

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021721\
 Data File : VW018191.D
 Acq On : 17 Feb 2021 17:00
 Operator : SY/VA
 Sample : M1371-16
 Misc : 3.93G/10ML/MSVOA_W/SOIL
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 DBH30

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: VW018191.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.355	66	74	88	rVB	70639	141576	11.51%	1.291%
2	2.775	141	143	150	rBV5	496	898	0.07%	0.008%
3	2.897	155	163	178	rVV	50796	131429	10.69%	1.199%
4	3.348	236	237	240	rVB	311	199	0.02%	0.002%
5	3.464	253	256	259	rBV3	239	343	0.03%	0.003%
6	3.531	260	267	275	rVV6	988	2323	0.19%	0.021%
7	3.592	275	277	278	rVV	219	187	0.02%	0.002%
8	3.854	318	320	324	rVB	182	185	0.02%	0.002%
9	4.025	334	348	360	rBV	129247	415919	33.82%	3.794%
10	4.251	383	385	388	rBV2	180	208	0.02%	0.002%
11	4.336	397	399	400	rVB	317	211	0.02%	0.002%
12	4.391	400	408	416	rBV3	1804	5655	0.46%	0.052%
13	4.726	460	463	465	rBV2	178	196	0.02%	0.002%
14	4.921	484	495	512	rBV2	11789	43505	3.54%	0.397%
15	5.055	514	517	520	rBV2	152	214	0.02%	0.002%
16	5.122	520	528	529	rBV3	680	1144	0.09%	0.010%
17	5.141	529	531	533	rBV3	681	832	0.07%	0.008%
18	5.427	574	578	579	rBV2	434	570	0.05%	0.005%
19	5.580	601	603	606	rVB	223	211	0.02%	0.002%
20	5.939	655	662	672	rVB7	1876	5499	0.45%	0.050%
21	6.201	703	705	707	rVV2	163	195	0.02%	0.002%
22	6.226	707	709	713	rVB2	272	220	0.02%	0.002%
23	6.384	727	735	748	rBV2	8466	26325	2.14%	0.240%
24	6.543	759	761	765	rBV2	203	220	0.02%	0.002%
25	6.781	797	800	801	rBV2	171	193	0.02%	0.002%
26	6.884	813	817	818	rBV2	410	426	0.03%	0.004%
27	6.903	818	820	822	rVV2	413	431	0.04%	0.004%
28	7.073	840	848	853	rBV	31273	79520	6.47%	0.725%
29	7.646	931	942	957	rBV	228950	542442	44.11%	4.948%
30	7.762	959	961	963	rVB	359	307	0.02%	0.003%
31	7.878	975	980	984	rVV3	445	698	0.06%	0.006%
32	7.951	985	992	998	rVB3	1185	2407	0.20%	0.022%
33	8.061	1006	1010	1011	rBV	434	351	0.03%	0.003%
34	8.274	1034	1045	1061	rBV2	427658	1229703	100.00%	11.218%
35	8.396	1062	1065	1069	rVB5	948	1484	0.12%	0.014%
36	8.695	1109	1114	1115	rBV3	459	711	0.06%	0.006%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Max Peaks: 100

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Peak Location: TOP

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

37	8.841	1129	1138	1151	rBV	435146	860201	69.95%	7.847%
38	9.024	1159	1168	1177	rBV	59935	126840	10.31%	1.157%
39	9.274	1200	1209	1224	rBV	276518	560296	45.56%	5.111%
40	9.414	1229	1232	1239	rVB3	1287	2648	0.22%	0.024%
41	9.500	1239	1246	1257	rBV	70117	126662	10.30%	1.155%
42	9.658	1267	1272	1279	rVB5	1543	3314	0.27%	0.030%
43	9.737	1282	1285	1286	rBV2	337	244	0.02%	0.002%
44	9.780	1288	1292	1294	rVB2	242	332	0.03%	0.003%
45	9.890	1307	1310	1311	rBV2	159	186	0.02%	0.002%
46	9.926	1314	1316	1320	rVB2	403	449	0.04%	0.004%
47	10.036	1326	1334	1344	rBV	151654	273468	22.24%	2.495%
48	10.115	1344	1347	1352	rBV5	579	1078	0.09%	0.010%
49	10.213	1357	1363	1372	rBV2	5797	10730	0.87%	0.098%
50	10.323	1372	1381	1389	rBV	649123	1126071	91.57%	10.272%
51	10.579	1415	1423	1433	rBV	90556	160726	13.07%	1.466%
52	10.707	1438	1444	1448	rBV	1949	3558	0.29%	0.032%
53	10.786	1454	1457	1459	rBV3	349	334	0.03%	0.003%
54	10.859	1462	1469	1473	rVB7	1405	3187	0.26%	0.029%
55	10.926	1473	1480	1485	rBV	120292	205895	16.74%	1.878%
56	10.975	1485	1488	1495	rVB	12869	21276	1.73%	0.194%
57	11.066	1496	1503	1517	rBV	178238	299731	24.37%	2.734%
58	11.274	1536	1537	1539	rVB	382	190	0.02%	0.002%
59	11.408	1557	1559	1563	rVB3	328	463	0.04%	0.004%
60	11.469	1563	1569	1573	rBV3	1957	3376	0.27%	0.031%
61	11.572	1581	1586	1588	rBV3	1413	2227	0.18%	0.020%
62	11.633	1588	1596	1604	rVV	587663	1010532	82.18%	9.218%
63	11.713	1604	1609	1618	rVB2	8481	17384	1.41%	0.159%
64	11.841	1625	1630	1634	rBV2	2978	6253	0.51%	0.057%
65	11.908	1638	1641	1645	rVB3	1067	1638	0.13%	0.015%
66	11.975	1650	1652	1653	rBV2	335	182	0.01%	0.002%
67	12.017	1657	1659	1663	rVB2	521	799	0.06%	0.007%
68	12.085	1664	1670	1676	rBV2	7245	13878	1.13%	0.127%
69	12.158	1676	1682	1685	rBV2	12561	21138	1.72%	0.193%
70	12.237	1692	1695	1706	rVB2	5054	10363	0.84%	0.095%
71	12.341	1707	1712	1722	rVB4	3887	7458	0.61%	0.068%
72	12.444	1723	1729	1738	rBV2	15075	32529	2.65%	0.297%
73	12.554	1742	1747	1752	rVV	15567	28072	2.28%	0.256%
74	12.609	1752	1756	1761	rVB	17536	27119	2.21%	0.247%
75	12.694	1762	1770	1776	rBV	218518	366543	29.81%	3.344%
76	12.840	1792	1794	1795	rBV2	338	244	0.02%	0.002%

