

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
 Data File : VW018167.D
 Acq On : 16 Feb 2021 17:03
 Operator : SY/VA
 Sample : M1414-09
 Misc : 5.45G/10ML/MSVOA_W/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 DBGX5

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\SOM2WLM012921S.M

Title : VOC Analysis

Signal : TIC: VW018167.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	69	73	81	rVB	11392	19206	6.64%	1.087%
2	2.891	155	162	172	rVB	14345	32714	11.31%	1.852%
3	3.275	223	225	229	rBV2	148	191	0.07%	0.011%
4	3.489	251	260	261	rBV3	283	642	0.22%	0.036%
5	3.617	279	281	285	rVB	129	160	0.06%	0.009%
6	3.678	288	291	294	rVB3	306	403	0.14%	0.023%
7	3.824	313	315	316	rVV	248	138	0.05%	0.008%
8	3.867	319	322	323	rVV	133	150	0.05%	0.008%
9	3.885	323	325	327	rVV	204	165	0.06%	0.009%
10	3.903	327	328	333	rVV2	118	171	0.06%	0.010%
11	4.019	337	347	358	rVV	31734	86452	29.88%	4.893%
12	4.129	358	365	371	rVV2	1431	3556	1.23%	0.201%
13	4.220	377	380	383	rVB2	146	189	0.07%	0.011%
14	4.452	416	418	423	rVB2	131	150	0.05%	0.008%
15	4.629	446	447	450	rVB	195	140	0.05%	0.008%
16	4.915	487	494	504	rVB3	3423	9316	3.22%	0.527%
17	5.318	558	560	563	rBB2	151	155	0.05%	0.009%
18	5.440	578	580	585	rVV	178	270	0.09%	0.015%
19	5.489	585	588	592	rVV2	117	209	0.07%	0.012%
20	5.940	657	662	669	rVB6	922	2254	0.78%	0.128%
21	6.074	682	684	685	rBV2	150	145	0.05%	0.008%
22	6.208	699	706	708	rBV2	165	311	0.11%	0.018%
23	6.598	769	770	774	rVB	184	206	0.07%	0.012%
24	6.635	774	776	779	rBV2	119	146	0.05%	0.008%
25	6.732	789	792	793	rBV	157	169	0.06%	0.010%
26	6.872	812	815	816	rVB2	160	139	0.05%	0.008%
27	6.897	816	819	821	rBV2	192	268	0.09%	0.015%
28	7.086	840	850	867	rBV3	8880	32959	11.39%	1.865%
29	7.470	909	913	915	rVV2	129	147	0.05%	0.008%
30	7.647	932	942	953	rBV	60463	149679	51.73%	8.472%
31	7.756	959	960	964	rVB2	222	149	0.05%	0.008%
32	7.793	964	966	967	rBV	184	155	0.05%	0.009%
33	7.951	985	992	997	rBV6	1119	2878	0.99%	0.163%
34	8.019	1000	1003	1004	rBV2	495	543	0.19%	0.031%
35	8.110	1016	1018	1021	rBV	247	184	0.06%	0.010%
36	8.275	1036	1045	1061	rBV2	96236	289321	100.00%	16.376%

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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\SOM2WLM012921S.M
 Title : VOC Analysis

37	8.397	1064	1065	1068	rVB	276	157	0.05%	0.009%
38	8.482	1070	1079	1086	rVV2	3697	9116	3.15%	0.516%
39	8.701	1112	1115	1117	rBV2	162	152	0.05%	0.009%
40	8.842	1130	1138	1147	rBV	94325	186960	64.62%	10.582%
41	9.031	1166	1169	1172	rBB2	175	149	0.05%	0.008%
42	9.274	1200	1209	1225	rVB	80100	162174	56.05%	9.179%
43	9.439	1232	1236	1237	rBV	492	470	0.16%	0.027%
44	9.579	1257	1259	1264	rBV2	122	163	0.06%	0.009%
45	9.701	1277	1279	1282	rBV2	142	198	0.07%	0.011%
46	9.762	1284	1289	1292	rVV	730	1242	0.43%	0.070%
47	9.793	1292	1294	1298	rVB	161	164	0.06%	0.009%
48	9.841	1298	1302	1305	rBV3	282	450	0.16%	0.025%
49	9.902	1307	1312	1314	rVB3	627	1063	0.37%	0.060%
50	10.036	1326	1334	1343	rBV2	21490	38132	13.18%	2.158%
51	10.177	1353	1357	1360	rBV2	103	197	0.07%	0.011%
52	10.256	1366	1370	1371	rBV3	149	196	0.07%	0.011%
53	10.323	1373	1381	1388	rBV	99664	171791	59.38%	9.723%
54	10.500	1406	1410	1411	rBV2	136	147	0.05%	0.008%
55	10.579	1417	1423	1429	rBV2	7887	14400	4.98%	0.815%
56	10.707	1439	1444	1450	rVV3	1201	2162	0.75%	0.122%
57	10.927	1474	1480	1488	rBV	32562	56076	19.38%	3.174%
58	11.372	1549	1553	1556	rBV2	117	179	0.06%	0.010%
59	11.463	1567	1568	1572	rVB2	270	280	0.10%	0.016%
60	11.518	1575	1577	1581	rVB2	195	188	0.06%	0.011%
61	11.628	1589	1595	1607	rVB	64489	111540	38.55%	6.313%
62	11.719	1607	1610	1611	rBV2	187	192	0.07%	0.011%
63	11.737	1611	1613	1616	rBV2	140	170	0.06%	0.010%
64	11.817	1622	1626	1628	rBV	811	1018	0.35%	0.058%
65	11.926	1643	1644	1648	rVB3	317	261	0.09%	0.015%
66	12.054	1663	1665	1669	rVB2	217	217	0.08%	0.012%
67	12.109	1672	1674	1676	rBV	216	188	0.06%	0.011%
68	12.158	1681	1682	1684	rBV	170	166	0.06%	0.009%
69	12.256	1695	1698	1702	rBV2	141	151	0.05%	0.009%
70	12.365	1714	1716	1721	rVB	467	319	0.11%	0.018%
71	12.463	1729	1732	1736	rVB	453	575	0.20%	0.033%
72	12.536	1738	1744	1749	rBV5	1100	1794	0.62%	0.102%
73	12.609	1750	1756	1763	rVB2	6530	10648	3.68%	0.603%
74	12.688	1763	1769	1779	rVB	63386	112208	38.78%	6.351%
75	12.774	1779	1783	1787	rBV4	1588	2640	0.91%	0.149%
76	12.804	1787	1788	1792	rBV2	245	224	0.08%	0.013%

