

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW061523\
 Data File : VW031534.D
 Acq On : 15 Jun 2023 16:23
 Operator : SY/MD
 Sample : 03249-12
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EXYT2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.M
 Title : VOC Analysis

Signal : TIC: VW031534.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.179	38	43	59	rBV	8819	12752	0.02%	0.009%
2	1.285	69	76	95	rBV2	311235	396163	0.59%	0.281%
3	1.552	148	159	183	rVB2	201726	335934	0.50%	0.238%
4	1.674	190	197	206	rBV	28045	36630	0.05%	0.026%
5	1.722	207	212	216	rBV	8871	10263	0.02%	0.007%
6	1.806	232	238	250	rVB	18638	24660	0.04%	0.017%
7	2.076	309	322	345	rBV3	552856	964616	1.45%	0.684%
8	2.262	373	380	386	rVV5	1964	2496	0.00%	0.002%
9	2.455	431	440	454	rBV2	13203	24848	0.04%	0.018%
10	2.703	505	517	549	rBV	237641	466819	0.70%	0.331%
11	2.979	592	603	617	rVB8	4765	10026	0.02%	0.007%
12	3.121	627	647	694	rBV	8053269	17201580	25.77%	12.189%
13	3.680	809	821	832	rBV6	3927	9412	0.01%	0.007%
14	3.735	832	838	842	rBV5	1354	1648	0.00%	0.001%
15	3.825	846	866	922	rBV	26602652	66748285	100.00%	47.298%
16	4.262	988	1002	1029	rVB2	139019	370402	0.55%	0.262%
17	4.526	1061	1084	1114	rBV	8819592	22240518	33.32%	15.760%
18	4.966	1204	1221	1234	rBV3	310625	837971	1.26%	0.594%
19	5.243	1300	1307	1318	rVB7	2547	5204	0.01%	0.004%
20	5.420	1353	1362	1371	rBV8	2650	5458	0.01%	0.004%
21	5.542	1386	1400	1426	rBV	241453	496765	0.74%	0.352%
22	5.841	1479	1493	1523	rBV	283692	570582	0.85%	0.404%
23	5.995	1529	1541	1555	rBV	191287	392037	0.59%	0.278%
24	6.297	1626	1635	1645	rBV10	2113	4127	0.01%	0.003%
25	6.558	1708	1716	1719	rBV4	1388	1814	0.00%	0.001%
26	6.850	1802	1807	1814	rVB4	1146	1675	0.00%	0.001%
27	6.921	1815	1829	1850	rBV	99715	188560	0.28%	0.134%
28	7.172	1896	1907	1921	rBV4	24290	58914	0.09%	0.042%
29	7.249	1921	1931	1943	rVV	339771	587784	0.88%	0.417%
30	7.320	1943	1953	1975	rVB	992542	1668746	2.50%	1.182%
31	7.558	2018	2027	2051	rVB	81669	156668	0.23%	0.111%
32	7.722	2071	2078	2087	rVB6	4250	6893	0.01%	0.005%
33	7.780	2087	2096	2110	rBV2	24836	43514	0.07%	0.031%
34	7.911	2121	2137	2156	rBV2	157462	273404	0.41%	0.194%
35	8.030	2164	2174	2195	rBV	223190	394568	0.59%	0.280%
36	8.230	2227	2236	2247	rVB3	15490	25329	0.04%	0.018%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Stop Thrs : 0

Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.M
 Title : VOC Analysis

37	8.288	2247	2254	2263	rVB4	6616	10623	0.02%	0.008%
38	8.529	2324	2329	2338	rVB4	1797	3101	0.00%	0.002%
39	8.612	2346	2355	2357	rBV5	1170	1219	0.00%	0.001%
40	8.657	2360	2369	2384	rVB5	4058	7419	0.01%	0.005%
41	8.789	2398	2410	2428	rBV	378217	620874	0.93%	0.440%
42	8.953	2449	2461	2476	rBV	529255	818669	1.23%	0.580%
43	9.079	2487	2500	2515	rBV	1310706	2087849	3.13%	1.479%
44	9.262	2547	2557	2565	rVB7	3742	6534	0.01%	0.005%
45	9.333	2572	2579	2591	rBV4	2984	5782	0.01%	0.004%
46	9.416	2592	2605	2614	rBV8	6628	14960	0.02%	0.011%
47	9.484	2614	2626	2643	rBV	1190218	1768149	2.65%	1.253%
48	9.644	2669	2676	2690	rVB3	19942	33264	0.05%	0.024%
49	9.741	2696	2706	2719	rBV8	14083	25933	0.04%	0.018%
50	9.873	2736	2747	2767	rBV	84686	135620	0.20%	0.096%
51	9.960	2767	2774	2788	rVB7	6896	10482	0.02%	0.007%
52	10.034	2788	2797	2801	rBV5	4968	7576	0.01%	0.005%
53	10.066	2803	2807	2819	rVB9	7913	12759	0.02%	0.009%
54	10.159	2824	2836	2864	rVB	307192	500593	0.75%	0.355%
55	10.297	2867	2879	2897	rBV	374043	591231	0.89%	0.419%
56	10.390	2897	2908	2926	rVV	990305	2071802	3.10%	1.468%
57	10.484	2926	2937	2957	rVB	896862	1328178	1.99%	0.941%
58	10.680	2988	2998	3016	rVB	938848	1408214	2.11%	0.998%
59	10.808	3032	3038	3041	rBV2	5054	5687	0.01%	0.004%
60	10.857	3041	3053	3071	rVV	3998008	5824464	8.73%	4.127%
61	10.998	3091	3097	3101	rBV	31442	43303	0.06%	0.031%
62	11.030	3103	3107	3118	rVB	71603	98826	0.15%	0.070%
63	11.140	3130	3141	3148	rBV	101687	154974	0.23%	0.110%
64	11.191	3148	3157	3171	rVB2	548568	828670	1.24%	0.587%
65	11.278	3172	3184	3200	rBV	1403991	2079317	3.12%	1.473%
66	11.397	3213	3221	3231	rBV2	18458	28539	0.04%	0.020%
67	11.471	3234	3244	3254	rBV2	425849	807997	1.21%	0.573%
68	11.570	3266	3275	3301	rVB6	626167	1430558	2.14%	1.014%
69	11.689	3302	3312	3325	rBV2	39328	59563	0.09%	0.042%
70	11.763	3325	3335	3350	rVB	116030	182061	0.27%	0.129%
71	11.863	3353	3366	3382	rBV2	565992	1229844	1.84%	0.871%
72	11.960	3388	3396	3417	rVB2	258786	472002	0.71%	0.334%
73	12.059	3417	3427	3433	rBV	88577	151924	0.23%	0.108%
74	12.204	3459	3472	3478	rBV6	18918	35188	0.05%	0.025%
75	12.259	3480	3489	3501	rVB3	79398	134355	0.20%	0.095%
76	12.358	3502	3520	3528	rBV	108878	196934	0.30%	0.140%

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 Title : VOC Analysis

77	12.410	3528	3536	3553	rVB	188685	292040	0.44%	0.207%
78	12.496	3555	3563	3569	rBV3	19060	28292	0.04%	0.020%
79	12.561	3580	3583	3588	rBV3	4749	4397	0.01%	0.003%
80	12.670	3610	3617	3631	rVB2	48953	75860	0.11%	0.054%
81	12.741	3631	3639	3646	rBV7	9061	12645	0.02%	0.009%
82	12.824	3647	3665	3678	rBV2	212658	356380	0.53%	0.253%
83	12.943	3696	3702	3708	rBV2	9755	12955	0.02%	0.009%
84	12.992	3709	3717	3729	rVB2	52664	83972	0.13%	0.060%
85	13.107	3745	3753	3762	rBV7	13569	21640	0.03%	0.015%
86	13.162	3762	3770	3777	rBV3	9342	13526	0.02%	0.010%
87	13.213	3777	3786	3793	rBV4	23962	38733	0.06%	0.027%
88	13.445	3846	3858	3869	rBV	154019	242335	0.36%	0.172%
89	13.561	3886	3894	3902	rBV8	4069	6496	0.01%	0.005%
90	13.664	3918	3926	3937	rVV8	3887	10318	0.02%	0.007%
91	13.712	3937	3941	3950	rVB5	2857	3870	0.01%	0.003%
92	13.850	3978	3984	4001	rVB5	4338	8891	0.01%	0.006%
93	13.985	4021	4026	4032	rBV5	1443	1686	0.00%	0.001%
94	14.049	4034	4046	4055	rVB10	7111	15606	0.02%	0.011%
95	14.178	4078	4086	4091	rBV5	2827	3843	0.01%	0.003%
96	14.239	4098	4105	4111	rBV7	1857	2586	0.00%	0.002%
97	14.377	4139	4148	4155	rBV5	4478	6769	0.01%	0.005%
98	14.577	4198	4210	4220	rBV2	25106	40111	0.06%	0.028%
99	14.760	4259	4267	4283	rVB	15189	25525	0.04%	0.018%
100	15.609	4524	4531	4537	rBV7	1749	3070	0.00%	0.002%