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Integration Parameters: LSCINT.P

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Title : VOC Analysis

77	12.889	1795	1802	1805	rBV6	1655	3693	0.30%	0.034%
78	12.969	1805	1815	1820	rBV	44313	71440	5.81%	0.652%
79	13.121	1836	1840	1841	rBV3	3218	3995	0.32%	0.036%
80	13.194	1849	1852	1855	rBV4	1879	2883	0.23%	0.026%
81	13.255	1858	1862	1863	rBV3	2123	2659	0.22%	0.024%
82	13.304	1865	1870	1879	rVB3	19110	39236	3.19%	0.358%
83	13.395	1881	1885	1890	rVB3	9401	15803	1.29%	0.144%
84	13.462	1891	1896	1901	rBV5	12046	24555	2.00%	0.224%
85	13.554	1905	1911	1918	rBV	663180	1058858	86.11%	9.659%
86	13.645	1921	1926	1933	rBV	146894	234406	19.06%	2.138%
87	13.853	1953	1960	1967	rBV	593605	993768	80.81%	9.065%
88	14.072	1991	1996	2005	rVB	20275	38476	3.13%	0.351%
89	14.194	2013	2016	2017	rBV3	2666	3294	0.27%	0.030%
90	14.322	2034	2037	2043	rVB	18618	28576	2.32%	0.261%
91	14.523	2063	2070	2082	rBV	167830	262226	21.32%	2.392%
92	14.627	2083	2087	2091	rBV7	4748	11262	0.92%	0.103%
93	15.169	2172	2176	2178	rBV4	11483	15932	1.30%	0.145%
94	15.206	2179	2182	2185	rVV	20114	34067	2.77%	0.311%
95	15.249	2185	2189	2198	rVB3	32321	61899	5.03%	0.565%
96	15.401	2211	2214	2218	rVB3	8225	11868	0.97%	0.108%
97	16.096	2323	2328	2335	rBV8	5876	13828	1.12%	0.126%
98	16.358	2368	2371	2376	rVB6	3441	5379	0.44%	0.049%
99	16.450	2377	2386	2398	rVB7	13098	41573	3.38%	0.379%
100	16.565	2402	2405	2413	rVB7	4397	7979	0.65%	0.073%

