LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

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\SFAMWLM101222SMA.M
Title : SFAM01.0

Signal : TIC: VW024925.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.349	67	73	82	rVB	100159	156903	3.23%	0.809%
2	2.623	116	118	120	rBV2	2151	2308	0.05%	0.012%
3	2.879	154	160	173	rVB	53995	119010	2.45%	0.614%
4	2.983	175	177	179	rBV3	1516	1599	0.03%	0.008%
5	3.373	239	241	245	rBV4	699	965	0.02%	0.005%
6	3.501	260	262	264	rVB2	699	636	0.01%	0.003%
7	3.538	264	268	271	rBV4	1027	1932	0.04%	0.010%
8	3.727	295	299	308	rVB6	952	1690	0.03%	0.009%
9	4.013	334	346	363	rBV2	151118	481919	9.92%	2.484%
10	4.397	408	409	414	rVB3	865	1025	0.02%	0.005%
11	4.611	439	444	445	rBV3	1125	1478	0.03%	0.008%
12	4.921	481	495	503	rBV2	6570	24393	0.50%	0.126%
13	5.184	536	538	541	rVB2	503	519	0.01%	0.003%
14	5.232	541	546	548	rBV3	452	755	0.02%	0.004%
15	5.440	577	580	582	rVV2	464	516	0.01%	0.003%
16	5.940	654	662	671	rBV4	2456	5854	0.12%	0.030%
17	7.080	838	849	868	rBV2	57772	195944	4.03%	1.010%
18	7.421	902	905	906	rBV2	648	624	0.01%	0.003%
19	7.647	931	942	959	rBV	288107	702792	14.46%	3.623%
20	7.817	968	970	973	rVB2	592	670	0.01%	0.003%
21	7.848	973	975	977	rBV	451	598	0.01%	0.003%
22	7.921	984	987	991	rVB2	433	621	0.01%	0.003%
23	8.012	999	1002	1006	rBV	431	568	0.01%	0.003%
24	8.275	1031	1045	1065	rBV2	536513	1625313	33.44%	8.379%
25	8.506	1082	1083	1091	rVB5	1459	2191	0.05%	0.011%
26	8.841	1129	1138	1150	rBV	615781	1200115	24.70%	6.187%
27	9.079	1172	1177	1183	rVB6	2038	3979	0.08%	0.021%
28	9.274	1200	1209	1219	rBV	389609	796535	16.39%	4.106%
29	9.390	1226	1228	1230	rVB2	705	531	0.01%	0.003%
30	9.469	1240	1241	1243	rBV2	674	559	0.01%	0.003%
31	9.597	1258	1262	1263	rBV2	737	782	0.02%	0.004%
32	9.665	1267	1273	1279	rBV6	1844	3945	0.08%	0.020%
33	9.878	1306	1308	1313	rVB2	342	521	0.01%	0.003%
34	9.933	1313	1317	1319	rBV2	424	648	0.01%	0.003%
35	10.030	1325	1333	1343	rBV	177638	329855	6.79%	1.700%
36	10.122	1346	1348	1356	rVB5	1606	2944	0.06%	0.015%

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Instrument :
MSVOA_W
ClientSampleId :
C0BL0

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\SFAMWLM101222SMA.M
Title : SFAM01.0

37	10.250	1364	1369	1373	rBV5	519	926	0.02%	0.005%
38	10.323	1373	1381	1388	rBV	708957	1268295	26.10%	6.538%
39	10.384	1388	1391	1397	rVV2	11481	19805	0.41%	0.102%
40	10.457	1400	1403	1409	rVB2	2120	2971	0.06%	0.015%
41	10.579	1416	1423	1433	rBV	105909	189442	3.90%	0.977%
42	10.701	1438	1443	1449	rVB4	4809	9027	0.19%	0.047%
43	10.786	1454	1457	1458	rVB2	742	534	0.01%	0.003%
44	10.920	1472	1479	1492	rBV	202324	365279	7.52%	1.883%
45	11.201	1522	1525	1526	rBV3	503	549	0.01%	0.003%
46	11.231	1529	1530	1533	rVV3	625	571	0.01%	0.003%
47	11.317	1540	1544	1546	rBV2	507	647	0.01%	0.003%
48	11.408	1555	1559	1563	rBV4	545	830	0.02%	0.004%
49	11.628	1587	1595	1607	rBV	832956	1363090	28.05%	7.027%
50	11.731	1607	1612	1618	rVB4	4099	7443	0.15%	0.038%
51	11.823	1620	1627	1634	rVB4	4847	11069	0.23%	0.057%
52	12.109	1671	1674	1678	rVB4	396	698	0.01%	0.004%
53	12.170	1678	1684	1692	rVB5	3871	8536	0.18%	0.044%
54	12.365	1702	1716	1725	rBV	230468	394933	8.13%	2.036%
55	12.469	1725	1733	1741	rBV	751327	1233821	25.39%	6.361%
56	12.603	1749	1755	1762	rBV2	20301	38287	0.79%	0.197%
57	12.719	1762	1774	1789	rVV2	2523494	4859684	100.00%	25.053%
58	12.859	1794	1797	1799	rBV2	1587	1558	0.03%	0.008%
59	12.951	1804	1812	1817	rVV4	33334	78306	1.61%	0.404%
60	13.024	1817	1824	1832	rVB	341518	565779	11.64%	2.917%
61	13.152	1843	1845	1849	rBV3	673	1002	0.02%	0.005%
62	13.243	1854	1860	1866	rBV2	12050	21116	0.43%	0.109%
63	13.359	1873	1879	1882	rBV4	11541	19135	0.39%	0.099%
64	13.408	1882	1887	1891	rVV5	38858	71642	1.47%	0.369%
65	13.463	1891	1896	1905	rVV2	231036	410303	8.44%	2.115%
66	13.554	1905	1911	1917	rVV	761449	1227223	25.25%	6.327%
67	13.615	1917	1921	1930	rVB3	73522	125696	2.59%	0.648%
68	13.713	1932	1937	1940	rBV2	10918	14924	0.31%	0.077%
69	13.755	1940	1944	1948	rVB2	9427	11495	0.24%	0.059%
70	13.847	1952	1959	1967	rBV	565355	958452	19.72%	4.941%
71	13.938	1973	1974	1978	rVB3	1773	1303	0.03%	0.007%
72	13.993	1978	1983	1988	rBV3	21872	38456	0.79%	0.198%
73	14.060	1990	1994	2004	rVV2	16647	34748	0.72%	0.179%
74	14.145	2005	2008	2011	rVV2	4708	6384	0.13%	0.033%
75	14.200	2014	2017	2024	rVB2	5855	9656	0.20%	0.050%
76	14.298	2030	2033	2034	rBV3	1115	1168	0.02%	0.006%

