

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
 Data File : VV027003.D
 Acq On : 27 Jul 2022 16:58
 Operator : SY/MD
 Sample : N3843-07DL 10X
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBUE7DL

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\SFAMVLM072222WMA.M

Title : VOC Analysis

Signal : TIC: VV027003.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.301	74	81	97	rVB	224969	238034	5.80%	0.973%
2	1.378	100	105	125	rVB	32390	43412	1.06%	0.177%
3	1.564	155	163	175	rBV	193522	224756	5.47%	0.918%
4	2.101	320	330	342	rBV	406171	572250	13.94%	2.339%
5	2.172	342	352	381	rVB	233648	548017	13.35%	2.240%
6	2.957	585	596	611	rBV3	5272	12751	0.31%	0.052%
7	3.458	746	752	757	rBV2	575	589	0.01%	0.002%
8	3.593	787	794	800	rBV2	418	696	0.02%	0.003%
9	3.732	826	837	842	rVV5	1643	3200	0.08%	0.013%
10	3.767	846	848	852	rVV2	751	667	0.02%	0.003%
11	3.883	873	884	905	rBV	177968	534666	13.02%	2.185%
12	4.342	1007	1027	1056	rBV	326805	831938	20.26%	3.400%
13	4.564	1094	1096	1100	rBV4	853	539	0.01%	0.002%
14	4.683	1128	1133	1135	rBV4	539	497	0.01%	0.002%
15	5.040	1227	1244	1285	rBV2	721894	1910208	46.53%	7.806%
16	5.336	1325	1336	1337	rBV5	2623	3384	0.08%	0.014%
17	5.535	1387	1398	1410	rBV2	11941	26940	0.66%	0.110%
18	5.613	1410	1422	1458	rVB	695091	1416106	34.49%	5.787%
19	5.905	1503	1513	1528	rBV3	11933	27344	0.67%	0.112%
20	6.066	1548	1563	1592	rBV	429956	892113	21.73%	3.646%
21	6.198	1595	1604	1627	rVB2	19659	51877	1.26%	0.212%
22	6.362	1649	1655	1666	rVB4	2900	5573	0.14%	0.023%
23	6.580	1713	1723	1736	rBV6	10646	28134	0.69%	0.115%
24	6.982	1837	1848	1881	rBV	206530	390758	9.52%	1.597%
25	7.236	1918	1927	1940	rBV3	17529	36199	0.88%	0.148%
26	7.313	1940	1951	1970	rBV	771614	1340005	32.64%	5.476%
27	7.619	2036	2046	2067	rBV	143495	255523	6.22%	1.044%
28	7.850	2115	2118	2120	rBV2	829	486	0.01%	0.002%
29	7.944	2139	2147	2149	rBV4	1893	2667	0.06%	0.011%
30	7.979	2149	2158	2168	rVV4	11349	19848	0.48%	0.081%
31	8.088	2180	2192	2234	rBV2	650323	1251945	30.49%	5.116%
32	8.728	2389	2391	2396	rVB5	1340	1137	0.03%	0.005%
33	8.850	2413	2429	2463	rBV	1057434	1750452	42.64%	7.153%
34	9.014	2470	2480	2498	rBV	53196	93625	2.28%	0.383%
35	9.143	2513	2520	2534	rVB2	16475	29932	0.73%	0.122%
36	9.252	2548	2554	2562	rVV5	3536	5936	0.14%	0.024%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072222WMA.M
 Title : VOC Analysis

37	9.397	2590	2599	2608	rBV7	4310	8098	0.20%	0.033%
38	9.436	2608	2611	2614	rVB4	1054	769	0.02%	0.003%
39	9.506	2627	2633	2637	rBV5	1924	2372	0.06%	0.010%
40	9.548	2638	2646	2659	rBV4	12218	24470	0.60%	0.100%
41	9.673	2678	2685	2687	rBV3	732	808	0.02%	0.003%
42	9.712	2688	2697	2706	rBV3	13416	21816	0.53%	0.089%
43	9.815	2717	2729	2738	rBV5	12974	28650	0.70%	0.117%
44	10.021	2783	2793	2807	rVV3	12777	23996	0.58%	0.098%
45	10.098	2811	2817	2819	rBV3	740	626	0.02%	0.003%
46	10.214	2839	2853	2875	rBV	749490	1160851	28.28%	4.744%
47	10.336	2889	2891	2893	rVV3	859	496	0.01%	0.002%
48	10.355	2893	2897	2903	rVV6	2457	3160	0.08%	0.013%
49	10.452	2916	2927	2941	rBV9	3566	7770	0.19%	0.032%
50	10.525	2947	2950	2951	rBV3	1352	664	0.02%	0.003%
51	10.693	2993	3002	3012	rBV4	5505	9532	0.23%	0.039%
52	10.776	3025	3028	3033	rBV4	704	692	0.02%	0.003%
53	10.828	3034	3044	3049	rBV4	4439	7015	0.17%	0.029%
54	10.918	3063	3072	3091	rBV4	11454	22804	0.56%	0.093%
55	11.021	3097	3104	3122	rVB2	16482	30756	0.75%	0.126%
56	11.172	3142	3151	3161	rBV	53092	87111	2.12%	0.356%
57	11.249	3163	3175	3194	rVB	1158589	1800389	43.85%	7.357%
58	11.490	3243	3250	3252	rBV3	3705	3996	0.10%	0.016%
59	11.529	3252	3262	3279	rVV2	185845	318928	7.77%	1.303%
60	11.622	3280	3291	3317	rVB	972075	1523783	37.12%	6.227%
61	11.744	3318	3329	3346	rBV	433275	665429	16.21%	2.719%
62	11.921	3376	3384	3396	rVB2	29851	50195	1.22%	0.205%
63	12.014	3410	3413	3417	rBV6	1965	1933	0.05%	0.008%
64	12.114	3437	3444	3451	rBV4	6827	9257	0.23%	0.038%
65	12.159	3456	3458	3464	rVB6	1070	653	0.02%	0.003%
66	12.197	3467	3470	3473	rVB4	771	524	0.01%	0.002%
67	12.239	3473	3483	3484	rBV5	2682	2449	0.06%	0.010%
68	12.355	3511	3519	3534	rVB6	19098	39564	0.96%	0.162%
69	12.468	3545	3554	3564	rBV3	27530	52254	1.27%	0.214%
70	12.683	3617	3621	3625	rVV3	812	812	0.02%	0.003%
71	12.725	3625	3634	3644	rVV4	26803	43747	1.07%	0.179%
72	12.786	3646	3653	3664	rVB5	12141	20256	0.49%	0.083%
73	12.882	3674	3683	3694	rBV5	26722	41960	1.02%	0.171%
74	12.950	3694	3704	3714	rVB7	9925	15795	0.38%	0.065%
75	13.046	3723	3734	3749	rBV3	69169	109644	2.67%	0.448%
76	13.156	3759	3768	3777	rBV7	10879	16796	0.41%	0.069%

