

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV040320\
 Data File : VV015398.D
 Acq On : 03 Apr 2020 16:52
 Operator : SY/MD
 Sample : L2065-10
 Misc : 5.0mL/MSVOA V/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBF06

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_V\METHOD\SOMVLM040120WMA.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	1.119	5	9	13	rVB3	7946	6646	0.01%	0.002%
2	1.229	39	43	46	rBV3	5144	5227	0.01%	0.001%
3	1.315	63	70	88	rVB	414176	434542	0.43%	0.103%
4	1.579	145	152	165	rVB	329419	398252	0.39%	0.095%
5	2.122	312	321	334	rVB	717595	1023972	1.01%	0.244%
6	2.518	441	444	447	rBV5	2344	2005	0.00%	0.000%
7	2.772	516	523	525	rBV7	3018	3640	0.00%	0.001%
8	2.852	545	548	552	rVB6	2557	1871	0.00%	0.000%
9	2.881	552	557	561	rBV7	2887	3436	0.00%	0.001%
10	2.997	589	593	598	rVB8	2409	2149	0.00%	0.001%
11	3.325	692	695	700	rVB5	2355	2325	0.00%	0.001%
12	3.351	700	703	706	rBV4	2586	1982	0.00%	0.000%
13	3.939	873	886	914	rBV2	110309	465883	0.46%	0.111%
14	4.174	956	959	965	rVB8	2387	2135	0.00%	0.001%
15	4.376	1007	1022	1047	rVV2	421466	1044376	1.03%	0.248%
16	4.675	1112	1115	1119	rVB6	2657	2157	0.00%	0.001%
17	4.836	1161	1165	1169	rVV6	3351	3023	0.00%	0.001%
18	4.859	1169	1172	1178	rVB7	2498	2698	0.00%	0.001%
19	5.074	1221	1239	1275	rVV2	1045634	2636247	2.61%	0.627%
20	5.643	1402	1416	1437	rBV	701217	1385529	1.37%	0.329%
21	5.878	1485	1489	1492	rBV6	2232	2033	0.00%	0.000%
22	5.923	1494	1503	1504	rBV9	7150	6447	0.01%	0.002%
23	6.093	1538	1556	1577	rBV	563397	1164214	1.15%	0.277%
24	6.338	1628	1632	1641	rBV9	2294	3405	0.00%	0.001%
25	6.585	1705	1709	1713	rBV5	2161	2390	0.00%	0.001%
26	7.010	1827	1841	1861	rBV	356042	647587	0.64%	0.154%
27	7.209	1899	1903	1908	rVB8	2240	2428	0.00%	0.001%
28	7.338	1930	1943	1956	rBV	1340152	2278930	2.26%	0.542%
29	7.412	1956	1966	1990	rVB	500819	880985	0.87%	0.210%
30	7.502	1990	1994	1997	rBV6	2253	2054	0.00%	0.000%
31	7.540	2002	2006	2007	rBV4	3382	2162	0.00%	0.001%
32	7.643	2027	2038	2056	rBV	252020	435161	0.43%	0.103%
33	7.955	2129	2135	2137	rBV7	6749	6570	0.01%	0.002%
34	8.003	2144	2150	2157	rVB10	5800	8864	0.01%	0.002%

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Integration Parameters: LSCINT.P

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 Title : VOC Analysis

35	8.058	2161	2167	2169	rBV7	2659	2405	0.00%	0.001%
36	8.116	2173	2185	2217	rBV2	688674	1330178	1.32%	0.316%
37	8.592	2329	2333	2338	rBV7	2487	2373	0.00%	0.001%
38	8.707	2363	2369	2372	rBV7	2286	2691	0.00%	0.001%
39	8.875	2407	2421	2438	rBV	1115526	1798794	1.78%	0.428%
40	9.032	2458	2470	2487	rBV	1068385	1644618	1.63%	0.391%
41	9.158	2497	2509	2533	rVB	2462715	3936670	3.90%	0.936%
42	9.334	2556	2564	2574	rVB8	11797	21844	0.02%	0.005%
43	9.383	2577	2579	2586	rVB7	3718	2968	0.00%	0.001%
44	9.415	2586	2589	2592	rBV4	3034	2827	0.00%	0.001%
45	9.482	2600	2610	2618	rBV9	15490	31458	0.03%	0.007%
46	9.569	2625	2637	2659	rVB3	1166172	2474975	2.45%	0.589%
47	9.740	2688	2690	2692	rBV2	3378	2076	0.00%	0.000%
48	9.775	2693	2701	2716	rVB4	27758	55566	0.05%	0.013%
49	9.900	2735	2740	2743	rVB6	2341	2402	0.00%	0.001%
50	9.955	2743	2757	2775	rBV	260477	421752	0.42%	0.100%
51	10.167	2811	2823	2827	rBV2	225611	370670	0.37%	0.088%
52	10.206	2828	2835	2840	rBV	513204	768713	0.76%	0.183%
53	10.299	2858	2864	2875	rVB2	35012	54178	0.05%	0.013%
54	10.376	2878	2888	2904	rBV	109780	187508	0.19%	0.045%
55	10.469	2905	2917	2935	rBV	3037741	6421941	6.35%	1.527%
56	10.563	2935	2946	2957	rBV2	4165972	6376745	6.31%	1.516%
57	10.627	2957	2966	2976	rBV	2307520	3799300	3.76%	0.904%
58	10.759	3000	3007	3013	rBV	482075	735912	0.73%	0.175%
59	10.939	3047	3063	3075	rBV	7741481	11804892	11.68%	2.807%
60	11.006	3075	3084	3093	rBV	2131495	3098933	3.07%	0.737%
61	11.058	3093	3100	3110	rVB	602218	852882	0.84%	0.203%
62	11.190	3128	3141	3155	rBV3	3062313	6830684	6.76%	1.624%
63	11.267	3155	3165	3183	rVV	13787678	21157379	20.94%	5.032%
64	11.360	3183	3194	3206	rVV	2346771	3664397	3.63%	0.871%
65	11.421	3206	3213	3224	rVB	192446	286031	0.28%	0.068%
66	11.553	3240	3254	3274	rBV	19146481	29669065	29.36%	7.056%
67	11.653	3274	3285	3301	rVB3	1709246	3411861	3.38%	0.811%
68	11.772	3309	3322	3338	rBV	29784322	46234017	45.75%	10.995%
69	11.852	3342	3347	3355	rVB2	58996	80906	0.08%	0.019%
70	11.926	3361	3370	3378	rBV3	744344	1209068	1.20%	0.288%
71	12.038	3391	3405	3412	rBV	696750	1135570	1.12%	0.270%

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72	12.138	3428	3436	3452	rVB	1445952	2197529	2.17%	0.523%
73	12.219	3453	3461	3470	rVB	306267	447743	0.44%	0.106%
74	12.276	3471	3479	3489	rVB10	85692	160036	0.16%	0.038%
75	12.341	3491	3499	3502	rBV2	81004	110262	0.11%	0.026%
76	12.386	3503	3513	3526	rBV2	2288752	5045875	4.99%	1.200%
77	12.489	3535	3545	3555	rBV2	5519042	9410230	9.31%	2.238%
78	12.614	3578	3584	3596	rBV3	159471	260387	0.26%	0.062%
79	12.749	3616	3626	3644	rVB	2298795	3496506	3.46%	0.832%
80	12.910	3656	3676	3686	rBV2	3544040	5979733	5.92%	1.422%
81	12.984	3686	3699	3717	rVV	16473028	28625341	28.33%	6.808%
82	13.080	3718	3729	3743	rVB2	3278401	5620281	5.56%	1.337%
83	13.193	3752	3764	3775	rVB2	1646751	2928700	2.90%	0.696%
84	13.662	3899	3910	3927	rBV	15133687	22135620	21.90%	5.264%
85	13.794	3946	3951	3960	rVB2	141497	196341	0.19%	0.047%
86	13.936	3988	3995	4005	rBV	156963	249293	0.25%	0.059%
87	14.119	4043	4052	4064	rBV3	161529	338926	0.34%	0.081%
88	14.318	4107	4114	4125	rVB2	168167	263472	0.26%	0.063%
89	14.604	4192	4203	4210	rBV	1645860	2567832	2.54%	0.611%
90	14.669	4210	4223	4245	rVV2	60765586	101053799	100.00%	24.032%
91	14.775	4246	4256	4266	rVV	1571451	2615642	2.59%	0.622%
92	14.849	4266	4279	4313	rVB	29524104	46601988	46.12%	11.083%
93	15.341	4426	4432	4436	rBV3	23180	31160	0.03%	0.007%
94	15.389	4437	4447	4462	rVV	1943442	2892665	2.86%	0.688%
95	15.579	4498	4506	4515	rBV	451107	664425	0.66%	0.158%
96	15.694	4532	4542	4563	rVB	458908	848495	0.84%	0.202%
97	15.839	4578	4587	4593	rBV	400291	624165	0.62%	0.148%
98	16.514	4786	4797	4814	rBV	1051167	1707263	1.69%	0.406%
99	16.813	4881	4890	4905	rVB	242302	372778	0.37%	0.089%
100	17.418	5070	5078	5091	rVB2	176916	293881	0.29%	0.070%

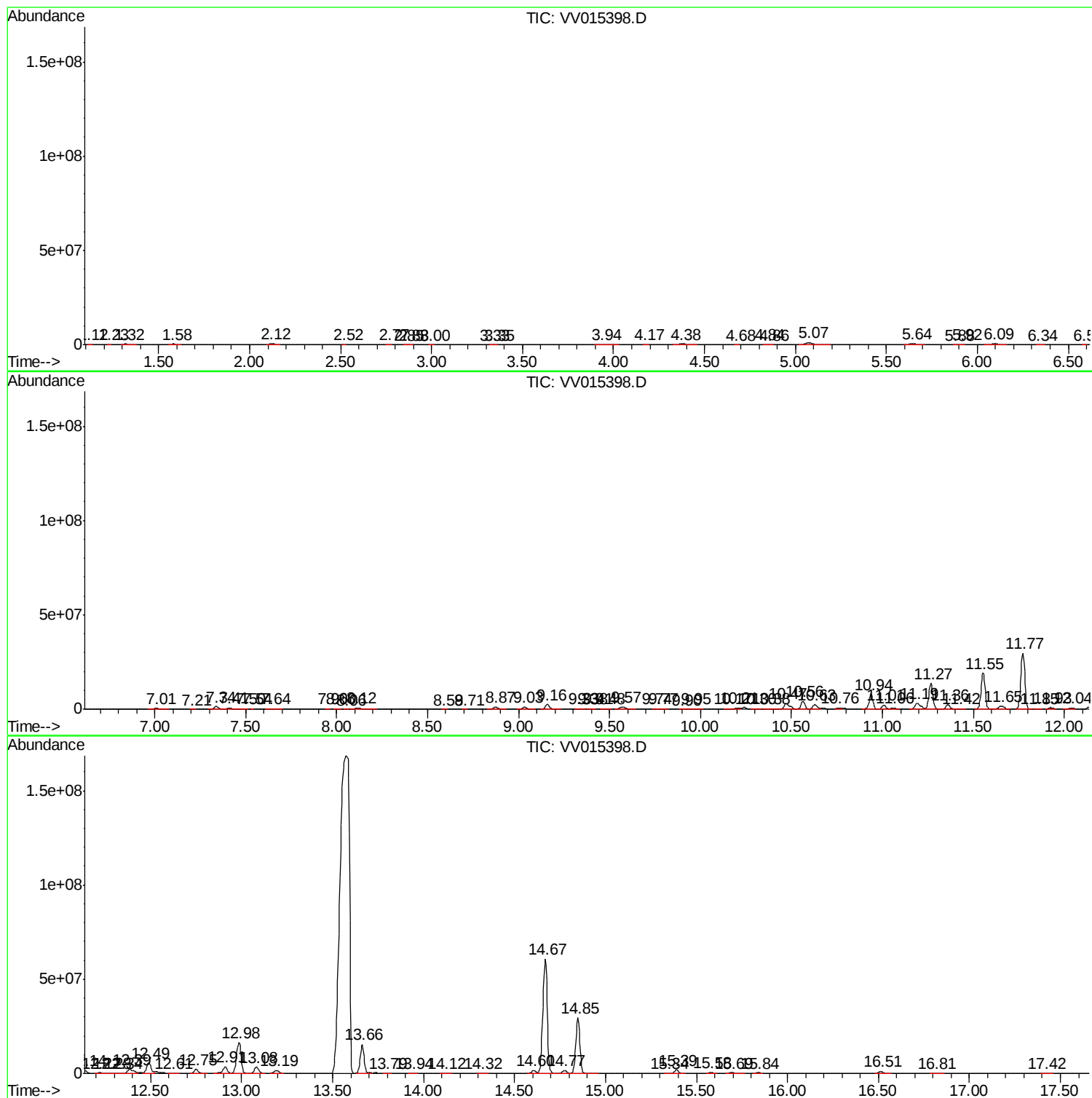
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM040120WMA.M
Quant Title : VOC Analysis

