LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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: VV021886.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	100	rVB	216735	219927	0.74%	0.308%
2	1.558	154	161	181	rVB	160963	194783	0.65%	0.272%
3	2.105	322	331	345	rVB	302801	426793	1.43%	0.597%
4	2.391	417	420	421	rBV2	455	247	0.00%	0.000%
5	2.770	528	538	555	rVB2	9444	19570	0.07%	0.027%
6	2.918	580	584	589	rBV3	336	318	0.00%	0.000%
7	2.979	600	603	607	rVB4	327	264	0.00%	0.000%
8	3.105	640	642	645	rBV	414	245	0.00%	0.000%
9	3.269	690	693	694	rBV	420	243	0.00%	0.000%
10	3.381	723	728	730	rBV2	486	387	0.00%	0.001%
11	3.407	731	736	741	rVB3	356	482	0.00%	0.001%
12	3.555	780	782	788	rVB3	426	333	0.00%	0.000%
13	3.587	788	792	796	rBV2	269	286	0.00%	0.000%
14	3.677	816	820	825	rVB2	314	265	0.00%	0.000%
15	3.716	829	832	834	rBV	421	257	0.00%	0.000%
16	3.741	837	840	845	rBV2	531	573	0.00%	0.001%
17	3.876	872	882	915	rBV2	121337	326559	1.09%	0.457%
18	4.349	1013	1029	1063	rBV	236434	585472	1.96%	0.819%
19	4.629	1113	1116	1120	rBV3	369	295	0.00%	0.000%
20	4.648	1120	1122	1124	rBV	359	182	0.00%	0.000%
21	4.664	1124	1127	1128	rBV2	537	244	0.00%	0.000%
22	4.770	1158	1160	1163	rVB	393	209	0.00%	0.000%
23	4.789	1163	1166	1172	rVB3	334	317	0.00%	0.000%
24	5.047	1228	1246	1283	rBV2	597788	1577923	5.29%	2.207%
25	5.413	1357	1360	1361	rBV2	520	251	0.00%	0.000%
26	5.481	1378	1381	1382	rBV	373	190	0.00%	0.000%
27	5.619	1412	1424	1451	rBV	497311	1002101	3.36%	1.401%
28	5.908	1504	1514	1529	rBV5	13591	29593	0.10%	0.041%
29	6.072	1551	1565	1588	rBV	325289	667199	2.24%	0.933%
30	6.310	1636	1639	1645	rBV3	314	269	0.00%	0.000%
31	6.378	1657	1660	1666	rBV2	286	221	0.00%	0.000%
32	6.436	1673	1678	1681	rBV2	289	272	0.00%	0.000%
33	6.513	1695	1702	1704	rBV3	427	393	0.00%	0.001%
34	6.571	1718	1720	1722	rBV2	315	175	0.00%	0.000%
35	6.641	1739	1742	1743	rBV	320	199	0.00%	0.000%
36	6.728	1764	1769	1770	rBV2	351	229	0.00%	0.000%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

37	6.876	1810	1815	1820	rBV2	248	302	0.00%	0.000%
38	6.989	1837	1850	1876	rBV	187663	342223	1.15%	0.479%
39	7.259	1930	1934	1936	rBV2	548	389	0.00%	0.001%
40	7.317	1940	1952	1966	rBV	726150	1258366	4.22%	1.760%
41	7.391	1967	1975	2000	rVB	211485	377866	1.27%	0.528%
42	7.625	2037	2048	2066	rBV	124242	218974	0.73%	0.306%
43	7.812	2104	2106	2110	rBV	363	180	0.00%	0.000%
44	7.834	2110	2113	2117	rVB4	733	512	0.00%	0.001%
45	7.866	2118	2123	2128	rVB3	333	361	0.00%	0.001%
46	7.892	2129	2131	2134	rVB2	416	185	0.00%	0.000%
47	7.944	2140	2147	2152	rBV4	1763	2475	0.01%	0.003%
48	8.088	2182	2192	2231	rBV	344163	622995	2.09%	0.871%
49	8.455	2303	2306	2308	rBV2	611	404	0.00%	0.001%
50	8.503	2317	2321	2324	rBV2	380	377	0.00%	0.001%
51	8.699	2375	2382	2385	rBV3	284	315	0.00%	0.000%
52	8.754	2396	2399	2402	rVB2	500	260	0.00%	0.000%
53	8.793	2407	2411	2413	rBV2	440	276	0.00%	0.000%
54	8.809	2413	2416	2418	rBV2	362	221	0.00%	0.000%
55	8.854	2418	2430	2457	rBV	746777	1215095	4.07%	1.699%
56	9.014	2468	2480	2504	rBV	887172	1404607	4.71%	1.964%
57	9.140	2508	2519	2539	rVB	423927	691045	2.32%	0.966%
58	9.461	2615	2619	2622	rBV3	434	307	0.00%	0.000%
59	9.503	2624	2632	2633	rBV3	2021	1659	0.01%	0.002%
60	9.548	2635	2646	2672	rVV4	307807	684664	2.29%	0.957%
61	9.719	2697	2699	2705	rVB3	672	531	0.00%	0.001%
62	9.934	2759	2766	2777	rBV3	7967	11857	0.04%	0.017%
63	10.030	2793	2796	2804	rBV6	907	1029	0.00%	0.001%
64	10.130	2818	2827	2837	rBV3	4147	6640	0.02%	0.009%
65	10.220	2842	2855	2871	rBV	466886	740066	2.48%	1.035%
66	10.358	2890	2898	2914	rVB	13647	25571	0.09%	0.036%
67	10.452	2916	2927	2945	rBV	76911	161596	0.54%	0.226%
68	10.542	2946	2955	2970	rVB	73424	115825	0.39%	0.162%
69	10.661	2990	2992	2993	rBV	560	213	0.00%	0.000%
70	10.741	3008	3017	3037	rVB	36758	71562	0.24%	0.100%
71	10.841	3043	3048	3051	rVB4	656	588	0.00%	0.001%
72	10.915	3057	3071	3085	rBV	233111	368399	1.23%	0.515%
73	10.989	3086	3094	3103	rBV	54059	83975	0.28%	0.117%
74	11.172	3139	3151	3165	rBV	599197	922517	3.09%	1.290%
75	11.249	3165	3175	3194	rVV	849238	1329753	4.46%	1.860%
76	11.339	3195	3203	3216	rVB	99899	158440	0.53%	0.222%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

