

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW011420\
 Data File : VW014634.D
 Acq On : 14 Jan 2020 15:47
 Operator : SY/VA
 Sample : L1109-14
 Misc : 5.04G/10ML/MSVOA W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 GASH3

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\SOM2WLM122719S.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	2.160	36	42	47	rBV2	12622	25066	0.79%	0.080%
2	2.349	67	73	84	rVB	280376	468466	14.83%	1.501%
3	2.885	154	161	174	rVB	177967	407867	12.91%	1.307%
4	3.531	265	267	272	rVB3	1695	2730	0.09%	0.009%
5	3.794	306	310	313	rBV3	545	1146	0.04%	0.004%
6	4.013	334	346	359	rBV	393972	1174680	37.19%	3.765%
7	4.672	451	454	457	rBV4	548	945	0.03%	0.003%
8	4.732	461	464	467	rVB2	805	845	0.03%	0.003%
9	4.915	482	494	509	rBV2	42837	151629	4.80%	0.486%
10	5.129	526	529	535	rVV4	697	1483	0.05%	0.005%
11	5.775	633	635	639	rVB3	1074	854	0.03%	0.003%
12	5.933	651	661	673	rVB6	5300	18004	0.57%	0.058%
13	6.074	682	684	687	rVB2	821	827	0.03%	0.003%
14	6.104	687	689	691	rBV2	1084	1163	0.04%	0.004%
15	6.129	691	693	697	rVB2	957	954	0.03%	0.003%
16	6.604	767	771	774	rBV3	688	900	0.03%	0.003%
17	6.708	786	788	791	rBV2	694	837	0.03%	0.003%
18	7.074	839	848	855	rBV	90454	262074	8.30%	0.840%
19	7.525	915	922	923	rBV4	1714	2511	0.08%	0.008%
20	7.647	930	942	961	rBV	503966	1238194	39.20%	3.968%
21	7.945	984	991	1002	rVB5	4936	10662	0.34%	0.034%
22	8.177	1027	1029	1031	rBV2	1040	1001	0.03%	0.003%
23	8.275	1033	1045	1060	rBV2	1068778	3158445	100.00%	10.123%
24	8.841	1128	1138	1155	rBV	1056867	2083175	65.96%	6.677%
25	9.073	1172	1176	1182	rVB6	3290	7091	0.22%	0.023%
26	9.159	1188	1190	1192	rVB	1117	940	0.03%	0.003%
27	9.274	1199	1209	1219	rBV	705336	1436679	45.49%	4.605%
28	9.396	1227	1229	1232	rVB3	1140	1012	0.03%	0.003%
29	9.476	1240	1242	1245	rVB2	1001	897	0.03%	0.003%
30	9.543	1250	1253	1258	rBV5	549	1016	0.03%	0.003%
31	9.665	1268	1273	1281	rVB8	3208	7706	0.24%	0.025%
32	9.725	1281	1283	1286	rBV	855	956	0.03%	0.003%
33	9.884	1304	1309	1312	rBV2	703	796	0.03%	0.003%
34	10.030	1325	1333	1342	rBV	364731	670772	21.24%	2.150%

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Integration Parameters: LSCINT.P

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 Title : VOC Analysis

35	10.164	1353	1355	1357	rVB3	1176	882	0.03%	0.003%
36	10.207	1357	1362	1367	rBV4	4455	7484	0.24%	0.024%
37	10.323	1372	1381	1387	rBV	1552404	2713750	85.92%	8.698%
38	10.384	1387	1391	1399	rVV	90965	168121	5.32%	0.539%
39	10.500	1408	1410	1413	rBV3	729	797	0.03%	0.003%
40	10.573	1415	1422	1435	rBV	233015	422021	13.36%	1.353%
41	10.701	1438	1443	1451	rVB2	14176	24733	0.78%	0.079%
42	10.829	1458	1464	1465	rBV3	831	1689	0.05%	0.005%
43	10.853	1465	1468	1473	rVV4	2264	3549	0.11%	0.011%
44	10.920	1473	1479	1494	rVV	342770	592366	18.75%	1.899%
45	11.061	1497	1502	1509	rVV2	25029	41879	1.33%	0.134%
46	11.170	1518	1520	1523	rBV4	997	1211	0.04%	0.004%
47	11.292	1535	1540	1543	rBV3	760	1293	0.04%	0.004%
48	11.347	1545	1549	1553	rBV3	746	1477	0.05%	0.005%
49	11.402	1555	1558	1560	rBV3	1153	1108	0.04%	0.004%
50	11.445	1560	1565	1568	rBV4	2991	4774	0.15%	0.015%
51	11.530	1572	1579	1583	rBV6	821	1782	0.06%	0.006%
52	11.628	1587	1595	1605	rBV	1429284	2388815	75.63%	7.656%
53	11.701	1605	1607	1609	rBV3	1193	1249	0.04%	0.004%
54	11.835	1622	1629	1639	rVV4	6786	15329	0.49%	0.049%
55	12.121	1673	1676	1678	rBV2	1516	1920	0.06%	0.006%
56	12.164	1678	1683	1693	rVV4	10817	22759	0.72%	0.073%
57	12.256	1693	1698	1702	rVV5	1216	2309	0.07%	0.007%
58	12.298	1702	1705	1709	rVV6	1309	1572	0.05%	0.005%
59	12.524	1739	1742	1744	rBV4	1154	1336	0.04%	0.004%
60	12.603	1750	1755	1762	rBV	11806	20808	0.66%	0.067%
61	12.688	1762	1769	1779	rBV	516469	856613	27.12%	2.745%
62	12.865	1792	1798	1800	rBV2	15535	25668	0.81%	0.082%
63	12.938	1805	1810	1817	rVV2	57605	104649	3.31%	0.335%
64	12.987	1817	1818	1823	rVB4	3509	3755	0.12%	0.012%
65	13.097	1832	1836	1845	rBV4	11497	22832	0.72%	0.073%
66	13.249	1854	1861	1866	rBV	165670	262305	8.30%	0.841%
67	13.298	1866	1869	1874	rVB4	8729	14707	0.47%	0.047%
68	13.408	1883	1887	1889	rBV5	3006	5183	0.16%	0.017%
69	13.469	1890	1897	1905	rBV7	15154	39423	1.25%	0.126%
70	13.554	1905	1911	1920	rBV2	1361126	2391177	75.71%	7.664%
71	13.621	1920	1922	1932	rVB2	39480	79845	2.53%	0.256%