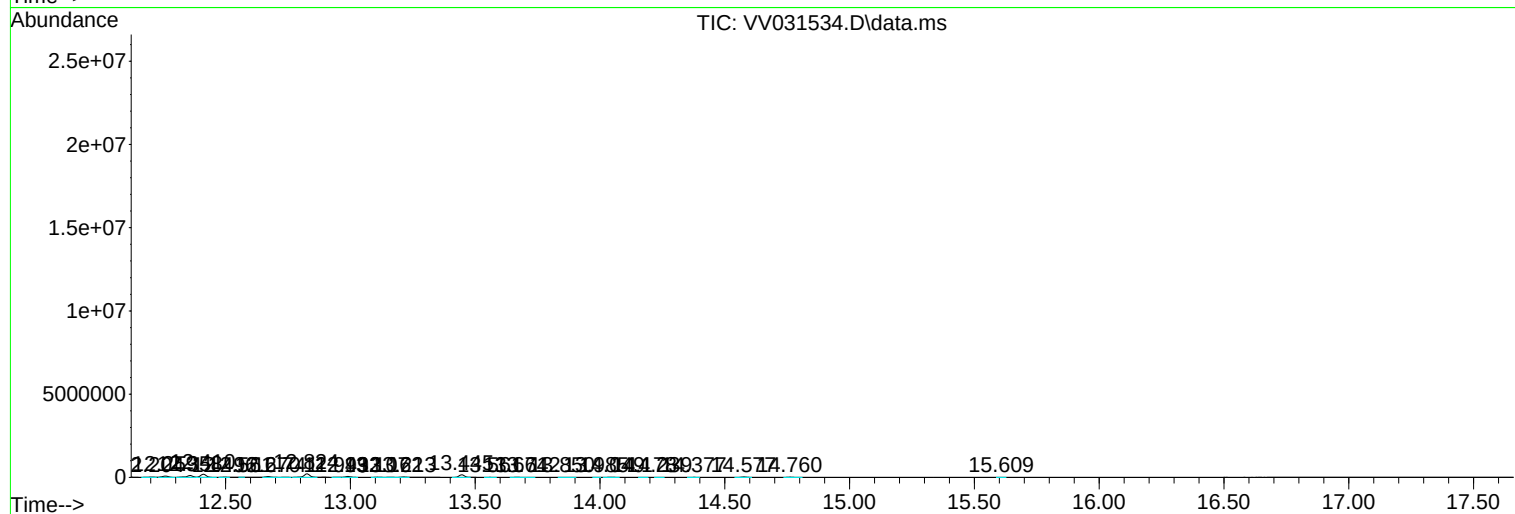
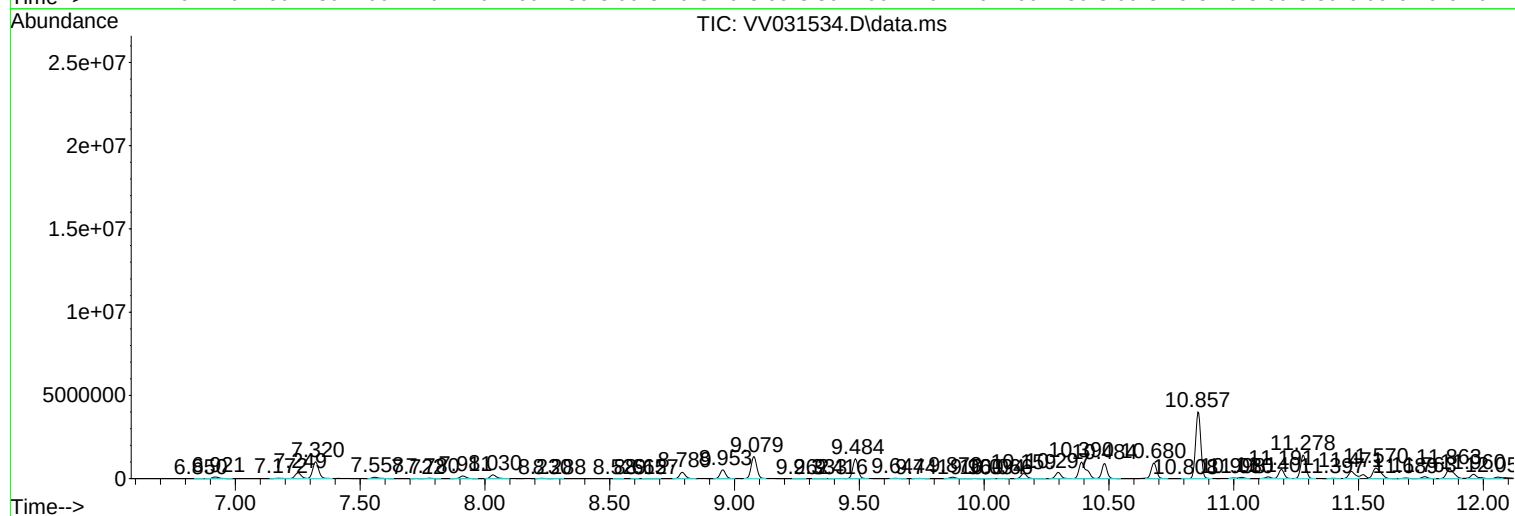
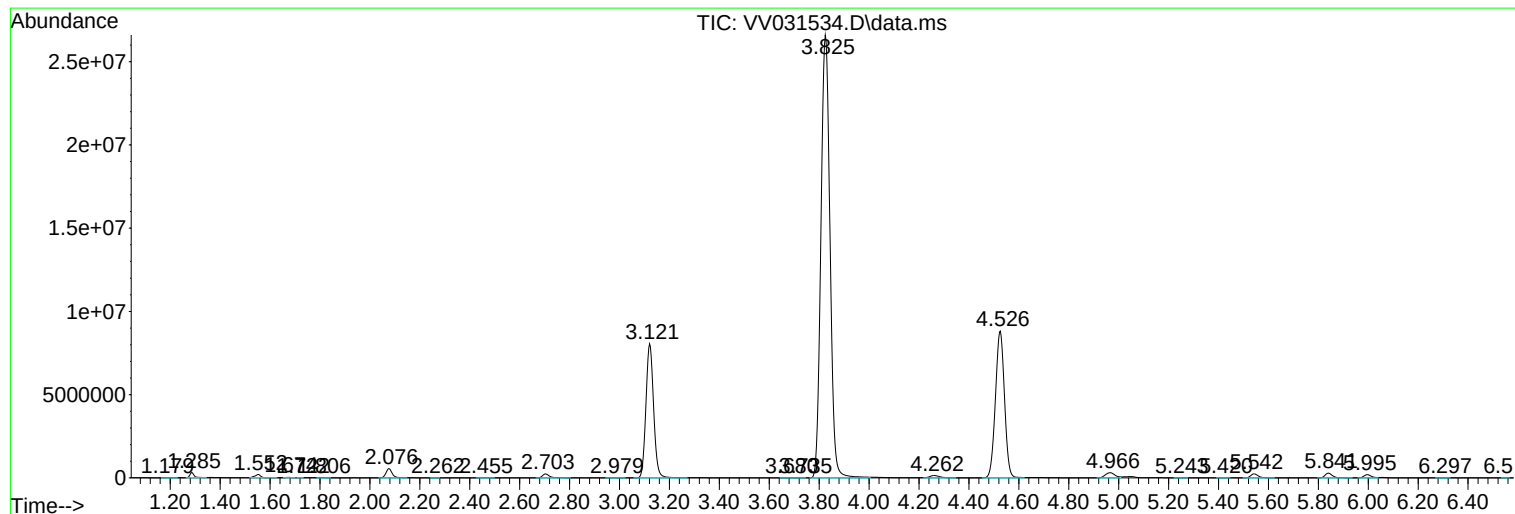
Sum of corrected areas: 141123648

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ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
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Instrument :
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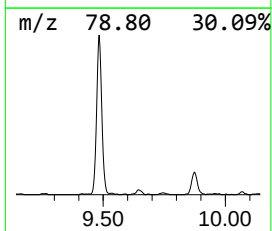
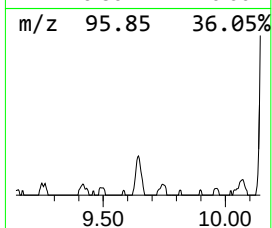
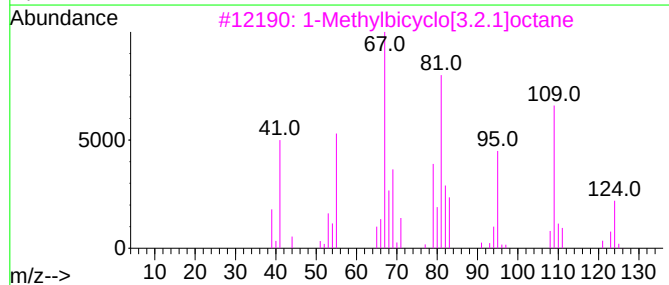
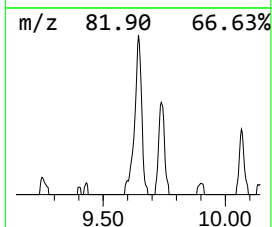
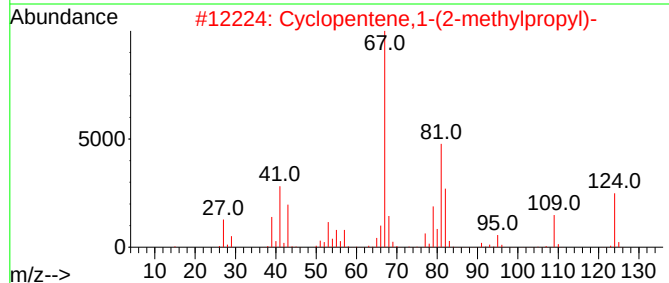
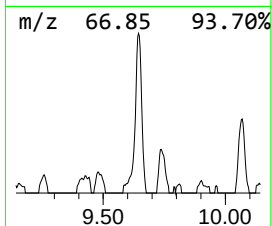
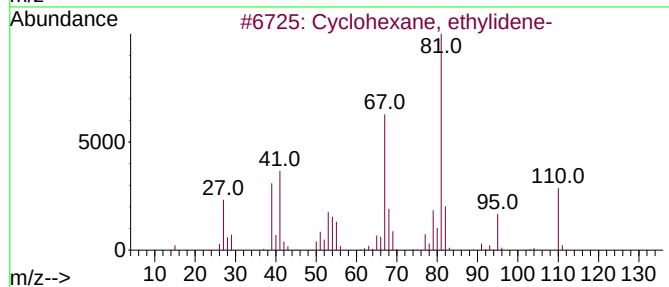
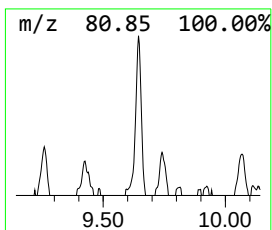
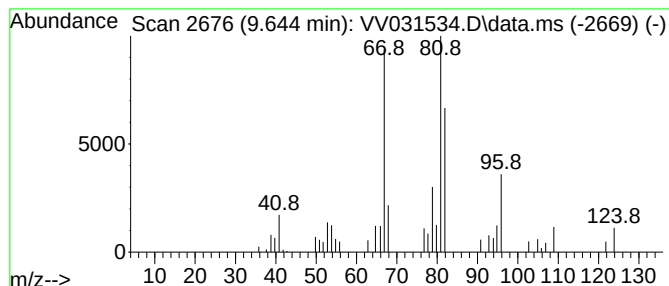
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 unknown-01 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.644	2.68 ug/L	33264	Chlorobenzene-d5	8.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethylidene-	110	C8H14	001003-64-1	47
2			Cyclopentene,1-(2-methylpropyl)-	124	C9H16	053098-47-8	46
3			1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	43
4			1H-Indene, octahydro-	124	C9H16	000496-10-6	43
5			1,3-Pentadiene, 2-methyl-, (E)-	82	C6H10	000926-54-5	38



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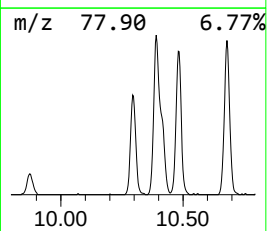
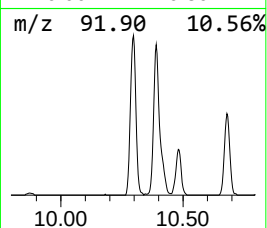
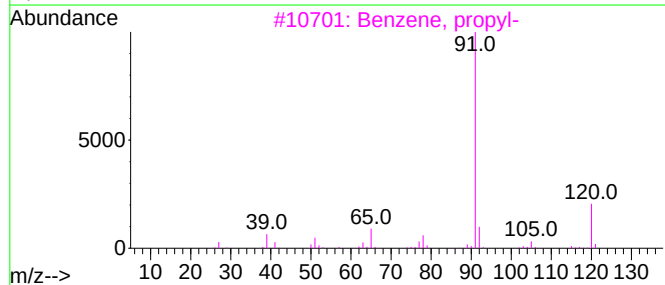
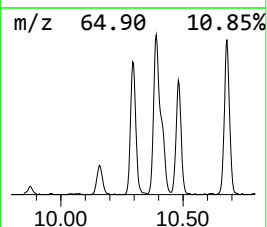
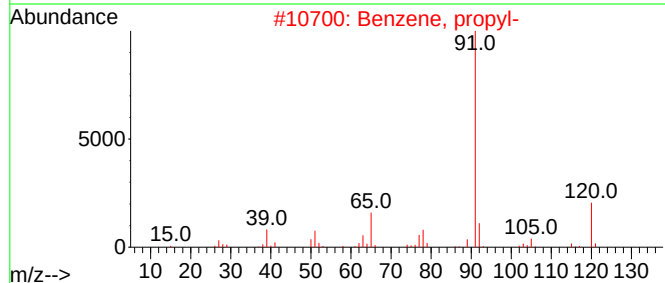
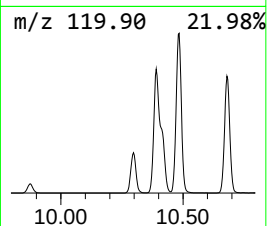
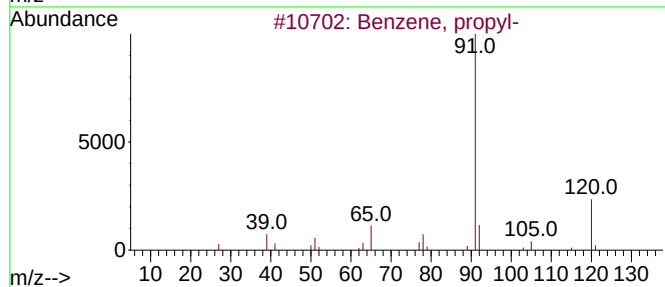
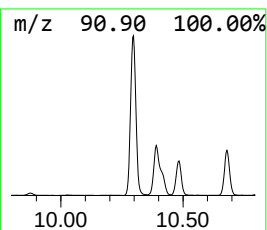
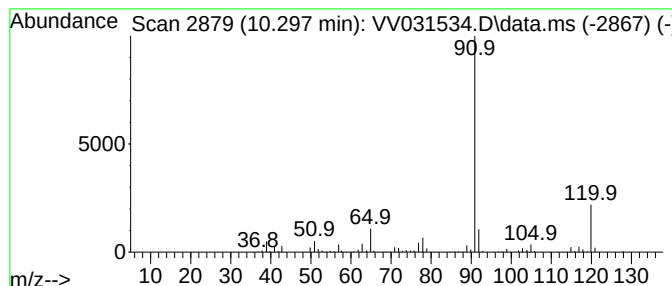
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.297	35.67 ug/L	591231	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	83
5			Benzene, propyl-	120	C9H12	000103-65-1	83



