

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071322\
 Data File : VW023656.D
 Acq On : 13 Jul 2022 18:53
 Operator : SY/VA
 Sample : N3676-08RE
 Misc : 6.26g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 F5B07RE

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\SFAMWLM071322SMA.M
 Title : SFAM01.0

Signal : TIC: VW023656.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.233	48	54	55	rBV2	12972	24921	6.74%	1.009%
2	2.337	68	71	81	rVB	56351	87190	23.58%	3.530%
3	3.715	295	297	300	rBV4	932	1159	0.31%	0.047%
4	3.873	322	323	325	rBV2	1123	1056	0.29%	0.043%
5	4.007	334	345	356	rBV2	31615	92500	25.01%	3.745%
6	4.129	356	365	374	rBV	16920	45781	12.38%	1.854%
7	4.330	394	398	399	rBV3	1358	2042	0.55%	0.083%
8	4.367	399	404	407	rBV3	5028	10965	2.96%	0.444%
9	4.556	431	435	437	rBV5	1459	2382	0.64%	0.096%
10	4.580	437	439	442	rVV4	817	1007	0.27%	0.041%
11	4.605	442	443	446	rVB2	1204	952	0.26%	0.039%
12	4.794	472	474	478	rBV5	1606	1616	0.44%	0.065%
13	4.885	487	489	490	rBV2	3186	2615	0.71%	0.106%
14	5.111	523	526	527	rBV3	2144	2253	0.61%	0.091%
15	5.233	544	546	550	rVB4	829	1141	0.31%	0.046%
16	5.416	570	576	578	rBV6	1889	3878	1.05%	0.157%
17	5.440	578	580	583	rVV4	2335	2471	0.67%	0.100%
18	5.672	616	618	621	rVB4	1097	965	0.26%	0.039%
19	5.885	651	653	654	rBV2	1714	1406	0.38%	0.057%
20	6.019	674	675	678	rVB3	1878	1226	0.33%	0.050%
21	6.062	680	682	686	rBV5	1543	1550	0.42%	0.063%
22	6.190	701	703	707	rVB5	1246	1440	0.39%	0.058%
23	6.464	746	748	751	rVB4	927	1070	0.29%	0.043%
24	6.531	757	759	761	rBV3	912	1042	0.28%	0.042%
25	6.684	782	784	789	rVB6	1672	2386	0.65%	0.097%
26	6.751	789	795	798	rBV8	1418	2729	0.74%	0.110%
27	6.812	802	805	807	rBV4	1014	1199	0.32%	0.049%
28	7.086	842	850	854	rBV	14248	38132	10.31%	1.544%
29	7.391	896	900	902	rBV5	833	1054	0.28%	0.043%
30	7.415	902	904	909	rVB6	1331	1928	0.52%	0.078%
31	7.488	913	916	918	rBV3	820	1017	0.27%	0.041%
32	7.549	923	926	929	rBV5	1020	1294	0.35%	0.052%
33	7.647	933	942	951	rBV	55975	142139	38.43%	5.755%
34	7.744	956	958	961	rBV3	955	1180	0.32%	0.048%
35	7.860	976	977	979	rBV2	1000	946	0.26%	0.038%
36	8.269	1033	1044	1060	rBV3	108922	369834	100.00%	14.974%

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Integrator: RTE

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
 Title : SFAM01.0

37	8.488	1077	1080	1081	rBV3	1446	1539	0.42%	0.062%
38	8.695	1107	1114	1122	rVB9	8076	18437	4.99%	0.746%
39	8.842	1130	1138	1146	rBV	100468	201132	54.38%	8.143%
40	8.994	1161	1163	1165	rBV3	1324	990	0.27%	0.040%
41	9.165	1189	1191	1193	rBV3	1106	1127	0.30%	0.046%
42	9.275	1201	1209	1220	rBV	74314	157567	42.60%	6.379%
43	9.451	1236	1238	1240	rBV3	1509	1280	0.35%	0.052%
44	9.506	1245	1247	1249	rBV3	1143	1048	0.28%	0.042%
45	9.646	1266	1270	1272	rVB6	1089	1348	0.36%	0.055%
46	9.665	1272	1273	1275	rBV2	1129	987	0.27%	0.040%
47	10.037	1327	1334	1343	rVV	23050	44909	12.14%	1.818%
48	10.128	1347	1349	1352	rVB4	2467	1999	0.54%	0.081%
49	10.159	1352	1354	1357	rBV4	790	1098	0.30%	0.044%
50	10.323	1373	1381	1388	rBV	124790	217801	58.89%	8.818%
51	10.476	1405	1406	1409	rBV3	1018	1060	0.29%	0.043%
52	10.579	1418	1423	1429	rVV2	17219	31523	8.52%	1.276%
53	10.701	1438	1443	1451	rVB2	6282	13387	3.62%	0.542%
54	10.768	1451	1454	1459	rBV6	1232	1779	0.48%	0.072%
55	10.841	1462	1466	1467	rBV4	2065	2204	0.60%	0.089%
56	10.927	1474	1480	1489	rBV	49416	90569	24.49%	3.667%
57	11.061	1498	1502	1511	rVB3	12278	21820	5.90%	0.883%
58	11.433	1560	1563	1570	rVB7	4555	7856	2.12%	0.318%
59	11.628	1583	1595	1605	rBV	130639	235670	63.72%	9.542%
60	12.030	1657	1661	1663	rBV5	1485	1928	0.52%	0.078%
61	12.109	1669	1674	1676	rBV6	1206	2175	0.59%	0.088%
62	12.158	1680	1682	1685	rBV4	1072	1301	0.35%	0.053%
63	12.433	1722	1727	1729	rBV5	1408	1807	0.49%	0.073%
64	12.524	1740	1742	1744	rVB3	1146	1026	0.28%	0.042%
65	12.603	1749	1755	1760	rVB5	7559	14237	3.85%	0.576%
66	12.689	1764	1769	1777	rVV	78777	126939	34.32%	5.139%
67	12.926	1804	1808	1813	rVB8	4510	7732	2.09%	0.313%
68	13.115	1836	1839	1840	rBV2	869	934	0.25%	0.038%
69	13.158	1845	1846	1848	rVV2	1457	1485	0.40%	0.060%
70	13.182	1848	1850	1855	rVB6	3399	4971	1.34%	0.201%
71	13.256	1860	1862	1865	rBV4	1269	985	0.27%	0.040%
72	13.298	1865	1869	1875	rVB7	3323	6618	1.79%	0.268%
73	13.402	1878	1886	1894	rBV7	6179	14223	3.85%	0.576%
74	13.463	1894	1896	1898	rBV3	1702	1182	0.32%	0.048%
75	13.487	1898	1900	1901	rBV2	1526	1220	0.33%	0.049%
76	13.554	1905	1911	1917	rBV	91858	154534	41.78%	6.257%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
 Title : SFAM01.0