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Instrument :
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ClientSampleId :
C0BL0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Start Thrs: 0.2 Max Peaks: 100
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Title : SFAM01.0

77	14.432	2048	2055	2063	rBV	77475	139857	2.88%	0.721%
78	14.517	2066	2069	2077	rVB4	6127	10474	0.22%	0.054%
79	14.688	2092	2097	2099	rBV4	4949	8330	0.17%	0.043%
80	14.712	2099	2101	2107	rVB2	6824	9068	0.19%	0.047%
81	14.767	2108	2110	2111	rBV2	866	644	0.01%	0.003%
82	14.804	2111	2116	2120	rVV3	6696	12042	0.25%	0.062%
83	14.847	2121	2123	2129	rVB6	2363	3178	0.07%	0.016%
84	14.950	2138	2140	2143	rVB3	719	607	0.01%	0.003%
85	15.029	2146	2153	2154	rBV4	4257	8045	0.17%	0.041%
86	15.066	2154	2159	2165	rVV2	22249	42161	0.87%	0.217%
87	15.200	2178	2181	2185	rBV3	1984	2493	0.05%	0.013%
88	15.243	2185	2188	2198	rVB9	5810	12320	0.25%	0.064%
89	15.359	2201	2207	2219	rVB	25475	50849	1.05%	0.262%
90	15.474	2222	2226	2235	rVB3	3505	8038	0.17%	0.041%
91	15.548	2236	2238	2243	rVB5	1433	2093	0.04%	0.011%
92	15.584	2243	2244	2246	rBV2	517	523	0.01%	0.003%
93	15.609	2246	2248	2252	rVB5	634	643	0.01%	0.003%
94	15.651	2252	2255	2257	rBV3	1214	1630	0.03%	0.008%
95	15.987	2308	2310	2312	rBV3	626	601	0.01%	0.003%
96	16.035	2312	2318	2323	rVB9	3129	6215	0.13%	0.032%
97	16.133	2331	2334	2337	rBV4	570	719	0.01%	0.004%
98	16.438	2375	2384	2390	rBV4	7591	17326	0.36%	0.089%
99	16.596	2406	2410	2411	rBV4	1124	1459	0.03%	0.008%
100	16.633	2411	2416	2424	rVV4	6016	12627	0.26%	0.065%

Sum of corrected areas: 19397962

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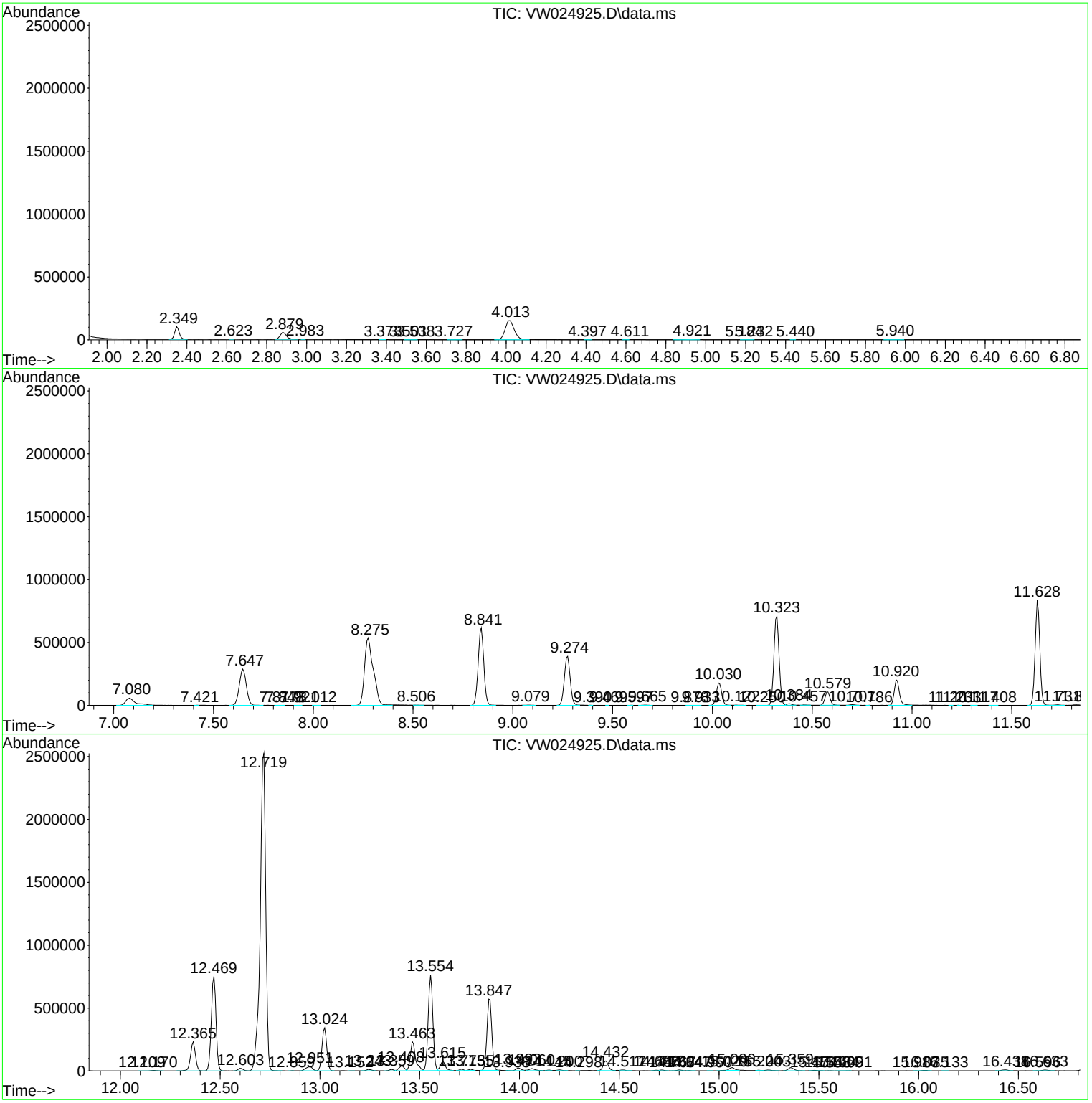
Instrument :
MSVOA_W
ClientSampleId :
C0BL0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
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Acq On : 14 Oct 2022 12:40
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Instrument :
MSVOA_W
ClientSampleId :
C0BL0

