

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
 Data File : VW022517.D
 Acq On : 12 Apr 2022 19:11
 Operator : SY/VA
 Sample : N2373-01
 Misc : 3.29g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BFFD2

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: VW022517.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.355	69	74	81	rBV	98781	159312	14.88%	2.207%
2	2.891	156	162	170	rVB	57722	131014	12.23%	1.815%
3	3.379	238	242	246	rBV7	2589	5340	0.50%	0.074%
4	3.458	254	255	258	rBV3	1005	1200	0.11%	0.017%
5	3.495	258	261	263	rBV3	1379	1745	0.16%	0.024%
6	3.543	263	269	279	rBV	13585	36245	3.38%	0.502%
7	3.909	327	329	331	rBV2	1294	1268	0.12%	0.018%
8	4.019	336	347	360	rBV2	91693	283741	26.49%	3.931%
9	4.403	407	410	411	rBV3	980	1005	0.09%	0.014%
10	4.610	441	444	449	rVV6	861	1569	0.15%	0.022%
11	4.696	453	458	463	rBV7	1416	2487	0.23%	0.034%
12	4.921	485	495	505	rVV3	13737	48737	4.55%	0.675%
13	5.007	507	509	513	rVV4	971	1113	0.10%	0.015%
14	5.153	531	533	537	rVB4	878	1184	0.11%	0.016%
15	5.196	537	540	545	rBV6	1060	1459	0.14%	0.020%
16	5.305	554	558	561	rVB5	1296	2072	0.19%	0.029%
17	5.336	561	563	565	rBV3	1194	1500	0.14%	0.021%
18	5.689	617	621	623	rVB5	1614	1712	0.16%	0.024%
19	5.769	630	634	635	rBV4	960	890	0.08%	0.012%
20	5.793	635	638	641	rBV4	1313	1417	0.13%	0.020%
21	5.848	644	647	648	rBV2	1065	909	0.08%	0.013%
22	5.945	653	663	671	rBV4	6967	21271	1.99%	0.295%
23	6.171	697	700	703	rBV4	1018	1500	0.14%	0.021%
24	6.311	720	723	728	rVB6	1659	2735	0.26%	0.038%
25	6.354	728	730	734	rBV3	877	1301	0.12%	0.018%
26	6.409	737	739	742	rVB2	881	895	0.08%	0.012%
27	6.488	749	752	754	rVB3	1096	1032	0.10%	0.014%
28	6.555	757	763	764	rBV4	1120	1984	0.19%	0.027%
29	6.604	766	771	775	rVB6	814	1470	0.14%	0.020%
30	6.665	778	781	784	rBV4	744	990	0.09%	0.014%
31	6.854	809	812	813	rBV2	948	992	0.09%	0.014%
32	7.079	840	849	864	rBV2	33071	110604	10.33%	1.532%
33	7.220	870	872	875	rVB4	1364	1309	0.12%	0.018%
34	7.311	883	887	889	rBV4	778	967	0.09%	0.013%
35	7.329	889	890	892	rBV2	1304	881	0.08%	0.012%
36	7.543	924	925	930	rVB6	1072	879	0.08%	0.012%

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

37	7.652	932	943	952	rBV	166280	412440	38.51%	5.714%
38	8.274	1035	1045	1061	rBV2	274481	884227	82.56%	12.251%
39	8.469	1075	1077	1079	rVB3	1915	1005	0.09%	0.014%
40	8.841	1130	1138	1148	rBV2	280554	550459	51.40%	7.626%
41	9.134	1184	1186	1187	rBV2	1136	916	0.09%	0.013%
42	9.274	1200	1209	1217	rBV	213181	436967	40.80%	6.054%
43	9.823	1296	1299	1300	rBV3	1290	1056	0.10%	0.015%
44	10.030	1327	1333	1345	rVB	76201	146979	13.72%	2.036%
45	10.323	1373	1381	1388	rBV	318851	565574	52.81%	7.836%
46	10.384	1389	1391	1396	rVB4	5099	7765	0.73%	0.108%
47	10.579	1417	1423	1431	rBV2	49493	90558	8.46%	1.255%
48	10.652	1434	1435	1437	rBV2	1058	884	0.08%	0.012%
49	10.926	1474	1480	1493	rBV	113902	215820	20.15%	2.990%
50	11.365	1548	1552	1555	rBV6	996	1572	0.15%	0.022%
51	11.426	1560	1562	1565	rVB4	1402	1299	0.12%	0.018%
52	11.457	1565	1567	1571	rVB4	1014	1010	0.09%	0.014%
53	11.548	1581	1582	1587	rVB5	1650	2308	0.22%	0.032%
54	11.627	1587	1595	1609	rBV	284116	494854	46.21%	6.856%
55	11.725	1609	1611	1619	rVV8	2025	4233	0.40%	0.059%
56	11.786	1619	1621	1625	rVV4	856	1135	0.11%	0.016%
57	11.847	1625	1631	1636	rVV10	2386	4658	0.43%	0.065%
58	11.975	1650	1652	1655	rBV3	886	998	0.09%	0.014%
59	12.005	1655	1657	1660	rVB3	1000	1177	0.11%	0.016%
60	12.139	1676	1679	1682	rBV4	1107	1727	0.16%	0.024%
61	12.170	1682	1684	1686	rVV3	1095	1171	0.11%	0.016%
62	12.200	1686	1689	1692	rVB4	1353	1284	0.12%	0.018%
63	12.249	1694	1697	1700	rVB4	1851	2612	0.24%	0.036%
64	12.280	1700	1702	1704	rBV3	1727	1911	0.18%	0.026%
65	12.334	1708	1711	1712	rBV2	1154	1255	0.12%	0.017%
66	12.359	1712	1715	1717	rVB3	1042	1066	0.10%	0.015%
67	12.389	1717	1720	1721	rVB3	1204	1178	0.11%	0.016%
68	12.493	1734	1737	1740	rBV4	1037	1504	0.14%	0.021%
69	12.548	1741	1746	1747	rBV5	1146	1673	0.16%	0.023%
70	12.603	1747	1755	1762	rVB6	10478	23133	2.16%	0.320%
71	12.688	1762	1769	1777	rBV	156640	264892	24.73%	3.670%
72	12.956	1805	1813	1817	rBV10	3304	9207	0.86%	0.128%
73	13.103	1831	1837	1839	rBV7	2287	4544	0.42%	0.063%
74	13.249	1843	1861	1878	rVV5	224110	1070985	100.00%	14.838%
75	13.407	1884	1887	1893	rVB5	10035	16351	1.53%	0.227%
76	13.554	1905	1911	1918	rBV	175389	286199	26.72%	3.965%