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Title : VOC Analysis

77	12.951	1801	1812	1820	rBV2	13687	26533	9.17%	1.502%
78	13.048	1820	1828	1838	rBV2	40540	71779	24.81%	4.063%
79	13.152	1844	1845	1847	rBV	275	267	0.09%	0.015%
80	13.201	1848	1853	1858	rVB3	3118	5524	1.91%	0.313%
81	13.249	1858	1861	1863	rBV2	162	253	0.09%	0.014%
82	13.353	1877	1878	1880	rBV	237	149	0.05%	0.008%
83	13.469	1892	1897	1904	rBV4	4022	7342	2.54%	0.416%
84	13.560	1906	1912	1916	rBV	17145	27345	9.45%	1.548%
85	13.761	1941	1945	1948	rBV3	820	1166	0.40%	0.066%
86	13.804	1948	1952	1953	rBV4	784	900	0.31%	0.051%
87	13.853	1954	1960	1966	rVV2	27680	45817	15.84%	2.593%
88	13.993	1979	1983	1985	rBV3	947	1516	0.52%	0.086%
89	14.066	1991	1995	2001	rVB3	2763	4888	1.69%	0.277%
90	14.200	2012	2017	2024	rVB5	1752	3863	1.34%	0.219%
91	14.341	2035	2040	2052	rBV6	4448	10759	3.72%	0.609%
92	14.493	2063	2065	2066	rBV2	611	515	0.18%	0.029%
93	14.530	2067	2071	2073	rVV3	2222	2874	0.99%	0.163%
94	14.572	2073	2078	2082	rVB	8703	12353	4.27%	0.699%
95	14.651	2086	2091	2096	rBV5	4968	9897	3.42%	0.560%
96	14.853	2123	2124	2127	rVB3	374	204	0.07%	0.012%
97	14.926	2132	2136	2141	rVB3	2238	3618	1.25%	0.205%
98	15.005	2147	2149	2152	rBV2	380	421	0.15%	0.024%
99	15.487	2224	2228	2234	rVB	3276	5900	2.04%	0.334%
100	15.535	2234	2236	2239	rBV2	430	361	0.12%	0.020%

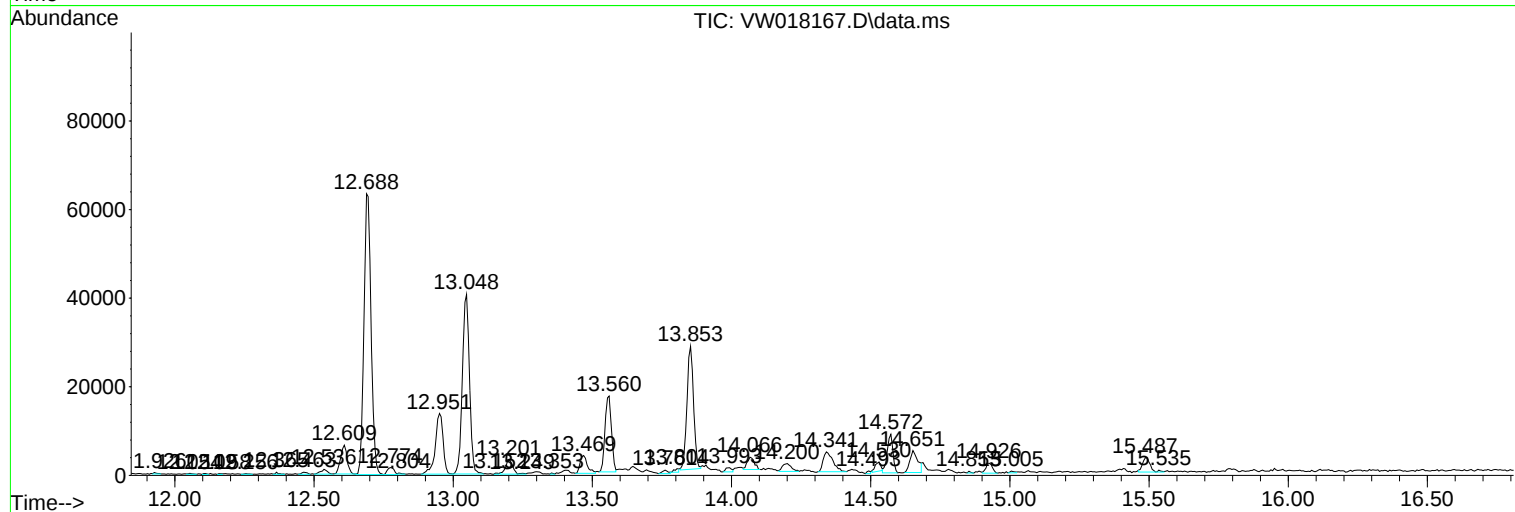
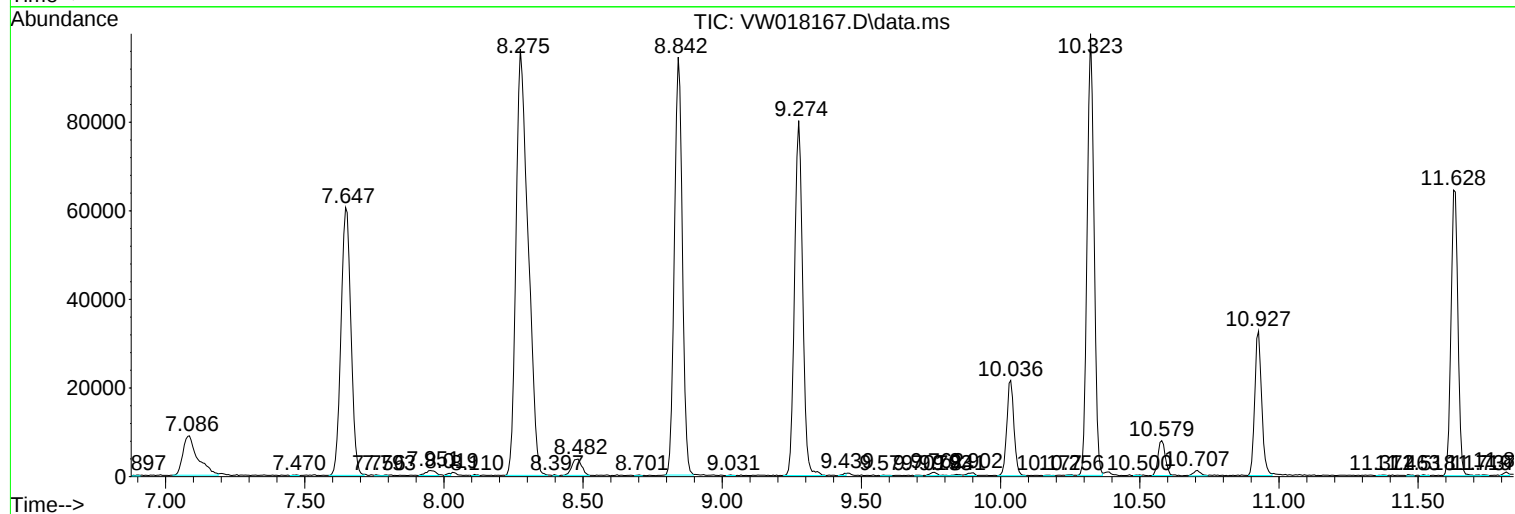
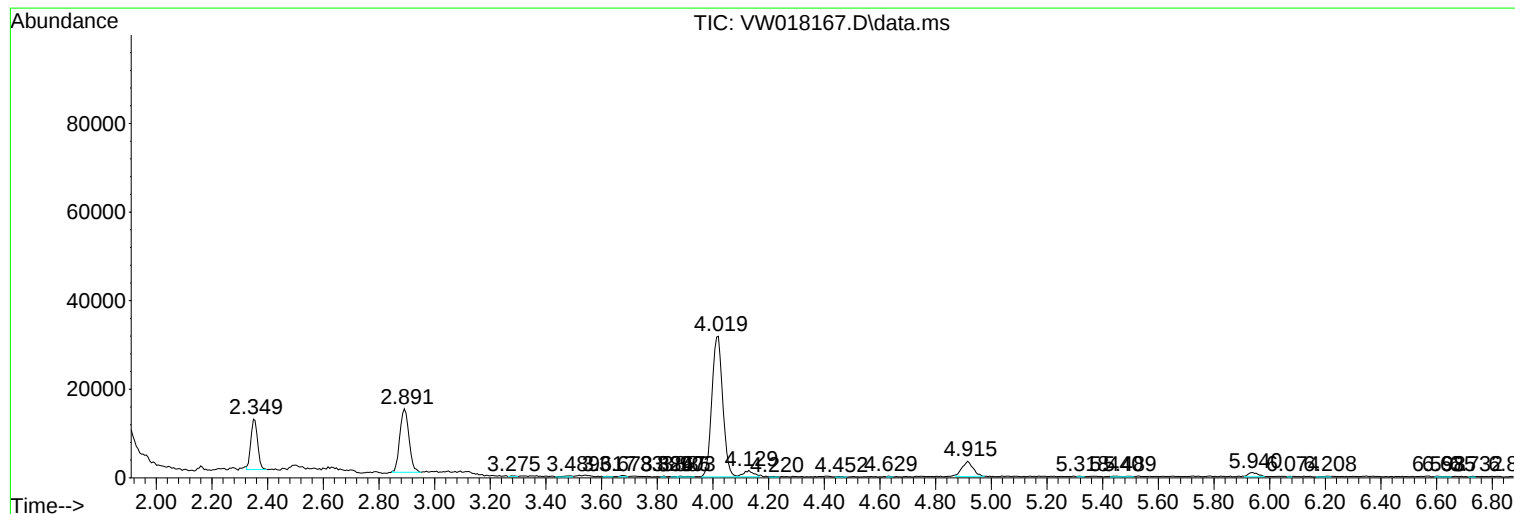
Sum of corrected areas: 1766791

