LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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: VV022061.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	76	83	97	rVB	201844	205129	0.62%	0.335%
2	1.564	156	163	175	rBV	157269	178669	0.54%	0.292%
3	2.108	322	332	358	rVB	328290	489579	1.48%	0.799%
4	2.404	422	424	430	rBV3	309	276	0.00%	0.000%
5	2.510	451	457	462	rBV7	1016	1484	0.00%	0.002%
6	2.574	472	477	480	rBV3	429	409	0.00%	0.001%
7	2.638	495	497	499	rVB2	477	180	0.00%	0.000%
8	2.661	501	504	506	rVB2	457	236	0.00%	0.000%
9	2.902	575	579	581	rBV2	289	244	0.00%	0.000%
10	3.005	609	611	615	rBV2	407	237	0.00%	0.000%
11	3.124	644	648	649	rBV	304	209	0.00%	0.000%
12	3.741	830	840	846	rBV5	999	1982	0.01%	0.003%
13	3.876	870	882	908	rBV	104658	271946	0.82%	0.444%
14	4.352	1013	1030	1064	rBV	209305	524925	1.59%	0.857%
15	4.629	1114	1116	1122	rVB4	346	223	0.00%	0.000%
16	4.661	1122	1126	1130	rBV3	310	345	0.00%	0.001%
17	4.709	1139	1141	1144	rBV	267	195	0.00%	0.000%
18	4.889	1194	1197	1202	rVB3	238	195	0.00%	0.000%
19	5.050	1228	1247	1287	rBV2	531278	1398050	4.23%	2.282%
20	5.371	1345	1347	1352	rBV2	476	243	0.00%	0.000%
21	5.484	1377	1382	1386	rBV3	414	463	0.00%	0.001%
22	5.516	1389	1392	1394	rBV3	337	202	0.00%	0.000%
23	5.545	1397	1401	1404	rBV	400	367	0.00%	0.001%
24	5.619	1412	1424	1454	rBV	342111	695582	2.10%	1.135%
25	5.911	1507	1515	1532	rVB5	5637	11940	0.04%	0.019%
26	6.072	1549	1565	1591	rBV	287826	596658	1.80%	0.974%
27	6.256	1620	1622	1626	rBV3	519	330	0.00%	0.001%
28	6.397	1664	1666	1669	rVB2	495	285	0.00%	0.000%
29	6.529	1705	1707	1711	rVB2	396	263	0.00%	0.000%
30	6.564	1712	1718	1722	rVB4	384	392	0.00%	0.001%
31	6.648	1742	1744	1747	rBV2	476	233	0.00%	0.000%
32	6.712	1762	1764	1768	rBV	269	206	0.00%	0.000%
33	6.879	1814	1816	1823	rVB2	462	270	0.00%	0.000%
34	6.989	1839	1850	1874	rBV	161855	296523	0.90%	0.484%
35	7.172	1905	1907	1912	rVB2	425	312	0.00%	0.001%
36	7.317	1940	1952	1968	rBV	659226	1137144	3.44%	1.856%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

37	7.625	2037	2048	2071	rBV	109731	193405	0.59%	0.316%
38	7.863	2120	2122	2127	rVB2	453	333	0.00%	0.001%
39	7.940	2137	2146	2153	rBV2	4322	6992	0.02%	0.011%
40	8.088	2182	2192	2224	rBV	338604	602981	1.82%	0.984%
41	8.853	2418	2430	2463	rBV	522223	860065	2.60%	1.404%
42	9.017	2470	2481	2498	rBV	85644	139603	0.42%	0.228%
43	9.088	2501	2503	2505	rBV2	502	195	0.00%	0.000%
44	9.143	2510	2520	2538	rVB	62383	105500	0.32%	0.172%
45	9.503	2628	2632	2634	rBV3	755	745	0.00%	0.001%
46	9.545	2634	2645	2667	rBV	217463	348005	1.05%	0.568%
47	9.648	2674	2677	2683	rVB3	852	808	0.00%	0.001%
48	9.673	2683	2685	2687	rBV2	557	278	0.00%	0.000%
49	9.773	2713	2716	2717	rBV2	383	201	0.00%	0.000%
50	9.853	2738	2741	2743	rBV2	361	181	0.00%	0.000%
51	9.873	2745	2747	2751	rBV3	395	230	0.00%	0.000%
52	9.934	2756	2766	2782	rBV2	25972	41659	0.13%	0.068%
53	10.008	2786	2789	2791	rVB2	378	204	0.00%	0.000%
54	10.034	2791	2797	2799	rBV	483	498	0.00%	0.001%
55	10.046	2799	2801	2806	rVB2	481	260	0.00%	0.000%
56	10.124	2818	2825	2835	rVV5	3942	6228	0.02%	0.010%
57	10.217	2842	2854	2874	rBV	423026	662305	2.00%	1.081%
58	10.358	2891	2898	2903	rBV3	6212	9846	0.03%	0.016%
59	10.452	2918	2927	2945	rBV	16417	37365	0.11%	0.061%
60	10.542	2945	2955	2971	rVB4	15116	27734	0.08%	0.045%
61	10.657	2987	2991	2993	rBV3	653	455	0.00%	0.001%
62	10.683	2997	2999	3005	rBV3	376	421	0.00%	0.001%
63	10.741	3006	3017	3028	rBV	51044	78471	0.24%	0.128%
64	10.812	3036	3039	3045	rVB2	771	670	0.00%	0.001%
65	10.837	3045	3047	3049	rBV2	576	292	0.00%	0.000%
66	10.918	3061	3072	3088	rVB	126181	198759	0.60%	0.324%
67	10.992	3089	3095	3102	rBV4	4168	6326	0.02%	0.010%
68	11.172	3139	3151	3165	rBV	382116	596374	1.80%	0.973%
69	11.249	3165	3175	3193	rVV	623136	986491	2.98%	1.610%
70	11.336	3194	3202	3214	rVB	142771	218025	0.66%	0.356%
71	11.529	3251	3262	3280	rBV	306111	487301	1.47%	0.795%
72	11.625	3281	3292	3313	rVB	706513	1122724	3.40%	1.833%
73	11.747	3319	3330	3347	rBV	596543	909707	2.75%	1.485%
74	11.914	3374	3382	3387	rBV2	15289	24100	0.07%	0.039%
75	12.021	3406	3415	3423	rBV	33847	52807	0.16%	0.086%
76	12.117	3434	3445	3467	rBV	139215	229443	0.69%	0.375%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

77	12.233	3478	3481	3482	rBV	834	458	0.00%	0.001%
78	12.361	3513	3521	3531	rBV	114051	171726	0.52%	0.280%
79	12.468	3544	3554	3566	rBV	296589	519945	1.57%	0.849%
80	12.728	3628	3635	3651	rVB	58428	91053	0.28%	0.149%
81	12.885	3676	3684	3697	rVB2	121779	186413	0.56%	0.304%
82	13.049	3725	3735	3751	rVB3	205792	320631	0.97%	0.523%
83	13.168	3760	3772	3779	rBV9	8069	16002	0.05%	0.026%
84	13.217	3780	3787	3796	rVB2	19688	29664	0.09%	0.048%
85	13.275	3797	3805	3811	rBV3	22308	34294	0.10%	0.056%
86	13.509	3863	3878	3901	rBV	20341624	33059963	100.00%	53.966%
87	13.622	3903	3913	3930	rVB	1044592	1607225	4.86%	2.624%
88	14.091	4049	4059	4074	rBV3	38096	85929	0.26%	0.140%
89	14.577	4200	4210	4218	rBV	146975	236166	0.71%	0.386%
90	14.635	4219	4228	4254	rVB	560800	931127	2.82%	1.520%
91	14.750	4256	4264	4273	rBV	53130	88468	0.27%	0.144%
92	14.818	4273	4285	4300	rVV	3131759	4944942	14.96%	8.072%
93	15.364	4432	4455	4467	rBV	706446	1083785	3.28%	1.769%
94	15.554	4507	4514	4521	rBV	150394	210499	0.64%	0.344%
95	15.670	4540	4550	4572	rVB	281841	516584	1.56%	0.843%
96	15.811	4586	4594	4601	rBV	404810	598381	1.81%	0.977%
97	16.184	4702	4710	4723	rVB	58601	90599	0.27%	0.148%
98	16.278	4730	4739	4755	rVB2	67608	140870	0.43%	0.230%
99	16.487	4793	4804	4825	rVB	1410658	2178628	6.59%	3.556%
100	16.786	4888	4897	4911	rVB	216914	344007	1.04%	0.562%

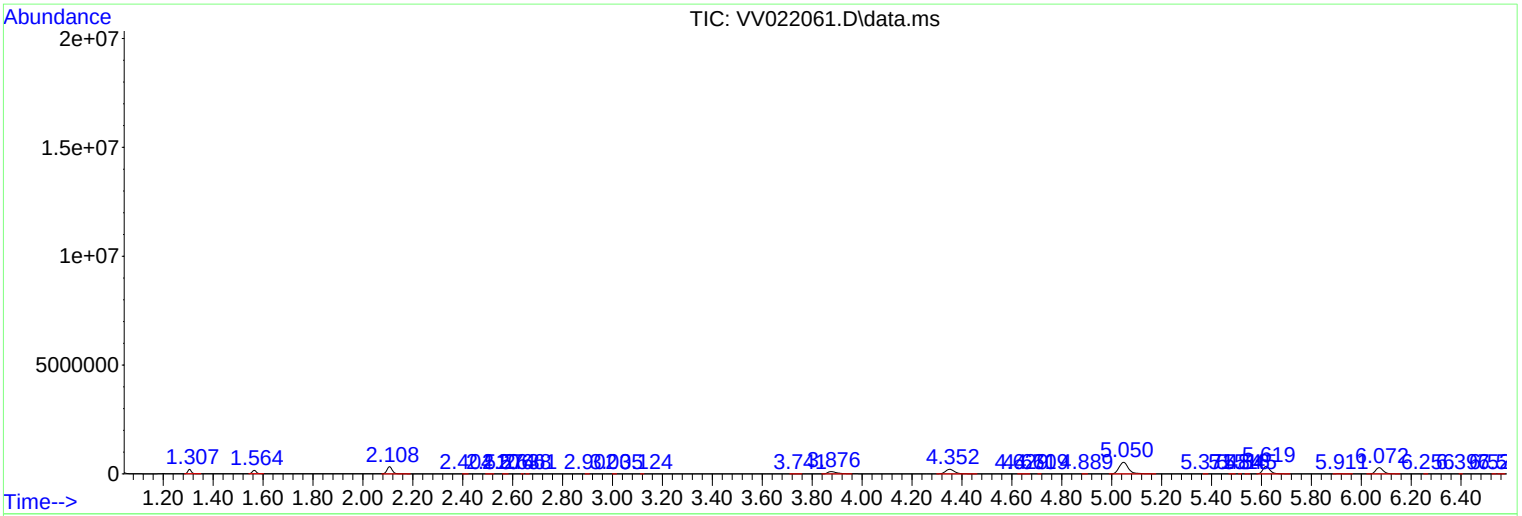
Sum of corrected areas: 61261272

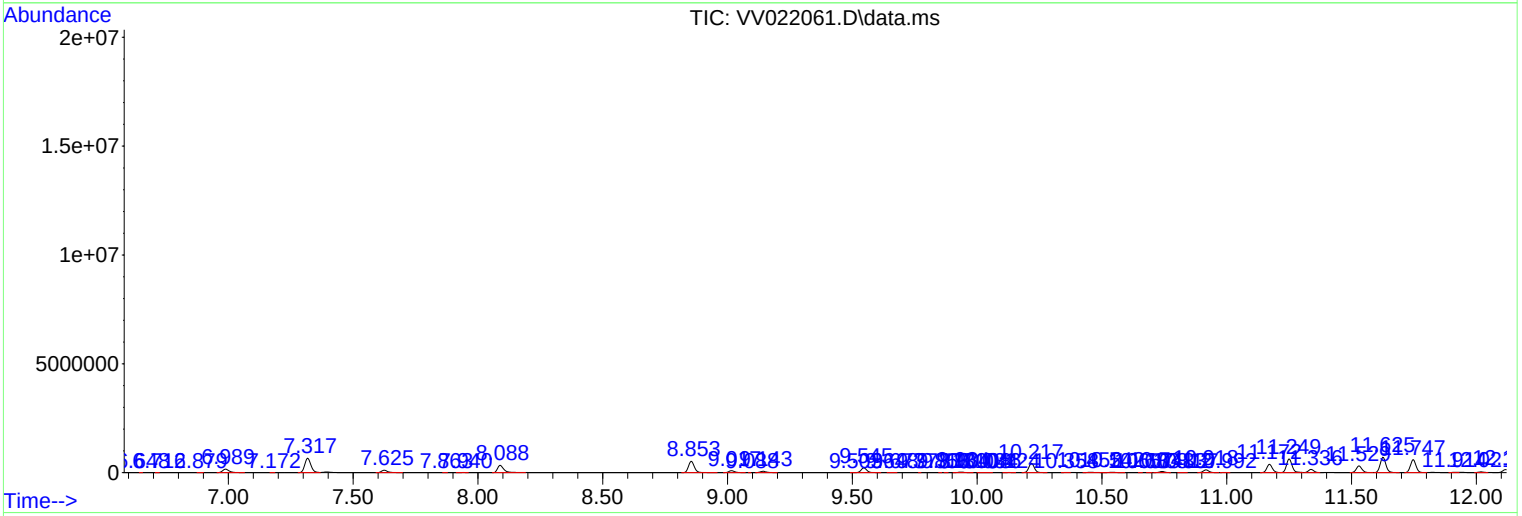
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

