LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKF3

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\SFAMVLM081621WMA.M
Title : VOC Analysis

Signal : TIC: VV021892.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.192	40	47	75	rBV	98036	163663	5.79%	0.874%
2	1.304	76	82	88	rBV	204786	205045	7.26%	1.094%
3	1.561	156	162	180	rVB	160510	184957	6.55%	0.987%
4	2.105	322	331	355	rVB	324003	491144	17.39%	2.622%
5	2.609	484	488	489	rBV3	726	455	0.02%	0.002%
6	2.764	532	536	543	rBV7	1373	1678	0.06%	0.009%
7	2.851	560	563	568	rVB2	360	251	0.01%	0.001%
8	2.925	584	586	589	rVB2	529	278	0.01%	0.001%
9	2.966	597	599	603	rBV3	344	212	0.01%	0.001%
10	3.117	644	646	651	rVB2	409	272	0.01%	0.001%
11	3.195	662	670	677	rBV4	527	787	0.03%	0.004%
12	3.243	681	685	688	rVB2	395	295	0.01%	0.002%
13	3.352	716	719	721	rBV2	324	205	0.01%	0.001%
14	3.388	727	730	732	rBV2	267	216	0.01%	0.001%
15	3.597	792	795	800	rVB2	337	220	0.01%	0.001%
16	3.638	804	808	813	rVB2	375	319	0.01%	0.002%
17	3.732	835	837	842	rVB2	434	301	0.01%	0.002%
18	3.757	842	845	848	rBV	352	207	0.01%	0.001%
19	3.780	848	852	857	rVB3	444	464	0.02%	0.002%
20	3.876	872	882	915	rBV	118841	322920	11.43%	1.724%
21	4.352	1013	1030	1057	rVV	234792	585668	20.74%	3.126%
22	4.789	1160	1166	1168	rBV2	349	300	0.01%	0.002%
23	5.047	1229	1246	1285	rBV2	589695	1508241	53.40%	8.050%
24	5.484	1378	1382	1383	rBV	496	335	0.01%	0.002%
25	5.494	1383	1385	1388	rVV2	706	380	0.01%	0.002%
26	5.532	1392	1397	1401	rVB3	564	537	0.02%	0.003%
27	5.619	1411	1424	1460	rBV	494770	996145	35.27%	5.317%
28	5.915	1505	1516	1528	rBV7	4697	10335	0.37%	0.055%
29	6.008	1542	1545	1551	rBV3	325	367	0.01%	0.002%
30	6.072	1551	1565	1594	rBV	329271	673981	23.86%	3.597%
31	6.320	1639	1642	1645	rVB3	755	423	0.01%	0.002%
32	6.378	1655	1660	1661	rBV3	367	270	0.01%	0.001%
33	6.503	1695	1699	1702	rBV	550	287	0.01%	0.002%
34	6.526	1702	1706	1709	rBV2	518	460	0.02%	0.002%
35	6.574	1717	1721	1725	rBV3	307	335	0.01%	0.002%
36	6.641	1736	1742	1747	rBV4	518	657	0.02%	0.004%

LSC Area Percent Report

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ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
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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 >
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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Title : VOC Analysis

37	6.989	1839	1850	1872	rBV	186244	343752	12.17%	1.835%
38	7.185	1909	1911	1913	rBV2	476	217	0.01%	0.001%
39	7.320	1940	1953	1970	rBV	729555	1269473	44.95%	6.776%
40	7.625	2036	2048	2065	rBV	123809	215616	7.63%	1.151%
41	7.802	2101	2103	2106	rBV2	451	287	0.01%	0.002%
42	7.870	2121	2124	2127	rVB3	427	281	0.01%	0.001%
43	7.892	2127	2131	2133	rBV3	560	417	0.01%	0.002%
44	7.944	2137	2147	2159	rVV4	4914	8945	0.32%	0.048%
45	8.088	2182	2192	2229	rBV	354702	629525	22.29%	3.360%
46	8.442	2300	2302	2305	rVB2	569	224	0.01%	0.001%
47	8.500	2316	2320	2324	rBV2	338	372	0.01%	0.002%
48	8.526	2327	2328	2331	rBV3	382	254	0.01%	0.001%
49	8.612	2353	2355	2357	rBV2	542	274	0.01%	0.001%
50	8.722	2385	2389	2390	rBV2	471	277	0.01%	0.001%
51	8.854	2418	2430	2461	rBV	727134	1200162	42.49%	6.406%
52	9.021	2474	2482	2503	rVB3	7268	12805	0.45%	0.068%
53	9.149	2514	2522	2534	rVB4	7104	12281	0.43%	0.066%
54	9.336	2575	2580	2582	rVV2	506	437	0.02%	0.002%
55	9.551	2637	2647	2656	rBV4	5210	8994	0.32%	0.048%
56	9.683	2681	2688	2691	rBV4	456	395	0.01%	0.002%
57	9.783	2715	2719	2725	rBV4	279	365	0.01%	0.002%
58	9.844	2736	2738	2743	rVB3	553	336	0.01%	0.002%
59	9.937	2758	2767	2777	rBV2	6059	10453	0.37%	0.056%
60	10.091	2813	2815	2818	rVV2	280	202	0.01%	0.001%
61	10.130	2818	2827	2836	rVV4	4021	6336	0.22%	0.034%
62	10.217	2842	2854	2875	rBV	489660	776588	27.49%	4.145%
63	10.455	2921	2928	2929	rBV4	1074	1158	0.04%	0.006%
64	10.519	2944	2948	2949	rBV	430	254	0.01%	0.001%
65	10.545	2949	2956	2961	rBV5	1717	2559	0.09%	0.014%
66	10.661	2990	2992	2999	rVB4	489	477	0.02%	0.003%
67	10.744	3008	3018	3031	rVB3	9298	14670	0.52%	0.078%
68	10.860	3052	3054	3057	rBV2	494	223	0.01%	0.001%
69	10.882	3059	3061	3064	rBV2	385	251	0.01%	0.001%
70	10.921	3064	3073	3084	rVV3	7451	11285	0.40%	0.060%
71	11.037	3102	3109	3119	rVB6	1276	2205	0.08%	0.012%
72	11.091	3123	3126	3132	rBV3	717	577	0.02%	0.003%
73	11.146	3138	3143	3146	rBV4	1698	2075	0.07%	0.011%
74	11.249	3163	3175	3197	rBV	795061	1243394	44.02%	6.637%
75	11.529	3246	3262	3279	rBV	1852246	2824485	100.00%	15.076%
76	11.625	3281	3292	3312	rVB	765332	1201049	42.52%	6.411%