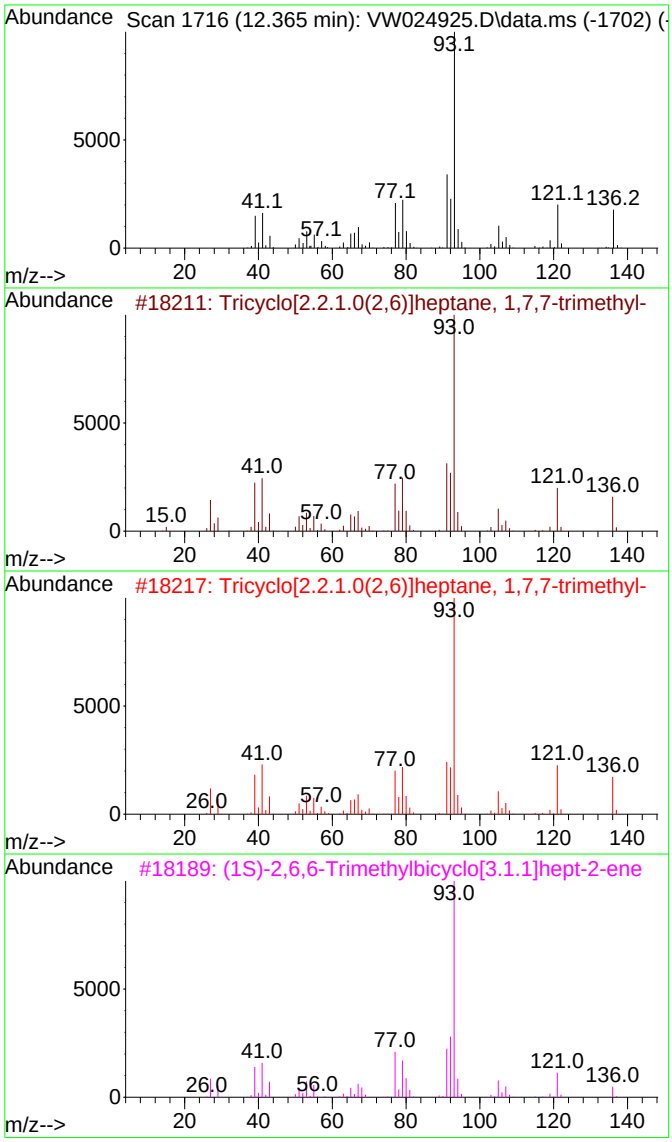
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

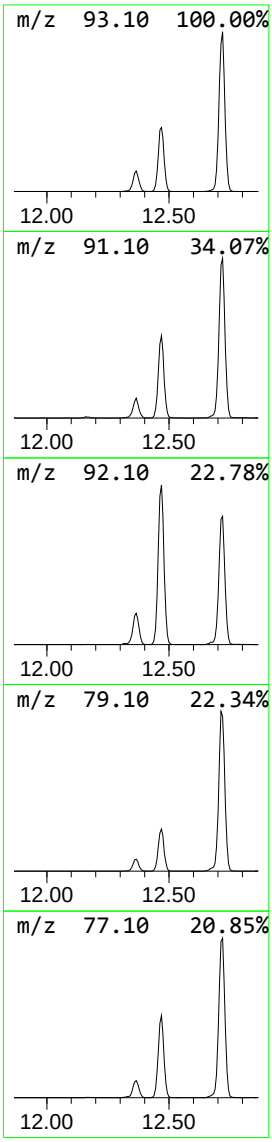
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Tricyclo[2.2.1.0(2,6)]hepta... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.365	7.24 ug/L	394933	Chlorobenzene-d5	11.628

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	96
2	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	96
3	(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	95
4	3-Carene	136	C10H16	013466-78-9	94
5	(+)-3-Carene	136	C10H16	000498-15-7	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

