

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
 Data File : VW022489.D
 Acq On : 11 Apr 2022 17:45
 Operator : SY/VA
 Sample : N2316-10
 Misc : 6.46g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNQ9

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: VW022489.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.958	4	9	29	rVB	768800	1142199	8.36%	3.422%
2	2.233	49	54	55	rBV5	3922	4960	0.04%	0.015%
3	2.355	67	74	84	rBV	141924	270468	1.98%	0.810%
4	2.501	93	98	106	rBV	20467	46090	0.34%	0.138%
5	2.629	114	119	125	rBV5	11843	33272	0.24%	0.100%
6	2.891	155	162	172	rVB	93387	204178	1.50%	0.612%
7	3.019	180	183	184	rBV3	3972	4277	0.03%	0.013%
8	3.391	237	244	249	rBV6	3167	8665	0.06%	0.026%
9	3.684	290	292	296	rVV5	2601	4379	0.03%	0.013%
10	3.903	322	328	330	rVV7	2070	4460	0.03%	0.013%
11	4.019	336	347	360	rVV	174523	546834	4.00%	1.639%
12	4.129	360	365	372	rVV	8796	24567	0.18%	0.074%
13	4.208	374	378	387	rVV	3883	12188	0.09%	0.037%
14	4.391	405	408	418	rVB3	7057	17060	0.12%	0.051%
15	5.190	530	539	547	rVB7	6329	22347	0.16%	0.067%
16	5.421	572	577	580	rBV5	2247	4645	0.03%	0.014%
17	5.561	592	600	602	rBV6	4941	10960	0.08%	0.033%
18	5.946	653	663	670	rBV5	5734	17990	0.13%	0.054%
19	6.122	687	692	699	rVB9	2426	6336	0.05%	0.019%
20	6.311	717	723	732	rVB9	2254	6577	0.05%	0.020%
21	7.000	827	836	841	rBV8	7736	24253	0.18%	0.073%
22	7.073	841	848	862	rBV2	76426	215294	1.58%	0.645%
23	7.646	932	942	958	rBV	307519	775454	5.68%	2.324%
24	7.951	989	992	1003	rVB7	6280	14299	0.10%	0.043%
25	8.152	1018	1025	1026	rBV4	3100	5021	0.04%	0.015%
26	8.274	1034	1045	1047	rBV	562899	1095909	8.03%	3.284%
27	8.323	1048	1053	1065	rVB	1351048	3132603	22.94%	9.386%
28	8.561	1088	1092	1101	rVB6	8839	21516	0.16%	0.064%
29	8.695	1108	1114	1122	rBV3	5556	11993	0.09%	0.036%
30	8.841	1126	1138	1148	rBV	550628	1080911	7.92%	3.239%
31	8.988	1156	1162	1167	rBV2	14049	27603	0.20%	0.083%
32	9.091	1167	1179	1188	rVB10	21864	63992	0.47%	0.192%
33	9.274	1199	1209	1216	rBV	421412	863470	6.32%	2.587%
34	9.335	1216	1219	1233	rVB5	20286	48541	0.36%	0.145%
35	9.506	1240	1247	1255	rBV2	19499	38008	0.28%	0.114%
36	9.664	1268	1273	1279	rBV10	2855	5236	0.04%	0.016%

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

37	9.902	1306	1312	1315	rBV6	3629	7386	0.05%	0.022%
38	9.963	1315	1322	1328	rBV2	103970	199502	1.46%	0.598%
39	10.036	1328	1334	1342	rVV	233923	424932	3.11%	1.273%
40	10.250	1365	1369	1373	rVV5	4805	7541	0.06%	0.023%

41	10.323	1373	1381	1387	rVV	816929	1443580	10.57%	4.325%
42	10.384	1387	1391	1399	rVV2	180085	324290	2.37%	0.972%
43	10.500	1401	1410	1415	rVB2	14067	37205	0.27%	0.111%
44	10.579	1415	1423	1431	rBV	157412	273031	2.00%	0.818%
45	10.676	1433	1439	1440	rVV5	3674	7115	0.05%	0.021%

46	10.719	1440	1446	1458	rVB3	44862	100754	0.74%	0.302%
47	10.926	1472	1480	1490	rBV2	393882	717303	5.25%	2.149%
48	11.060	1497	1502	1507	rBV9	7306	14295	0.10%	0.043%
49	11.176	1517	1521	1527	rVB8	4458	8418	0.06%	0.025%
50	11.292	1534	1540	1545	rVV2	20018	39952	0.29%	0.120%

51	11.481	1566	1571	1577	rVB5	10557	19068	0.14%	0.057%
52	11.560	1577	1584	1588	rVB7	10013	19042	0.14%	0.057%
53	11.658	1588	1600	1607	rBV	7313943	13654824	100.00%	40.915%
54	11.725	1607	1611	1619	rVV2	50485	108179	0.79%	0.324%
55	11.841	1621	1630	1638	rVB	42571	86626	0.63%	0.260%

56	11.969	1646	1651	1654	rBV7	4170	8370	0.06%	0.025%
57	12.127	1673	1677	1678	rBV4	3656	4928	0.04%	0.015%
58	12.164	1678	1683	1690	rVB	45437	80337	0.59%	0.241%
59	12.347	1708	1713	1717	rVV8	3131	6761	0.05%	0.020%
60	12.463	1727	1732	1737	rVB3	10971	20613	0.15%	0.062%

61	12.603	1750	1755	1759	rBV3	19011	33282	0.24%	0.100%
62	12.688	1759	1769	1774	rBV	442449	816856	5.98%	2.448%
63	12.743	1774	1778	1784	rVB	284599	450790	3.30%	1.351%
64	12.798	1784	1787	1793	rVB	13931	20418	0.15%	0.061%
65	12.889	1793	1802	1807	rBV2	91655	186290	1.36%	0.558%

