

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
 Data File : VU047647.D
 Acq On : 24 Mar 2022 21:08
 Operator : SY/MD
 Sample : N2081-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EW5X5

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_U\Method\SFAMULM032422WMA.M
 Title : VOC Analysis

Signal : TIC: VU047647.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.607	74	83	105	rVB2	367370	456519	8.31%	1.433%
2	1.919	171	180	182	rBV	81080	88938	1.62%	0.279%
3	2.478	351	354	363	rVB3	1153	1516	0.03%	0.005%
4	2.575	371	384	407	rBV3	189818	348636	6.35%	1.094%
5	3.359	615	628	643	rBV4	18992	39484	0.72%	0.124%
6	3.703	720	735	748	rVB4	4792	10115	0.18%	0.032%
7	3.874	767	788	820	rBV	729030	1738544	31.66%	5.458%
8	4.539	984	995	1008	rVB4	3337	7761	0.14%	0.024%
9	4.671	1013	1036	1079	rBV	1648282	3890955	70.85%	12.215%
10	5.066	1144	1159	1179	rBV	149510	326621	5.95%	1.025%
11	5.320	1219	1238	1256	rBV	594596	1332566	24.26%	4.183%
12	5.591	1316	1322	1331	rBV6	1393	2284	0.04%	0.007%
13	5.729	1343	1365	1382	rBV2	297772	805301	14.66%	2.528%
14	5.976	1437	1442	1443	rBV2	1802	1364	0.02%	0.004%
15	6.137	1483	1492	1503	rBV7	2851	5897	0.11%	0.019%
16	6.250	1511	1527	1547	rBV	206441	389293	7.09%	1.222%
17	6.546	1605	1619	1641	rVB2	99767	189702	3.45%	0.596%
18	6.693	1651	1665	1678	rBV	180713	353948	6.45%	1.111%
19	6.764	1680	1687	1703	rVB2	33526	64087	1.17%	0.201%
20	6.980	1749	1754	1765	rVB7	1892	3180	0.06%	0.010%
21	7.237	1827	1834	1843	rVB6	1842	3128	0.06%	0.010%
22	7.568	1923	1937	1956	rBV	102284	183986	3.35%	0.578%
23	7.655	1961	1964	1970	rVB5	749	870	0.02%	0.003%
24	7.793	2000	2007	2017	rVV4	4997	8937	0.16%	0.028%
25	7.857	2017	2027	2029	rBV4	11563	16361	0.30%	0.051%
26	7.899	2029	2040	2052	rVV	376893	655474	11.94%	2.058%
27	7.970	2053	2062	2075	rVB	127485	215588	3.93%	0.677%
28	8.179	2117	2127	2139	rBV	72109	119572	2.18%	0.375%
29	8.365	2177	2185	2193	rBV5	5907	11026	0.20%	0.035%
30	8.471	2211	2218	2231	rVB3	3455	6154	0.11%	0.019%
31	8.555	2231	2244	2258	rBV	303568	516146	9.40%	1.620%
32	8.635	2258	2269	2304	rVB	264639	496246	9.04%	1.558%
33	8.809	2313	2323	2331	rBV7	6644	12089	0.22%	0.038%
34	8.864	2332	2340	2350	rVV2	22904	37307	0.68%	0.117%
35	8.925	2350	2359	2369	rVB3	9406	15705	0.29%	0.049%
36	9.070	2399	2404	2412	rVB4	998	1258	0.02%	0.004%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
 Title : VOC Analysis

37	9.166	2427	2434	2442	rVB7	2722	3675	0.07%	0.012%
38	9.279	2460	2469	2472	rBV4	1180	1731	0.03%	0.005%
39	9.333	2483	2486	2495	rVB2	813	848	0.02%	0.003%
40	9.417	2499	2512	2533	rBV	306481	520612	9.48%	1.634%
41	9.510	2534	2541	2549	rVV2	9475	14623	0.27%	0.046%
42	9.571	2549	2560	2579	rVV	223843	363076	6.61%	1.140%
43	9.693	2585	2598	2615	rBV	229468	389959	7.10%	1.224%
44	9.889	2651	2659	2669	rVV4	4580	7336	0.13%	0.023%
45	10.037	2690	2705	2711	rBV6	6675	13946	0.25%	0.044%
46	10.102	2713	2725	2751	rVB	256271	429073	7.81%	1.347%
47	10.237	2757	2767	2770	rBV8	4538	6843	0.12%	0.021%
48	10.272	2771	2778	2790	rVB3	24775	41364	0.75%	0.130%
49	10.362	2795	2806	2819	rVB8	11933	24624	0.45%	0.077%
50	10.442	2825	2831	2833	rBV3	1123	957	0.02%	0.003%
51	10.484	2833	2844	2855	rVV	74873	117987	2.15%	0.370%
52	10.574	2864	2872	2880	rVB2	3308	4597	0.08%	0.014%
53	10.645	2887	2894	2897	rBV7	1530	2212	0.04%	0.007%
54	10.696	2901	2910	2918	rVB5	11131	17122	0.31%	0.054%
55	10.758	2918	2929	2941	rBV	277368	450295	8.20%	1.414%
56	10.815	2941	2947	2957	rVB9	8506	15044	0.27%	0.047%
57	10.905	2957	2975	2992	rBV	328907	517189	9.42%	1.624%
58	10.999	2992	3004	3009	rBV	319875	536034	9.76%	1.683%
59	11.089	3021	3032	3050	rVB	866166	1324520	24.12%	4.158%
60	11.230	3070	3076	3082	rBV4	2786	4356	0.08%	0.014%
61	11.291	3082	3095	3116	rVB	849677	1301780	23.70%	4.087%
62	11.420	3126	3135	3137	rBV3	8794	10903	0.20%	0.034%
63	11.468	3137	3150	3175	rVB	3639954	5491789	100.00%	17.240%
64	11.610	3183	3194	3198	rBV	42468	65067	1.18%	0.204%
65	11.642	3199	3204	3218	rVB	80507	119175	2.17%	0.374%
66	11.741	3225	3235	3244	rBV	119017	182979	3.33%	0.574%
67	11.812	3244	3257	3270	rVB2	412609	728113	13.26%	2.286%
68	11.893	3270	3282	3297	rBV	1025603	1600222	29.14%	5.024%
69	12.008	3310	3318	3331	rVB2	22521	35810	0.65%	0.112%
70	12.095	3331	3345	3351	rBV3	408405	688310	12.53%	2.161%
71	12.192	3363	3375	3392	rVB4	758070	1393918	25.38%	4.376%
72	12.301	3399	3409	3421	rBV2	39228	62484	1.14%	0.196%
73	12.375	3421	3432	3445	rVB	130328	199896	3.64%	0.628%
74	12.465	3448	3460	3464	rBV	174956	271069	4.94%	0.851%
75	12.571	3483	3493	3501	rBV	252034	377173	6.87%	1.184%
76	12.684	3518	3528	3548	rBV4	112657	303271	5.52%	0.952%

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77	12.809	3561	3567	3575	rVB3	17728	24994	0.46%	0.078%
78	12.873	3576	3587	3599	rVB3	97133	164360	2.99%	0.516%
79	12.973	3599	3618	3626	rBV	122256	206221	3.76%	0.647%
80	13.024	3626	3634	3649	rVB	208553	317420	5.78%	0.996%
81	13.108	3650	3660	3666	rBV3	21040	36923	0.67%	0.116%
82	13.253	3699	3705	3710	rBV3	12326	17610	0.32%	0.055%
83	13.291	3711	3717	3726	rVB3	40716	58069	1.06%	0.182%
84	13.404	3744	3752	3756	rBV3	16756	24802	0.45%	0.078%
85	13.449	3757	3766	3781	rVB3	222299	377199	6.87%	1.184%
86	13.622	3811	3820	3837	rVB	71196	117033	2.13%	0.367%
87	13.732	3845	3854	3862	rBV5	9406	15888	0.29%	0.050%
88	13.783	3863	3870	3877	rVV3	8660	12727	0.23%	0.040%
89	13.838	3878	3887	3902	rVV7	23138	48559	0.88%	0.152%
90	14.085	3949	3964	3983	rBV	187646	311594	5.67%	0.978%
91	14.188	3990	3996	4007	rVB8	4194	6990	0.13%	0.022%
92	14.256	4011	4017	4021	rBV6	1646	2183	0.04%	0.007%
93	14.291	4021	4028	4037	rVV8	5157	8412	0.15%	0.026%
94	14.487	4082	4089	4094	rBV5	3461	4597	0.08%	0.014%
95	14.683	4137	4150	4160	rBV8	8225	17708	0.32%	0.056%
96	14.806	4184	4188	4196	rVB7	2122	2451	0.04%	0.008%
97	14.873	4205	4209	4217	rVB7	2069	3118	0.06%	0.010%
98	15.015	4246	4253	4265	rVB6	4867	7757	0.14%	0.024%
99	15.224	4307	4318	4328	rBV	25225	40392	0.74%	0.127%
100	15.410	4367	4376	4389	rBV	19205	30980	0.56%	0.097%

Sum of corrected areas: 31854527

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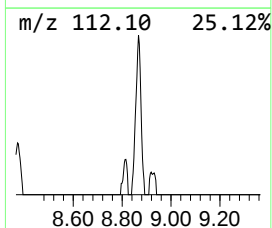
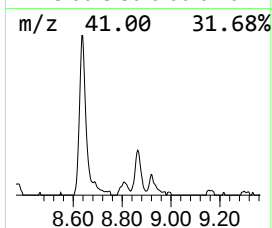
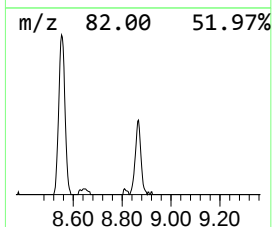
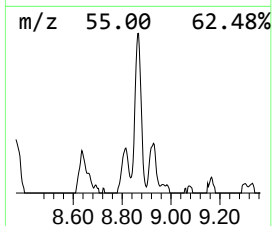
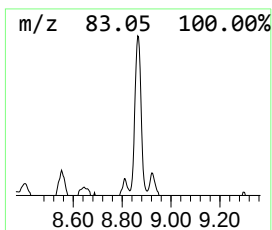
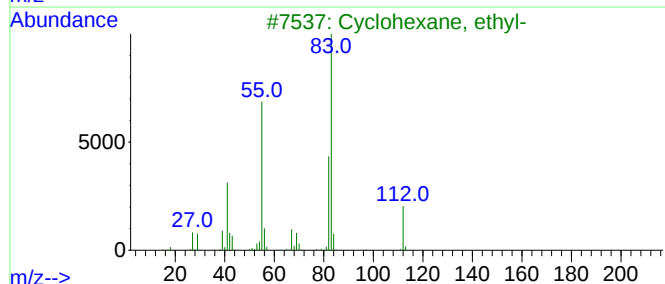
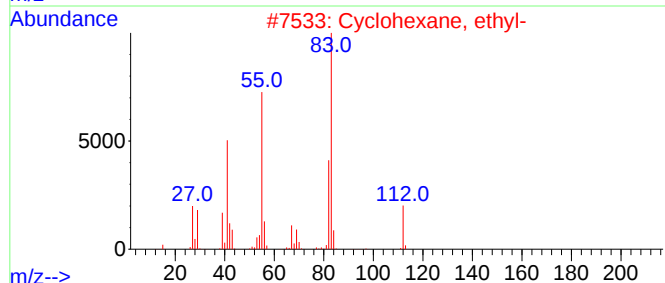
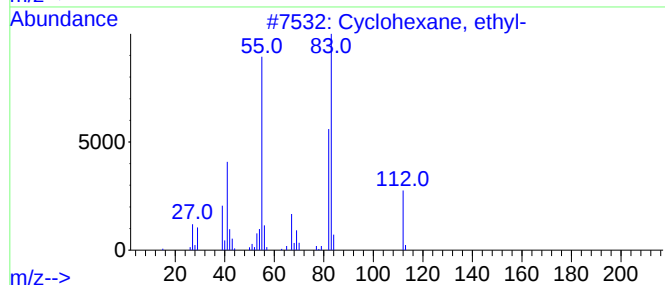
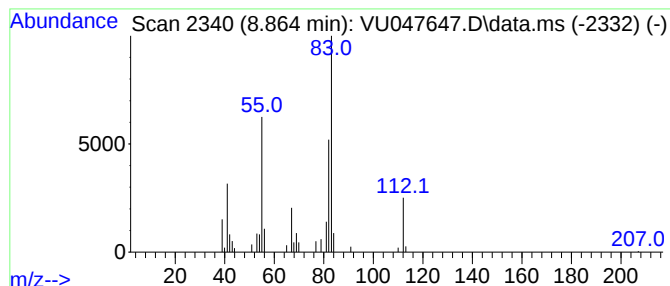
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Cyclic8.864 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.864	3.58 ug/L	37307	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	93
2			Cyclohexane, ethyl-	112	C8H16	001678-91-7	90
3			Cyclohexane, ethyl-	112	C8H16	001678-91-7	87
4			Cyclohexane, ethyl-	112	C8H16	001678-91-7	87
5			Cyclohexane, ethyl-	112	C8H16	001678-91-7	87