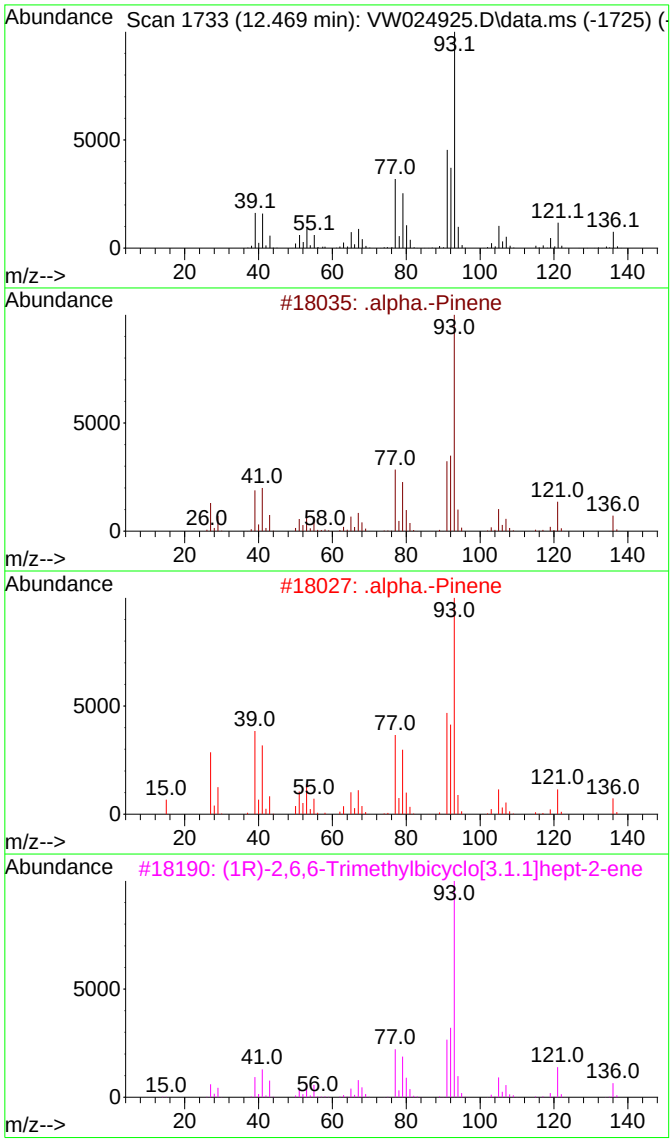
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

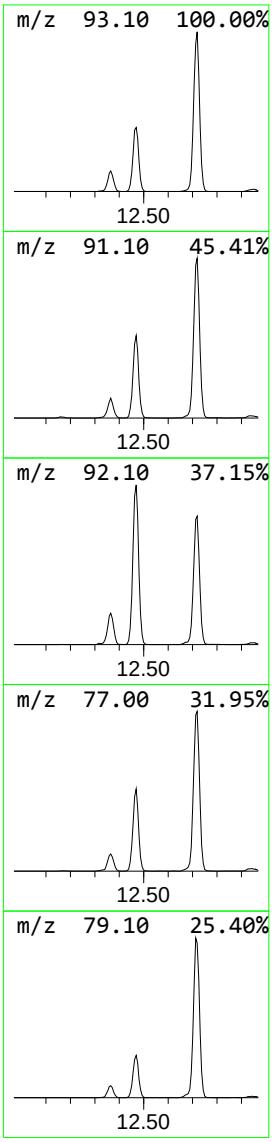
****

Peak Number 3 .alpha.-Pinene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	22.63 ug/L	1233820	Chlorobenzene-d5	11.628

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	97
2	.alpha.-Pinene	136	C10H16	000080-56-8	95
3	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	94
4	(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	94
5	.alpha.-Pinene	136	C10H16	000080-56-8	91





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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

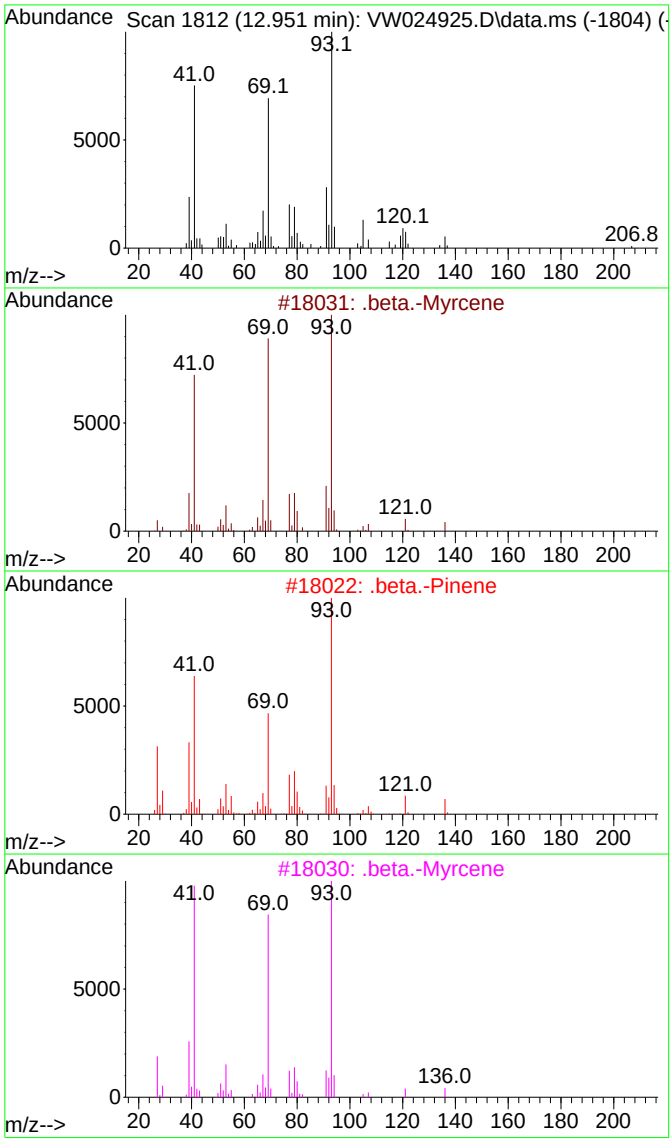
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

