

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
 Data File : VU059323.D
 Acq On : 18 Jun 2024 19:55
 Operator : MD/SY
 Sample : P2897-14 200X
 Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0T61

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\SFAMULM052924WMA.M
 Title : VOC Analysis

Signal : TIC: VU059323.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.595	88	95	113	rBV	52022	62952	3.24%	0.793%
2	1.907	185	192	204	rBV	46878	63080	3.25%	0.795%
3	2.560	384	395	414	rBV	85499	138937	7.16%	1.751%
4	2.901	495	501	502	rBV	219	198	0.01%	0.002%
5	2.949	512	516	517	rBV2	191	171	0.01%	0.002%
6	3.039	533	544	557	rVB2	6516	11962	0.62%	0.151%
7	3.145	574	577	580	rBV	350	273	0.01%	0.003%
8	3.373	644	648	651	rBV	205	218	0.01%	0.003%
9	3.660	731	737	739	rBV	239	249	0.01%	0.003%
10	3.730	756	759	762	rBV	166	152	0.01%	0.002%
11	3.769	769	771	772	rBV	175	94	0.00%	0.001%
12	3.840	790	793	797	rBV2	198	167	0.01%	0.002%
13	3.907	811	814	816	rBV	196	124	0.01%	0.002%
14	3.949	824	827	830	rBV	243	185	0.01%	0.002%
15	3.968	830	833	837	rVV	214	165	0.01%	0.002%
16	4.055	857	860	862	rBV	256	195	0.01%	0.002%
17	4.193	898	903	909	rBV	321	453	0.02%	0.006%
18	4.357	951	954	957	rBV	287	247	0.01%	0.003%
19	4.544	1009	1012	1014	rBV	166	121	0.01%	0.002%
20	4.615	1024	1034	1065	rBV	49148	125247	6.45%	1.578%
21	5.055	1156	1171	1197	rBV2	84889	189748	9.78%	2.391%
22	5.412	1279	1282	1286	rVB	260	214	0.01%	0.003%
23	5.496	1305	1308	1311	rBV2	152	118	0.01%	0.001%
24	5.602	1337	1341	1344	rBV2	259	242	0.01%	0.003%
25	5.721	1358	1378	1401	rBV2	165313	463884	23.90%	5.846%
26	5.994	1459	1463	1467	rVB3	366	302	0.02%	0.004%
27	6.058	1478	1483	1484	rBV	248	203	0.01%	0.003%
28	6.245	1529	1541	1560	rBV	195338	371288	19.13%	4.679%
29	6.479	1612	1614	1621	rVB3	298	383	0.02%	0.005%
30	6.531	1621	1630	1641	rBV4	4553	8317	0.43%	0.105%
31	6.618	1654	1657	1659	rVB3	193	108	0.01%	0.001%
32	6.685	1665	1678	1697	rBV	110487	213791	11.01%	2.694%
33	6.804	1713	1715	1718	rBV	305	204	0.01%	0.003%
34	6.955	1759	1762	1765	rBV2	232	198	0.01%	0.002%
35	7.004	1775	1777	1780	rBV	195	131	0.01%	0.002%
36	7.061	1792	1795	1802	rBV2	232	243	0.01%	0.003%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
 Title : VOC Analysis

37	7.090	1802	1804	1808	rBV2	275	185	0.01%	0.002%
38	7.110	1808	1810	1815	rBV	226	219	0.01%	0.003%
39	7.190	1831	1835	1837	rBV3	794	626	0.03%	0.008%
40	7.370	1885	1891	1893	rBV	158	170	0.01%	0.002%
41	7.438	1908	1912	1920	rVB	288	455	0.02%	0.006%
42	7.473	1920	1923	1925	rBV	147	134	0.01%	0.002%
43	7.563	1940	1951	1967	rVV	49412	86779	4.47%	1.094%
44	7.708	1992	1996	2001	rBV4	737	844	0.04%	0.011%
45	7.807	2024	2027	2031	rVB2	308	236	0.01%	0.003%
46	7.840	2031	2037	2041	rBV2	285	418	0.02%	0.005%
47	7.894	2041	2054	2070	rBV	184515	329064	16.95%	4.147%
48	8.177	2132	2142	2161	rBV	34206	58252	3.00%	0.734%
49	8.290	2174	2177	2180	rVB2	283	195	0.01%	0.002%
50	8.357	2194	2198	2201	rBV	225	219	0.01%	0.003%
51	8.402	2209	2212	2214	rBV2	259	211	0.01%	0.003%
52	8.550	2249	2258	2270	rVB4	7362	12036	0.62%	0.152%
53	8.627	2272	2282	2318	rBV	120855	244450	12.59%	3.080%
54	9.029	2399	2407	2410	rBV2	310	433	0.02%	0.005%
55	9.071	2417	2420	2422	rBV	240	166	0.01%	0.002%
56	9.116	2431	2434	2436	rBV	162	140	0.01%	0.002%
57	9.258	2474	2478	2482	rBV2	312	345	0.02%	0.004%
58	9.412	2514	2526	2562	rBV	284802	488556	25.17%	6.156%
59	9.566	2565	2574	2590	rVB	63372	103417	5.33%	1.303%
60	9.692	2603	2613	2634	rVB	40654	70603	3.64%	0.890%
61	9.827	2652	2655	2660	rBV2	251	175	0.01%	0.002%
62	10.061	2725	2728	2730	rBV	195	120	0.01%	0.002%
63	10.100	2730	2740	2756	rVV	22811	37596	1.94%	0.474%
64	10.283	2794	2797	2799	rBV	187	138	0.01%	0.002%
65	10.405	2832	2835	2838	rVB2	173	98	0.01%	0.001%
66	10.483	2849	2859	2873	rBV2	7154	11363	0.59%	0.143%
67	10.749	2931	2942	2964	rVB	168492	272207	14.02%	3.430%
68	10.859	2972	2976	2978	rBV2	181	171	0.01%	0.002%
69	10.997	3008	3019	3038	rBV	47765	112237	5.78%	1.414%
70	11.084	3039	3046	3062	rVB3	12442	19515	1.01%	0.246%
71	11.290	3103	3110	3124	rVB5	3143	5710	0.29%	0.072%
72	11.409	3143	3147	3151	rBV2	245	282	0.01%	0.004%
73	11.466	3152	3165	3178	rVV2	39714	63332	3.26%	0.798%
74	11.540	3178	3188	3202	rVB2	9585	15372	0.79%	0.194%
75	11.589	3202	3203	3206	rBV2	199	126	0.01%	0.002%
76	11.637	3216	3218	3226	rVB4	584	391	0.02%	0.005%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
 Title : VOC Analysis