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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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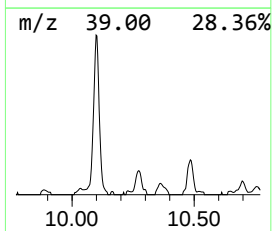
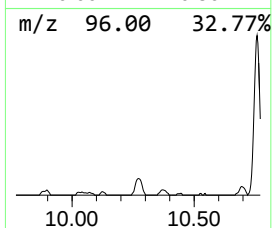
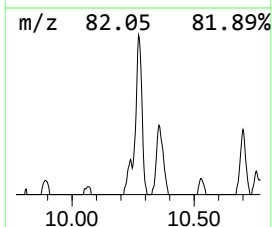
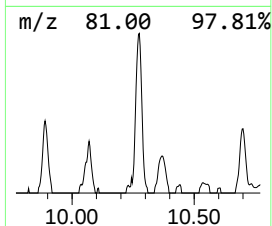
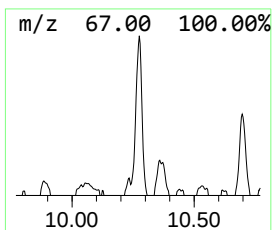
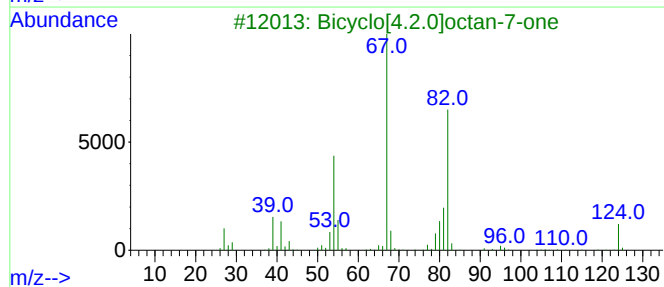
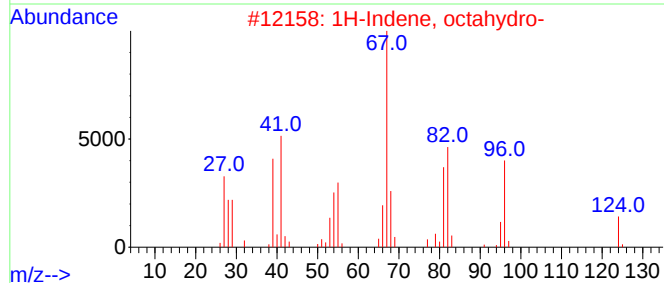
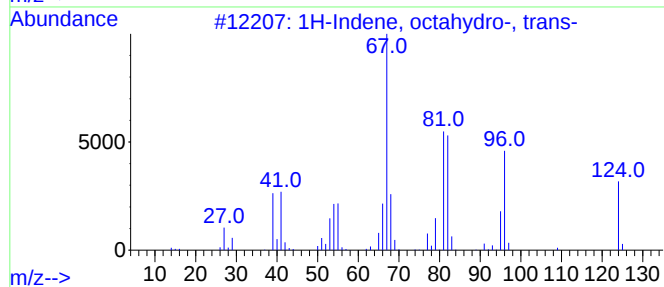
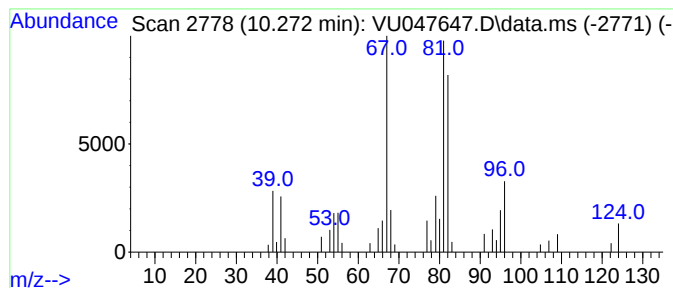
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.272	3.97 ug/L	41364	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-, trans-			124	C9H16	003296-50-2	47
2	1H-Indene, octahydro-			124	C9H16	000496-10-6	46
3	Bicyclo[4.2.0]octan-7-one			124	C8H12O	054211-18-6	43
4	Cyclopropane, trimethylmethylene-			96	C7H12	034462-28-7	43
5	2,4-Hexadiene			82	C6H10	000592-46-1	43



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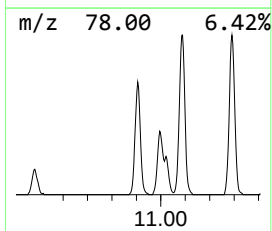
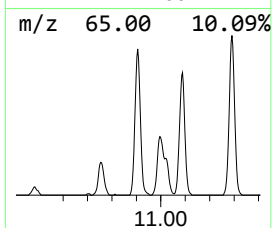
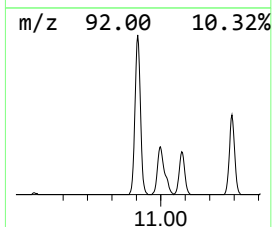
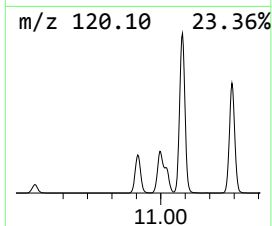
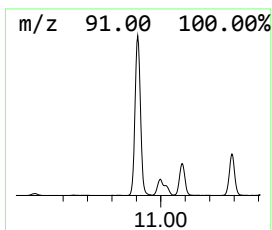
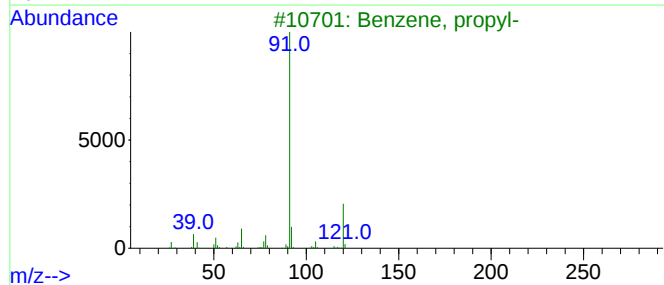
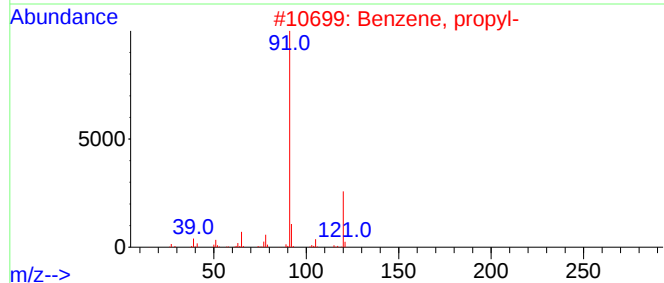
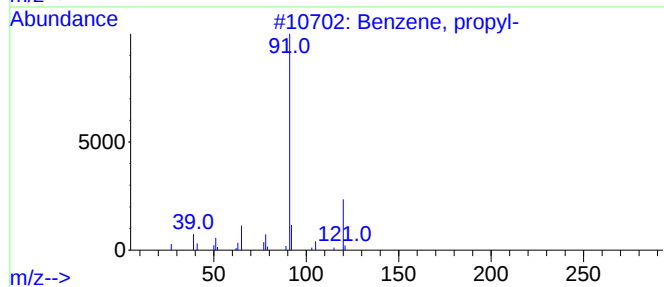
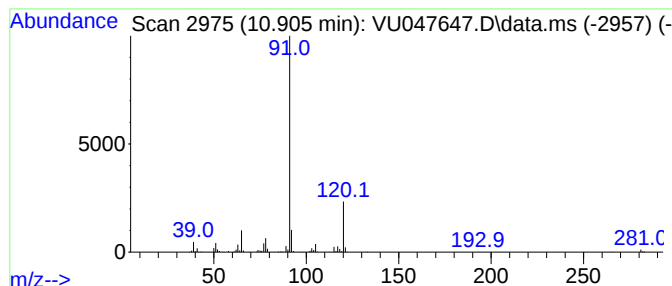
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Benzene, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	35.52 ug/L	517189	1,4-Dichlorobenzene-d4	11.812

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	87
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			Benzene, propyl-	120	C9H12	000103-65-1	87
5			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64



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ClientSampleId :
EW5X5

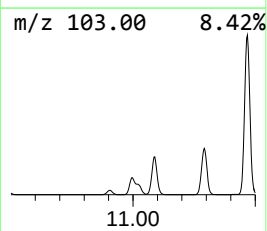
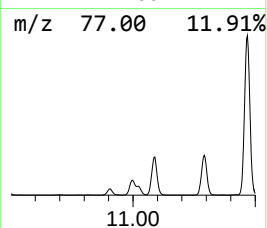
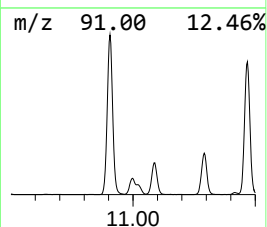
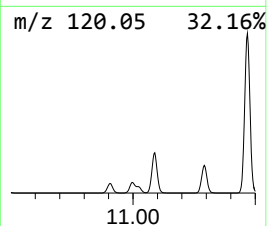
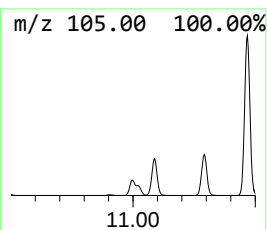
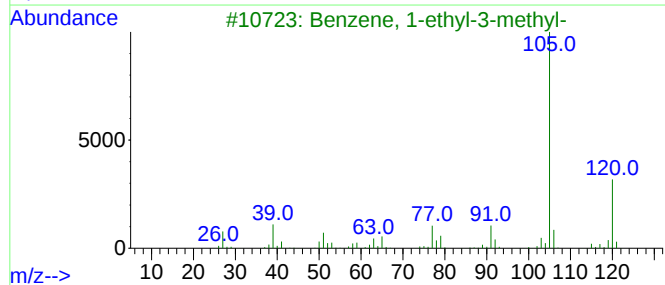
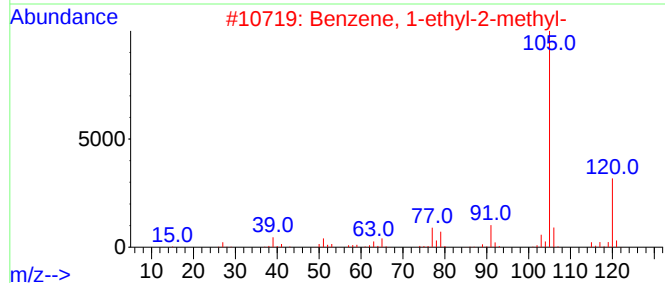
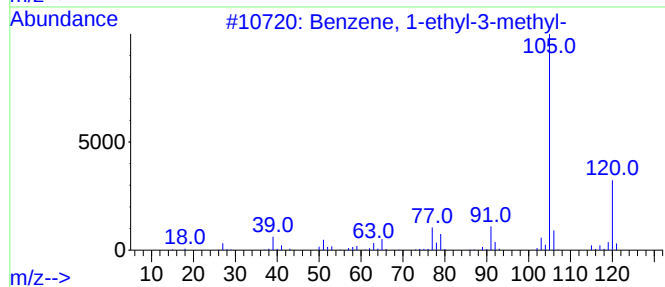
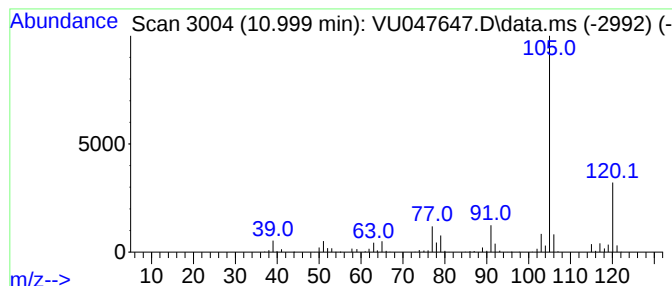
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	36.81 ug/L	536034	1,4-Dichlorobenzene-d4	11.812

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-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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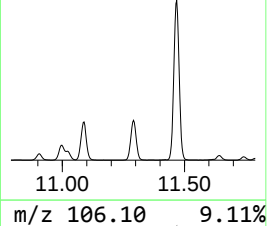
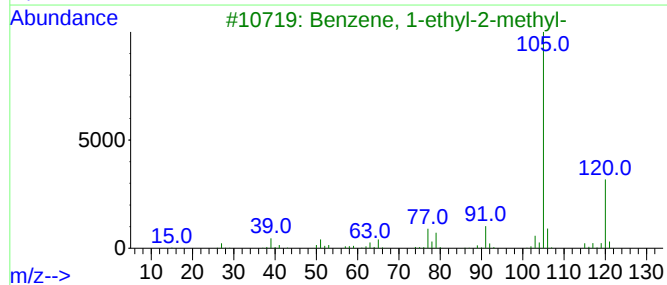
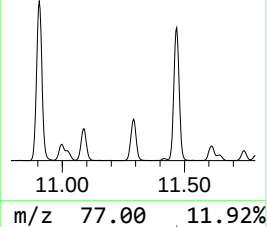
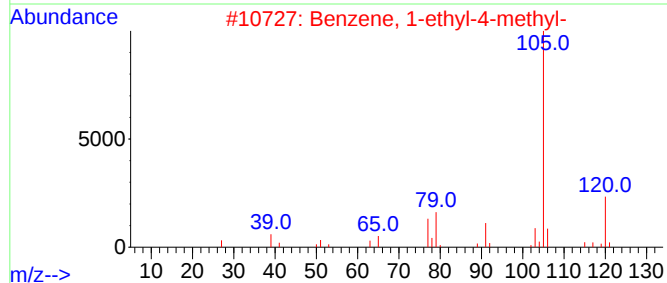
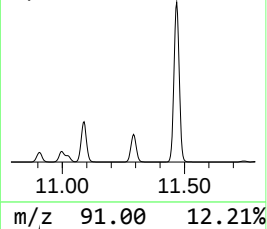
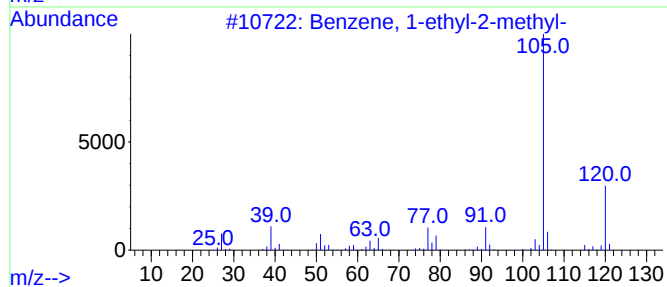
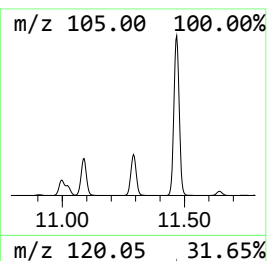
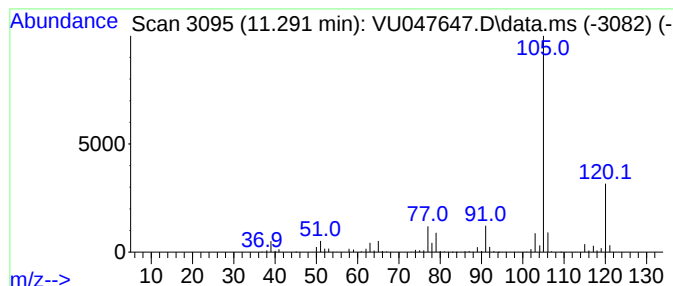
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	89.39 ug/L	1301780	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4			Mesitylene	120	C9H12	000108-67-8	90
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