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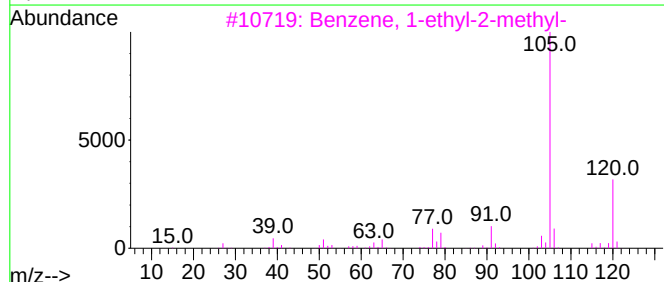
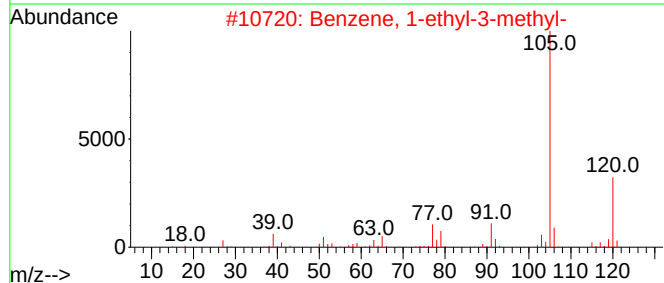
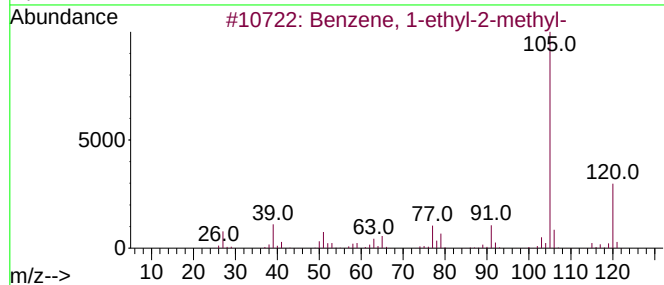
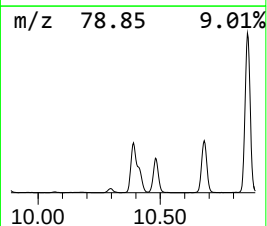
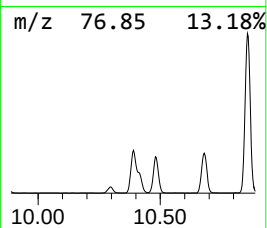
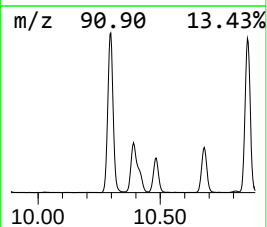
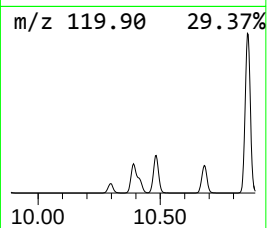
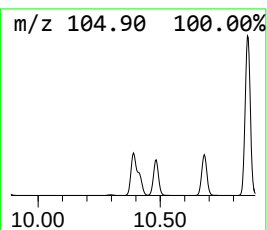
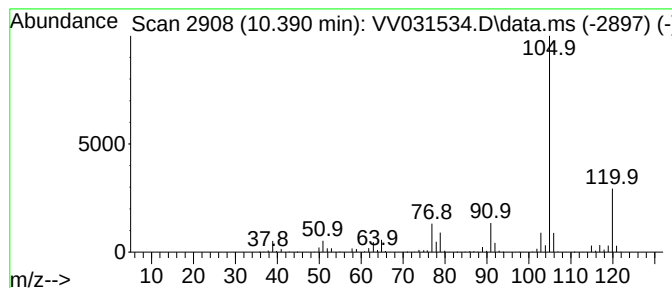
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Quant Title : VOC Analysis

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.390	125.01 ug/L	2071800	1,4-Dichlorobenzene-d4	11.191

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-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

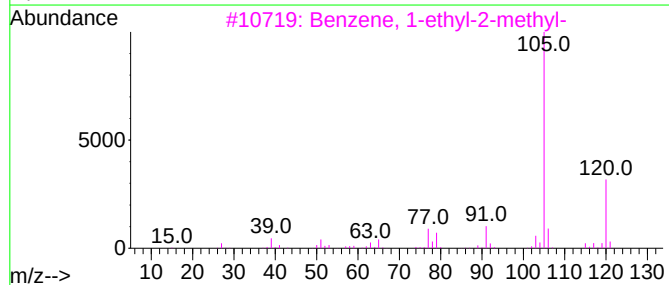
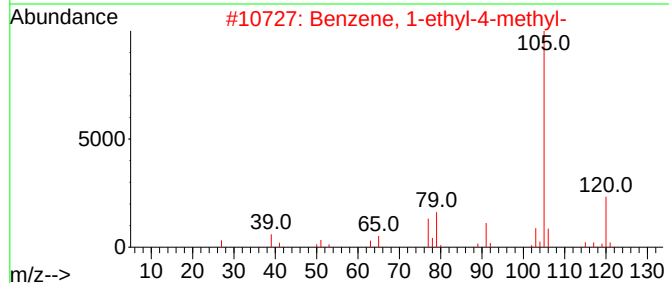
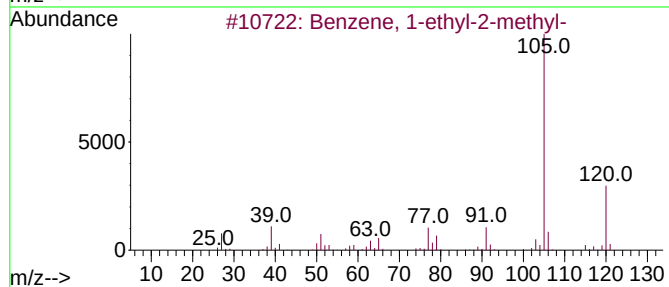
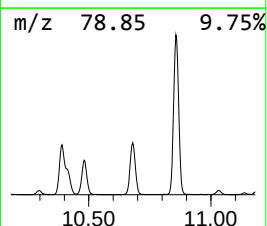
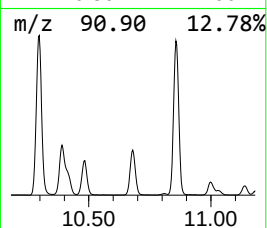
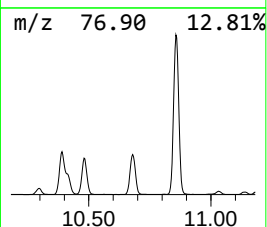
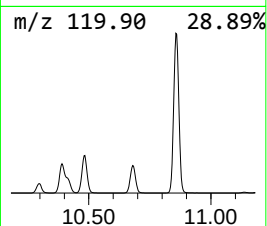
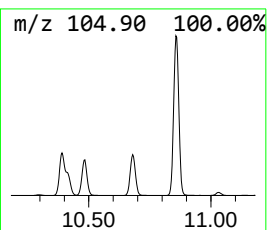
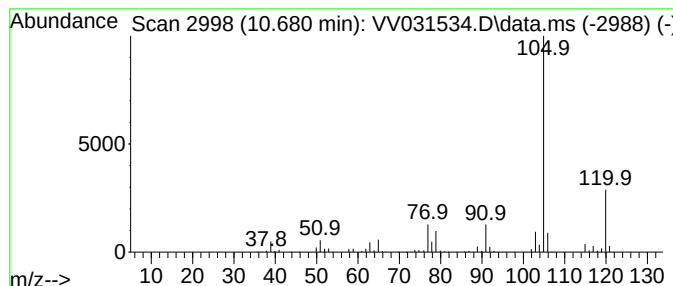
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Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.680	84.97 ug/L	1408210	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

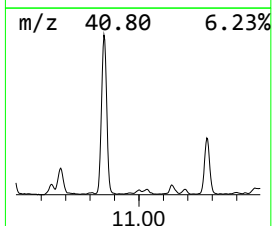
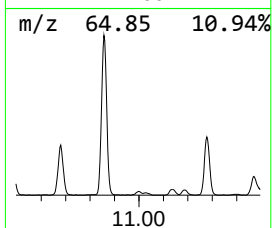
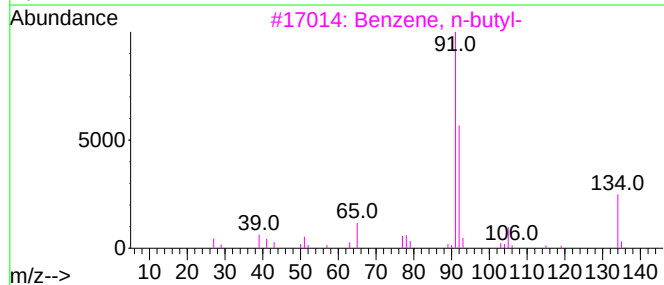
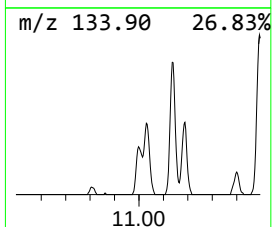
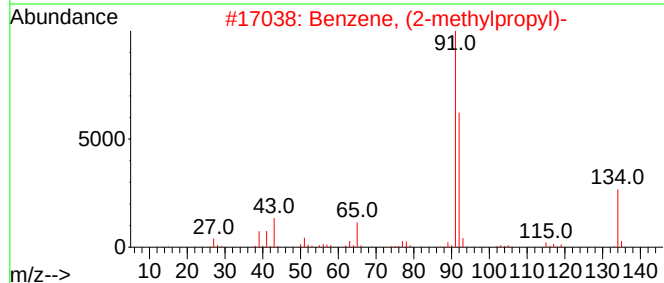
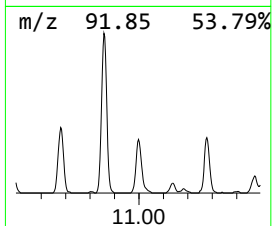
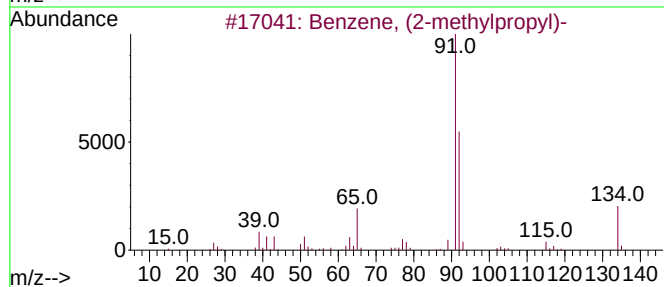
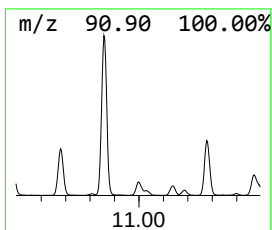
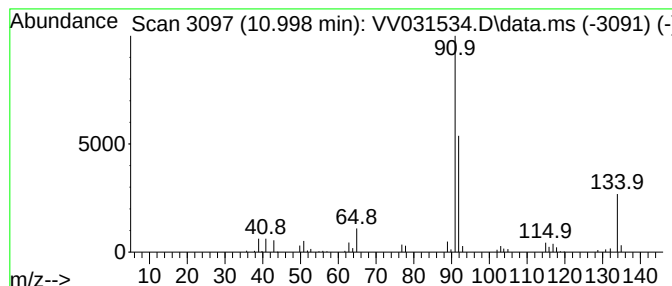
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, (2-methylpropyl)- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.998	2.61 ug/L	43303	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87
2			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	80
3			Benzene, n-butyl-	134	C10H14	000104-51-8	80
4			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	80
5			Benzene, n-butyl-	134	C10H14	000104-51-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