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 Title : SFAM01.0

77	13.767	1940	1946	1950	rBV8	2784	4933	0.46%	0.068%
78	13.852	1952	1960	1970	rVV	138520	241110	22.51%	3.340%
79	14.066	1983	1995	2000	rBV	7568	16361	1.53%	0.227%
80	14.212	2014	2019	2025	rBV8	3309	6514	0.61%	0.090%
81	14.328	2034	2038	2041	rVB5	1968	2563	0.24%	0.036%
82	14.468	2054	2061	2067	rBV2	18561	32182	3.00%	0.446%
83	14.535	2071	2072	2075	rBV3	1471	1105	0.10%	0.015%
84	14.682	2093	2096	2098	rBV4	1606	1449	0.14%	0.020%
85	14.730	2098	2104	2113	rVV2	67638	119696	11.18%	1.658%
86	14.816	2115	2118	2126	rVB9	3952	9278	0.87%	0.129%
87	15.005	2141	2149	2154	rBV10	4901	11373	1.06%	0.158%
88	15.053	2154	2157	2158	rBV3	1617	1849	0.17%	0.026%
89	15.090	2160	2163	2168	rVB7	1443	2663	0.25%	0.037%
90	15.206	2175	2182	2183	rBV5	4323	6691	0.62%	0.093%
91	15.334	2193	2203	2212	rBV5	18288	56393	5.27%	0.781%
92	15.547	2222	2238	2244	rBV5	63894	177953	16.62%	2.465%
93	15.846	2284	2287	2288	rBV3	2311	2257	0.21%	0.031%
94	15.962	2304	2306	2311	rVB6	2548	3403	0.32%	0.047%
95	16.181	2341	2342	2346	rVB4	1281	1154	0.11%	0.016%
96	16.224	2346	2349	2353	rBV6	1213	2199	0.21%	0.030%
97	16.395	2374	2377	2380	rBV5	1088	1545	0.14%	0.021%
98	16.498	2380	2394	2399	rVV6	16754	49464	4.62%	0.685%
99	16.553	2401	2403	2408	rVB6	1406	2138	0.20%	0.030%
100	16.730	2423	2432	2441	rVV3	49092	110242	10.29%	1.527%

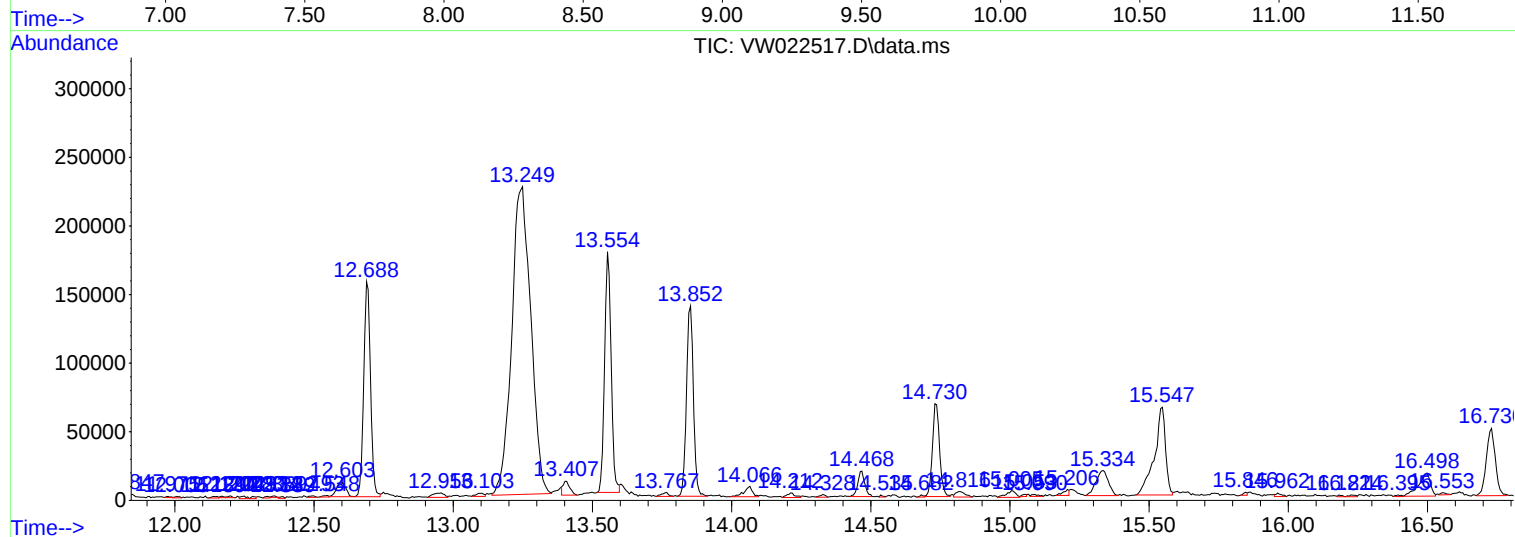
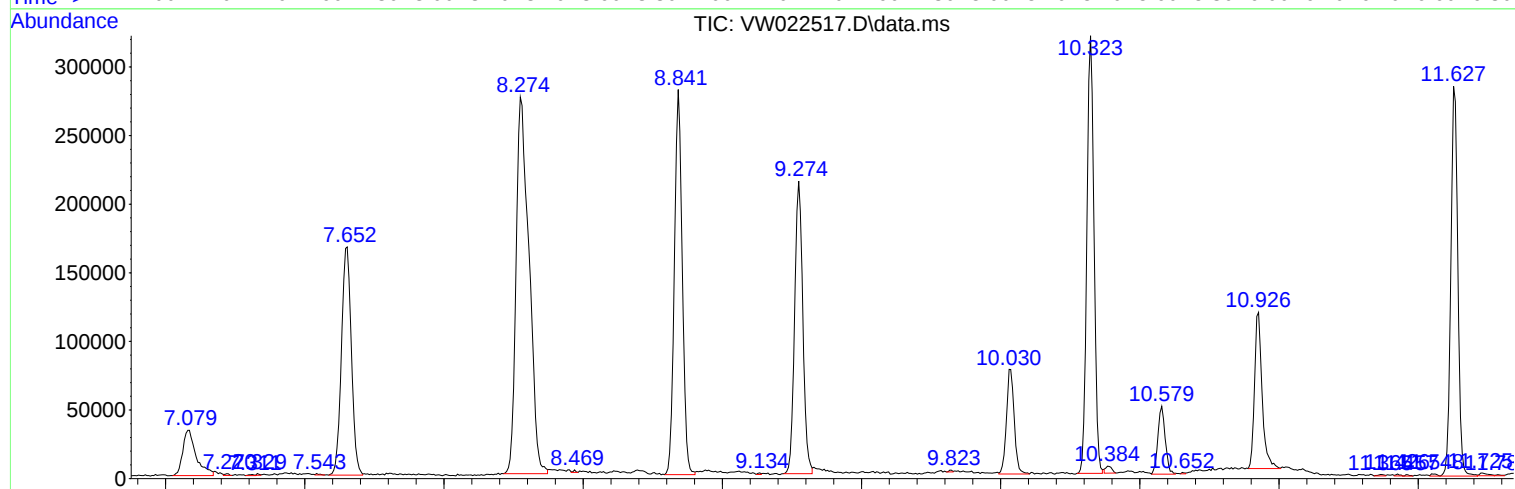
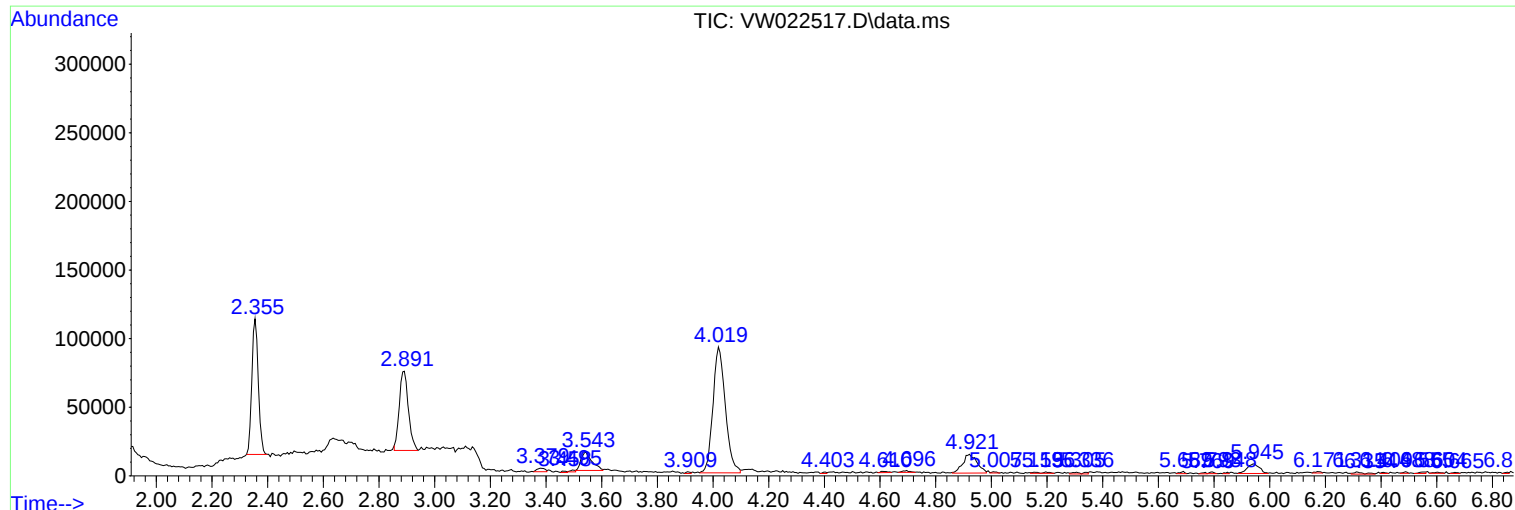
Sum of corrected areas: 7217830