77	11.740	3241	3250	3258	rBV6	4309	6435	0.33%	0.081%
78	11.807	3259	3271	3288	rVV	298858	488504	25.17%	6.156%
79	11.894	3291	3298	3307	rVV3	14169	22282	1.15%	0.281%
80	11.949	3310	3315	3328	rVB2	7585	11826	0.61%	0.149%
81	12.093	3350	3360	3377	rVB4	21097	35349	1.82%	0.445%
82	12.190	3377	3390	3409	rBV	212102	348854	17.97%	4.396%
83	12.309	3417	3427	3442	rBV	120705	200813	10.35%	2.531%
84	12.569	3499	3508	3518	rBV	14972	23138	1.19%	0.292%
85	12.749	3554	3564	3575	rVV5	14244	22161	1.14%	0.279%
86	12.804	3576	3581	3594	rVB5	3901	5941	0.31%	0.075%
87	13.026	3641	3650	3663	rBV2	7053	11775	0.61%	0.148%
88	13.293	3727	3733	3741	rVB4	4396	6183	0.32%	0.078%
89	13.444	3770	3780	3790	rVB6	8182	16531	0.85%	0.208%
90	13.518	3793	3803	3815	rBV2	29542	46052	2.37%	0.580%
91	13.624	3827	3836	3850	rVB3	34998	58592	3.02%	0.738%
92	13.727	3861	3868	3875	rBV5	5529	7089	0.37%	0.089%
93	14.080	3966	3978	4004	rBV	1226375	1941091	100.00%	24.460%
94	14.958	4249	4251	4253	rBV	697	439	0.02%	0.006%
95	15.219	4321	4332	4353	rVB	336421	544755	28.06%	6.865%
96	15.405	4379	4390	4412	rBV	233444	385684	19.87%	4.860%
97	15.952	4553	4560	4571	rVB3	15923	24434	1.26%	0.308%
98	16.145	4613	4620	4628	rBV2	28531	40247	2.07%	0.507%
99	16.260	4649	4656	4675	rVB	25870	47022	2.42%	0.593%
100	16.405	4694	4701	4708	rBV2	32436	48923	2.52%	0.616%

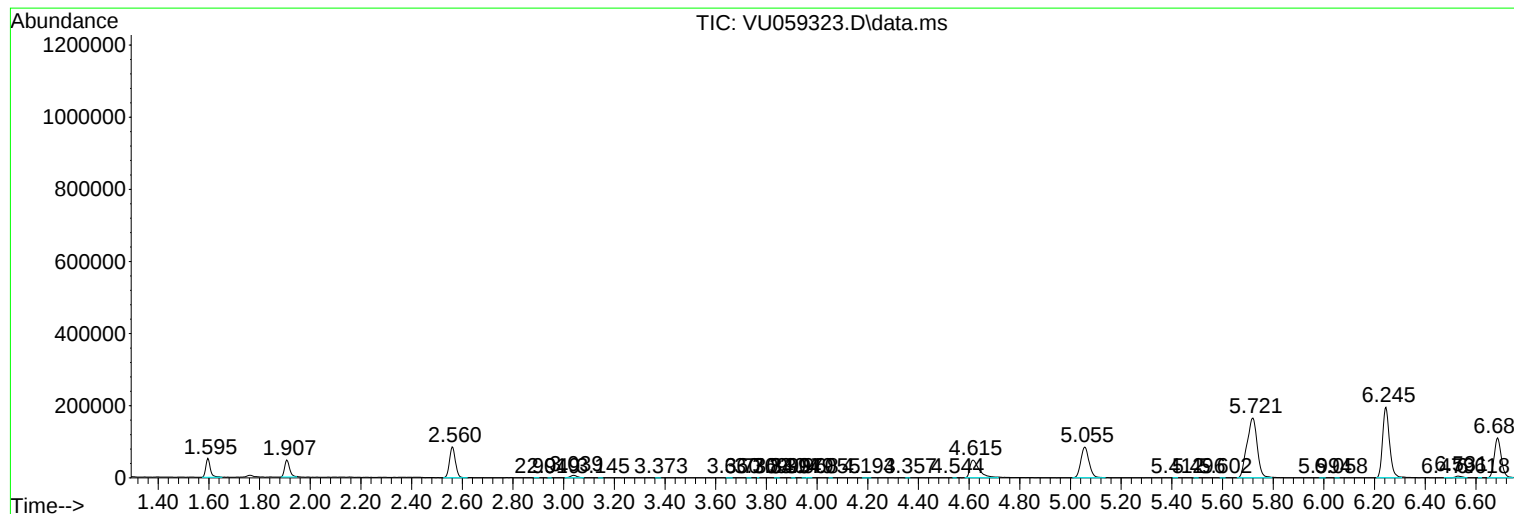
Sum of corrected areas: 7935670

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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
Quant Title : VOC Analysis

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



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Instrument :
MSVOA_U
ClientSampleId :
C0T61