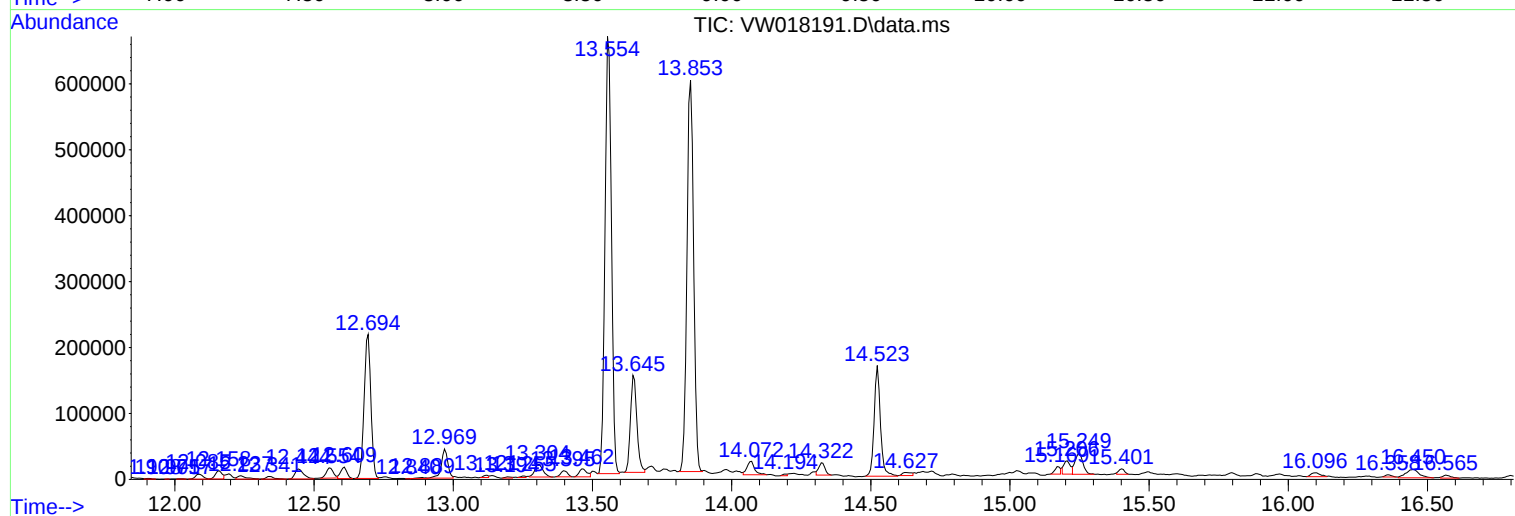
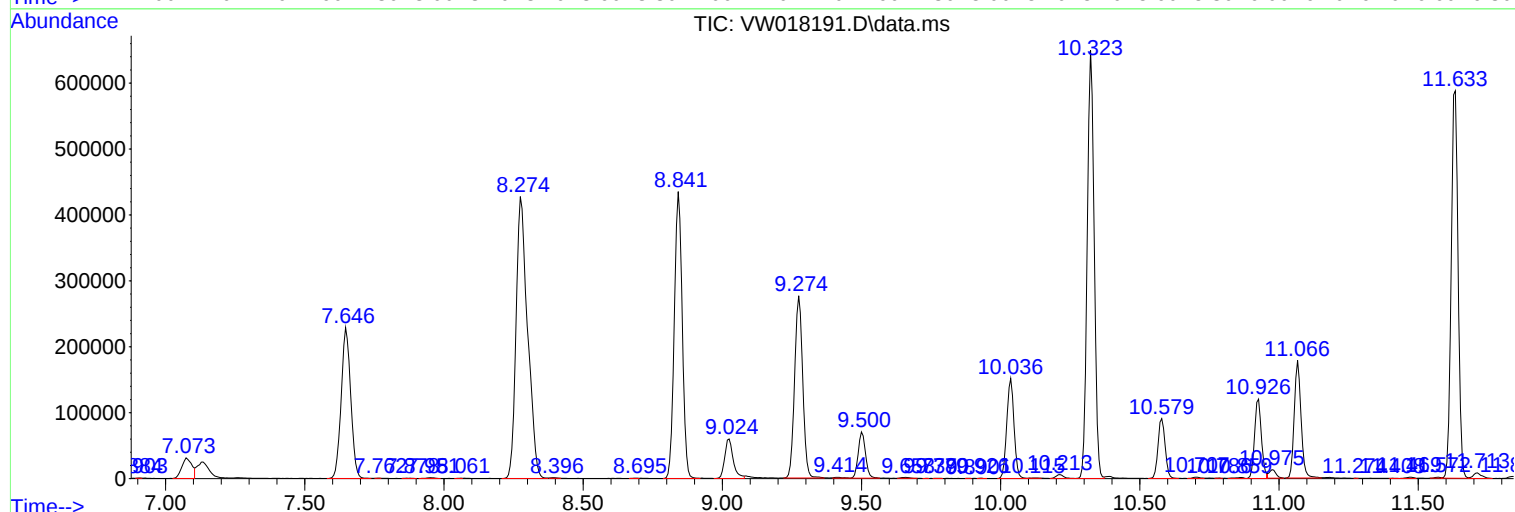
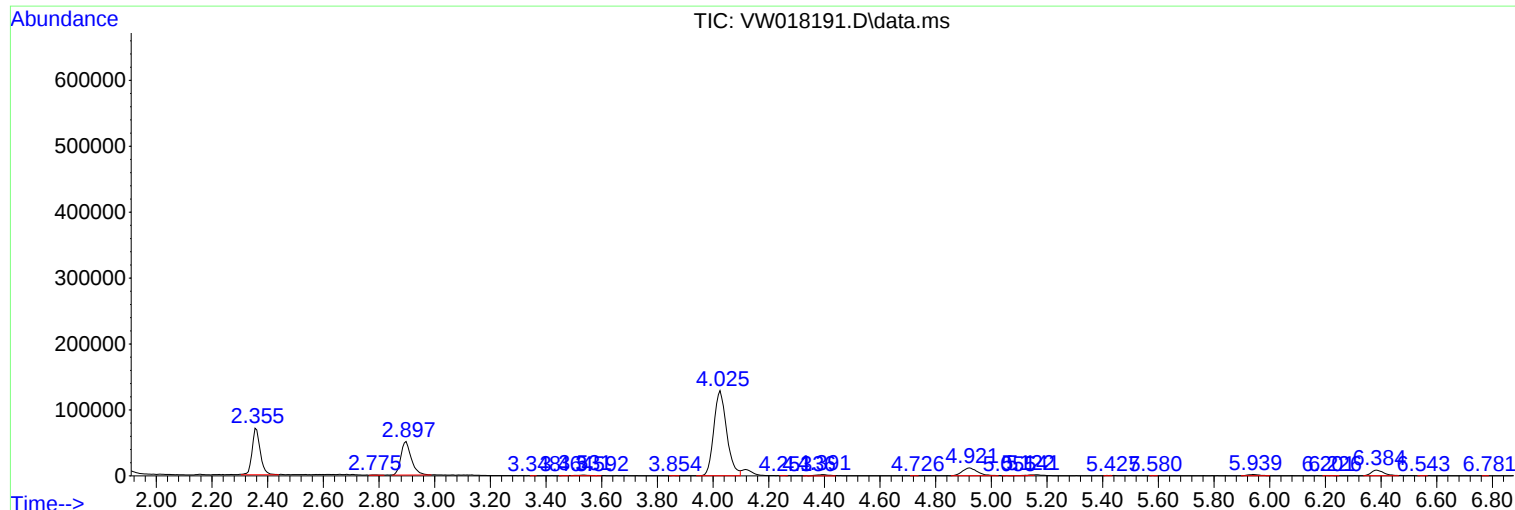
Sum of corrected areas: 10962206

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Instrument :
MSVOA_W
ClientSampleId :
DBH30

