

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
 Data File : VW022470.D
 Acq On : 08 Apr 2022 17:43
 Operator : SY/VA
 Sample : N2293-18
 Misc : 6.67g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNP6

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: VW022470.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	36	rVB	3465382	5495398	12.78%	7.982%
2	2.355	68	74	84	rBV	149619	257134	0.60%	0.373%
3	2.495	92	97	104	rBV	27159	59608	0.14%	0.087%
4	2.623	114	118	119	rBV4	12572	14118	0.03%	0.021%
5	2.885	155	161	171	rVB	90759	196445	0.46%	0.285%
6	4.019	337	347	359	rVV2	158685	482318	1.12%	0.701%
7	4.123	359	364	373	rVV2	14923	43107	0.10%	0.063%
8	4.214	373	379	387	rVV10	6728	18856	0.04%	0.027%
9	4.391	399	408	417	rBV2	11155	36077	0.08%	0.052%
10	4.916	483	494	495	rBV6	6796	13749	0.03%	0.020%
11	5.190	527	539	553	rVB4	15656	56376	0.13%	0.082%
12	5.556	588	599	607	rBV8	6330	21319	0.05%	0.031%
13	5.946	652	663	676	rVB3	9780	31379	0.07%	0.046%
14	6.604	762	771	776	rBV8	4223	11460	0.03%	0.017%
15	6.988	823	834	841	rBV10	12514	39921	0.09%	0.058%
16	7.074	841	848	861	rVV	44446	148497	0.35%	0.216%
17	7.171	861	864	875	rVB2	10462	26350	0.06%	0.038%
18	7.537	912	924	926	rBV8	3562	12162	0.03%	0.018%
19	7.653	930	943	954	rBV	273981	676094	1.57%	0.982%
20	7.915	978	986	989	rBV8	5973	14534	0.03%	0.021%
21	8.159	1019	1026	1031	rBV7	5591	13153	0.03%	0.019%
22	8.281	1031	1046	1047	rBV	502694	1003465	2.33%	1.458%
23	8.323	1047	1053	1064	rVB	1936505	4557865	10.60%	6.620%
24	8.567	1089	1093	1101	rVB5	9080	21003	0.05%	0.031%
25	8.695	1108	1114	1127	rVB4	15126	39334	0.09%	0.057%
26	8.842	1129	1138	1152	rBV	485528	970219	2.26%	1.409%
27	9.049	1164	1172	1179	rBV3	23345	61683	0.14%	0.090%
28	9.275	1200	1209	1217	rBV	389966	814517	1.89%	1.183%
29	9.506	1240	1247	1256	rVB	19538	44493	0.10%	0.065%
30	9.811	1292	1297	1302	rBV6	8524	13246	0.03%	0.019%
31	9.963	1315	1322	1328	rBV	350389	652670	1.52%	0.948%
32	10.037	1328	1334	1341	rVV	201653	373318	0.87%	0.542%
33	10.091	1341	1343	1354	rVB7	9237	18250	0.04%	0.027%
34	10.207	1354	1362	1365	rBV5	10661	20890	0.05%	0.030%
35	10.244	1365	1368	1373	rVB3	9353	15905	0.04%	0.023%
36	10.323	1373	1381	1386	rBV	730748	1303813	3.03%	1.894%

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

37	10.384	1386	1391	1401	rVV2	740506	1335941	3.11%	1.940%
38	10.469	1401	1405	1415	rVV5	24835	73670	0.17%	0.107%
39	10.579	1415	1423	1430	rVV	126551	217311	0.51%	0.316%
40	10.719	1438	1446	1459	rVB	275879	513390	1.19%	0.746%
41	10.854	1460	1468	1473	rBV3	10977	22912	0.05%	0.033%
42	10.927	1473	1480	1492	rBV2	202618	383469	0.89%	0.557%
43	11.061	1495	1502	1515	rVB2	36207	70313	0.16%	0.102%
44	11.177	1515	1521	1527	rVB2	8959	14242	0.03%	0.021%
45	11.292	1533	1540	1548	rBV	102385	191767	0.45%	0.279%
46	11.481	1566	1571	1576	rVB7	6270	10215	0.02%	0.015%
47	11.555	1579	1583	1588	rBV6	9617	15625	0.04%	0.023%
48	11.658	1588	1600	1609	rBV	26848119	42996909	100.00%	62.452%
49	11.725	1609	1611	1619	rVB	30473	46866	0.11%	0.068%
50	11.835	1619	1629	1642	rVB	53744	127691	0.30%	0.185%
51	11.969	1642	1651	1653	rBV7	5488	13520	0.03%	0.020%
52	12.164	1677	1683	1691	rVB	50867	94706	0.22%	0.138%
53	12.341	1703	1712	1718	rVB	5159	11315	0.03%	0.016%
54	12.463	1728	1732	1738	rVB5	9179	16075	0.04%	0.023%
55	12.603	1747	1755	1758	rBV5	12169	23480	0.05%	0.034%
56	12.658	1758	1764	1766	rBV	150614	249106	0.58%	0.362%
57	12.689	1766	1769	1775	rVV	286028	454887	1.06%	0.661%
58	12.743	1775	1778	1784	rVB2	20859	31632	0.07%	0.046%
59	12.798	1784	1787	1791	rVB	7534	11076	0.03%	0.016%
60	12.890	1794	1802	1807	rBV	142227	267539	0.62%	0.389%
61	12.987	1814	1818	1823	rVB	21709	34932	0.08%	0.051%
62	13.109	1832	1838	1845	rBV4	30754	64594	0.15%	0.094%
63	13.182	1847	1850	1855	rVB5	6556	9616	0.02%	0.014%
64	13.243	1855	1860	1865	rBV	26545	45504	0.11%	0.066%
65	13.292	1865	1868	1875	rVB9	9013	16480	0.04%	0.024%
66	13.365	1875	1880	1886	rBV4	29292	60158	0.14%	0.087%
67	13.493	1895	1901	1904	rBV8	8289	17005	0.04%	0.025%
68	13.554	1904	1911	1920	rVV	755629	1305722	3.04%	1.897%
69	13.640	1920	1925	1934	rVB2	126651	221154	0.51%	0.321%
70	13.713	1934	1937	1941	rBV5	8075	13762	0.03%	0.020%
71	13.749	1941	1943	1947	rVB5	6599	9027	0.02%	0.013%
72	13.798	1947	1951	1953	rBV3	6462	9399	0.02%	0.014%
73	13.853	1953	1960	1970	rVB3	790150	1484203	3.45%	2.156%
74	13.944	1970	1975	1979	rBV3	14550	22797	0.05%	0.033%
75	13.999	1979	1984	1988	rBV6	10147	20285	0.05%	0.029%
76	14.200	2009	2017	2026	rBV6	20460	47252	0.11%	0.069%

