

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW091923\
 Data File : VW027144.D
 Acq On : 19 Sep 2023 12:21
 Operator : SY/MD
 Sample : 04472-04RE
 Misc : 3.67g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EZYK8RE

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\SFAMWLM091823SMA.M
 Title : SFAM01.0

Signal : TIC: VW027144.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.357	72	77	86	rVB	34069	61542	12.11%	2.623%
2	2.820	149	153	154	rBV3	17019	22839	4.49%	0.974%
3	3.277	225	228	229	rBV3	1677	1762	0.35%	0.075%
4	3.667	291	292	296	rBV4	1317	1477	0.29%	0.063%
5	3.832	317	319	321	rBV3	1199	783	0.15%	0.033%
6	3.936	334	336	337	rBV2	1567	1173	0.23%	0.050%
7	4.021	341	350	361	rVV	56248	150903	29.70%	6.432%
8	4.131	366	368	373	rVV4	1758	2868	0.56%	0.122%
9	4.173	373	375	377	rVV3	1302	1324	0.26%	0.056%
10	4.234	380	385	387	rBV6	1053	2027	0.40%	0.086%
11	4.332	398	401	402	rBV3	1265	834	0.16%	0.036%
12	4.368	404	407	408	rBV3	1596	1863	0.37%	0.079%
13	4.533	432	434	436	rVB3	1730	1273	0.25%	0.054%
14	4.685	456	459	461	rVB3	1028	1072	0.21%	0.046%
15	4.740	466	468	471	rVB4	818	922	0.18%	0.039%
16	4.771	471	473	476	rBV3	1897	1915	0.38%	0.082%
17	4.862	485	488	489	rBV2	1026	1176	0.23%	0.050%
18	4.917	489	497	498	rBV3	11137	22991	4.52%	0.980%
19	5.063	519	521	525	rVB5	1214	995	0.20%	0.042%
20	5.173	537	539	541	rBV3	1646	1528	0.30%	0.065%
21	5.307	558	561	562	rBV2	1067	1214	0.24%	0.052%
22	5.447	582	584	585	rBV2	1459	906	0.18%	0.039%
23	5.490	590	591	594	rBV3	1407	1432	0.28%	0.061%
24	5.594	606	608	610	rBV3	851	753	0.15%	0.032%
25	5.655	617	618	621	rBV2	971	1026	0.20%	0.044%
26	5.704	624	626	628	rBV3	1369	871	0.17%	0.037%
27	5.722	628	629	631	rBV2	1164	957	0.19%	0.041%
28	5.752	631	634	636	rBV3	1779	1817	0.36%	0.077%
29	5.905	655	659	660	rBV4	1352	1722	0.34%	0.073%
30	5.929	660	663	664	rBV3	1996	2127	0.42%	0.091%
31	6.149	698	699	702	rBV2	1379	1367	0.27%	0.058%
32	6.240	712	714	716	rVV3	1181	873	0.17%	0.037%
33	6.283	718	721	722	rVB3	945	839	0.17%	0.036%
34	6.392	737	739	742	rVB4	1595	1671	0.33%	0.071%
35	6.423	742	744	745	rBV2	1567	1097	0.22%	0.047%
36	6.441	745	747	750	rVB3	1181	1111	0.22%	0.047%

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Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM091823SMA.M
 Title : SFAM01.0

37	6.496	754	756	758	rVB3	1320	1287	0.25%	0.055%
38	6.551	764	765	768	rVB3	1301	941	0.19%	0.040%
39	6.697	786	789	790	rBV3	926	1073	0.21%	0.046%
40	6.770	799	801	804	rVB4	928	880	0.17%	0.038%
41	6.850	810	814	817	rBV6	1487	2071	0.41%	0.088%
42	6.898	820	822	824	rBV3	954	942	0.19%	0.040%
43	6.929	824	827	829	rBV3	1203	1558	0.31%	0.066%
44	7.081	846	852	858	rBV2	12151	32921	6.48%	1.403%
45	7.283	882	885	886	rBV2	960	783	0.15%	0.033%
46	7.343	891	895	898	rVB5	1481	1754	0.35%	0.075%
47	7.411	903	906	910	rBV6	1422	1900	0.37%	0.081%
48	7.514	922	923	926	rVB3	1291	1060	0.21%	0.045%
49	7.545	926	928	930	rBV3	1153	1418	0.28%	0.060%
50	7.648	937	945	958	rVV2	98277	246713	48.55%	10.516%
51	7.764	962	964	967	rBV4	1067	969	0.19%	0.041%
52	7.910	985	988	989	rBV3	1911	2224	0.44%	0.095%
53	8.026	1005	1007	1009	rBV2	1750	1109	0.22%	0.047%
54	8.136	1022	1025	1026	rBV3	1086	975	0.19%	0.042%
55	8.276	1039	1048	1063	rBV2	173482	508155	100.00%	21.660%
56	8.532	1087	1090	1093	rBV5	1596	2250	0.44%	0.096%
57	8.679	1112	1114	1116	rBV3	2258	2648	0.52%	0.113%
58	8.843	1133	1141	1151	rVV	122351	250678	49.33%	10.685%
59	9.087	1176	1181	1188	rBV	2976	6864	1.35%	0.293%
60	9.276	1203	1212	1220	rBV	125325	268540	52.85%	11.447%
61	9.392	1228	1231	1234	rBV5	1194	1854	0.36%	0.079%
62	9.575	1258	1261	1262	rBV3	1249	1429	0.28%	0.061%
63	9.642	1267	1272	1273	rBV4	1622	1742	0.34%	0.074%
64	9.745	1287	1289	1291	rBV2	1325	1180	0.23%	0.050%
65	9.800	1295	1298	1301	rBV5	1877	2273	0.45%	0.097%
66	9.849	1304	1306	1309	rBV4	1359	1489	0.29%	0.063%
67	9.886	1310	1312	1313	rBV2	1512	847	0.17%	0.036%
68	9.971	1323	1326	1328	rBV3	1725	1556	0.31%	0.066%
69	10.032	1331	1336	1342	rVV2	28816	52156	10.26%	2.223%
70	10.123	1349	1351	1353	rBV	1730	1060	0.21%	0.045%
71	10.270	1370	1375	1376	rBV5	1251	1764	0.35%	0.075%
72	10.325	1376	1384	1390	rBV	148079	263862	51.93%	11.247%
73	10.483	1408	1410	1412	rBV3	1670	1414	0.28%	0.060%
74	10.508	1412	1414	1418	rVB5	2213	2615	0.51%	0.111%
75	10.581	1419	1426	1428	rBV	13914	22308	4.39%	0.951%
76	11.288	1537	1542	1548	rBV10	17854	41682	8.20%	1.777%