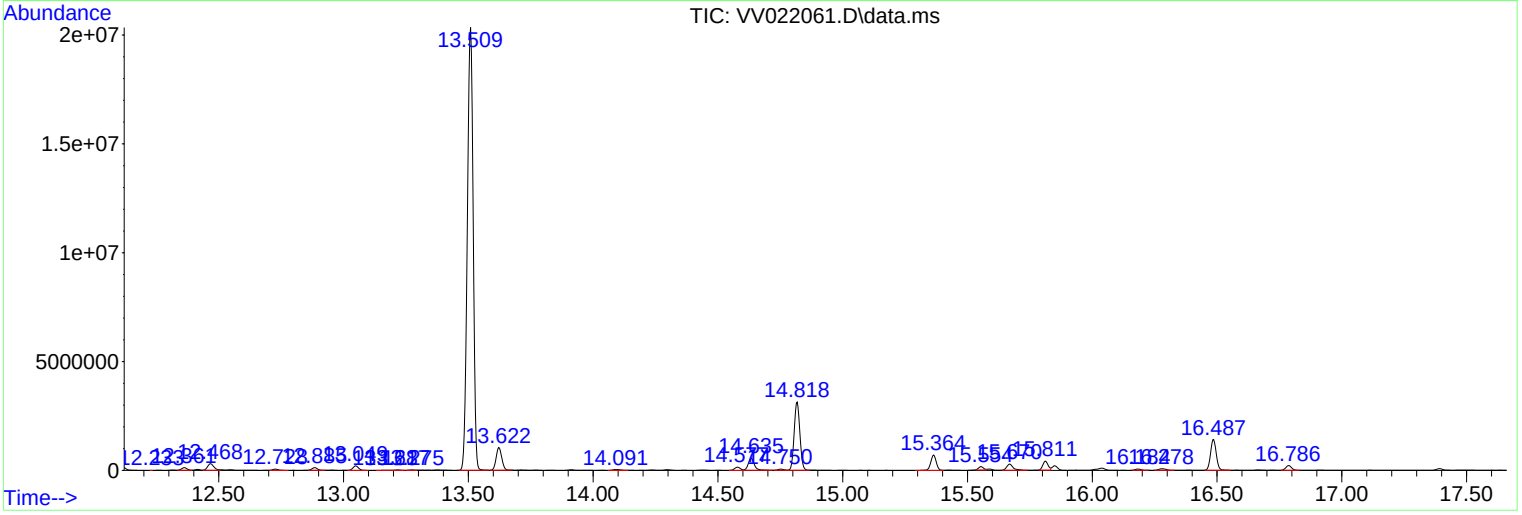
Instrument :
MSVOA_V
ClientSampleId :
DBKJ9

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

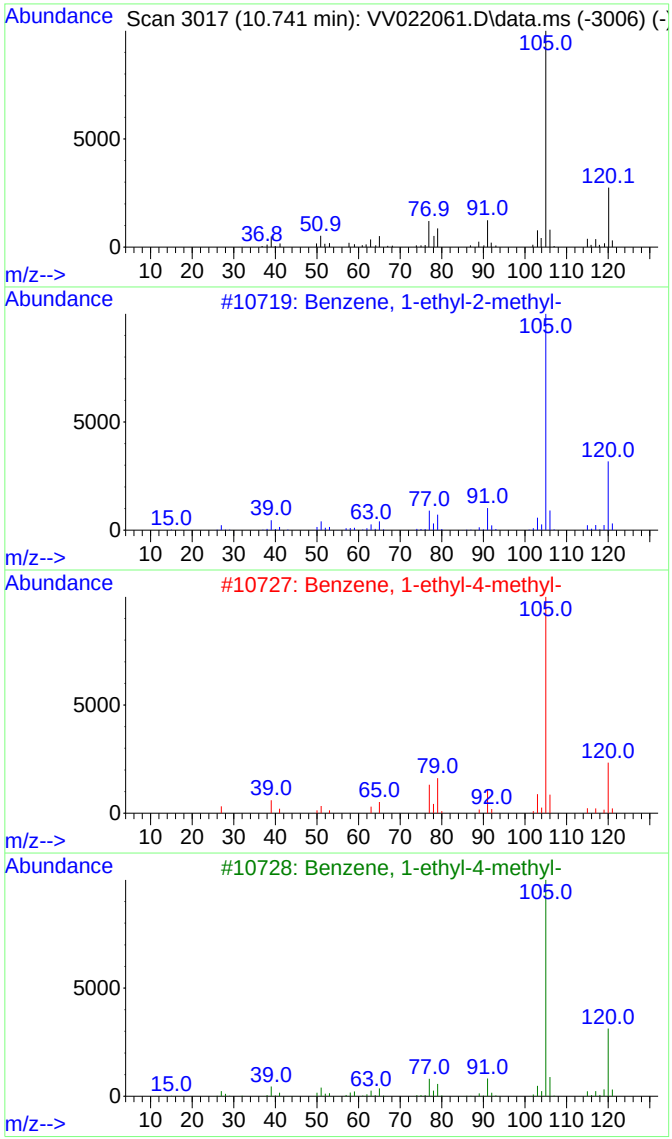
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

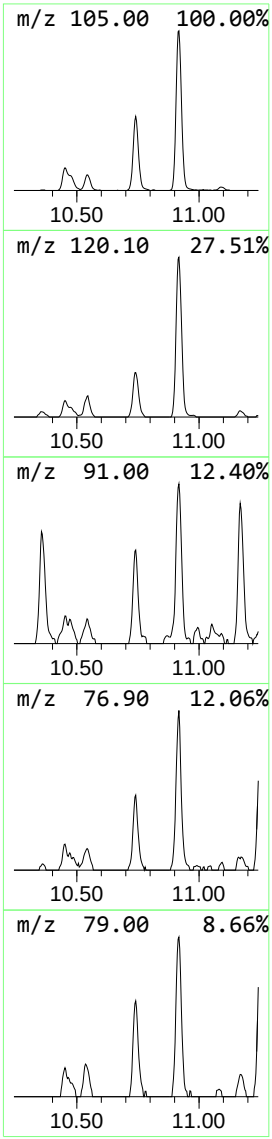
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.741	3.98 ug/L	78471	1,4-Dichlorobenzene-d4	11.252

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

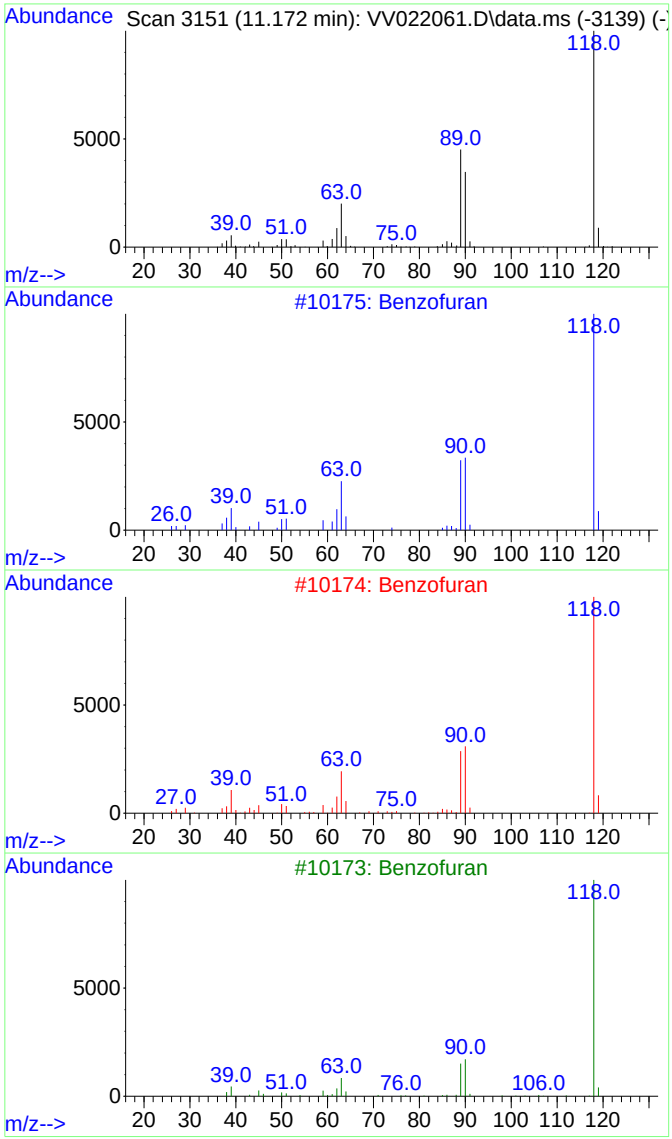
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

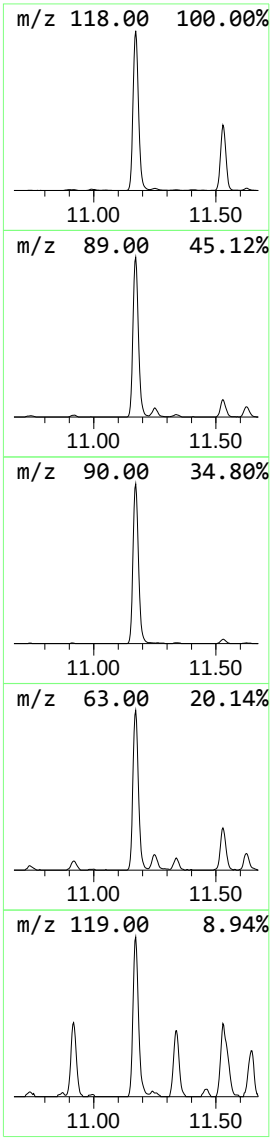
****

Peak Number 3 Benzofuran Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.172	30.23 ug/L	596374	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	96
2	Benzofuran	118	C8H6O	000271-89-6	91
3	Benzofuran	118	C8H6O	000271-89-6	90
4	Benzofuran	118	C8H6O	000271-89-6	87
5	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

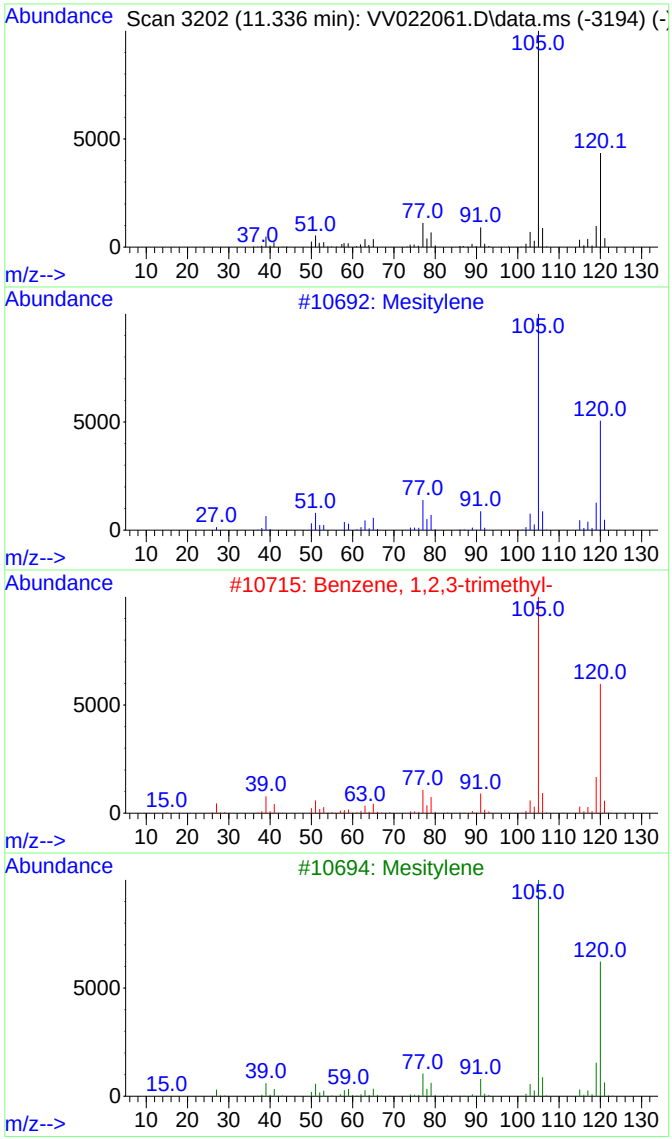
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

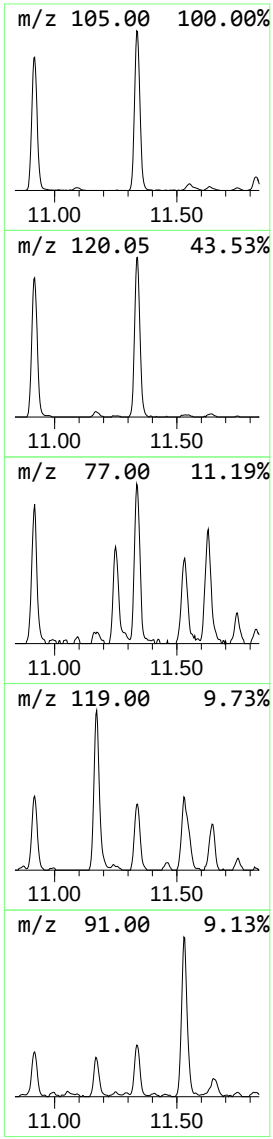
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	11.05 ug/L	218025	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Mesitylene	120	C9H12	000108-67-8	93
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

