

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051922\
 Data File : VW023181.D
 Acq On : 19 May 2022 18:33
 Operator : SY/VA
 Sample : N2929-01
 Misc : 6.89g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 DBTF7

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\SFAMWLM050222SMA.M
 Title : SFAM01.0

Signal : TIC: VW023181.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.172	42	44	49	rBV	6375	6960	0.22%	0.052%
2	2.349	68	73	82	rVB	161842	273751	8.71%	2.050%
3	2.885	154	161	170	rBV	95261	203766	6.49%	1.526%
4	3.020	178	183	194	rVB2	23531	54570	1.74%	0.409%
5	3.294	224	228	235	rVB9	2681	5135	0.16%	0.038%
6	3.385	235	243	250	rBV3	9393	20569	0.65%	0.154%
7	4.019	336	347	358	rBV2	149371	454333	14.46%	3.403%
8	4.129	360	365	372	rVB	11581	29187	0.93%	0.219%
9	4.385	405	407	408	rVV2	1423	1287	0.04%	0.010%
10	4.909	483	493	494	rBV5	5544	11904	0.38%	0.089%
11	5.354	562	566	568	rBV4	1048	1394	0.04%	0.010%
12	5.385	568	571	575	rBV4	1250	2581	0.08%	0.019%
13	5.623	605	610	613	rBV4	1077	1998	0.06%	0.015%
14	5.897	650	655	656	rVV3	1105	1239	0.04%	0.009%
15	5.946	656	663	671	rVB6	4166	12444	0.40%	0.093%
16	6.019	671	675	678	rVB6	1520	1727	0.05%	0.013%
17	6.299	715	721	723	rBV6	1348	2004	0.06%	0.015%
18	6.610	770	772	776	rVB4	1859	1792	0.06%	0.013%
19	6.976	829	832	833	rBV2	2279	2136	0.07%	0.016%
20	7.080	840	849	858	rBV	44043	130595	4.16%	0.978%
21	7.409	900	903	905	rBV4	1107	1333	0.04%	0.010%
22	7.482	913	915	916	rBV2	1340	1225	0.04%	0.009%
23	7.647	931	942	955	rBV	226366	553032	17.60%	4.142%
24	7.830	969	972	975	rBV4	1191	1414	0.05%	0.011%
25	8.275	1033	1045	1061	rBV2	432802	1350389	42.98%	10.114%
26	8.641	1103	1105	1108	rVB4	1245	1214	0.04%	0.009%
27	8.701	1108	1115	1120	rBV10	3392	9063	0.29%	0.068%
28	8.842	1127	1138	1147	rBV	435259	847607	26.98%	6.349%
29	9.086	1172	1178	1180	rBV6	2854	5639	0.18%	0.042%
30	9.275	1196	1209	1226	rBV	295376	626384	19.94%	4.692%
31	9.439	1231	1236	1238	rBV4	1913	2807	0.09%	0.021%
32	9.500	1245	1246	1249	rBV3	1268	1540	0.05%	0.012%
33	9.604	1260	1263	1267	rVV6	1200	1816	0.06%	0.014%
34	9.659	1267	1272	1273	rBV5	1462	1937	0.06%	0.015%
35	9.756	1284	1288	1292	rBV4	1188	2438	0.08%	0.018%
36	9.957	1317	1321	1324	rVB5	1222	1421	0.05%	0.011%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
 Title : SFAM01.0

37	10.037	1326	1334	1345	rVV	155400	286812	9.13%	2.148%
38	10.134	1347	1350	1352	rVV4	1974	2633	0.08%	0.020%
39	10.152	1352	1353	1357	rVV4	1540	1793	0.06%	0.013%
40	10.207	1357	1362	1367	rVV8	2173	4233	0.13%	0.032%

41	10.323	1372	1381	1387	rBV	610803	1083462	34.49%	8.115%
42	10.384	1387	1391	1402	rVB	26097	54431	1.73%	0.408%
43	10.488	1403	1408	1409	rBV4	2485	2894	0.09%	0.022%
44	10.579	1416	1423	1430	rBV	98698	173830	5.53%	1.302%
45	10.652	1433	1435	1438	rVV4	1331	1194	0.04%	0.009%

46	10.683	1438	1440	1441	rVV	1649	1484	0.05%	0.011%
47	10.701	1441	1443	1449	rVB5	4418	5725	0.18%	0.043%
48	10.786	1453	1457	1458	rVB4	1258	1427	0.05%	0.011%
49	10.860	1463	1469	1474	rBV2	31268	53973	1.72%	0.404%
50	10.927	1474	1480	1496	rVB	178132	338505	10.77%	2.535%

51	11.225	1528	1529	1534	rVB5	1456	1761	0.06%	0.013%
52	11.542	1578	1581	1587	rVB7	1458	1718	0.05%	0.013%
53	11.628	1587	1595	1605	rBV	530864	922204	29.35%	6.907%
54	11.731	1605	1612	1619	rVV	45052	83627	2.66%	0.626%
55	11.835	1623	1629	1639	rVB8	7668	20699	0.66%	0.155%

56	11.914	1639	1642	1647	rBV6	1034	2063	0.07%	0.015%
57	12.067	1665	1667	1670	rVB3	2020	1597	0.05%	0.012%
58	12.164	1678	1683	1684	rBV4	4603	5954	0.19%	0.045%
59	12.329	1707	1710	1712	rBV4	1454	2007	0.06%	0.015%
60	12.469	1725	1733	1741	rBV2	26208	51246	1.63%	0.384%

61	12.603	1746	1755	1760	rBV2	48499	80483	2.56%	0.603%
62	12.689	1762	1769	1779	rBV	235864	456329	14.52%	3.418%
63	12.804	1785	1788	1792	rVB5	2863	4316	0.14%	0.032%
64	12.859	1795	1797	1799	rVV3	1739	1407	0.04%	0.011%
65	12.932	1805	1809	1817	rVB9	8371	14486	0.46%	0.109%