77	11.529	3251	3262	3275	rBV	874765	1334679	4.47%	1.866%
78	11.625	3282	3292	3311	rVB	781760	1245102	4.17%	1.741%
79	11.744	3317	3329	3349	rBV	7140220	10856286	36.38%	15.182%
80	12.117	3436	3445	3465	rVB	79510	129791	0.43%	0.182%
81	12.361	3513	3521	3531	rBV	117786	180930	0.61%	0.253%
82	12.464	3543	3553	3563	rBV2	340576	617867	2.07%	0.864%
83	12.728	3625	3635	3651	rVB	141269	217192	0.73%	0.304%
84	12.886	3665	3684	3695	rBV2	243750	383827	1.29%	0.537%
85	12.950	3696	3704	3714	rVB2	66024	101217	0.34%	0.142%
86	13.050	3723	3735	3756	rVB2	761720	1210920	4.06%	1.693%
87	13.156	3759	3768	3779	rBV6	19247	37220	0.12%	0.052%
88	13.506	3862	3877	3902	rBV	18449797	29838733	100.00%	41.728%
89	13.622	3903	3913	3933	rVB	935692	1452490	4.87%	2.031%
90	13.911	3995	4003	4019	rBV2	17188	28678	0.10%	0.040%
91	14.095	4051	4060	4061	rBV3	19866	23614	0.08%	0.033%
92	14.580	4203	4211	4217	rBV	74215	104779	0.35%	0.147%
93	14.635	4217	4228	4254	rVB	2569601	3921674	13.14%	5.484%
94	14.747	4255	4263	4273	rBV	73163	115301	0.39%	0.161%
95	14.818	4273	4285	4300	rBV	1520640	2336428	7.83%	3.267%
96	15.365	4447	4455	4468	rVB	212560	312367	1.05%	0.437%
97	15.554	4507	4514	4521	rBV	44794	67323	0.23%	0.094%
98	15.670	4541	4550	4571	rVB	78586	154687	0.52%	0.216%
99	15.812	4586	4594	4601	rBV	112134	174695	0.59%	0.244%
100	16.487	4793	4804	4821	rBV	494393	782614	2.62%	1.094%