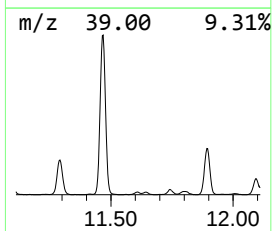
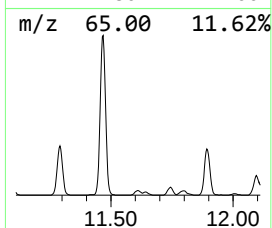
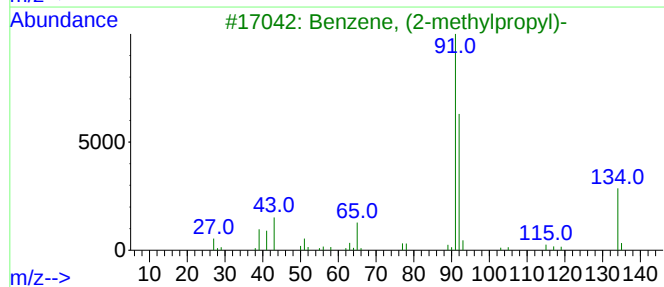
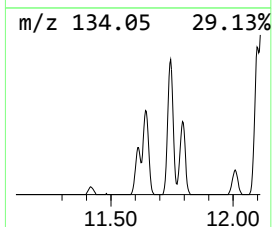
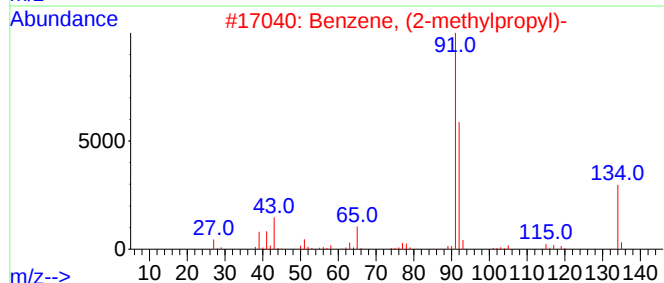
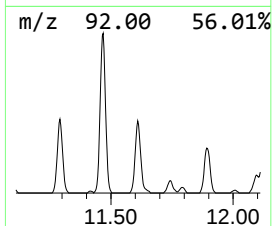
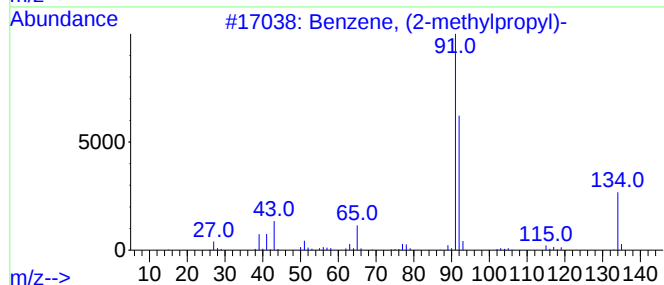
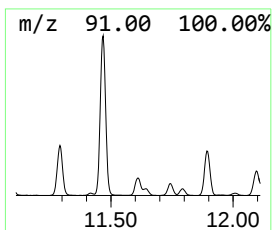
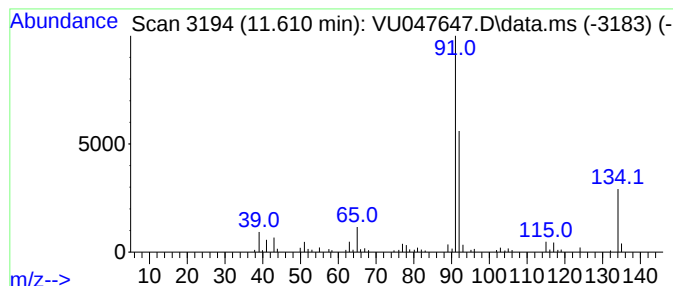
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Quant Title : VOC Analysis

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

Peak Number 7 Benzene, (2-methylpropyl)- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	4.47 ug/L	65067	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
4			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
5			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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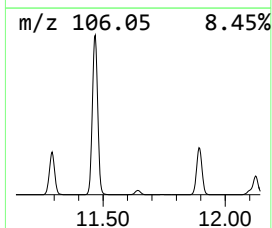
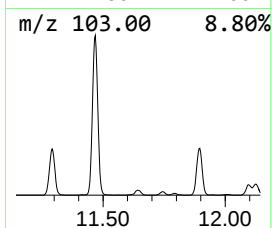
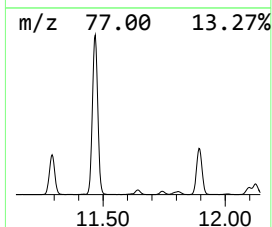
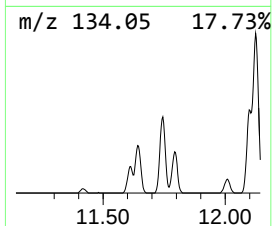
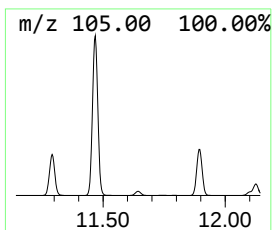
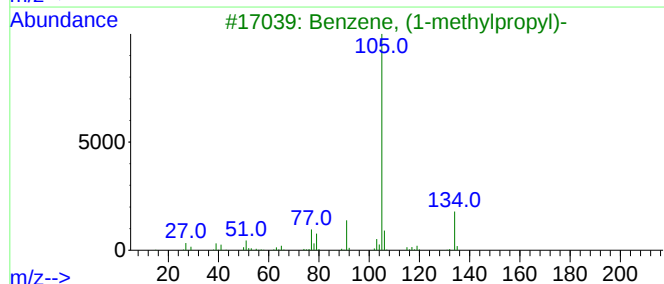
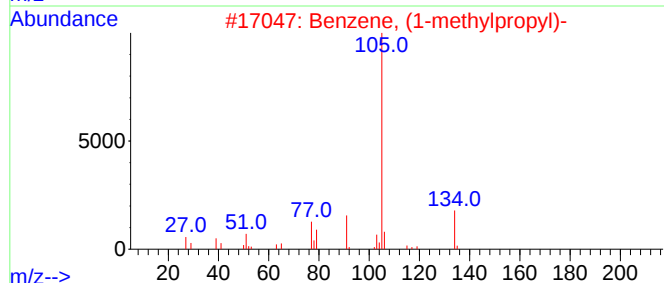
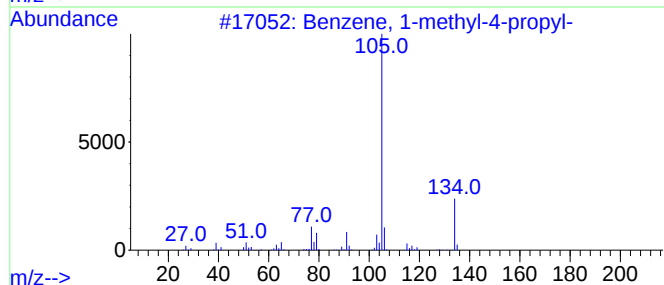
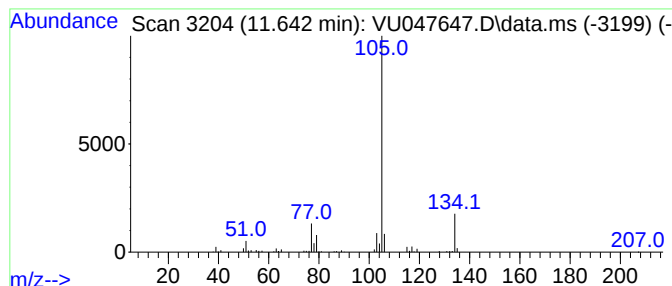
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Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1-methyl-4-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.642	8.18 ug/L	119175	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

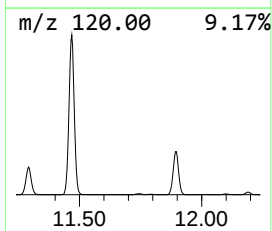
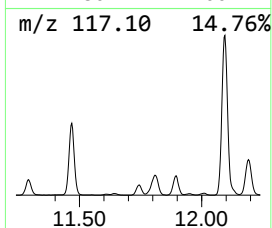
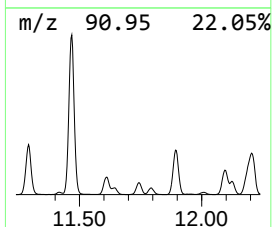
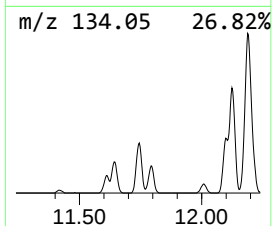
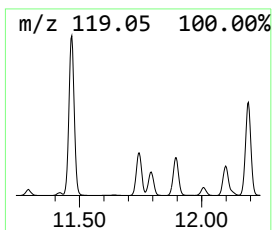
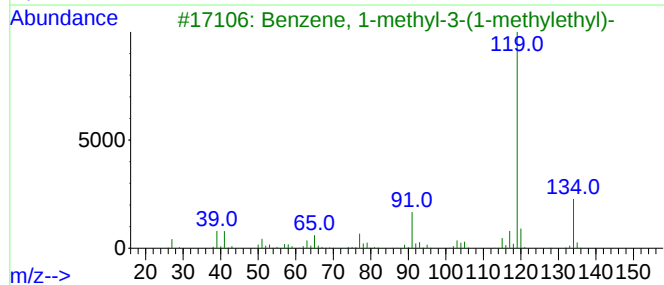
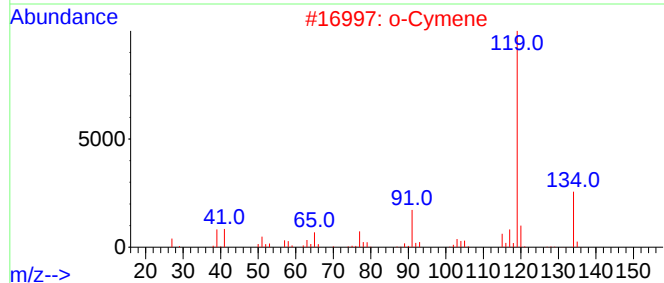
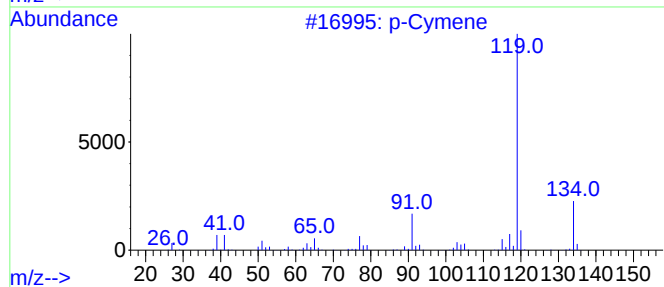
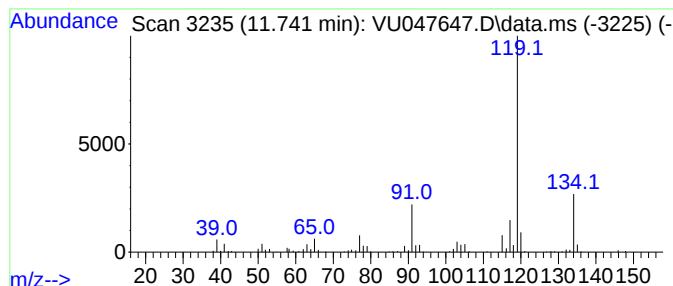
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Quant Title : VOC Analysis

