

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW021721\
 Data File : VW020342.D
 Acq On : 17 Feb 2021 13:08
 Operator : SY/MD
 Sample : M1414-15ME 10X
 Misc : 5.64g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBH01ME

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_V\Method\SOMVLM020521WMA.M

Title : VOC Analysis

Signal : TIC: VW020342.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	77	83	101	rVB	255063	253458	9.27%	1.272%
2	1.568	157	164	175	rBV	217029	244313	8.94%	1.227%
3	2.108	322	332	351	rBV	439885	669196	24.48%	3.360%
4	2.433	429	433	434	rBV	702	430	0.02%	0.002%
5	2.507	452	456	470	rVB7	2287	3517	0.13%	0.018%
6	2.561	470	473	475	rBV	448	282	0.01%	0.001%
7	2.606	483	487	488	rBV3	766	548	0.02%	0.003%
8	2.677	506	509	514	rVB2	314	202	0.01%	0.001%
9	2.741	526	529	531	rBV3	318	190	0.01%	0.001%
10	2.812	549	551	552	rBV	211	95	0.00%	0.000%
11	2.880	569	572	578	rBB	226	176	0.01%	0.001%
12	2.986	603	605	608	rVB2	270	88	0.00%	0.000%
13	3.069	628	631	635	rBV	427	306	0.01%	0.002%
14	3.137	648	652	655	rBV2	171	150	0.01%	0.001%
15	3.188	664	668	673	rBV2	227	251	0.01%	0.001%
16	3.214	674	676	679	rVB	157	72	0.00%	0.000%
17	3.298	700	702	705	rVB	266	123	0.00%	0.001%
18	3.320	705	709	713	rVB2	186	143	0.01%	0.001%
19	3.381	725	728	731	rBV2	197	157	0.01%	0.001%
20	3.462	751	753	756	rVB2	352	200	0.01%	0.001%
21	3.503	762	766	769	rBV2	269	222	0.01%	0.001%
22	3.519	769	771	772	rBV	231	95	0.00%	0.000%
23	3.600	791	796	797	rBV2	250	185	0.01%	0.001%
24	3.651	809	812	814	rBV2	125	85	0.00%	0.000%
25	3.664	814	816	818	rVB2	199	90	0.00%	0.000%
26	3.706	827	829	833	rBV	362	215	0.01%	0.001%
27	3.741	837	840	843	rVV2	159	105	0.00%	0.001%
28	3.780	849	852	857	rVB	257	226	0.01%	0.001%
29	3.889	875	886	920	rBV	144093	423047	15.47%	2.124%
30	4.352	1011	1030	1056	rBV	304444	755921	27.65%	3.795%
31	4.709	1138	1141	1144	rBV3	288	208	0.01%	0.001%
32	4.783	1162	1164	1168	rBV3	276	229	0.01%	0.001%
33	4.812	1170	1173	1177	rBV3	257	299	0.01%	0.002%
34	4.851	1183	1185	1186	rBV3	310	111	0.00%	0.001%
35	4.870	1187	1191	1196	rVV3	353	375	0.01%	0.002%
36	4.889	1196	1197	1200	rVV2	255	112	0.00%	0.001%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
 Title : VOC Analysis

37	4.908	1200	1203	1205	rVV	262	143	0.01%	0.001%
38	4.918	1205	1206	1209	rVV2	243	111	0.00%	0.001%
39	4.934	1209	1211	1213	rVV	171	78	0.00%	0.000%
40	5.050	1229	1247	1278	rBV2	740453	1891262	69.18%	9.495%
41	5.391	1351	1353	1354	rBV	323	109	0.00%	0.001%
42	5.619	1412	1424	1453	rBV	577168	1146477	41.93%	5.756%
43	5.908	1502	1514	1530	rBV3	21726	45581	1.67%	0.229%
44	6.072	1549	1565	1590	rBV	417190	854120	31.24%	4.288%
45	6.333	1643	1646	1649	rBV3	448	307	0.01%	0.002%
46	6.387	1661	1663	1664	rBV3	179	86	0.00%	0.000%
47	6.445	1679	1681	1684	rVB2	368	158	0.01%	0.001%
48	6.484	1689	1693	1696	rBV	342	241	0.01%	0.001%
49	6.510	1699	1701	1703	rBV	174	75	0.00%	0.000%
50	6.574	1720	1721	1722	rBV	324	95	0.00%	0.000%
51	6.609	1729	1732	1733	rBV	135	77	0.00%	0.000%
52	6.645	1738	1743	1746	rBV4	698	758	0.03%	0.004%
53	6.760	1775	1779	1782	rVB2	341	254	0.01%	0.001%
54	6.818	1795	1797	1799	rVB	455	141	0.01%	0.001%
55	6.886	1813	1818	1822	rBV2	147	136	0.00%	0.001%
56	6.908	1822	1825	1826	rBV	204	128	0.00%	0.001%
57	6.944	1833	1836	1837	rBV	145	65	0.00%	0.000%
58	6.989	1839	1850	1876	rBV	216945	396197	14.49%	1.989%
59	7.320	1940	1953	1968	rBV	877831	1509949	55.23%	7.580%
60	7.625	2037	2048	2071	rBV	154446	269135	9.84%	1.351%
61	7.857	2118	2120	2123	rBV2	442	183	0.01%	0.001%
62	7.924	2139	2141	2144	rBV2	313	158	0.01%	0.001%
63	7.982	2158	2159	2160	rBV2	341	97	0.00%	0.000%
64	8.095	2183	2194	2240	rBV	491281	986750	36.09%	4.954%
65	8.564	2337	2340	2343	rBV2	489	309	0.01%	0.002%
66	8.670	2371	2373	2376	rBB3	253	112	0.00%	0.001%
67	8.706	2380	2384	2388	rBV4	510	401	0.01%	0.002%
68	8.812	2415	2417	2419	rBV	278	109	0.00%	0.001%
69	8.857	2419	2431	2457	rBV	888858	1434726	52.48%	7.203%
70	9.149	2516	2522	2532	rVB4	3653	5612	0.21%	0.028%
71	9.288	2562	2565	2567	rBV	211	143	0.01%	0.001%
72	9.352	2582	2585	2587	rBV	303	189	0.01%	0.001%
73	9.391	2595	2597	2598	rBV	227	74	0.00%	0.000%
74	9.419	2604	2606	2609	rBV2	326	188	0.01%	0.001%
75	9.574	2642	2654	2665	rBV6	3695	8097	0.30%	0.041%
76	9.754	2702	2710	2711	rBV3	6045	5549	0.20%	0.028%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Filtering: 5