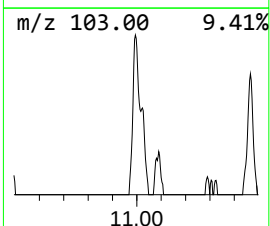
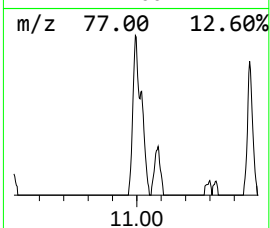
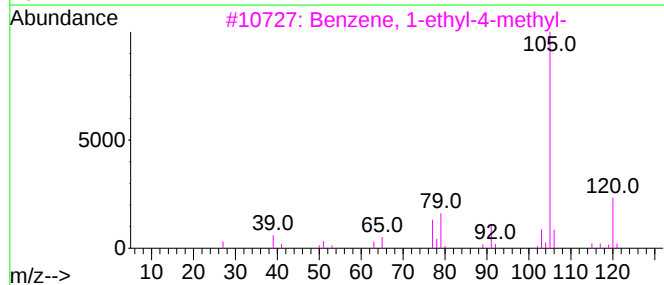
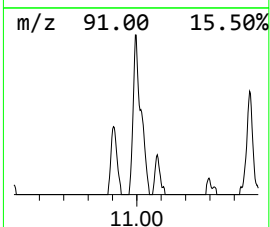
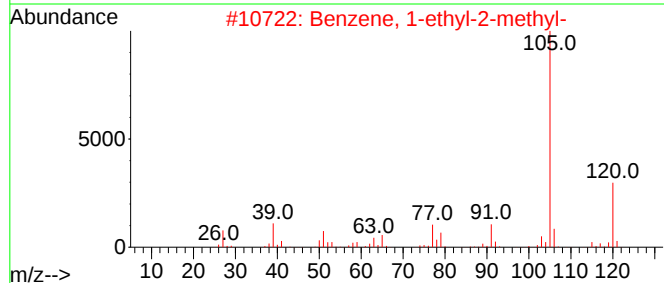
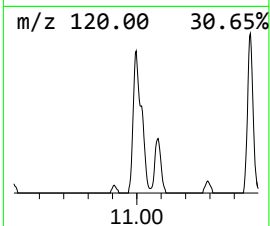
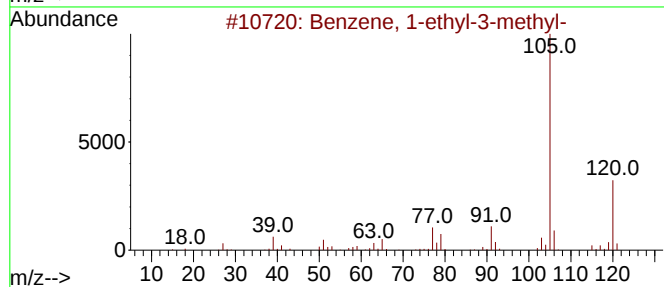
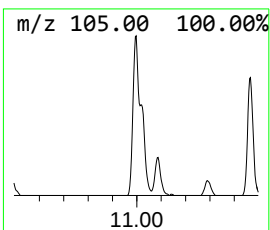
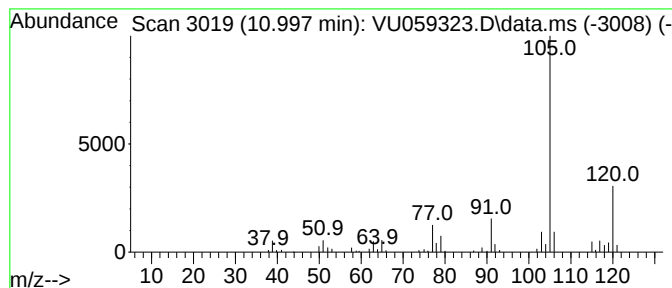
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.997	11.49 ug/L	112237	1,4-Dichlorobenzene-d4	11.807

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



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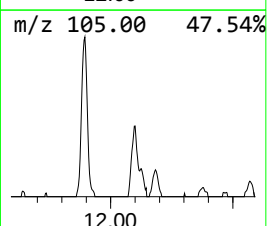
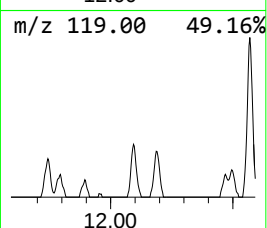
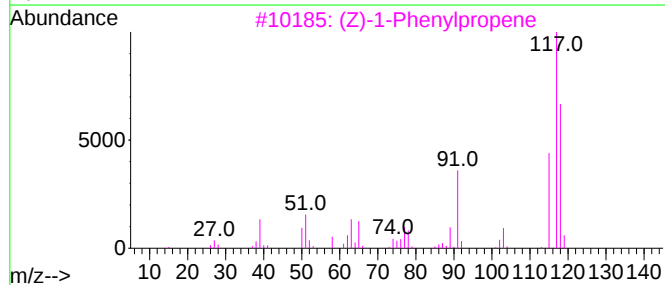
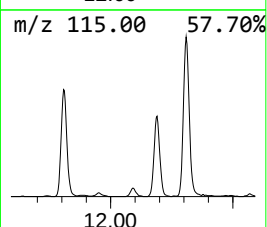
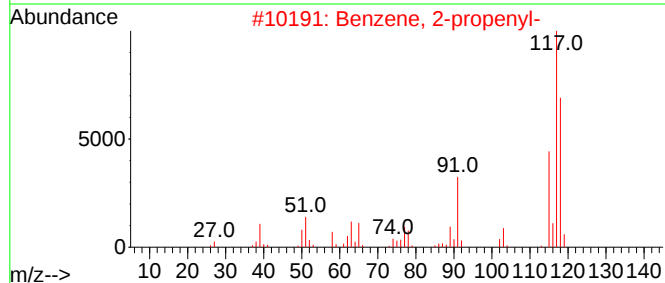
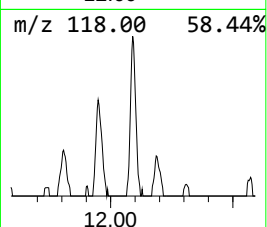
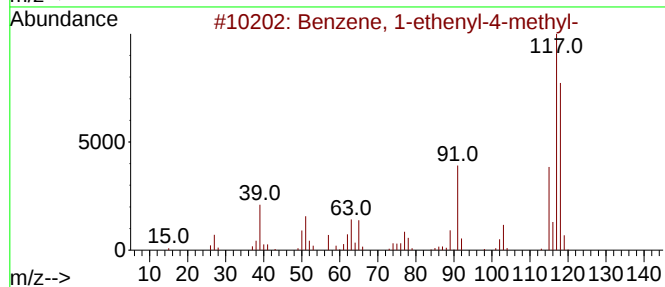
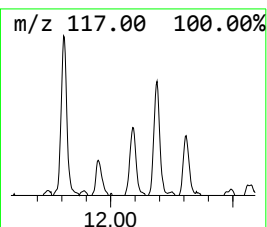
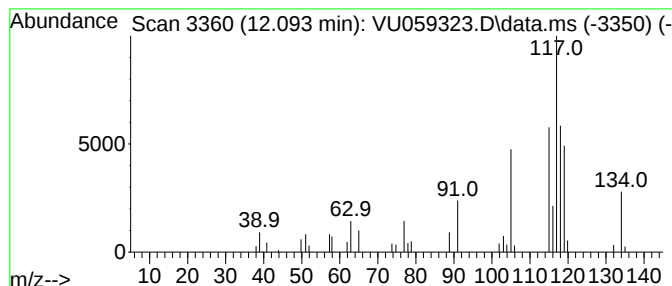
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1-ethenyl-4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.093	3.62 ug/L	35349	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	50
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	50
3			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	50
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	43
5			Diglycolic acid, isobutyl 3-phen...	308	C17H24O5	1010382-17-1	43



