

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
 Data File : VW022501.D
 Acq On : 12 Apr 2022 12:45
 Operator : SY/VA
 Sample : N2316-07
 Misc : 6.39g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNQ6

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: VW022501.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	24	rBV	1037652	1583174	58.01%	8.841%
2	2.355	69	74	84	rVB	132347	238959	8.76%	1.334%
3	2.495	93	97	105	rBV	22075	46294	1.70%	0.259%
4	2.885	154	161	172	rBV	91697	198634	7.28%	1.109%
5	3.373	236	241	247	rBV4	2537	6001	0.22%	0.034%
6	3.909	323	329	330	rBV6	3314	5000	0.18%	0.028%
7	4.019	335	347	359	rVV	150741	481377	17.64%	2.688%
8	4.123	359	364	371	rVV	17353	50495	1.85%	0.282%
9	4.208	371	378	391	rVV2	32370	103670	3.80%	0.579%
10	4.385	400	407	420	rVV2	7420	25576	0.94%	0.143%
11	4.921	486	495	498	rBV7	4992	13602	0.50%	0.076%
12	5.177	531	537	538	rBV5	2384	3466	0.13%	0.019%
13	5.562	594	600	607	rVB5	8641	21978	0.81%	0.123%
14	5.793	630	638	639	rVV4	2945	5199	0.19%	0.029%
15	5.940	657	662	671	rVB7	2833	7518	0.28%	0.042%
16	6.116	689	691	694	rVV4	1795	2659	0.10%	0.015%
17	6.622	766	774	779	rBV9	2174	6720	0.25%	0.038%
18	7.006	826	837	842	rVV5	21279	66169	2.42%	0.370%
19	7.080	843	849	859	rVV	45337	132374	4.85%	0.739%
20	7.177	863	865	875	rVB	7402	14058	0.52%	0.079%
21	7.647	931	942	955	rBV	274306	664272	24.34%	3.710%
22	7.890	975	982	985	rBV6	2788	6150	0.23%	0.034%
23	8.153	1021	1025	1033	rVB9	1623	3627	0.13%	0.020%
24	8.274	1033	1045	1048	rBV	492190	1106804	40.56%	6.181%
25	8.695	1107	1114	1124	rBV3	8347	19180	0.70%	0.107%
26	8.841	1129	1138	1149	rBV	475133	932251	34.16%	5.206%
27	9.079	1165	1177	1193	rBV9	17135	73751	2.70%	0.412%
28	9.274	1200	1209	1216	rBV	355404	739221	27.09%	4.128%
29	9.506	1239	1247	1256	rVV2	79116	169125	6.20%	0.944%
30	9.610	1259	1264	1269	rVB9	2581	5649	0.21%	0.032%
31	9.756	1282	1288	1292	rBV8	3206	6393	0.23%	0.036%
32	9.902	1305	1312	1313	rBV7	3454	5494	0.20%	0.031%
33	9.963	1315	1322	1328	rVV	121970	222569	8.16%	1.243%
34	10.036	1328	1334	1340	rVV	191578	333460	12.22%	1.862%
35	10.207	1357	1362	1365	rBV5	4035	8619	0.32%	0.048%
36	10.323	1373	1381	1387	rBV	723602	1311946	48.07%	7.327%

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

37	10.384	1387	1391	1399	rVB2	174180	301714	11.06%	1.685%
38	10.481	1402	1407	1408	rBV5	8854	14155	0.52%	0.079%
39	10.579	1416	1423	1430	rVB	122012	208397	7.64%	1.164%
40	10.664	1432	1437	1440	rVV4	7170	12861	0.47%	0.072%
41	10.719	1440	1446	1460	rVB3	60226	129853	4.76%	0.725%
42	10.859	1466	1469	1473	rVB5	3141	4564	0.17%	0.025%
43	10.926	1473	1480	1490	rBV2	366393	663936	24.33%	3.708%
44	11.061	1498	1502	1506	rVB5	6812	9962	0.37%	0.056%
45	11.176	1513	1521	1527	rVB5	9933	19767	0.72%	0.110%
46	11.292	1533	1540	1547	rVV	26563	48601	1.78%	0.271%
47	11.475	1566	1570	1577	rVB4	5331	9674	0.35%	0.054%
48	11.542	1577	1581	1587	rVB9	1463	3066	0.11%	0.017%
49	11.652	1587	1599	1606	rBV2	980340	2729021	100.00%	15.240%
50	11.725	1606	1611	1617	rVV2	49642	106939	3.92%	0.597%
51	11.841	1620	1630	1641	rVB2	22000	55963	2.05%	0.313%
52	11.963	1645	1650	1653	rBV6	2871	5760	0.21%	0.032%
53	12.164	1673	1683	1693	rVB	41594	74582	2.73%	0.417%
54	12.341	1708	1712	1717	rVB7	3548	6243	0.23%	0.035%
55	12.463	1727	1732	1737	rVB3	9133	14960	0.55%	0.084%
56	12.530	1738	1743	1748	rBV4	3836	6623	0.24%	0.037%
57	12.603	1750	1755	1760	rBV3	11697	21062	0.77%	0.118%
58	12.688	1760	1769	1775	rBV	288202	501111	18.36%	2.798%
59	12.743	1775	1778	1783	rVV	40296	60575	2.22%	0.338%
60	12.804	1783	1788	1793	rVB	12038	20664	0.76%	0.115%
61	12.889	1793	1802	1811	rBV	297073	498140	18.25%	2.782%
62	12.987	1812	1818	1827	rVB	172895	277812	10.18%	1.551%
63	13.103	1831	1837	1843	rBV	34343	58114	2.13%	0.325%
64	13.188	1845	1851	1854	rBV7	8951	14375	0.53%	0.080%
65	13.249	1856	1861	1865	rBV2	14651	20835	0.76%	0.116%
66	13.371	1876	1881	1886	rBV8	5667	12987	0.48%	0.073%
67	13.493	1896	1901	1903	rBV5	5569	9267	0.34%	0.052%
68	13.554	1904	1911	1922	rVB	793461	1334130	48.89%	7.450%
69	13.645	1923	1926	1932	rVB8	6972	12151	0.45%	0.068%
70	13.719	1934	1938	1939	rBV4	5805	7677	0.28%	0.043%
71	13.853	1953	1960	1970	rVB	725381	1229577	45.06%	6.867%
72	13.993	1971	1983	1988	rBV	100106	236282	8.66%	1.320%
73	14.090	1996	1999	2005	rVB3	16982	26970	0.99%	0.151%
74	14.200	2011	2017	2022	rVB5	13387	26378	0.97%	0.147%
75	14.273	2022	2029	2030	rBV5	4837	8920	0.33%	0.050%
76	14.328	2035	2038	2045	rVB5	11218	19674	0.72%	0.110%

