

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
 Data File : VW021909.D
 Acq On : 14 Jan 2022 13:25
 Operator : SY/VA
 Sample : N1115-12
 Misc : 6.26g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BGLS2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
 Title : SFAM01.0

Signal : TIC: VW021909.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.952	4	8	10	rBV3	1775	2449	0.00%	0.002%
2	1.989	10	14	20	rVB	6518	10758	0.01%	0.009%
3	2.148	36	40	41	rBV3	303	466	0.00%	0.000%
4	2.196	46	48	49	rBV2	652	589	0.00%	0.001%
5	2.349	63	73	81	rBV	55665	94467	0.11%	0.081%
6	2.410	81	83	84	rVB2	516	297	0.00%	0.000%
7	2.440	85	88	89	rBV2	423	474	0.00%	0.000%
8	2.495	91	97	104	rVV	2884	5665	0.01%	0.005%
9	2.617	115	117	118	rBV2	818	560	0.00%	0.000%
10	2.666	118	125	127	rBV6	1696	3328	0.00%	0.003%
11	2.800	145	147	148	rBV	572	402	0.00%	0.000%
12	2.879	153	160	172	rVV	38671	83976	0.10%	0.072%
13	3.019	176	183	186	rVV7	1321	3498	0.00%	0.003%
14	3.050	186	188	190	rVV3	1026	1183	0.00%	0.001%
15	3.093	193	195	198	rVV3	1162	1897	0.00%	0.002%
16	3.141	201	203	208	rVB3	1134	1480	0.00%	0.001%
17	3.257	219	222	225	rBV	202	282	0.00%	0.000%
18	3.300	225	229	230	rBV2	1517	1491	0.00%	0.001%
19	3.312	230	231	237	rVB2	1618	2247	0.00%	0.002%
20	3.373	237	241	244	rBV2	521	858	0.00%	0.001%
21	3.446	246	253	265	rVB2	5666	13049	0.02%	0.011%
22	3.617	277	281	283	rVB3	347	492	0.00%	0.000%
23	3.684	283	292	298	rBV3	2244	5650	0.01%	0.005%
24	3.781	298	308	319	rVB2	6359	17458	0.02%	0.015%
25	3.903	326	328	330	rBV	282	294	0.00%	0.000%
26	4.037	333	350	381	rBV	1494062	4414479	5.27%	3.804%
27	4.300	391	393	396	rVB	513	477	0.00%	0.000%
28	4.379	396	406	416	rBV2	5318	16703	0.02%	0.014%
29	4.452	416	418	422	rVB	263	268	0.00%	0.000%