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

Peak Number 9 p-Cymene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.741	12.57 ug/L	182979	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			p-Cymene	134	C10H14	000099-87-6	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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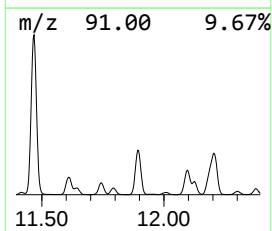
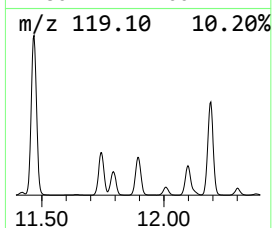
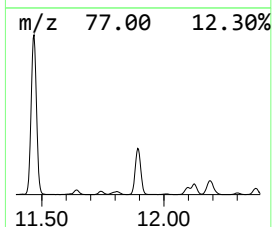
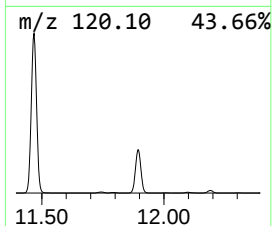
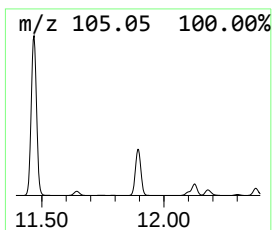
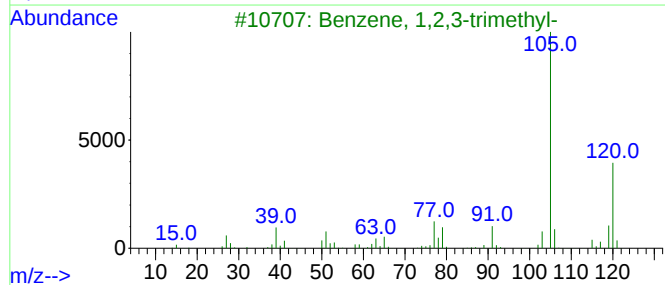
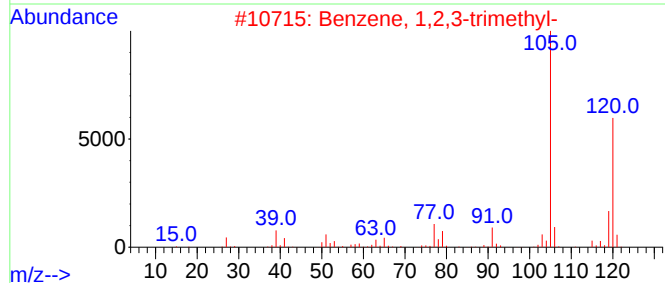
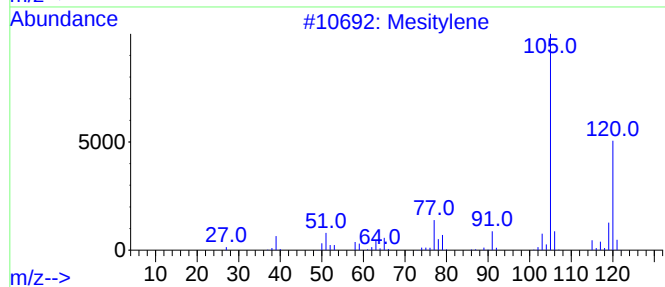
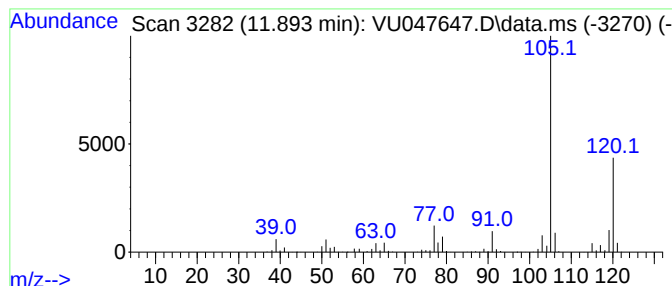
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Quant Title : VOC Analysis

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

Peak Number 10 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.893	109.89 ug/L	1600220	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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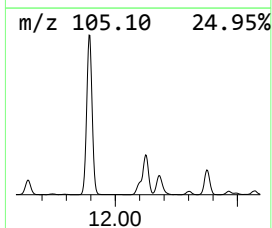
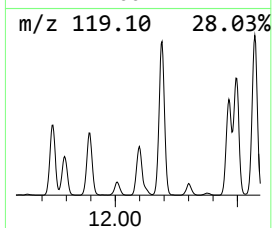
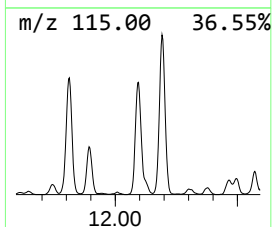
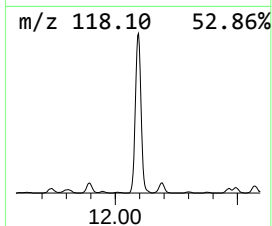
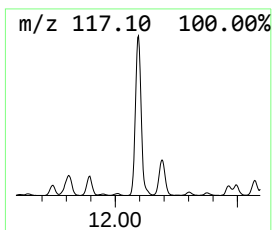
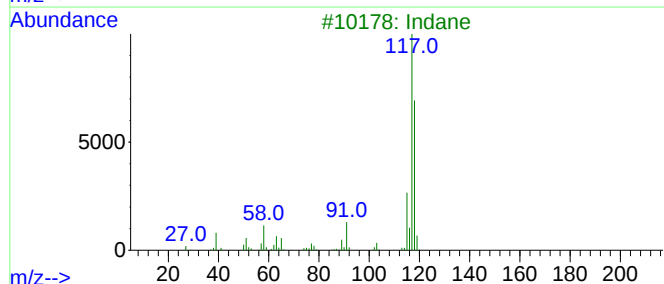
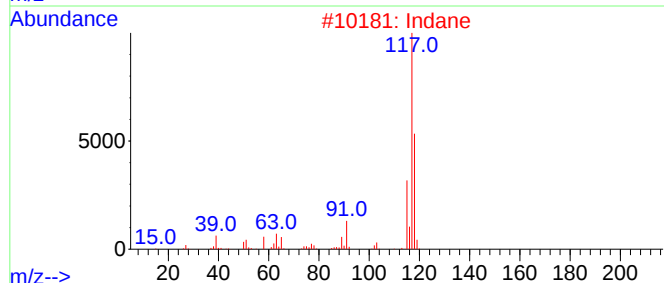
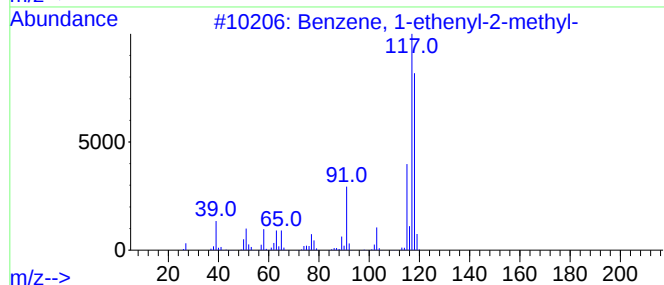
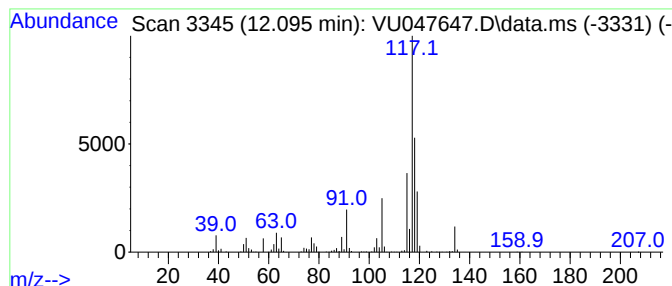
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Quant Title : VOC Analysis

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

Peak Number 11 Benzene, 1-ethenyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	47.27 ug/L	688310	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	83
2			Indane	118	C9H10	000496-11-7	64
3			Indane	118	C9H10	000496-11-7	64
4			Deltacyclene	118	C9H10	007785-10-6	64
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

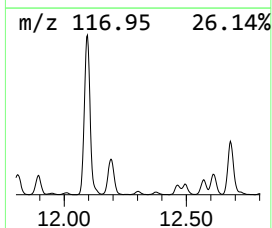
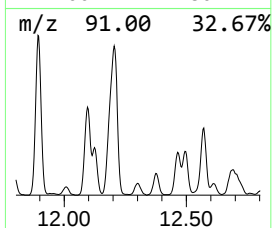
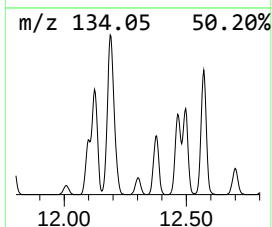
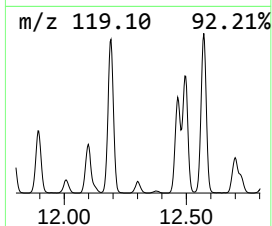
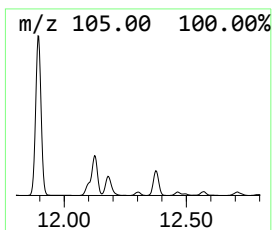
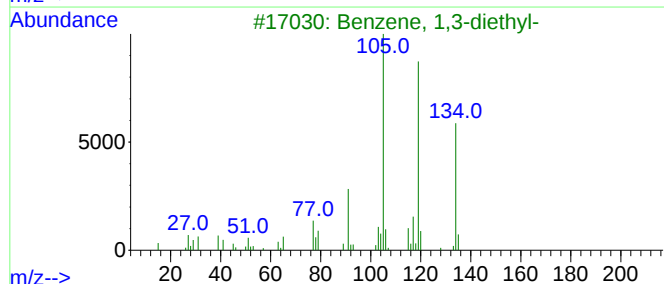
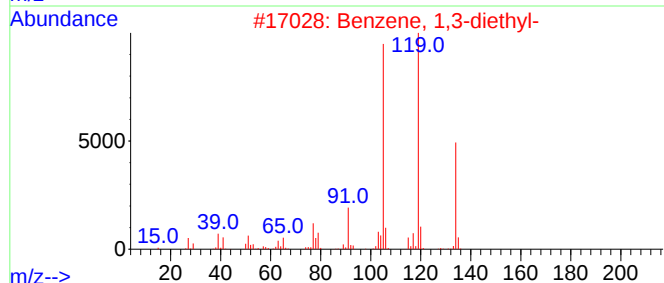
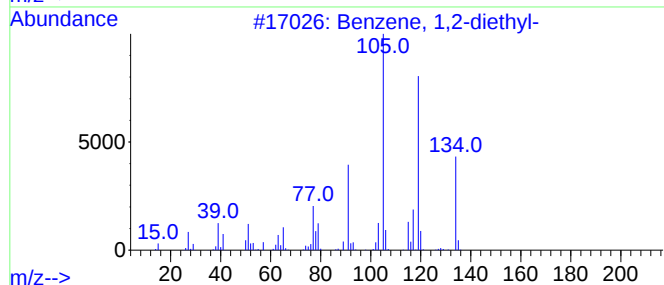
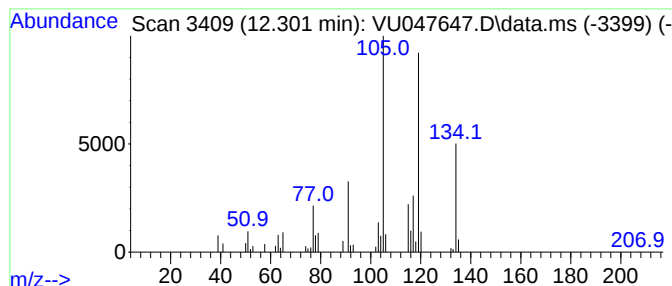
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, 1,2-diethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.301	4.29 ug/L	62484	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	90
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	81
4			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	81
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

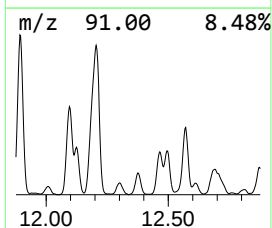
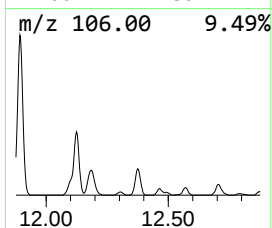
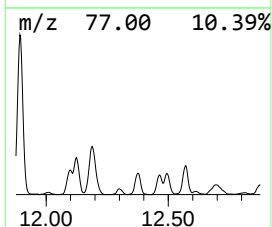
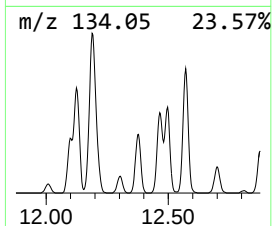
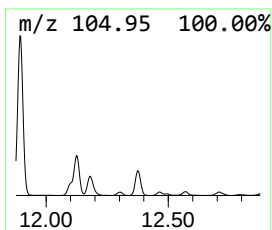
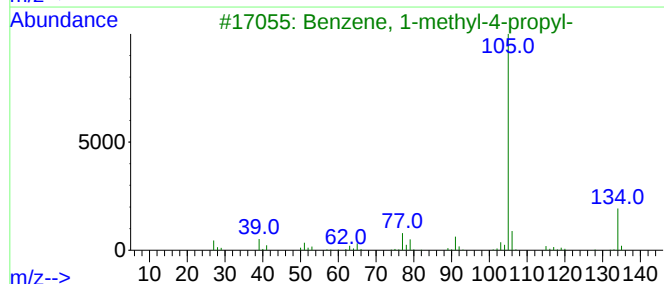
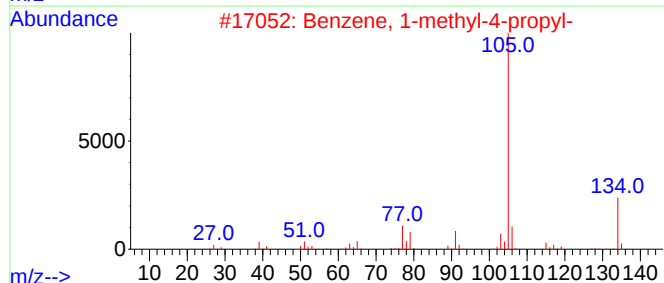
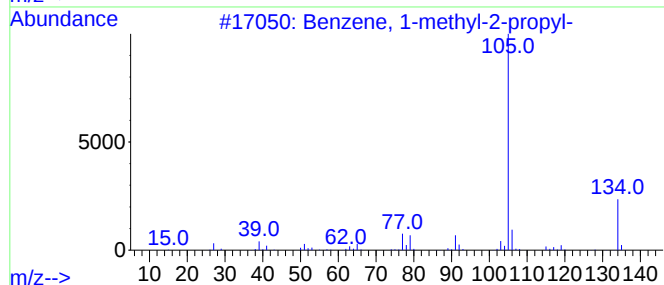
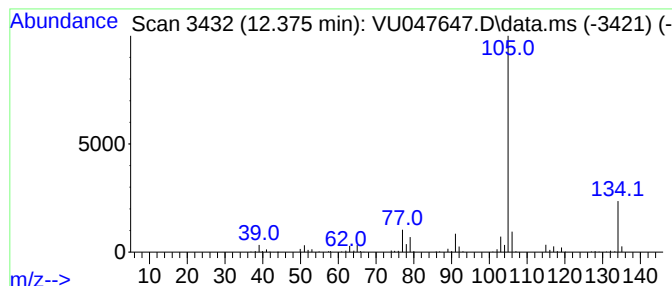
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Quant Title : VOC Analysis

