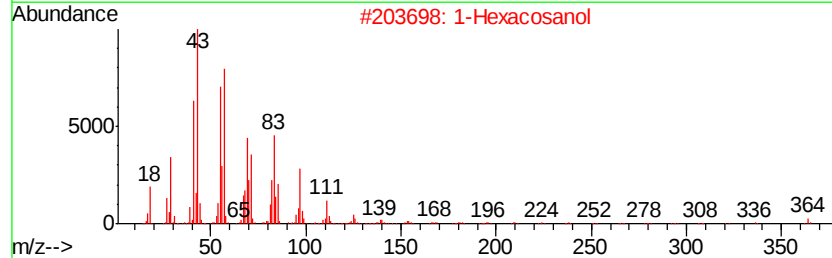
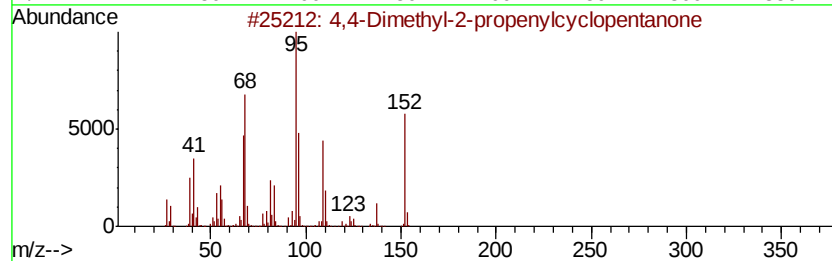
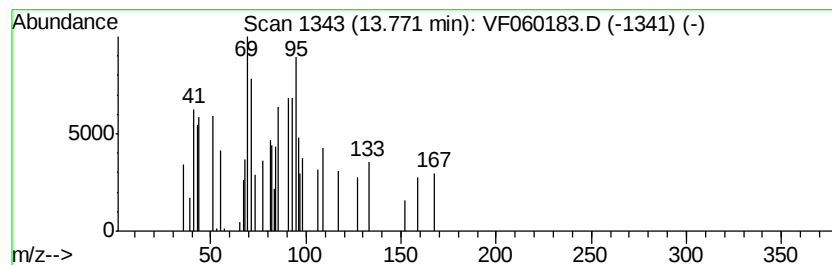
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Quant Title : SW846 8260

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

Peak Number 8 unknown13.77 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	13.39 ug/l	13137	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4,4-Dimethyl-2-propenylcyclopent...	152	C10H16O	068261-88-1	25
2		1-Hexacosanol	382	C26H54O	000506-52-5	17
3		Octadecane, 1-(ethenyl)-	296	C20H40	000930-02-9	17
4		Carbonic acid, ethyl octadecyl e...	342	C21H42O3	1000314-59-8	12
5		Bicyclo[4.1.0]heptan-2-one, 3,5,...	152	C10H16O	029750-24-1	12



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA F/SOIL
ALS Vial : 46 Sample Multiplier: 1

