

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
 Data File : VW022465.D
 Acq On : 08 Apr 2022 15:43
 Operator : SY/VA
 Sample : N2293-13
 Misc : 6.32g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNP1

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\SFAMWLM040722SMA.M
 Title : SFAM01.0

Signal : TIC: VW022465.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.965	4	10	43	rVB	3667593	5790158	40.22%	16.295%
2	2.355	69	74	82	rBV	125052	207442	1.44%	0.584%
3	2.495	93	97	104	rBV	21136	42380	0.29%	0.119%
4	2.885	156	161	170	rVB	73741	157195	1.09%	0.442%
5	3.690	287	293	297	rBV5	3289	8465	0.06%	0.024%
6	3.897	323	327	328	rBV4	1809	2399	0.02%	0.007%
7	4.019	335	347	359	rBV	128577	381148	2.65%	1.073%
8	4.129	360	365	372	rVV2	8930	25196	0.18%	0.071%
9	4.208	376	378	389	rVB10	4144	9218	0.06%	0.026%
10	4.391	398	408	417	rBV3	9000	30473	0.21%	0.086%
11	4.916	486	494	506	rBV7	8394	29192	0.20%	0.082%
12	5.086	520	522	525	rBV3	1340	1692	0.01%	0.005%
13	5.361	563	567	569	rVV4	1571	1920	0.01%	0.005%
14	5.403	572	574	577	rVV4	1330	1830	0.01%	0.005%
15	5.440	577	580	582	rVV4	1459	1915	0.01%	0.005%
16	5.574	592	602	607	rVV10	2581	8174	0.06%	0.023%
17	5.800	636	639	644	rVB5	1692	2705	0.02%	0.008%
18	5.952	653	664	669	rBV10	5008	14041	0.10%	0.040%
19	6.068	681	683	685	rBV2	1514	1985	0.01%	0.006%
20	6.226	708	709	715	rVB5	1503	1981	0.01%	0.006%
21	6.299	715	721	722	rBV4	1851	2277	0.02%	0.006%
22	6.324	722	725	730	rVB5	1872	3004	0.02%	0.008%
23	6.610	767	772	774	rBV6	1531	2875	0.02%	0.008%
24	6.720	786	790	793	rBV6	1384	2125	0.01%	0.006%
25	6.982	830	833	834	rBV3	2321	1968	0.01%	0.006%
26	7.074	842	848	860	rBV	36683	107153	0.74%	0.302%
27	7.171	860	864	874	rVB	9736	25331	0.18%	0.071%
28	7.531	915	923	925	rBV7	2233	4833	0.03%	0.014%
29	7.647	933	942	952	rBV	216648	531610	3.69%	1.496%
30	7.903	975	984	988	rBV10	5985	18178	0.13%	0.051%
31	8.159	1024	1026	1032	rVB7	2684	3900	0.03%	0.011%
32	8.275	1035	1045	1047	rBV	397194	780785	5.42%	2.197%
33	8.323	1048	1053	1063	rVB	985333	2233241	15.51%	6.285%
34	8.567	1089	1093	1097	rVB5	3526	5722	0.04%	0.016%
35	8.701	1109	1115	1126	rVB6	6966	15866	0.11%	0.045%
36	8.842	1129	1138	1149	rBV	408182	802965	5.58%	2.260%

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Peak Location: TOP

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
 Title : SFAM01.0

37	9.043	1165	1171	1183	rBV3	18776	49027	0.34%	0.138%
38	9.275	1200	1209	1219	rBV	310353	657910	4.57%	1.852%
39	9.500	1243	1246	1253	rVB7	4098	8386	0.06%	0.024%
40	9.610	1260	1264	1268	rVB7	2063	3378	0.02%	0.010%
41	9.659	1268	1272	1275	rBV6	1984	3340	0.02%	0.009%
42	9.701	1276	1279	1284	rVV6	3525	5883	0.04%	0.017%
43	9.768	1288	1290	1291	rVV2	2039	1868	0.01%	0.005%
44	9.817	1292	1298	1305	rVB7	7558	17536	0.12%	0.049%
45	9.963	1314	1322	1328	rBV	346351	653169	4.54%	1.838%
46	10.037	1328	1334	1341	rVB	159334	292710	2.03%	0.824%
47	10.213	1358	1363	1366	rBV	7332	12004	0.08%	0.034%
48	10.323	1373	1381	1386	rBV	596409	1046226	7.27%	2.944%
49	10.384	1386	1391	1400	rVV2	462872	840359	5.84%	2.365%
50	10.469	1400	1405	1415	rVB5	25660	71966	0.50%	0.203%
51	10.579	1416	1423	1430	rVV	100805	178080	1.24%	0.501%
52	10.713	1438	1445	1459	rVB	222799	417732	2.90%	1.176%
53	10.860	1465	1469	1473	rVB	11888	15782	0.11%	0.044%
54	10.921	1473	1479	1485	rBV	177302	308710	2.14%	0.869%
55	11.061	1497	1502	1514	rVB2	32727	59648	0.41%	0.168%
56	11.177	1517	1521	1527	rVB5	6094	9082	0.06%	0.026%
57	11.292	1533	1540	1550	rBV	90151	167503	1.16%	0.471%
58	11.372	1550	1553	1556	rBV3	4659	6741	0.05%	0.019%
59	11.475	1566	1570	1574	rBV5	3099	5334	0.04%	0.015%
60	11.555	1579	1583	1587	rVB6	4174	6978	0.05%	0.020%
61	11.658	1587	1600	1608	rBV	7924011	14396928	100.00%	40.517%
62	11.835	1620	1629	1638	rVB3	23815	59827	0.42%	0.168%
63	12.158	1678	1682	1690	rVB2	17958	34062	0.24%	0.096%
64	12.347	1708	1713	1716	rBV6	5517	9845	0.07%	0.028%
65	12.463	1727	1732	1737	rVB6	5658	9907	0.07%	0.028%
66	12.597	1749	1754	1759	rBV3	9166	19493	0.14%	0.055%
67	12.658	1759	1764	1765	rVV	93297	127178	0.88%	0.358%
68	12.689	1765	1769	1775	rVB	258251	440538	3.06%	1.240%
69	12.890	1794	1802	1807	rBV	93940	181140	1.26%	0.510%
70	13.115	1833	1839	1846	rBV4	71085	135932	0.94%	0.383%
71	13.182	1846	1850	1856	rVB5	14201	23594	0.16%	0.066%
72	13.249	1856	1861	1864	rBV	16316	27385	0.19%	0.077%
73	13.292	1864	1868	1874	rVB6	19145	42600	0.30%	0.120%
74	13.365	1875	1880	1885	rBV2	32622	61076	0.42%	0.172%
75	13.457	1891	1895	1898	rBV6	3264	5748	0.04%	0.016%
76	13.499	1898	1902	1904	rVV5	7817	12064	0.08%	0.034%