77	13.762	1941	1945	1949	rBV6	4297	8564	2.32%	0.347%
78	13.847	1952	1959	1970	rVV	93543	163993	44.34%	6.640%
79	13.944	1974	1975	1984	rVB9	1884	3129	0.85%	0.127%
80	14.060	1992	1994	1998	rBV5	2639	4804	1.30%	0.195%
81	14.426	2051	2054	2058	rBV5	916	1561	0.42%	0.063%
82	14.463	2058	2060	2062	rBV3	1217	1043	0.28%	0.042%
83	14.542	2071	2073	2076	rBV4	1226	1545	0.42%	0.063%
84	14.719	2098	2102	2107	rVB7	2638	4216	1.14%	0.171%
85	14.828	2118	2120	2124	rVB5	1194	1285	0.35%	0.052%
86	15.085	2160	2162	2165	rVV4	1091	982	0.27%	0.040%
87	15.371	2208	2209	2213	rBV5	1032	1464	0.40%	0.059%
88	15.402	2213	2214	2216	rVV2	1616	1219	0.33%	0.049%
89	15.438	2216	2220	2223	rVB5	2154	3216	0.87%	0.130%
90	15.475	2223	2226	2230	rBV6	1357	2043	0.55%	0.083%
91	15.536	2233	2236	2237	rBV3	956	1085	0.29%	0.044%
92	15.761	2271	2273	2274	rBV2	1181	1150	0.31%	0.047%
93	15.847	2286	2287	2292	rVB5	776	1227	0.33%	0.050%
94	15.938	2300	2302	2304	rVB3	1701	1285	0.35%	0.052%
95	15.975	2304	2308	2313	rBV7	1437	3023	0.82%	0.122%
96	16.279	2354	2358	2359	rBV4	998	1057	0.29%	0.043%
97	16.328	2363	2366	2368	rVB3	1125	992	0.27%	0.040%
98	16.353	2368	2370	2373	rVB4	935	987	0.27%	0.040%
99	16.383	2373	2375	2378	rBV5	1148	1500	0.41%	0.061%
100	16.468	2387	2389	2393	rVB5	1358	1651	0.45%	0.067%