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72	13.719	1933	1938	1939	rBV4	8412	13678	0.43%	0.044%
73	13.755	1939	1944	1947	rBV2	52926	89407	2.83%	0.287%
74	13.853	1953	1960	1968	rVB	1275579	2125952	67.31%	6.814%
75	13.938	1969	1974	1981	rVV2	71170	123396	3.91%	0.395%
76	14.011	1981	1986	1987	rVV	28011	36219	1.15%	0.116%
77	14.036	1988	1990	1994	rVV	37038	59654	1.89%	0.191%
78	14.097	1996	2000	2005	rVB	59228	95452	3.02%	0.306%
79	14.200	2013	2017	2028	rBV5	27889	64652	2.05%	0.207%
80	14.334	2034	2039	2043	rBV2	28776	50598	1.60%	0.162%
81	14.389	2043	2048	2050	rBV	97044	149327	4.73%	0.479%
82	14.456	2055	2059	2062	rVV	105900	172121	5.45%	0.552%
83	14.493	2062	2065	2073	rVB4	73071	180829	5.73%	0.580%
84	14.688	2092	2097	2100	rBV	59674	95821	3.03%	0.307%
85	14.718	2100	2102	2108	rVB2	49572	72445	2.29%	0.232%
86	14.792	2108	2114	2122	rBV2	184178	405260	12.83%	1.299%
87	14.871	2123	2127	2134	rVB2	48322	86670	2.74%	0.278%
88	14.950	2134	2140	2148	rBV	180504	319537	10.12%	1.024%
89	15.066	2154	2159	2163	rBV4	35724	66641	2.11%	0.214%
90	15.170	2172	2176	2184	rVB3	38858	79963	2.53%	0.256%
91	15.285	2190	2195	2199	rBV	81245	160989	5.10%	0.516%
92	15.365	2202	2208	2216	rVB	900595	1566445	49.60%	5.020%
93	15.450	2217	2222	2232	rBV5	54127	187138	5.93%	0.600%
94	15.657	2251	2256	2262	rBV4	21205	43116	1.37%	0.138%
95	15.730	2264	2268	2274	rVB6	19137	36772	1.16%	0.118%
96	15.962	2301	2306	2312	rBV4	19162	47912	1.52%	0.154%
97	16.377	2365	2374	2378	rBV4	36592	103951	3.29%	0.333%
98	16.444	2378	2385	2399	rVB	496713	1082043	34.26%	3.468%
99	16.566	2401	2405	2409	rBV7	14160	29221	0.93%	0.094%
100	16.645	2409	2418	2429	rVB	1005142	2262760	71.64%	7.252%

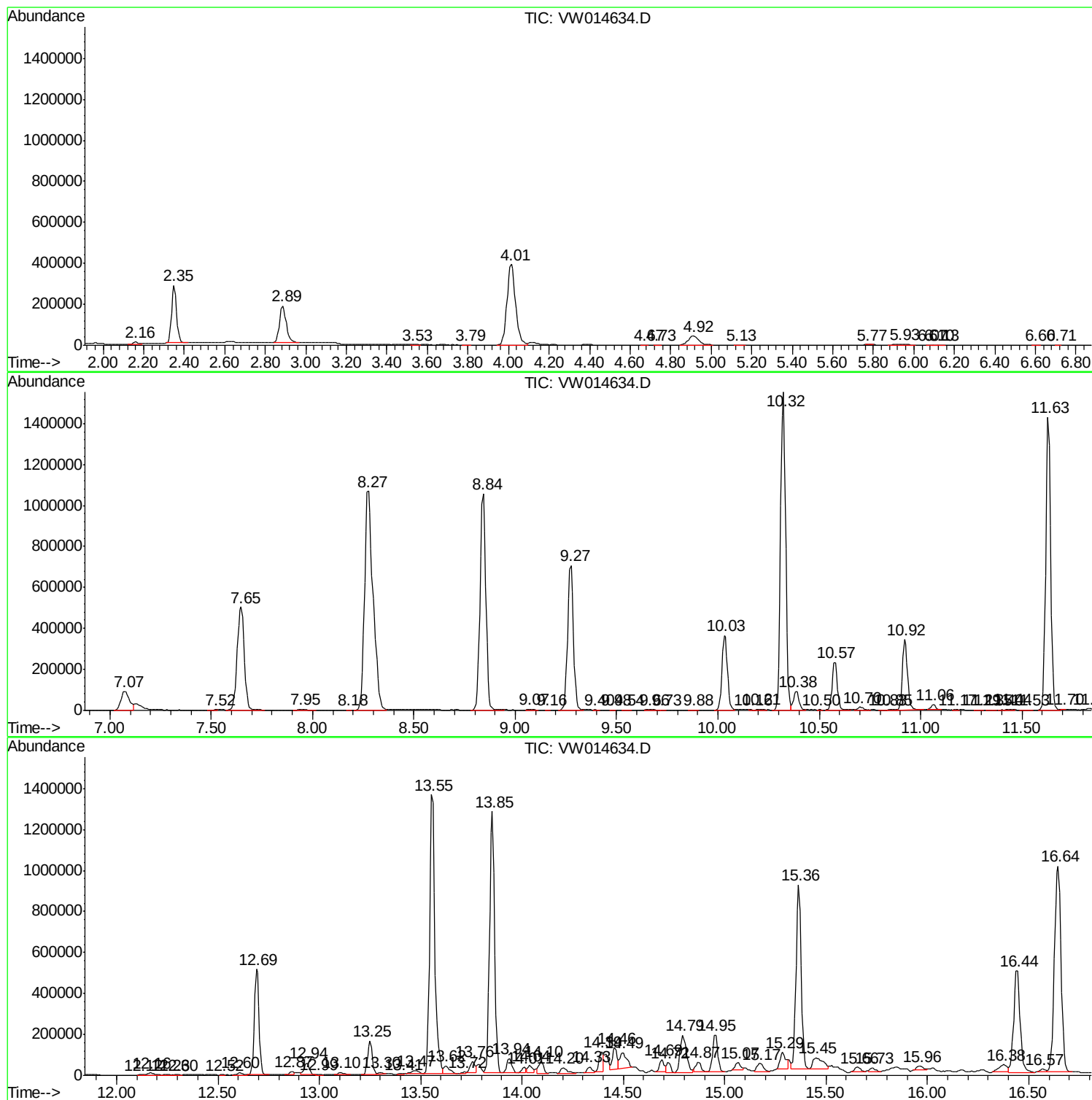
Sum of corrected areas: 31201462

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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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