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

Peak Number 13 Benzene, 1-methyl-2-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.375	13.73 ug/L	199896	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

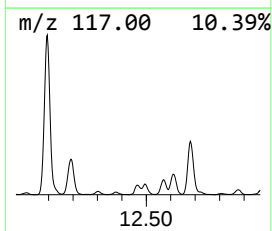
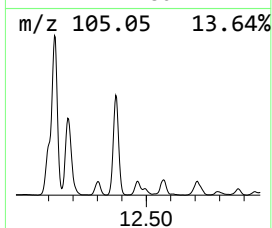
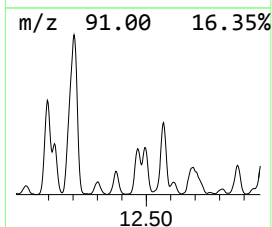
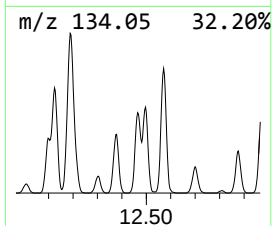
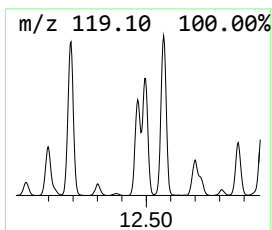
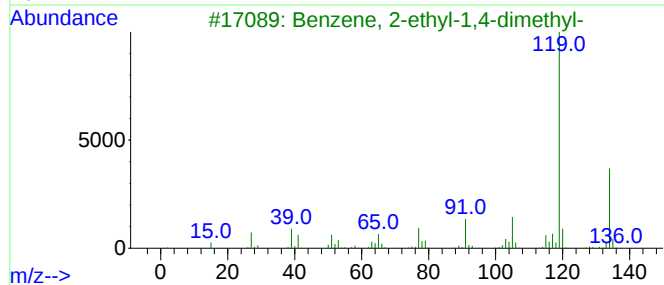
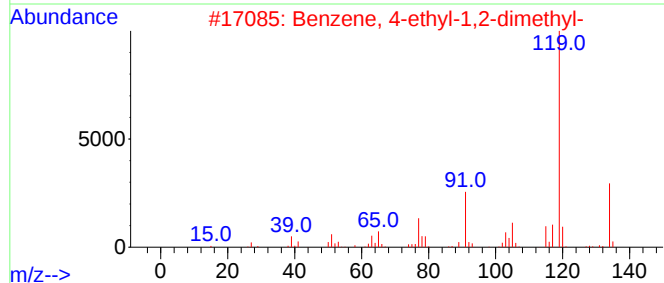
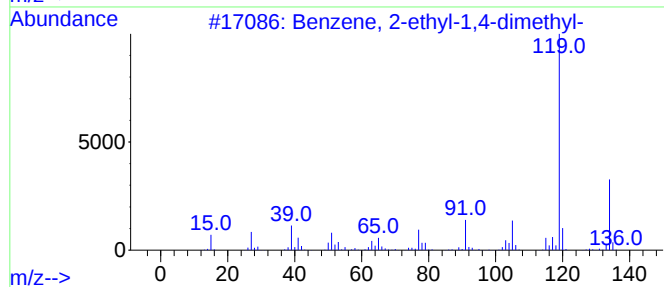
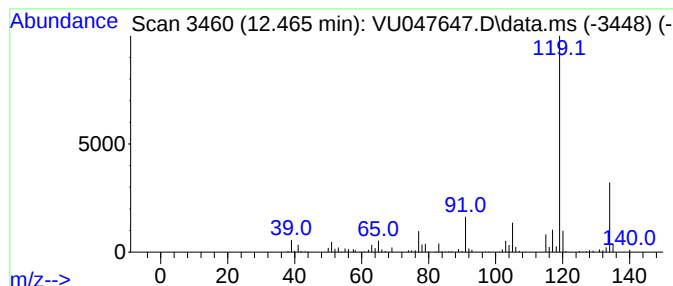
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.465	18.61 ug/L	271069	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

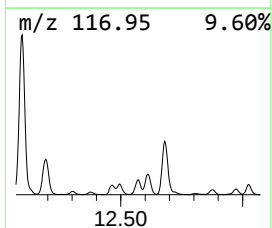
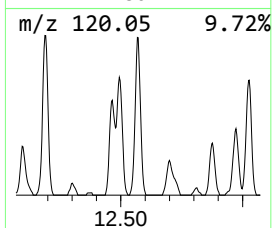
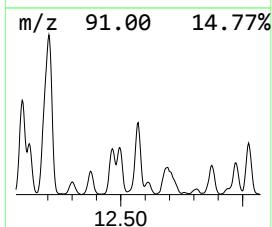
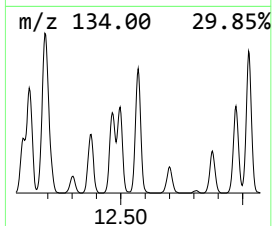
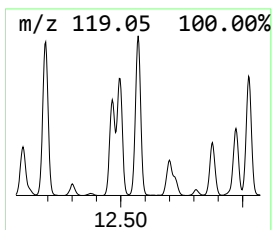
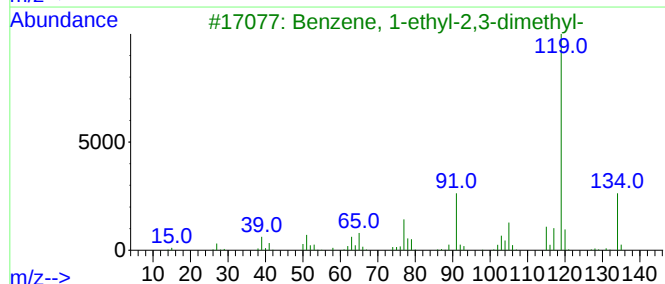
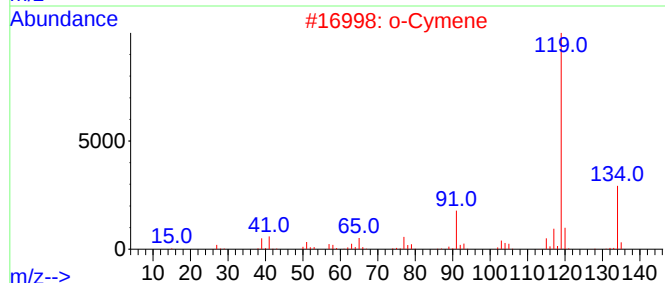
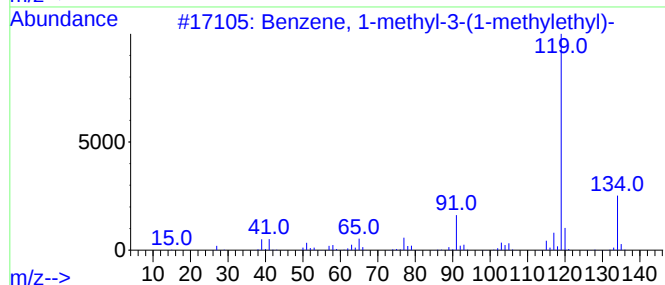
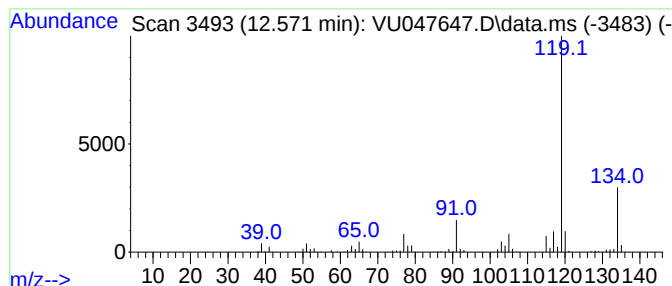
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Quant Title : VOC Analysis

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

Peak Number 15 Benzene, 1-methyl-3-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	25.90 ug/L	377173	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

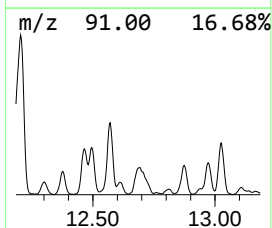
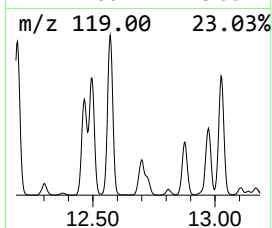
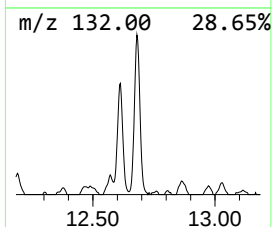
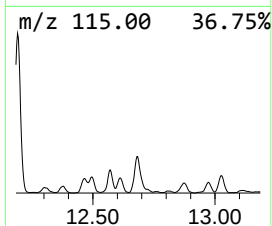
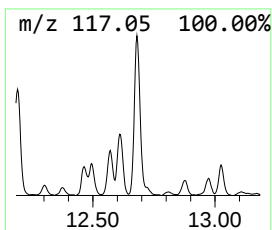
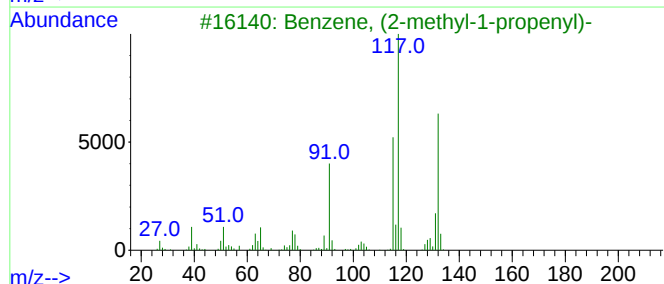
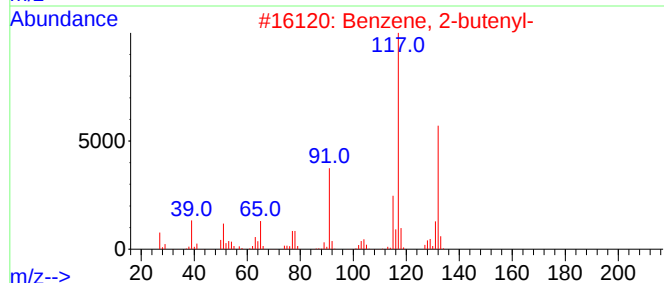
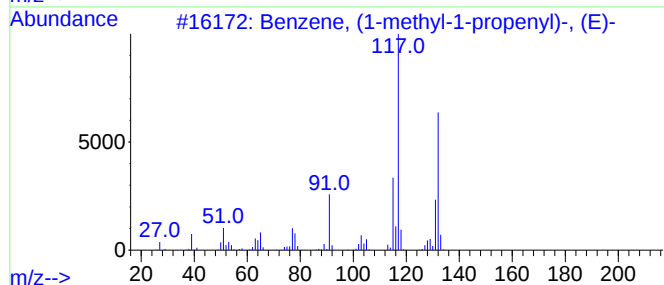
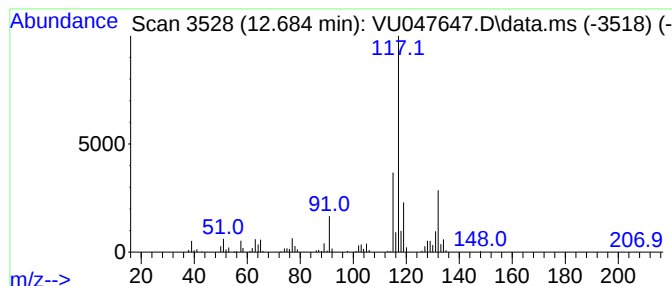
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Quant Title : VOC Analysis

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

Peak Number 16 Benzene, (1-methyl-1-propen... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.684	20.83 ug/L	303271	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
4			Indan, 1-methyl-	132	C10H12	000767-58-8	81
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