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Integration Parameters: LSCINT.P

Integrator: RTE

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 Title : VOC Analysis

77	13.210	3783	3785	3792	rVB5	1383	1312	0.03%	0.005%
78	13.265	3794	3802	3812	rBV8	4054	5893	0.14%	0.024%
79	13.371	3826	3835	3836	rBV4	8828	10606	0.26%	0.043%
80	13.500	3857	3875	3902	rBV	2702484	4105520	100.00%	16.777%
81	13.622	3902	3913	3928	rVB	123026	201891	4.92%	0.825%
82	13.696	3934	3936	3937	rBV2	1287	591	0.01%	0.002%
83	13.812	3968	3972	3974	rBV3	2208	1857	0.05%	0.008%
84	13.908	3996	4002	4008	rBV6	2056	2656	0.06%	0.011%
85	14.104	4054	4063	4071	rVB9	2144	3625	0.09%	0.015%
86	14.139	4071	4074	4080	rBV6	970	1028	0.03%	0.004%
87	14.291	4114	4121	4130	rBV9	4528	6899	0.17%	0.028%
88	14.429	4159	4164	4166	rBV4	1556	1368	0.03%	0.006%
89	14.577	4197	4210	4216	rBV	7790	14026	0.34%	0.057%
90	14.631	4216	4227	4245	rVV	320672	506165	12.33%	2.068%
91	14.747	4253	4263	4273	rBV2	10626	18805	0.46%	0.077%
92	14.815	4273	4284	4300	rVV	159736	261722	6.37%	1.070%
93	15.365	4444	4455	4466	rVB	55830	86343	2.10%	0.353%
94	15.554	4505	4514	4520	rBV2	10696	16318	0.40%	0.067%
95	15.667	4540	4549	4558	rBV3	23140	39722	0.97%	0.162%
96	15.811	4586	4594	4600	rBV2	29173	44420	1.08%	0.182%
97	16.274	4729	4738	4752	rBV2	29024	51645	1.26%	0.211%
98	16.483	4792	4803	4818	rBV	161247	258471	6.30%	1.056%
99	16.786	4889	4897	4907	rVB	49238	76714	1.87%	0.313%
100	17.387	5079	5084	5098	rVB2	26813	39855	0.97%	0.163%