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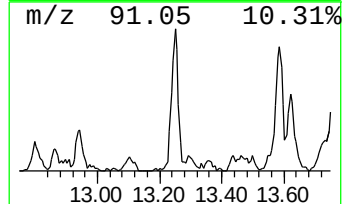
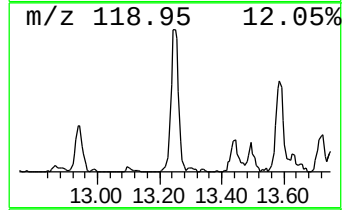
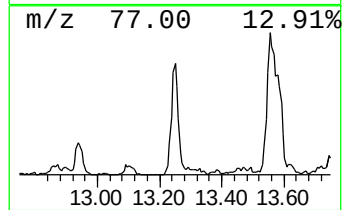
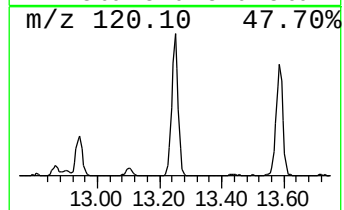
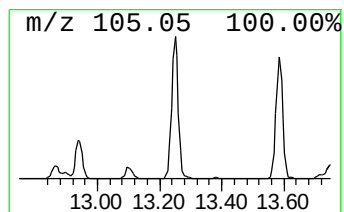
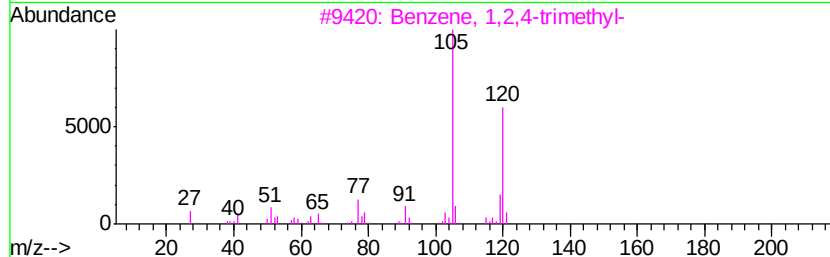
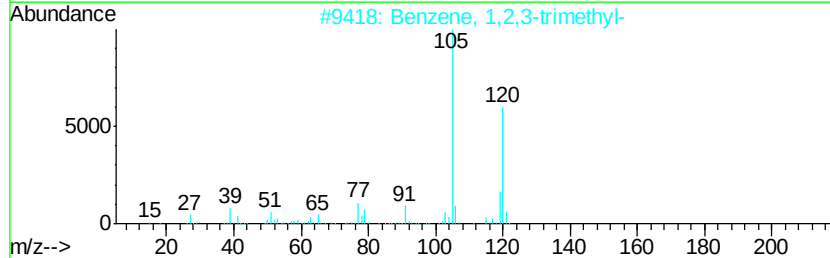
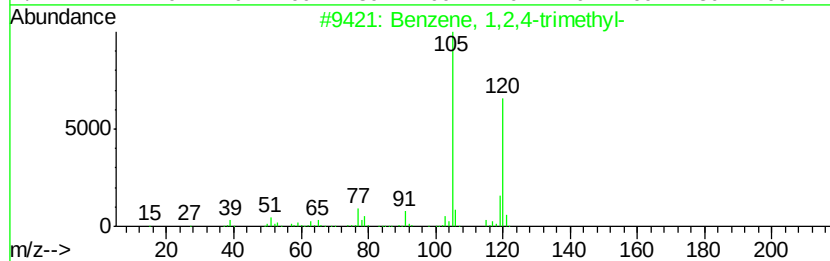
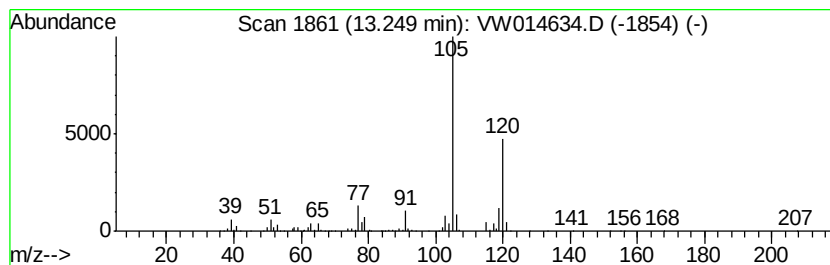
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.25	2.74 ug/L	262305	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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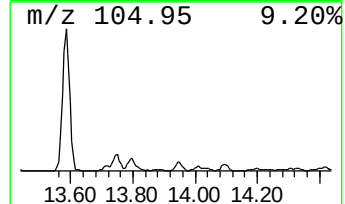
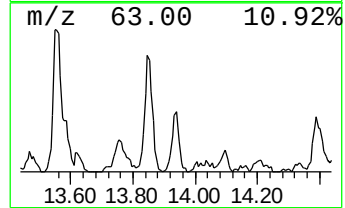
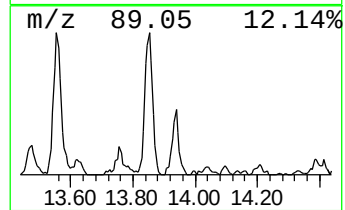
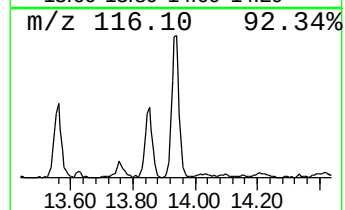
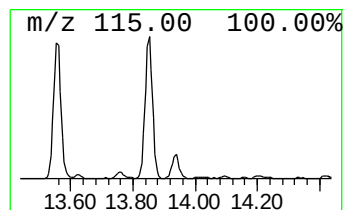
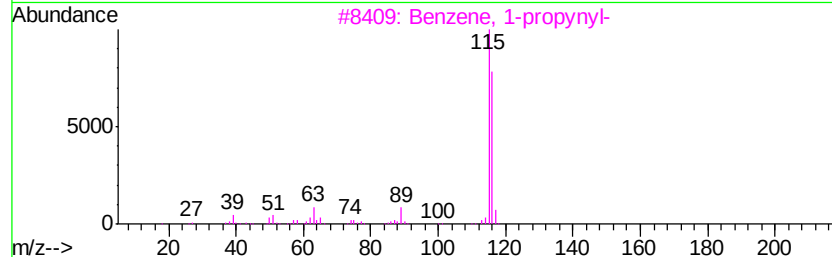
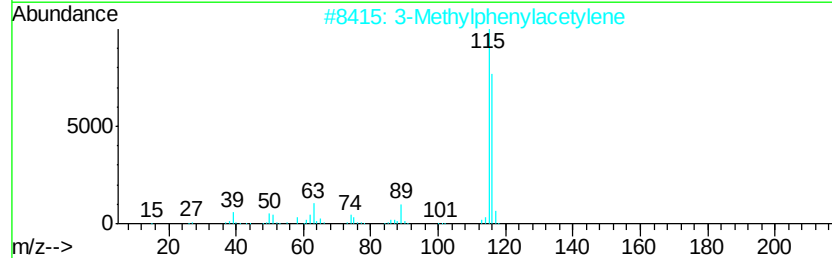
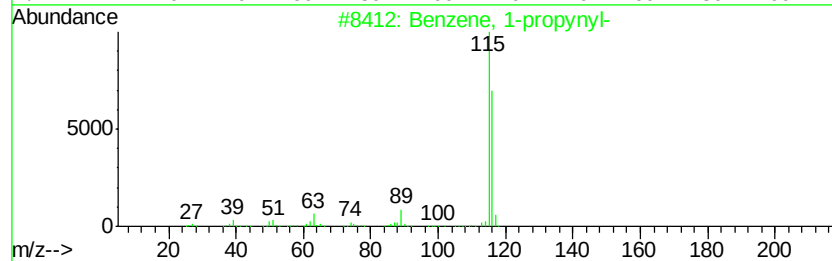
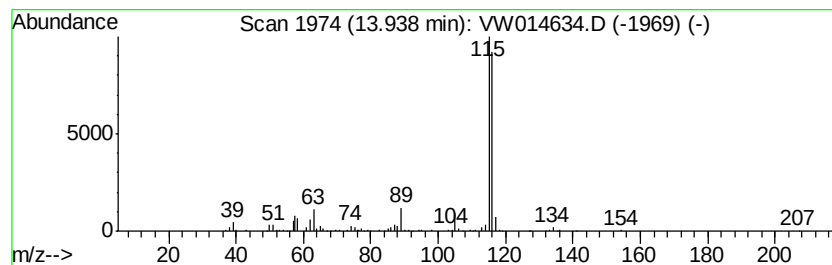
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1-propynyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	1.29 ug/L	123396	1,4-Dichlorobenzene-d4	13.55