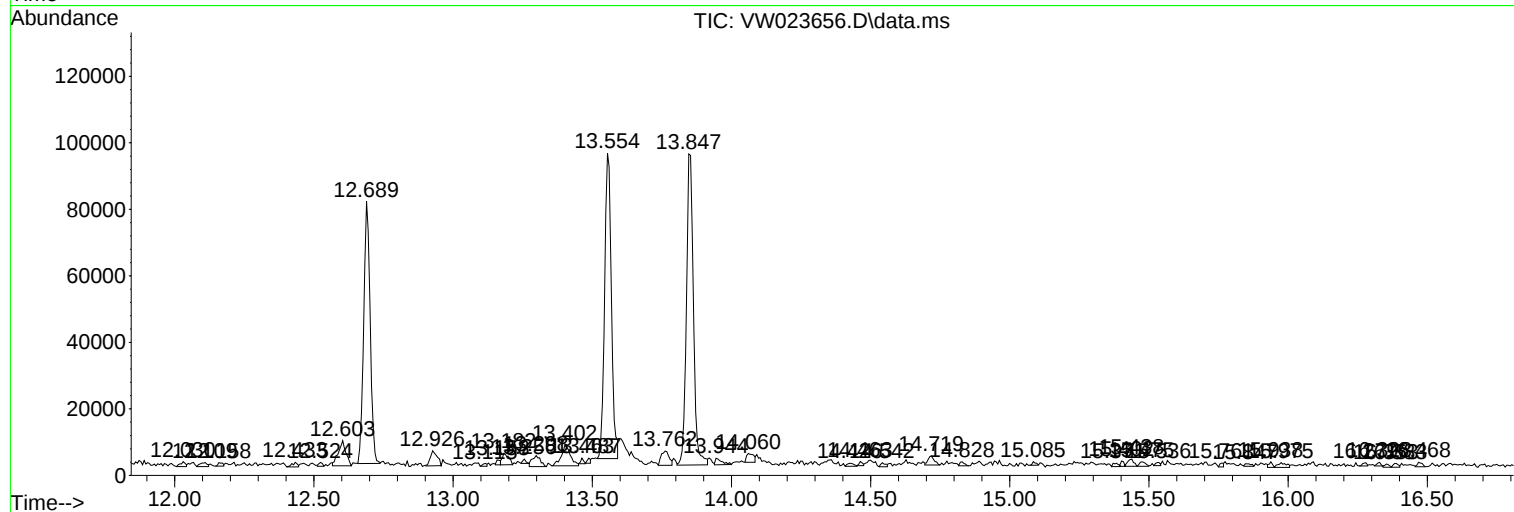
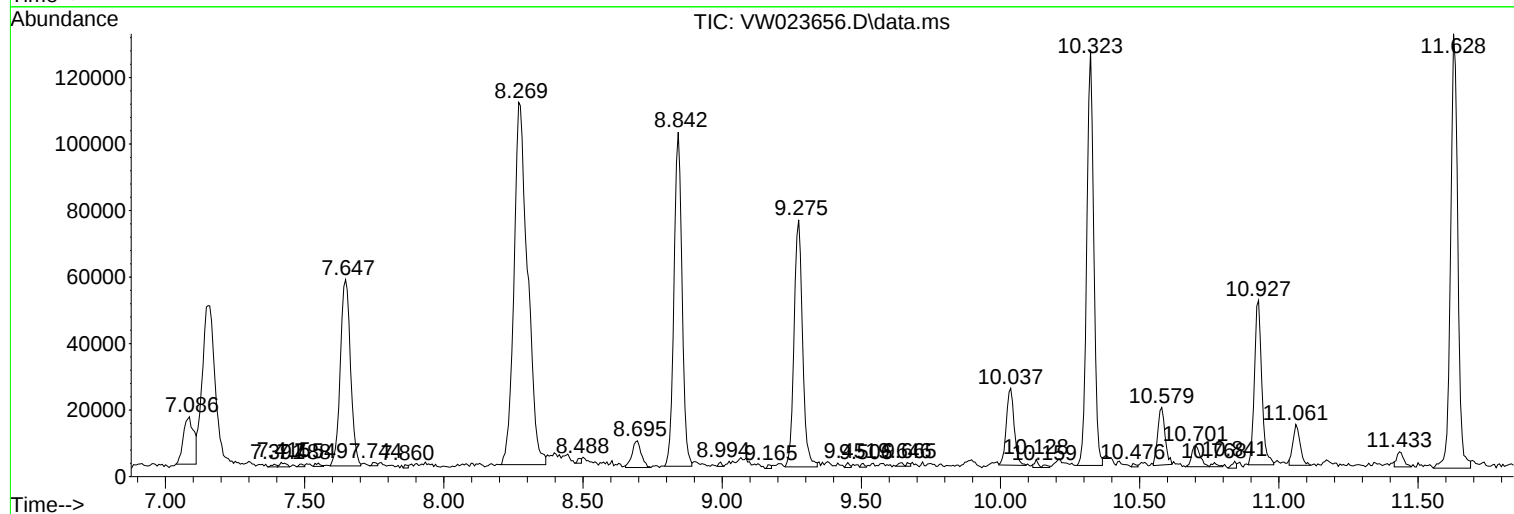
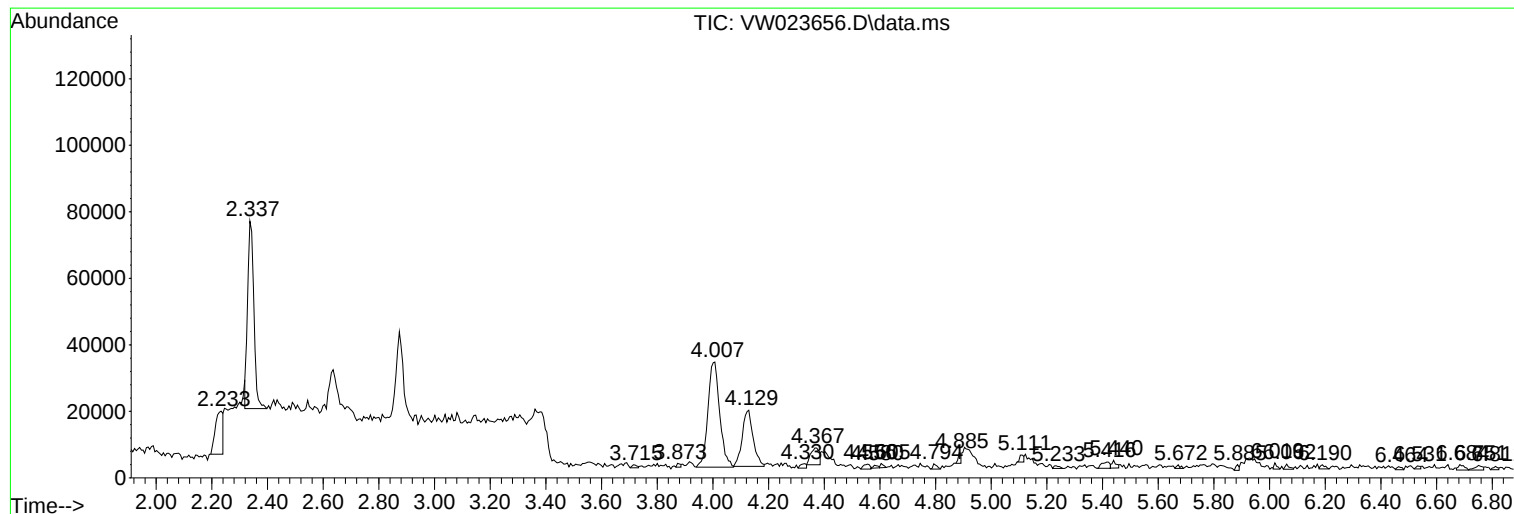
Sum of corrected areas: 2469914

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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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



