

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
 Data File : VV022111.D
 Acq On : 07 Sep 2021 19:25
 Operator : SY/MD
 Sample : M3651-05
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBKM1

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\SFAMVLM082721WMA.M

Title : VOC Analysis

Signal : TIC: VV022111.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	76	82	99	rVB	124293	129133	2.66%	0.745%
2	1.565	157	163	179	rVB	115112	127277	2.62%	0.734%
3	2.105	321	331	346	rBV	231625	339594	6.99%	1.959%
4	2.423	427	430	433	rBV3	530	408	0.01%	0.002%
5	2.693	509	514	519	rBV2	518	693	0.01%	0.004%
6	2.831	554	557	558	rVB2	320	145	0.00%	0.001%
7	2.886	572	574	579	rVB	351	234	0.00%	0.001%
8	2.957	589	596	601	rBV3	398	603	0.01%	0.003%
9	3.124	645	648	652	rBV	316	265	0.01%	0.002%
10	3.224	677	679	683	rVV	267	199	0.00%	0.001%
11	3.503	764	766	768	rBV	295	148	0.00%	0.001%
12	3.777	849	851	856	rBV	312	202	0.00%	0.001%
13	3.799	856	858	860	rBV	337	154	0.00%	0.001%
14	3.851	868	874	877	rBV2	312	334	0.01%	0.002%
15	3.896	877	888	921	rBV	85735	271399	5.59%	1.566%
16	4.352	1013	1030	1059	rBV	169666	426219	8.77%	2.459%
17	4.648	1120	1122	1129	rVB3	539	308	0.01%	0.002%
18	4.699	1134	1138	1140	rBV2	460	385	0.01%	0.002%
19	4.754	1151	1155	1156	rBV	326	188	0.00%	0.001%
20	5.047	1230	1246	1285	rBV2	417340	1095054	22.54%	6.317%
21	5.314	1327	1329	1332	rVB	622	298	0.01%	0.002%
22	5.333	1332	1335	1337	rBV	505	287	0.01%	0.002%
23	5.416	1357	1361	1369	rVB5	632	781	0.02%	0.005%
24	5.494	1380	1385	1387	rBV2	436	308	0.01%	0.002%
25	5.535	1397	1398	1402	rVB2	344	171	0.00%	0.001%
26	5.619	1412	1424	1446	rBV	377464	755114	15.54%	4.356%
27	5.908	1505	1514	1530	rBV5	10629	22325	0.46%	0.129%
28	6.072	1551	1565	1591	rBV	240410	491139	10.11%	2.833%
29	6.288	1630	1632	1634	rVB2	521	210	0.00%	0.001%
30	6.362	1653	1655	1658	rVB2	530	224	0.00%	0.001%
31	6.388	1660	1663	1665	rBV2	373	191	0.00%	0.001%
32	6.516	1697	1703	1705	rBV2	262	270	0.01%	0.002%
33	6.561	1713	1717	1718	rBV2	362	293	0.01%	0.002%
34	6.593	1725	1727	1730	rVB2	364	173	0.00%	0.001%
35	6.863	1809	1811	1816	rVB2	423	232	0.00%	0.001%
36	6.905	1819	1824	1825	rBV2	244	190	0.00%	0.001%

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
 Title : VOC Analysis

37	6.989	1839	1850	1882	rBV	122203	228194	4.70%	1.316%
38	7.317	1941	1952	1967	rBV	484807	841068	17.31%	4.852%
39	7.625	2038	2048	2070	rBV	81735	146729	3.02%	0.846%
40	7.847	2113	2117	2120	rBV2	271	168	0.00%	0.001%
41	7.902	2132	2134	2138	rBV2	254	185	0.00%	0.001%
42	7.947	2141	2148	2153	rVB3	980	1579	0.03%	0.009%
43	8.043	2175	2178	2180	rVB2	273	156	0.00%	0.001%
44	8.092	2180	2193	2230	rBV	306586	570184	11.74%	3.289%
45	8.580	2342	2345	2350	rVB2	462	331	0.01%	0.002%
46	8.654	2365	2368	2371	rVB2	438	261	0.01%	0.002%
47	8.719	2385	2388	2389	rBV	279	144	0.00%	0.001%
48	8.789	2407	2410	2413	rVB2	257	163	0.00%	0.001%
49	8.854	2418	2430	2450	rBV	561345	902208	18.57%	5.204%
50	9.014	2470	2480	2501	rBV	165392	262847	5.41%	1.516%
51	9.140	2509	2519	2535	rBV	70726	117403	2.42%	0.677%
52	9.249	2551	2553	2561	rVB3	768	810	0.02%	0.005%
53	9.297	2566	2568	2570	rBV	323	186	0.00%	0.001%
54	9.310	2570	2572	2574	rBV2	606	361	0.01%	0.002%
55	9.381	2591	2594	2598	rVB2	413	296	0.01%	0.002%
56	9.423	2605	2607	2611	rVB	244	143	0.00%	0.001%
57	9.477	2621	2624	2625	rBV	292	153	0.00%	0.001%
58	9.548	2636	2646	2667	rBV3	50554	98902	2.04%	0.571%
59	9.699	2691	2693	2698	rBV2	294	198	0.00%	0.001%
60	9.770	2708	2715	2718	rBV3	458	523	0.01%	0.003%
61	9.786	2718	2720	2725	rVV2	402	290	0.01%	0.002%
62	9.863	2740	2744	2745	rBV	329	179	0.00%	0.001%
63	9.918	2759	2761	2762	rBV	426	170	0.00%	0.001%
64	9.937	2764	2767	2773	rVB	1547	1387	0.03%	0.008%
65	10.124	2820	2825	2826	rBV2	1500	1197	0.02%	0.007%
66	10.217	2841	2854	2874	rBV	356155	561378	11.55%	3.238%
67	10.301	2879	2880	2882	rBV	362	181	0.00%	0.001%
68	10.452	2919	2927	2945	rVB	10400	20708	0.43%	0.119%
69	10.542	2949	2955	2969	rVB3	8589	12955	0.27%	0.075%
70	10.744	3009	3018	3025	rBV2	5870	7936	0.16%	0.046%
71	10.815	3038	3040	3043	rVB2	335	171	0.00%	0.001%
72	10.918	3062	3072	3085	rVB2	29940	47125	0.97%	0.272%
73	10.992	3089	3095	3101	rBV4	3919	5321	0.11%	0.031%
74	11.137	3137	3140	3141	rBV	585	298	0.01%	0.002%
75	11.172	3141	3151	3165	rVV	121238	197214	4.06%	1.138%
76	11.249	3165	3175	3194	rVV	587928	916538	18.86%	5.287%