66	12.987	1813	1818	1828	rVB2	34484	65264	0.48%	0.196%
67	13.103	1832	1837	1845	rVB3	28144	57179	0.42%	0.171%
68	13.188	1847	1851	1855	rBV6	5496	10851	0.08%	0.033%
69	13.249	1855	1861	1864	rBV	27045	46881	0.34%	0.140%
70	13.554	1904	1911	1920	rBV	912758	1515541	11.10%	4.541%

71	13.639	1920	1925	1934	rVB3	39105	74914	0.55%	0.224%
72	13.712	1934	1937	1941	rBV5	8862	17442	0.13%	0.052%
73	13.749	1941	1943	1947	rVB5	8032	9392	0.07%	0.028%
74	13.853	1953	1960	1970	rVB	826147	1427816	10.46%	4.278%
75	13.956	1971	1977	1980	rVV7	16441	39669	0.29%	0.119%

76	13.987	1980	1982	1988	rVV2	24193	39283	0.29%	0.118%
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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
 Title : SFAM01.0

77	14.066	1991	1995	2005	rVB2	17670	42558	0.31%	0.128%
78	14.200	2008	2017	2027	rBV7	21169	54096	0.40%	0.162%
79	14.334	2032	2039	2045	rVB4	12390	30480	0.22%	0.091%
80	14.468	2054	2061	2067	rBV	183579	309980	2.27%	0.929%
81	14.523	2067	2070	2072	rVB4	4872	5757	0.04%	0.017%
82	14.627	2084	2087	2092	rVB7	6390	10444	0.08%	0.031%
83	14.682	2093	2096	2100	rVB6	5524	6571	0.05%	0.020%
84	14.737	2101	2105	2109	rBV7	9104	13987	0.10%	0.042%
85	14.846	2120	2123	2125	rBV3	5437	7671	0.06%	0.023%
86	14.950	2138	2140	2148	rVB9	8655	13165	0.10%	0.039%
87	15.084	2152	2162	2164	rBV10	16651	48453	0.35%	0.145%
88	15.249	2186	2189	2196	rVB6	18976	32946	0.24%	0.099%
89	15.334	2197	2203	2204	rBV6	8471	13640	0.10%	0.041%
90	15.419	2213	2217	2220	rBV6	5768	10495	0.08%	0.031%
91	15.487	2220	2228	2235	rVB2	55377	123829	0.91%	0.371%
92	15.615	2239	2249	2261	rVB7	21564	84719	0.62%	0.254%
93	15.736	2262	2269	2279	rBV2	76669	168782	1.24%	0.506%
94	16.114	2321	2331	2337	rBV10	14057	48038	0.35%	0.144%
95	16.267	2350	2356	2365	rBV6	27948	58728	0.43%	0.176%
96	16.352	2365	2370	2371	rBV5	4691	6688	0.05%	0.020%
97	16.529	2396	2399	2404	rVB7	8226	14675	0.11%	0.044%
98	16.590	2404	2409	2410	rBV4	6000	7214	0.05%	0.022%
99	16.633	2413	2416	2417	rBV3	4007	4325	0.03%	0.013%
100	16.730	2427	2432	2437	rBV8	7725	14165	0.10%	0.042%

