

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072922\
 Data File : VV027092.D
 Acq On : 29 Jul 2022 18:09
 Operator : SY/MD
 Sample : N3899-04 10X
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBUR3

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\SFAMVLM072822WMA.M
 Title : VOC Analysis

Signal : TIC: VV027092.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.304	74	82	98	rVB	322442	343476	5.80%	1.293%
2	1.565	155	163	178	rVB	276834	310885	5.25%	1.170%
3	2.101	319	330	342	rBV	563859	800277	13.51%	3.013%
4	2.201	343	361	380	rVV2	229299	462916	7.81%	1.743%
5	3.828	858	867	871	rVB5	562	970	0.02%	0.004%
6	3.886	871	885	909	rBV2	112023	409040	6.90%	1.540%
7	4.343	1011	1027	1050	rBV	361115	892084	15.06%	3.359%
8	4.847	1181	1184	1192	rVB5	388	543	0.01%	0.002%
9	5.040	1227	1244	1298	rBV2	858503	2315406	39.08%	8.717%
10	5.326	1330	1333	1337	rVV2	821	631	0.01%	0.002%
11	5.352	1339	1341	1344	rVV2	703	518	0.01%	0.002%
12	5.381	1347	1350	1354	rVB2	1254	802	0.01%	0.003%
13	5.416	1354	1361	1363	rBV5	552	618	0.01%	0.002%
14	5.487	1378	1383	1386	rBV5	693	773	0.01%	0.003%
15	5.548	1394	1402	1403	rBV4	2239	2341	0.04%	0.009%
16	5.613	1410	1422	1458	rBV	571889	1210892	20.44%	4.559%
17	5.908	1503	1514	1534	rVB3	13930	30056	0.51%	0.113%
18	6.066	1547	1563	1596	rBV	487597	1028589	17.36%	3.872%
19	6.384	1658	1662	1668	rVB6	1327	1365	0.02%	0.005%
20	6.413	1668	1671	1677	rVB6	836	834	0.01%	0.003%
21	6.567	1711	1719	1720	rBV2	1310	1388	0.02%	0.005%
22	6.658	1740	1747	1749	rVB5	1004	1250	0.02%	0.005%
23	6.986	1837	1849	1879	rBV	223358	434318	7.33%	1.635%
24	7.240	1919	1928	1939	rBV6	9396	22183	0.37%	0.084%
25	7.313	1940	1951	1969	rBV	946328	1673289	28.24%	6.300%
26	7.619	2037	2046	2074	rBV	159387	295627	4.99%	1.113%
27	7.863	2119	2122	2125	rBV3	881	666	0.01%	0.003%
28	7.985	2152	2160	2164	rVV7	1463	1904	0.03%	0.007%
29	8.088	2181	2192	2224	rBV	465635	1098126	18.53%	4.134%
30	8.600	2347	2351	2355	rBV4	648	489	0.01%	0.002%
31	8.850	2411	2429	2461	rBV	906716	1510519	25.49%	5.687%
32	9.014	2470	2480	2496	rBV	73910	120506	2.03%	0.454%
33	9.140	2510	2519	2536	rVB2	26312	46747	0.79%	0.176%
34	9.223	2542	2545	2550	rVB	767	511	0.01%	0.002%
35	9.262	2550	2557	2559	rVV4	972	1236	0.02%	0.005%
36	9.291	2563	2566	2572	rVB4	883	841	0.01%	0.003%

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072822WMA.M
 Title : VOC Analysis

37	9.320	2572	2575	2579	rBV3	683	677	0.01%	0.003%
38	9.545	2635	2645	2662	rBV2	20435	37461	0.63%	0.141%
39	9.722	2694	2700	2701	rBV3	1765	1388	0.02%	0.005%
40	9.796	2720	2723	2725	rBV2	1001	656	0.01%	0.002%
41	10.034	2792	2797	2799	rBV3	897	753	0.01%	0.003%
42	10.214	2840	2853	2878	rBV	747310	1181505	19.94%	4.448%
43	10.362	2897	2899	2905	rVB5	906	613	0.01%	0.002%
44	10.452	2919	2927	2944	rVV3	8260	16588	0.28%	0.062%
45	10.542	2946	2955	2963	rBV2	8936	13220	0.22%	0.050%
46	10.696	2996	3003	3009	rBV4	2929	4152	0.07%	0.016%
47	10.741	3012	3017	3026	rVB3	4880	6401	0.11%	0.024%
48	10.915	3059	3071	3083	rBV	25169	40118	0.68%	0.151%
49	11.021	3091	3104	3112	rBV4	7982	13235	0.22%	0.050%
50	11.136	3137	3140	3142	rBV2	751	525	0.01%	0.002%
51	11.172	3142	3151	3162	rVV	38459	65716	1.11%	0.247%
52	11.249	3163	3175	3194	rVV	944686	1479490	24.97%	5.570%
53	11.336	3195	3202	3217	rVB2	18248	31710	0.54%	0.119%
54	11.413	3223	3226	3234	rVB7	1244	1144	0.02%	0.004%
55	11.529	3248	3262	3280	rBV2	161762	272500	4.60%	1.026%
56	11.622	3280	3291	3318	rVB	1009869	1585209	26.75%	5.968%
57	11.744	3319	3329	3346	rBV	179330	281787	4.76%	1.061%
58	11.924	3375	3385	3399	rBV10	5793	12963	0.22%	0.049%
59	12.017	3407	3414	3418	rBV3	6586	8377	0.14%	0.032%
60	12.114	3434	3444	3457	rBV2	18906	30283	0.51%	0.114%
61	12.162	3457	3459	3462	rVB4	1340	677	0.01%	0.003%
62	12.252	3482	3487	3492	rBV7	1942	2352	0.04%	0.009%
63	12.291	3496	3499	3500	rBV3	1055	579	0.01%	0.002%
64	12.361	3512	3521	3531	rBV5	17582	33883	0.57%	0.128%
65	12.468	3544	3554	3566	rBV2	31004	58857	0.99%	0.222%
66	12.725	3623	3634	3647	rVV	33659	52889	0.89%	0.199%
67	12.789	3647	3654	3662	rVB7	4223	6430	0.11%	0.024%
68	12.882	3671	3683	3694	rBV2	62372	97792	1.65%	0.368%
69	12.950	3698	3704	3713	rVB5	7765	11693	0.20%	0.044%
70	13.046	3723	3734	3748	rBV2	156321	243971	4.12%	0.919%
71	13.159	3760	3769	3779	rVB9	9454	16462	0.28%	0.062%
72	13.207	3779	3784	3785	rBV3	3506	2335	0.04%	0.009%
73	13.268	3796	3803	3811	rBV10	4941	7641	0.13%	0.029%
74	13.378	3829	3837	3854	rVB10	6177	14302	0.24%	0.054%
75	13.500	3858	3875	3902	rBV	4133877	5925475	100.00%	22.308%
76	13.622	3903	3913	3927	rVV	164018	261518	4.41%	0.985%