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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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



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Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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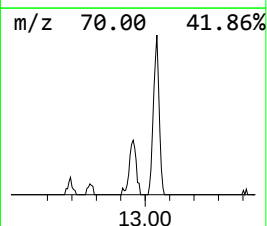
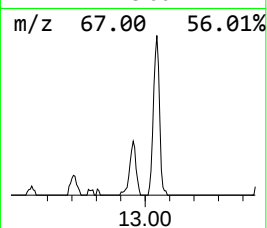
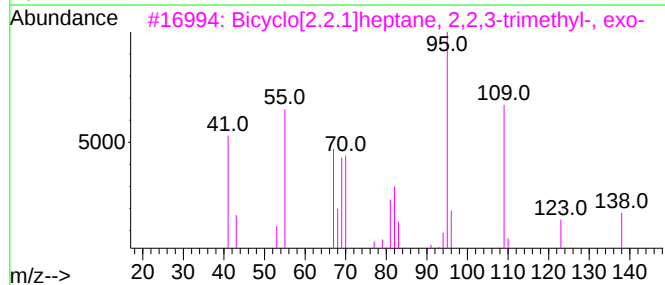
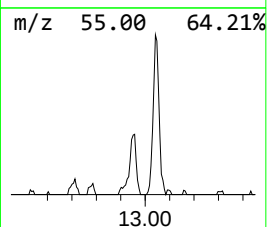
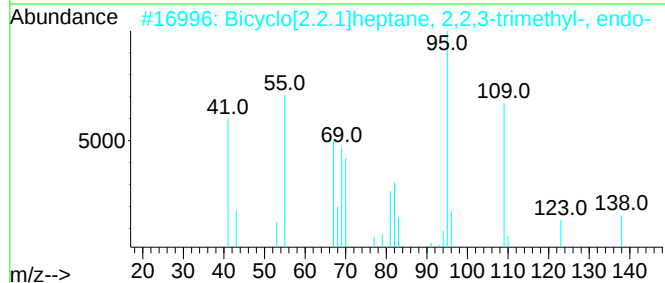
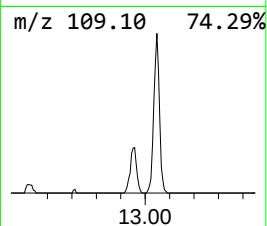
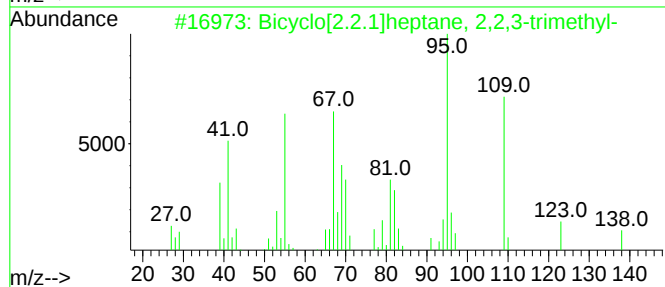
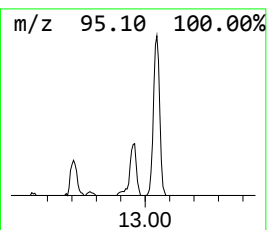
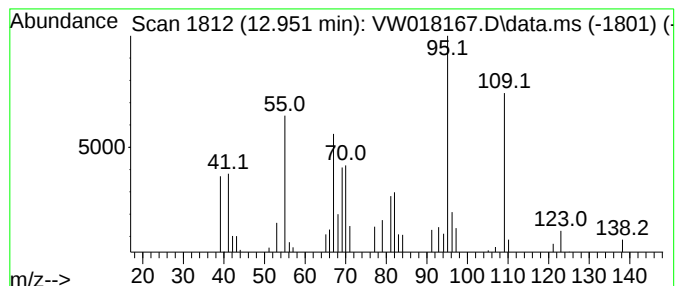
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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

Peak Number 4 Bicyclo[2.2.1]heptane, 2,2,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.951	24.26 ug/L	26533	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	95
2			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	87
3			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-41-8	80
4			Cyclopentane, (3-methylbutylidene)-	138	C10H18	053366-51-1	38
5			Bicyclo[2.2.1]heptane, 1,7,7-tri...	138	C10H18	000464-15-3	35



