

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041322\
 Data File : VW022536.D
 Acq On : 13 Apr 2022 17:00
 Operator : SY/VA
 Sample : N2374-07RE
 Misc : 5.27g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BFFE3RE

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: VW022536.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.117	33	35	38	rBV4	1317	2015	0.01%	0.007%
2	2.233	49	54	56	rBV5	3896	7702	0.05%	0.026%
3	2.355	68	74	82	rVB	121010	203720	1.32%	0.695%
4	2.891	155	162	169	rBV2	83676	187741	1.21%	0.640%
5	3.257	219	222	225	rVB4	859	1126	0.01%	0.004%
6	3.306	228	230	233	rVB3	1424	1338	0.01%	0.005%
7	3.471	254	257	259	rBV3	1023	1108	0.01%	0.004%
8	3.538	261	268	277	rVV	18953	52230	0.34%	0.178%
9	4.013	333	346	362	rBV	124162	392897	2.54%	1.340%
10	4.574	435	438	440	rBV3	947	1124	0.01%	0.004%
11	4.916	483	494	514	rVB3	16409	62655	0.40%	0.214%
12	5.050	514	516	521	rBV6	881	1223	0.01%	0.004%
13	5.178	534	537	540	rVB4	815	1245	0.01%	0.004%
14	5.257	548	550	554	rBV5	1455	1947	0.01%	0.007%
15	5.306	554	558	561	rBV5	755	1517	0.01%	0.005%
16	5.617	606	609	614	rBV6	792	1223	0.01%	0.004%
17	5.745	628	630	632	rBV2	852	1041	0.01%	0.004%
18	5.769	632	634	636	rVB3	1128	1136	0.01%	0.004%
19	5.934	651	661	668	rBV2	7602	24267	0.16%	0.083%
20	6.805	801	804	806	rBV3	996	1501	0.01%	0.005%
21	7.080	841	849	871	rVB	35690	117389	0.76%	0.400%
22	7.336	889	891	895	rVB4	925	1204	0.01%	0.004%
23	7.488	914	916	921	rVB4	1575	1882	0.01%	0.006%
24	7.543	921	925	927	rBV4	860	1047	0.01%	0.004%
25	7.647	932	942	957	rBV	232138	576269	3.72%	1.966%
26	7.933	986	989	991	rBV4	1609	1421	0.01%	0.005%
27	7.958	991	993	996	rVB4	1533	1650	0.01%	0.006%
28	8.128	1014	1021	1025	rBV9	1320	3542	0.02%	0.012%
29	8.208	1030	1034	1036	rBV4	1704	2834	0.02%	0.010%
30	8.275	1036	1045	1063	rVV2	415525	1273923	8.23%	4.345%
31	8.506	1081	1083	1086	rVB4	1664	1810	0.01%	0.006%
32	8.762	1120	1125	1129	rBV7	2281	3384	0.02%	0.012%
33	8.842	1129	1138	1147	rBV	393744	782845	5.06%	2.670%
34	9.049	1167	1172	1174	rBV5	2032	3635	0.02%	0.012%
35	9.183	1192	1194	1197	rVB4	1134	1055	0.01%	0.004%
36	9.274	1200	1209	1226	rVV	305763	639059	4.13%	2.180%

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

37	9.470	1239	1241	1243	rBV3	2074	1790	0.01%	0.006%
38	9.713	1279	1281	1283	rBV3	1214	1247	0.01%	0.004%
39	10.037	1326	1334	1348	rVB	138002	254943	1.65%	0.870%
40	10.213	1360	1363	1365	rBV3	1869	2365	0.02%	0.008%
41	10.323	1374	1381	1388	rBV	558492	952055	6.15%	3.247%
42	10.390	1388	1392	1397	rVB2	8399	14617	0.09%	0.050%
43	10.579	1415	1423	1433	rVB	85739	154009	0.99%	0.525%
44	10.701	1437	1443	1450	rBV2	15606	31347	0.20%	0.107%
45	10.847	1464	1467	1469	rBV4	1209	1431	0.01%	0.005%
46	10.921	1473	1479	1493	rBV	135891	259648	1.68%	0.886%
47	11.164	1517	1519	1522	rVB3	970	1224	0.01%	0.004%
48	11.207	1522	1526	1527	rBV4	1151	1280	0.01%	0.004%
49	11.396	1555	1557	1560	rVB4	1477	1490	0.01%	0.005%
50	11.439	1560	1564	1568	rBV6	1423	2341	0.02%	0.008%
51	11.567	1583	1585	1587	rBV2	1482	1206	0.01%	0.004%
52	11.628	1587	1595	1604	rVV	501455	844834	5.46%	2.882%
53	11.707	1604	1608	1617	rVB8	4617	10062	0.06%	0.034%
54	11.805	1621	1624	1626	rVV4	1127	1296	0.01%	0.004%
55	11.841	1626	1630	1633	rVV6	2647	4157	0.03%	0.014%
56	12.073	1663	1668	1672	rBV7	1854	3355	0.02%	0.011%
57	12.128	1672	1677	1680	rVB5	1198	2250	0.01%	0.008%
58	12.170	1680	1684	1686	rBV5	1304	1999	0.01%	0.007%
59	12.329	1706	1710	1713	rBV5	8286	14238	0.09%	0.049%
60	12.365	1713	1716	1722	rVB5	9447	14640	0.09%	0.050%
61	12.469	1725	1733	1745	rBV	9645995	15480591	100.00%	52.804%
62	12.603	1749	1755	1762	rVB2	29985	55122	0.36%	0.188%
63	12.689	1762	1769	1772	rBV	265837	511844	3.31%	1.746%
64	12.823	1787	1791	1793	rVB5	1101	1517	0.01%	0.005%
65	12.957	1803	1813	1817	rBV	210417	386348	2.50%	1.318%
66	13.024	1817	1824	1836	rVB	1882086	3089492	19.96%	10.538%
67	13.195	1849	1852	1855	rVB4	3131	4080	0.03%	0.014%
68	13.243	1855	1860	1866	rVV3	13479	24003	0.16%	0.082%
69	13.298	1866	1869	1873	rVB4	2626	3287	0.02%	0.011%
70	13.359	1873	1879	1884	rBV2	36442	59326	0.38%	0.202%
71	13.402	1884	1886	1889	rVV3	5810	6981	0.05%	0.024%
72	13.463	1890	1896	1905	rVV	677320	1074401	6.94%	3.665%
73	13.554	1905	1911	1925	rVV	497077	810070	5.23%	2.763%
74	13.713	1929	1937	1943	rBV3	22459	38840	0.25%	0.132%
75	13.755	1943	1944	1946	rVV	2519	1993	0.01%	0.007%
76	13.853	1949	1960	1967	rBV	342321	589477	3.81%	2.011%