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 Title : SFAM01.0

77	13.554	1904	1911	1919	rVV	697814	1153749	8.01%	3.247%
78	13.640	1920	1925	1934	rVB2	76571	138608	0.96%	0.390%
79	13.853	1953	1960	1971	rVV3	1025989	1979703	13.75%	5.571%
80	13.944	1972	1975	1980	rVB3	8918	11638	0.08%	0.033%
81	14.200	2011	2017	2023	rBV6	11158	22676	0.16%	0.064%
82	14.426	2050	2054	2056	rBV4	5620	8878	0.06%	0.025%
83	14.463	2056	2060	2065	rVV5	15977	25411	0.18%	0.072%
84	14.694	2093	2098	2100	rBV5	7796	12917	0.09%	0.036%
85	14.737	2100	2105	2109	rVV7	12496	23585	0.16%	0.066%
86	14.786	2109	2113	2121	rVV5	12076	25279	0.18%	0.071%
87	14.956	2136	2141	2146	rBV7	5901	12187	0.08%	0.034%
88	15.060	2154	2158	2161	rBV6	11418	19394	0.13%	0.055%
89	15.316	2196	2200	2203	rBV6	3545	5699	0.04%	0.016%
90	15.359	2203	2207	2212	rVV2	15743	28736	0.20%	0.081%
91	15.420	2212	2217	2221	rVB3	11070	18093	0.13%	0.051%
92	15.737	2261	2269	2281	rBV3	98211	222927	1.55%	0.627%
93	16.170	2333	2340	2348	rBV3	8834	22416	0.16%	0.063%
94	16.267	2348	2356	2363	rBV9	12600	29324	0.20%	0.083%
95	16.365	2369	2372	2379	rVB9	5972	11105	0.08%	0.031%
96	16.438	2379	2384	2389	rBV9	9915	22814	0.16%	0.064%
97	16.535	2398	2400	2404	rVB5	2227	2909	0.02%	0.008%
98	16.578	2404	2407	2409	rBV3	1958	2542	0.02%	0.007%
99	16.639	2410	2417	2423	rVV8	9343	24124	0.17%	0.068%
100	16.730	2427	2432	2438	rVB9	8723	20218	0.14%	0.057%

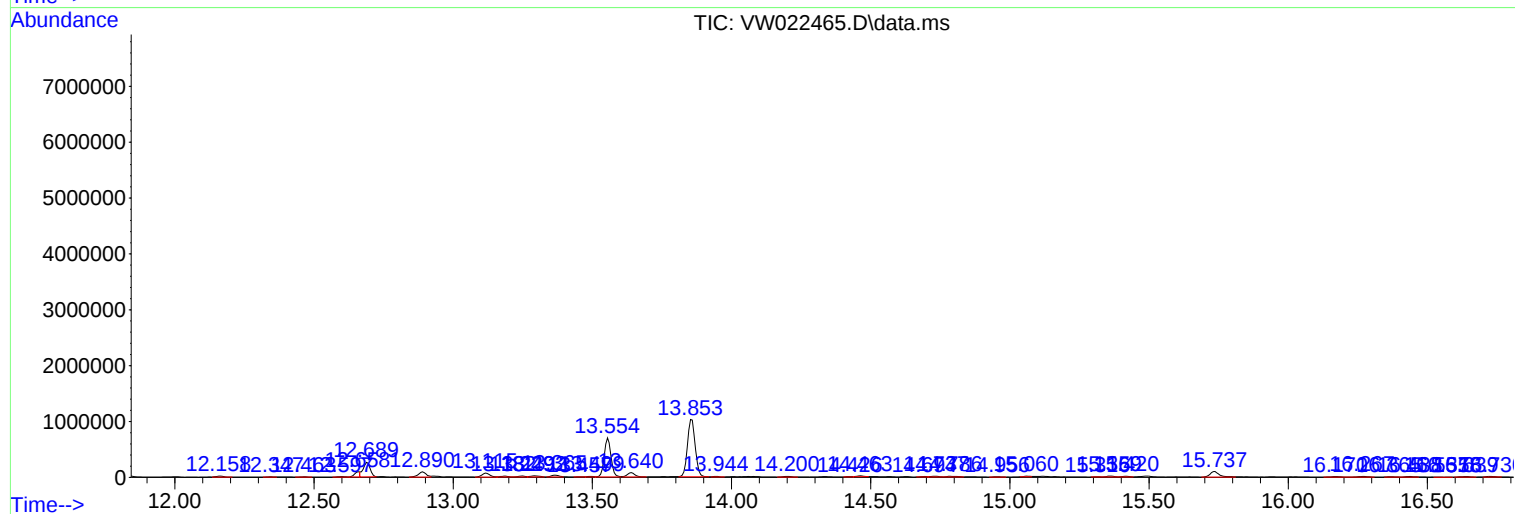
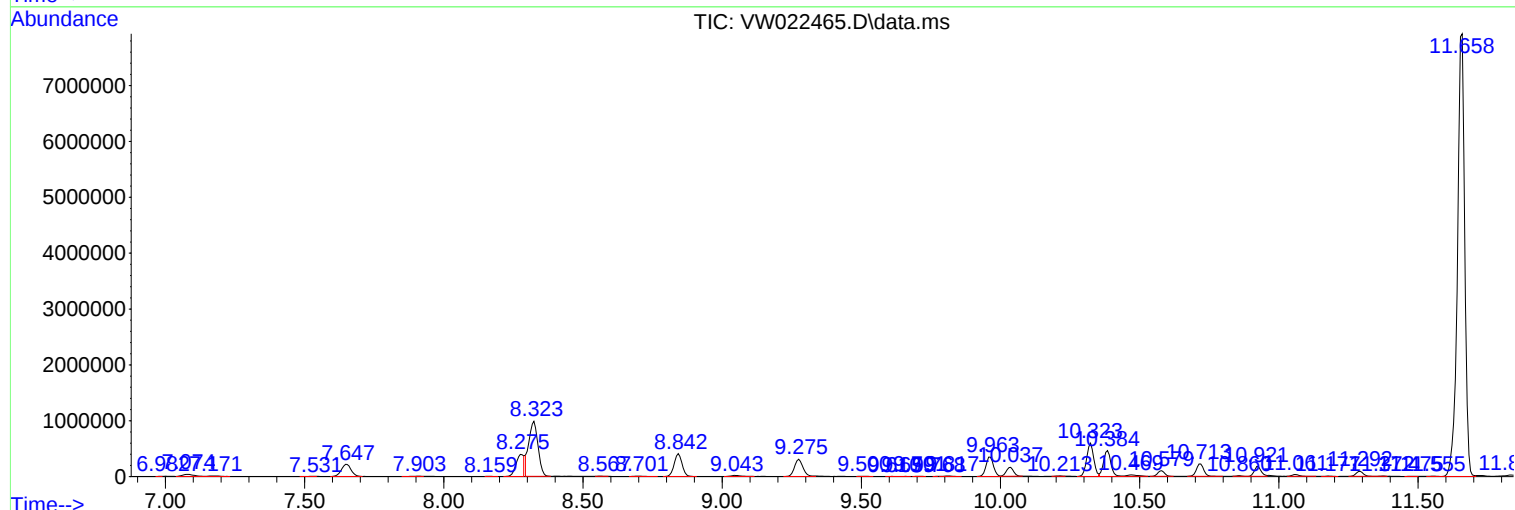
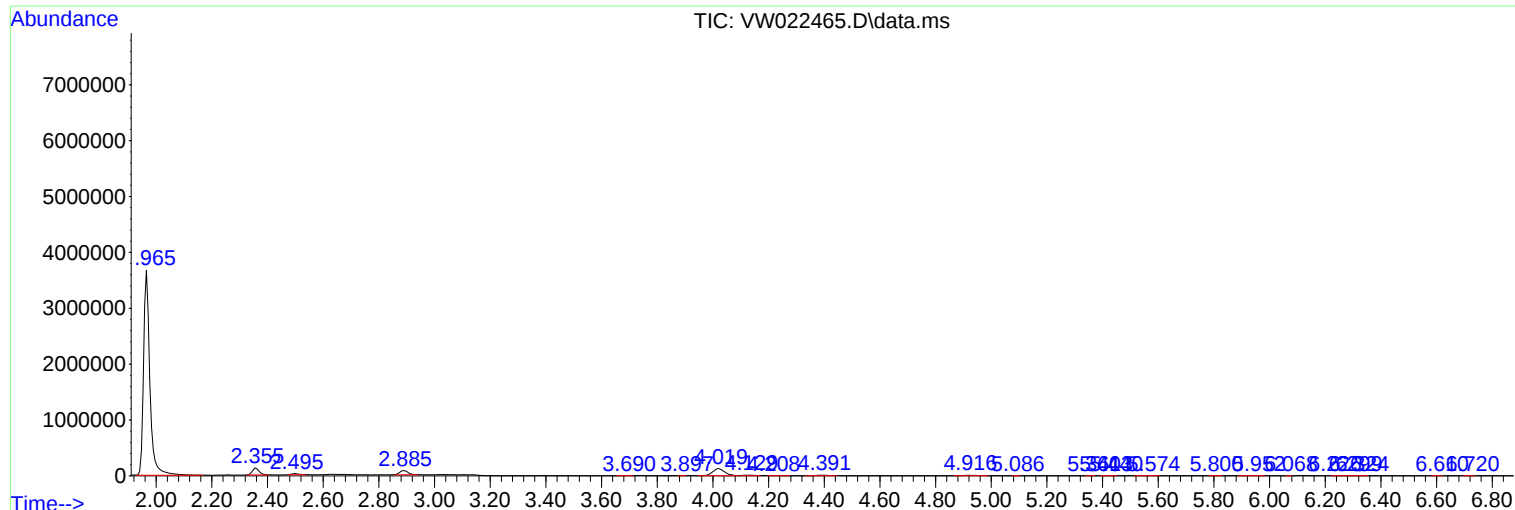
Sum of corrected areas: 35532943