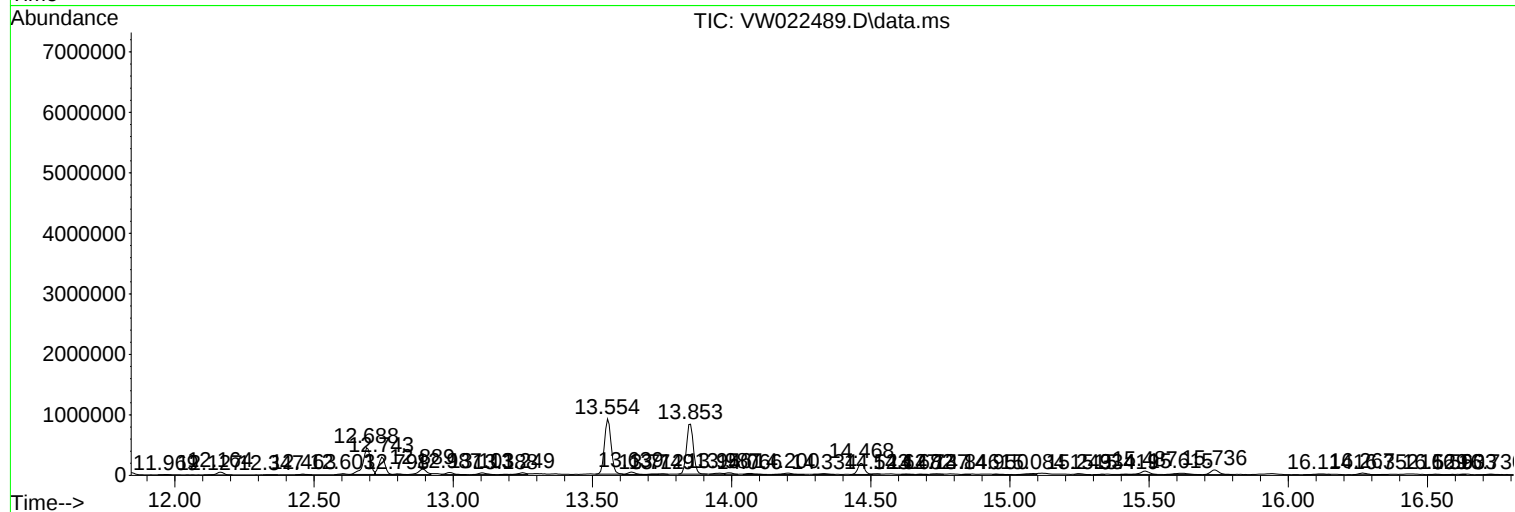
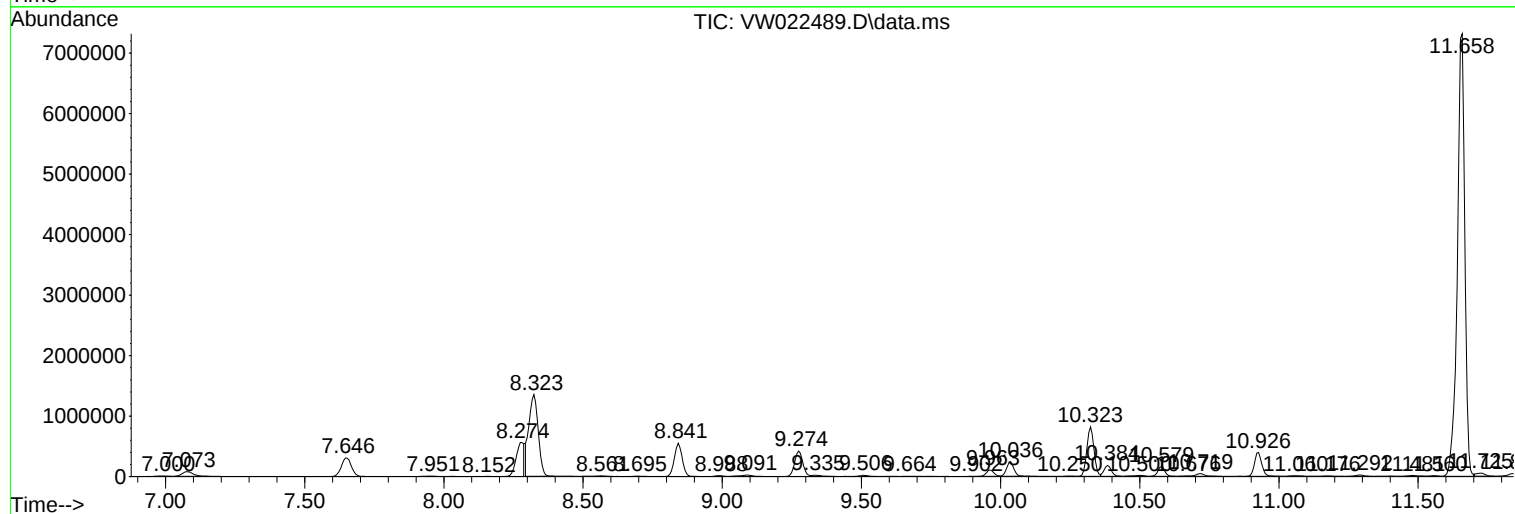
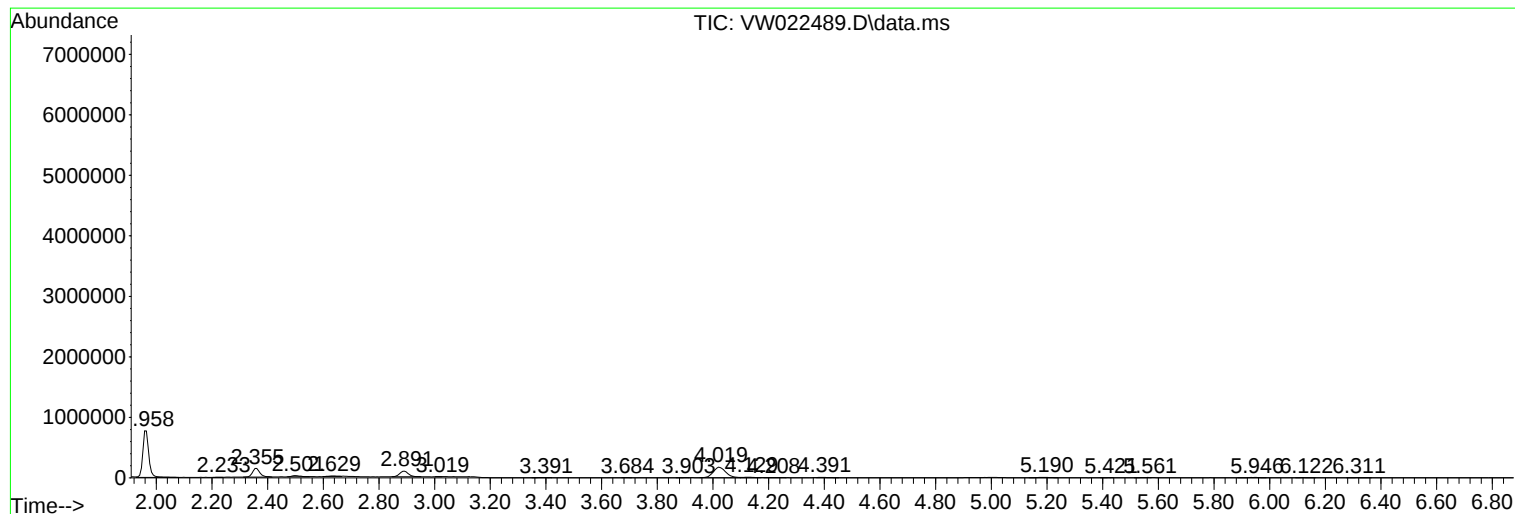
Sum of corrected areas: 33373881