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



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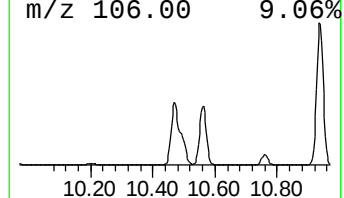
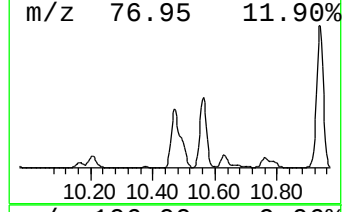
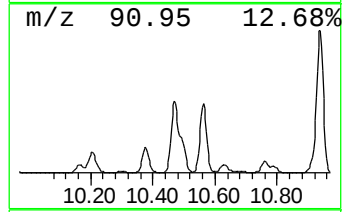
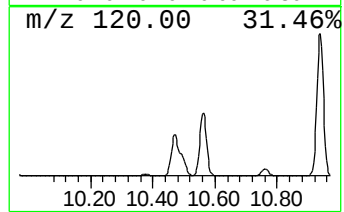
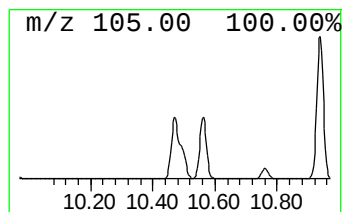
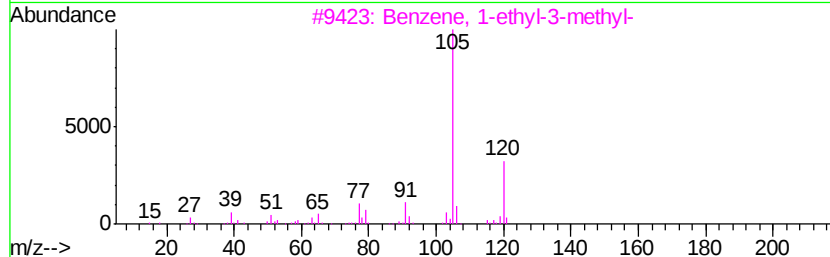
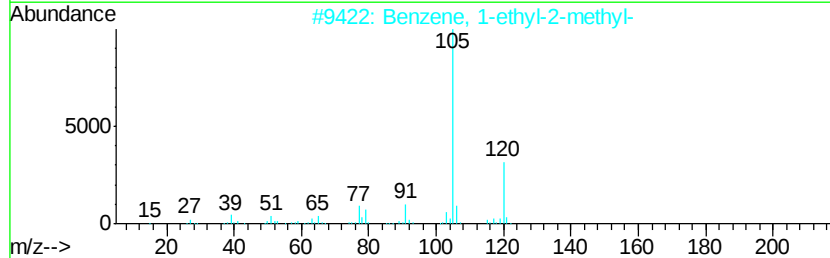
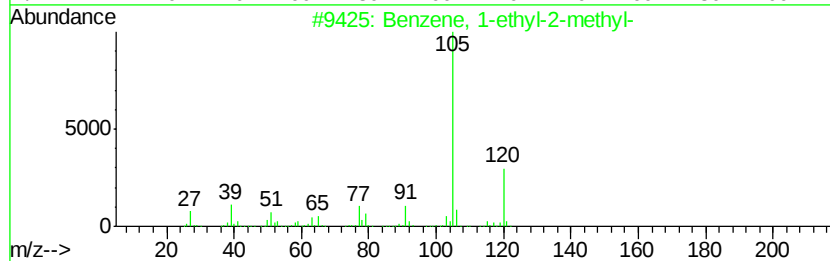
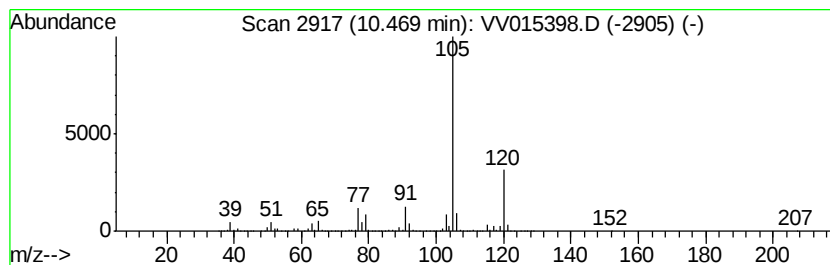
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM040120WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	15.18 ug/L	6421940	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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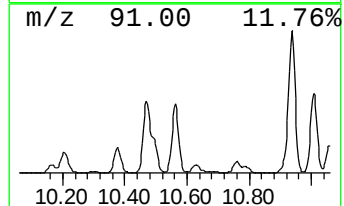
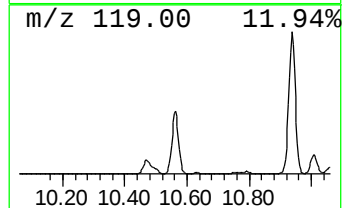
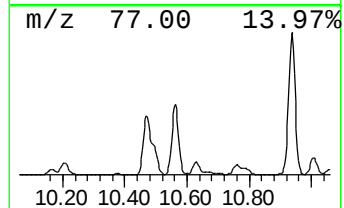
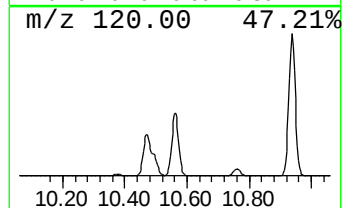
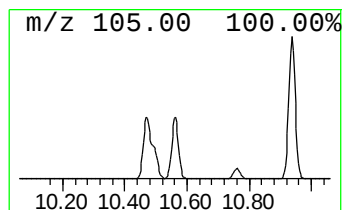
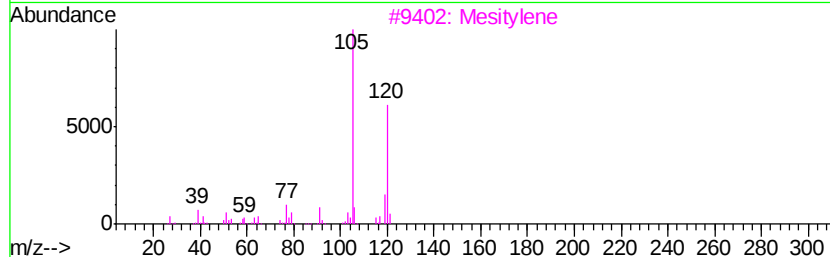
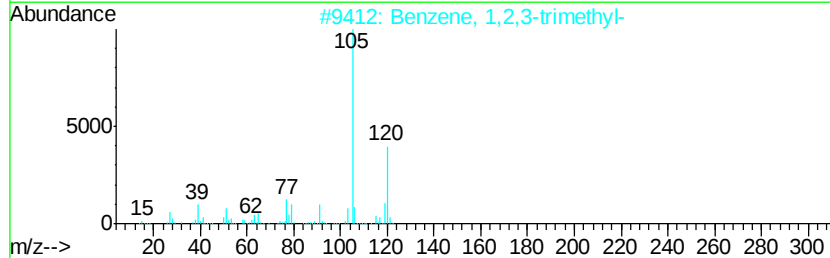
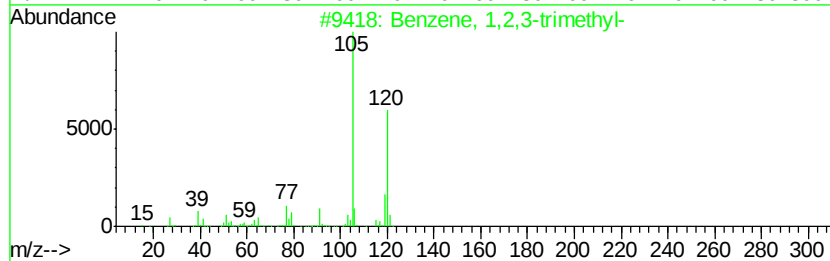
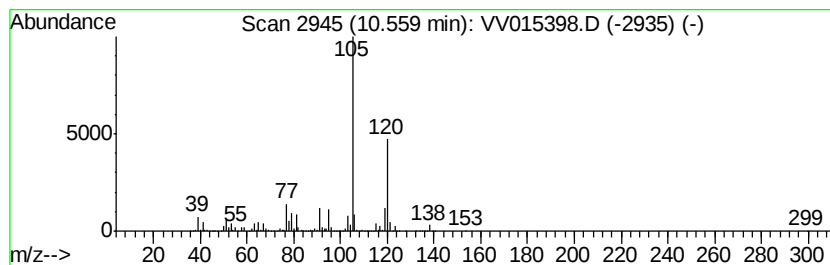
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Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	15.07 ug/L	6376750	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Mesitylene	120	C9H12	000108-67-8	94
4		Mesitylene	120	C9H12	000108-67-8	93
5		Mesitylene	120	C9H12	000108-67-8	93



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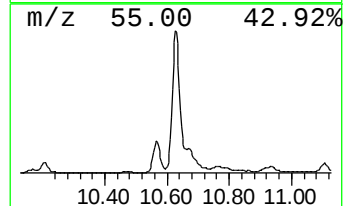
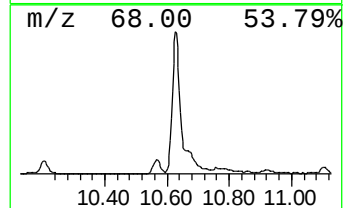
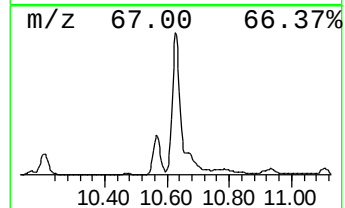
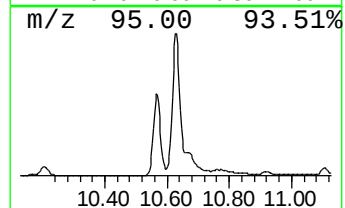
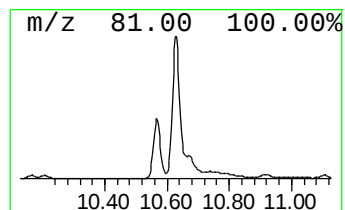
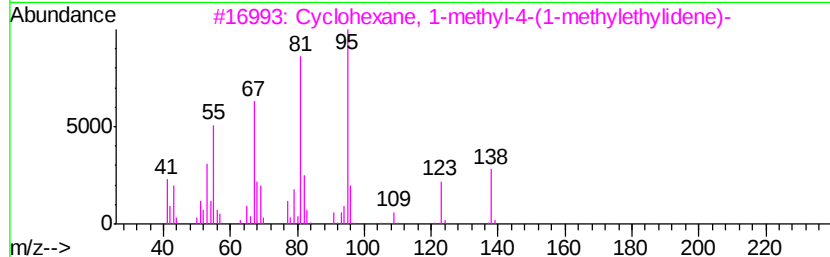
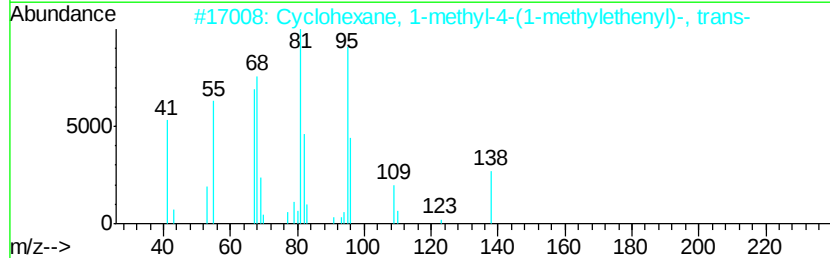
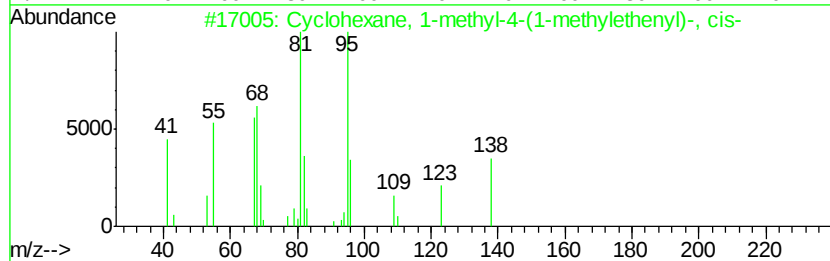
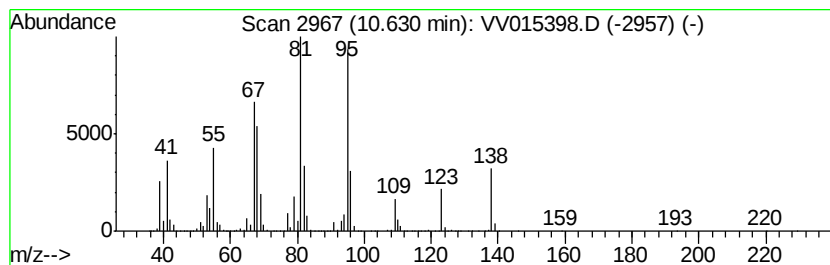
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Quant Title : VOC Analysis

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

Peak Number 4 Cyclohexane, 1-methyl-4-(1-... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.63	8.98 ug/L	3799300	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-methyl-4-(1-methyl-4-(1-methylethenyl)-, cis-	138	C10H18	001879-07-8	98
2		Cyclohexane, 1-methyl-4-(1-methyl-4-(1-methylethenyl)-, trans-	138	C10H18	001124-25-0	87
3		Cyclohexane, 1-methyl-4-(1-methyl-4-(1-methylethenyl)-, trans-	138	C10H18	001124-27-2	87
4		Cyclohexene, 4-methyl-1-(1-methyl-4-(1-methylethenyl)-, trans-	138	C10H18	000500-00-5	81
5		Bicyclo[4.1.0]heptane, 3,7,7-trimethyl-	138	C10H18	002778-68-9	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

