

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
 Data File : VW028814.D
 Acq On : 26 May 2024 02:04
 Operator : SY/MD
 Sample : P2590-18
 Misc : 6.11g/10mL/MSVOA_W/SOIL/A
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BHBM2

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\SFAMWLM050224SMA.M
 Title : SFAM01.0

Signal : TIC: VW028814.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.710	134	135	138	rBV3	668	657	2.58%	0.373%
2	2.753	140	142	143	rVB2	758	436	1.71%	0.247%
3	2.899	162	166	172	rVB5	2021	4228	16.59%	2.397%
4	3.137	201	205	207	rBV5	843	1006	3.95%	0.570%
5	3.222	217	219	222	rVB3	692	420	1.65%	0.238%
6	3.838	319	320	323	rVB2	652	543	2.13%	0.308%
7	4.009	345	348	349	rBV2	1723	1623	6.37%	0.920%
8	4.179	374	376	378	rBV3	769	785	3.08%	0.445%
9	4.326	398	400	403	rBV4	515	529	2.08%	0.300%
10	4.387	408	410	412	rVB2	530	444	1.74%	0.252%
11	4.454	419	421	423	rBV3	562	501	1.97%	0.284%
12	4.533	433	434	437	rBV3	602	614	2.41%	0.348%
13	4.600	443	445	448	rBV3	464	465	1.82%	0.264%
14	4.771	472	473	475	rBV2	708	716	2.81%	0.406%
15	5.283	553	557	558	rVB3	572	706	2.77%	0.400%
16	5.313	560	562	565	rVV3	773	667	2.62%	0.378%
17	5.453	582	585	588	rBV3	719	1074	4.21%	0.609%
18	5.582	604	606	609	rVV3	575	568	2.23%	0.322%
19	5.673	619	621	623	rVV2	709	460	1.80%	0.261%
20	5.722	627	629	630	rVB2	781	431	1.69%	0.244%
21	5.935	662	664	666	rBV3	607	598	2.35%	0.339%
22	6.051	681	683	687	rVB3	884	703	2.76%	0.399%
23	6.088	687	689	692	rBV3	654	538	2.11%	0.305%
24	6.222	706	711	715	rBV5	909	1235	4.85%	0.700%
25	6.264	716	718	722	rBV3	637	832	3.26%	0.472%
26	6.362	731	734	735	rBV2	456	474	1.86%	0.269%
27	6.459	747	750	754	rBV3	635	592	2.32%	0.336%
28	6.618	774	776	778	rVB2	704	472	1.85%	0.268%
29	6.697	784	789	790	rBV4	785	804	3.15%	0.456%
30	6.831	809	811	814	rBV4	690	782	3.07%	0.443%
31	6.941	826	829	831	rBV3	552	473	1.86%	0.268%
32	7.063	846	849	851	rBV3	382	421	1.65%	0.239%
33	7.087	851	853	855	rBV2	1031	1188	4.66%	0.674%
34	7.209	871	873	875	rVB3	909	657	2.58%	0.373%
35	7.410	905	906	909	rVB3	742	505	1.98%	0.286%
36	7.453	909	913	917	rBV4	569	971	3.81%	0.551%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050224SMA.M
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37	7.502	919	921	923	rBV2	595	468	1.84%	0.265%
38	7.532	924	926	929	rVB3	703	606	2.38%	0.344%
39	7.563	929	931	932	rBV2	628	456	1.79%	0.259%
40	7.648	941	945	952	rVB2	4526	9797	38.44%	5.555%
41	7.727	955	958	961	rBV4	639	949	3.72%	0.538%
42	7.758	961	963	967	rVB2	704	708	2.78%	0.401%
43	7.965	995	997	1000	rBV3	615	666	2.61%	0.378%
44	8.020	1002	1006	1008	rBV3	594	819	3.21%	0.464%
45	8.215	1035	1038	1041	rBV3	624	806	3.16%	0.457%
46	8.282	1041	1049	1061	rBV4	7071	25485	100.00%	14.451%
47	8.721	1119	1121	1123	rBV3	782	647	2.54%	0.367%
48	8.764	1126	1128	1129	rBV2	659	420	1.65%	0.238%
49	8.843	1134	1141	1149	rVV2	5751	13251	52.00%	7.514%
50	9.056	1173	1176	1177	rVV3	559	452	1.77%	0.256%
51	9.111	1184	1185	1187	rBV2	573	513	2.01%	0.291%
52	9.227	1201	1204	1205	rBV2	623	482	1.89%	0.273%
53	9.276	1205	1212	1222	rVB2	6578	15467	60.69%	8.770%
54	9.495	1246	1248	1250	rBV3	908	881	3.46%	0.500%
55	9.855	1305	1307	1309	rBV2	492	569	2.23%	0.323%
56	9.971	1322	1326	1328	rBV3	586	917	3.60%	0.520%
57	10.093	1344	1346	1347	rBV	793	425	1.67%	0.241%
58	10.325	1379	1384	1392	rVB3	6513	12451	48.86%	7.060%
59	10.453	1403	1405	1407	rBV2	780	645	2.53%	0.366%
60	10.581	1424	1426	1428	rBV3	709	592	2.32%	0.336%
61	10.666	1438	1440	1443	rBV2	954	1115	4.38%	0.632%
62	10.751	1452	1454	1456	rBV3	705	484	1.90%	0.274%
63	11.129	1515	1516	1518	rBV2	767	613	2.41%	0.348%
64	11.233	1531	1533	1535	rBV3	551	566	2.22%	0.321%
65	11.300	1542	1544	1547	rVB4	627	507	1.99%	0.287%
66	11.373	1554	1556	1563	rVB5	846	1547	6.07%	0.877%
67	11.629	1594	1598	1603	rVV3	5389	9882	38.78%	5.603%
68	11.733	1613	1615	1619	rBV3	931	1159	4.55%	0.657%
69	11.818	1627	1629	1631	rVB2	673	551	2.16%	0.312%
70	12.330	1711	1713	1715	rBV3	554	426	1.67%	0.242%
71	12.599	1755	1757	1759	rBV2	757	457	1.79%	0.259%
72	12.647	1764	1765	1767	rBV2	450	447	1.75%	0.253%
73	12.690	1767	1772	1778	rVB2	6035	10409	40.84%	5.902%
74	12.964	1814	1817	1818	rVB2	443	451	1.77%	0.256%
75	12.995	1818	1822	1823	rBV3	422	465	1.82%	0.264%
76	13.068	1832	1834	1837	rBV3	648	834	3.27%	0.473%