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 Title : SFAM01.0

77	14.335	2034	2039	2045	rBV3	16741	28858	0.07%	0.042%
78	14.420	2048	2053	2055	rVV5	9836	15254	0.04%	0.022%
79	14.457	2055	2059	2069	rVV7	18872	50710	0.12%	0.074%
80	14.572	2072	2078	2083	rVV7	14162	37178	0.09%	0.054%
81	14.627	2083	2087	2092	rVB8	8898	15753	0.04%	0.023%
82	14.682	2092	2096	2099	rBV6	7566	12835	0.03%	0.019%
83	14.731	2099	2104	2109	rVV5	23149	46421	0.11%	0.067%
84	14.786	2109	2113	2120	rVV2	19484	37328	0.09%	0.054%
85	15.060	2150	2158	2165	rBV7	16280	43110	0.10%	0.063%
86	15.365	2203	2208	2212	rVV2	8695	16480	0.04%	0.024%
87	15.414	2212	2216	2220	rVV4	7617	12423	0.03%	0.018%
88	15.487	2220	2228	2236	rVB8	15625	42150	0.10%	0.061%
89	15.603	2241	2247	2251	rBV8	12215	23631	0.05%	0.034%
90	15.645	2251	2254	2262	rVB10	6976	13727	0.03%	0.020%
91	15.731	2262	2268	2277	rBV4	43265	99631	0.23%	0.145%
92	15.816	2277	2282	2288	rVV6	6909	19713	0.05%	0.029%
93	15.938	2298	2302	2310	rVB6	7049	12507	0.03%	0.018%
94	16.170	2335	2340	2347	rVB10	6326	14174	0.03%	0.021%
95	16.273	2351	2357	2361	rVV8	3849	9833	0.02%	0.014%
96	16.365	2370	2372	2379	rVV8	4530	8992	0.02%	0.013%
97	16.438	2379	2384	2386	rVV5	6430	12891	0.03%	0.019%
98	16.474	2388	2390	2395	rVV6	5314	9196	0.02%	0.013%
99	16.633	2409	2416	2424	rBV7	14155	35716	0.08%	0.052%
100	16.724	2424	2431	2441	rVB2	31117	74642	0.17%	0.108%

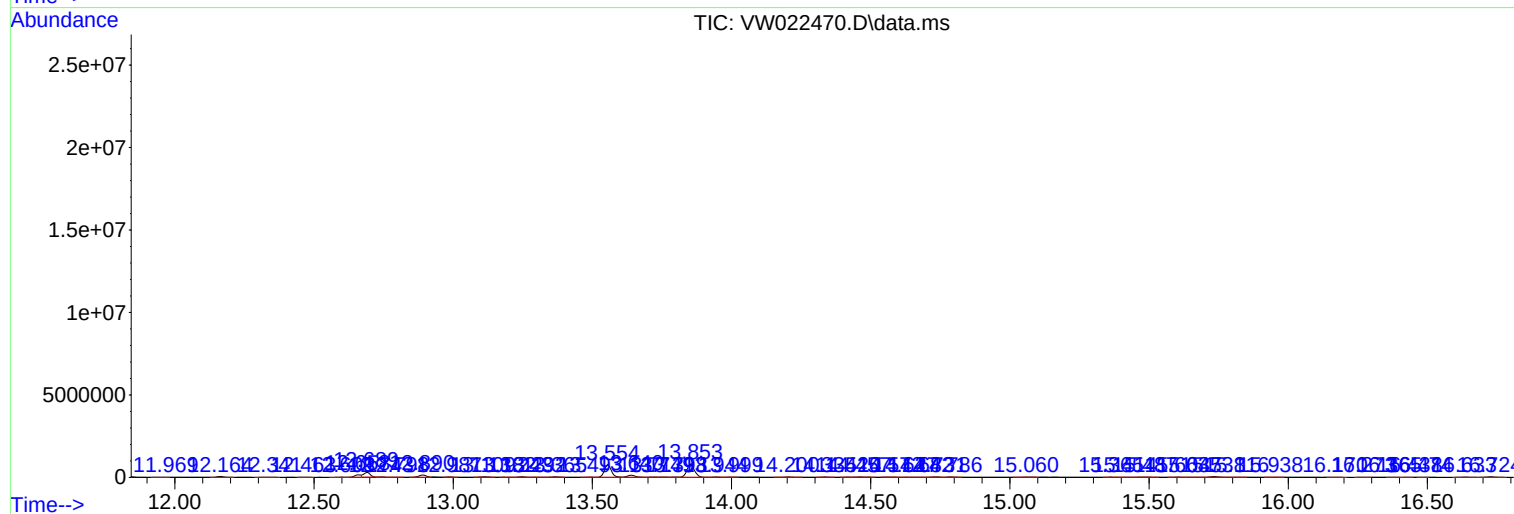
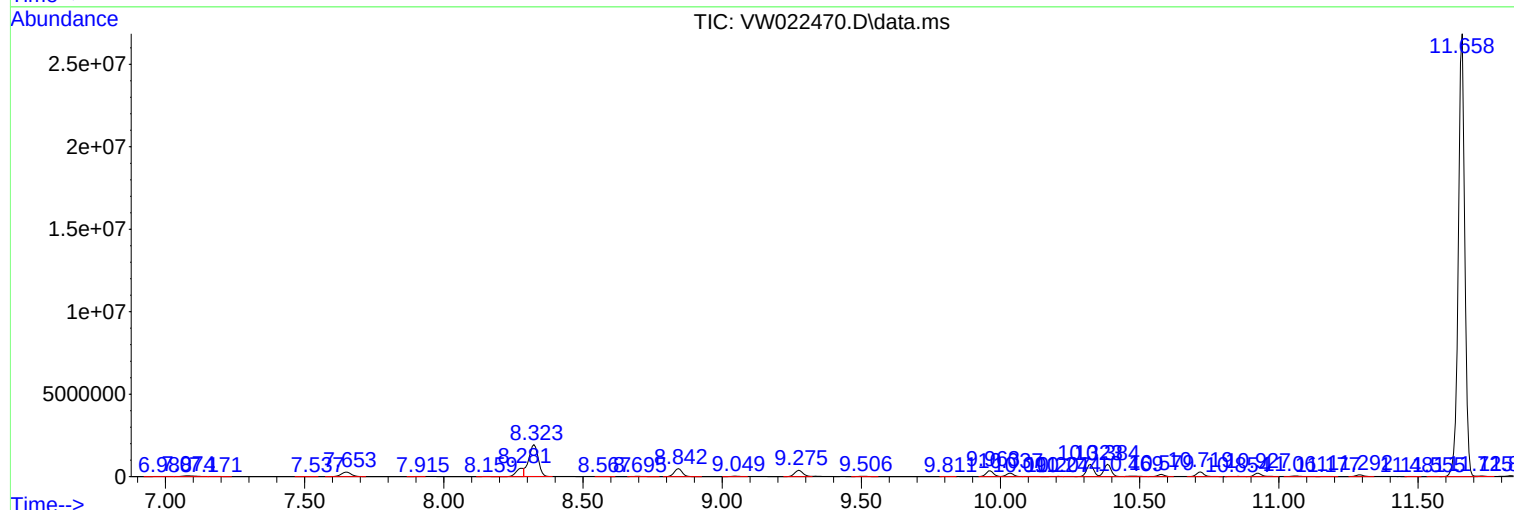
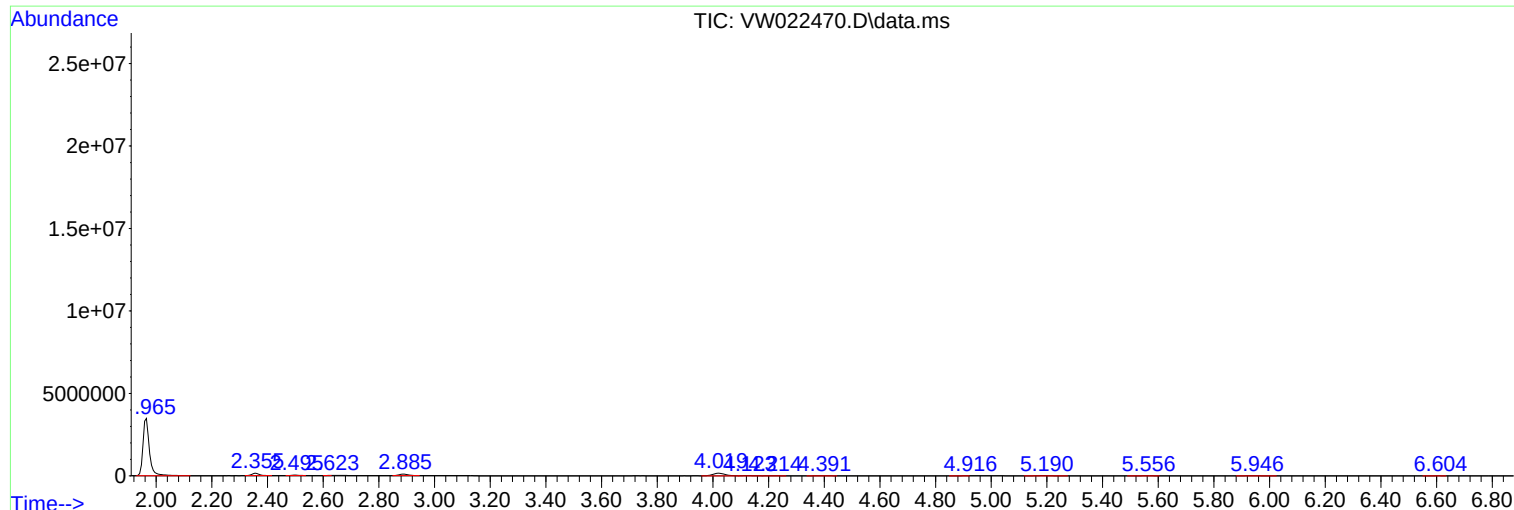
Sum of corrected areas: 68847426