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77	14.414	2049	2052	2055	rBV	7996	8881	0.33%	0.050%
78	14.456	2055	2059	2063	rBV3	12477	20454	0.75%	0.114%
79	14.578	2074	2079	2083	rBV3	18891	32349	1.19%	0.181%
80	14.627	2083	2087	2092	rVB5	14777	27682	1.01%	0.155%
81	14.682	2092	2096	2099	rBV5	8789	13983	0.51%	0.078%
82	14.785	2108	2113	2120	rBV2	22847	47907	1.76%	0.268%
83	14.853	2121	2124	2129	rVV4	9654	15534	0.57%	0.087%
84	14.907	2131	2133	2137	rVB4	4909	7287	0.27%	0.041%
85	15.060	2153	2158	2164	rBV4	16791	31996	1.17%	0.179%
86	15.127	2165	2169	2172	rBV4	12102	19276	0.71%	0.108%
87	15.310	2194	2199	2203	rVB6	4142	6894	0.25%	0.038%
88	15.365	2203	2208	2211	rBV2	5631	9076	0.33%	0.051%
89	15.420	2211	2217	2220	rVV3	12210	20352	0.75%	0.114%
90	15.487	2220	2228	2236	rVB4	17737	48526	1.78%	0.271%
91	15.578	2239	2243	2249	rVB8	4470	8412	0.31%	0.047%
92	15.645	2249	2254	2260	rBV6	7352	16444	0.60%	0.092%
93	15.743	2264	2270	2271	rBV5	2615	4709	0.17%	0.026%
94	15.889	2289	2294	2297	rBV6	5249	9179	0.34%	0.051%
95	15.938	2298	2302	2309	rVB2	11159	19033	0.70%	0.106%
96	16.163	2332	2339	2346	rBV7	9433	22384	0.82%	0.125%
97	16.273	2349	2357	2364	rBV8	9658	23031	0.84%	0.129%
98	16.340	2364	2368	2369	rBV4	2830	3801	0.14%	0.021%
99	16.425	2378	2382	2389	rBV8	3774	9764	0.36%	0.055%
100	16.639	2408	2417	2424	rVV6	13074	35027	1.28%	0.196%

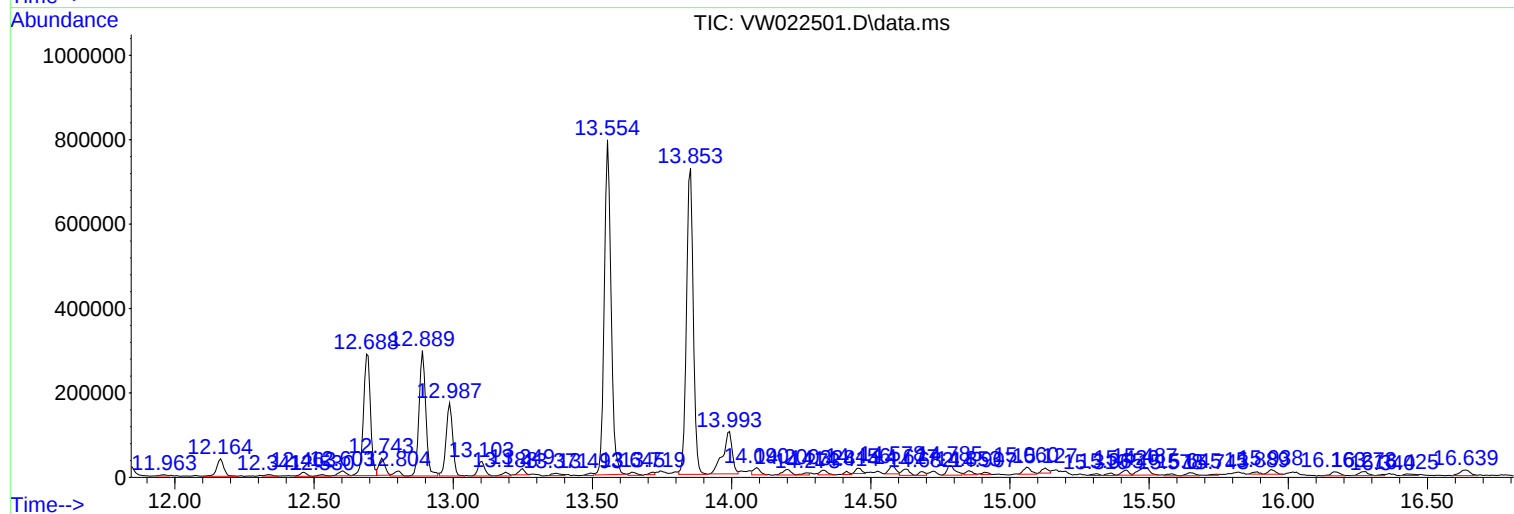
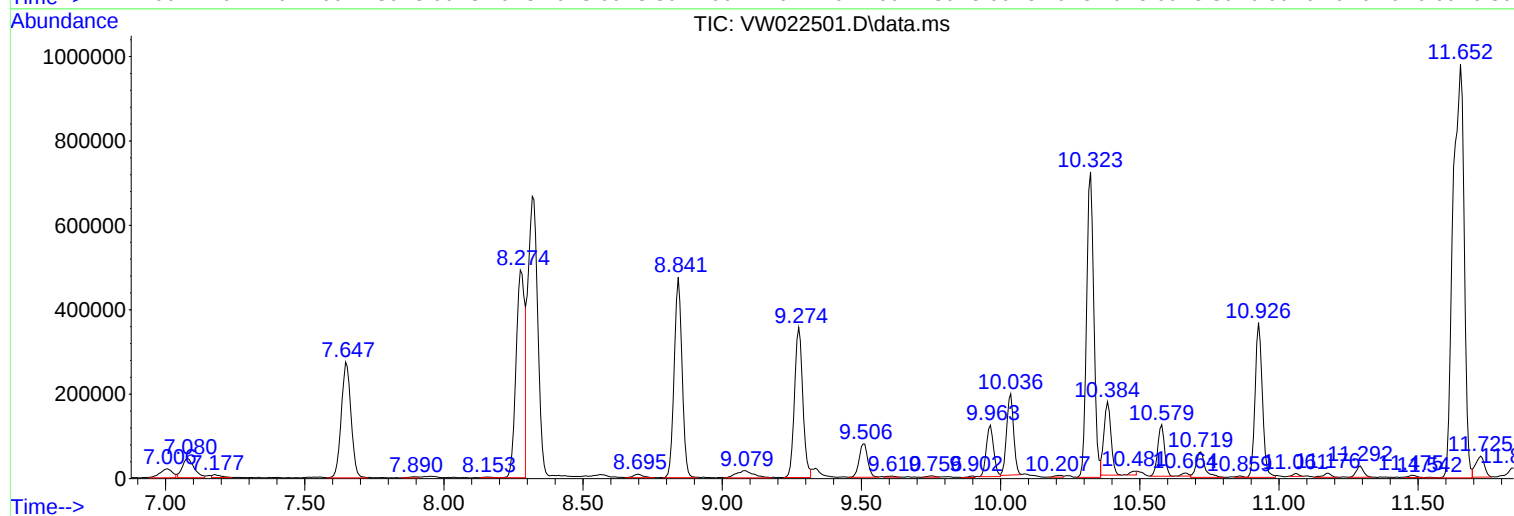
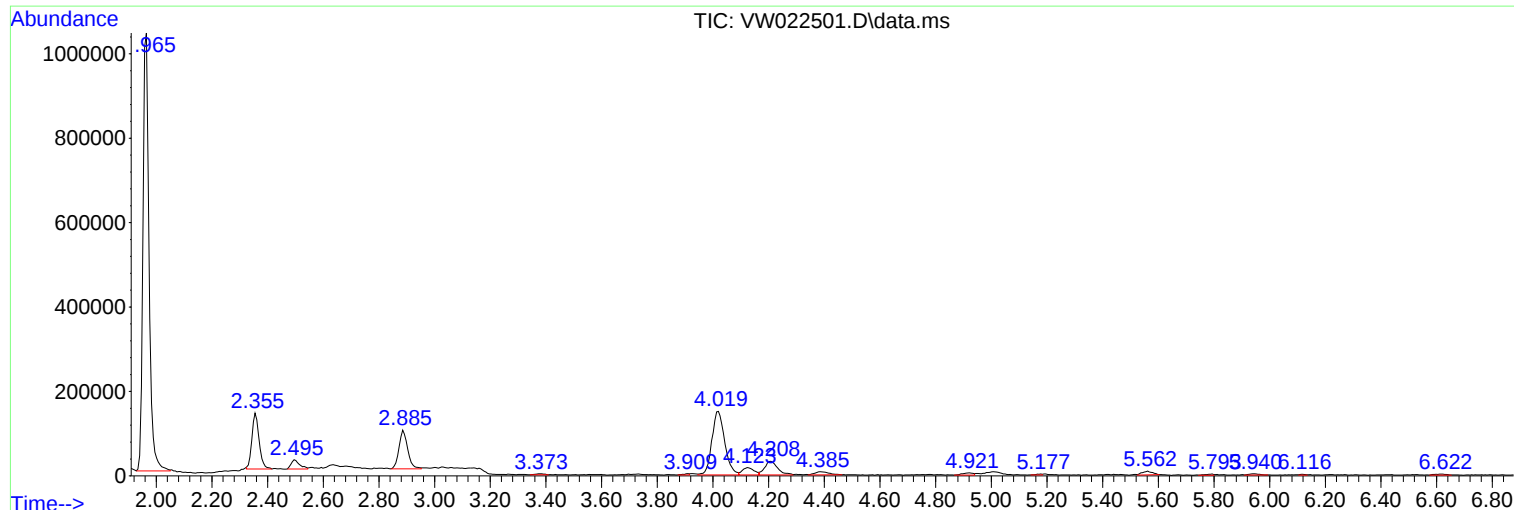
Sum of corrected areas: 17906757