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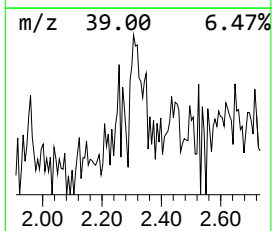
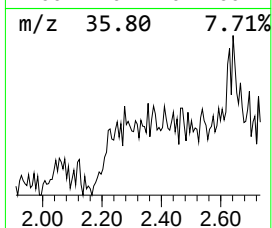
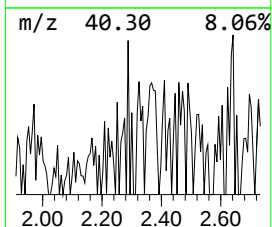
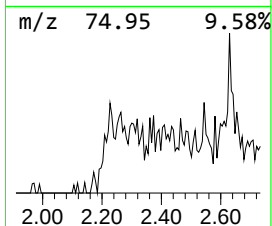
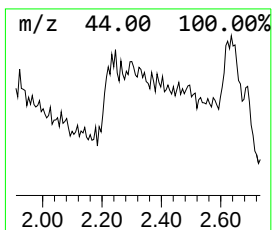
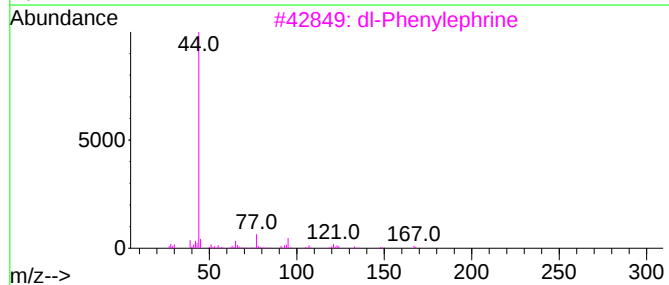
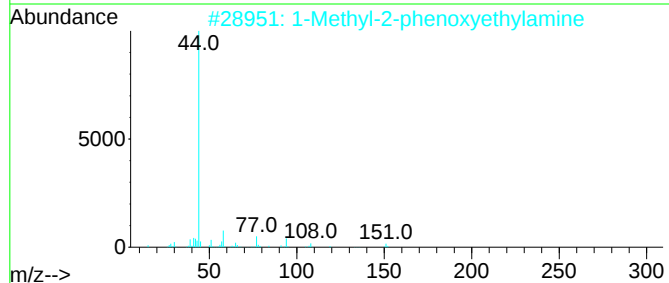
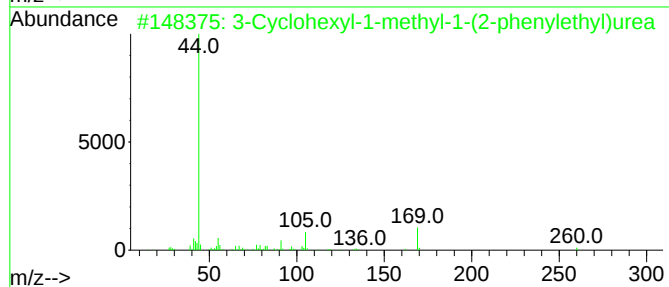
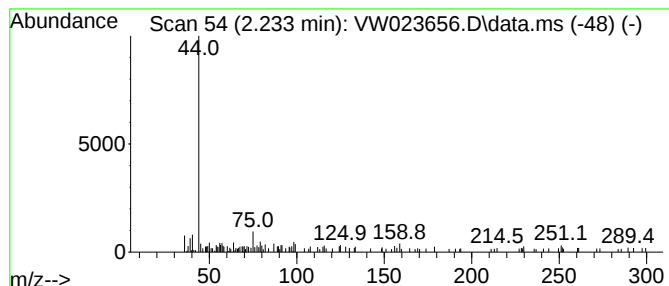
Instrument :
MSVOA_W
ClientSampleId :
F5B07RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 3-Cyclohexyl-1-methyl-1-(2-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.233	3.10 ug/L	24921	1,4-Difluorobenzene	8.842	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Cyclohexyl-1-methyl-1-(2-pheny...	260	C16H24N2O	1024470-80-1	64
2	1-Methyl-2-phenoxyethylamine	151	C9H13NO	035205-54-0	59
3	dl-Phenylephrine	167	C9H13NO2	001477-63-0	59
4	Ortetamine	149	C10H15N	005580-32-5	59
5	1-Guanidinosuccinimide	141	C5H7N3O2	1000130-20-8	59



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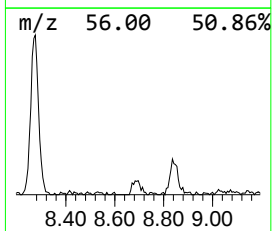
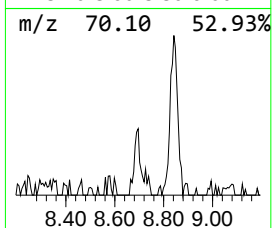
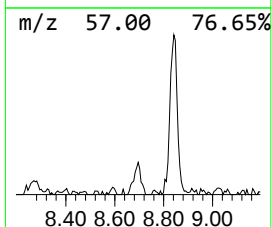
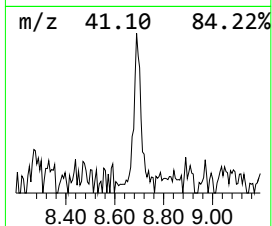
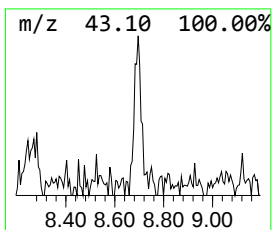
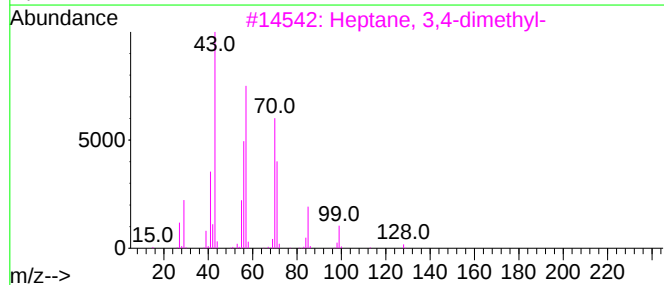
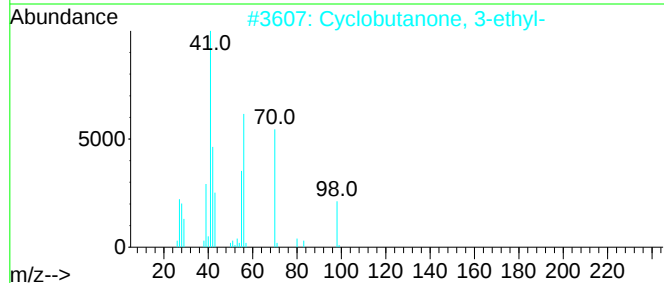
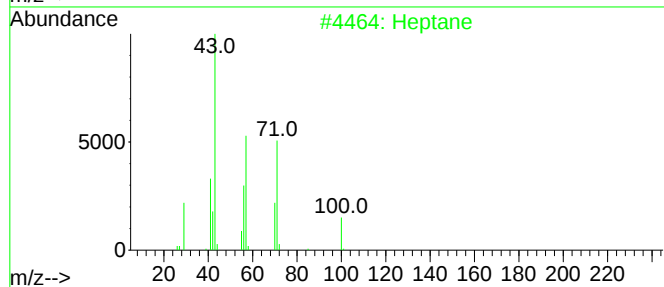
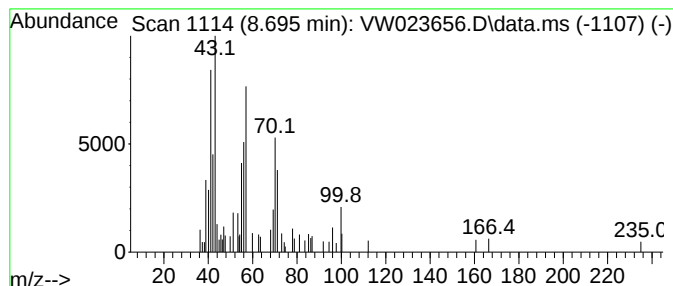
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.695	2.29 ug/L	18437	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	47
2			Cyclobutanone, 3-ethyl-	98	C6H10O	056335-73-0	43
3			Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	43
4			Pentaerythritol	136	C5H12O4	000115-77-5	43
5			Pentaerythritol	136	C5H12O4	000115-77-5	38