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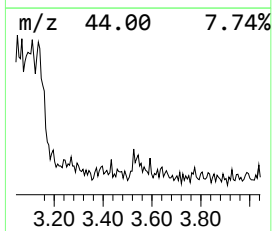
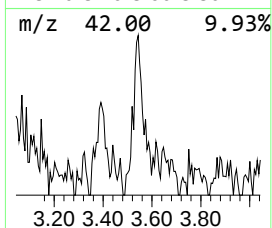
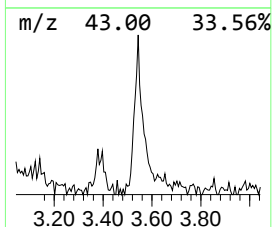
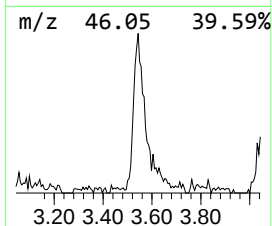
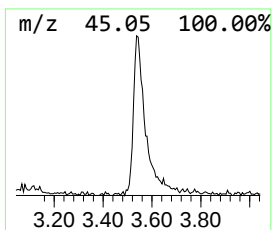
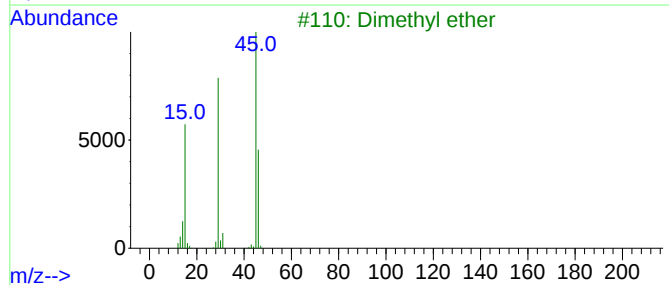
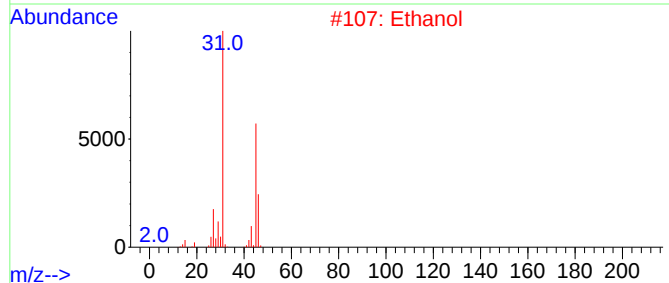
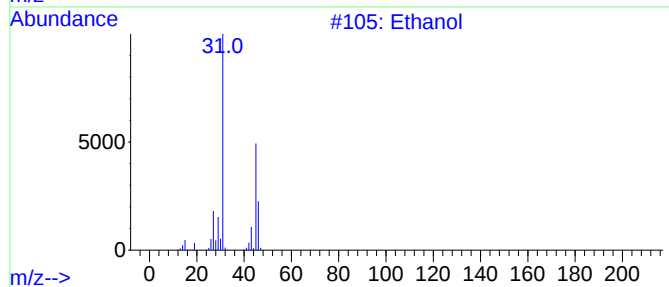
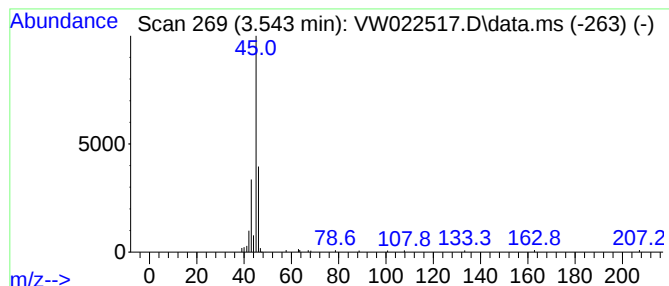
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 Ethanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.543	1.65 ug/L	36245	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol			46	C2H6O	000064-17-5	64
2	Ethanol			46	C2H6O	000064-17-5	64
3	Dimethyl ether			46	C2H6O	000115-10-6	9
4	Ethanol			46	C2H6O	000064-17-5	9
5	Methane, nitroso-			45	CH3NO	000865-40-7	7