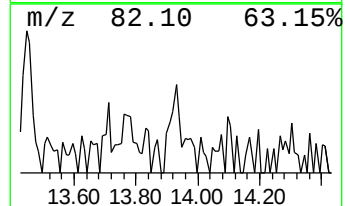
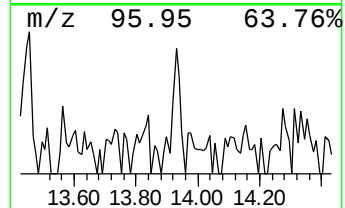
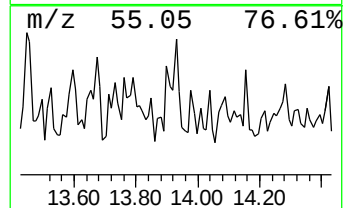
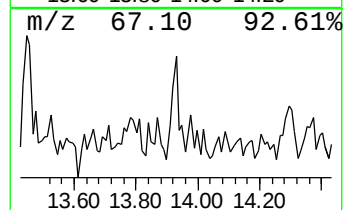
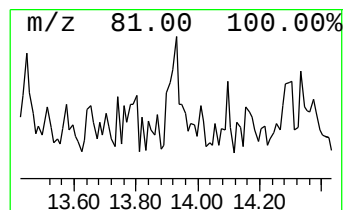
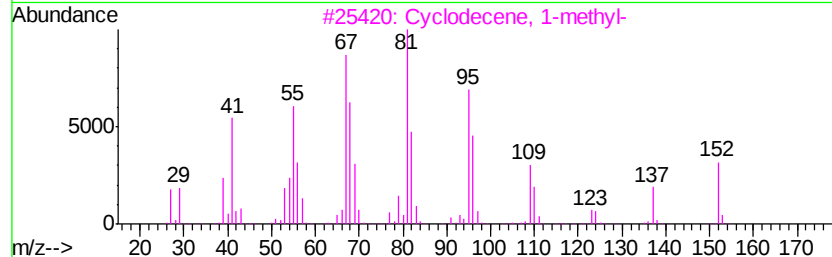
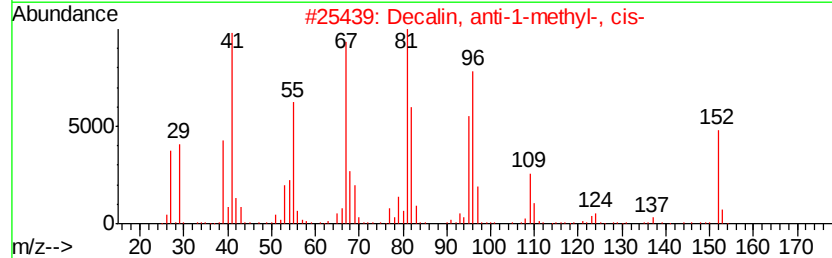
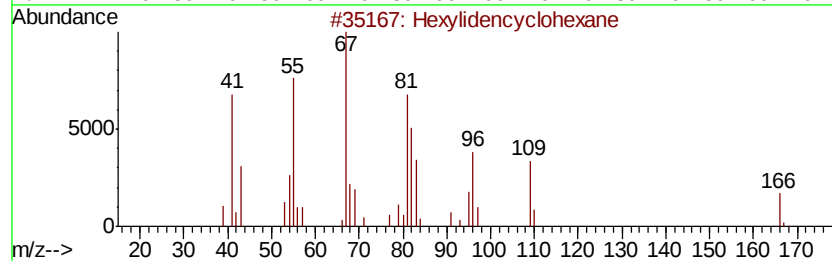
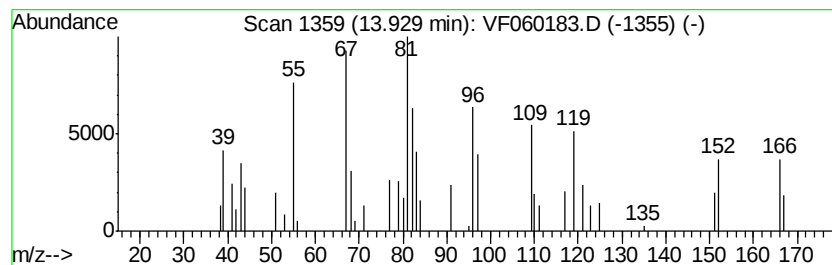
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ClientSampleId :
OR-01-082718-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.M
Quant Title : SW846 8260

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

Peak Number 9 Hexylidencyclohexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	16.27 ug/l	15962	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Hexylidencyclohexane	166 C12H22	010201-62-4 64
2		Decalin, anti-1-methyl-, cis-	152 C11H20	1000158-89-0 60
3		Cyclodecene, 1-methyl-	152 C11H20	066633-38-3 53
4		Bicyclo[3.3.2]decan-9-one	152 C10H16O	028054-91-3 53
5		Cyclohexene,1-hexyl-	166 C12H22	003964-66-7 47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA F/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-01-082718-A

