

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW012720\
 Data File : VW014776.D
 Acq On : 27 Jan 2020 20:05
 Operator : SY/VA
 Sample : L1271-14
 Misc : 5.00G/10ML/MSVOA W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 GAT05

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM012720S.M
 Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.154	37	41	45	rBV2	4014	7790	0.03%	0.010%
2	2.349	67	73	82	rVB	328656	558458	2.14%	0.739%
3	2.885	154	161	180	rVB	192387	482468	1.85%	0.639%
4	3.434	249	251	255	rVB3	810	1048	0.00%	0.001%
5	3.495	259	261	263	rBV2	926	1069	0.00%	0.001%
6	3.666	284	289	290	rBV3	1484	2171	0.01%	0.003%
7	3.684	290	292	295	rVB3	1245	1210	0.00%	0.002%
8	4.013	332	346	359	rBV	329995	976237	3.73%	1.292%
9	4.397	401	409	416	rBV5	935	2778	0.01%	0.004%
10	4.769	466	470	474	rBV3	876	1075	0.00%	0.001%
11	4.909	483	493	508	rBV2	28134	97394	0.37%	0.129%
12	5.068	515	519	520	rBV2	1194	1303	0.00%	0.002%
13	5.159	533	534	538	rVB4	1224	1083	0.00%	0.001%
14	5.361	565	567	569	rBV2	919	1007	0.00%	0.001%
15	5.422	574	577	578	rBV	648	789	0.00%	0.001%
16	5.513	589	592	595	rVB2	597	840	0.00%	0.001%
17	5.763	626	633	635	rBV5	1284	2285	0.01%	0.003%
18	5.934	654	661	669	rVV3	7181	20601	0.08%	0.027%
19	6.056	680	681	686	rVB3	615	916	0.00%	0.001%
20	6.141	691	695	697	rBV2	562	801	0.00%	0.001%
21	6.367	729	732	737	rBV4	606	1119	0.00%	0.001%
22	6.799	801	803	809	rVB2	808	1023	0.00%	0.001%
23	6.879	813	816	819	rVB2	718	876	0.00%	0.001%
24	6.921	819	823	827	rBV4	912	1377	0.01%	0.002%
25	7.074	839	848	855	rBV	100538	291748	1.12%	0.386%
26	7.244	869	876	889	rVB	111057	277765	1.06%	0.368%
27	7.513	919	920	923	rBV3	1317	1655	0.01%	0.002%
28	7.647	931	942	955	rBV	563881	1350136	5.17%	1.787%
29	7.946	986	991	999	rVB6	3844	9878	0.04%	0.013%
30	8.061	1007	1010	1014	rBV4	932	1368	0.01%	0.002%
31	8.275	1032	1045	1062	rBV2	1122968	3346631	12.80%	4.430%
32	8.689	1108	1113	1114	rBV3	858	1498	0.01%	0.002%
33	8.842	1125	1138	1152	rBV	1185243	2324014	8.89%	3.076%
34	9.073	1172	1176	1184	rVB6	3734	7960	0.03%	0.011%

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35	9.275	1199	1209	1218	rBV	742781	1531766	5.86%	2.027%
36	9.464	1236	1240	1245	rVB7	812	1316	0.01%	0.002%
37	9.567	1253	1257	1259	rBV3	886	1134	0.00%	0.002%
38	9.659	1267	1272	1281	rVB6	2494	5443	0.02%	0.007%
39	9.750	1283	1287	1291	rBV5	873	1305	0.00%	0.002%
40	9.896	1303	1311	1315	rVB5	3274	6856	0.03%	0.009%
41	10.031	1325	1333	1348	rBV	374130	692954	2.65%	0.917%
42	10.213	1360	1363	1365	rBV3	1486	1629	0.01%	0.002%
43	10.323	1373	1381	1388	rBV	1561741	2748827	10.52%	3.638%
44	10.384	1388	1391	1403	rVB	43803	76122	0.29%	0.101%
45	10.573	1415	1422	1437	rBV	248772	441039	1.69%	0.584%
46	10.701	1437	1443	1449	rVB	27628	49213	0.19%	0.065%
47	10.921	1472	1479	1495	rBV	402453	683271	2.61%	0.904%
48	11.061	1497	1502	1509	rVV	30658	49667	0.19%	0.066%
49	11.171	1518	1520	1527	rVB7	2463	3154	0.01%	0.004%
50	11.439	1559	1564	1572	rVB2	6960	13113	0.05%	0.017%
51	11.628	1587	1595	1604	rBV	1605621	2641099	10.10%	3.496%
52	11.725	1608	1611	1616	rVB6	2739	4406	0.02%	0.006%
53	11.841	1623	1630	1632	rBV5	3392	5564	0.02%	0.007%
54	11.902	1639	1640	1645	rVB3	1807	1632	0.01%	0.002%
55	12.122	1674	1676	1678	rBV2	1416	1477	0.01%	0.002%
56	12.170	1678	1684	1687	rVV5	4754	8679	0.03%	0.011%
57	12.292	1701	1704	1709	rVB5	1901	3404	0.01%	0.005%
58	12.463	1727	1732	1736	rBV5	1592	3169	0.01%	0.004%
59	12.609	1749	1756	1761	rVB2	8872	15842	0.06%	0.021%
60	12.689	1761	1769	1778	rBV	554202	913554	3.49%	1.209%
61	12.853	1793	1796	1797	rBV3	2136	2274	0.01%	0.003%
62	12.932	1804	1809	1814	rVB5	9335	16373	0.06%	0.022%
63	12.981	1815	1817	1821	rVB4	2069	2798	0.01%	0.004%
64	13.067	1824	1831	1833	rBV5	1069	2355	0.01%	0.003%
65	13.115	1835	1839	1840	rBV3	1041	1221	0.00%	0.002%
66	13.243	1855	1860	1865	rBV5	8465	15981	0.06%	0.021%
67	13.298	1865	1869	1873	rVB3	10769	15374	0.06%	0.020%
68	13.335	1873	1875	1876	rBV2	1171	918	0.00%	0.001%
69	13.396	1883	1885	1890	rBV6	1945	2422	0.01%	0.003%
70	13.469	1893	1897	1903	rBV7	3908	6134	0.02%	0.008%
71	13.554	1904	1911	1919	rBV	1406191	2265843	8.67%	2.999%