66	13.024	1818	1824	1831	rVB2	57221	96070	3.06%	0.720%
67	13.164	1843	1847	1850	rBV5	5145	9025	0.29%	0.068%
68	13.243	1850	1860	1866	rVV2	51346	100384	3.20%	0.752%
69	13.365	1876	1880	1882	rVV5	1940	3042	0.10%	0.023%
70	13.493	1892	1901	1906	rBV	214009	346213	11.02%	2.593%

71	13.554	1906	1911	1921	rVB	389842	624338	19.87%	4.676%
72	13.755	1941	1944	1948	rVB5	2329	3472	0.11%	0.026%
73	13.853	1952	1960	1969	rBV	292598	491918	15.66%	3.684%
74	13.999	1980	1984	1985	rBV4	2416	2587	0.08%	0.019%
75	14.066	1989	1995	2003	rVB2	47305	80463	2.56%	0.603%

76	14.188	2012	2015	2017	rBV4	1629	2422	0.08%	0.018%
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 Title : SFAM01.0

77	14.322	2030	2037	2042	rBV9	4146	8221	0.26%	0.062%
78	14.365	2042	2044	2050	rVB6	1411	1826	0.06%	0.014%
79	14.585	2076	2080	2081	rBV4	1533	1517	0.05%	0.011%
80	14.627	2081	2087	2092	rVB6	4065	8322	0.26%	0.062%
81	14.688	2094	2097	2098	rVV3	2328	2694	0.09%	0.020%
82	14.719	2099	2102	2105	rVB3	7495	9405	0.30%	0.070%
83	14.944	2136	2139	2141	rBV3	2199	2311	0.07%	0.017%
84	15.072	2152	2160	2161	rBV5	4028	7483	0.24%	0.056%
85	15.091	2161	2163	2166	rVB4	1220	1230	0.04%	0.009%
86	15.127	2166	2169	2177	rVB8	4089	7034	0.22%	0.053%
87	15.225	2182	2185	2189	rVB6	3100	4651	0.15%	0.035%
88	15.298	2189	2197	2204	rBV	2026911	3141748	100.00%	23.532%
89	15.414	2211	2216	2222	rVB2	26123	45168	1.44%	0.338%
90	15.487	2223	2228	2232	rVV2	9524	19035	0.61%	0.143%
91	15.554	2234	2239	2242	rVV3	7146	12591	0.40%	0.094%
92	15.676	2255	2259	2265	rBV9	1397	2971	0.09%	0.022%
93	15.761	2268	2273	2276	rBV7	2297	3468	0.11%	0.026%
94	15.792	2276	2278	2282	rVV5	1841	1977	0.06%	0.015%
95	15.834	2284	2285	2290	rVB5	1374	1821	0.06%	0.014%
96	16.036	2316	2318	2323	rVB6	1783	2215	0.07%	0.017%
97	16.340	2366	2368	2372	rVB5	1708	1516	0.05%	0.011%
98	16.462	2385	2388	2390	rVV4	1591	2203	0.07%	0.017%
99	16.487	2390	2392	2396	rVB4	2511	2806	0.09%	0.021%
100	16.566	2400	2405	2412	rVV10	3587	8078	0.26%	0.061%

