

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
 Data File : VV024258.D
 Acq On : 31 Dec 2021 02:34
 Operator : SY/MD
 Sample : M5213-06DL 200X
 Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGLD8DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

Signal : TIC: VV024258.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	77	83	99	rVB	140582	144286	14.88%	1.431%
2	1.568	157	164	177	rBV	110991	126132	13.01%	1.251%
3	2.111	322	333	350	rBV	244358	361212	37.25%	3.583%
4	2.301	385	392	395	rBV2	1087	1265	0.13%	0.013%
5	2.429	429	432	437	rBV2	1037	1137	0.12%	0.011%
6	2.709	517	519	522	rBV	571	396	0.04%	0.004%
7	2.741	526	529	533	rVV2	619	473	0.05%	0.005%
8	2.834	554	558	561	rBV	497	456	0.05%	0.005%
9	2.854	561	564	566	rBV	595	359	0.04%	0.004%
10	2.889	572	575	579	rBV	637	477	0.05%	0.005%
11	3.233	679	682	686	rVB	590	436	0.04%	0.004%
12	3.252	686	688	691	rBV	347	220	0.02%	0.002%
13	3.323	707	710	715	rVB2	492	454	0.05%	0.005%
14	3.371	721	725	729	rBV	706	679	0.07%	0.007%
15	3.442	744	747	751	rVB	598	405	0.04%	0.004%
16	3.516	766	770	773	rBV2	485	428	0.04%	0.004%
17	3.648	805	811	816	rBV2	886	972	0.10%	0.010%
18	3.683	817	822	825	rVV	643	634	0.07%	0.006%
19	3.706	825	829	833	rVV	461	418	0.04%	0.004%
20	3.754	840	844	848	rBV2	762	709	0.07%	0.007%
21	3.802	856	859	861	rBV	699	479	0.05%	0.005%
22	3.841	866	871	873	rBV	304	250	0.03%	0.002%
23	3.889	875	886	911	rBV	54315	163829	16.89%	1.625%
24	4.236	992	994	996	rBV	379	225	0.02%	0.002%
25	4.352	1013	1030	1056	rBV	152175	372962	38.46%	3.700%
26	4.571	1095	1098	1102	rBV2	551	448	0.05%	0.004%