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-propynyl-	116	C9H8	000673-32-5	93
2		3-Methylphenylacetylene	116	C9H8	000766-82-5	90
3		Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
4		Indene	116	C9H8	000095-13-6	87
5		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	83



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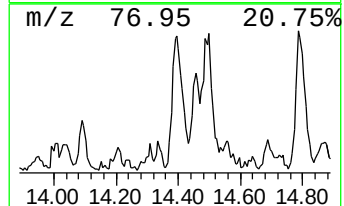
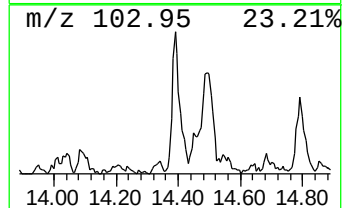
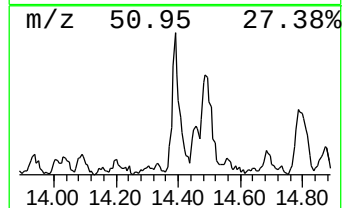
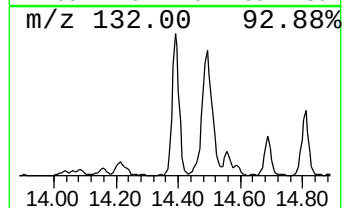
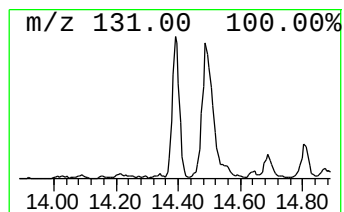
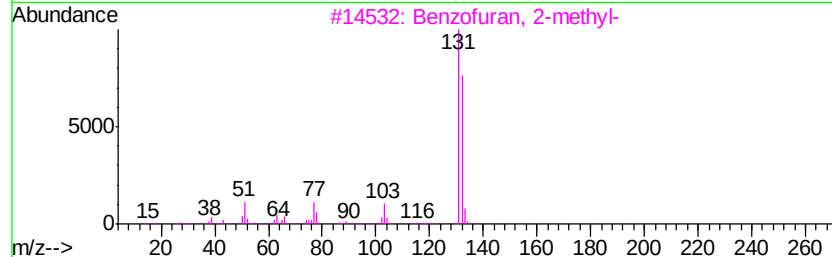
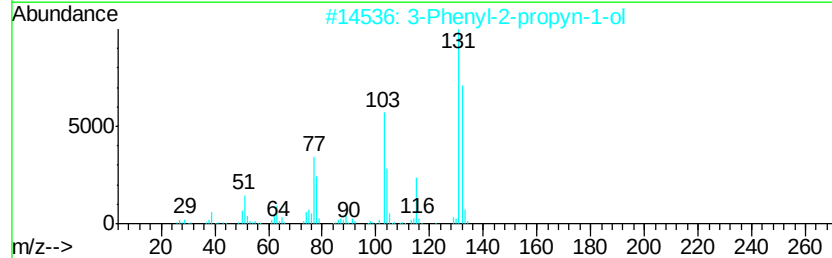
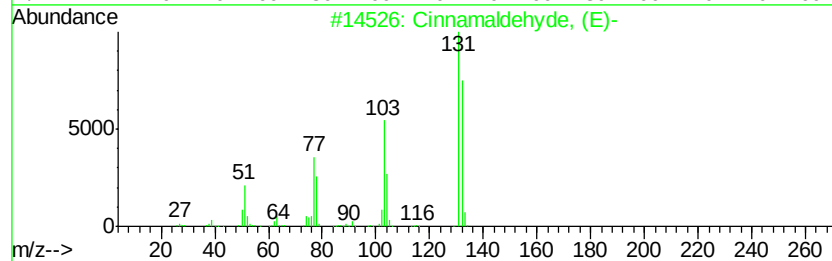
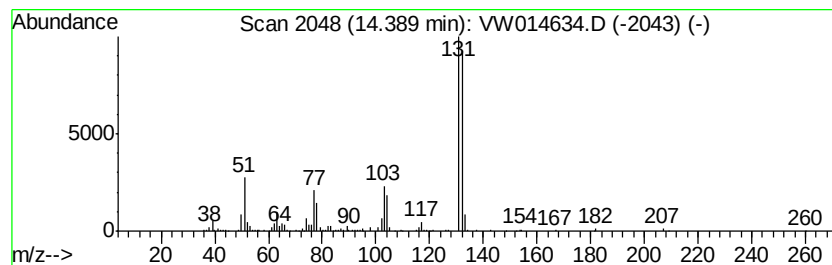
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Quant Title : VOC Analysis

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Peak Number 4 Cinnamaldehyde, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	1.56 ug/L	149327	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cinnamaldehyde, (E)-	132 C9H8O	014371-10-9 91
2		3-Phenyl-2-propyn-1-ol	132 C9H8O	001504-58-1 91
3		Benzofuran, 2-methyl-	132 C9H8O	004265-25-2 90
4		Benzofuran, 2-methyl-	132 C9H8O	004265-25-2 90
5		Cinnamaldehyde, (E)-	132 C9H8O	014371-10-9 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
Misc : 5.04G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GASH3