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Instrument :
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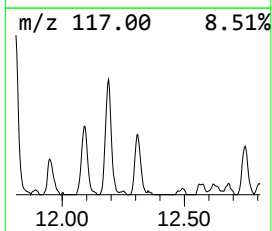
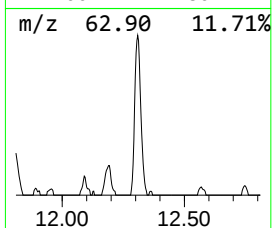
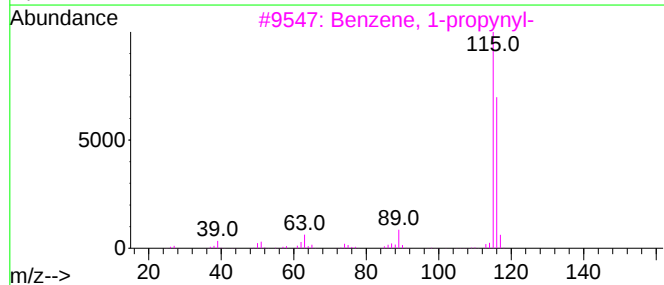
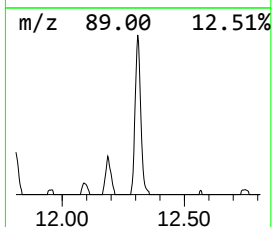
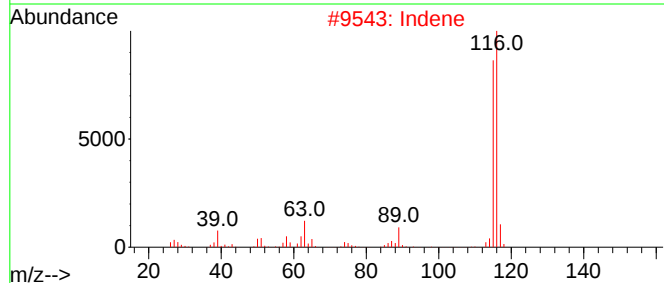
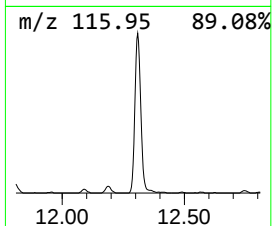
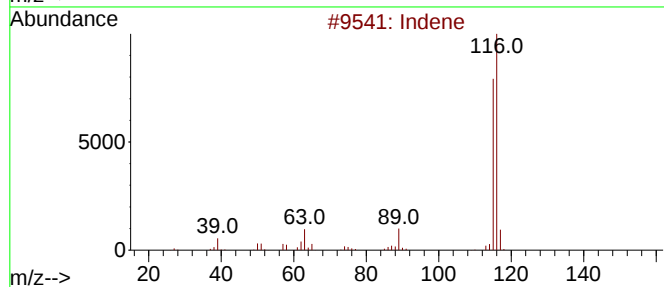
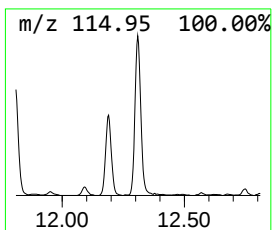
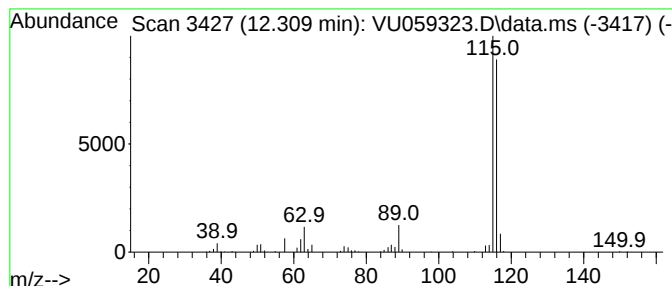
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.309	20.55 ug/L	200813	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Indene	116	C9H8	000095-13-6	95
3			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
4			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5			3-Methylphenylacetylene	116	C9H8	000766-82-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
COT61

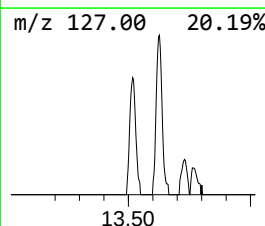
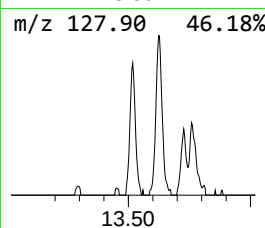
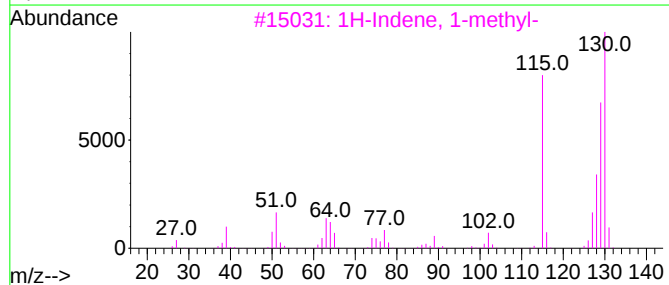
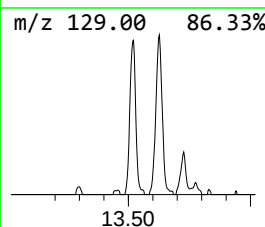
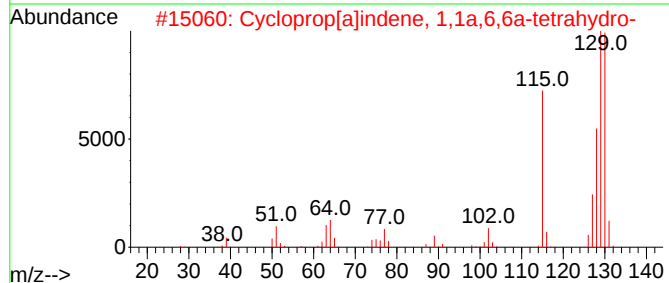
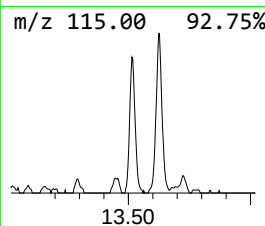
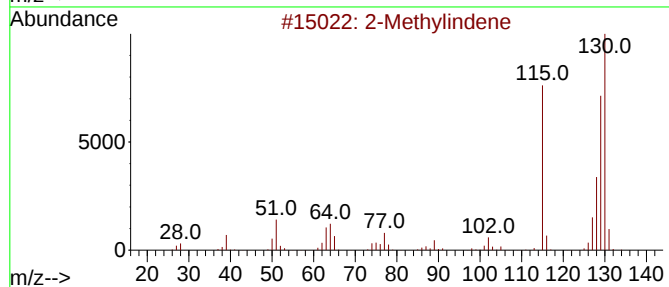
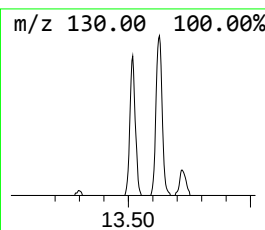
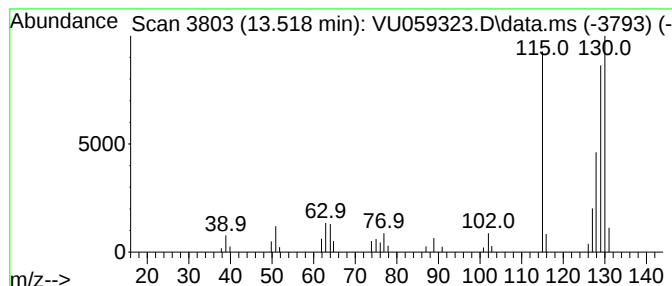
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Quant Title : VOC Analysis