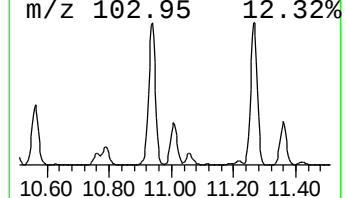
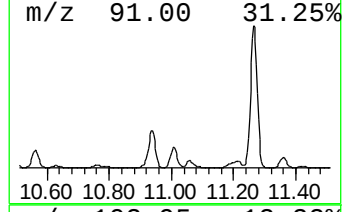
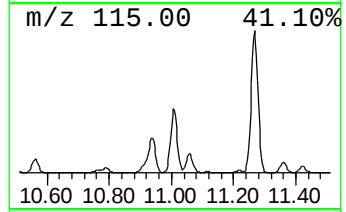
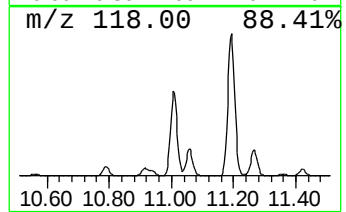
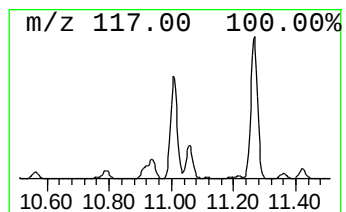
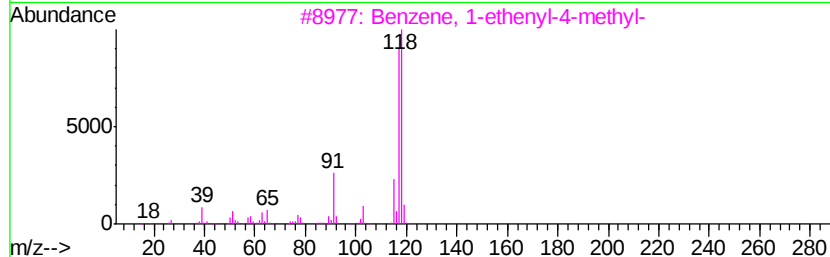
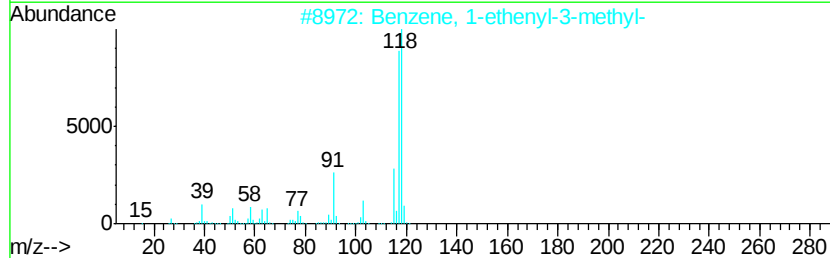
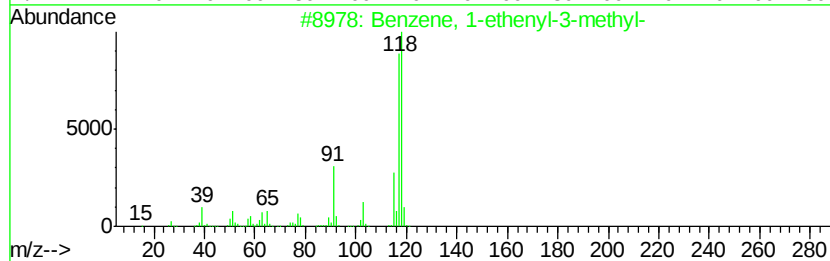
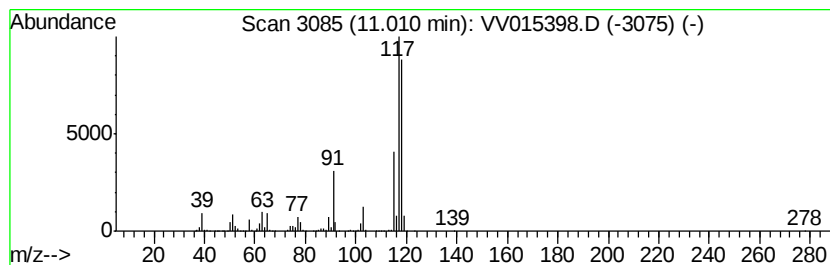
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-ethenyl-3-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.01	7.32 ug/L	3098930	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	95
4		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	94
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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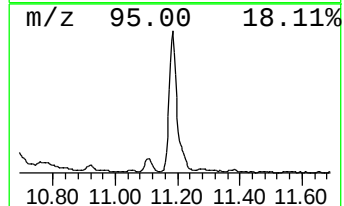
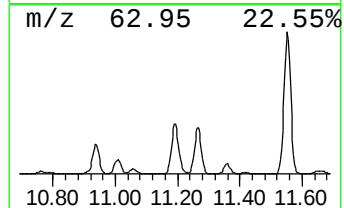
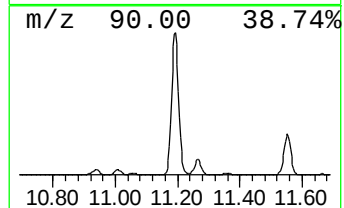
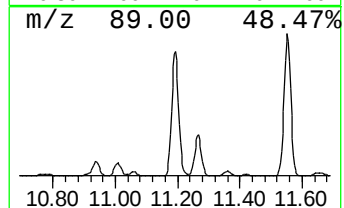
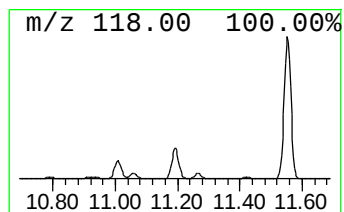
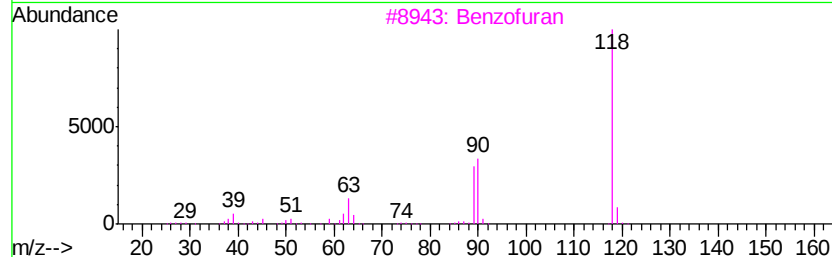
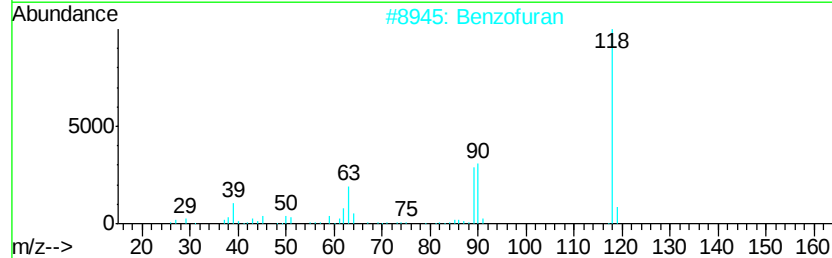
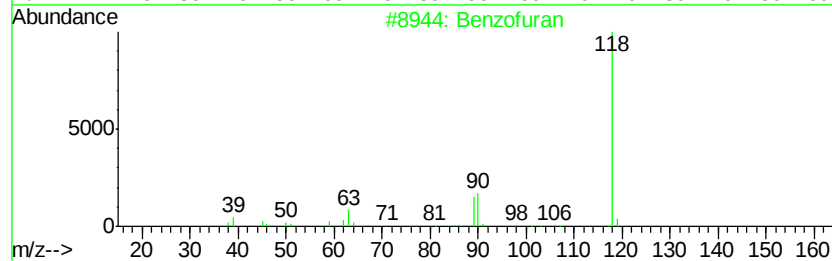
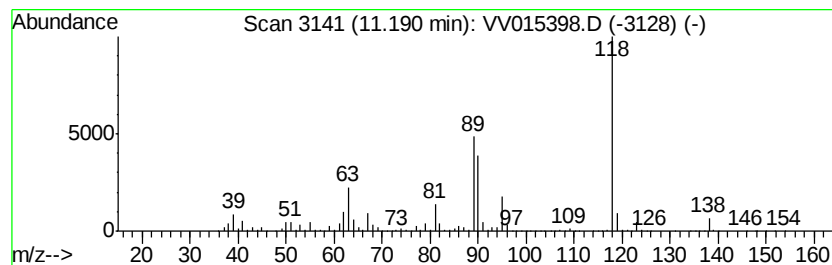
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Quant Title : VOC Analysis

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

Peak Number 7 Benzofuran Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	16.14 ug/L	6830680	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran	118	C8H6O	000271-89-6	83
2		Benzofuran	118	C8H6O	000271-89-6	81
3		Benzofuran	118	C8H6O	000271-89-6	76
4		3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	64
5		2H-Cyclopenta[d]pyridazine	118	C7H6N2	000270-64-4	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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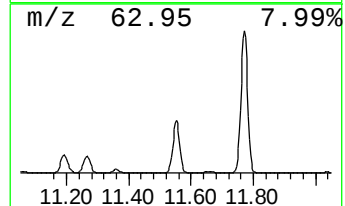
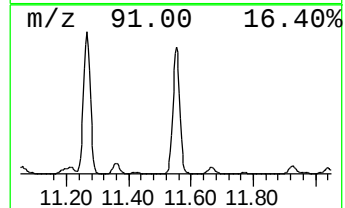
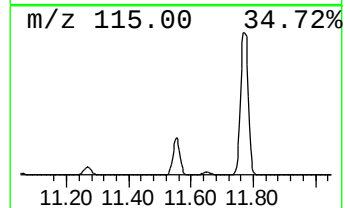
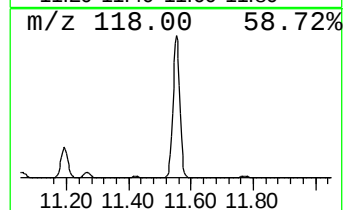
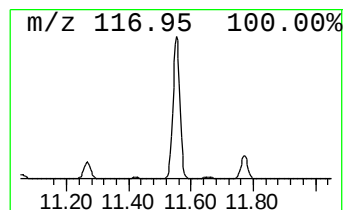
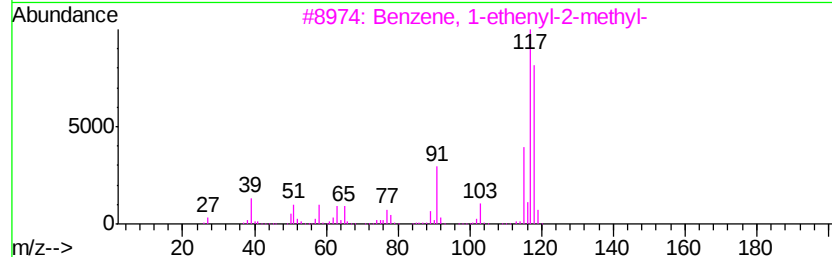
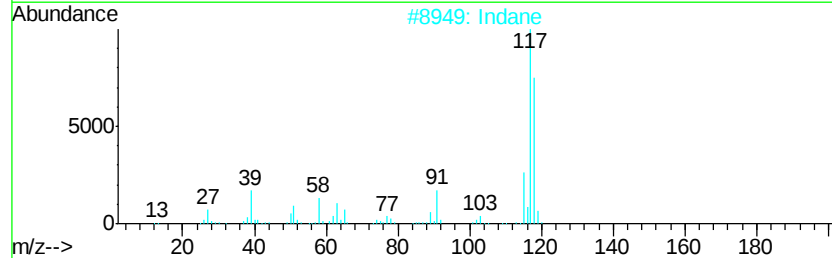
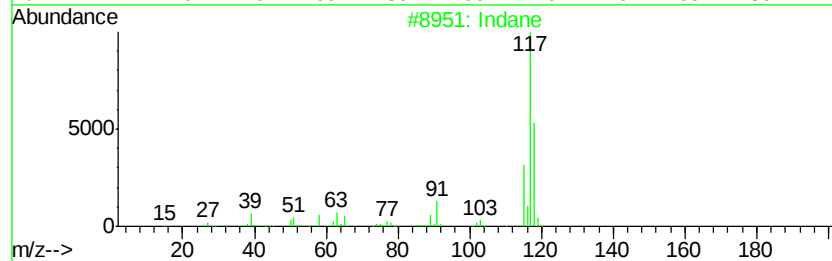
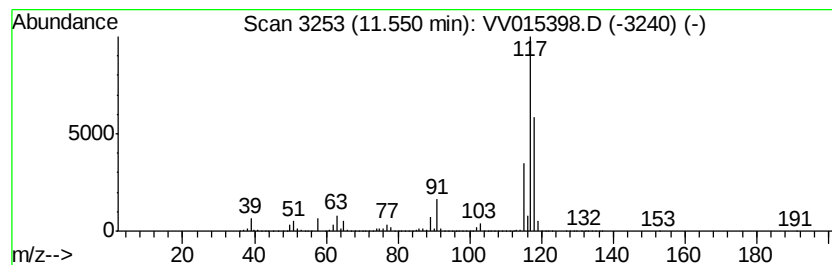
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Quant Title : VOC Analysis