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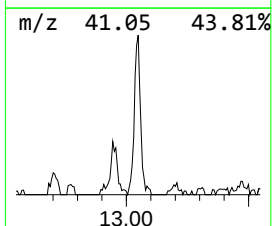
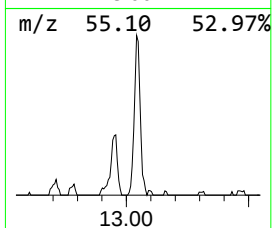
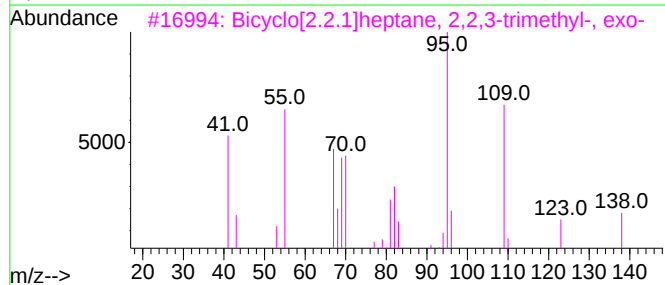
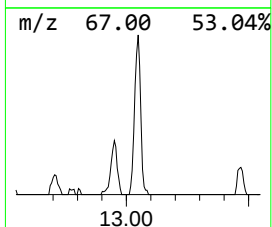
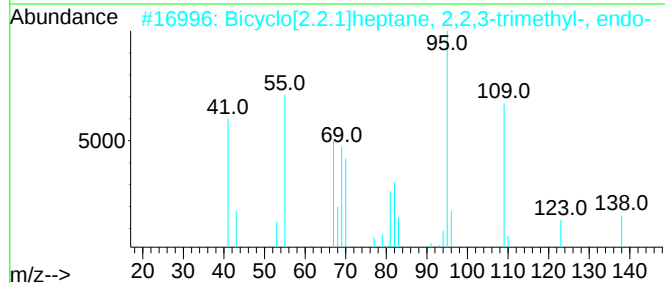
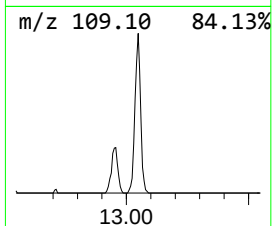
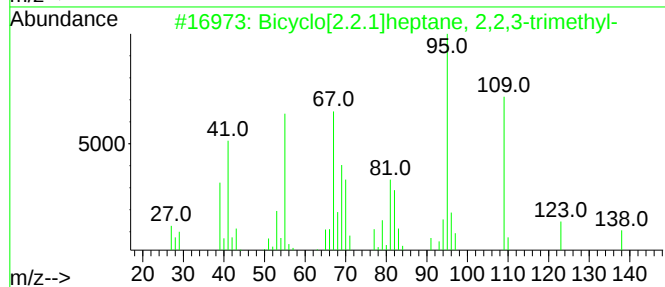
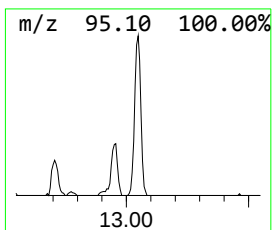
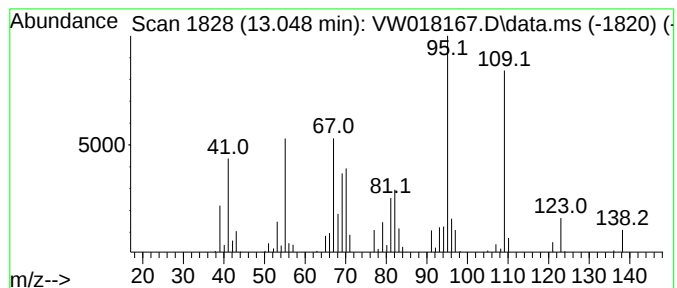
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Bicyclo[2.2.1]heptane, 2,2,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.048	65.62 ug/L	71779	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	96
2			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	93
3			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-41-8	53
4			Bicyclo[2.2.1]heptane, 1,7,7-tri...	138	C10H18	000464-15-3	43
5			Cyclopentene, 1,4-dimethyl-5-(1-...	138	C10H18	061142-33-4	38



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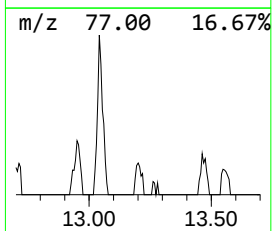
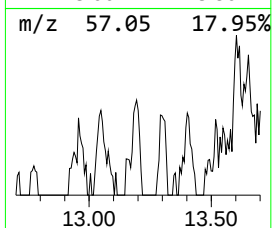
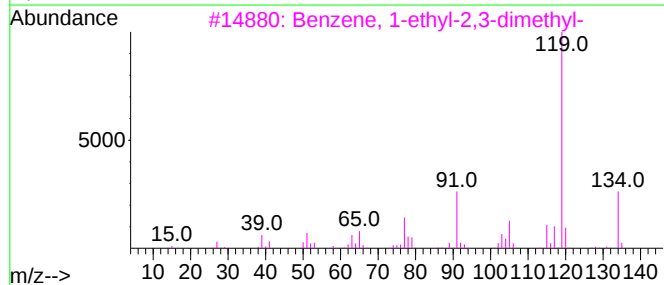
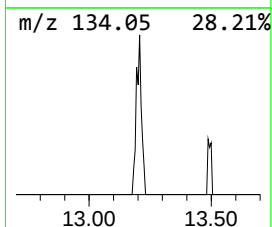
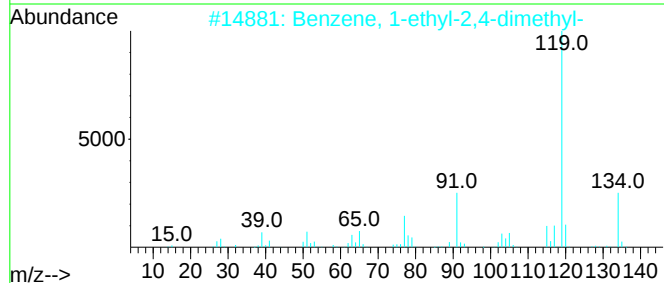
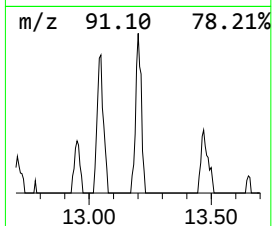
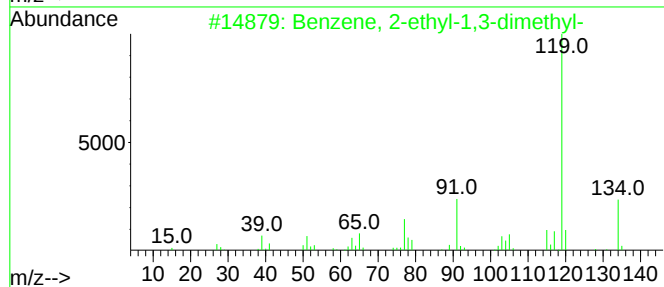
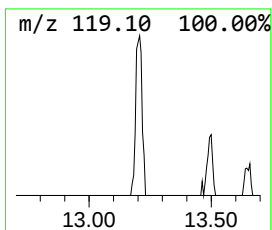
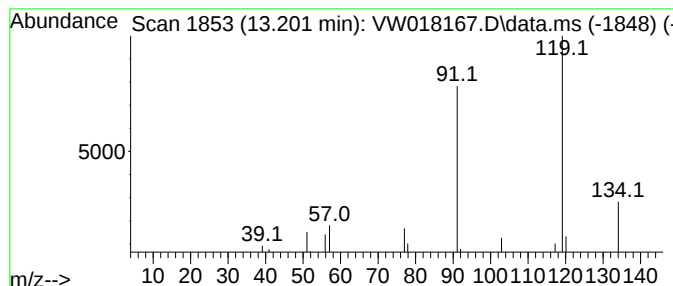
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.201	5.05 ug/L	5524	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	72
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	72
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	72
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