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 Title : SFAM01.0

77	13.147	1845	1847	1850	rBV2	700	581	2.28%	0.329%
78	13.415	1889	1891	1896	rVB4	455	610	2.39%	0.346%
79	13.556	1906	1914	1920	rBV6	3106	7480	29.35%	4.241%
80	13.604	1920	1922	1924	rVV	705	497	1.95%	0.282%
81	13.629	1924	1926	1929	rVB2	590	646	2.53%	0.366%
82	13.665	1929	1932	1934	rBV3	660	879	3.45%	0.498%
83	13.708	1937	1939	1944	rVB5	658	675	2.65%	0.383%
84	13.763	1946	1948	1950	rVB	670	598	2.35%	0.339%
85	13.793	1950	1953	1955	rVB2	359	446	1.75%	0.253%
86	13.848	1955	1962	1967	rBV5	4094	7505	29.45%	4.256%
87	13.928	1973	1975	1977	rVB3	697	523	2.05%	0.297%
88	14.251	2026	2028	2031	rBV3	608	671	2.63%	0.380%
89	14.440	2056	2059	2061	rVB3	435	537	2.11%	0.304%
90	14.684	2096	2099	2102	rVB2	865	715	2.81%	0.405%
91	14.732	2102	2107	2109	rBV4	630	1114	4.37%	0.632%
92	14.915	2135	2137	2139	rBV2	437	456	1.79%	0.259%
93	14.964	2143	2145	2148	rBV3	633	626	2.46%	0.355%
94	15.116	2168	2170	2172	rVB2	689	563	2.21%	0.319%
95	15.446	2223	2224	2227	rBV3	784	534	2.10%	0.303%
96	15.500	2229	2233	2238	rVB8	907	1527	5.99%	0.866%
97	15.824	2284	2286	2288	rBV	635	569	2.23%	0.323%
98	16.074	2325	2327	2328	rBV	1051	673	2.64%	0.382%
99	16.330	2368	2369	2371	rBV2	796	689	2.70%	0.391%
100	16.653	2420	2422	2424	rBV3	655	742	2.91%	0.421%

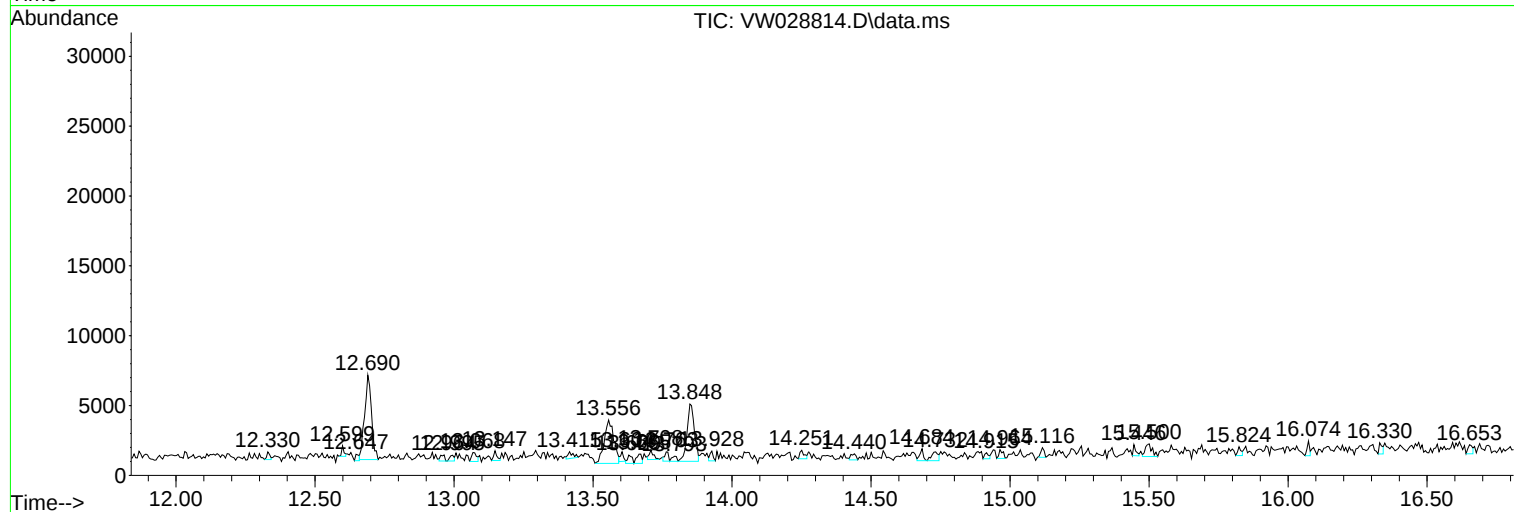
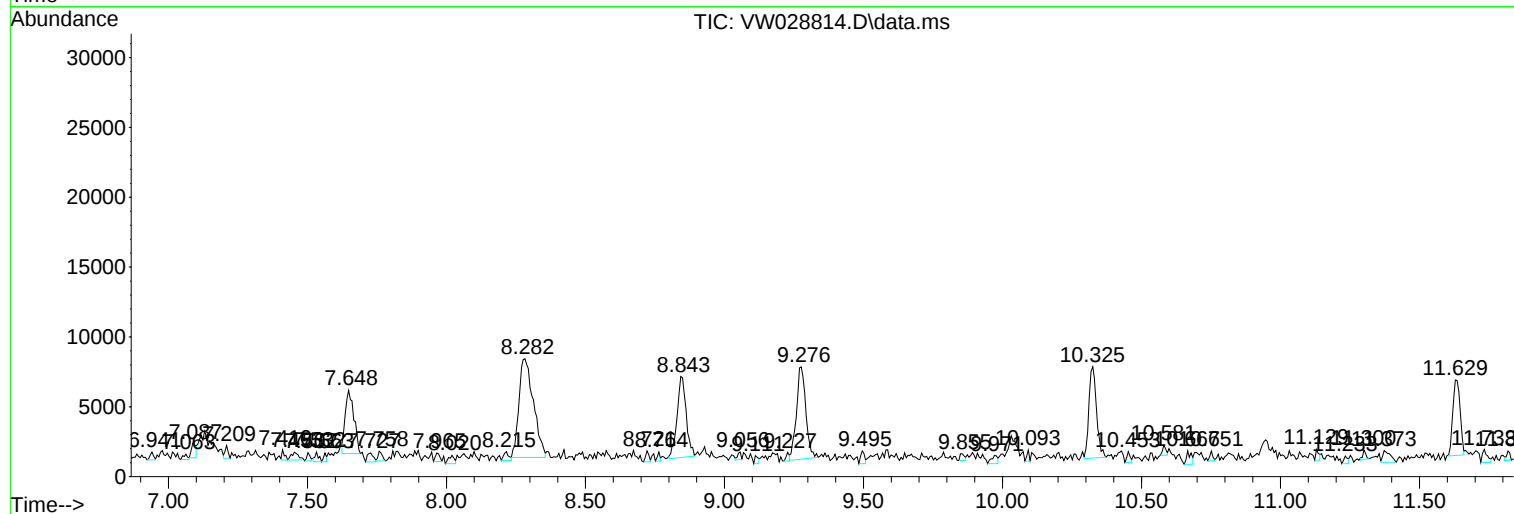
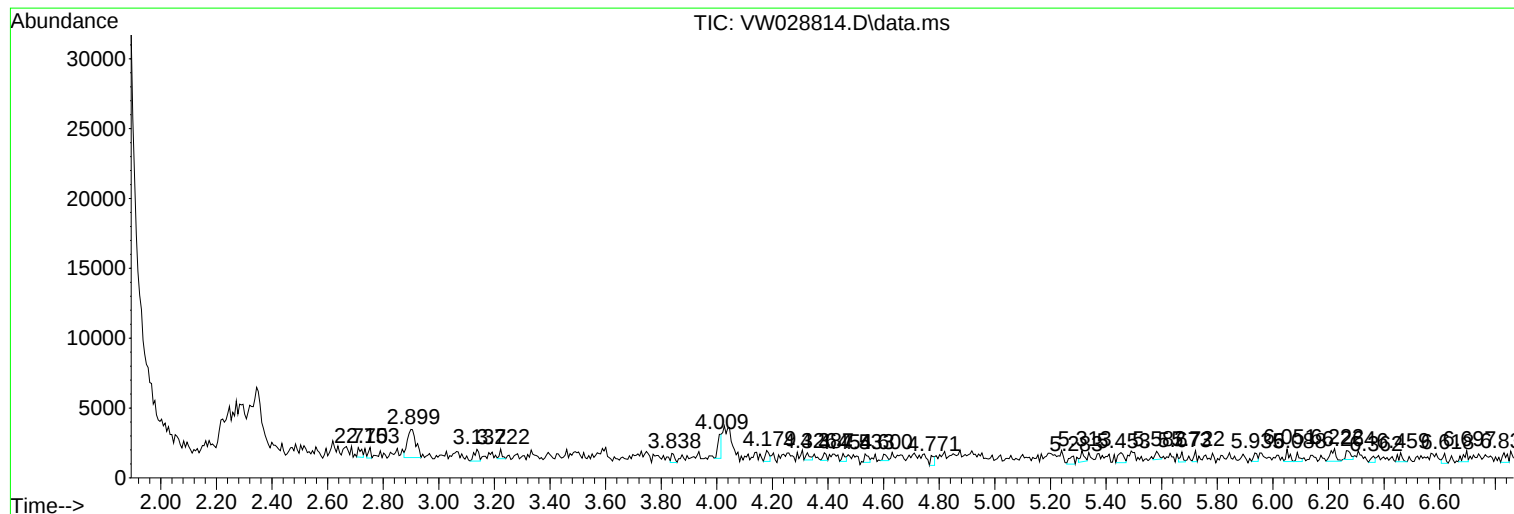
Sum of corrected areas: 176359