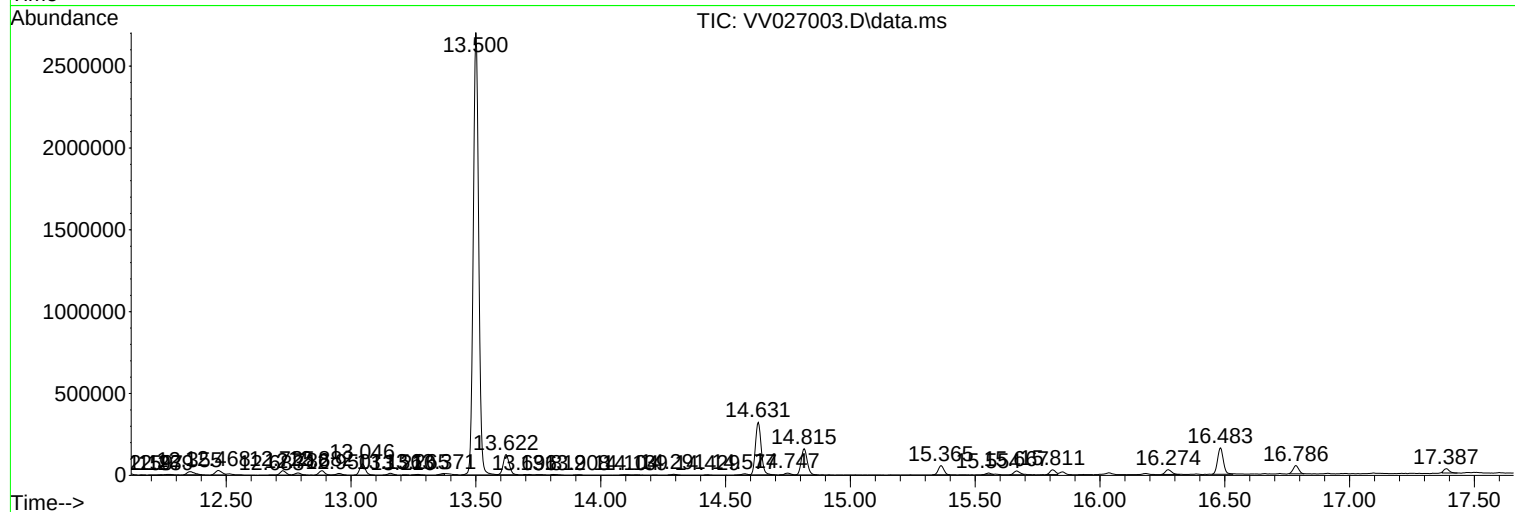
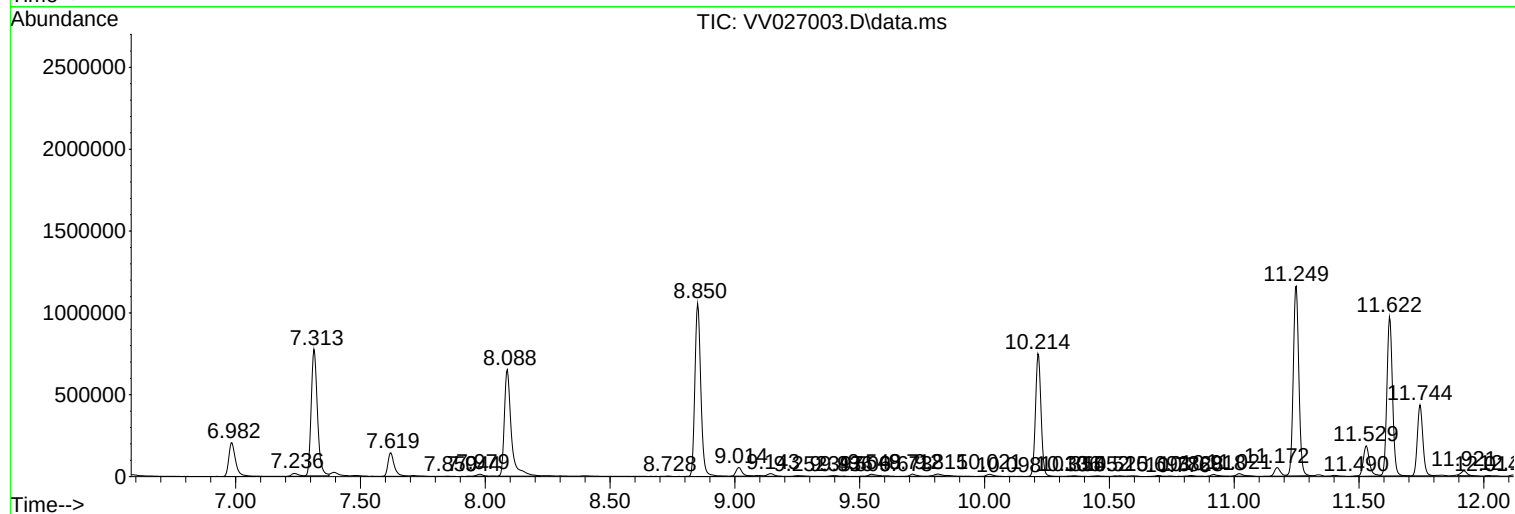
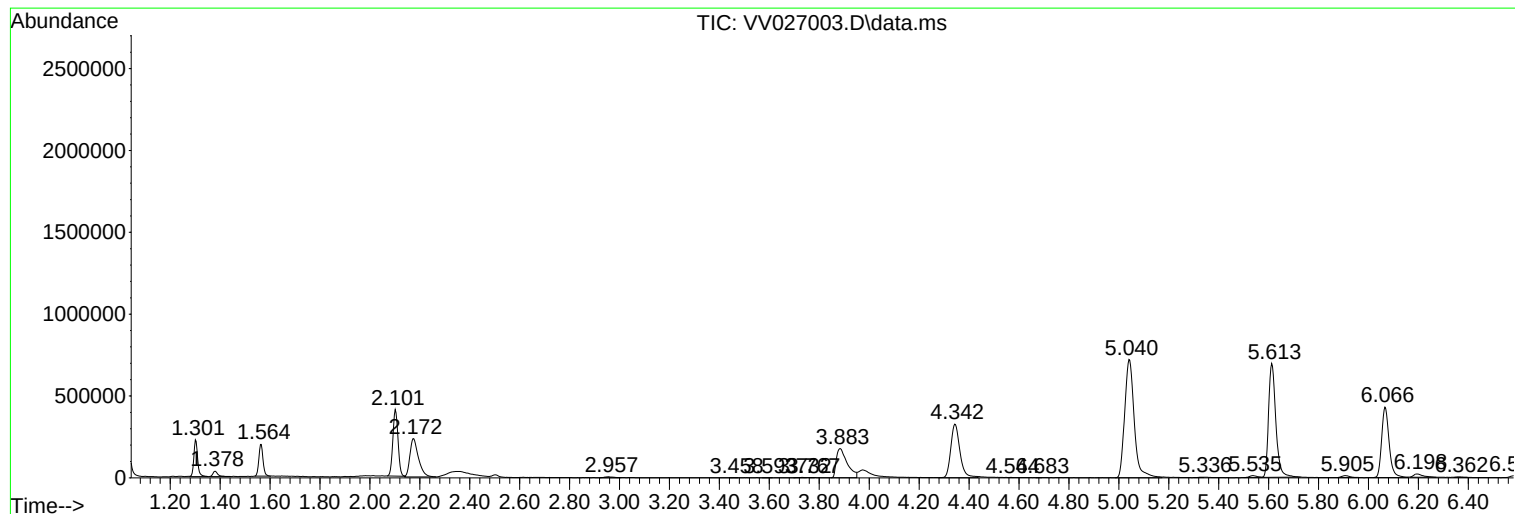
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072222WMA.M
Quant Title : VOC Analysis