27	4.703	1135	1139	1145	rBV	828	734	0.08%	0.007%
28	4.741	1148	1151	1156	rVB2	409	318	0.03%	0.003%
29	4.799	1162	1169	1170	rBV	556	533	0.05%	0.005%
30	4.892	1195	1198	1199	rBV	282	162	0.02%	0.002%
31	4.966	1215	1221	1225	rBV2	448	464	0.05%	0.005%
32	5.047	1229	1246	1278	rVV2	340109	887793	91.55%	8.807%
33	5.265	1311	1314	1317	rBV2	556	387	0.04%	0.004%
34	5.313	1327	1329	1334	rVB2	469	349	0.04%	0.003%
35	5.349	1334	1340	1343	rBV2	550	465	0.05%	0.005%
36	5.381	1347	1350	1352	rVB	437	238	0.02%	0.002%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Peak Location: TOP

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

37	5.407	1352	1358	1361	rBV2	631	601	0.06%	0.006%
38	5.461	1371	1375	1378	rVV	546	412	0.04%	0.004%
39	5.555	1401	1404	1407	rBV	275	176	0.02%	0.002%
40	5.616	1411	1423	1450	rBV	272150	541564	55.85%	5.372%
41	5.857	1495	1498	1500	rBV2	432	269	0.03%	0.003%
42	5.905	1505	1513	1527	rVV2	7873	15893	1.64%	0.158%
43	6.014	1540	1547	1548	rBV	550	636	0.07%	0.006%
44	6.069	1551	1564	1587	rVV	188779	382680	39.46%	3.796%
45	6.207	1604	1607	1611	rBV3	1188	1048	0.11%	0.010%
46	6.255	1616	1622	1625	rBV2	627	679	0.07%	0.007%
47	6.304	1635	1637	1643	rVB	732	532	0.05%	0.005%
48	6.342	1646	1649	1654	rBV2	609	475	0.05%	0.005%
49	6.432	1675	1677	1682	rVB	409	275	0.03%	0.003%
50	6.468	1682	1688	1690	rBV	559	541	0.06%	0.005%
51	6.571	1714	1720	1722	rBV2	571	623	0.06%	0.006%
52	6.603	1727	1730	1732	rBV	487	364	0.04%	0.004%
53	6.641	1737	1742	1746	rBV2	526	536	0.06%	0.005%
54	6.664	1746	1749	1751	rBV2	374	164	0.02%	0.002%
55	6.680	1751	1754	1756	rBV	351	264	0.03%	0.003%