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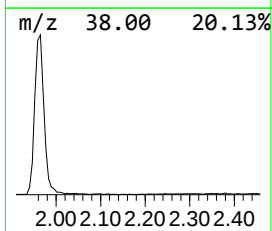
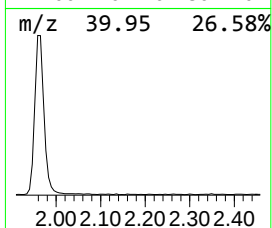
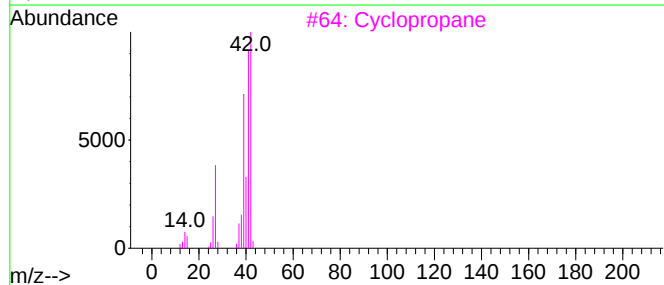
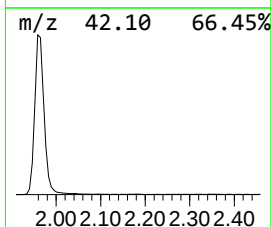
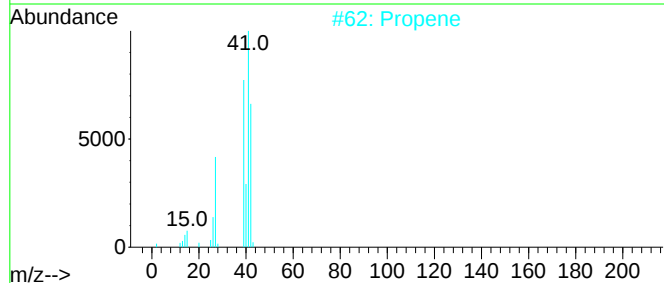
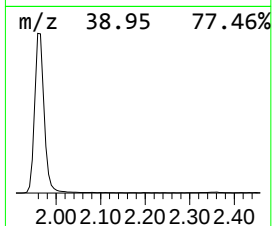
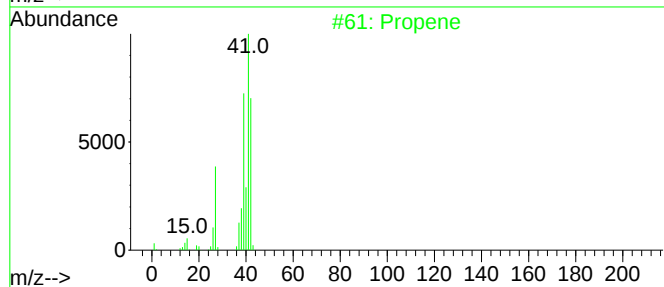
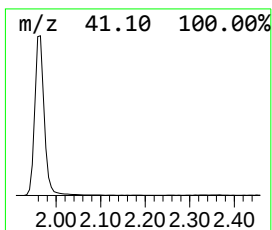
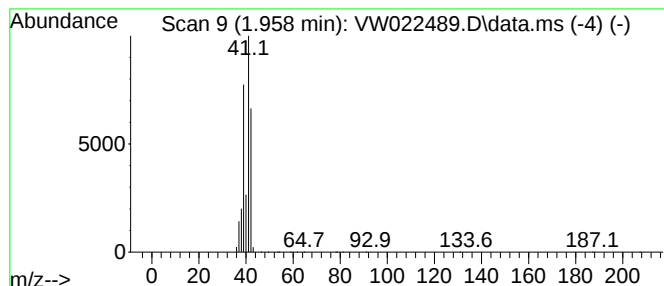
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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.
1.958	26.42 ug/L	1142200	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	64
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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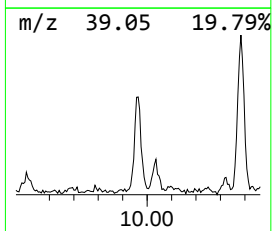
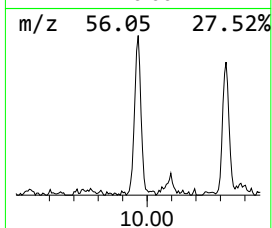
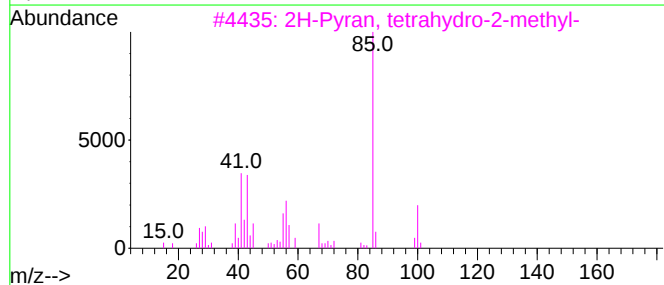
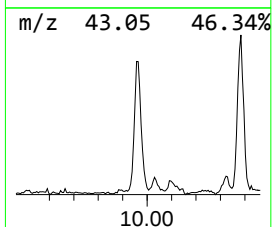
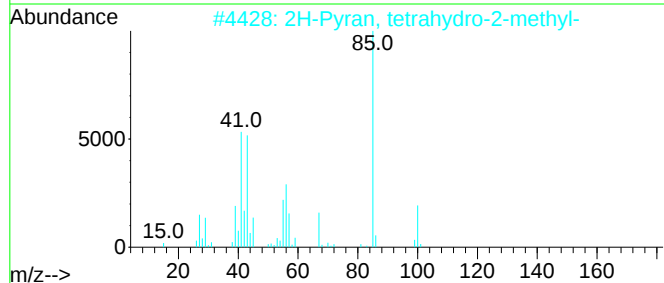
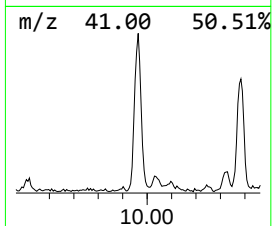
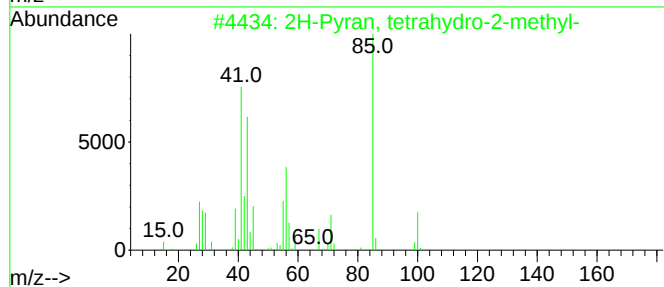
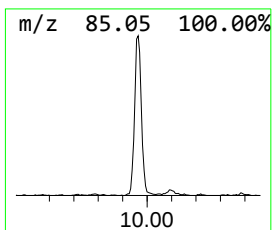
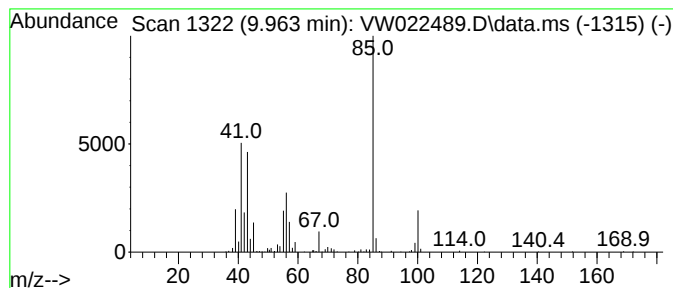
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	4.61 ug/L	199502	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	91
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	87
3	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	80
4	2-Piperazinecarboxamide			129	C5H11N3O	084501-64-4	59
5	1,3-Butadiene, 1-(methylthio)-			100	C5H8S	010574-97-7	58