30	4.903	481	492	502	rBV8	2807	9935	0.01%	0.009%
31	5.037	511	514	516	rBV2	247	320	0.00%	0.000%
32	5.196	537	540	544	rVB2	179	263	0.00%	0.000%
33	5.421	568	577	587	rBV2	5396	15950	0.02%	0.014%
34	5.763	631	633	634	rBV	272	273	0.00%	0.000%
35	5.787	634	637	642	rVB5	272	365	0.00%	0.000%
36	6.214	691	707	724	rBV	818020	2492079	2.97%	2.148%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
 Title : SFAM01.0

37	6.903	815	820	823	rBV4	325	481	0.00%	0.000%
38	7.080	840	849	853	rBV	13011	34518	0.04%	0.030%
39	7.165	853	863	875	rVB2	200719	535273	0.64%	0.461%
40	7.256	875	878	879	rBV3	700	825	0.00%	0.001%
41	7.482	912	915	919	rBV2	197	309	0.00%	0.000%
42	7.647	929	942	955	rBV	99108	246289	0.29%	0.212%
43	7.854	974	976	978	rBV	232	254	0.00%	0.000%
44	7.915	982	986	991	rBV4	597	1124	0.00%	0.001%
45	8.110	1015	1018	1021	rBV	262	385	0.00%	0.000%
46	8.274	1035	1045	1059	rBV2	174501	548810	0.65%	0.473%
47	8.396	1059	1065	1075	rVB4	6667	16103	0.02%	0.014%
48	8.561	1090	1092	1096	rVB4	601	748	0.00%	0.001%
49	8.616	1100	1101	1107	rVB3	347	469	0.00%	0.000%
50	8.689	1110	1113	1115	rBV2	483	447	0.00%	0.000%
51	8.713	1115	1117	1122	rVB3	650	813	0.00%	0.001%
52	8.841	1128	1138	1148	rBV	160499	320736	0.38%	0.276%
53	8.909	1148	1149	1154	rVB2	565	439	0.00%	0.000%
54	9.091	1168	1179	1195	rBV	11870835	20877335	24.92%	17.991%
55	9.201	1195	1197	1199	rVV3	1856	2232	0.00%	0.002%
56	9.274	1199	1209	1220	rVV	113432	237013	0.28%	0.204%
57	9.347	1220	1221	1226	rVV4	1253	1241	0.00%	0.001%
58	9.384	1226	1227	1229	rVV2	507	421	0.00%	0.000%
59	9.402	1229	1230	1232	rVV2	635	627	0.00%	0.001%
60	9.433	1232	1235	1236	rVV3	1035	1102	0.00%	0.001%
61	9.476	1240	1242	1244	rVV2	397	355	0.00%	0.000%