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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



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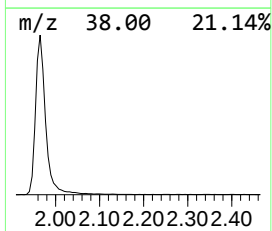
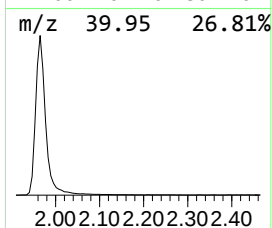
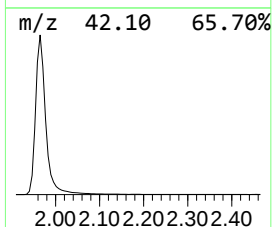
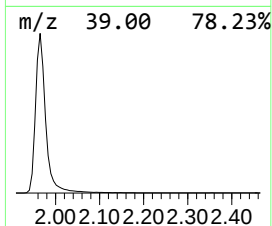
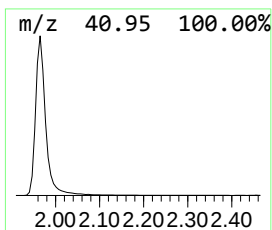
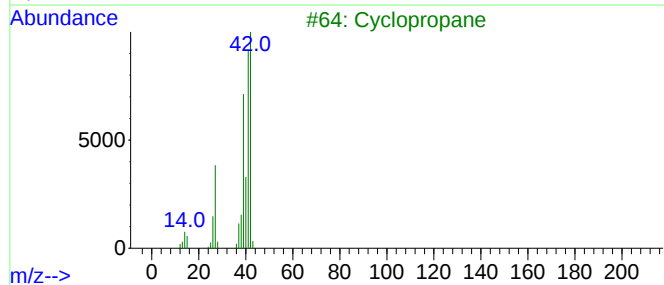
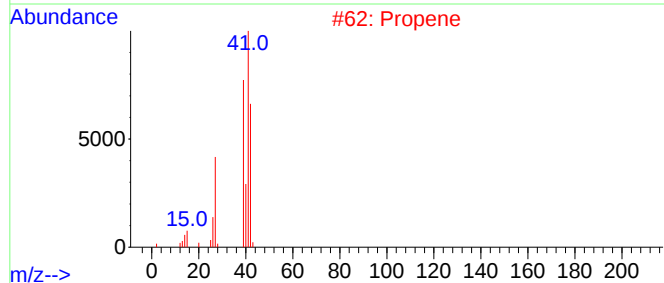
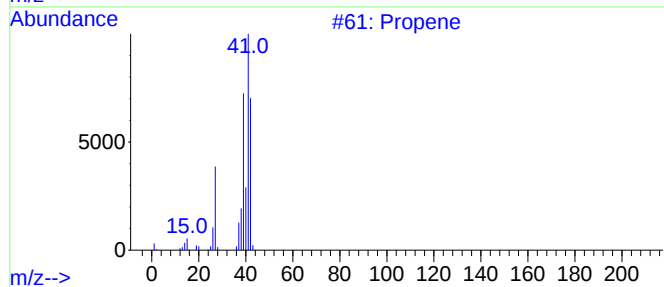
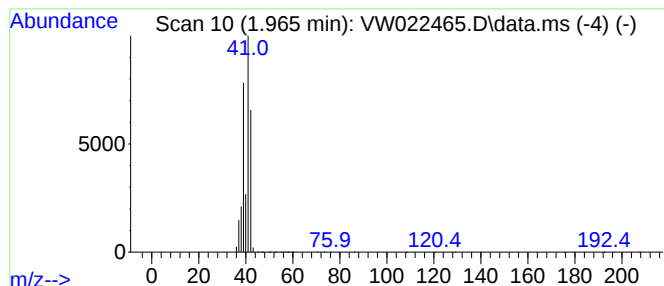
Instrument :
MSVOA_W
ClientSampleId :
ETNP1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.965	180.27 ug/L	5790160	1,4-Difluorobenzene	8.842	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Propene	42	C3H6	000115-07-1	59
3	Cyclopropane	42	C3H6	000075-19-4	10
4	Cyclopropene	40	C3H4	002781-85-3	10
5	Cyclopropane	42	C3H6	000075-19-4	7