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

Peak Number 5 2-Methylindene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.518	4.71 ug/L	46052	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methylindene	130	C10H10	002177-47-1	96
2			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93
3			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93
4			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	92
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
COT61

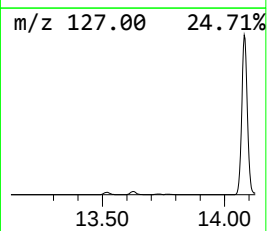
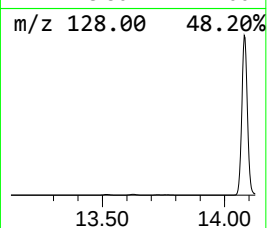
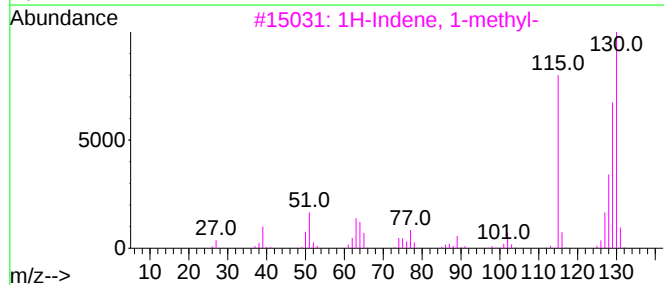
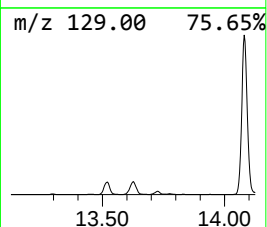
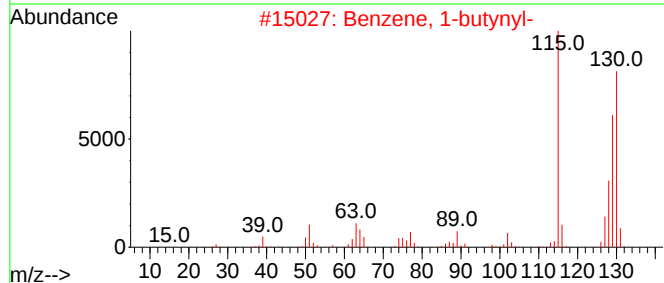
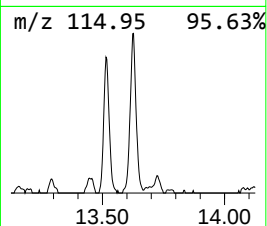
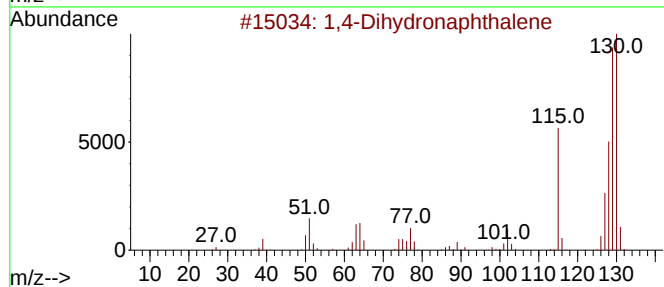
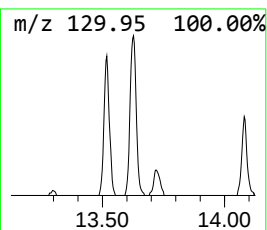
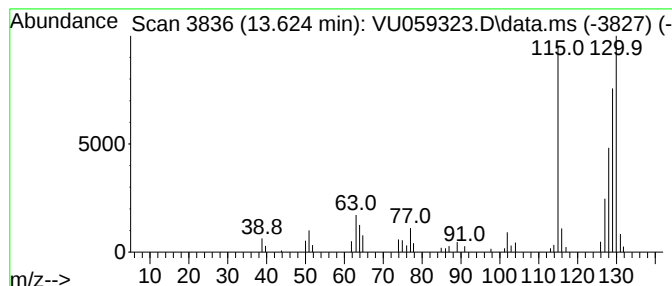
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Quant Title : VOC Analysis