62	9.634	1266	1268	1269	rBV2	558	494	0.00%	0.000%
63	9.665	1269	1273	1278	rVV4	1021	1450	0.00%	0.001%
64	9.732	1278	1284	1291	rVB8	2178	4755	0.01%	0.004%
65	9.780	1291	1292	1295	rBV3	291	292	0.00%	0.000%
66	9.847	1300	1303	1304	rBV3	481	503	0.00%	0.000%
67	9.878	1304	1308	1316	rVB3	4265	9190	0.01%	0.008%
68	9.939	1316	1318	1319	rBV2	459	444	0.00%	0.000%
69	10.030	1327	1333	1344	rBV2	48735	89270	0.11%	0.077%
70	10.116	1345	1347	1349	rVV3	616	582	0.00%	0.001%
71	10.134	1349	1350	1353	rVB3	633	356	0.00%	0.000%
72	10.238	1365	1367	1369	rVB3	514	295	0.00%	0.000%
73	10.323	1373	1381	1387	rBV	220980	400505	0.48%	0.345%
74	10.384	1387	1391	1400	rVB	40921	70426	0.08%	0.061%
75	10.445	1400	1401	1403	rVB2	623	387	0.00%	0.000%
76	10.494	1405	1409	1411	rBV4	752	1045	0.00%	0.001%

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 Title : SFAM01.0

77	10.579	1414	1423	1429	rBV	25312	46349	0.06%	0.040%
78	10.701	1438	1443	1447	rBV2	2545	4560	0.01%	0.004%
79	10.756	1447	1452	1455	rBV2	2605	4145	0.00%	0.004%
80	10.792	1455	1458	1460	rBV3	1730	1727	0.00%	0.001%
81	10.872	1461	1471	1479	rBV3	34512051	83791175	100.00%	72.207%
82	11.286	1534	1539	1542	rBV7	3149	6964	0.01%	0.006%
83	11.323	1543	1545	1547	rVV3	1983	1833	0.00%	0.002%
84	11.628	1588	1595	1606	rVB	172341	289443	0.35%	0.249%
85	11.725	1606	1611	1620	rVB	7256	13421	0.02%	0.012%
86	11.835	1624	1629	1634	rBV2	8368	15043	0.02%	0.013%
87	12.164	1678	1683	1690	rVB2	19715	33534	0.04%	0.029%
88	12.365	1712	1716	1719	rBV5	4035	6710	0.01%	0.006%
89	12.457	1728	1731	1737	rBV5	4379	8870	0.01%	0.008%
90	12.603	1750	1755	1764	rVV	18618	32615	0.04%	0.028%
91	12.688	1764	1769	1775	rVB	67363	109691	0.13%	0.095%
92	12.871	1783	1799	1816	rBV3	44019	265170	0.32%	0.229%