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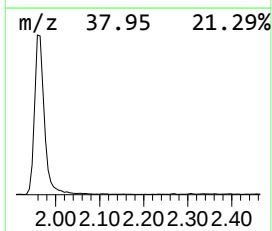
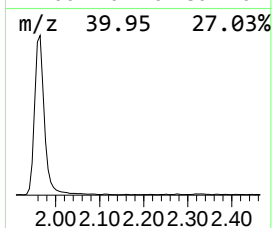
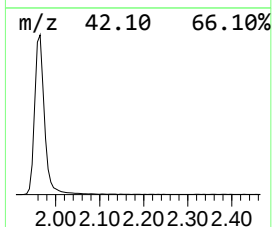
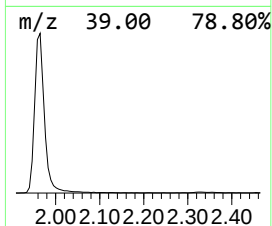
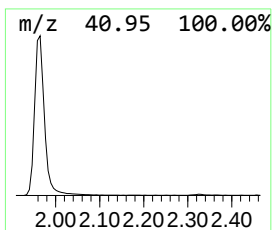
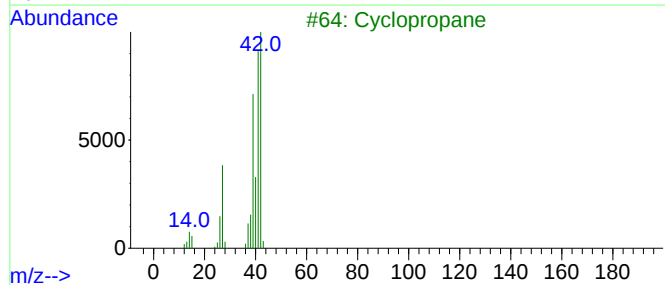
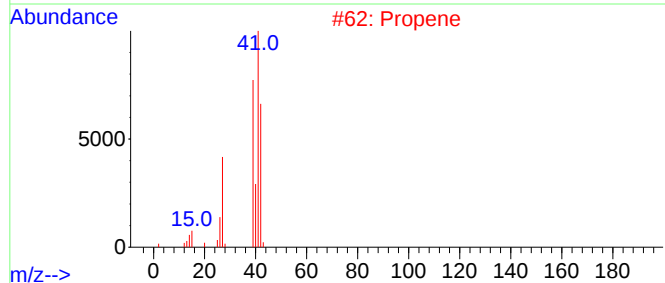
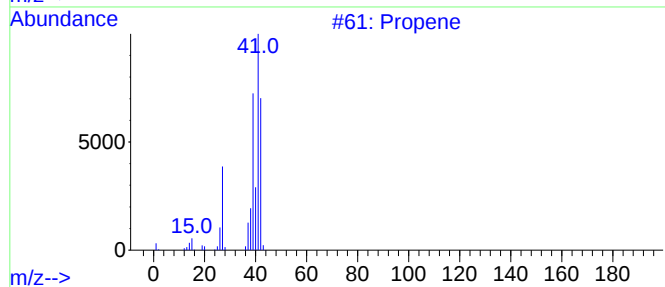
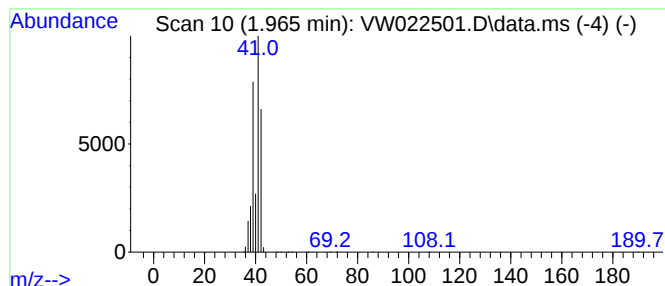
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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	42.46 ug/L	1583170	1,4-Difluorobenzene	8.841

