

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
 Data File : VV020344.D
 Acq On : 17 Feb 2021 14:18
 Operator : SY/MD
 Sample : M1414-14ME 10X
 Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBGZ9ME

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: VV020344.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	76	83	100	rVB	261255	267477	9.95%	1.068%
2	1.568	157	164	177	rBV	230552	260412	9.69%	1.040%
3	1.976	284	291	302	rVB	19103	26286	0.98%	0.105%
4	2.105	320	331	348	rBV	461014	704906	26.23%	2.815%
5	2.342	402	405	409	rVV3	519	467	0.02%	0.002%
6	2.429	428	432	433	rBV2	677	461	0.02%	0.002%
7	2.449	434	438	442	rVV4	964	1210	0.05%	0.005%
8	2.513	451	458	467	rVB4	2297	3914	0.15%	0.016%
9	2.574	476	477	481	rVB2	550	327	0.01%	0.001%
10	2.600	482	485	488	rBV4	1010	889	0.03%	0.004%
11	2.667	502	506	508	rBV2	404	327	0.01%	0.001%
12	2.719	518	522	527	rVB2	412	382	0.01%	0.002%
13	2.757	527	534	537	rBV4	707	781	0.03%	0.003%
14	3.522	767	772	776	rVV5	832	889	0.03%	0.004%
15	3.548	776	780	783	rVB	563	442	0.02%	0.002%
16	3.822	860	865	869	rBV2	333	339	0.01%	0.001%
17	3.892	875	887	919	rBV	154727	472391	17.58%	1.887%
18	4.352	1012	1030	1064	rBV	323976	816696	30.39%	3.262%
19	4.947	1206	1215	1216	rBV	264	328	0.01%	0.001%
20	4.960	1216	1219	1225	rVB3	357	333	0.01%	0.001%
21	5.050	1228	1247	1276	rBV2	801192	2051324	76.34%	8.193%
22	5.413	1357	1360	1366	rBV5	585	395	0.01%	0.002%
23	5.555	1391	1404	1411	rBV5	3473	8900	0.33%	0.036%
24	5.619	1411	1424	1450	rVV	691812	1379091	51.32%	5.508%
25	5.911	1504	1515	1531	rBV	23358	50104	1.86%	0.200%
26	6.072	1551	1565	1602	rVV	451166	926580	34.48%	3.701%
27	6.210	1606	1608	1610	rVV2	1144	793	0.03%	0.003%
28	6.236	1614	1616	1622	rVB5	1158	1117	0.04%	0.004%
29	6.645	1736	1743	1744	rBV3	1031	1016	0.04%	0.004%
30	6.686	1754	1756	1762	rVB4	538	378	0.01%	0.002%
31	6.764	1775	1780	1782	rBV2	370	332	0.01%	0.001%
32	6.908	1821	1825	1831	rBV3	316	341	0.01%	0.001%
33	6.989	1839	1850	1874	rBV	234680	432139	16.08%	1.726%
34	7.239	1921	1928	1940	rBV4	5833	10243	0.38%	0.041%
35	7.320	1940	1953	1967	rBV	938129	1616963	60.17%	6.458%
36	7.394	1967	1976	2000	rVB	228339	412834	15.36%	1.649%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Filtering: 5

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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

37	7.545	2021	2023	2025	rBV3	756	336	0.01%	0.001%
38	7.625	2037	2048	2070	rBV	167804	292779	10.90%	1.169%
39	7.879	2123	2127	2130	rBV3	560	483	0.02%	0.002%
40	7.944	2141	2147	2149	rBV3	339	383	0.01%	0.002%

41	8.095	2180	2194	2238	rBV	569118	1109616	41.29%	4.432%
42	8.413	2290	2293	2294	rBV	686	335	0.01%	0.001%
43	8.571	2339	2342	2347	rVB4	390	347	0.01%	0.001%
44	8.673	2371	2374	2377	rVB2	463	330	0.01%	0.001%
45	8.696	2377	2381	2382	rBV3	615	479	0.02%	0.002%

46	8.857	2415	2431	2458	rBV	1061320	1712433	63.73%	6.840%
47	9.024	2476	2483	2494	rVB2	9429	15056	0.56%	0.060%
48	9.079	2498	2500	2502	rBV3	615	362	0.01%	0.001%
49	9.146	2512	2521	2537	rBV	14340	25506	0.95%	0.102%
50	9.204	2537	2539	2542	rVV4	700	398	0.01%	0.002%

51	9.570	2638	2653	2665	rBV5	11313	27384	1.02%	0.109%
52	9.632	2669	2672	2683	rVV6	2278	3336	0.12%	0.013%
53	9.689	2683	2690	2699	rVV4	1400	2238	0.08%	0.009%
54	9.757	2703	2711	2721	rVB5	6857	10315	0.38%	0.041%
55	9.911	2751	2759	2763	rBV6	1425	1760	0.07%	0.007%

56	9.940	2763	2768	2776	rVB3	1163	1578	0.06%	0.006%
57	10.011	2786	2790	2793	rVB4	418	341	0.01%	0.001%
58	10.107	2813	2820	2824	rBV5	1444	2120	0.08%	0.008%
59	10.146	2825	2832	2835	rVV5	5316	6721	0.25%	0.027%
60	10.223	2837	2856	2881	rVB	651247	1094200	40.72%	4.370%

61	10.384	2900	2906	2911	rVB5	1371	1661	0.06%	0.007%
62	10.458	2918	2929	2942	rBV2	24079	38667	1.44%	0.154%
63	10.551	2949	2958	2966	rBV8	7145	10529	0.39%	0.042%
64	10.615	2971	2978	2987	rVV6	6812	10798	0.40%	0.043%
65	10.673	2991	2996	3003	rVV3	2224	2105	0.08%	0.008%

66	10.702	3003	3005	3010	rVB2	515	356	0.01%	0.001%
67	10.763	3022	3024	3026	rVV2	607	402	0.01%	0.002%
68	10.786	3026	3031	3040	rVB5	2408	2943	0.11%	0.012%
69	10.860	3048	3054	3055	rBV4	538	426	0.02%	0.002%
70	10.918	3058	3072	3093	rBV5	14312	34592	1.29%	0.138%

71	11.091	3108	3126	3144	rBV	960026	1635443	60.86%	6.532%
72	11.197	3148	3159	3167	rVV	709222	1042878	38.81%	4.165%
73	11.252	3167	3176	3197	rVB3	1691357	2687109	100.00%	10.733%
74	11.358	3197	3209	3226	rBV	740680	1177284	43.81%	4.702%
75	11.464	3237	3242	3252	rVB6	7456	13710	0.51%	0.055%