93	13.158	1843	1846	1850	rBV5	4950	9648	0.01%	0.008%
94	13.304	1857	1870	1881	rBV4	64185	292809	0.35%	0.252%
95	13.438	1883	1892	1903	rVV6	23324	91719	0.11%	0.079%
96	13.554	1906	1911	1917	rVB	95960	152220	0.18%	0.131%
97	13.853	1954	1960	1966	rBV	88608	166795	0.20%	0.144%
98	15.066	2154	2159	2167	rVB4	21770	55058	0.07%	0.047%
99	15.322	2197	2201	2205	rBV4	9579	19225	0.02%	0.017%
100	16.273	2354	2357	2362	rBV5	6177	11160	0.01%	0.010%

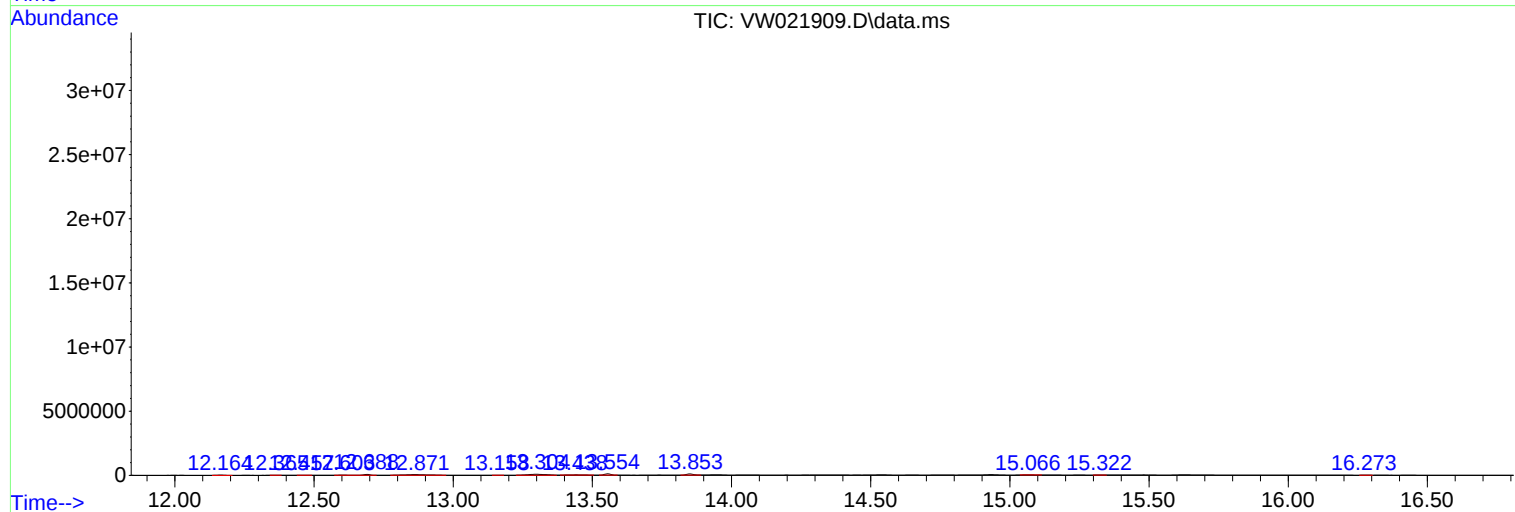
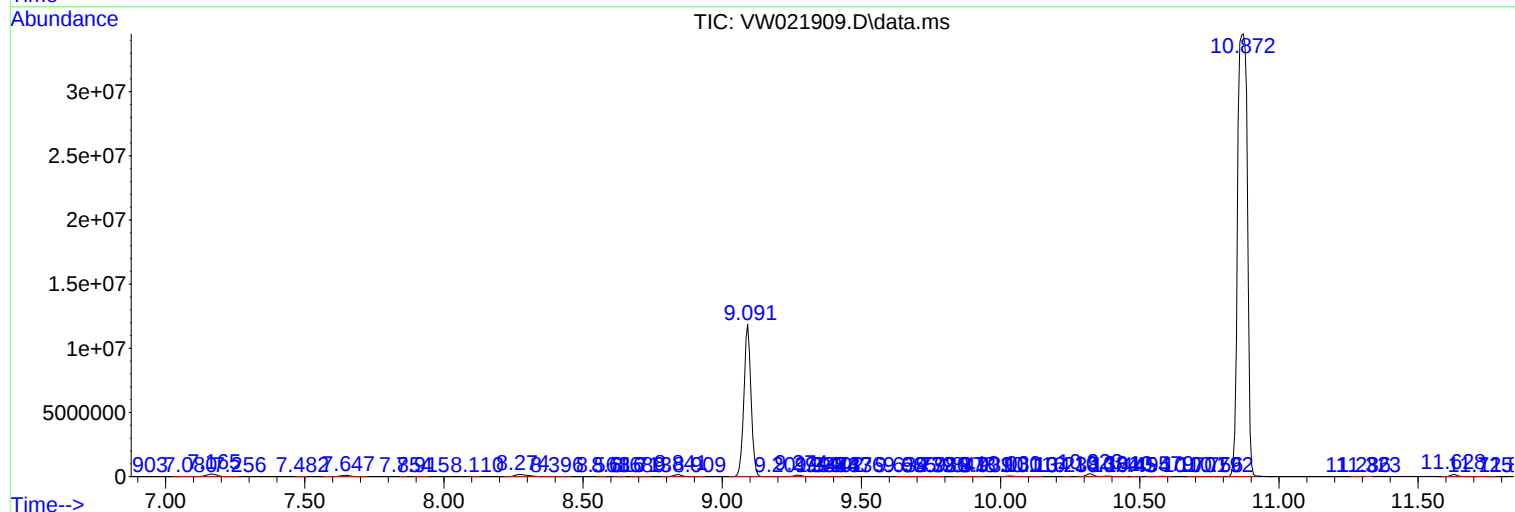
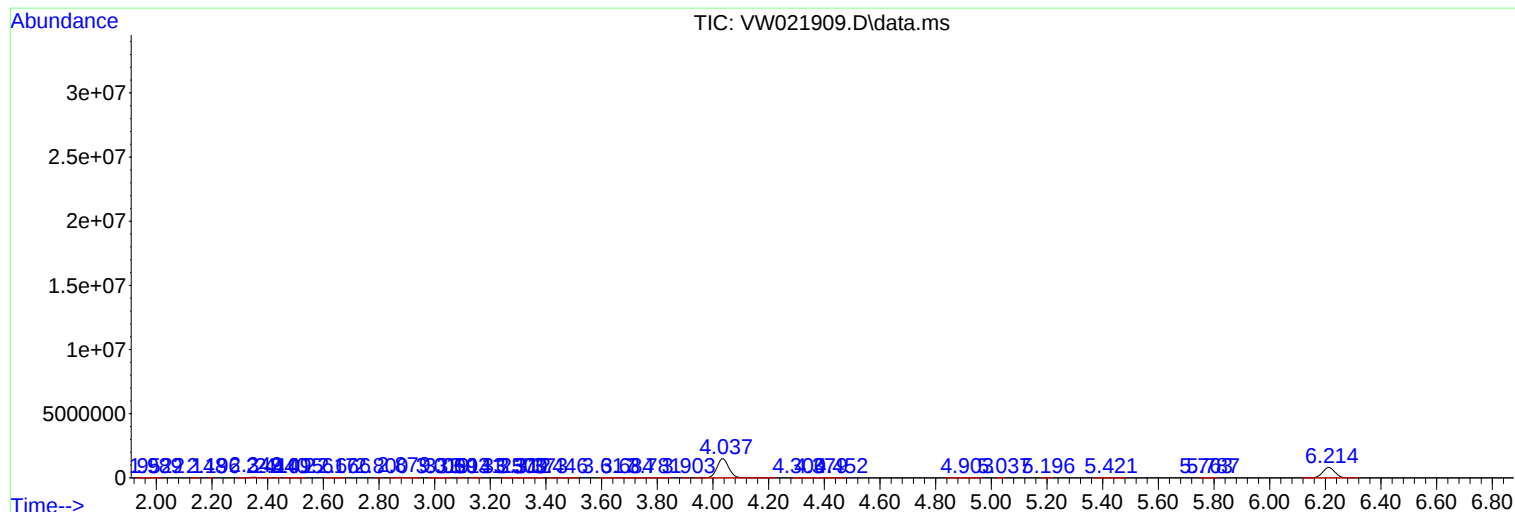
Sum of corrected areas: 116042649

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MSVOA_W
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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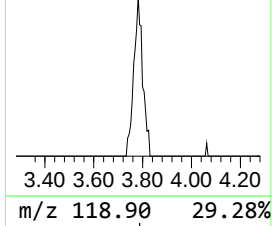
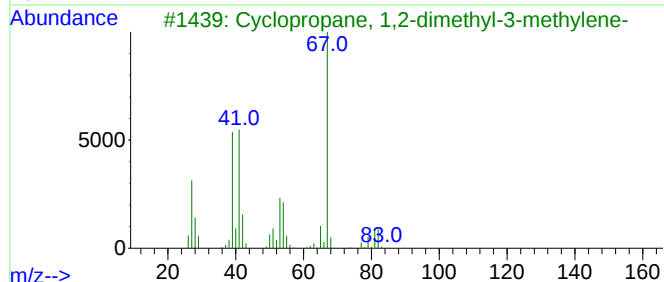
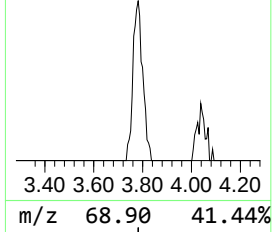
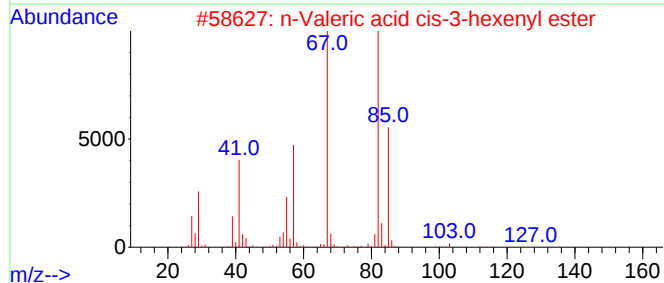
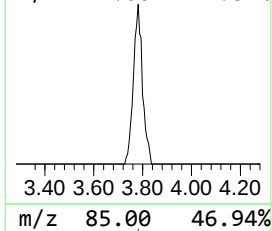
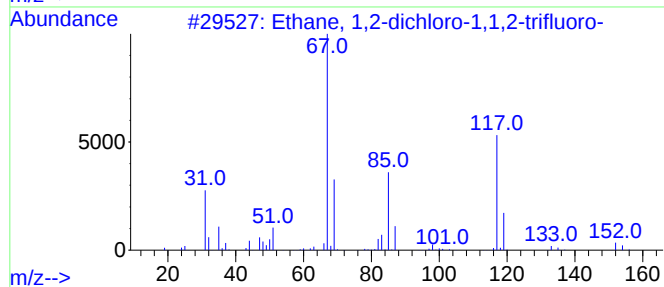
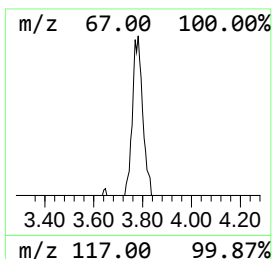
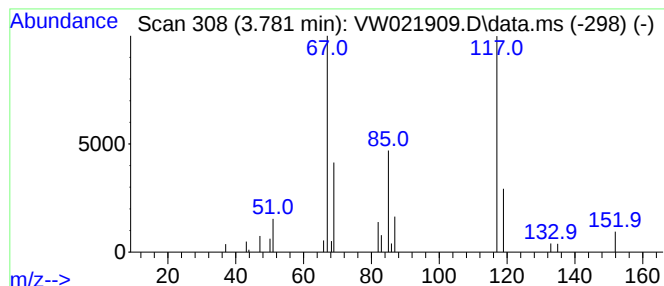
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethane, 1,2-dichloro-1,1,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.781	1.36 ug/L	17458	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,2-dichloro-1,1,2-trifl...	152	C2HCl2F3	000354-23-4	52
2			n-Valeric acid cis-3-hexenyl ester	184	C11H20O2	035852-46-1	17
3			Cyclopropane, 1,2-dimethyl-3-met...	82	C6H10	062338-02-7	14
4			2-Pentyn-4-one	82	C5H6O	007299-55-0	10
5			Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	10



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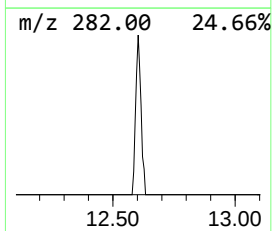
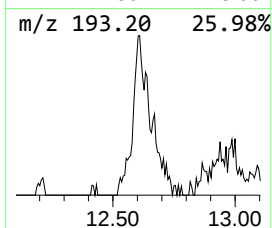