56	6.709	1760	1763	1766	rBV	562	422	0.04%	0.004%
57	6.796	1787	1790	1796	rBV2	477	455	0.05%	0.005%
58	6.821	1796	1798	1803	rVB	480	358	0.04%	0.004%
59	6.847	1803	1806	1810	rBV	464	375	0.04%	0.004%
60	6.889	1816	1819	1820	rBV	423	285	0.03%	0.003%
61	6.985	1836	1849	1870	rBV	111379	201366	20.76%	1.998%
62	7.136	1894	1896	1899	rBV2	307	227	0.02%	0.002%
63	7.268	1934	1937	1940	rBV3	474	353	0.04%	0.004%
64	7.317	1940	1952	1970	rBV	419076	723818	74.64%	7.180%
65	7.622	2037	2047	2069	rBV	81696	137971	14.23%	1.369%
66	7.757	2085	2089	2094	rBV3	649	706	0.07%	0.007%
67	7.834	2110	2113	2115	rBV2	368	254	0.03%	0.003%
68	7.870	2122	2124	2132	rVB2	522	434	0.04%	0.004%
69	7.918	2137	2139	2142	rBV2	439	240	0.02%	0.002%
70	8.037	2172	2176	2177	rBV	384	280	0.03%	0.003%
71	8.088	2182	2192	2223	rBV	209437	381780	39.37%	3.787%
72	8.278	2249	2251	2252	rBV2	1066	435	0.04%	0.004%
73	8.480	2312	2314	2318	rVB	584	380	0.04%	0.004%
74	8.503	2318	2321	2326	rBV3	953	679	0.07%	0.007%
75	8.538	2326	2332	2334	rBV	471	406	0.04%	0.004%
76	8.667	2366	2372	2384	rVB6	3978	6414	0.66%	0.064%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Peak Location: TOP

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 Title : VOC Analysis

77	8.792	2402	2411	2418	rBV4	4060	6502	0.67%	0.065%
78	8.850	2418	2429	2448	rBV	401912	650134	67.04%	6.449%
79	9.014	2469	2480	2496	rBV	161170	251398	25.92%	2.494%
80	9.136	2507	2518	2531	rBV	613846	969748	100.00%	9.620%
81	9.477	2609	2624	2631	rBV9	10221	21997	2.27%	0.218%
82	9.545	2635	2645	2662	rVB	149999	242654	25.02%	2.407%