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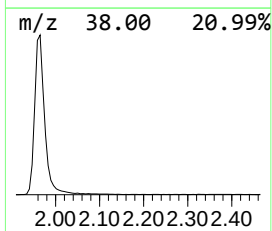
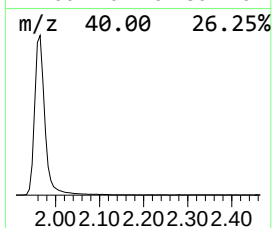
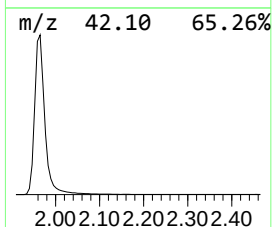
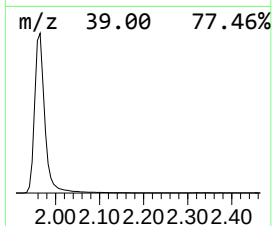
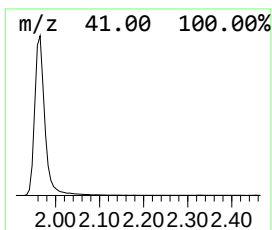
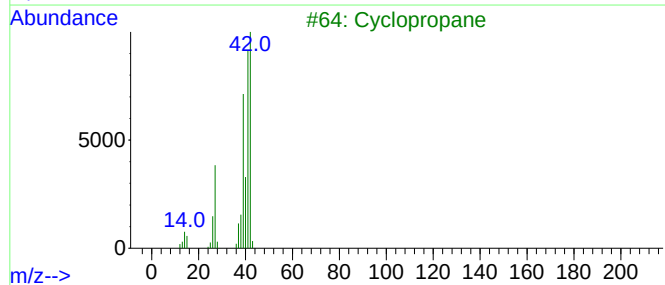
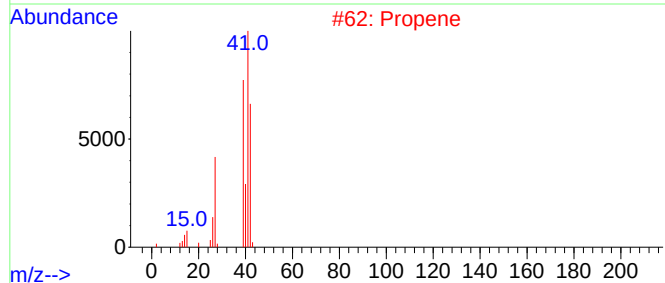
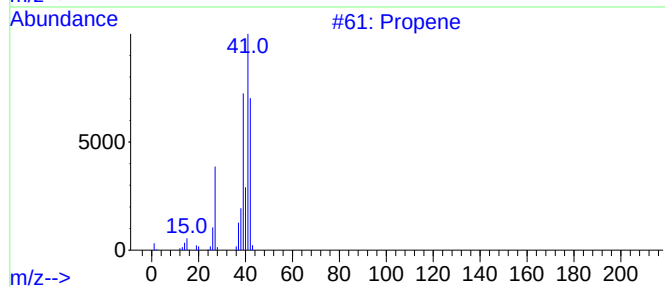
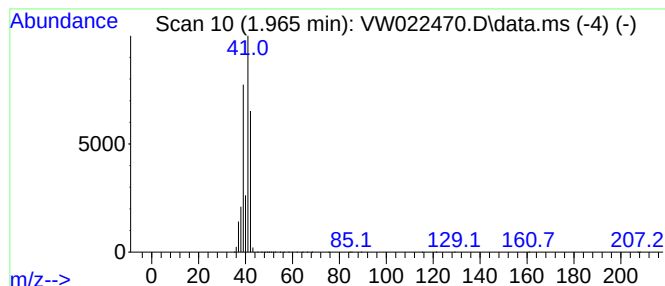
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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	141.60 ug/L	5495400	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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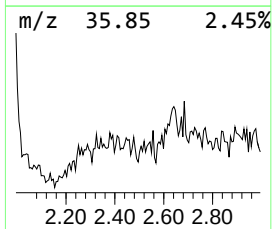
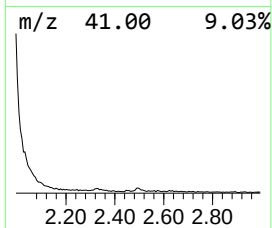
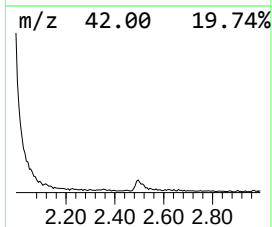
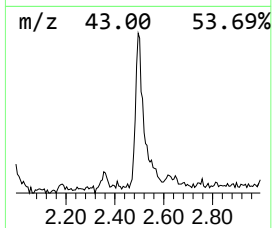
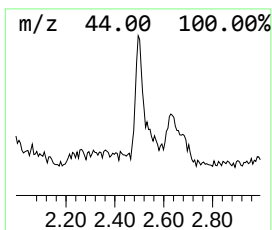
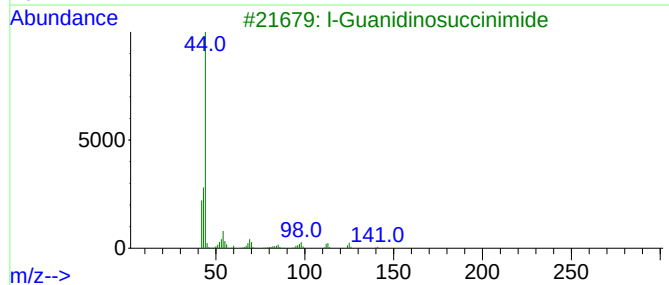
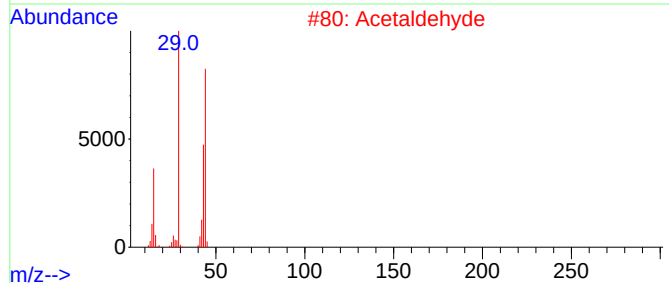
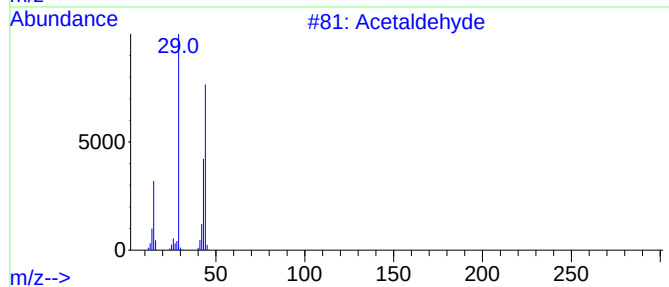
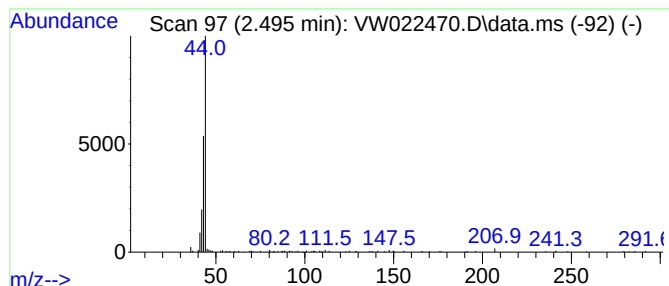
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.495	1.54 ug/L	59608	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetaldehyde	44	C2H4O	000075-07-0	9
2			Acetaldehyde	44	C2H4O	000075-07-0	9
3			1-Guanidinosuccinimide	141	C5H7N3O2	1000130-20-8	9
4			2-Propanamine	59	C3H9N	000075-31-0	9
5			Ethylamine, 1-(2-thenyl)-, hydro...	141	C7H11NS	053632-92-1	7