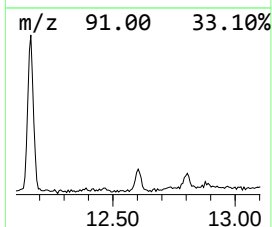
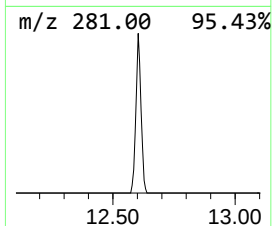
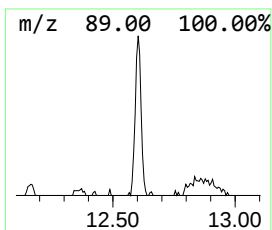
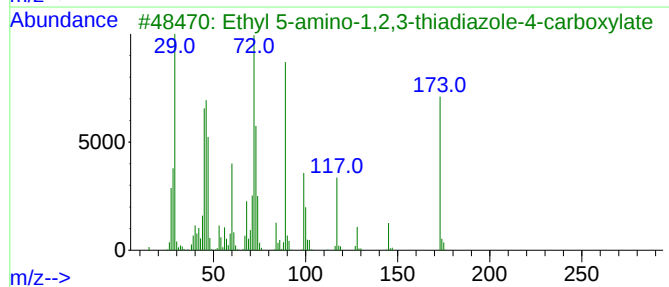
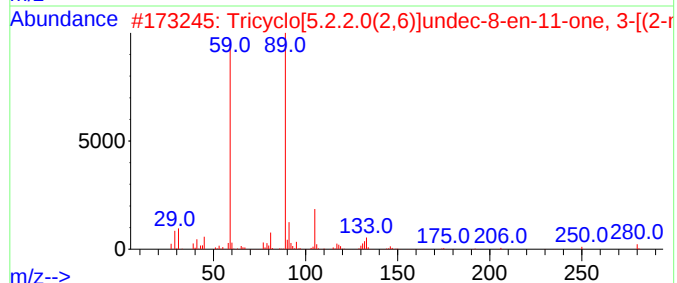
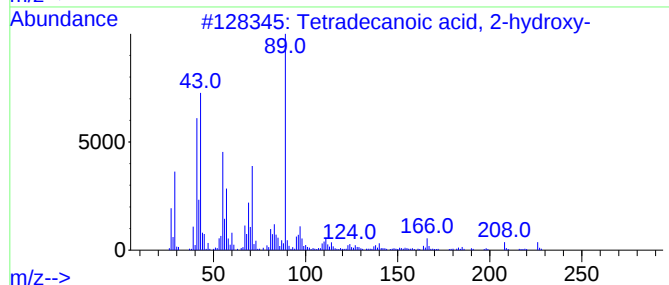
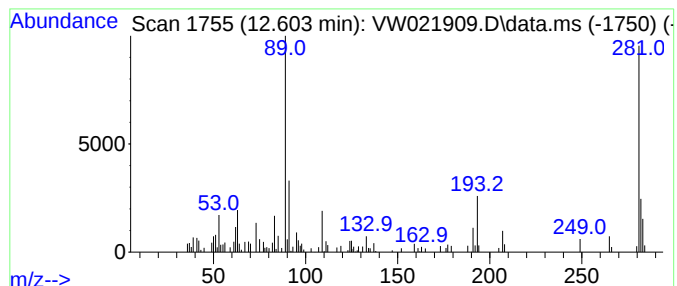
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.603	5.36 ug/L	32615	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid, 2-hydroxy-	244	C14H28O3	002507-55-3	47
2			Tricyclo[5.2.2.0(2,6)]undec-8-en...	280	C16H24O4	1000154-01-3	47
3			Ethyl 5-amino-1,2,3-thiadiazole-	173	C5H7N3O2S	006440-02-4	47
4			(2R,5S)-5-Methyl-2-(2-(methylthi...	200	C11H20O5	032637-94-8	38
5			2,2-Dimethoxypropionamide	133	C5H11NO3	1000237-69-7	38



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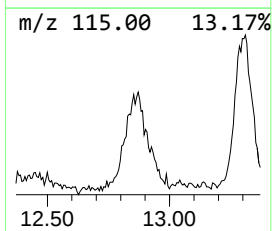
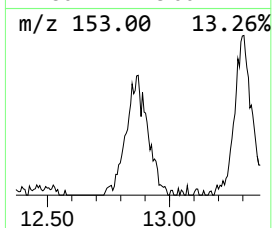
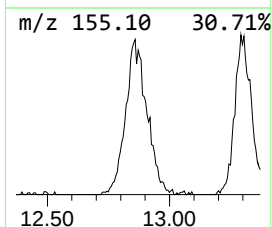
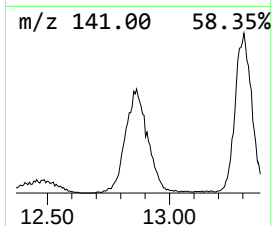
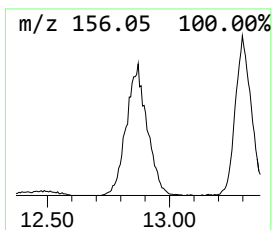
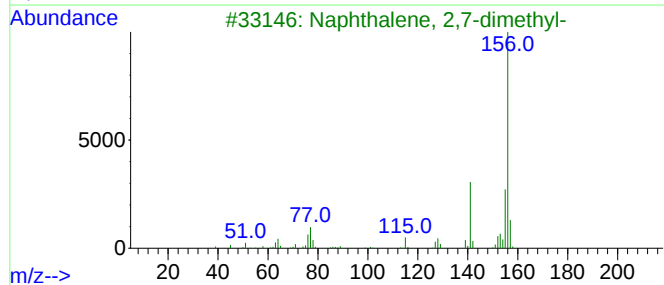
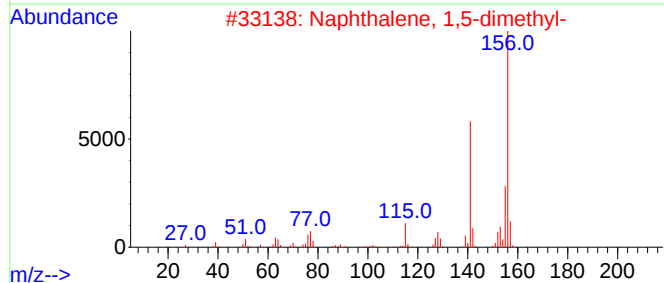
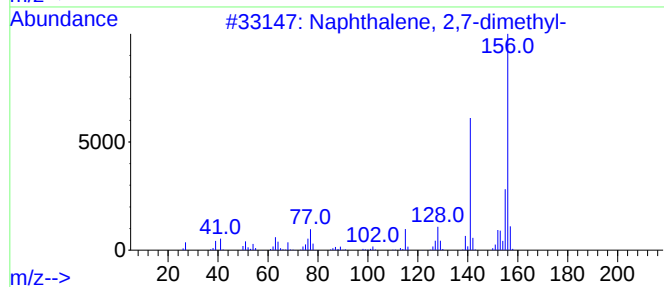
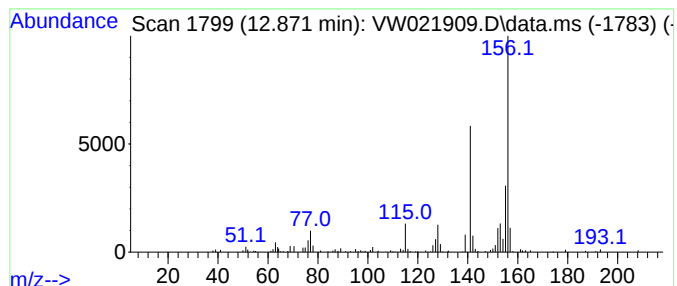
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 2,7-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.871	43.55 ug/L	265170	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

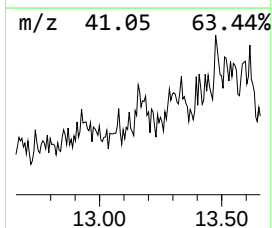
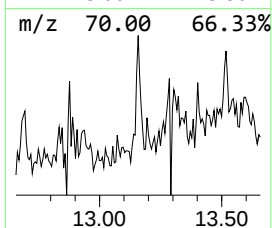
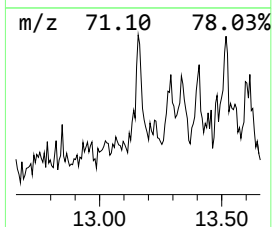
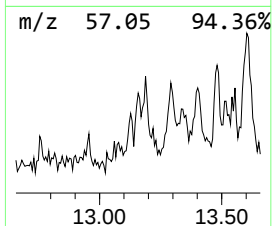
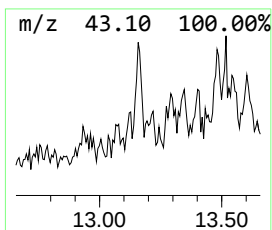
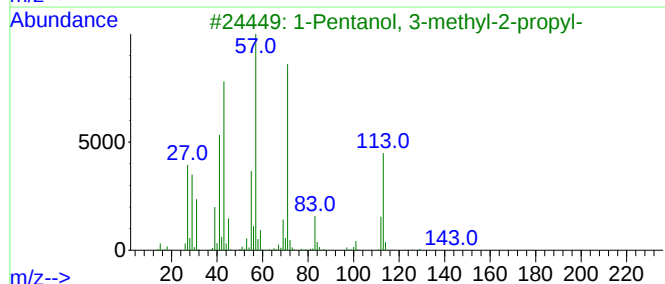
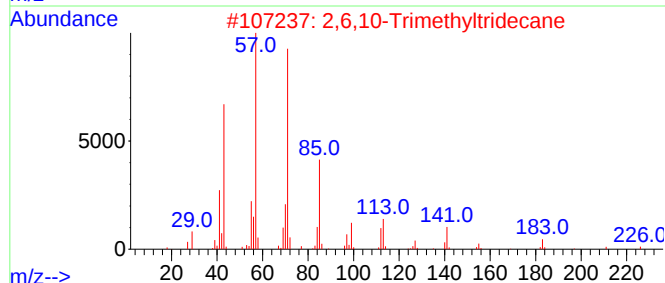
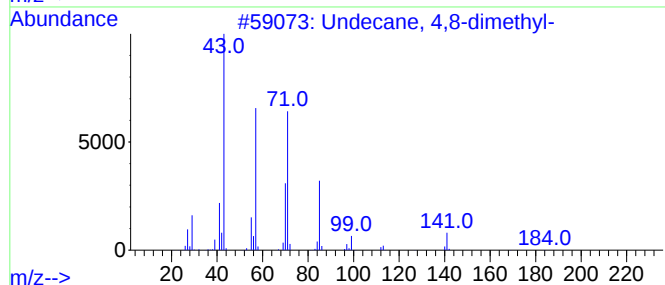
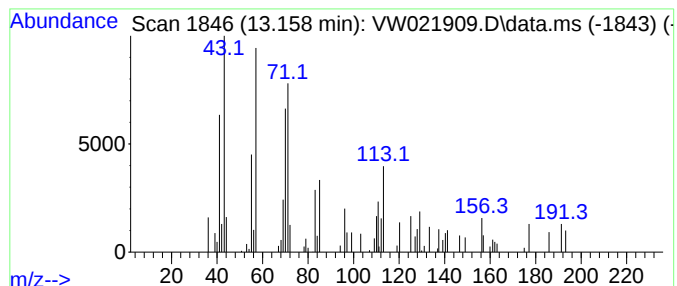