LSC Area Percent Report

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ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKF3

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Title : VOC Analysis

77	12.024	3410	3416	3417	rVV2	1574	1548	0.05%	0.008%
78	12.050	3418	3424	3431	rVV5	8676	12414	0.44%	0.066%
79	12.159	3456	3458	3461	rVB3	1230	496	0.02%	0.003%
80	12.233	3479	3481	3482	rBV	567	234	0.01%	0.001%
81	12.300	3498	3502	3503	rBV4	1063	745	0.03%	0.004%
82	12.365	3514	3522	3532	rBV4	8681	14754	0.52%	0.079%
83	12.590	3590	3592	3593	rBV	572	271	0.01%	0.001%
84	12.654	3608	3612	3613	rBV2	627	383	0.01%	0.002%
85	12.725	3625	3634	3646	rBV2	17886	27542	0.98%	0.147%
86	12.886	3674	3684	3697	rBV3	48065	74813	2.65%	0.399%
87	13.050	3726	3735	3747	rBV3	20595	33159	1.17%	0.177%
88	13.159	3760	3769	3772	rBV5	1235	1564	0.06%	0.008%
89	13.503	3865	3876	3903	rVB	1222136	1878876	66.52%	10.029%
90	14.294	4115	4122	4133	rVB3	10369	16097	0.57%	0.086%
91	14.580	4205	4211	4218	rVB2	18416	23297	0.82%	0.124%
92	14.635	4218	4228	4243	rBV	233597	358065	12.68%	1.911%
93	14.818	4275	4285	4300	rBV	193981	311123	11.02%	1.661%
94	15.059	4358	4360	4363	rBV2	992	632	0.02%	0.003%
95	15.365	4446	4455	4466	rVB	64171	100555	3.56%	0.537%
96	15.670	4542	4550	4570	rVB3	31668	63836	2.26%	0.341%
97	15.812	4586	4594	4601	rBV2	45763	71590	2.53%	0.382%
98	16.278	4730	4739	4752	rVB	37433	70150	2.48%	0.374%
99	16.487	4793	4804	4825	rBV	369506	583202	20.65%	3.113%
100	16.789	4890	4898	4918	rVB	82922	132869	4.70%	0.709%

Sum of corrected areas: 18734755

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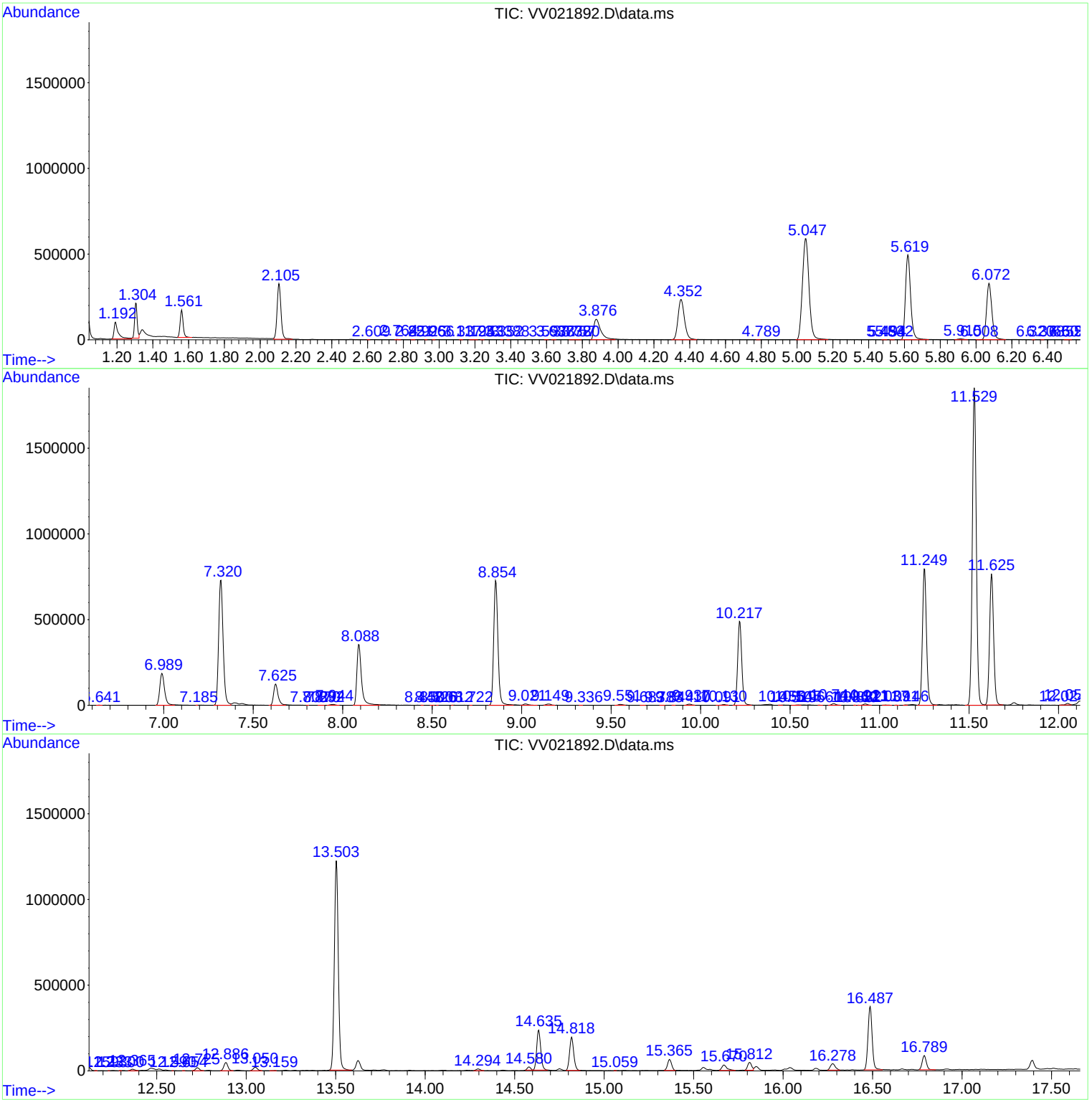
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