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 Title : VOC Analysis

77	13.767	3953	3958	3968	rVB8	4392	6890	0.12%	0.026%
78	13.818	3970	3974	3980	rVB7	922	982	0.02%	0.004%
79	13.905	3996	4001	4018	rVB4	5748	9539	0.16%	0.036%
80	13.972	4018	4022	4026	rVB3	795	707	0.01%	0.003%
81	14.008	4027	4033	4036	rBV4	983	980	0.02%	0.004%
82	14.095	4049	4060	4071	rBV7	7674	17553	0.30%	0.066%
83	14.149	4073	4077	4081	rVV5	1031	975	0.02%	0.004%
84	14.188	4085	4089	4092	rBV4	748	678	0.01%	0.003%
85	14.294	4114	4122	4133	rVB8	6789	11336	0.19%	0.043%
86	14.358	4139	4142	4144	rBV3	784	540	0.01%	0.002%
87	14.390	4149	4152	4156	rBV4	805	489	0.01%	0.002%
88	14.429	4158	4164	4176	rBV10	4851	8140	0.14%	0.031%
89	14.503	4185	4187	4192	rVB3	994	561	0.01%	0.002%
90	14.577	4200	4210	4216	rBV2	16686	24997	0.42%	0.094%
91	14.631	4216	4227	4244	rVV	456865	718356	12.12%	2.704%
92	14.750	4256	4264	4272	rVB3	11396	18477	0.31%	0.070%
93	14.815	4272	4284	4299	rBV	385163	619272	10.45%	2.331%
94	15.361	4446	4454	4467	rVB	34147	55091	0.93%	0.207%
95	15.461	4483	4485	4488	rBV3	837	580	0.01%	0.002%
96	15.554	4507	4514	4520	rBV	12544	16174	0.27%	0.061%
97	15.667	4540	4549	4569	rVB5	20327	41683	0.70%	0.157%
98	15.808	4585	4593	4601	rBV2	33905	50979	0.86%	0.192%
99	16.011	4651	4656	4657	rBV4	3137	2691	0.05%	0.010%
100	16.483	4793	4803	4821	rBV	68504	108154	1.83%	0.407%

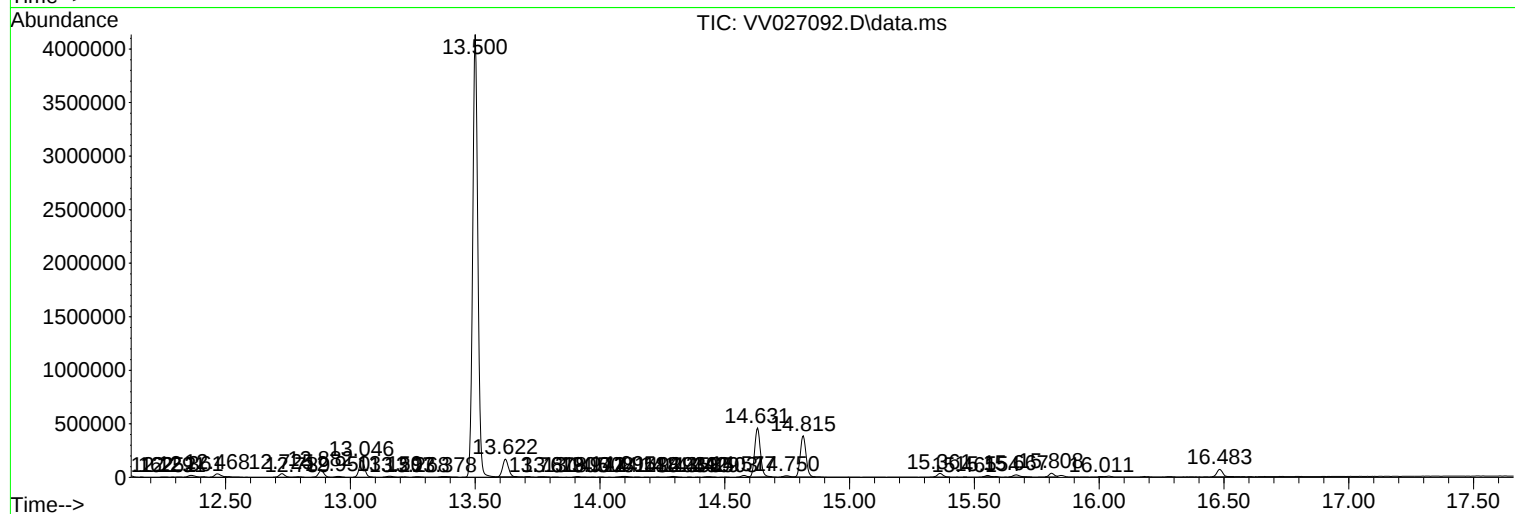
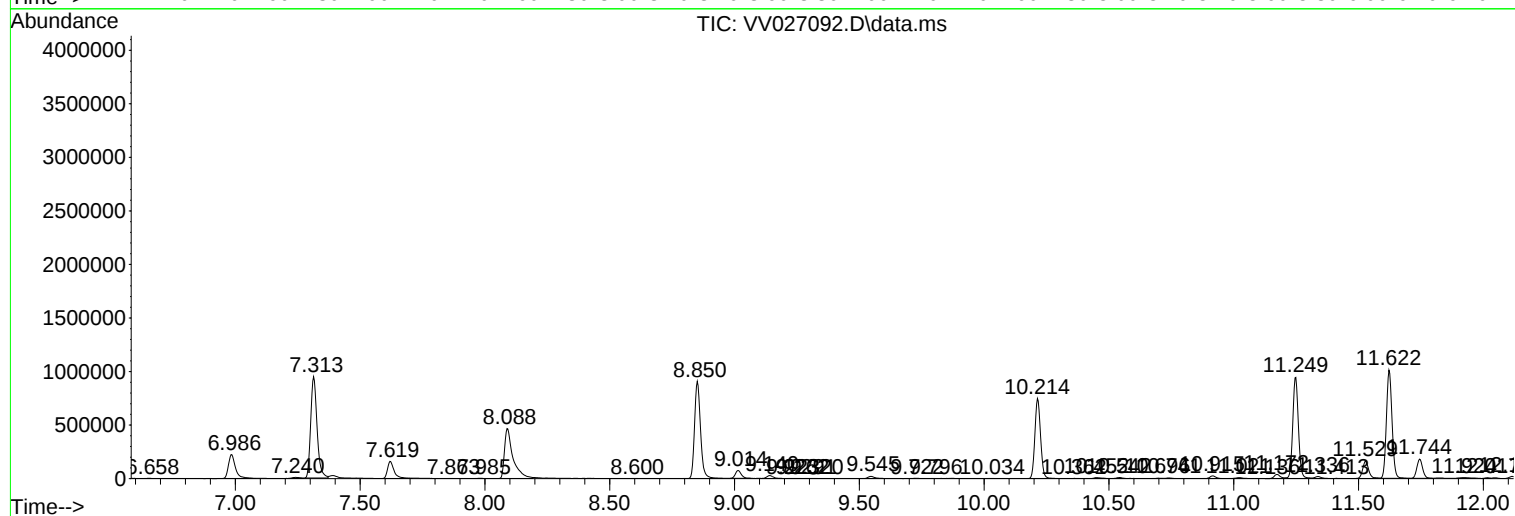
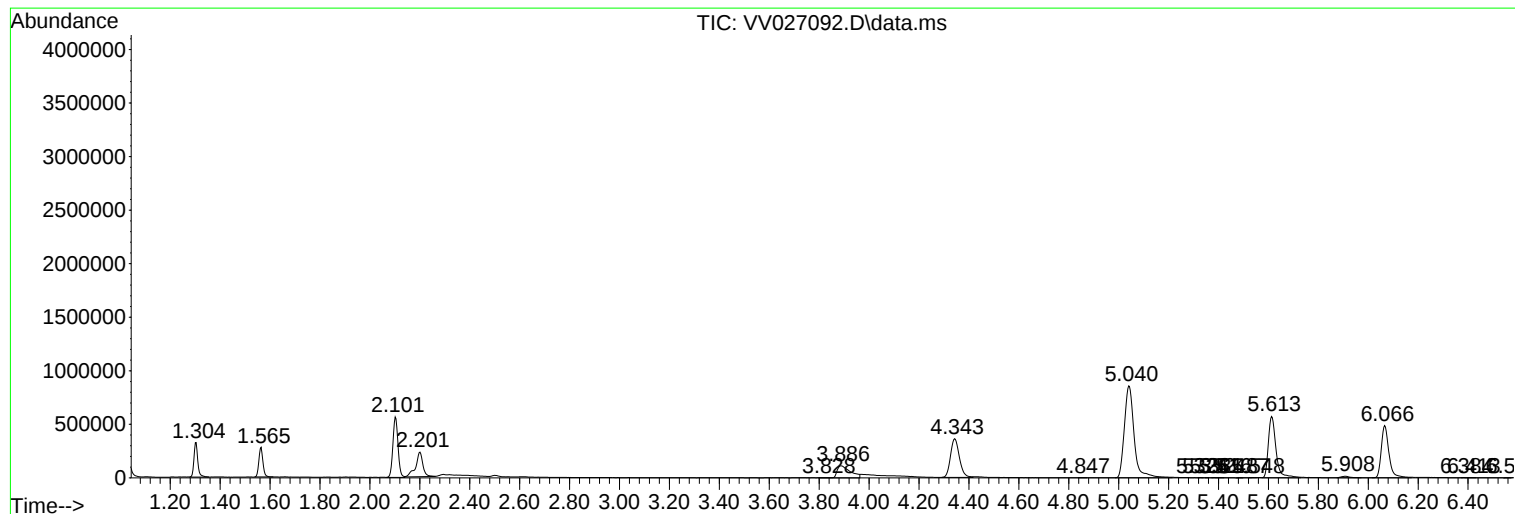
Sum of corrected areas: 26561767