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050224SMA.M
Quant Title : SFAM01.0

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



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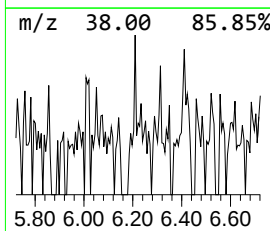
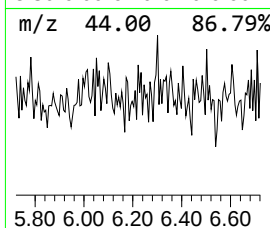
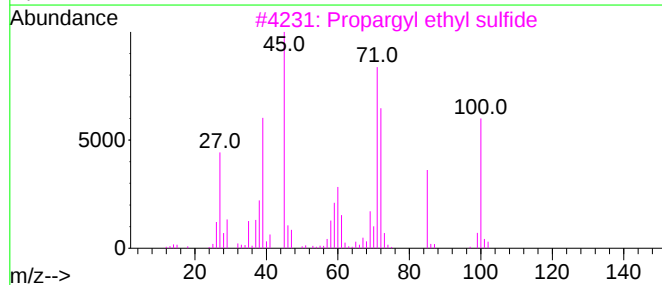
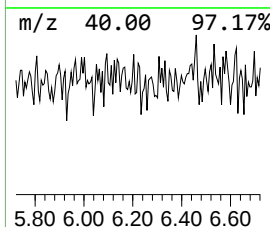
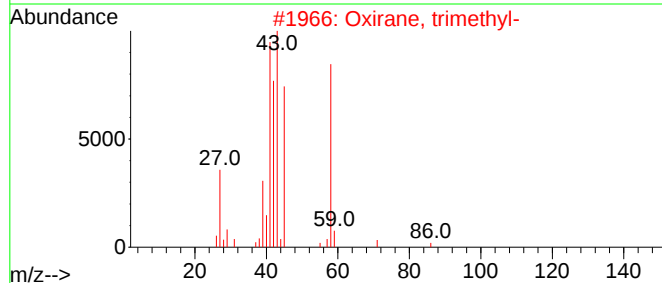
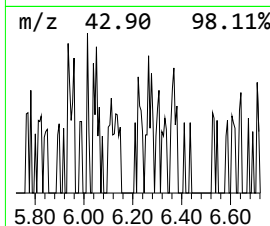
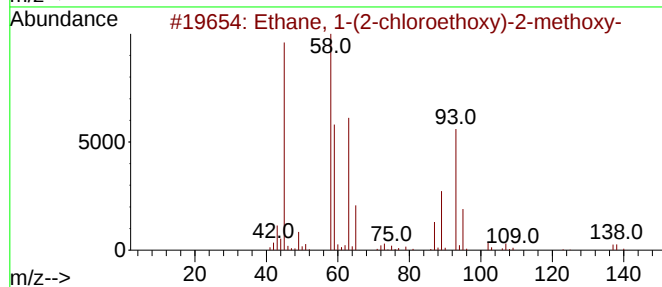
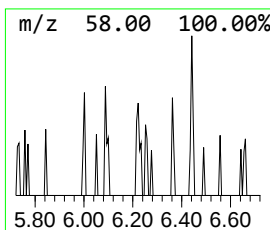
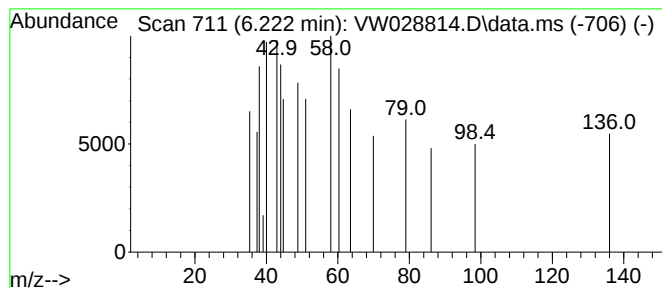
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050224SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.222	2.33 ug/L	1235	1,4-Difluorobenzene	8.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1-(2-chloroethoxy)-2-met...	138	C5H11ClO2	052808-36-3	9
2			Oxirane, trimethyl-	86	C5H10O	005076-19-7	7
3			Propargyl ethyl sulfide	100	C5H8S	007310-92-1	4
4			2-Hexynoic acid	112	C6H8O2	000764-33-0	4
5			Butanedioic acid, mercapto-	150	C4H6O4S	000070-49-5	4