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Title : VOC Analysis

77	9.889	2748	2752	2754	rBV2	414	274	0.01%	0.001%
78	10.011	2789	2790	2793	rVB2	436	162	0.01%	0.001%
79	10.056	2801	2804	2810	rBV2	365	385	0.01%	0.002%
80	10.104	2813	2819	2825	rBV4	2200	2761	0.10%	0.014%
81	10.223	2838	2856	2879	rVB	590255	1003872	36.72%	5.040%
82	10.355	2895	2897	2898	rBV	355	152	0.01%	0.001%
83	10.458	2921	2929	2945	rVB6	9123	15228	0.56%	0.076%
84	10.551	2949	2958	2963	rBV7	5690	8439	0.31%	0.042%
85	10.783	3024	3030	3039	rBV5	958	1232	0.05%	0.006%
86	10.915	3060	3071	3081	rBV3	17929	31677	1.16%	0.159%
87	11.091	3108	3126	3142	rBV2	372362	700796	25.63%	3.518%
88	11.198	3149	3159	3167	rBV	734483	1078449	39.45%	5.414%
89	11.252	3167	3176	3196	rVB3	1707708	2733971	100.00%	13.725%
90	11.358	3198	3209	3230	rVB	442266	717080	26.23%	3.600%
91	11.542	3258	3266	3281	rVB3	36487	53694	1.96%	0.270%
92	11.632	3282	3294	3310	rBV	1004115	1563613	57.19%	7.850%
93	11.908	3369	3380	3387	rBV	200781	308569	11.29%	1.549%
94	12.095	3428	3438	3453	rBV	190733	287412	10.51%	1.443%
95	12.230	3478	3480	3482	rBV	654	338	0.01%	0.002%
96	12.358	3512	3520	3522	rBV6	3510	4107	0.15%	0.021%
97	12.738	3628	3638	3652	rBV2	103855	188191	6.88%	0.945%
98	12.963	3699	3708	3713	rBV	40960	63556	2.32%	0.319%
99	13.374	3826	3836	3859	rBV2	101804	221962	8.12%	1.114%
100	14.349	4131	4139	4150	rBV5	12228	19320	0.71%	0.097%