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MSVOA_W
ClientSampleId :
BGLS2

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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.158	1.58 ug/L	9648	1,4-Dichlorobenzene-d4		13.554	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane, 4,8-dimethyl-		184	C13H28	017301-33-6	35
2	2,6,10-Trimethyltridecane		226	C16H34	003891-99-4	35
3	1-Pentanol, 3-methyl-2-propyl-		144	C9H20O	054004-40-9	35
4	Dodecane, 4-methyl-		184	C13H28	006117-97-1	35
5	Hexadecane, 7-methyl-		240	C17H36	026730-20-1	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGLS2

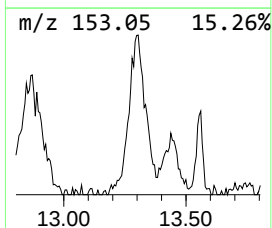
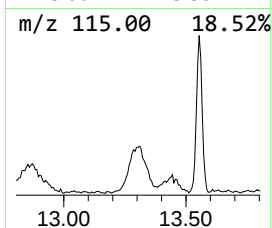
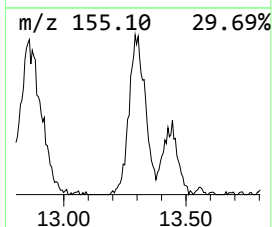
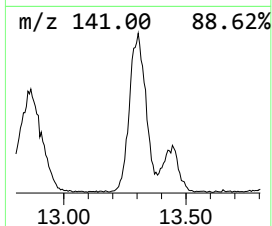
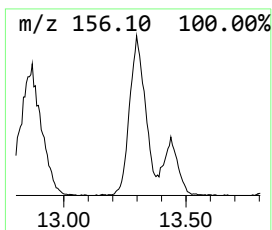
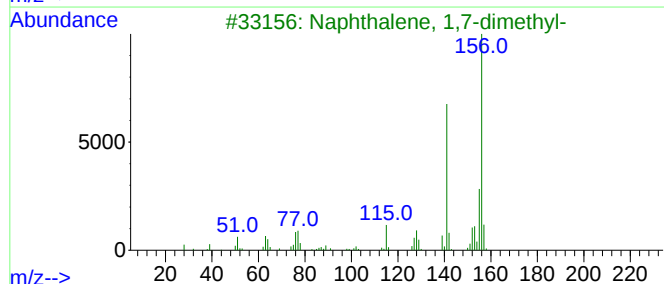
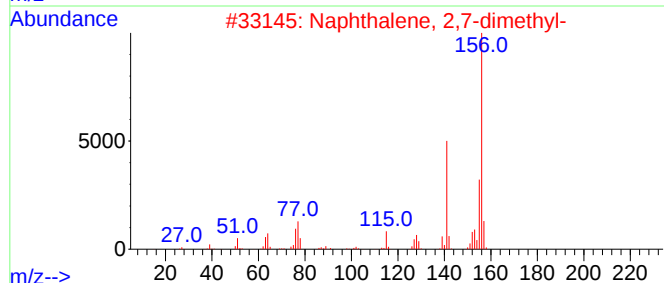
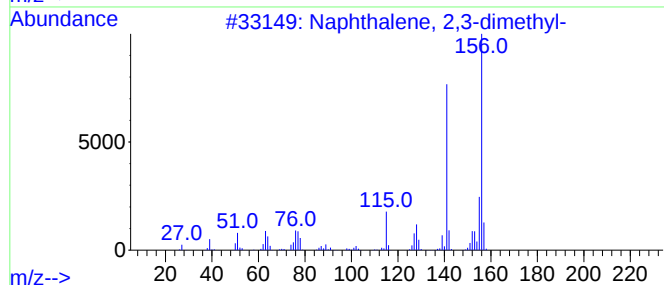
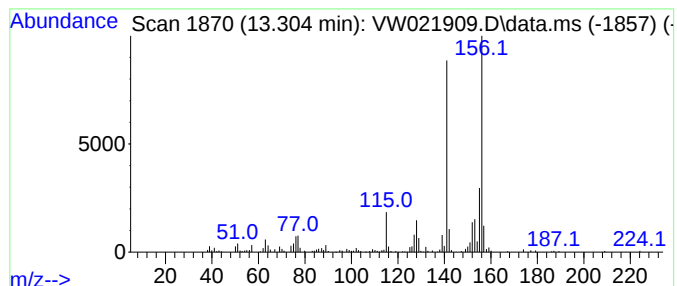
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.304	48.09 ug/L	292809	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGLS2

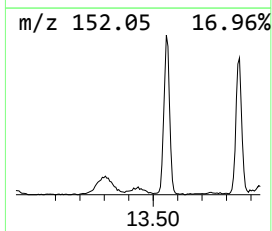
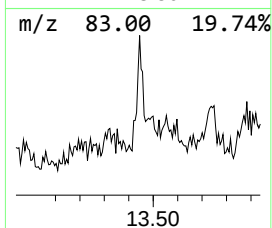
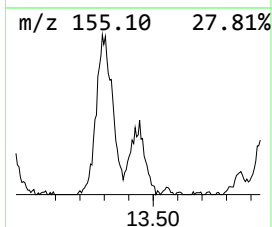
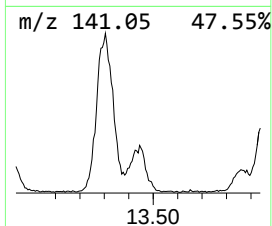
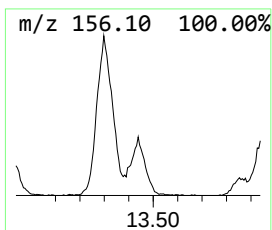
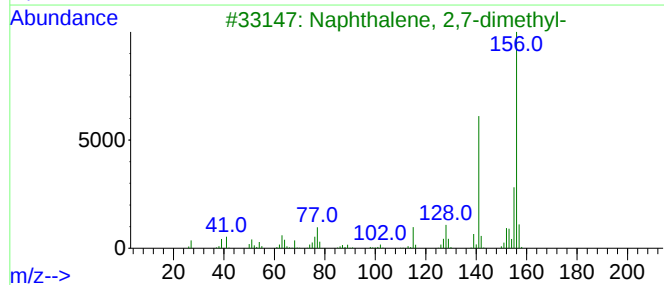
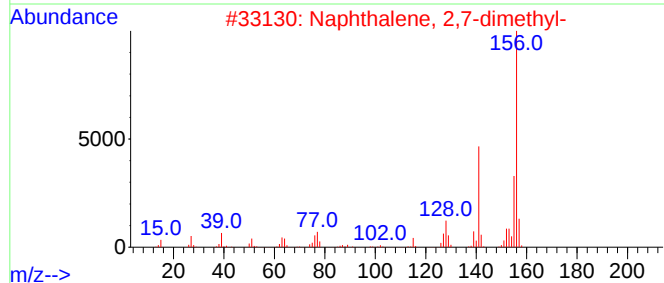
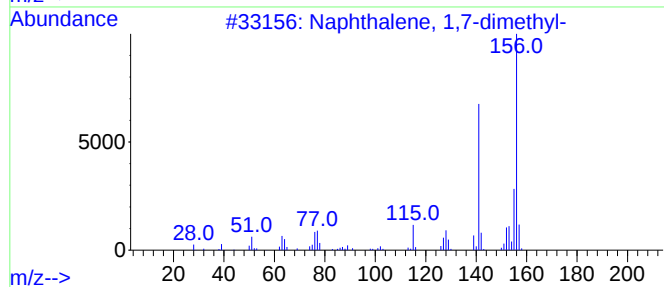
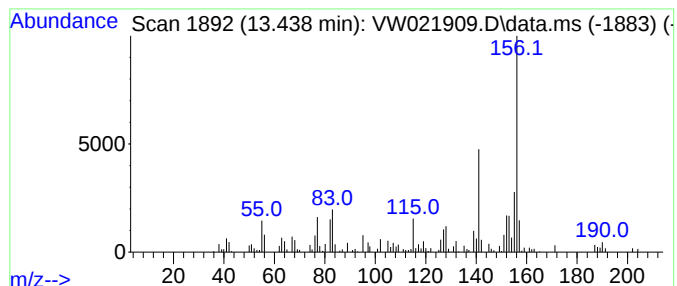
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1,7-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.438	15.06 ug/L	91719	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	91
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	91