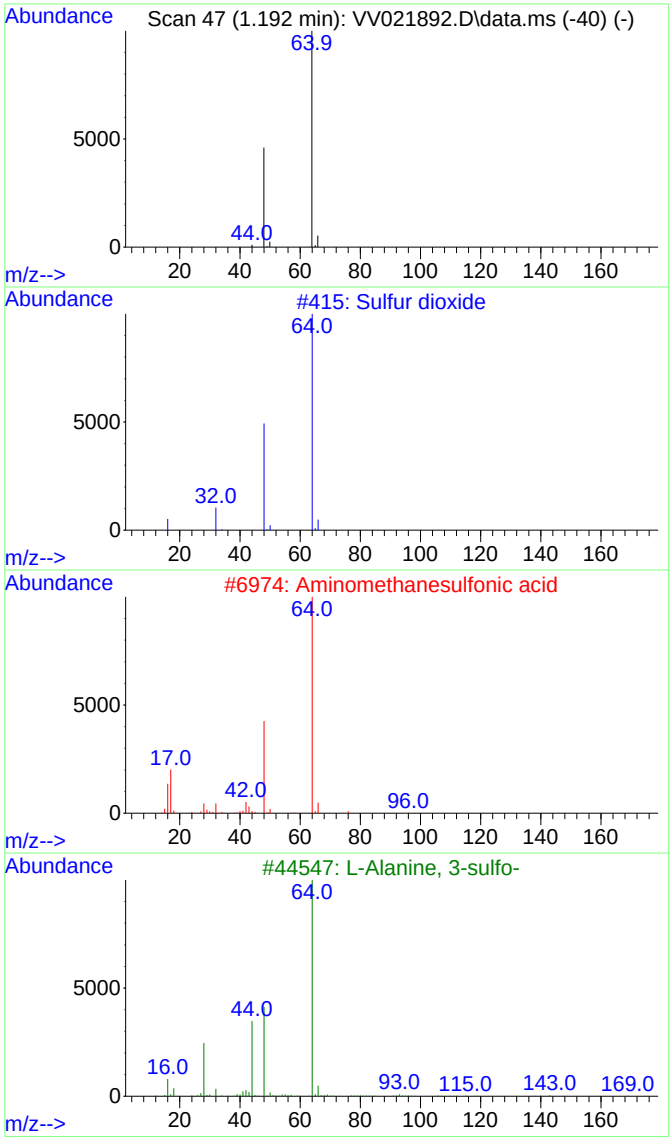
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

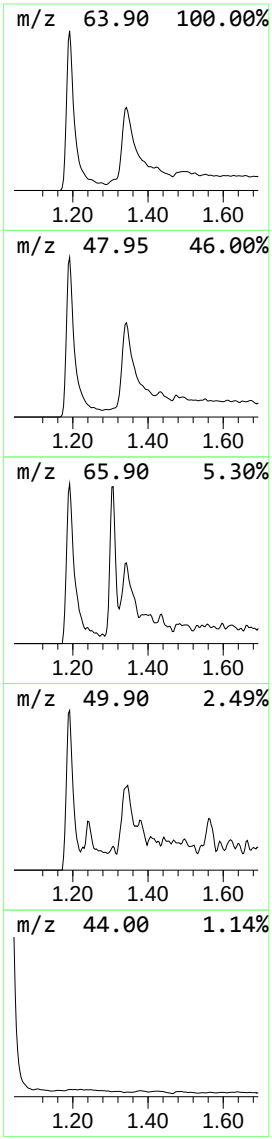
****

Peak Number 1 Sulfur dioxide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.192	8.21 ug/L	163663	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
5	Sulfur dioxide	64	O2S	007446-09-5	5





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

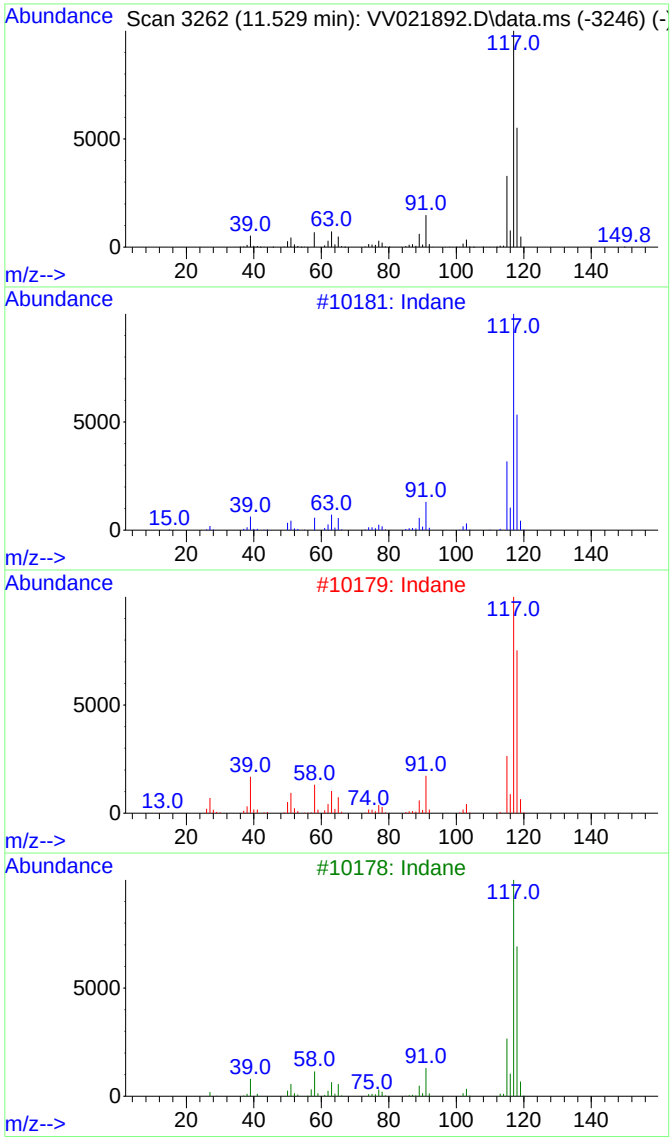
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