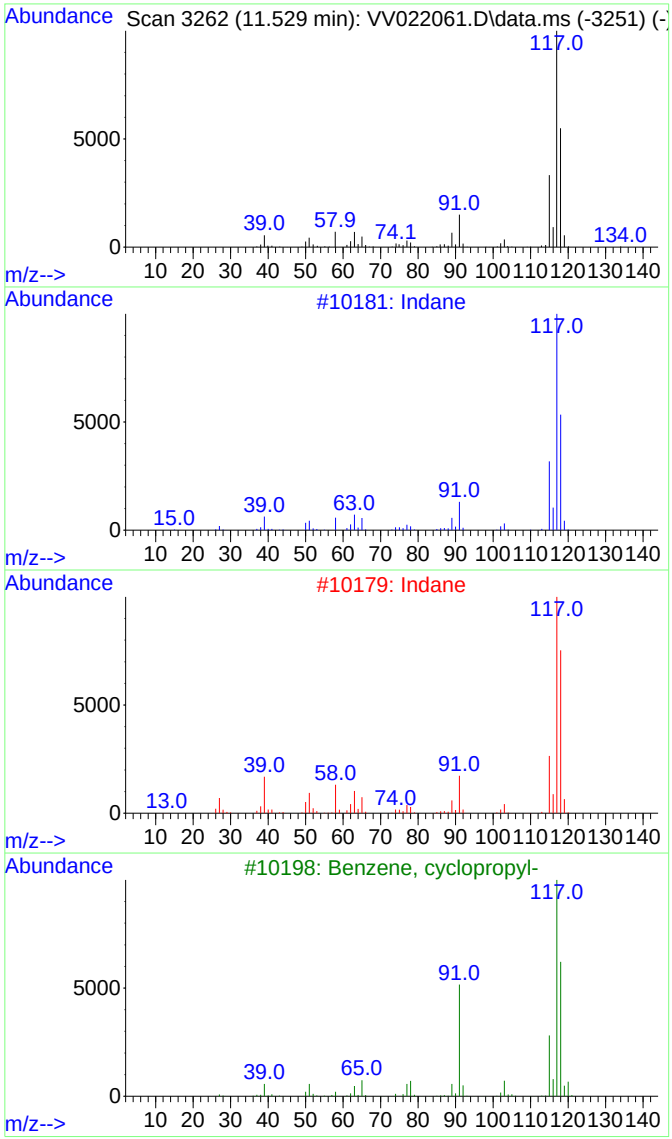
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

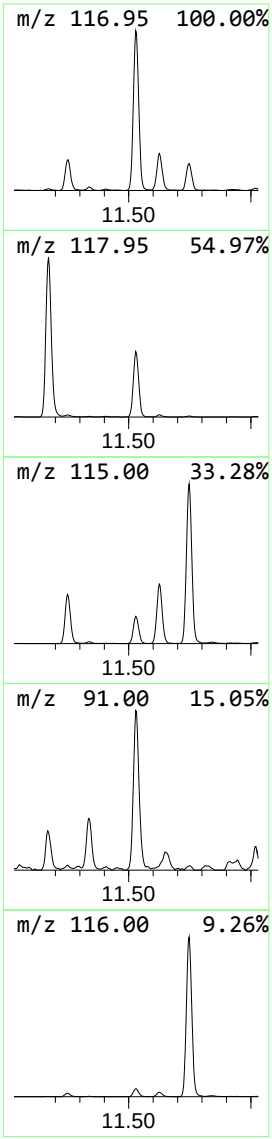
****

Peak Number 5 Indane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	24.70 ug/L	487301	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	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	74
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

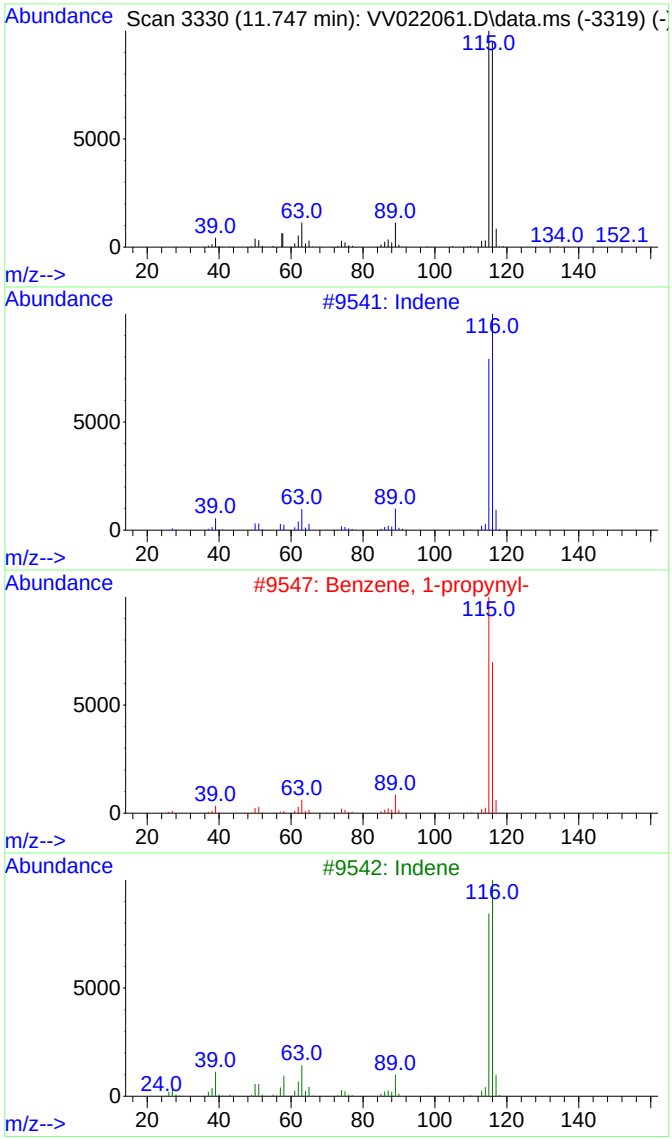
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

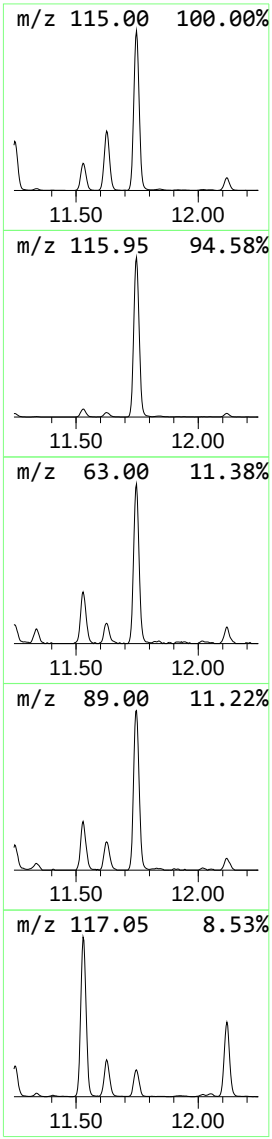
****

Peak Number 6 Indene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.747	46.11 ug/L	909707	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\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

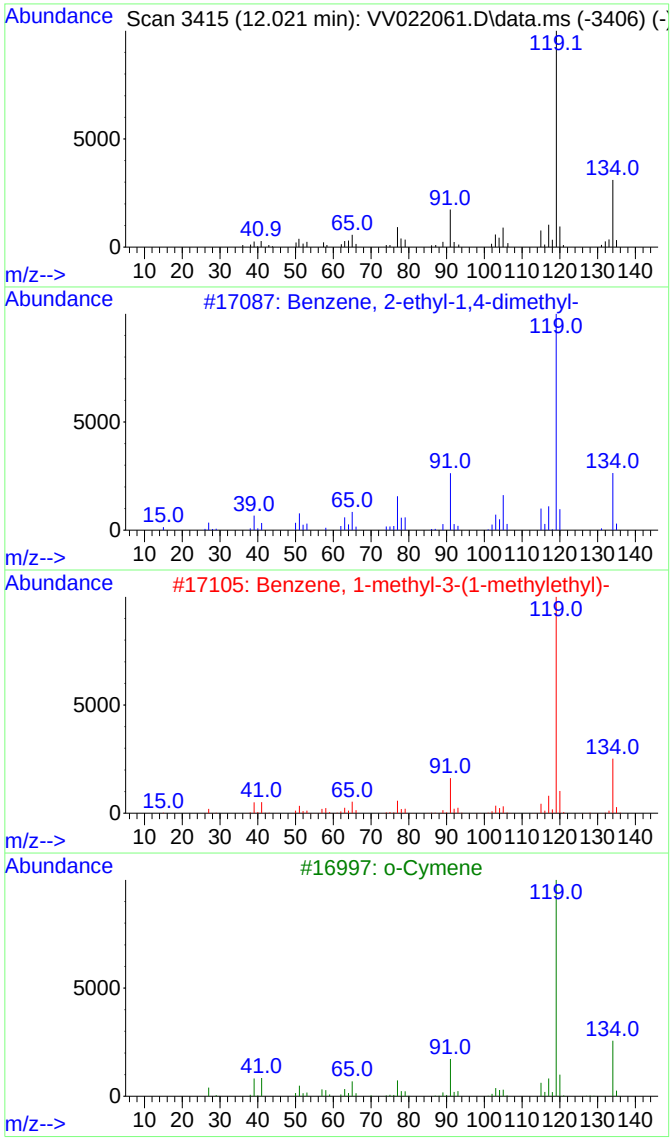
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

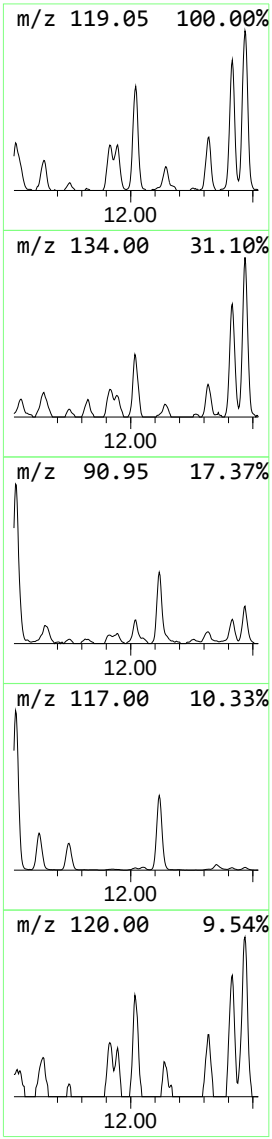
****

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.021	2.68 ug/L	52807	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
3	o-Cymene	134	C10H14	000527-84-4	94
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	p-Cymene	134	C10H14	000099-87-6	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

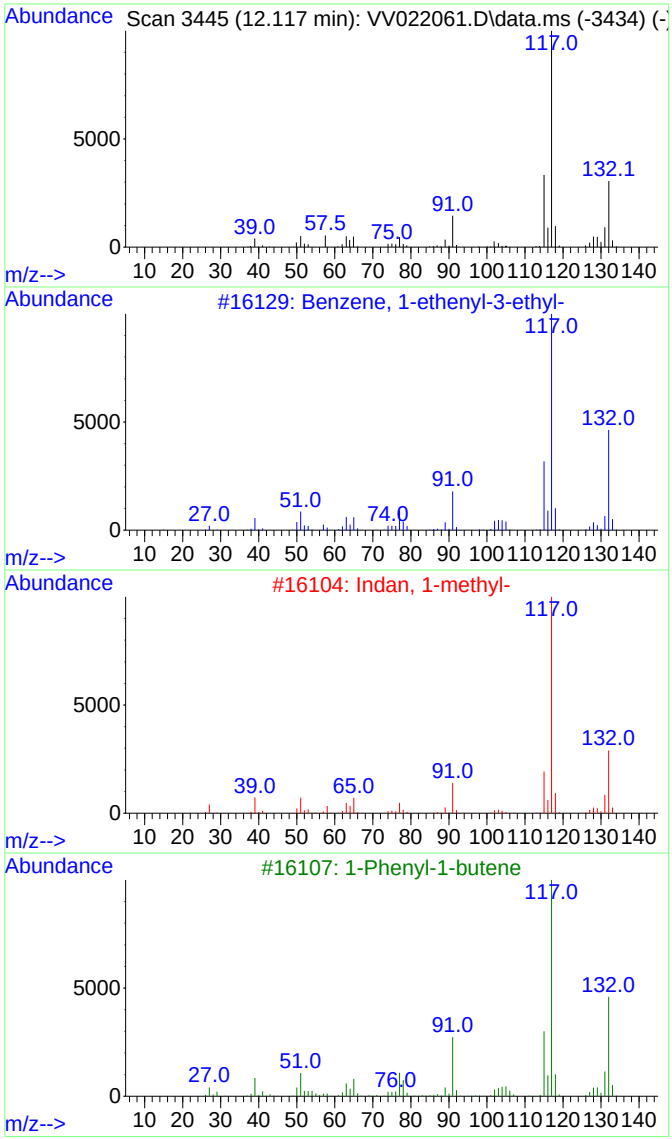
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

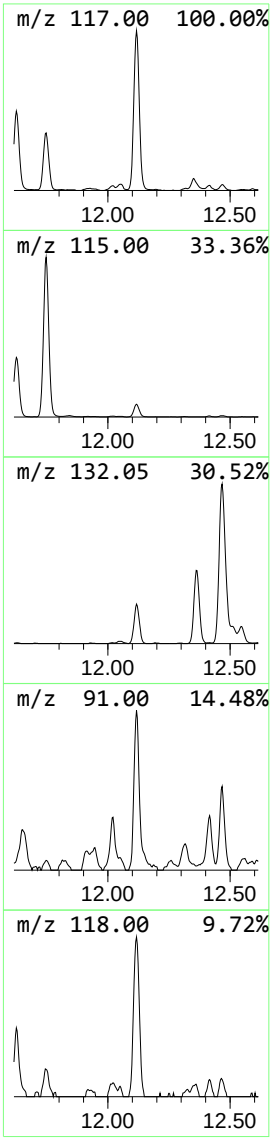
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	11.63 ug/L	229443	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	91
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	86