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 Title : VOC Analysis

77	11.339	3197	3203	3214	rVB	16848	25735	0.53%	0.148%
78	11.474	3240	3245	3249	rVB3	539	372	0.01%	0.002%
79	11.529	3250	3262	3280	rBV	268106	412670	8.49%	2.380%
80	11.625	3281	3292	3314	rVB	537201	841528	17.32%	4.854%
81	11.744	3318	3329	3351	rBV	853120	1317872	27.12%	7.602%
82	12.117	3438	3445	3461	rVB	12134	19073	0.39%	0.110%
83	12.239	3482	3483	3484	rBV	497	182	0.00%	0.001%
84	12.365	3513	3522	3533	rBV2	22532	34301	0.71%	0.198%
85	12.468	3544	3554	3565	rBV2	55374	103057	2.12%	0.594%
86	12.651	3606	3611	3613	rBV3	874	811	0.02%	0.005%
87	12.728	3627	3635	3647	rVB	17811	25483	0.52%	0.147%
88	12.886	3675	3684	3697	rVB2	34557	52025	1.07%	0.300%
89	12.953	3697	3705	3715	rVB6	4528	6834	0.14%	0.039%
90	13.050	3723	3735	3752	rBV2	111681	178011	3.66%	1.027%
91	13.198	3780	3781	3782	rBV	465	160	0.00%	0.001%
92	13.271	3798	3804	3810	rBV5	2175	2861	0.06%	0.017%
93	13.416	3846	3849	3852	rBV4	1251	794	0.02%	0.005%
94	13.503	3863	3876	3902	rBV	3243287	4858756	100.00%	28.027%
95	13.622	3904	3913	3928	rVB	163299	252686	5.20%	1.458%
96	14.580	4205	4211	4218	rBV3	9186	12430	0.26%	0.072%
97	14.635	4218	4228	4248	rVB	131166	210704	4.34%	1.215%
98	14.818	4275	4285	4299	rBV	183154	292511	6.02%	1.687%
99	15.815	4588	4595	4601	rBV5	8282	11997	0.25%	0.069%
100	16.487	4796	4804	4816	rVB2	41721	63578	1.31%	0.367%

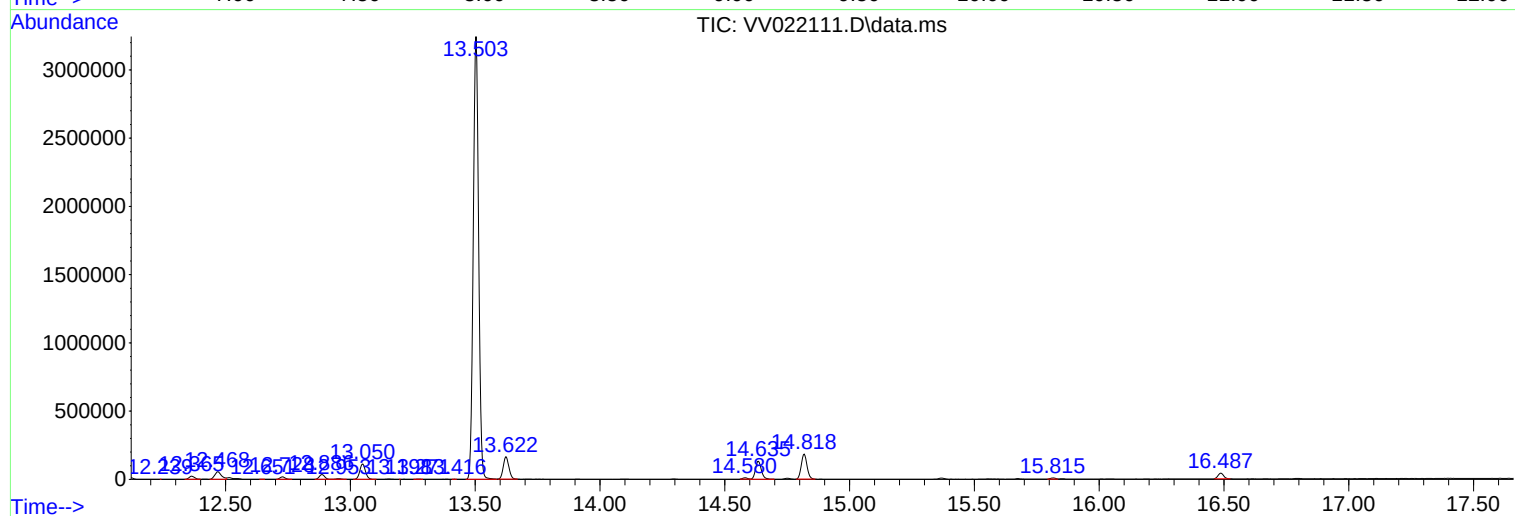
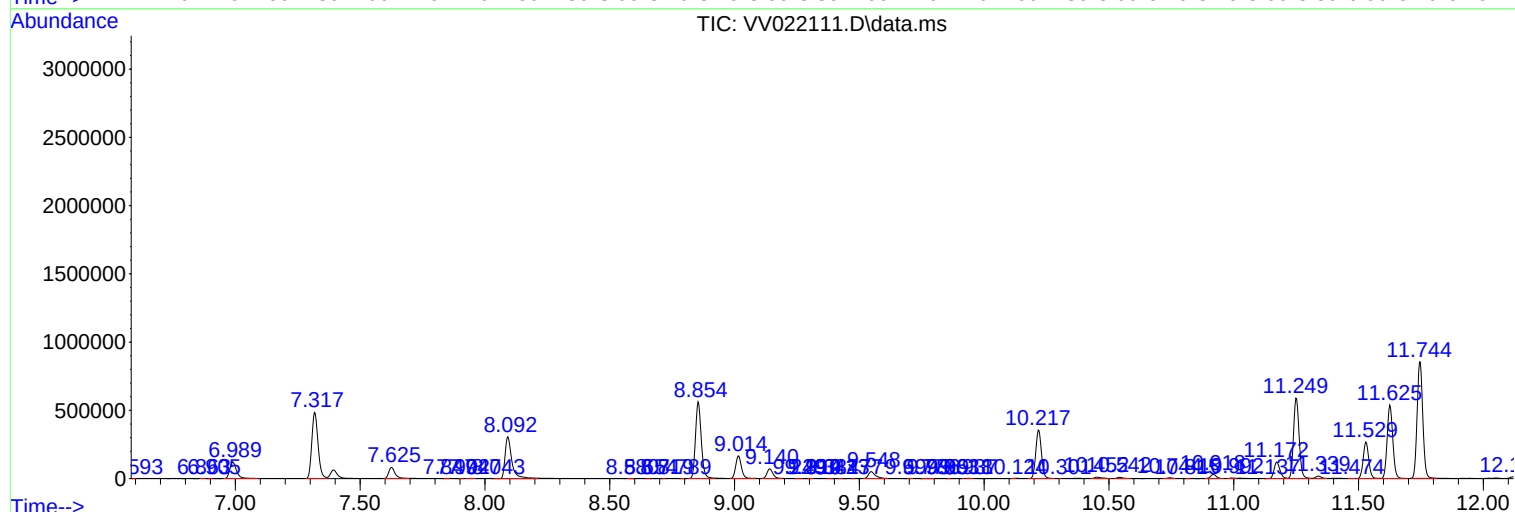
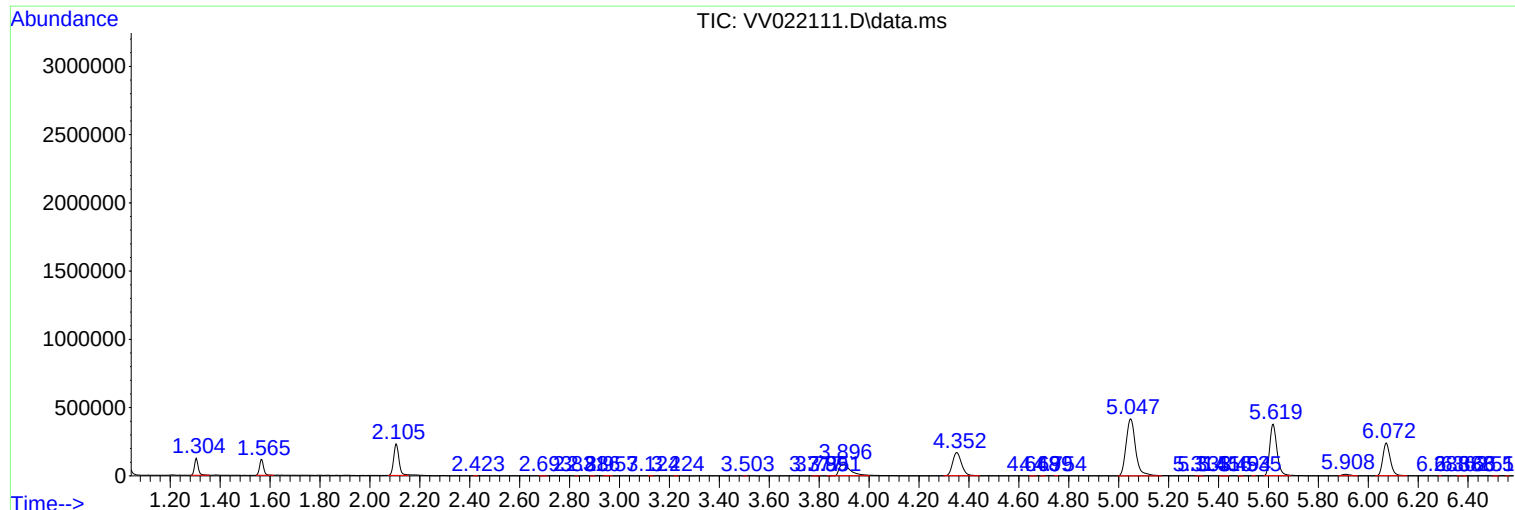
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