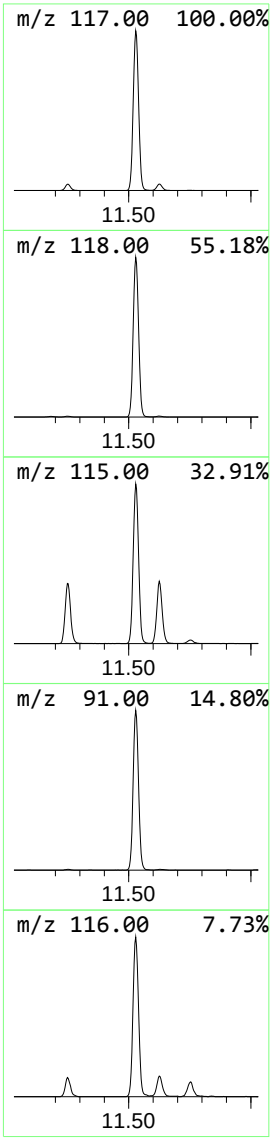
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	113.58 ug/L	2824490	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	91
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64





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ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

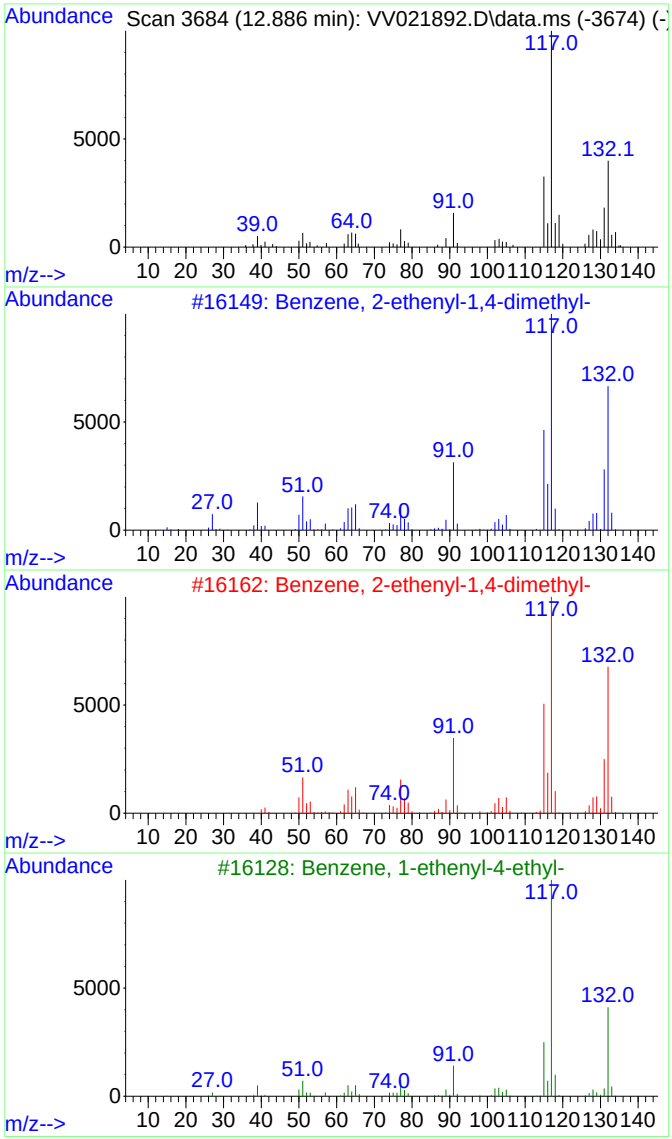
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

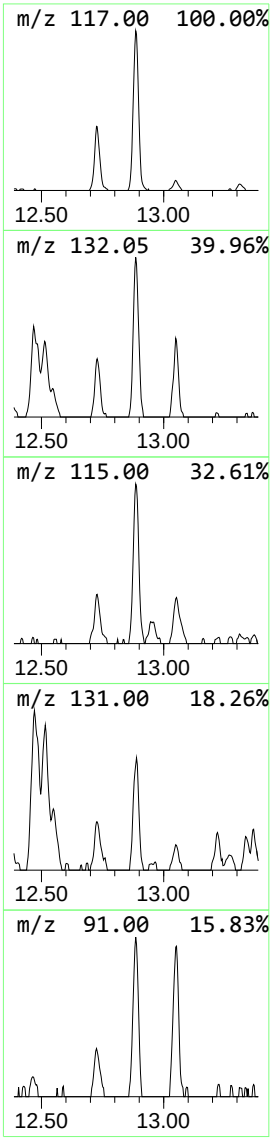
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	3.01 ug/L	74813	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, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





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Data File : VV021892.D
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ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

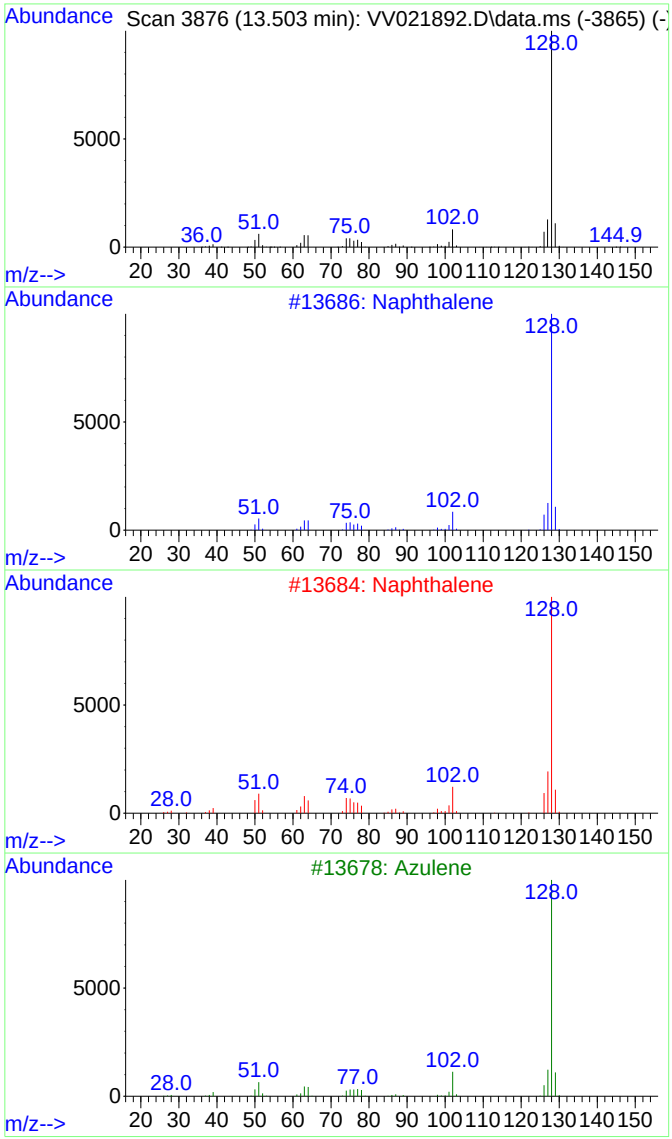
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