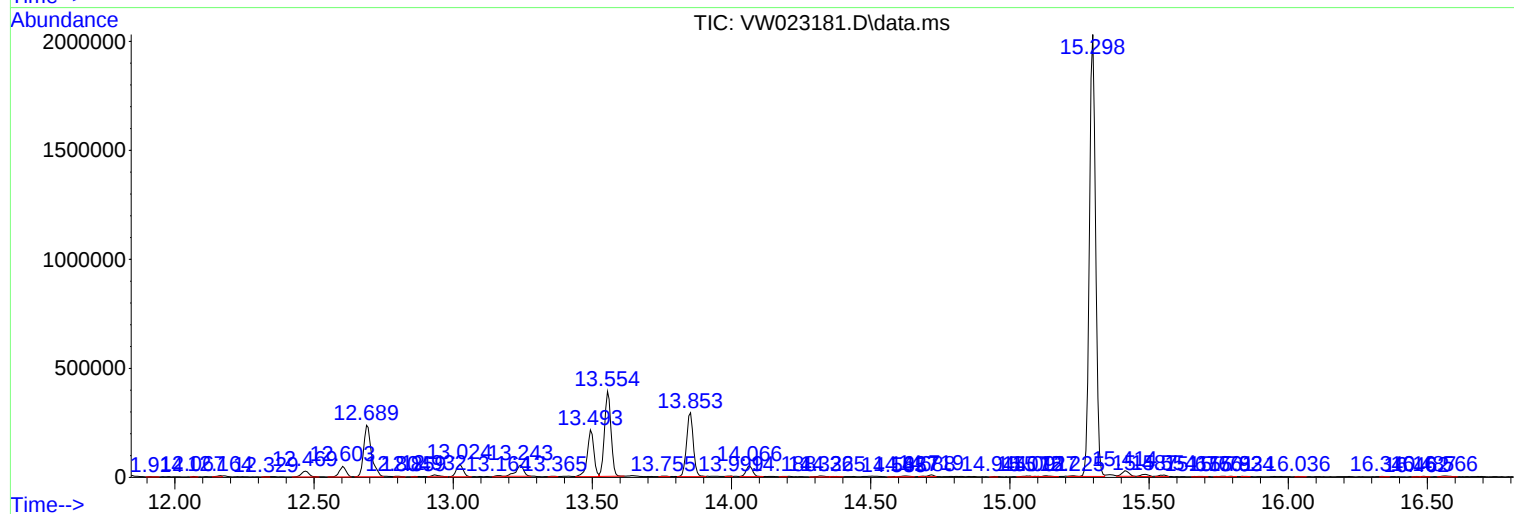
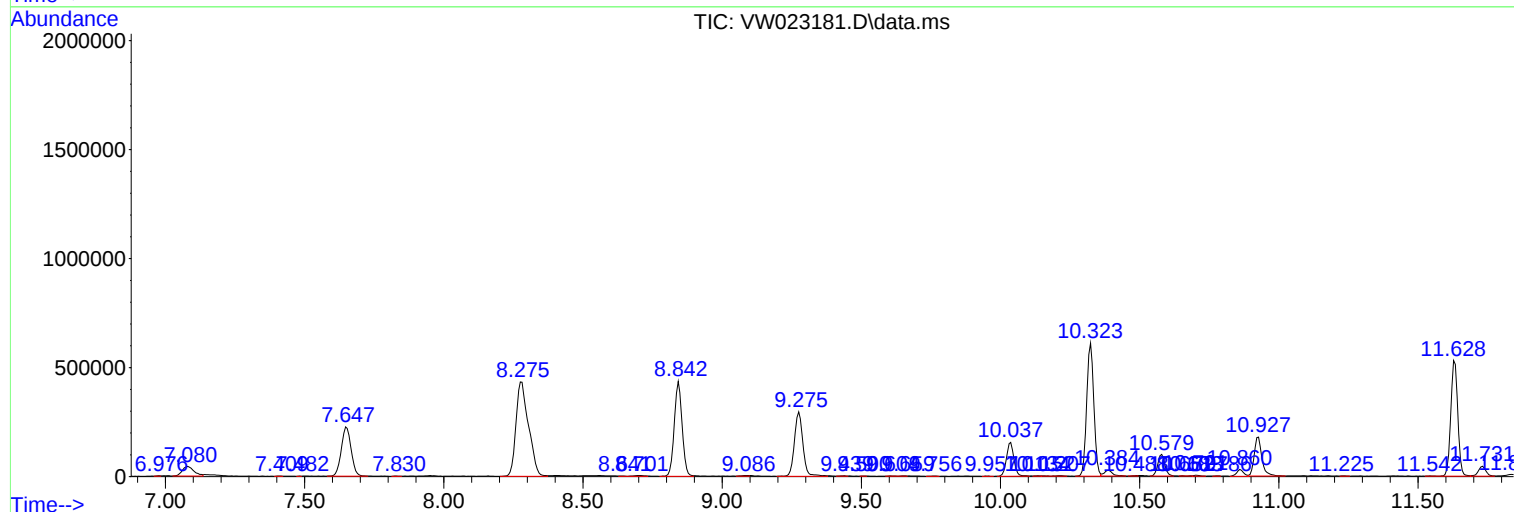
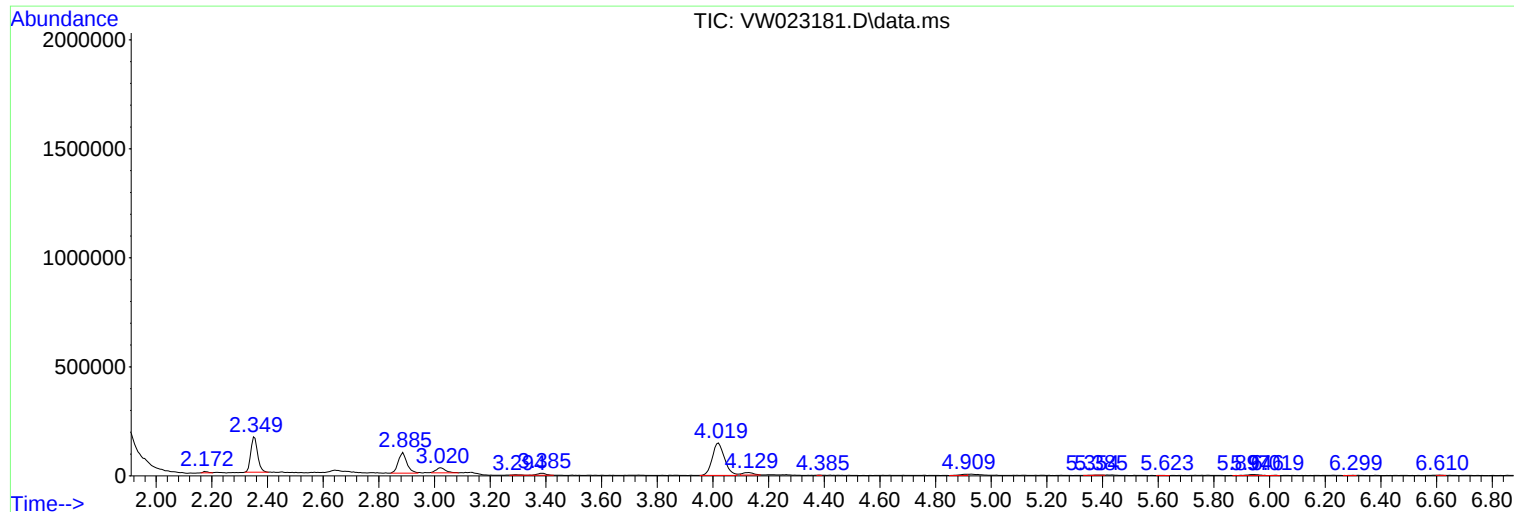
Sum of corrected areas: 13351149