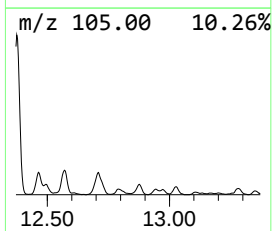
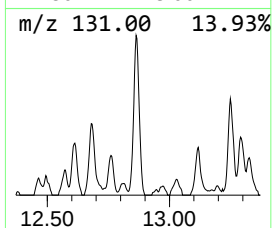
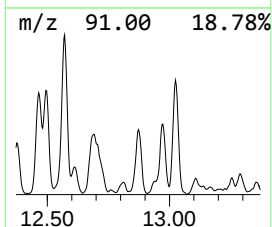
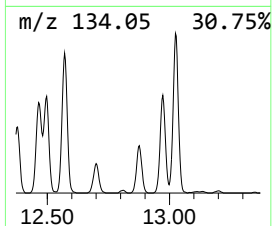
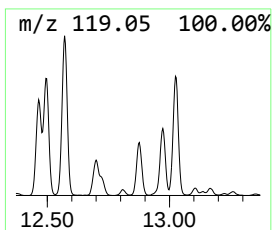
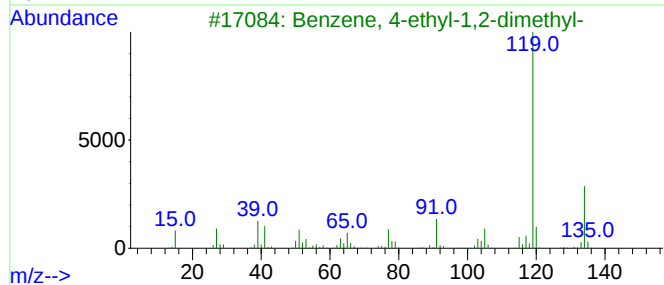
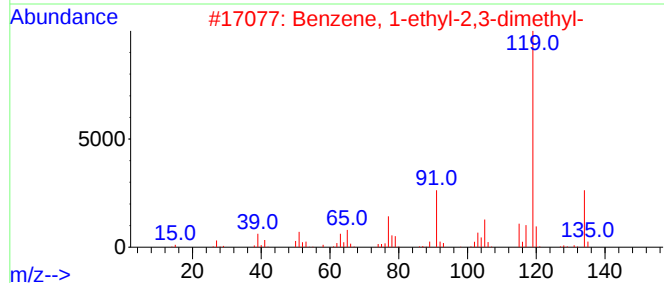
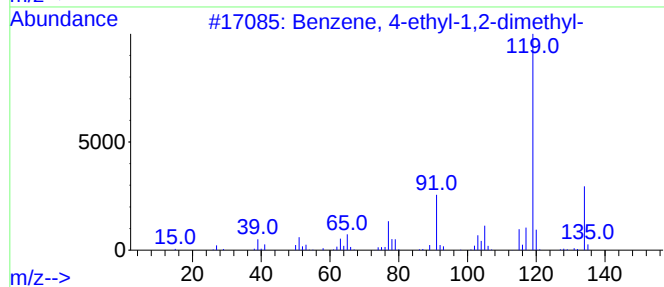
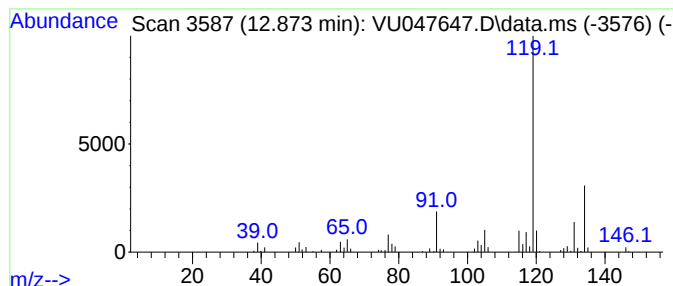
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Quant Title : VOC Analysis

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

Peak Number 17 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.873	11.29 ug/L	164360	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	91
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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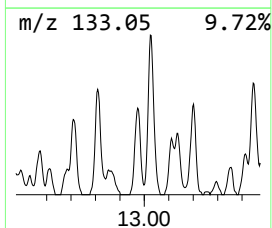
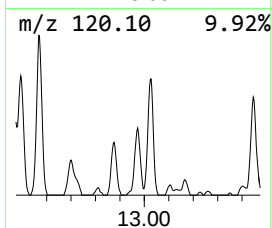
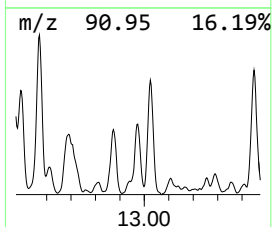
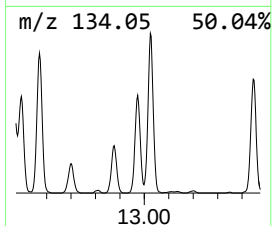
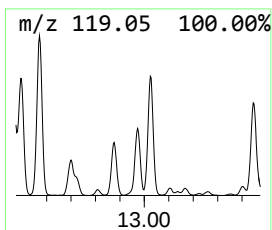
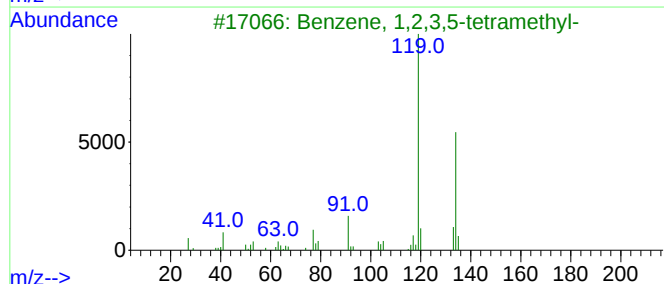
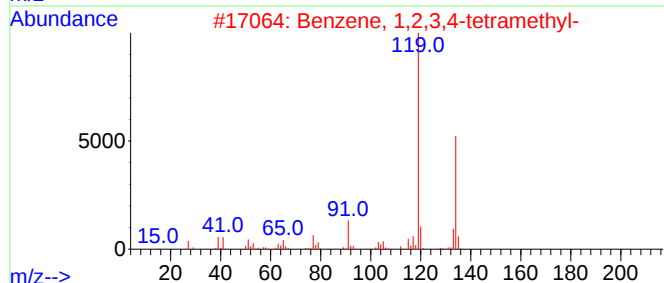
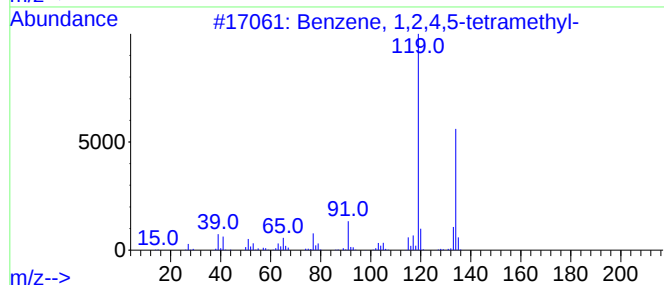
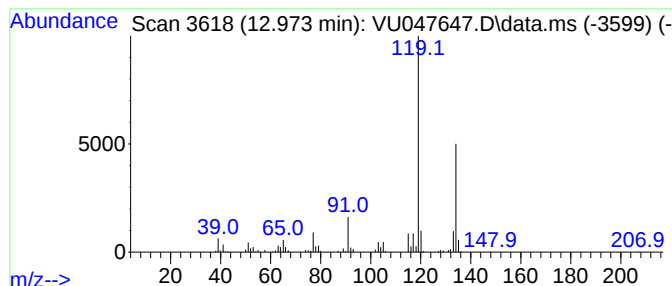
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Quant Title : VOC Analysis

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

Peak Number 18 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.973	14.16 ug/L	206221	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

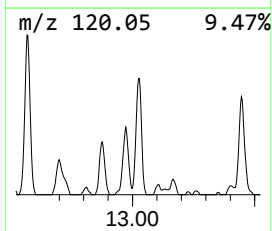
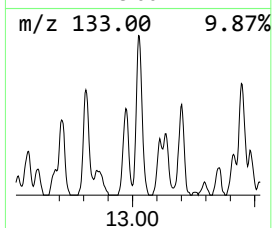
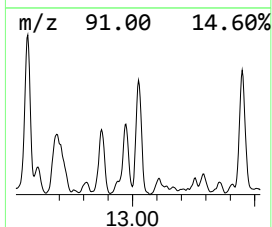
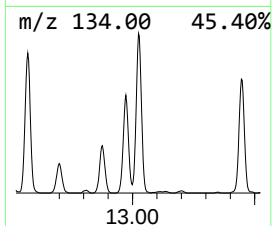
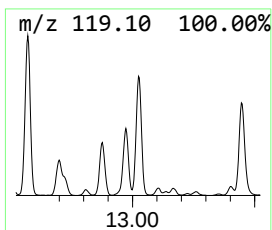
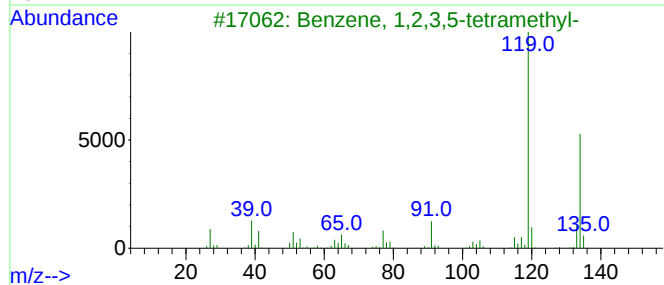
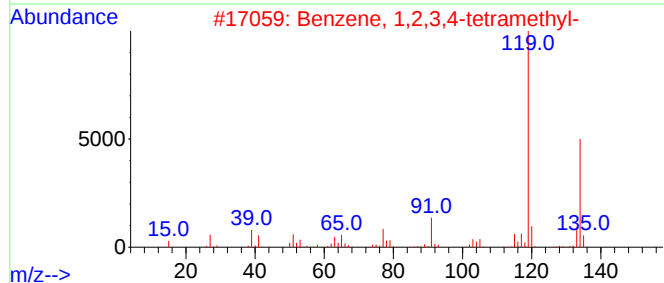
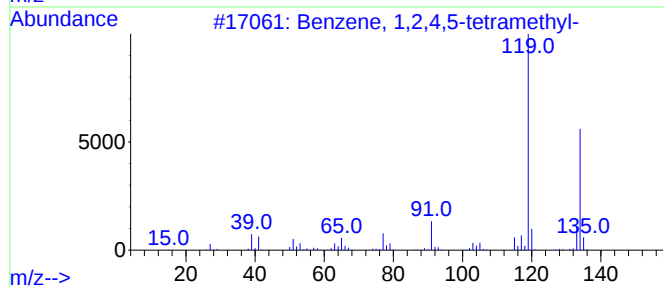
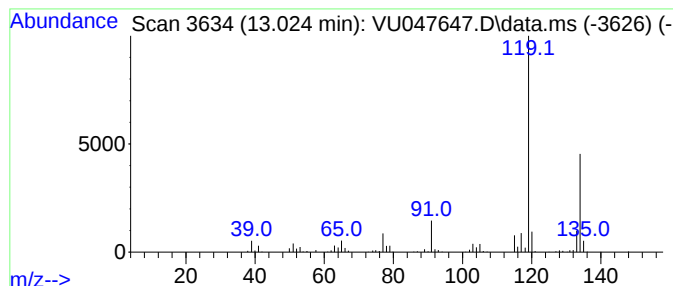
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Quant Title : VOC Analysis

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

Peak Number 19 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.024	21.80 ug/L	317420	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

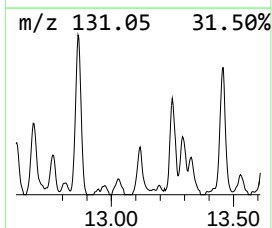
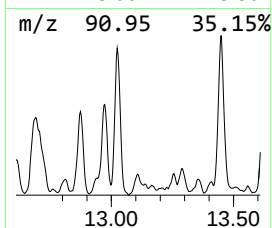
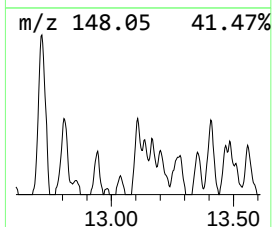
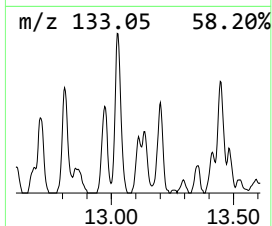
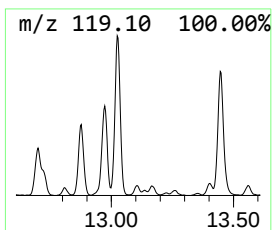
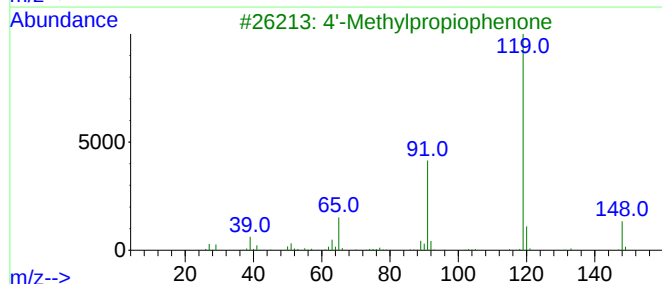
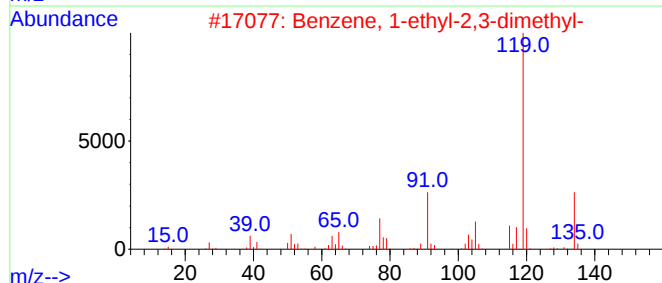
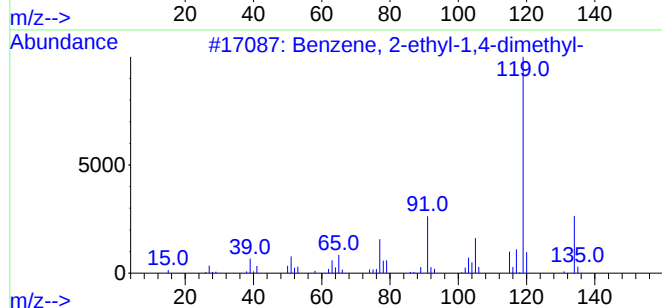
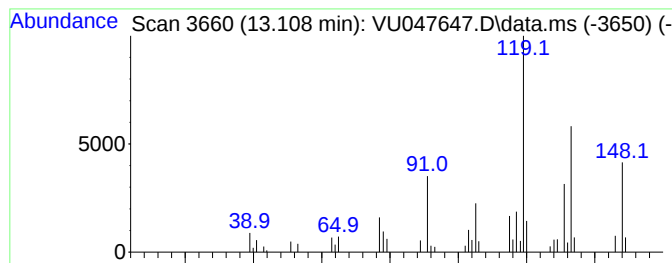
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Quant Title : VOC Analysis