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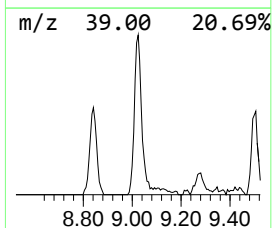
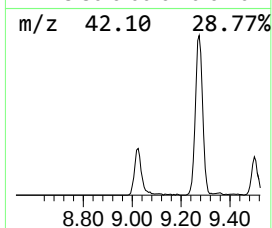
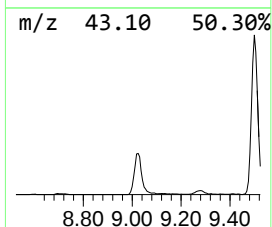
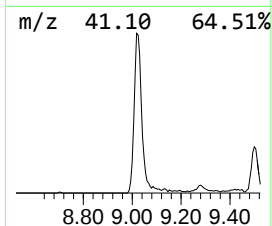
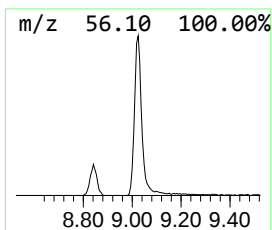
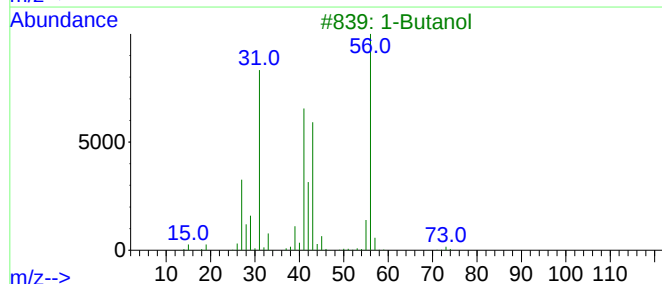
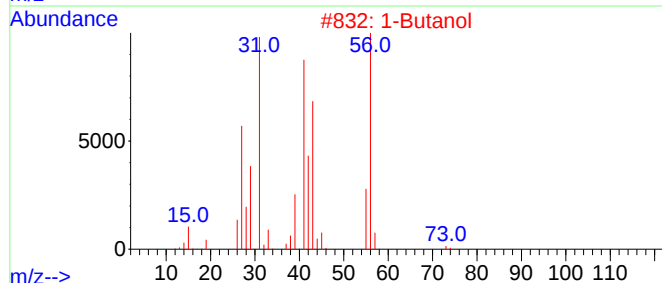
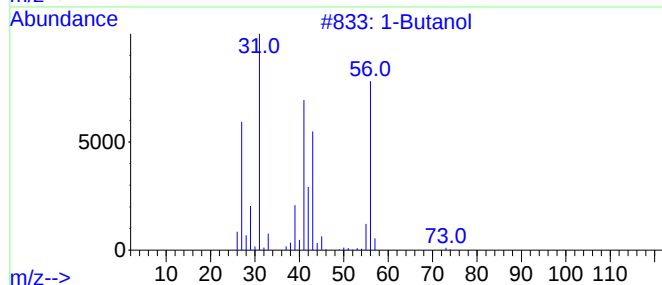
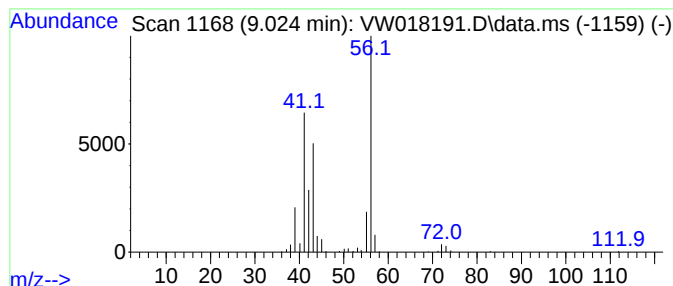
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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

Peak Number 1 1-Butanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.024	3.69 ug/L	126840	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol			74	C4H10O	000071-36-3	86
2	1-Butanol			74	C4H10O	000071-36-3	86
3	1-Butanol			74	C4H10O	000071-36-3	86
4	1-Butanol			74	C4H10O	000071-36-3	78
5	1-Butanol			74	C4H10O	000071-36-3	47



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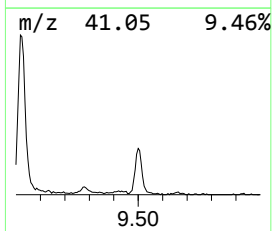
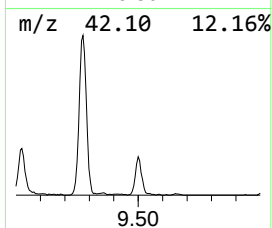
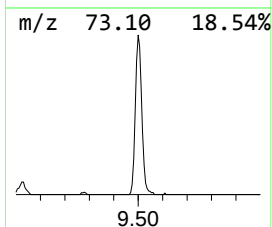
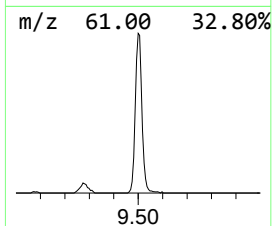
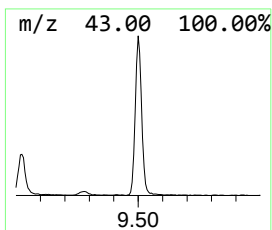
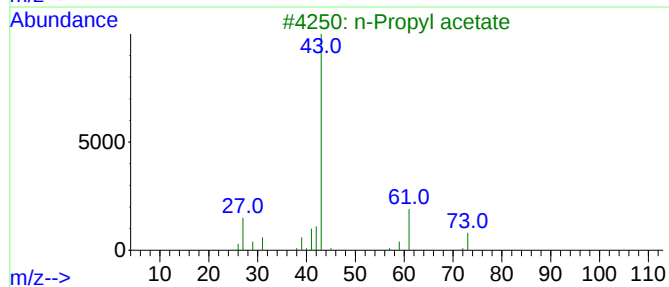
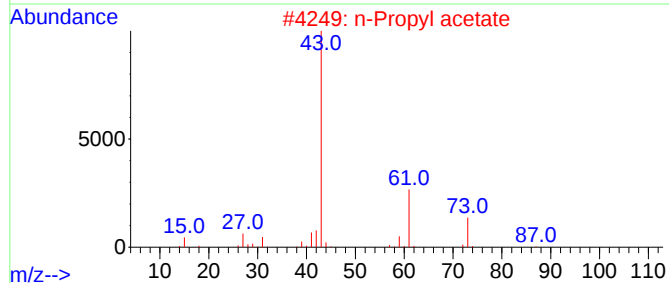
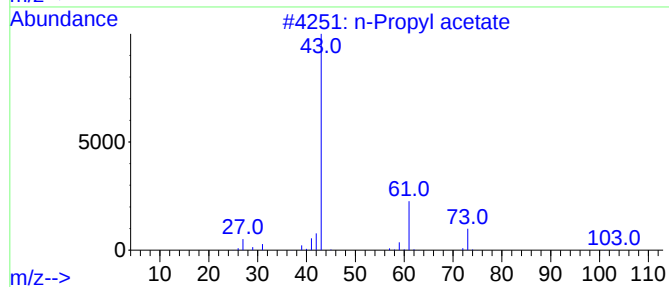
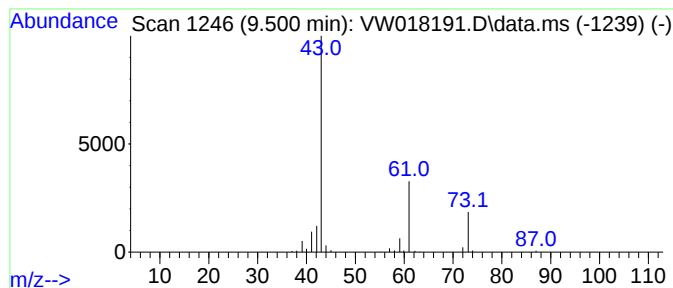
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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