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

77	13.902	1967	1968	1972	rVV4	3120	3211	0.02%	0.011%
78	13.950	1973	1976	1978	rVV3	1736	2056	0.01%	0.007%
79	13.993	1978	1983	1988	rVV	62218	99141	0.64%	0.338%
80	14.060	1988	1994	2005	rVV2	34118	64284	0.42%	0.219%
81	14.146	2005	2008	2015	rVB8	2988	6207	0.04%	0.021%
82	14.298	2030	2033	2035	rBV3	1128	1374	0.01%	0.005%
83	14.414	2046	2052	2054	rBV5	1144	1730	0.01%	0.006%
84	14.585	2077	2080	2085	rVB4	1558	2086	0.01%	0.007%
85	14.627	2085	2087	2089	rBV3	1273	1263	0.01%	0.004%
86	14.694	2095	2098	2101	rBV5	1522	1722	0.01%	0.006%
87	14.725	2101	2103	2106	rVV4	1625	1656	0.01%	0.006%
88	14.950	2137	2140	2141	rBV3	1254	1043	0.01%	0.004%
89	15.023	2149	2152	2155	rBV3	1031	1573	0.01%	0.005%
90	15.243	2185	2188	2194	rVV7	3166	5730	0.04%	0.020%
91	15.365	2204	2208	2210	rBV5	1349	2029	0.01%	0.007%
92	15.401	2212	2214	2216	rVB3	1621	1283	0.01%	0.004%
93	15.481	2222	2227	2232	rVV	9757	15955	0.10%	0.054%
94	15.786	2271	2277	2285	rBV8	2953	7802	0.05%	0.027%
95	15.889	2291	2294	2298	rVB6	2737	2899	0.02%	0.010%
96	16.017	2312	2315	2318	rBV5	1530	2075	0.01%	0.007%
97	16.042	2318	2319	2323	rVB4	1566	1668	0.01%	0.006%
98	16.413	2376	2380	2388	rVB4	4120	8104	0.05%	0.028%
99	16.517	2394	2397	2399	rBV4	1053	1552	0.01%	0.005%
100	16.761	2432	2437	2439	rBV6	2259	3481	0.02%	0.012%