76	11.541	3256	3266	3276	rVB2	43375	63871	2.38%	0.255%
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 ClientSampleId :
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Integration Parameters: LSCINT.P

Integrator: RTE

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Title : VOC Analysis

77	11.631	3282	3294	3312	rBV	1064614	1653337	61.53%	6.604%
78	11.747	3329	3330	3338	rVB4	1063	882	0.03%	0.004%
79	11.792	3338	3344	3347	rBV5	1791	1971	0.07%	0.008%
80	11.905	3366	3379	3387	rBV2	219184	354558	13.19%	1.416%
81	12.094	3427	3438	3454	rVV	190985	289103	10.76%	1.155%
82	12.255	3481	3488	3494	rVV5	1432	1907	0.07%	0.008%
83	12.390	3512	3530	3548	rBV2	36086	68661	2.56%	0.274%
84	12.522	3564	3571	3579	rBV5	4275	5738	0.21%	0.023%
85	12.738	3626	3638	3651	rBV2	163918	299558	11.15%	1.196%
86	12.808	3653	3660	3668	rVB2	60653	93176	3.47%	0.372%
87	12.963	3696	3708	3713	rBV2	98238	159944	5.95%	0.639%
88	13.001	3713	3720	3737	rVB	261167	415933	15.48%	1.661%
89	13.168	3764	3772	3788	rVB2	75704	130313	4.85%	0.520%
90	13.246	3790	3796	3799	rBV6	13621	17859	0.66%	0.071%
91	13.374	3825	3836	3858	rBV	186544	404199	15.04%	1.614%
92	13.480	3860	3869	3880	rVV	313756	505411	18.81%	2.019%
93	13.538	3882	3887	3905	rVB3	67711	110618	4.12%	0.442%
94	14.204	4089	4094	4097	rBV6	1619	1579	0.06%	0.006%
95	14.345	4131	4138	4151	rVB7	10128	16052	0.60%	0.064%
96	14.834	4287	4290	4291	rBV2	873	426	0.02%	0.002%
97	15.149	4381	4388	4393	rVB9	2097	2430	0.09%	0.010%
98	15.316	4436	4440	4451	rVB9	4664	6053	0.23%	0.024%
99	15.512	4494	4501	4508	rBV10	9221	13071	0.49%	0.052%
100	15.728	4566	4568	4573	rVB4	1187	719	0.03%	0.003%

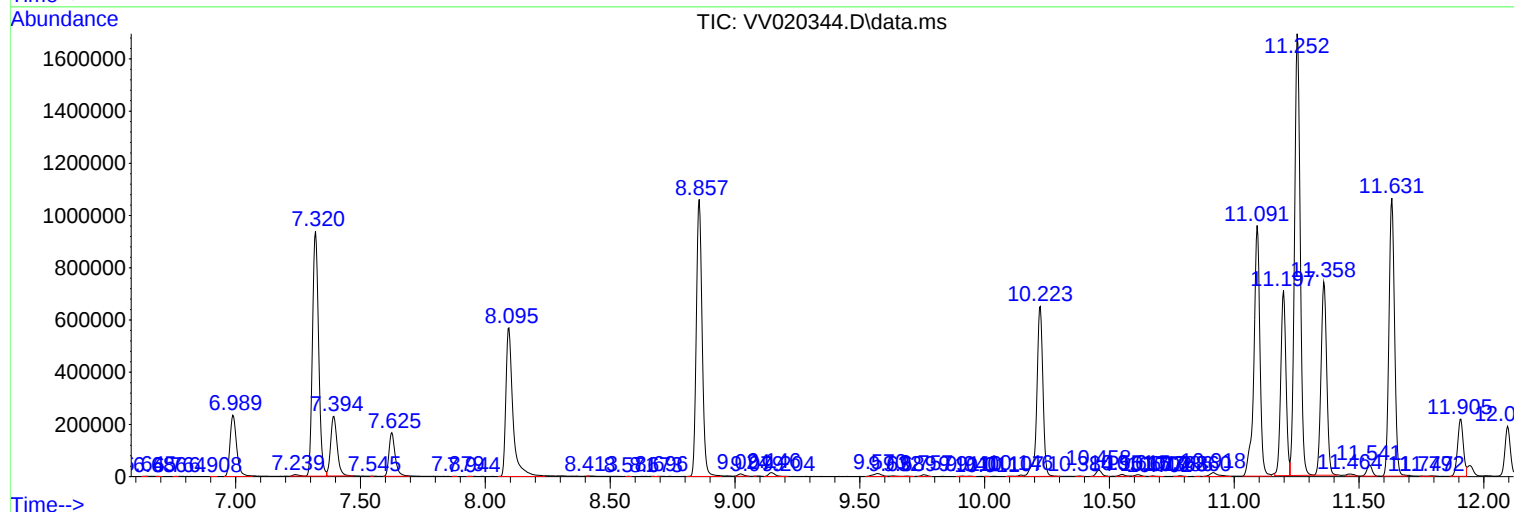
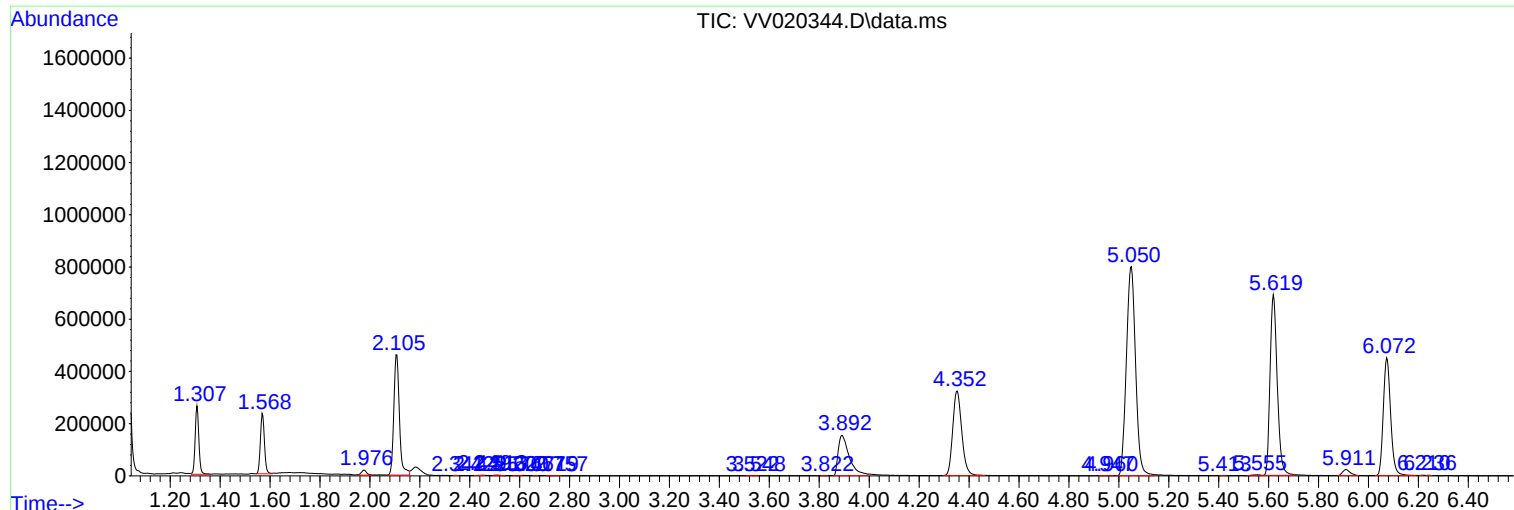
Sum of corrected areas: 25036915