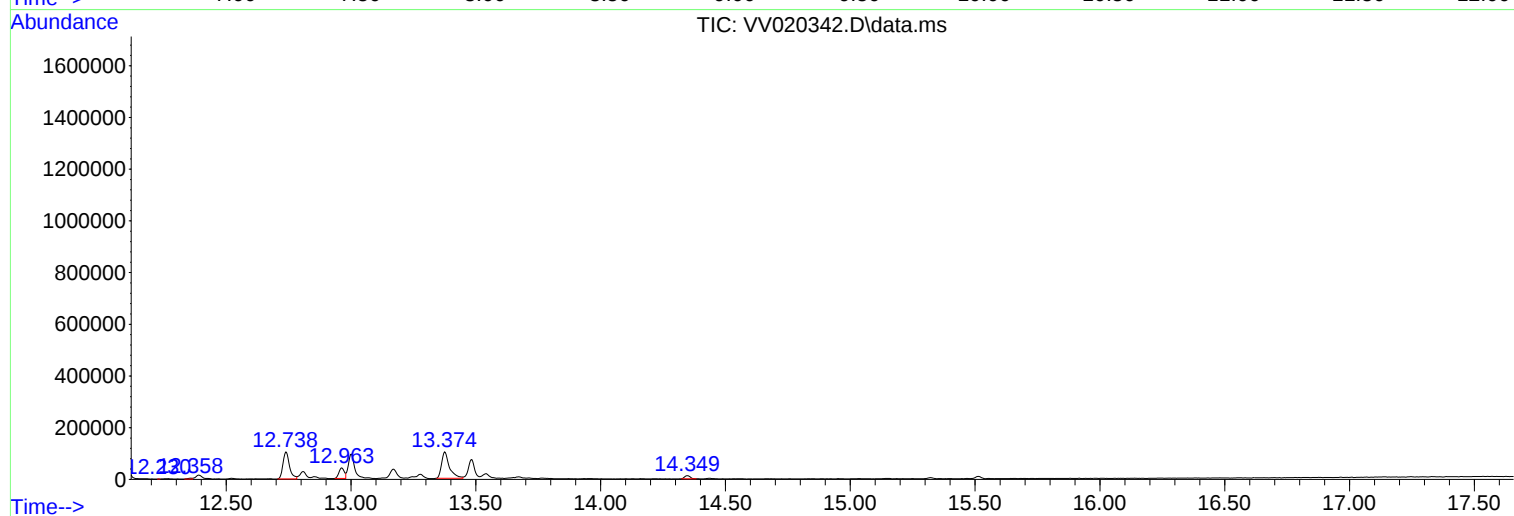
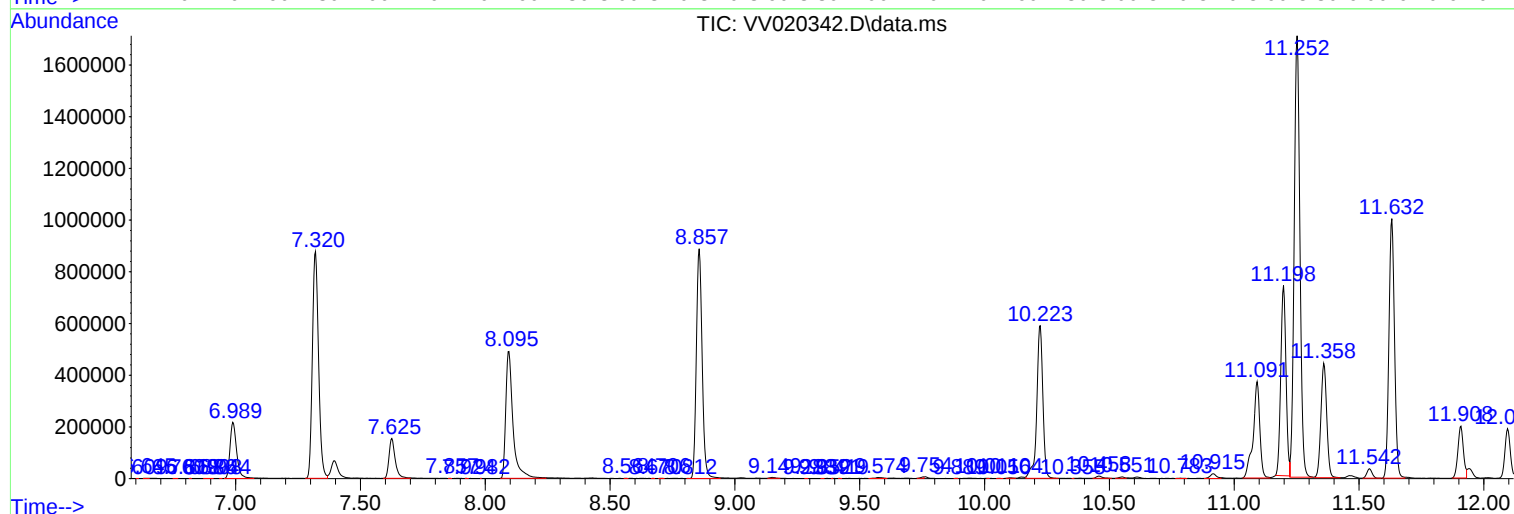
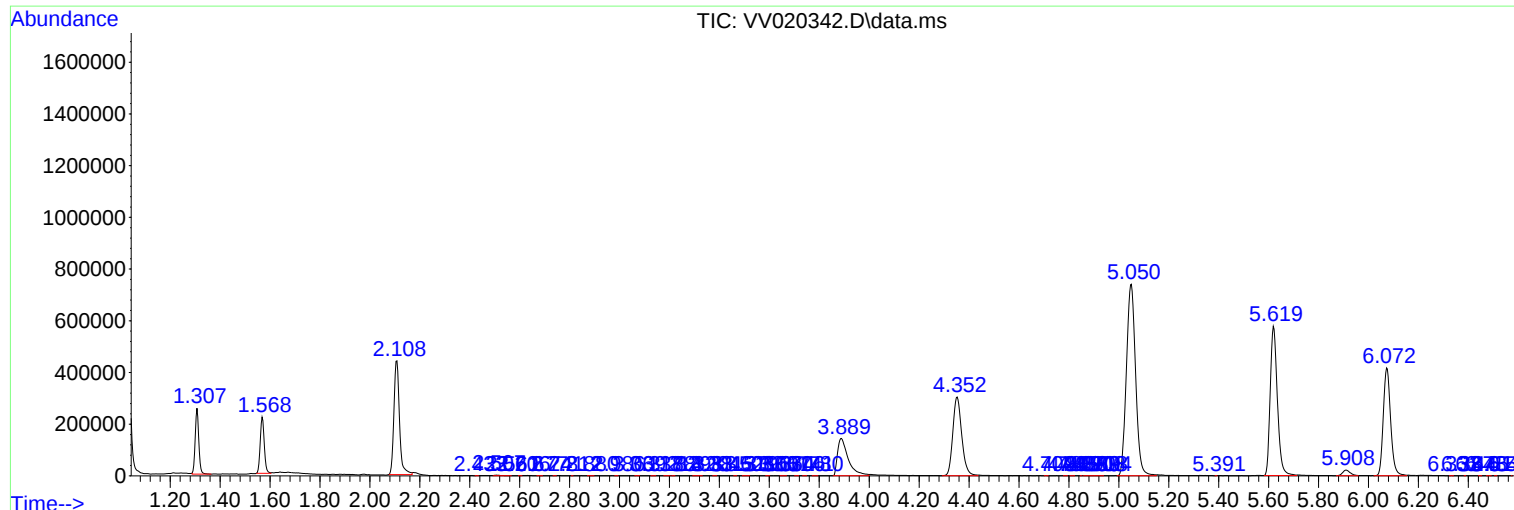
Sum of corrected areas: 19919042