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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072822WMA.M
Quant Title : VOC Analysis

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



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Instrument :
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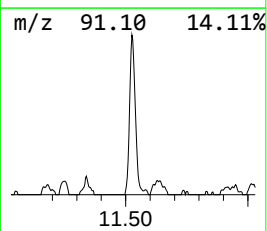
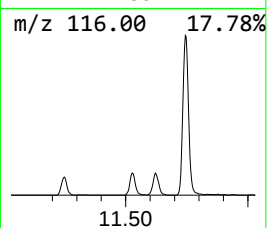
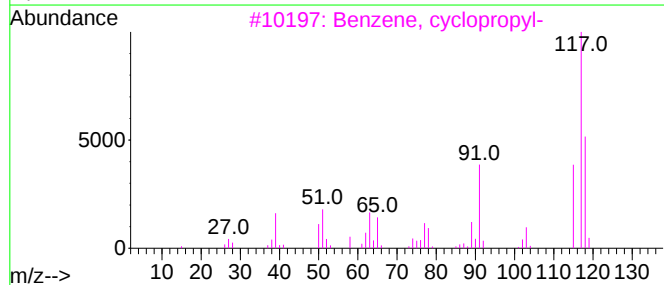
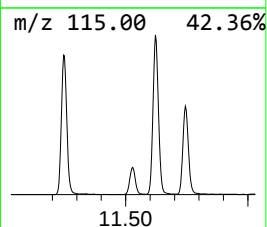
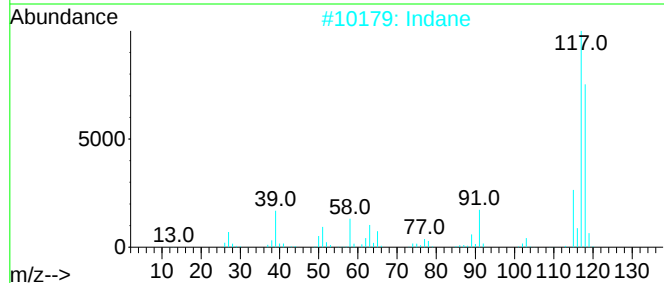
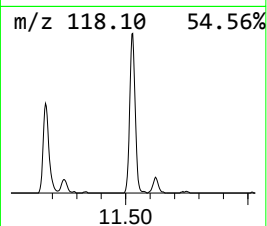
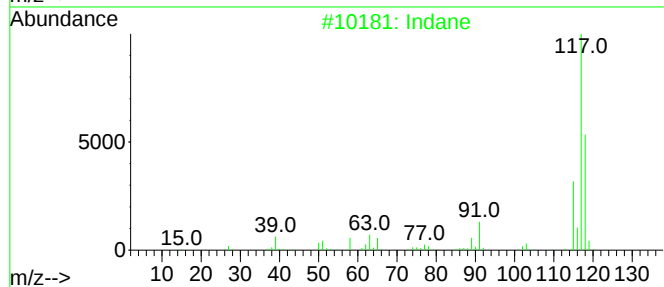
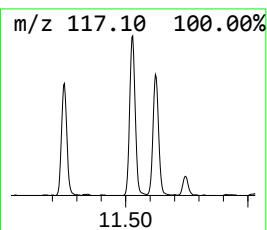
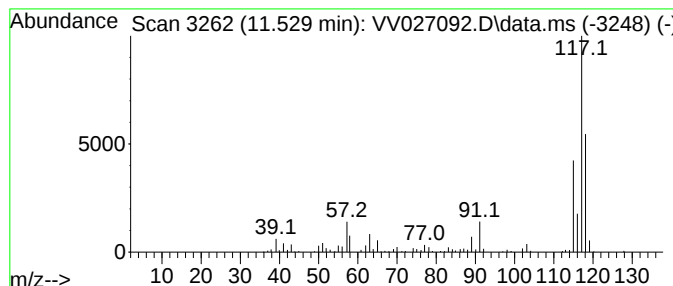
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072822WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	9.21 ug/L	272500	1,4-Dichlorobenzene-d4	11.249