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 Title : SFAM01.0

77	12.690	1768	1772	1781	rVB	98363	162554	31.99%	6.929%
78	13.214	1856	1858	1861	rBV4	11685	9640	1.90%	0.411%
79	13.355	1880	1881	1883	rVB2	1495	811	0.16%	0.035%
80	13.397	1883	1888	1889	rBV4	1752	2453	0.48%	0.105%
81	13.434	1891	1894	1896	rBV4	1064	1319	0.26%	0.056%
82	13.507	1903	1906	1908	rBV4	2429	2314	0.46%	0.099%
83	13.556	1909	1914	1920	rVV	24213	40663	8.00%	1.733%
84	13.769	1946	1949	1954	rVB6	1399	2265	0.45%	0.097%
85	13.848	1955	1962	1969	rBV2	34221	64594	12.71%	2.753%
86	13.983	1982	1984	1985	rBV2	1841	1870	0.37%	0.080%
87	14.263	2029	2030	2031	rBV	1630	791	0.16%	0.034%
88	14.452	2059	2061	2064	rVB4	1027	952	0.19%	0.041%
89	14.629	2088	2090	2091	rBV2	1377	1226	0.24%	0.052%
90	14.714	2102	2104	2111	rVB7	2786	3695	0.73%	0.157%
91	14.860	2126	2128	2131	rBV4	1181	941	0.19%	0.040%
92	14.958	2142	2144	2147	rVB4	1626	1698	0.33%	0.072%
93	14.988	2147	2149	2151	rBV3	934	753	0.15%	0.032%
94	15.013	2151	2153	2155	rBV3	1565	1110	0.22%	0.047%
95	15.129	2171	2172	2175	rBV3	1460	1338	0.26%	0.057%
96	15.190	2180	2182	2183	rBV2	1342	816	0.16%	0.035%
97	15.525	2235	2237	2238	rBV2	1874	1368	0.27%	0.058%
98	15.921	2300	2302	2303	rBV2	1454	772	0.15%	0.033%
99	16.086	2328	2329	2331	rBV2	1246	784	0.15%	0.033%
100	16.269	2357	2359	2361	rBV3	1604	1315	0.26%	0.056%