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72	13.707	1934	1936	1939	rBV4	2060	2137	0.01%	0.003%
73	13.755	1939	1944	1946	rBV6	4599	7283	0.03%	0.010%
74	13.780	1946	1948	1952	rVB4	4299	5416	0.02%	0.007%
75	13.853	1952	1960	1968	rBV	1217231	2029276	7.76%	2.686%
76	13.938	1968	1974	1982	rVB	57019	100285	0.38%	0.133%
77	14.042	1982	1991	1992	rBV6	9527	19398	0.07%	0.026%
78	14.140	2003	2007	2013	rVB2	11999	23021	0.09%	0.030%
79	14.231	2014	2022	2027	rBV2	64100	104596	0.40%	0.138%
80	14.280	2027	2030	2034	rBV2	11284	15476	0.06%	0.020%
81	14.329	2036	2038	2044	rVB2	12483	15922	0.06%	0.021%
82	14.414	2048	2052	2055	rBV	23117	32954	0.13%	0.044%
83	14.457	2055	2059	2064	rBV2	54638	91176	0.35%	0.121%
84	14.554	2070	2075	2080	rBV2	71231	103498	0.40%	0.137%
85	14.688	2092	2097	2099	rBV	26173	39162	0.15%	0.052%
86	14.719	2100	2102	2108	rVB3	31369	42113	0.16%	0.056%
87	14.792	2108	2114	2120	rBV2	89711	189904	0.73%	0.251%
88	14.871	2121	2127	2134	rVV	233275	404685	1.55%	0.536%
89	14.950	2134	2140	2148	rVB	481091	819605	3.14%	1.085%
90	15.054	2153	2157	2162	rBV3	35705	64907	0.25%	0.086%
91	15.176	2172	2177	2183	rBV5	29785	73972	0.28%	0.098%
92	15.365	2201	2208	2218	rVB	5459141	9529277	36.46%	12.613%
93	15.469	2220	2225	2232	rVV3	60836	136800	0.52%	0.181%
94	15.554	2235	2239	2246	rVB	48898	84086	0.32%	0.111%
95	15.658	2251	2256	2260	rBV3	34589	73243	0.28%	0.097%
96	15.847	2283	2287	2292	rVB3	57483	88058	0.34%	0.117%
97	15.968	2299	2307	2312	rBV4	81859	221192	0.85%	0.293%
98	16.377	2367	2374	2376	rBV	70976	145649	0.56%	0.193%
99	16.438	2376	2384	2398	rVB	12594924	26138899	100.00%	34.598%
100	16.645	2409	2418	2432	rVB	5894037	12932869	49.48%	17.118%

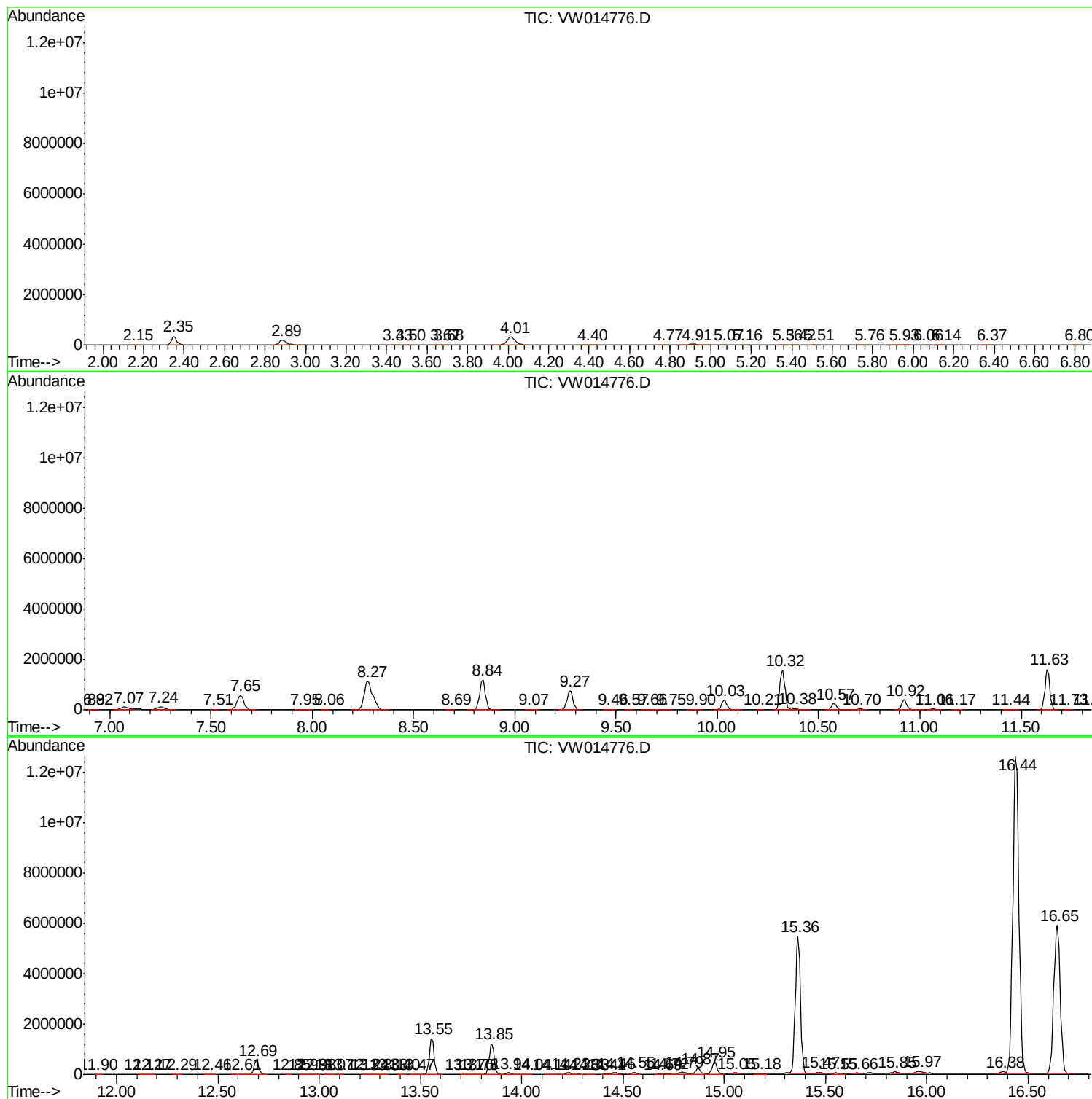
Sum of corrected areas: 75550558

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM012720S.M
Quant Title : VOC Analysis