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

Peak Number 9 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	70.12 ug/L	29669100	1,4-Dichlorobenzene-d4	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Indane		118	C9H10	000496-11-7	91
3	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	87
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	83
5	Indane		118	C9H10	000496-11-7	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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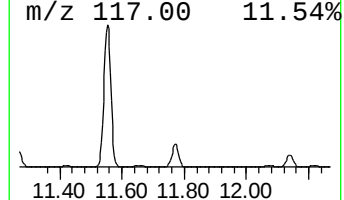
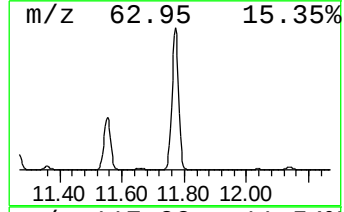
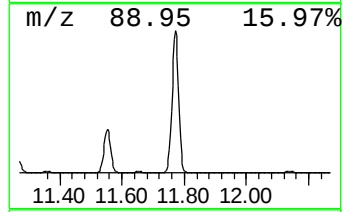
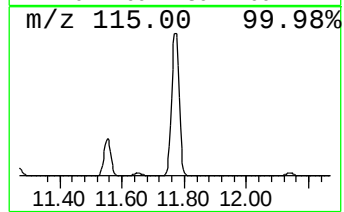
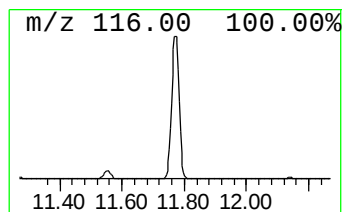
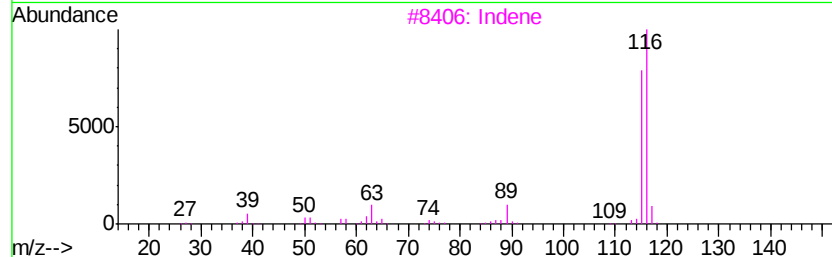
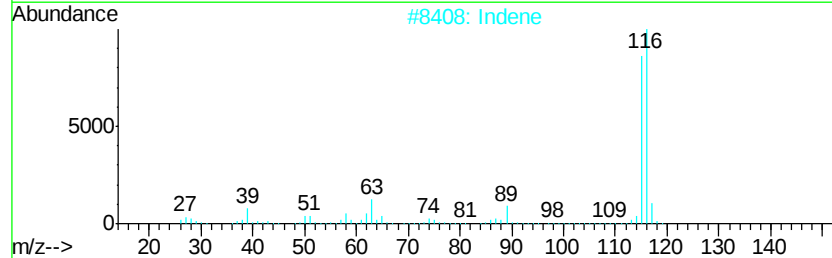
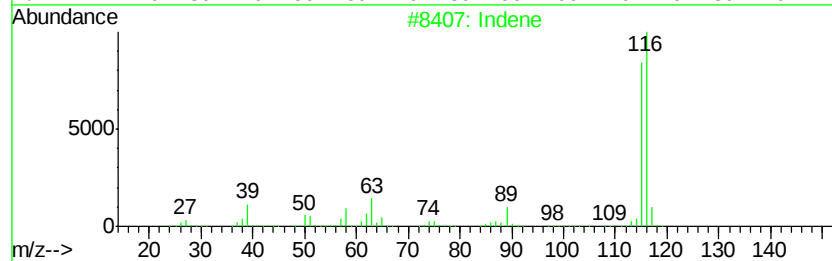
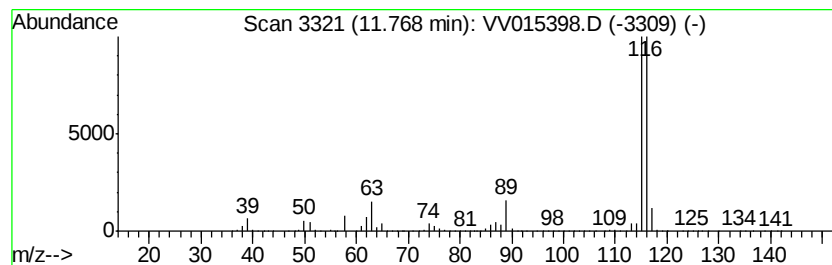
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Quant Title : VOC Analysis

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

Peak Number 10 Indene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	109.26 ug/L	46234000	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	97
2		Indene	116	C9H8	000095-13-6	96
3		Indene	116	C9H8	000095-13-6	93
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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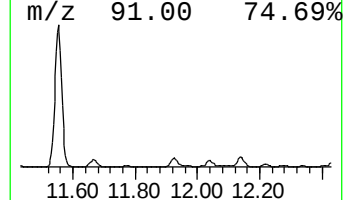
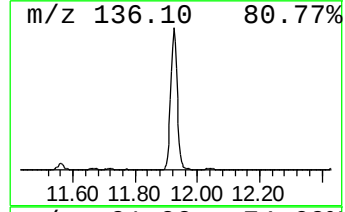
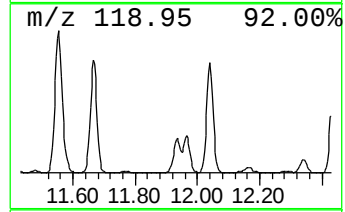
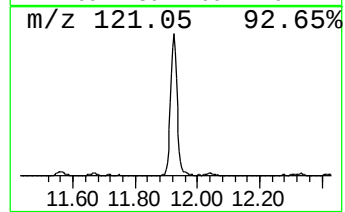
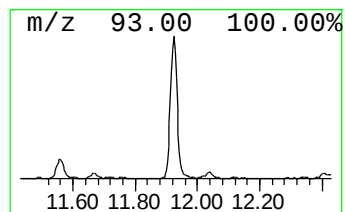
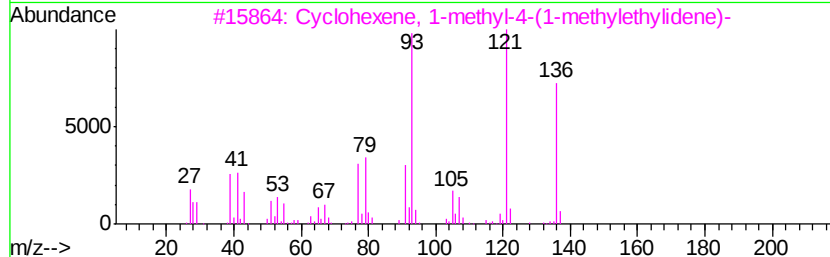
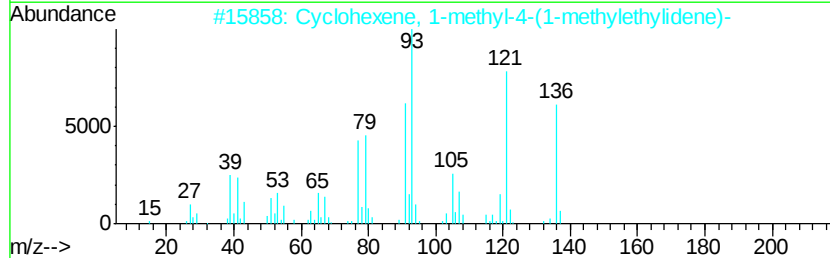
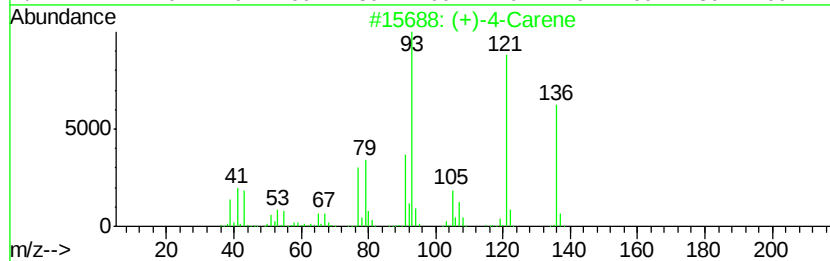
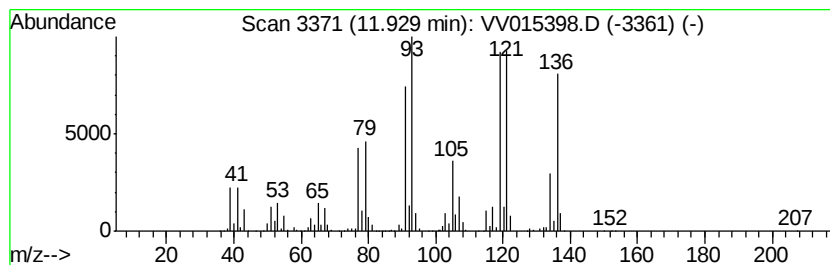
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Quant Title : VOC Analysis

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

Peak Number 11 (+)-4-Carene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	2.86 ug/L	1209070	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(+)-4-Carene	136	C10H16	029050-33-7	96
2			Cyclohexene, 1-methyl-4-(1-methyl-4-ethyldiene)-	136	C10H16	000586-62-9	96
3			Cyclohexene, 1-methyl-4-(1-methyl-4-ethyldiene)-	136	C10H16	000586-62-9	96
4			1,3-Cyclohexadiene, 1-methyl-4-(1-methyl-4-ethyldiene)-	136	C10H16	000099-86-5	93
5			Cyclohexene, 1-methyl-4-(1-methyl-4-ethyldiene)-	136	C10H16	000586-62-9	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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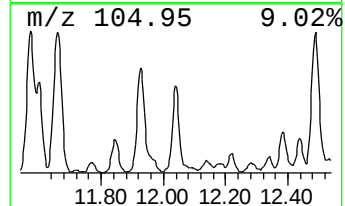
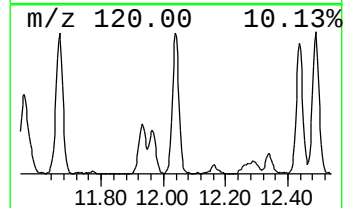
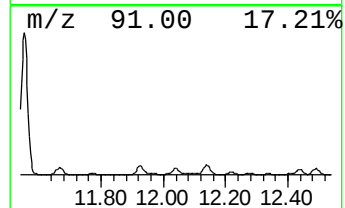
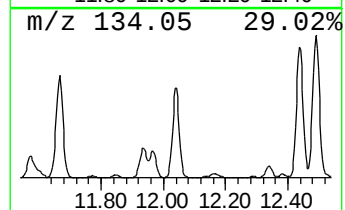
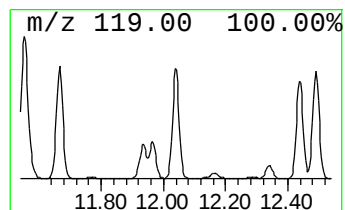
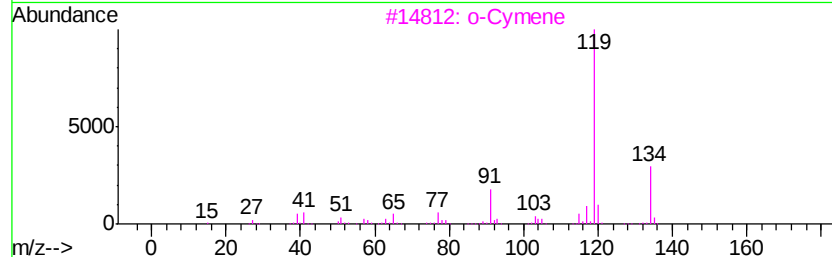
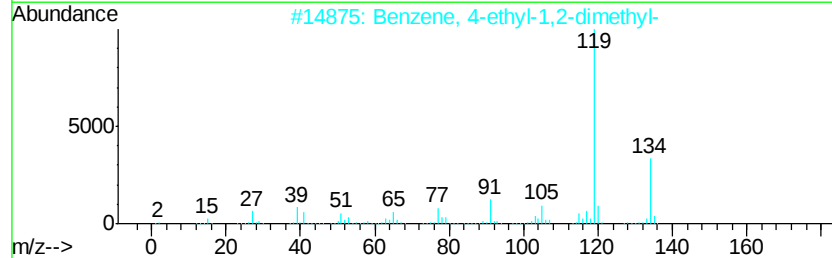
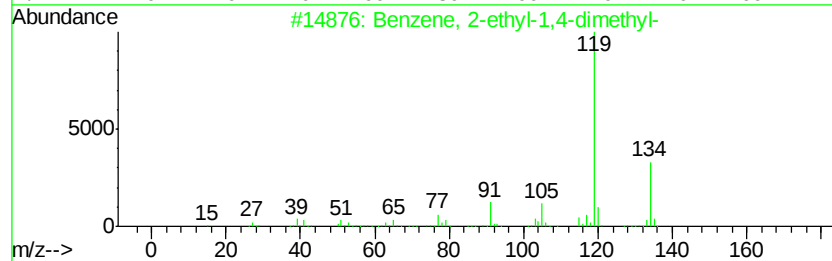
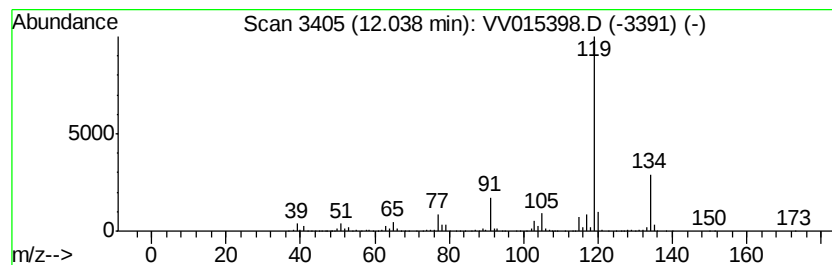
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	2.68 ug/L	1135570	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
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ALS Vial : 1 Sample Multiplier: 1