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

Peak Number 20 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.108	2.54 ug/L	36923	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	55
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	50
3			4'-Methylpropiophenone	148	C10H12O	005337-93-9	46
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	45
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

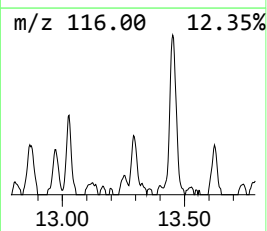
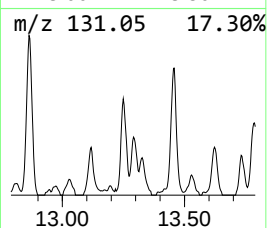
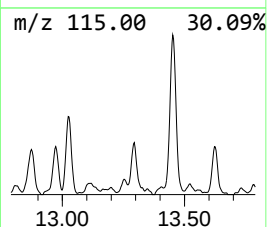
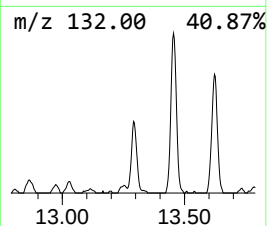
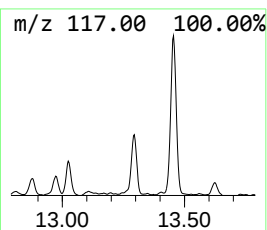
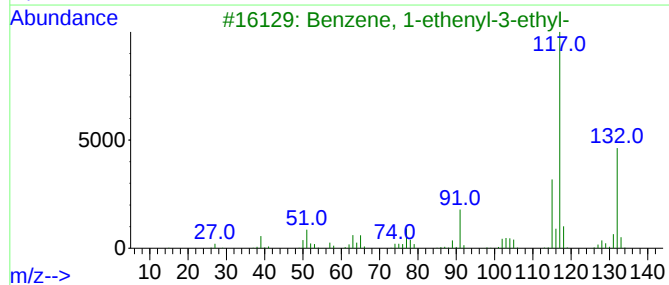
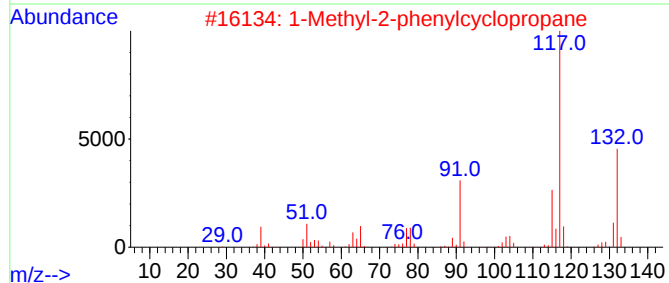
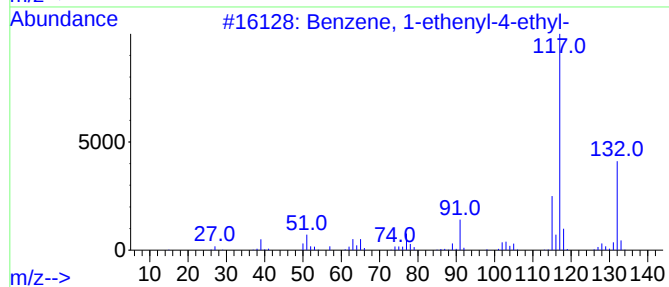
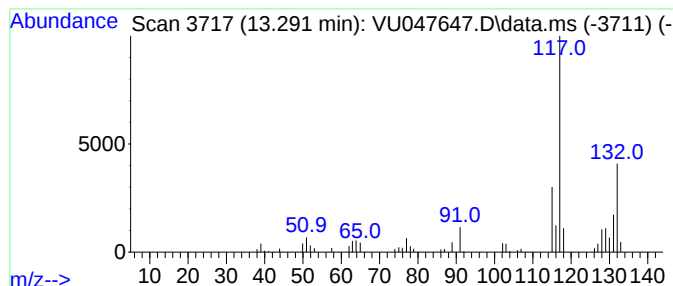
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Quant Title : VOC Analysis

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

Peak Number 21 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.291	3.99 ug/L	58069	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

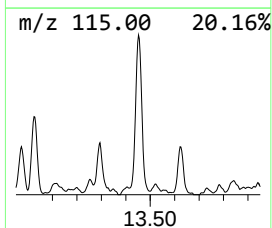
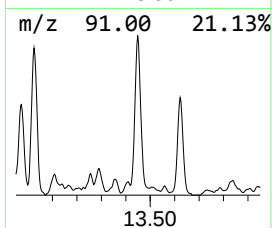
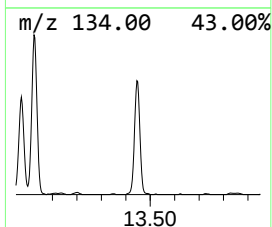
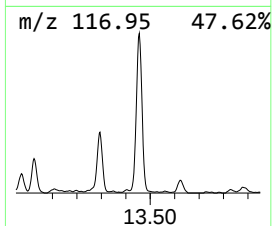
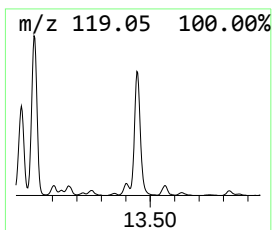
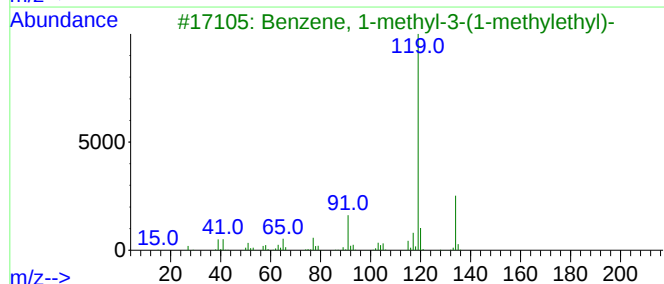
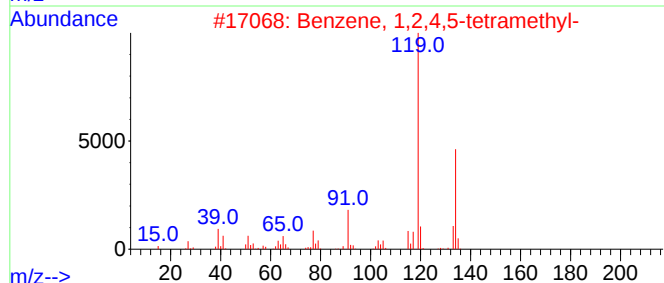
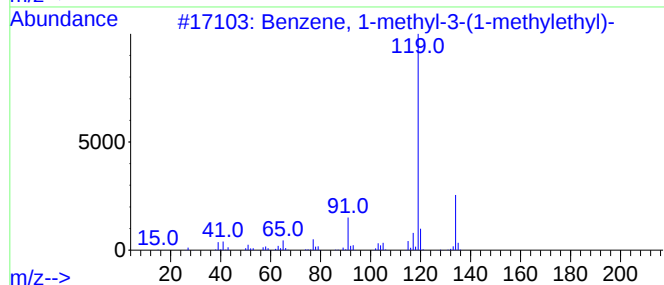
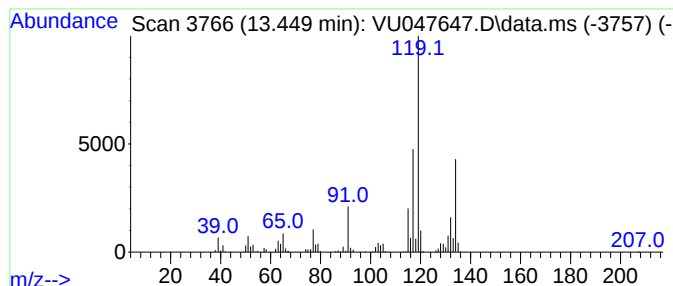
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Quant Title : VOC Analysis

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

Peak Number 22 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.449	25.90 ug/L	377199	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	64
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	64
4			p-Cymene	134	C10H14	000099-87-6	64
5			o-Cymene	134	C10H14	000527-84-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

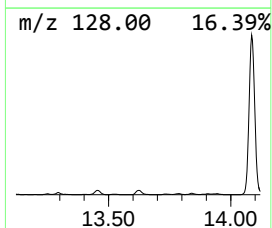
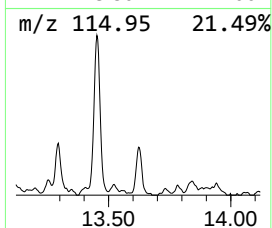
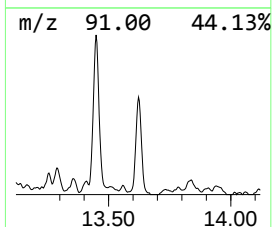
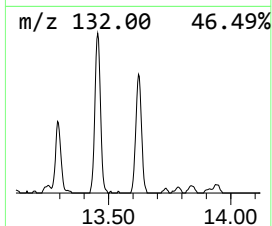
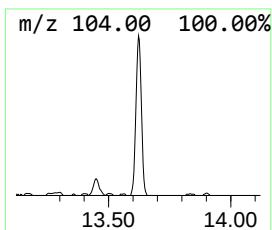
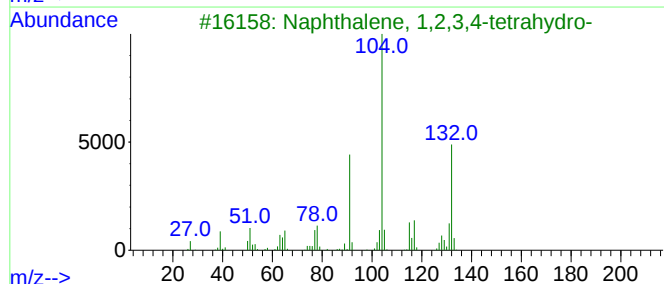
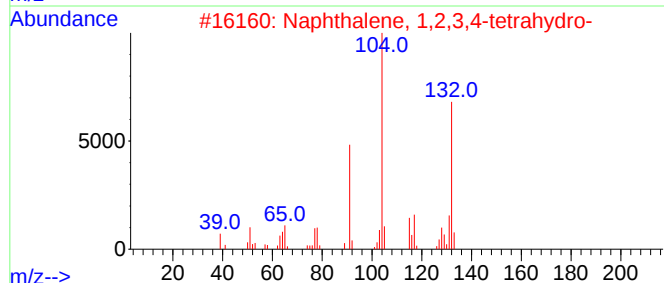
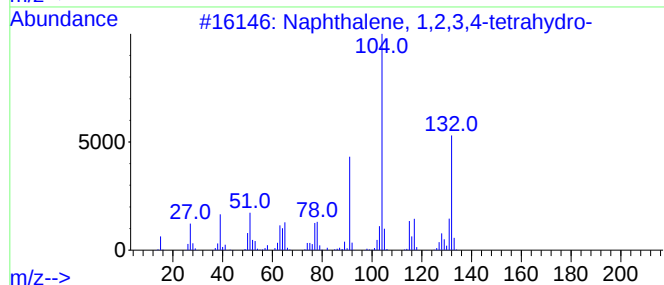
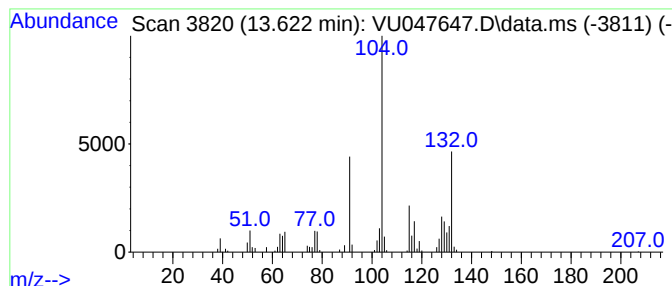
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Quant Title : VOC Analysis

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

Peak Number 23 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	8.04 ug/L	117033	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	89
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