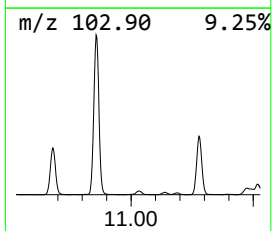
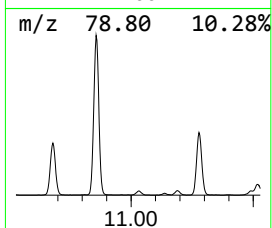
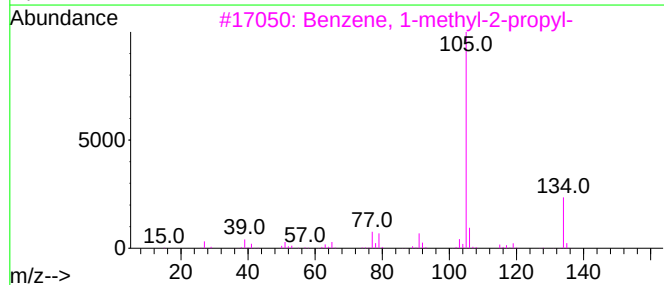
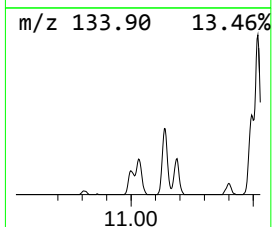
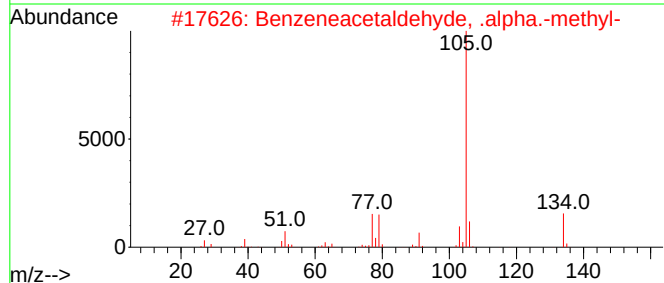
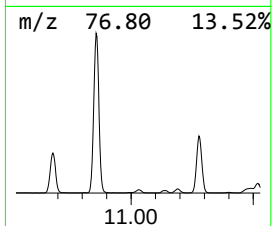
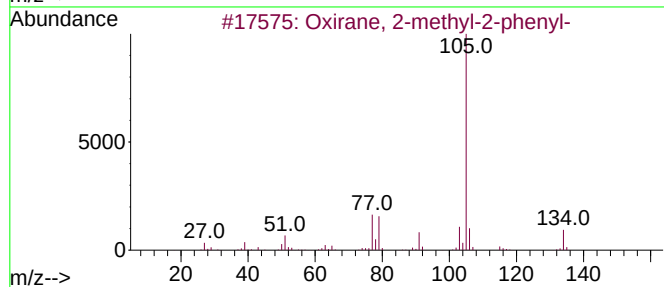
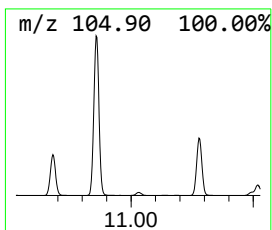
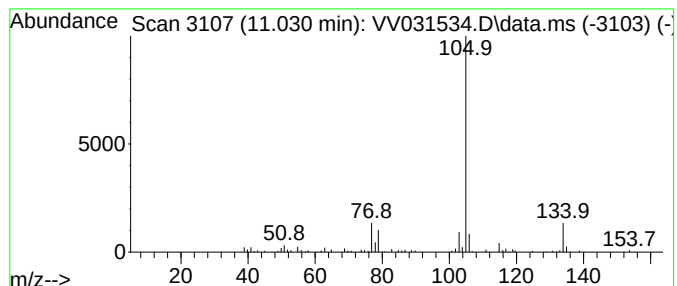
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Quant Title : VOC Analysis

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

Peak Number 7 Oxirane, 2-methyl-2-phenyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.030	5.96 ug/L	98826	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	87
2			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	87
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	86
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	83
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

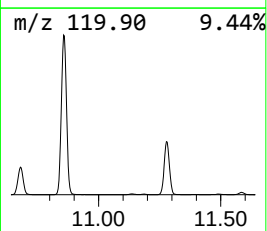
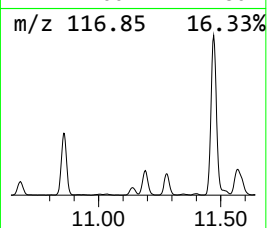
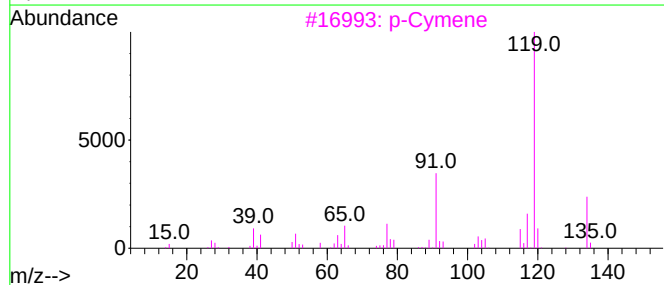
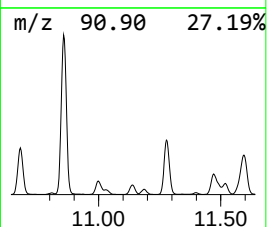
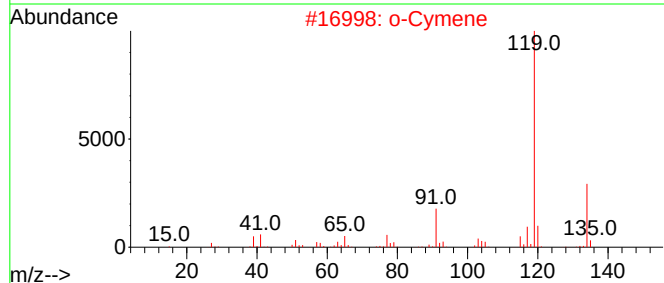
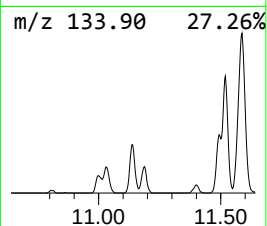
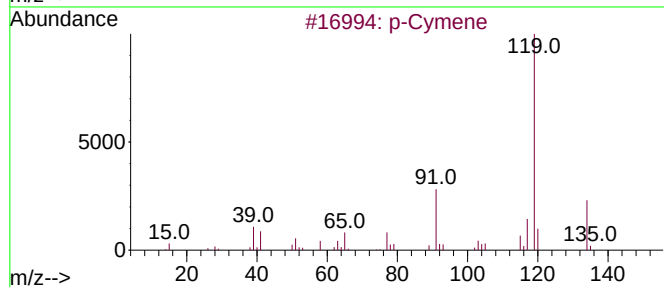
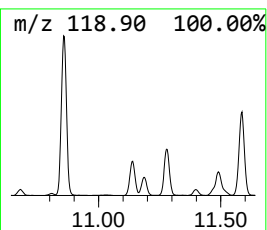
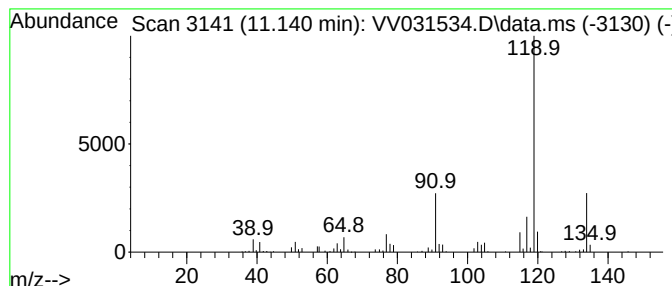
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Quant Title : VOC Analysis

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

Peak Number 8 p-Cymene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.140	9.35 ug/L	154974	1,4-Dichlorobenzene-d4	11.191

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	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
5			1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

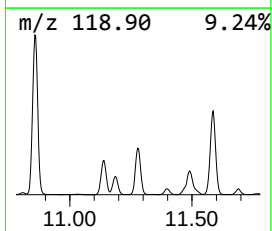
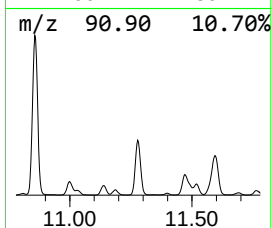
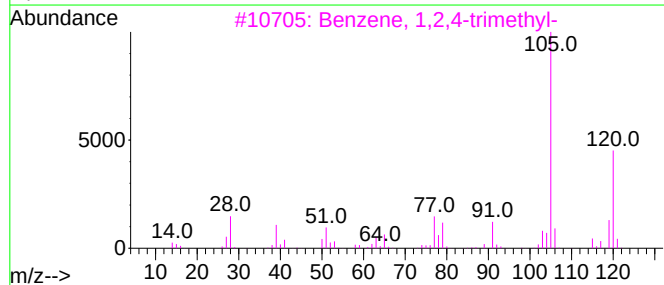
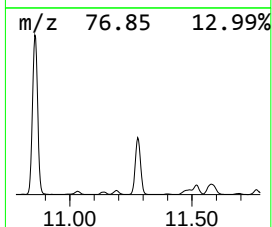
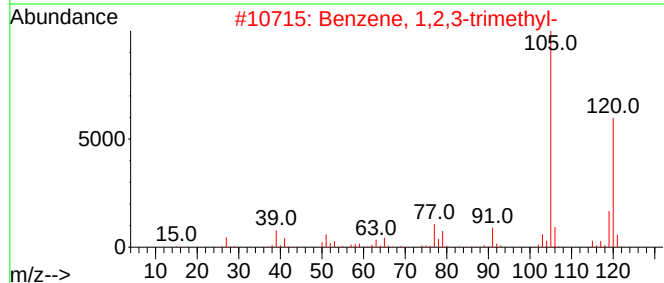
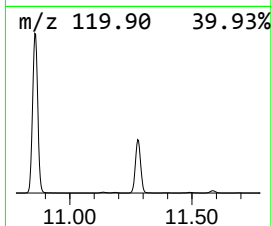
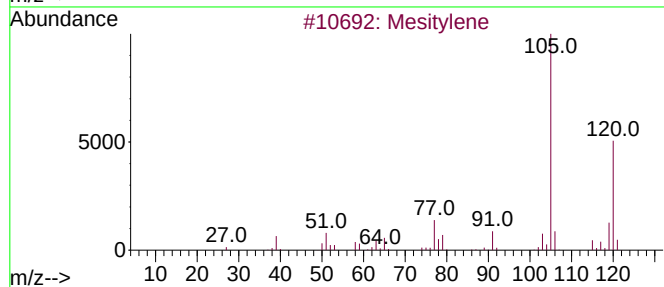
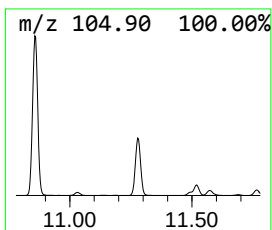
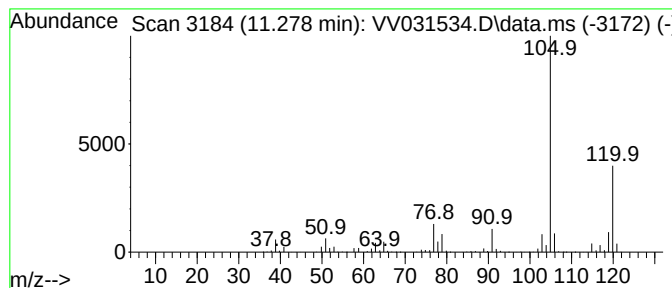
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.278	125.46 ug/L	2079320	1,4-Dichlorobenzene-d4	11.191

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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

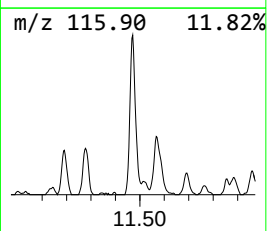
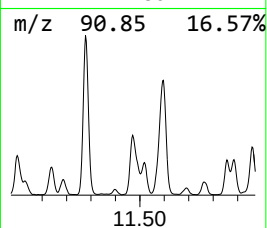
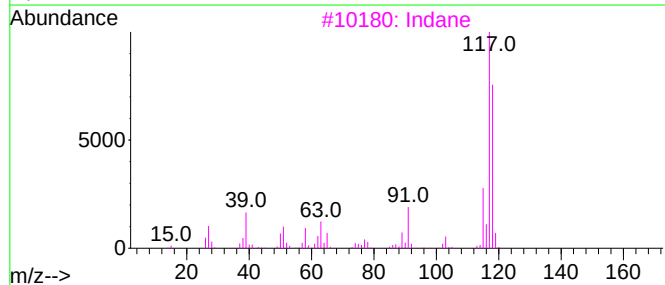
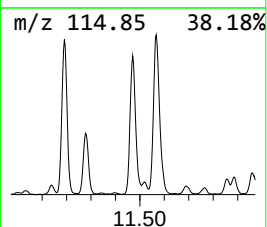
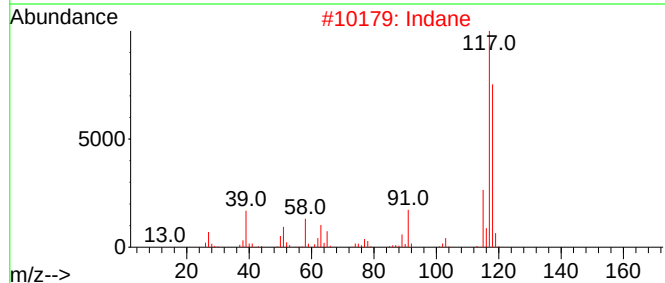
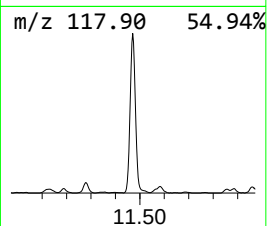
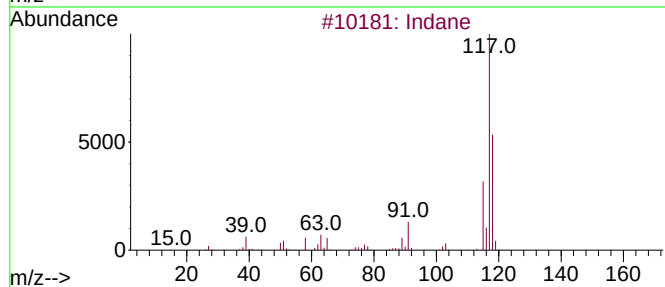
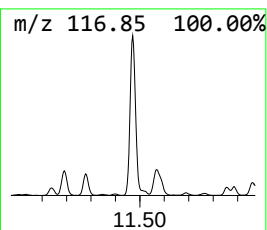
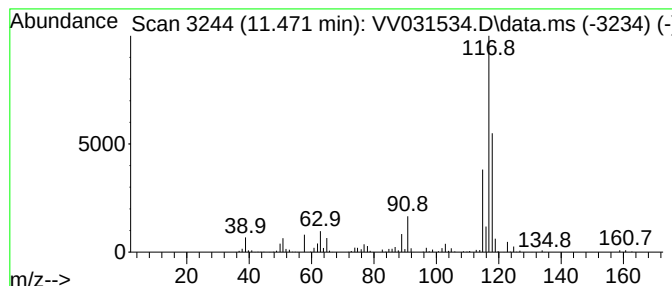
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Quant Title : VOC Analysis