Peak Number 2 n-Propyl acetate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.500	3.68 ug/L	126662	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Propyl acetate			102	C5H10O2	000109-60-4	83
2	n-Propyl acetate			102	C5H10O2	000109-60-4	83
3	n-Propyl acetate			102	C5H10O2	000109-60-4	78
4	Isopropyl acetate			102	C5H10O2	000108-21-4	36
5	Acetamide, N-acetyl-N-methyl-			115	C5H9NO2	001113-68-4	9



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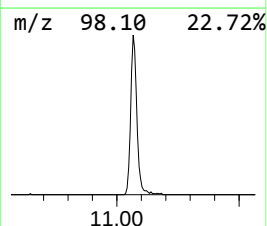
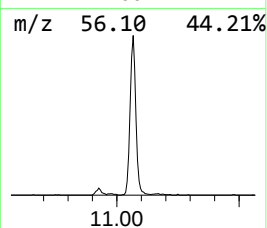
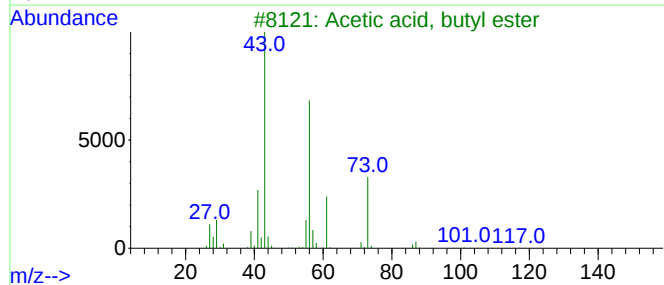
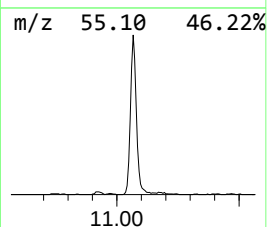
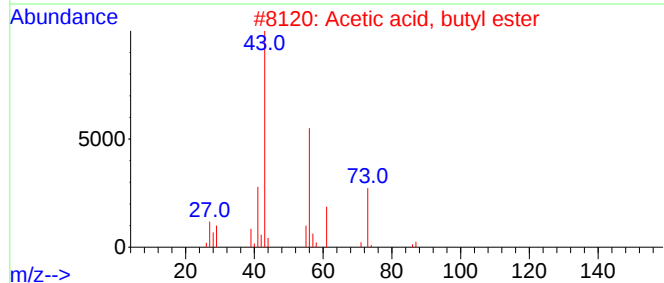
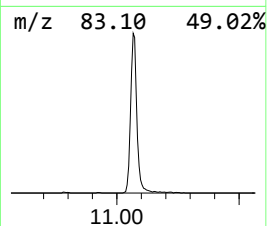
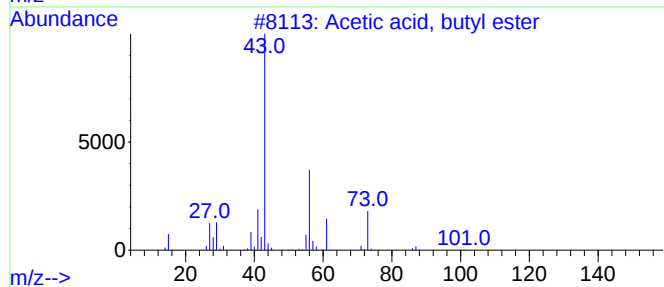
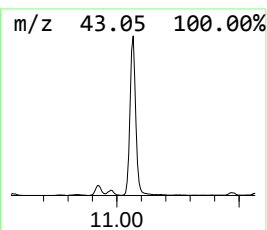
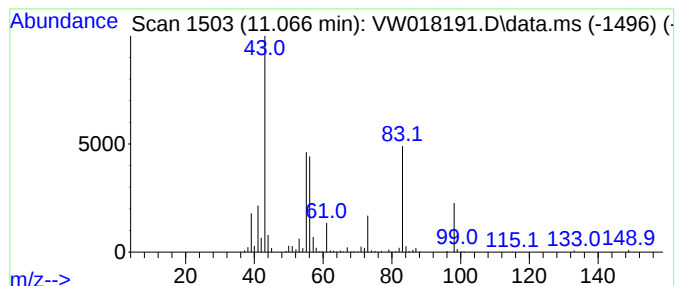
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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

Peak Number 4 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.066	7.42 ug/L	299731	Chlorobenzene-d5	11.633

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	43
2			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	38
3			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	38
4			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	38
5			Isobutyl acetate	116	C6H12O2	000110-19-0	32



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021721\
Data File : VW018191.D
Acq On : 17 Feb 2021 17:00
Operator : SY/VA
Sample : M1371-16
Misc : 3.93G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBH30