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

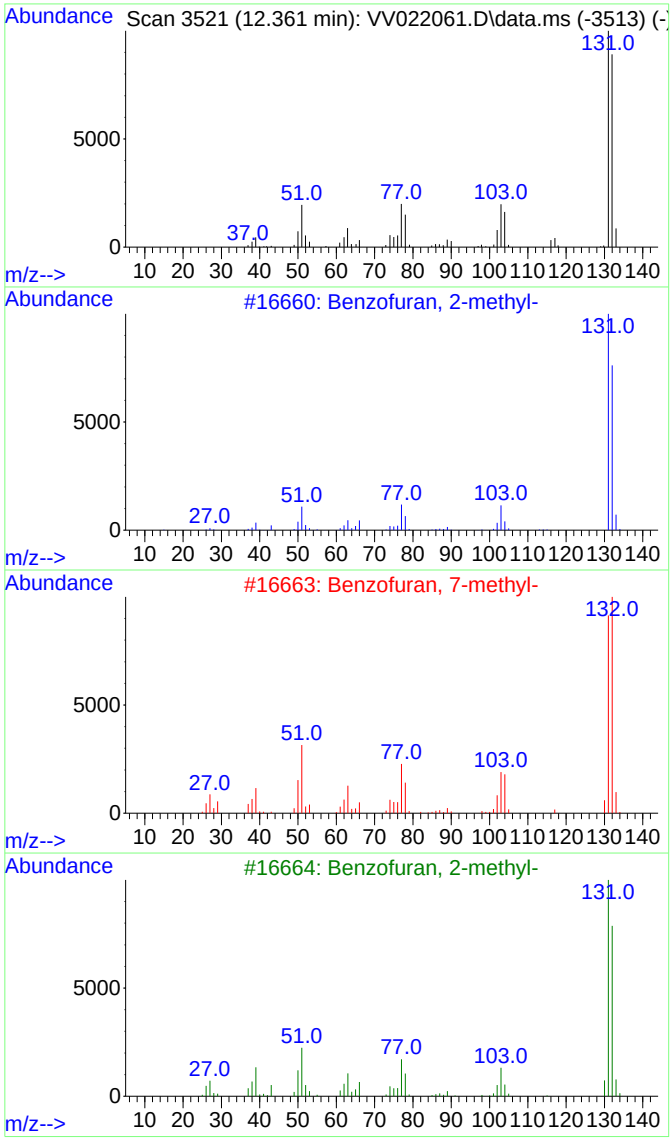
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

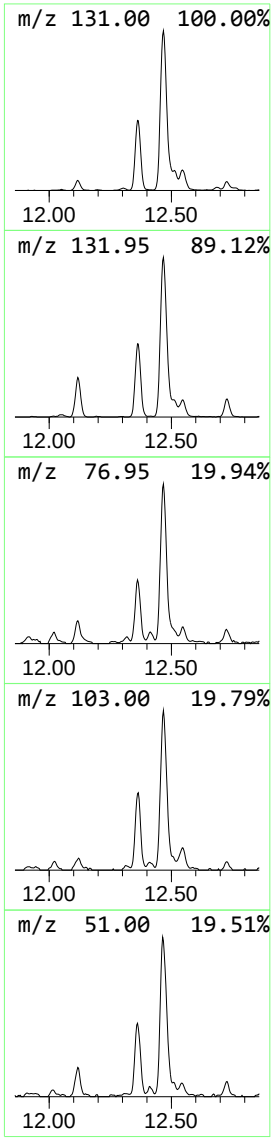
****

Peak Number 9 Benzofuran, 2-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.361	8.70 ug/L	171726	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
4	1H-Indenol	132	C9H8O	056631-57-3	91
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

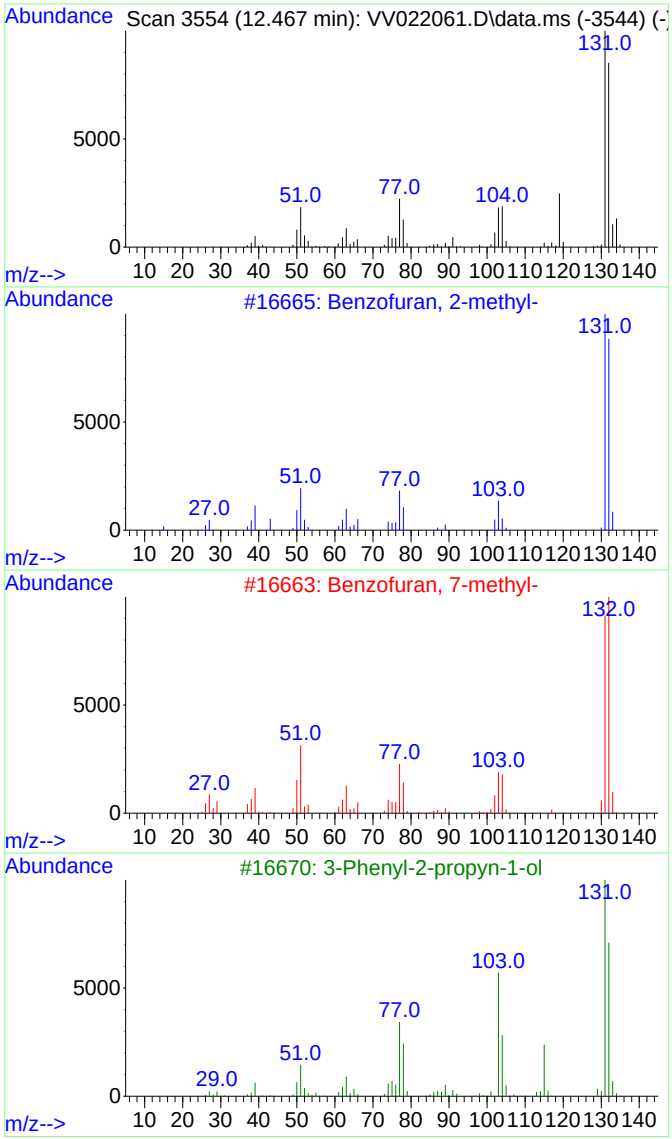
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

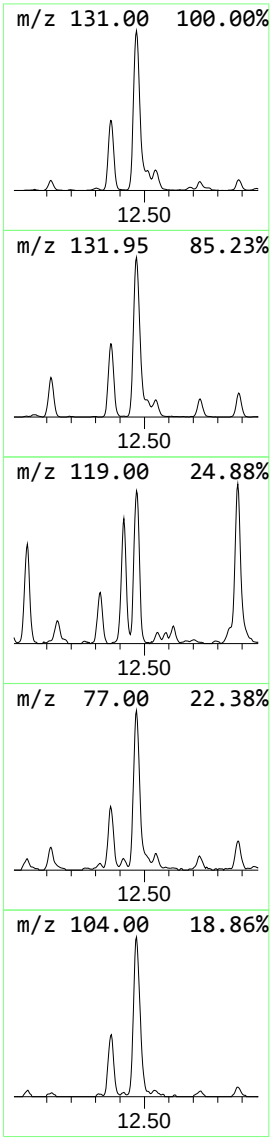
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.467	26.35 ug/L	519945	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	91
3	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	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\VV090121\
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Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

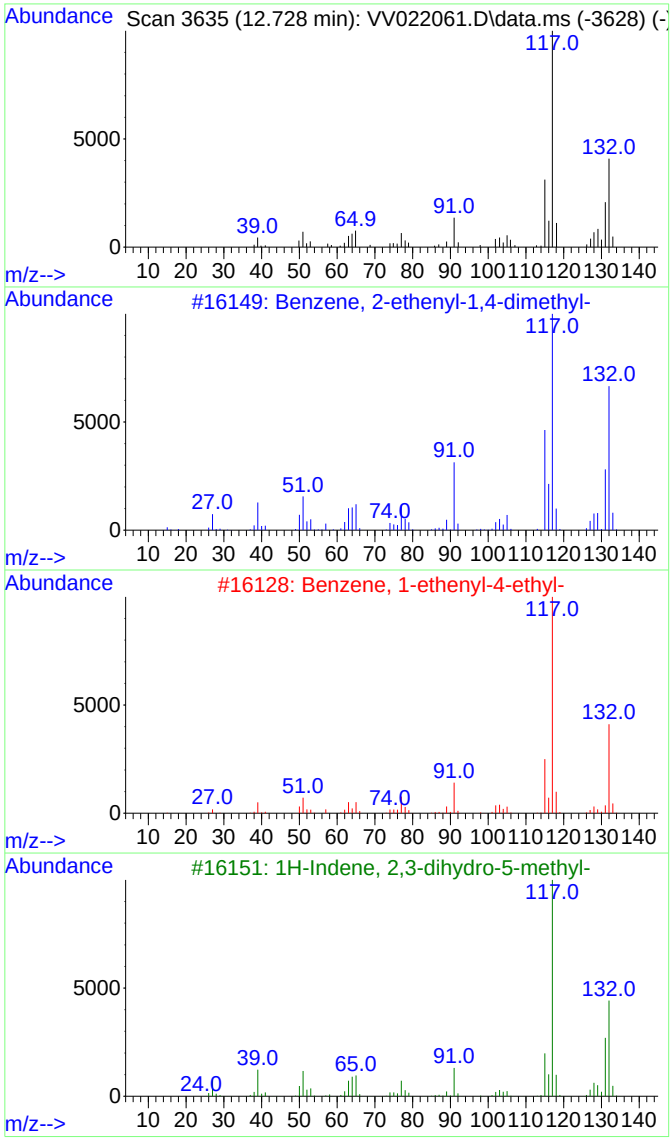
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

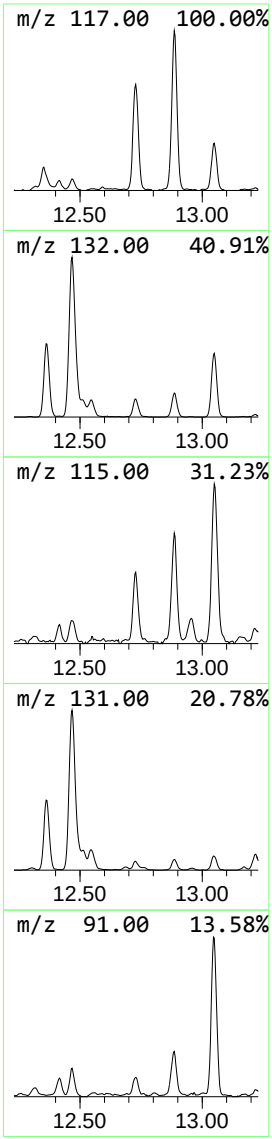
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.728	4.61 ug/L	91053	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

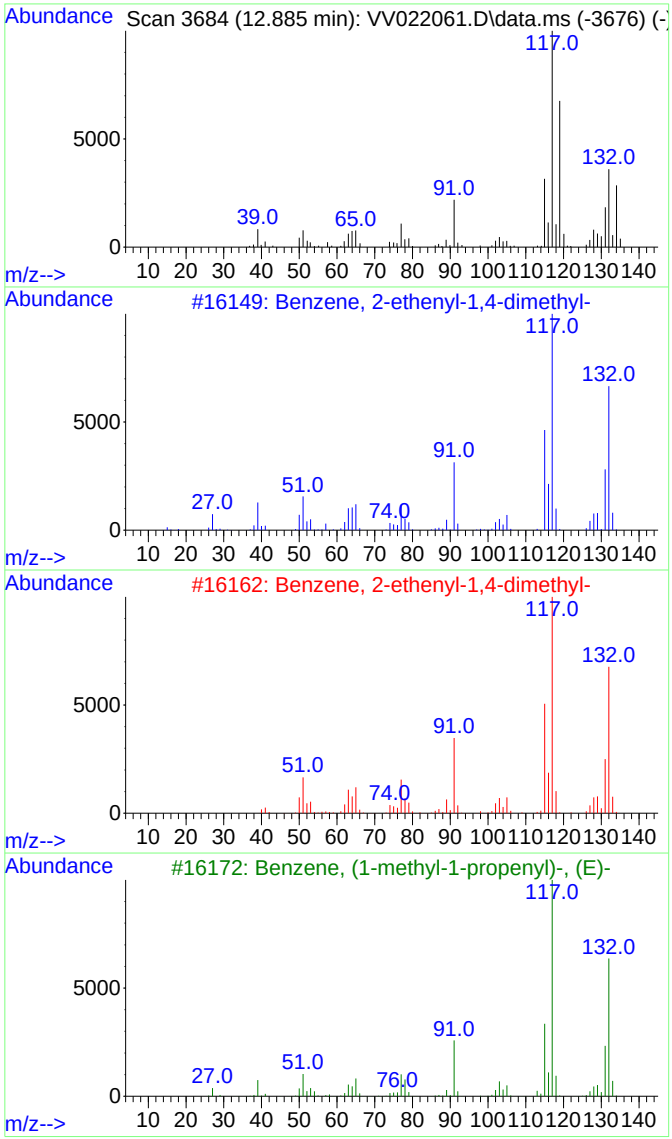
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