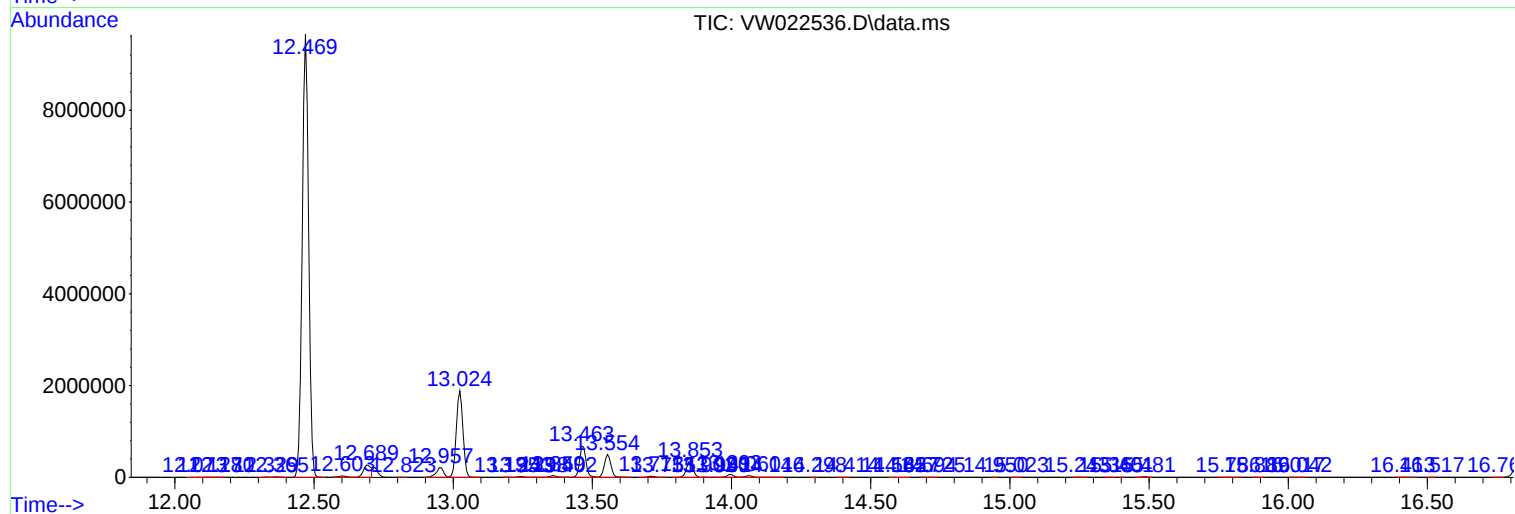
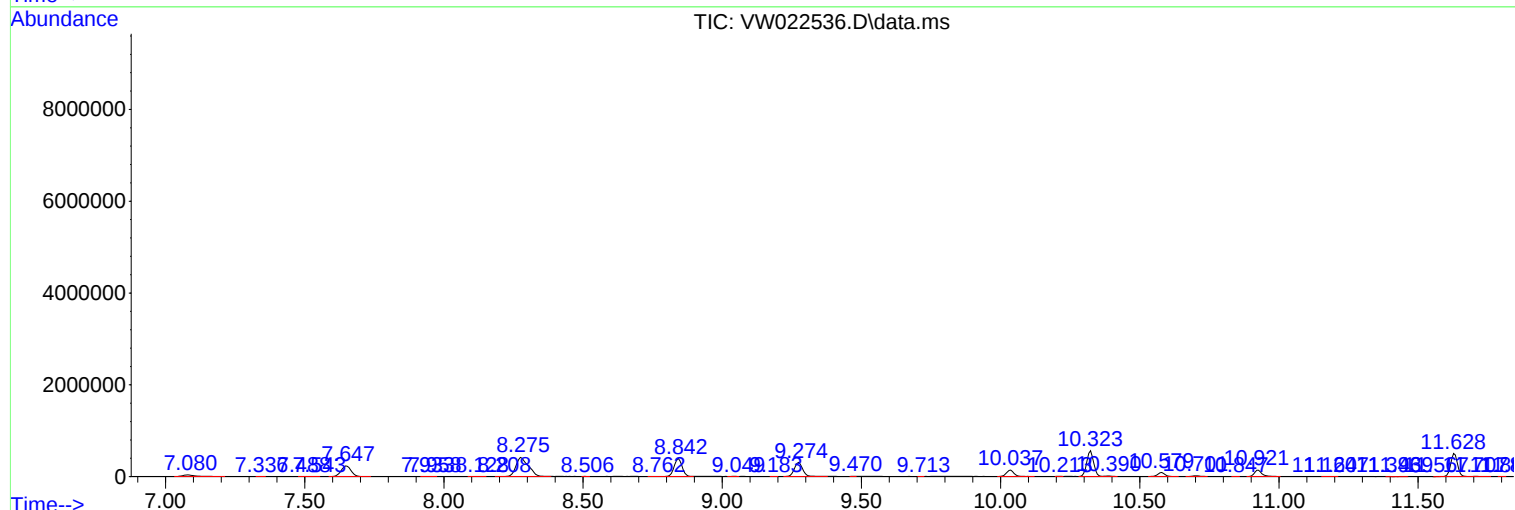
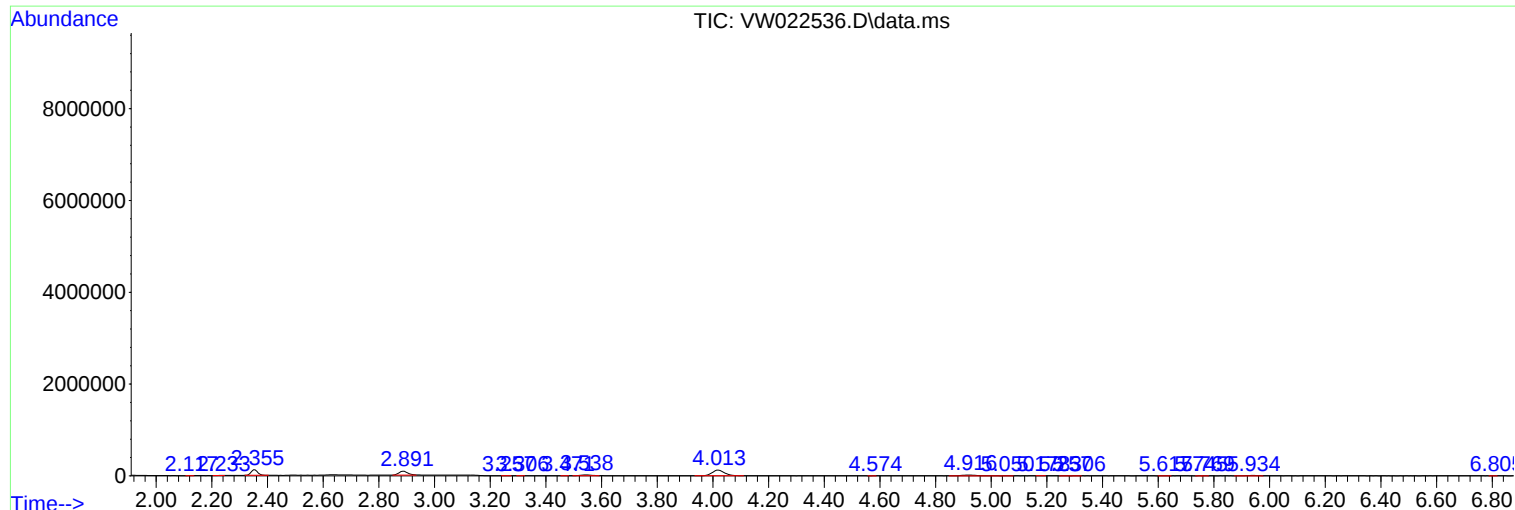
Sum of corrected areas: 29316825