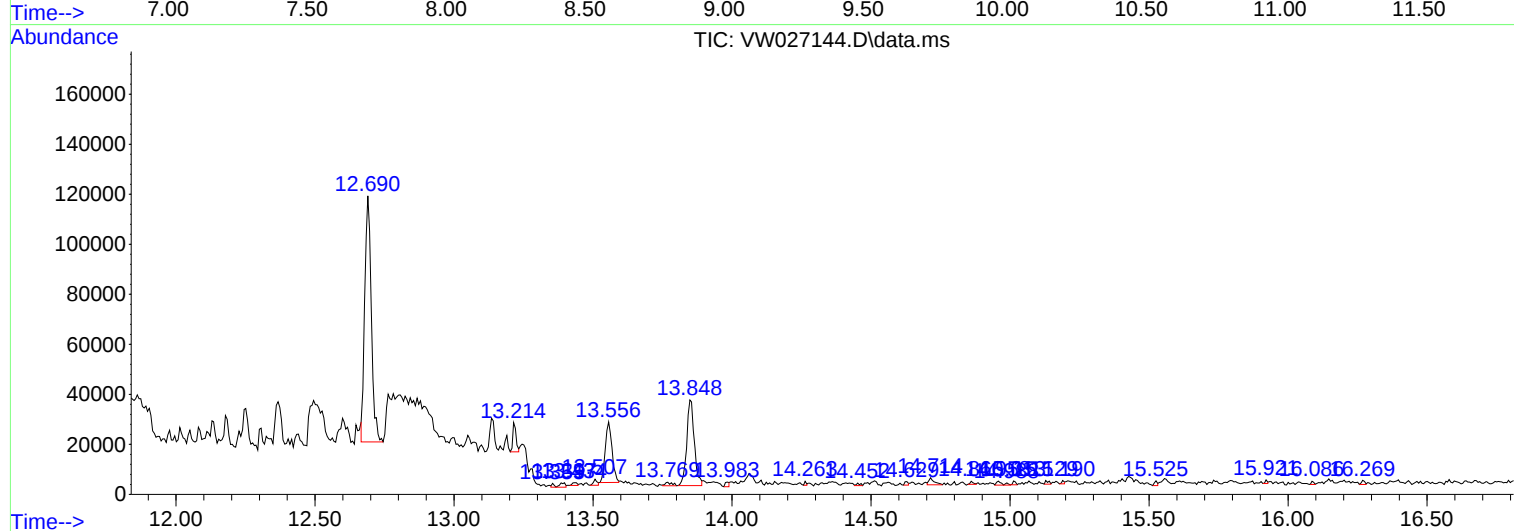
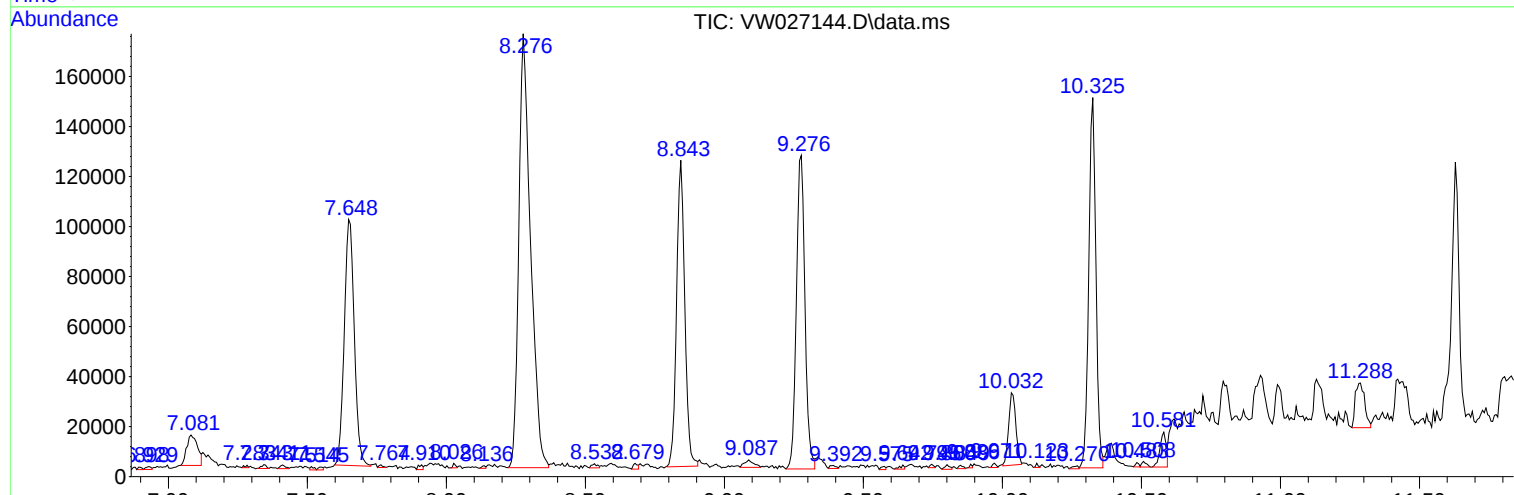
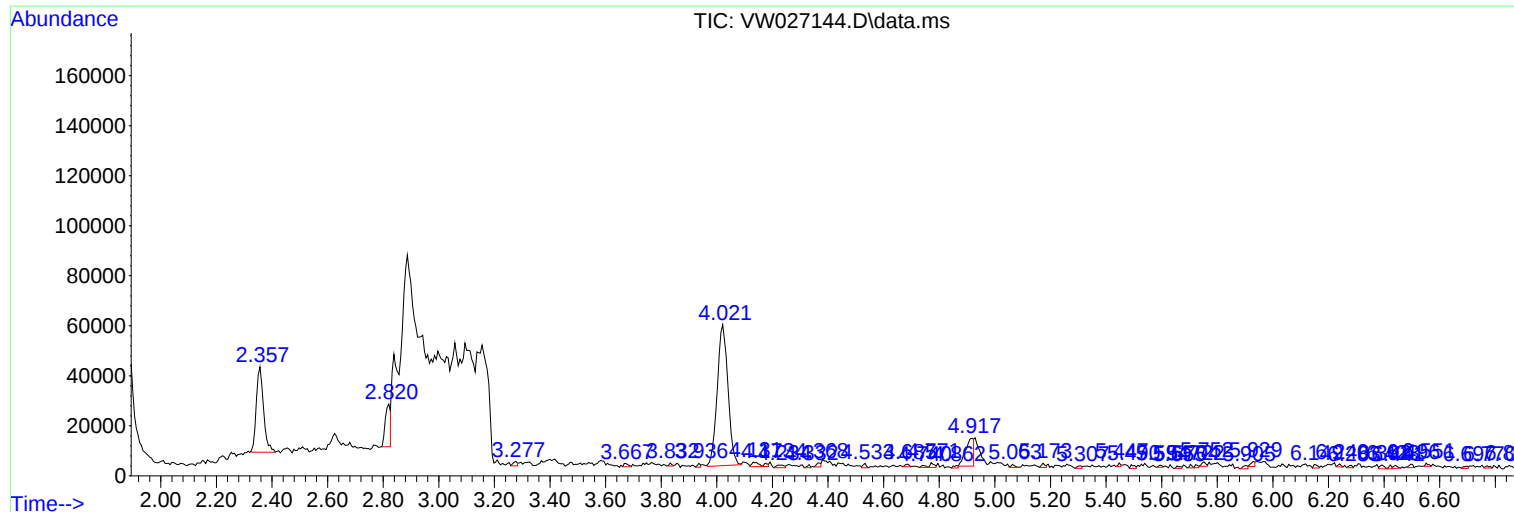
Sum of corrected areas: 2346036