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

Peak Number 10 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.471	48.75 ug/L	807997	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	95
2	Indane			118	C9H10	000496-11-7	89
3	Indane			118	C9H10	000496-11-7	87
4	Indane			118	C9H10	000496-11-7	70
5	Benzene, 1-ethenyl-4-methyl-			118	C9H10	000622-97-9	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

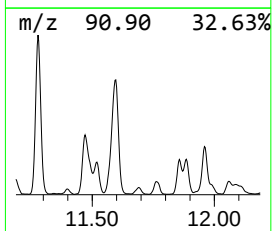
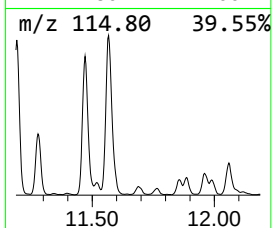
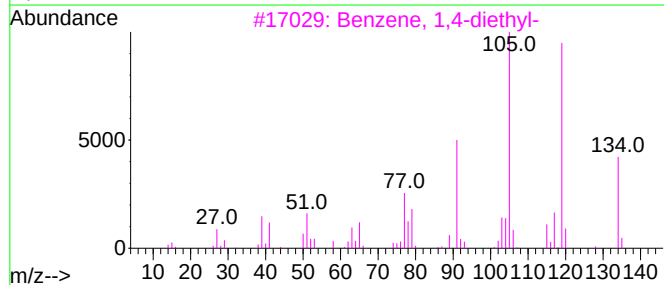
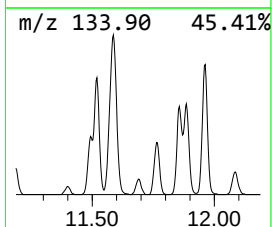
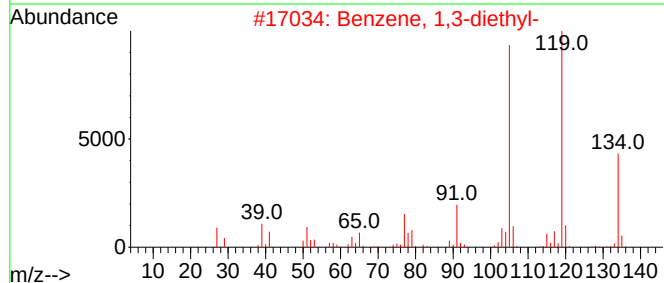
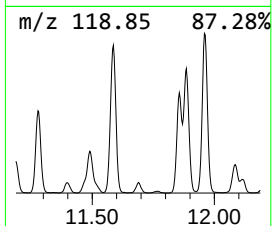
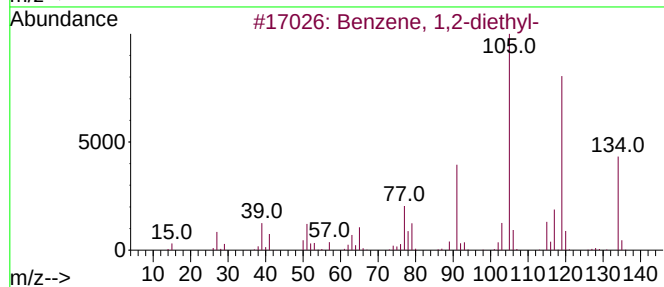
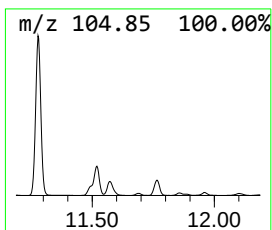
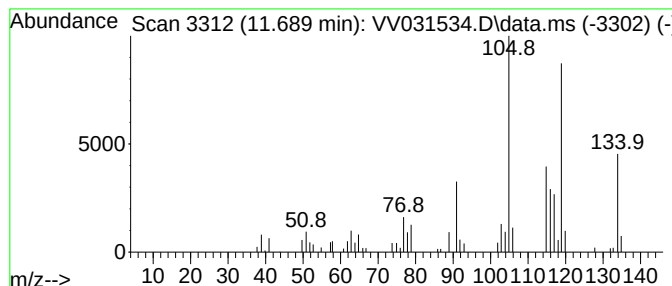
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Quant Title : VOC Analysis

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

Peak Number 11 Benzene, 1,2-diethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.689	3.59 ug/L	59563	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

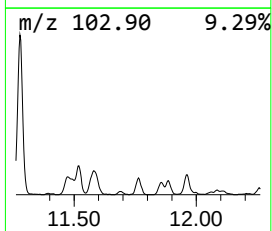
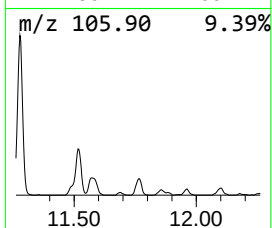
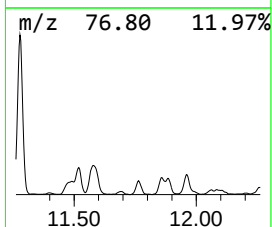
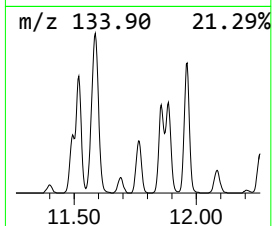
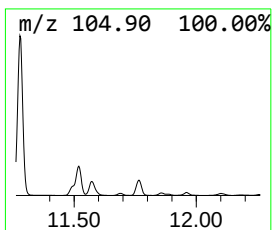
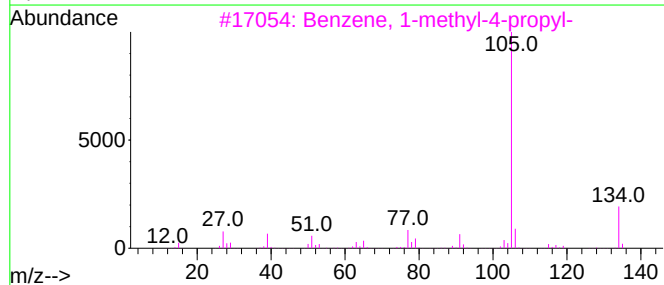
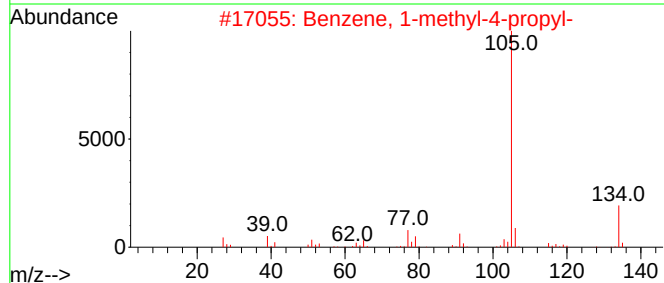
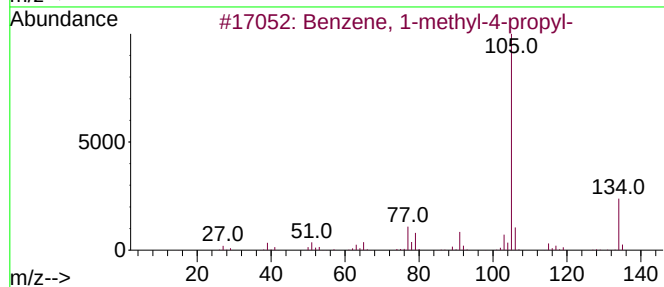
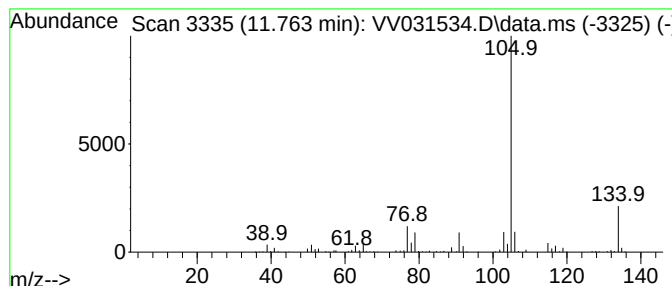
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, 1-methyl-4-propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.763	10.99 ug/L	182061	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	91
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

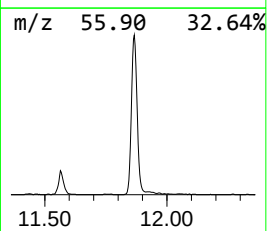
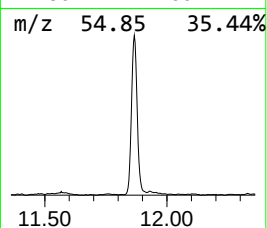
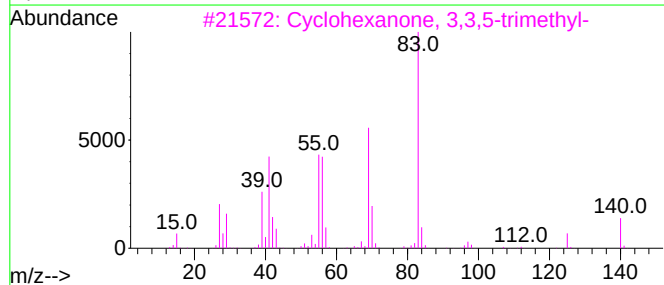
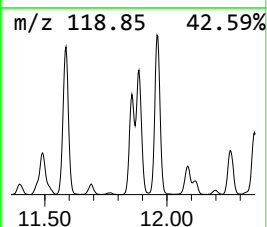
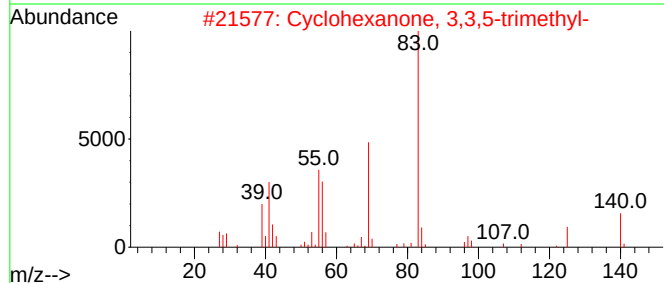
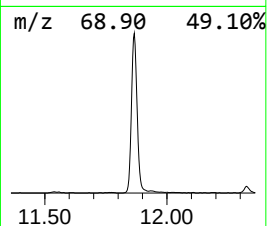
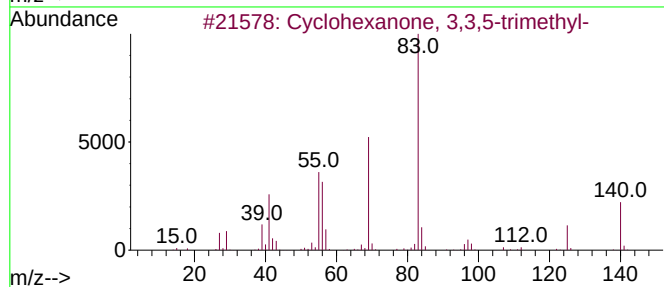
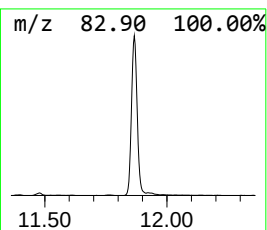
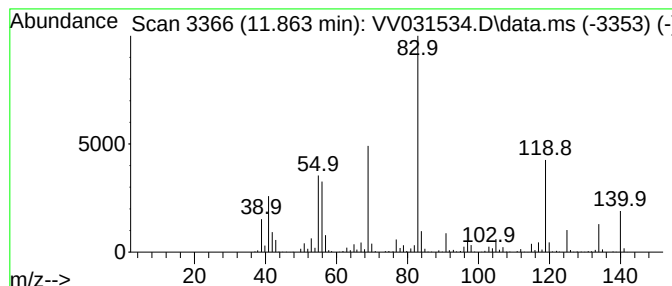
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Quant Title : VOC Analysis