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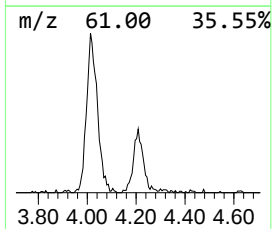
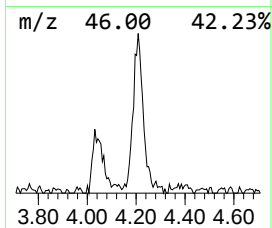
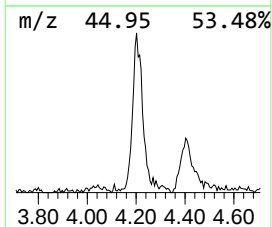
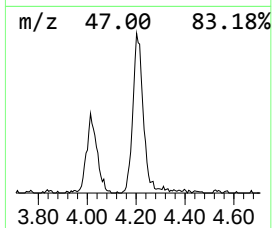
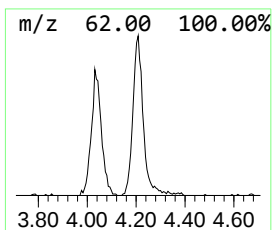
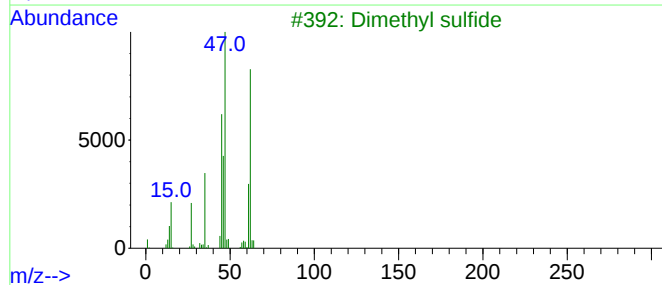
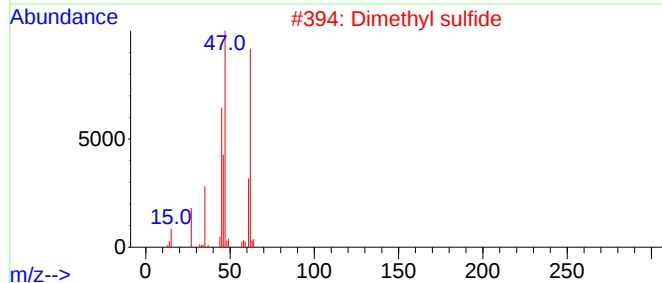
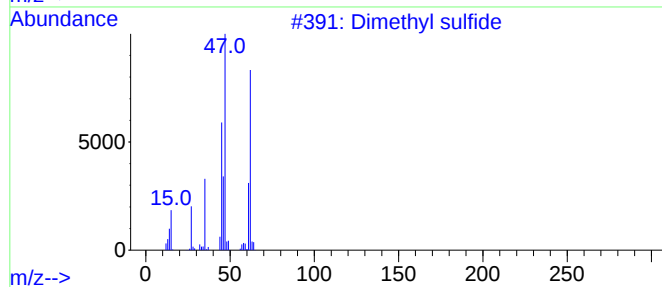
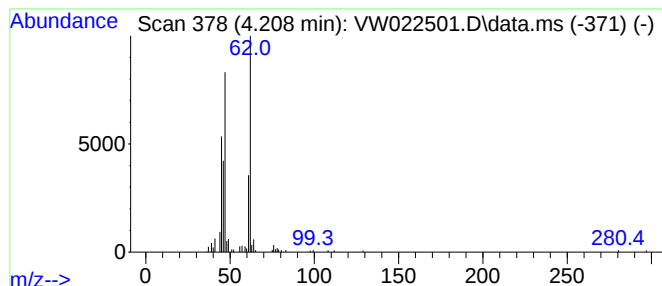
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 Dimethyl sulfide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.208	2.78 ug/L	103670	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dimethyl sulfide	62	C2H6S	000075-18-3	94
2			Dimethyl sulfide	62	C2H6S	000075-18-3	91
3			Dimethyl sulfide	62	C2H6S	000075-18-3	91
4			Dimethyl sulfide	62	C2H6S	000075-18-3	87
5			Dimethyl sulfide	62	C2H6S	000075-18-3	86



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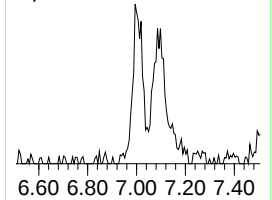
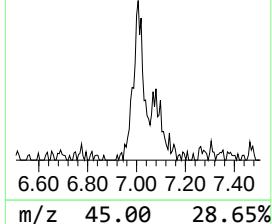
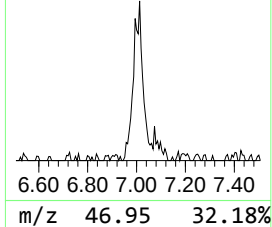
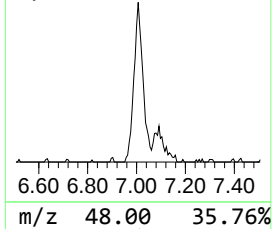
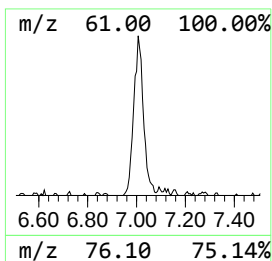
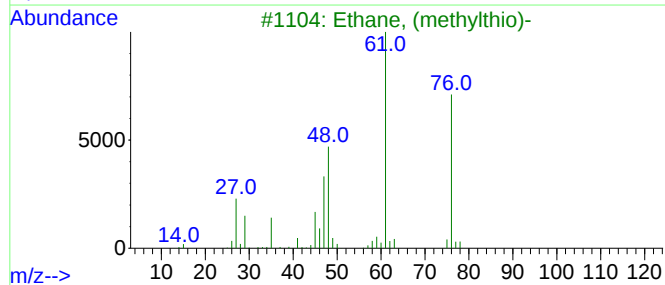
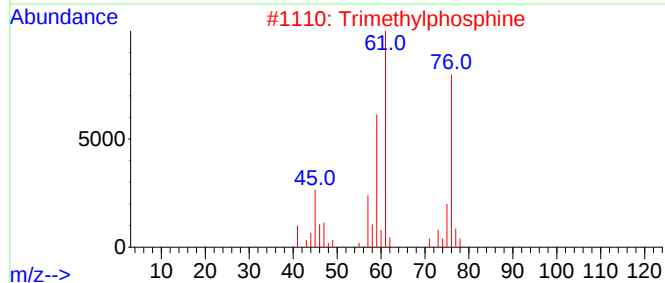
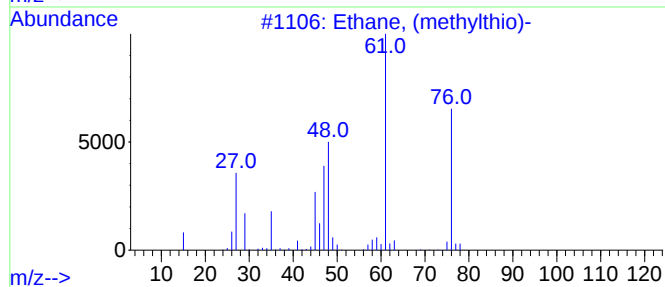
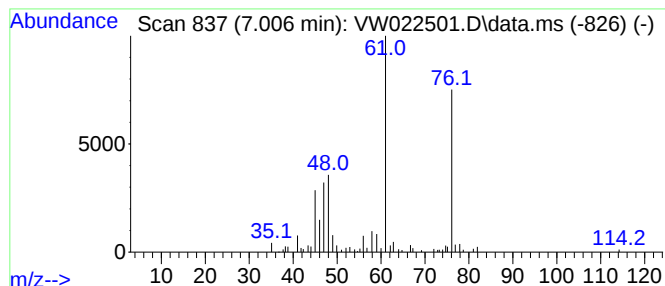
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Ethane, (methylthio)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.006	1.77 ug/L	66169	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, (methylthio)-	76	C3H8S	000624-89-5	91
2			Trimethylphosphine	76	C3H9P	000594-09-2	86
3			Ethane, (methylthio)-	76	C3H8S	000624-89-5	86
4			Ethane, (methylthio)-	76	C3H8S	000624-89-5	83
5			Ethane, (methylthio)-	76	C3H8S	000624-89-5	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022501.D
Acq On : 12 Apr 2022 12:45
Operator : SY/VA
Sample : N2316-07
Misc : 6.39g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ6