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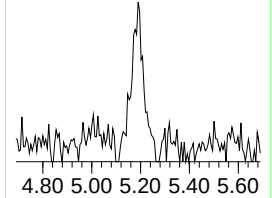
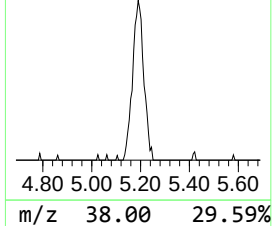
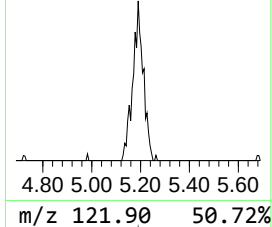
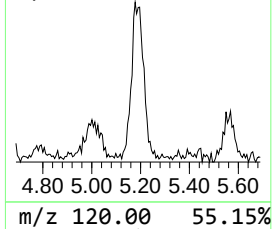
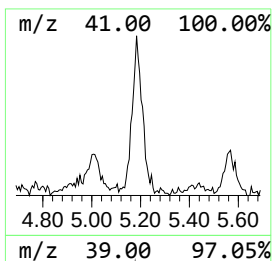
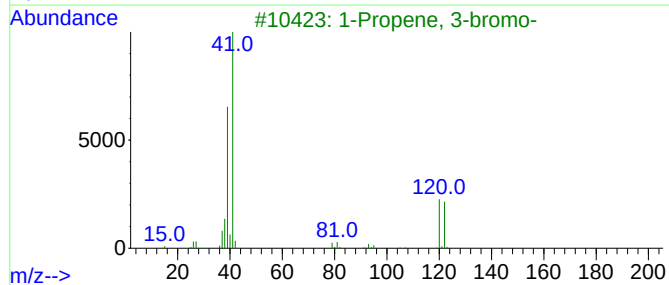
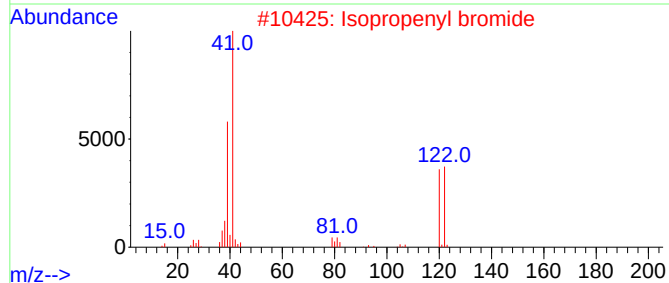
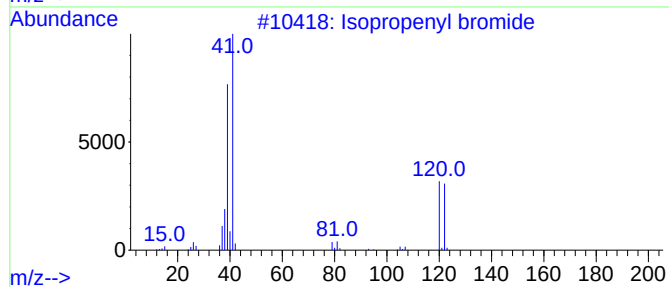
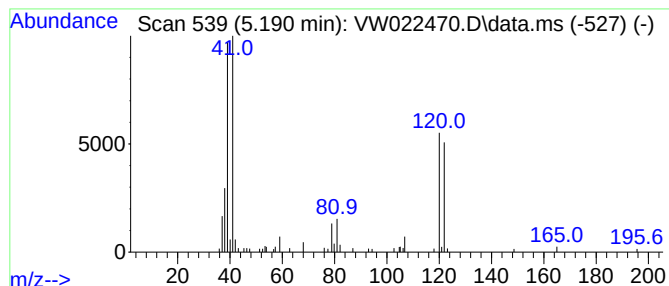
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Isopropenyl bromide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.190	1.45 ug/L	56376	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropenyl bromide	120	C3H5Br	000557-93-7	90
2			Isopropenyl bromide	120	C3H5Br	000557-93-7	80
3			1-Propene, 3-bromo-	120	C3H5Br	000106-95-6	72
4			1-Methoxy-1-buten-3-yne	82	C5H6O	002798-73-4	64
5			Chloro(prop-2-yn-1-yloxy)methanone	118	C4H3ClO2	035718-08-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022470.D
Acq On : 08 Apr 2022 17:43
Operator : SY/VA
Sample : N2293-18
Misc : 6.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP6