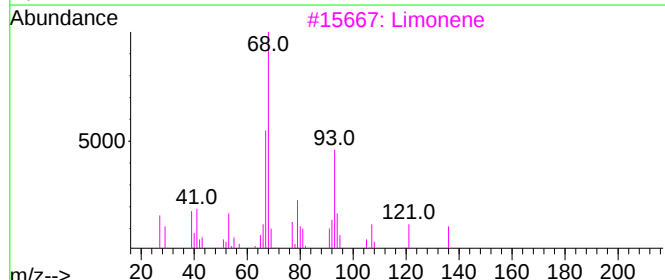
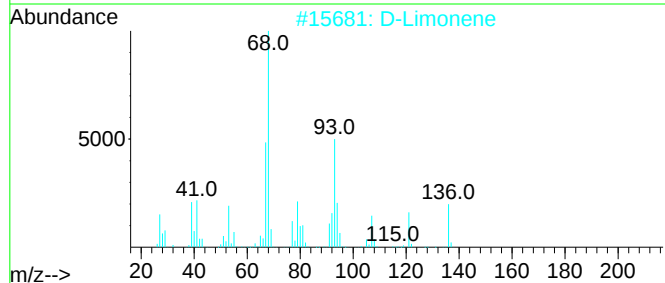
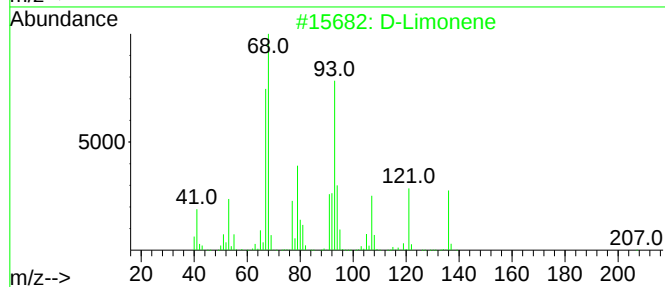
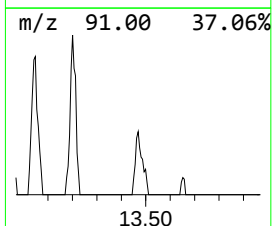
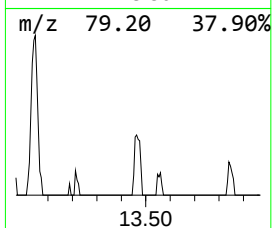
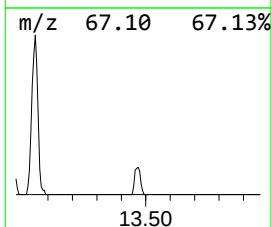
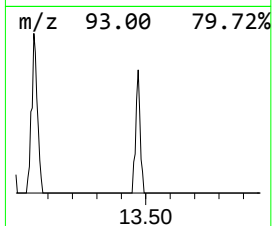
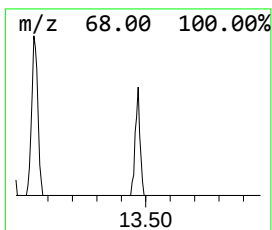
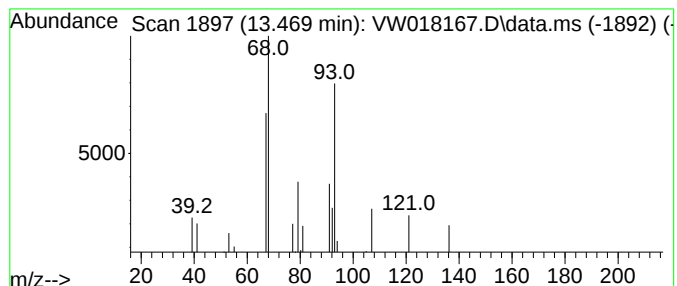
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Quant Title : VOC Analysis

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

Peak Number 7 D-Limonene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.469	6.71 ug/L	7342	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Limonene			136	C10H16	005989-27-5	91
2	D-Limonene			136	C10H16	005989-27-5	90
3	Limonene			136	C10H16	000138-86-3	87
4	D-Limonene			136	C10H16	005989-27-5	86
5	D-Limonene			136	C10H16	005989-27-5	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