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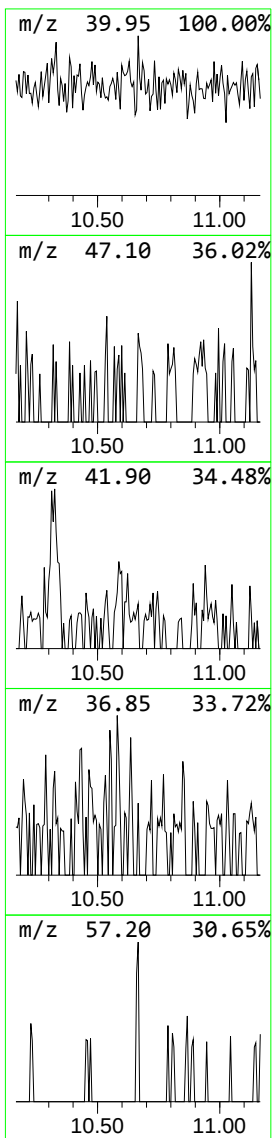
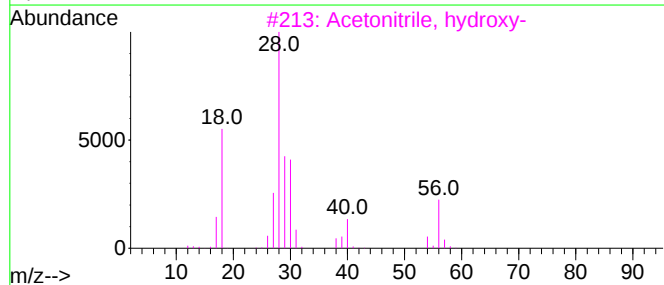
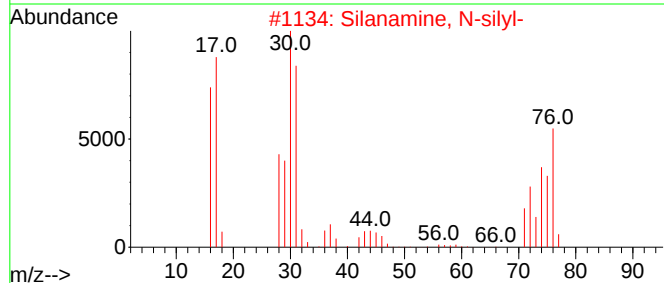
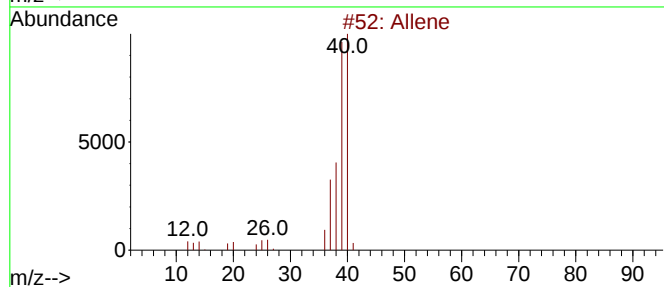
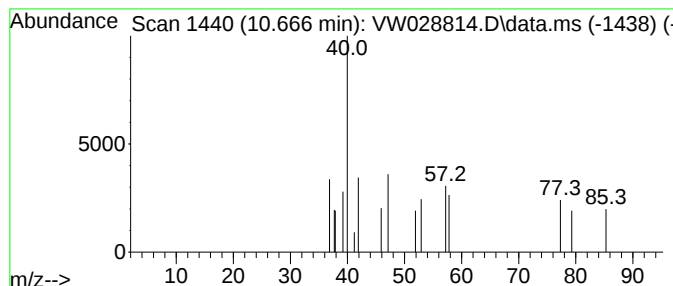
Instrument :
MSVOA_W
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050224SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.666	2.82 ug/L	1115	Chlorobenzene-d5		11.635	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Allene		40	C3H4	000463-49-0	23
2	Silanamine, N-silyl-		77	H7NSi2	005702-11-4	7
3	Acetonitrile, hydroxy-		57	C2H3NO	000107-16-4	5
4	Difluoramine		53	F2HN	010405-27-3	5
5	Propyne		40	C3H4	000074-99-7	5



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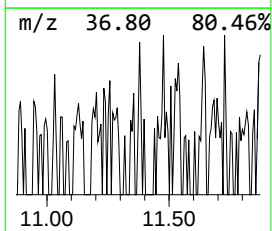
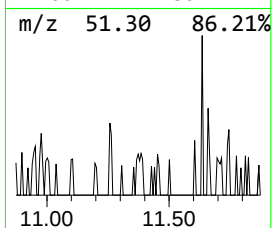
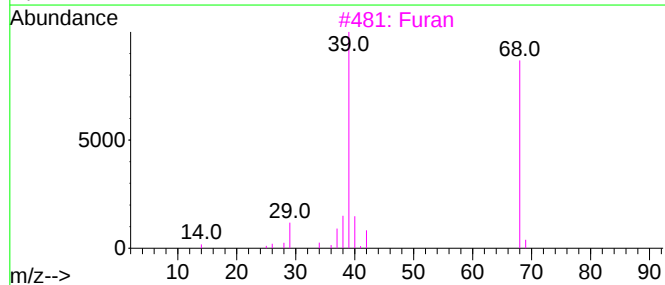
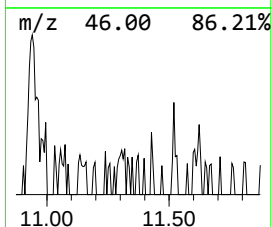
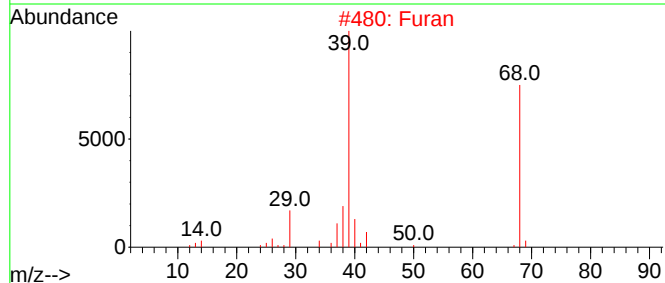
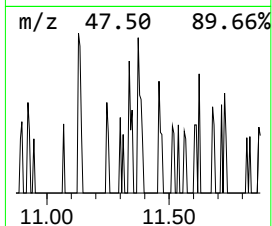
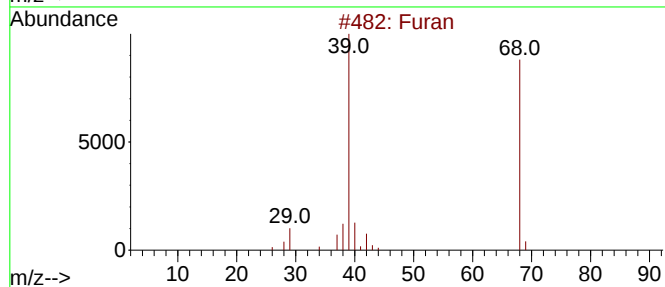
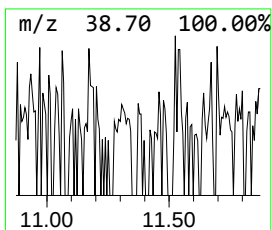
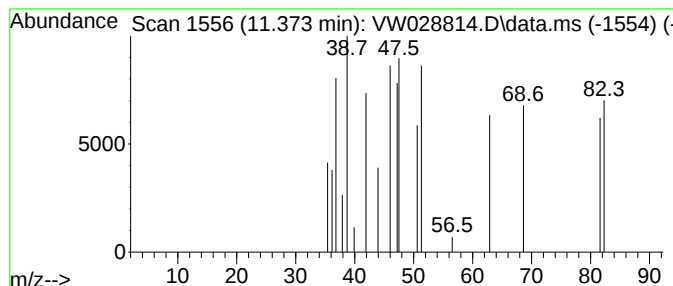
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050224SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.373	3.91 ug/L	1547	Chlorobenzene-d5	11.635

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan			68	C4H4O	000110-00-9	9
2	Furan			68	C4H4O	000110-00-9	9
3	Furan			68	C4H4O	000110-00-9	9
4	Furan, 2,5-dihydro-			70	C4H6O	001708-29-8	9
5	2-Butynal			68	C4H4O	001119-19-3	4