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



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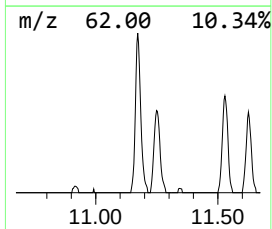
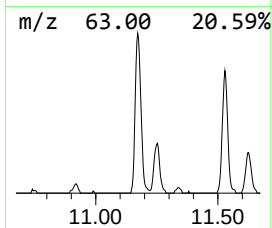
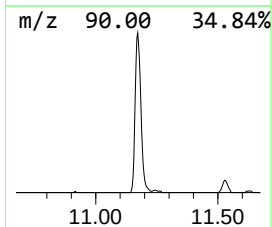
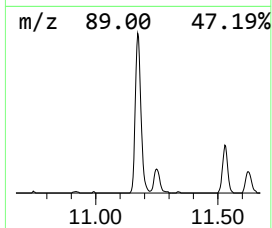
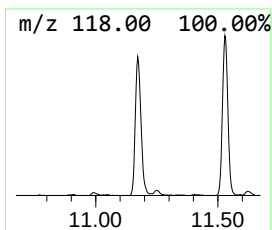
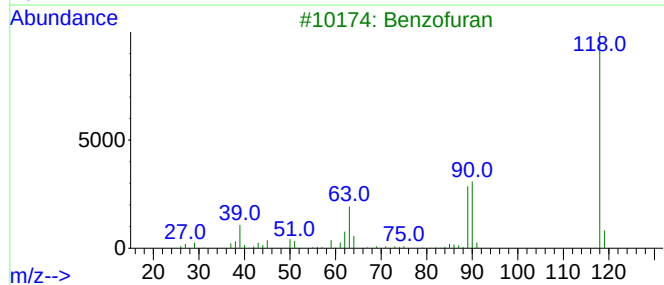
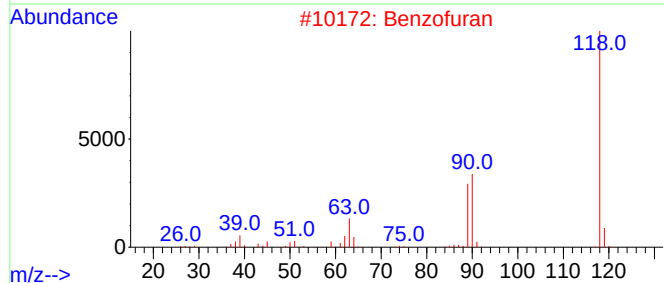
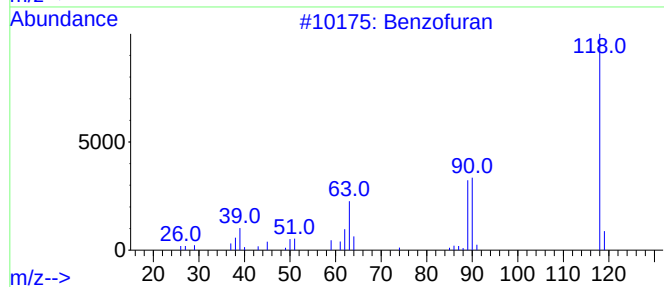
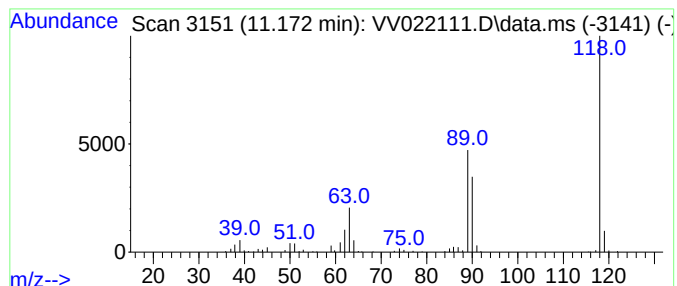
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzofuran Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.172	10.76 ug/L	197214	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran	118	C8H6O	000271-89-6	94
2			Benzofuran	118	C8H6O	000271-89-6	87
3			Benzofuran	118	C8H6O	000271-89-6	86
4			3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	80
5			Benzofuran	118	C8H6O	000271-89-6	74