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM091823SMA.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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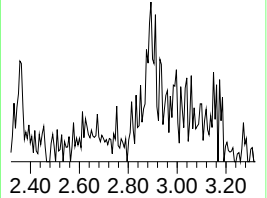
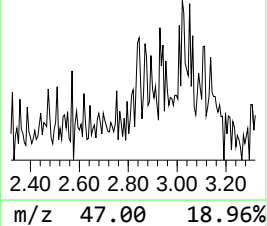
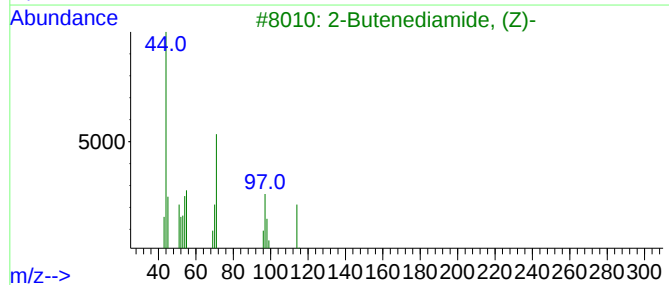
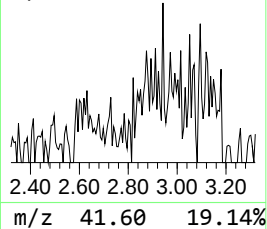
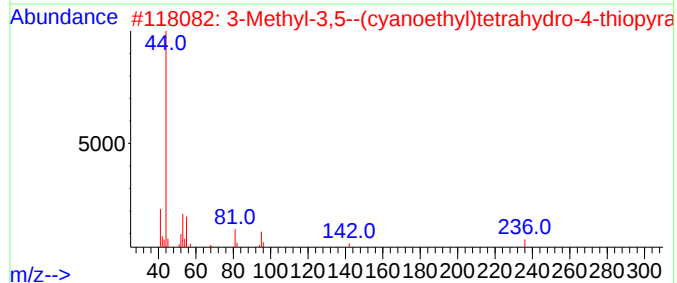
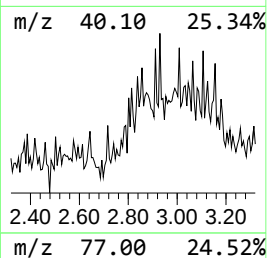
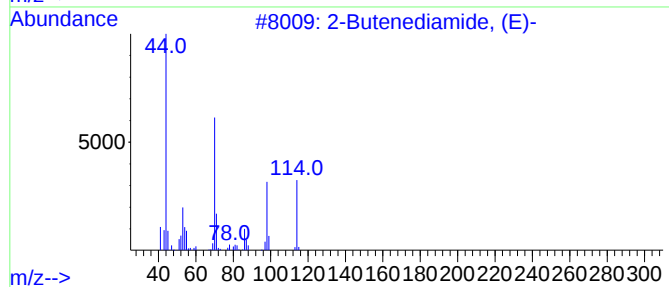
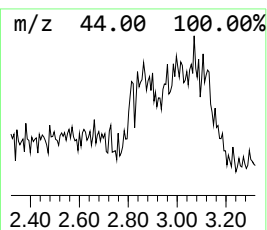
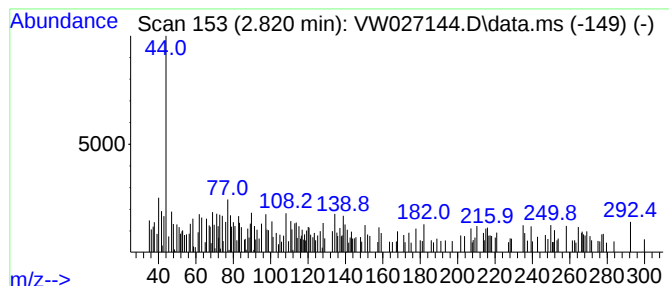
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM091823SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.820	2.28 ug/L	22839	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butenediamide, (E)-		114	C4H6N2O2		000627-64-5	38
2	3-Methyl-3,5--(cyanoethyl)tetrah...		236	C12H16N2OS		1000140-32-3	38
3	2-Butenediamide, (Z)-		114	C4H6N2O2		000928-01-8	38
4	Benzeneethanamine, 5-benzyloxy-3...		275	C16H18FN02		1000116-56-1	27
5	1-[.alpha.-(1-Adamantyl)benzylid...		313	C18H23N3S		1000222-83-3	27



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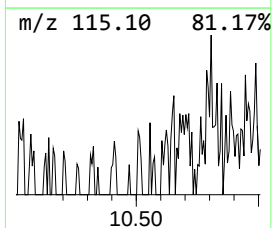
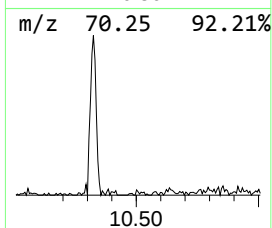
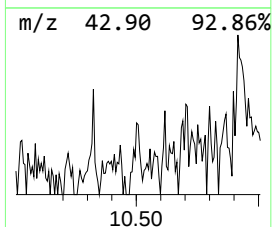
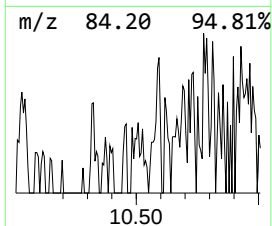
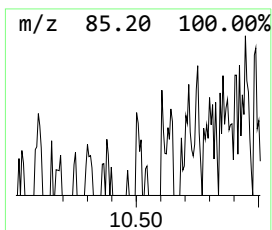
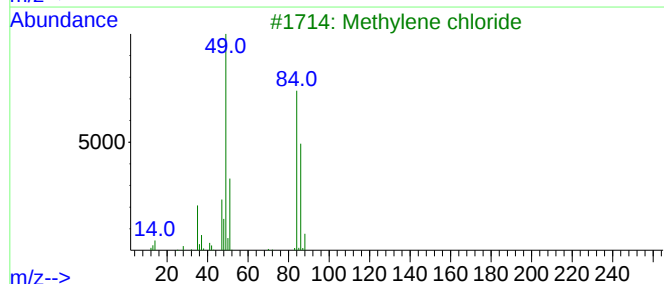
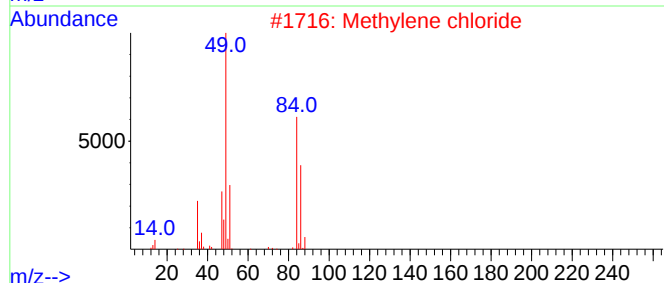
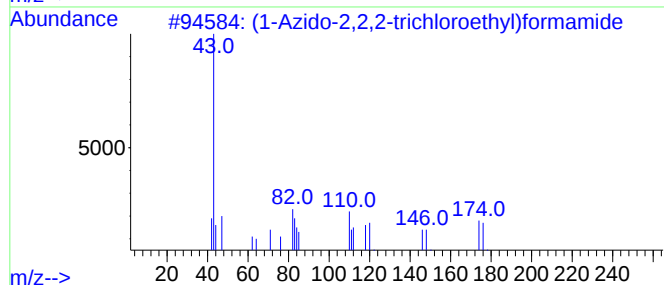
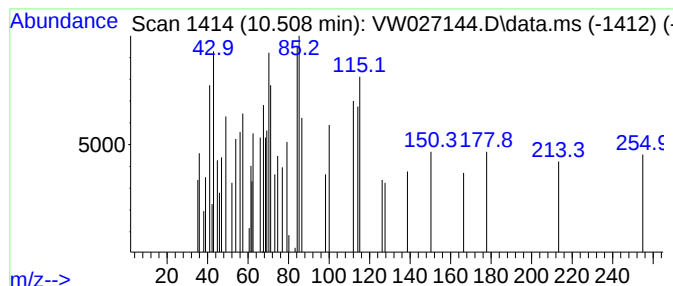
Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM091823SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.508	1.57 ug/L	2615	Chlorobenzene-d5		11.629	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	(1-Azido-2,2,2-trichloroethyl)fo...		216	C3H3Cl3N4O	035530-17-7	23
2	Methylene chloride		84	CH2Cl2	000075-09-2	18
3	Methylene chloride		84	CH2Cl2	000075-09-2	14
4	Dimethyl sulfide		62	C2H6S	000075-18-3	12
5	Heptane, 4-methyl-		114	C8H18	000589-53-7	9