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BHBM2

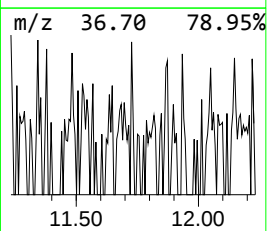
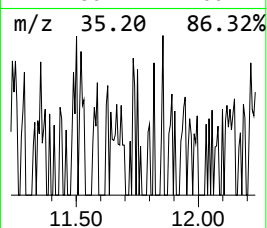
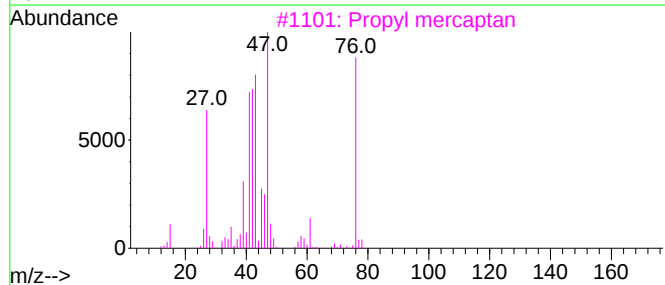
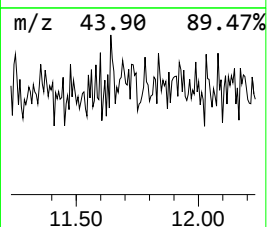
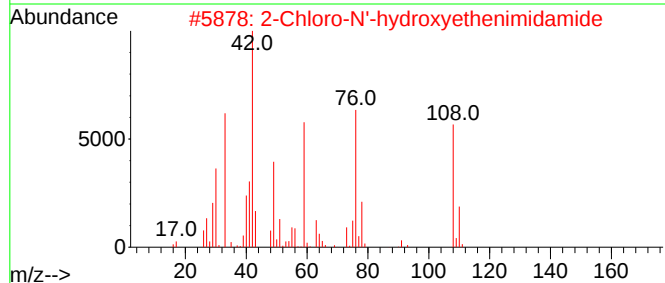
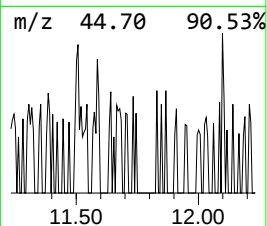
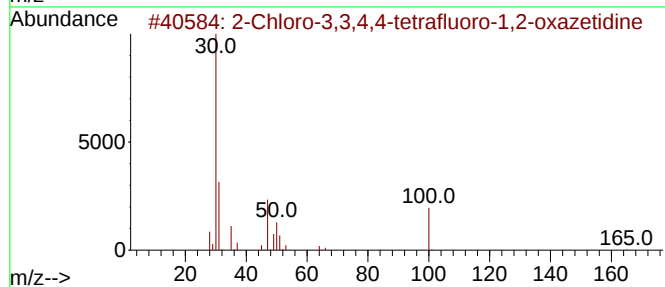
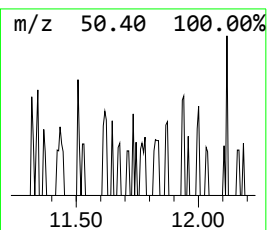
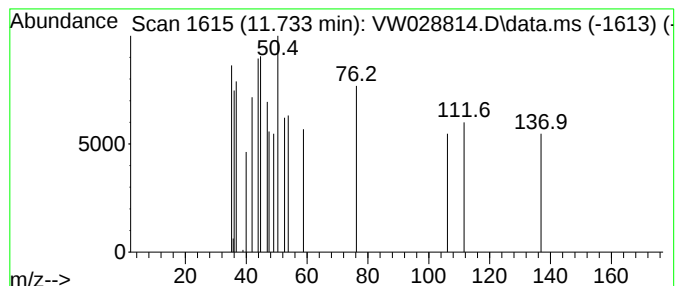
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Quant Title : SFAM01.0

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

Peak Number 4 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.733	2.93 ug/L	1159	Chlorobenzene-d5	11.635

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-3,3,4,4-tetrafluoro-1,2...	165	C2ClF4NO	1000305-82-6	9
2			2-Chloro-N'-hydroxyethenimidamide	108	C2H5ClN2O	1010461-06-2	9
3			Propyl mercaptan	76	C3H8S	000107-03-9	9
4			Propyl mercaptan	76	C3H8S	000107-03-9	9
5			1,3-Butadiyne	50	C4H2	000460-12-8	7



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BHBM2

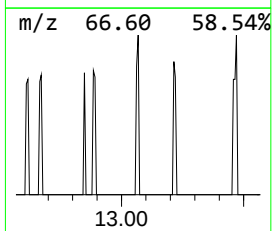
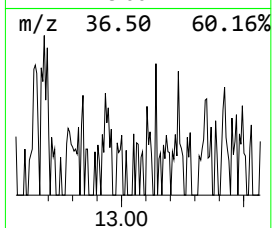
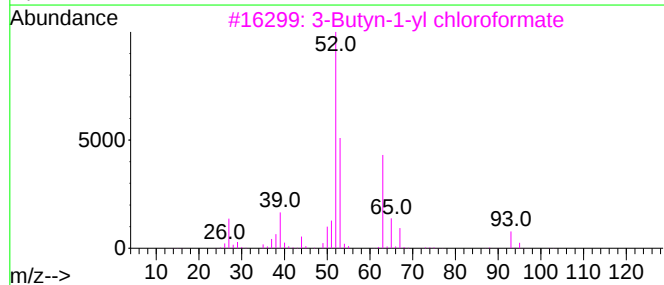
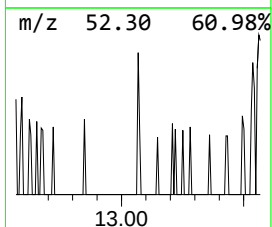
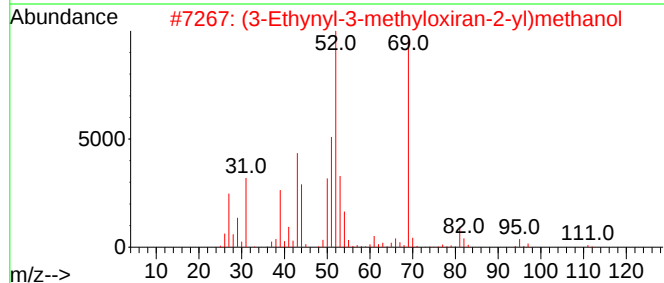
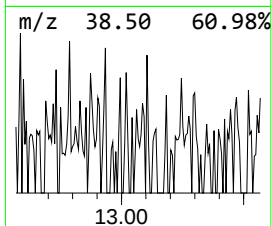
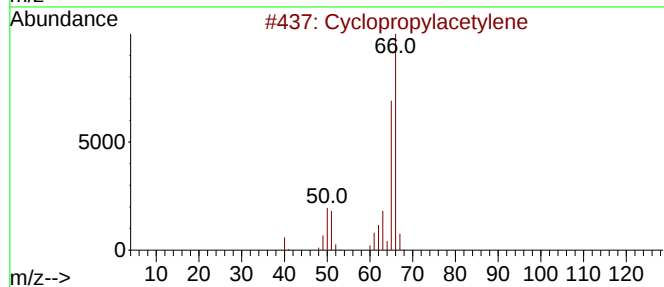
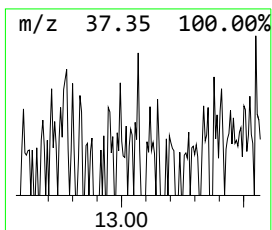
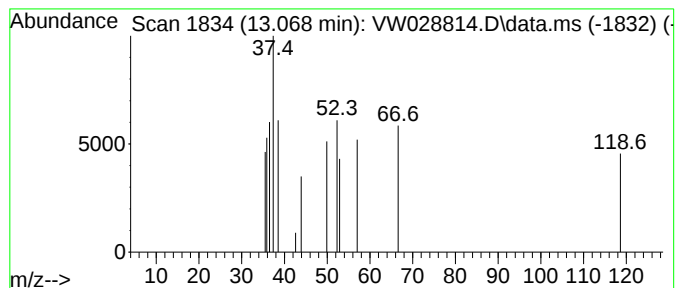
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.068	2.79 ug/L	834	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropylacetylene	66	C5H6	006746-94-7	3
2			(3-Ethynyl-3-methyloxiran-2-yl)m...	112	C6H8O2	1000191-14-3	2
3			3-Butyn-1-yl chloroformate	132	C5H5ClO2	190774-49-3	2
4			2-Butyne-1,4-diol, diformate	142	C6H6O4	036677-73-3	2
5			Pentanedinitrile, 2-methylene-	106	C6H6N2	001572-52-7	2