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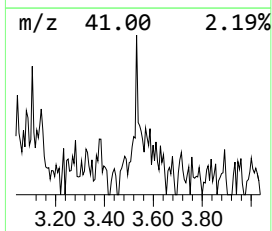
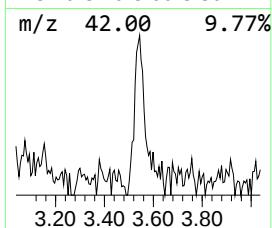
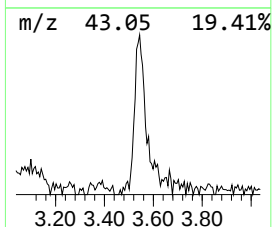
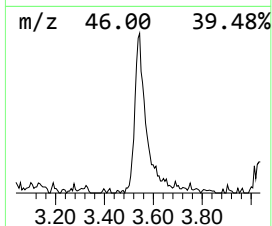
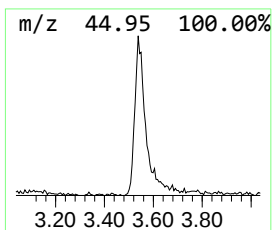
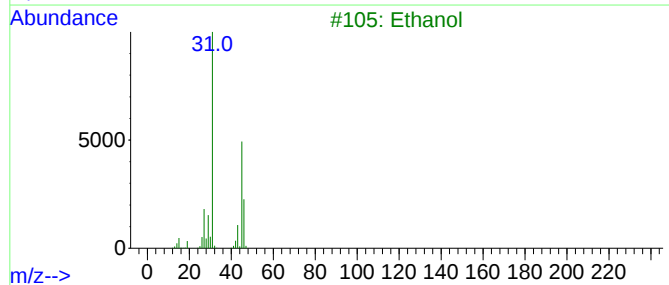
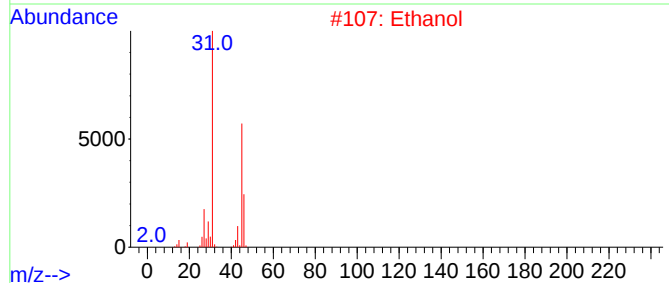
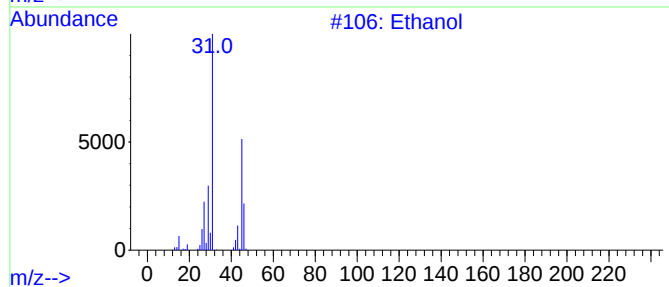
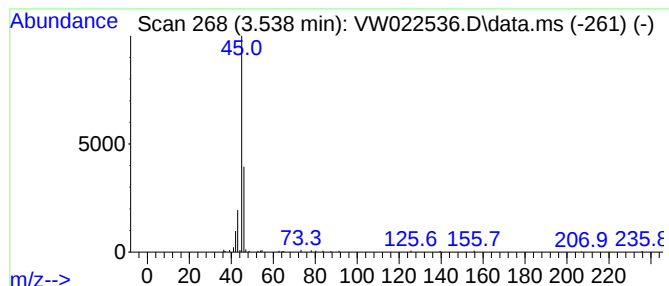
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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.
3.538	1.67 ug/L	52230	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol			46	C2H6O	000064-17-5	86
2	Ethanol			46	C2H6O	000064-17-5	64
3	Ethanol			46	C2H6O	000064-17-5	40
4	Dimethyl ether			46	C2H6O	000115-10-6	9
5	Ethanol			46	C2H6O	000064-17-5	9



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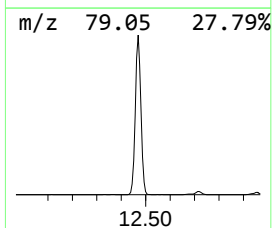
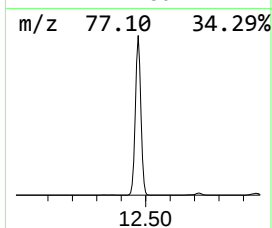
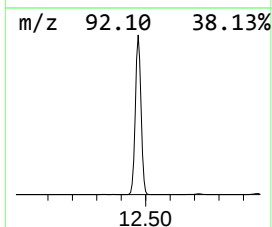
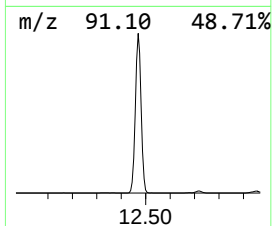
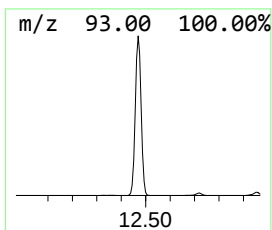
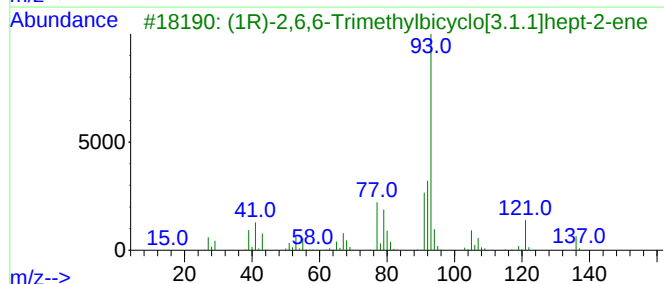
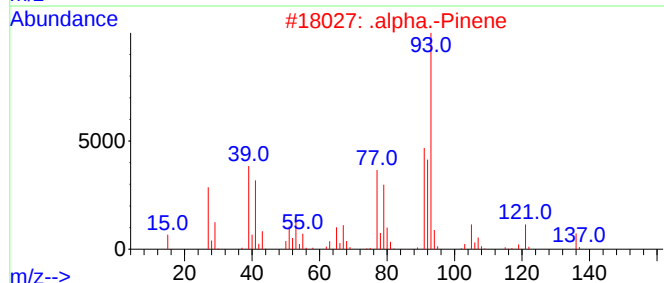
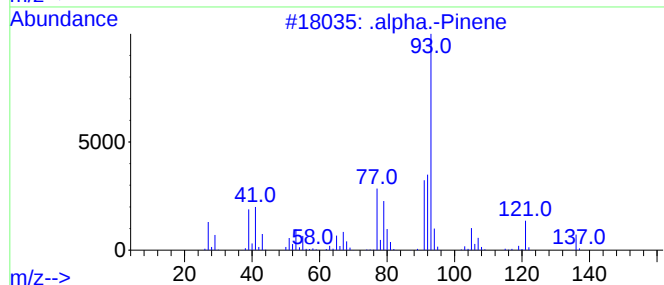
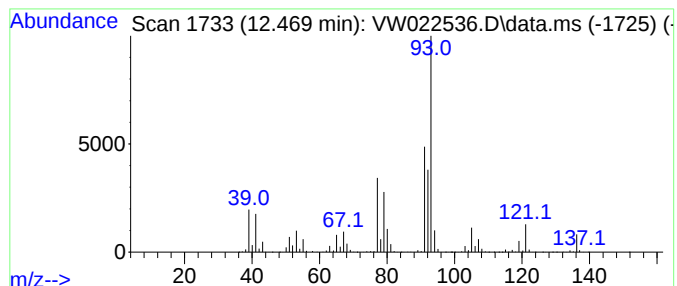
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 .alpha.-Pinene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	458.10 ug/L	15480600	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	.gamma.-Terpinene			136	C10H16	000099-85-4	93
5	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...			136	C10H16	004889-83-2	91