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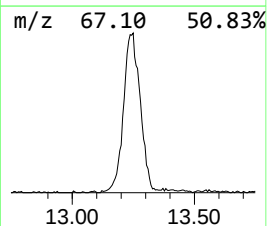
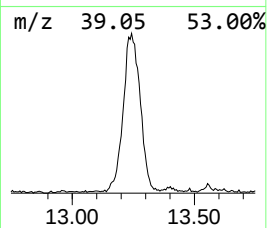
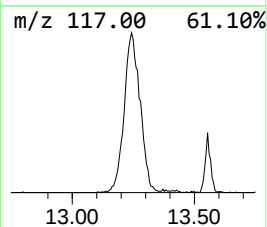
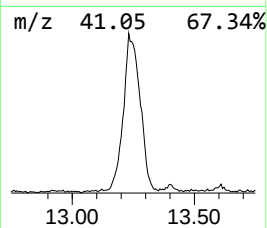
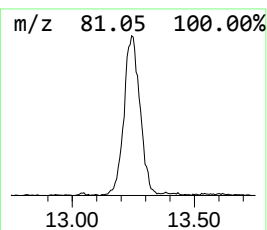
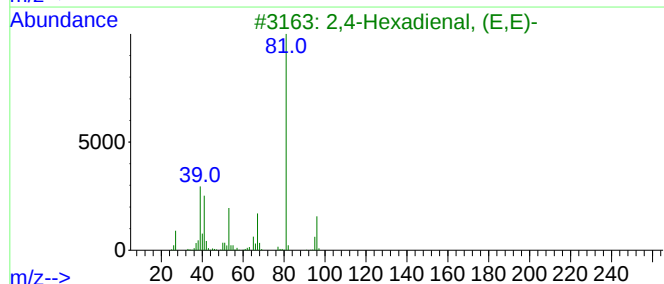
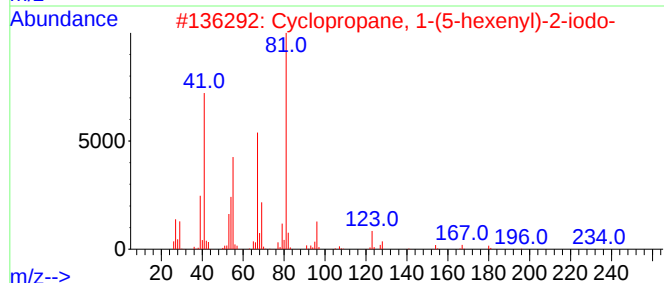
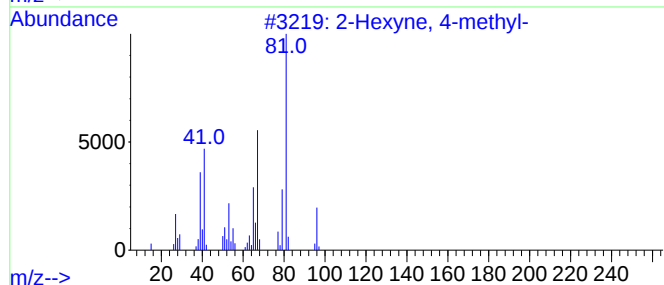
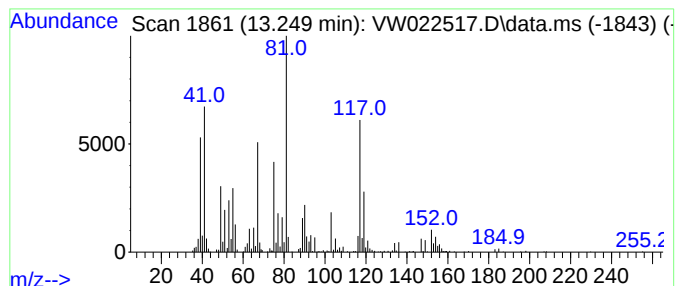
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.249	93.55 ug/L	1070990	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hexyne, 4-methyl-	96	C7H12	020198-49-6	12
2			Cyclopropane, 1-(5-hexenyl)-2-iodo-	250	C9H15I	074685-40-8	12
3			2,4-Hexadienal, (E,E)-	96	C6H8O	000142-83-6	10
4			Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	10
5			Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	10



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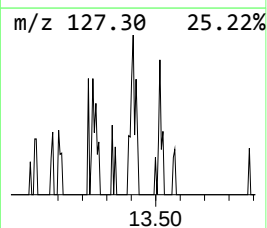
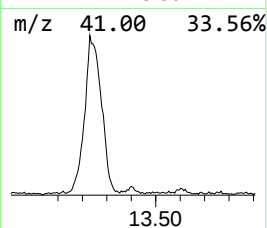
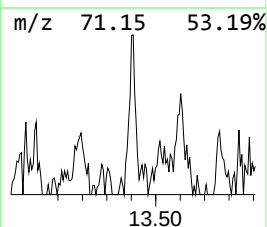
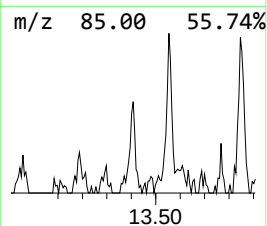
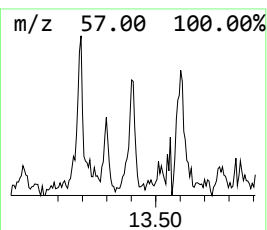
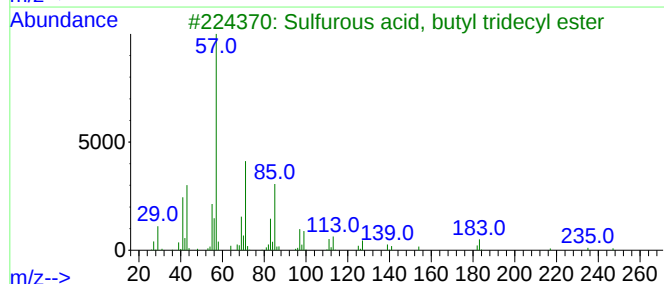
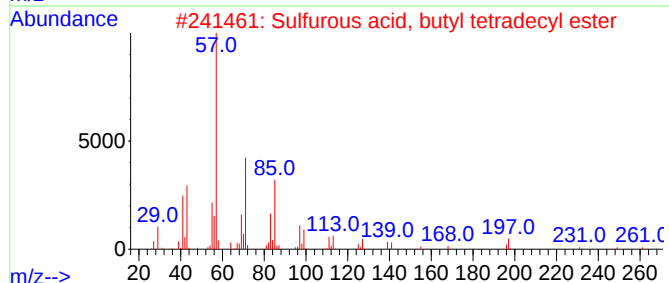
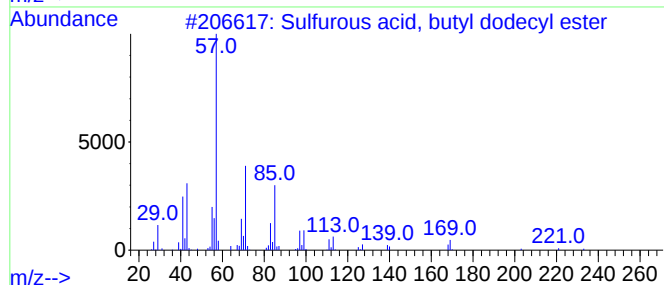
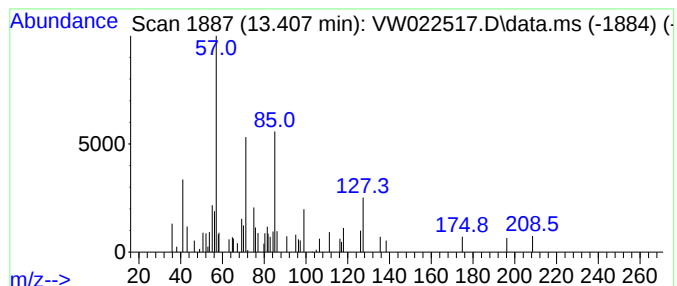
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Quant Title : SFAM01.0