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



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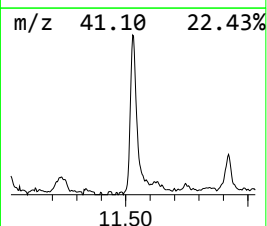
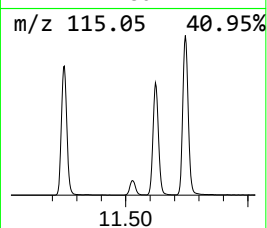
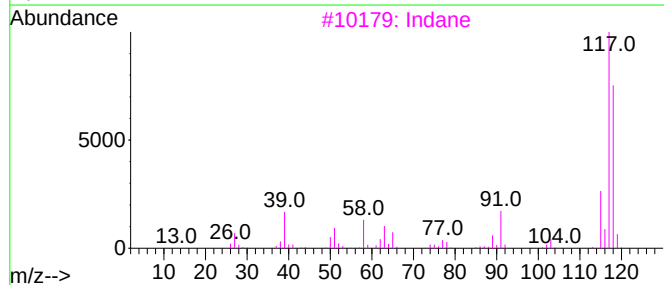
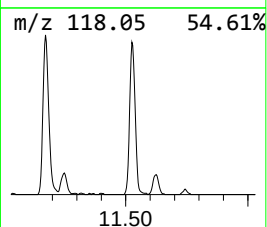
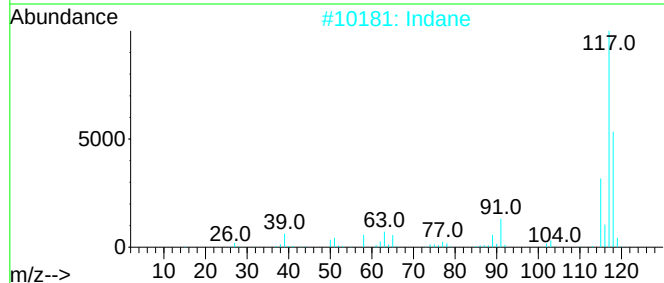
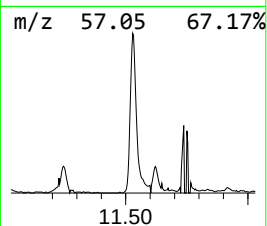
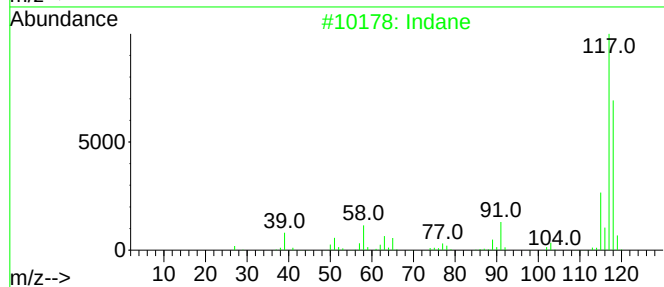
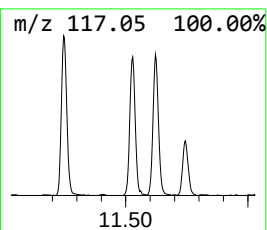
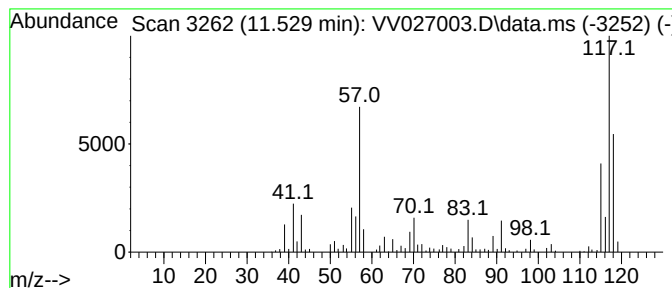
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072222WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	8.86 ug/L	318928	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	76
3	Indane			118	C9H10	000496-11-7	70
4	Benzene, cyclopropyl-			118	C9H10	000873-49-4	52
5	Benzene, 2-propenyl-			118	C9H10	000300-57-2	50



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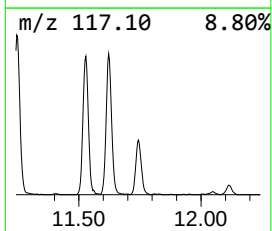
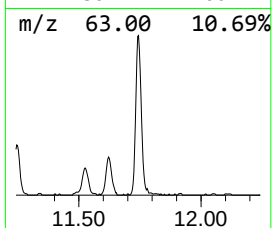
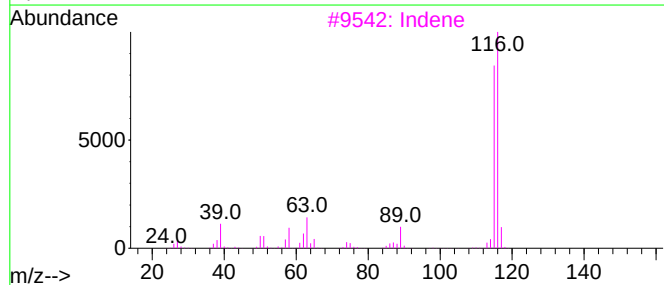
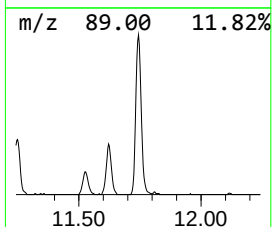
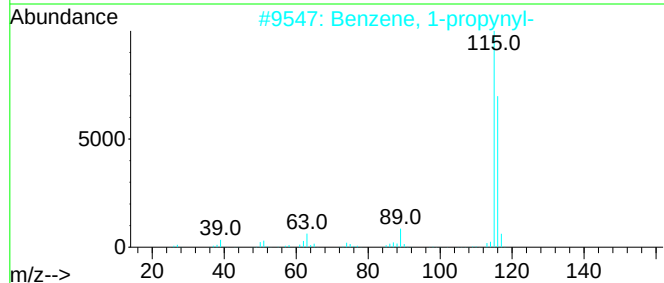
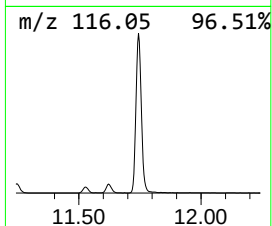
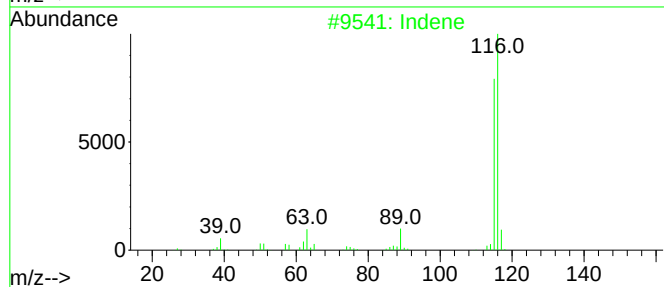
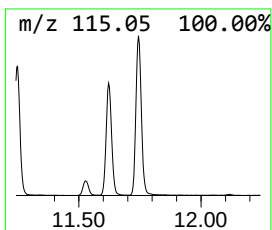
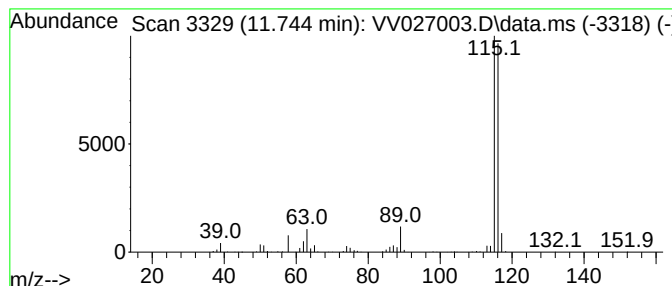
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072222WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.744	18.48 ug/L	665429	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			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91