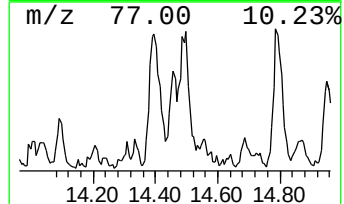
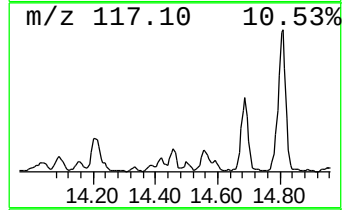
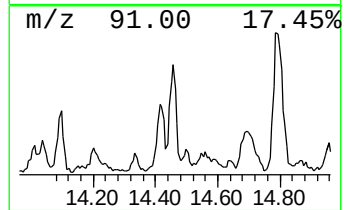
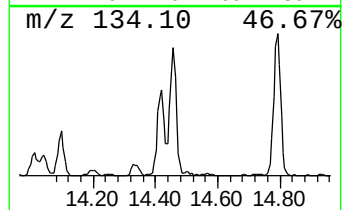
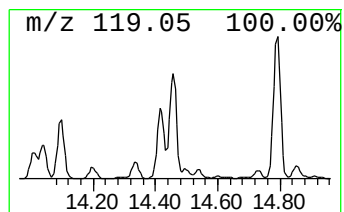
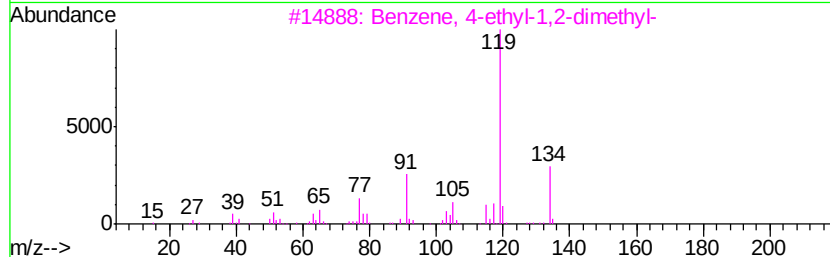
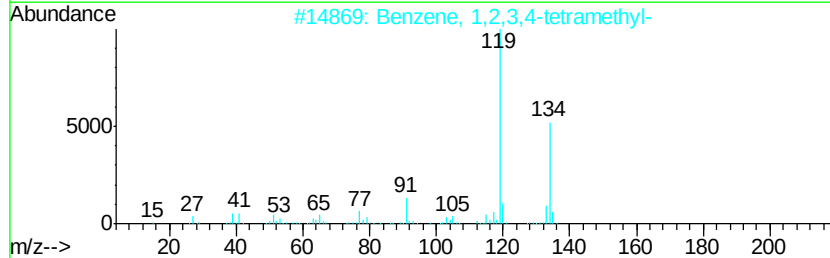
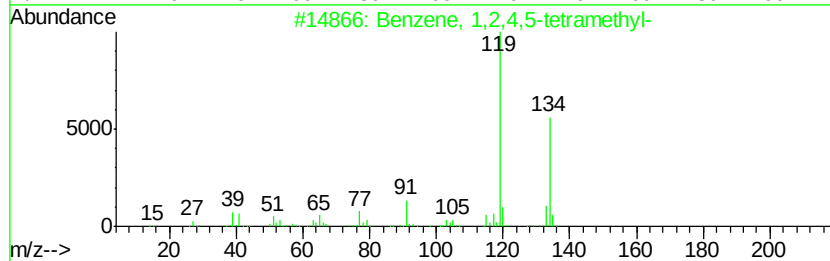
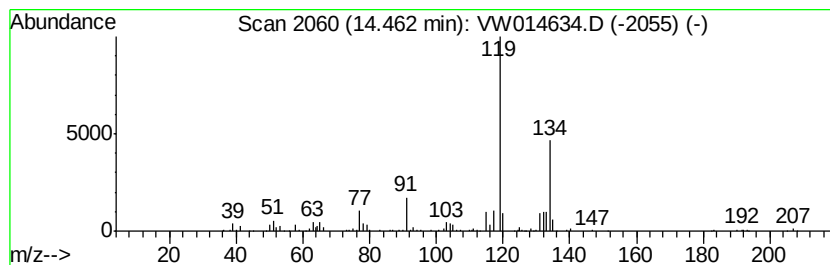
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	1.80 ug/L	172121	1,4-Dichlorobenzene-d4	13.55

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
Misc : 5.04G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

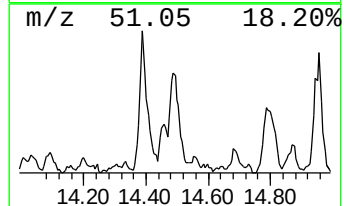
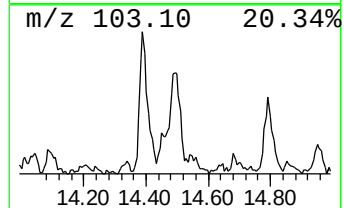
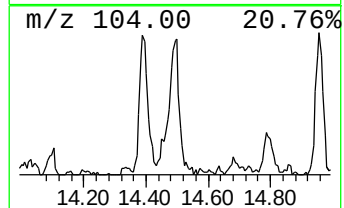
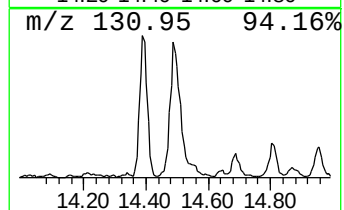
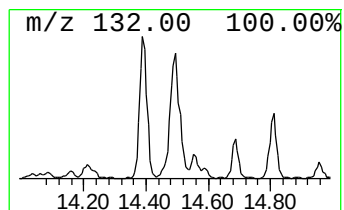
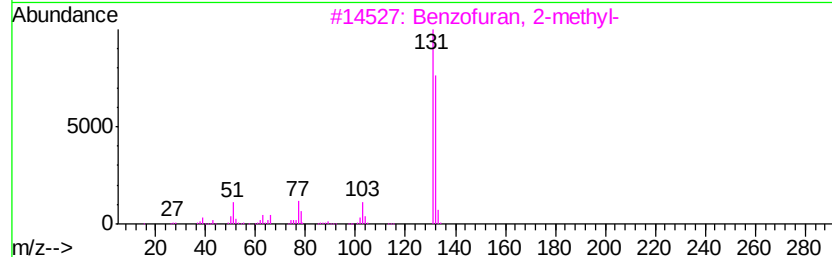
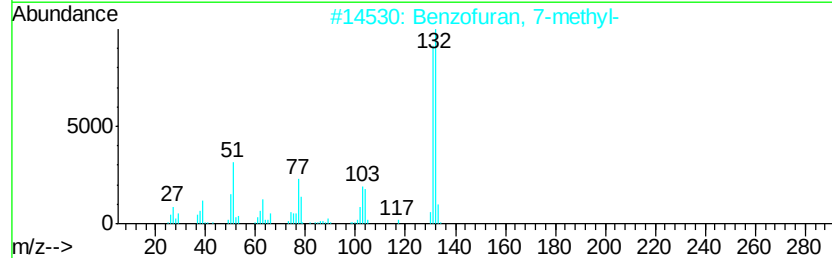
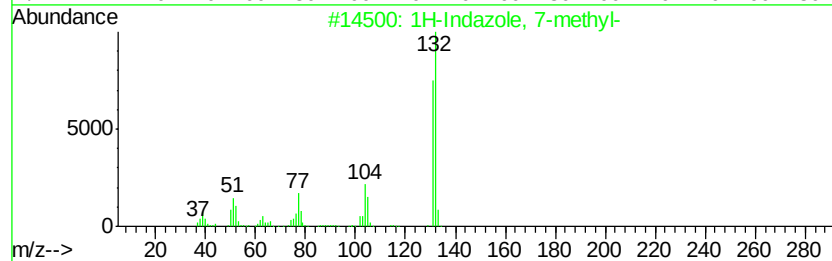
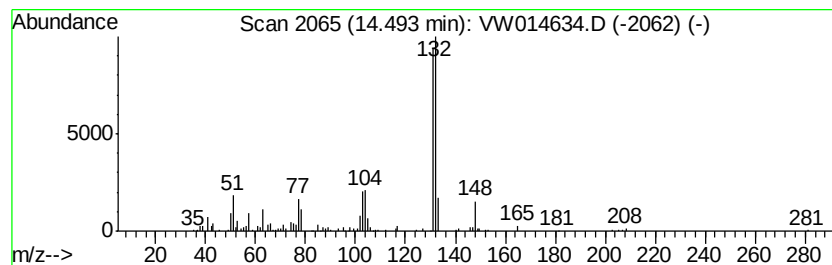
Instrument :
MSVOA_W
ClientSampleId :
GASH3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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