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ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

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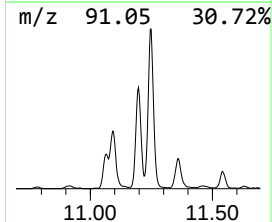
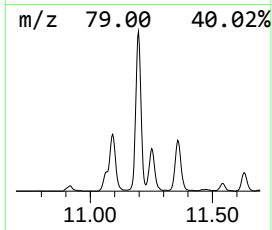
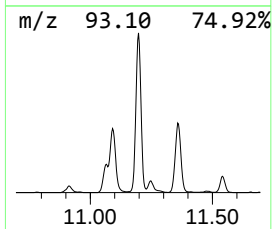
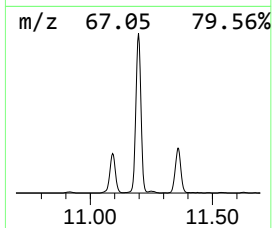
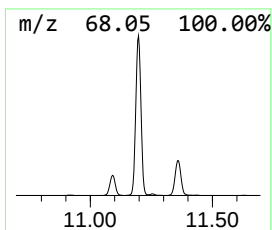
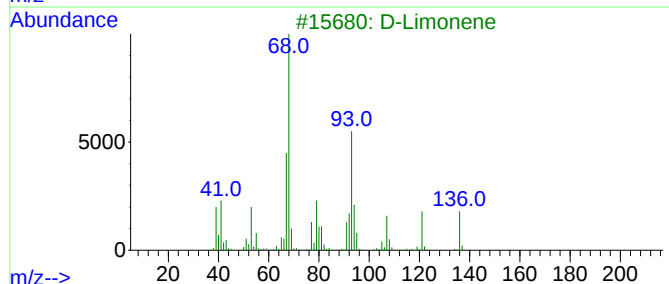
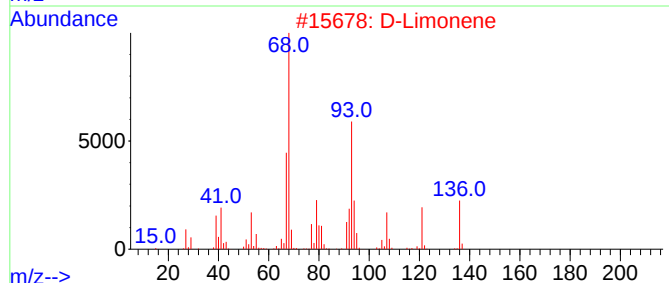
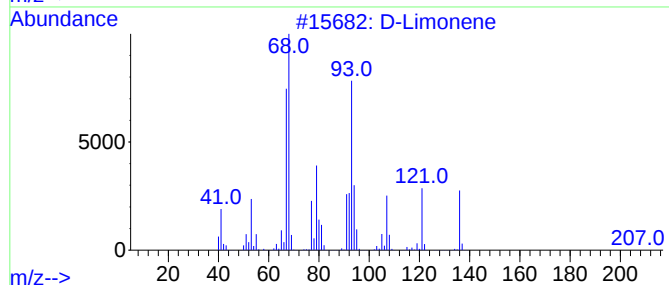
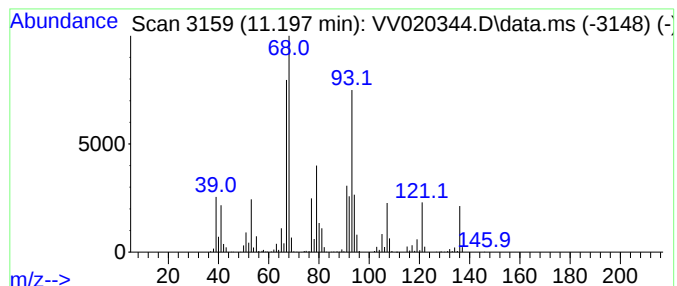
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 D-Limonene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.197	19.41 ug/L	1042880	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Limonene			136	C10H16	005989-27-5	99
2	D-Limonene			136	C10H16	005989-27-5	94
3	D-Limonene			136	C10H16	005989-27-5	94
4	D-Limonene			136	C10H16	005989-27-5	93
5	Limonene			136	C10H16	000138-86-3	93