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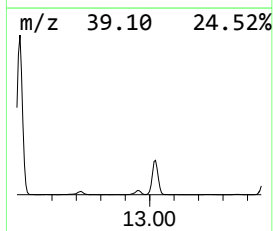
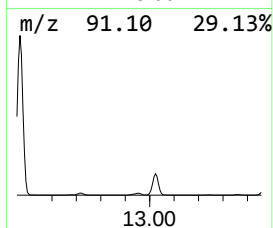
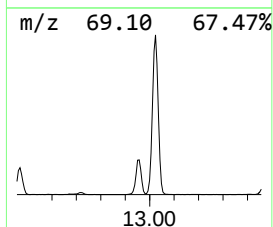
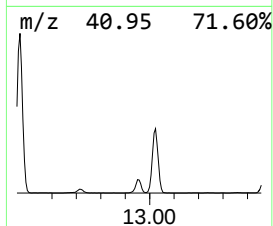
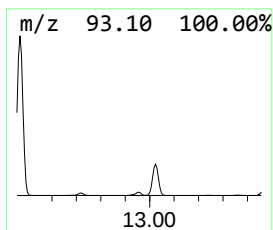
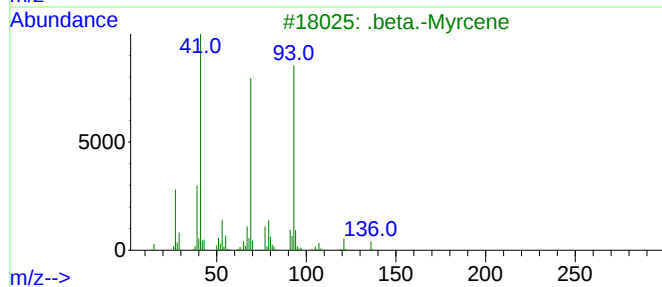
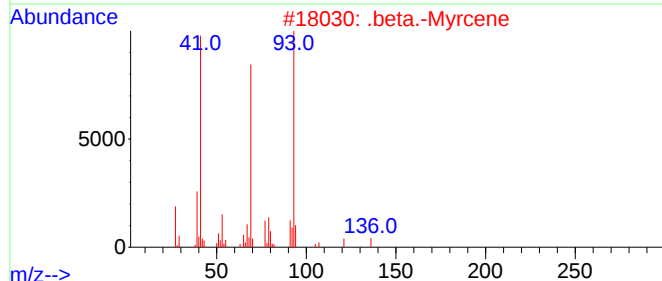
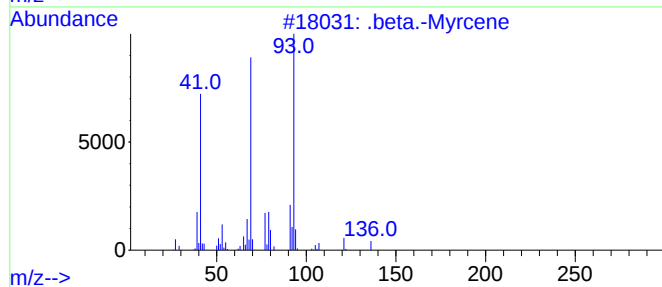
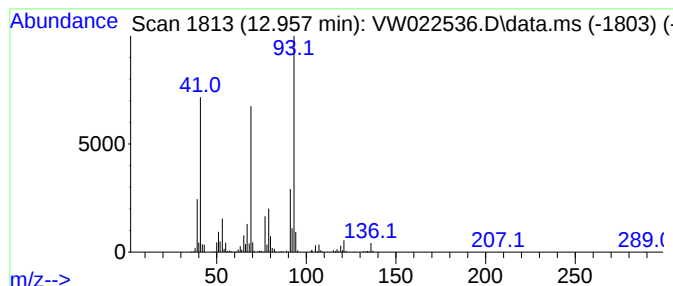
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Quant Title : SFAM01.0

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

Peak Number 6 .beta.-Myrcene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.957	11.92 ug/L	386348	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.-Myrcene	136	C10H16	000123-35-3	90	
3	.	beta.-Myrcene	136	C10H16	000123-35-3	90	
4	.	beta.-Myrcene	136	C10H16	000123-35-3	64	
5	Pyridine, 2-propyl-	121	C8H11N	000622-39-9	58		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041322\
Data File : VW022536.D
Acq On : 13 Apr 2022 17:00
Operator : SY/VA
Sample : N2374-07RE
Misc : 5.27g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFE3RE