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indane	118	C9H10	000496-11-7	95
2		Indane	118	C9H10	000496-11-7	70
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	58
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	58



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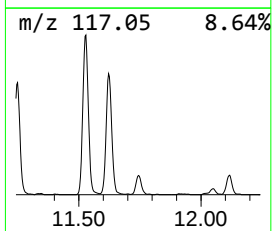
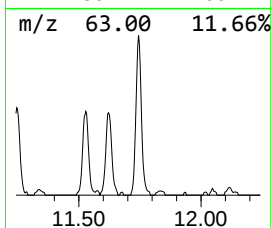
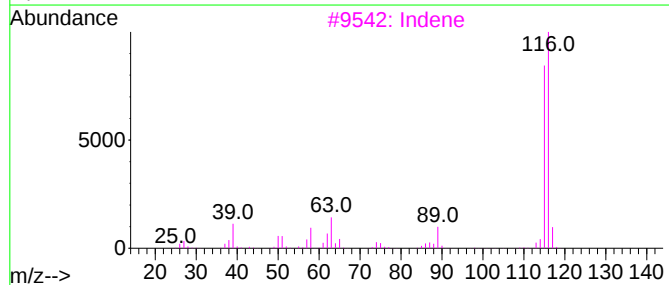
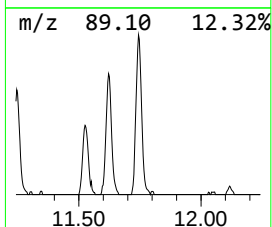
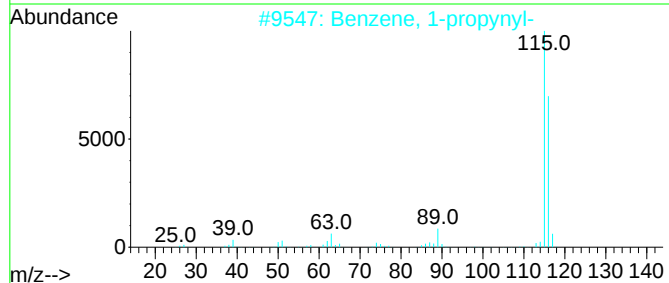
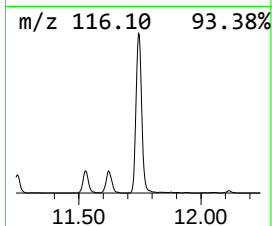
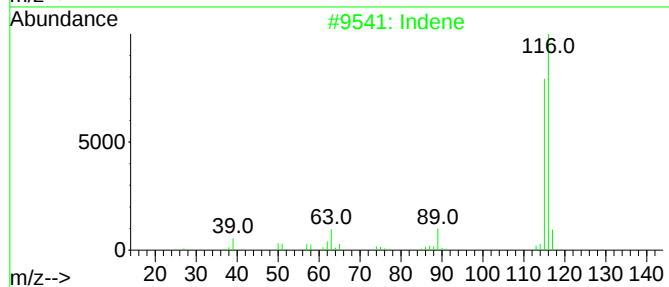
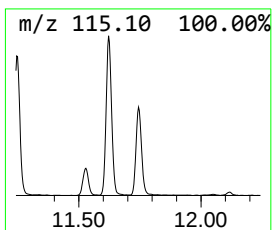
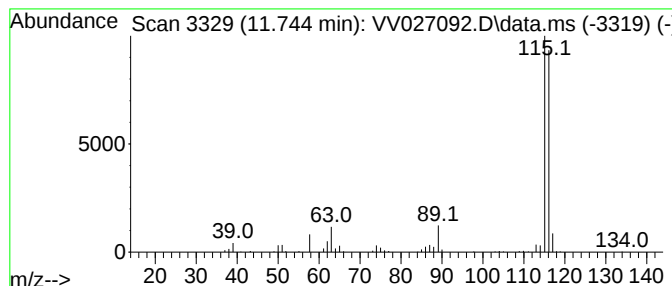
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072822WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Indene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.744	9.52 ug/L	281787	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3			Indene	116	C9H8	000095-13-6	94
4			Indene	116	C9H8	000095-13-6	93
5			3-Methylphenylacetylene	116	C9H8	000766-82-5	91