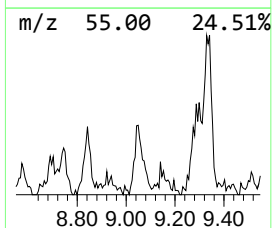
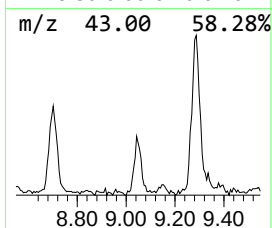
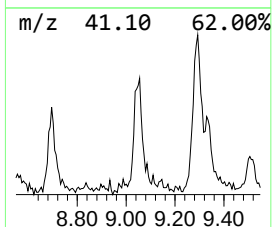
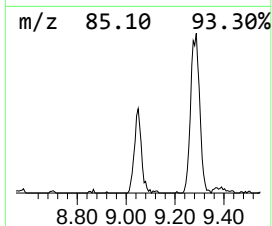
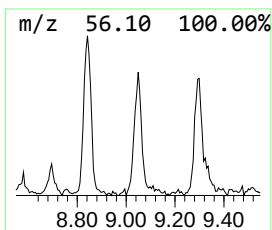
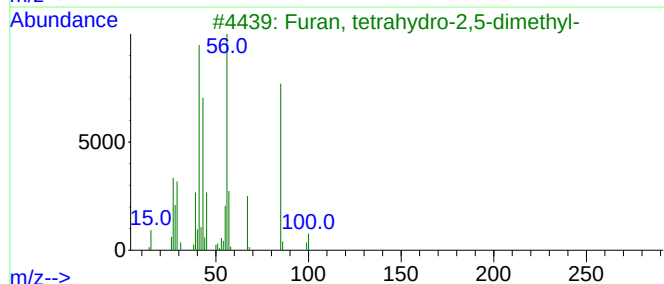
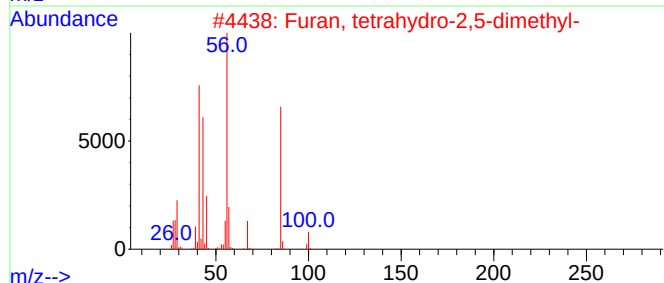
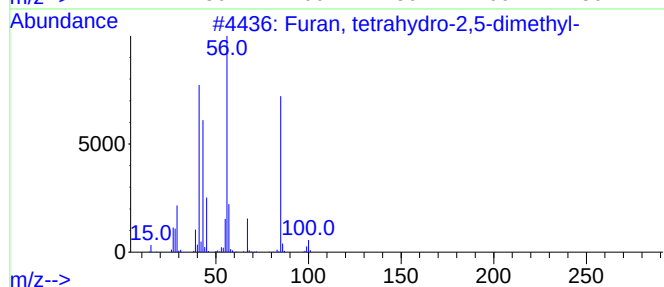
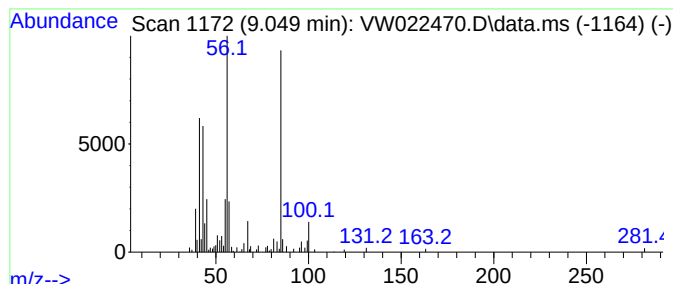
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Quant Title : SFAM01.0