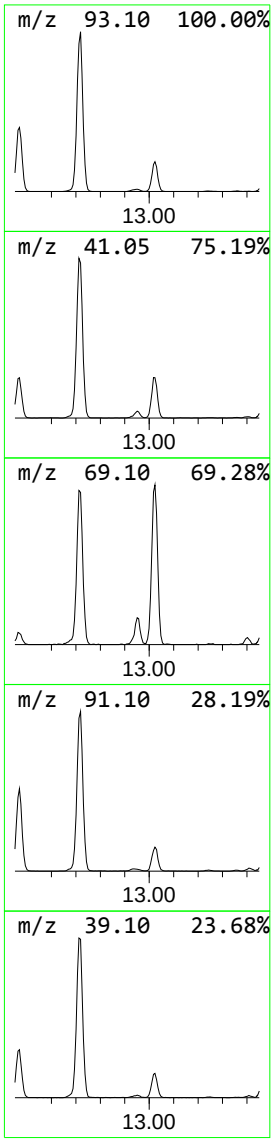
****

Peak Number 4 .beta.-Myrcene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.951	1.60 ug/L	78306	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Myrcene	136	C10H16	000123-35-3	93
2	.beta.-Pinene	136	C10H16	000127-91-3	83
3	.beta.-Myrcene	136	C10H16	000123-35-3	74
4	Bicyclo[3.1.0]hex-2-ene, 4-methy...	136	C10H16	028634-89-1	72
5	.beta.-Pinene	136	C10H16	000127-91-3	64





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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

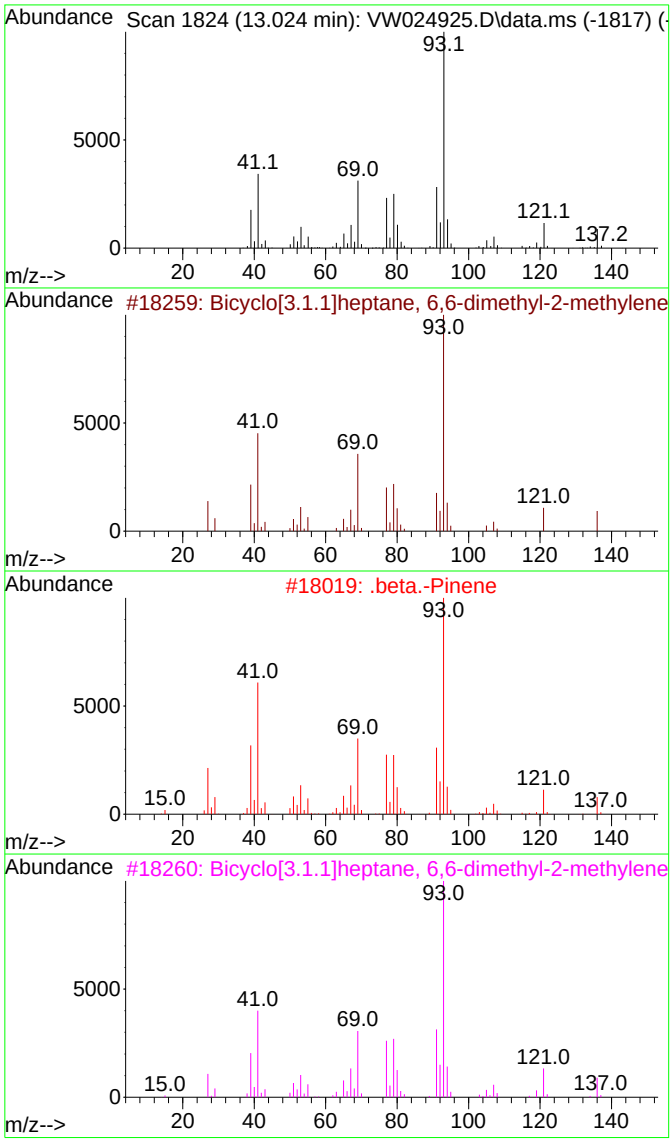
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

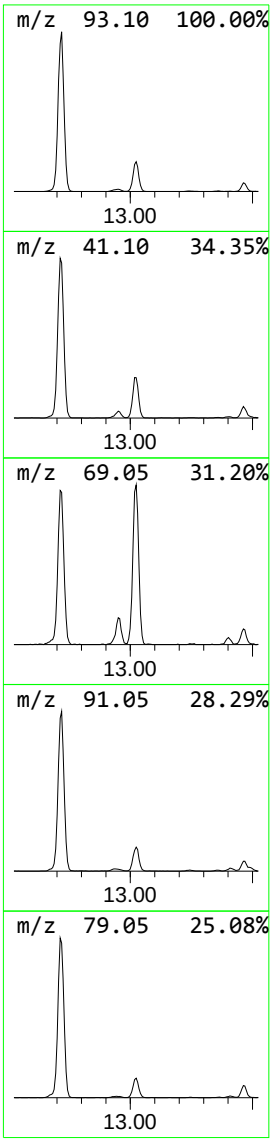
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Bicyclo[3.1.1]heptane, 6,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.024	11.53 ug/L	565779	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	97
2	.beta.-Pinene	136	C10H16	000127-91-3	96
3	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	96
4	.beta.-Pinene	136	C10H16	000127-91-3	94
5	Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