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BHBM2

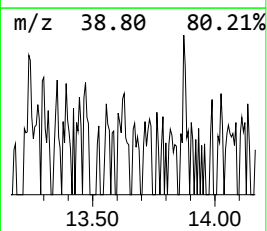
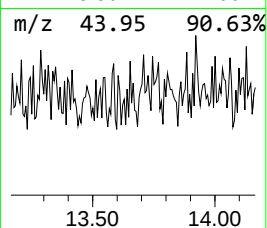
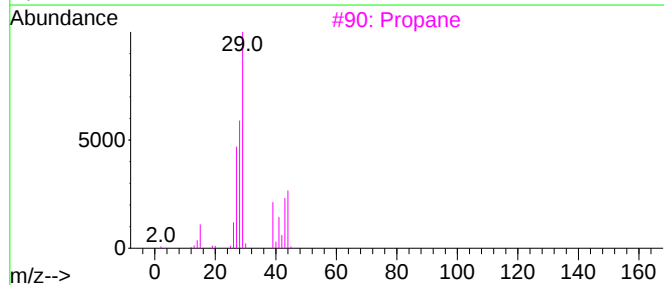
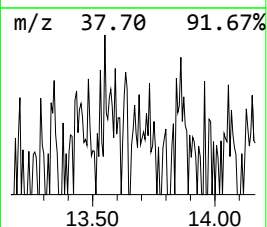
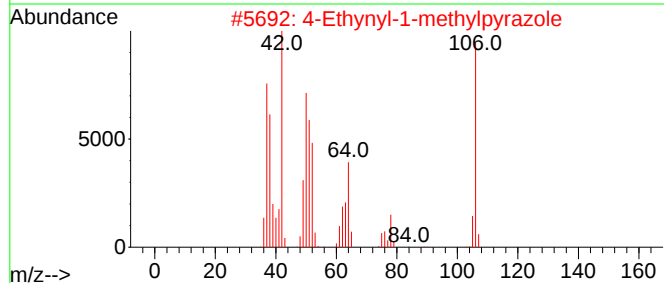
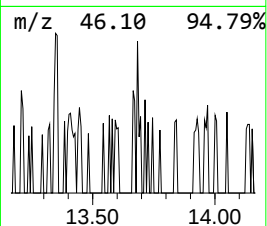
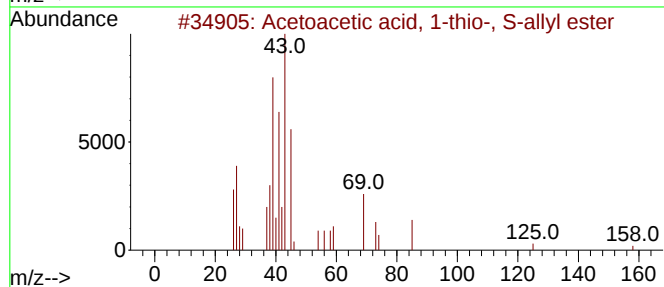
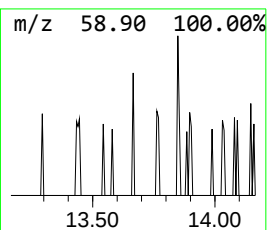
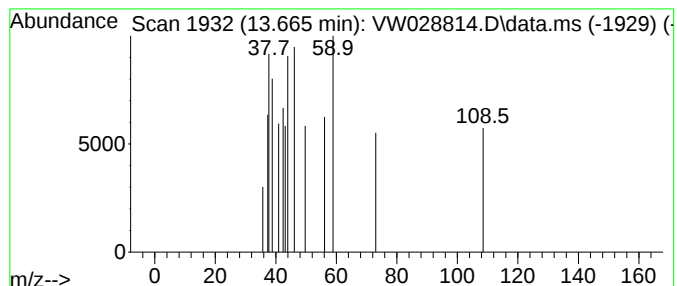
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Quant Title : SFAM01.0

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

Peak Number 6 unknown-06 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.665	2.94 ug/L	879	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetoacetic acid, 1-thio-, S-all...	158	C7H10O2S	015780-65-1	12
2			4-Ethynyl-1-methylpyrazole	106	C6H6N2	039806-89-8	9
3			Propane	44	C3H8	000074-98-6	4
4			Propene	42	C3H6	000115-07-1	4
5			1-Propene, 2-methyl-	56	C4H8	000115-11-7	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BHBM2

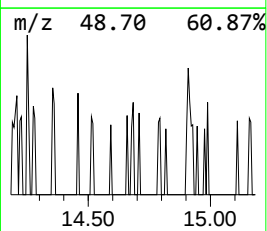
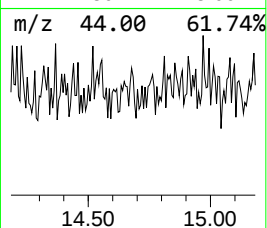
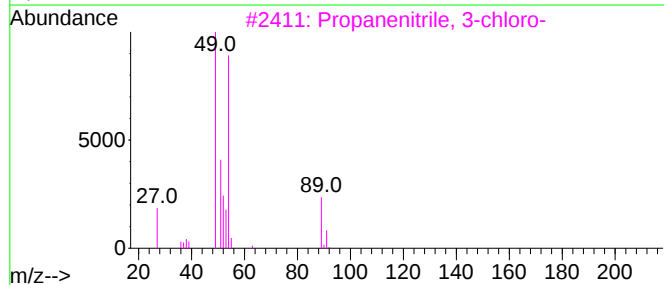
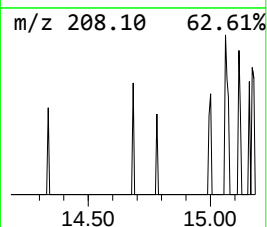
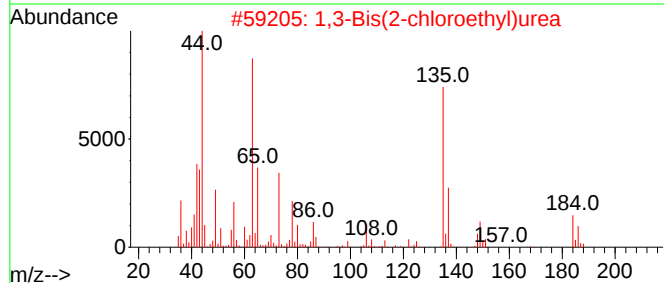
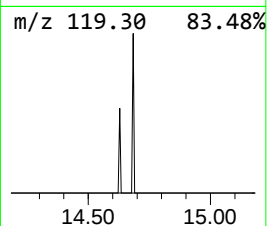
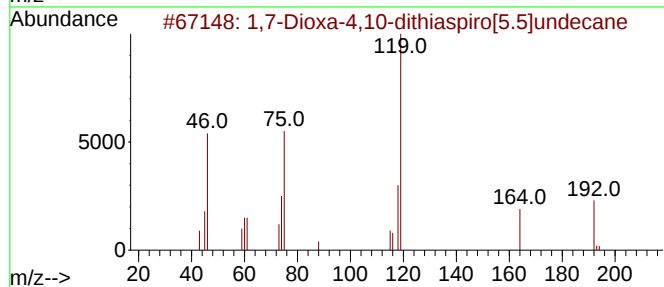
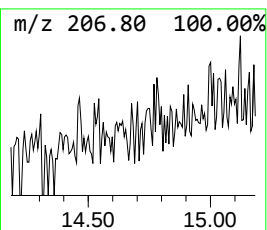
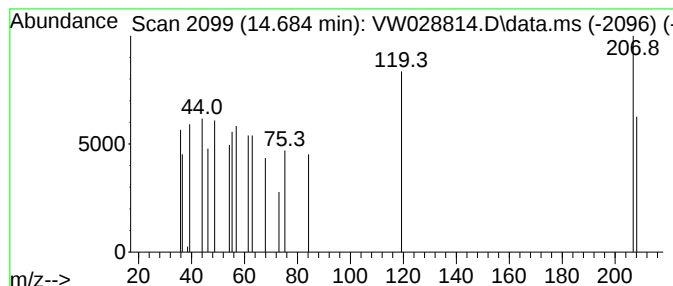
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Quant Title : SFAM01.0