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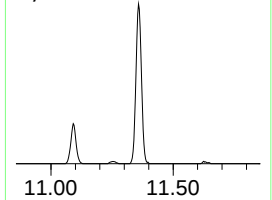
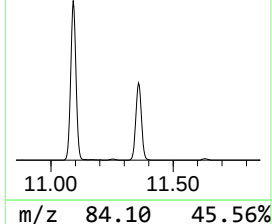
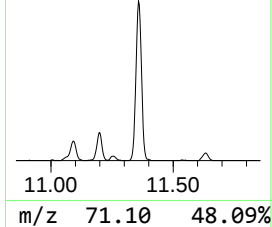
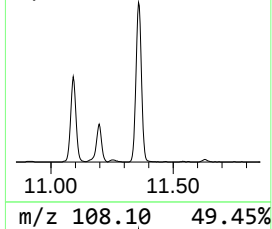
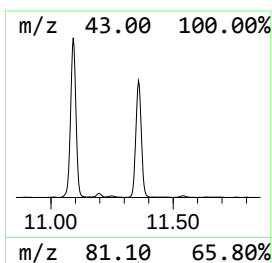
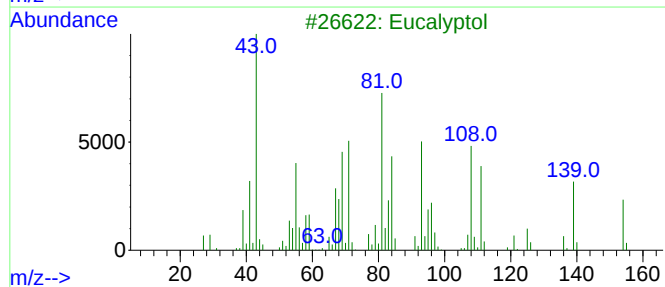
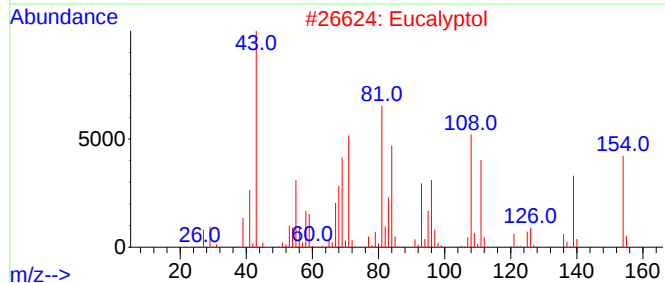
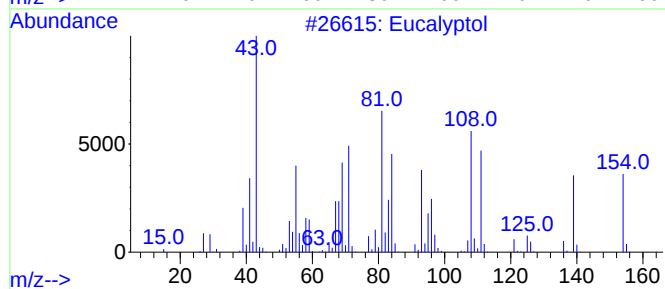
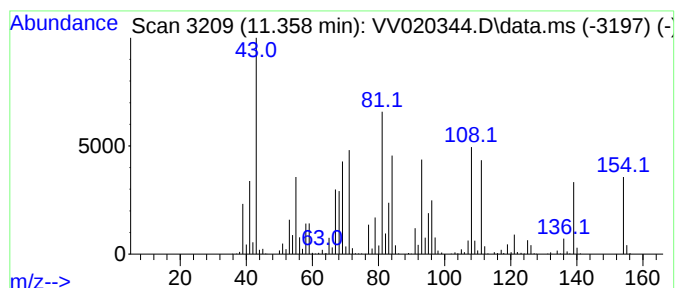
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM020521WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Eucalyptol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.358	21.91 ug/L	1177280	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	98
3	Eucalyptol			154	C10H18O	000470-82-6	95
4	Eucalyptol			154	C10H18O	000470-82-6	94
5	Trifluoroacetyl-.alpha.-terpineol			250	C12H17F3O2	1000058-17-6	83



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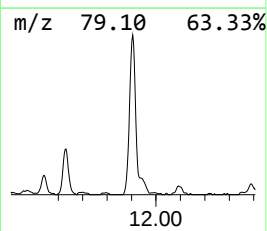
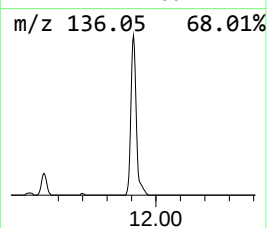
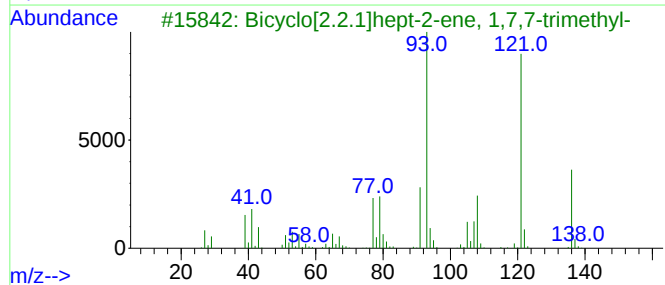
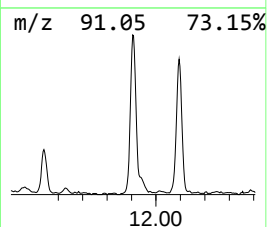
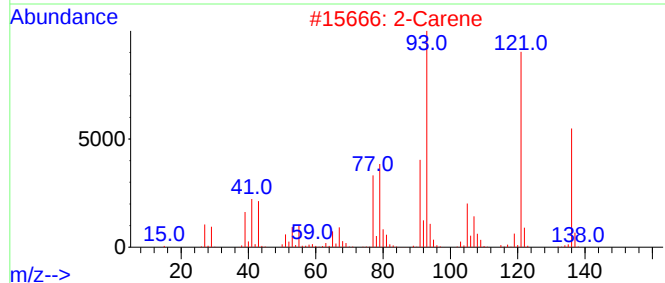
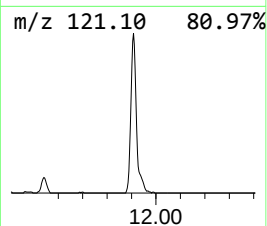
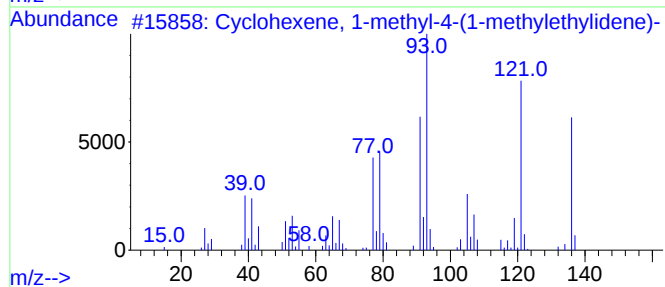
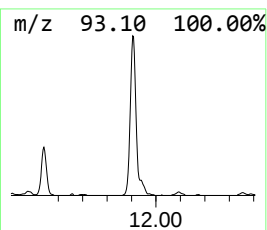
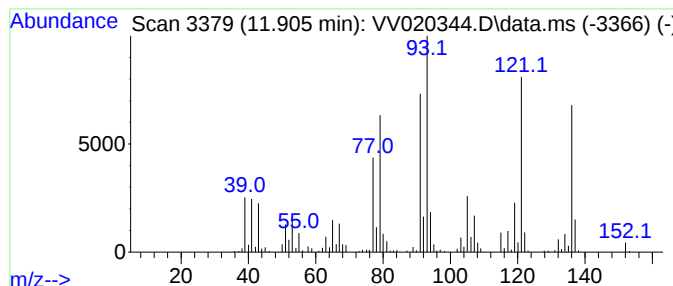
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Quant Title : VOC Analysis

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

Peak Number 5 Cyclohexene, 1-methyl-4-(1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.905	6.60 ug/L	354558	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	95
2			2-Carene	136	C10H16	000554-61-0	91
3			Bicyclo[2.2.1]hept-2-ene, 1,7,7-...	136	C10H16	000464-17-5	90
4			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	90
5			1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