83	9.709	2679	2696	2708	rBV7	18153	38107	3.93%	0.378%
84	10.217	2843	2854	2873	rVB	285897	461065	47.54%	4.574%
85	10.355	2889	2897	2905	rBV	37223	57372	5.92%	0.569%
86	10.448	2917	2926	2941	rBV	141066	297466	30.67%	2.951%
87	10.914	3064	3071	3093	rVB	335449	518425	53.46%	5.143%
88	11.249	3165	3175	3186	rBV	479892	725460	74.81%	7.197%
89	11.336	3195	3202	3209	rBV	95377	131033	13.51%	1.300%
90	11.529	3254	3262	3270	rBV	68593	116779	12.04%	1.158%
91	11.625	3283	3292	3311	rVB	513058	887726	91.54%	8.806%
92	12.371	3515	3524	3530	rBV9	12258	22554	2.33%	0.224%
93	12.464	3547	3553	3562	rBV2	27189	38978	4.02%	0.387%
94	13.265	3796	3802	3809	rBV7	3463	5481	0.57%	0.054%
95	13.503	3869	3876	3886	rVB2	47106	68915	7.11%	0.684%
96	13.628	3910	3915	3916	rBV2	2765	2238	0.23%	0.022%
97	13.975	4022	4023	4024	rBV	507	187	0.02%	0.002%
98	14.168	4080	4083	4085	rBV2	645	430	0.04%	0.004%
99	14.631	4218	4227	4241	rBV	43675	65365	6.74%	0.648%
100	14.818	4278	4285	4300	rVB	13830	23833	2.46%	0.236%

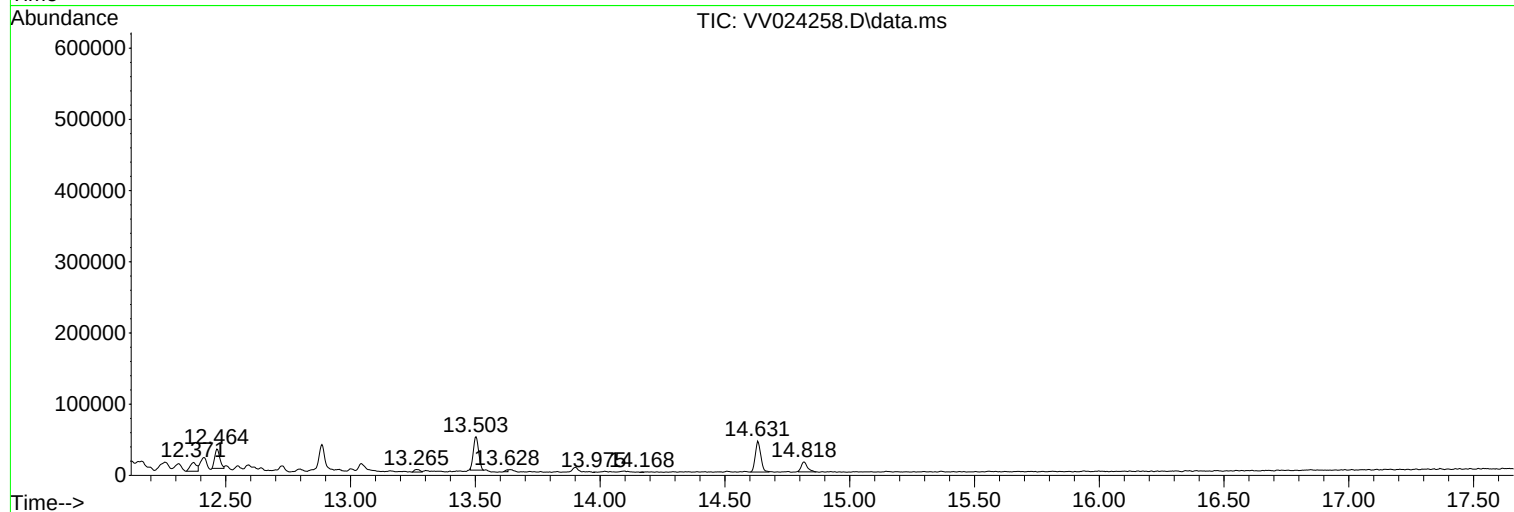
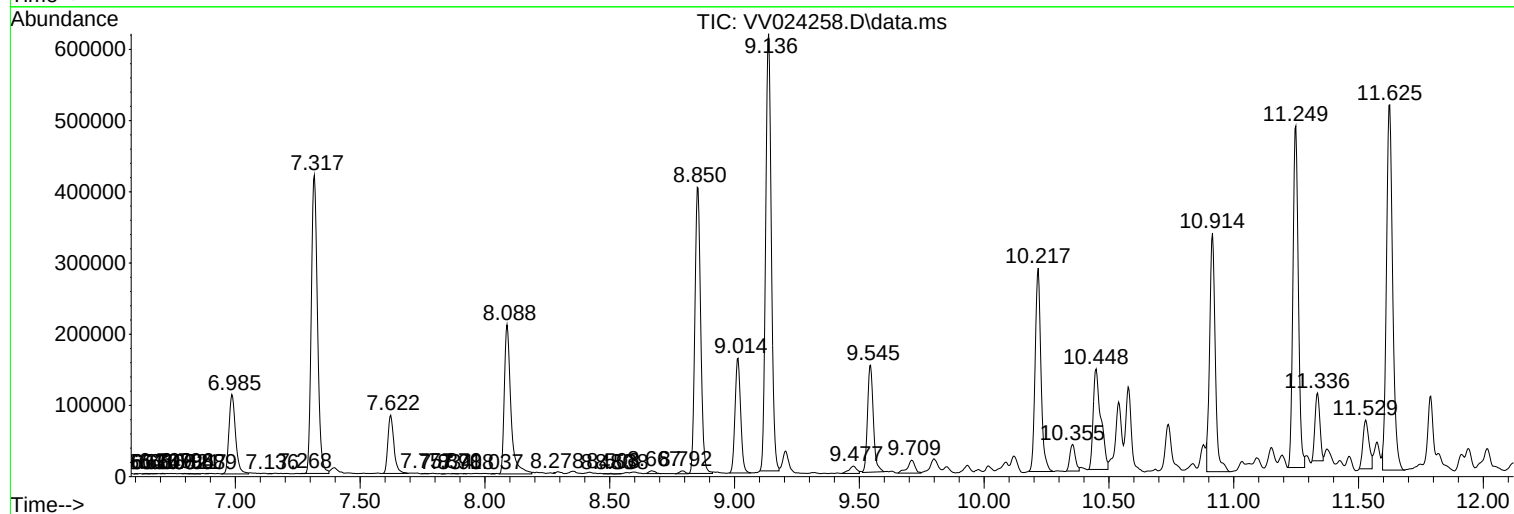
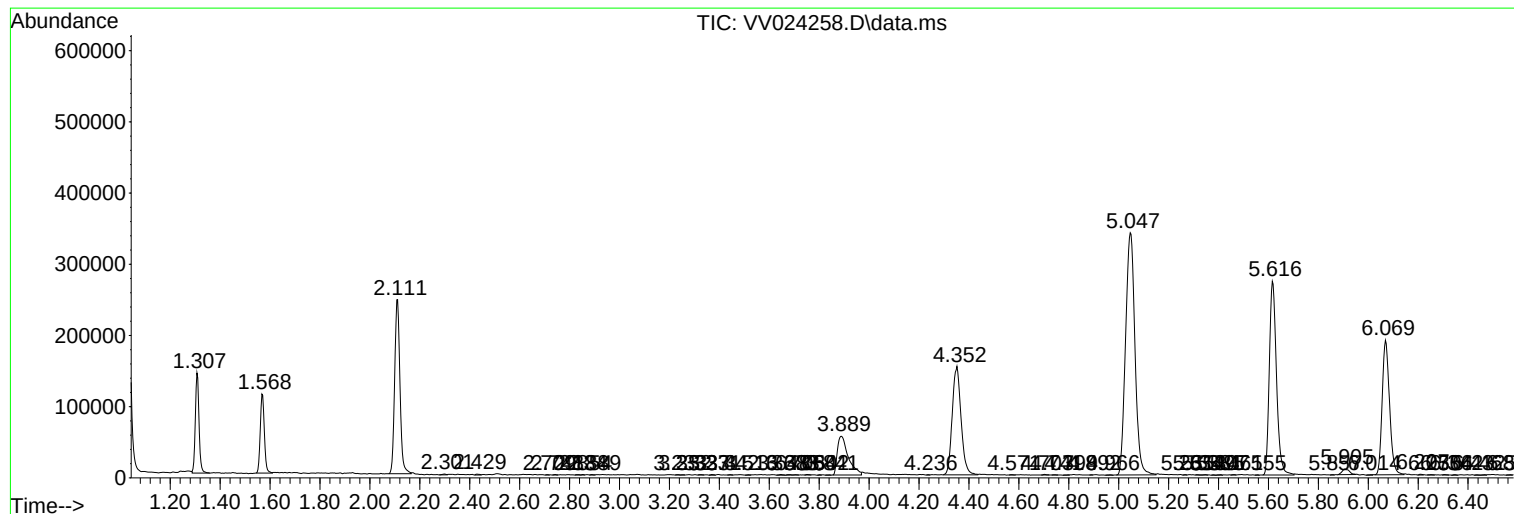
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLD8DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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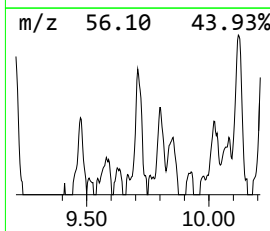
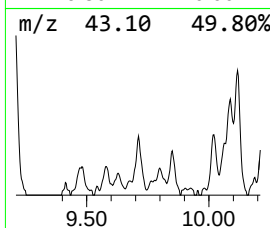
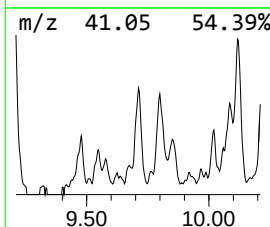
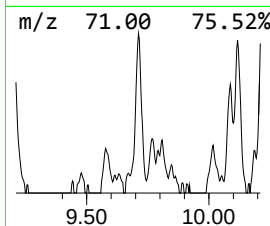
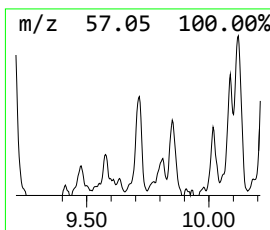
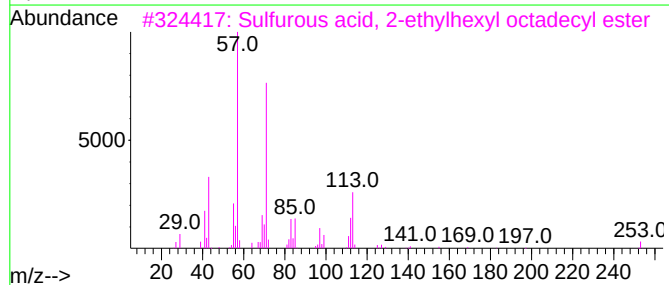
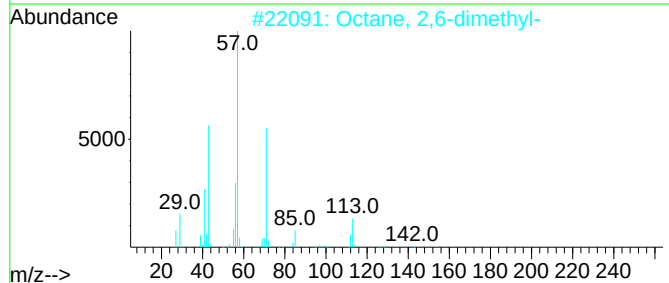