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

Peak Number 7 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	2.39 ug/L	715	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,7-Dioxa-4,10-dithiaspiro[5.5]u...	192	C7H12O2S2	000180-86-9	9		
2	1,3-Bis(2-chloroethyl)urea	184	C5H10Cl2N2O	002214-72-4	9		
3	Propanenitrile, 3-chloro-	89	C3H4ClN	000542-76-7	9		
4	2-(Chloromethyl)-5-ethyl-1,3,4-o...	146	C5H7ClN2O	003914-45-2	9		
5	Methylene chloride	84	CH2Cl2	000075-09-2	7		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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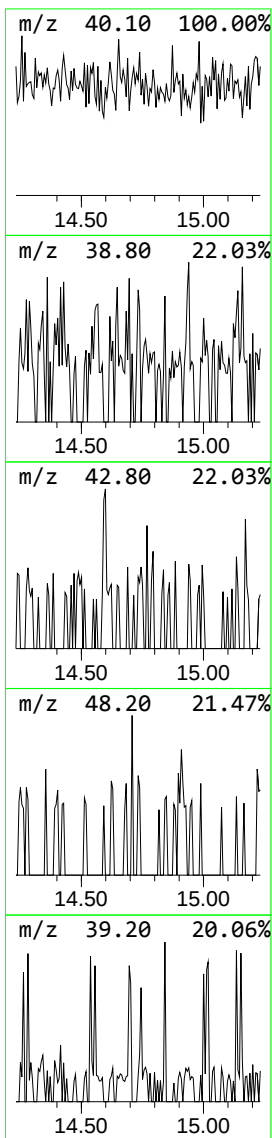
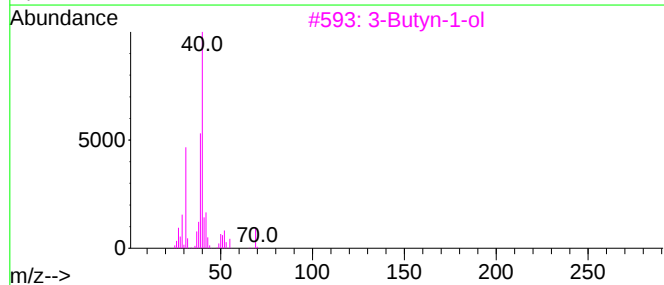
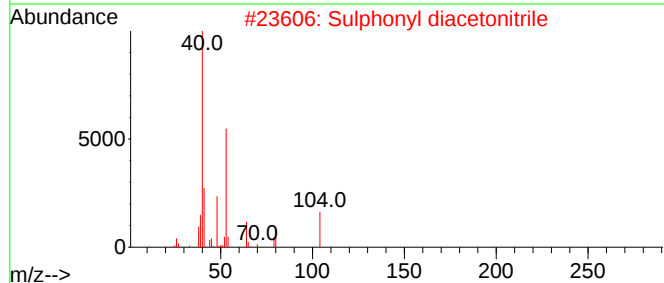
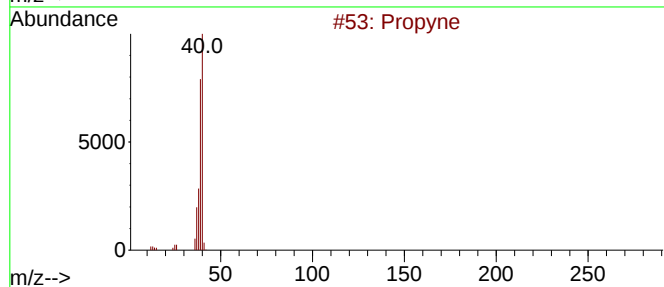
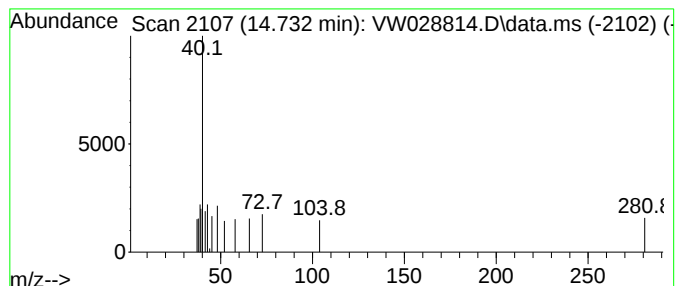
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Quant Title : SFAM01.0

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

Peak Number 8 unknown-08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.732	3.72 ug/L	1114	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propyne	40	C3H4	000074-99-7	9
2			Sulphonyl diacetonitrile	144	C4H4N2O2S	1000342-33-0	9
3			3-Butyn-1-ol	70	C4H6O	000927-74-2	9
4			Pyridine, 2-amino-3,5-dinitro-	184	C5H4N4O4	003073-30-1	4
5			cis-Aconitic anhydride	156	C6H4O5	006318-55-4	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BHBM2