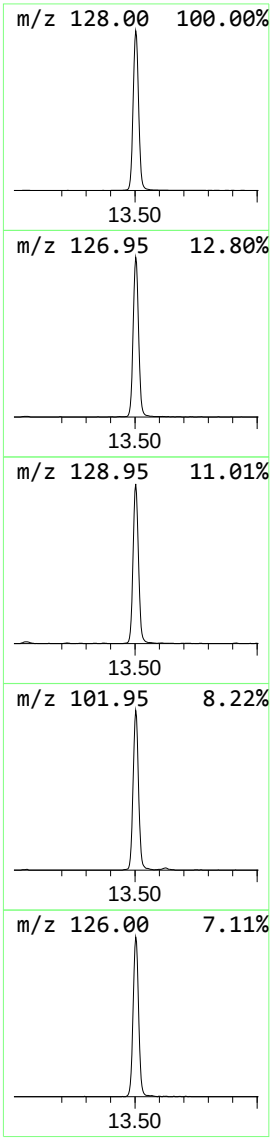
****

Peak Number 5 Azulene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.503	75.55 ug/L	1878880	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





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Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

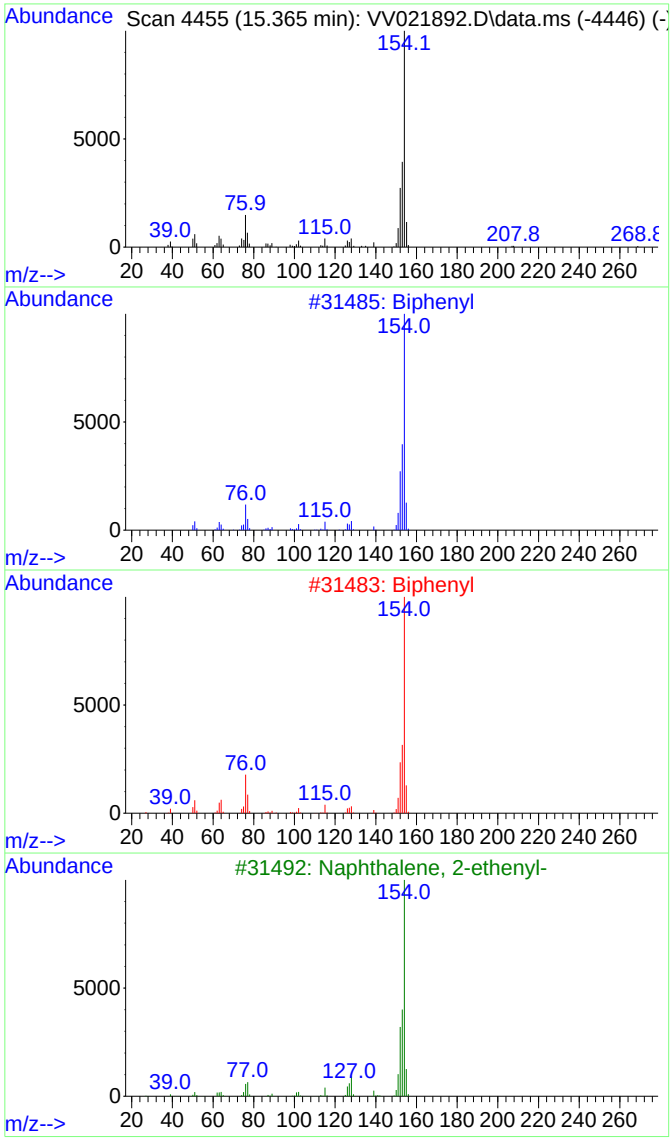
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

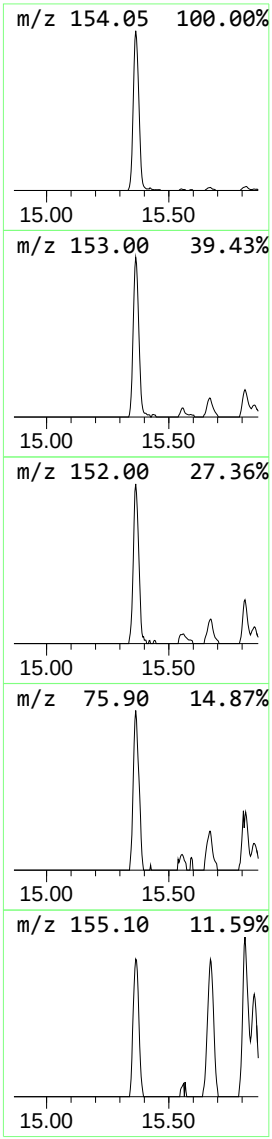
****

Peak Number 8 Biphenyl Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.365	4.04 ug/L	100555	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl	154	C12H10	000092-52-4	94
2	Biphenyl	154	C12H10	000092-52-4	87
3	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	81
4	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	81
5	Biphenyl	154	C12H10	000092-52-4	76





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Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