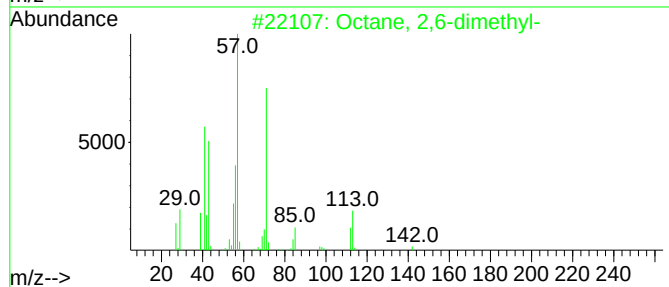
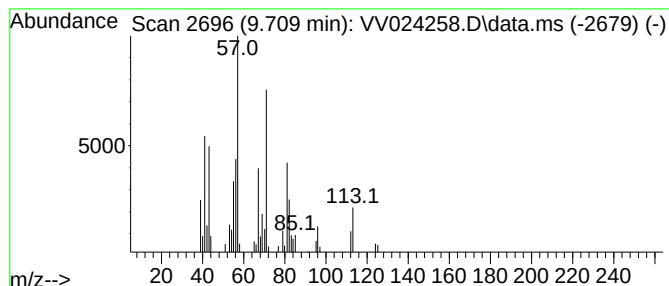
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.709	2.93 ug/L	38107	Chlorobenzene-d5	8.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	58
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	50
3			Sulfurous acid, 2-ethylhexyl oct...	446	C26H54O3S	1000309-20-1	47
4			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	47
5			1,5,7-Octatrien-3-ol, 3,7-dimethyl-	152	C10H16O	029957-43-5	43



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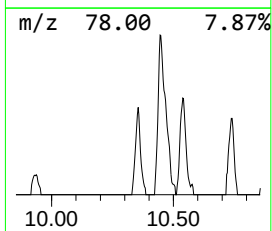
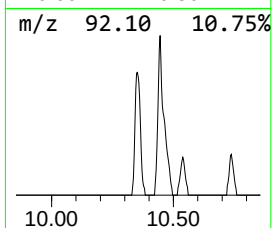
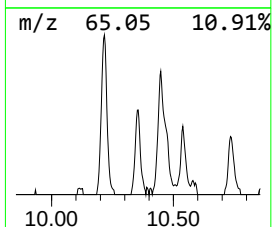
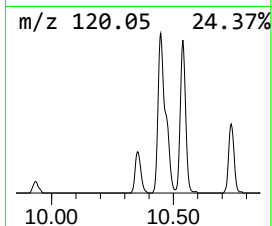
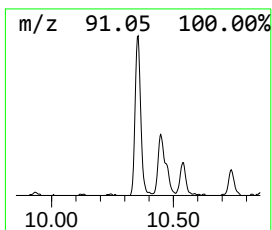
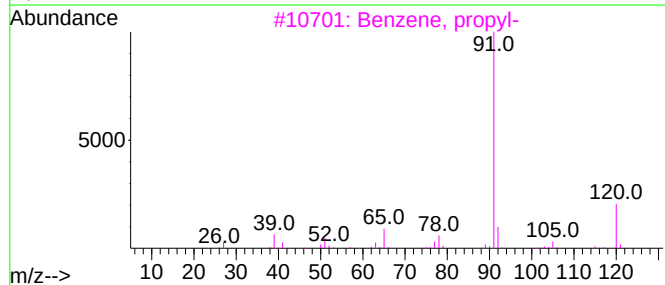
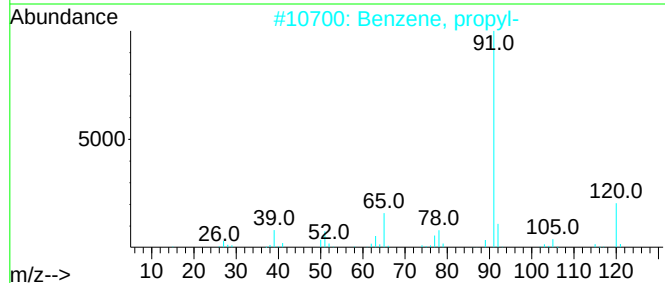
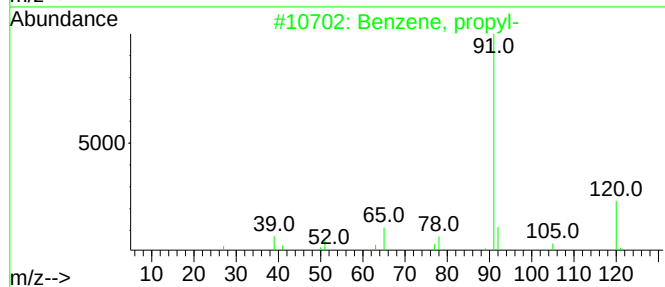
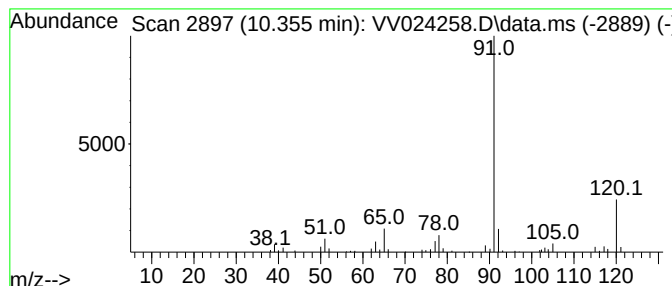
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.355	3.95 ug/L	57372	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			Benzene, propyl-	120	C9H12	000103-65-1	80
5			1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	72



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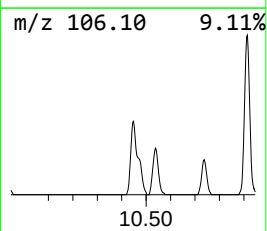
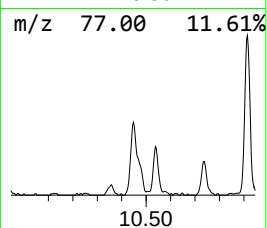
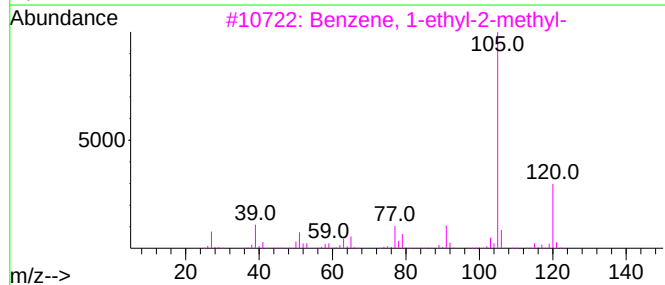
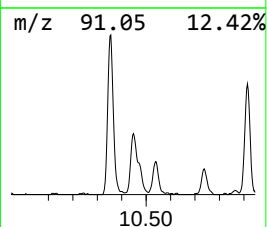
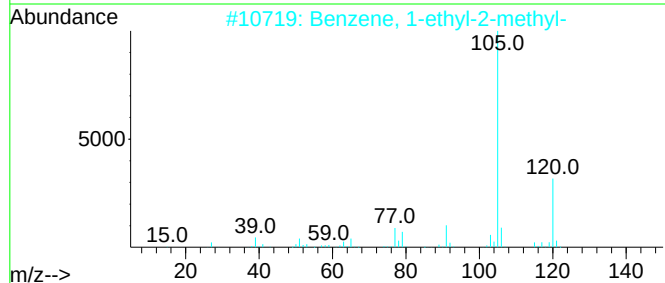
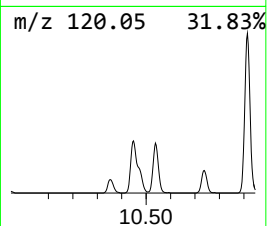
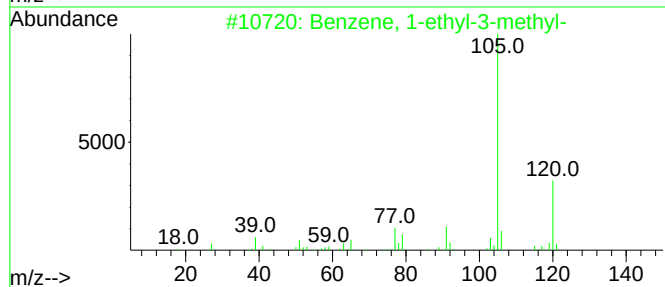
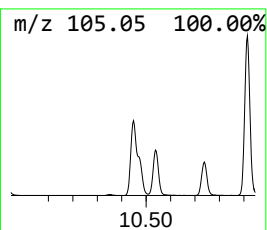
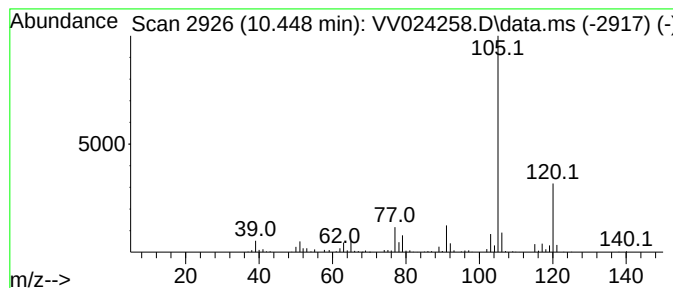