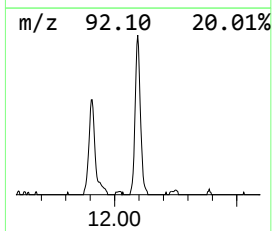
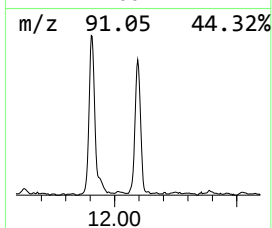
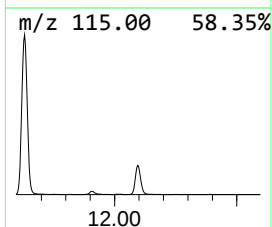
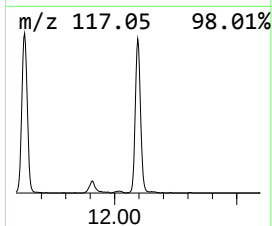
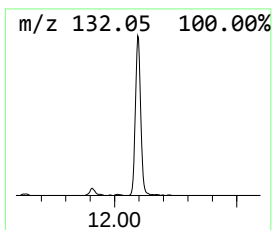
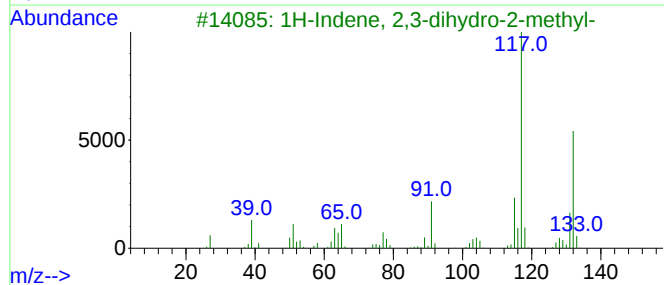
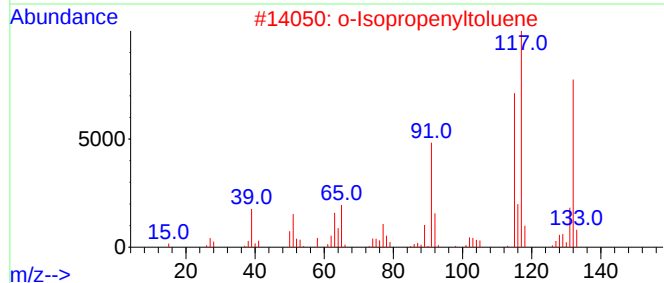
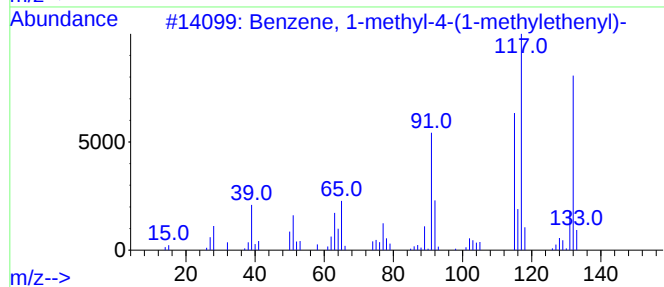
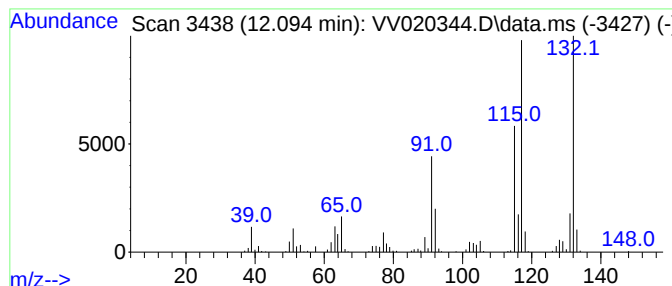
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-methyl-4-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	5.38 ug/L	289103	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	96
2			o-Isopropenyltoluene	132	C10H12	007399-49-7	95
3			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	90
4			Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	90
5			Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

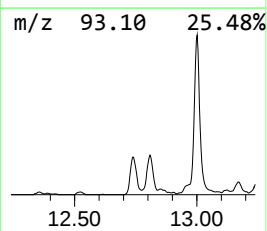
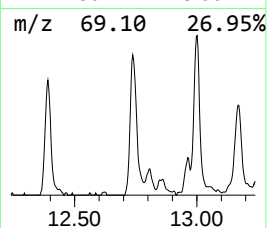
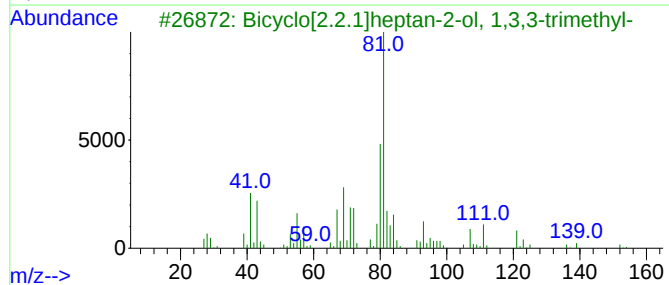
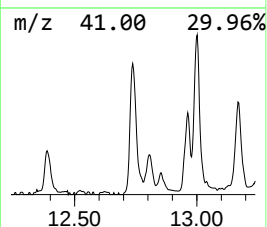
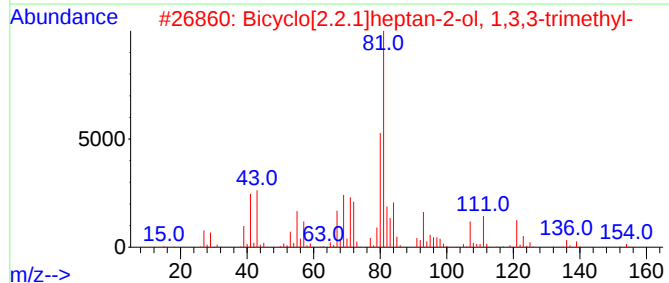
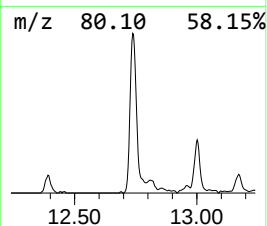
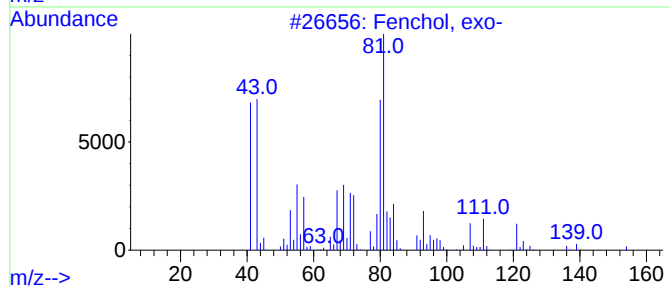
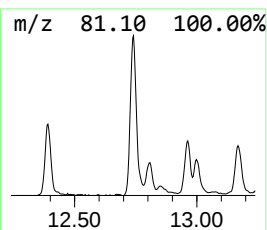
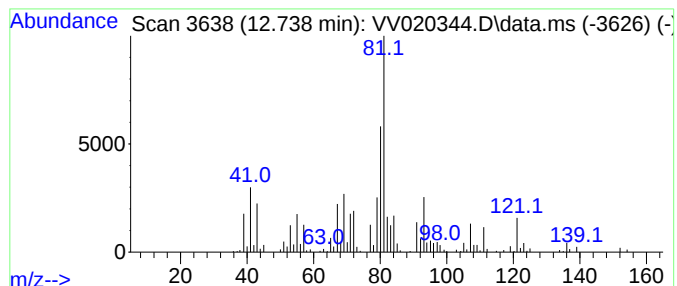
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Quant Title : VOC Analysis