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ClientSampleId :
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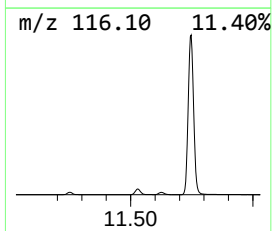
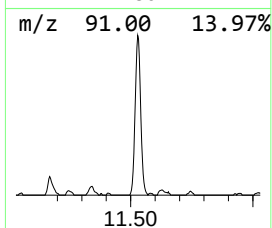
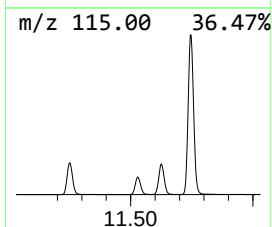
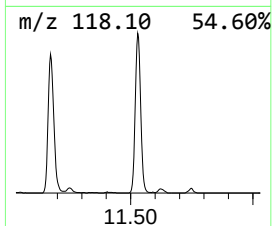
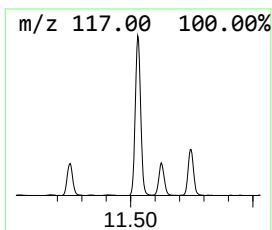
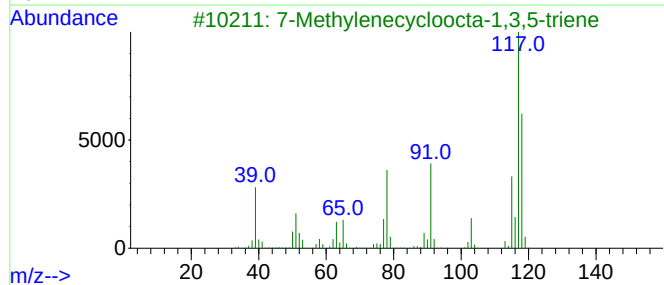
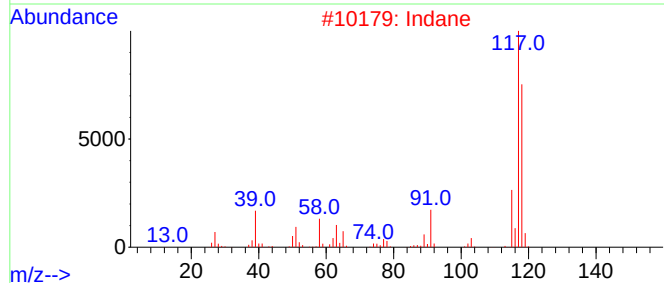
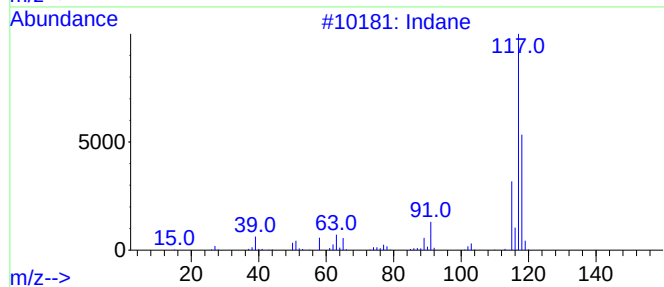
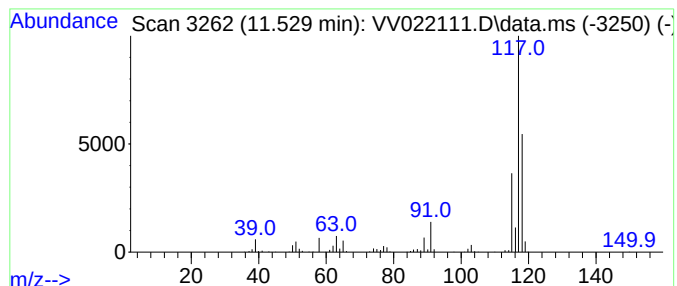
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	22.51 ug/L	412670	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	95
2			Indane	118	C9H10	000496-11-7	91
3			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	83
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
5			Deltacyclene	118	C9H10	007785-10-6	80



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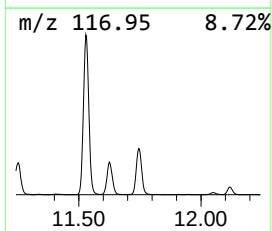
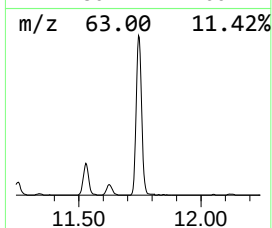
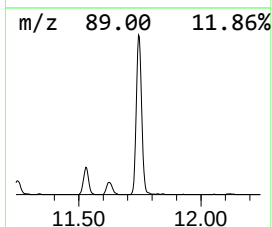
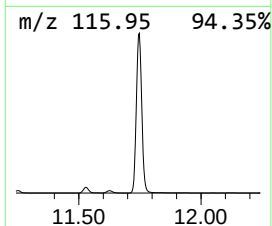
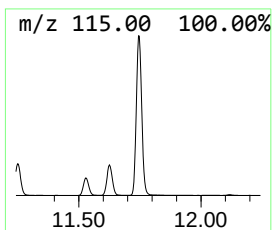
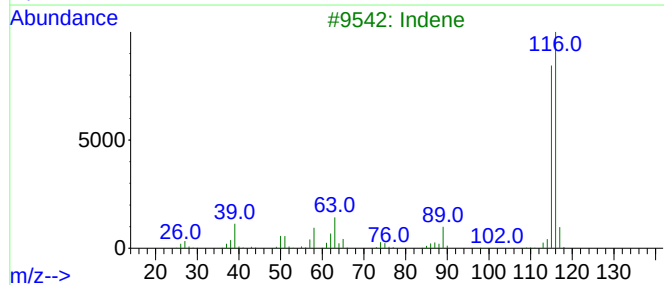
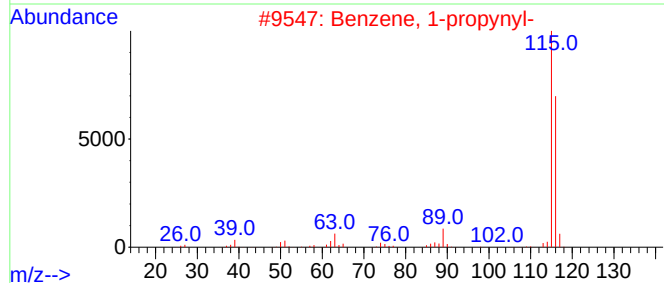
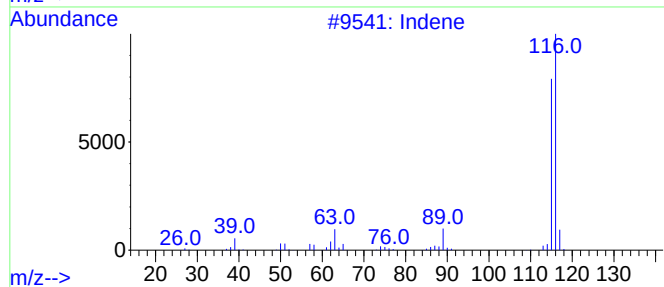
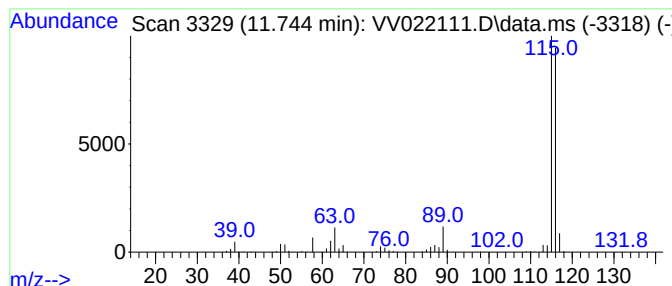
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.744	71.89 ug/L	1317870	1,4-Dichlorobenzene-d4	11.252