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

Peak Number 6 1,4-Dihydronaphthalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.624	6.00 ug/L	58592	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95
2			Benzene, 1-butynyl-	130	C10H10	000622-76-4	91
3			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	91
4			1,4-Dihydronaphthalene	130	C10H10	000612-17-9	91
5			Tetracyclo[5.3.0.0<2,6>.0<3,10>]...	130	C10H10	034324-40-8	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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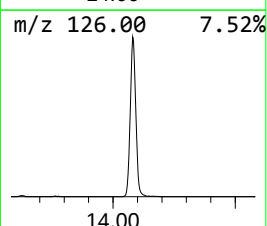
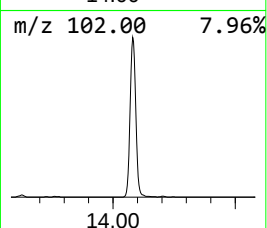
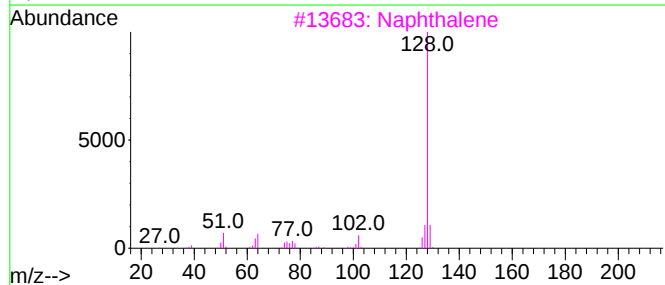
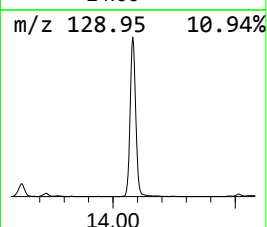
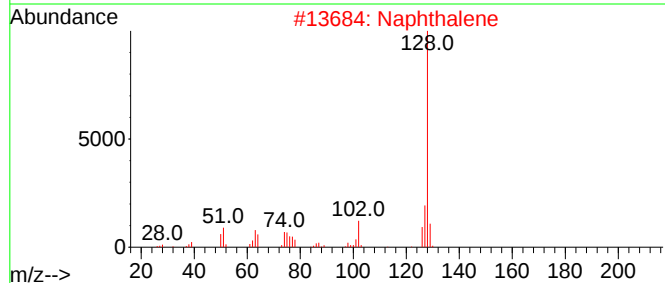
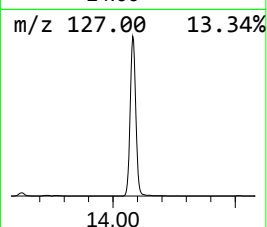
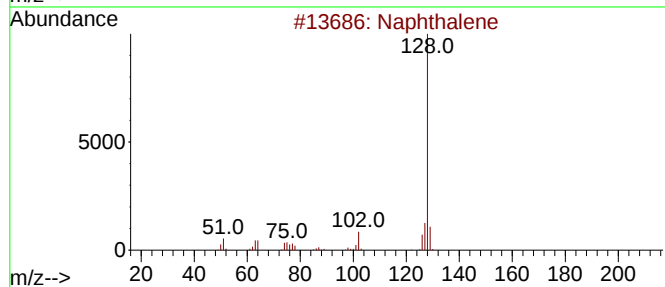
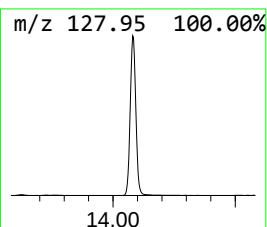
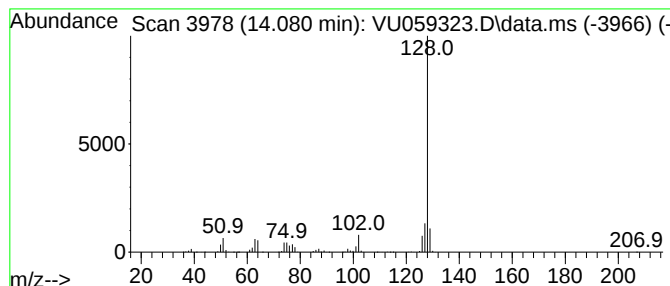
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Quant Title : VOC Analysis

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

Peak Number 7 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	198.68 ug/L	1941090	1,4-Dichlorobenzene-d4	11.807

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\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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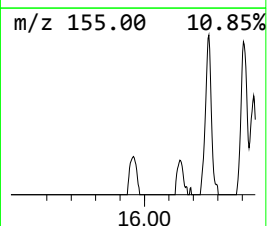
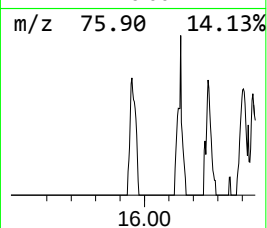
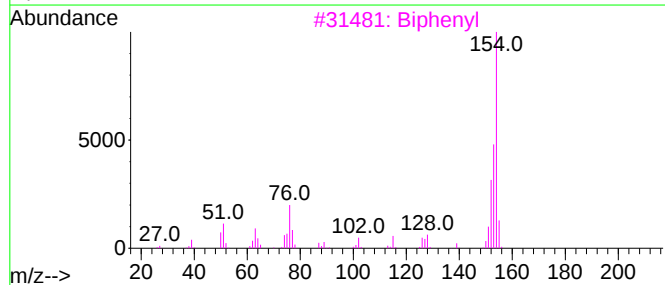
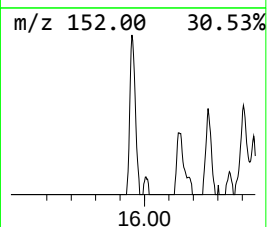
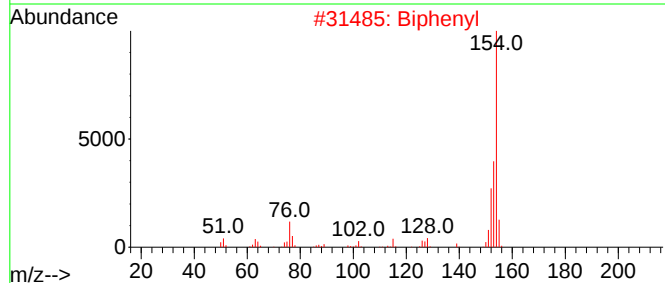
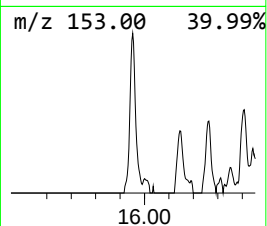
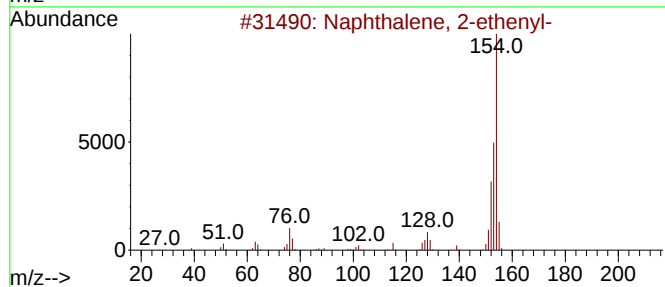
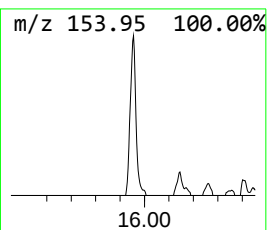
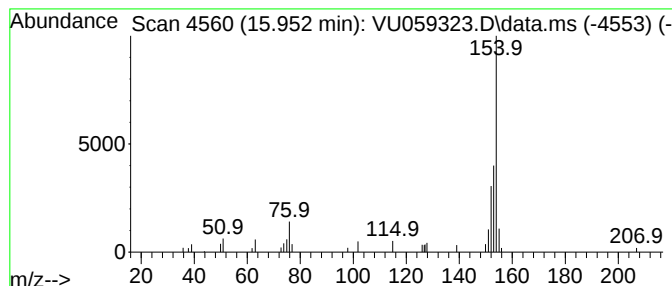
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Quant Title : VOC Analysis