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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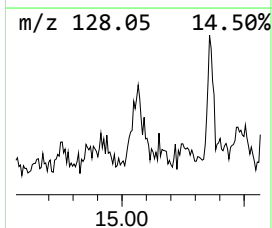
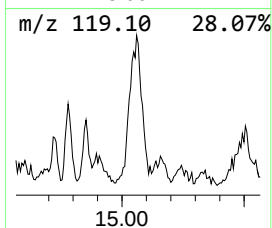
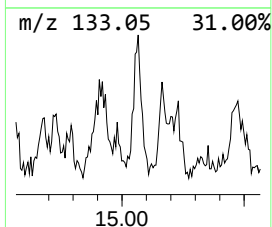
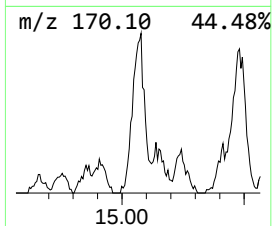
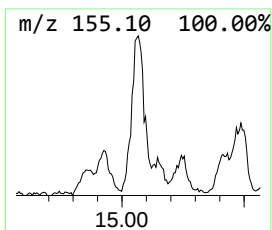
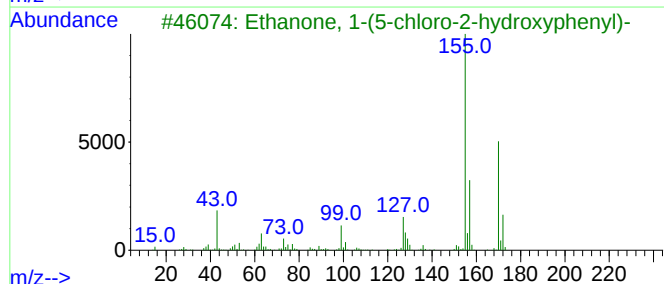
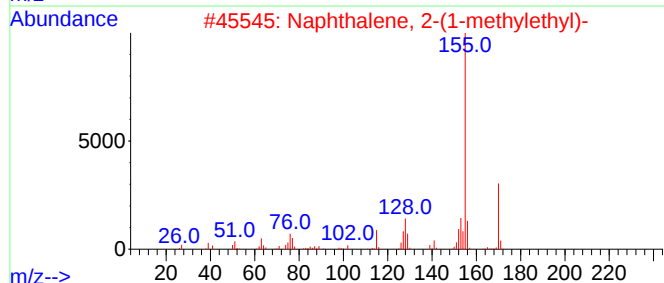
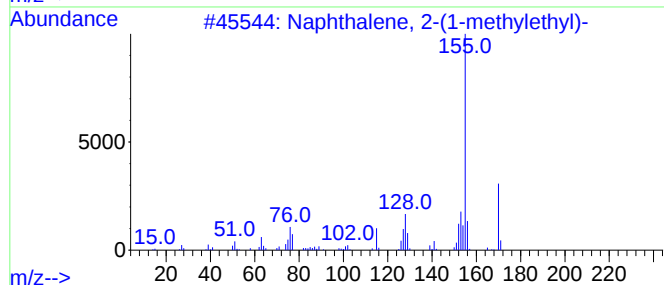
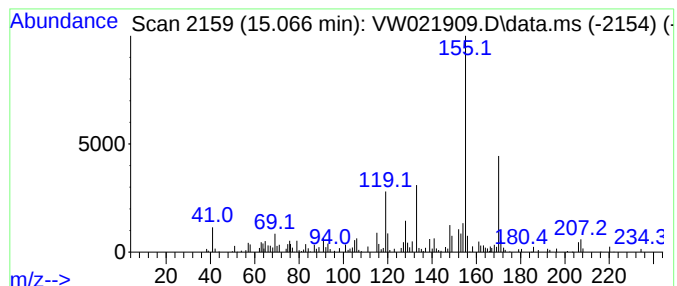
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, 2-(1-methyleth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.066	9.04 ug/L	55058	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	58
2			Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	53
3			Ethanone, 1-(5-chloro-2-hydroxyp...	170	C8H7ClO2	001450-74-4	49
4			2-Naphthyl methyl ketone	170	C12H10O	000093-08-3	49
5			Ethanone, 1-(5-chloro-2-hydroxyp...	170	C8H7ClO2	001450-74-4	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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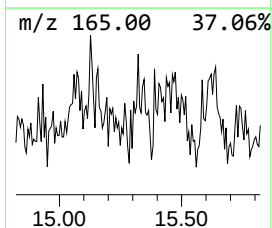
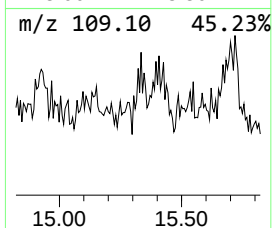
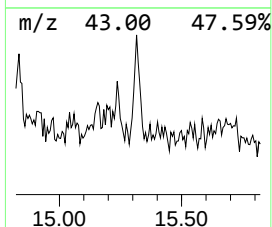
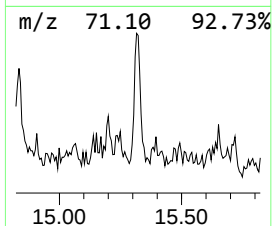
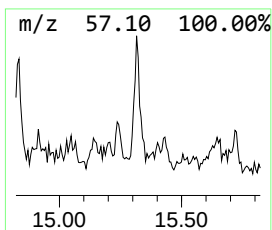
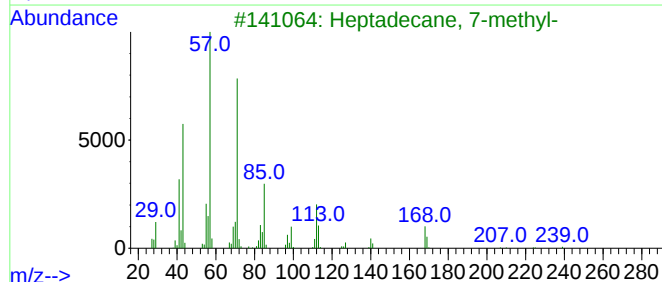
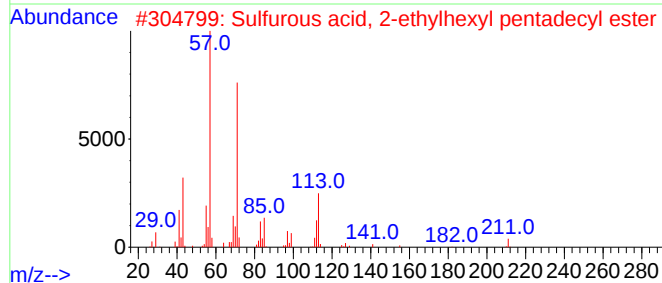
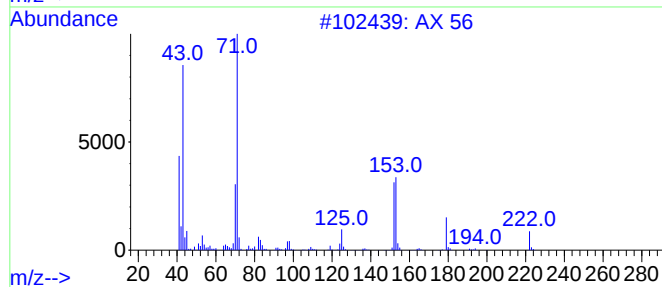
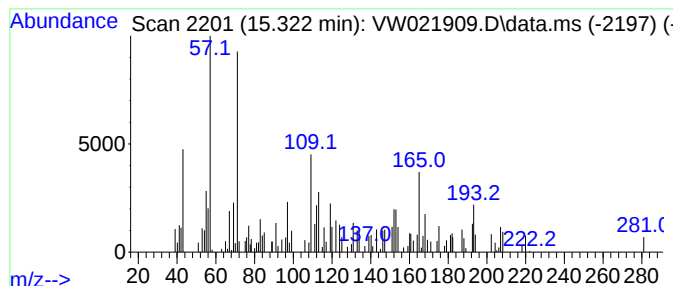
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.322	3.16 ug/L	19225	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	AX	56		222	C9H10N4OS	042204-09-1	35
2	Sulfurous acid, 2-ethylhexyl pen...			404	C23H48O3S	1000309-19-8	22
3	Heptadecane, 7-methyl-			254	C18H38	020959-33-5	22