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

Peak Number 7 Fenchol, exo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.738	5.57 ug/L	299558	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchol, exo-	154	C10H18O	022627-95-8	98
2			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	96
3			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	95
4			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	002217-02-9	55
5			Cyclohexane, 1,2-dichloro-, trans-	152	C6H10Cl2	000822-86-6	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

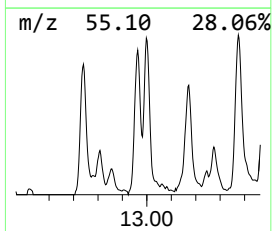
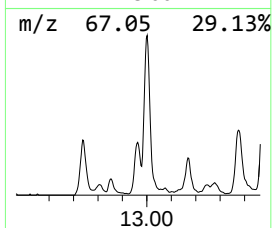
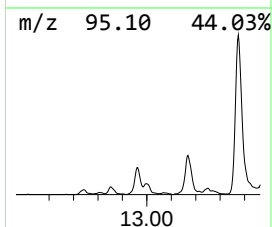
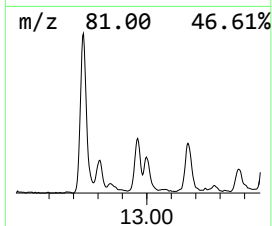
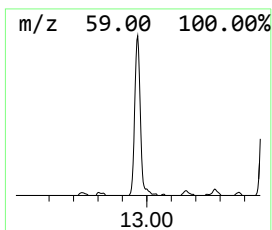
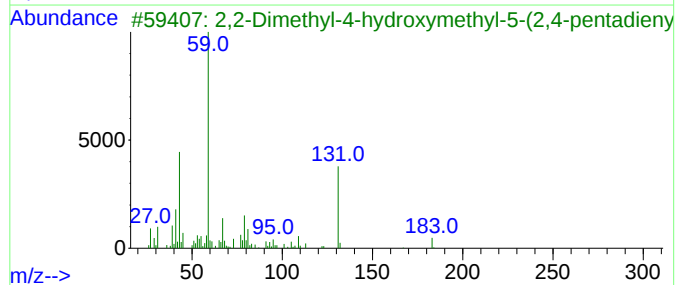
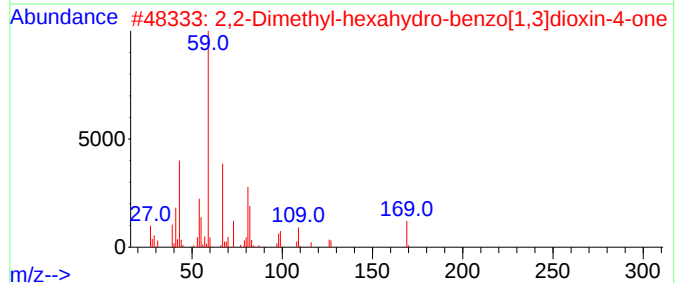
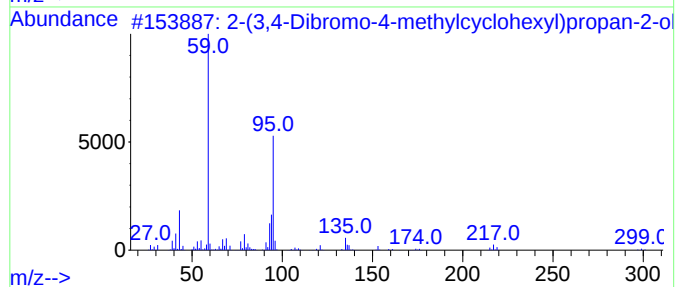
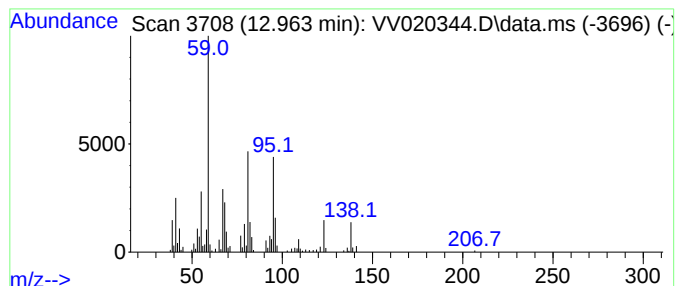
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Quant Title : VOC Analysis

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

Peak Number 8 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	2.98 ug/L	159944	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(3,4-Dibromo-4-methylcyclohexy...	312	C10H18Br2O	110202-13-6	47		
2	2,2-Dimethyl-hexahydro-benzo[1,3...	184	C10H16O3	1000187-24-1	37		
3	2,2-Dimethyl-4-hydroxymethyl-5-(...	198	C11H18O3	1000284-41-5	35		
4	trans-2,7-Dimethyl-4,6-octadien-...	154	C10H18O	1000281-69-5	25		
5	2-Hydroxymethyl-5-(1-hydroxy-1-i...	184	C10H16O3	097762-06-6	25		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