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.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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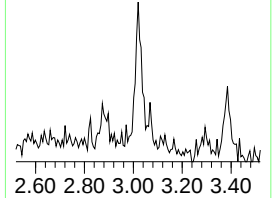
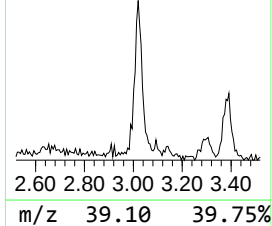
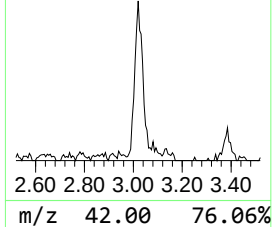
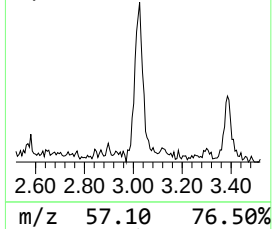
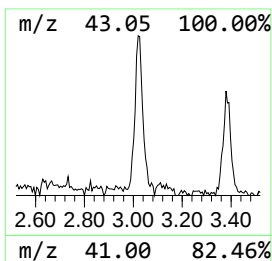
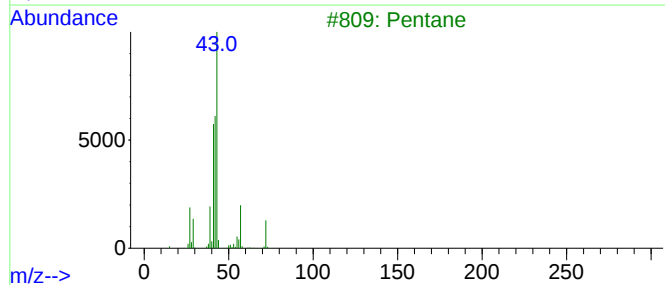
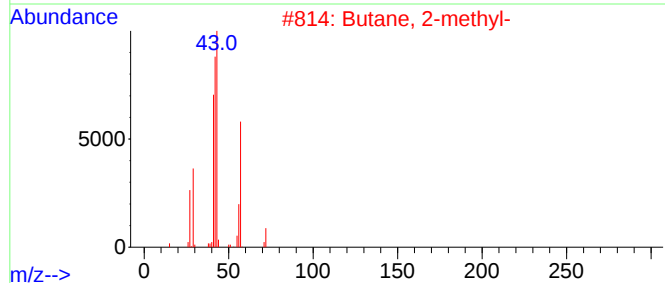
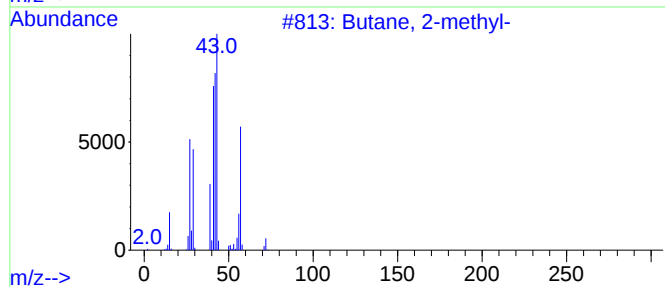
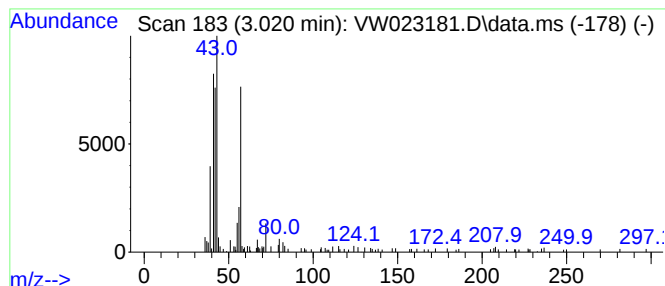
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.020	1.61 ug/L	54570	1,4-Difluorobenzene	8.842

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	53
2		Butane, 2-methyl-	72	C5H12	000078-78-4	53
3		Pentane	72	C5H12	000109-66-0	43
4		Pentane	72	C5H12	000109-66-0	43
5		3-Buten-1-ol	72	C4H8O	000627-27-0	42



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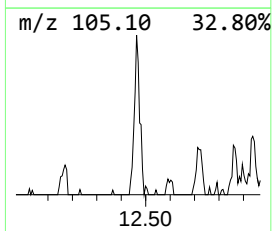
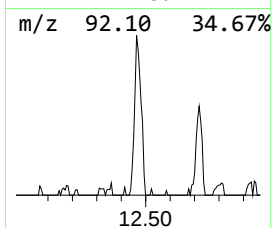
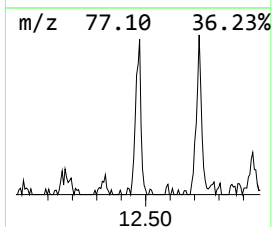
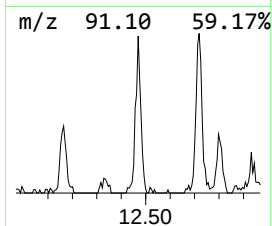
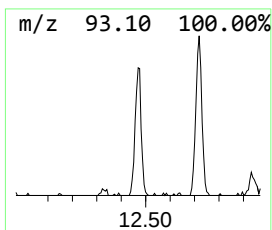
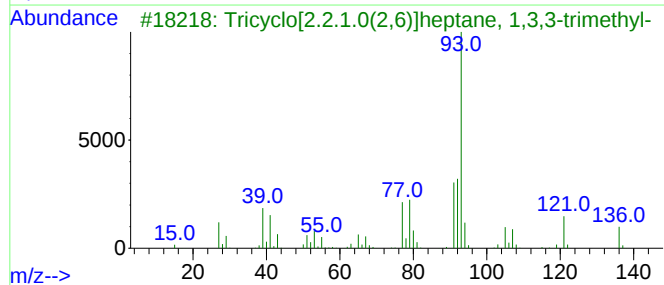
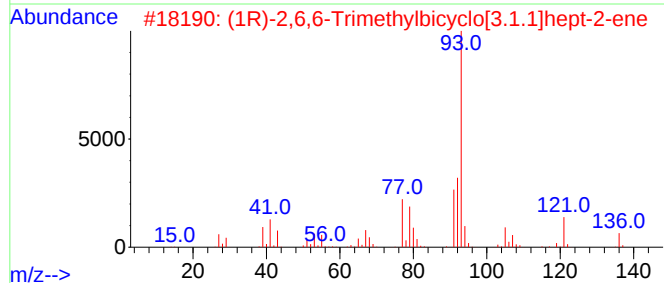
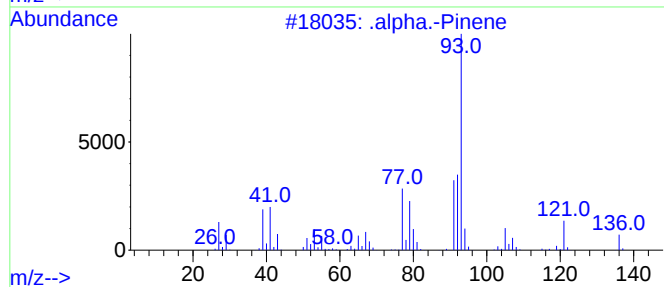
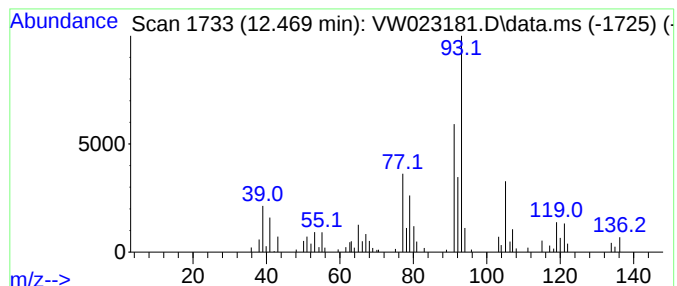
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 .alpha.-Pinene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	1.39 ug/L	51246	Chlorobenzene-d5	11.628