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

Peak Number 13 Cyclohexanone, 3,3,5-trimet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.863	74.21 ug/L	1229840	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
5			Verbenyl angelate, cis-	234	C15H22O2	1000383-56-3	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

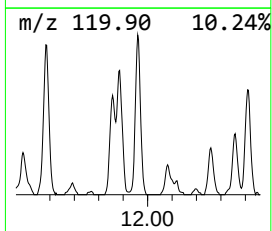
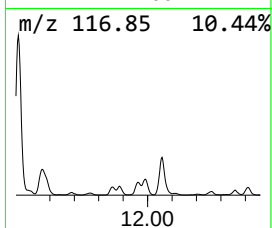
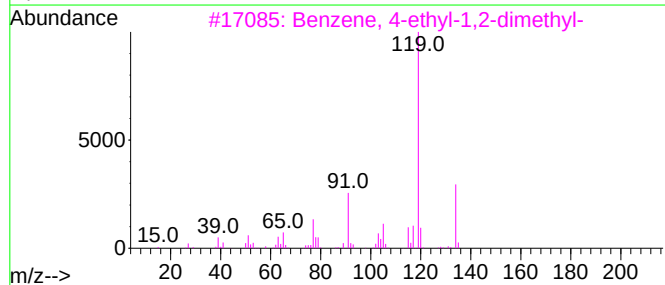
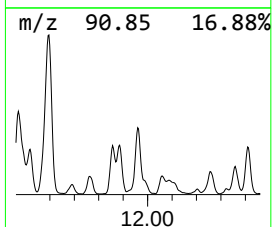
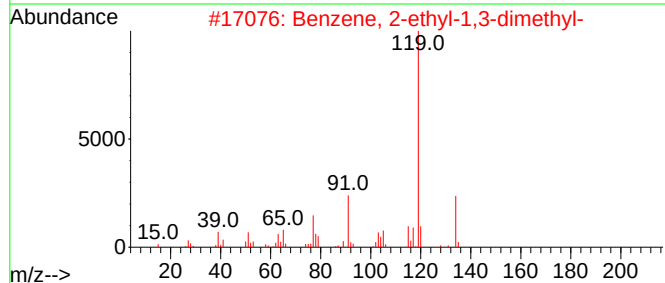
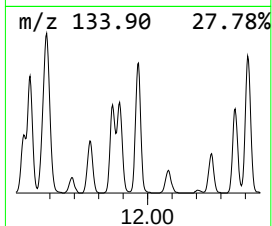
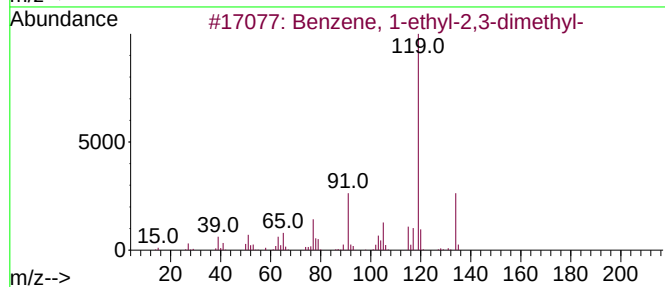
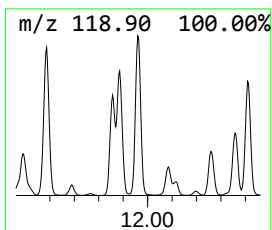
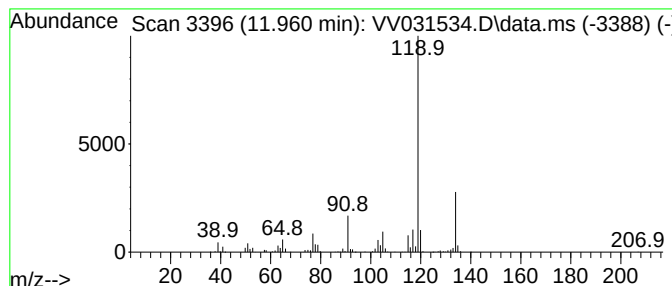
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.960	28.48 ug/L	472002	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

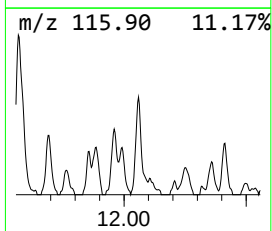
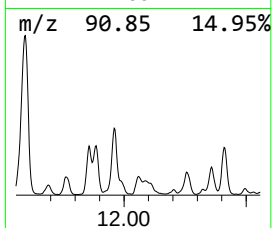
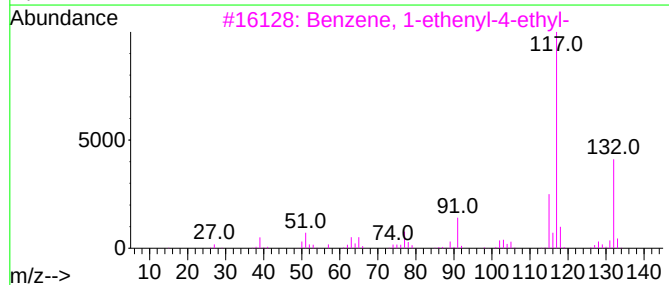
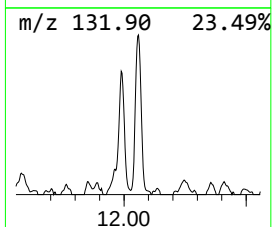
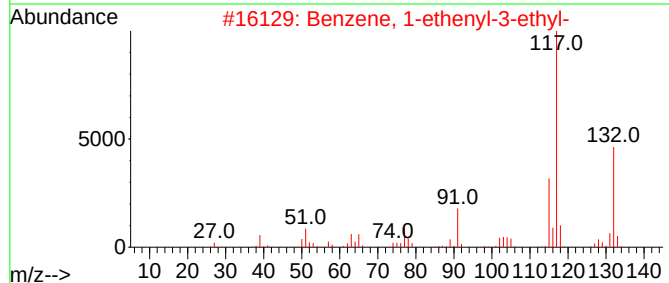
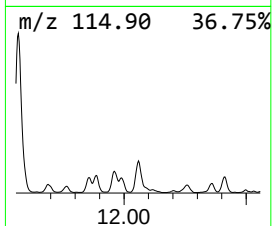
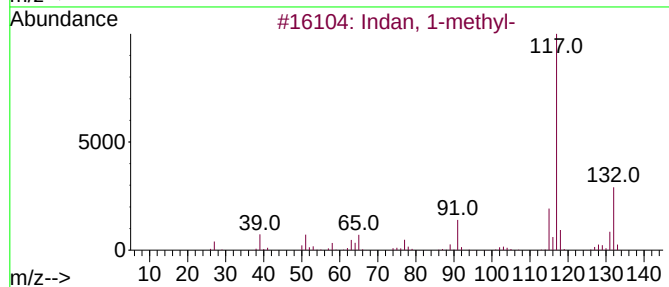
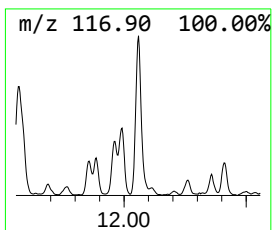
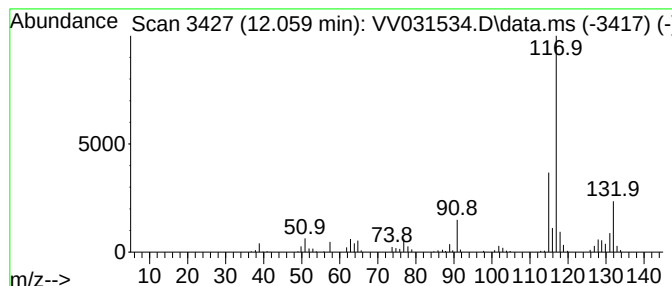
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Quant Title : VOC Analysis

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

Peak Number 15 Indan, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.059	9.17 ug/L	151924	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

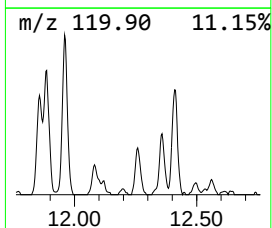
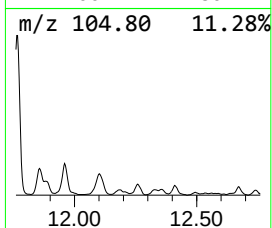
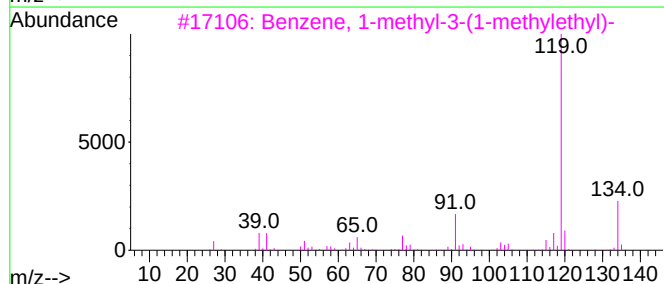
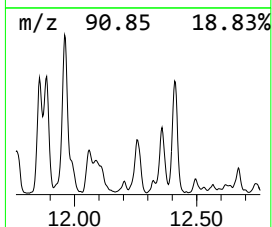
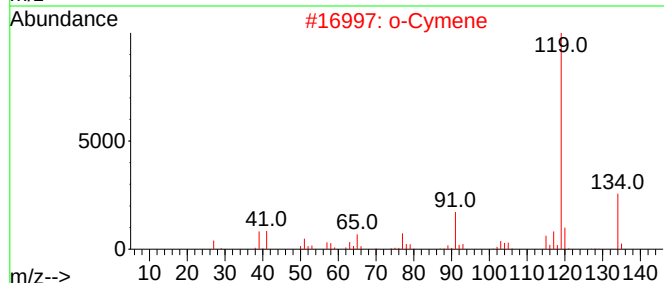
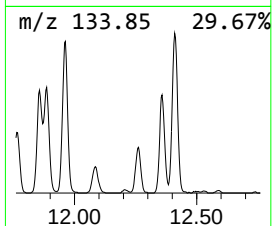
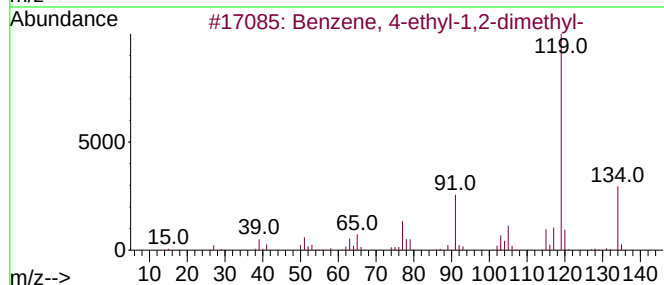
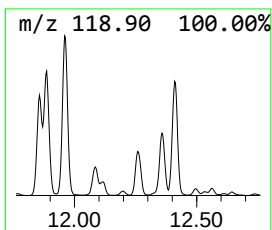
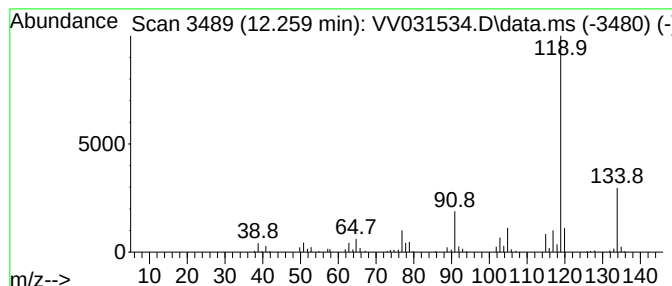
Instrument :
MSVOA_V
ClientSampleId :
EXYT2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.M
Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.259	8.11 ug/L	134355	1,4-Dichlorobenzene-d4			11.191
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-		134	C10H14	000934-80-5	95
2	o-Cymene		134	C10H14	000527-84-4	94
3	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	94
4	p-Cymene		134	C10H14	000099-87-6	93
5	o-Cymene		134	C10H14	000527-84-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