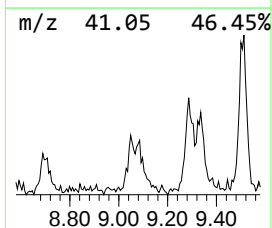
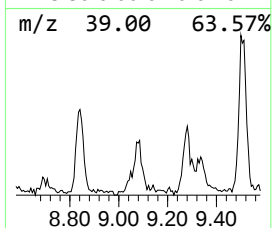
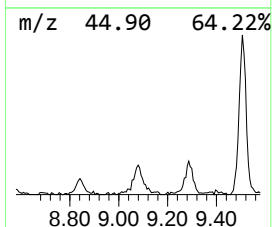
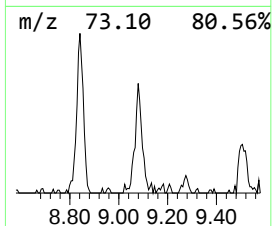
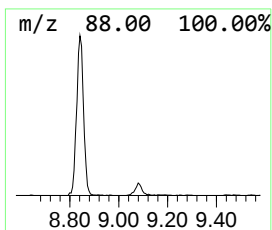
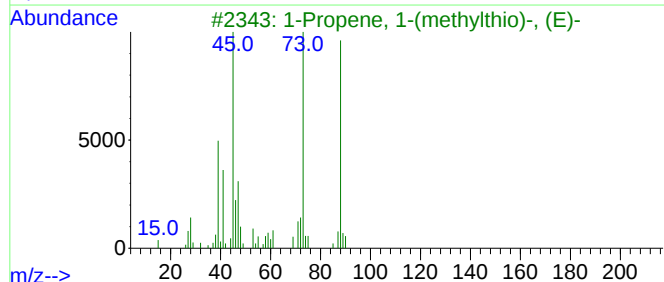
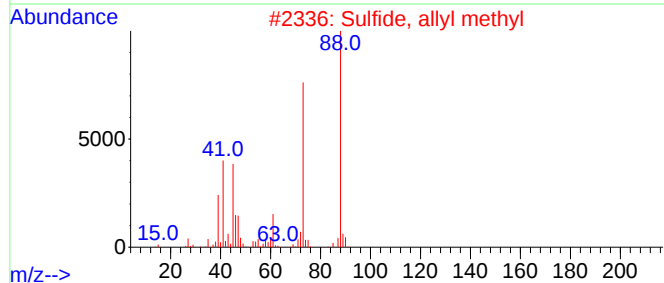
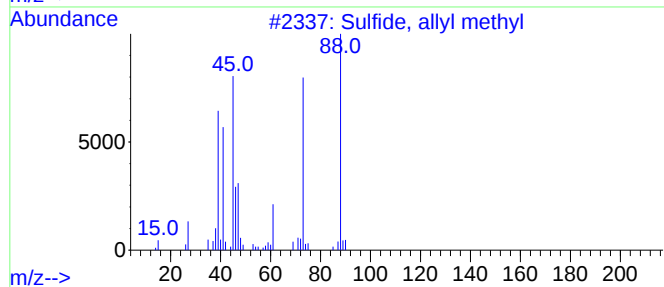
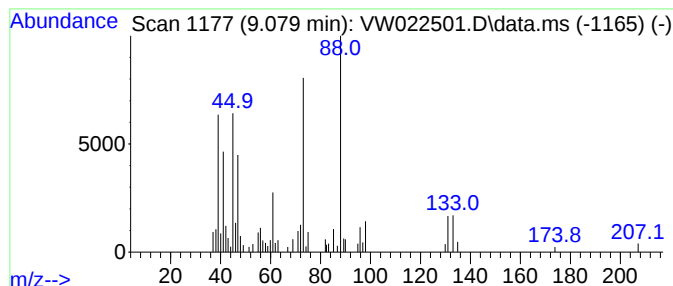
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Quant Title : SFAM01.0

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

Peak Number 4 Sulfide, allyl methyl Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.079	1.98 ug/L	73751	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfide, allyl methyl	88	C4H8S	010152-76-8	60
2			Sulfide, allyl methyl	88	C4H8S	010152-76-8	55
3			1-Propene, 1-(methylthio)-, (E)-	88	C4H8S	042848-06-6	53
4			Sulfide, allyl methyl	88	C4H8S	010152-76-8	40
5			1,2-Dimethoxy-ethene	88	C4H8O2	1000194-22-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022501.D
Acq On : 12 Apr 2022 12:45
Operator : SY/VA
Sample : N2316-07
Misc : 6.39g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ6