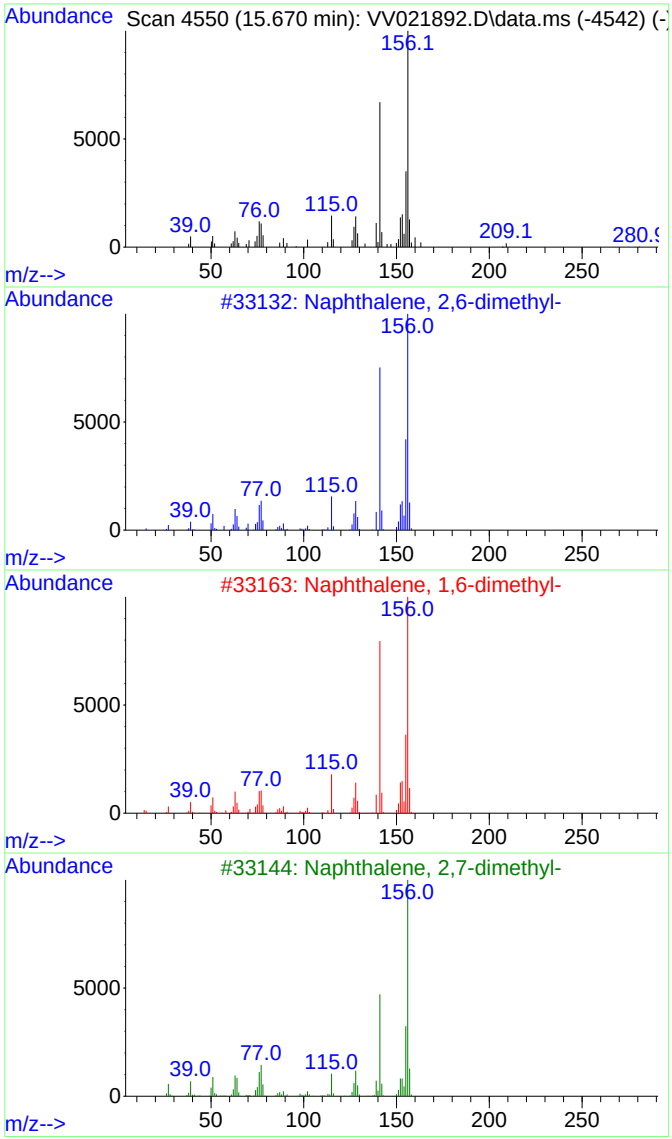
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

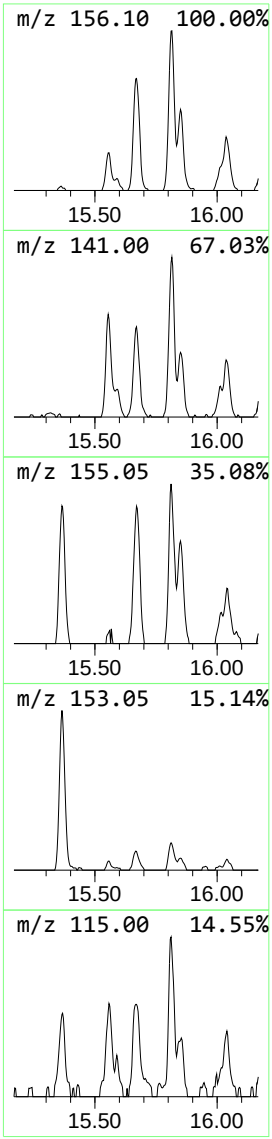
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 2,6-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.670	2.57 ug/L	63836	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, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKF3

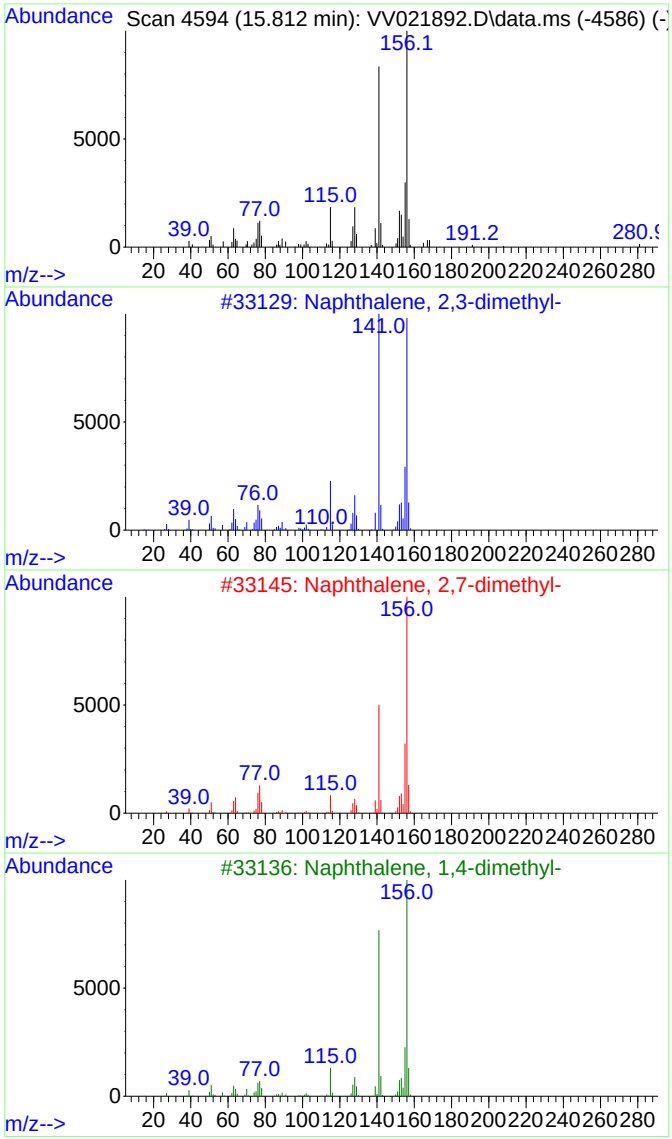
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

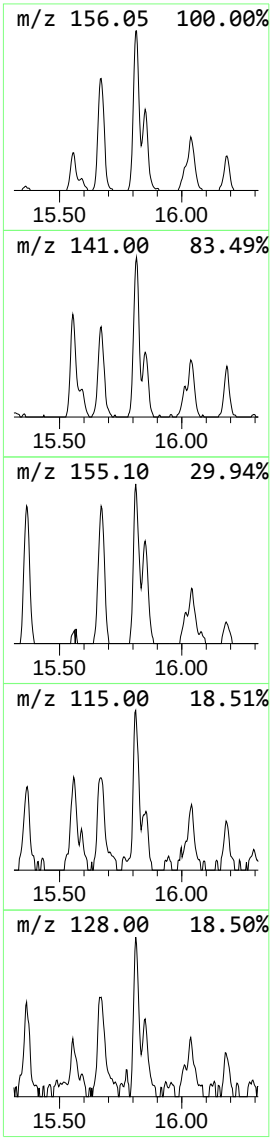
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.812	2.88 ug/L	71590	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	98
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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