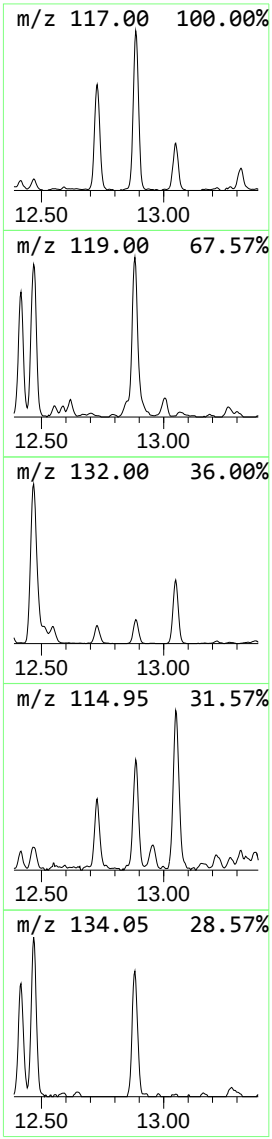
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, (1-methyl-1-propen... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	9.45 ug/L	186413	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

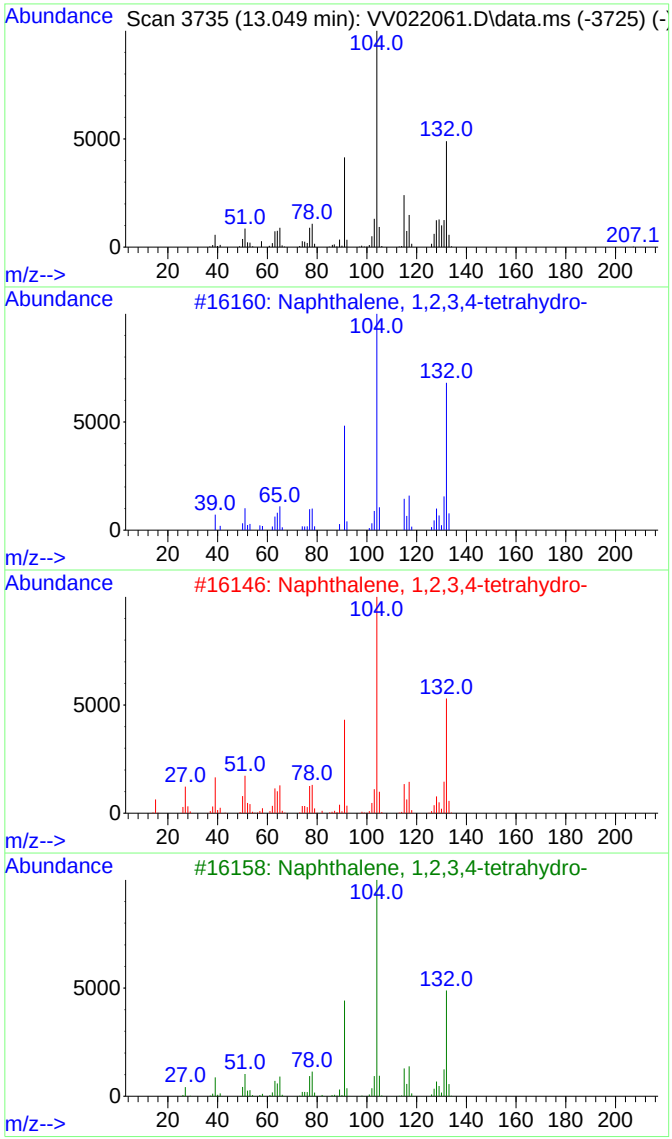
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

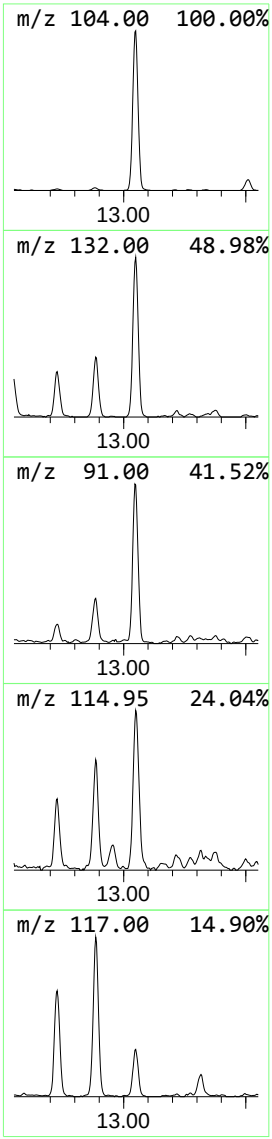
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.050	16.25 ug/L	320631	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	93





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

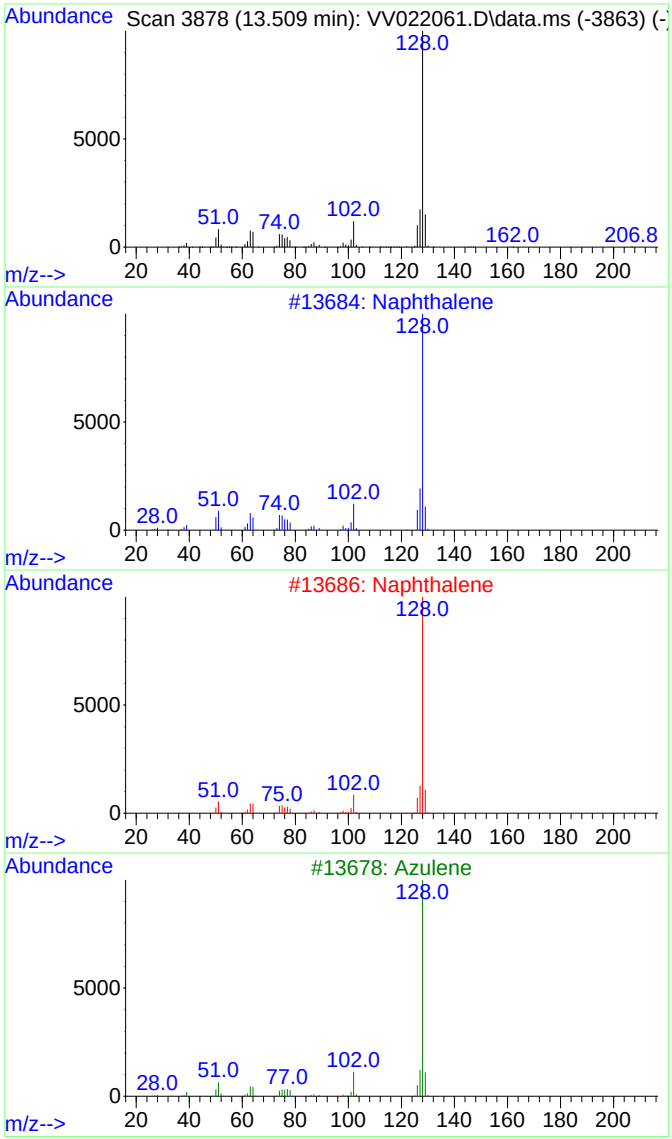
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

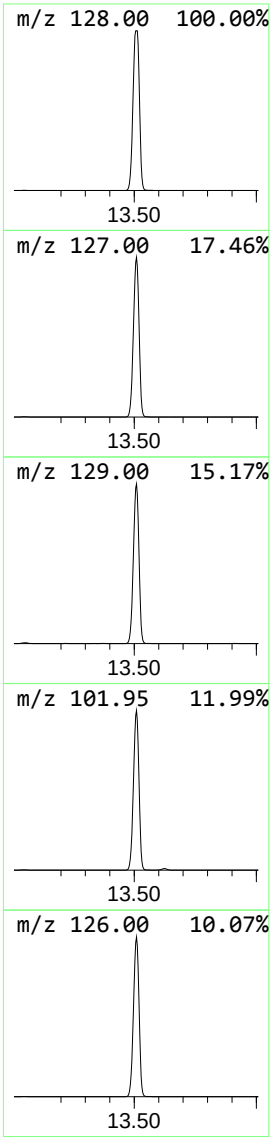
****

Peak Number 14 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.509	1675.63 ug/L	33060000	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	97
2	Naphthalene	128	C10H8	000091-20-3	95
3	Azulene	128	C10H8	000275-51-4	94
4	Azulene	128	C10H8	000275-51-4	91
5	Naphthalene	128	C10H8	000091-20-3	90





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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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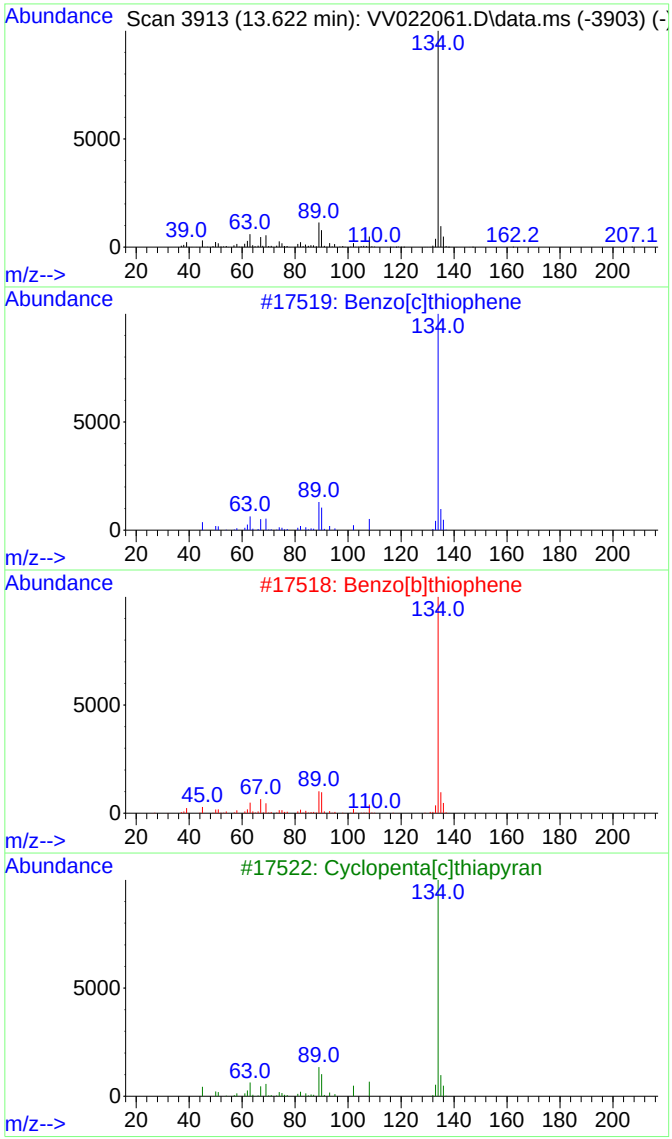
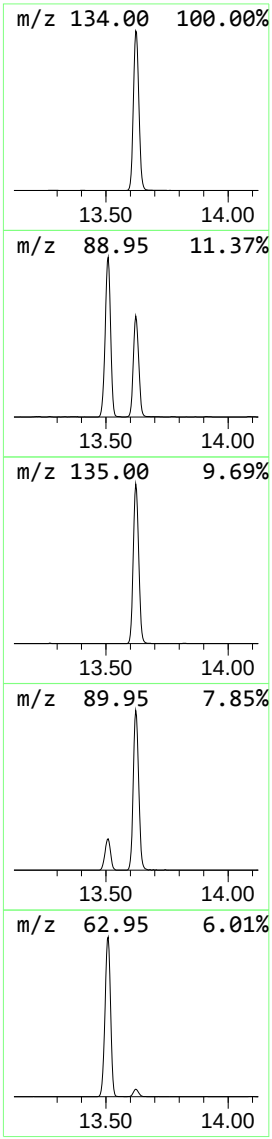
Peak Number 15 Benzo[c]thiophene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	81.46 ug/L	1607230	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	94
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	91





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

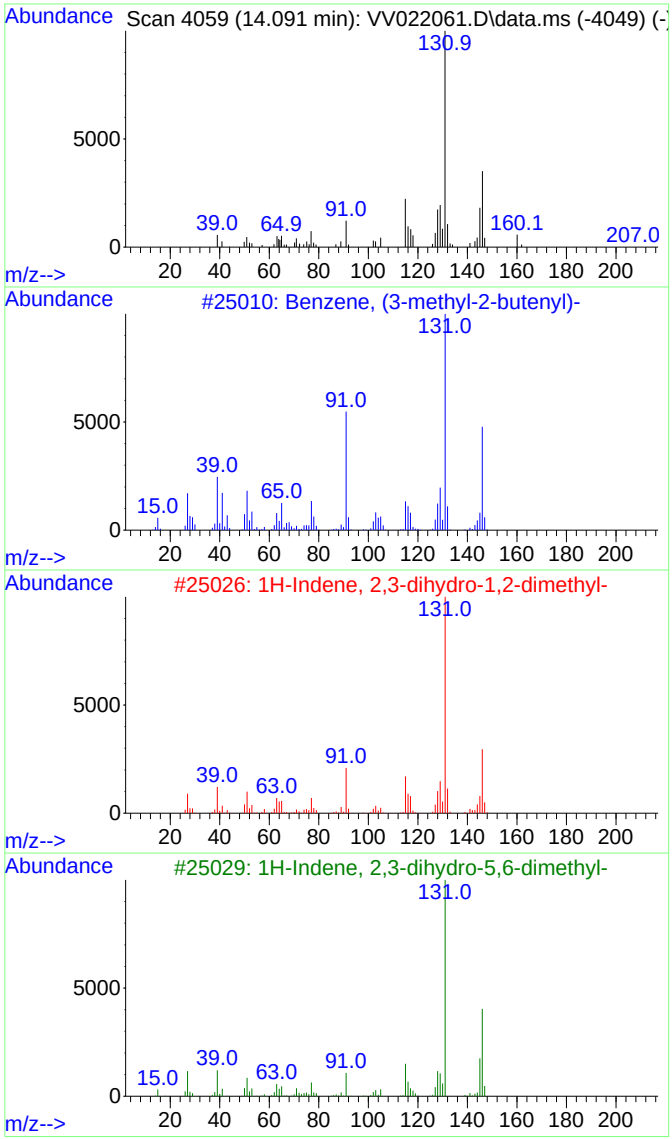
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