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ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBH01ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
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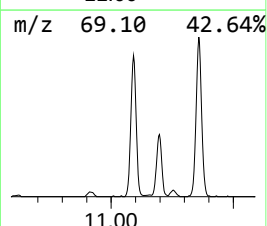
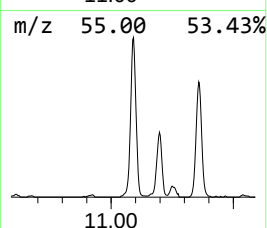
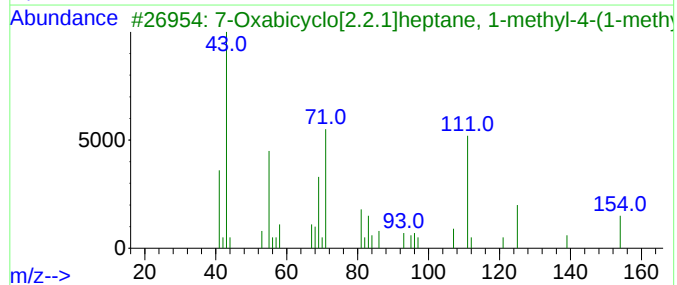
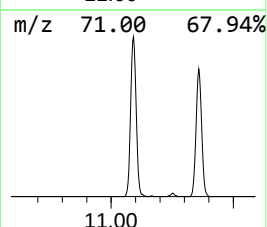
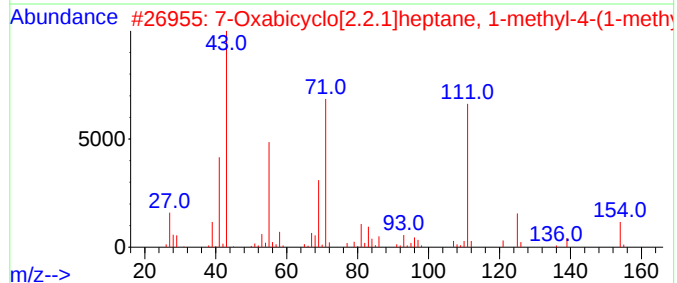
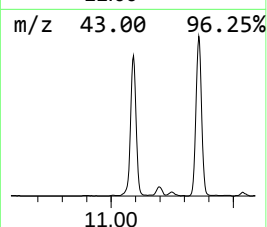
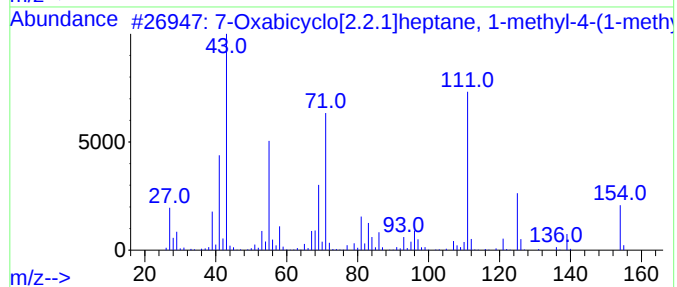
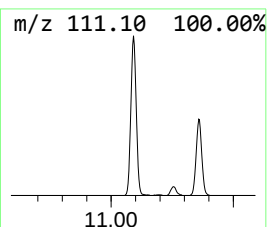
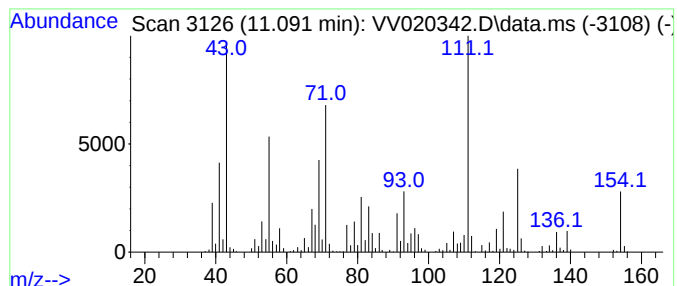
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 7-Oxabicyclo[2.2.1]heptane,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.091	12.82 ug/L	700796	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	93
2			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	76
3			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	64
4			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	43
5			Terpinen-4-ol	154	C10H18O	000562-74-3	38



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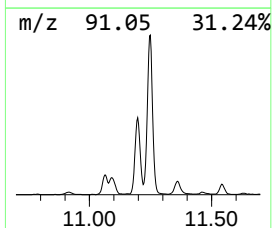
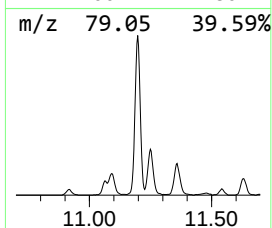
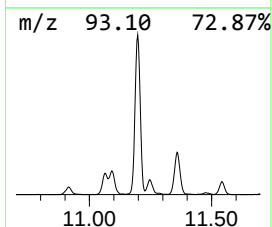
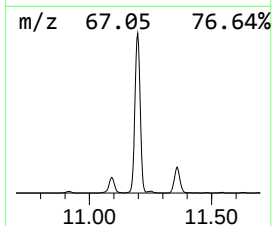
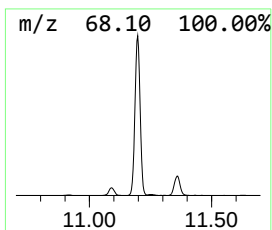
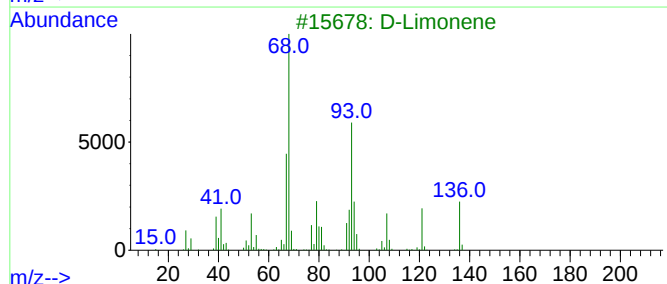
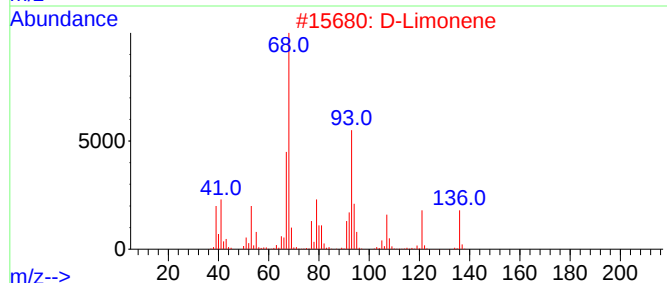
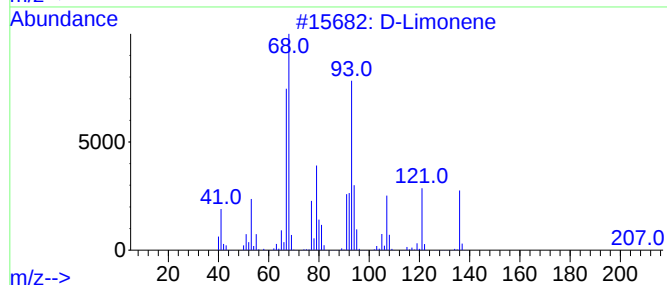
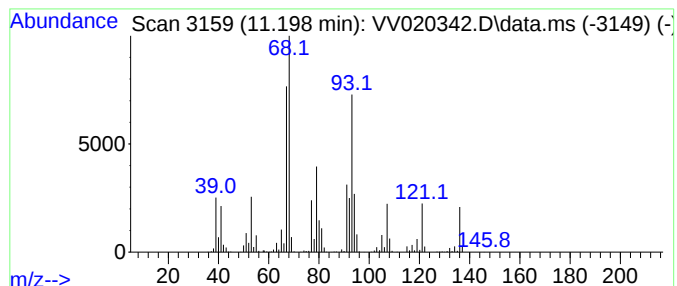
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 D-Limonene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.198	19.72 ug/L	1078450	1,4-Dichlorobenzene-d4	11.255