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	20.50 ug/L	297466	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
Data File : VV024258.D
Acq On : 31 Dec 2021 02:34
Operator : SY/MD
Sample : M5213-06DL 200X
Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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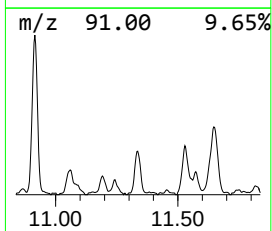
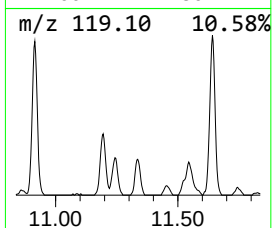
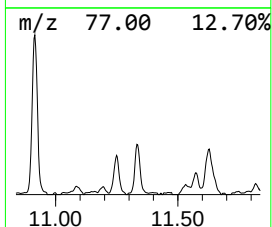
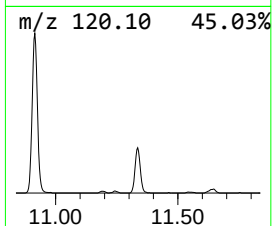
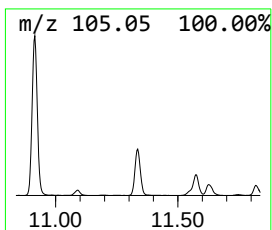
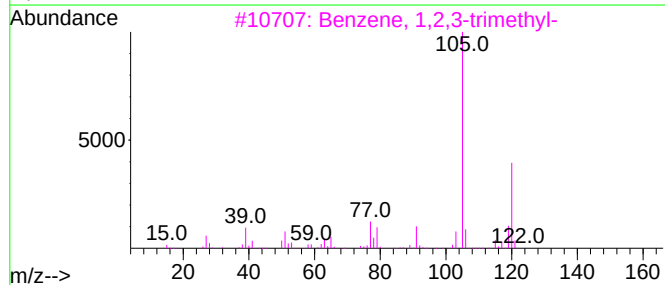
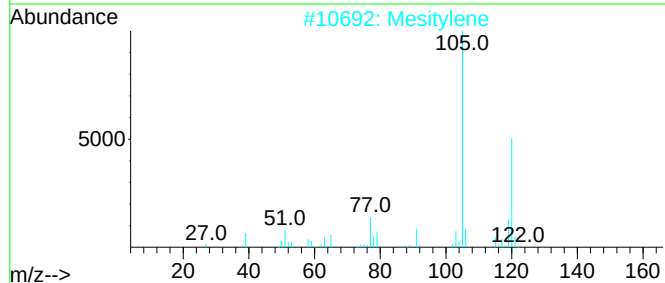
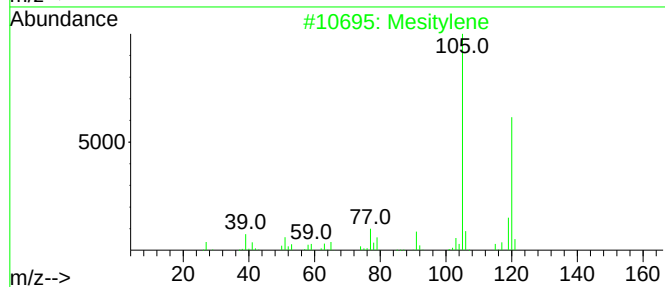
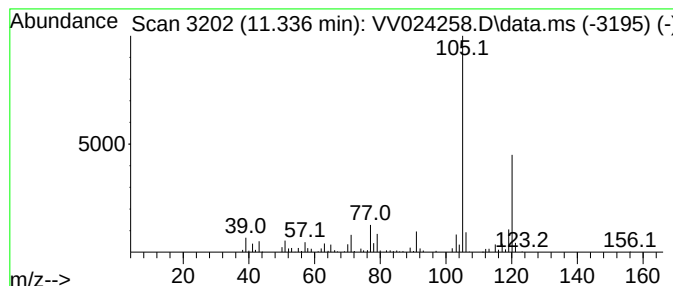
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	9.03 ug/L	131033	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	95
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
Data File : VV024258.D
Acq On : 31 Dec 2021 02:34
Operator : SY/MD
Sample : M5213-06DL 200X
Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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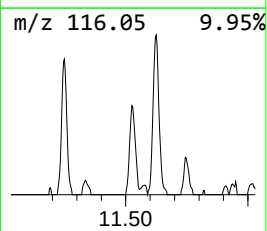
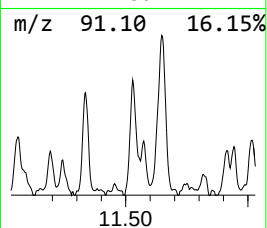
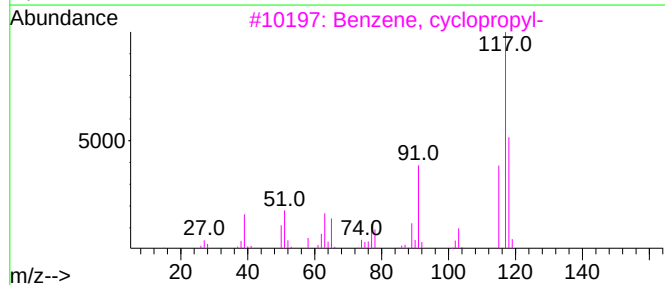
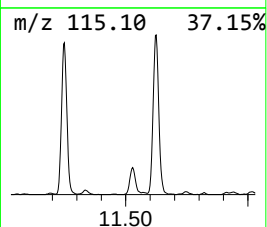
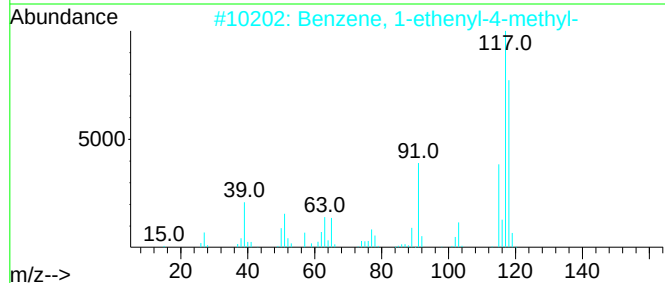
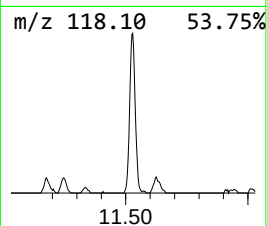
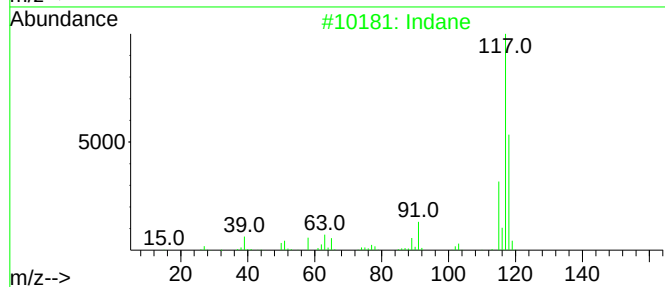
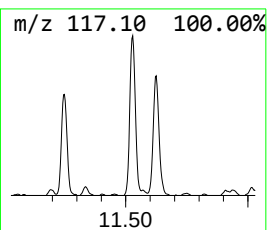
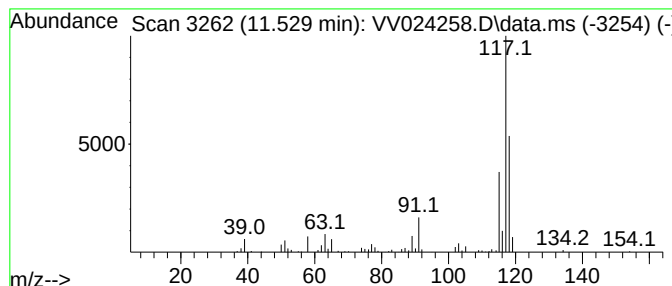
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	8.05 ug/L	116779	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	95