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Instrument :
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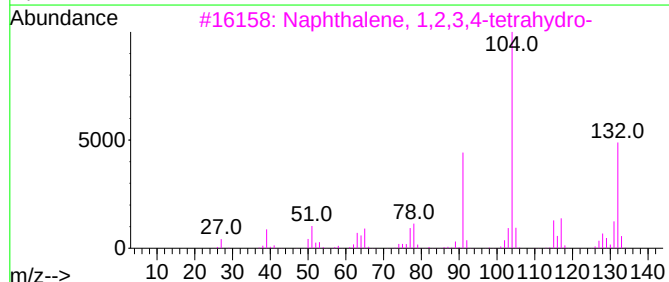
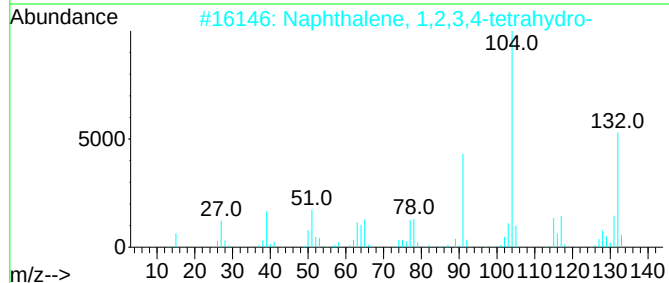
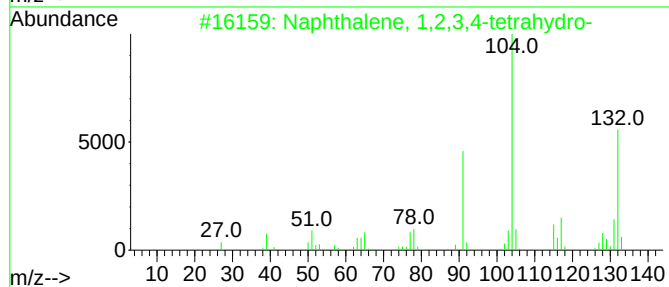
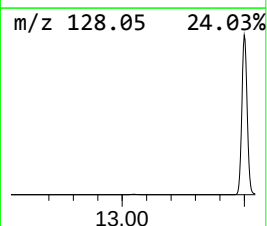
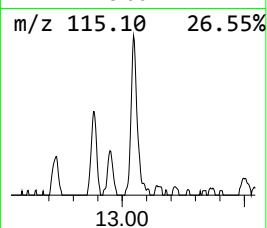
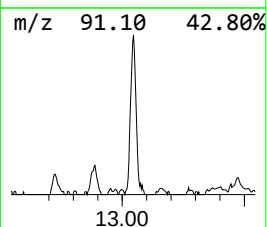
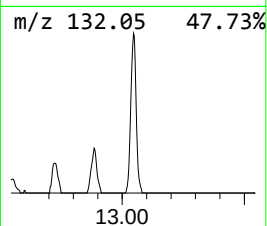
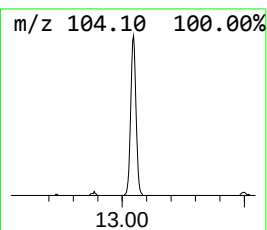
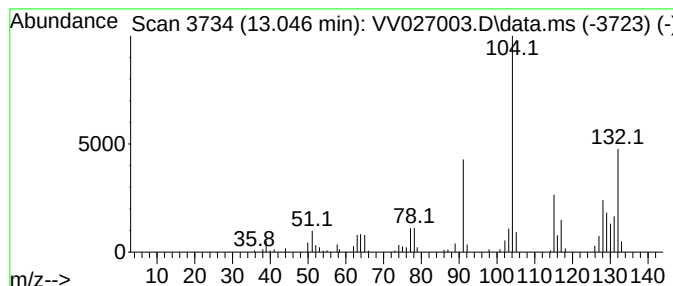
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072222WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.046	3.05 ug/L	109644	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	94
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	94
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
Data File : VV027003.D
Acq On : 27 Jul 2022 16:58
Operator : SY/MD
Sample : N3843-07DL 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUE7DL

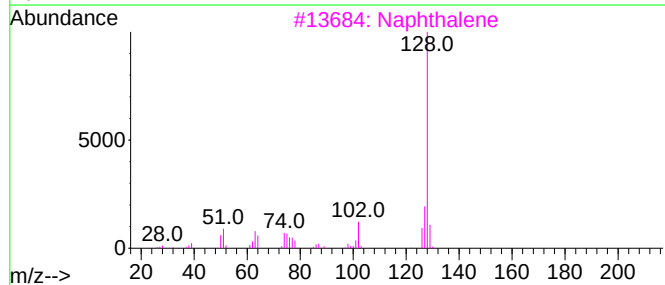
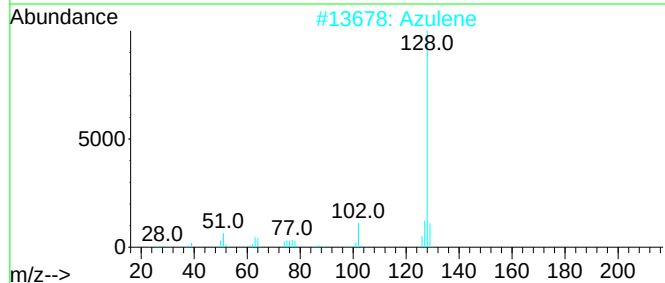
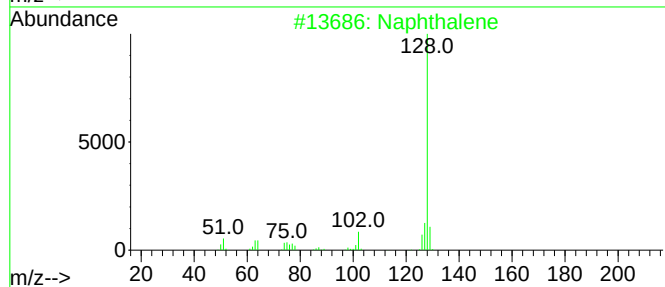
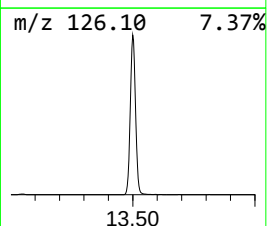
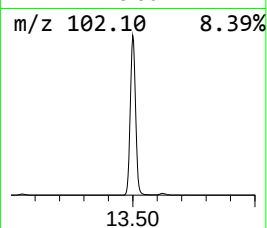
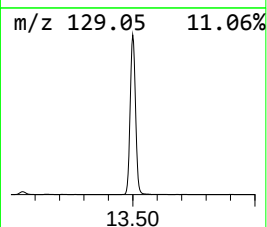
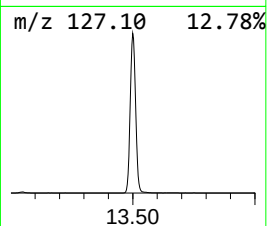
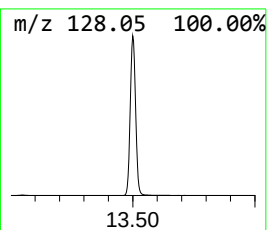
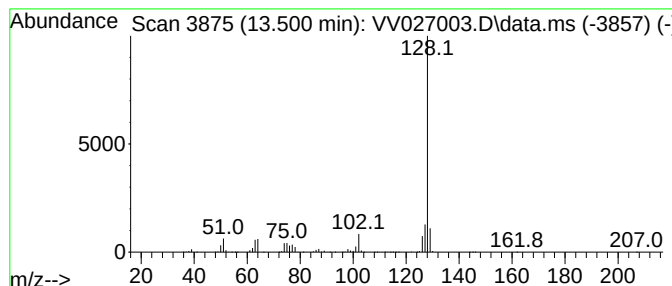
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Quant Title : VOC Analysis