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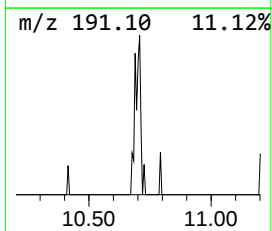
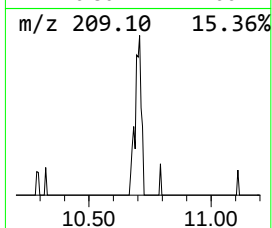
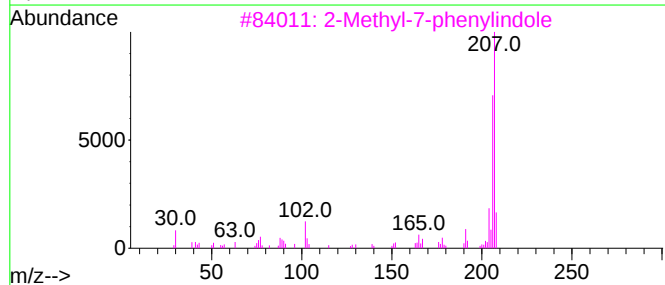
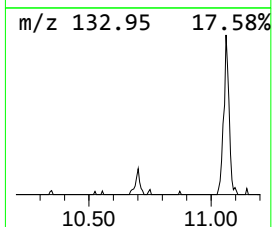
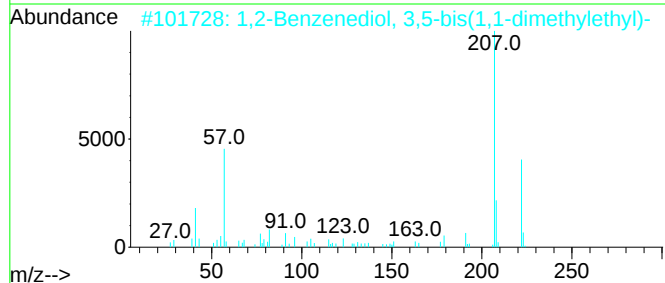
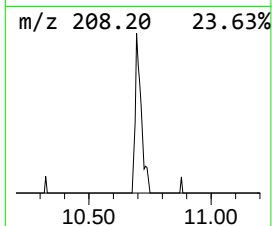
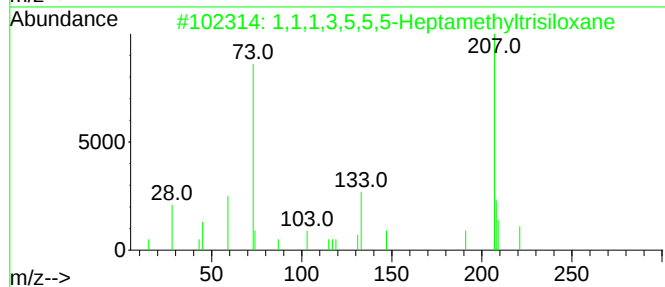
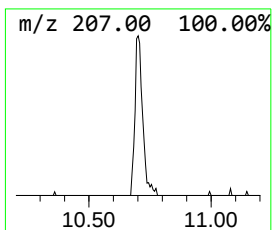
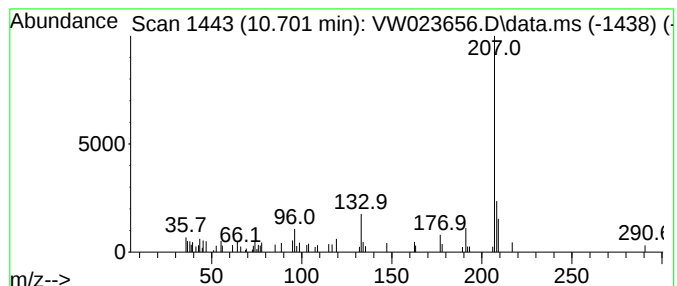
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Quant Title : SFAM01.0

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

Peak Number 4 1,1,1,3,5,5,5-Heptamethyltr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.701	1.42 ug/L	13387	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,1,3,5,5,5-Heptamethyltrisilo...	222	C7H22O2Si3	001873-88-7	64		
2	1,2-Benzenediol, 3,5-bis(1,1-dim...	222	C14H22O2	001020-31-1	59		
3	2-Methyl-7-phenylindole	207	C15H13N	001140-08-5	52		
4	Thymol, TMS derivative	222	C13H22OSi	055012-80-1	50		
5	Arsenous acid, tris(trimethylsil...	342	C9H27AsO3Si3	055429-29-3	50		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071322\
Data File : VW023656.D
Acq On : 13 Jul 2022 18:53
Operator : SY/VA
Sample : N3676-08RE
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5B07RE