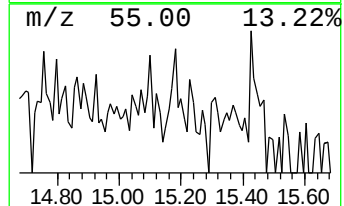
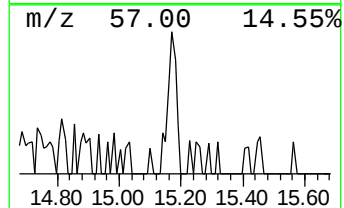
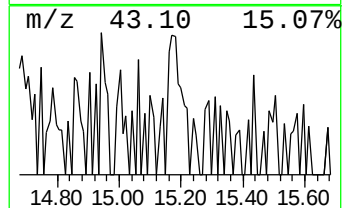
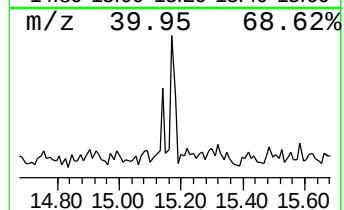
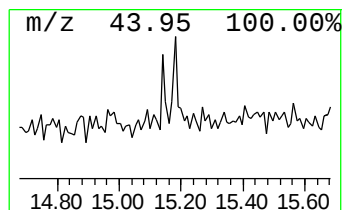
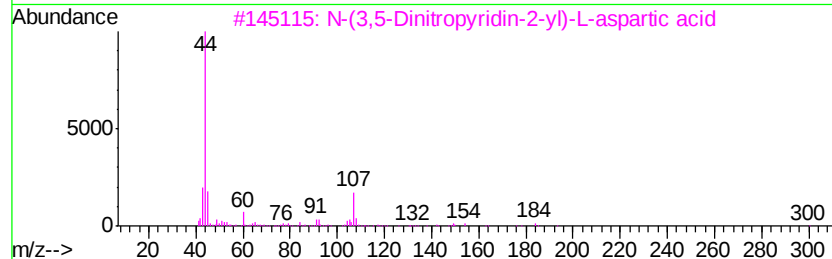
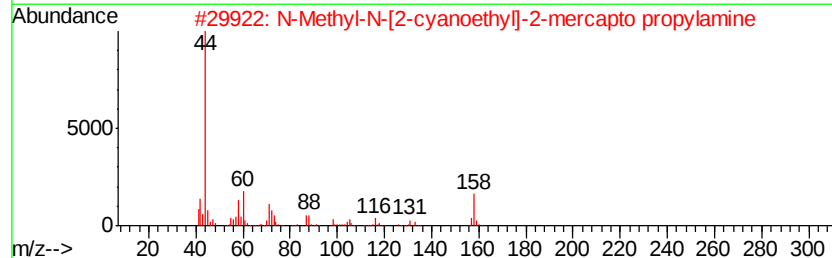
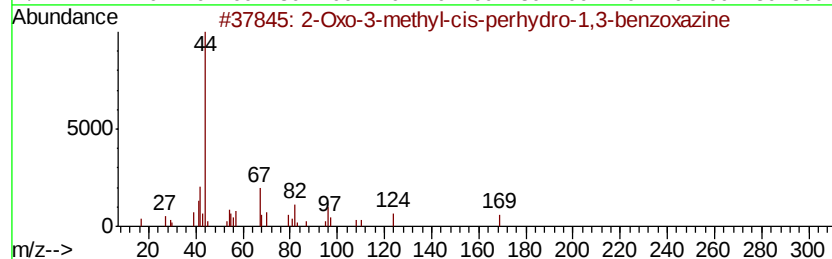
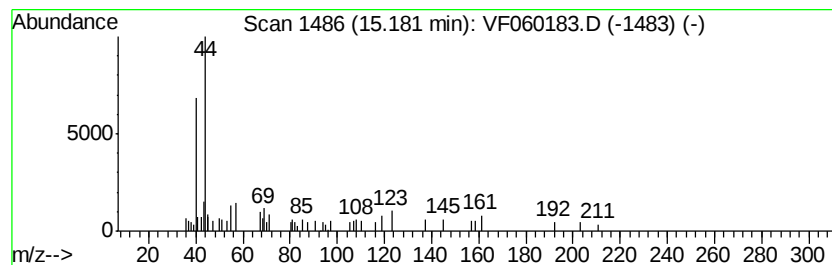
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Quant Title : SW846 8260

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

Peak Number 10 unknown15.18 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	12.75 ug/l	12501	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Oxo-3-methyl-cis-perhydro-1,3-...	169	C9H15N02	1000138-84-6	38
2		N-Methyl-N-[2-cyanoethyl]-2-merc...	158	C7H14N2S	1000257-11-6	36
3		N-(3,5-Dinitropyridin-2-yl)-L-as...	300	C9H8N4O8	035899-60-6	28
4		4-Amino-2-oxo-furazan-3-carboxyl...	145	C3H3N3O4	1000300-51-9	9
5		p-Hydroxyamphetamine	151	C9H13NO	000103-86-6	9



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF082918\
Data File : VF060183.D
Acq On : 29 Aug 2018 8:44
Operator : VA/AP
Sample : J4277-09
Misc : 6.48g/5mL/MSVOA_F/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-01-082718-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F082918S.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
unknown12.69	12.69	12.4	ug/l	12206	4	12.63	49039	50.0
unknown12.92	12.92	12.1	ug/l	11838	4	12.63	49039	50.0
unknown13.18	13.18	11.5	ug/l	11273	4	12.63	49039	50.0
unknown13.27	13.27	23.6	ug/l	23143	4	12.63	49039	50.0
Adamantane, 1-iso...	13.41	20.6	ug/l	20237	4	12.63	49039	50.0
1-Methyldecahydro...	13.45	17.7	ug/l	17360	4	12.63	49039	50.0
unknown13.73	13.73	11.9	ug/l	11622	4	12.63	49039	50.0
unknown13.77	13.77	13.4	ug/l	13137	4	12.63	49039	50.0
Hexylidencyclohexane	13.93	16.3	ug/l	15962	4	12.63	49039	50.0
unknown15.18	15.18	12.8	ug/l	12501	4	12.63	49039	50.0