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		.alpha.-Pinene	136	C10H16	000080-56-8	93
2		(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	90
3		Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000488-97-1	86
4		.alpha.-Pinene	136	C10H16	000080-56-8	83
5		.alpha.-Pinene	136	C10H16	000080-56-8	83



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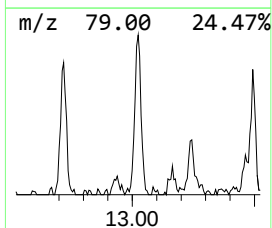
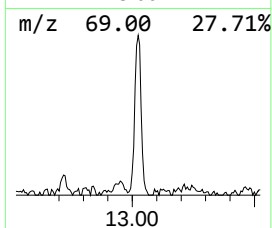
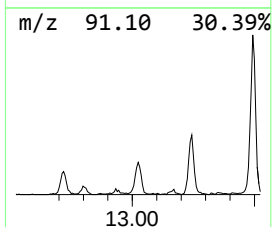
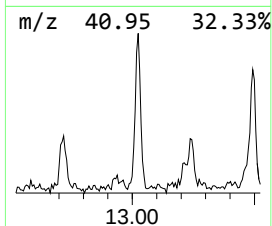
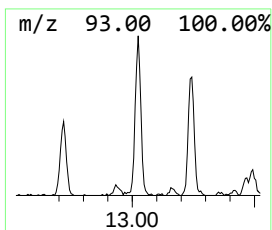
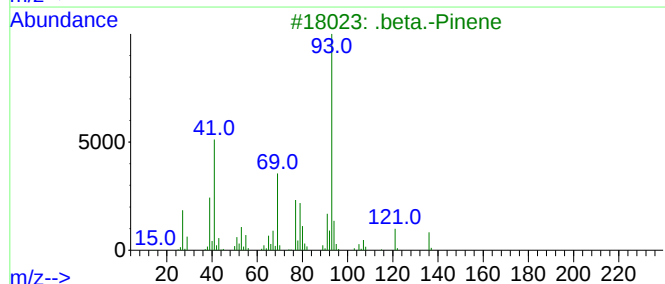
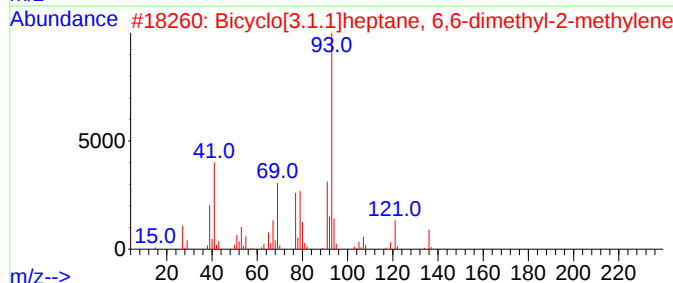
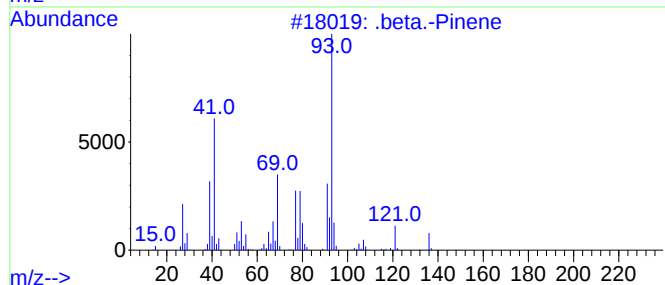
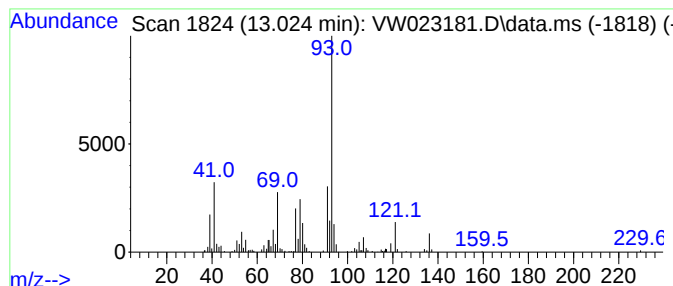
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Quant Title : SFAM01.0

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

Peak Number 5 .beta.-Pinene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.024	3.85 ug/L	96070	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-Pinene	136	C10H16	000127-91-3	97
2			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	96
3			.beta.-Pinene	136	C10H16	000127-91-3	95
4			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	94
5			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051922\
Data File : VW023181.D
Acq On : 19 May 2022 18:33
Operator : SY/VA
Sample : N2929-01
Misc : 6.89g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBTF7