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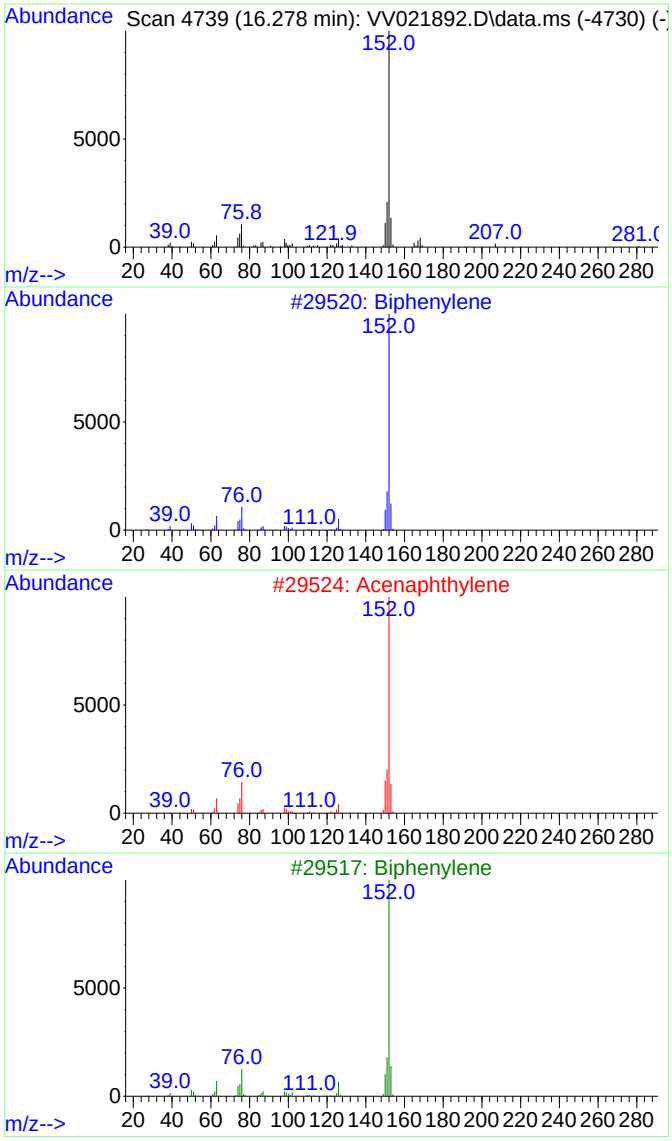
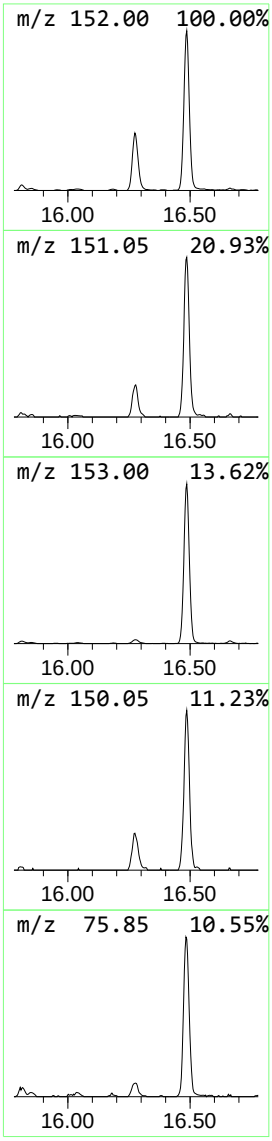
Peak Number 11 Biphenylene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.278	2.82 ug/L	70150	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenylene	152	C12H8	000259-79-0	93
2	Acenaphthylene	152	C12H8	000208-96-8	87
3	Biphenylene	152	C12H8	000259-79-0	87
4	Acenaphthylene	152	C12H8	000208-96-8	87
5	Biphenylene	152	C12H8	000259-79-0	62





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

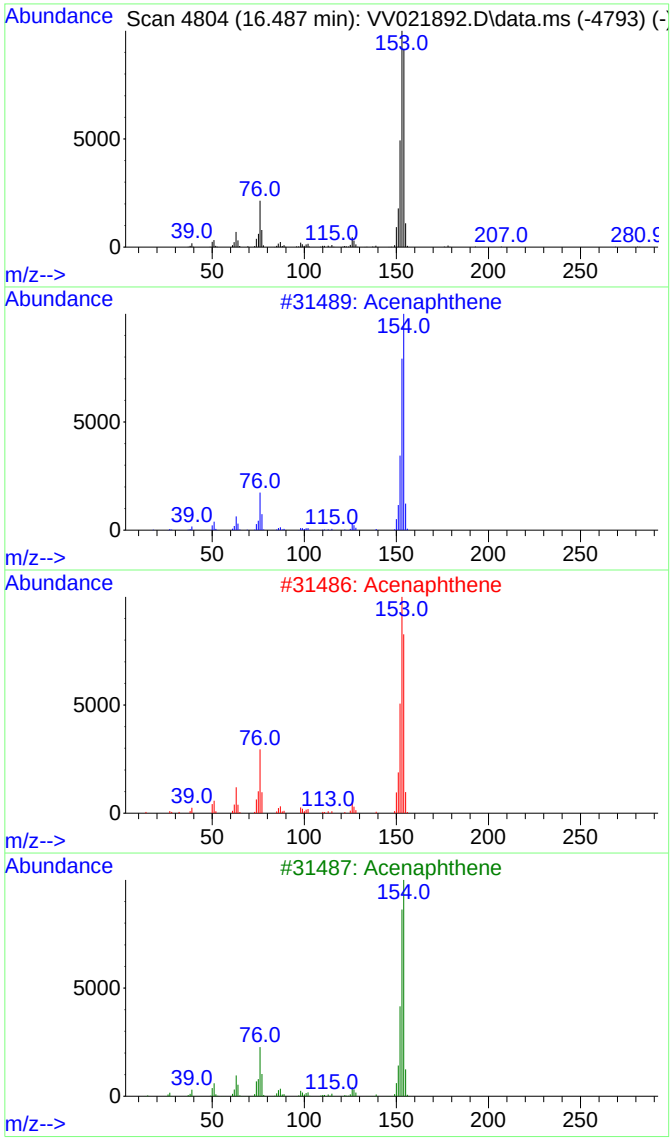
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