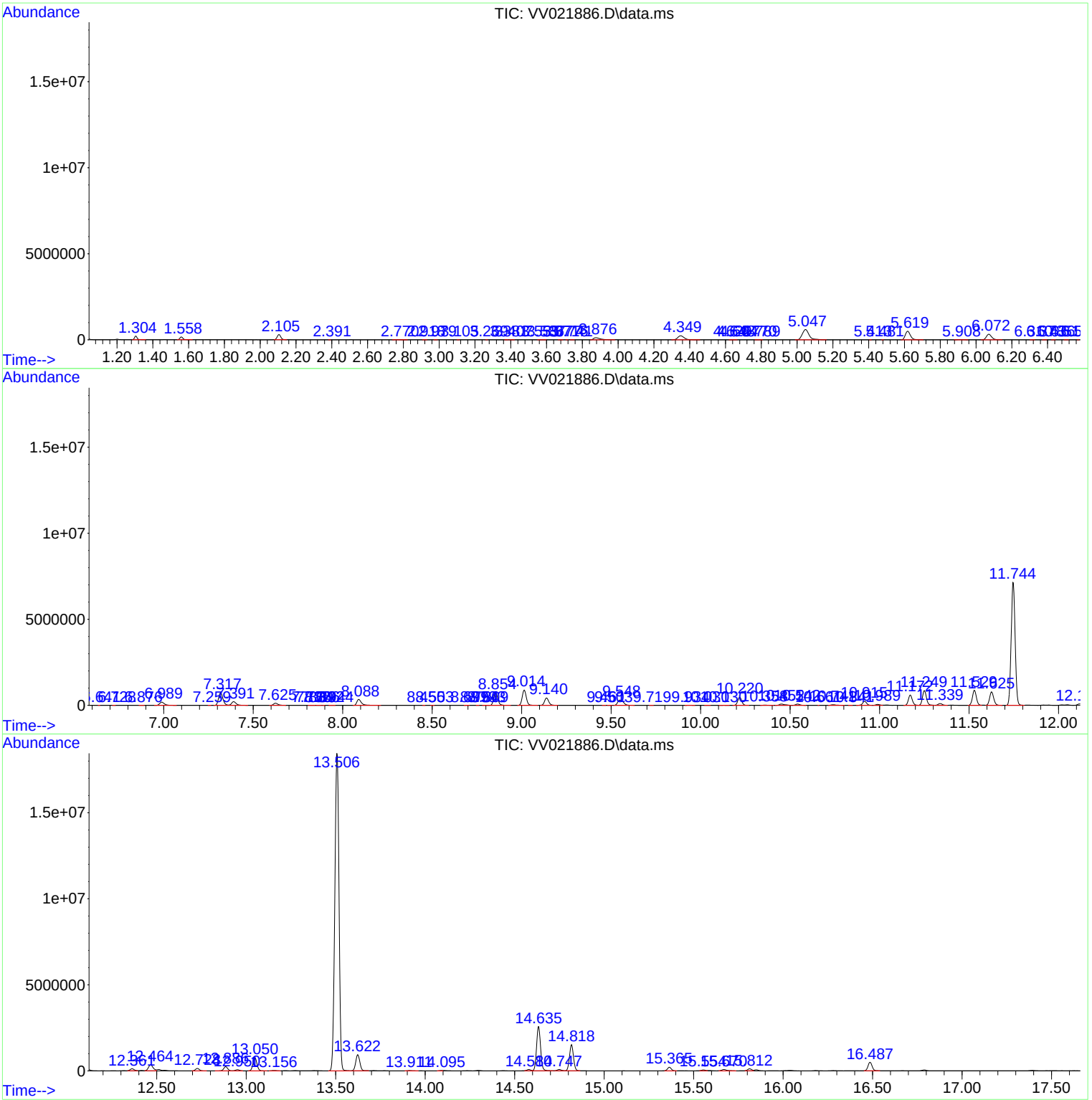
Sum of corrected areas: 71508300

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

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 : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

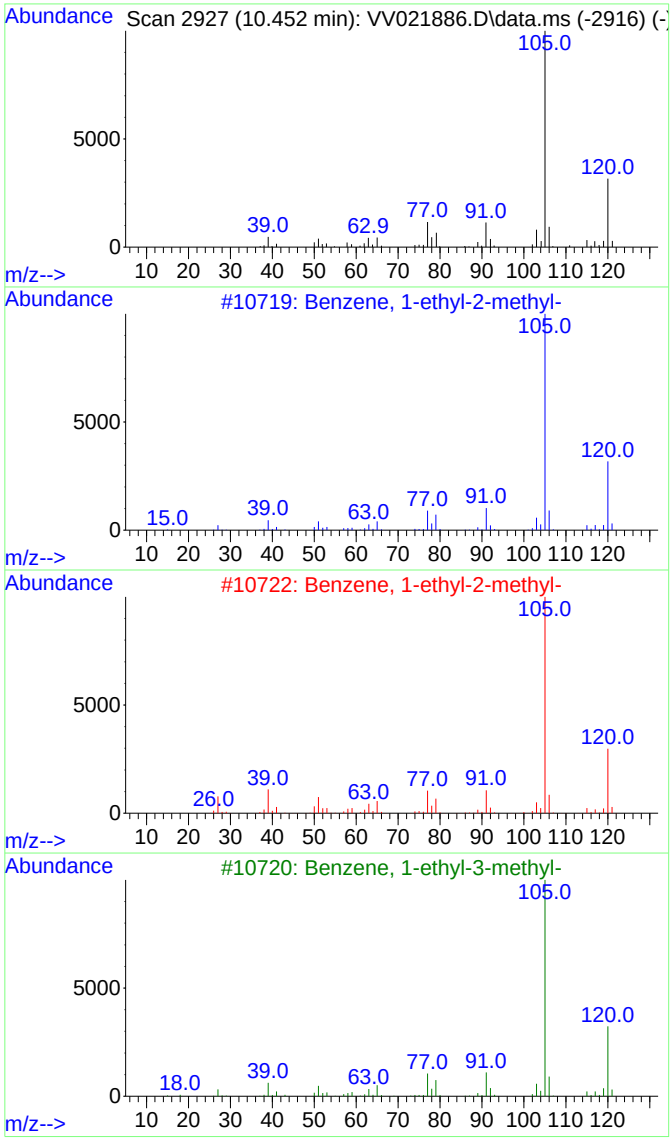
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

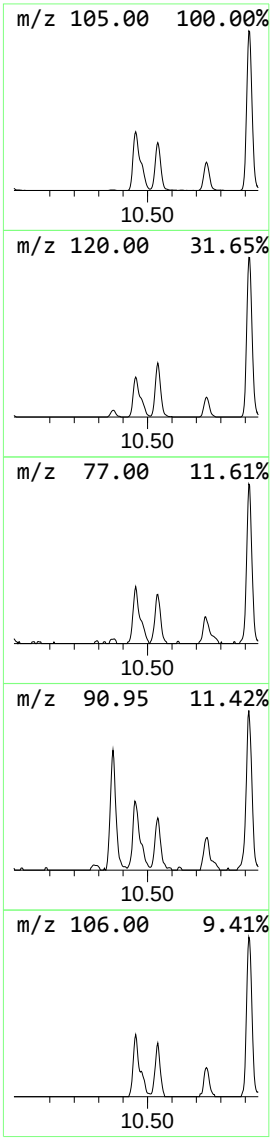
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.452	6.08 ug/L	161596	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	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