Peak Number 6 1H-Indazole, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	1.89 ug/L	180829	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indazole, 7-methyl-	132 C8H8N2	1000316-00-7 87
2		Benzofuran, 7-methyl-	132 C9H8O	017059-52-8 87
3		Benzofuran, 2-methyl-	132 C9H8O	004265-25-2 87
4		1H-Indazole, 3-methyl-	132 C8H8N2	1000316-00-2 83
5		Benzofuran, 2-methyl-	132 C9H8O	004265-25-2 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
Misc : 5.04G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

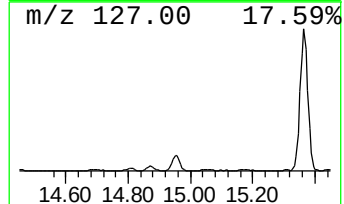
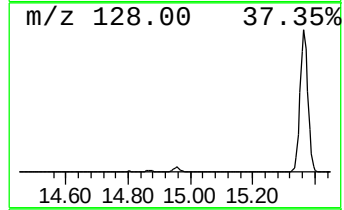
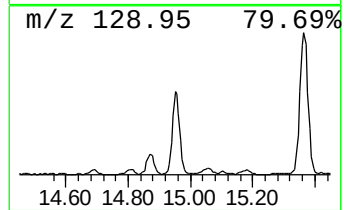
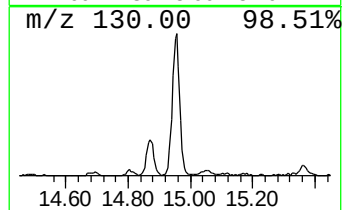
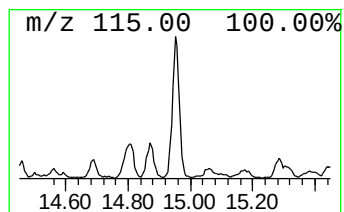
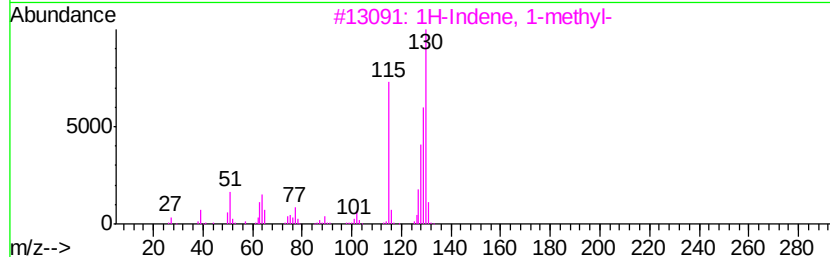
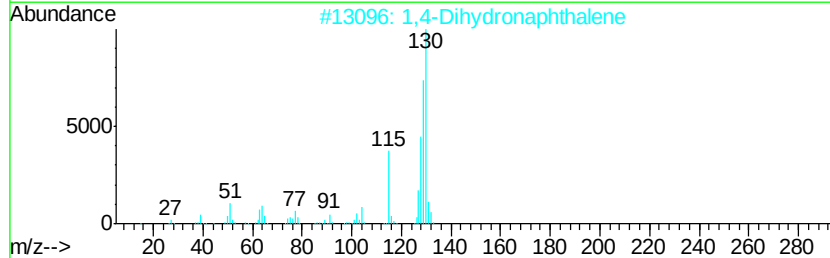
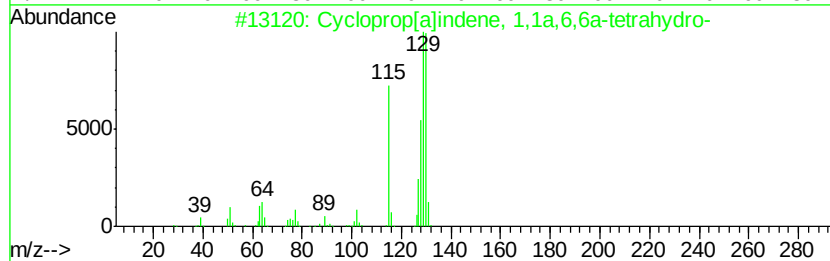
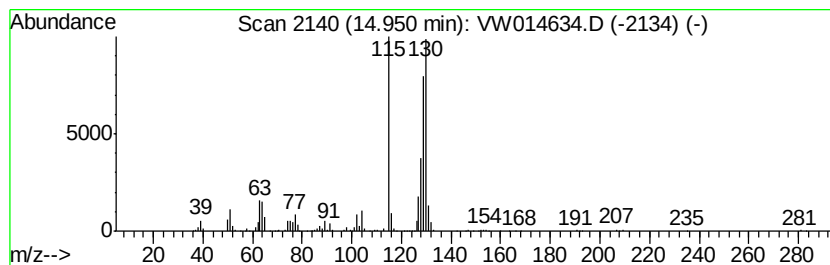
Instrument :
MSVOA_W
ClientSampleId :
GASH3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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

Peak Number 8 Cycloprop[a]indene, 1,1a,6,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	3.34 ug/L	319537	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cycloprop[a]indene, 1,1a,6,6a-te...	130 C10H10	015677-15-3 93
2		1,4-Dihydronaphthalene	130 C10H10	000612-17-9 93
3		1H-Indene, 1-methyl-	130 C10H10	000767-59-9 93
4		2-Methylindene	130 C10H10	002177-47-1 93
5		1H-Indene, 1-methyl-	130 C10H10	000767-59-9 93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
Misc : 5.04G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