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ClientSampleId :
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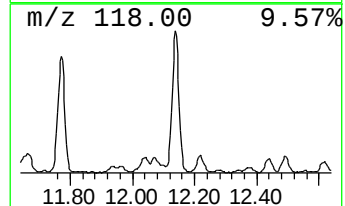
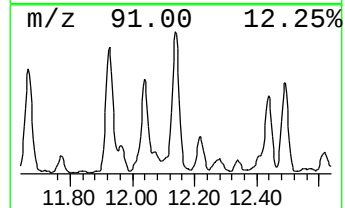
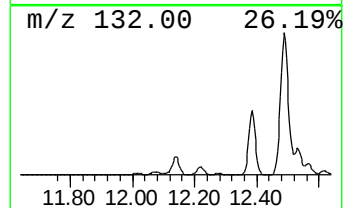
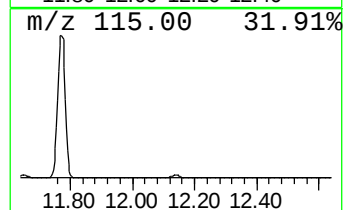
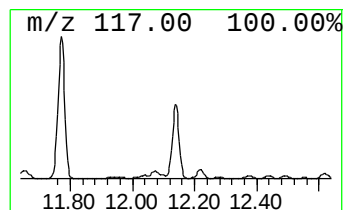
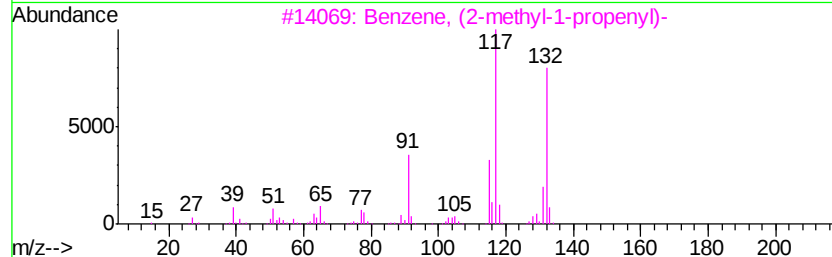
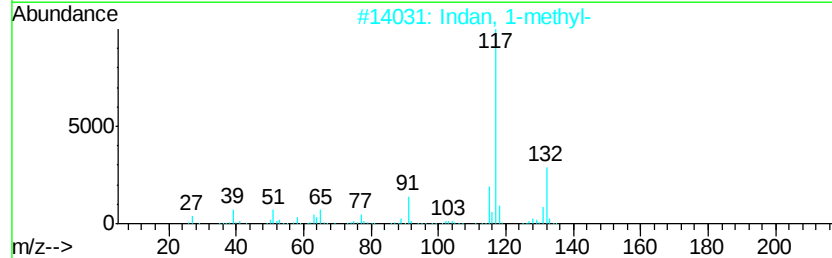
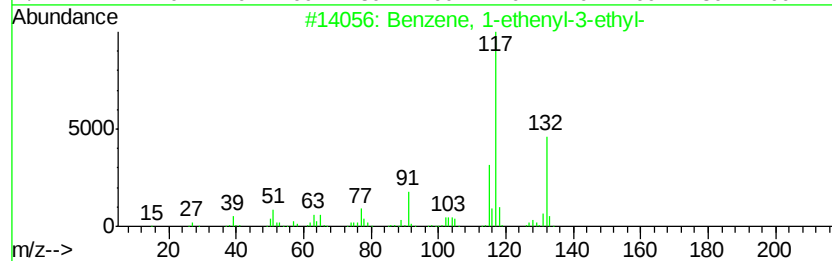
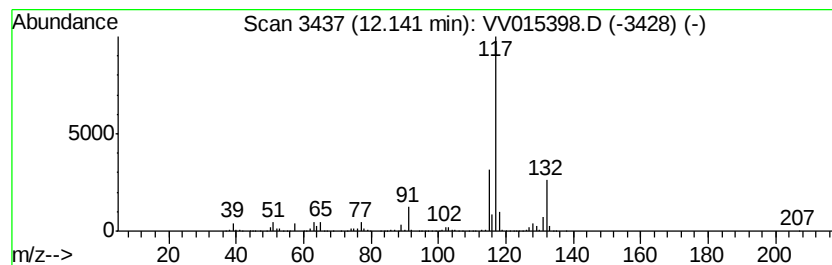
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Quant Title : VOC Analysis

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

Peak Number 13 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	5.19 ug/L	2197530	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

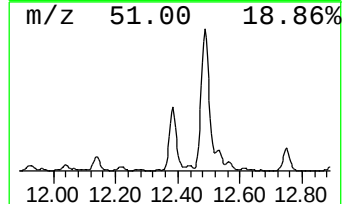
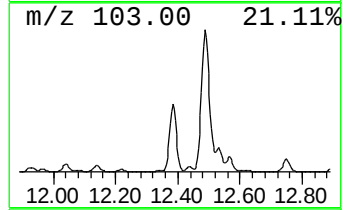
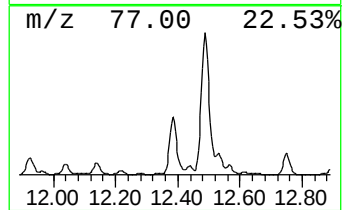
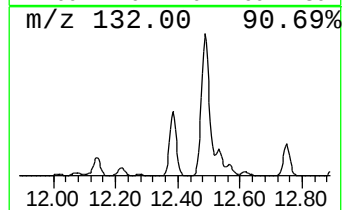
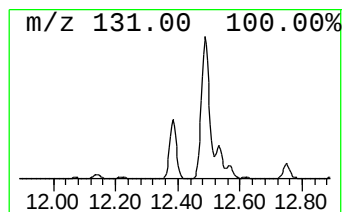
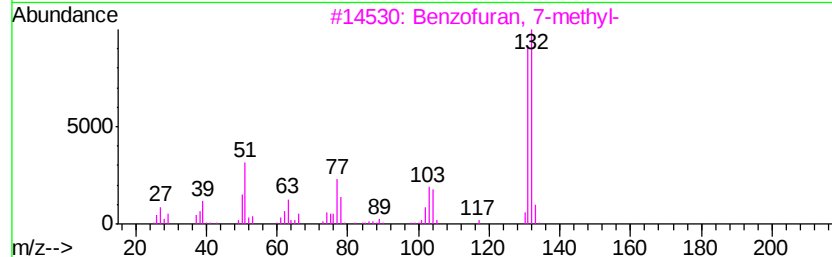
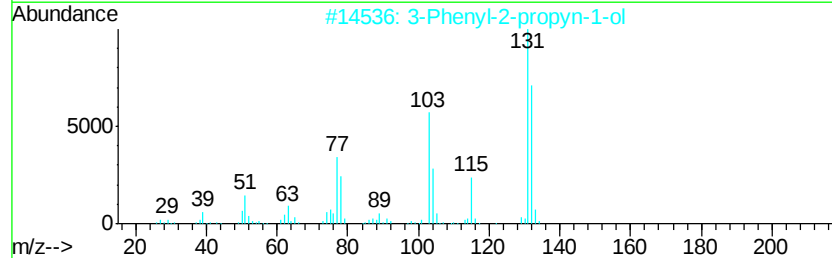
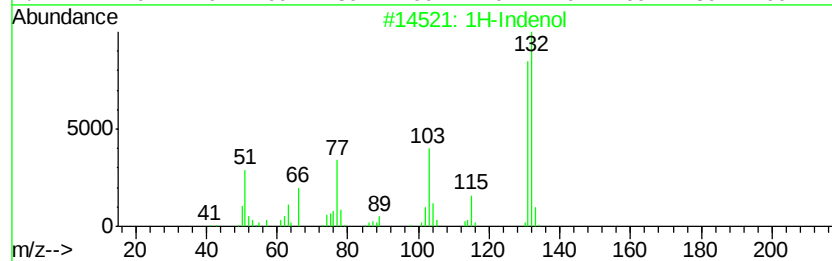
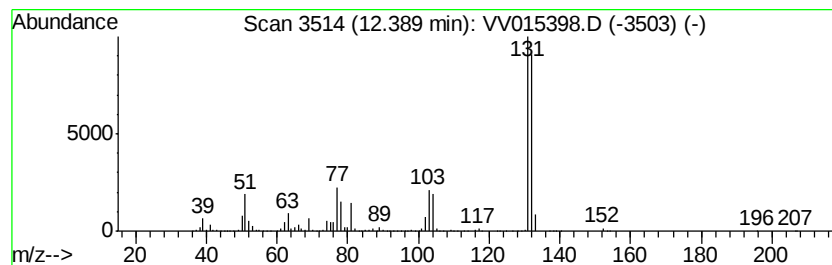
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Quant Title : VOC Analysis

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

Peak Number 14 1H-Indenol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.39	11.92 ug/L	5045880	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indenol	132	C9H8O	056631-57-3	91
2			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	91
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

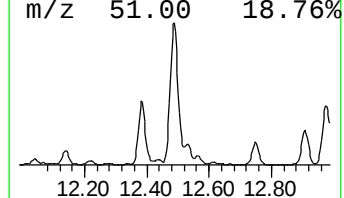
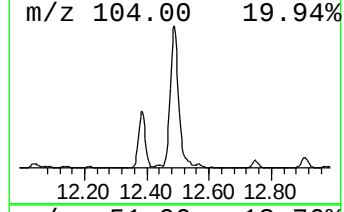
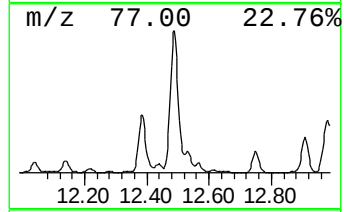
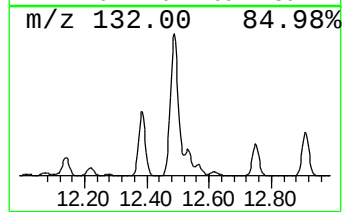
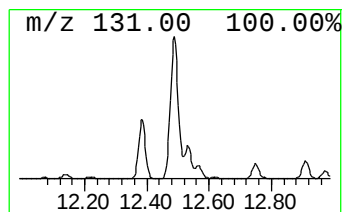
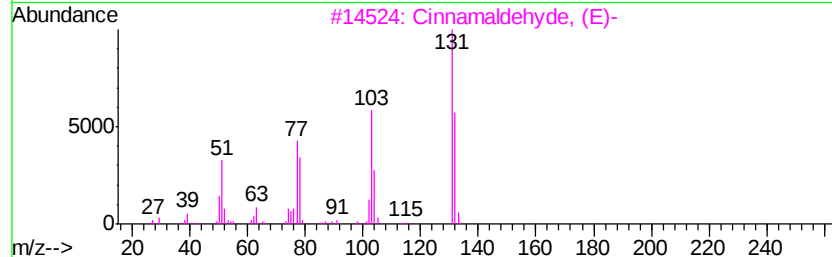
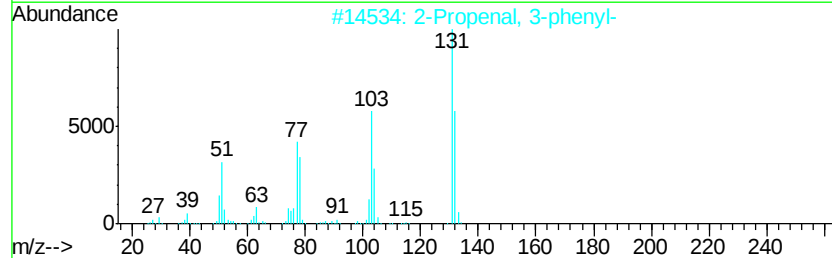
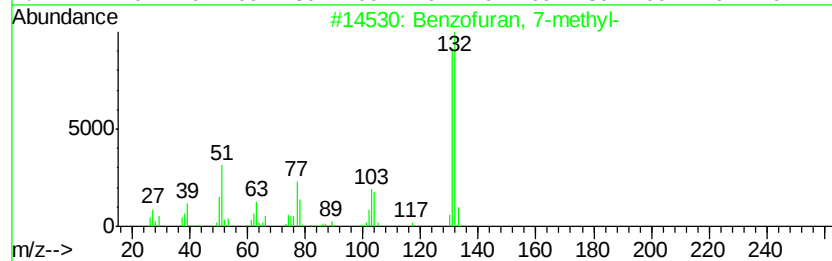
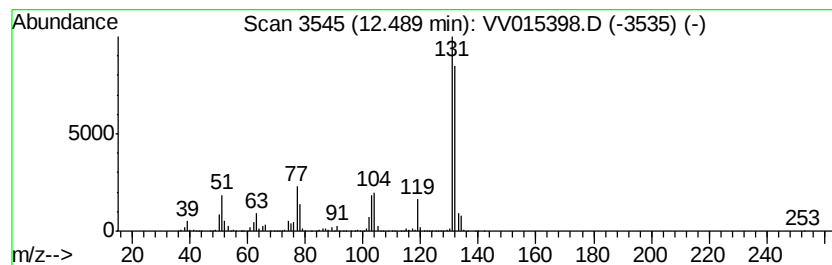
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Quant Title : VOC Analysis

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

Peak Number 15 Benzofuran, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	22.24 ug/L	9410230	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
2		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90
3		Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	90
4		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
5		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