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

Peak Number 10 Naphthalene, 2-ethenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.952	2.50 ug/L	24434	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	87
2			Biphenyl	154	C12H10	000092-52-4	87
3			Biphenyl	154	C12H10	000092-52-4	87
4			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	83
5			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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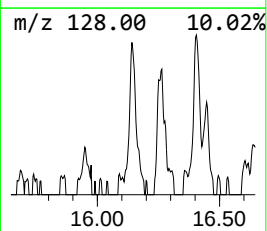
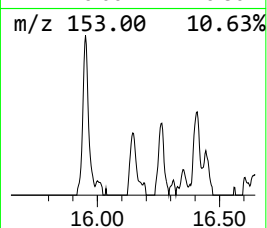
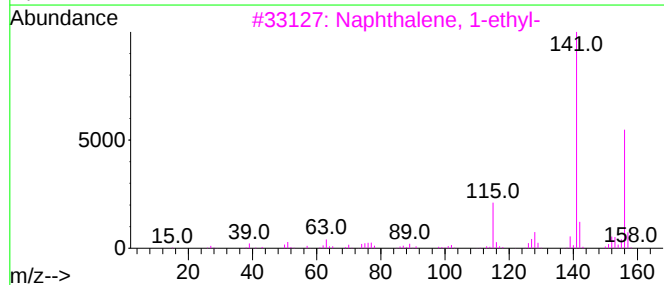
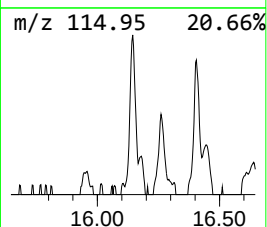
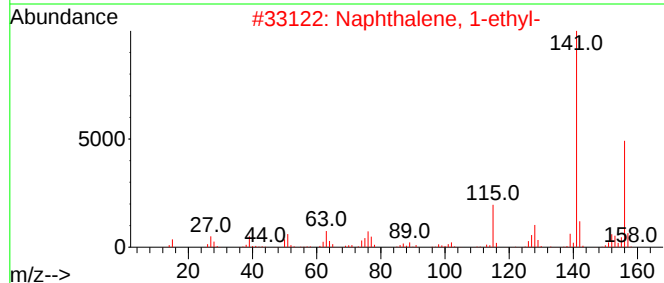
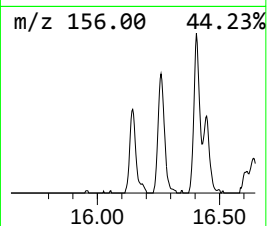
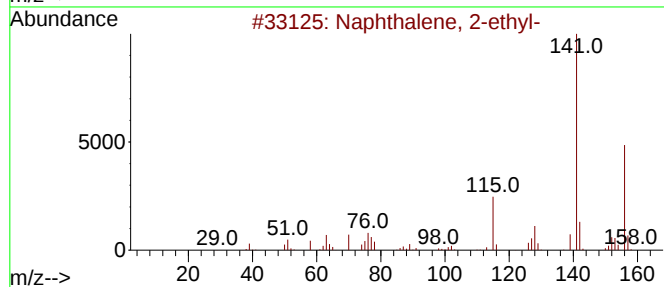
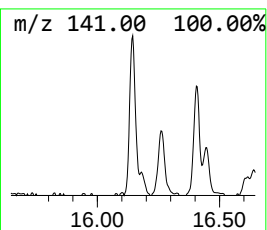
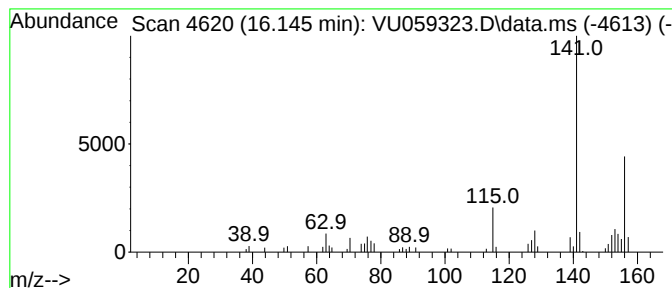
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Quant Title : VOC Analysis

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

Peak Number 11 Naphthalene, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.145	4.12 ug/L	40247	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
2			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
3			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	93
4			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	90
5			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
COT61

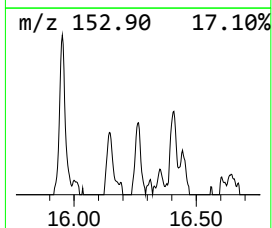
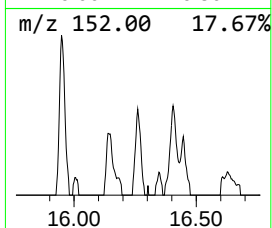
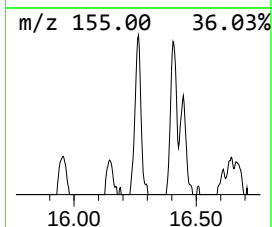
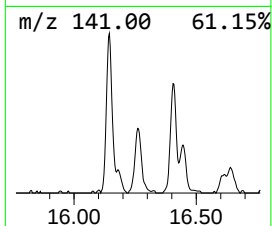
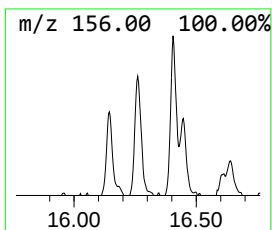
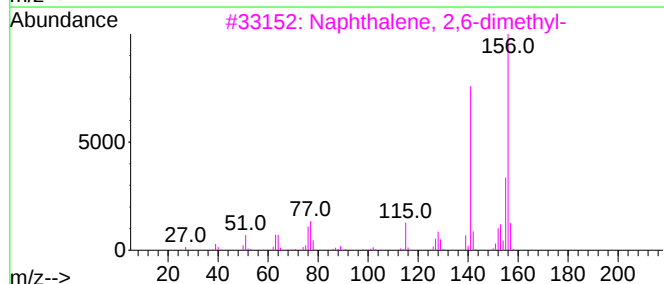
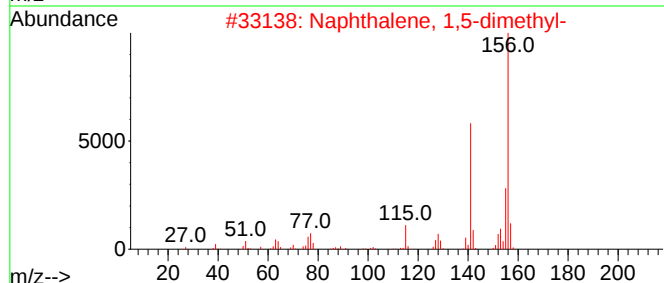
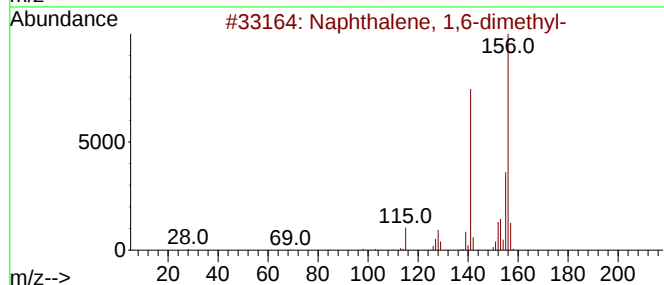
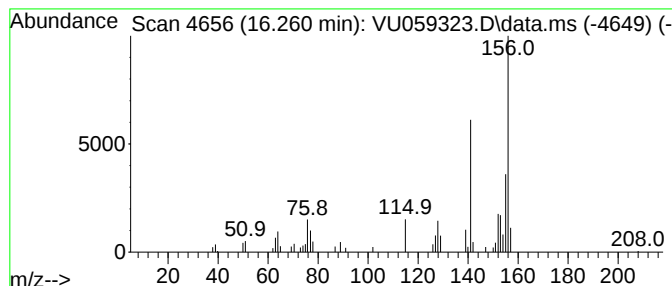
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Quant Title : VOC Analysis