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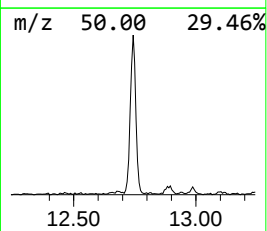
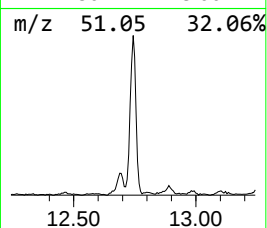
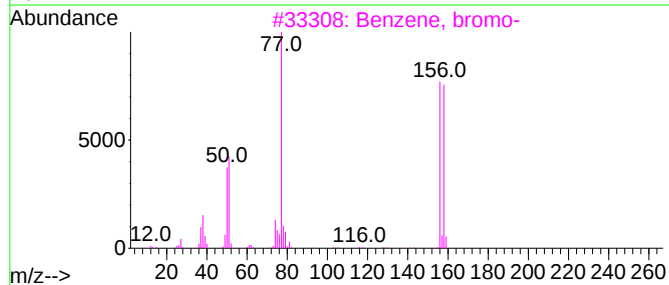
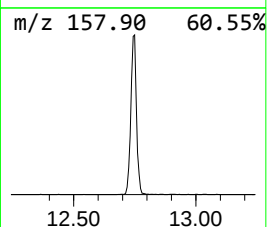
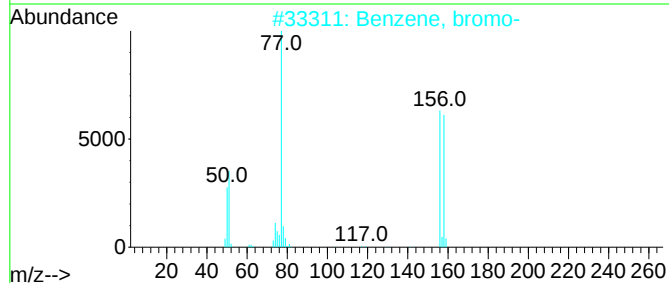
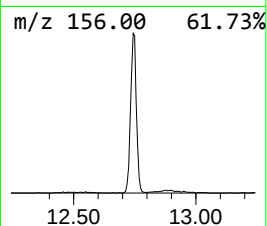
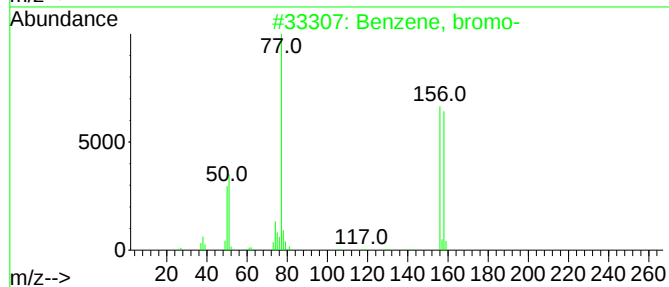
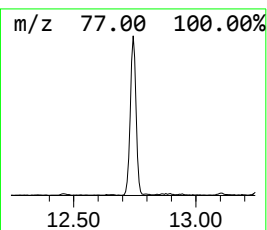
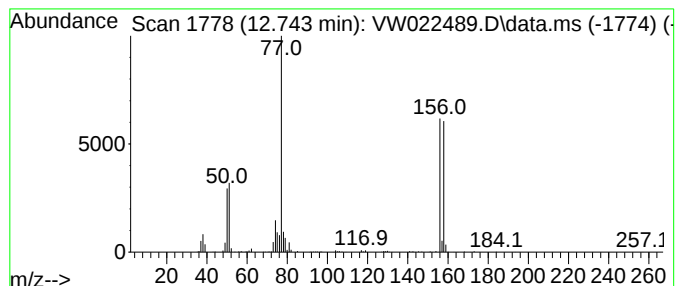
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Quant Title : SFAM01.0

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

Peak Number 4 Benzene, bromo- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.743	7.44 ug/L	450790	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, bromo-	156	C6H5Br	000108-86-1	97
2			Benzene, bromo-	156	C6H5Br	000108-86-1	96
3			Benzene, bromo-	156	C6H5Br	000108-86-1	95
4			Benzene, bromo-	156	C6H5Br	000108-86-1	94
5			Benzene, bromo-	156	C6H5Br	000108-86-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022489.D
Acq On : 11 Apr 2022 17:45
Operator : SY/VA
Sample : N2316-10
Misc : 6.46g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ9