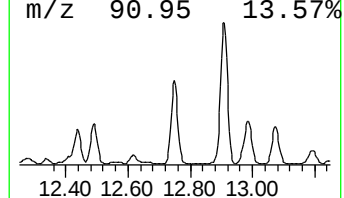
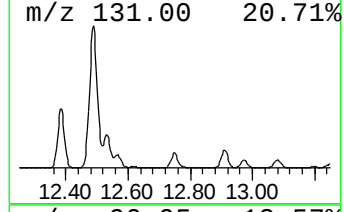
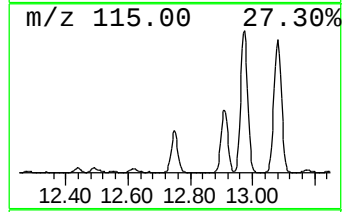
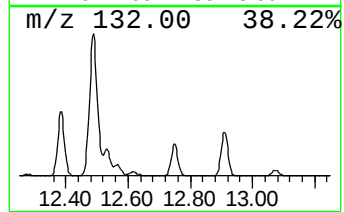
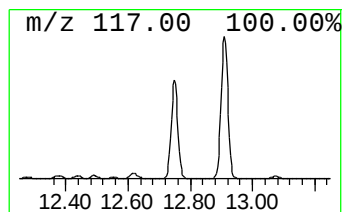
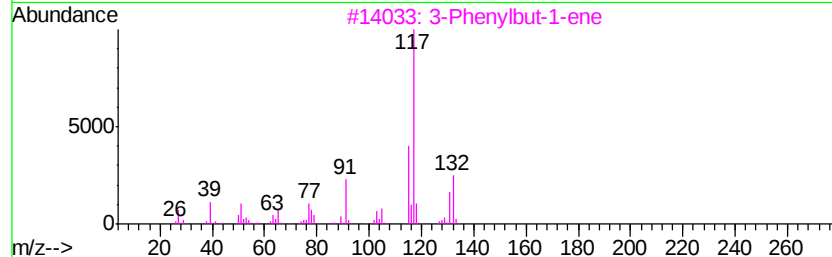
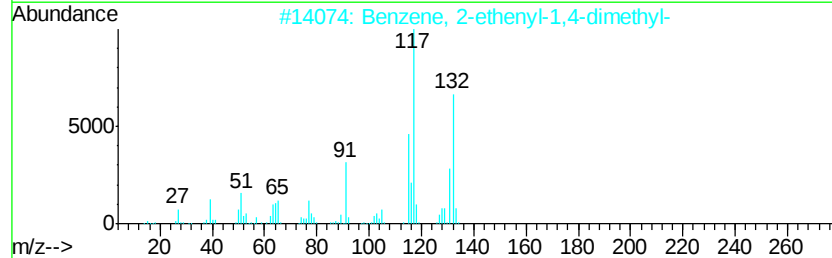
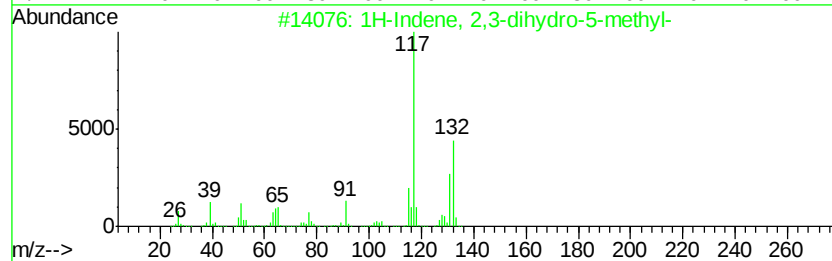
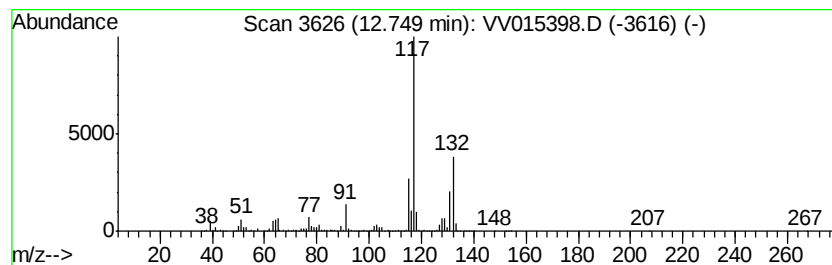
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Quant Title : VOC Analysis

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

Peak Number 16 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	8.26 ug/L	3496510	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

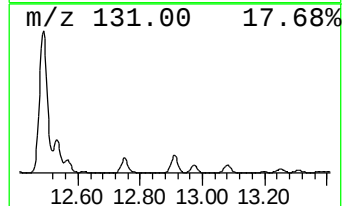
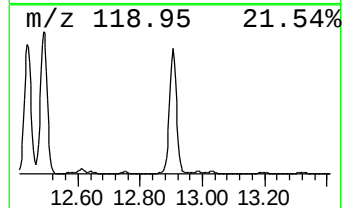
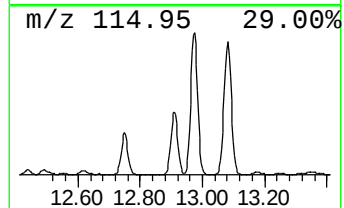
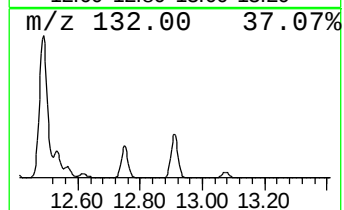
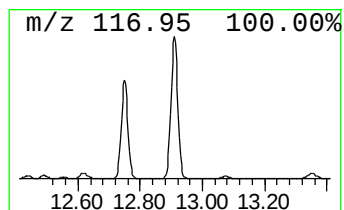
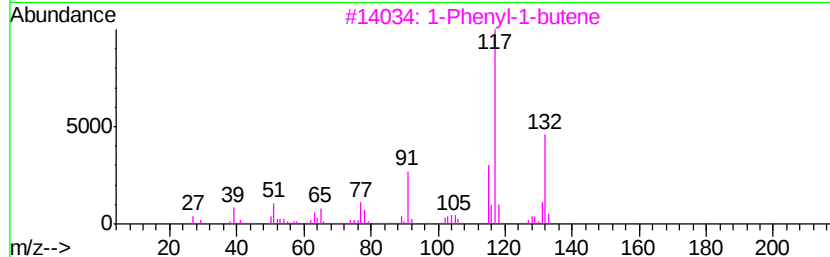
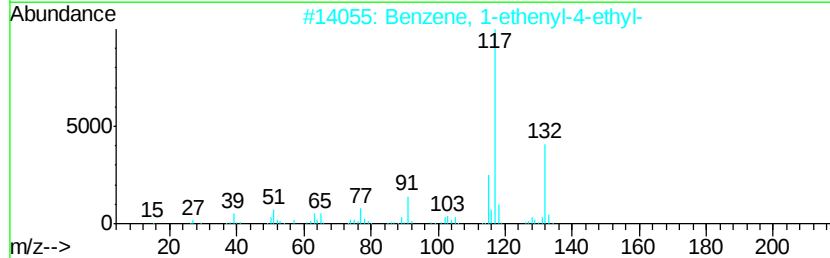
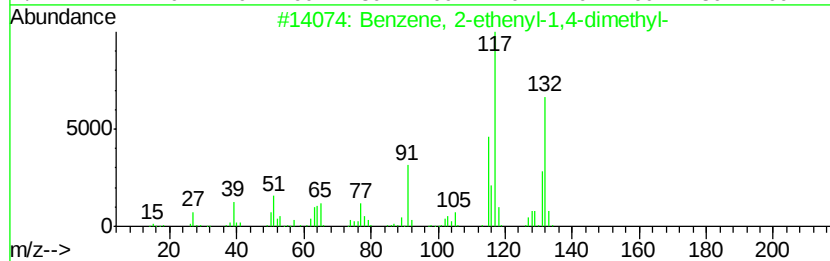
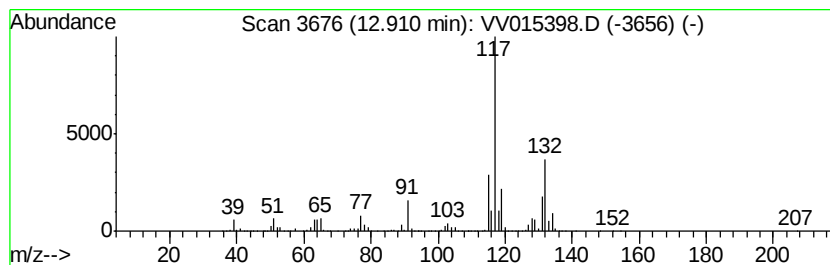
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Quant Title : VOC Analysis

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

Peak Number 17 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.91	14.13 ug/L	5979730	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

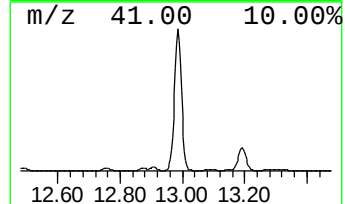
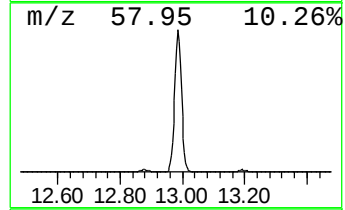
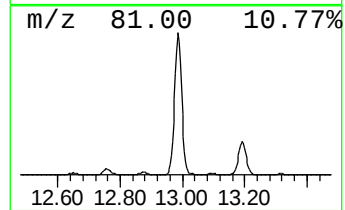
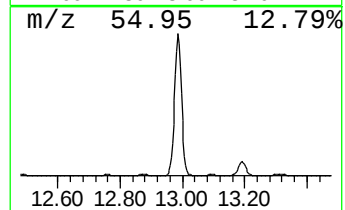
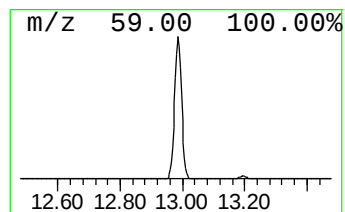
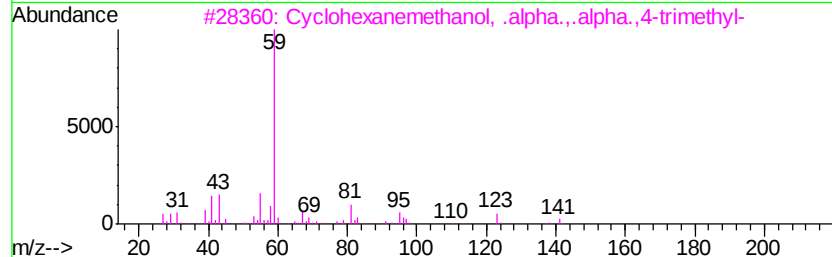
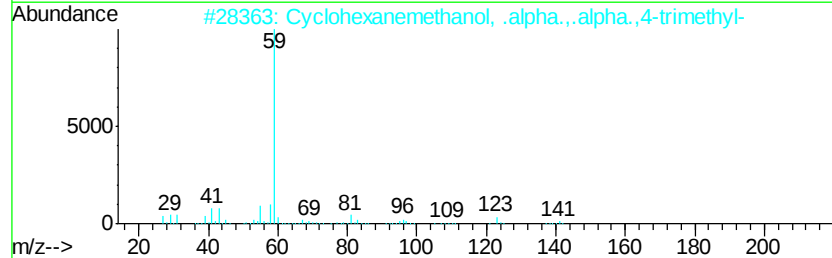
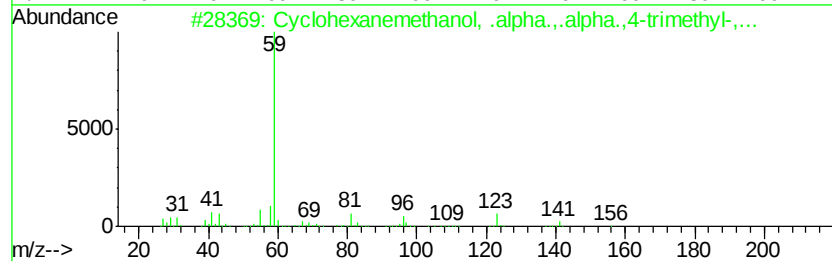
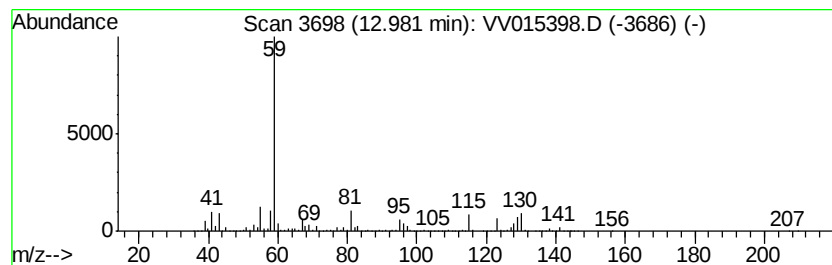
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Quant Title : VOC Analysis

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

Peak Number 18 Cyclohexanemethanol, .alpha... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	67.65 ug/L	28625300	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanemethanol, .alpha...al...	156	C10H20O	005114-00-1	78
2		Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	78
3		Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	78
4		Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	78
5		Cyclohexanemethanol, .alpha...al...	156	C10H20O	007322-63-6	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

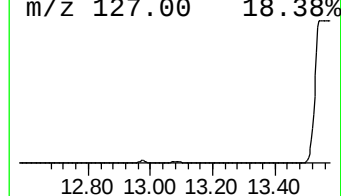
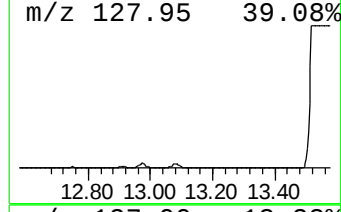
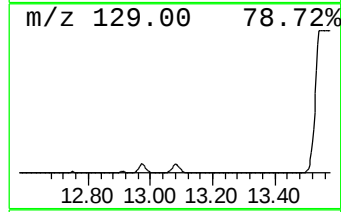
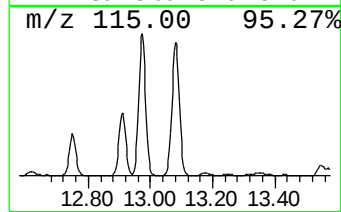
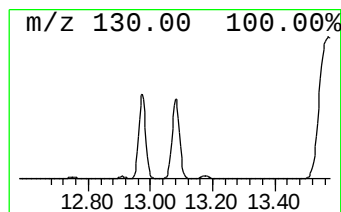
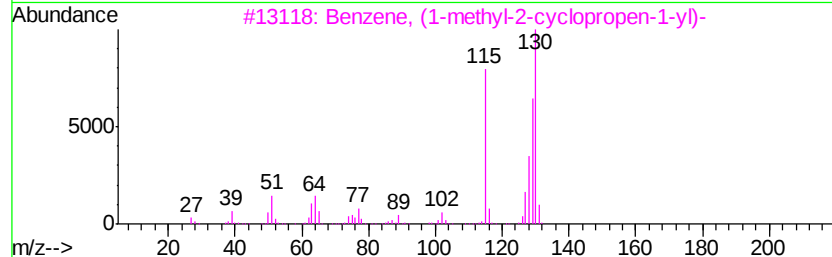
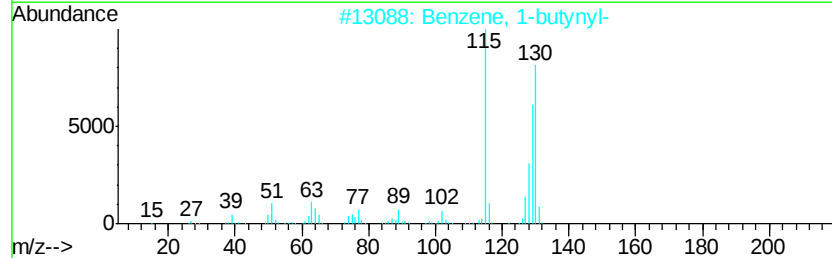
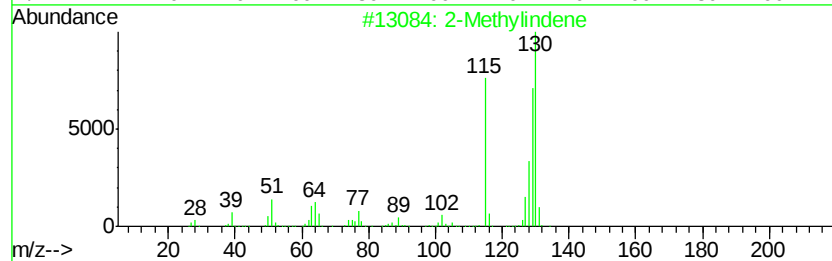
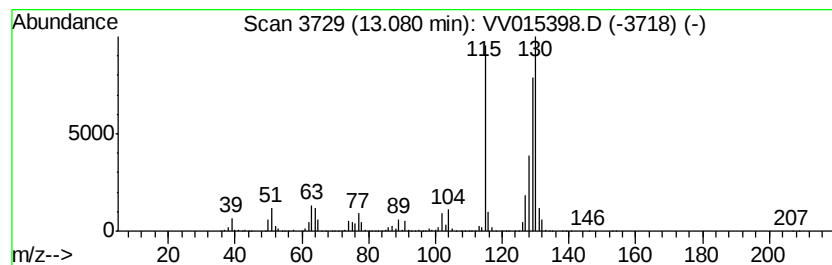
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Quant Title : VOC Analysis