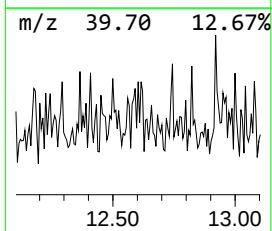
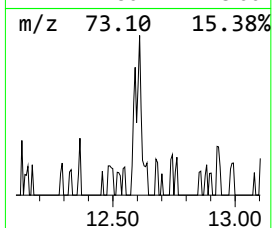
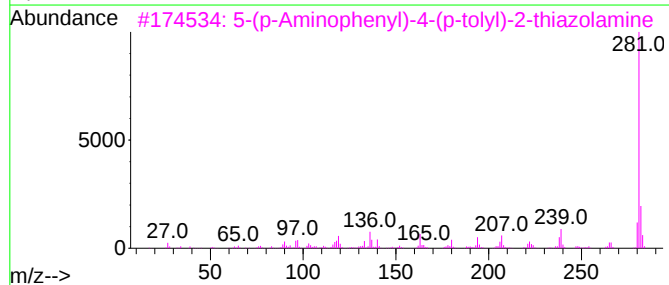
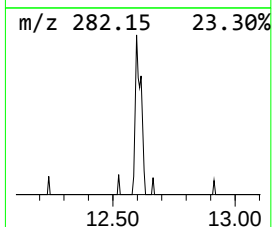
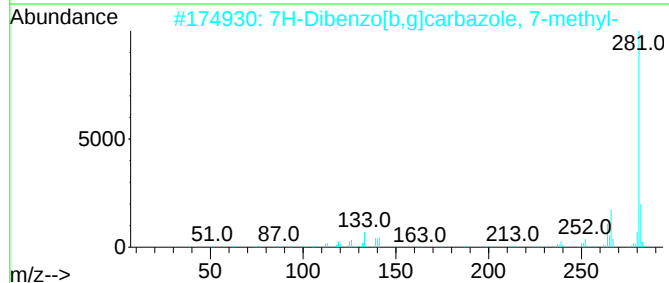
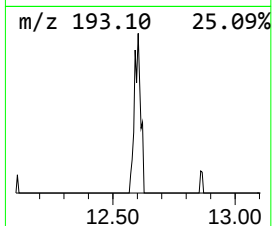
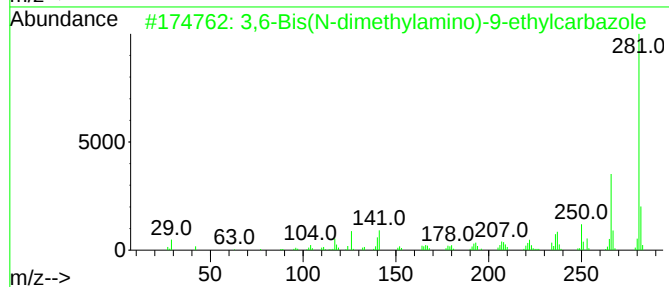
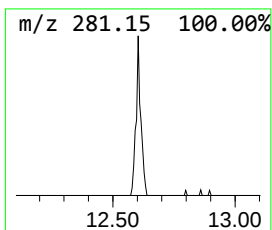
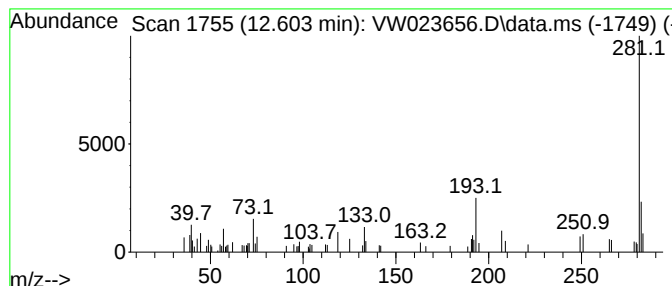
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Quant Title : SFAM01.0

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

Peak Number 6 3,6-Bis(N-dimethylamino)-9-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.603	2.30 ug/L	14237	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,6-Bis(N-dimethylamino)-9-ethyl...	281	C18H23N3	057103-04-5	53
2			7H-Dibenzo[b,g]carbazole, 7-methyl-	281	C21H15N	003557-49-1	53
3			5-(p-Aminophenyl)-4-(p-tolyl)-2-...	281	C16H15N3S	094460-46-5	53
4			3-Phenyl oxindole, BSTFA derivative	281	C17H19NOSi	1000458-51-5	50
5			2-(Benzthiazol-2-yl)-6-methoxybe...	281	C16H11NO2S	1000284-45-0	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071322\
Data File : VW023656.D
Acq On : 13 Jul 2022 18:53
Operator : SY/VA
Sample : N3676-08RE
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5B07RE