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			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

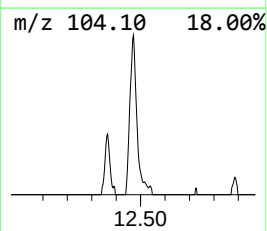
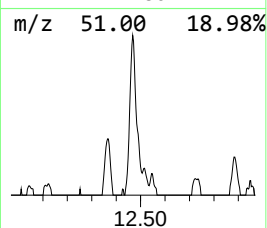
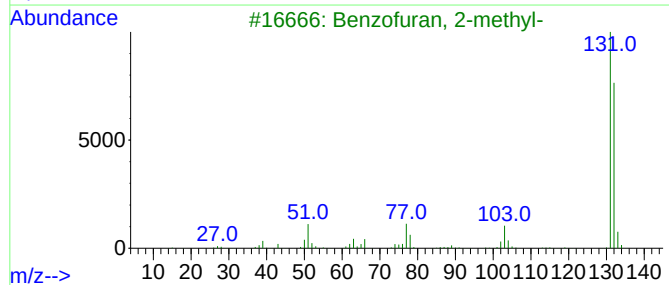
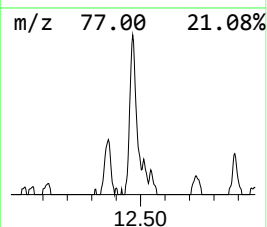
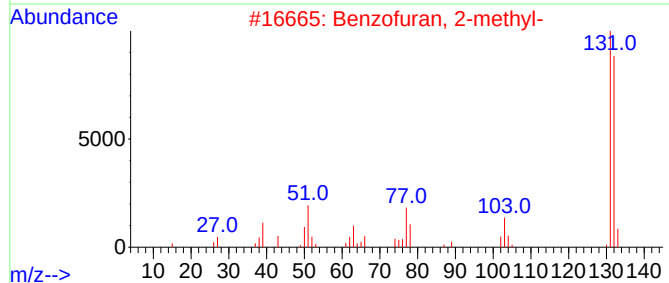
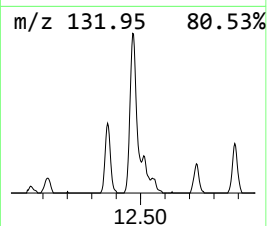
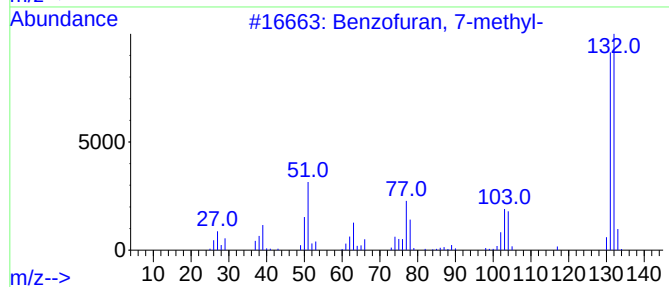
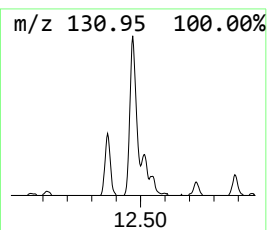
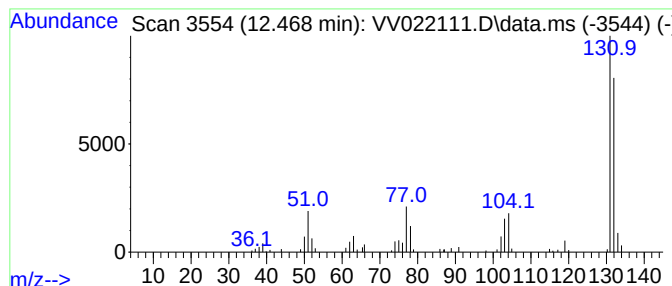
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Quant Title : VOC Analysis

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

Peak Number 5 Benzofuran, 7-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.468	5.62 ug/L	103057	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