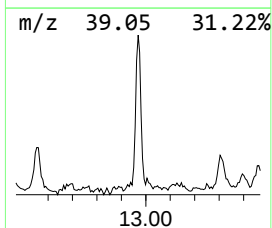
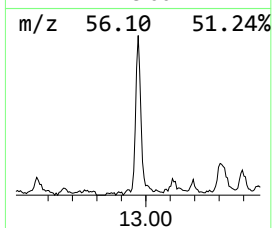
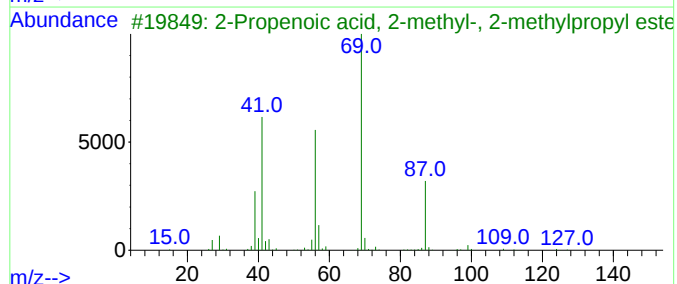
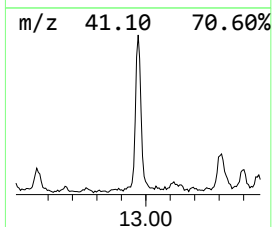
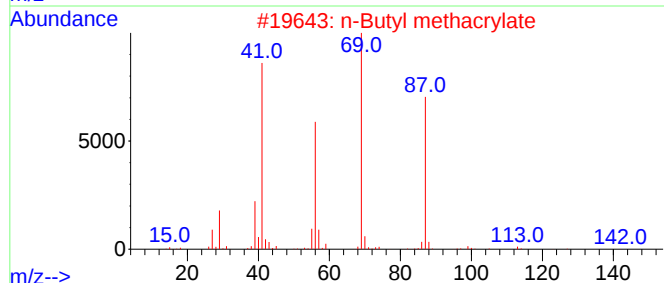
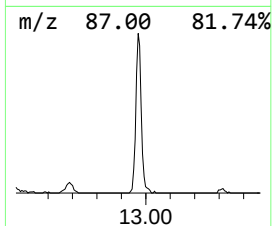
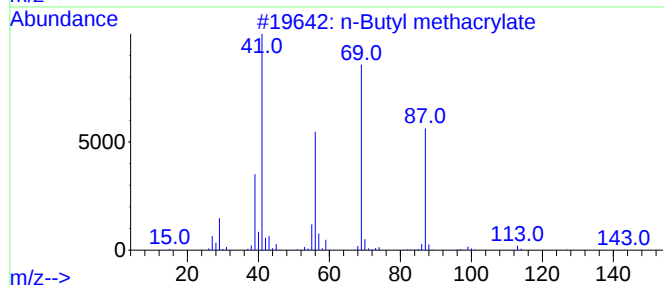
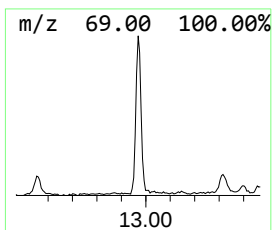
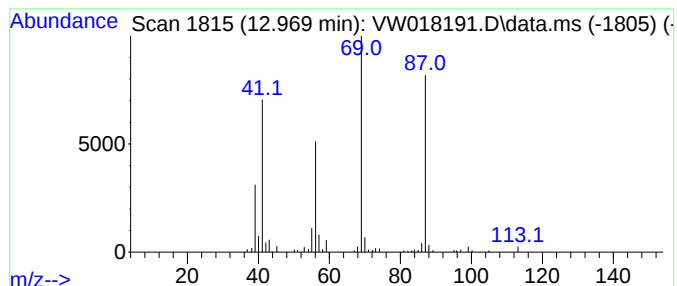
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Quant Title : VOC Analysis

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

Peak Number 5 n-Butyl methacrylate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.969	1.69 ug/L	71440	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Butyl methacrylate	142	C8H14O2	000097-88-1	83
2			n-Butyl methacrylate	142	C8H14O2	000097-88-1	83
3			2-Propenoic acid, 2-methyl-, 2-m...	142	C8H14O2	000097-86-9	74
4			2-Propenoic acid, 2-methyl-, hex...	170	C10H18O2	000142-09-6	74
5			2-Propenoic acid, 2-methyl-, hex...	170	C10H18O2	000142-09-6	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021721\
Data File : VW018191.D
Acq On : 17 Feb 2021 17:00
Operator : SY/VA
Sample : M1371-16
Misc : 3.93G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBH30

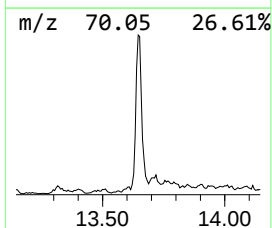
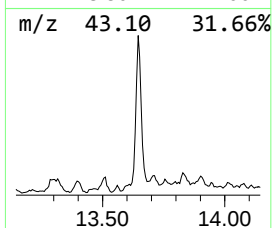
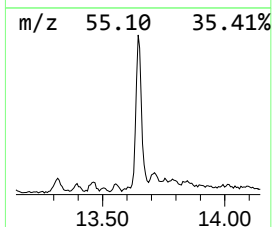
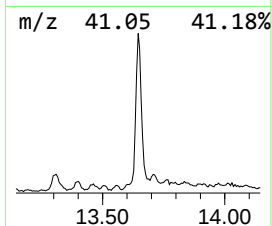
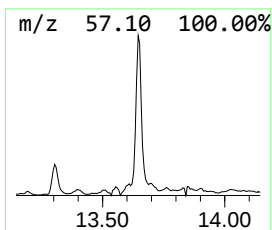
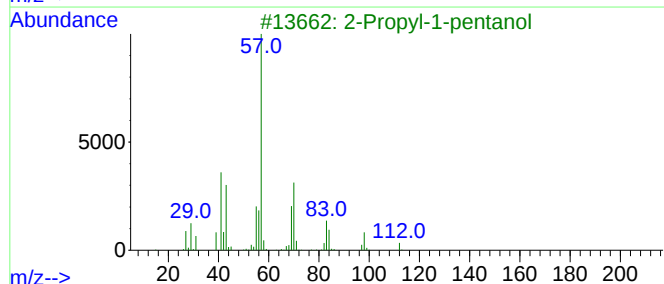
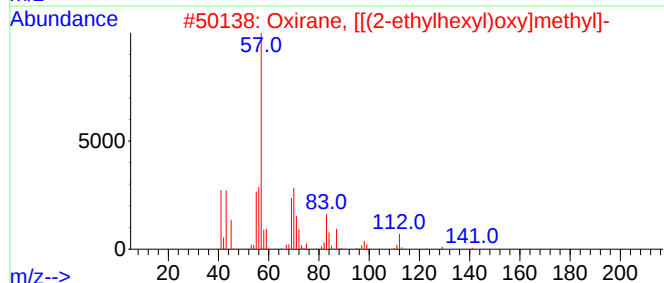
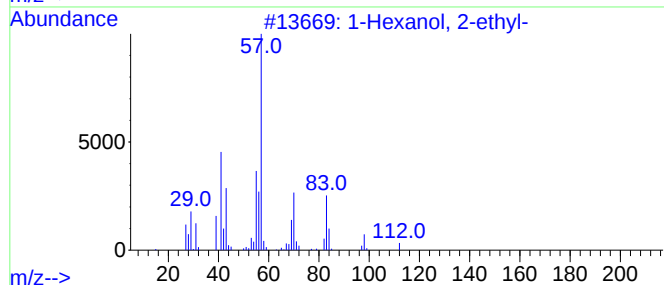
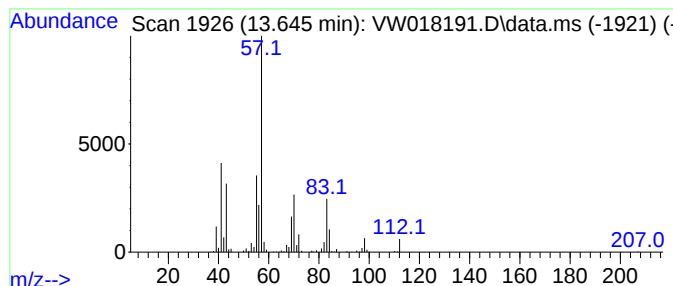
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Quant Title : VOC Analysis