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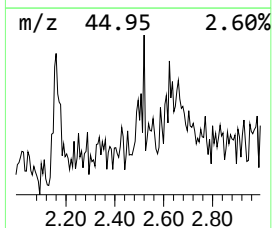
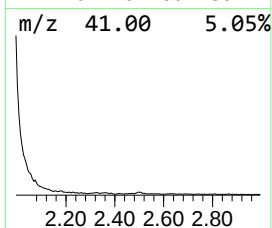
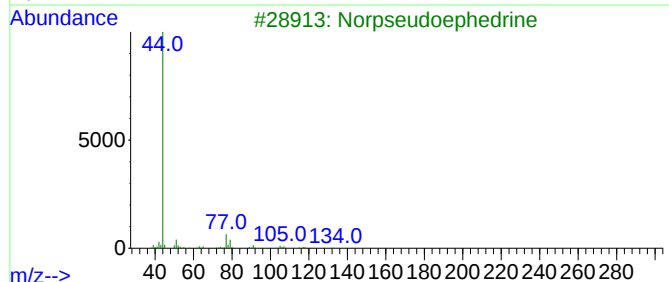
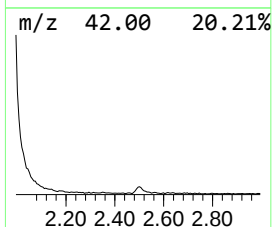
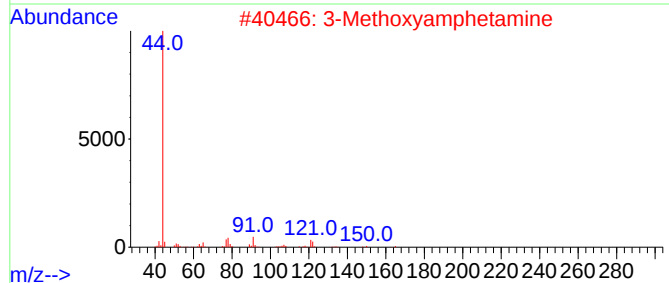
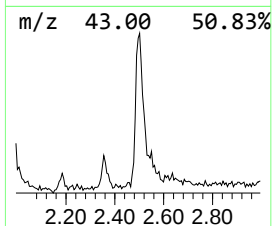
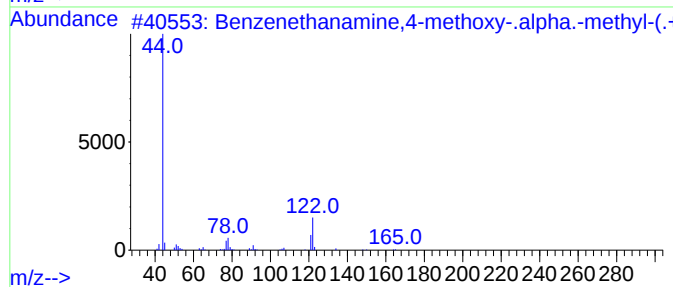
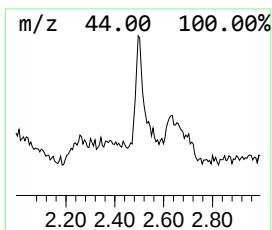
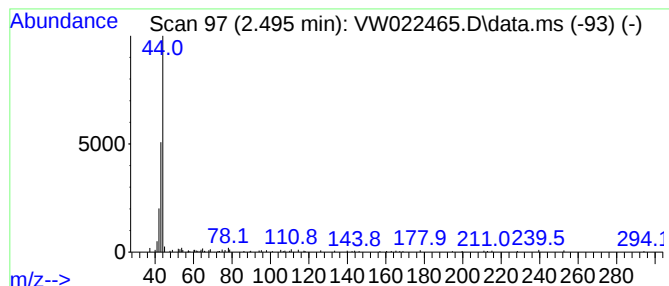
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 Benzenethanamine,4-methoxy-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.495	1.32 ug/L	42380	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenethanamine,4-methoxy-.alph...	165	C10H15NO	023239-32-9	50
2			3-Methoxyamphetamine	165	C10H15NO	017862-85-0	40
3			Norpseudoephedrine	151	C9H13NO	036393-56-3	33
4			Brolamfetamine	273	C11H16BrNO2	064638-07-9	33
5			dl-Alanine	89	C3H7NO2	000302-72-7	9



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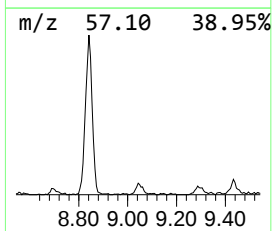
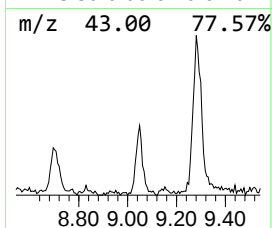
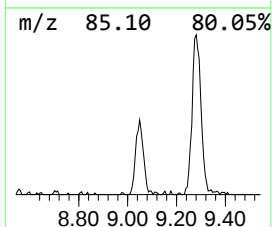
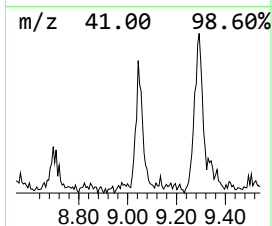
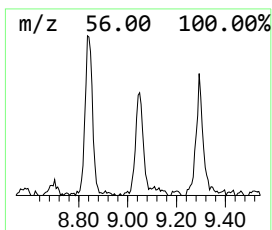
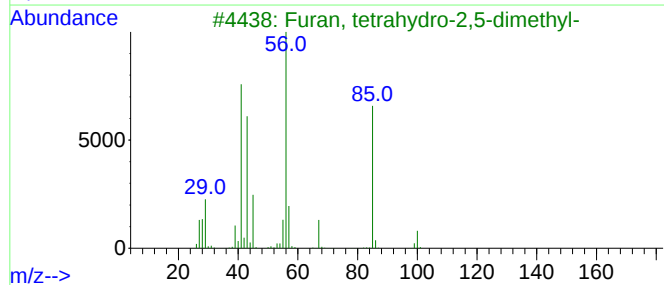
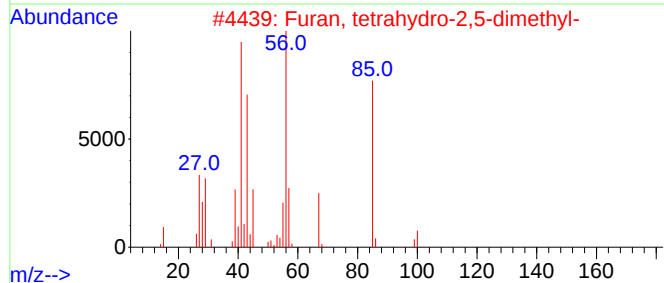
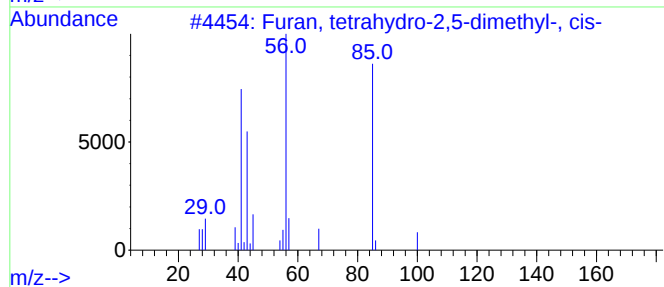
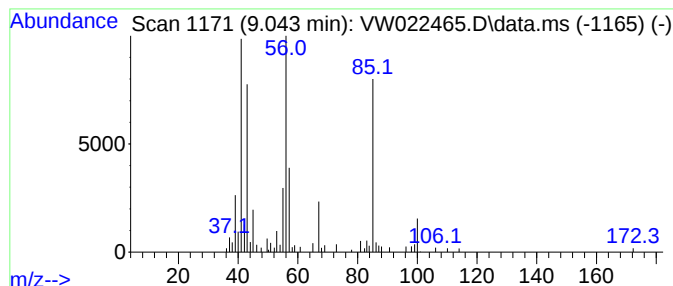
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Furan, tetrahydro-2,5-dimet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.043	1.53 ug/L	49027	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	002144-41-4	86
2			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	83
3			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	72
4			2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	72
5			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	038484-59-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