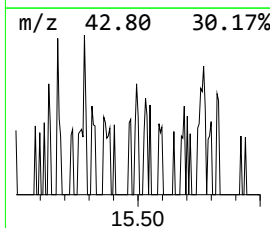
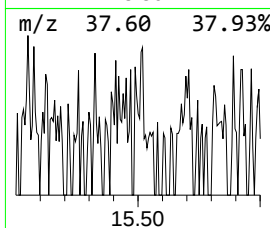
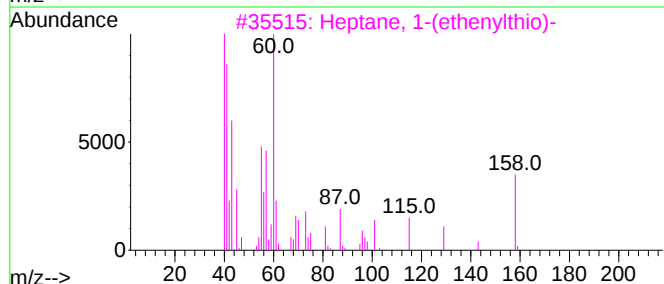
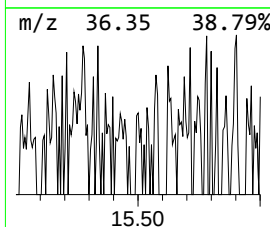
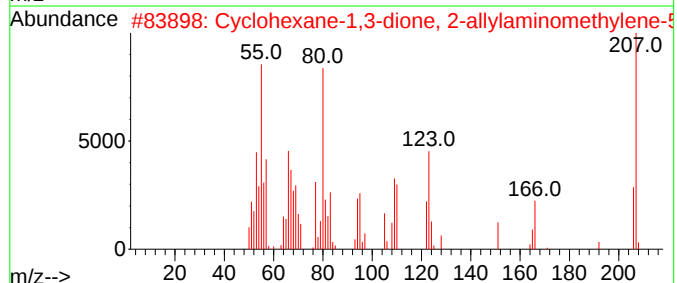
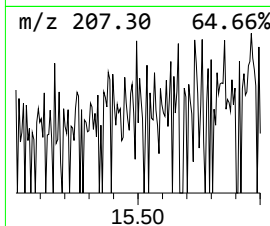
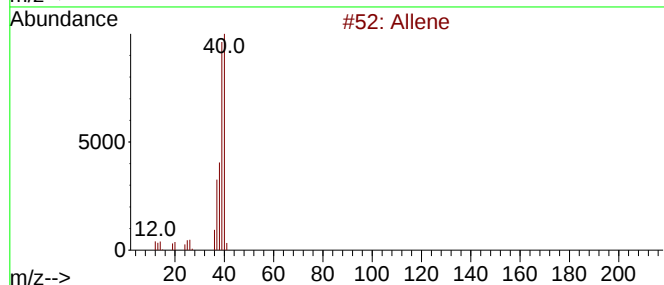
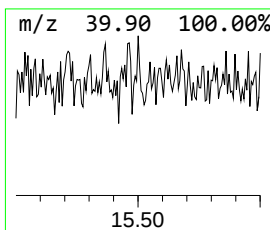
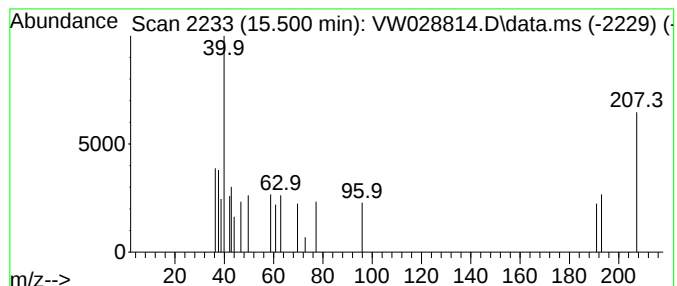
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.501	5.10 ug/L	1527	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Allene	40	C3H4	000463-49-0	5
2			Cyclohexane-1,3-dione, 2-allylam...	207	C12H17NO2	104926-37-6	4
3			Heptane, 1-(ethenylthio)-	158	C9H18S	021961-05-7	4
4			Silamine, N-silyl-	77	H7NSi2	005702-11-4	4
5			6-Nitro-8-methoxy-2H-chromene	207	C10H9NO4	062063-07-4	2



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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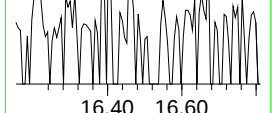
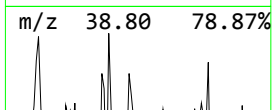
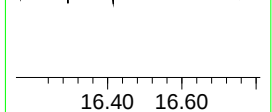
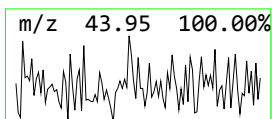
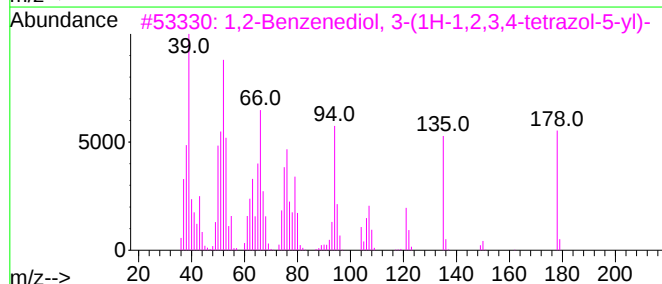
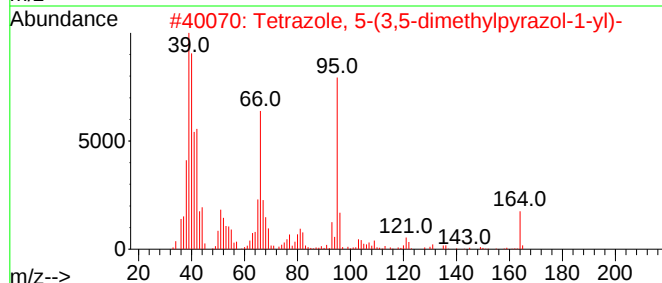
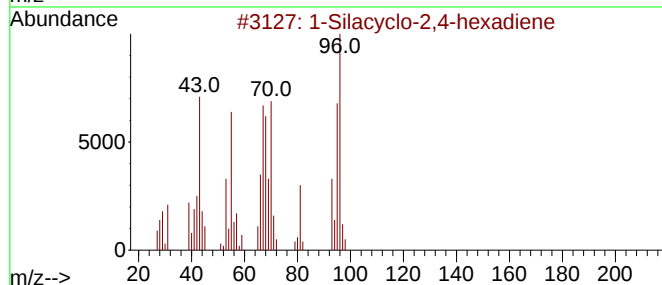
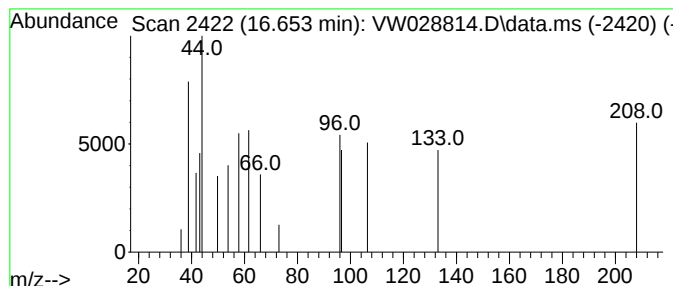
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-10 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.653	2.48 ug/L	742	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Silacyclo-2,4-hexadiene	96	C5H8Si	052023-17-3	9
2			Tetrazole, 5-(3,5-dimethylpyrazo...	164	C6H8N6	297763-48-5	9
3			1,2-Benzenediol, 3-(1H-1,2,3,4-t...	178	C7H6N4O2	1000362-63-3	9
4			4-Chloro-N-(1,2-dihydro-2-pyridy...	268	C11H9ClN2O2S	001213-38-3	9
5			Hydrazinecarboxamide, 2-(2,6-cyc...	179	C9H13N3O	068344-44-5	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW052424\
Data File : VW028814.D
Acq On : 26 May 2024 02:04
Operator : SY/MD
Sample : P2590-18
Misc : 6.11g/10mL/MSVOA_W/SOIL/A
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BHBM2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050224SMA.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
unknown-01	6.222	2.3	ug/L	1235	1	8.843	13251	25.0
unknown-02	10.666	2.8	ug/L	1115	2	11.635	9882	25.0
unknown-03	11.373	3.9	ug/L	1547	2	11.635	9882	25.0
unknown-04	11.733	2.9	ug/L	1159	2	11.635	9882	25.0
unknown-05	13.068	2.8	ug/L	834	3	13.556	7480	25.0
unknown-06	13.665	2.9	ug/L	879	3	13.556	7480	25.0
unknown-07	14.684	2.4	ug/L	715	3	13.556	7480	25.0
unknown-08	14.732	3.7	ug/L	1114	3	13.556	7480	25.0
unknown-09	15.501	5.1	ug/L	1527	3	13.556	7480	25.0
unknown-10	16.653	2.5	ug/L	742	3	13.556	7480	25.0