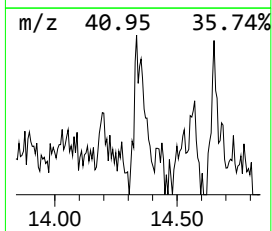
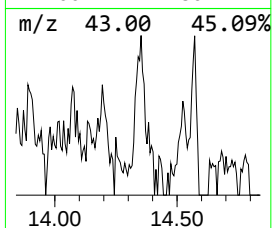
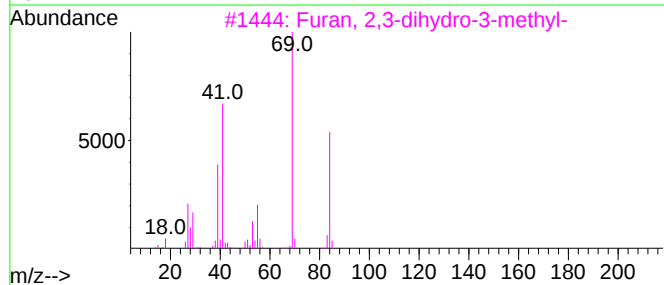
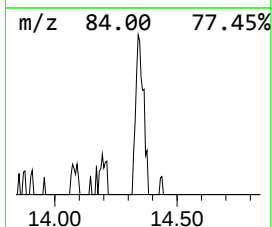
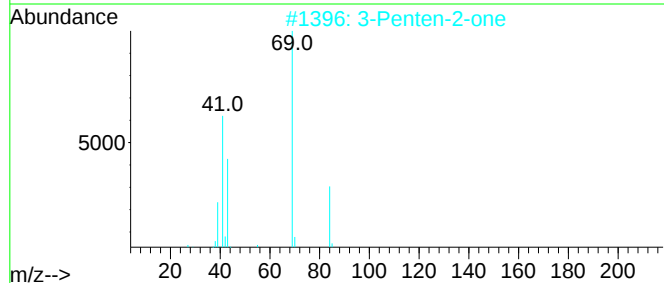
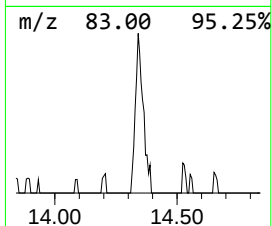
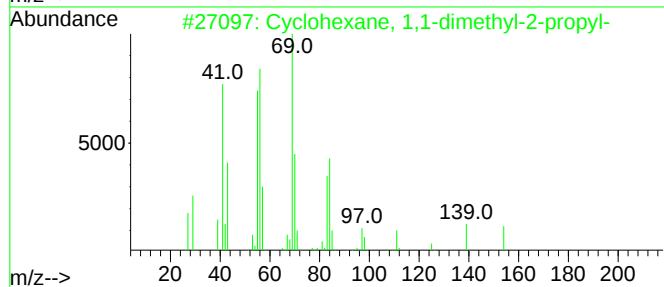
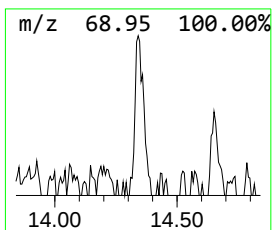
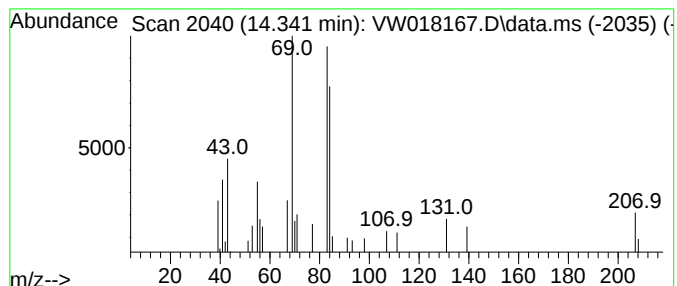
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Quant Title : VOC Analysis

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

Peak Number 11 (DEL) Alkane: Cyclic14.341 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.341	9.84 ug/L	10759	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1-dimethyl-2-propyl-	154	C11H22	081983-71-3	38
2			3-Penten-2-one	84	C5H8O	000625-33-2	38
3			Furan, 2,3-dihydro-3-methyl-	84	C5H8O	001708-27-6	38
4			3-Ethyl-2-methyl-1-heptene	140	C10H20	019780-60-0	32
5			2-Azetidinone, 3,3,4,4-tetramethyl-	127	C7H13NO	013423-22-8	23



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

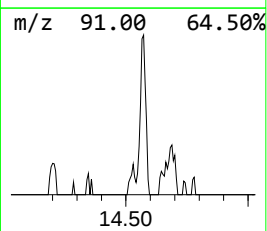
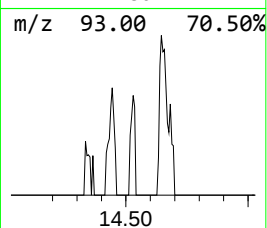
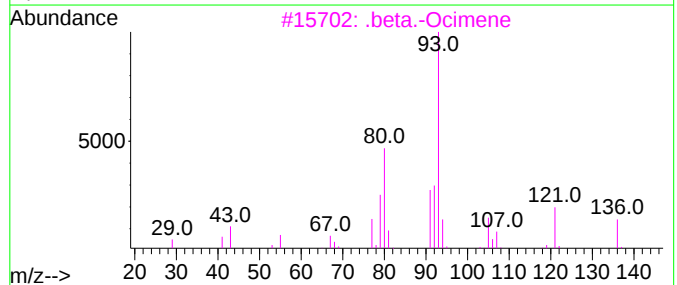
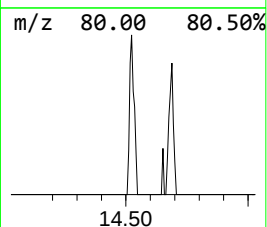
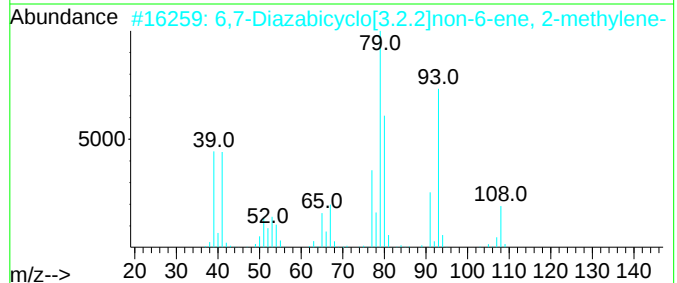
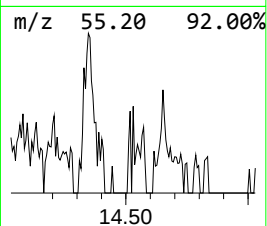
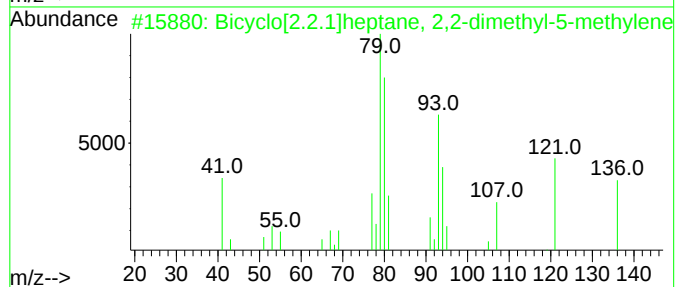
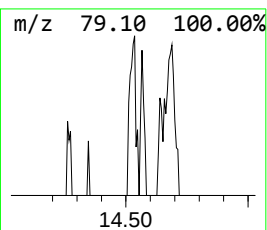
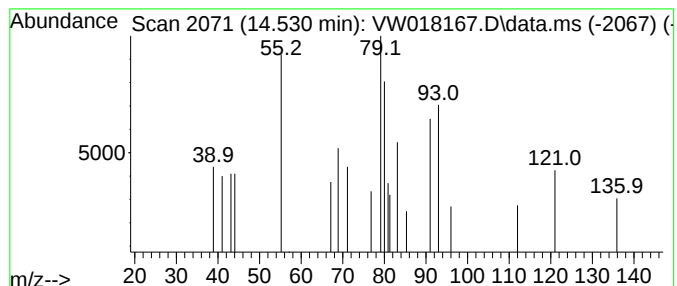
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Quant Title : VOC Analysis

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

Peak Number 12 6,7-Diazabicyclo[3.2.2]non-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.530	2.63 ug/L	2874	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	000497-32-5	58
2			6,7-Diazabicyclo[3.2.2]non-6-ene...	136	C8H12N2	097423-85-3	50
3			.beta.-Ocimene	136	C10H16	013877-91-3	47
4			3-Penten-1-yne, 3-methyl-	80	C6H8	001574-33-0	47
5			Tricyclo[4.2.1.0(2,5)]nonane	122	C9H14	1000195-06-2	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