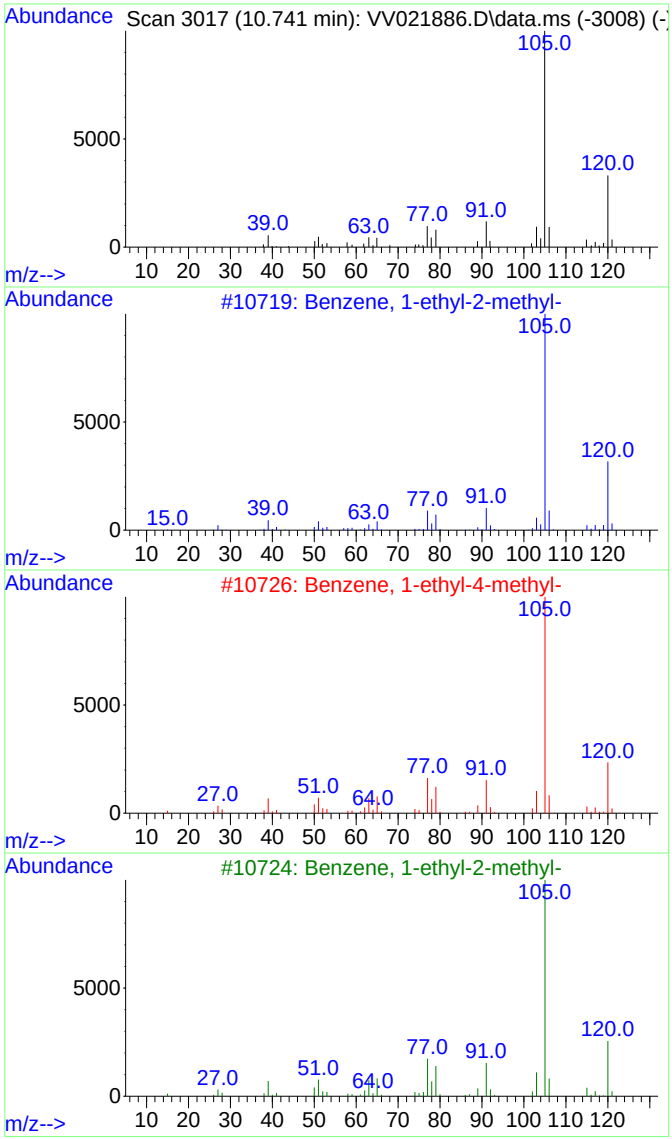
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

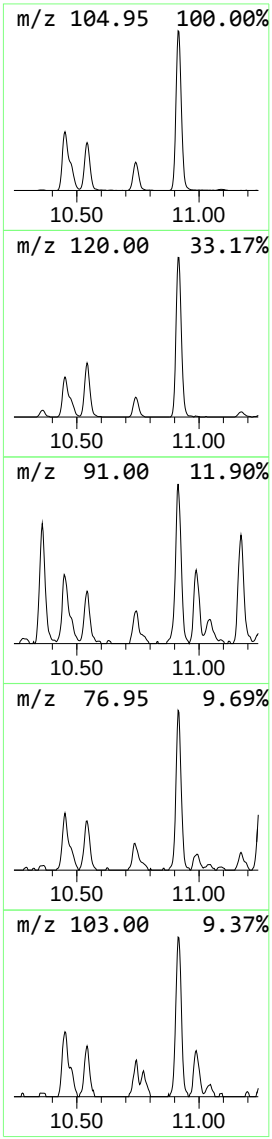
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.741	2.69 ug/L	71562	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	93
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
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\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

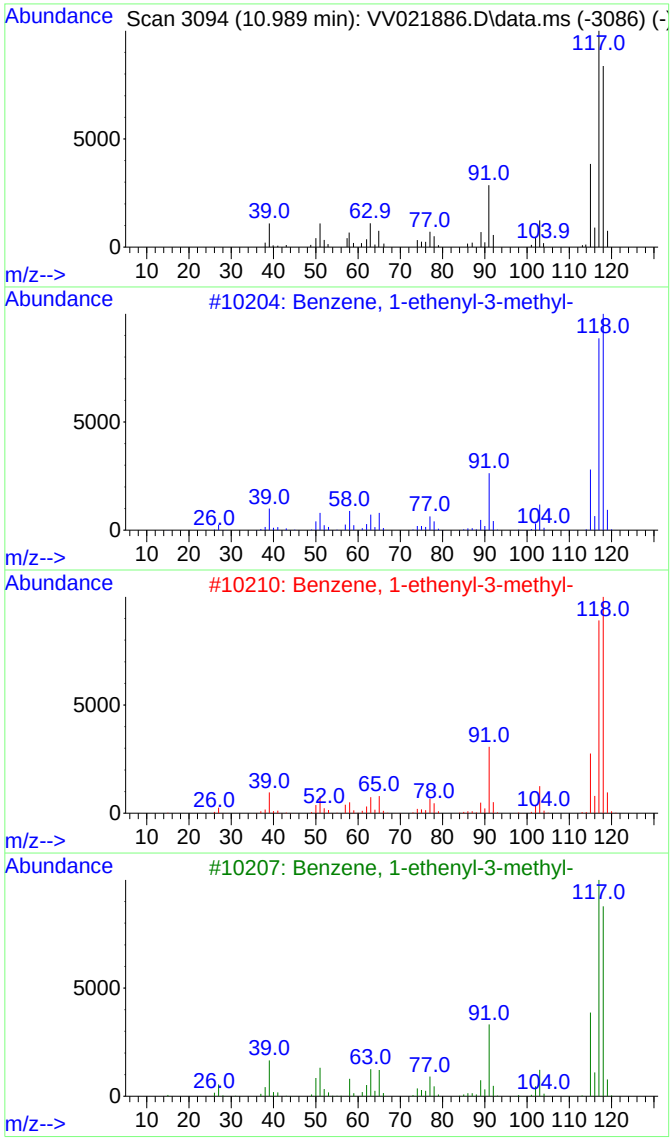
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