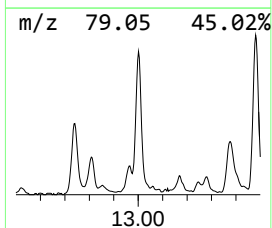
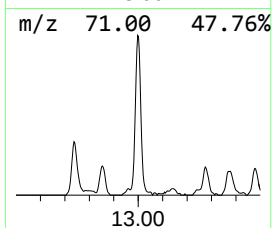
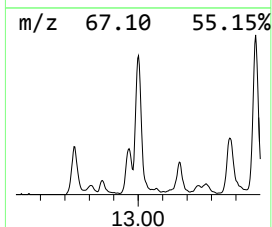
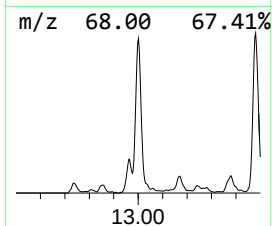
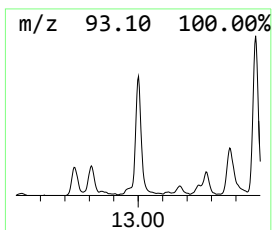
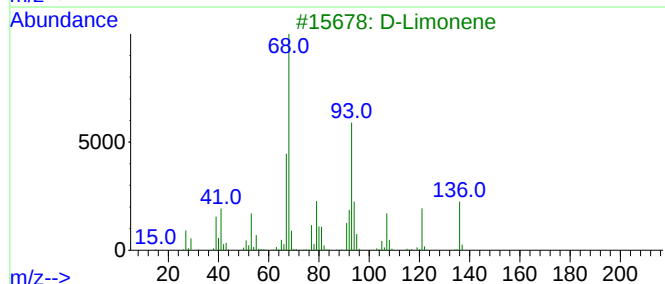
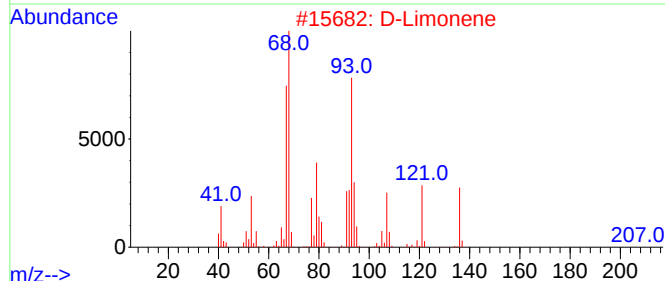
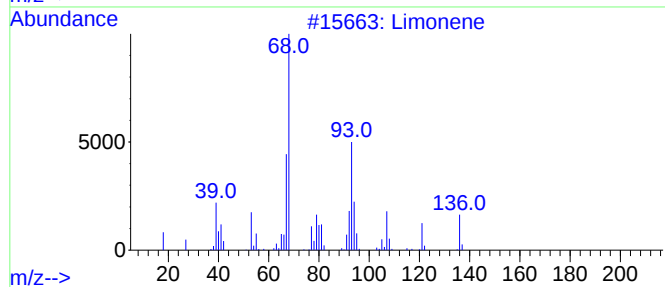
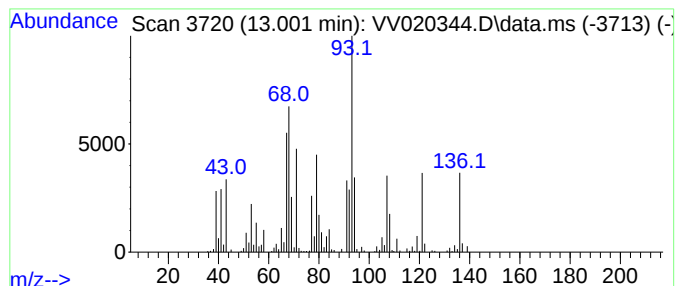
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Quant Title : VOC Analysis

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

Peak Number 9 Limonene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.001	7.74 ug/L	415933	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Limonene	136	C10H16	000138-86-3	62
2			D-Limonene	136	C10H16	005989-27-5	60
3			D-Limonene	136	C10H16	005989-27-5	52
4			Limonene	136	C10H16	000138-86-3	49
5			3-Carene	136	C10H16	013466-78-9	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

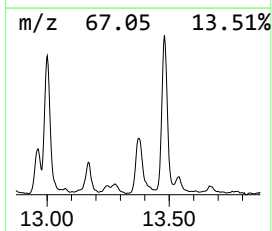
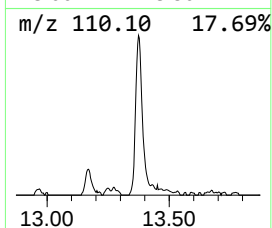
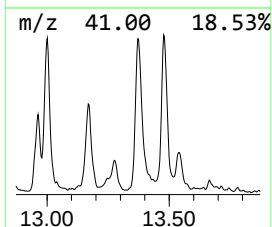
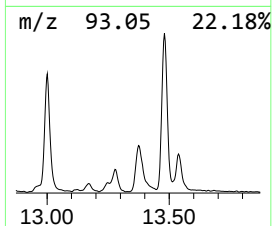
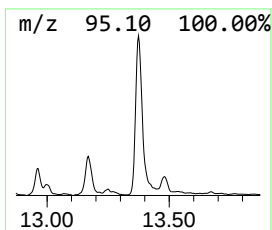
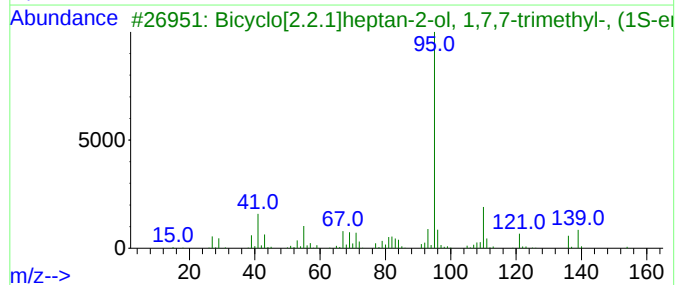
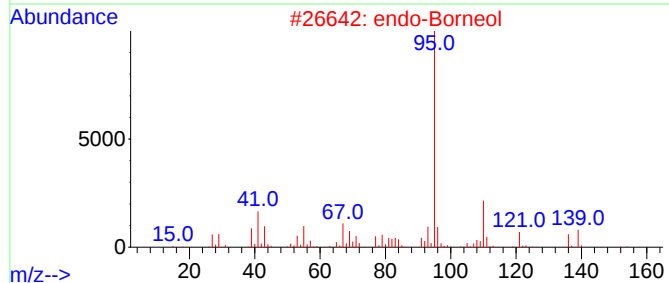
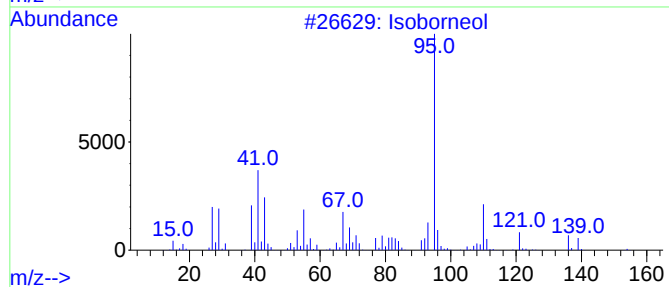
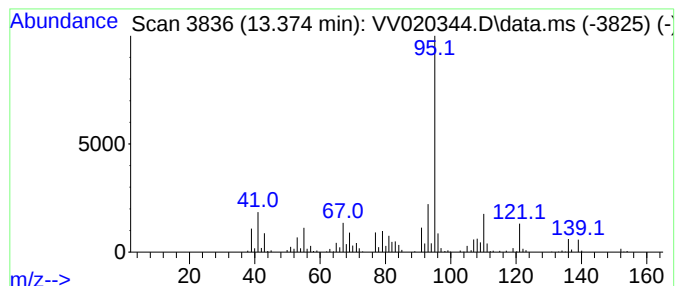
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Quant Title : VOC Analysis