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	D-Limonene			136	C10H16	005989-27-5	94
4	Limonene			136	C10H16	000138-86-3	94
5	D-Limonene			136	C10H16	005989-27-5	93



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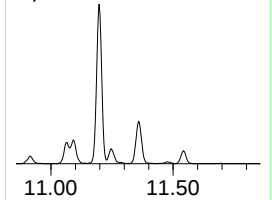
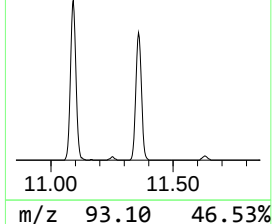
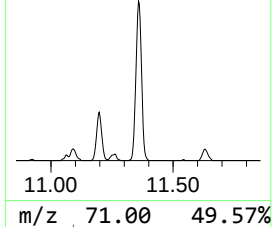
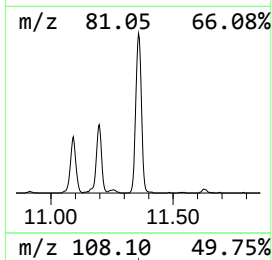
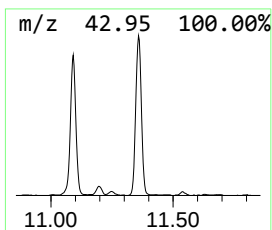
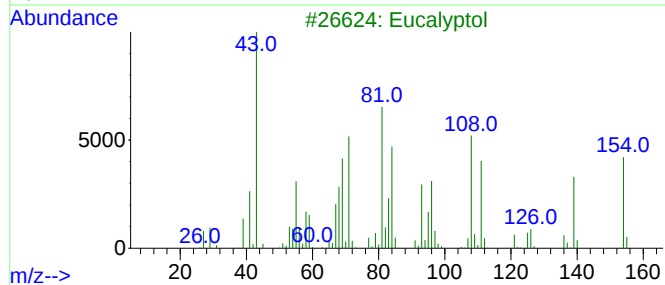
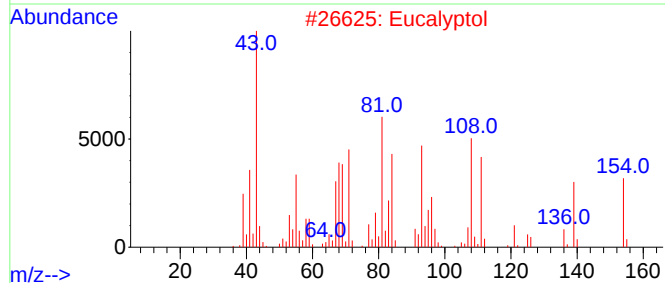
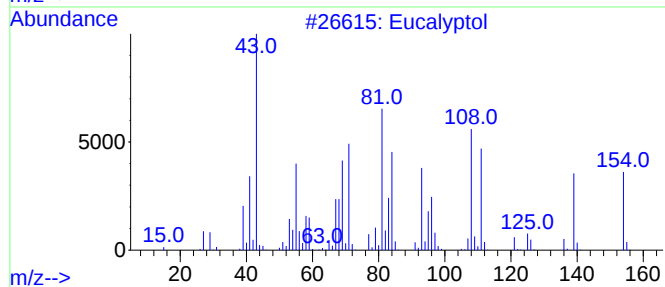
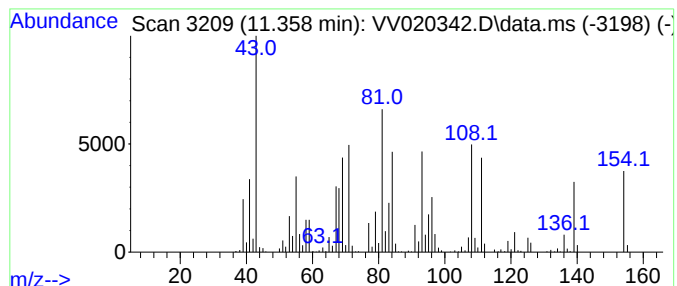
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Eucalyptol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.358	13.11 ug/L	717080	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eucalyptol	154	C10H18O	000470-82-6	99
2			Eucalyptol	154	C10H18O	000470-82-6	96
3			Eucalyptol	154	C10H18O	000470-82-6	95
4			Eucalyptol	154	C10H18O	000470-82-6	93
5			Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020342.D
Acq On : 17 Feb 2021 13:08
Operator : SY/MD
Sample : M1414-15ME 10X
Misc : 5.64g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBH01ME