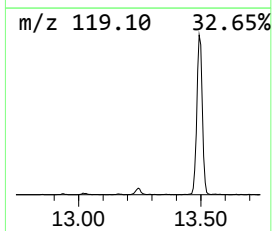
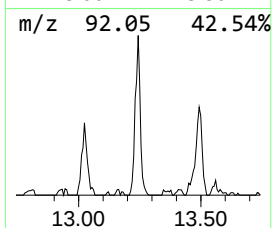
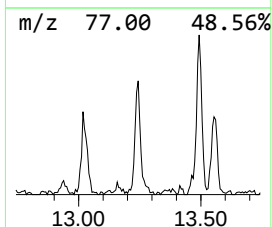
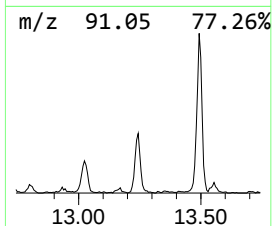
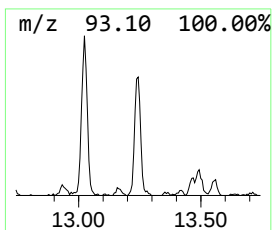
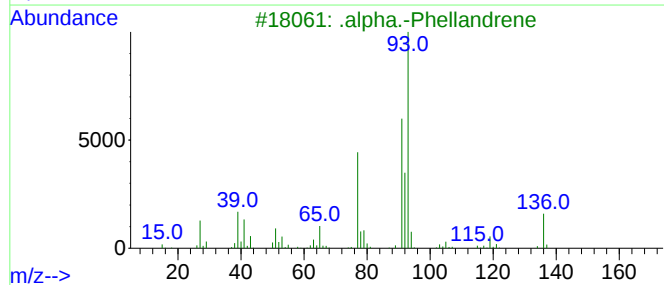
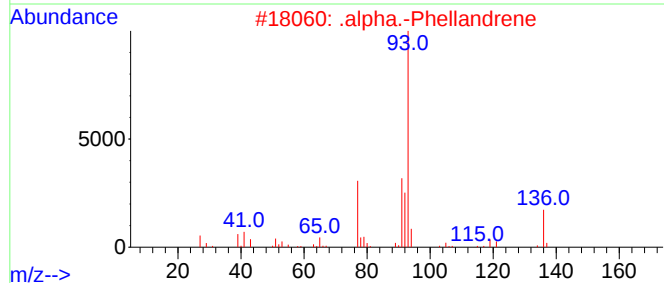
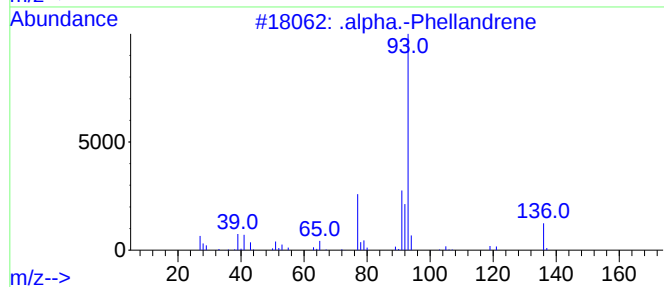
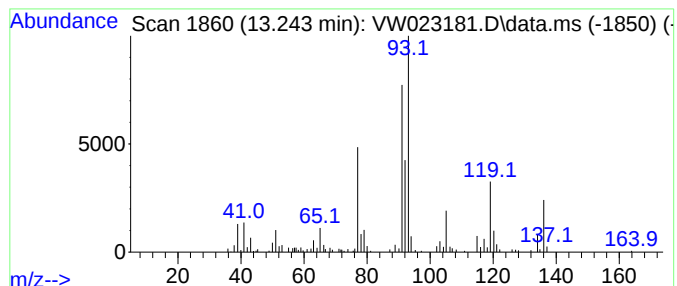
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Quant Title : SFAM01.0

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

Peak Number 6 .alpha.-Phellandrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.243	4.02 ug/L	100384	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	alpha.-Phellandrene	136	C10H16	000099-83-2	64	
2	.	alpha.-Phellandrene	136	C10H16	000099-83-2	64	
3	.	alpha.-Phellandrene	136	C10H16	000099-83-2	64	
4	Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	58		
5	.	alpha.-Phellandrene	136	C10H16	000099-83-2	45	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051922\
Data File : VW023181.D
Acq On : 19 May 2022 18:33
Operator : SY/VA
Sample : N2929-01
Misc : 6.89g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBTF7

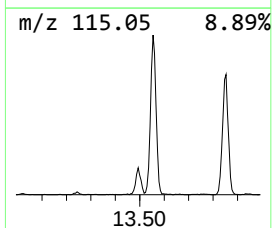
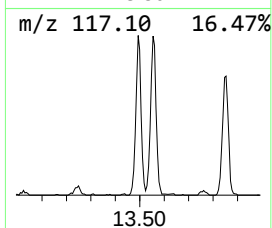
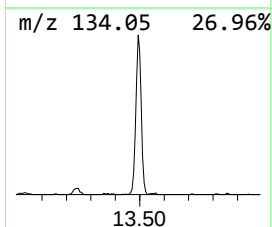
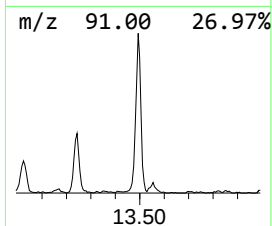
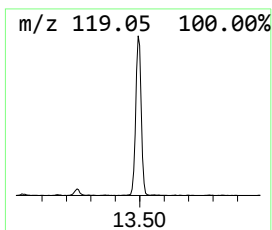
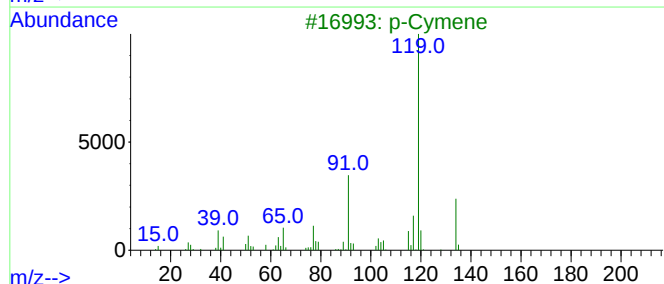
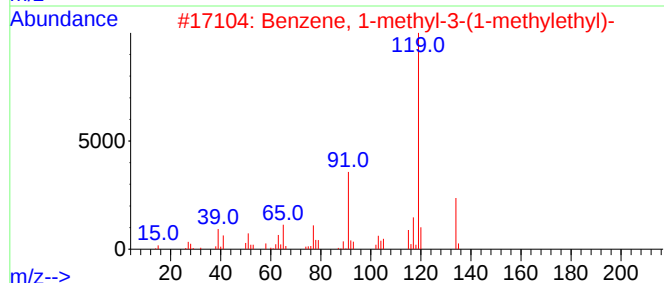
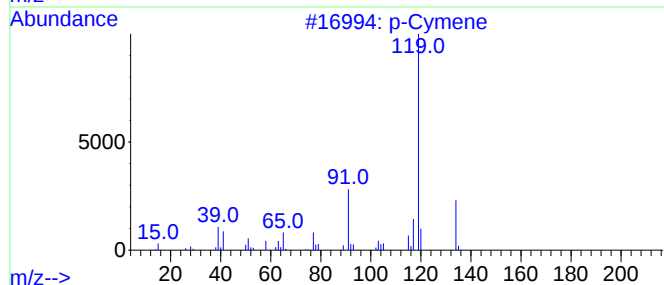
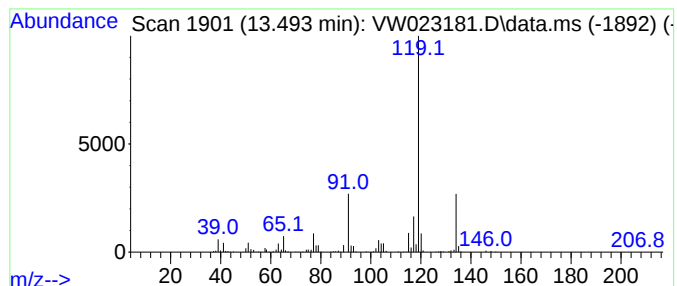
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Quant Title : SFAM01.0