2			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	81
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	80
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
Data File : VV024258.D
Acq On : 31 Dec 2021 02:34
Operator : SY/MD
Sample : M5213-06DL 200X
Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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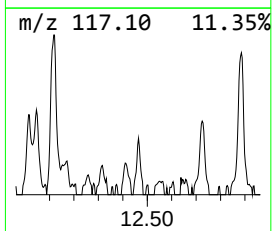
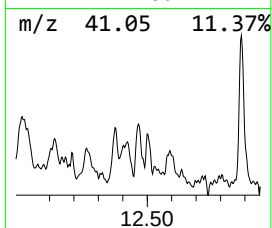
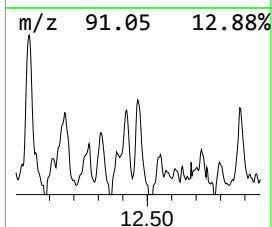
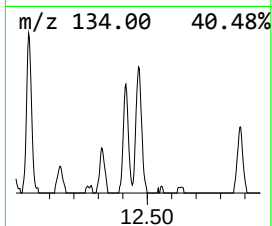
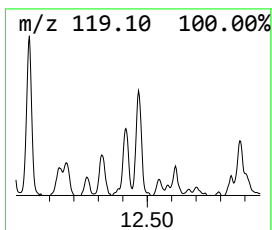
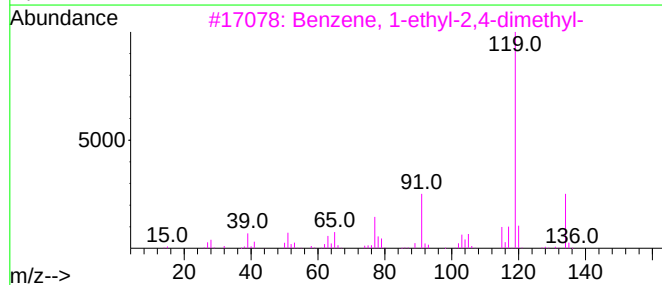
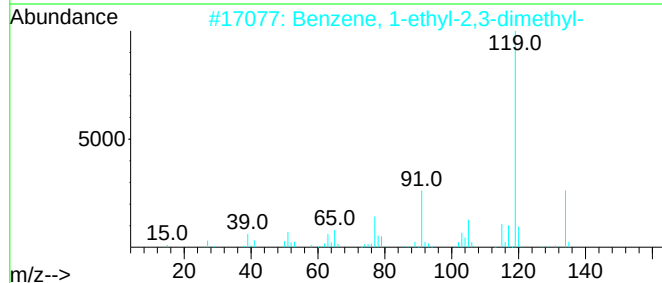
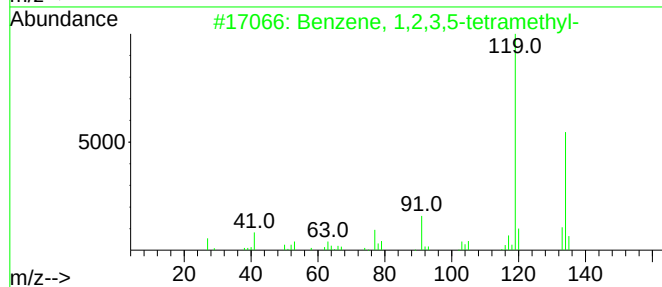
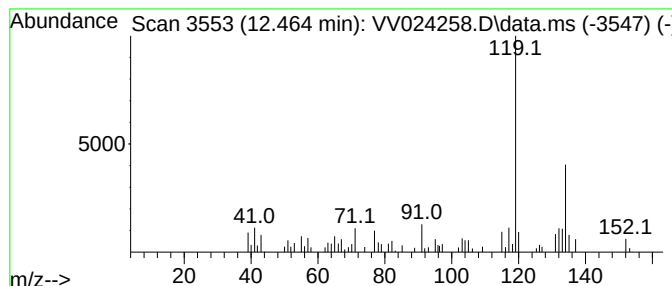
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	2.69 ug/L	38978	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
Data File : VV024258.D
Acq On : 31 Dec 2021 02:34
Operator : SY/MD
Sample : M5213-06DL 200X
Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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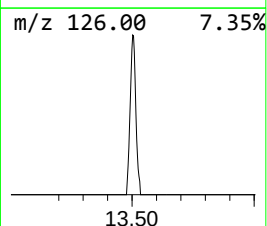
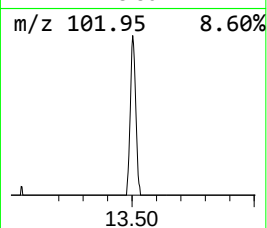
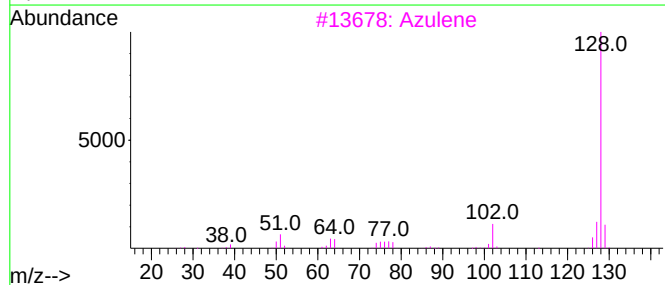
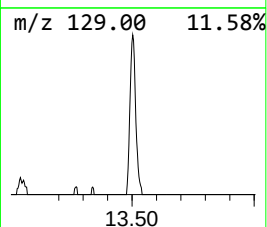
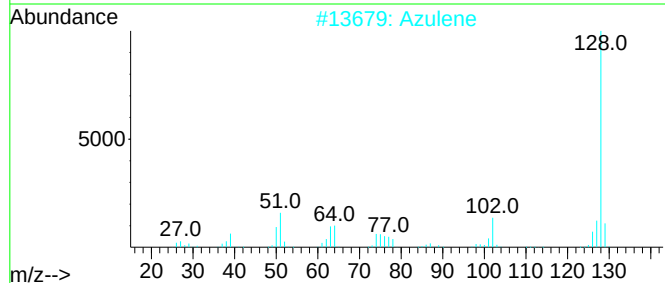
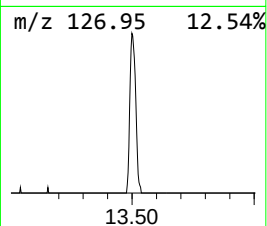
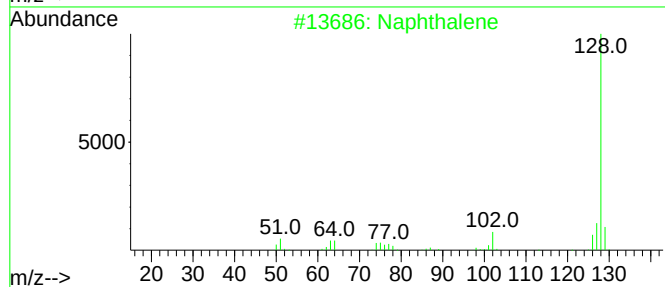
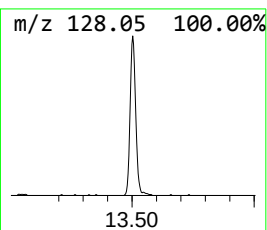
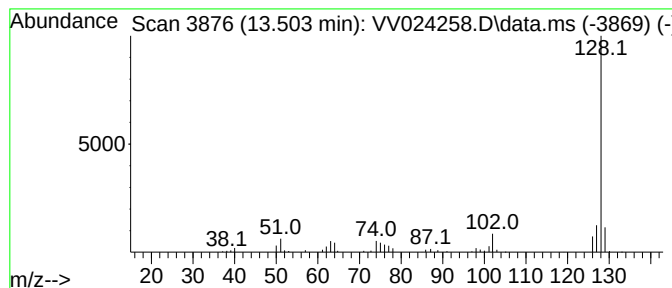
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Azulene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.503	4.75 ug/L	68915	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	91