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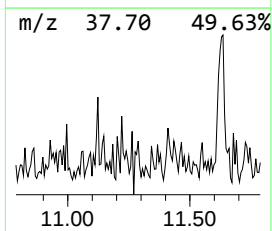
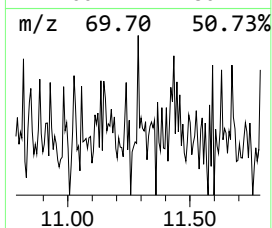
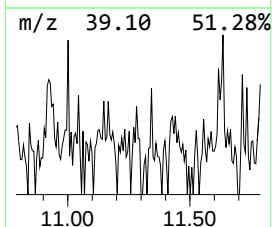
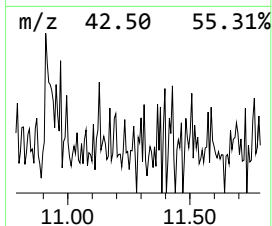
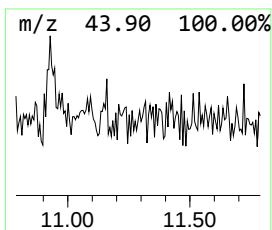
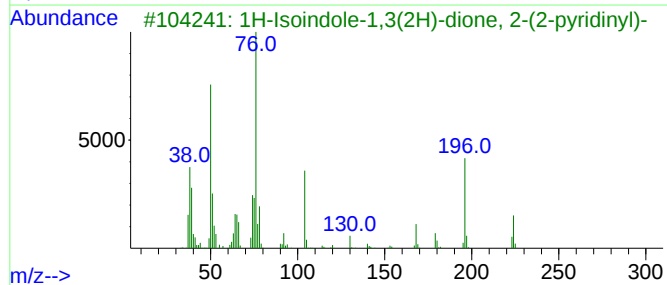
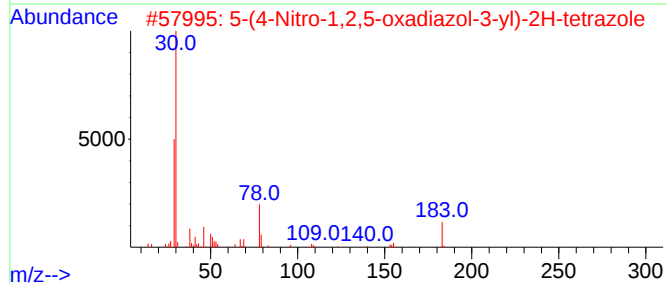
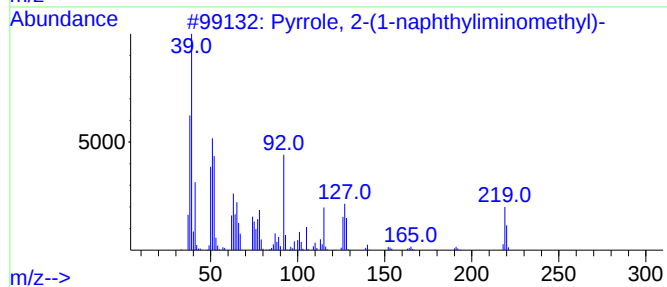
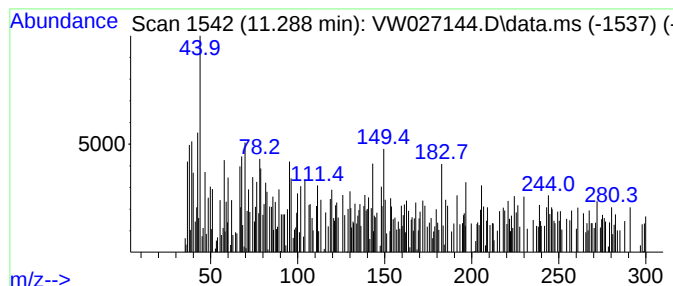
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Quant Title : SFAM01.0

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

Peak Number 4 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.288	25.00 ug/L	41682	Chlorobenzene-d5	11.629

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrrole, 2-(1-naphthyliminomethyl)-	220	C15H12N2	212759-16-5	47
2			5-(4-Nitro-1,2,5-oxadiazol-3-yl)...	183	C3HN7O3	355142-29-9	38
3			1H-Isoindole-1,3(2H)-dione, 2-(2...	224	C13H8N2O2	059208-49-0	35
4			Benzenesulfonamide, N-methyl-3-n...	216	C7H8N2O4S	058955-78-5	25
5			2,6-Dinitrobenzonitrile	193	C7H3N3O4	035213-00-4	16



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW091923\
Data File : VW027144.D
Acq On : 19 Sep 2023 12:21
Operator : SY/MD
Sample : 04472-04RE
Misc : 3.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYK8RE