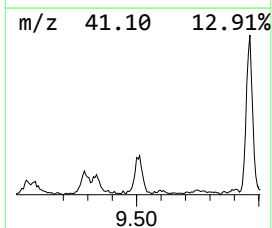
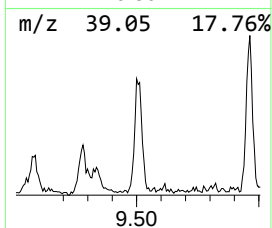
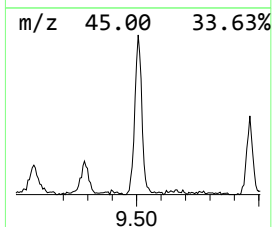
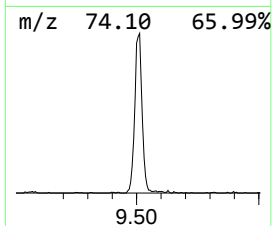
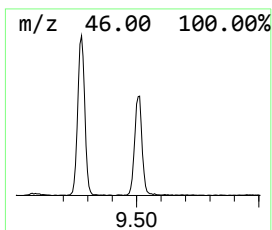
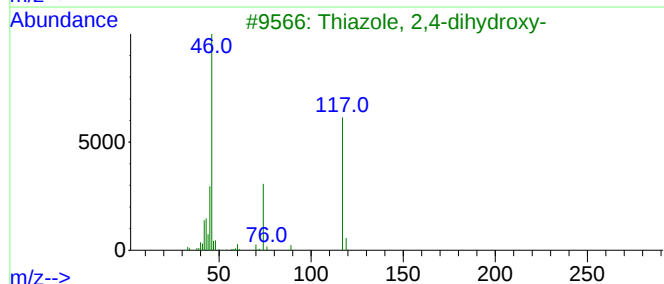
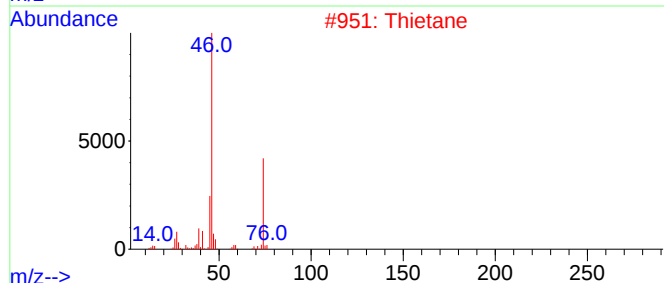
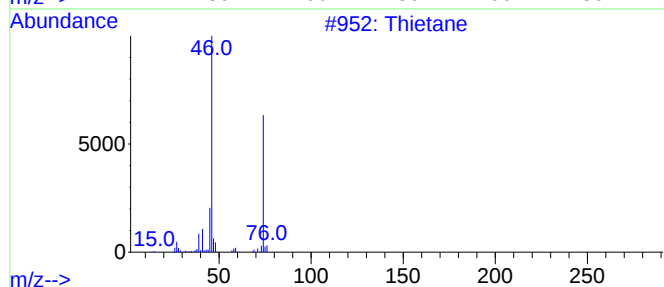
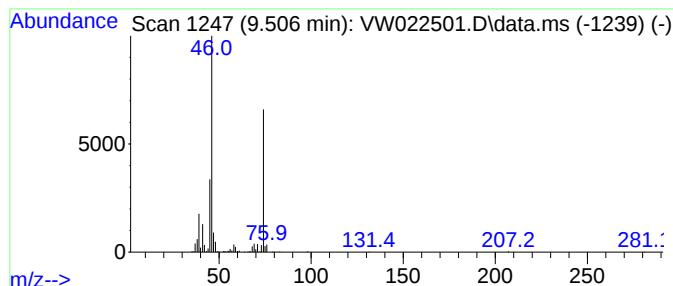
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Quant Title : SFAM01.0

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

Peak Number 5 Thietane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.506	4.54 ug/L	169125	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thietane	74	C3H6S	000287-27-4	90
2			Thietane	74	C3H6S	000287-27-4	72
3			Thiazole, 2,4-dihydroxy-	117	C3H3NO2S	000625-85-4	64
4			Thiirane, methyl-	74	C3H6S	001072-43-1	37
5			Dihydro-3-(2H)-thiophenone	102	C4H6OS	001003-04-9	33



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022501.D
Acq On : 12 Apr 2022 12:45
Operator : SY/VA
Sample : N2316-07
Misc : 6.39g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ6

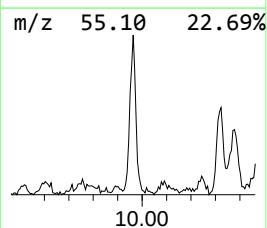
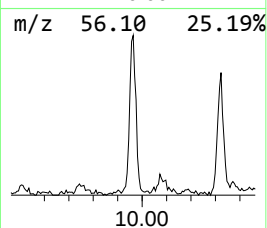
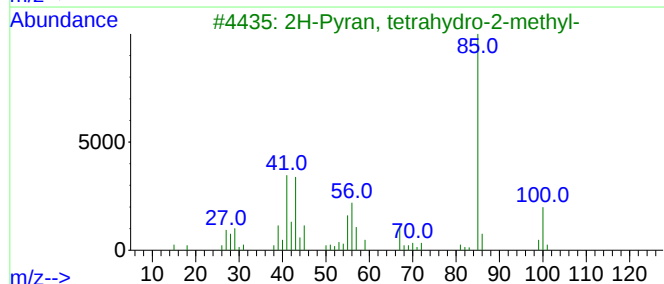
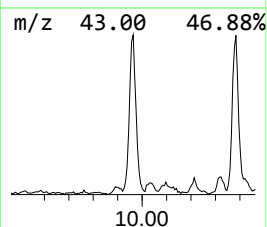
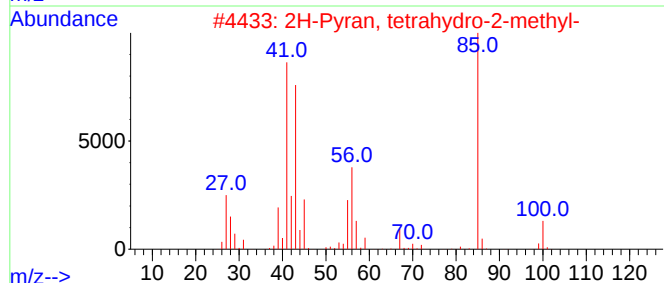
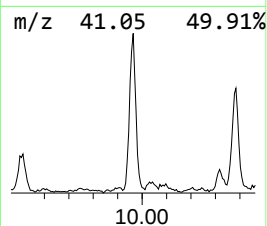
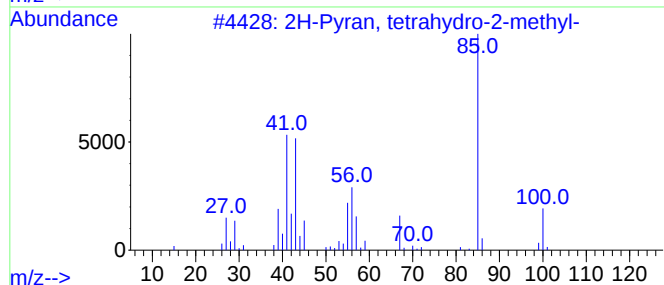
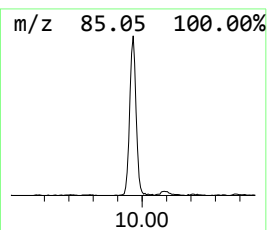
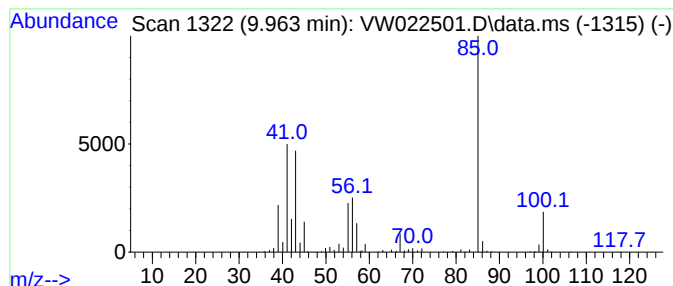
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Quant Title : SFAM01.0