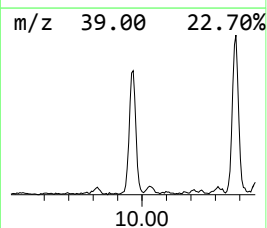
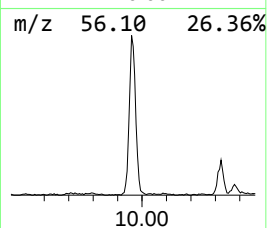
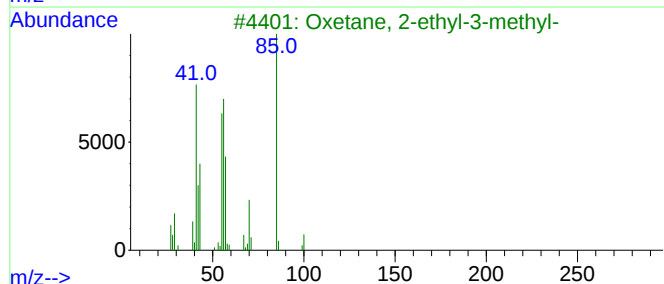
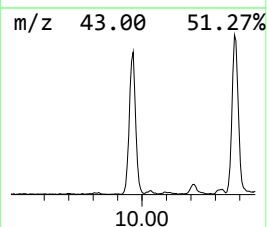
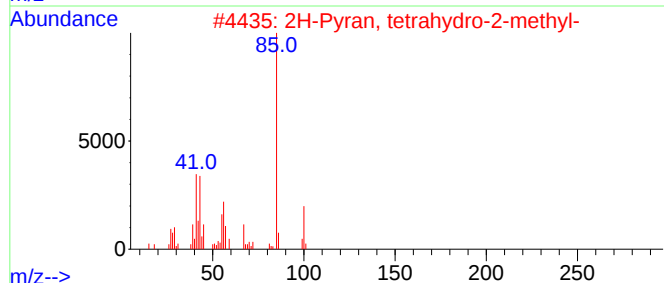
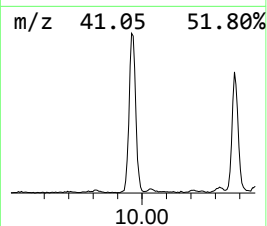
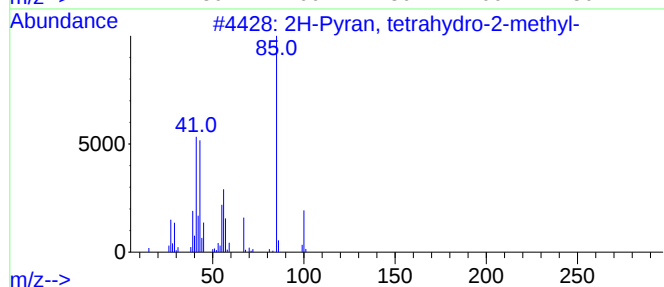
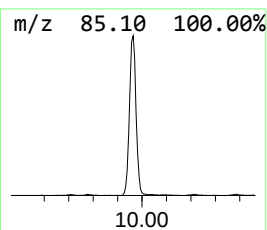
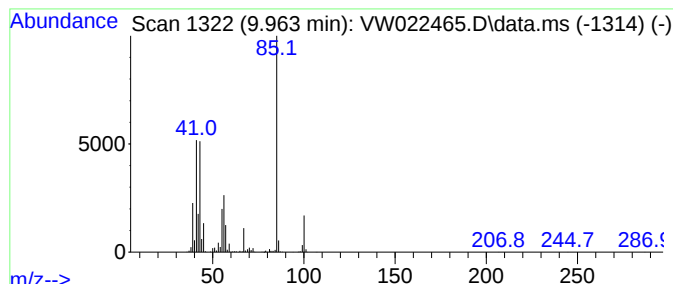
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Quant Title : SFAM01.0

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

Peak Number 4 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	20.34 ug/L	653169	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	95
2			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	90
3			Oxetane, 2-ethyl-3-methyl-	100	C6H12O	053778-62-4	72
4			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	59
5			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