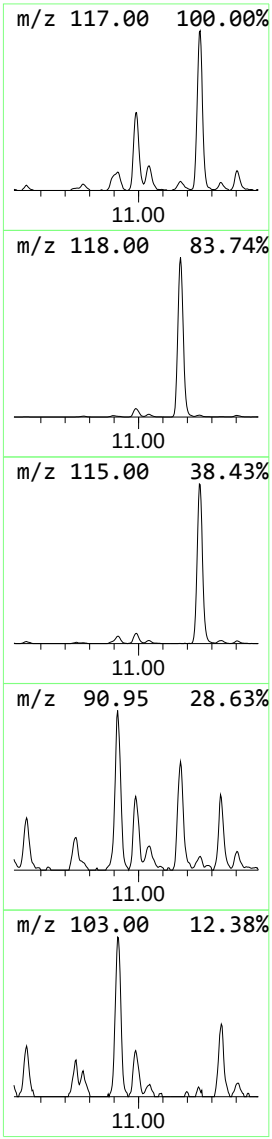
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethenyl-3-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.989	3.16 ug/L	83975	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	95
3	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	95
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

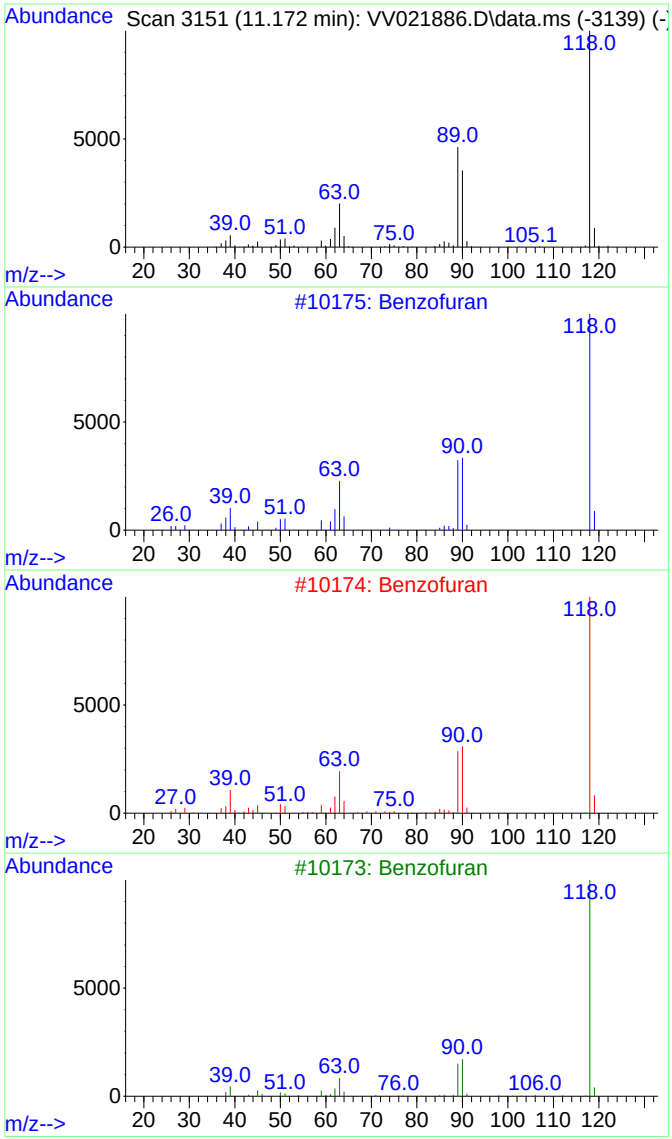
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

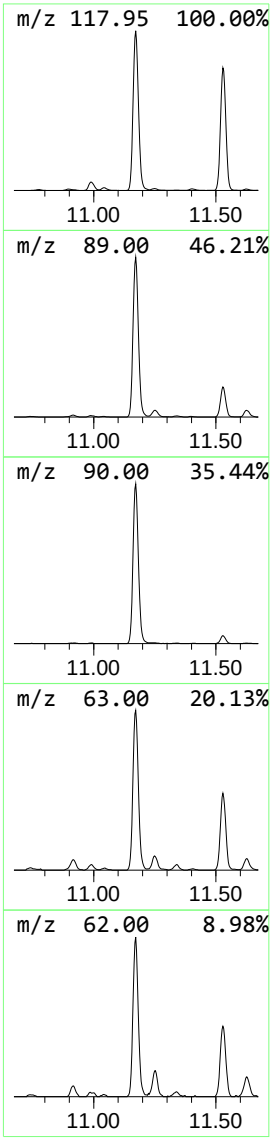
****

Peak Number 5 Benzofuran Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.172	34.69 ug/L	922517	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	90
5	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