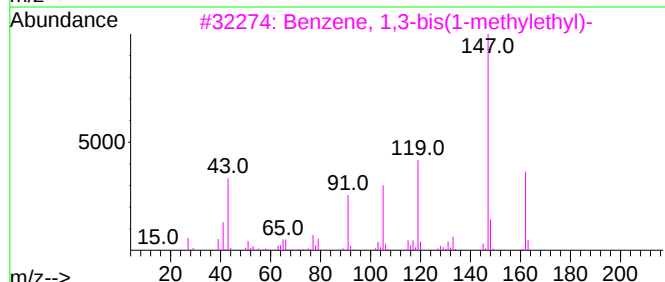
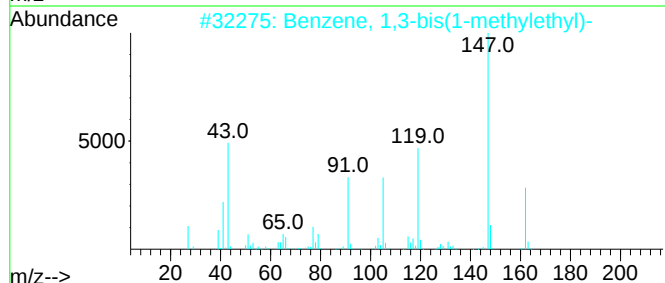
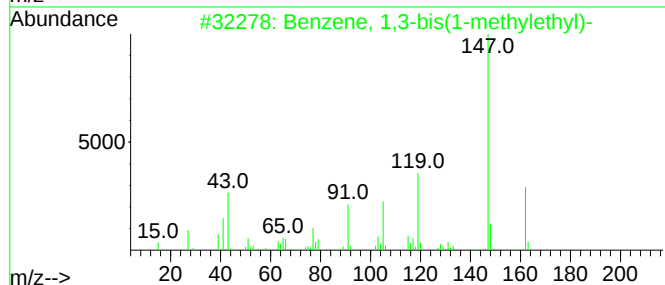
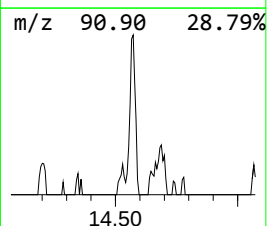
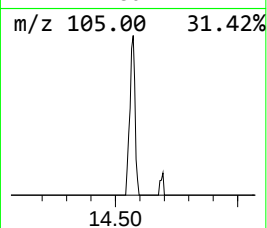
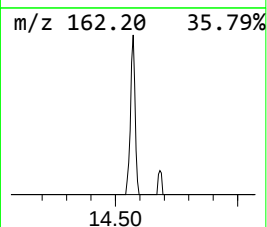
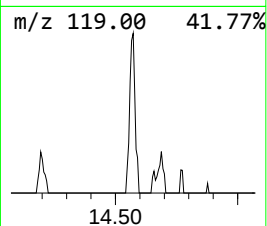
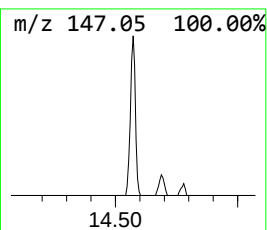
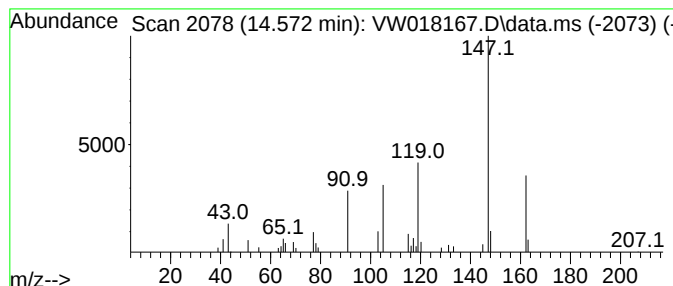
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Quant Title : VOC Analysis

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

Peak Number 13 Benzene, 1,3-bis(1-methylet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.572	11.29 ug/L	12353	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	93
2			Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	91
3			Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	90
4			Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	87
5			Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

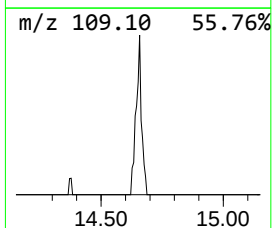
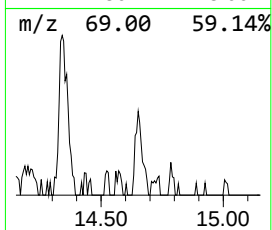
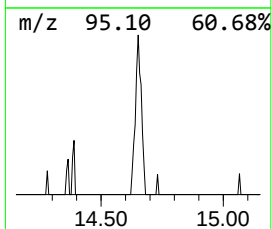
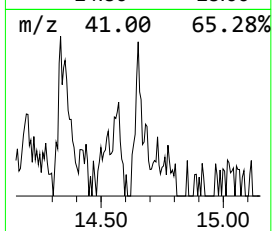
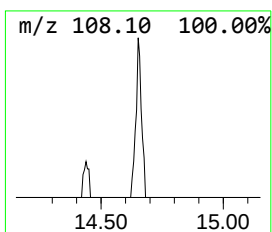
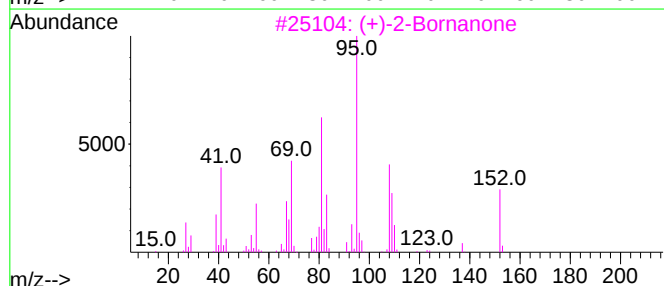
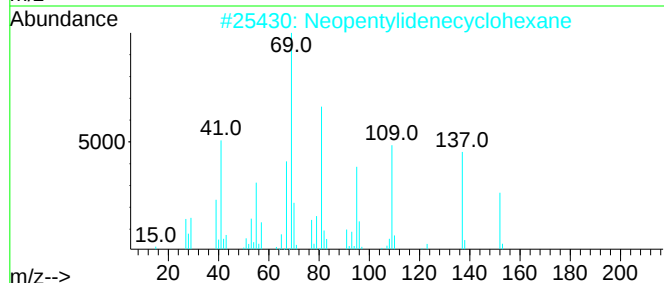
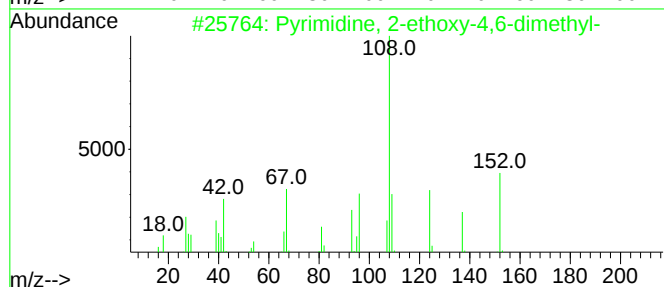
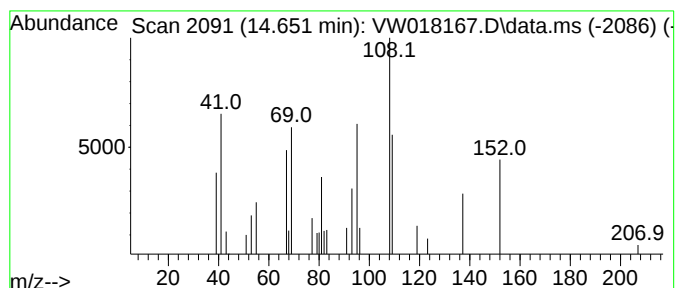
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Quant Title : VOC Analysis