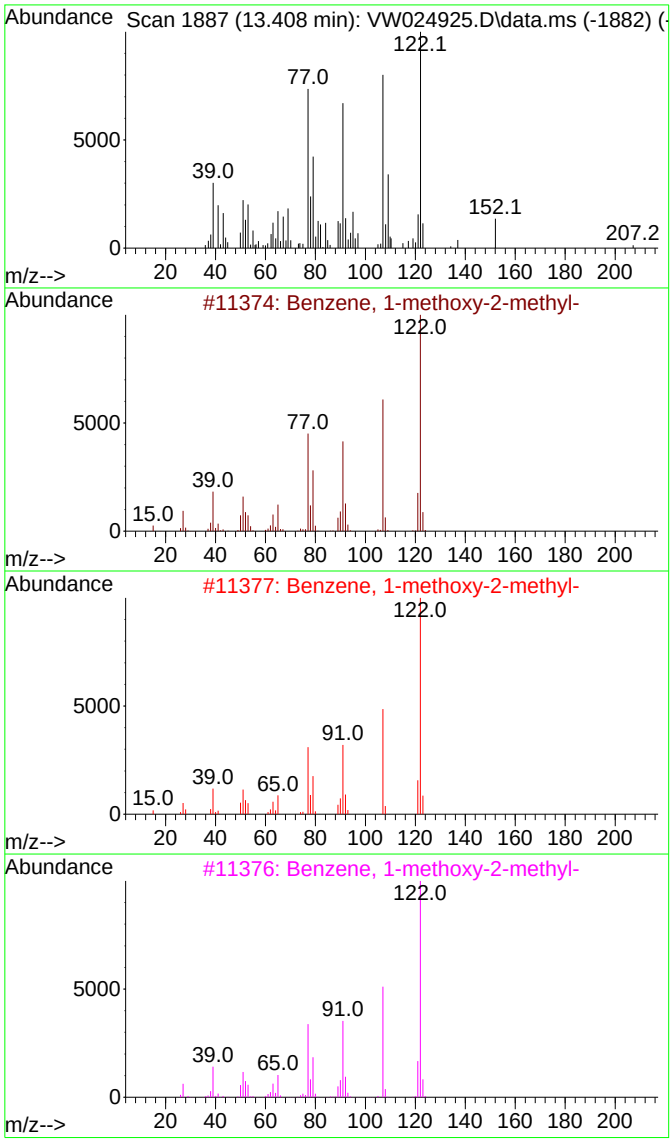
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

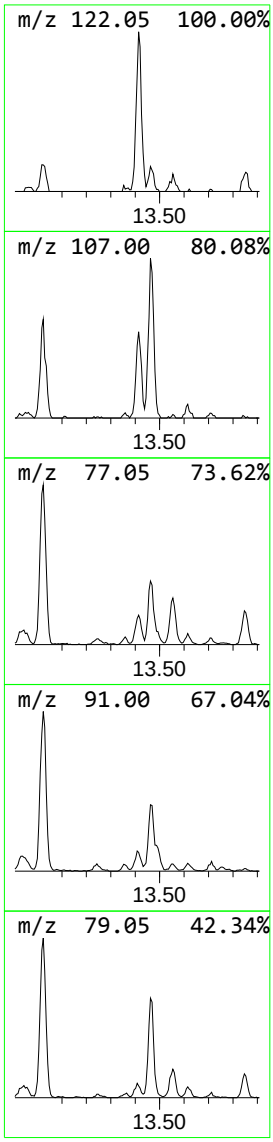
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.408	1.46 ug/L	71642	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methoxy-2-methyl-	122	C8H10O	000578-58-5	95
2	Benzene, 1-methoxy-2-methyl-	122	C8H10O	000578-58-5	94
3	Benzene, 1-methoxy-2-methyl-	122	C8H10O	000578-58-5	94
4	Benzene, 1-methoxy-2-methyl-	122	C8H10O	000578-58-5	93
5	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

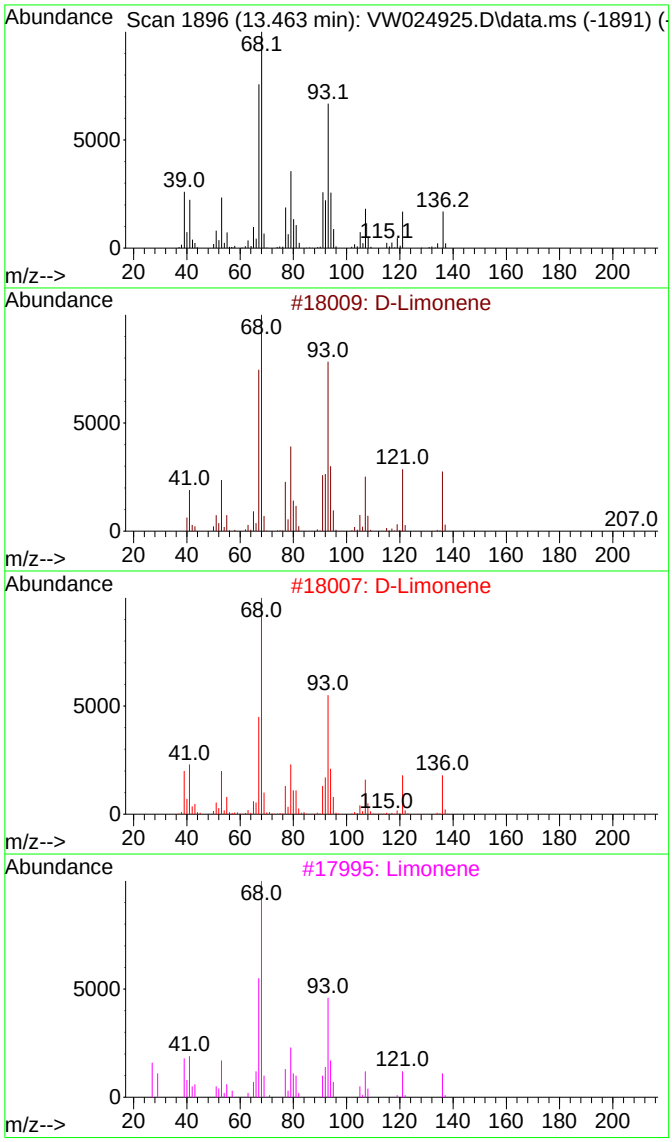
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