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

Peak Number 5 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	114.02 ug/L	4105520	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\VV072722\
Data File : VV027003.D
Acq On : 27 Jul 2022 16:58
Operator : SY/MD
Sample : N3843-07DL 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUE7DL

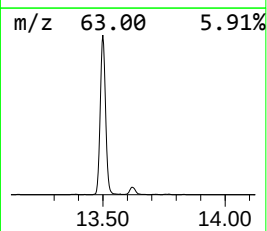
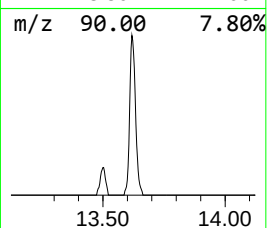
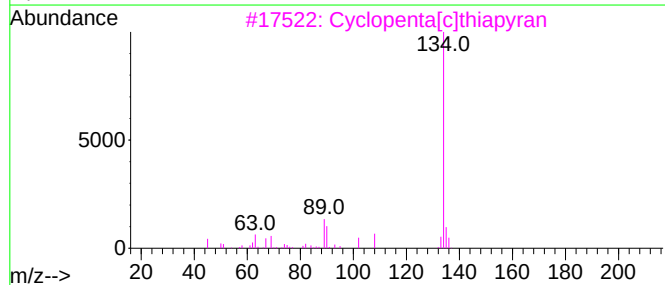
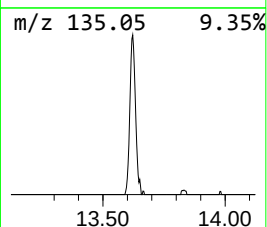
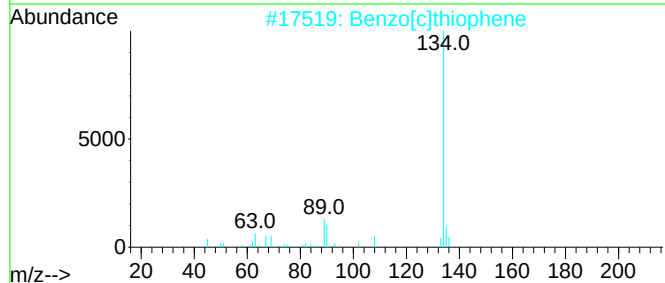
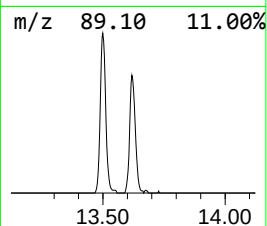
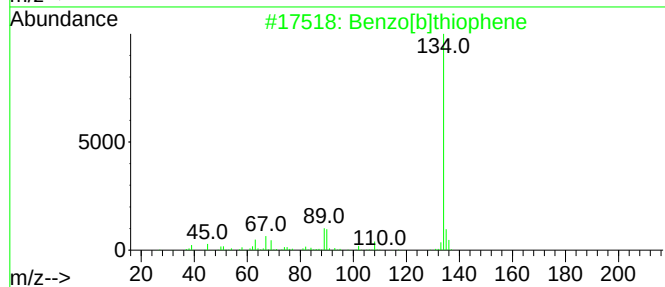
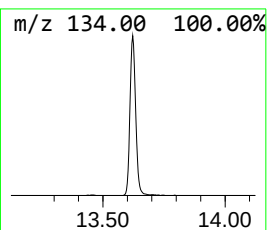
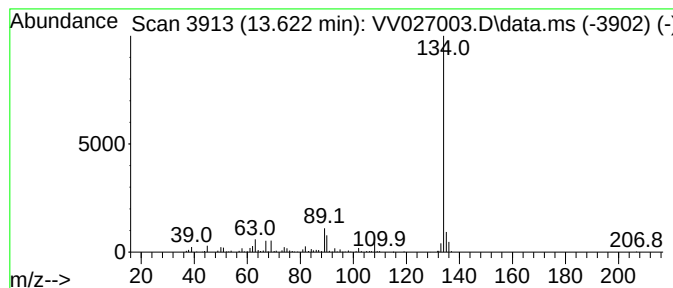
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Quant Title : VOC Analysis