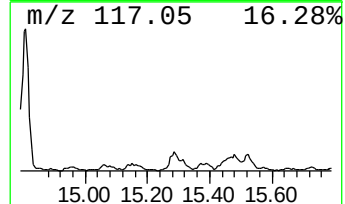
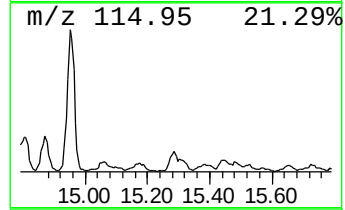
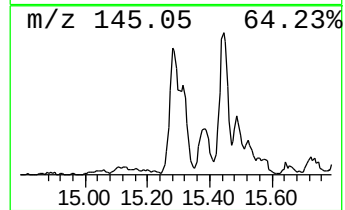
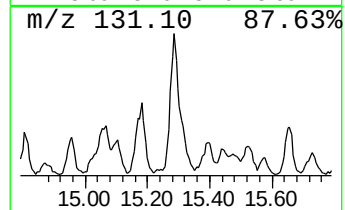
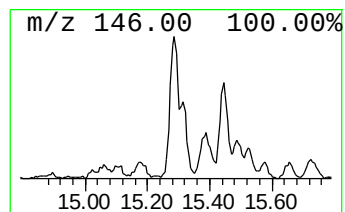
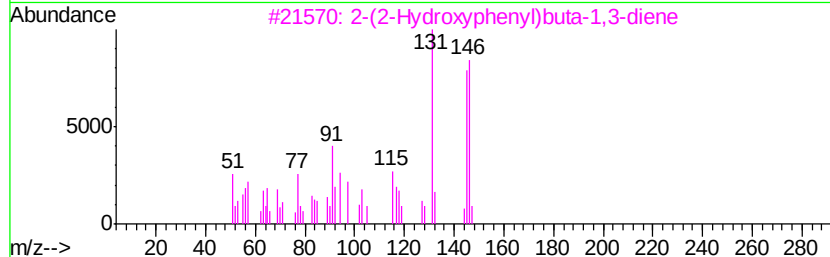
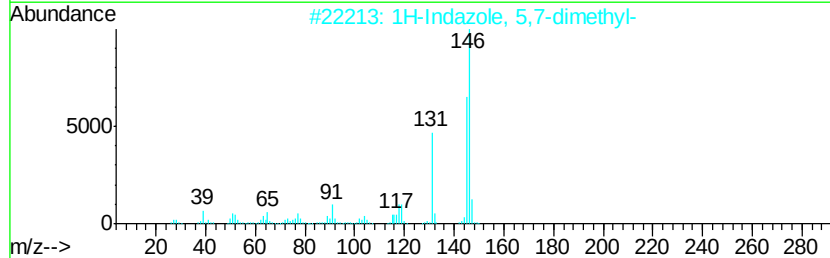
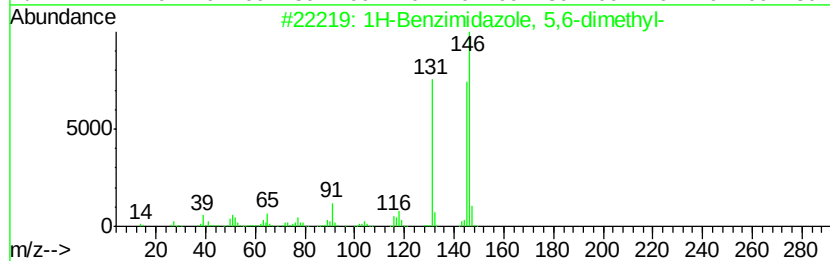
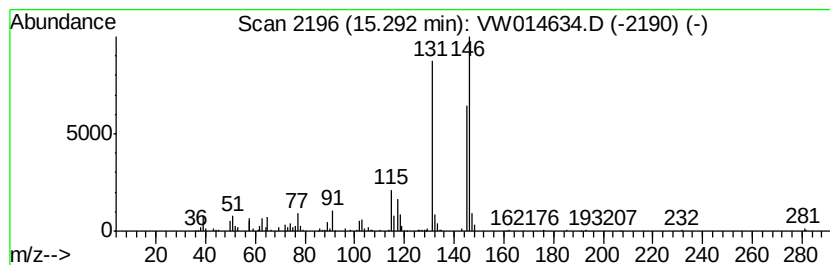
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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

Peak Number 9 1H-Benzimidazole, 5,6-dimet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.29	1.68 ug/L	160989	1,4-Dichlorobenzene-d4	13.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	90
2	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	87
3	2-(2-Hydroxyphenyl)buta-1,3-diene	146	C10H10O	038865-47-3	86
4	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	80
5	Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
Misc : 5.04G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

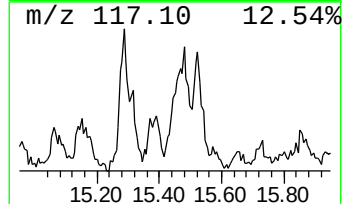
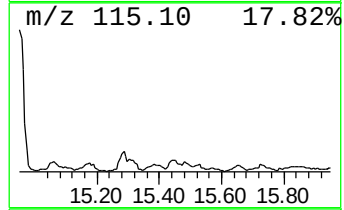
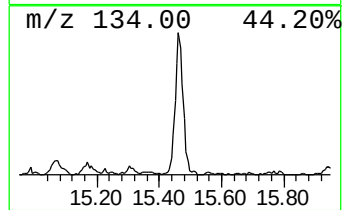
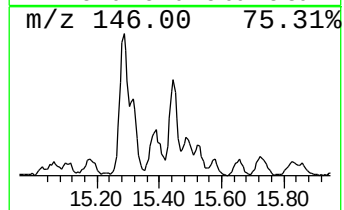
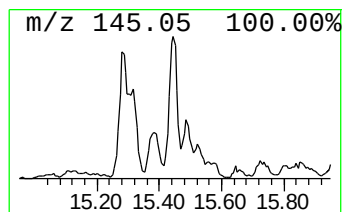
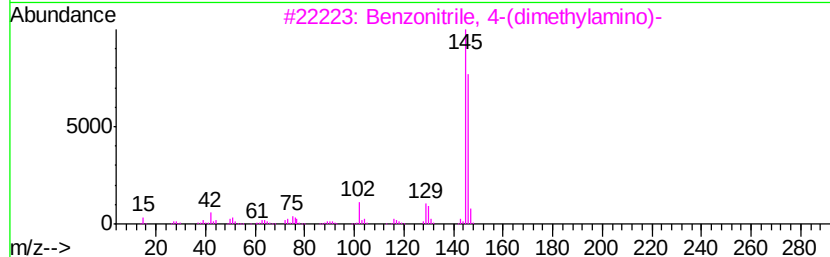
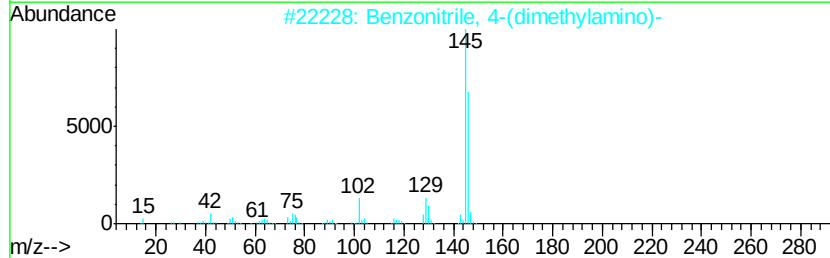
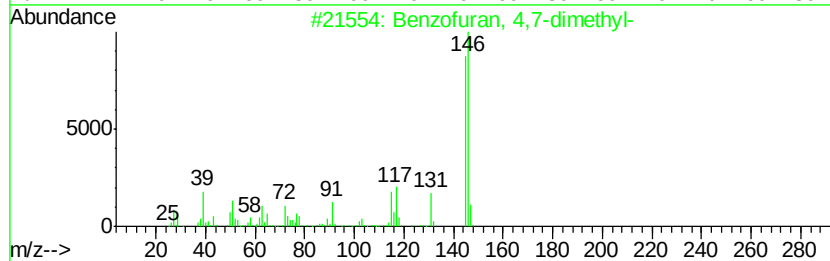
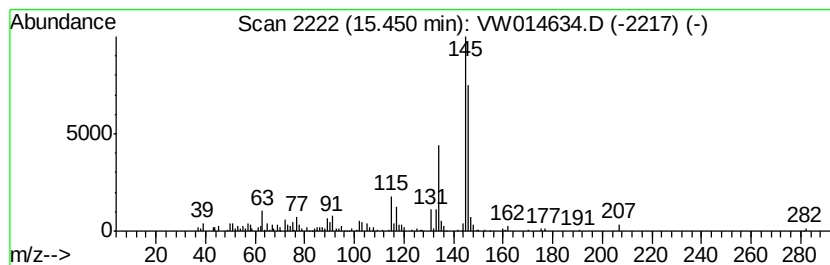
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.M
Quant Title : VOC Analysis