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

Peak Number 6 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	5.97 ug/L	222569	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	94
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	83
3	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	80
4	Oxetane, 2-ethyl-3-methyl-			100	C6H12O	053778-62-4	72
5	2H-Pyran-2-ol, tetrahydro-			102	C5H10O2	000694-54-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022501.D
Acq On : 12 Apr 2022 12:45
Operator : SY/VA
Sample : N2316-07
Misc : 6.39g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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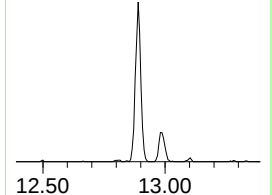
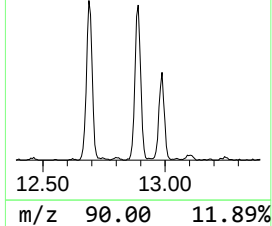
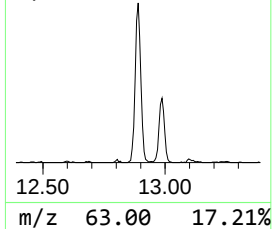
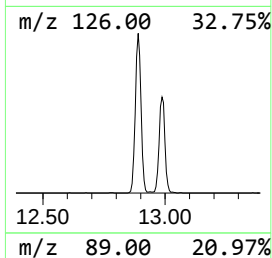
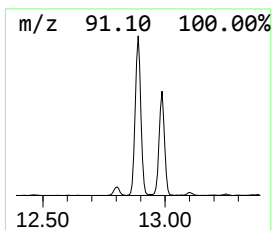
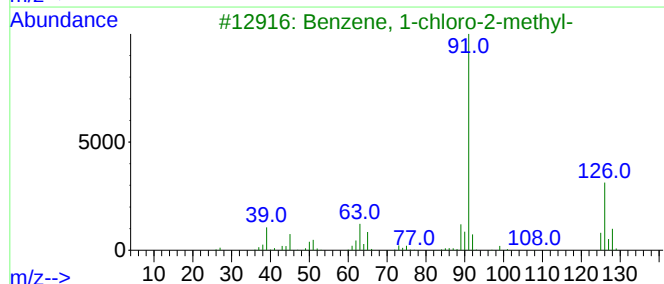
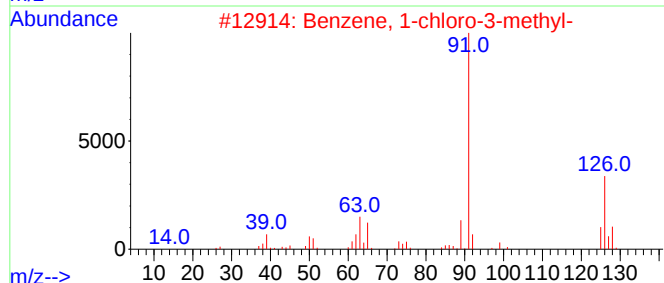
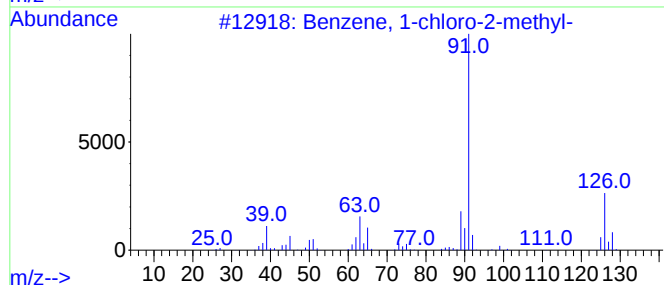
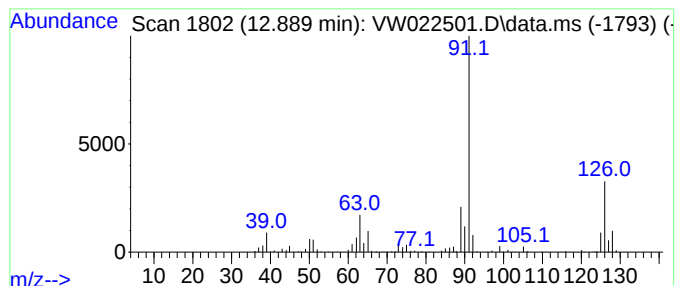
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Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-chloro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	9.33 ug/L	498140	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	96
2			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	95
3			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	94
4			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	93
5			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022501.D
Acq On : 12 Apr 2022 12:45
Operator : SY/VA
Sample : N2316-07
Misc : 6.39g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ6

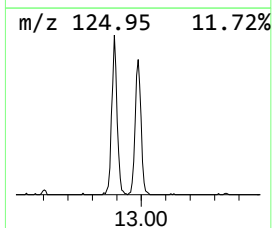
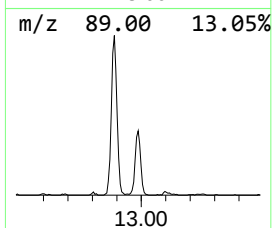
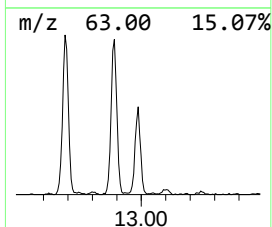
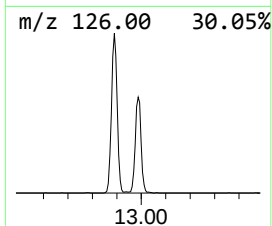
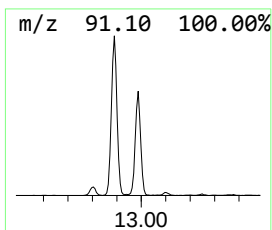
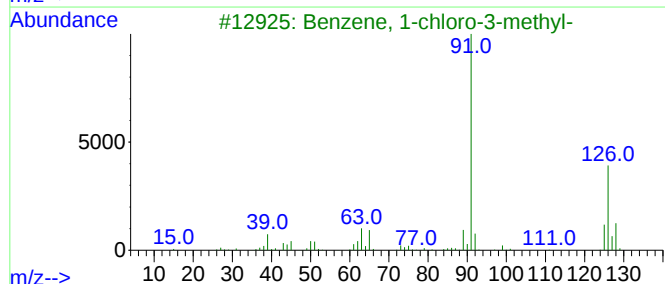
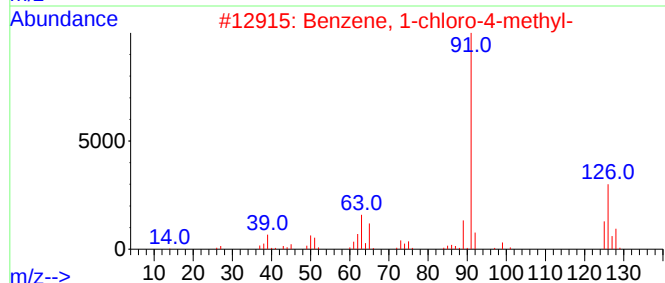
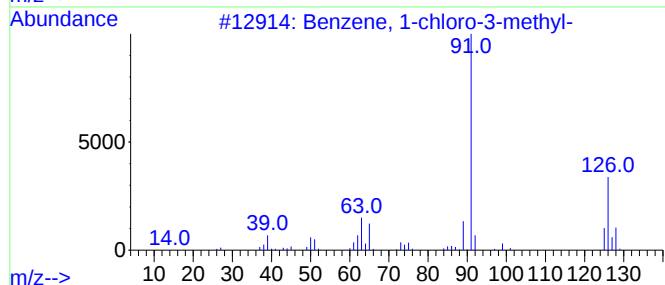
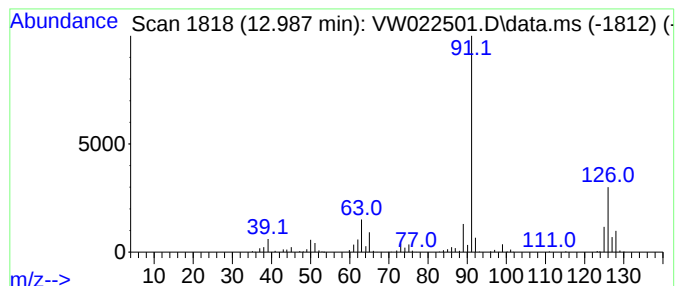
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Quant Title : SFAM01.0