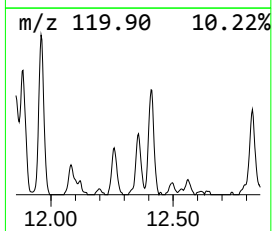
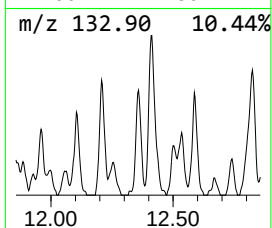
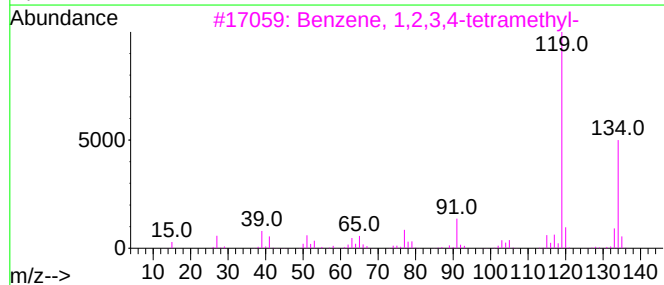
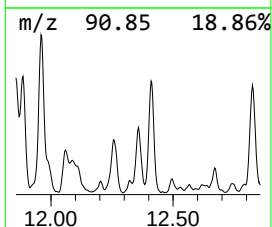
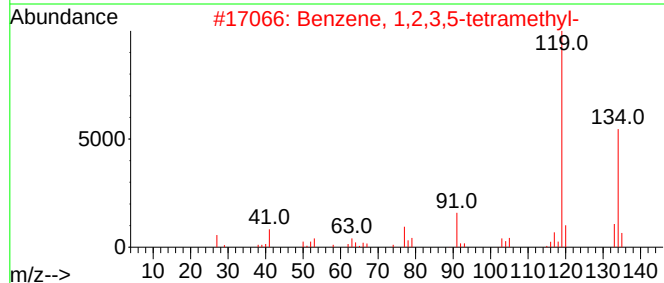
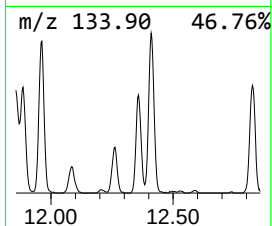
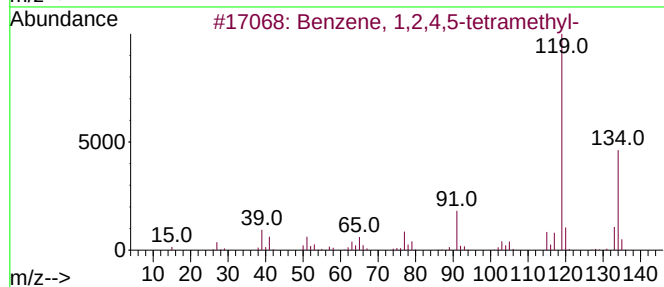
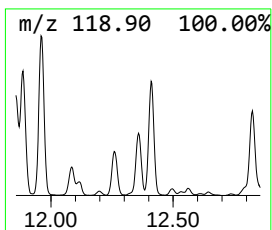
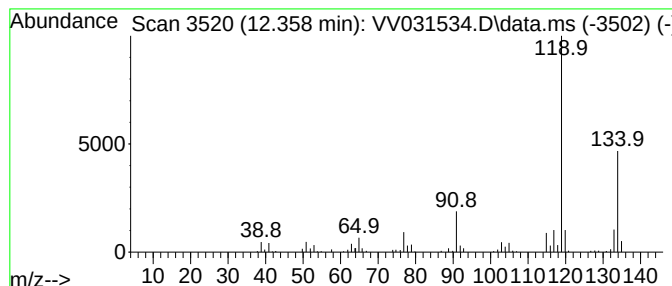
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Quant Title : VOC Analysis

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

Peak Number 17 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.358	11.88 ug/L	196934	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

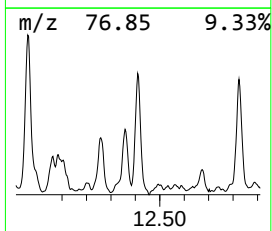
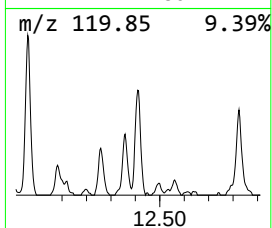
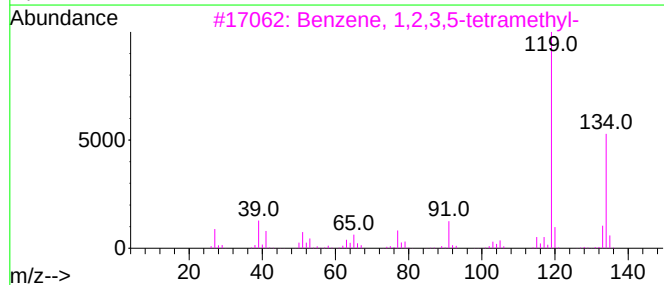
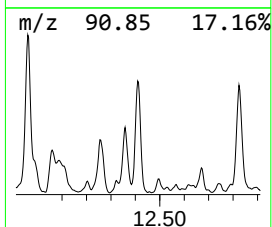
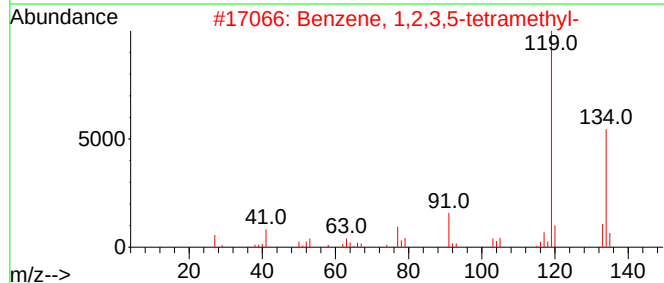
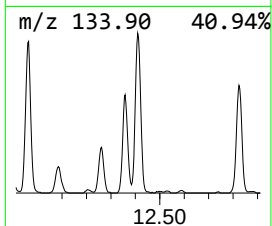
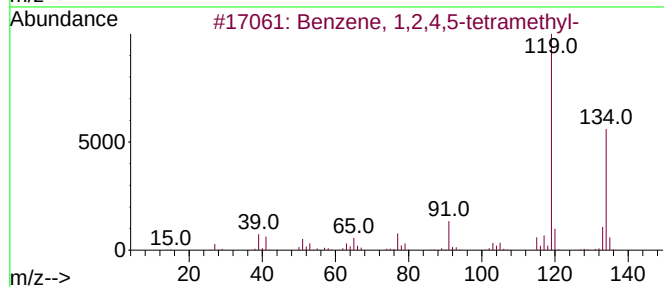
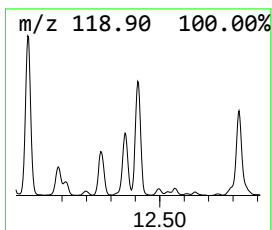
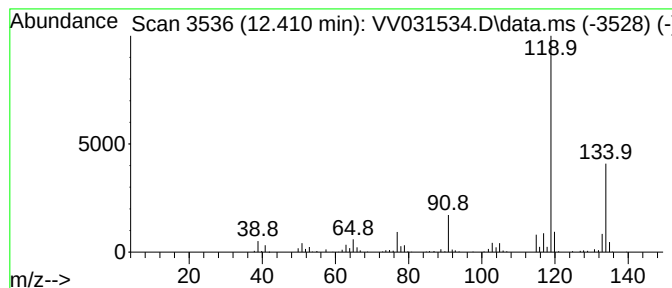
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Quant Title : VOC Analysis

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

Peak Number 18 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.410	17.62 ug/L	292040	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

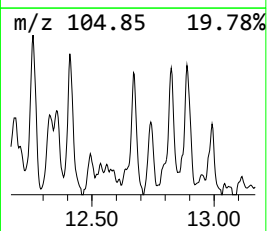
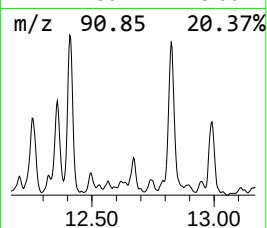
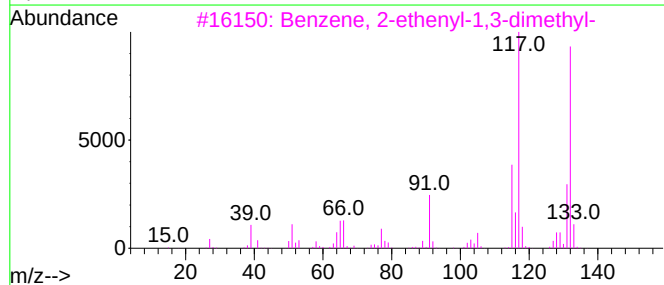
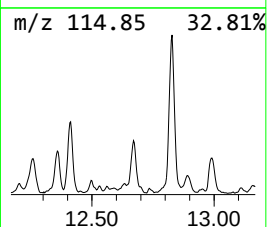
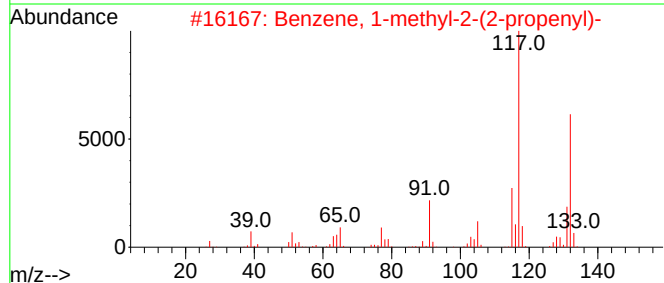
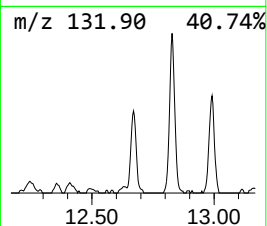
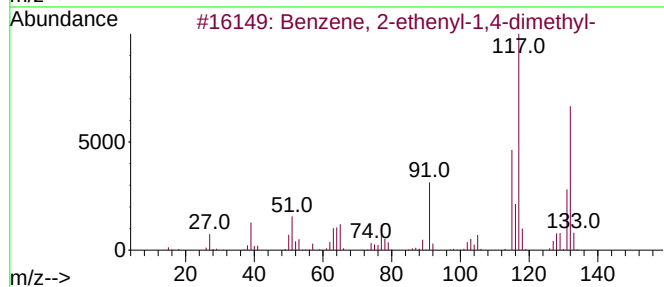
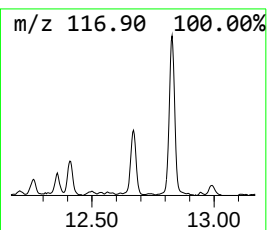
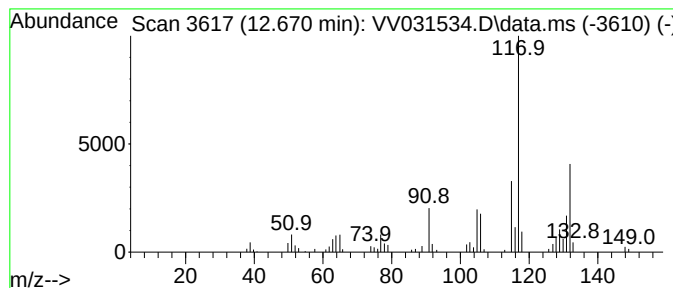
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Quant Title : VOC Analysis

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

Peak Number 19 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.670	4.58 ug/L	75860	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
3			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	81
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