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

Peak Number 7 p-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.493	13.86 ug/L	346213	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			p-Cymene	134	C10H14	000099-87-6	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051922\
Data File : VW023181.D
Acq On : 19 May 2022 18:33
Operator : SY/VA
Sample : N2929-01
Misc : 6.89g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBTF7

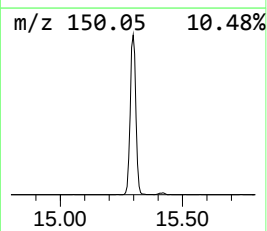
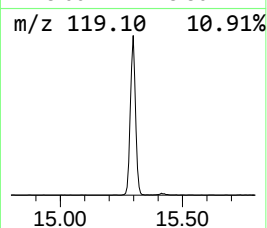
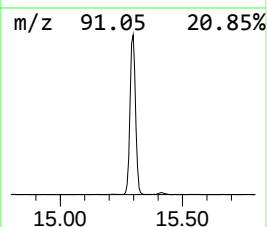
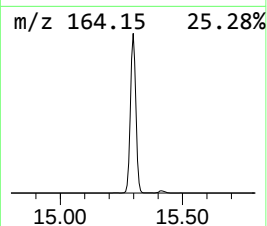
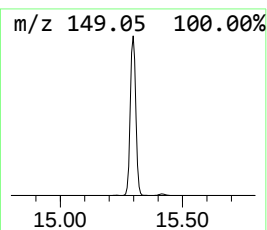
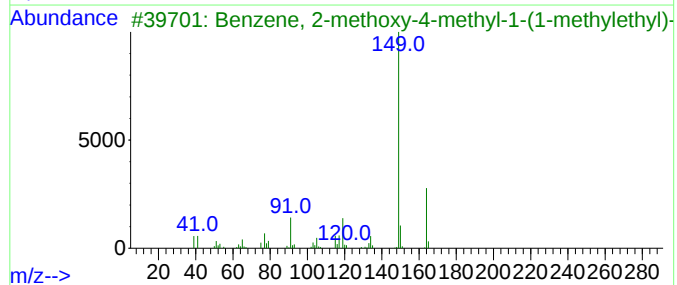
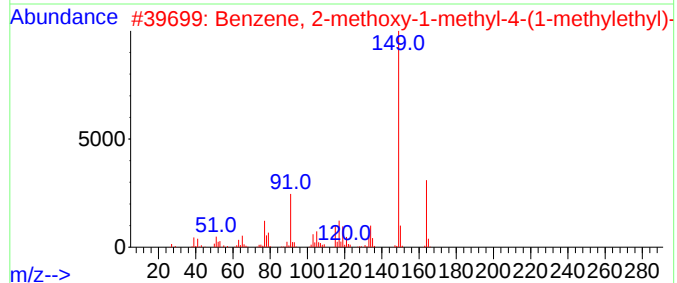
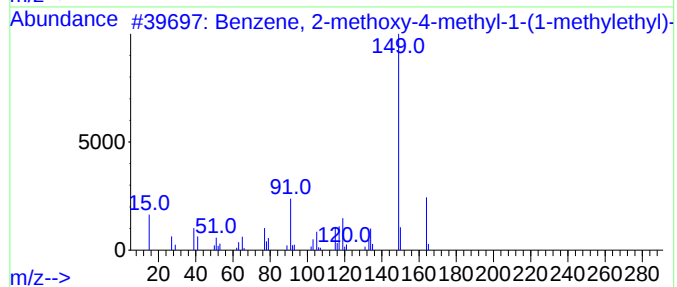
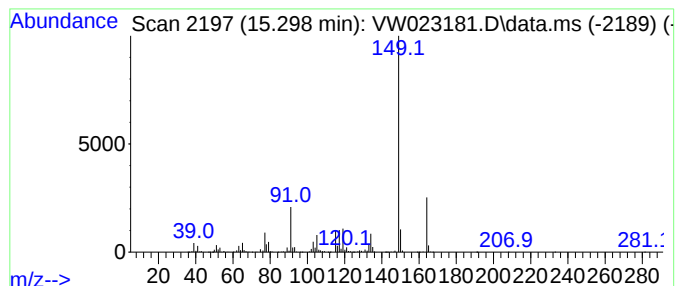
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Quant Title : SFAM01.0