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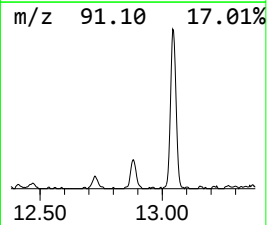
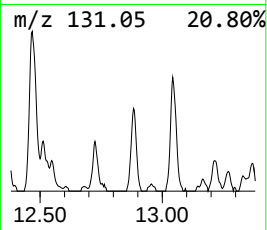
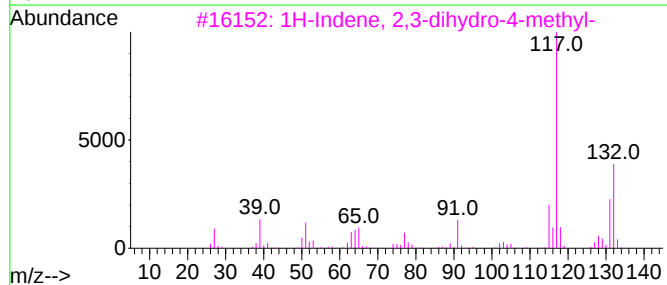
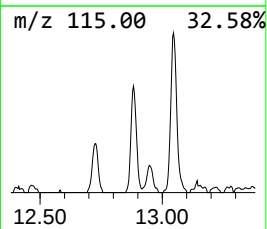
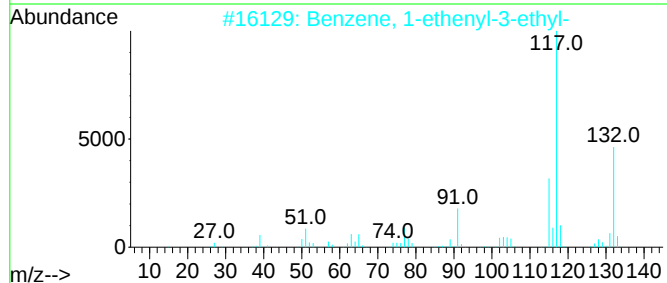
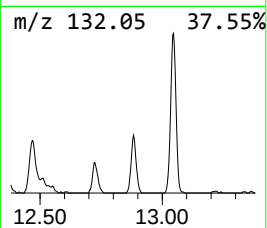
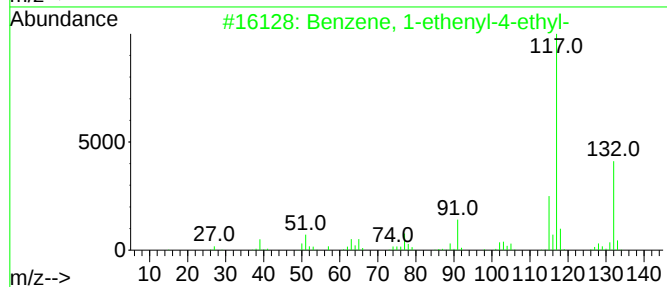
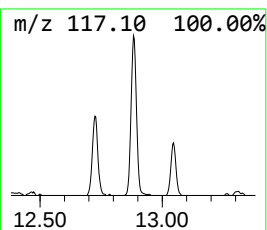
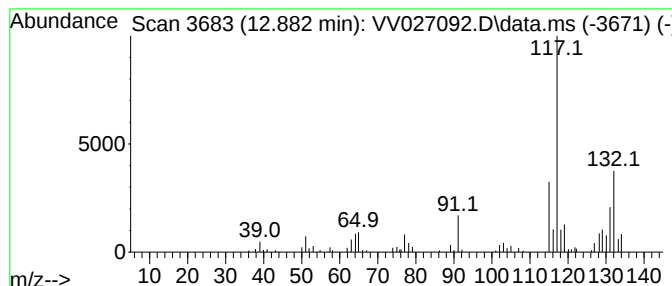
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Quant Title : VOC Analysis

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

Peak Number 4 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.882	3.30 ug/L	97792	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072922\
Data File : VV072092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

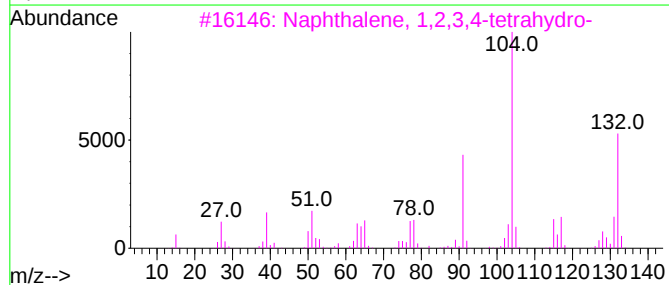
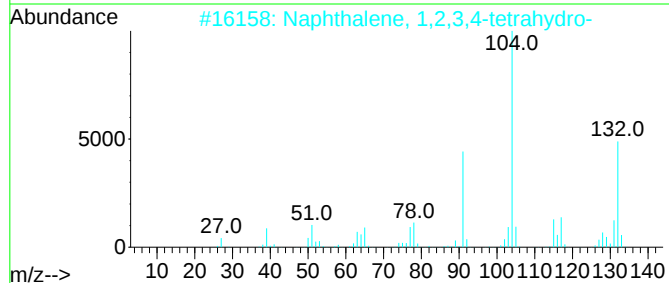
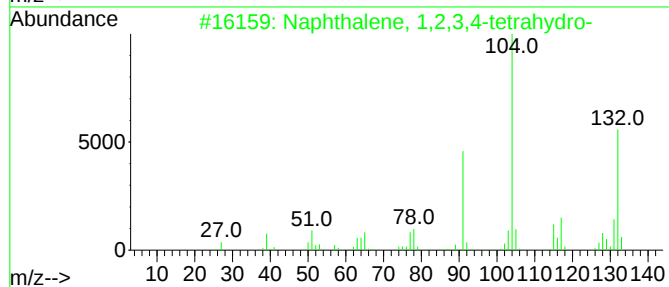
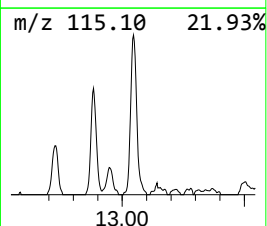
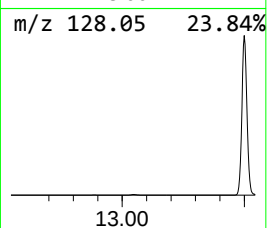
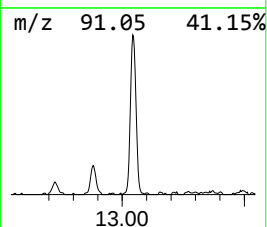
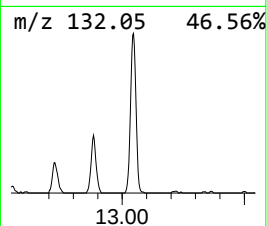
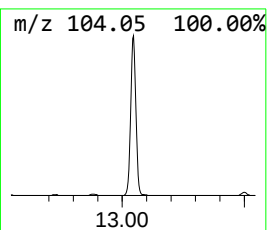
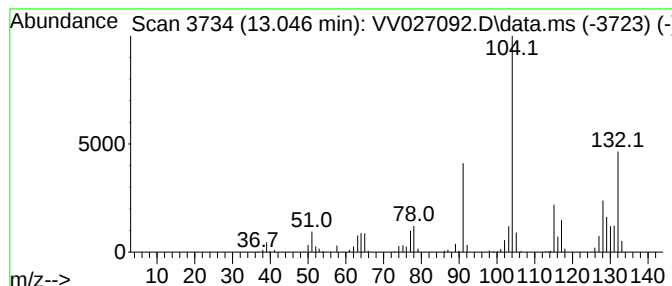
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Quant Title : VOC Analysis

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.046	8.25 ug/L	243971	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072922\
Data File : VV027092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