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

Peak Number 6 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.645	5.53 ug/L	234406	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2			Oxirane, [[(2-ethylhexyl)oxy]met...	186	C11H22O2	002461-15-6	59
3			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	59
4			2-Ethylhexyl hydrogen maleate	228	C12H20O4	002370-71-0	59
5			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021721\
Data File : VW018191.D
Acq On : 17 Feb 2021 17:00
Operator : SY/VA
Sample : M1371-16
Misc : 3.93G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBH30

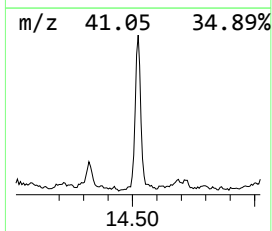
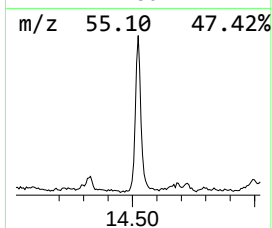
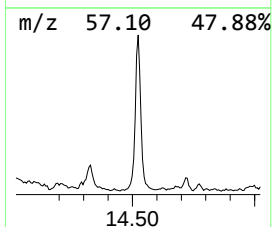
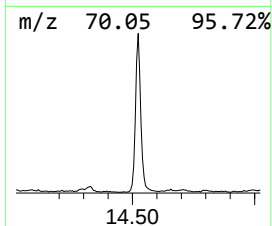
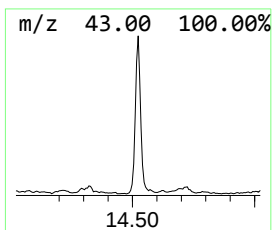
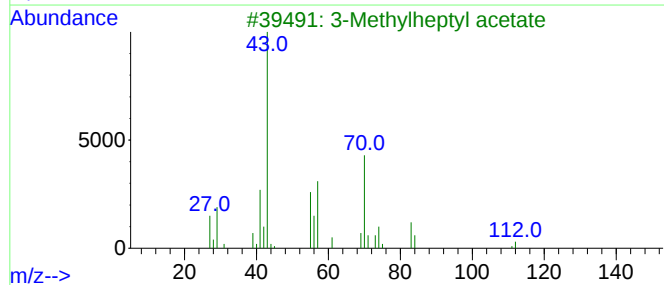
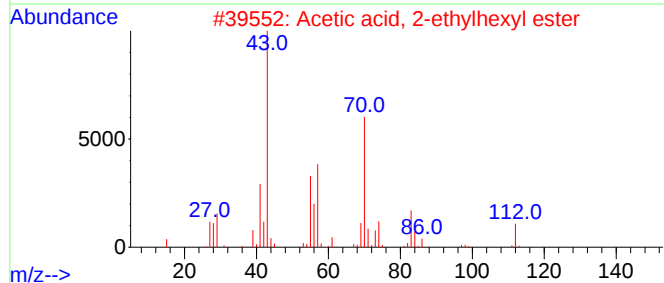
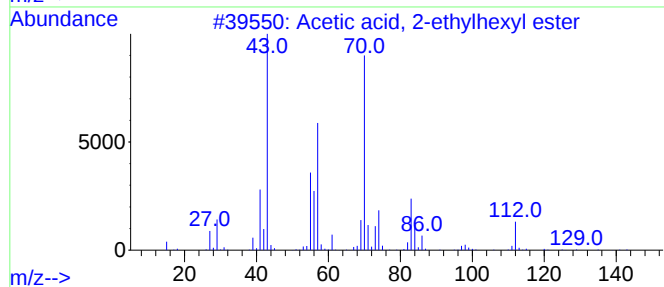
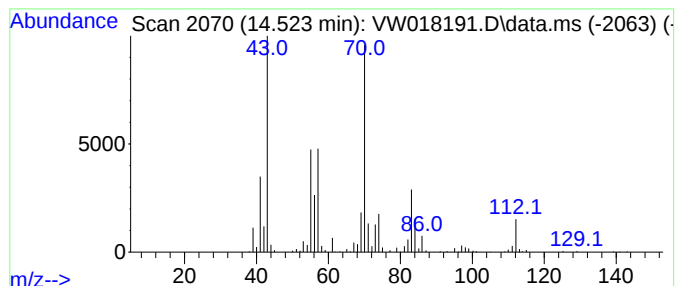
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Quant Title : VOC Analysis