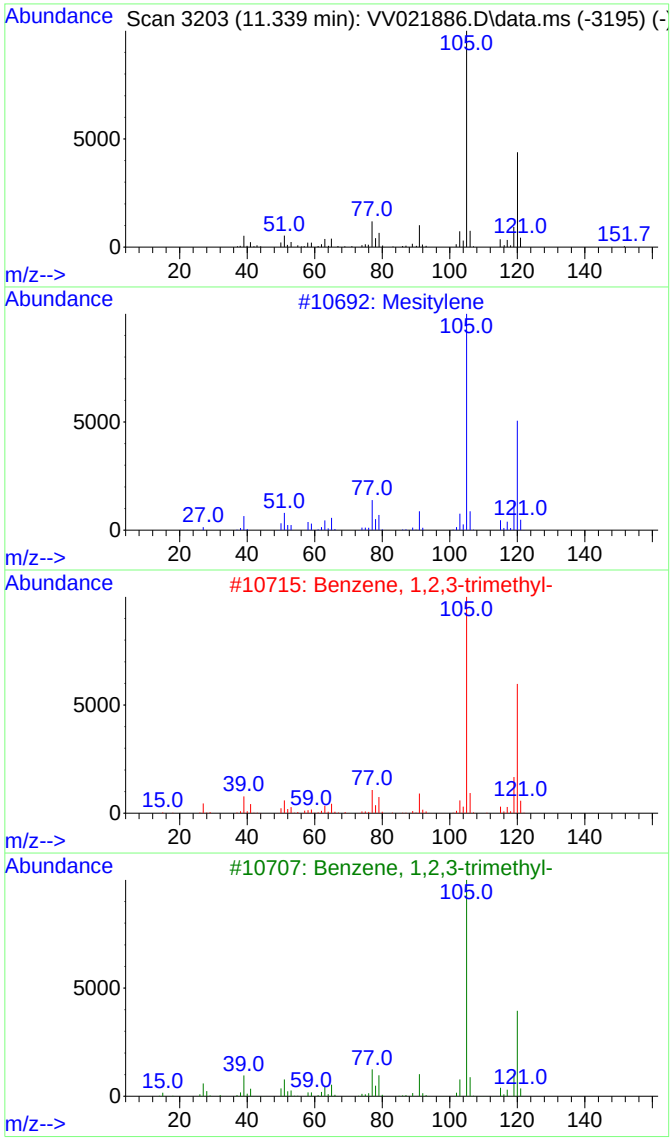
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

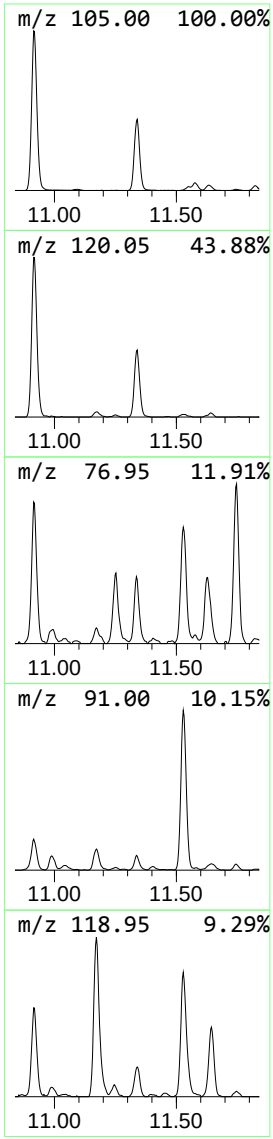
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.339	5.96 ug/L	158440	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	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

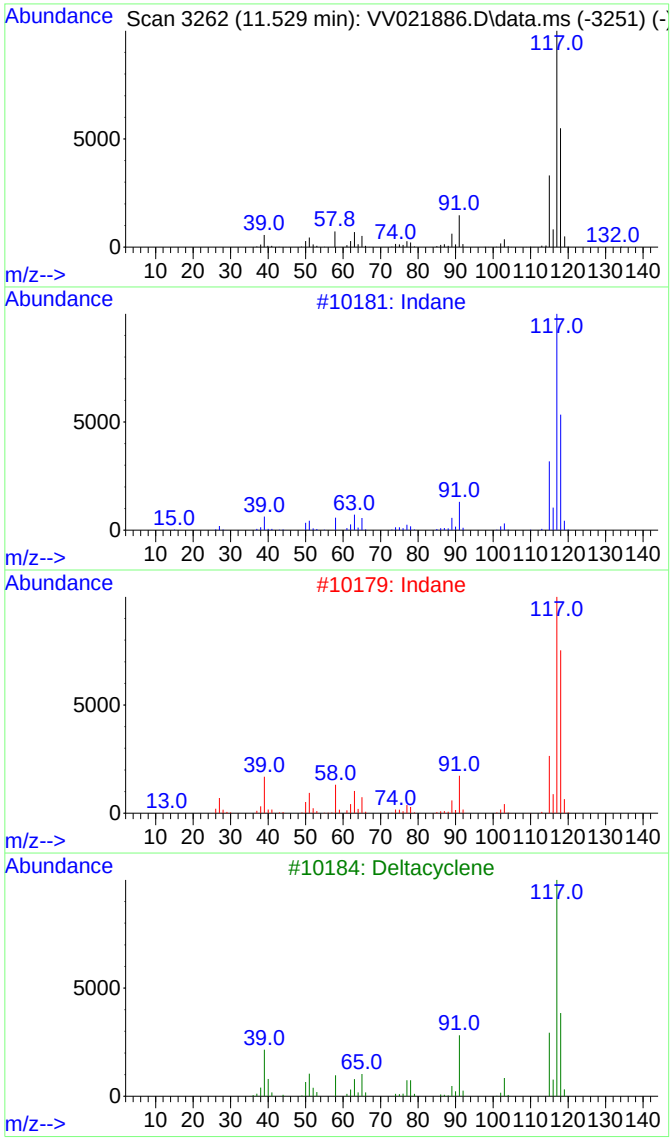
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