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

Peak Number 9 Benzene, 2-methoxy-4-methyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.298	125.80 ug/L	3141750	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-methoxy-4-methyl-1-(1-methylethyl)-	164	C11H16O	001076-56-8	97
2			Benzene, 2-methoxy-1-methyl-4-(1-methylethyl)-	164	C11H16O	006379-73-3	91
3			Benzene, 2-methoxy-4-methyl-1-(1-methylethyl)-	164	C11H16O	001076-56-8	87
4			Benzene, 2-methoxy-4-methyl-1-(1-methylethyl)-	164	C11H16O	001076-56-8	87
5			Phenol, 2-(1,1-dimethylethyl)-4-	164	C11H16O	002409-55-4	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051922\
Data File : VW023181.D
Acq On : 19 May 2022 18:33
Operator : SY/VA
Sample : N2929-01
Misc : 6.89g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBTF7

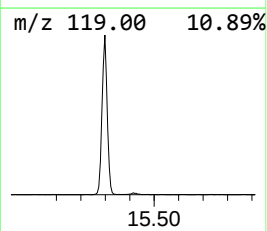
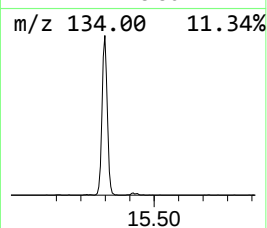
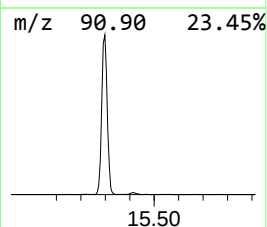
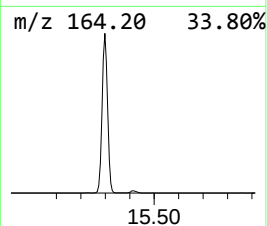
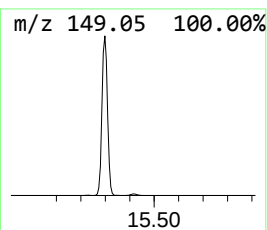
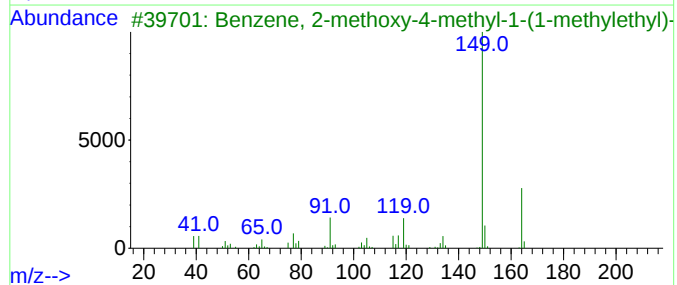
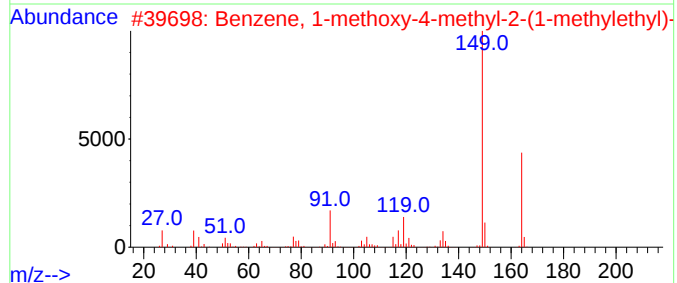
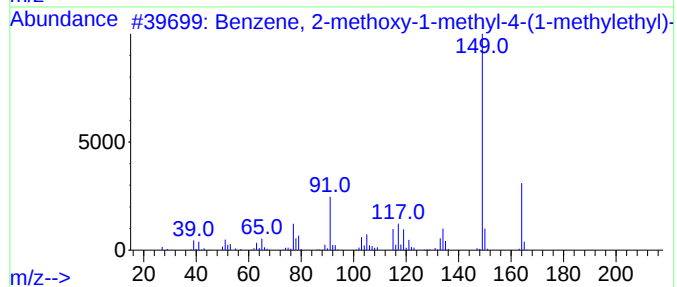
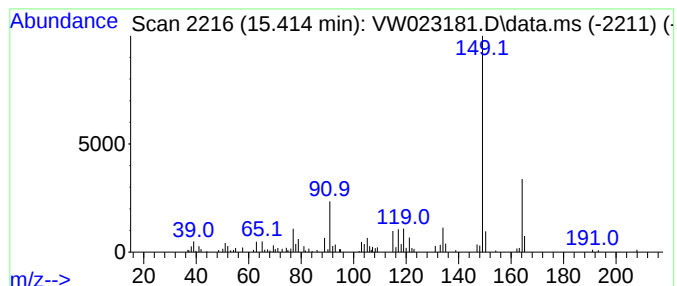
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Quant Title : SFAM01.0

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

Peak Number 10 Benzene, 2-methoxy-1-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.414	1.81 ug/L	45168	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-methoxy-1-methyl-4-(1...	164	C11H16O	006379-73-3	94
2			Benzene, 1-methoxy-4-methyl-2-(1...	164	C11H16O	031574-44-4	91
3			Benzene, 2-methoxy-4-methyl-1-(1...	164	C11H16O	001076-56-8	87
4			Benzene, 2-methoxy-4-methyl-1-(1...	164	C11H16O	001076-56-8	87
5			2'-Hydroxy-4',5'-dimethylacetoph...	164	C10H12O2	036436-65-4	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051922\
Data File : VW023181.D
Acq On : 19 May 2022 18:33
Operator : SY/VA
Sample : N2929-01
Misc : 6.89g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBTF7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.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...	3.020	1.6	ug/L	54570	1	8.842	847607	25.0
.alpha.-Pinene	12.469	1.4	ug/L	51246	2	11.628	922204	25.0
.beta.-Pinene	13.024	3.9	ug/L	96070	3	13.554	624338	25.0
.alpha.-Phellan...	13.243	4.0	ug/L	100384	3	13.554	624338	25.0
p-Cymene	13.493	13.9	ug/L	346213	3	13.554	624338	25.0
Benzene, 2-meth...	15.298	125.8	ug/L	3141750	3	13.554	624338	25.0
Benzene, 2-meth...	15.414	1.8	ug/L	45168	3	13.554	624338	25.0