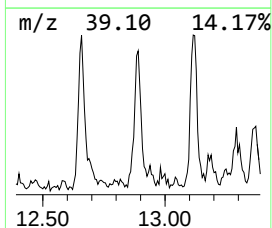
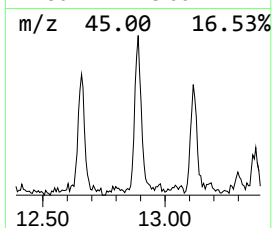
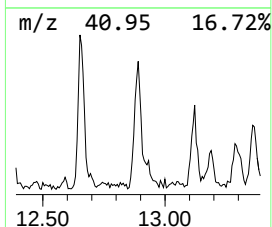
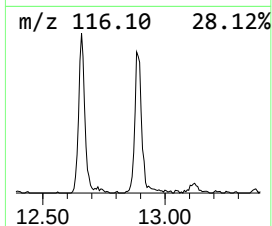
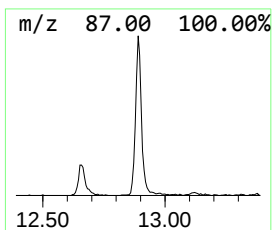
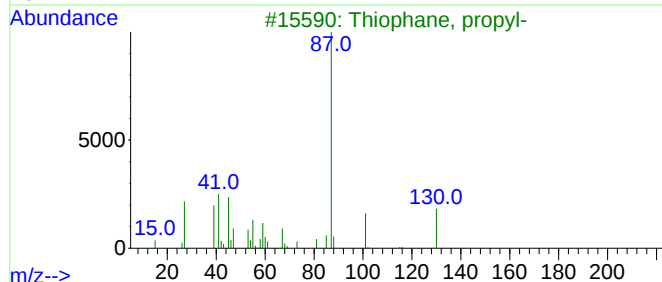
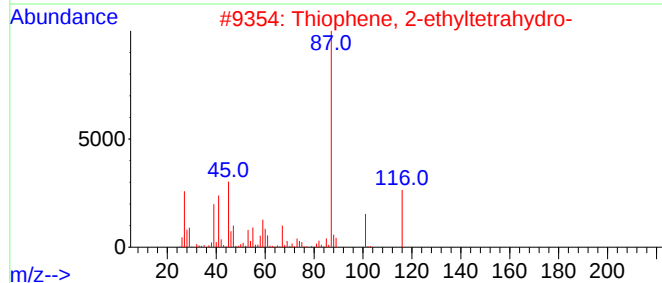
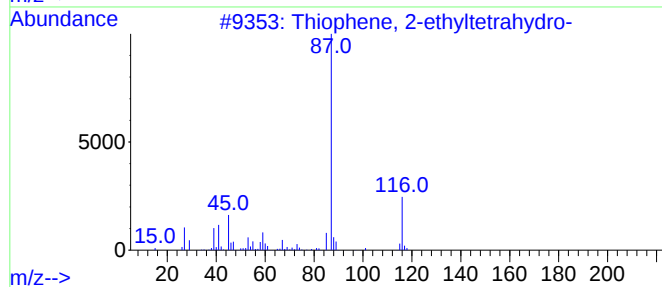
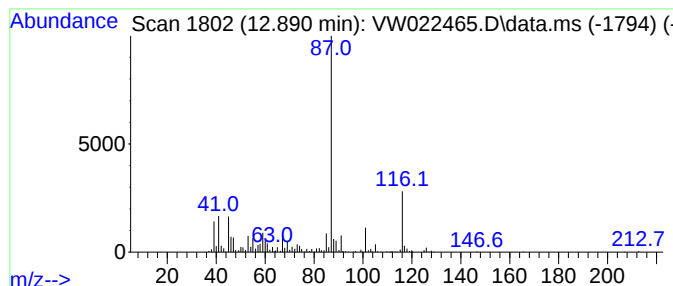
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Quant Title : SFAM01.0

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

Peak Number 6 Thiophene, 2-ethyltetrahydro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	3.93 ug/L	181140	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiophene, 2-ethyltetrahydro-	116	C6H12S	001551-32-2	94
2			Thiophene, 2-ethyltetrahydro-	116	C6H12S	001551-32-2	90
3			Thiophane, propyl-	130	C7H14S	001551-34-4	72
4			Hydrazine, 1-ethyl-1-(1-methylpr...	116	C6H16N2	020325-97-7	64
5			Thiophene, 2-butyltetrahydro-	144	C8H16S	001613-49-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

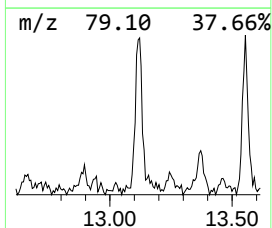
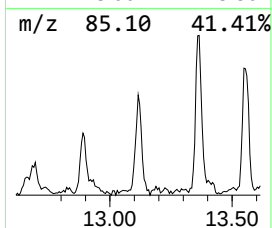
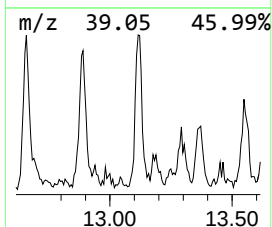
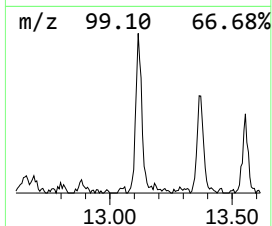
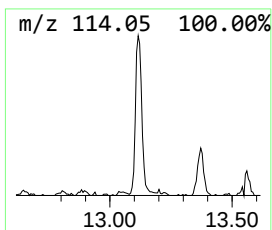
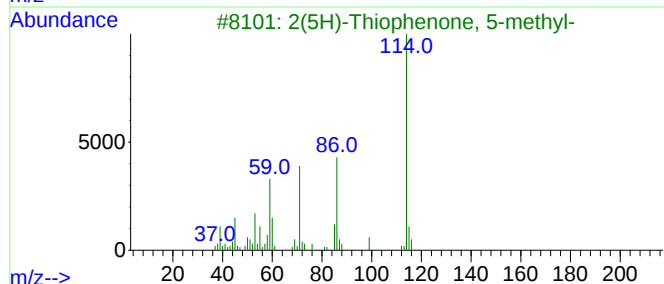
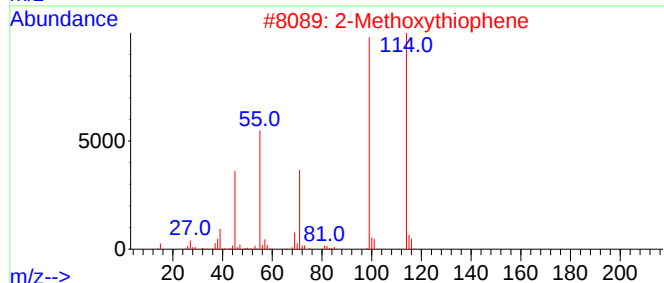
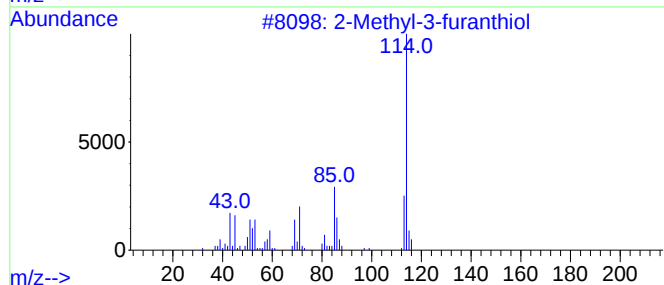
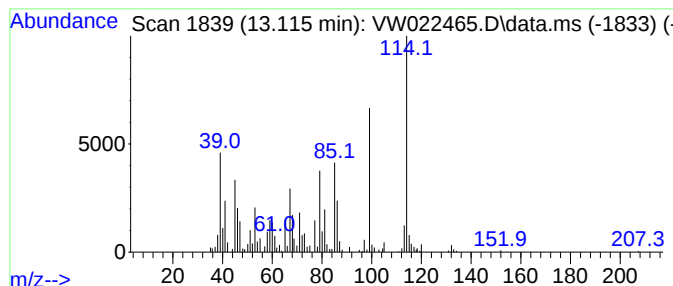
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.115	2.95 ug/L	135932	1,4-Dichlorobenzene-d4	13.554

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methyl-3-furanthiol	114	C5H6OS	028588-74-1	43
2		2-Methoxythiophene	114	C5H6OS	016839-97-7	41
3		2(5H)-Thiophenone, 5-methyl-	114	C5H6OS	007210-64-2	38
4		1-Propene, 1,1'-thiobis-	114	C6H10S	033922-80-4	38
5		Thiophene, 3-methoxy-	114	C5H6OS	017573-92-1	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