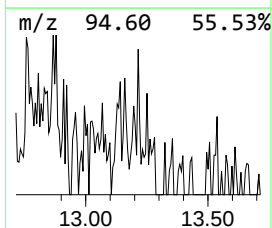
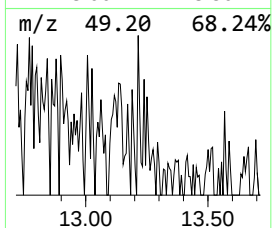
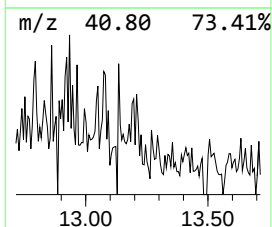
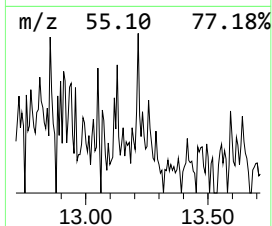
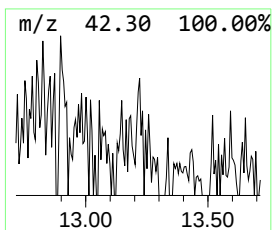
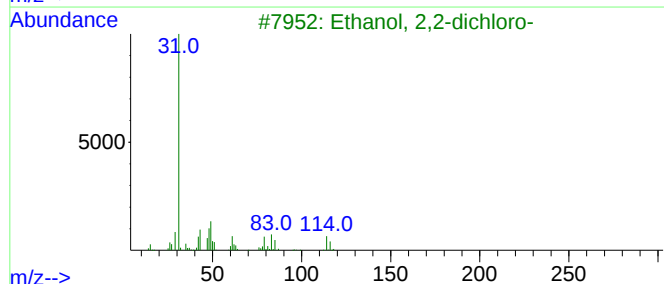
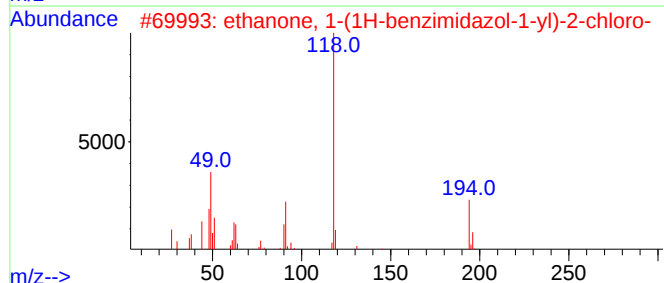
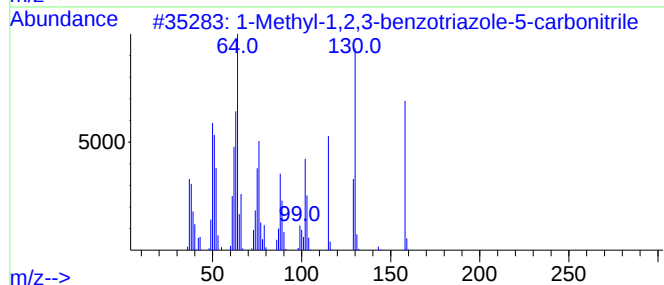
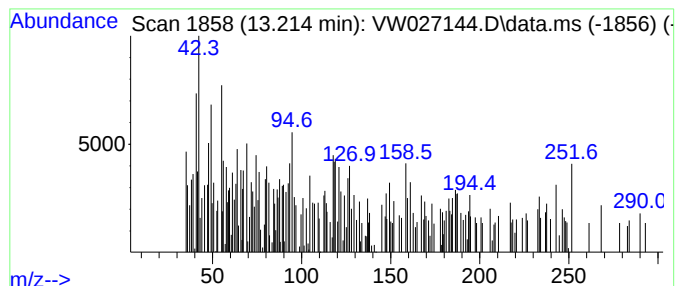
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.214	5.93 ug/L	9640	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-1,2,3-benzotriazole-5-c...	158	C8H6N4	1000435-82-2	25
2			ethanone, 1-(1H-benzimidazol-1-y...	194	C9H7C1N2O	1000400-31-1	15
3			Ethanol, 2,2-dichloro-	114	C2H4Cl2O	000598-38-9	11
4			tert-Butyl methyl sulfoxide	120	C5H12OS	014094-11-2	10
5			N,N'-Trimethylenebis[s-3-aminopr...	382	C9H22N2O6S4	035871-60-4	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW091923\
Data File : VW027144.D
Acq On : 19 Sep 2023 12:21
Operator : SY/MD
Sample : 04472-04RE
Misc : 3.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYK8RE

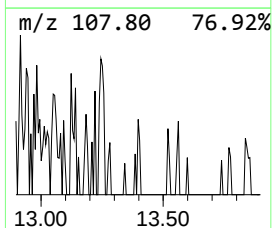
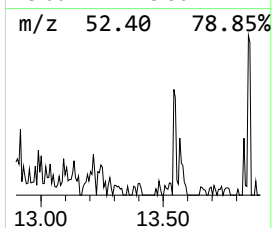
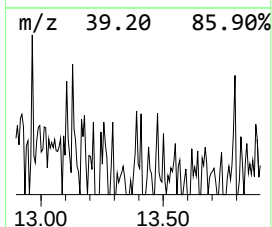
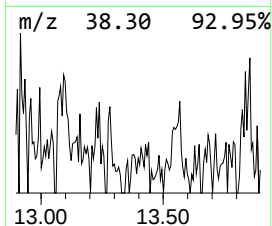
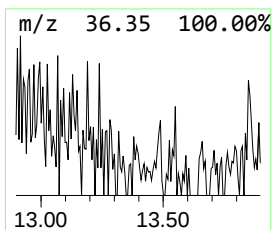
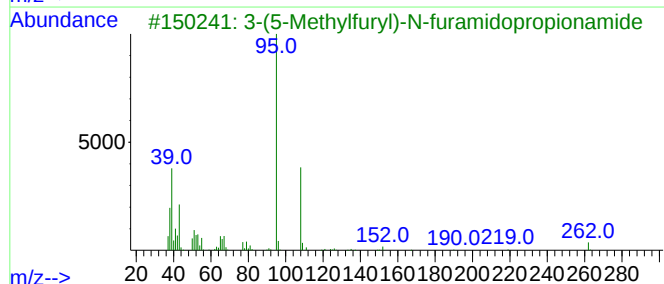
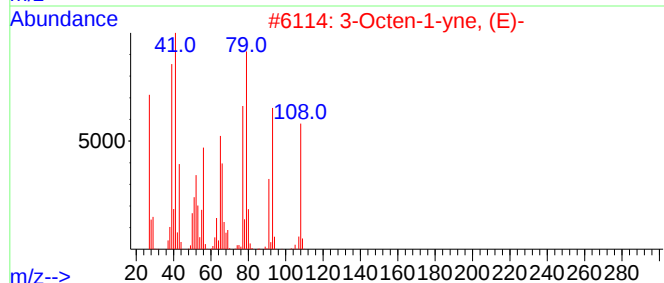
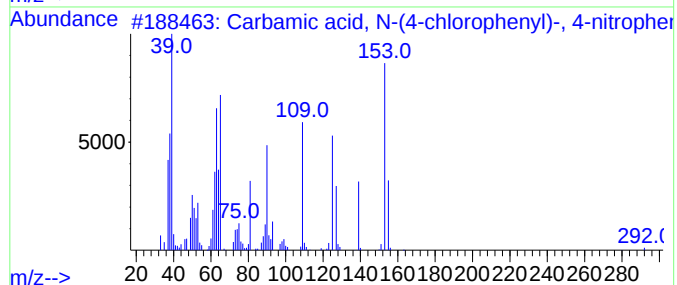
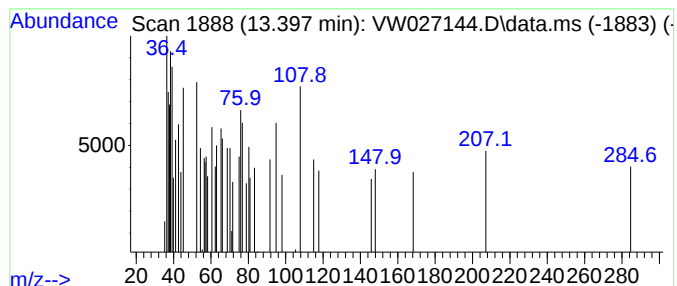
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Quant Title : SFAM01.0