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

Peak Number 4 Furan, tetrahydro-2,5-dimet... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.049	1.59 ug/L	61683	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	91
2			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	91
3			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	87
4			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	002144-41-4	87
5			2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022470.D
Acq On : 08 Apr 2022 17:43
Operator : SY/VA
Sample : N2293-18
Misc : 6.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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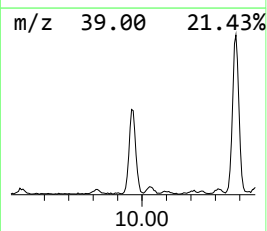
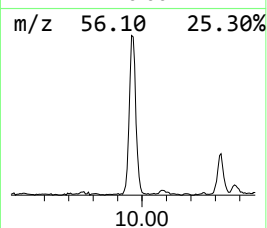
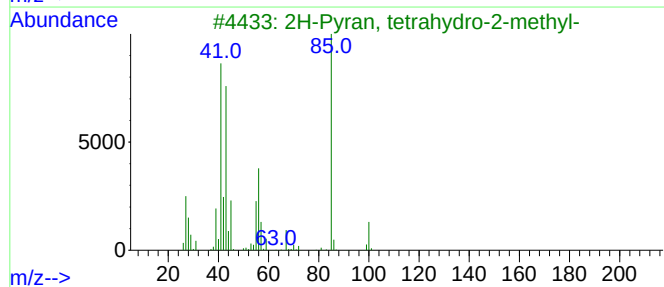
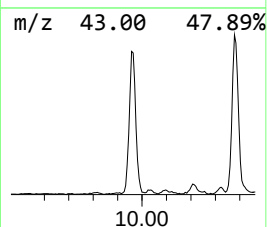
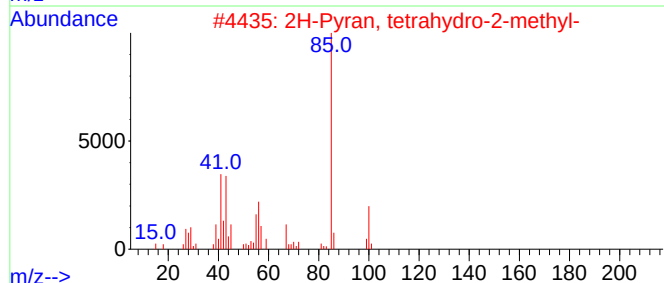
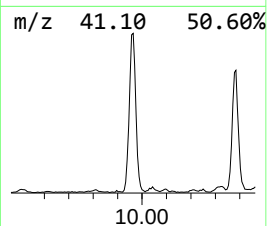
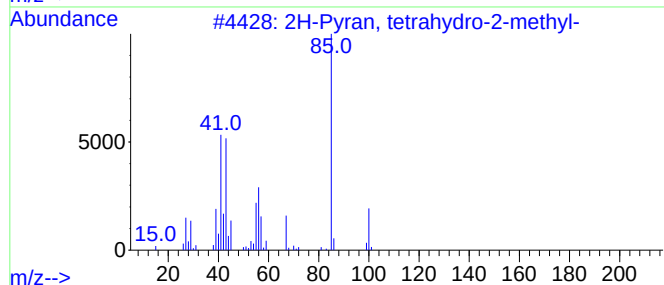
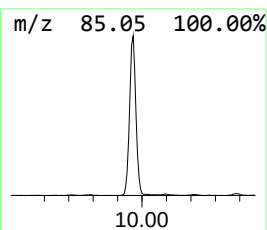
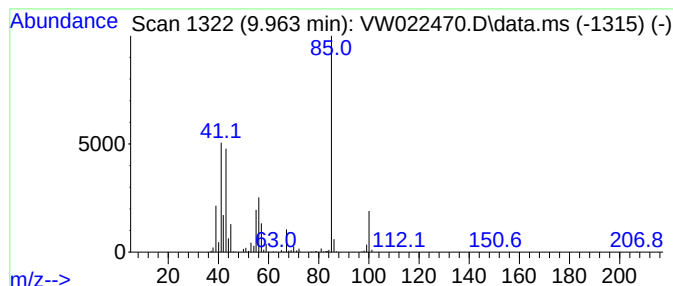
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Quant Title : SFAM01.0