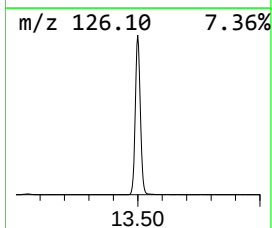
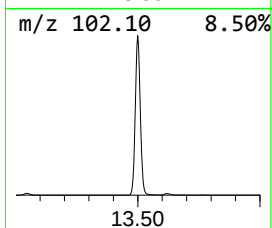
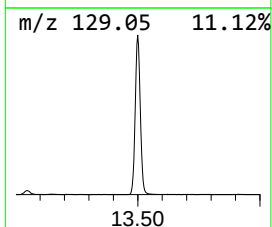
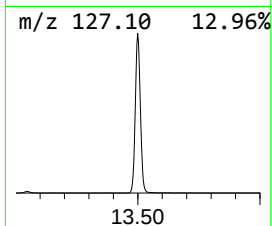
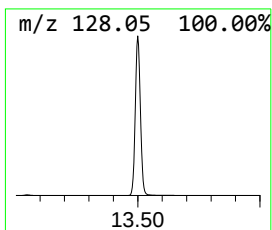
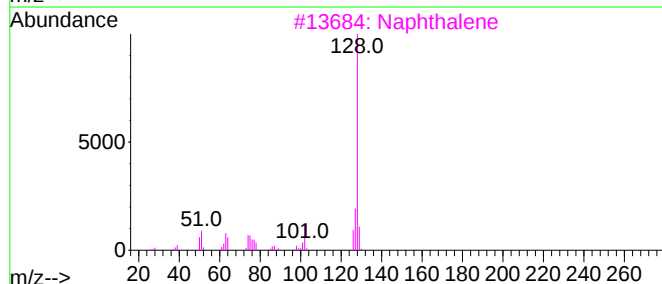
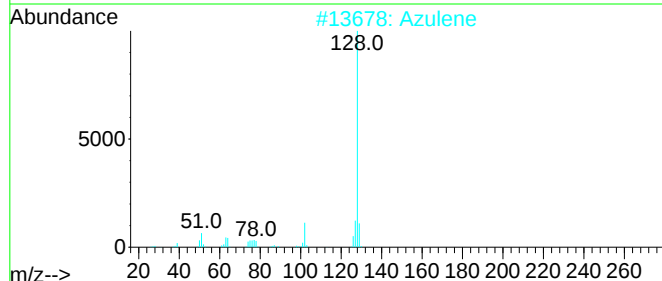
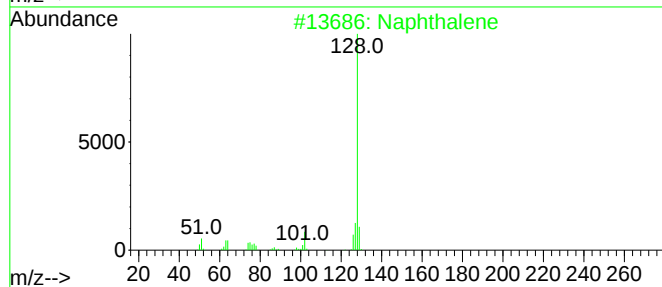
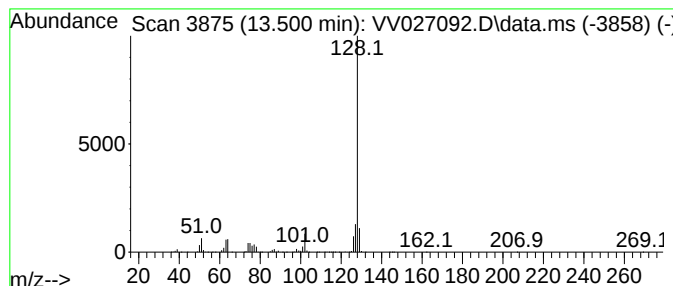
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Quant Title : VOC Analysis

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

Peak Number 6 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	200.25 ug/L	5925480	1,4-Dichlorobenzene-d4	11.249

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072922\
Data File : VV027092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

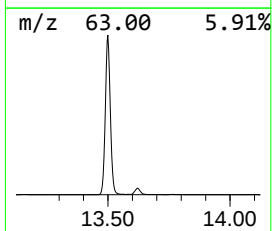
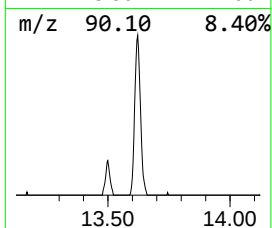
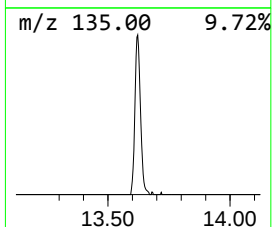
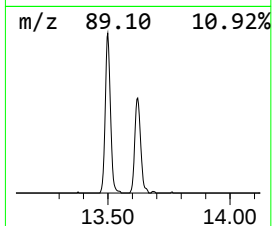
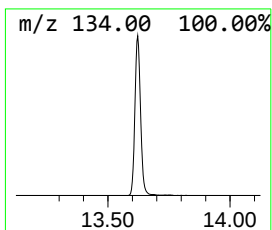
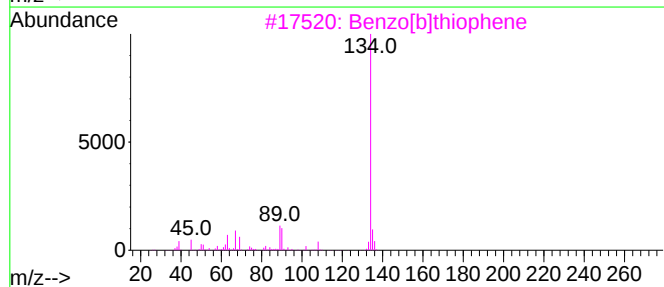
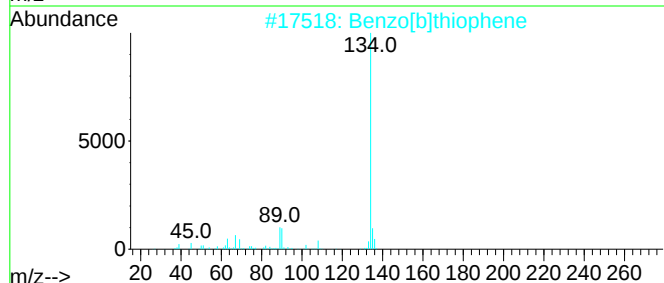
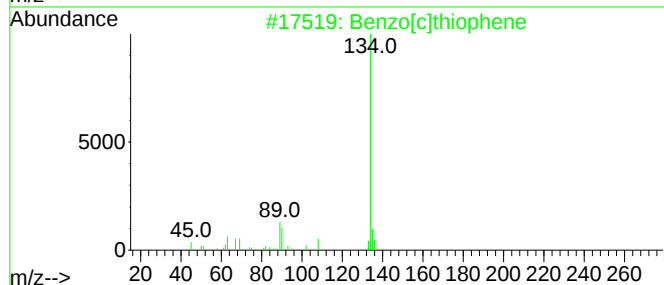
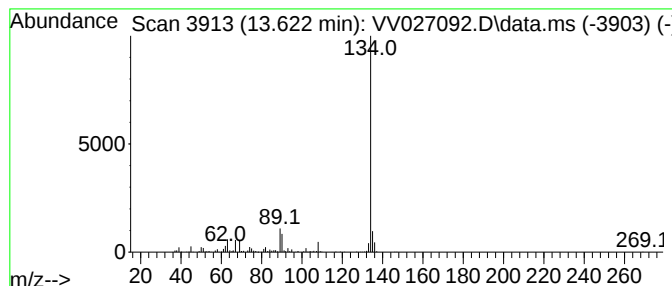
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Quant Title : VOC Analysis