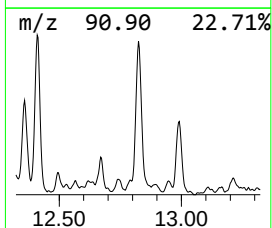
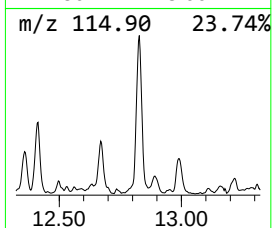
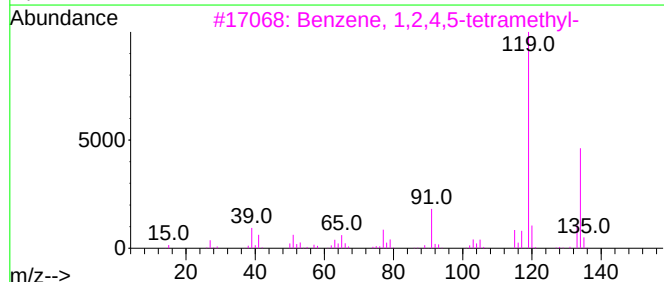
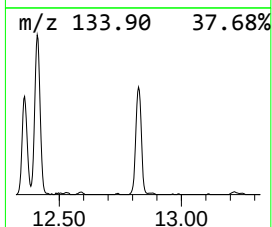
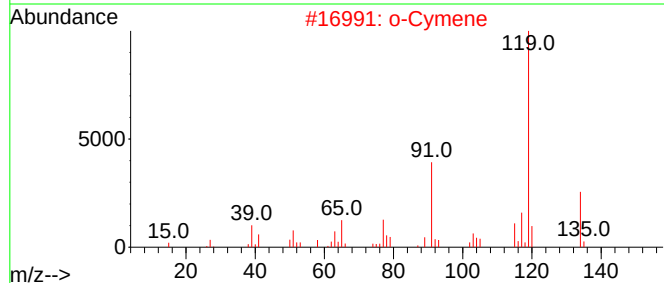
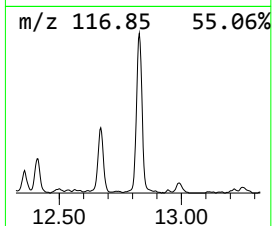
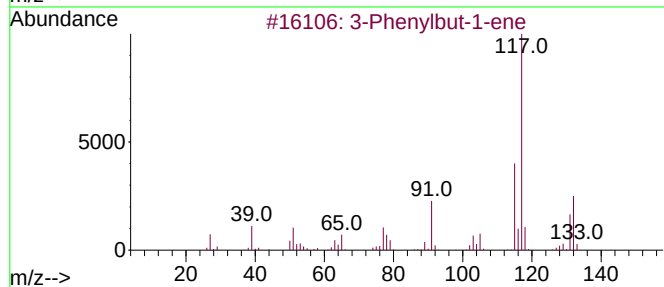
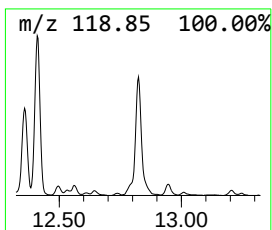
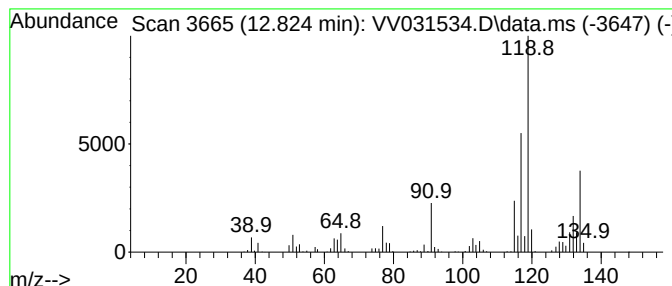
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Quant Title : VOC Analysis

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

Peak Number 20 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.824	21.50 ug/L	356380	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	80
2			o-Cymene	134	C10H14	000527-84-4	64
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	64
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

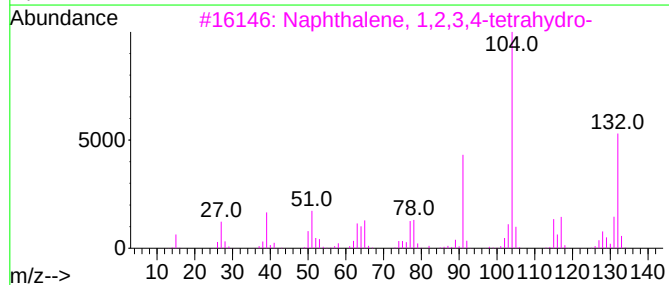
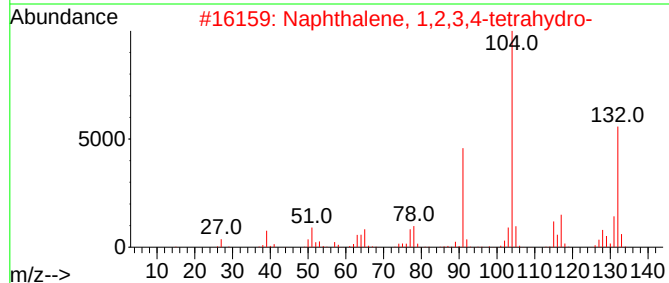
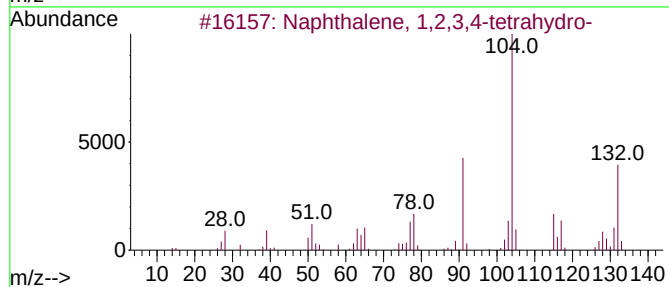
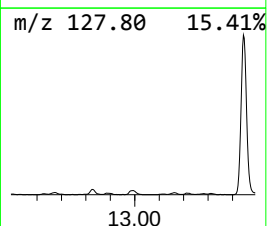
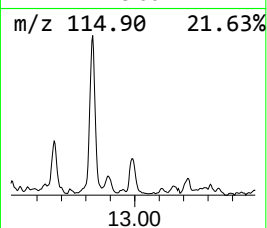
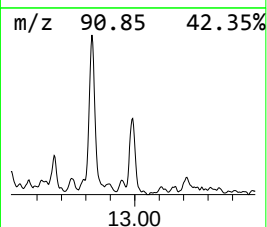
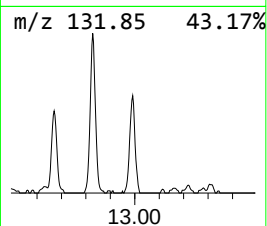
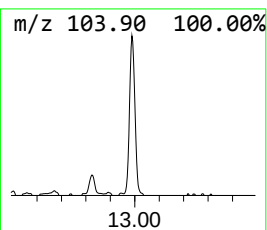
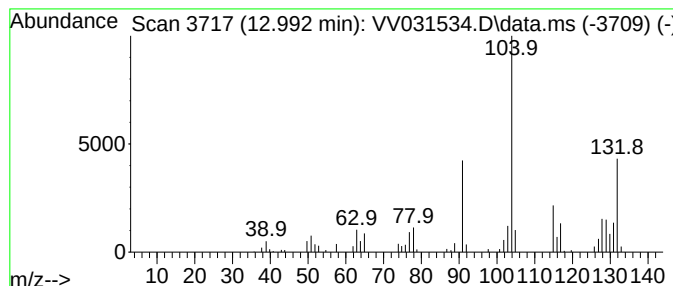
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Quant Title : VOC Analysis

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

Peak Number 21 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.992	5.07 ug/L	83972	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
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	90
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV061523\
Data File : VV031534.D
Acq On : 15 Jun 2023 16:23
Operator : SY/MD
Sample : 03249-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXYT2

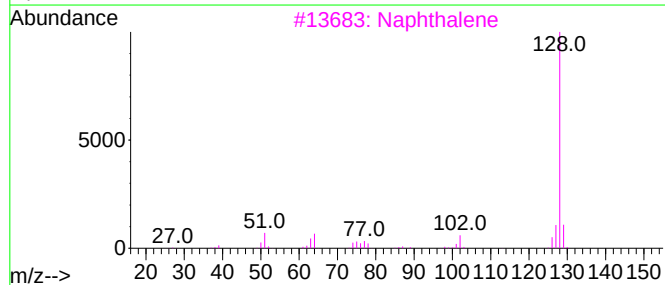
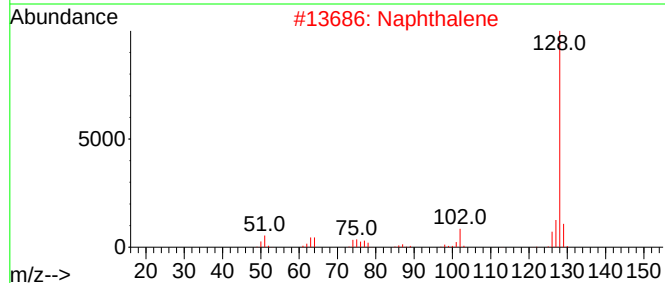
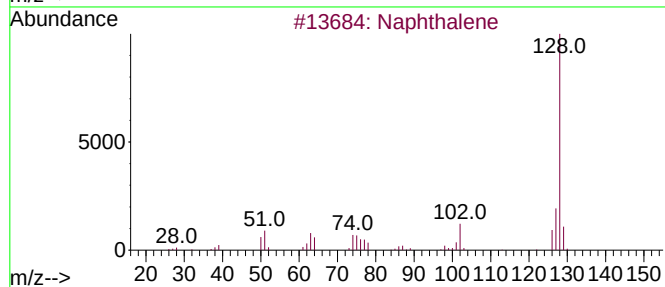
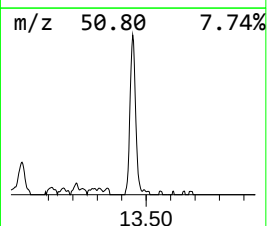
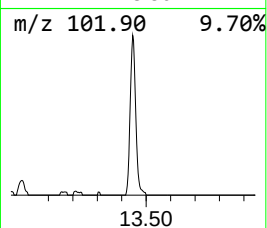
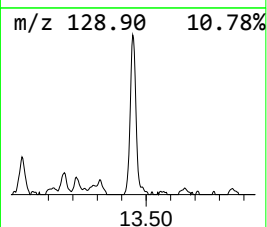
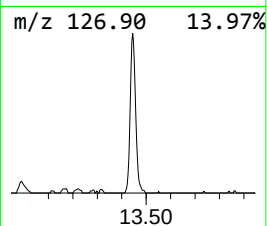
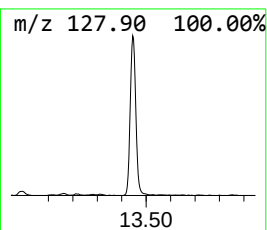
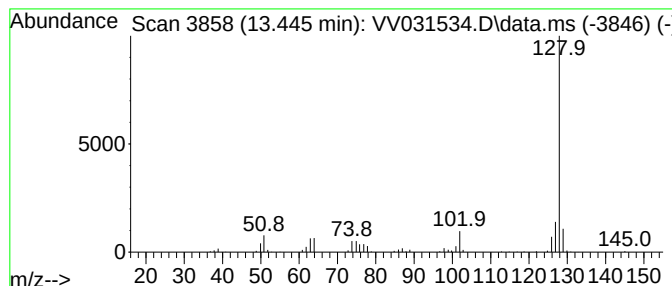
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Quant Title : VOC Analysis

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

Peak Number 22 Azulene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.445	14.62 ug/L	242335	1,4-Dichlorobenzene-d4	11.191

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	97
2			Naphthalene	128	C10H8	000091-20-3	95
3			Naphthalene	128	C10H8	000091-20-3	95
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW061523\
 Data File : VW031534.D
 Acq On : 15 Jun 2023 16:23
 Operator : SY/MD
 Sample : 03249-12
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EXYT2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM061423WMA.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
unknown-01	9.644	2.7	ug/L	33264	2	8.789	620874	50.0
Benzene, propyl-	10.297	35.7	ug/L	591231	3	11.191	828670	50.0
Benzene, 1-ethy...	10.390	125.0	ug/L	2071800	3	11.191	828670	50.0
Benzene, 1-ethy...	10.680	85.0	ug/L	1408210	3	11.191	828670	50.0
Benzene, (2-met...	10.998	2.6	ug/L	43303	3	11.191	828670	50.0
Oxirane, 2-meth...	11.030	6.0	ug/L	98826	3	11.191	828670	50.0
p-Cymene	11.140	9.3	ug/L	154974	3	11.191	828670	50.0
Benzene, 1,2,3-...	11.278	125.5	ug/L	2079320	3	11.191	828670	50.0
Indane	11.471	48.8	ug/L	807997	3	11.191	828670	50.0
Benzene, 1,2-di...	11.689	3.6	ug/L	59563	3	11.191	828670	50.0
Benzene, 1-meth...	11.763	11.0	ug/L	182061	3	11.191	828670	50.0
Cyclohexanone, ...	11.863	74.2	ug/L	1229840	3	11.191	828670	50.0
Benzene, 1-ethy...	11.960	28.5	ug/L	472002	3	11.191	828670	50.0
Indan, 1-methyl-	12.059	9.2	ug/L	151924	3	11.191	828670	50.0
Benzene, 4-ethy...	12.259	8.1	ug/L	134355	3	11.191	828670	50.0
Benzene, 1,2,4,...	12.358	11.9	ug/L	196934	3	11.191	828670	50.0
Benzene, 1,2,3,...	12.410	17.6	ug/L	292040	3	11.191	828670	50.0
Benzene, 2-ethe...	12.670	4.6	ug/L	75860	3	11.191	828670	50.0
3-Phenylbut-1-ene	12.824	21.5	ug/L	356380	3	11.191	828670	50.0
Naphthalene, 1,...	12.992	5.1	ug/L	83972	3	11.191	828670	50.0
Azulene	13.445	14.6	ug/L	242335	3	11.191	828670	50.0