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

Peak Number 5 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	16.82 ug/L	652670	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	93		
2	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	90		
3	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	83		
4	2(3H)-Furanone, dihydro-5-propyl-	128	C7H12O2	000105-21-5	59		
5	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	59		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022470.D
Acq On : 08 Apr 2022 17:43
Operator : SY/VA
Sample : N2293-18
Misc : 6.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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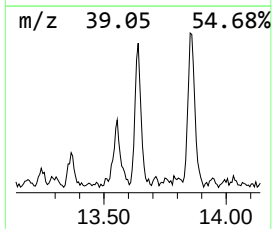
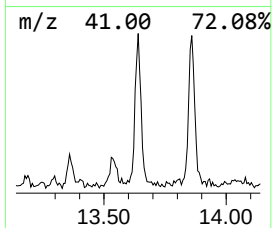
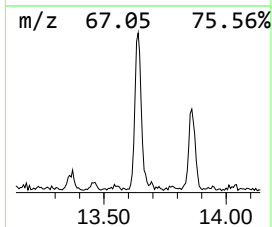
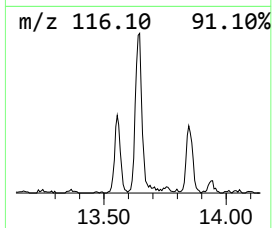
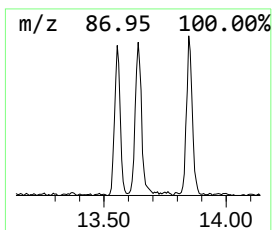
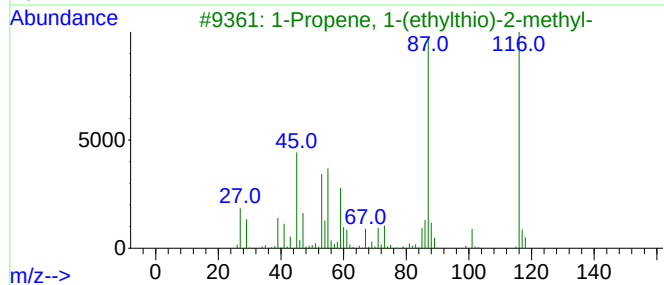
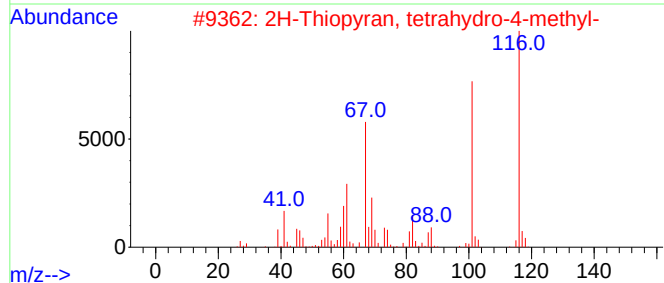
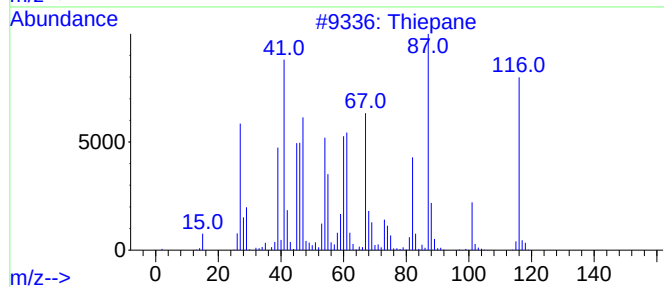
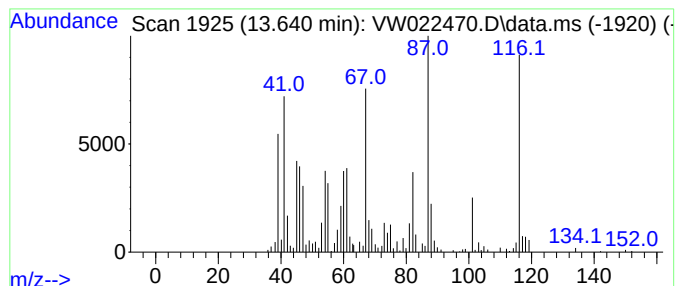
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Quant Title : SFAM01.0

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

Peak Number 7 Thiepane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	4.23 ug/L	221154	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiepane	116	C6H12S	004753-80-4	81
2			2H-Thiopyran, tetrahydro-4-methyl-	116	C6H12S	005161-17-1	43
3			1-Propene, 1-(ethylthio)-2-methyl-	116	C6H12S	027482-14-0	43
4			1,6-Hexanedithiol	150	C6H14S2	001191-43-1	43
5			4H-Thiopyran-4-one, tetrahydro-	116	C5H8OS	001072-72-6	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022470.D
Acq On : 08 Apr 2022 17:43
Operator : SY/VA
Sample : N2293-18
Misc : 6.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP6

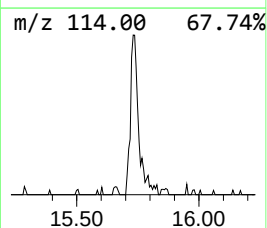
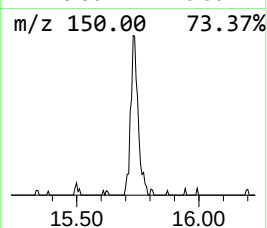
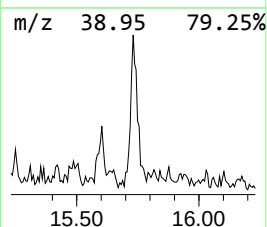
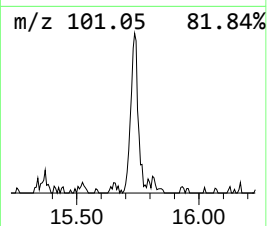
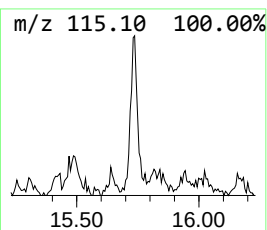
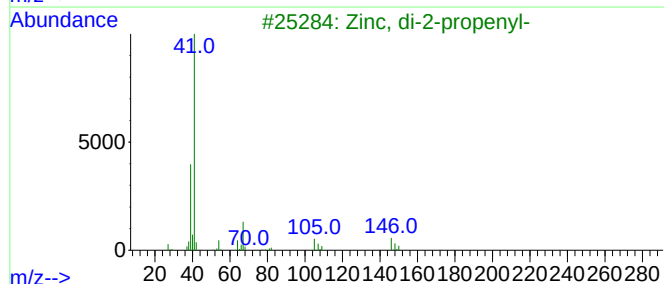
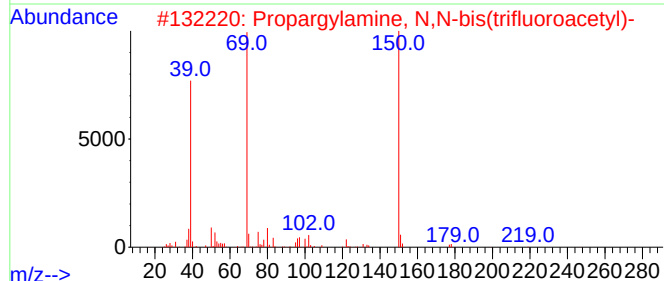
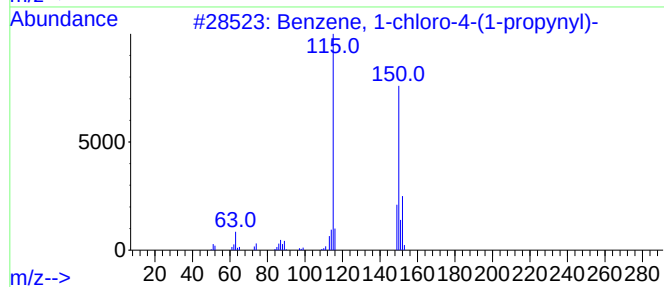
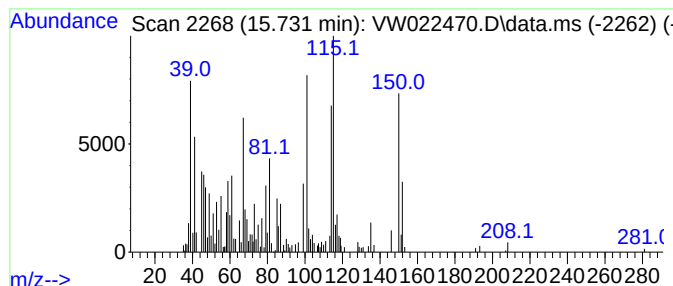
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Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-chloro-4-(1-prop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.731	1.91 ug/L	99631	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	53
2			Propargylamine, N,N-bis(trifluor...	247	C7H3F6N2	1000417-80-9	32
3			Zinc, di-2-propenyl-	146	C6H10Zn	001802-55-7	30
4			3-Hexenoic acid, (E)-	114	C6H10O2	001577-18-0	25
5			S-Methyl-1-propenethiosulfonate	152	C4H8O2S2	1000322-29-0	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022470.D
Acq On : 08 Apr 2022 17:43
Operator : SY/VA
Sample : N2293-18
Misc : 6.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP6