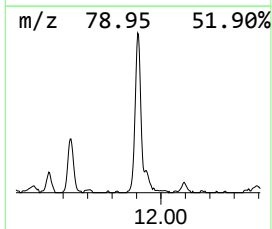
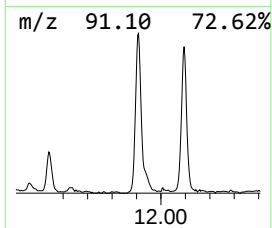
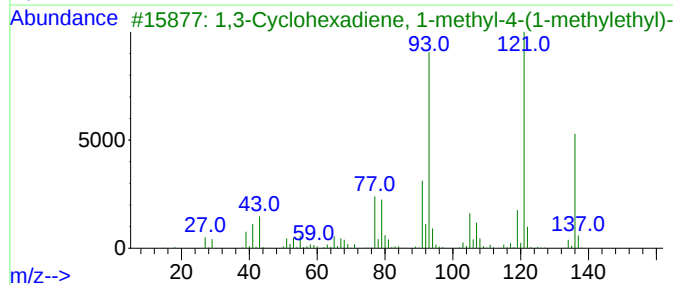
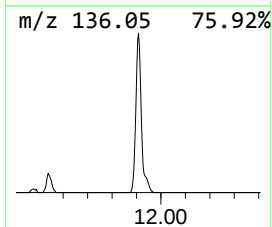
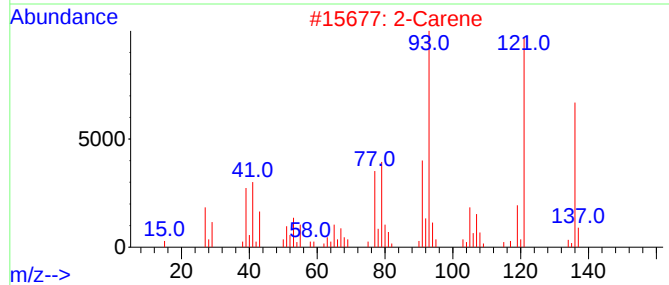
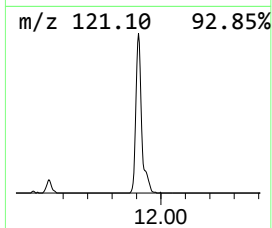
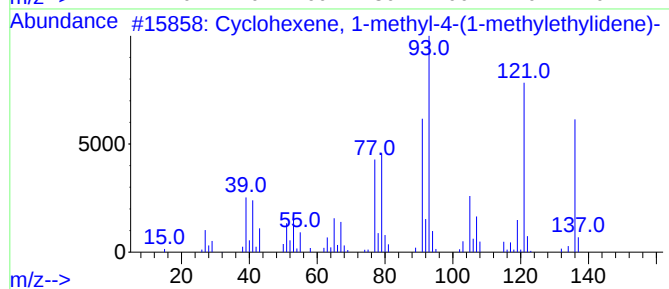
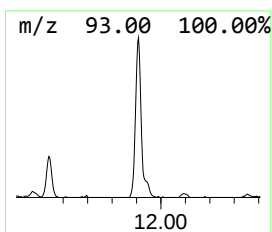
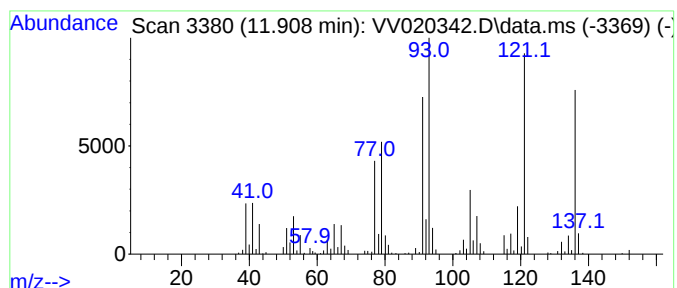
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclohexene, 1-methyl-4-(1-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.908	5.64 ug/L	308569	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	98
2			2-Carene	136	C10H16	000554-61-0	94
3			1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	93
4			(+)-4-Carene	136	C10H16	029050-33-7	93
5			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020342.D
Acq On : 17 Feb 2021 13:08
Operator : SY/MD
Sample : M1414-15ME 10X
Misc : 5.64g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBH01ME