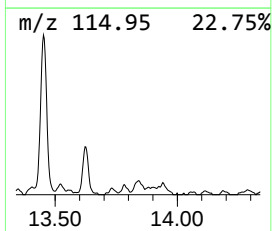
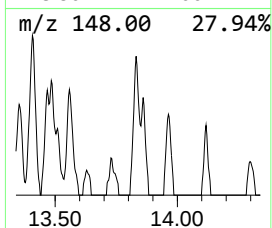
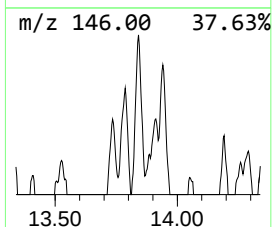
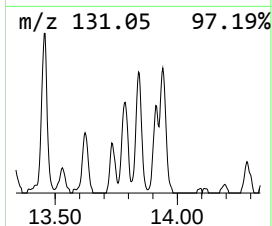
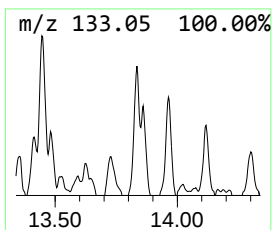
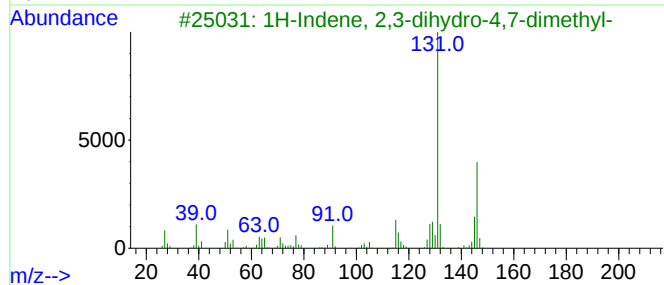
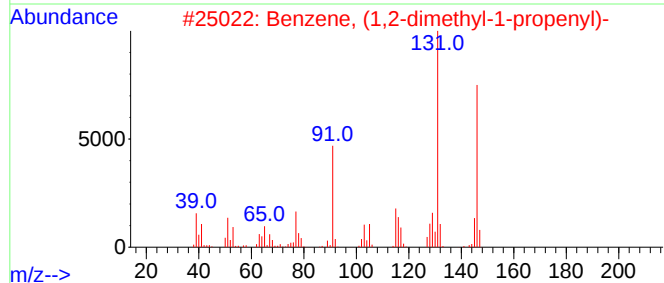
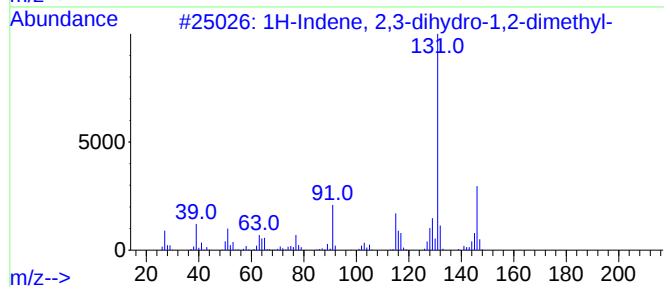
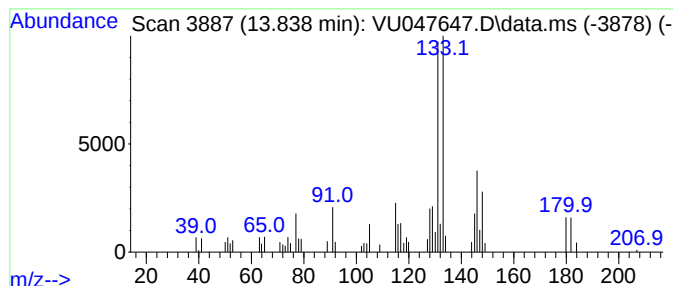
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Quant Title : VOC Analysis

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

Peak Number 24 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.838	3.33 ug/L	48559	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	50
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	46
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	46
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	45
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

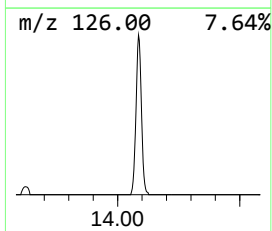
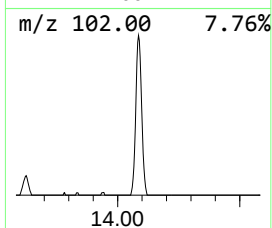
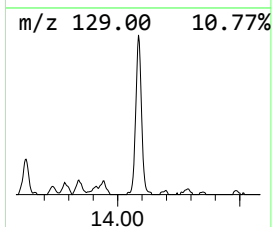
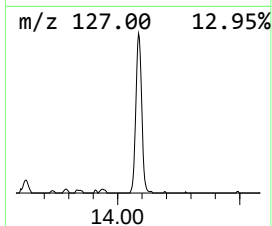
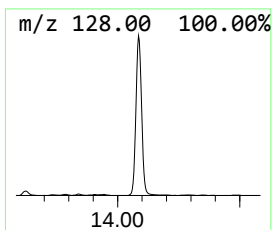
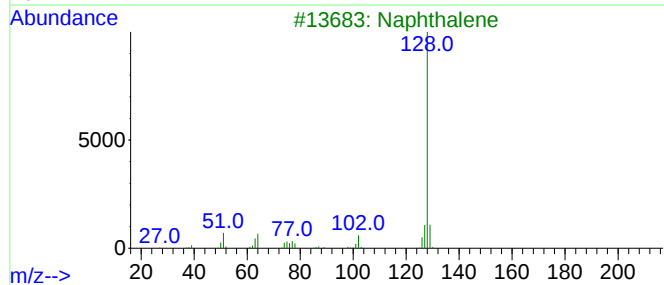
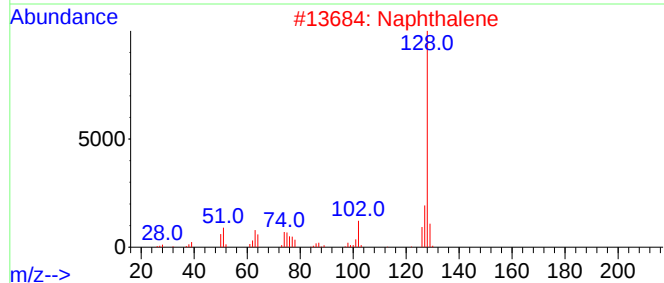
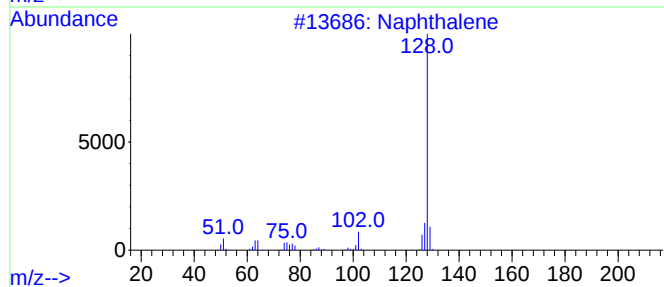
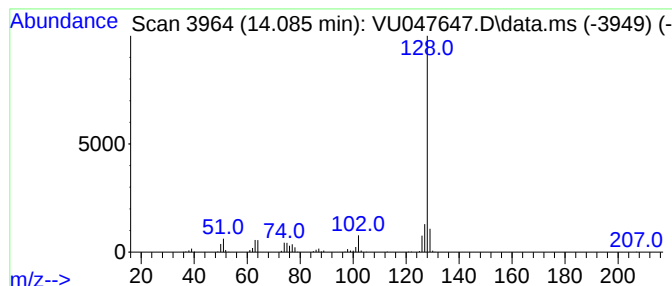
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Quant Title : VOC Analysis

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

Peak Number 25 Azulene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.085	21.40 ug/L	311594	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			Naphthalene	128	C10H8	000091-20-3	93
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
Data File : VU047647.D
Acq On : 24 Mar 2022 21:08
Operator : SY/MD
Sample : N2081-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5X5

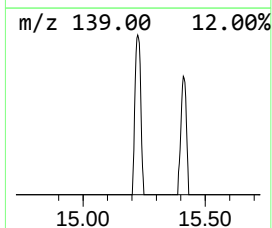
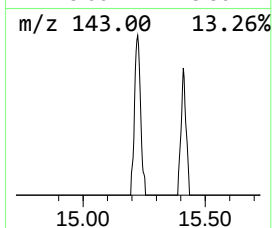
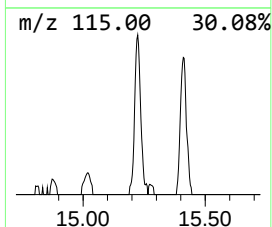
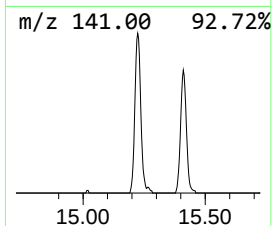
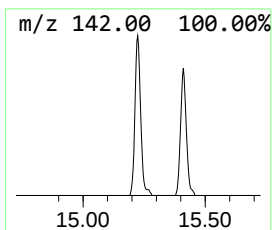
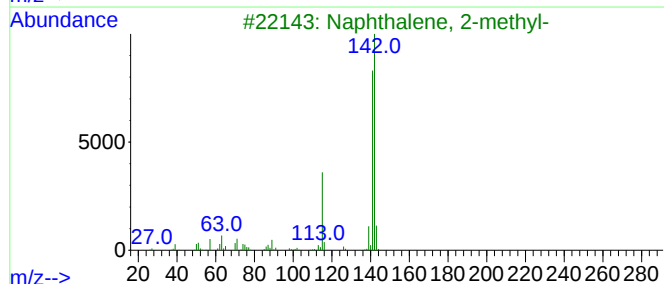
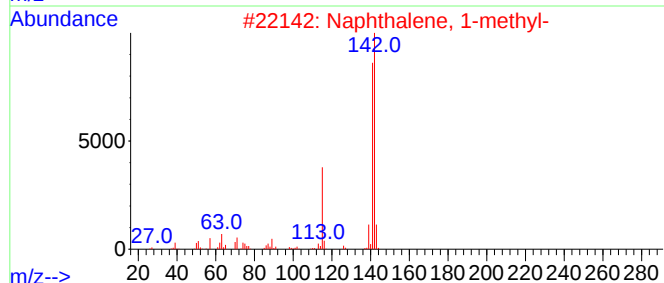
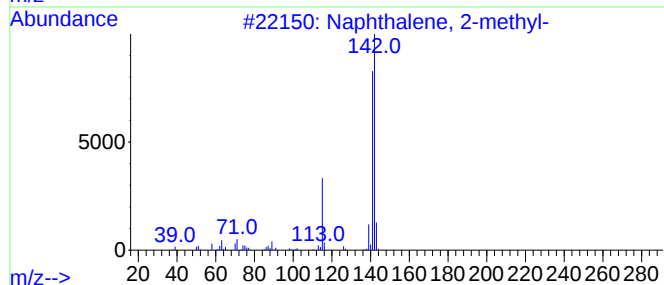
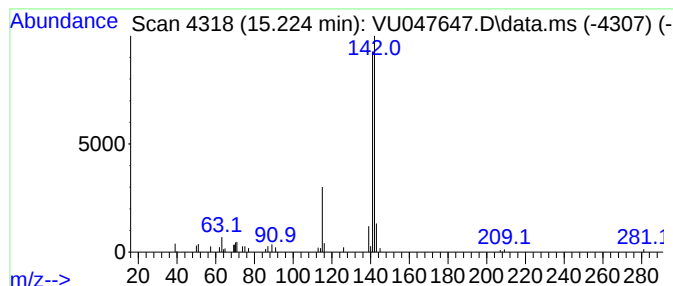
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

Peak Number 26 Naphthalene, 2-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.224	2.77 ug/L	40392	1,4-Dichlorobenzene-d4	11.812

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032422\
 Data File : VU047647.D
 Acq On : 24 Mar 2022 21:08
 Operator : SY/MD
 Sample : N2081-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EW5X5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.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
(DEL) Alkane: C...	8.864	3.6	ug/L	37307	2	9.417	520612	50.0
unknown-01	10.272	4.0	ug/L	41364	2	9.417	520612	50.0
Benzene, propyl-	10.906	35.5	ug/L	517189	3	11.812	728113	50.0
Benzene, 1-ethy...	10.999	36.8	ug/L	536034	3	11.812	728113	50.0
Benzene, 1-ethy...	11.291	89.4	ug/L	1301780	3	11.812	728113	50.0
Benzene, (2-met...	11.610	4.5	ug/L	65067	3	11.812	728113	50.0
Benzene, 1-meth...	11.642	8.2	ug/L	119175	3	11.812	728113	50.0
p-Cymene	11.741	12.6	ug/L	182979	3	11.812	728113	50.0
Benzene, 1,2,3-...	11.893	109.9	ug/L	1600220	3	11.812	728113	50.0
Benzene, 1-ethe...	12.095	47.3	ug/L	688310	3	11.812	728113	50.0
Benzene, 1,2-di...	12.301	4.3	ug/L	62484	3	11.812	728113	50.0
Benzene, 1-meth...	12.375	13.7	ug/L	199896	3	11.812	728113	50.0
Benzene, 2-ethy...	12.465	18.6	ug/L	271069	3	11.812	728113	50.0
Benzene, 1-meth...	12.571	25.9	ug/L	377173	3	11.812	728113	50.0
Benzene, (1-met...	12.684	20.8	ug/L	303271	3	11.812	728113	50.0
Benzene, 4-ethy...	12.873	11.3	ug/L	164360	3	11.812	728113	50.0
Benzene, 1,2,4,...	12.973	14.2	ug/L	206221	3	11.812	728113	50.0
Benzene, 1,2,3,...	13.024	21.8	ug/L	317420	3	11.812	728113	50.0
Benzene, 1-ethy...	13.108	2.5	ug/L	36923	3	11.812	728113	50.0
Benzene, 1-ethe...	13.291	4.0	ug/L	58069	3	11.812	728113	50.0
o-Cymene	13.449	25.9	ug/L	377199	3	11.812	728113	50.0
Naphthalene, 1,...	13.622	8.0	ug/L	117033	3	11.812	728113	50.0
1H-Indene, 2,3-...	13.838	3.3	ug/L	48559	3	11.812	728113	50.0
Azulene	14.085	21.4	ug/L	311594	3	11.812	728113	50.0
Naphthalene, 2-...	15.224	2.8	ug/L	40392	3	11.812	728113	50.0