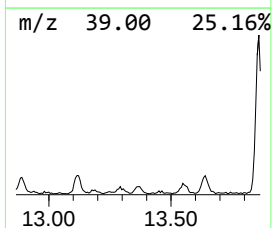
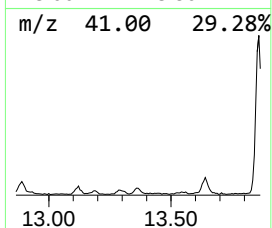
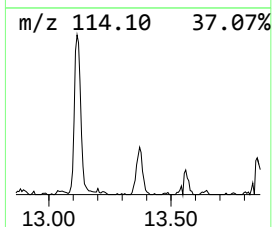
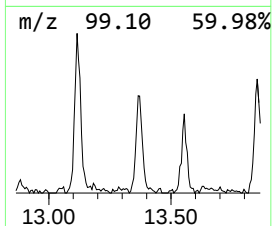
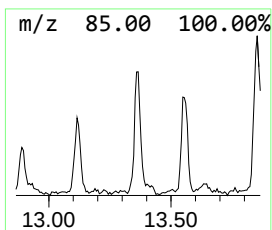
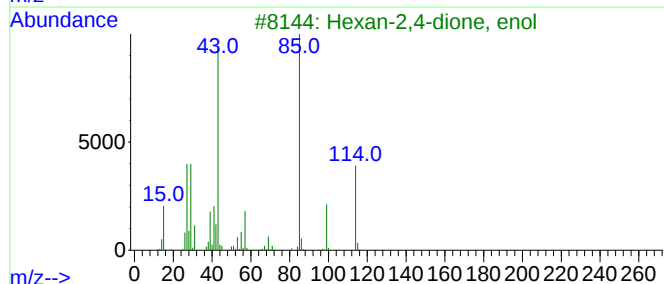
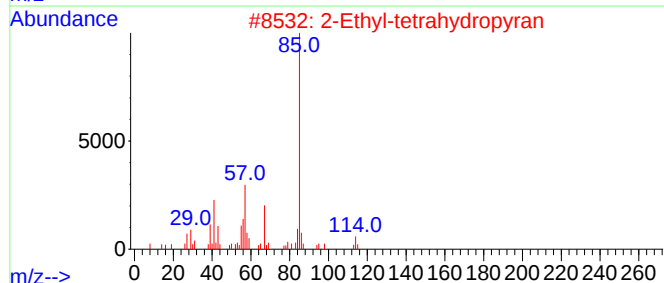
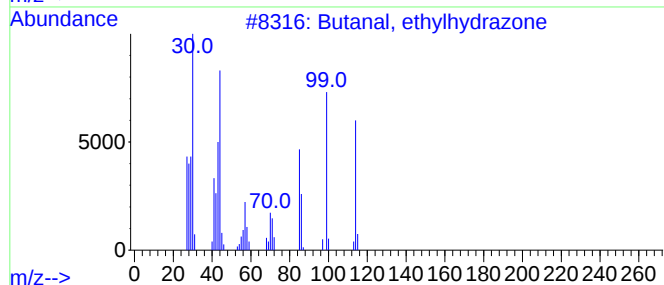
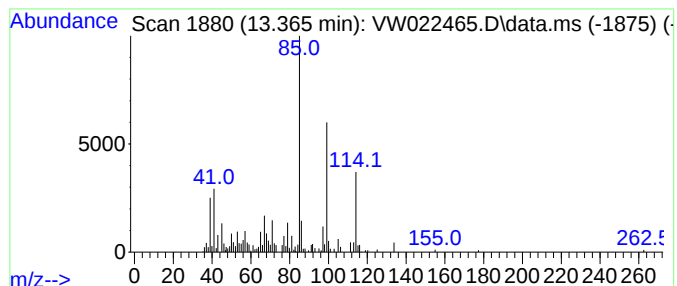
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.365	1.32 ug/L	61076	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanal, ethylhydrazone	114	C6H14N2	051576-30-8	49
2			2-Ethyl-tetrahydropyran	114	C7H14O	1000386-40-1	47
3			Hexan-2,4-dione, enol	114	C6H10O2	1000202-23-9	47
4			3-Ethoxy-2-methylacroleine	114	C6H10O2	042588-57-8	47
5			Propanal, propylhydrazone	114	C6H14N2	019718-39-9	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

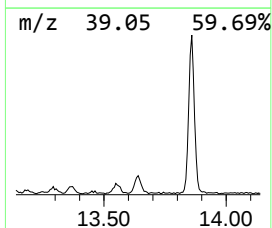
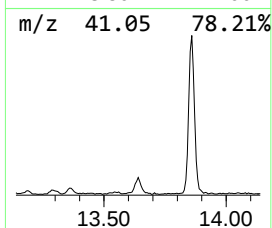
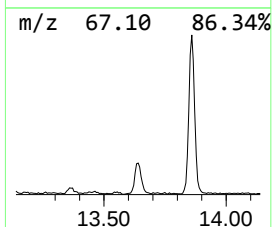
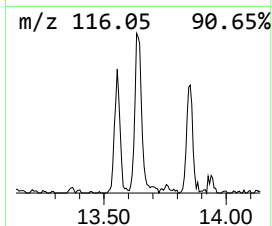
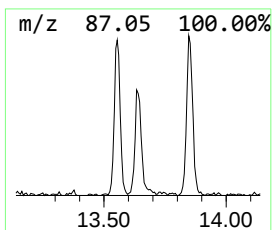
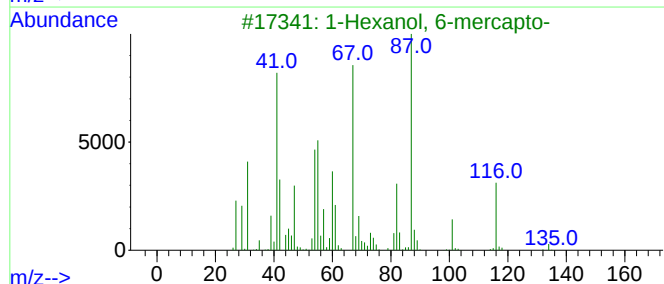
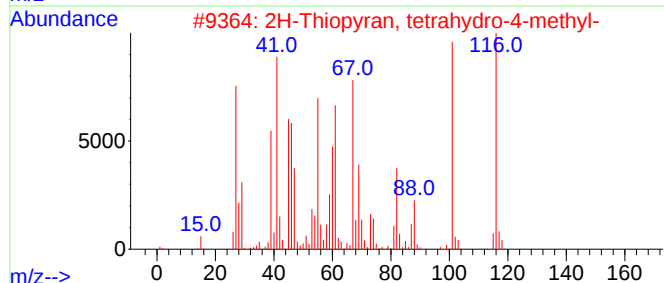
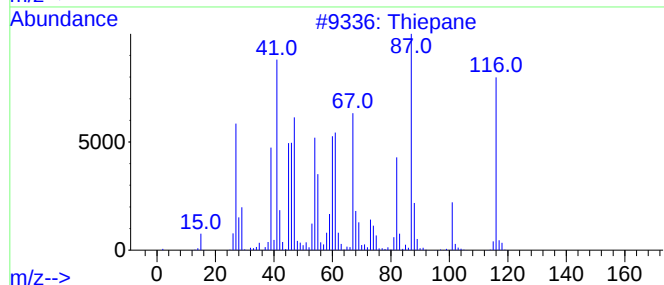
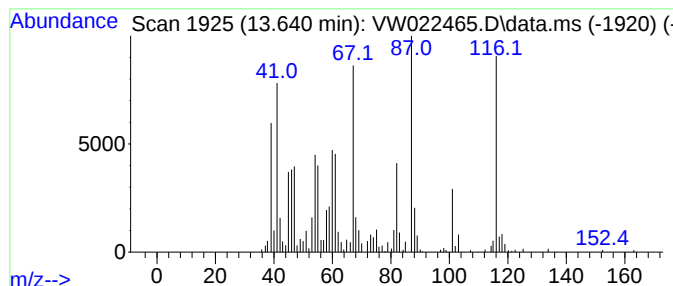
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Quant Title : SFAM01.0