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

Peak Number 7 Acetic acid, 2-ethylhexyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.523	6.19 ug/L	262226	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 2-ethylhexyl ester	172	C10H20O2	000103-09-3	91
2			Acetic acid, 2-ethylhexyl ester	172	C10H20O2	000103-09-3	91
3			3-Methylheptyl acetate	172	C10H20O2	072218-58-7	83
4			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	68
5			Bromoacetic acid, 2-ethylhexyl e...	250	C10H19BrO2	068144-73-0	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021721\
Data File : VW018191.D
Acq On : 17 Feb 2021 17:00
Operator : SY/VA
Sample : M1371-16
Misc : 3.93G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBH30

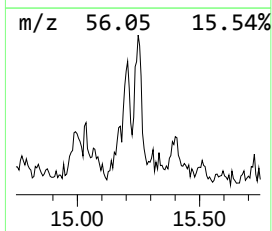
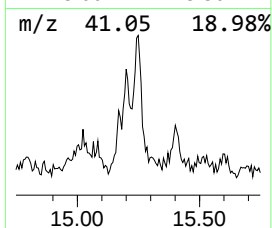
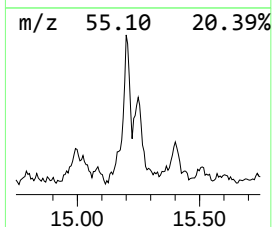
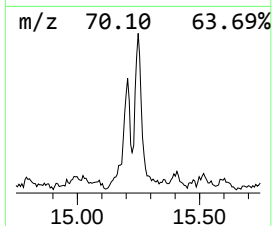
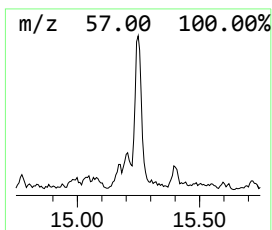
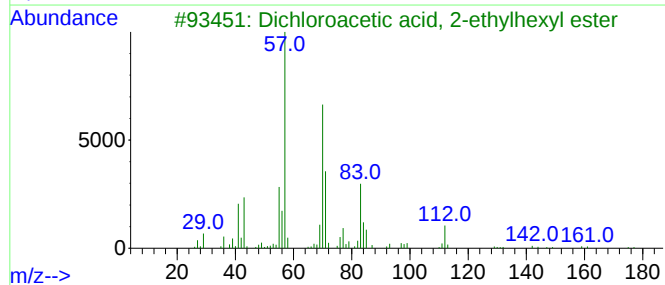
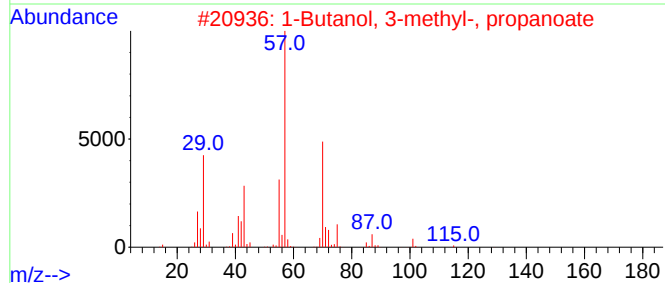
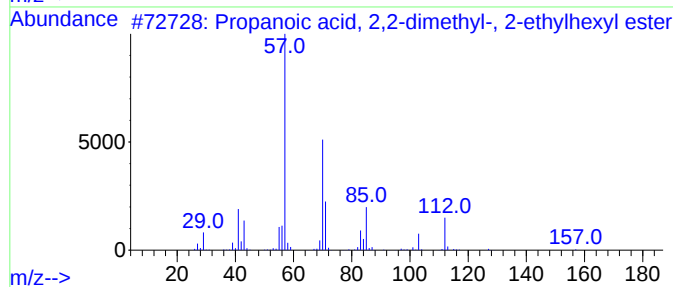
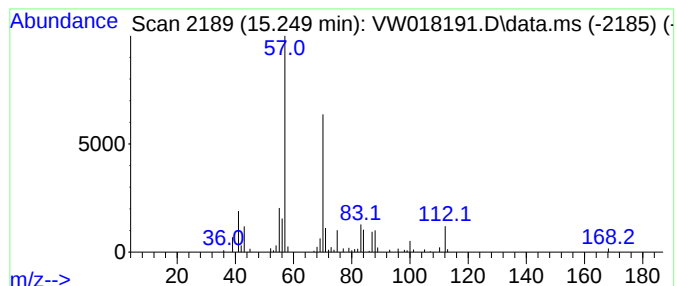
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SOM2WLM012921S.M
Quant Title : VOC Analysis

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

Peak Number 8 Propanoic acid, 2,2-dimethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.249	1.46 ug/L	61899	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2,2-dimethyl-, 2...	214	C13H26O2	016387-18-1	50
2			1-Butanol, 3-methyl-, propanoate	144	C8H16O2	000105-68-0	50
3			Dichloroacetic acid, 2-ethylhexy...	240	C10H18Cl2O2	068144-72-9	47
4			4-Nitrobenzoic acid, cyclobutyl ...	221	C11H11NO4	070335-00-1	47
5			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	015890-40-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW021721\
Data File : VW018191.D
Acq On : 17 Feb 2021 17:00
Operator : SY/VA
Sample : M1371-16
Misc : 3.93G/10ML/MSVOA_W/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBH30

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
1-Butanol	9.024	3.7	ug/L	126840	1	8.841	860201	25.0
n-Propyl acetate	9.500	3.7	ug/L	126662	1	8.841	860201	25.0
unknown-01	11.066	7.4	ug/L	299731	2	11.633	1010530	25.0
n-Butyl methacr...	12.969	1.7	ug/L	71440	3	13.554	1058860	25.0
1-Hexanol, 2-et...	13.645	5.5	ug/L	234406	3	13.554	1058860	25.0
Acetic acid, 2-...	14.523	6.2	ug/L	262226	3	13.554	1058860	25.0
Propanoic acid,...	15.249	1.5	ug/L	61899	3	13.554	1058860	25.0