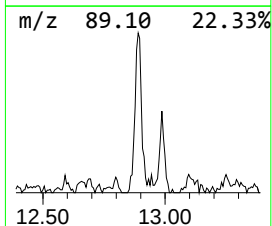
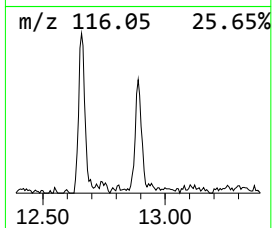
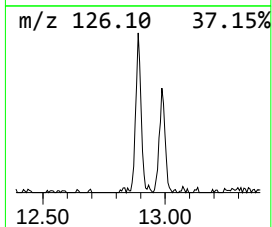
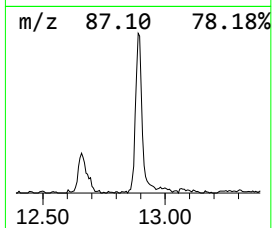
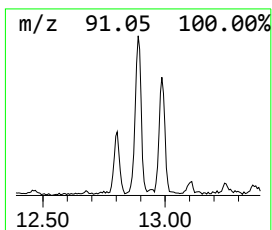
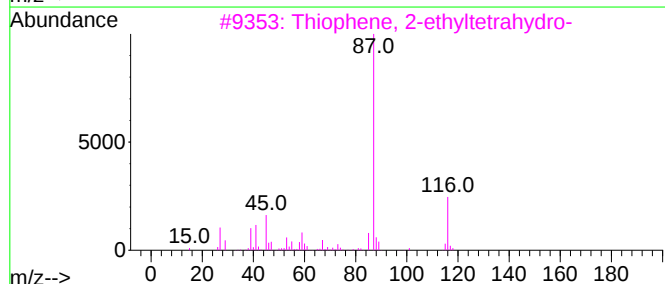
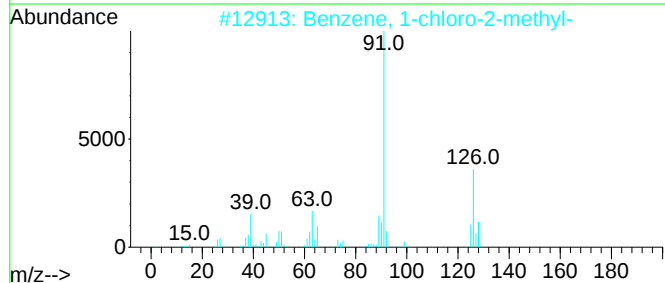
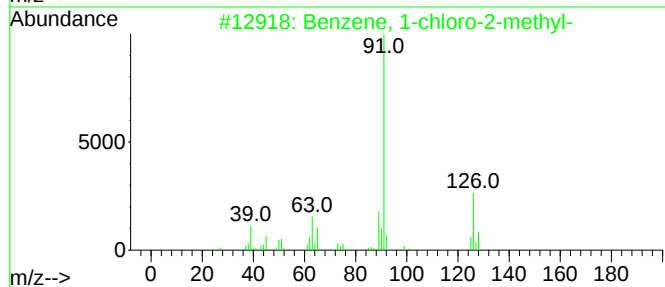
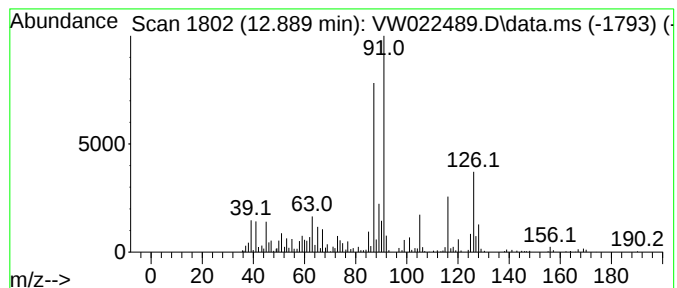
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Quant Title : SFAM01.0

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

Peak Number 5 Benzene, 1-chloro-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.889	3.07 ug/L	186290	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	80
2			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	72
3			Thiophene, 2-ethyltetrahydro-	116	C6H12S	001551-32-2	53
4			2-Propenenitrile, 2-chloro-	87	C3H2ClN	000920-37-6	43
5			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022489.D
Acq On : 11 Apr 2022 17:45
Operator : SY/VA
Sample : N2316-10
Misc : 6.46g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ9