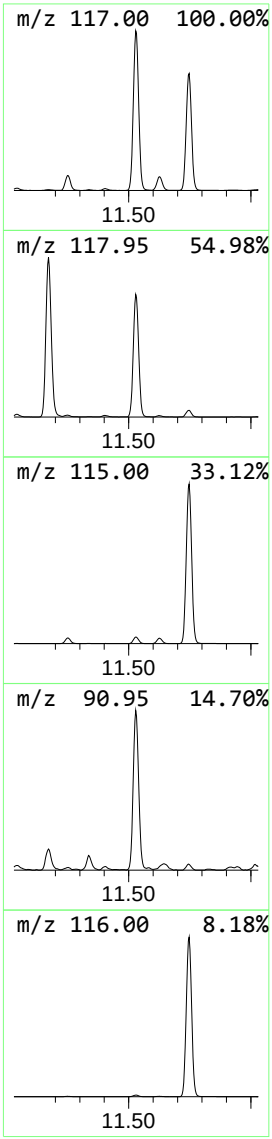
****

Peak Number 7 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	50.19 ug/L	1334680	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	Deltacyclene	118	C9H10	007785-10-6	72
4	Indane	118	C9H10	000496-11-7	64
5	Indane	118	C9H10	000496-11-7	64





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

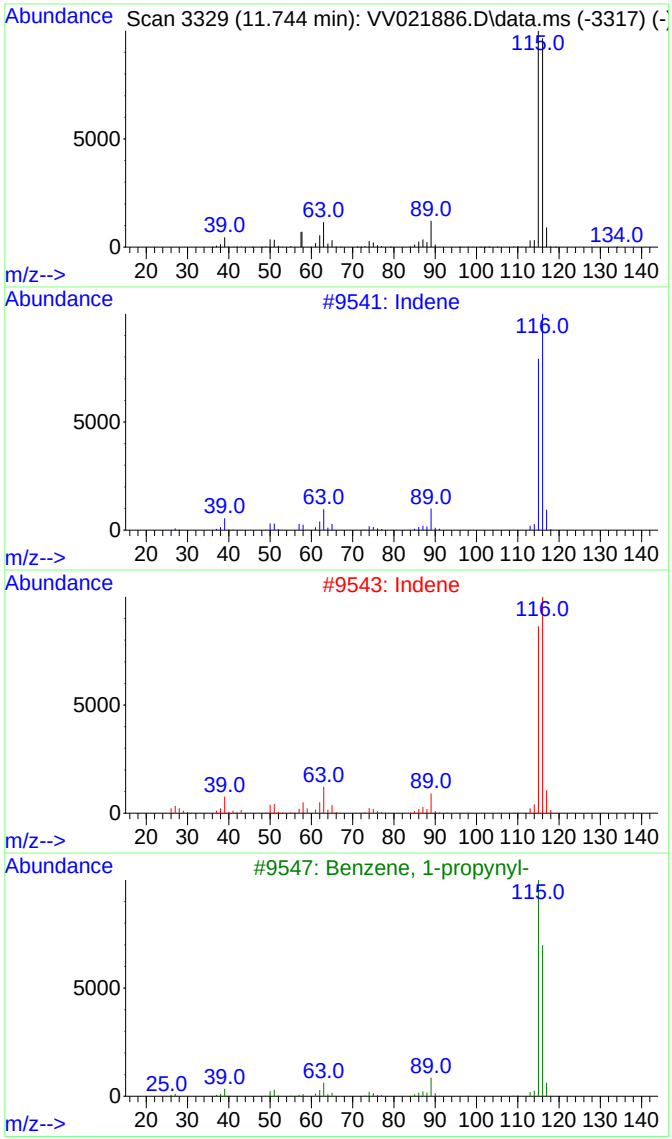
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

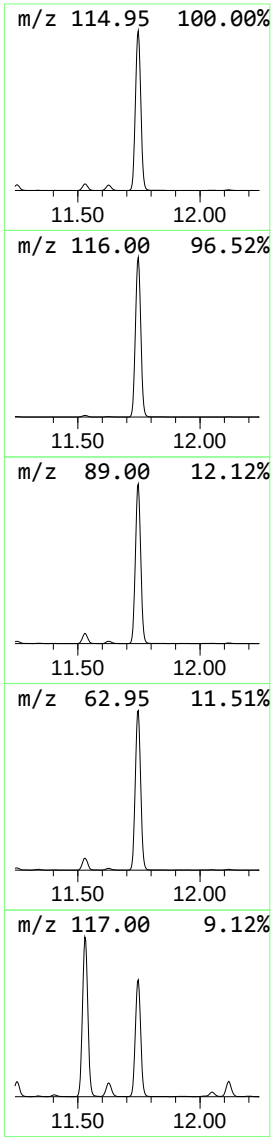
****

Peak Number 8 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.744	408.21 ug/L	10856300	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Indene	116	C9H8	000095-13-6	95
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
4	Indene	116	C9H8	000095-13-6	94
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