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

Peak Number 12 Naphthalene, 1,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.261	4.81 ug/L	47022	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	94
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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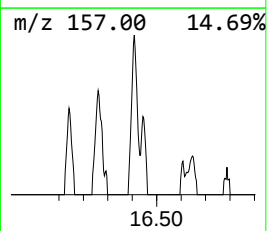
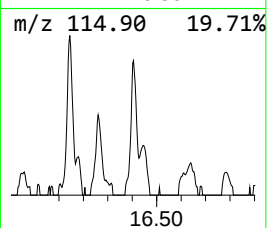
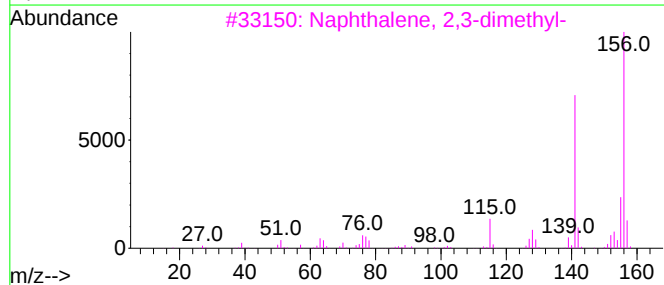
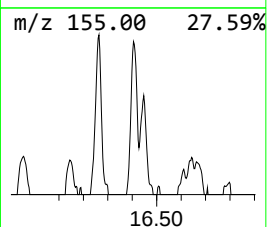
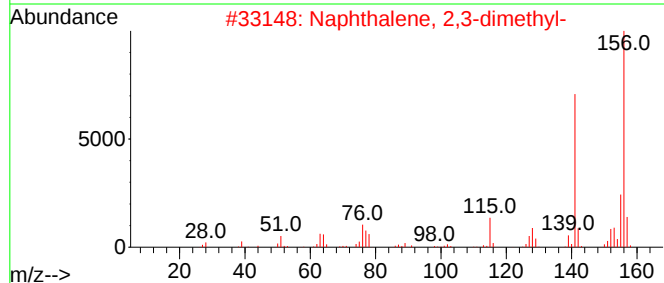
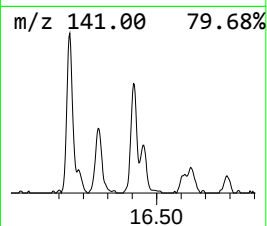
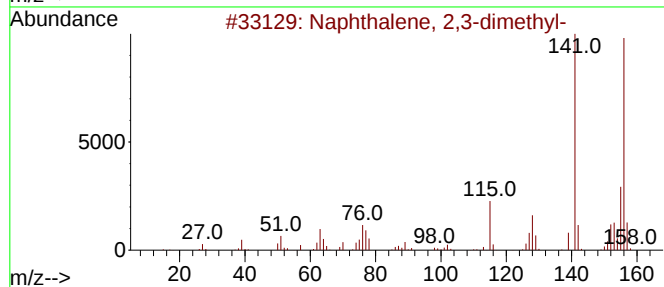
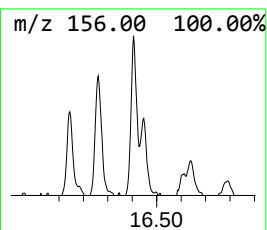
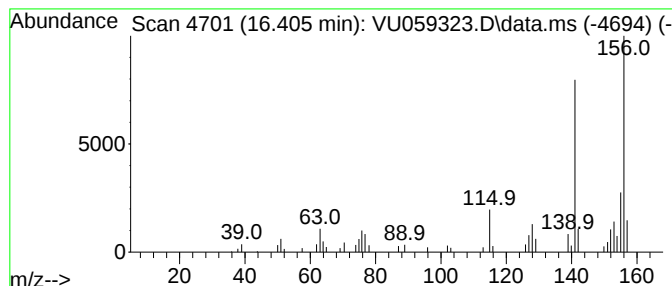
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.M
Quant Title : VOC Analysis

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

Peak Number 13 Naphthalene, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.405	5.01 ug/L	48923	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061824\
Data File : VU059323.D
Acq On : 18 Jun 2024 19:55
Operator : MD/SY
Sample : P2897-14 200X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0T61

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM052924WMA.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
Benzene, 1-ethy...	10.997	11.5	ug/L	112237	3	11.807	488504	50.0
Benzene, 1-ethe...	12.093	3.6	ug/L	35349	3	11.807	488504	50.0
Indene	12.309	20.6	ug/L	200813	3	11.807	488504	50.0
2-Methylindene	13.518	4.7	ug/L	46052	3	11.807	488504	50.0
1,4-Dihydronaph...	13.624	6.0	ug/L	58592	3	11.807	488504	50.0
Azulene	14.081	198.7	ug/L	1941090	3	11.807	488504	50.0
Naphthalene, 2-...	15.952	2.5	ug/L	24434	3	11.807	488504	50.0
Naphthalene, 2-...	16.145	4.1	ug/L	40247	3	11.807	488504	50.0
Naphthalene, 1,...	16.261	4.8	ug/L	47022	3	11.807	488504	50.0
Naphthalene, 2,...	16.405	5.0	ug/L	48923	3	11.807	488504	50.0