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

Peak Number 10 Isoborneol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.374	7.52 ug/L	404199	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isoborneol	154	C10H18O	000124-76-5	94
2			endo-Borneol	154	C10H18O	000507-70-0	81
3			Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	72
4			(+)-Borneol	154	C10H18O	1000150-76-3	64
5			1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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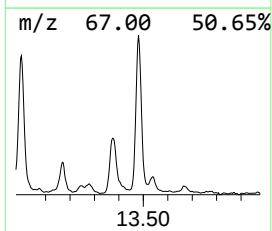
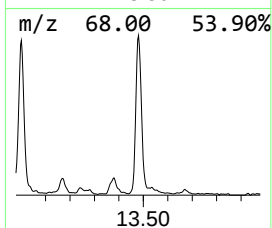
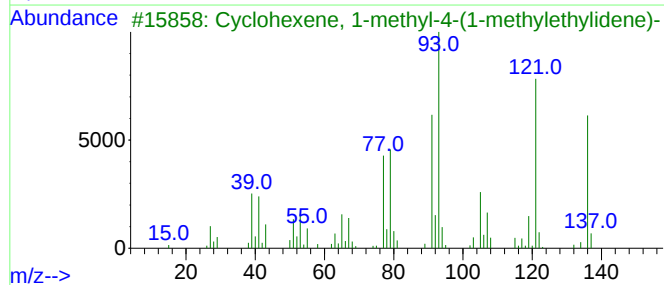
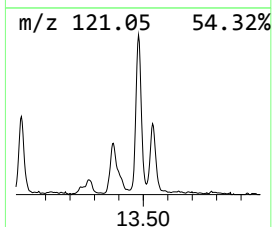
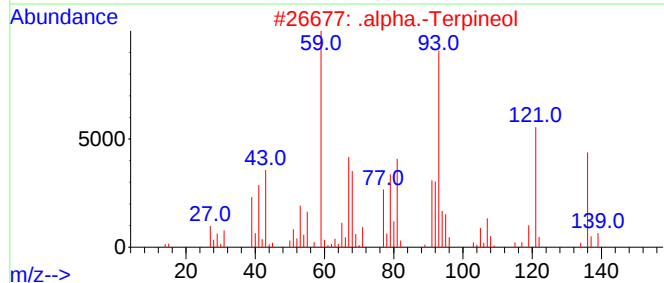
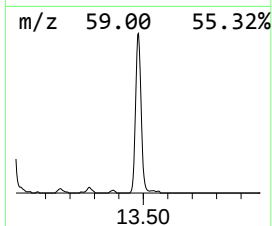
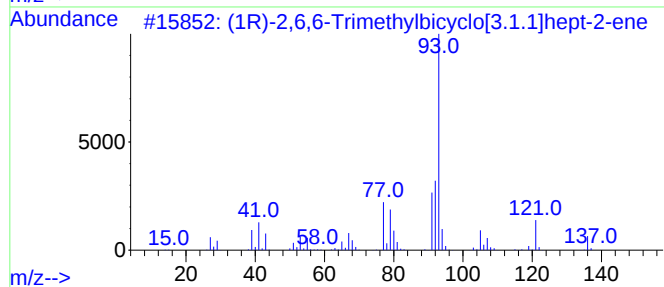
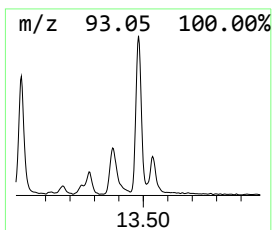
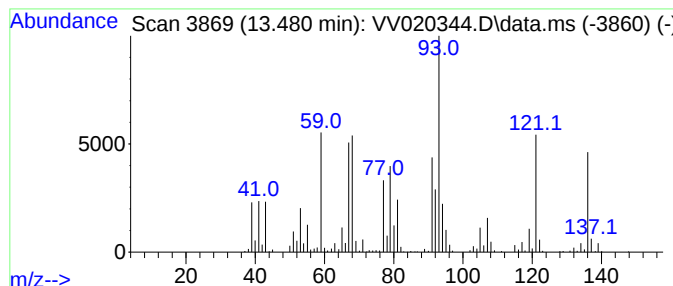
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Quant Title : VOC Analysis

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

Peak Number 11 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.480	9.40 ug/L	505411	1,4-Dichlorobenzene-d4	11.255

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	90		
2	.alpha.-Terpineol	154	C10H18O	000098-55-5	72		
3	Cyclohexene, 1-methyl-4-(1-methylethylidene)-	136	C10H16	000586-62-9	70		
4	(+)-4-Carene	136	C10H16	029050-33-7	64		
5	(+)-3-Carene	136	C10H16	000498-15-7	60		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV021721\
Data File : VV020344.D
Acq On : 17 Feb 2021 14:18
Operator : SY/MD
Sample : M1414-14ME 10X
Misc : 5.03g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBGZ9ME

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	30.4	ug/L	1635440	3	11.255	2687110	50.0
D-Limonene	11.197	19.4	ug/L	1042880	3	11.255	2687110	50.0
Eucalyptol	11.358	21.9	ug/L	1177280	3	11.255	2687110	50.0
Cyclohexene, 1-...	11.905	6.6	ug/L	354558	3	11.255	2687110	50.0
Benzene, 1-meth...	12.095	5.4	ug/L	289103	3	11.255	2687110	50.0
Fenchol, exo-	12.738	5.6	ug/L	299558	3	11.255	2687110	50.0
unknown-01	12.963	3.0	ug/L	159944	3	11.255	2687110	50.0
Limonene	13.001	7.7	ug/L	415933	3	11.255	2687110	50.0
Isoborneol	13.374	7.5	ug/L	404199	3	11.255	2687110	50.0
(1R)-2,6,6-Trim...	13.480	9.4	ug/L	505411	3	11.255	2687110	50.0