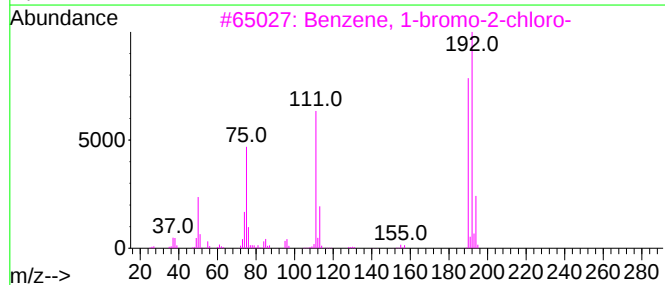
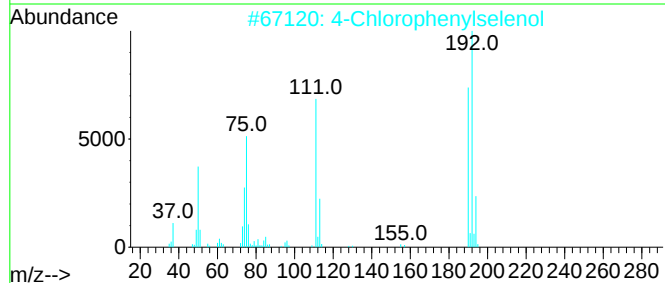
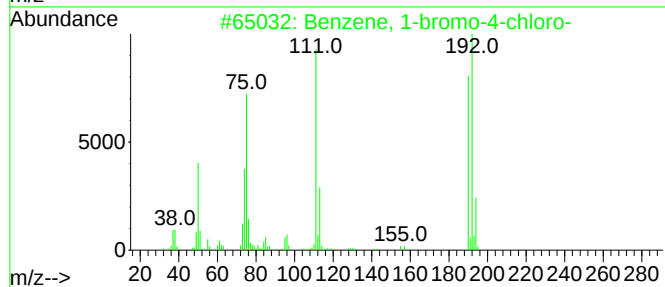
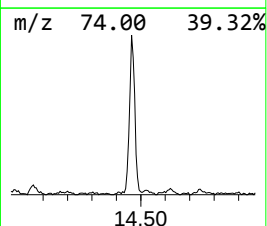
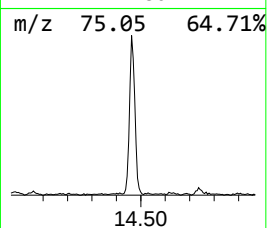
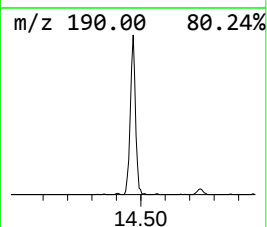
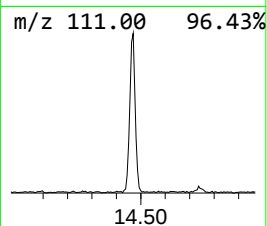
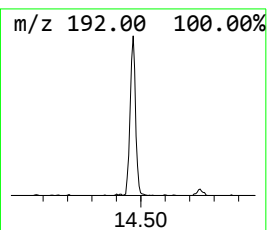
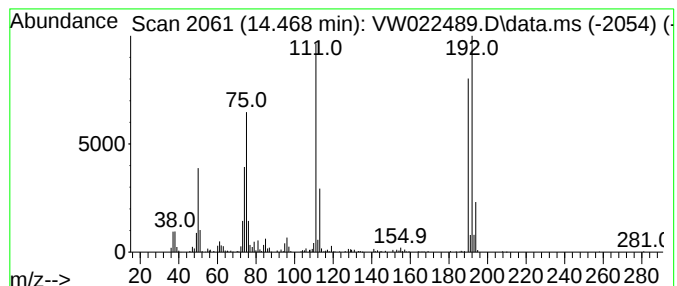
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Quant Title : SFAM01.0

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

Peak Number 6 Benzene, 1-bromo-4-chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.468	5.11 ug/L	309980	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	99
2			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	98
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	97
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022489.D
Acq On : 11 Apr 2022 17:45
Operator : SY/VA
Sample : N2316-10
Misc : 6.46g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ9

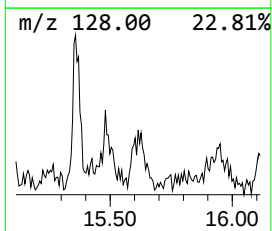
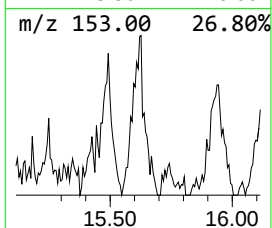
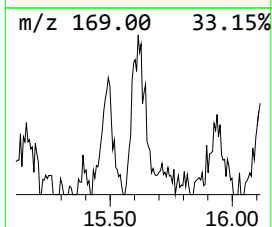
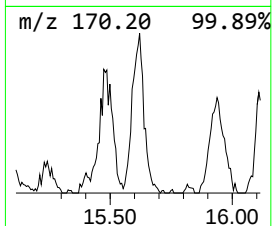
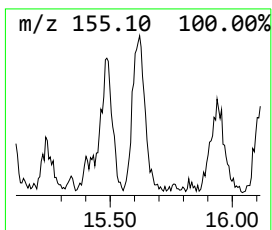
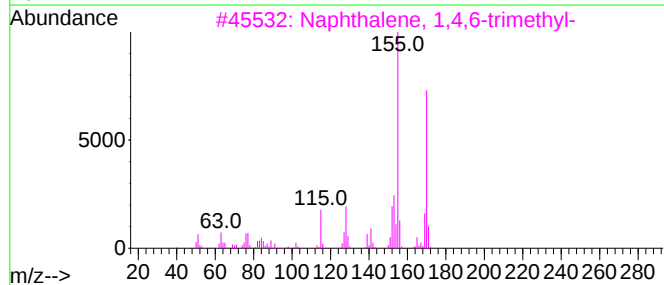
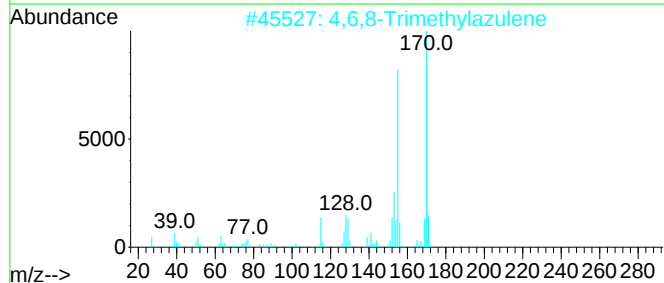
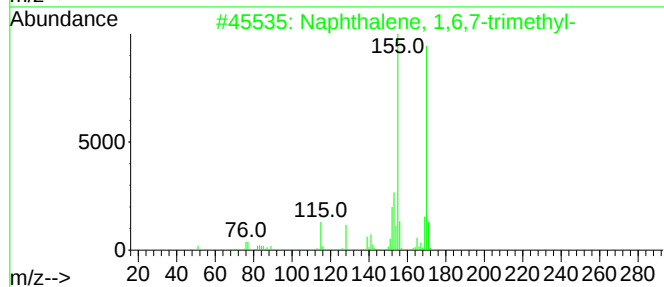
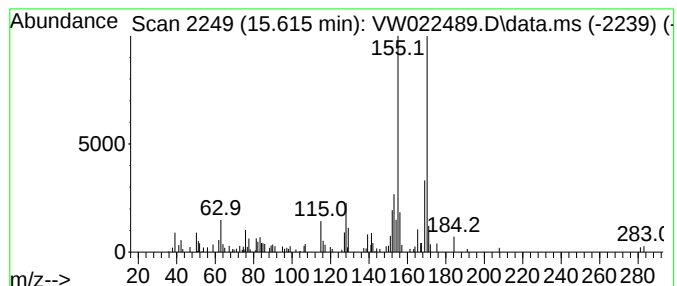
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Quant Title : SFAM01.0