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

Peak Number 6 Sulfurous acid, butyl dodec... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.407	1.43 ug/L	16351	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	53
2			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	50
3			Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	50
4			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	50
5			Heptadecane	240	C17H36	000629-78-7	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

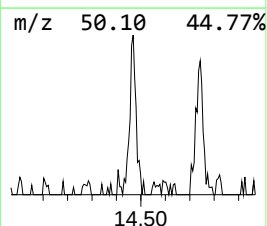
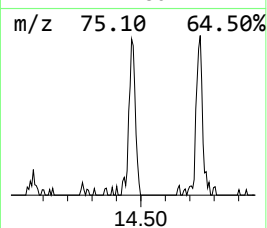
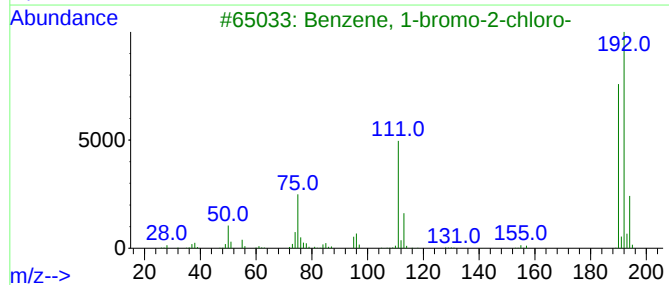
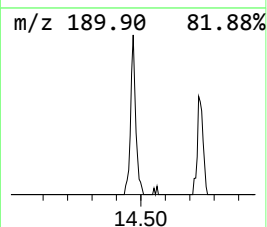
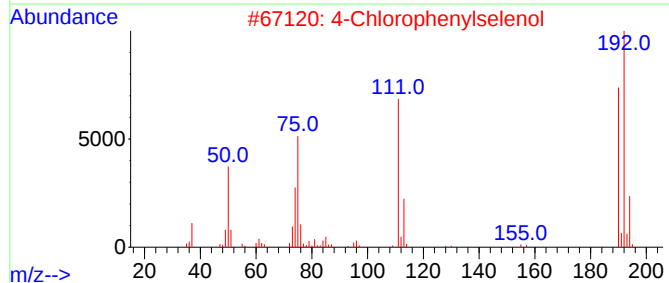
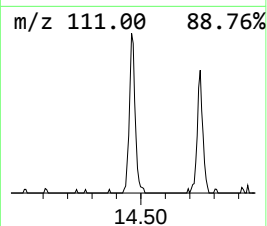
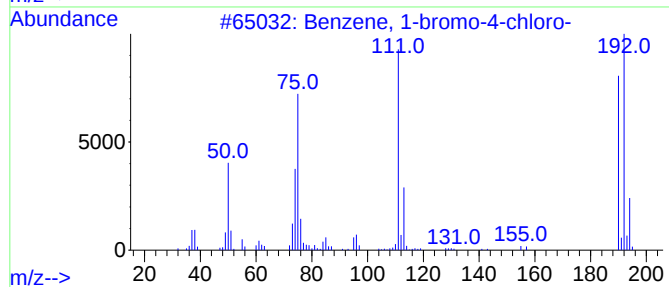
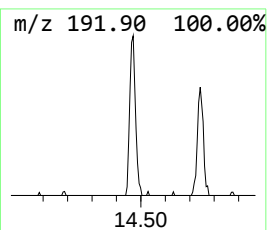
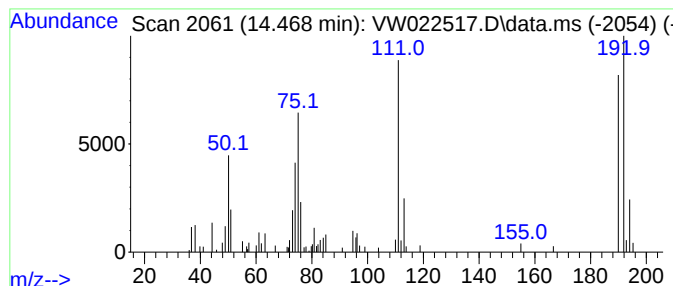
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Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-bromo-4-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.468	2.81 ug/L	32182	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
2			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