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



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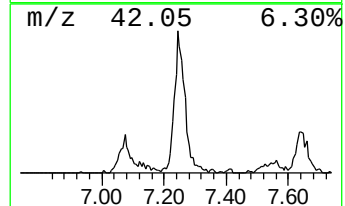
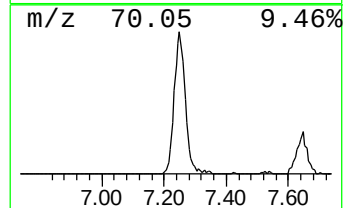
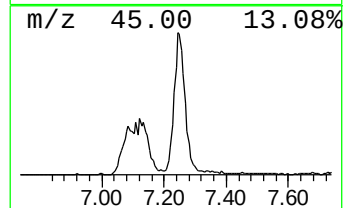
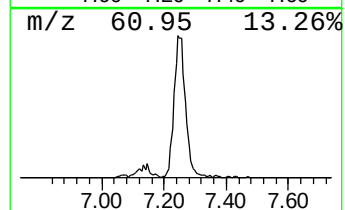
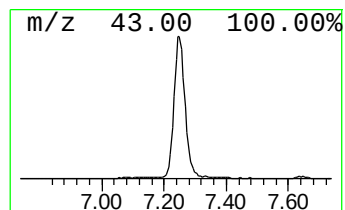
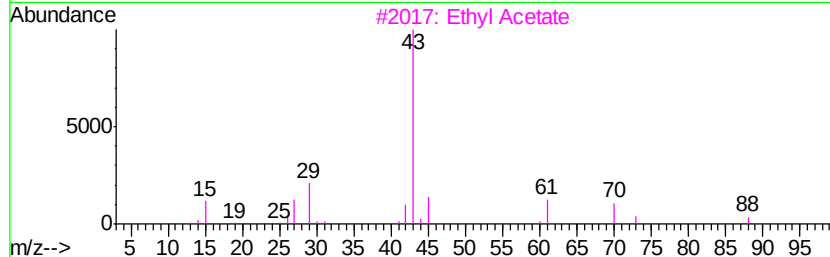
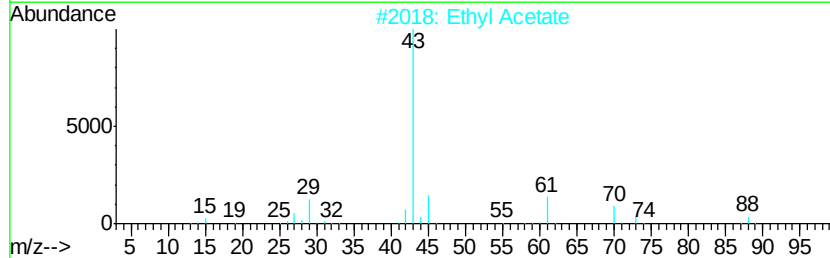
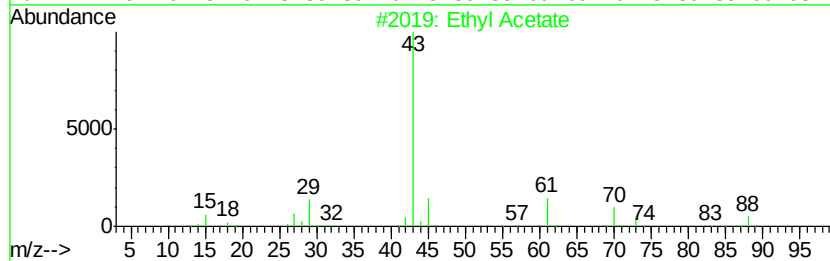
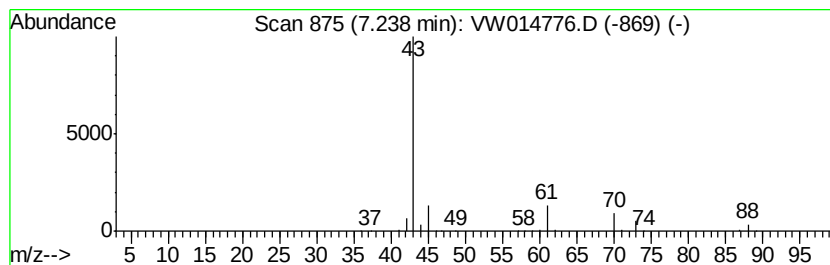
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM012720S.M
Quant Title : VOC Analysis

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

Peak Number 1 Ethyl Acetate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.24	2.99 ug/L	277765	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethyl Acetate	88	C4H8O2	000141-78-6	90
2		Ethyl Acetate	88	C4H8O2	000141-78-6	86
3		Ethyl Acetate	88	C4H8O2	000141-78-6	86
4		CH3C(O)CH2CH2OH	88	C4H8O2	000590-90-9	53
5		1,4-Butanediol, diacetate	174	C8H14O4	000628-67-1	28