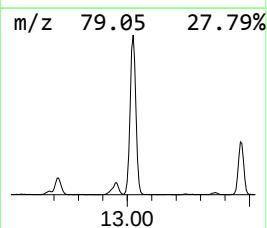
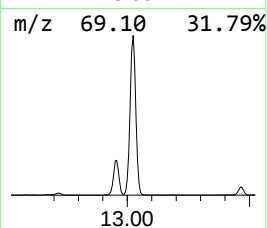
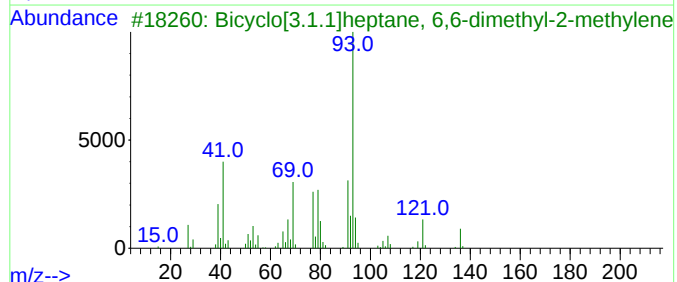
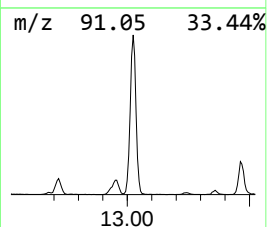
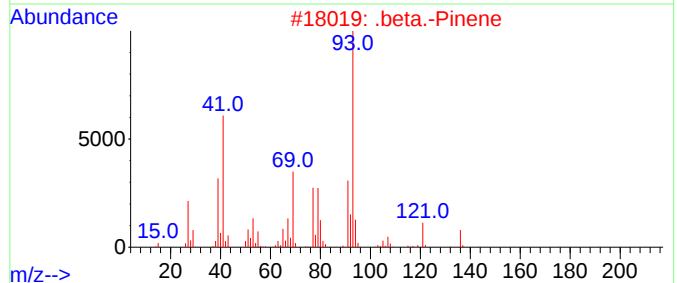
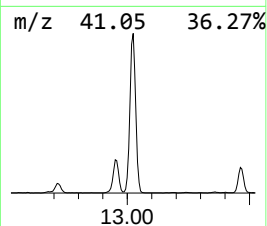
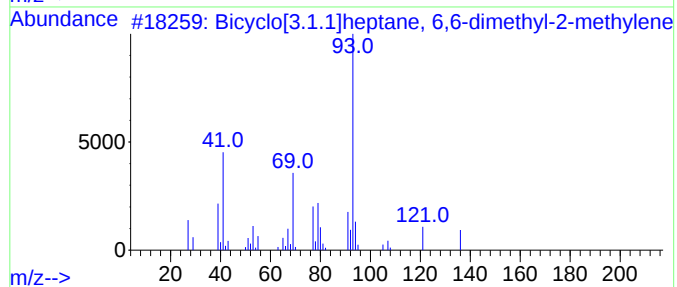
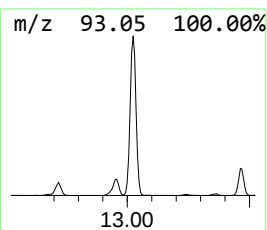
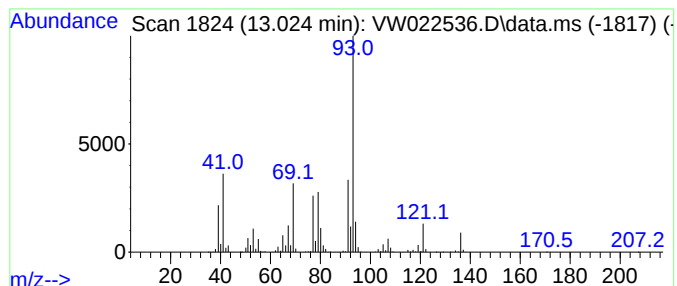
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Quant Title : SFAM01.0

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

Peak Number 7 Bicyclo[3.1.1]heptane, 6,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.024	95.35 ug/L	3089490	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	95
5			Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041322\
Data File : VW022536.D
Acq On : 13 Apr 2022 17:00
Operator : SY/VA
Sample : N2374-07RE
Misc : 5.27g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFE3RE