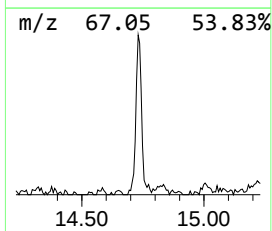
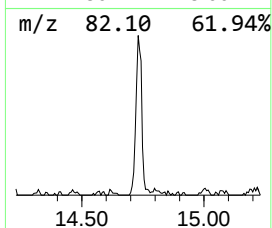
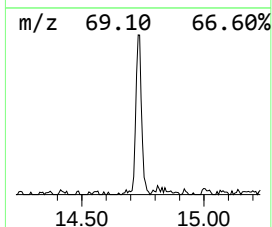
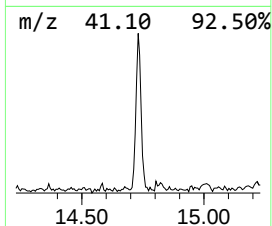
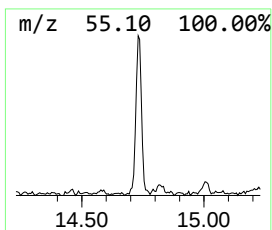
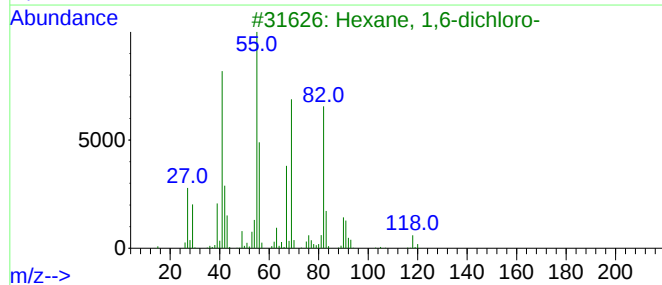
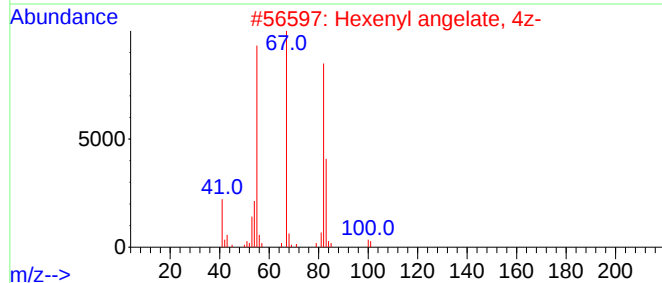
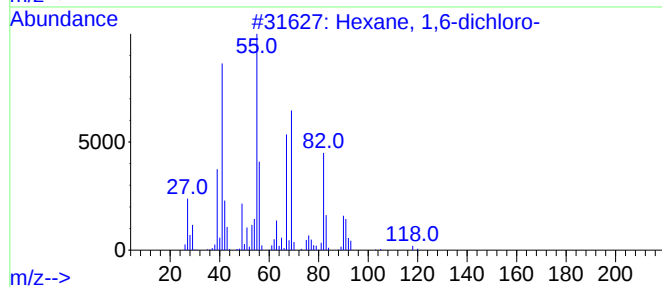
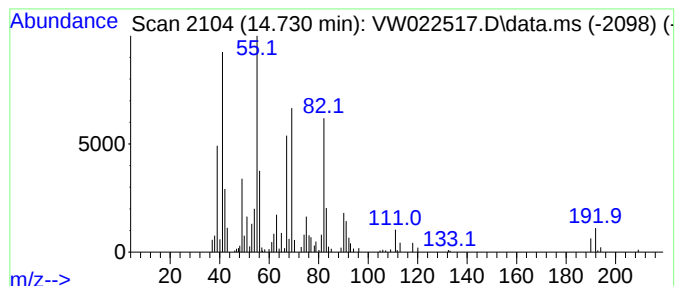
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Quant Title : SFAM01.0

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

Peak Number 9 Hexane, 1,6-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.730	10.46 ug/L	119696	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	80
2			Hexenyl angelate, 4z-	182	C11H18O2	1000383-63-7	38
3			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	38
4			1,1'-Bicyclohexyl	166	C12H22	000092-51-3	38
5			Cyclohexane, chloro-	118	C6H11Cl	000542-18-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

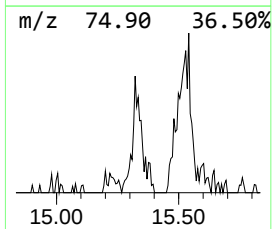
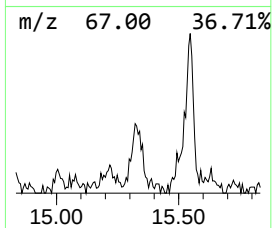
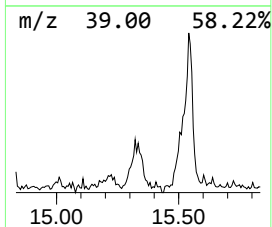
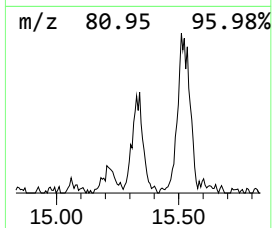
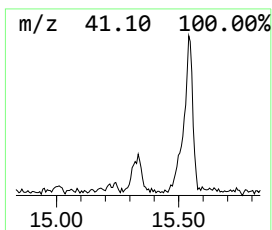
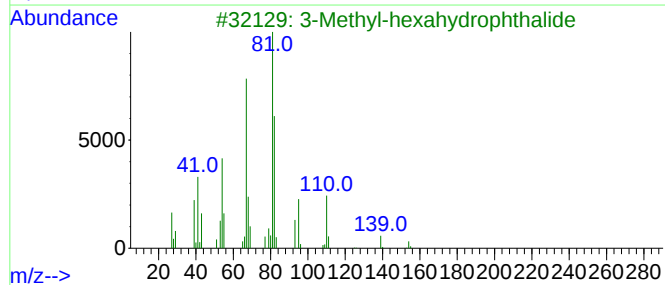
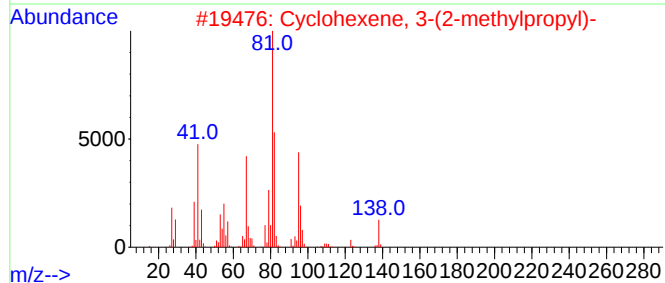
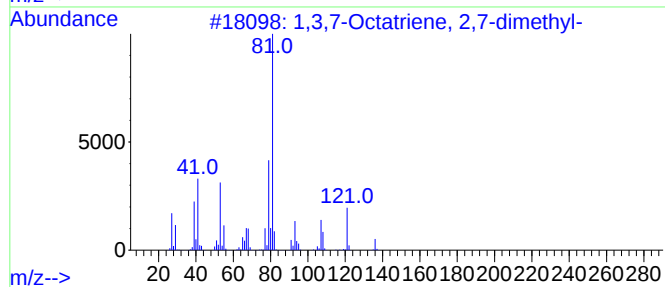
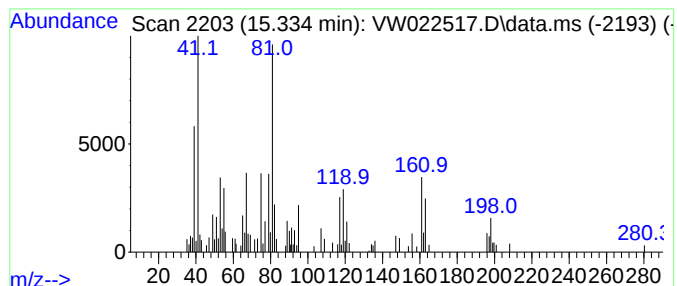
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.334	4.93 ug/L	56393	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,3,7-Octatriene, 2,7-dimethyl-	136	C10H16	036638-38-7	38
2		Cyclohexene, 3-(2-methylpropyl)-	138	C10H18	004104-56-7	35
3		3-Methyl-hexahydrophthalide	154	C9H14O2	002205-25-6	35
4		Cyclohexene, 3-(2-methylpropyl)-	138	C10H18	004104-56-7	35
5		Avocadynofuran	246	C17H26O	024708-33-6	32