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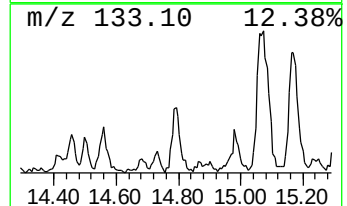
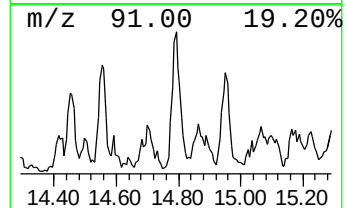
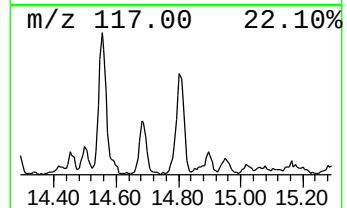
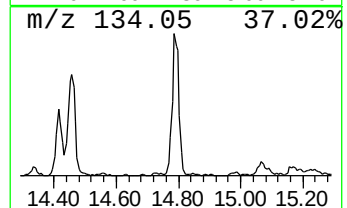
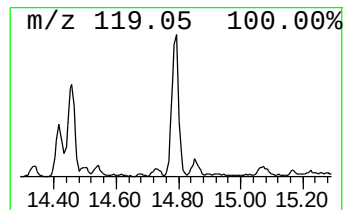
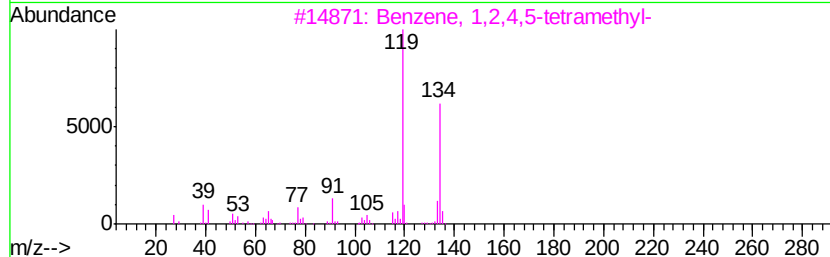
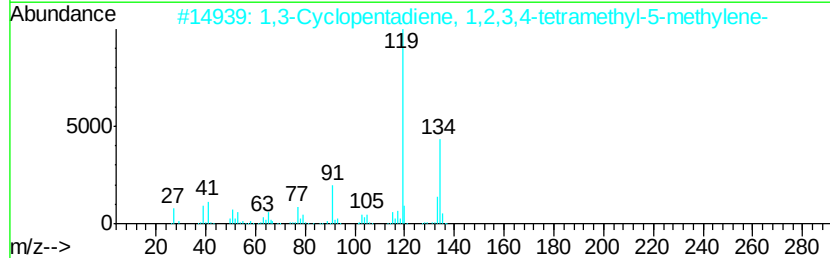
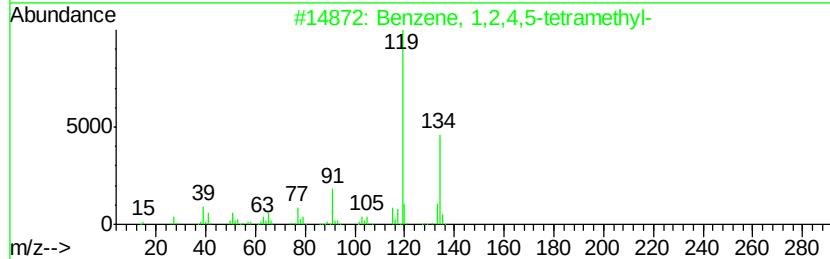
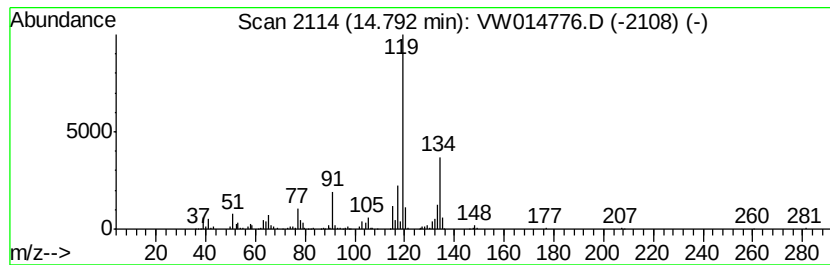
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Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	2.10 ug/L	189904	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	94
2	1,3-Cyclopentadiene, 1,2,3,4-tet...	134 C10H14	076089-59-3	93
3	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	87
4	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	87
5	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	87



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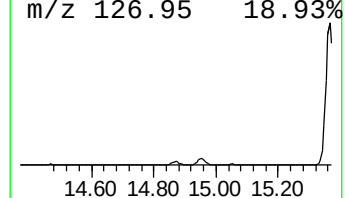
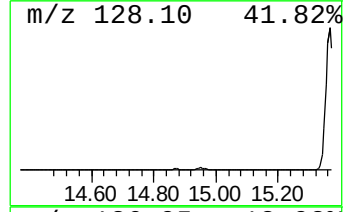
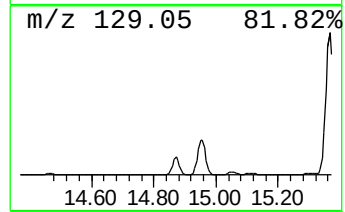
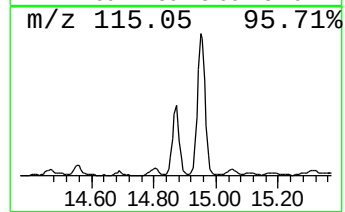
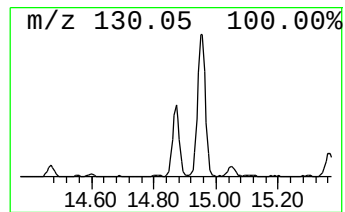
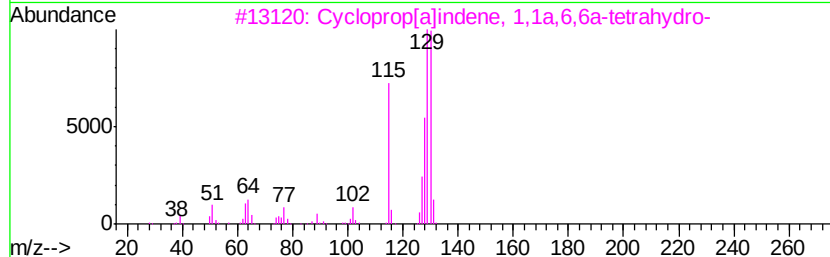
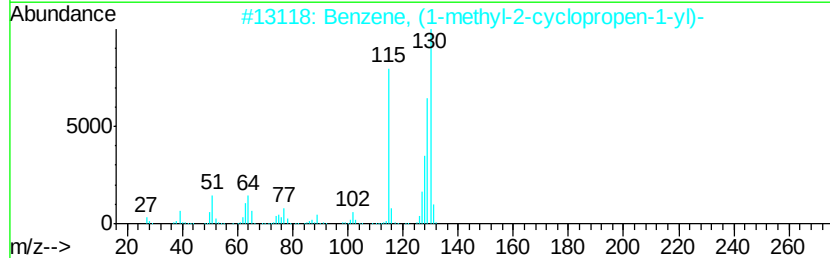
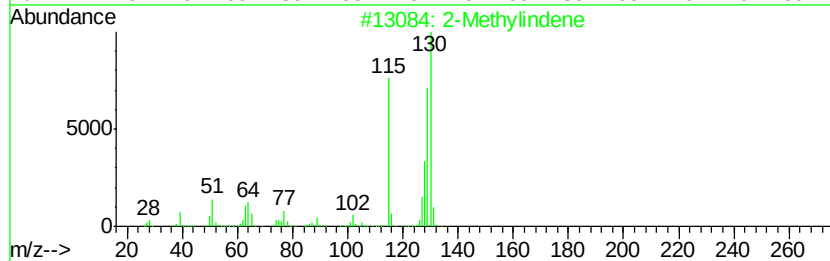
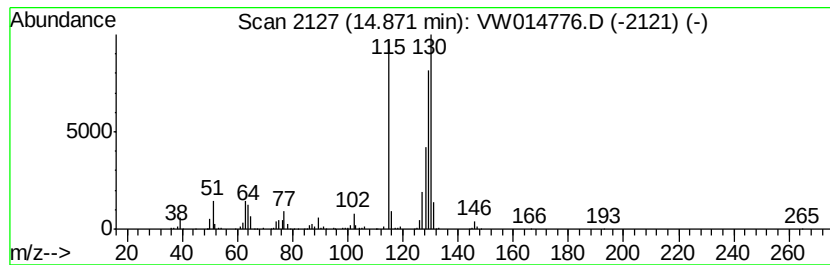
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Quant Title : VOC Analysis