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

Peak Number 6 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.397	1.51 ug/L	2453	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbamic acid, N-(4-chlorophenyl...	292	C13H9ClN2O4	003848-41-7	32
2			3-Octen-1-yne, (E)-	108	C8H12	042104-42-7	30
3			3-(5-Methylfuryl)-N-furamidoprop...	262	C13H14N2O4	331958-18-0	28
4			4-Amino-6-chloroquinolin	178	C9H7ClN2	020028-60-8	25
5			9,10-Diazatetracyclo[6.2.0(2,7)...	148	C8H8N2O	034098-80-1	16



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW091923\
Data File : VW027144.D
Acq On : 19 Sep 2023 12:21
Operator : SY/MD
Sample : 04472-04RE
Misc : 3.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYK8RE

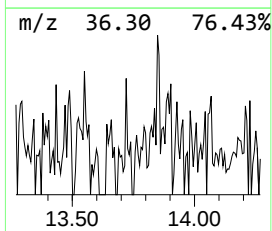
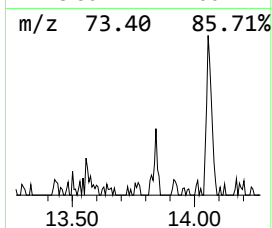
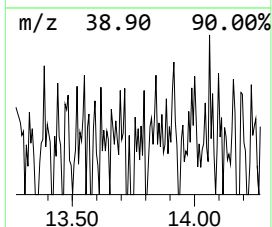
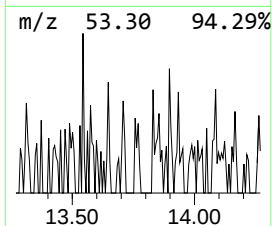
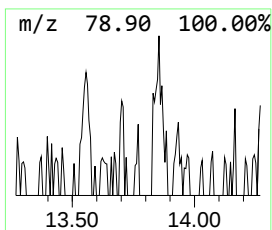
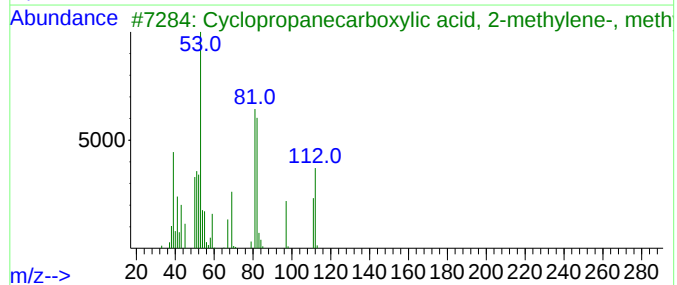
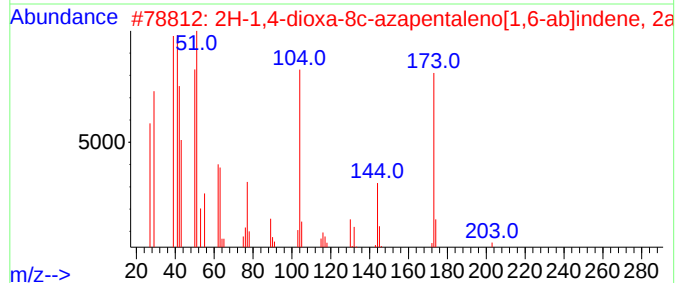
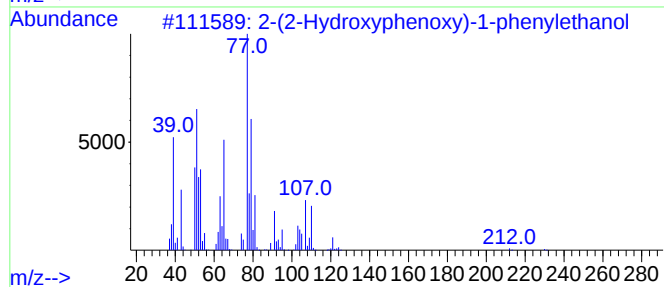
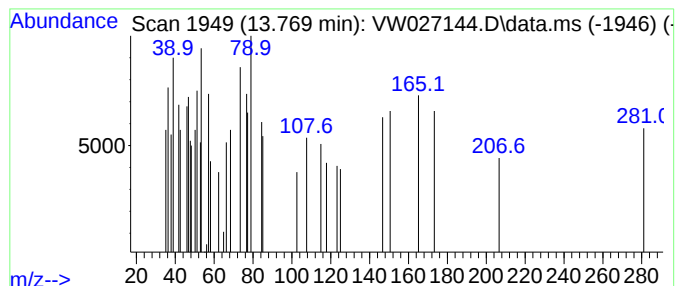
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.769	1.39 ug/L	2265	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2-Hydroxyphenoxy)-1-phenyleth...	230	C14H14O3	328104-89-8	32		
2	2H-1,4-dioxo-8c-azapentaleno[1,6...	203	C12H13NO2	1000400-04-0	27		
3	Cyclopropanecarboxylic acid, 2-m...	112	C6H8O2	088787-23-9	25		
4	1-Methyl-4-[2-nitroethenyl]-1H-p...	153	C6H7N3O2	003156-58-9	25		
5	4-Pentyn-1-ol	84	C5H8O	005390-04-5	23		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW091923\
Data File : VW027144.D
Acq On : 19 Sep 2023 12:21
Operator : SY/MD
Sample : 04472-04RE
Misc : 3.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYK8RE