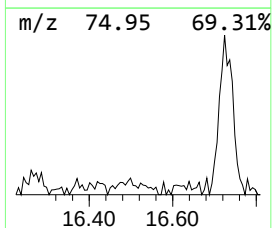
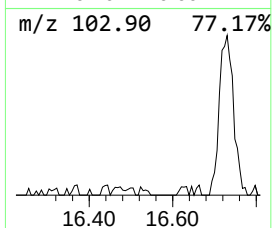
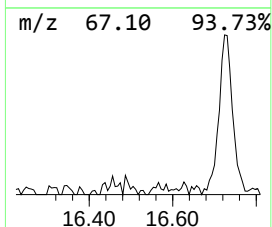
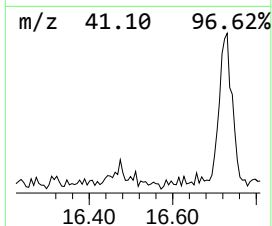
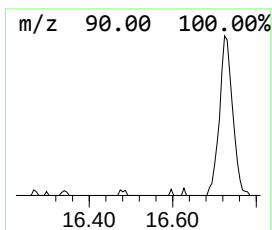
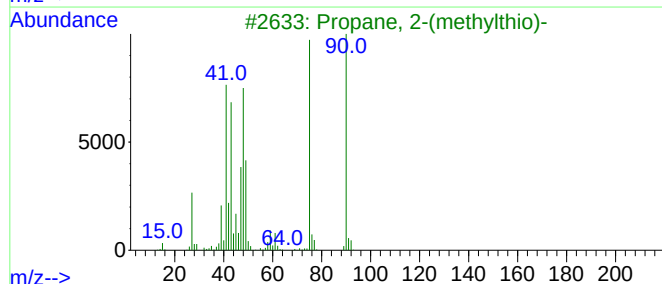
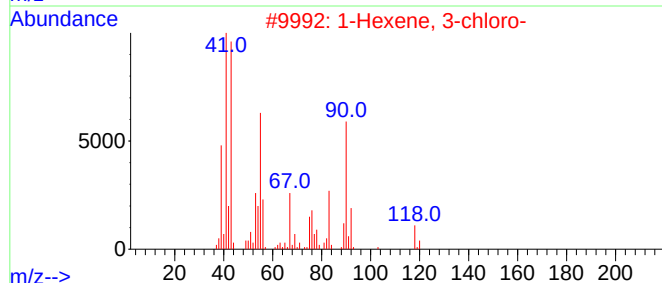
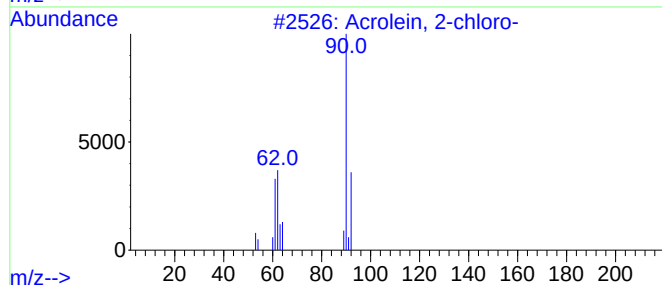
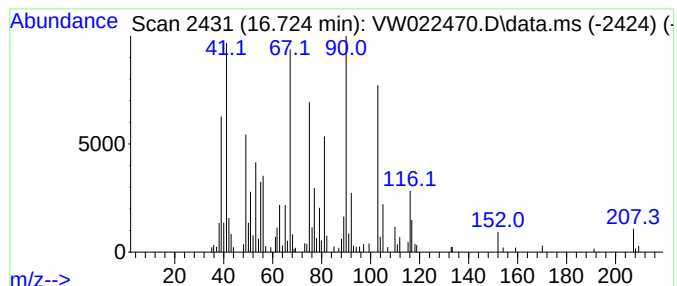
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.724	1.43 ug/L	74642	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	38
2			1-Hexene, 3-chloro-	118	C6H11Cl	053101-38-5	37
3			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	25
4			Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	22
5			Benzocyclobuten-1(2H)-one	118	C8H6O	003469-06-5	16



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW040822\
Data File : VW022470.D
Acq On : 08 Apr 2022 17:43
Operator : SY/VA
Sample : N2293-18
Misc : 6.67g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNP6

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Quant Title : SFAM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown-01	2.495	1.5	ug/L	59608	1	8.842	970219	25.0
Isopropenyl bro...	5.190	1.5	ug/L	56376	1	8.842	970219	25.0
Furan, tetrahyd...	9.049	1.6	ug/L	61683	1	8.842	970219	25.0
2H-Pyran, tetra...	9.963	16.8	ug/L	652670	1	8.842	970219	25.0
Thiepane	13.640	4.2	ug/L	221154	3	13.554	1305720	25.0
Benzene, 1-chlo...	15.731	1.9	ug/L	99631	3	13.554	1305720	25.0
unknown-02	16.724	1.4	ug/L	74642	3	13.554	1305720	25.0