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Peak Number 4 2-Methylindene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	4.47 ug/L	404685	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Methylindene	130 C10H10	002177-47-1	97
2	Benzene, (1-methyl-2-cyclopropen...	130 C10H10	065051-83-4	95
3	Cycloprop[alindene, 1,1a,6,6a-te...	130 C10H10	015677-15-3	93
4	1H-Indene, 1-methyl-	130 C10H10	000767-59-9	93
5	Benzene, 1-butyryl-	130 C10H10	000622-76-4	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012720\
Data File : VW014776.D
Acq On : 27 Jan 2020 20:05
Operator : SY/VA
Sample : L1271-14
Misc : 5.00G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GAT05

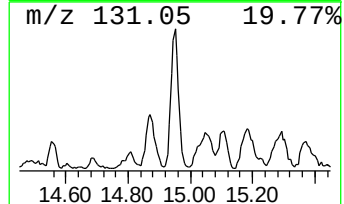
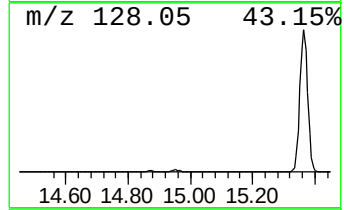
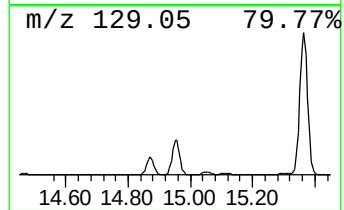
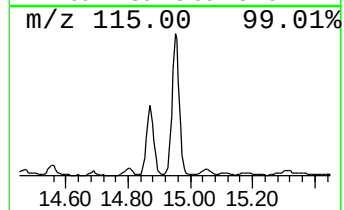
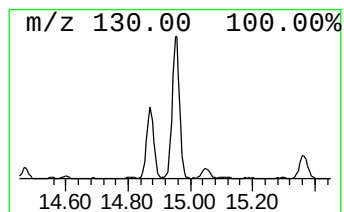
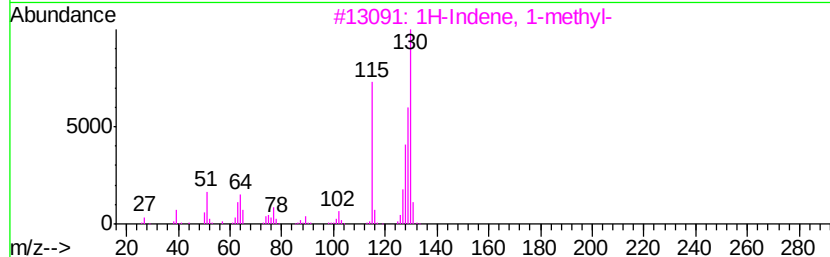
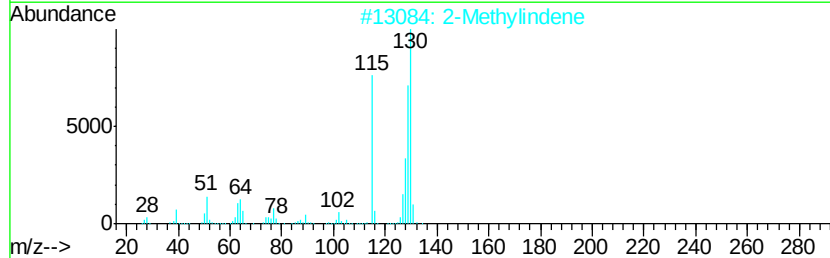
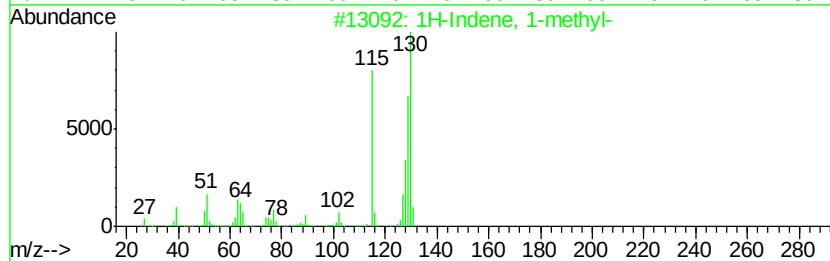
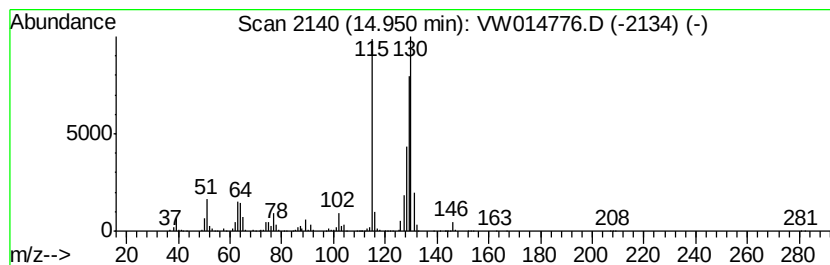
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Quant Title : VOC Analysis

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

Peak Number 5 1H-Indene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	9.04 ug/L	819605	1,4-Dichlorobenzene-d4	13.55

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2	2-Methylindene	130	C10H10	002177-47-1	96
3	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
4	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	95
5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012720\
Data File : VW014776.D
Acq On : 27 Jan 2020 20:05
Operator : SY/VA
Sample : L1271-14
Misc : 5.00G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