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

Peak Number 9 Thiepane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	3.00 ug/L	138608	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiepane	116	C6H12S	004753-80-4	95
2			2H-Thiopyran, tetrahydro-4-methyl-	116	C6H12S	005161-17-1	49
3			1-Hexanol, 6-mercapto-	134	C6H14OS	001633-78-9	47
4			2H-Thiopyran, tetrahydro-4-methyl-	116	C6H12S	005161-17-1	46
5			4H-Thiopyran-4-one, tetrahydro-	116	C5H8OS	001072-72-6	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

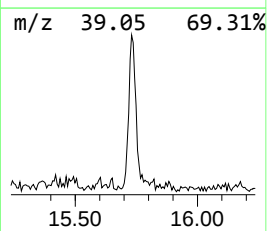
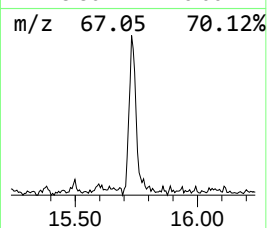
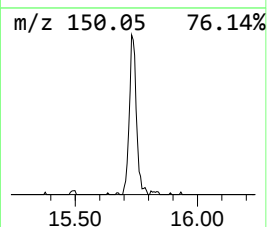
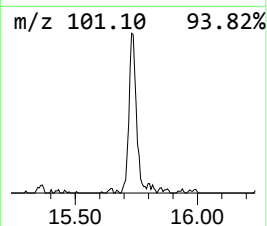
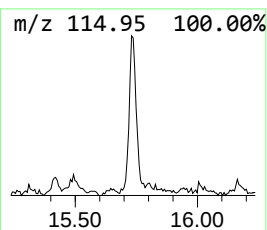
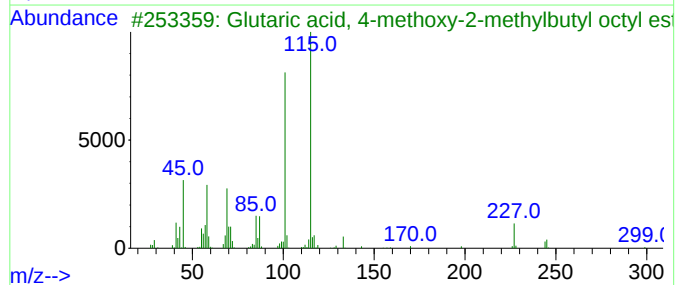
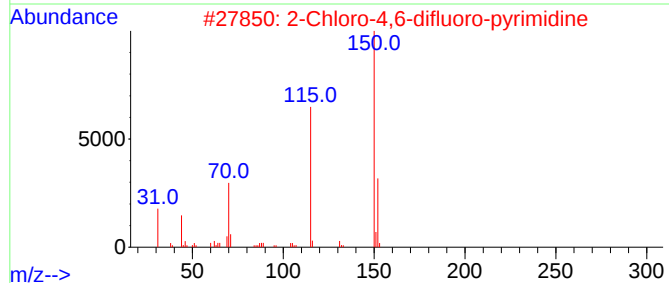
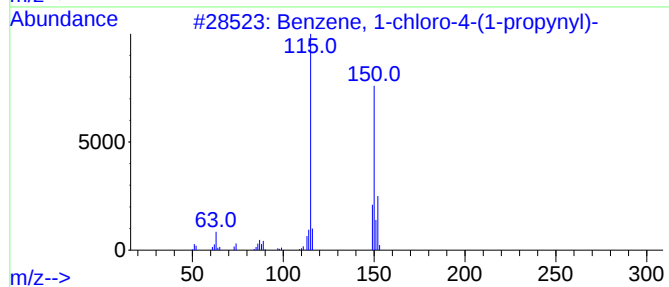
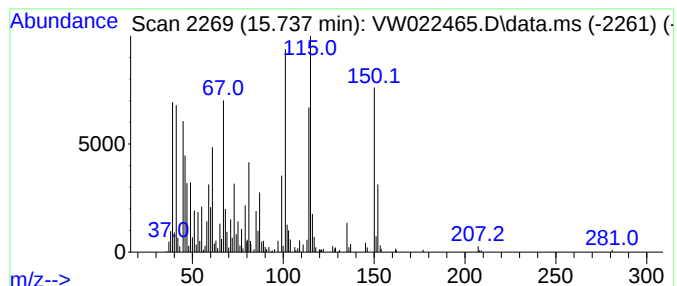
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.737	4.83 ug/L	222927	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-4-(1-propynyl)-	150	C9H7Cl	002809-65-6	43
2			2-Chloro-4,6-difluoro-pyrimidine	150	C4HClF2N2	038953-30-9	22
3			Glutaric acid, 4-methoxy-2-methy...	344	C19H36O5	1000359-41-7	14
4			Glutaric acid, isobutyl 2-(2-nit...	337	C17H23NO6	1000377-52-1	14
5			Glutaric acid, 2-(2-nitrophenyl)...	323	C16H21NO6	1000377-52-0	14



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022465.D
Acq On : 08 Apr 2022 15:43
Operator : SY/VA
Sample : N2293-13
Misc : 6.32g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.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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Benzenethanamin...	2.495	1.3	ug/L	42380	1	8.842	802965	25.0
Furan, tetrahyd...	9.043	1.5	ug/L	49027	1	8.842	802965	25.0
2H-Pyran, tetra...	9.963	20.3	ug/L	653169	1	8.842	802965	25.0
Thiophene, 2-et...	12.890	3.9	ug/L	181140	3	13.554	1153750	25.0
unknown-01	13.115	3.0	ug/L	135932	3	13.554	1153750	25.0
unknown-02	13.365	1.3	ug/L	61076	3	13.554	1153750	25.0
Thiepane	13.640	3.0	ug/L	138608	3	13.554	1153750	25.0
unknown-03	15.737	4.8	ug/L	222927	3	13.554	1153750	25.0