4	Undecane, 6,6-dimethyl-			184	C13H28	017312-76-4	22
5	Sulfurous acid, decyl 2-ethylhex...			334	C18H38O3S	1000309-19-3	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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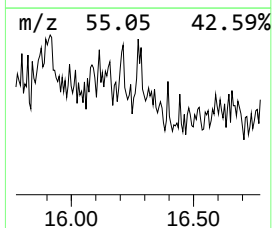
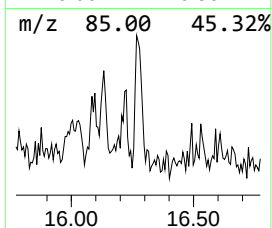
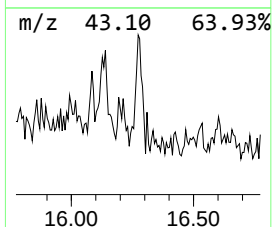
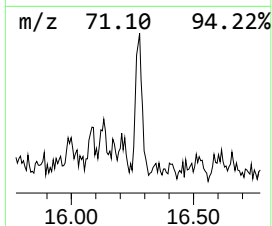
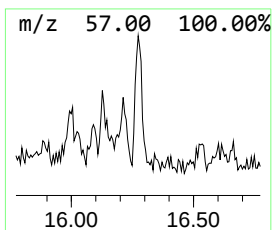
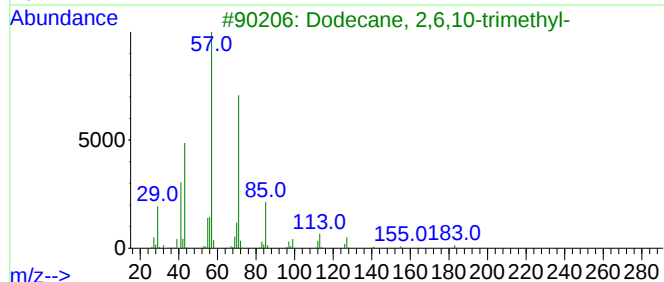
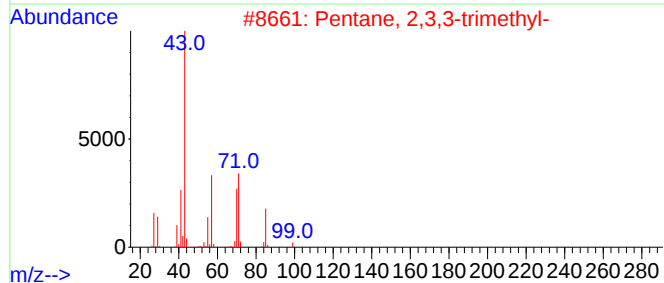
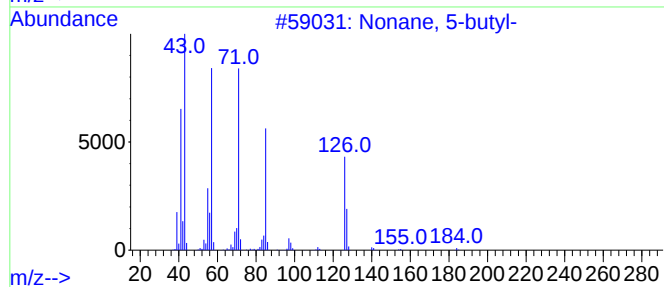
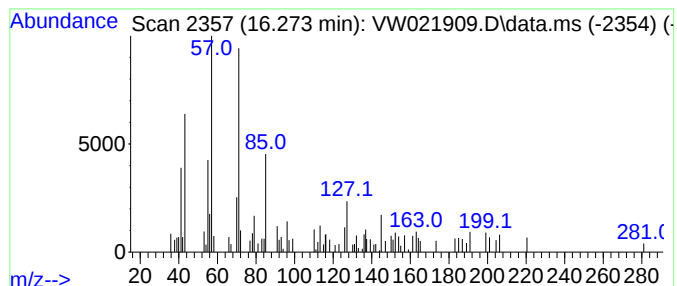
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.273	1.83 ug/L	11160	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 5-butyl-	184	C13H28	017312-63-9	50
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	47
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	38
4			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	38
5			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW011422\
Data File : VW021909.D
Acq On : 14 Jan 2022 13:25
Operator : SY/VA
Sample : N1115-12
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGLS2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM011122SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown-01	12.603	5.4	ug/L	32615	3	13.554	152220	25.0
Naphthalene, 2,...	12.871	43.5	ug/L	265170	3	13.554	152220	25.0
(DEL) Alkane: S...	13.158	1.6	ug/L	9648	3	13.554	152220	25.0
Naphthalene, 2,...	13.304	48.1	ug/L	292809	3	13.554	152220	25.0
Naphthalene, 1,...	13.438	15.1	ug/L	91719	3	13.554	152220	25.0
Naphthalene, 2-...	15.066	9.0	ug/L	55058	3	13.554	152220	25.0
unknown-02	15.322	3.2	ug/L	19225	3	13.554	152220	25.0
(DEL) Alkane: S...	16.273	1.8	ug/L	11160	3	13.554	152220	25.0