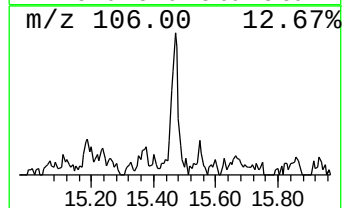
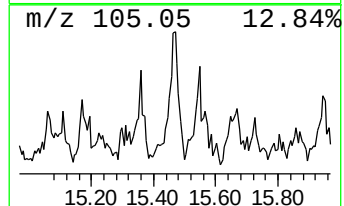
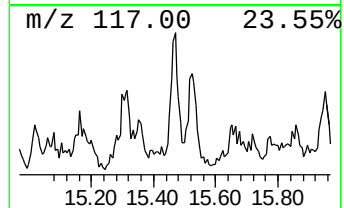
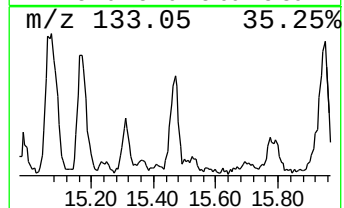
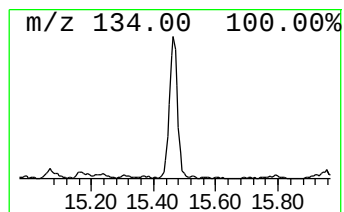
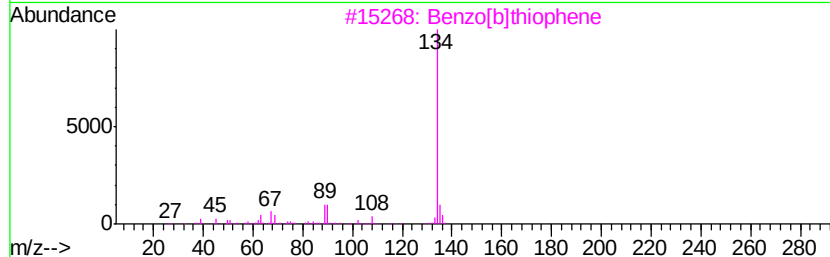
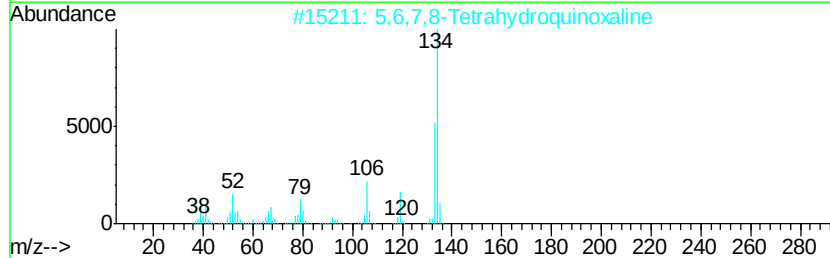
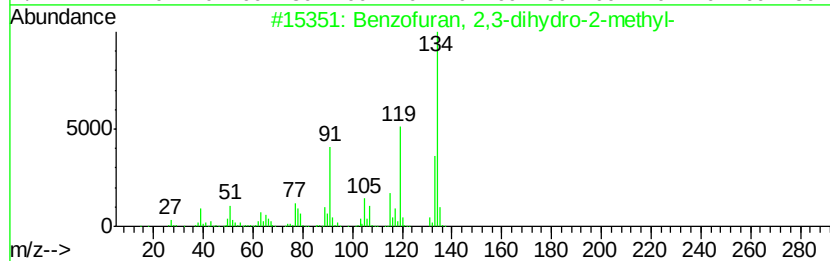
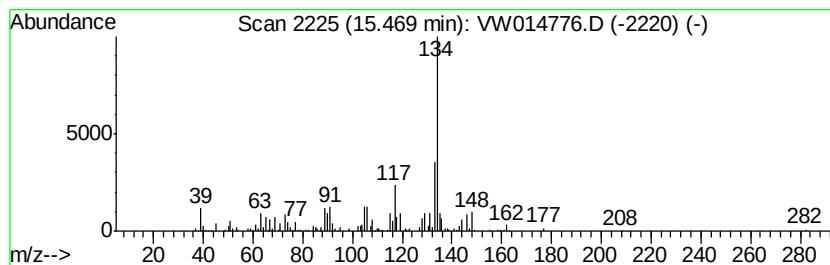
Instrument :
MSVOA_W
ClientSampleId :
GAT05

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM012720S.M
Quant Title : VOC Analysis

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

Peak Number 7 Benzofuran, 2,3-dihydro-2-m... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.47	1.51 ug/L	136800	1,4-Dichlorobenzene-d4	13.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2,3-dihydro-2-methyl-	134	C9H10O	001746-11-8	58
2	5,6,7,8-Tetrahydroquinoxaline	134	C8H10N2	034413-35-9	53
3	Benzo[blthiophene	134	C8H6S	000095-15-8	50
4	2-(1-Cyclopentenyl)furan	134	C9H10O	115754-78-4	50
5	2-Allylphenol	134	C9H10O	001745-81-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012720\
Data File : VW014776.D
Acq On : 27 Jan 2020 20:05
Operator : SY/VA
Sample : L1271-14
Misc : 5.00G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