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

Peak Number 7 Benzo[c]thiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	8.84 ug/L	261518	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2			Benzo[b]thiophene	134	C8H6S	000095-15-8	95
3			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
4			Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072922\
Data File : VV072092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

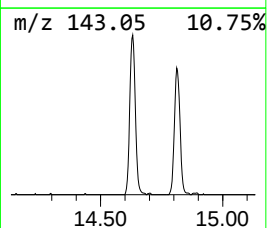
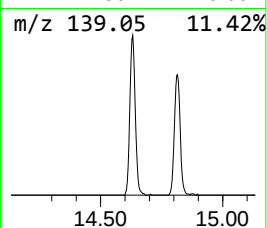
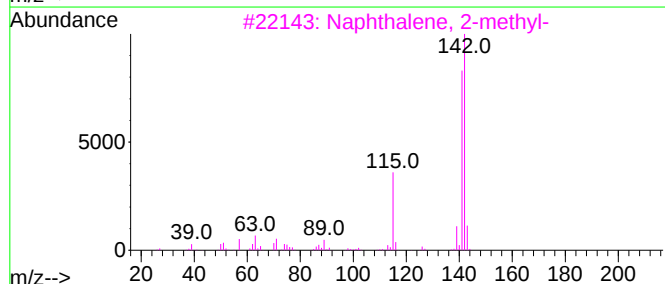
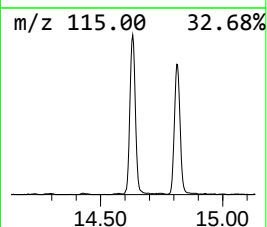
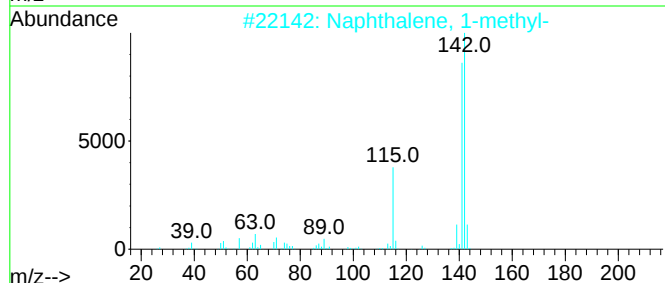
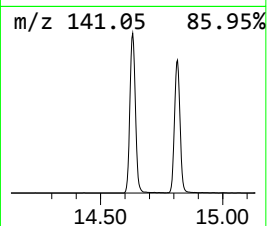
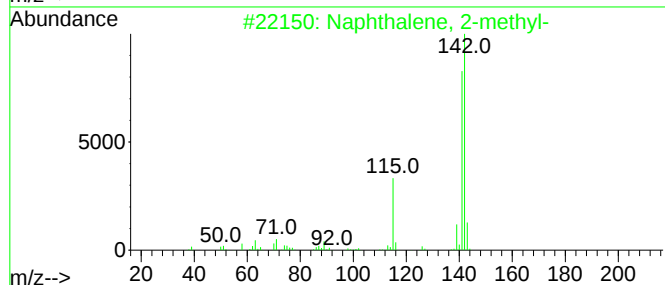
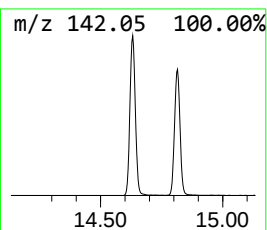
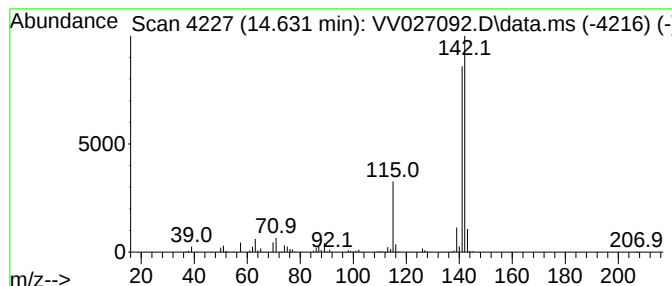
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Quant Title : VOC Analysis

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

Peak Number 8 Naphthalene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.632	24.28 ug/L	718356	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
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_V\Data\VV072922\
Data File : VV072092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

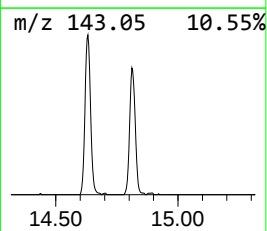
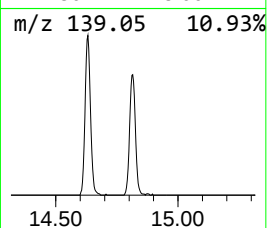
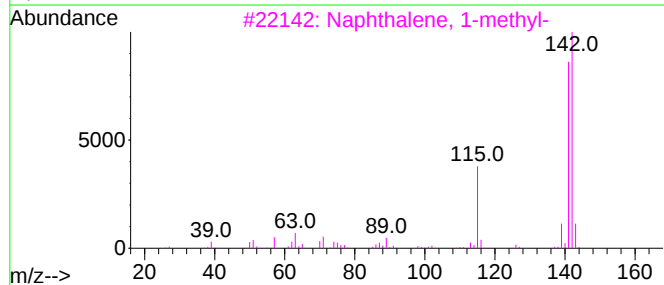
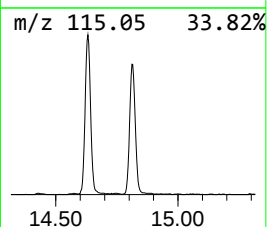
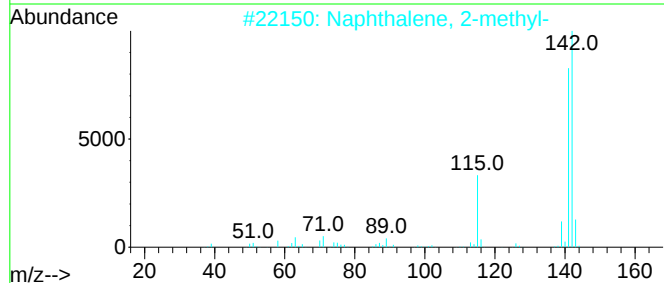
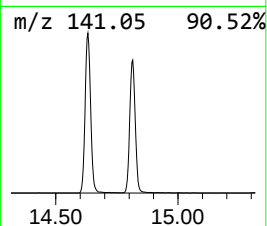
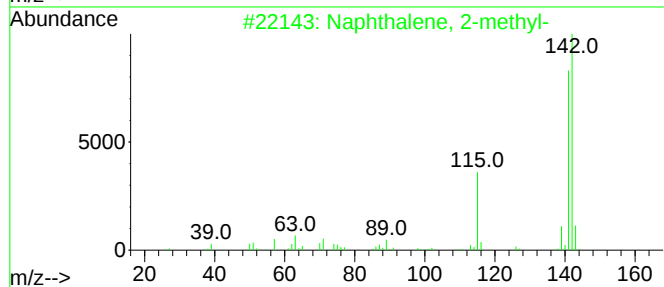
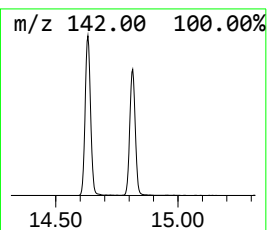
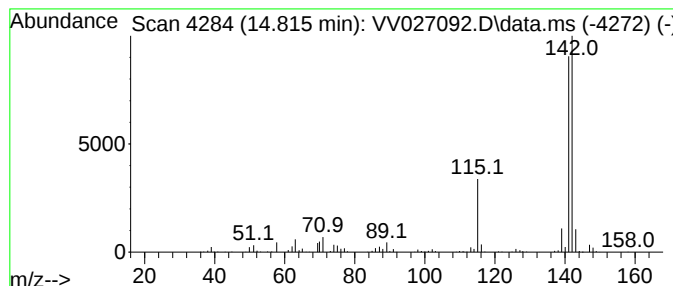
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Quant Title : VOC Analysis