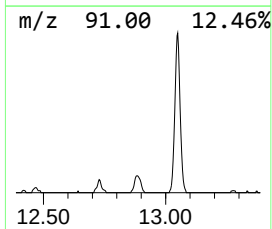
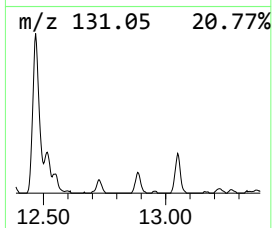
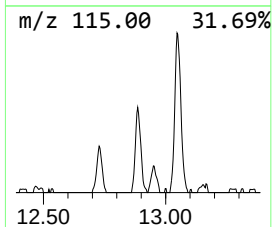
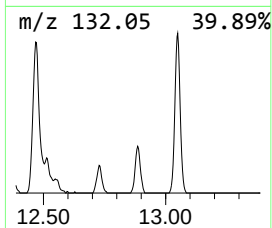
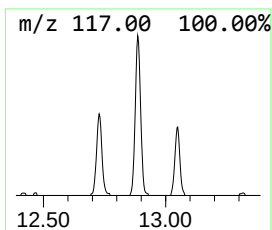
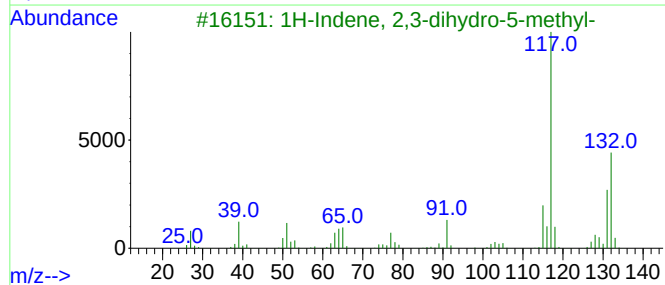
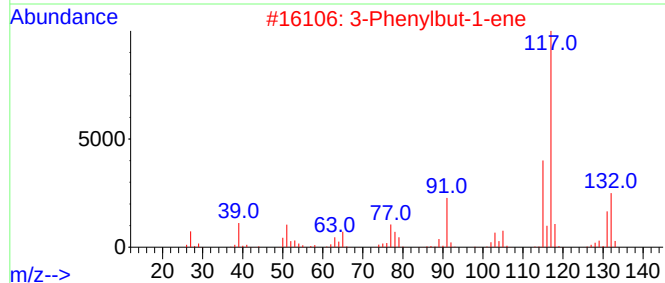
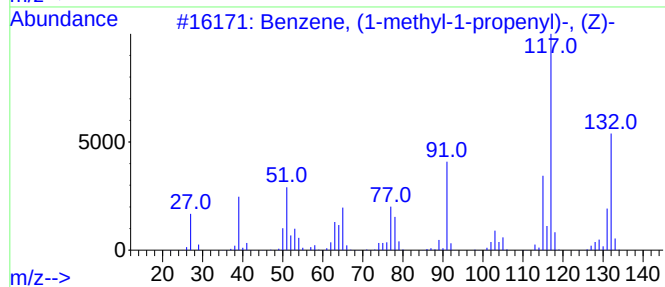
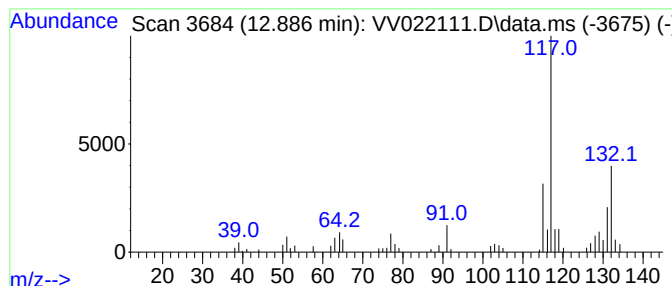
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, (1-methyl-1-propen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	2.84 ug/L	52025	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	90
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

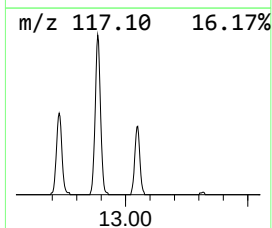
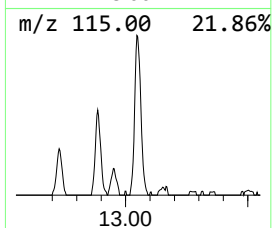
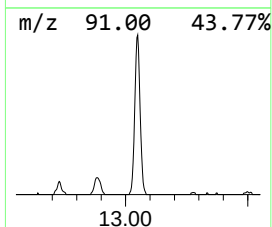
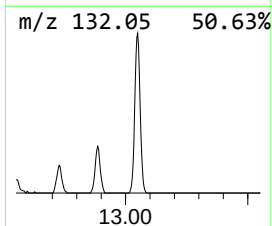
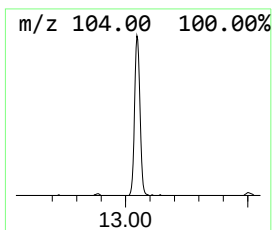
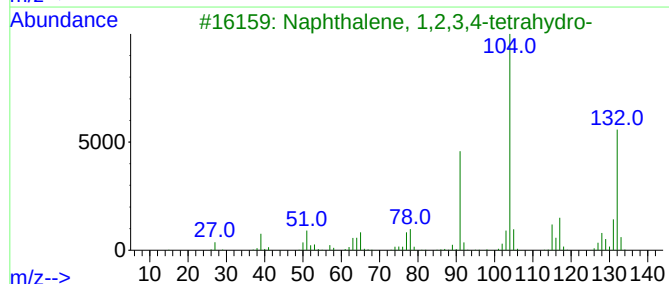
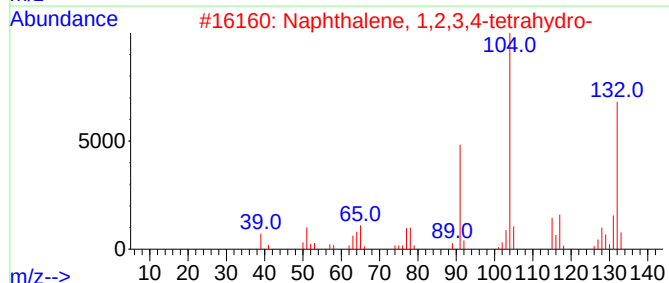
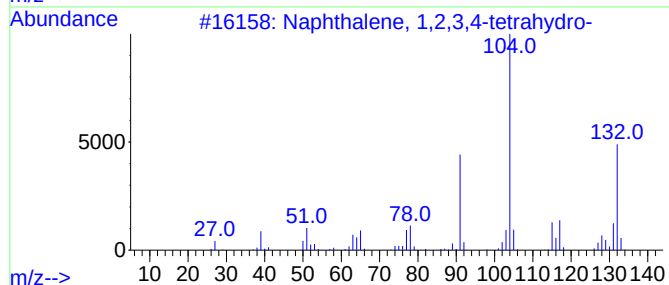
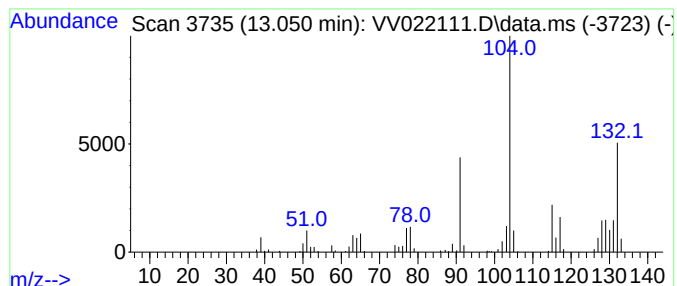
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Quant Title : VOC Analysis