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

Peak Number 19 2-Methylindene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	13.28 ug/L	5620280	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylindene	130	C10H10	002177-47-1	96
2		Benzene, 1-butynyl-	130	C10H10	000622-76-4	96
3		Benzene, (1-methyl-2-cyclopropen-1-yl)-	130	C10H10	065051-83-4	96
4		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95
5		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

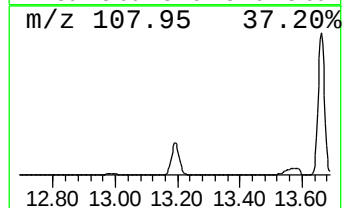
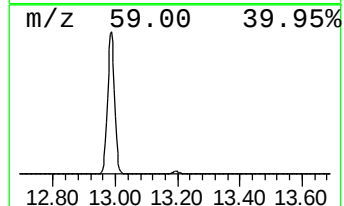
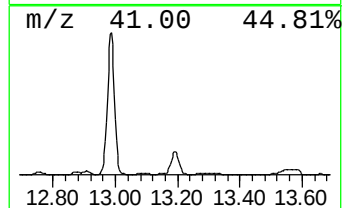
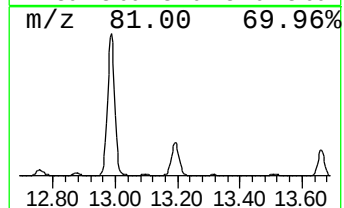
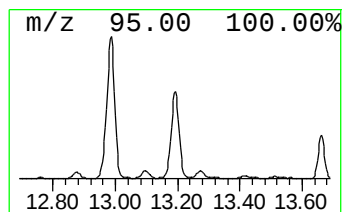
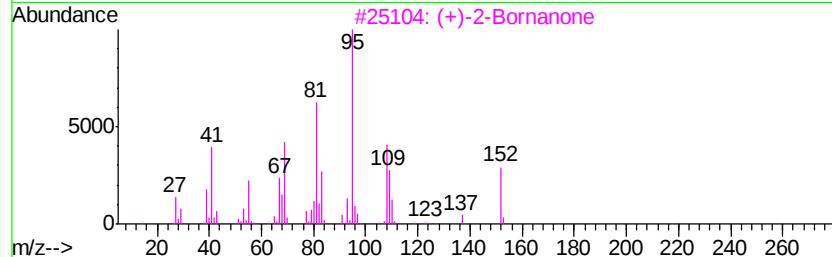
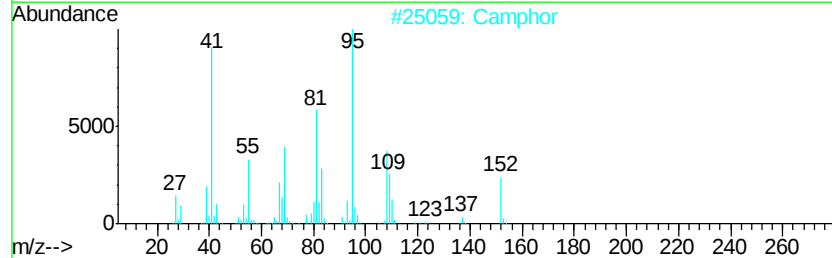
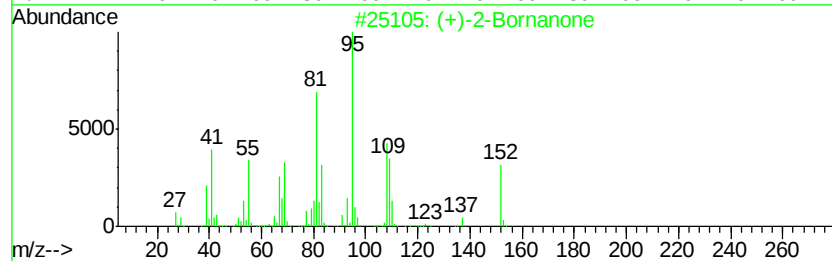
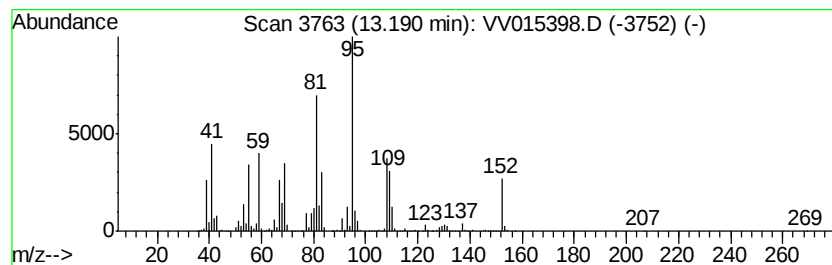
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Quant Title : VOC Analysis

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

Peak Number 20 (+)-2-Bornanone Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	6.92 ug/L	2928700	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(+)-2-Bornanone	152	C10H16O	000464-49-3	98
2			Camphor	152	C10H16O	000076-22-2	98
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	98
4			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	98
5			Camphor	152	C10H16O	000076-22-2	98



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

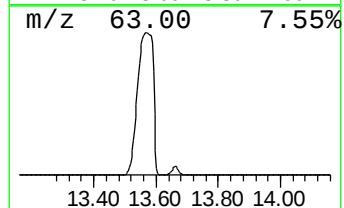
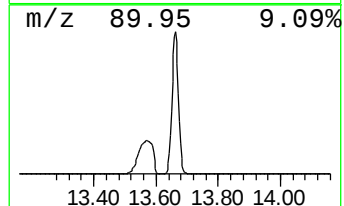
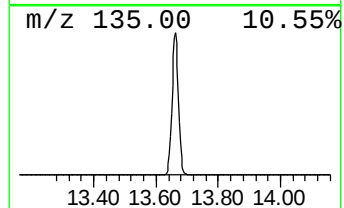
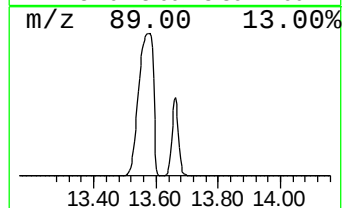
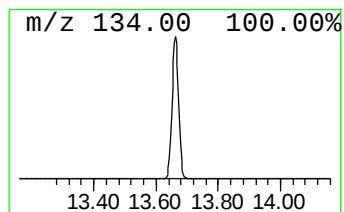
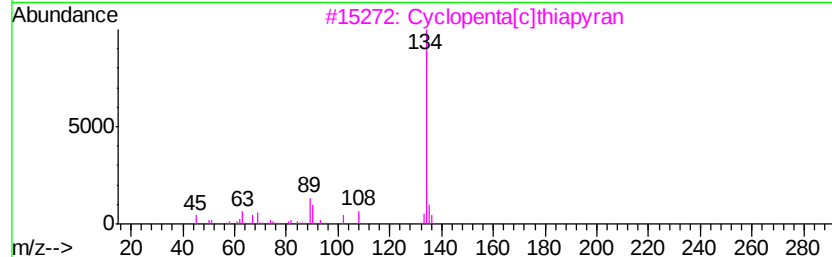
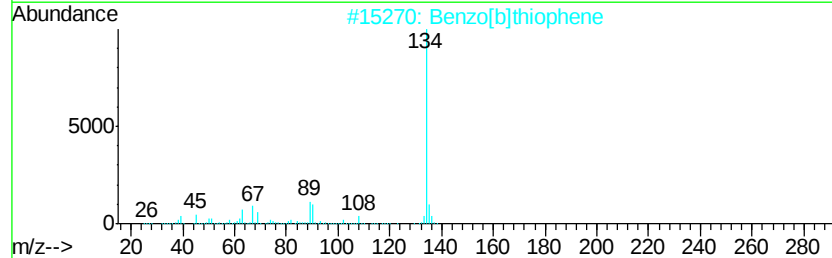
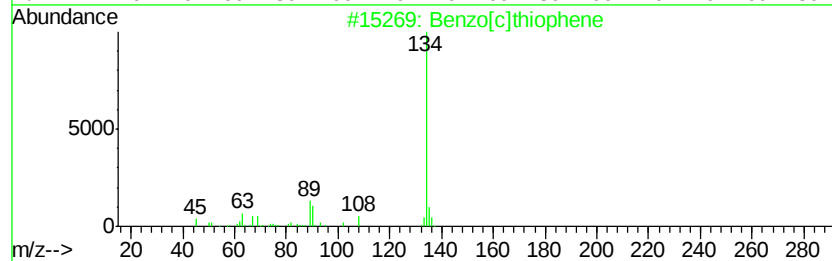
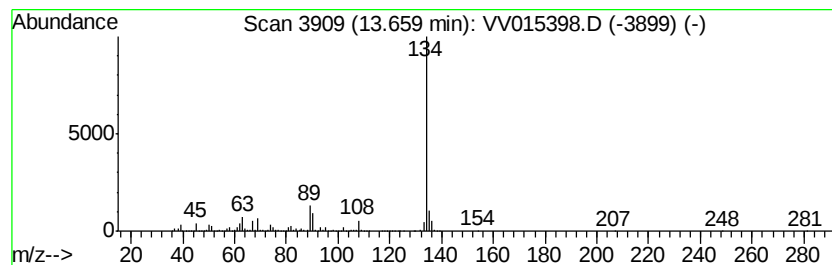
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Quant Title : VOC Analysis

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

Peak Number 21 Benzo[c]thiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	52.31 ug/L	22135600	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	94
3		Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
4		Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5		Benzo[b]thiophene	134	C8H6S	000095-15-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

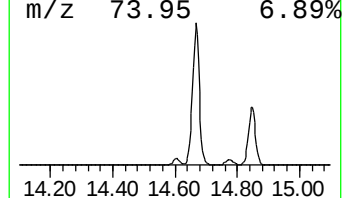
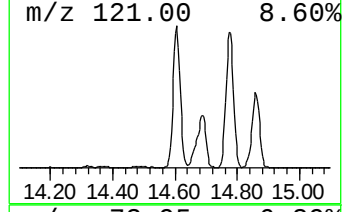
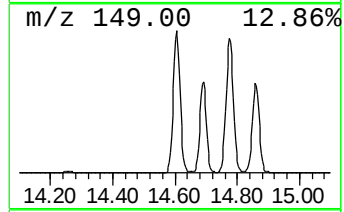
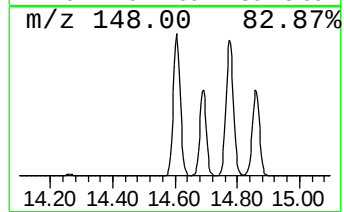
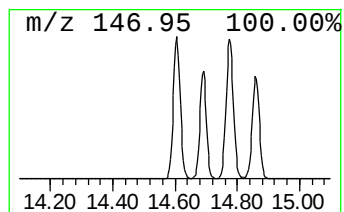
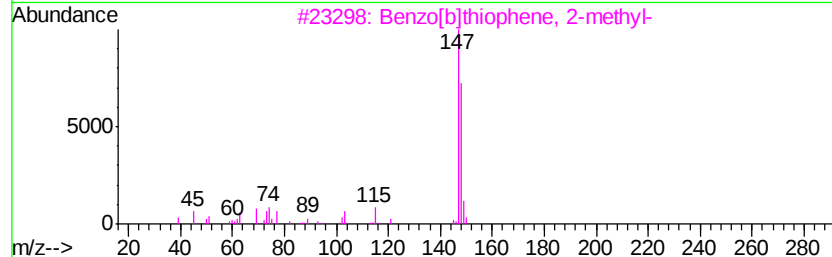
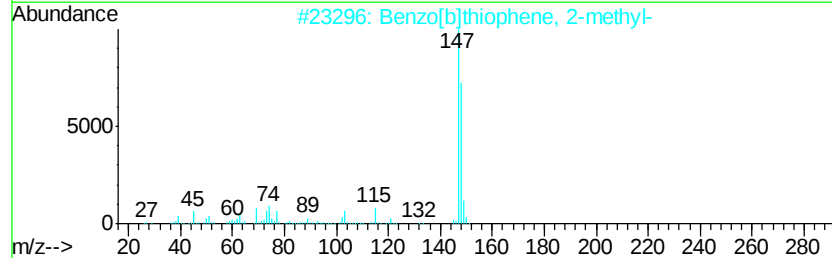
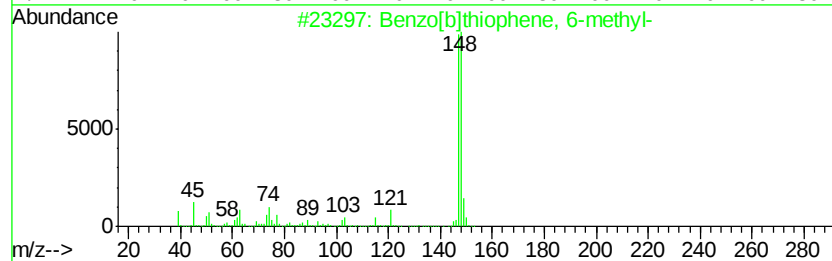
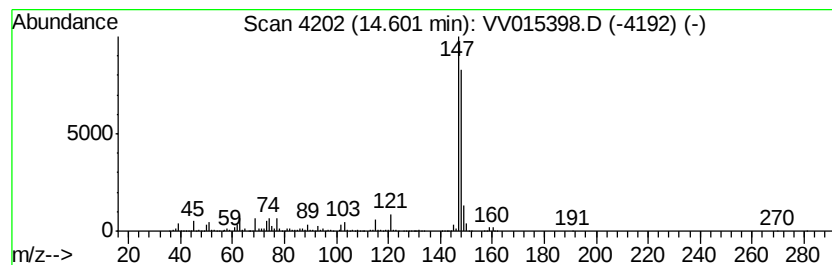
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MSVOA_V
ClientSampleId :
DBF06