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

Peak Number 6 Benzo[b]thiophene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	5.61 ug/L	201891	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene	134	C8H6S	000095-15-8	95
2			Benzo[c]thiophene	134	C8H6S	000270-82-6	95
3			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
4			Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5			Benzo[b]thiophene	134	C8H6S	000095-15-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
Data File : VV027003.D
Acq On : 27 Jul 2022 16:58
Operator : SY/MD
Sample : N3843-07DL 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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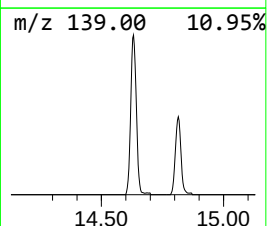
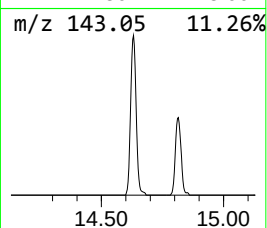
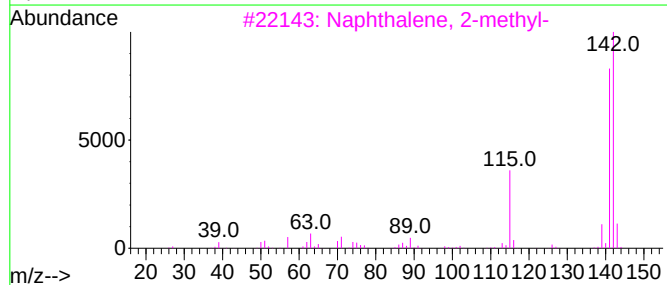
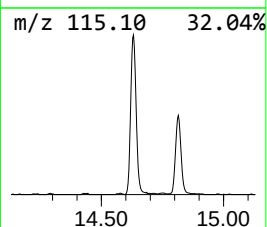
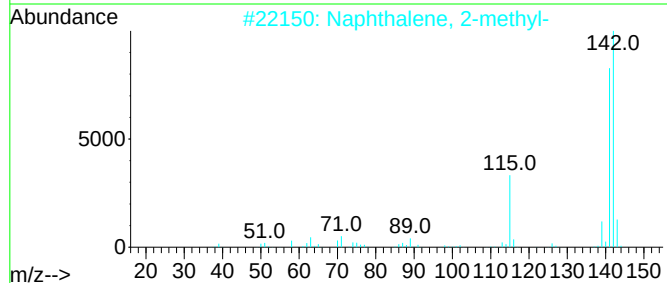
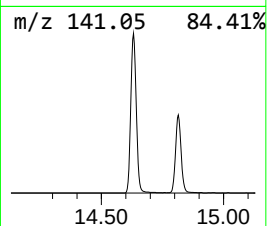
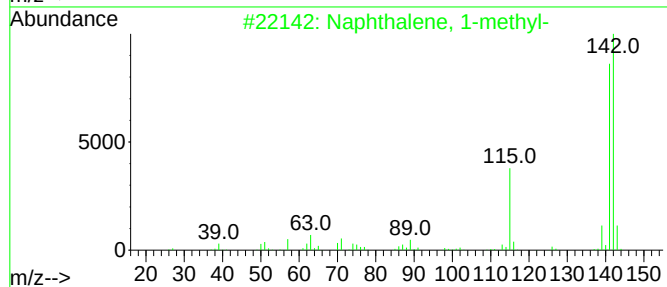
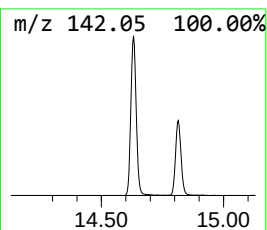
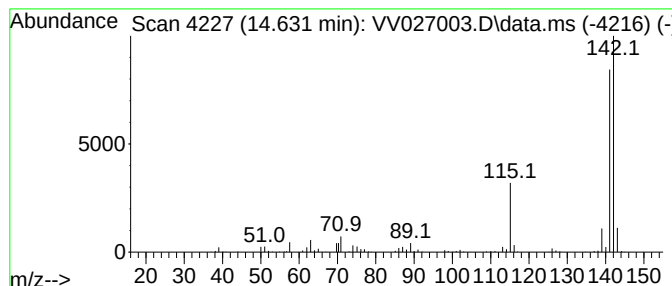
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Quant Title : VOC Analysis

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

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.631	14.06 ug/L	506165	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
Data File : VV027003.D
Acq On : 27 Jul 2022 16:58
Operator : SY/MD
Sample : N3843-07DL 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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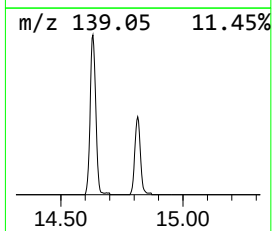
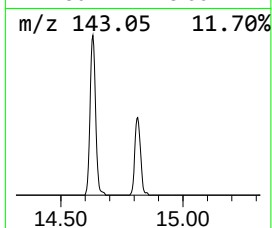
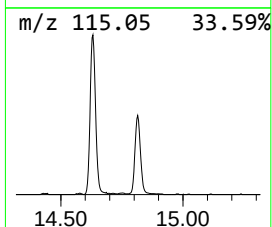
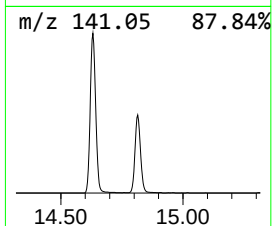
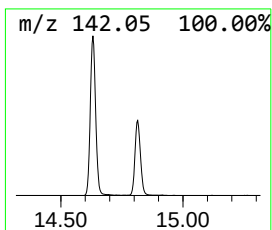
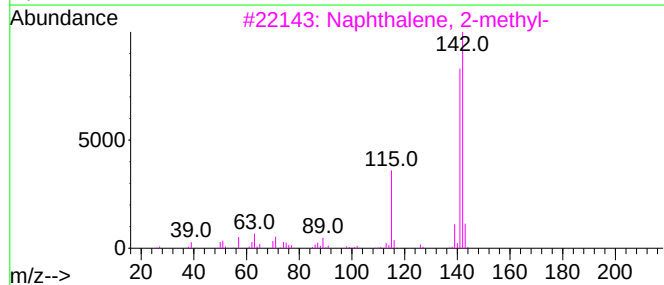
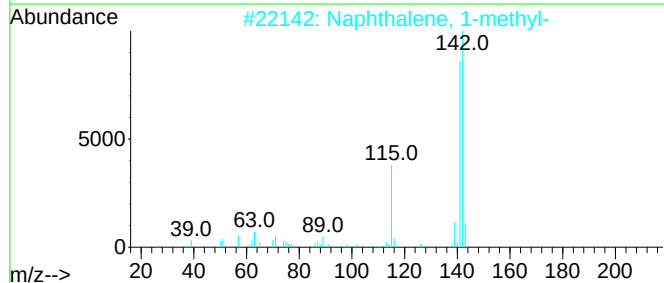
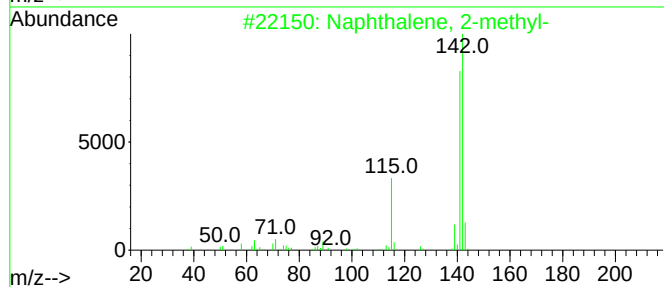
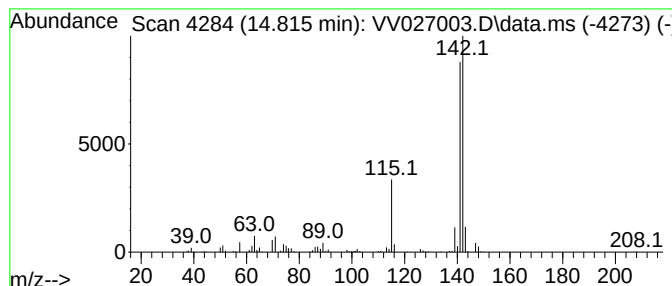
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Quant Title : VOC Analysis