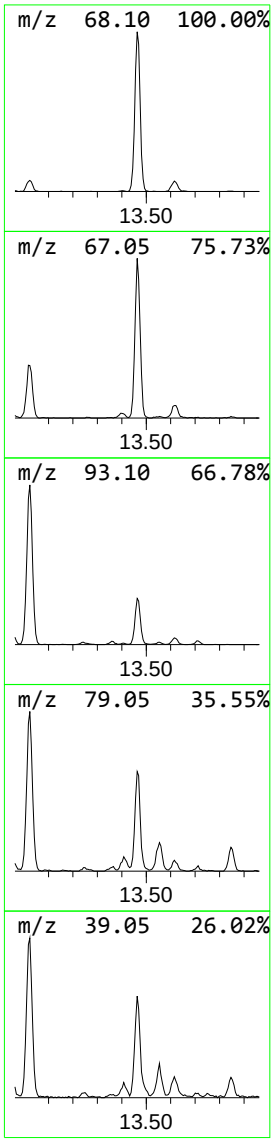
****

Peak Number 7 D-Limonene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.463	8.36 ug/L	410303	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Limonene	136	C10H16	005989-27-5	98
2	D-Limonene	136	C10H16	005989-27-5	95
3	Limonene	136	C10H16	000138-86-3	91
4	D-Limonene	136	C10H16	005989-27-5	87
5	Limonene	136	C10H16	000138-86-3	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

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

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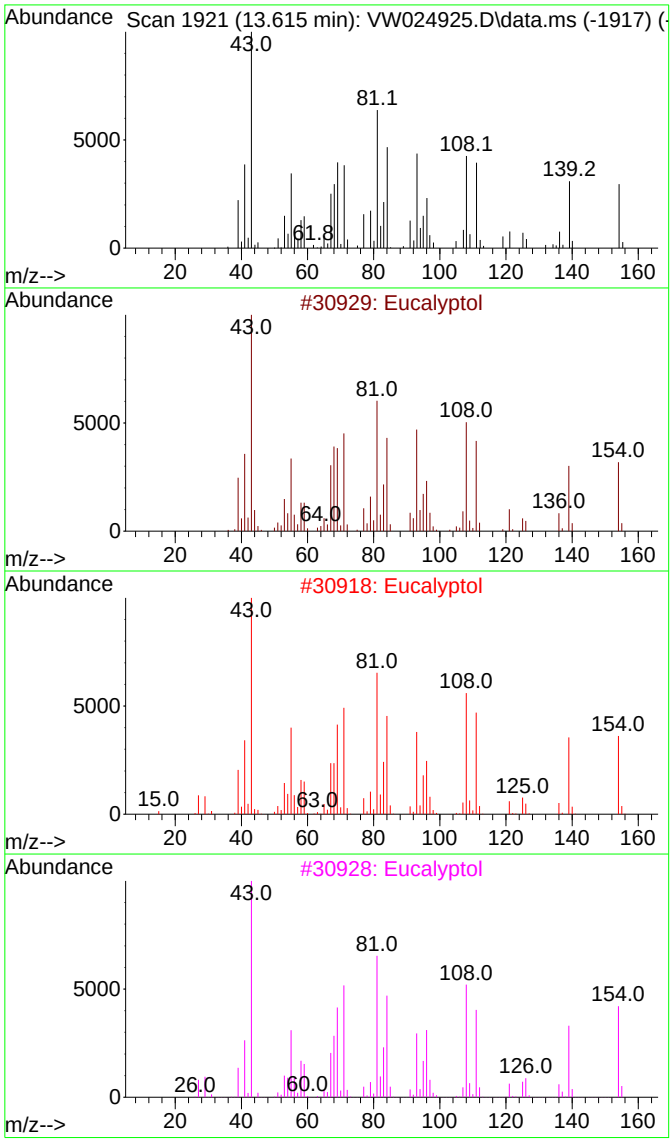
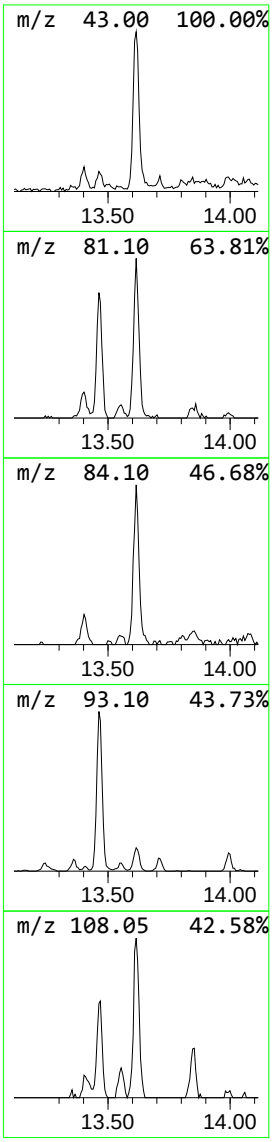
Peak Number 8 Eucalyptol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.615	2.56 ug/L	125696	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	99
2	Eucalyptol	154	C10H18O	000470-82-6	98
3	Eucalyptol	154	C10H18O	000470-82-6	94
4	Eucalyptol	154	C10H18O	000470-82-6	90
5	Eucalyptol	154	C10H18O	000470-82-6	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