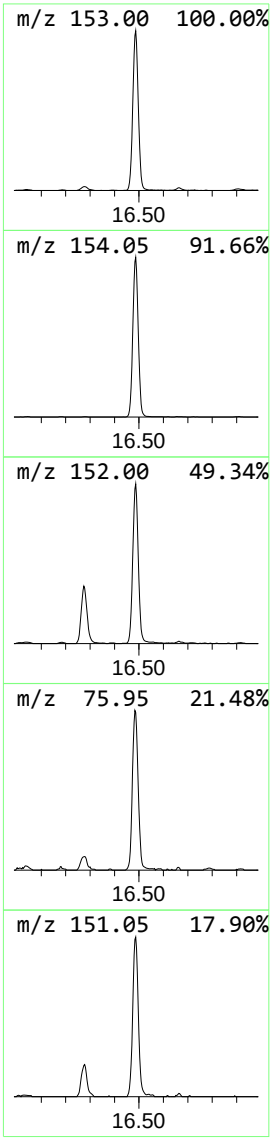
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Naphthalene, 2-ethenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.487	23.45 ug/L	583202	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	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	64
5	1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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

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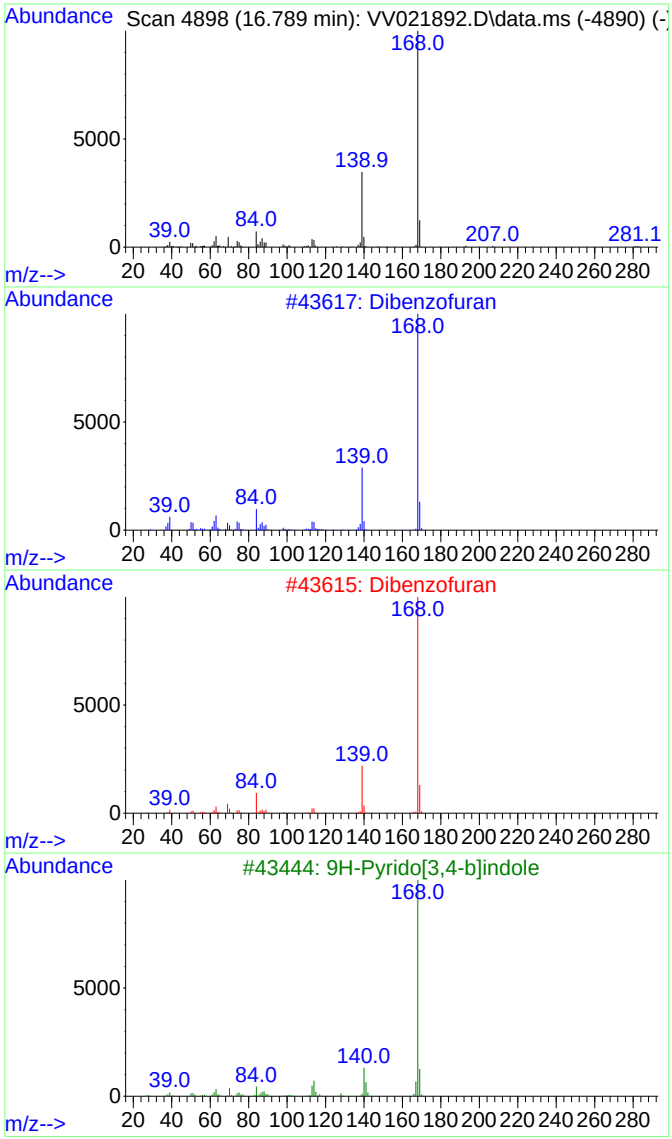
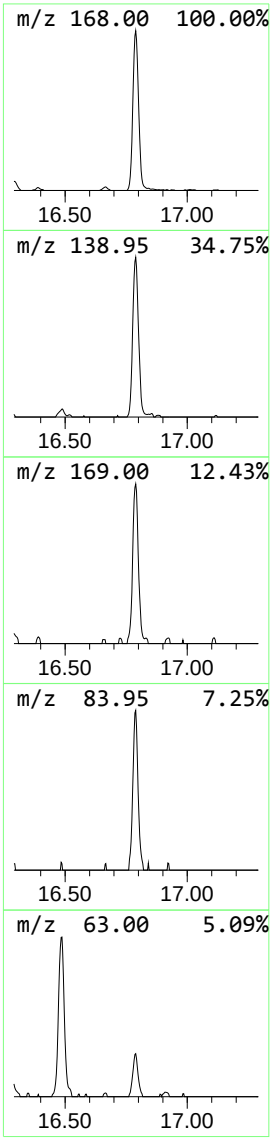
Peak Number 13 Dibenzofuran Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.789	5.34 ug/L	132869	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	80
3	9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	59
4	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	59
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\VV081921\
Data File : VV021892.D
Acq On : 19 Aug 2021 15:35
Operator : SY/MD
Sample : M3448-14
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKF3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.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
Sulfur dioxide	1.192	8.2	ug/L	163663	1	5.619	996145	50.0
Indane	11.529	113.6	ug/L	2824490	3	11.252	1243390	50.0
Benzene, 2-ethe...	12.886	3.0	ug/L	74813	3	11.252	1243390	50.0
Azulene	13.503	75.5	ug/L	1878880	3	11.252	1243390	50.0
Biphenyl	15.365	4.0	ug/L	100555	3	11.252	1243390	50.0
Naphthalene, 2,...	15.670	2.6	ug/L	63836	3	11.252	1243390	50.0
Naphthalene, 2,...	15.812	2.9	ug/L	71590	3	11.252	1243390	50.0
Biphenylene	16.278	2.8	ug/L	70150	3	11.252	1243390	50.0
Naphthalene, 2-...	16.487	23.4	ug/L	583202	3	11.252	1243390	50.0
Dibenzofuran	16.789	5.3	ug/L	132869	3	11.252	1243390	50.0