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM040120WMA.M
Quant Title : VOC Analysis

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

Peak Number 22 Benzo[b]thiophene, 6-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	6.07 ug/L	2567830	1,4-Dichlorobenzene-d4	11.27
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzo[b]thiophene, 6-methyl-	148 C9H8S	016587-47-6 94
2		Benzo[b]thiophene, 2-methyl-	148 C9H8S	001195-14-8 91
3		Benzo[b]thiophene, 2-methyl-	148 C9H8S	001195-14-8 91
4		Benzo[b]thiophene, 5-methyl-	148 C9H8S	014315-14-1 91
5		Benzo[b]thiophene, 4-methyl-	148 C9H8S	014315-11-8 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBF06

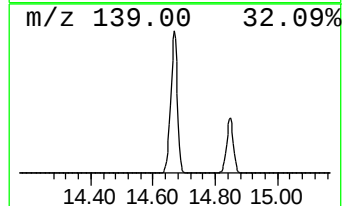
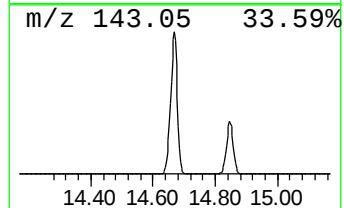
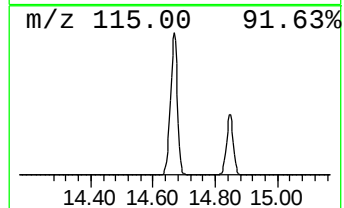
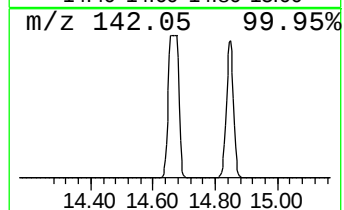
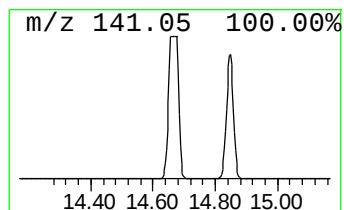
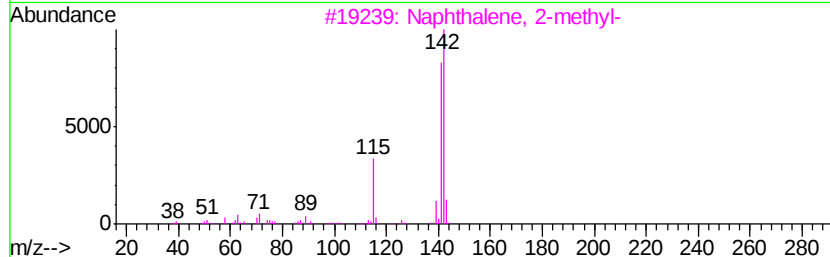
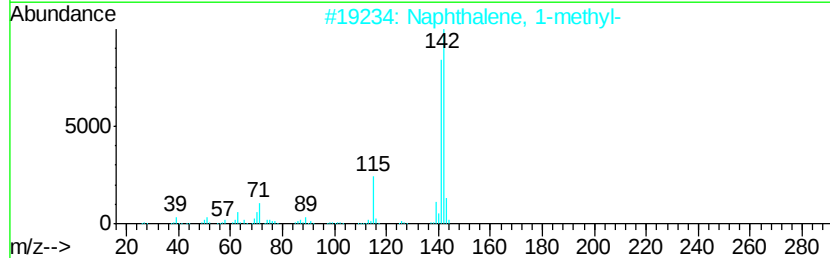
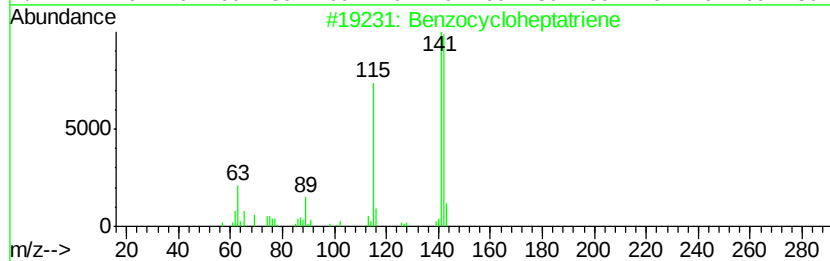
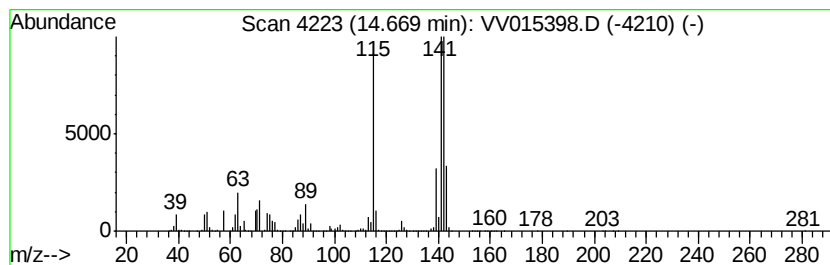
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Quant Title : VOC Analysis

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

Peak Number 23 Benzocycloheptatriene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	238.82 ug/L	101054000	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzocycloheptatriene	142	C11H10	000264-09-5	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
4		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	90
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV040320\
Data File : VV015398.D
Acq On : 03 Apr 2020 16:52
Operator : SY/MD
Sample : L2065-10
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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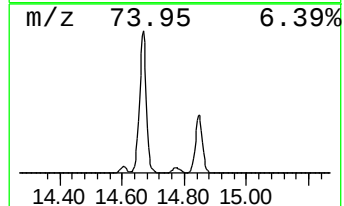
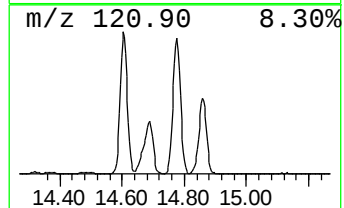
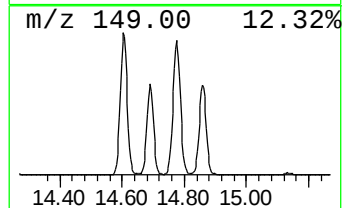
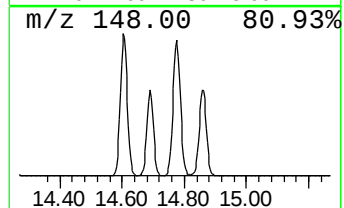
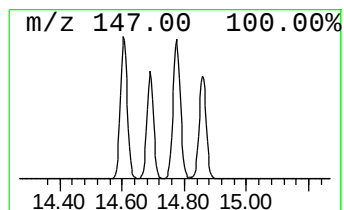
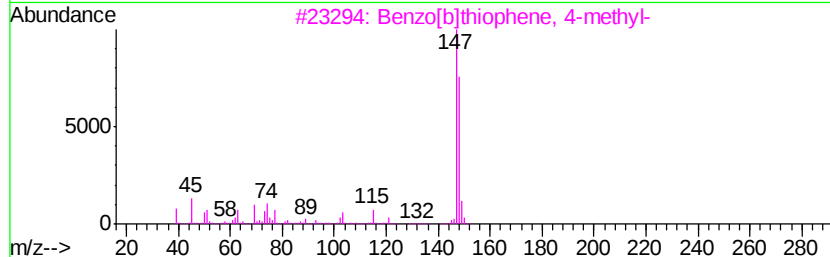
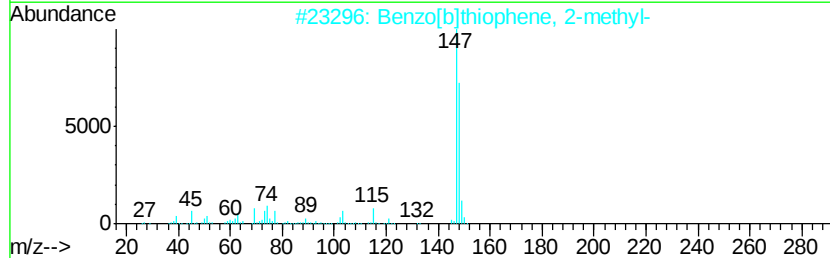
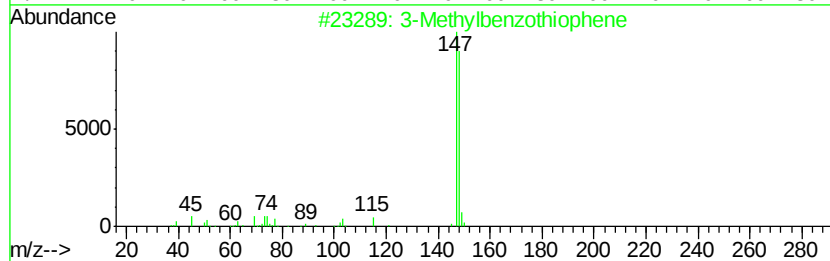
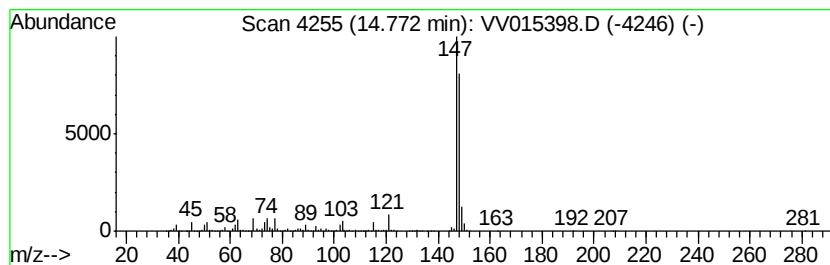
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Quant Title : VOC Analysis

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

Peak Number 24 3-Methylbenzothiophene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	6.18 ug/L	2615640	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Methylbenzothiophene	148	C9H8S	001455-18-1	94
2		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
3		Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	91
4		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
5		Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV040320\
 Data File : VV015398.D
 Acq On : 03 Apr 2020 16:52
 Operator : SY/MD
 Sample : L2065-10
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBF06

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM040120WMA.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-ethyl-...	10.47	15.2	ug/L	6421940	3	11.27	21157400	50.0
Benzene, 1,2,3-tr...	10.56	15.1	ug/L	6376750	3	11.27	21157400	50.0
Cyclohexane, 1-me...	10.63	9.0	ug/L	3799300	3	11.27	21157400	50.0
Benzene, 1-etheny...	11.01	7.3	ug/L	3098930	3	11.27	21157400	50.0
Benzofuran	11.19	16.1	ug/L	6830680	3	11.27	21157400	50.0
Indane	11.55	70.1	ug/L	29669100	3	11.27	21157400	50.0
Indene	11.77	109.3	ug/L	46234000	3	11.27	21157400	50.0
(+)-4-Carene	11.93	2.9	ug/L	1209070	3	11.27	21157400	50.0
Benzene, 2-ethyl-...	12.04	2.7	ug/L	1135570	3	11.27	21157400	50.0
Benzene, 1-etheny...	12.14	5.2	ug/L	2197530	3	11.27	21157400	50.0
1H-Indenol	12.39	11.9	ug/L	5045880	3	11.27	21157400	50.0
Benzofuran, 7-met...	12.49	22.2	ug/L	9410230	3	11.27	21157400	50.0
1H-Indene, 2,3-di...	12.75	8.3	ug/L	3496510	3	11.27	21157400	50.0
Benzene, 2-etheny...	12.91	14.1	ug/L	5979730	3	11.27	21157400	50.0
Cyclohexanemethan...	12.98	67.7	ug/L	28625300	3	11.27	21157400	50.0
2-Methylindene	13.08	13.3	ug/L	5620280	3	11.27	21157400	50.0
(+)-2-Bornanone	13.19	6.9	ug/L	2928700	3	11.27	21157400	50.0
Benzo[c]thiophene	13.66	52.3	ug/L	22135600	3	11.27	21157400	50.0
Benzo[b]thiophene...	14.60	6.1	ug/L	2567830	3	11.27	21157400	50.0
Benzocycloheptatr...	14.67	238.8	ug/L	101054000	3	11.27	21157400	50.0
3-Methylbenzothio...	14.77	6.2	ug/L	2615640	3	11.27	21157400	50.0