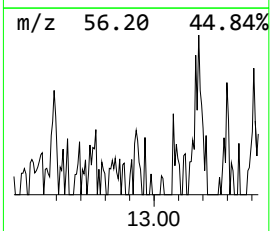
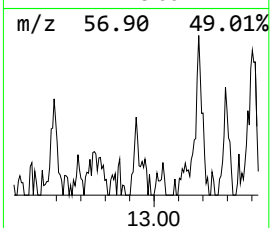
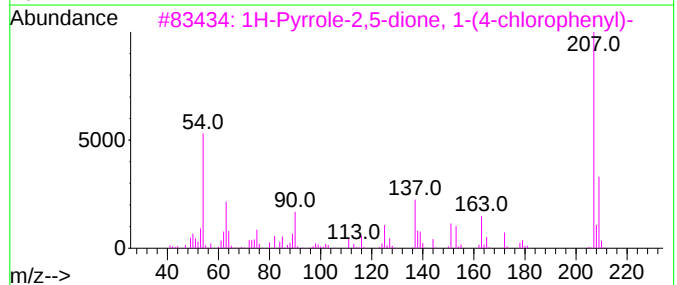
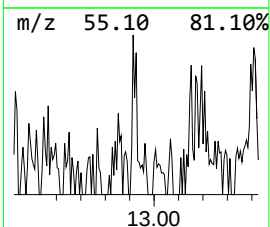
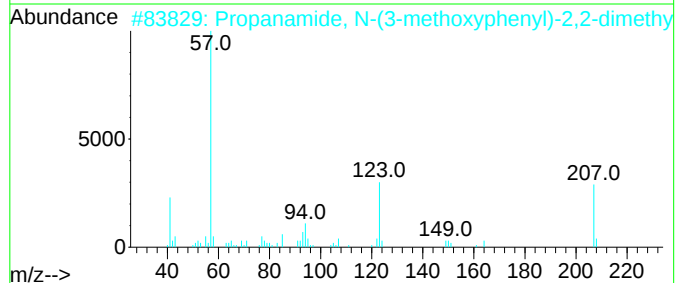
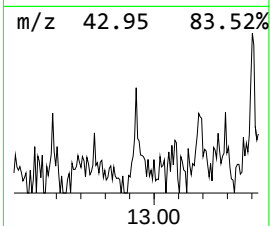
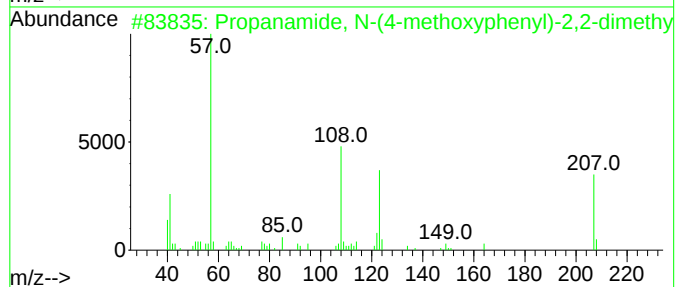
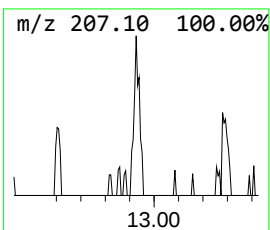
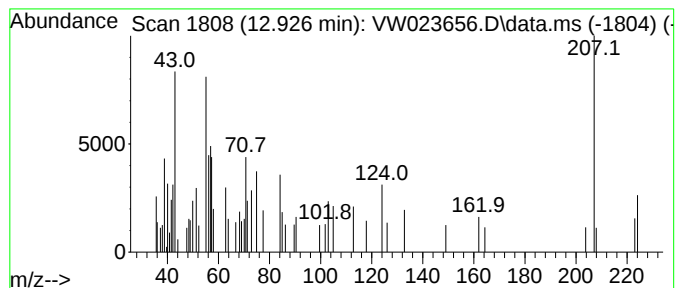
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.926	1.25 ug/L	7732	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanamide, N-(4-methoxyphenyl)...	207	C12H17NO2	056619-94-4	38
2			Propanamide, N-(3-methoxyphenyl)...	207	C12H17NO2	056619-93-3	35
3			1H-Pyrrole-2,5-dione, 1-(4-chlor...	207	C10H6ClNO2	001631-29-4	35
4			Benzofuran-2-one, 2,3-dihydro-3,...	207	C10H9NO4	1000129-53-4	35
5			N-Methyl-1-adamantaneacetamide	207	C13H21NO	031897-93-5	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071322\
Data File : VW023656.D
Acq On : 13 Jul 2022 18:53
Operator : SY/VA
Sample : N3676-08RE
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

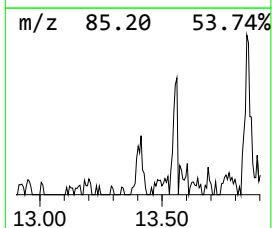
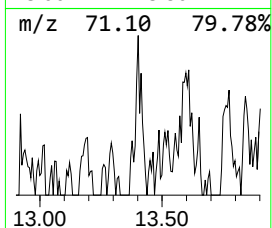
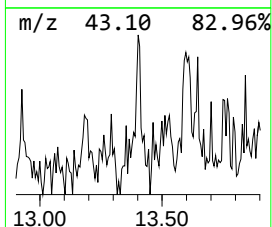
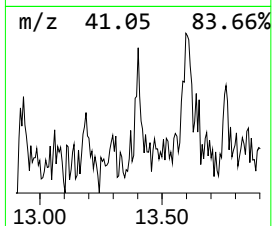
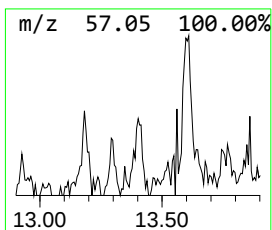
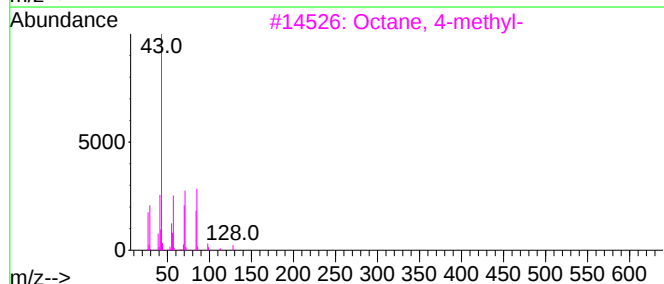
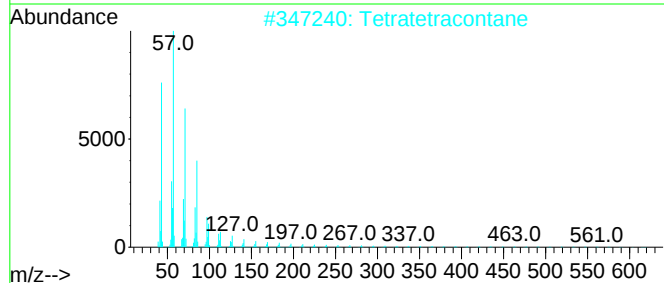
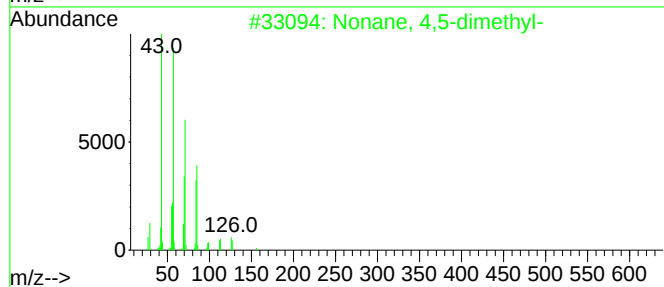
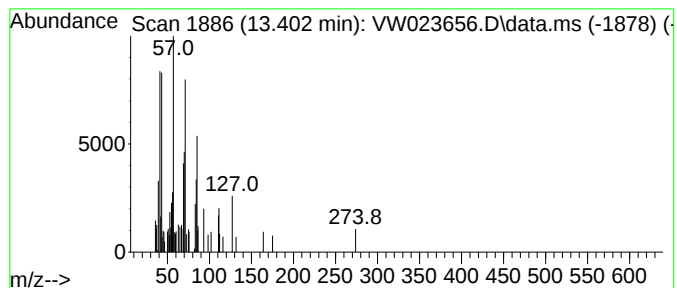
Instrument :
MSVOA_W
ClientSampleId :
F5B07RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.M
Quant Title : SFAM01.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.402	2.30 ug/L	14223	1,4-Dichlorobenzene-d4		13.560	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 4,5-dimethyl-		156	C11H24	017302-23-7	43
2	Tetratetracontane		619	C44H90	007098-22-8	43
3	Octane, 4-methyl-		128	C9H20	002216-34-4	38
4	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	38
5	1-Heptanol, 2-propyl-		158	C10H22O	010042-59-8	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071322\
Data File : VW023656.D
Acq On : 13 Jul 2022 18:53
Operator : SY/VA
Sample : N3676-08RE
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5B07RE