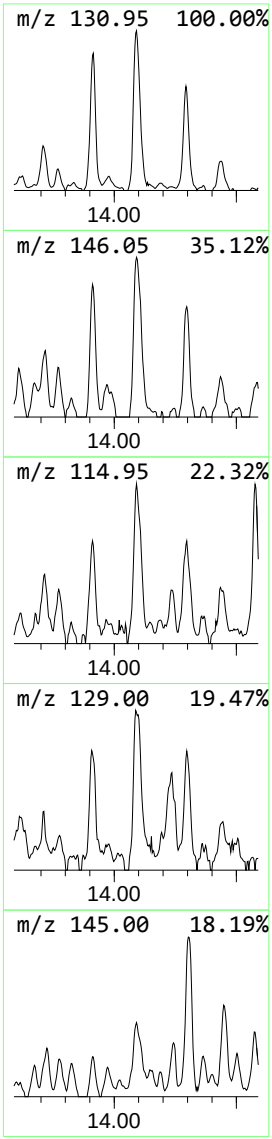
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, (3-methyl-2-butenyl)- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.091	4.36 ug/L	85929	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	87
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	87





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Data File : VV022061.D
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Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

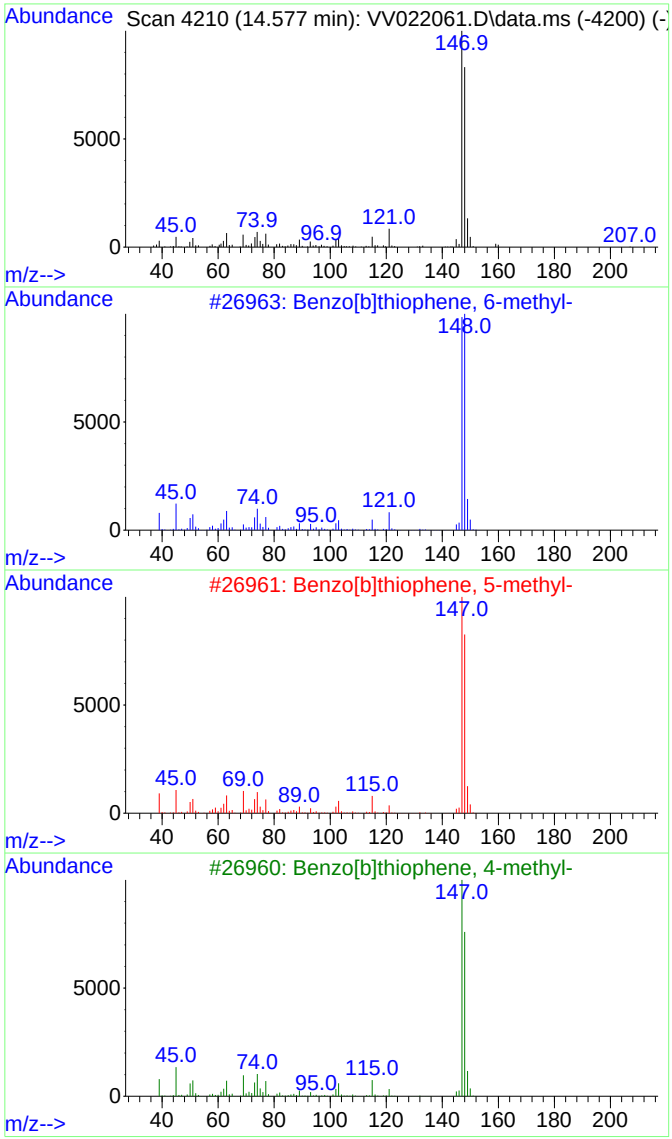
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

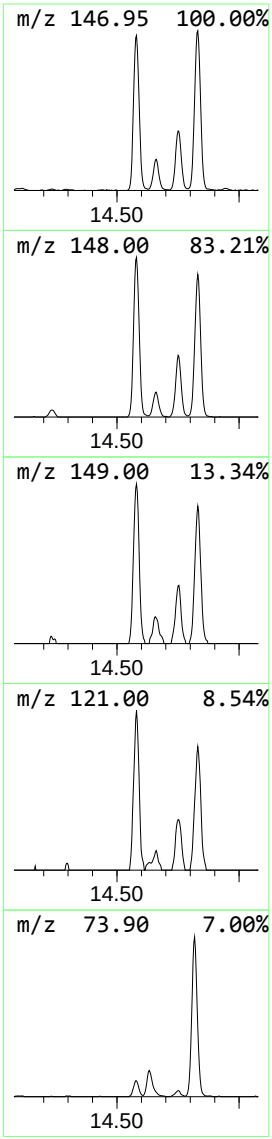
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzo[b]thiophene, 6-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.577	11.97 ug/L	236166	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	94
2	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	91
3	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	91
4	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
5	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

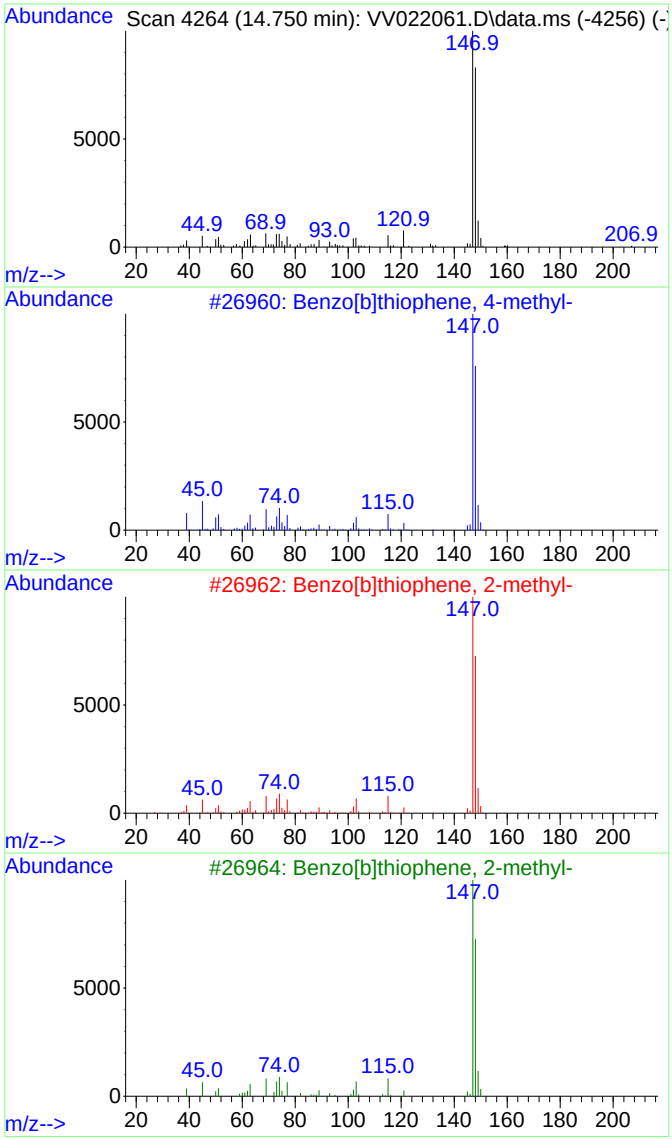
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

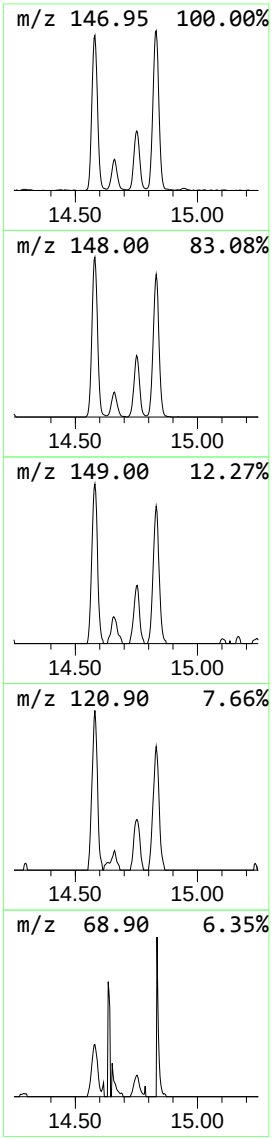
****

Peak Number 19 Benzo[b]thiophene, 4-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.750	4.48 ug/L	88468	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	91
2	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
3	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
4	3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
5	Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

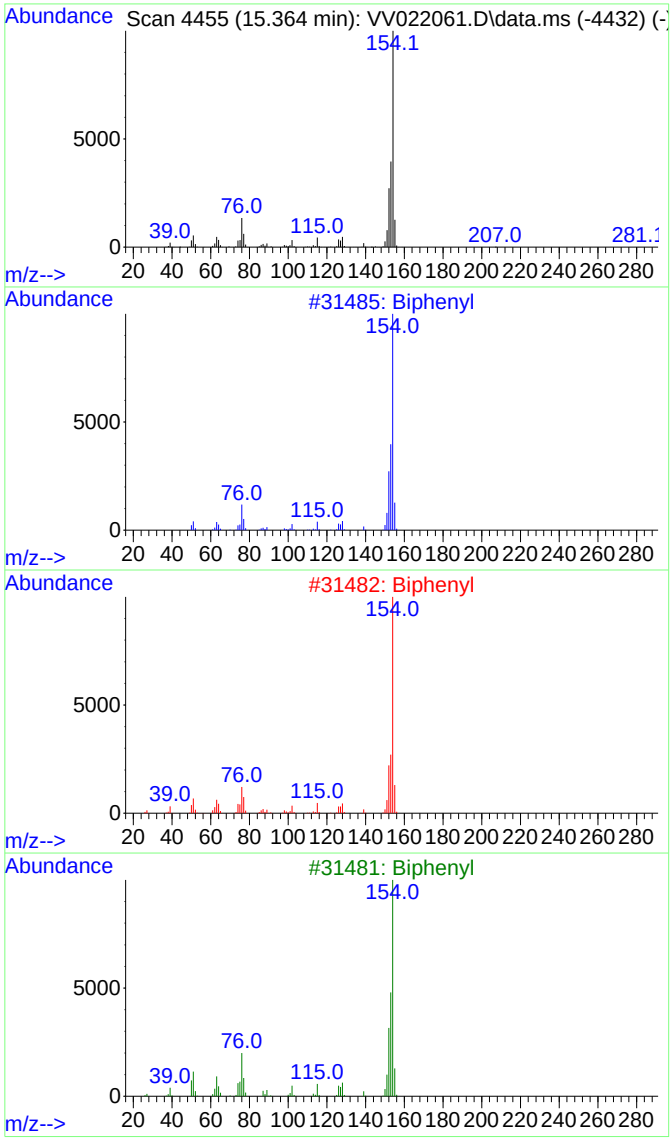
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

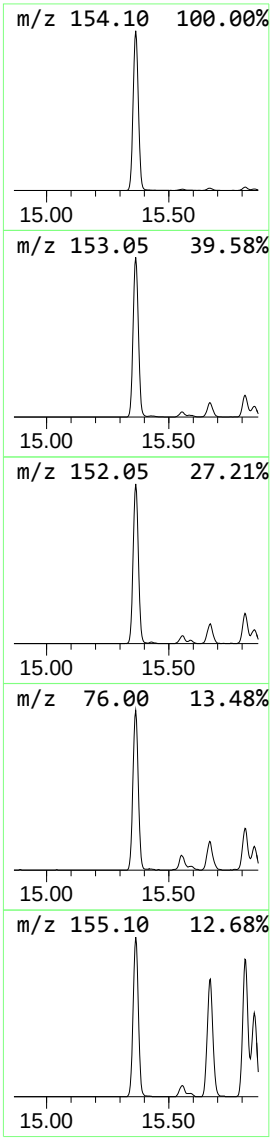
****

Peak Number 21 Biphenyl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.365	54.93 ug/L	1083790	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl	154	C12H10	000092-52-4	95
2	Biphenyl	154	C12H10	000092-52-4	94
3	Biphenyl	154	C12H10	000092-52-4	93
4	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	87
5	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	83