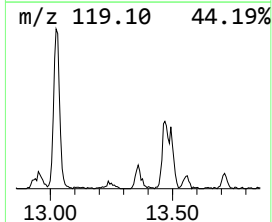
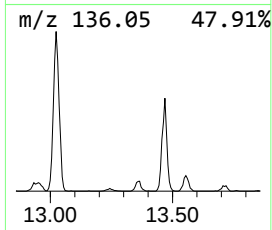
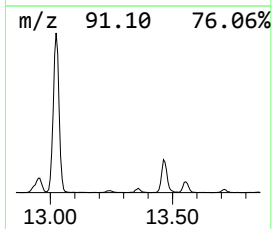
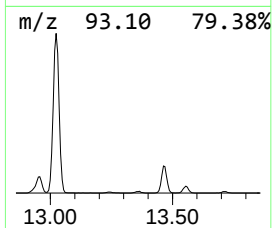
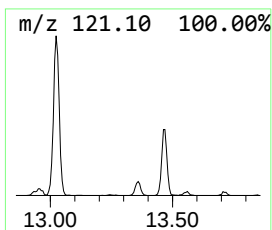
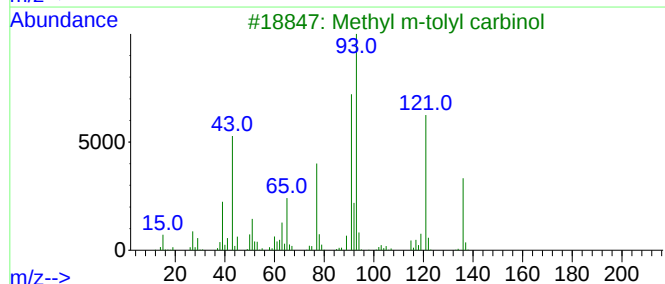
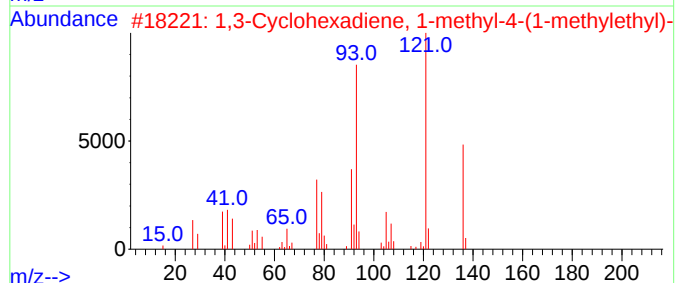
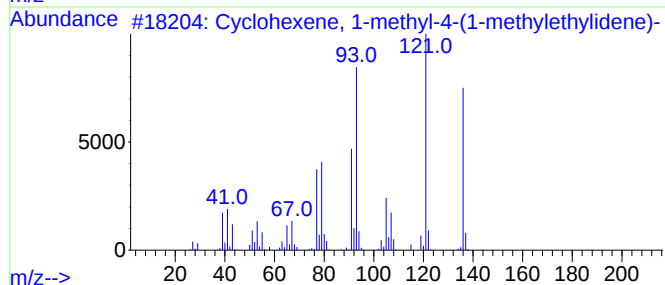
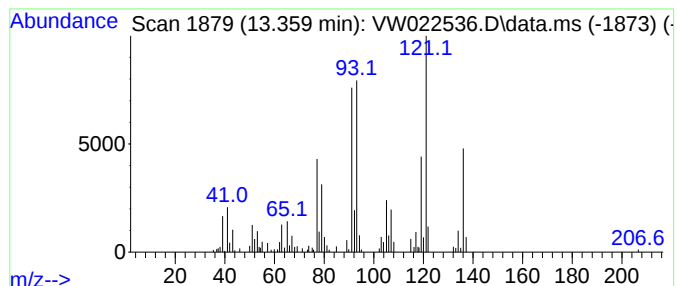
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Quant Title : SFAM01.0

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

Peak Number 8 Cyclohexene, 1-methyl-4-(1-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.359	1.83 ug/L	59326	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	95
2			1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	95
3			Methyl m-tolyl carbinol	136	C9H12O	1010342-29-9	93
4			(+)-4-Carene	136	C10H16	029050-33-7	93
5			Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041322\
Data File : VW022536.D
Acq On : 13 Apr 2022 17:00
Operator : SY/VA
Sample : N2374-07RE
Misc : 5.27g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFE3RE

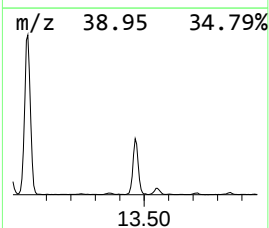
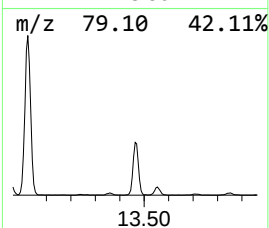
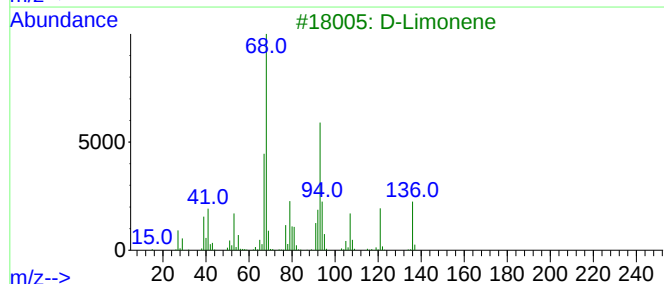
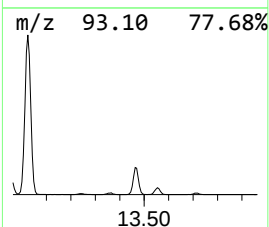
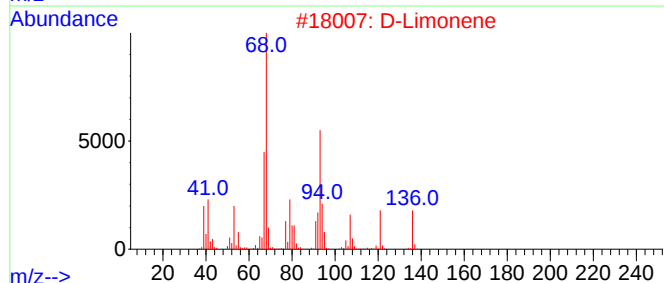
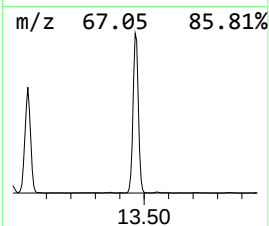
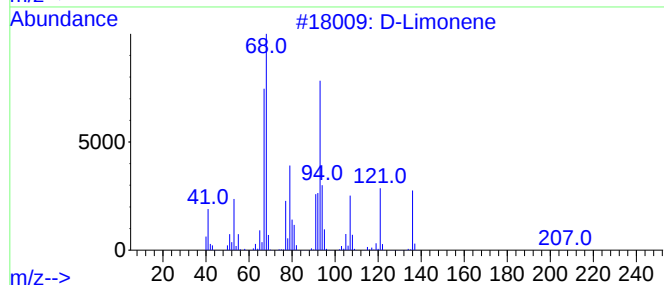
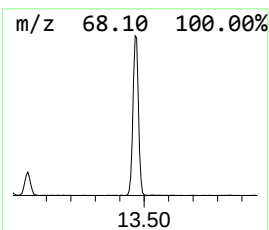
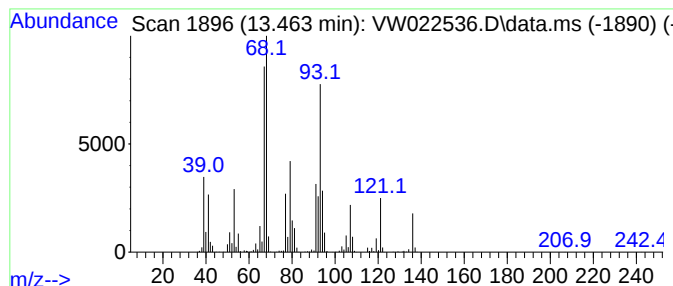
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Quant Title : SFAM01.0