3			Azulene	128	C10H8	000275-51-4	91
4			Naphthalene	128	C10H8	000091-20-3	90
5			Naphthalene	128	C10H8	000091-20-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
Data File : VV024258.D
Acq On : 31 Dec 2021 02:34
Operator : SY/MD
Sample : M5213-06DL 200X
Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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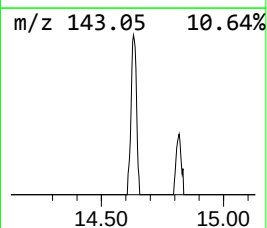
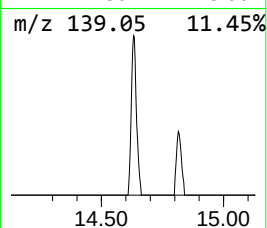
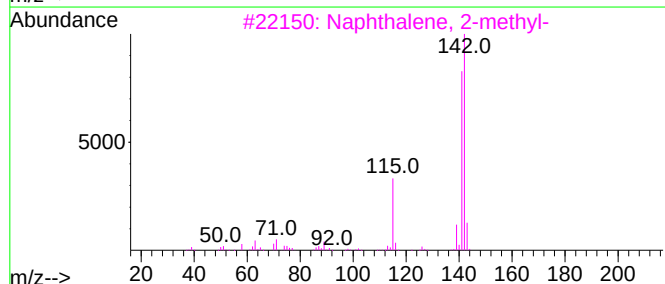
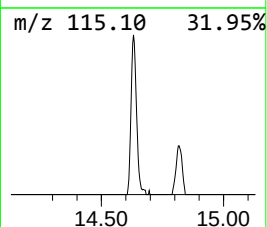
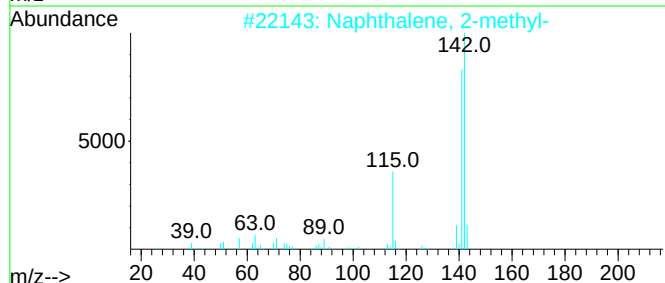
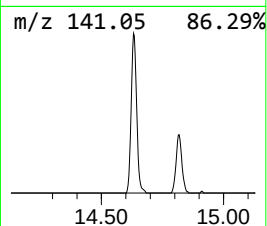
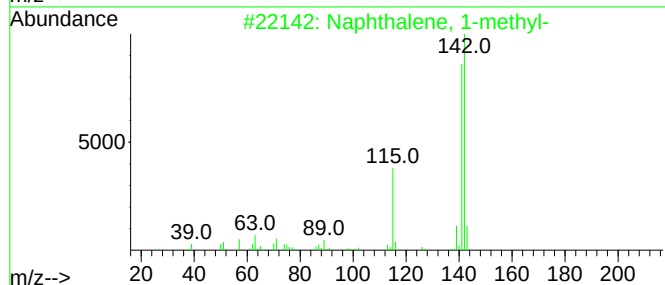
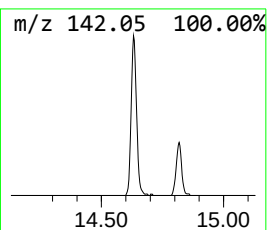
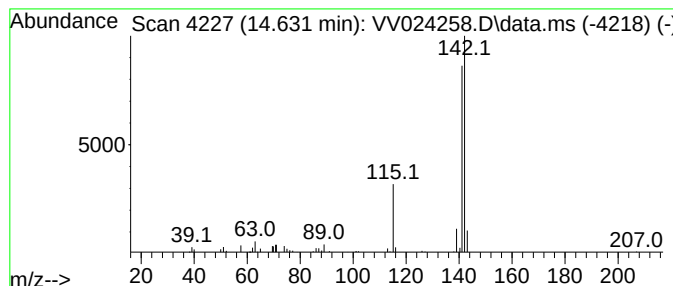
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.631	4.51 ug/L	65365	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV123121\
 Data File : VV024258.D
 Acq On : 31 Dec 2021 02:34
 Operator : SY/MD
 Sample : M5213-06DL 200X
 Misc : 6.00G/5.0mL/100uL//5.0mL/MSVOA_V/MEOH
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGLD8DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, propyl-	10.355	4.0	ug/L	57372	3	11.249	725460	50.0
Benzene, 1-ethy...	10.448	20.5	ug/L	297466	3	11.249	725460	50.0
Benzene, 1,2,3-...	11.336	9.0	ug/L	131033	3	11.249	725460	50.0
Indane	11.529	8.1	ug/L	116779	3	11.249	725460	50.0
Benzene, 1,2,3,...	12.464	2.7	ug/L	38978	3	11.249	725460	50.0
Azulene	13.503	4.8	ug/L	68915	3	11.249	725460	50.0
Naphthalene, 1-...	14.631	4.5	ug/L	65365	3	11.249	725460	50.0