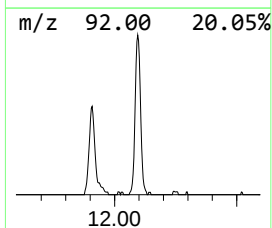
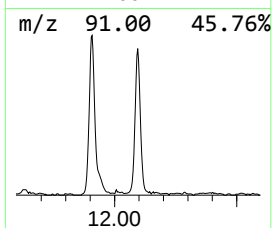
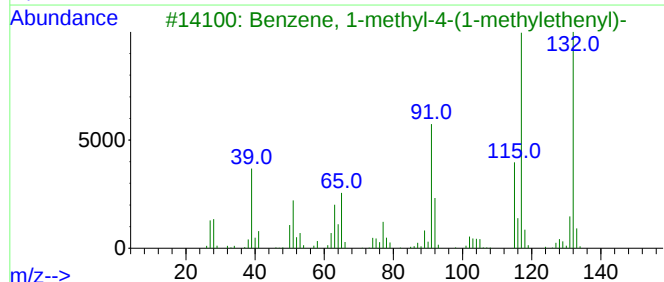
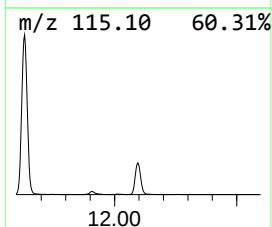
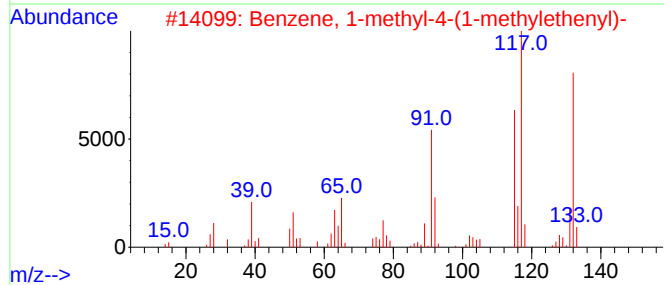
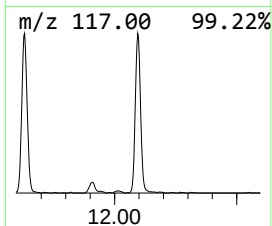
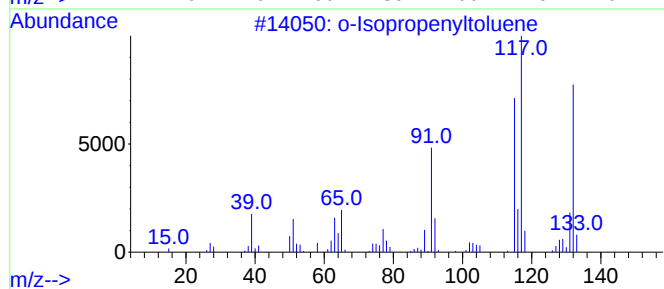
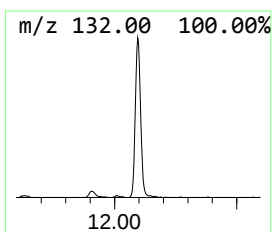
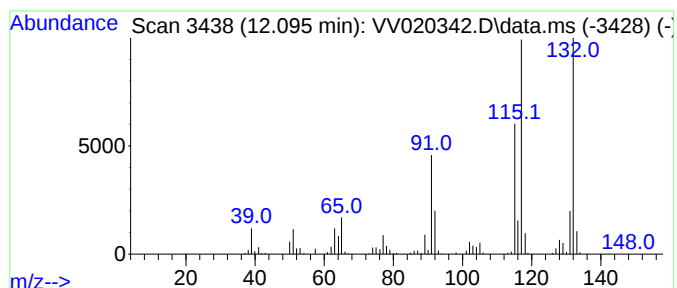
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 o-Isopropenyltoluene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	5.26 ug/L	287412	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Isopropenyltoluene	132	C10H12	007399-49-7	97
2			Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	96
3			Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	95
4			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	93
5			Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020342.D
Acq On : 17 Feb 2021 13:08
Operator : SY/MD
Sample : M1414-15ME 10X
Misc : 5.64g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBH01ME

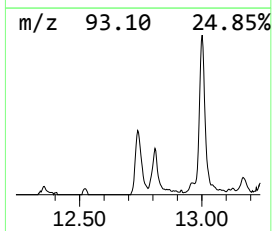
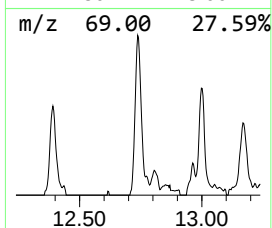
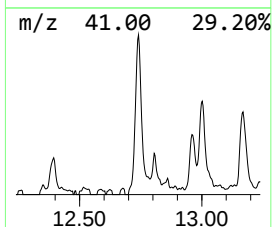
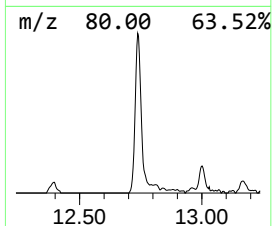
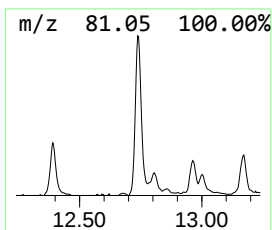
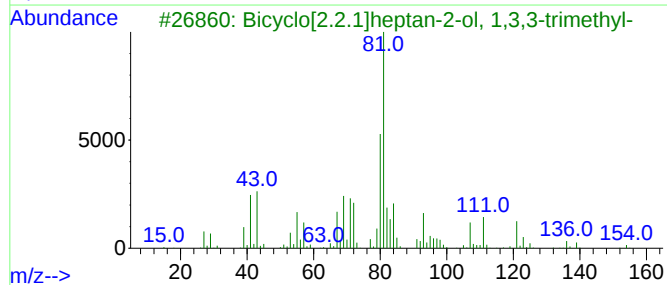
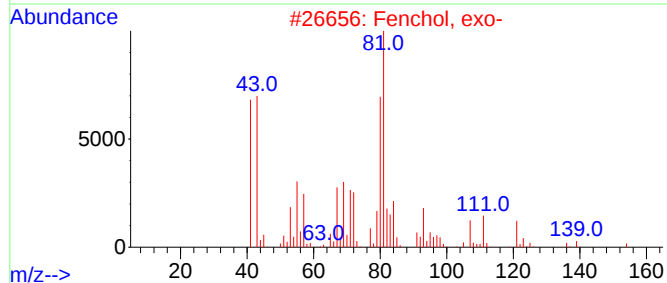
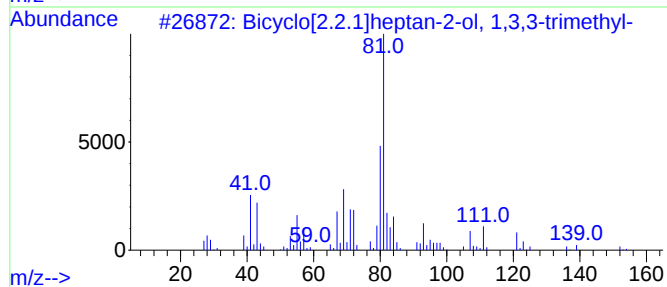
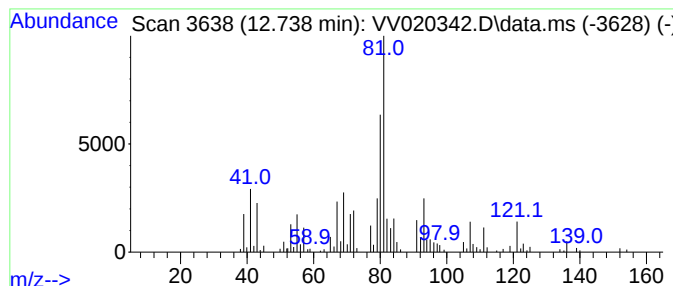
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.738	3.44 ug/L	188191	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	95
2			Fenchol, exo-	154	C10H18O	022627-95-8	95
3			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	94
4			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	002217-02-9	91
5			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020342.D
Acq On : 17 Feb 2021 13:08
Operator : SY/MD
Sample : M1414-15ME 10X
Misc : 5.64g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBH01ME