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

Peak Number 9 D-Limonene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.463	33.16 ug/L	1074400	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	94
3	D-Limonene			136	C10H16	005989-27-5	93
4	D-Limonene			136	C10H16	005989-27-5	93
5	Limonene			136	C10H16	000138-86-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041322\
Data File : VW022536.D
Acq On : 13 Apr 2022 17:00
Operator : SY/VA
Sample : N2374-07RE
Misc : 5.27g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFE3RE

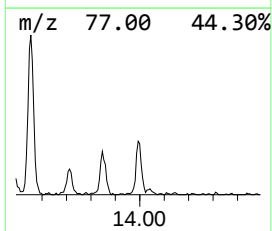
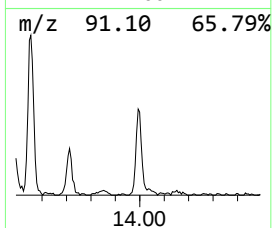
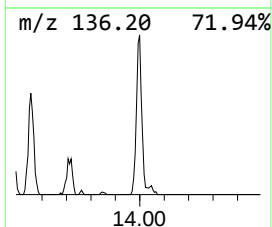
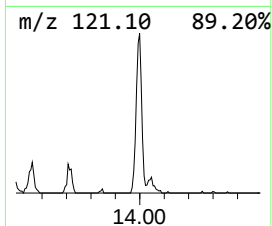
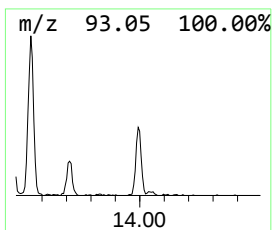
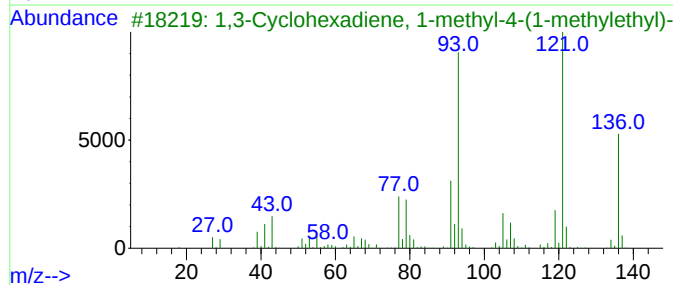
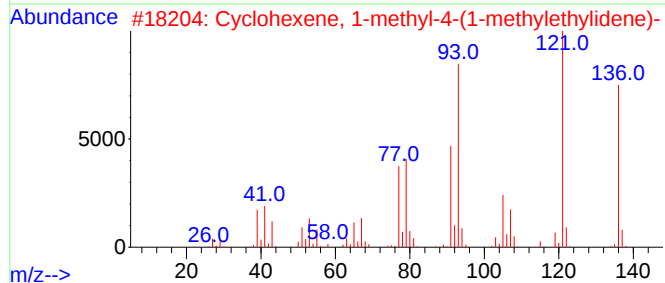
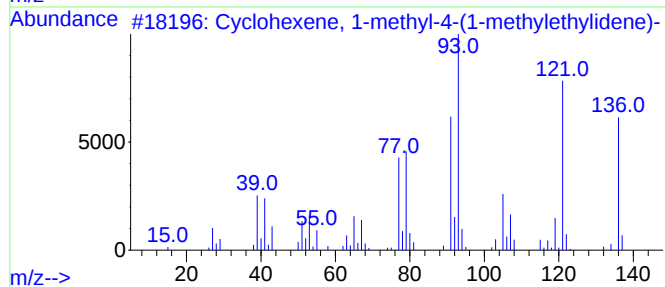
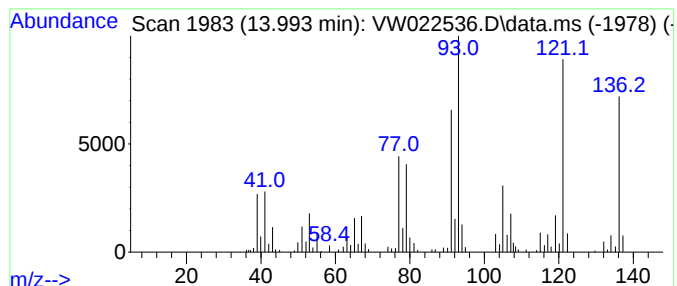
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Quant Title : SFAM01.0

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

Peak Number 10 1,3-Cyclohexadiene, 1-methy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.993	3.06 ug/L	99141	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	98
2			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	96
3			1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	96
4			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	95
5			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041322\
Data File : VW022536.D
Acq On : 13 Apr 2022 17:00
Operator : SY/VA
Sample : N2374-07RE
Misc : 5.27g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFFE3RE

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	3.538	1.7	ug/L	52230	1	8.842	782845	25.0
.alpha.-Pinene	12.469	458.1	ug/L	15480600	2	11.628	844834	25.0
.beta.-Myrcene	12.957	11.9	ug/L	386348	3	13.554	810070	25.0
Bicyclo[3.1.1]h...	13.024	95.3	ug/L	3089490	3	13.554	810070	25.0
Cyclohexene, 1-...	13.359	1.8	ug/L	59326	3	13.554	810070	25.0
D-Limonene	13.463	33.2	ug/L	1074400	3	13.554	810070	25.0
1,3-Cyclohexadi...	13.993	3.1	ug/L	99141	3	13.554	810070	25.0