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

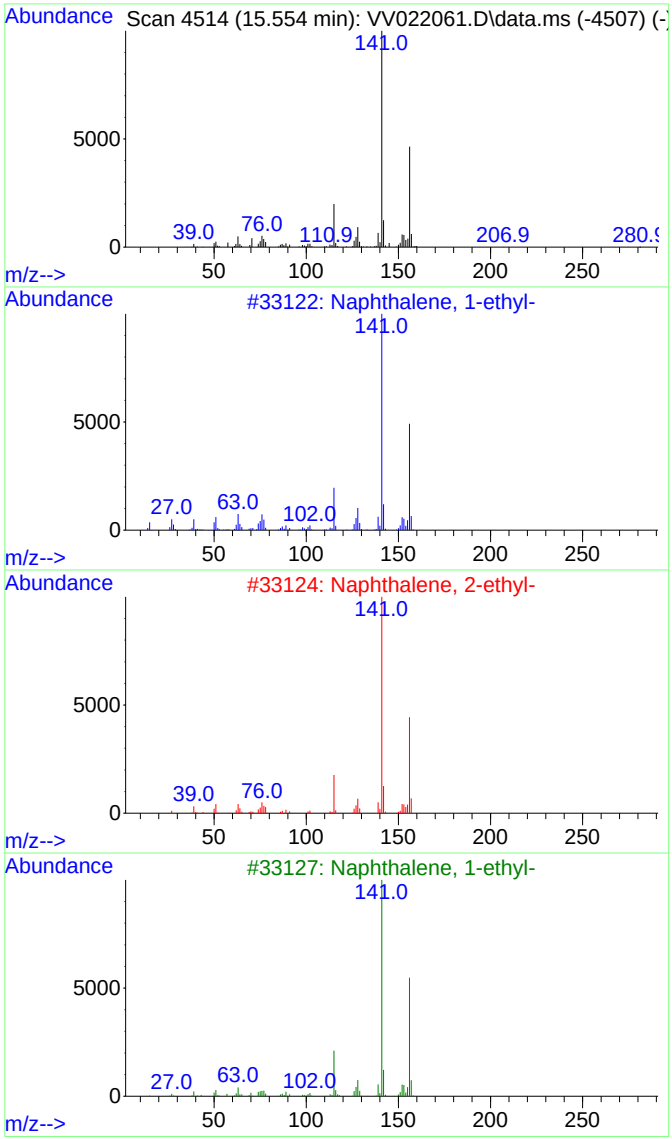
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

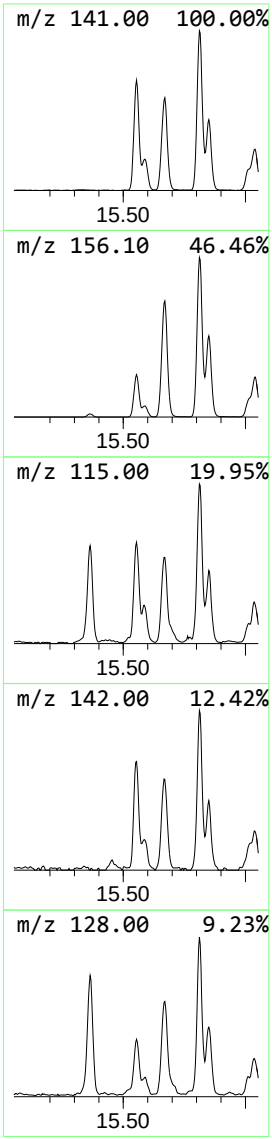
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Naphthalene, 1-ethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.554	10.67 ug/L	210499	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	97
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
4	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

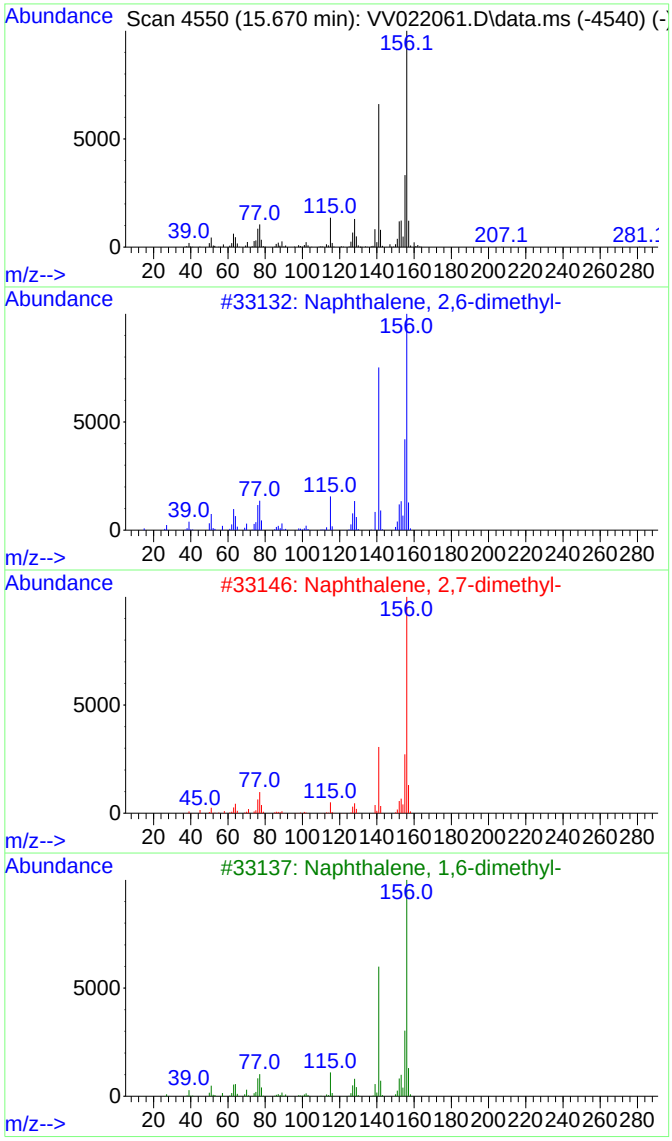
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

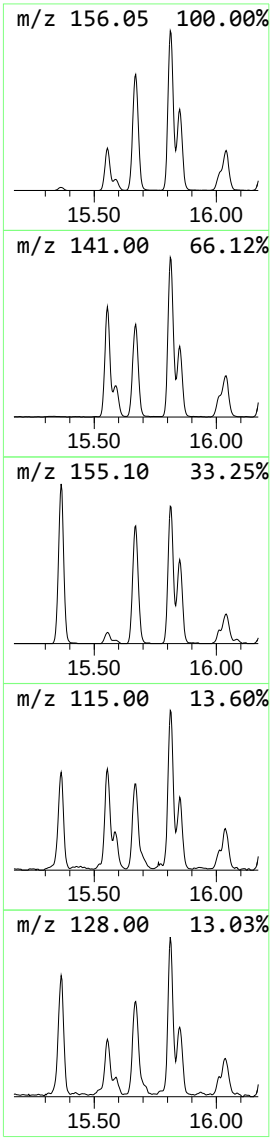
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Naphthalene, 2,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.670	26.18 ug/L	516584	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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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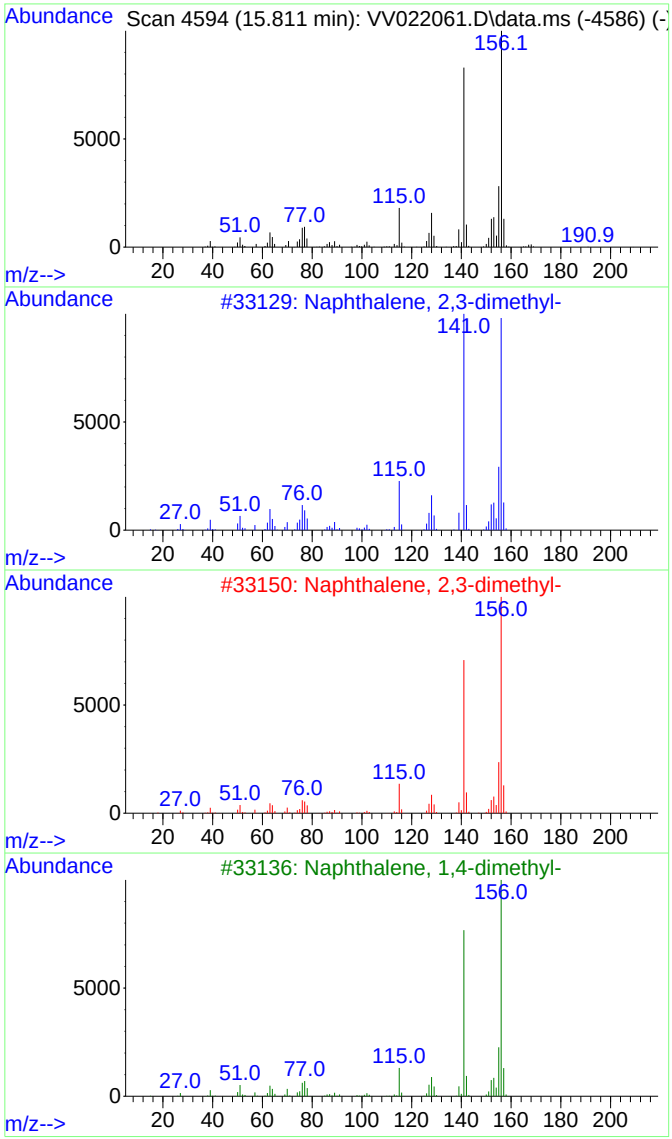
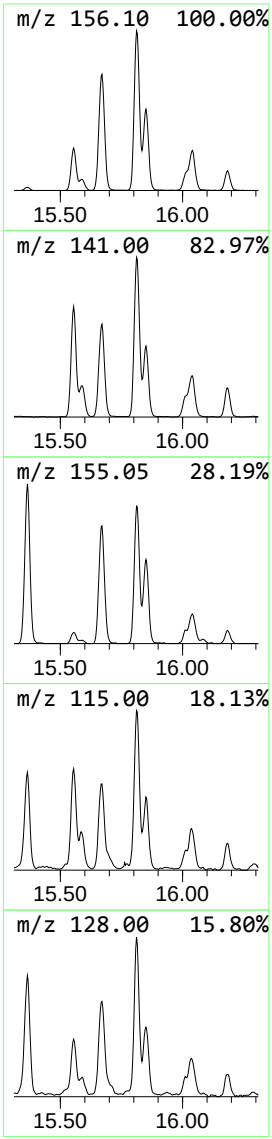
Peak Number 24 Naphthalene, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.811	30.33 ug/L	598381	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

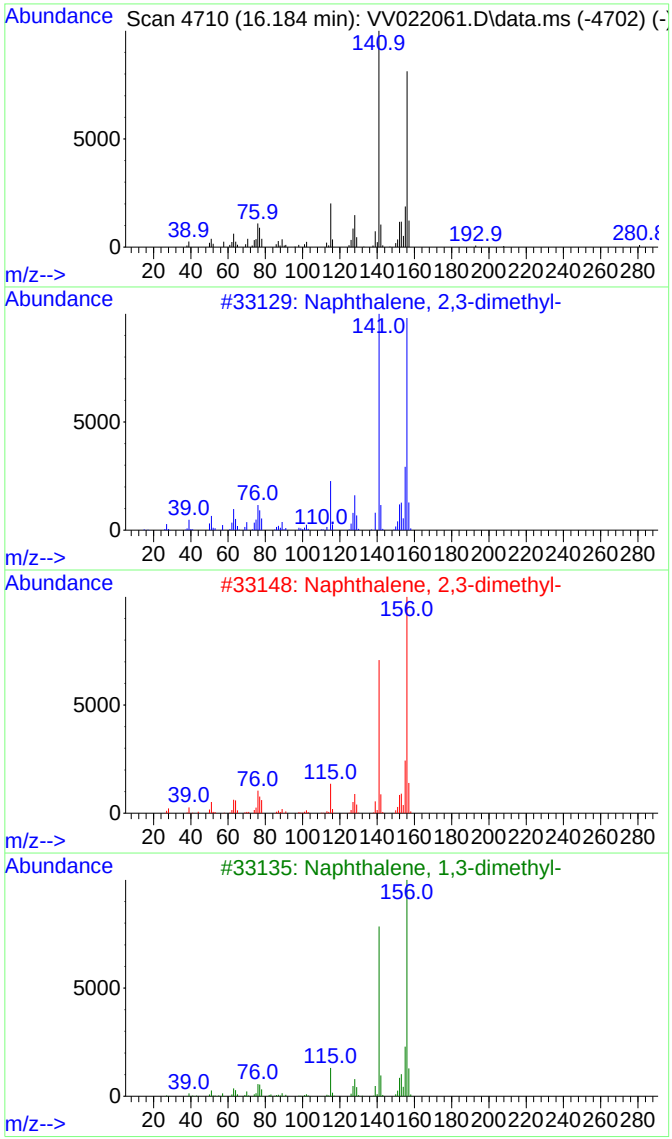
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