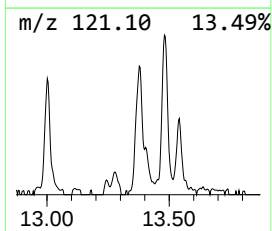
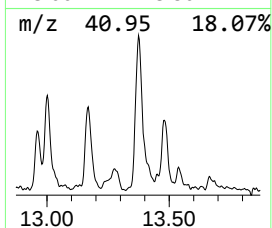
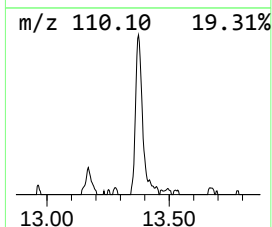
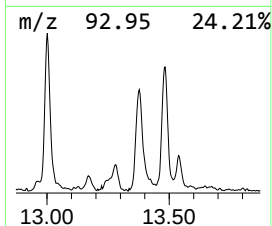
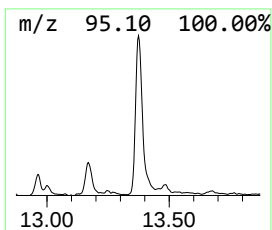
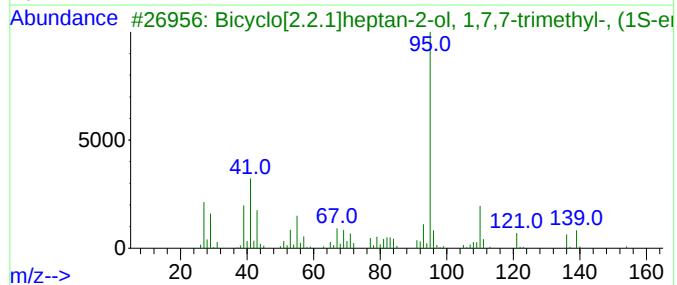
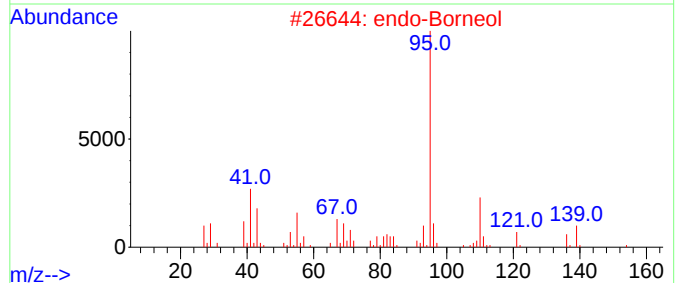
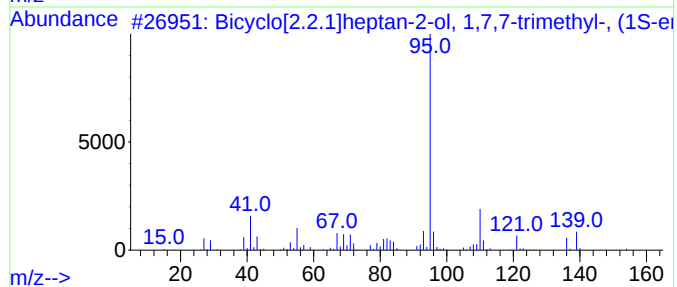
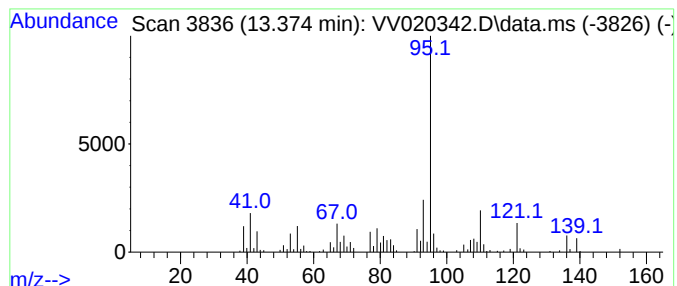
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.374	4.06 ug/L	221962	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	72
2			endo-Borneol	154	C10H18O	000507-70-0	64
3			Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	64
4			Isoborneol	154	C10H18O	000124-76-5	64
5			Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW021721\
Data File : VW020342.D
Acq On : 17 Feb 2021 13:08
Operator : SY/MD
Sample : M1414-15ME 10X
Misc : 5.64g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBH01ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
7-Oxabicyclo[2....	11.091	12.8	ug/L	700796	3	11.255	2733970	50.0
D-Limonene	11.198	19.7	ug/L	1078450	3	11.255	2733970	50.0
Eucalyptol	11.358	13.1	ug/L	717080	3	11.255	2733970	50.0
Cyclohexene, 1-...	11.908	5.6	ug/L	308569	3	11.255	2733970	50.0
o-Isopropenylto...	12.095	5.3	ug/L	287412	3	11.255	2733970	50.0
Bicyclo[2.2.1]h...	12.738	3.4	ug/L	188191	3	11.255	2733970	50.0
Bicyclo[2.2.1]h...	13.374	4.1	ug/L	221962	3	11.255	2733970	50.0