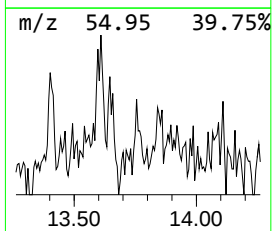
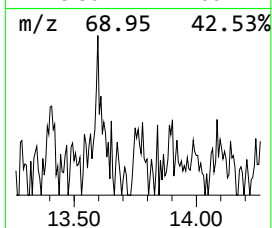
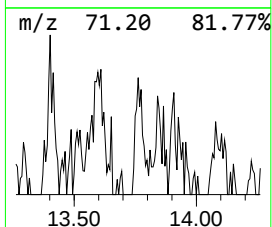
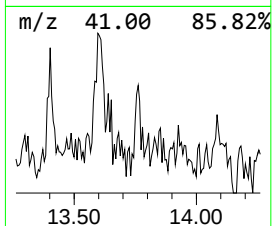
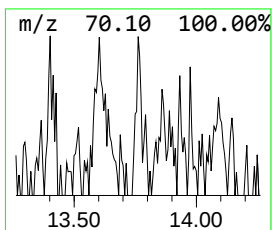
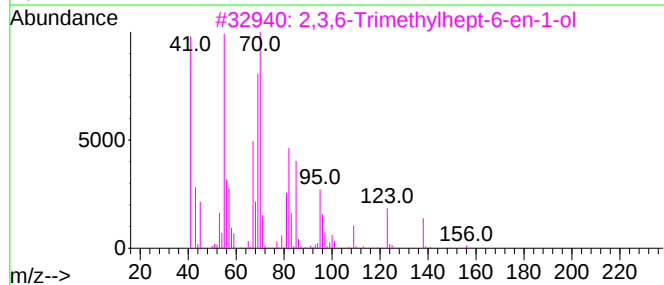
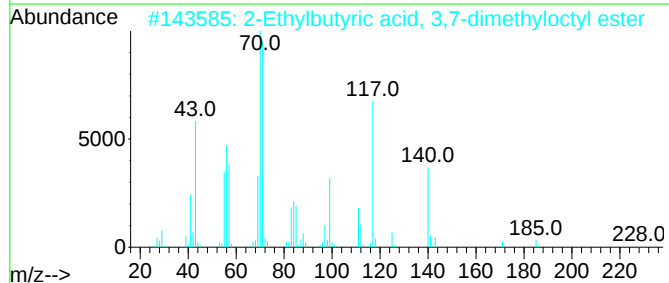
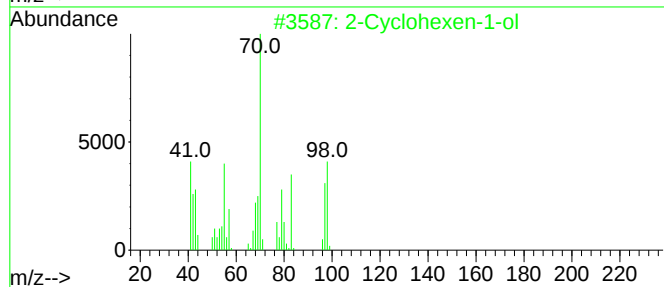
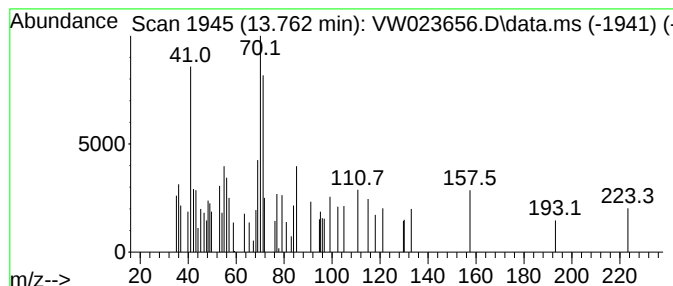
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.762	1.39 ug/L	8564	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-ol	98	C6H10O	000822-67-3	10
2			2-Ethylbutyric acid, 3,7-dimethy...	256	C16H32O2	1000369-41-9	10
3			2,3,6-Trimethylhept-6-en-1-ol	156	C10H20O	1000111-57-0	10
4			2-Propen-1-amine, N-ethyl-	85	C5H11N	002424-02-4	9
5			Maleamic acid	115	C4H5NO3	000557-24-4	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071322\
Data File : VW023656.D
Acq On : 13 Jul 2022 18:53
Operator : SY/VA
Sample : N3676-08RE
Misc : 6.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
F5B07RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM071322SMA.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
3-Cyclohexyl-1-...	2.233	3.1	ug/L	24921	1	8.842	201132	25.0
(DEL) Alkane: S...	8.695	2.3	ug/L	18437	1	8.842	201132	25.0
1,1,1,3,5,5,5-H...	10.701	1.4	ug/L	13387	2	11.634	235670	25.0
3,6-Bis(N-dimet...	12.603	2.3	ug/L	14237	3	13.560	154534	25.0
unknown-01	12.926	1.3	ug/L	7732	3	13.560	154534	25.0
(DEL) Alkane: S...	13.402	2.3	ug/L	14223	3	13.560	154534	25.0
unknown-02	13.762	1.4	ug/L	8564	3	13.560	154534	25.0