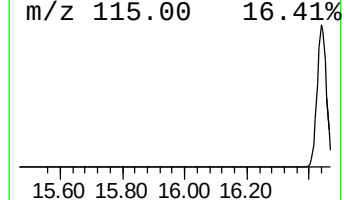
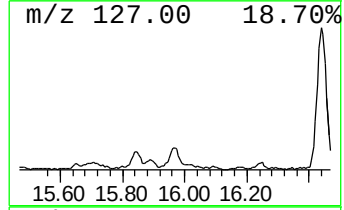
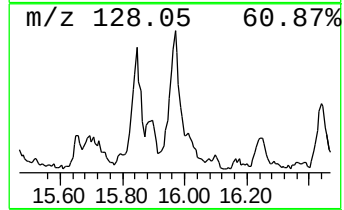
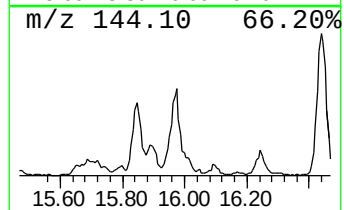
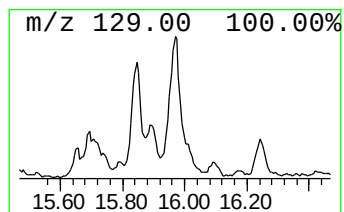
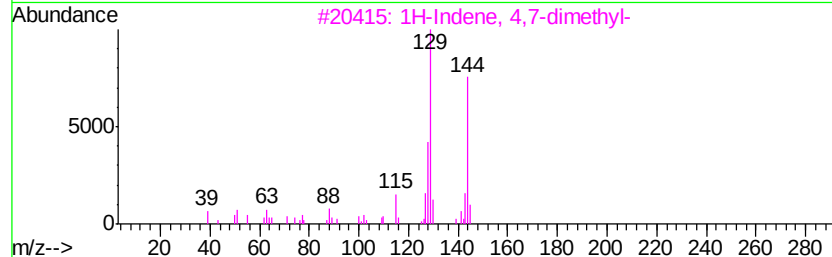
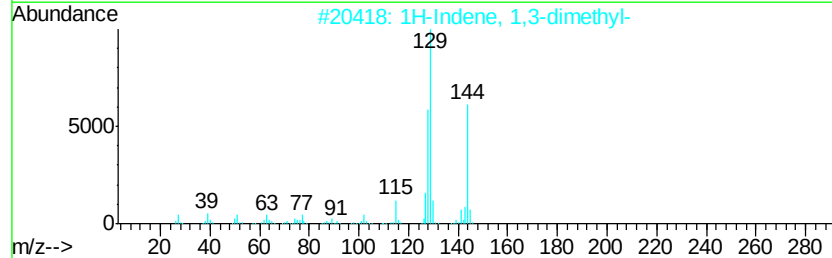
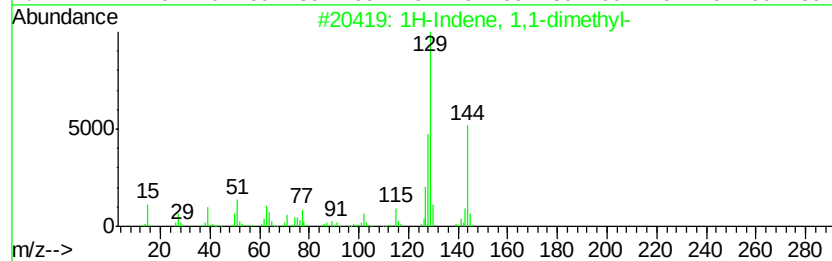
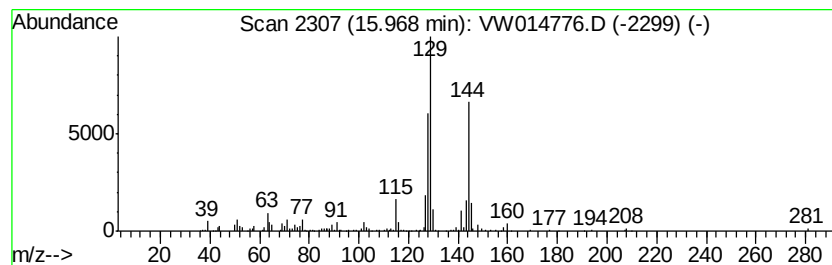
Instrument :
MSVOA_W
ClientSampleId :
GAT05

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM012720S.M
Quant Title : VOC Analysis

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

Peak Number 8 1H-Indene, 1,1-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	2.44 ug/L	221192	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 1,1-dimethyl-	144 C11H12	018636-55-0	91
2	1H-Indene, 1,3-dimethyl-	144 C11H12	002177-48-2	87
3	1H-Indene, 4,7-dimethyl-	144 C11H12	006974-97-6	87
4	(1-Methylbuta-1,3-dienyl)benzene	144 C11H12	054758-36-0	87
5	Benzene, 1-cyclopenten-1-yl-	144 C11H12	000825-54-7	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012720\
Data File : VW014776.D
Acq On : 27 Jan 2020 20:05
Operator : SY/VA
Sample : L1271-14
Misc : 5.00G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GAT05

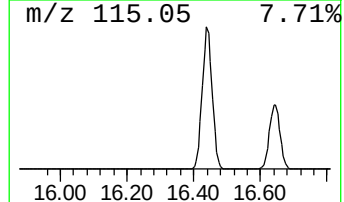
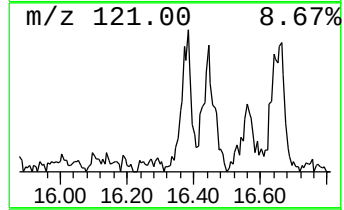
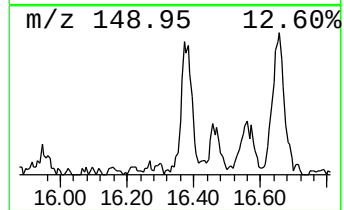
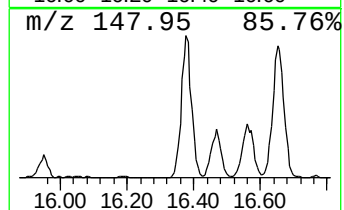
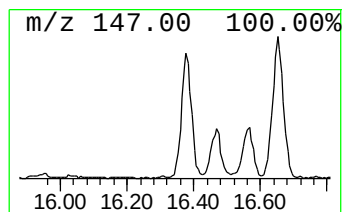
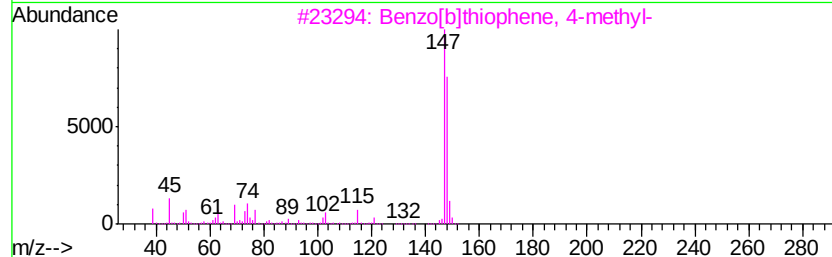
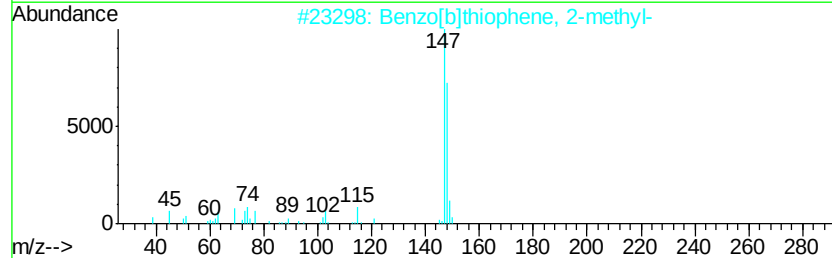
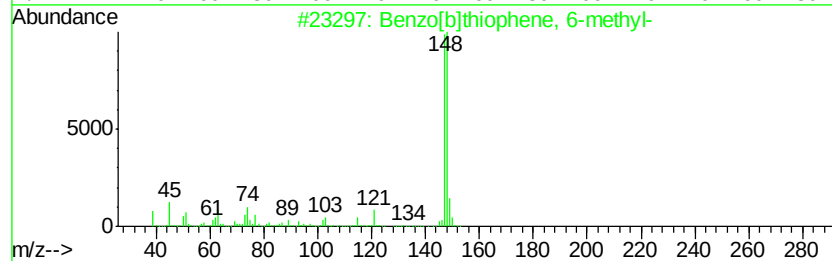
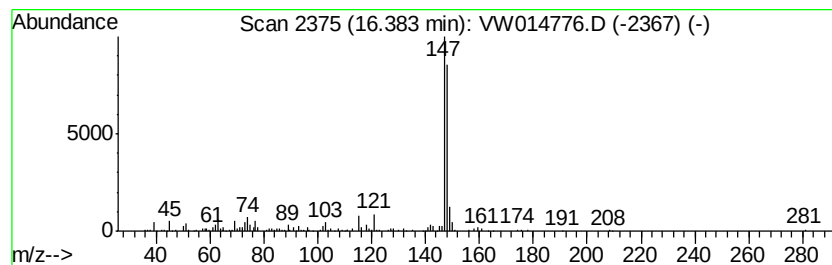
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Quant Title : VOC Analysis