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

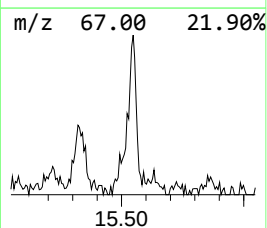
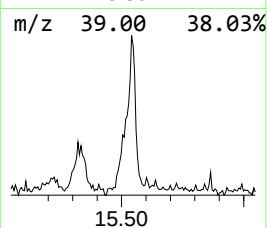
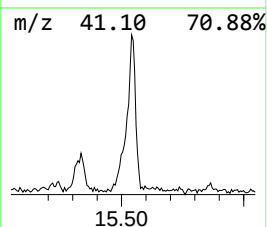
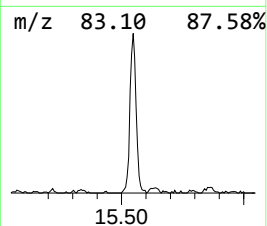
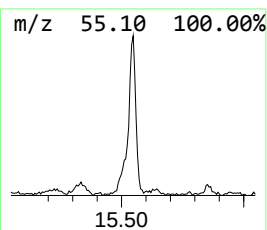
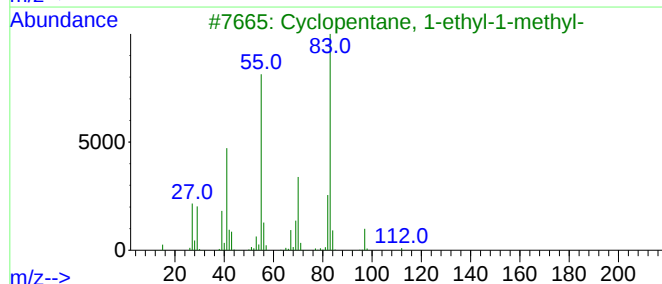
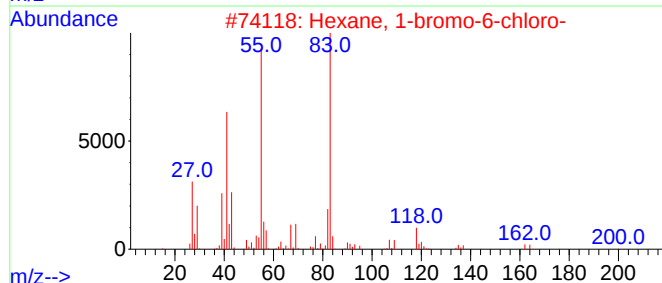
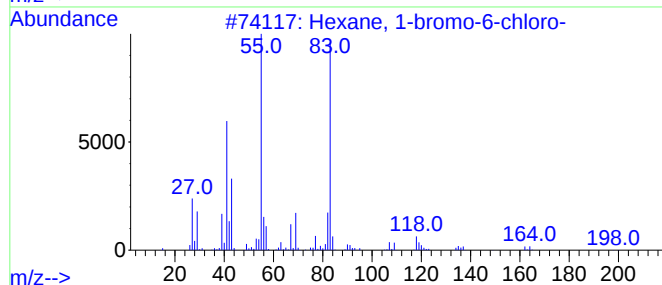
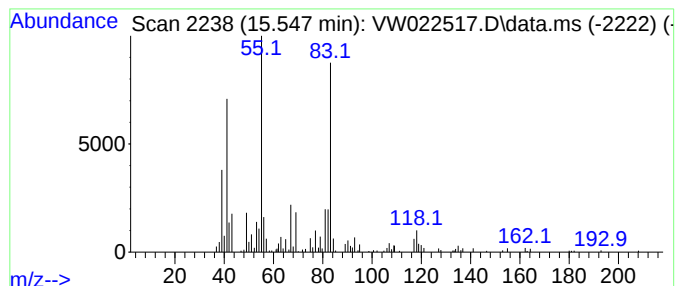
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Quant Title : SFAM01.0

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

Peak Number 11 Hexane, 1-bromo-6-chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.547	15.54 ug/L	177953	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	80
2			Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	72
3			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	53
4			Cyclohexane, pentyl-	154	C11H22	004292-92-6	50
5			Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

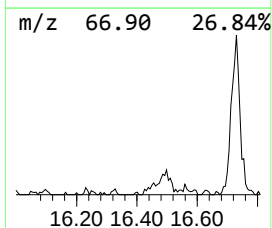
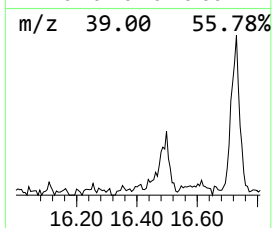
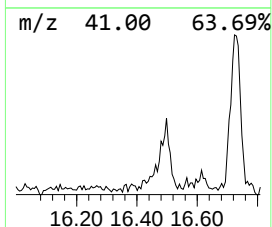
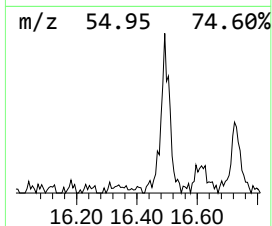
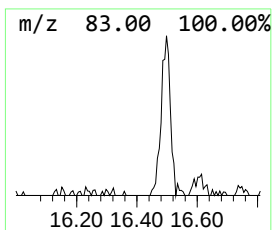
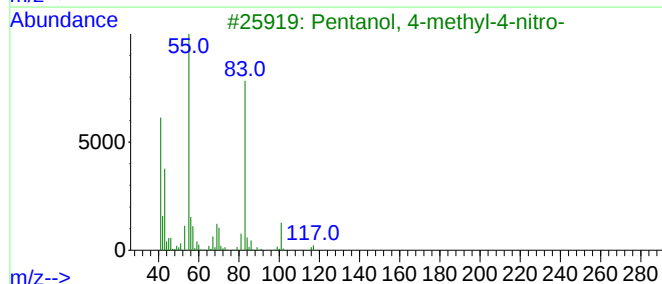
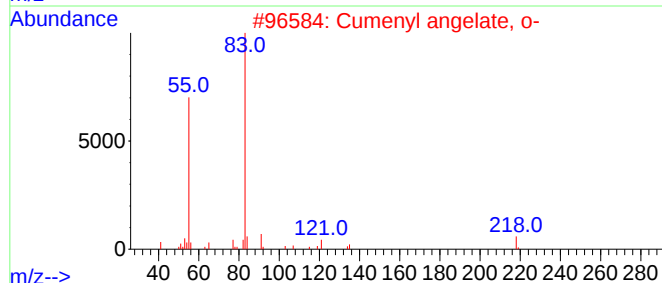
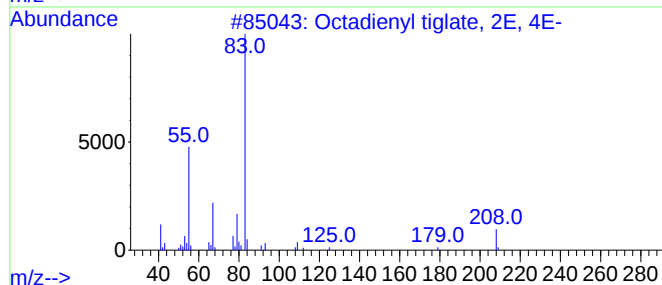
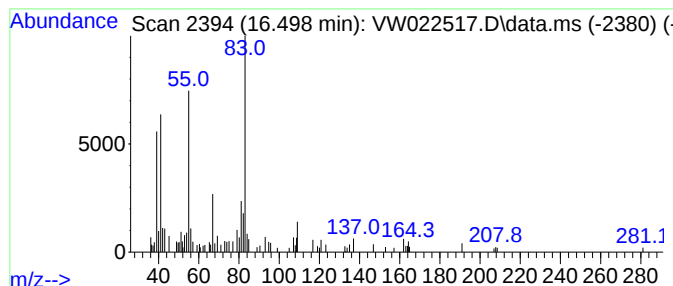
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Quant Title : SFAM01.0