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.815	20.93 ug/L	619272	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
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_V\Data\VV072922\
Data File : VV027092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

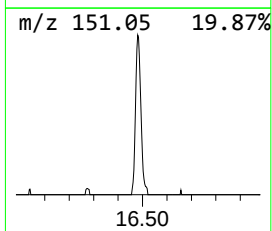
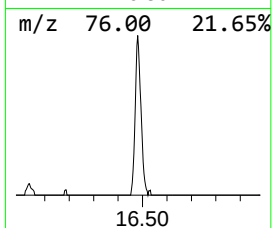
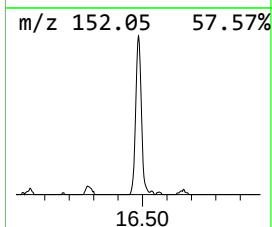
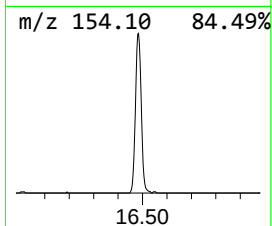
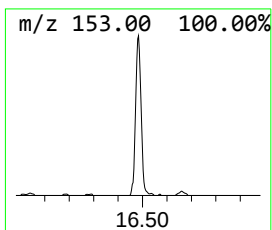
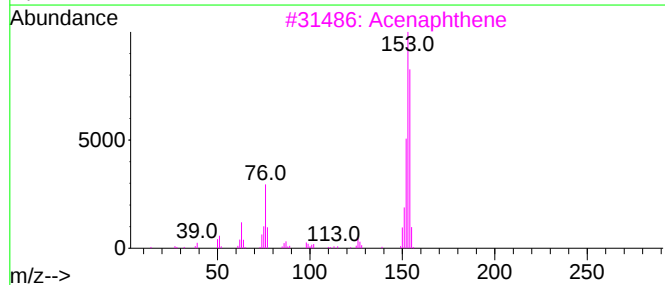
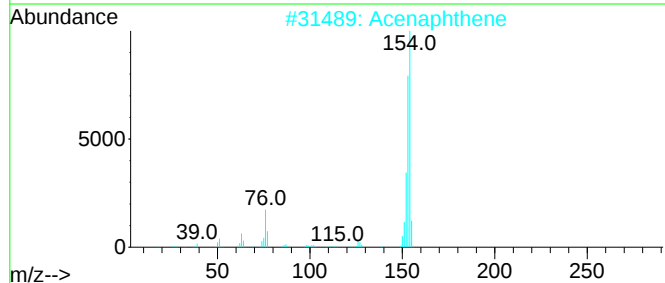
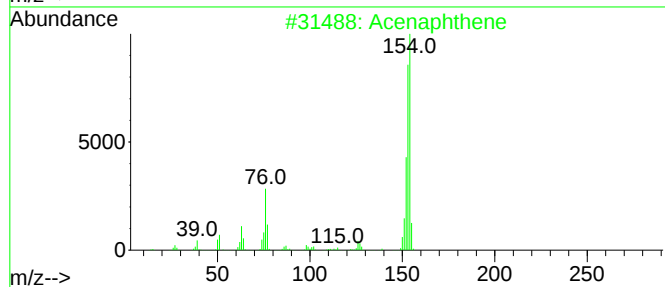
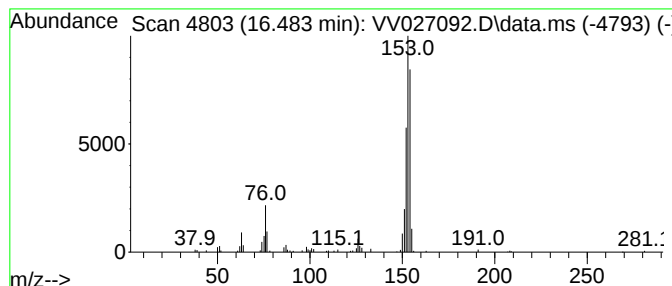
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Quant Title : VOC Analysis

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

Peak Number 10 Acenaphthene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.483	3.66 ug/L	108154	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	81
2			Acenaphthene	154	C12H10	000083-32-9	81
3			Acenaphthene	154	C12H10	000083-32-9	76
4			Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	58
5			1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072922\
Data File : VV027092.D
Acq On : 29 Jul 2022 18:09
Operator : SY/MD
Sample : N3899-04 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUR3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072822WMA.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
Indane	11.529	9.2	ug/L	272500	3	11.249	1479490	50.0
Indene	11.744	9.5	ug/L	281787	3	11.249	1479490	50.0
Benzene, 1-ethe...	12.882	3.3	ug/L	97792	3	11.249	1479490	50.0
Naphthalene, 1,...	13.046	8.3	ug/L	243971	3	11.249	1479490	50.0
Azulene	13.500	200.3	ug/L	5925480	3	11.249	1479490	50.0
Benzo[c]thiophene	13.622	8.8	ug/L	261518	3	11.249	1479490	50.0
Naphthalene, 2-...	14.632	24.3	ug/L	718356	3	11.249	1479490	50.0
Naphthalene, 1-...	14.815	20.9	ug/L	619272	3	11.249	1479490	50.0
Acenaphthene	16.483	3.7	ug/L	108154	3	11.249	1479490	50.0