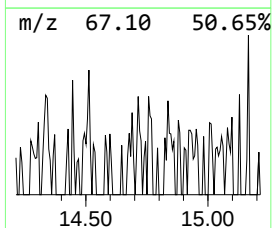
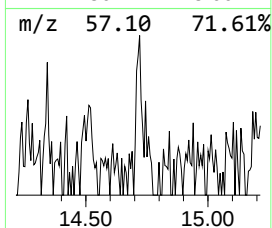
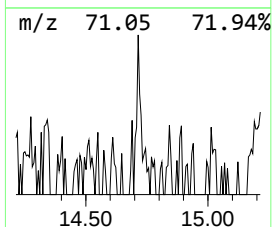
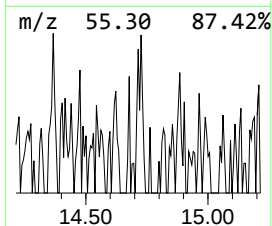
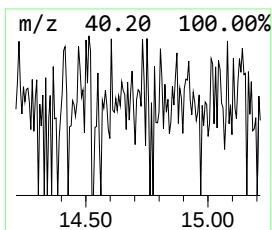
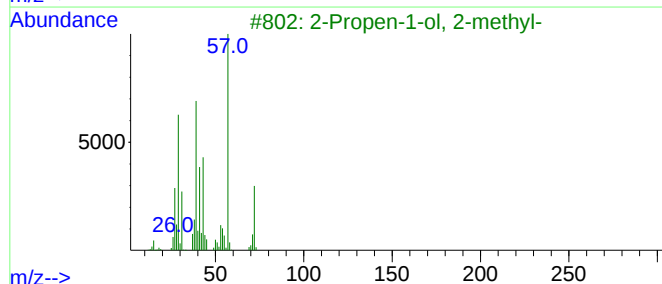
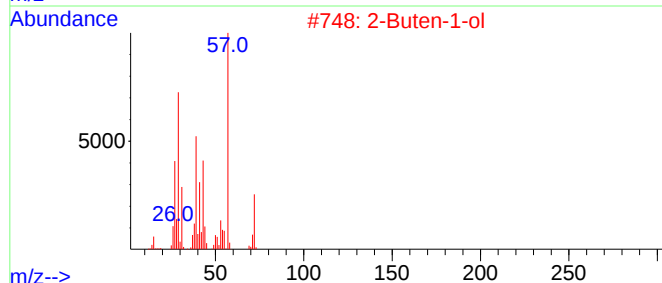
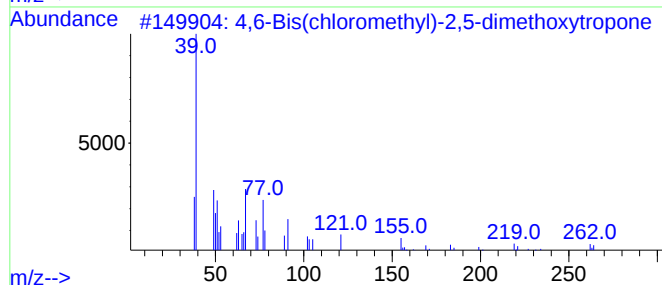
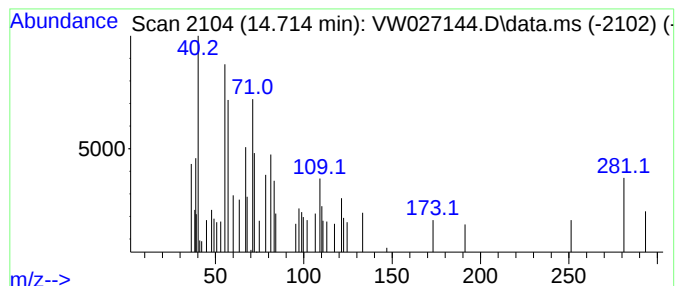
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-07 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.714	2.27 ug/L	3695	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,6-Bis(chloromethyl)-2,5-dimeth...	262	C11H12Cl2O3	1000433-38-3	17		
2	2-Buten-1-ol	72	C4H8O	006117-91-5	10		
3	2-Propen-1-ol, 2-methyl-	72	C4H8O	000513-42-8	9		
4	1-Ethyl-3,cis-(1,1-dimethylethyl...	214	C13H26O2	1000215-62-8	9		
5	1-(3-Acetamidophenyl)-3-(2,2,2-t...	438	C16H21Cl3N4O2S	294658-29-0	9		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW091923\
Data File : VW027144.D
Acq On : 19 Sep 2023 12:21
Operator : SY/MD
Sample : 04472-04RE
Misc : 3.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EZYK8RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM091823SMA.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-02	10.508	1.6	ug/L	2615	2	11.629	41682	25.0
unknown-03	11.288	25.0	ug/L	41682	2	11.629	41682	25.0
unknown-04	13.214	5.9	ug/L	9640	3	13.556	40663	25.0
unknown-05	13.397	1.5	ug/L	2453	3	13.556	40663	25.0
unknown-06	13.769	1.4	ug/L	2265	3	13.556	40663	25.0
unknown-07	14.714	2.3	ug/L	3695	3	13.556	40663	25.0