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

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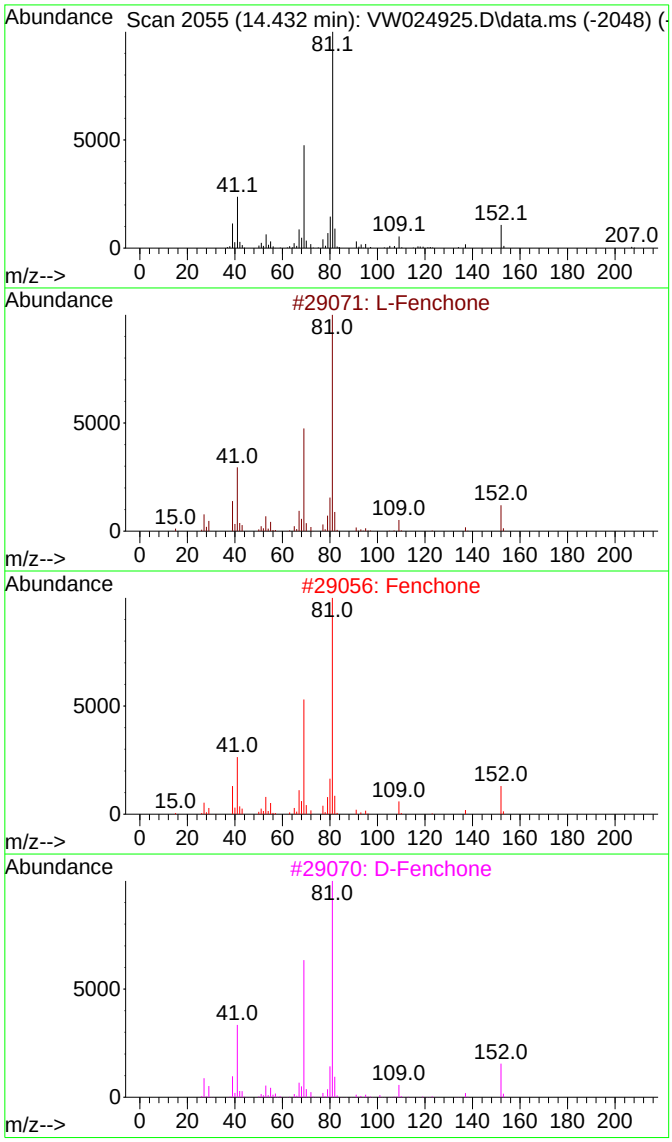
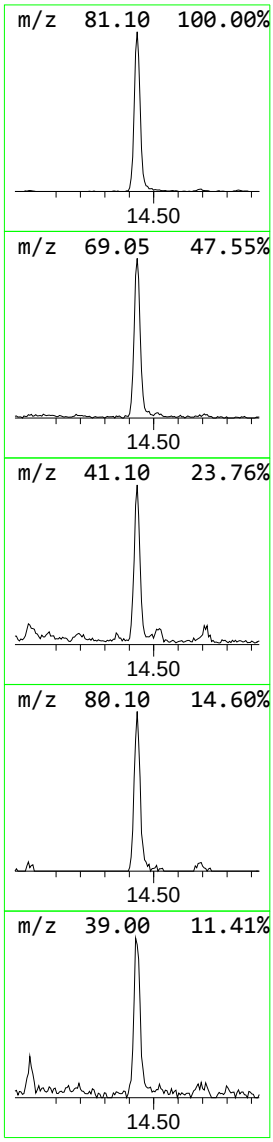
Peak Number 9 L-Fenchone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.432	2.85 ug/L	139857	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	L-Fenchone	152	C10H16O	007787-20-4	95
2	Fenchone	152	C10H16O	001195-79-5	95
3	D-Fenchone	152	C10H16O	004695-62-9	91
4	Fenchone	152	C10H16O	001195-79-5	91
5	L-Fenchone	152	C10H16O	007787-20-4	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101422\
Data File : VW024925.D
Acq On : 14 Oct 2022 12:40
Operator : SY/VA
Sample : N5144-04
Misc : 3.86g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0BL0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM101222SMA.M
Quant Title : SFAM01.0

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Tricyclo[2.2.1....	12.365	7.2	ug/L	394933	2	11.628	1363090	25.0
.alpha.-Pinene	12.469	22.6	ug/L	1233820	2	11.628	1363090	25.0
.beta.-Myrcene	12.951	1.6	ug/L	78306	3	13.554	1227220	25.0
Bicyclo[3.1.1]h...	13.024	11.5	ug/L	565779	3	13.554	1227220	25.0
Benzene, 1-meth...	13.408	1.5	ug/L	71642	3	13.554	1227220	25.0
D-Limonene	13.463	8.4	ug/L	410303	3	13.554	1227220	25.0
Eucalyptol	13.615	2.6	ug/L	125696	3	13.554	1227220	25.0
L-Fenchone	14.432	2.9	ug/L	139857	3	13.554	1227220	25.0