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.050	9.71 ug/L	178011	1,4-Dichlorobenzene-d4	11.252

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\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

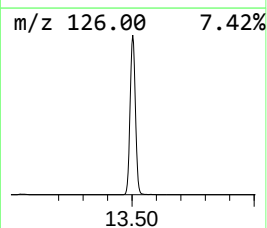
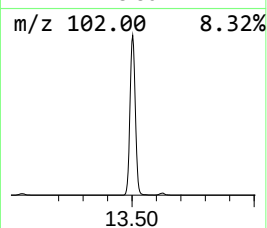
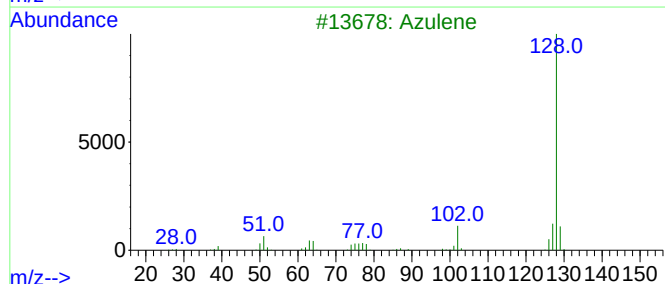
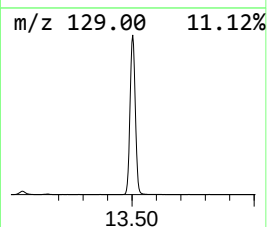
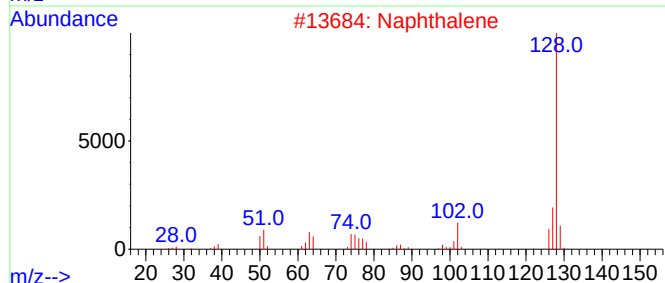
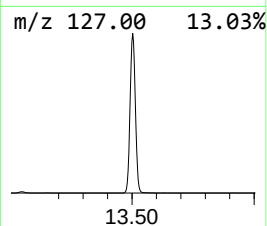
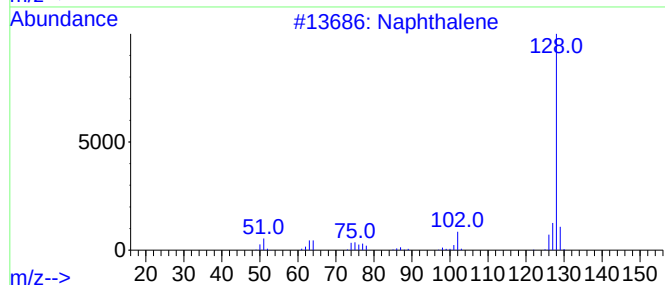
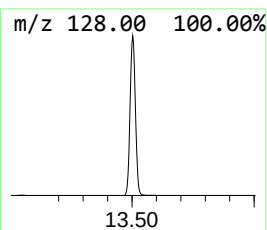
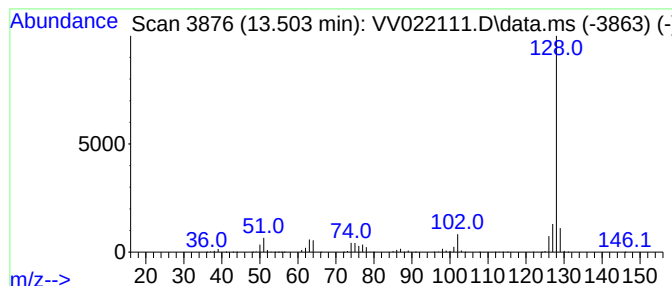
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Quant Title : VOC Analysis

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

Peak Number 8 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.503	265.06 ug/L	4858760	1,4-Dichlorobenzene-d4	11.252

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			Azulene	128	C10H8	000275-51-4	93
4			Naphthalene	128	C10H8	000091-20-3	93
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

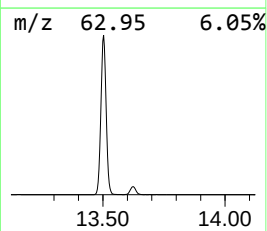
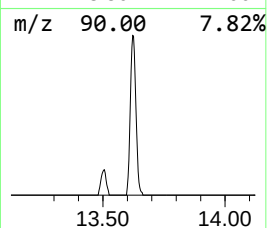
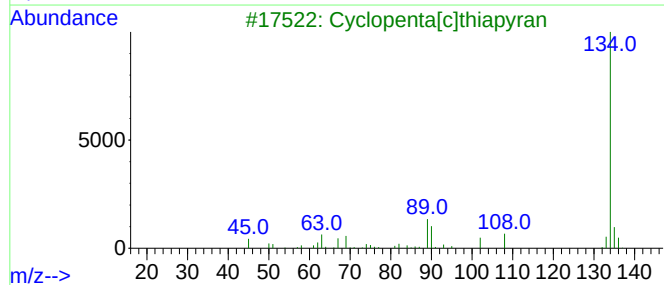
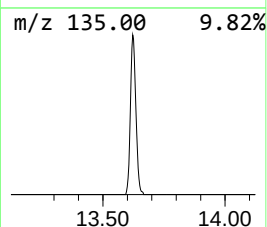
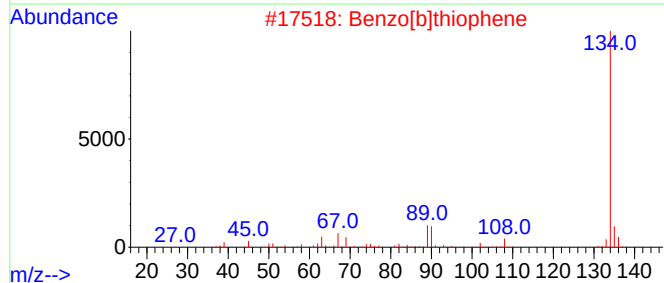
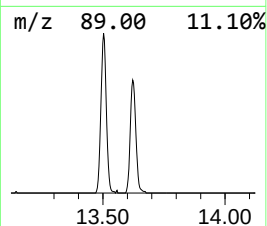
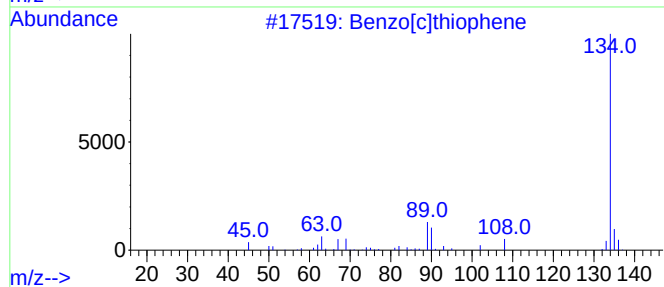
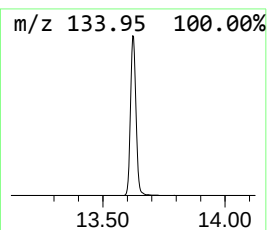
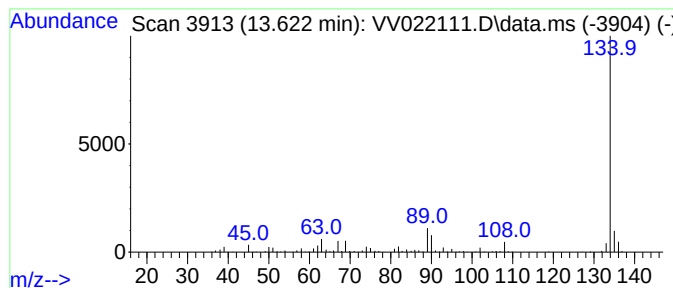
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Quant Title : VOC Analysis