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

Peak Number 8 Naphthalene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.815	7.27 ug/L	261722	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	95
5			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
Data File : VV027003.D
Acq On : 27 Jul 2022 16:58
Operator : SY/MD
Sample : N3843-07DL 10X
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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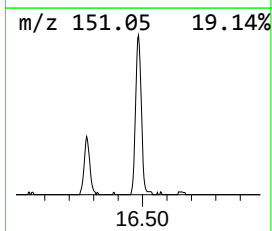
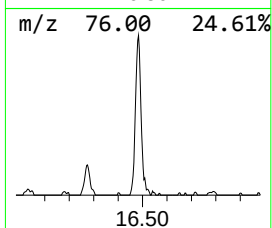
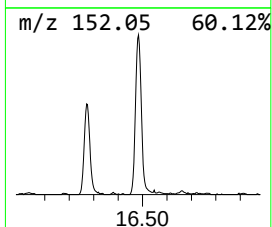
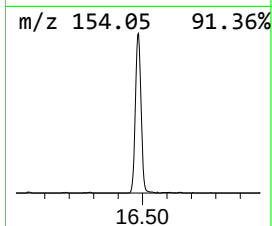
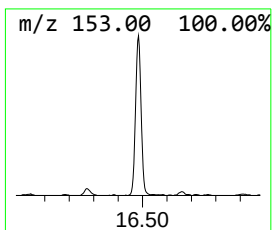
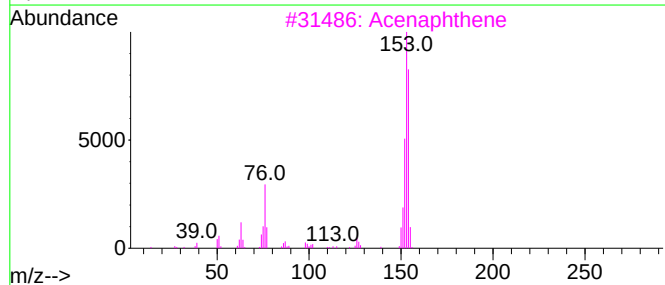
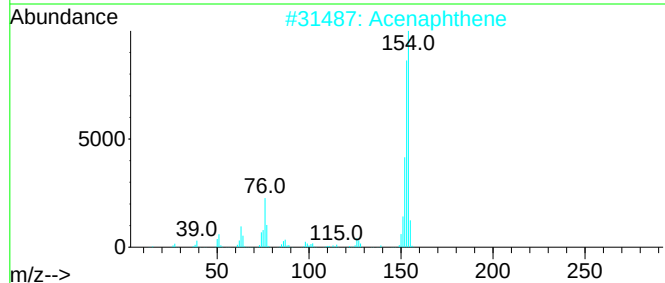
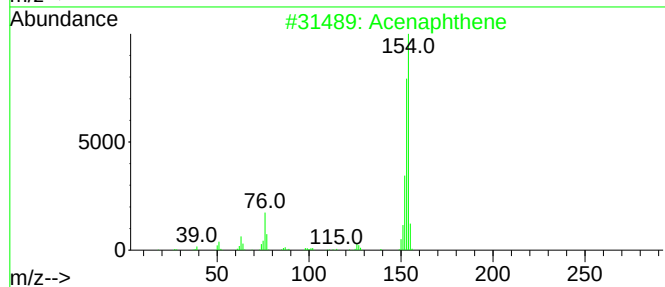
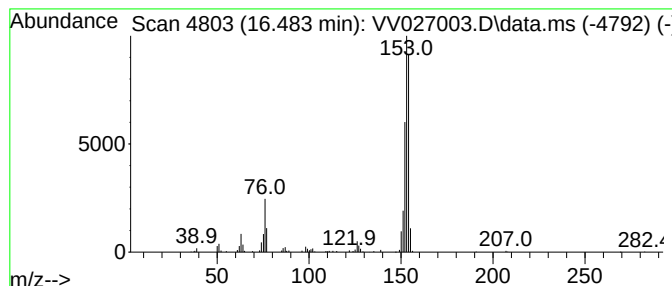
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Quant Title : VOC Analysis

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

Peak Number 9 Acenaphthene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.483	7.18 ug/L	258471	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	93
2			Acenaphthene	154	C12H10	000083-32-9	90
3			Acenaphthene	154	C12H10	000083-32-9	89
4			Acenaphthene	154	C12H10	000083-32-9	64
5			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072722\
Data File : VV027003.D
Acq On : 27 Jul 2022 16:58
Operator : SY/MD
Sample : N3843-07DL 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUE7DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM072222WMA.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
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Indene	11.744	18.5	ug/L	665429	3	11.249	1800390	50.0
Naphthalene, 1-...	13.046	3.0	ug/L	109644	3	11.249	1800390	50.0
Azulene	13.500	114.0	ug/L	4105520	3	11.249	1800390	50.0
Benzo[b]thiophene	13.622	5.6	ug/L	201891	3	11.249	1800390	50.0
Naphthalene, 1-...	14.631	14.1	ug/L	506165	3	11.249	1800390	50.0
Naphthalene, 2-...	14.815	7.3	ug/L	261722	3	11.249	1800390	50.0
Acenaphthene	16.483	7.2	ug/L	258471	3	11.249	1800390	50.0