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

Peak Number 8 Naphthalene, 1,6,7-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	1.40 ug/L	84719	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
2			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	93
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022489.D
Acq On : 11 Apr 2022 17:45
Operator : SY/VA
Sample : N2316-10
Misc : 6.46g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ9

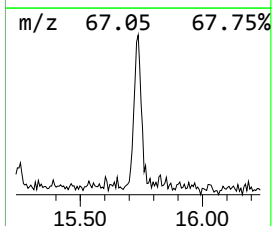
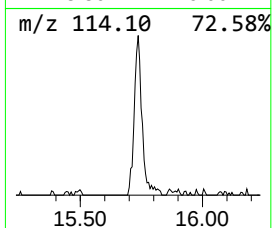
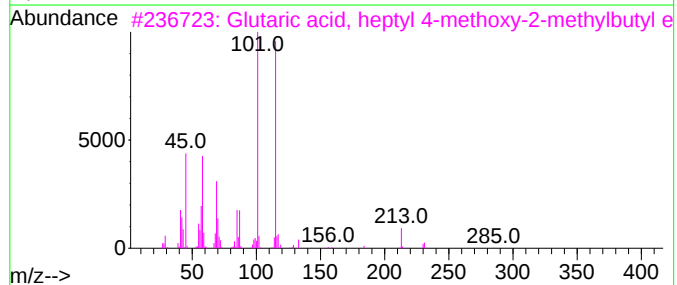
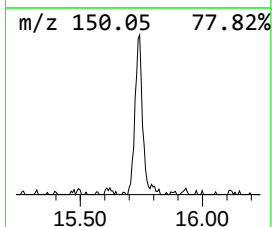
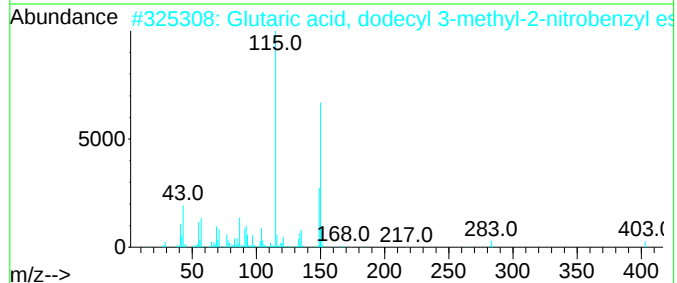
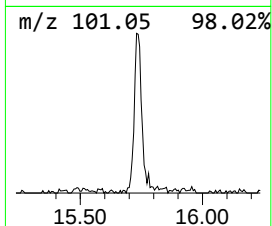
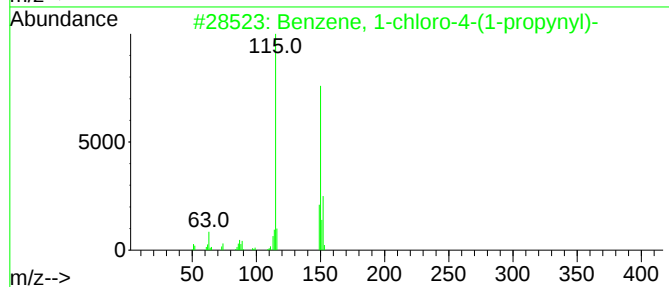
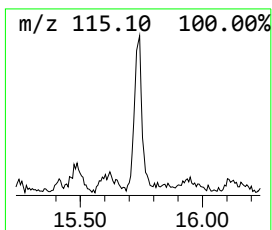
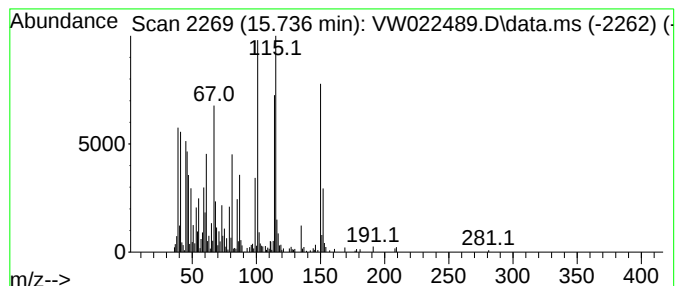
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Quant Title : SFAM01.0

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

Peak Number 9 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.736	2.78 ug/L	168782	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	38
2			Glutaric acid, dodecyl 3-methyl-...	449	C25H39NO6	1000376-74-2	14
3			Glutaric acid, heptyl 4-methoxy-...	330	C18H34O5	1000359-41-6	14
4			Glutaric acid, isobutyl 4-methox...	288	C15H28O5	1000359-41-1	12
5			Glutaric acid, 4-methoxy-2-methy...	302	C16H30O5	1000359-41-3	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041122\
Data File : VW022489.D
Acq On : 11 Apr 2022 17:45
Operator : SY/VA
Sample : N2316-10
Misc : 6.46g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNQ9

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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2H-Pyran, tetra...	9.963	4.6	ug/L	199502	1	8.841	1080910	25.0
Benzene, bromo-	12.743	7.4	ug/L	450790	3	13.554	1515540	25.0
Benzene, 1-chlo...	12.889	3.1	ug/L	186290	3	13.554	1515540	25.0
Benzene, 1-brom...	14.468	5.1	ug/L	309980	3	13.554	1515540	25.0
Naphthalene, 1,...	15.615	1.4	ug/L	84719	3	13.554	1515540	25.0
unknown-01	15.736	2.8	ug/L	168782	3	13.554	1515540	25.0