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

Peak Number 12 Octadienyl tiglate, 2E, 4E- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.498	4.32 ug/L	49464	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadienyl tiglate, 2E, 4E-	208	C13H20O2	1000383-60-3	59
2			Cumenyl angelate, o-	218	C14H18O2	1000383-67-2	47
3			Pentanol, 4-methyl-4-nitro-	147	C6H13NO3	005215-92-9	47
4			Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	43
5			Acetaldehyde 2E-hexenyl 3Z-hexen...	226	C14H26O2	1000431-45-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

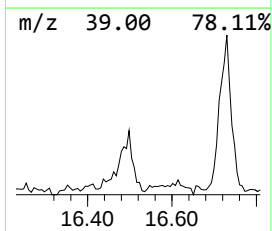
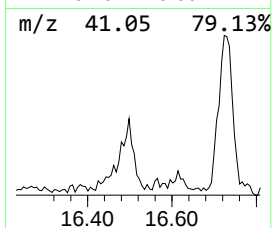
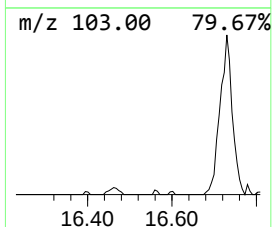
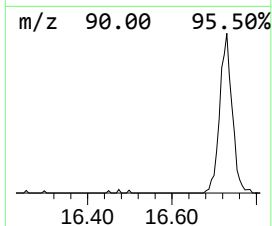
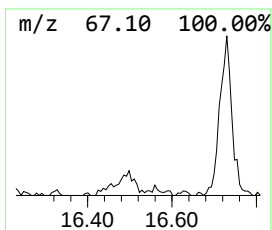
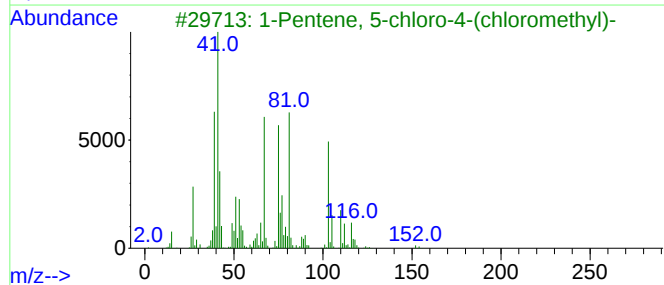
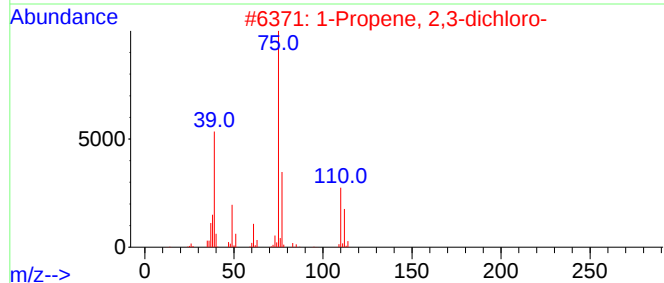
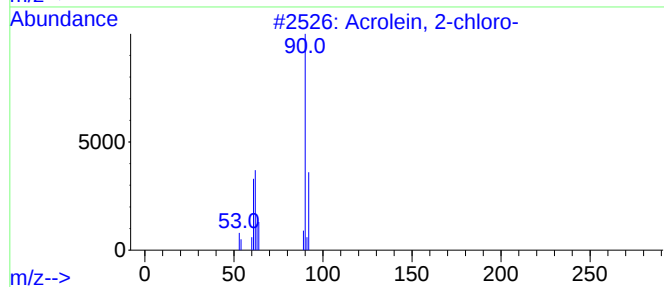
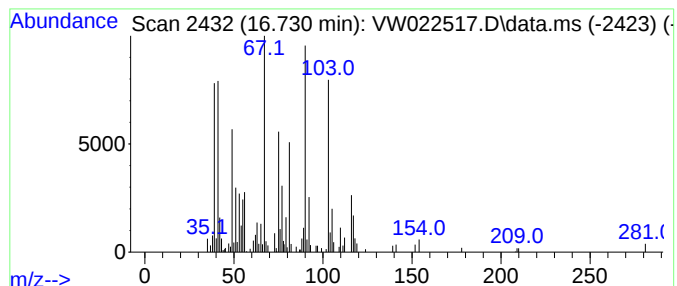
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Quant Title : SFAM01.0

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

Peak Number 13 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.730	9.63 ug/L	110242	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	47
2			1-Propene, 2,3-dichloro-	110	C3H4Cl2	000078-88-6	27
3			1-Pentene, 5-chloro-4-(chloromet...	152	C6H10Cl2	027990-74-5	25
4			1,3-Butadiene, 2,3-dimethyl-	82	C6H10	000513-81-5	22
5			1-Propene, 1,3-dichloro-, (Z)-	110	C3H4Cl2	010061-01-5	16



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022517.D
Acq On : 12 Apr 2022 19:11
Operator : SY/VA
Sample : N2373-01
Misc : 3.29g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFD2

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

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					#	RT	Resp	Conc
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unknown-01	13.249	93.5	ug/L	1070990	3	13.560	286199	25.0
Sulfurous acid,...	13.407	1.4	ug/L	16351	3	13.560	286199	25.0
Benzene, 1-brom...	14.468	2.8	ug/L	32182	3	13.560	286199	25.0
Hexane, 1,6-dic...	14.730	10.5	ug/L	119696	3	13.560	286199	25.0
unknown-02	15.334	4.9	ug/L	56393	3	13.560	286199	25.0
Hexane, 1-bromo...	15.547	15.5	ug/L	177953	3	13.560	286199	25.0
Octadienyl tigl...	16.498	4.3	ug/L	49464	3	13.560	286199	25.0
unknown-03	16.730	9.6	ug/L	110242	3	13.560	286199	25.0