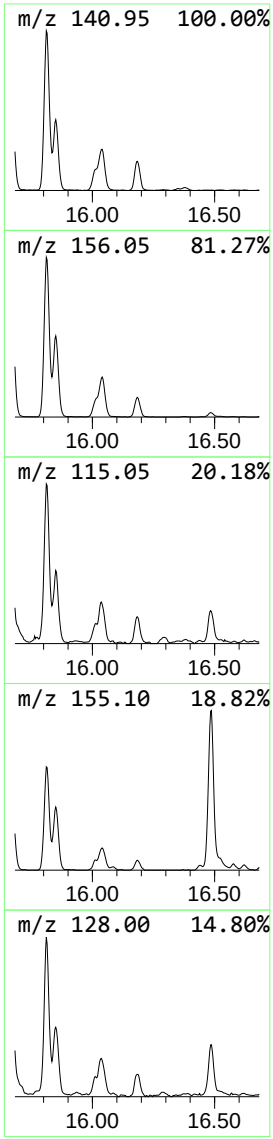
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Naphthalene, 1,3-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.184	4.59 ug/L	90599	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
4	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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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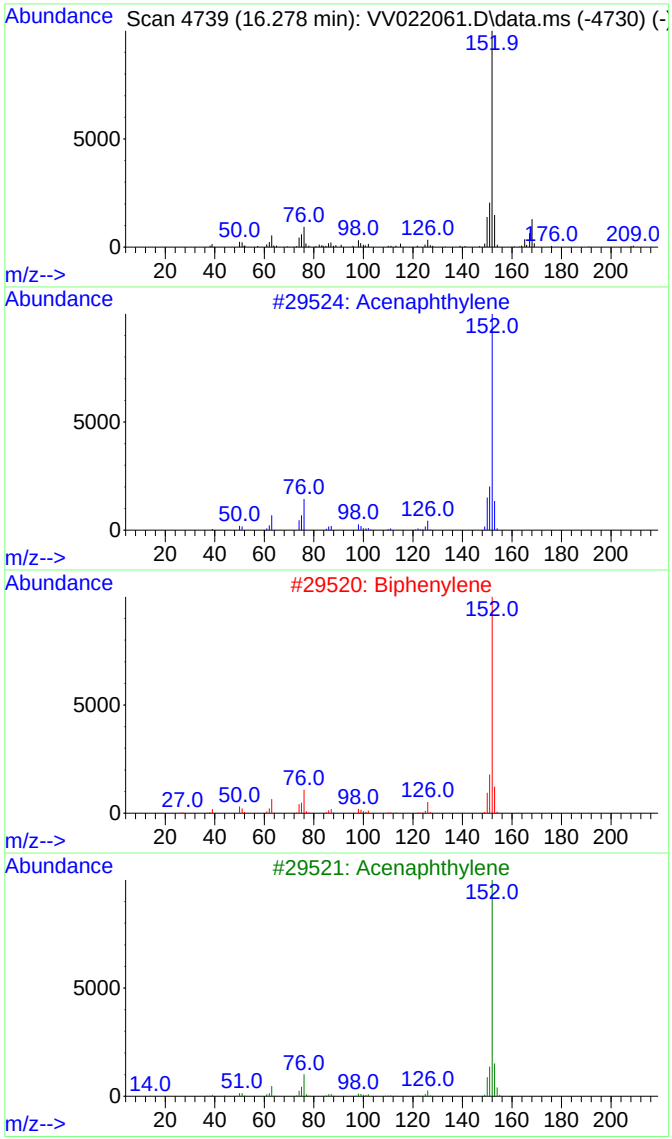
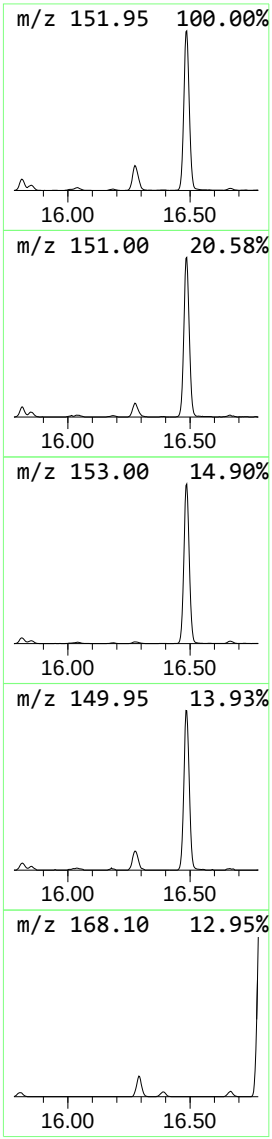
Peak Number 26 Biphenylene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.278	7.14 ug/L	140870	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthylene	152	C12H8	000208-96-8	90
2	Biphenylene	152	C12H8	000259-79-0	90
3	Acenaphthylene	152	C12H8	000208-96-8	81
4	Biphenylene	152	C12H8	000259-79-0	76
5	(6Balpha,7beta,11alpha,11aalpha)...	334	C25H18O	1000241-69-8	64





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Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

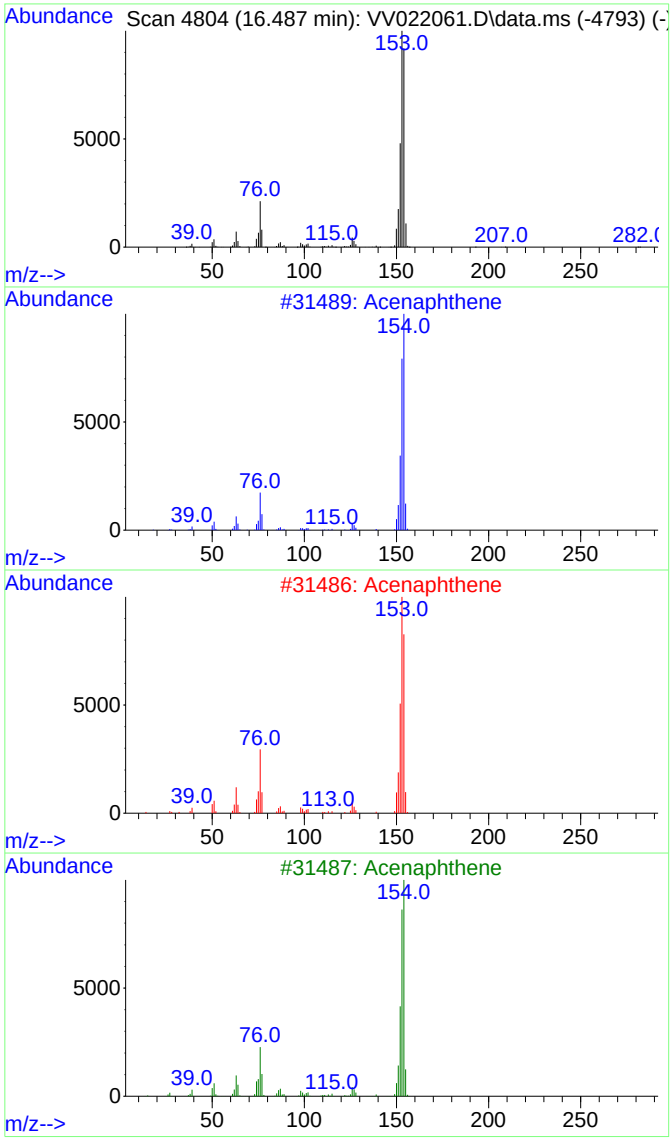
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis

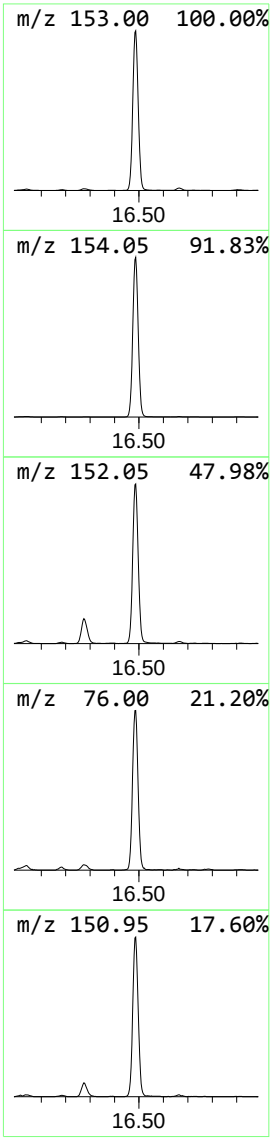
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 2,4,6-Trihydroxybenzaldehyde Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.487	110.42 ug/L	2178630	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthene	154	C12H10	000083-32-9	94
2	Acenaphthene	154	C12H10	000083-32-9	89
3	Acenaphthene	154	C12H10	000083-32-9	72
4	Acenaphthene	154	C12H10	000083-32-9	59
5	2,4,6-Trihydroxybenzaldehyde	154	C7H6O4	000487-70-7	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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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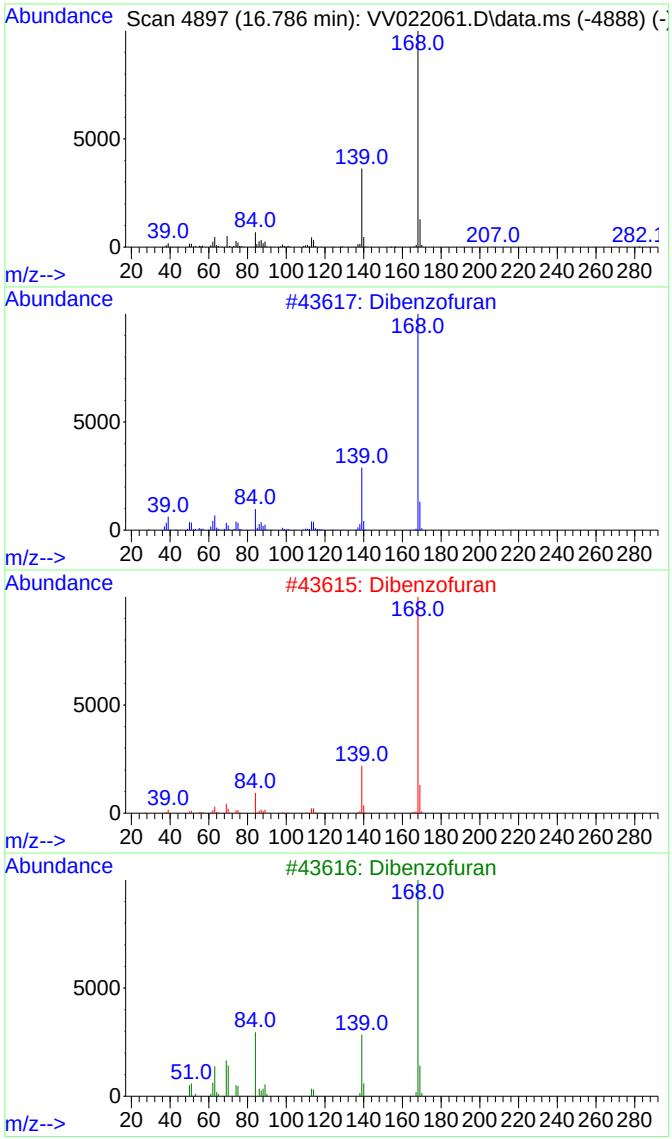
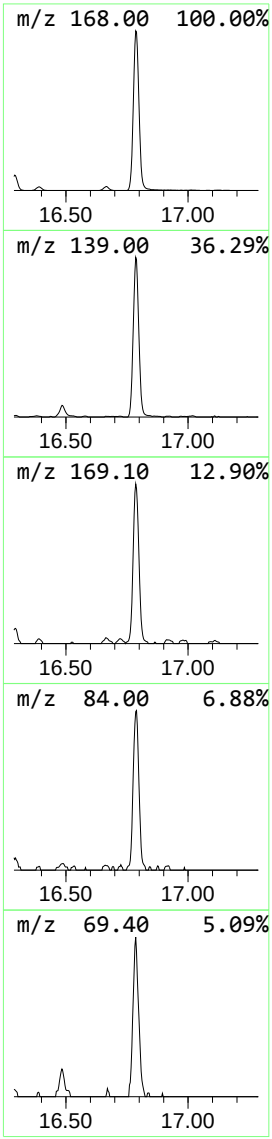
Peak Number 28 Dibenzofuran Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.786	17.44 ug/L	344007	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	91
2	Dibenzofuran	168	C12H8O	000132-64-9	87
3	Dibenzofuran	168	C12H8O	000132-64-9	64
4	3,5-Dimethoxybenzyl alcohol	168	C9H12O3	000705-76-0	56
5	Naphtho[2,1-d]imidazole	168	C11H8N2	000233-53-4	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090121\
Data File : VV022061.D
Acq On : 01 Sep 2021 12:47
Operator : SY/MD
Sample : M3608-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKJ9

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-ethy...	10.741	4.0	ug/L	78471	3	11.252	986491	50.0
Benzofuran	11.172	30.2	ug/L	596374	3	11.252	986491	50.0
Benzene, 1,2,3-...	11.336	11.1	ug/L	218025	3	11.252	986491	50.0
Indane	11.529	24.7	ug/L	487301	3	11.252	986491	50.0
Indene	11.747	46.1	ug/L	909707	3	11.252	986491	50.0
Benzene, 2-ethy...	12.021	2.7	ug/L	52807	3	11.252	986491	50.0
Benzene, 1-ethe...	12.117	11.6	ug/L	229443	3	11.252	986491	50.0
Benzofuran, 2-m...	12.361	8.7	ug/L	171726	3	11.252	986491	50.0
Benzofuran, 7-m...	12.467	26.4	ug/L	519945	3	11.252	986491	50.0
Benzene, 2-ethe...	12.728	4.6	ug/L	91053	3	11.252	986491	50.0
Benzene, (1-met...	12.886	9.4	ug/L	186413	3	11.252	986491	50.0
Naphthalene, 1,...	13.050	16.3	ug/L	320631	3	11.252	986491	50.0
Azulene	13.509	1675.6	ug/L	33060000	3	11.252	986491	50.0
Benzo[c]thiophene	13.622	81.5	ug/L	1607230	3	11.252	986491	50.0
Benzene, (3-met...	14.091	4.4	ug/L	85929	3	11.252	986491	50.0
Benzo[b]thiophe...	14.577	12.0	ug/L	236166	3	11.252	986491	50.0
Benzo[b]thiophe...	14.750	4.5	ug/L	88468	3	11.252	986491	50.0
Biphenyl	15.365	54.9	ug/L	1083790	3	11.252	986491	50.0
Naphthalene, 1-...	15.554	10.7	ug/L	210499	3	11.252	986491	50.0
Naphthalene, 2,...	15.670	26.2	ug/L	516584	3	11.252	986491	50.0
Naphthalene, 2,...	15.811	30.3	ug/L	598381	3	11.252	986491	50.0
Naphthalene, 1,...	16.184	4.6	ug/L	90599	3	11.252	986491	50.0
Biphenylene	16.278	7.1	ug/L	140870	3	11.252	986491	50.0
2,4,6-Trihydrox...	16.487	110.4	ug/L	2178630	3	11.252	986491	50.0
Dibenzofuran	16.786	17.4	ug/L	344007	3	11.252	986491	50.0