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

Peak Number 9 Benzene, 1-chloro-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.987	5.21 ug/L	277812	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	96
2			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	96
3			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	95
4			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	95
5			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022501.D
Acq On : 12 Apr 2022 12:45
Operator : SY/VA
Sample : N2316-07
Misc : 6.39g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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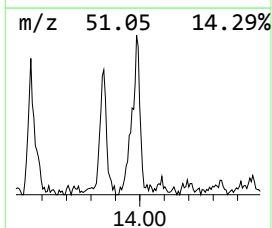
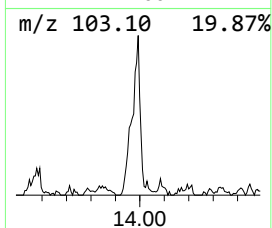
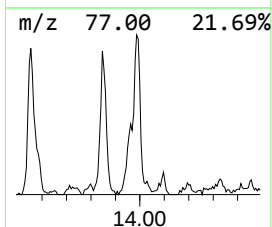
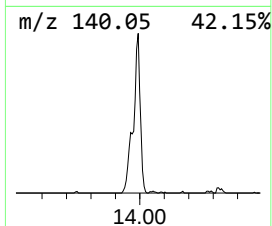
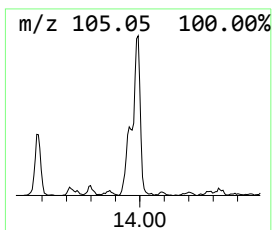
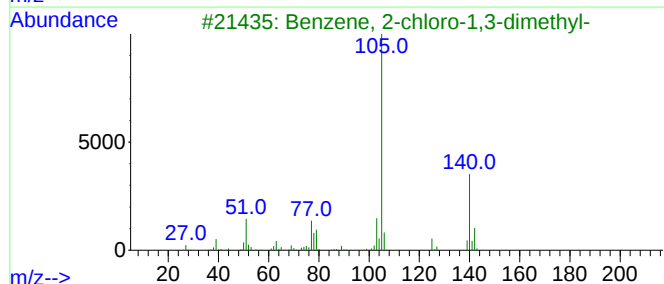
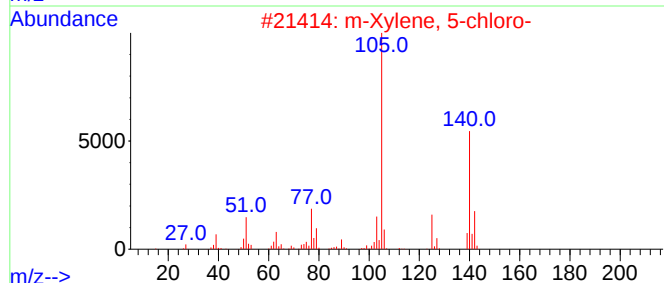
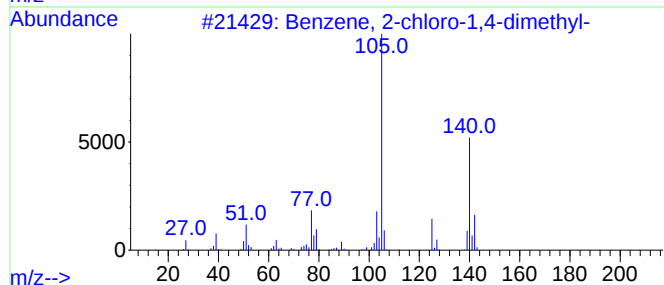
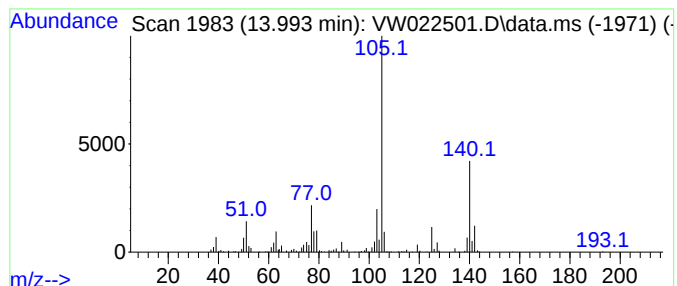
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Quant Title : SFAM01.0

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

Peak Number 10 Benzene, 2-chloro-1,4-dimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.993	4.43 ug/L	236282	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-chloro-1,4-dimethyl-	140	C8H9Cl	000095-72-7	97
2			m-Xylene, 5-chloro-	140	C8H9Cl	000556-97-8	94
3			Benzene, 2-chloro-1,3-dimethyl-	140	C8H9Cl	006781-98-2	87
4			Benzene, 2-chloro-1,4-dimethyl-	140	C8H9Cl	000095-72-7	87
5			Benzene, 1-chloro-2,3-dimethyl-	140	C8H9Cl	000608-23-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
 Data File : VW022501.D
 Acq On : 12 Apr 2022 12:45
 Operator : SY/VA
 Sample : N2316-07
 Misc : 6.39g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNQ6

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
(DEL) Alkane: S...	1.965	42.5	ug/L	1583170	1	8.841	932251	25.0
Dimethyl sulfide	4.208	2.8	ug/L	103670	1	8.841	932251	25.0
Ethane, (methyl...	7.006	1.8	ug/L	66169	1	8.841	932251	25.0
Sulfide, allyl ...	9.079	2.0	ug/L	73751	1	8.841	932251	25.0
Thietane	9.506	4.5	ug/L	169125	1	8.841	932251	25.0
2H-Pyran, tetra...	9.963	6.0	ug/L	222569	1	8.841	932251	25.0
Benzene, 1-chlo...	12.890	9.3	ug/L	498140	3	13.554	1334130	25.0
Benzene, 1-chlo...	12.987	5.2	ug/L	277812	3	13.554	1334130	25.0
Benzene, 2-chlo...	13.993	4.4	ug/L	236282	3	13.554	1334130	25.0