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

Peak Number 11 Benzofuran, 4,7-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	1.96 ug/L	187138	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzofuran, 4,7-dimethyl-	146 C10H10O	028715-26-6 55
2		Benzonitrile, 4-(dimethylamino)-	146 C9H10N2	001197-19-9 53
3		Benzonitrile, 4-(dimethylamino)-	146 C9H10N2	001197-19-9 50
4		2-Thiophenecarboxaldehyde, 5-chl...	146 C5H3ClOS	007283-96-7 50
5		1H-Benzimidazole, 2-ethyl-	146 C9H10N2	001848-84-6 50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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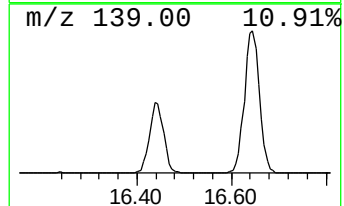
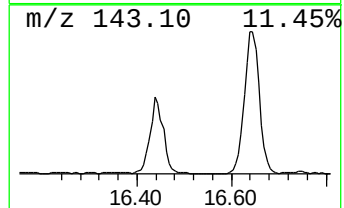
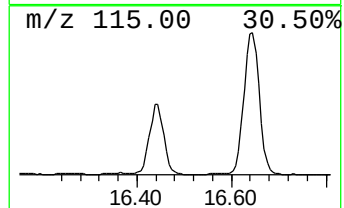
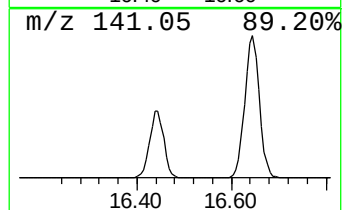
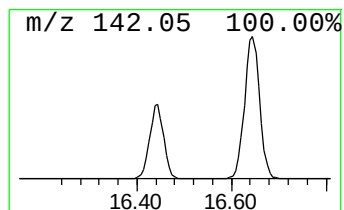
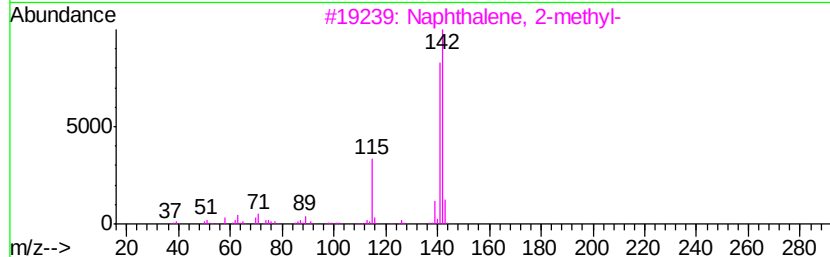
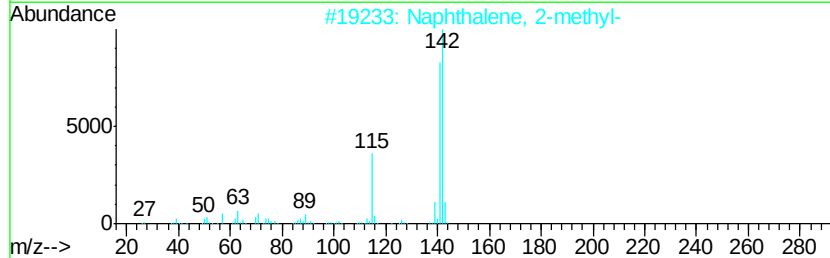
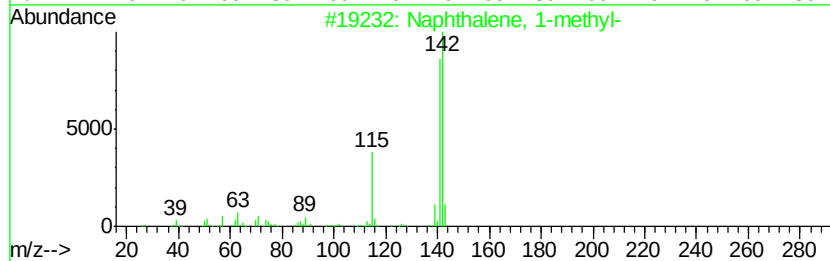
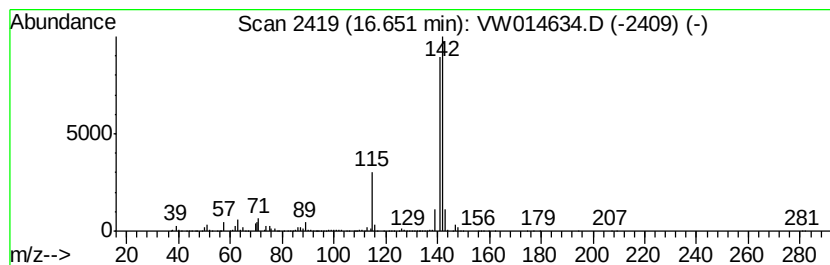
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Quant Title : VOC Analysis

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

Peak Number 13 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	23.66 ug/L	2262760	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW011420\
Data File : VW014634.D
Acq On : 14 Jan 2020 15:47
Operator : SY/VA
Sample : L1109-14
Misc : 5.04G/10ML/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GASH3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM122719S.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
Benzene, 1,2,4-tr...	13.25	2.7	ug/L	262305	3	13.55	2391180	25.0
Benzene, 1-propynyl-	13.94	1.3	ug/L	123396	3	13.55	2391180	25.0
Cinnamaldehyde, (E)-	14.39	1.6	ug/L	149327	3	13.55	2391180	25.0
Benzene, 1,2,4,5-...	14.46	1.8	ug/L	172121	3	13.55	2391180	25.0
1H-Indazole, 7-me...	14.49	1.9	ug/L	180829	3	13.55	2391180	25.0
Cycloprop[a]inden...	14.95	3.3	ug/L	319537	3	13.55	2391180	25.0
1H-Benzimidazole,...	15.29	1.7	ug/L	160989	3	13.55	2391180	25.0
Benzofuran, 4,7-d...	15.45	2.0	ug/L	187138	3	13.55	2391180	25.0
Naphthalene, 1-me...	16.65	23.7	ug/L	2262760	3	13.55	2391180	25.0