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

Peak Number 9 Benzo[c]thiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	13.78 ug/L	252686	1,4-Dichlorobenzene-d4	11.252

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			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\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

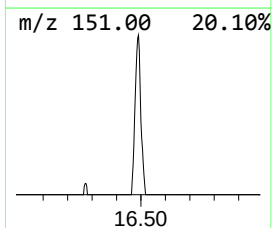
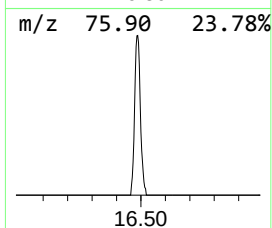
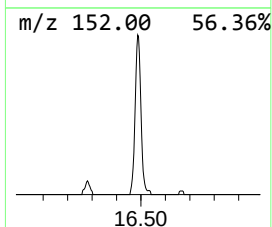
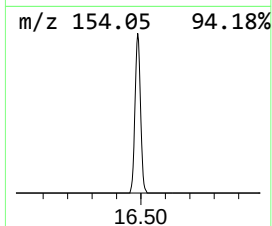
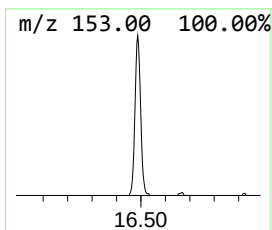
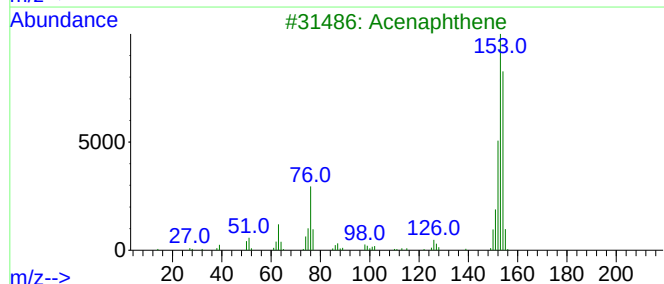
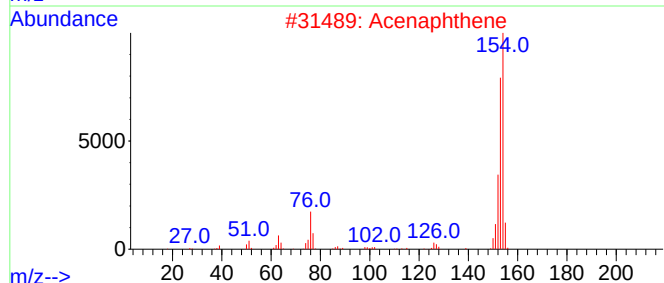
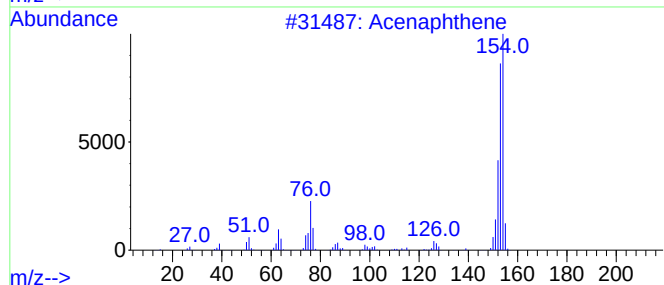
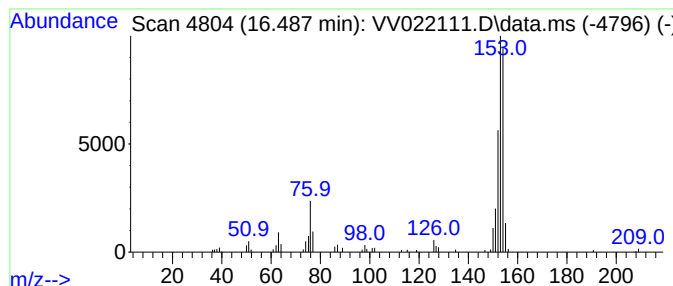
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Quant Title : VOC Analysis

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

Peak Number 12 Hexa-2,4-diyn-1-ylbenzene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.487	3.47 ug/L	63578	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	90
2			Acenaphthene	154	C12H10	000083-32-9	76
3			Acenaphthene	154	C12H10	000083-32-9	70
4			Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	64
5			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
Data File : VV022111.D
Acq On : 07 Sep 2021 19:25
Operator : SY/MD
Sample : M3651-05
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKM1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.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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Indane	11.529	22.5	ug/L	412670	3	11.252	916538	50.0
Indene	11.744	71.9	ug/L	1317870	3	11.252	916538	50.0
Benzofuran, 7-m...	12.468	5.6	ug/L	103057	3	11.252	916538	50.0
Benzene, (1-met...	12.886	2.8	ug/L	52025	3	11.252	916538	50.0
Naphthalene, 1,...	13.050	9.7	ug/L	178011	3	11.252	916538	50.0
Azulene	13.503	265.1	ug/L	4858760	3	11.252	916538	50.0
Benzo[c]thiophene	13.622	13.8	ug/L	252686	3	11.252	916538	50.0
Hexa-2,4-diyne-1...	16.487	3.5	ug/L	63578	3	11.252	916538	50.0