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

Peak Number 14 Pyrimidine, 2-ethoxy-4,6-di... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.652	9.05 ug/L	9897	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrimidine, 2-ethoxy-4,6-dimethyl-	152	C8H12N2O	007781-21-7	59
2			Neopentylidenecyclohexane	152	C11H20	039546-80-0	43
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	38
4			Camphor	152	C10H16O	000076-22-2	35
5			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

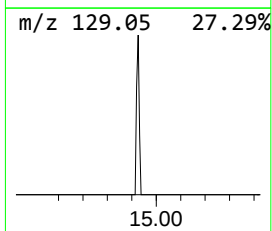
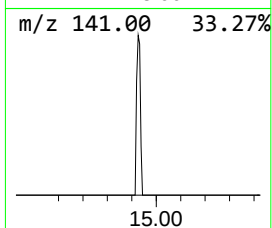
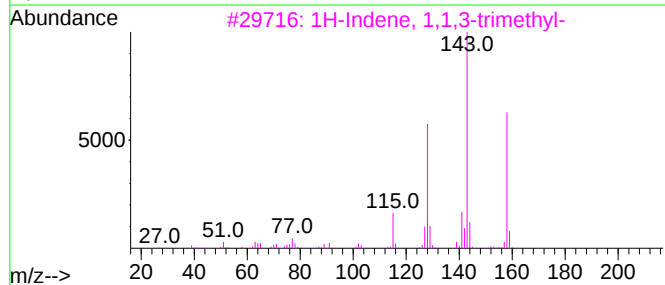
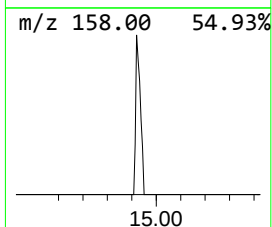
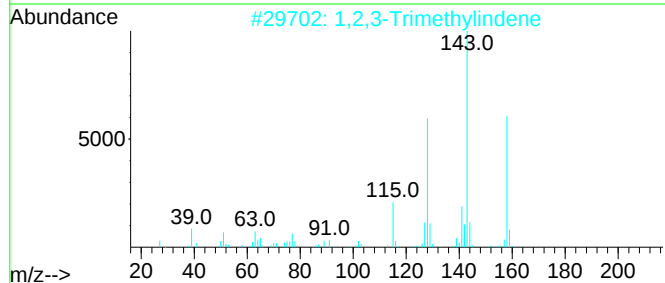
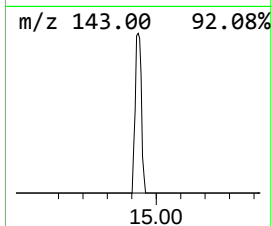
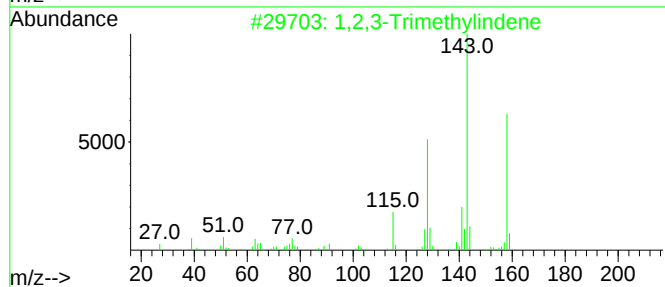
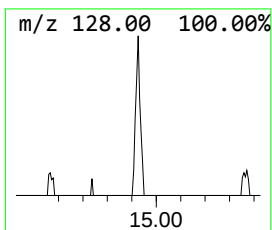
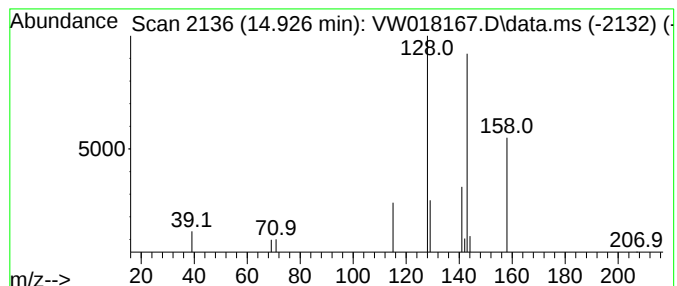
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Quant Title : VOC Analysis

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

Peak Number 15 1,2,3-Trimethylindene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.926	3.31 ug/L	3618	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,3-Trimethylindene	158	C12H14	004773-83-5	59
2			1,2,3-Trimethylindene	158	C12H14	004773-83-5	59
3			1H-Indene, 1,1,3-trimethyl-	158	C12H14	002177-45-9	59
4			(1-Methylpenta-1,3-dienyl)benzene	158	C12H14	116669-49-9	59
5			Benzene, (1,2,3-trimethyl-2-cycl...	158	C12H14	006393-13-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021621\
Data File : VW018167.D
Acq On : 16 Feb 2021 17:03
Operator : SY/VA
Sample : M1414-09
Misc : 5.45G/10ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBGX5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.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
Bicyclo[2.2.1]h...	12.951	24.3	ug/L	26533	3	13.560	27345	25.0
Bicyclo[2.2.1]h...	13.048	65.6	ug/L	71779	3	13.560	27345	25.0
Benzene, 2-ethy...	13.201	5.0	ug/L	5524	3	13.560	27345	25.0
D-Limonene	13.469	6.7	ug/L	7342	3	13.560	27345	25.0
(DEL) Alkane: C...	14.341	9.8	ug/L	10759	3	13.560	27345	25.0
6,7-Diazabicycl...	14.530	2.6	ug/L	2874	3	13.560	27345	25.0
Benzene, 1,3-bi...	14.572	11.3	ug/L	12353	3	13.560	27345	25.0
Pyrimidine, 2-e...	14.652	9.1	ug/L	9897	3	13.560	27345	25.0
1,2,3-Trimethyl...	14.926	3.3	ug/L	3618	3	13.560	27345	25.0