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

Peak Number 9 Benzo[b]thiophene, 6-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	1.61 ug/L	145649	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	91
2		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
3		Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	90
4		3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
5		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012720\
Data File : VW014776.D
Acq On : 27 Jan 2020 20:05
Operator : SY/VA
Sample : L1271-14
Misc : 5.00G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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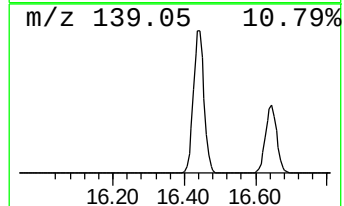
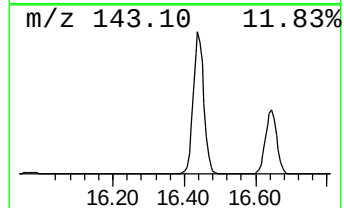
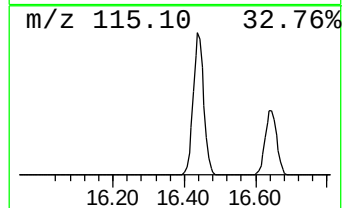
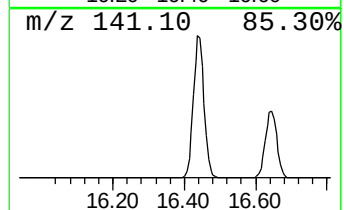
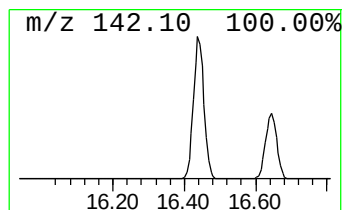
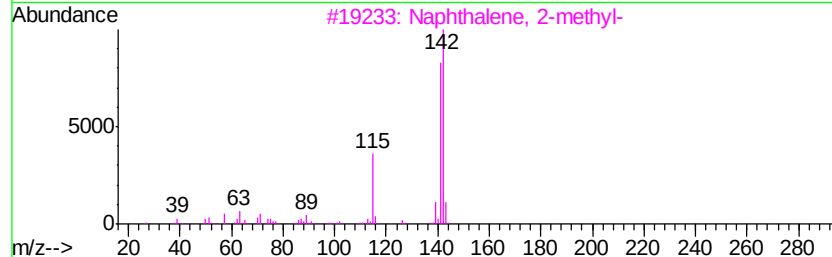
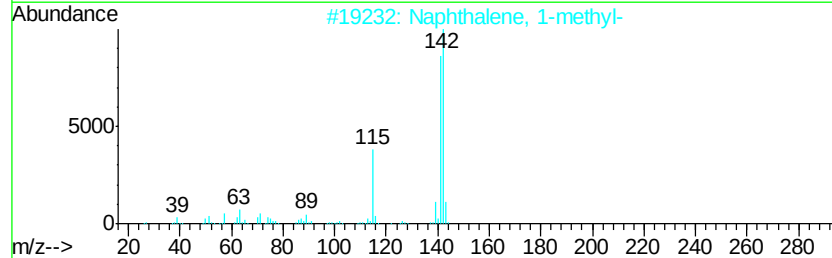
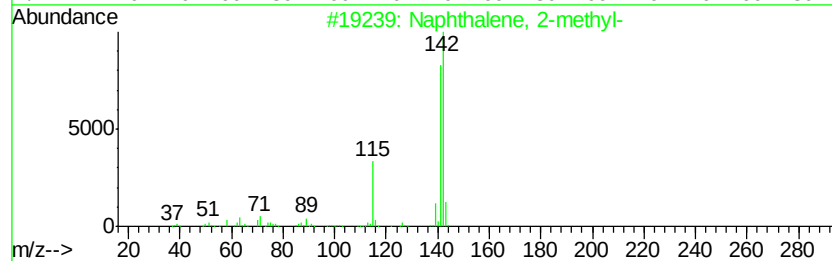
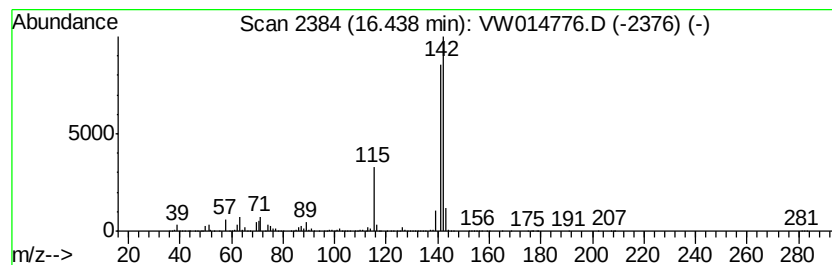
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Quant Title : VOC Analysis

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	288.40 ug/L	26138900	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW012720\
 Data File : VW014776.D
 Acq On : 27 Jan 2020 20:05
 Operator : SY/VA
 Sample : L1271-14
 Misc : 5.00G/10ML/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 GAT05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM012720S.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
Ethyl Acetate	7.24	3.0	ug/L	277765	1	8.84	2324010	25.0
Benzene, 1,2,4,5-...	14.79	2.1	ug/L	189904	3	13.55	2265840	25.0
2-Methyldene	14.87	4.5	ug/L	404685	3	13.55	2265840	25.0
1H-Indene, 1-methyl-	14.95	9.0	ug/L	819605	3	13.55	2265840	25.0
Benzofuran, 2,3-d...	15.47	1.5	ug/L	136800	3	13.55	2265840	25.0
1H-Indene, 1,1-di...	15.97	2.4	ug/L	221192	3	13.55	2265840	25.0
Benzo[b]thiophene...	16.38	1.6	ug/L	145649	3	13.55	2265840	25.0
Naphthalene, 1-me...	16.44	288.4	ug/L	26138900	3	13.55	2265840	25.0