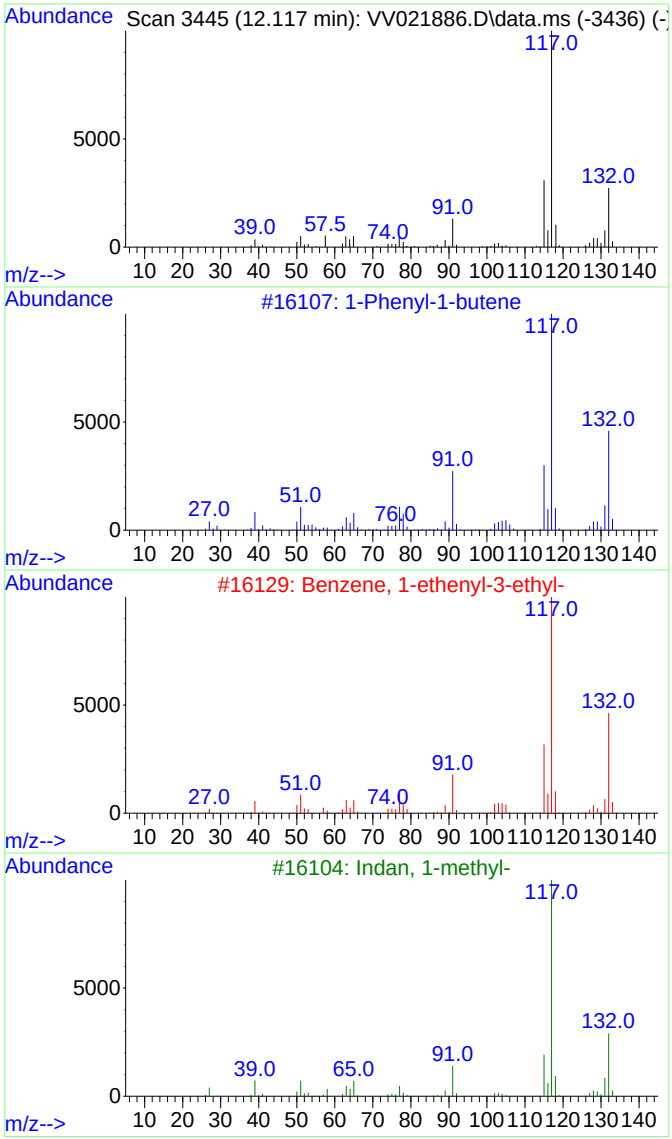
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

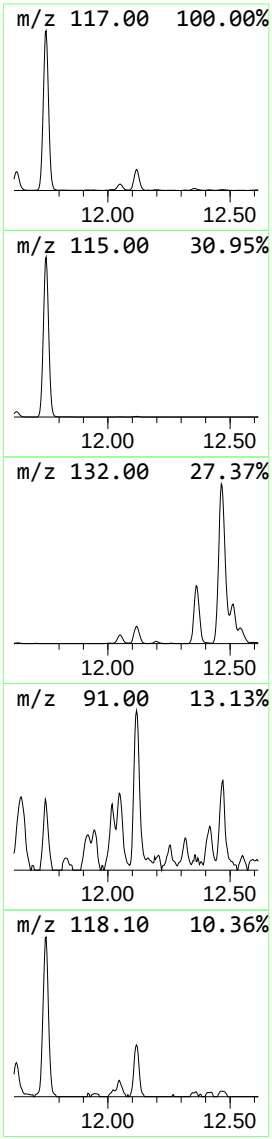
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1-Phenyl-1-butene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	4.88 ug/L	129791	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

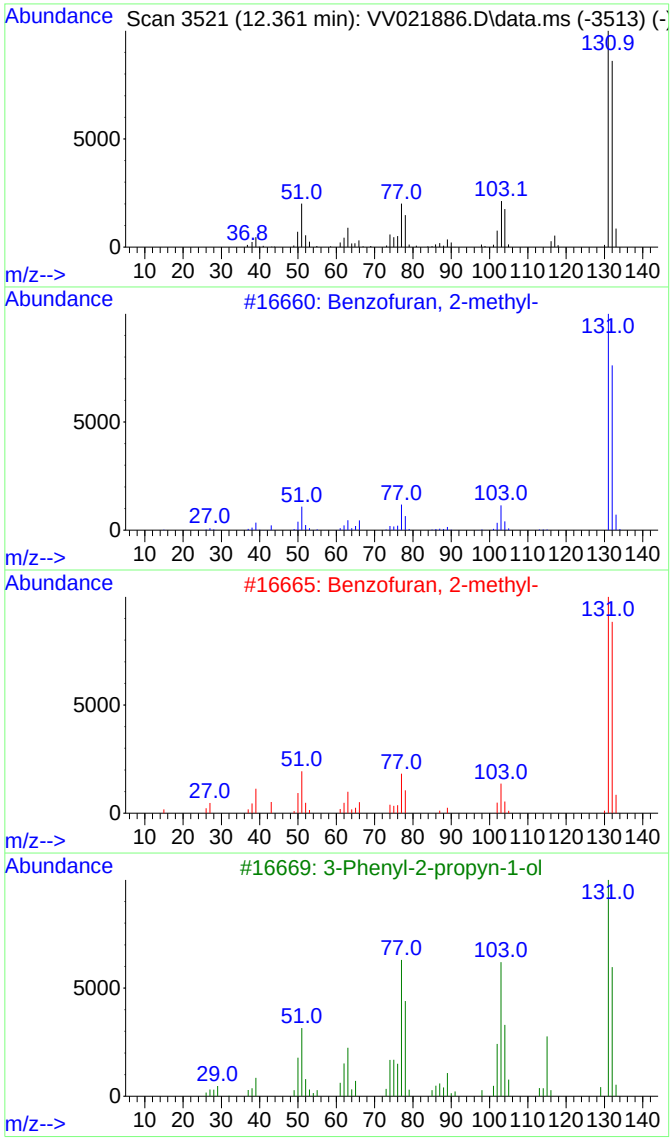
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

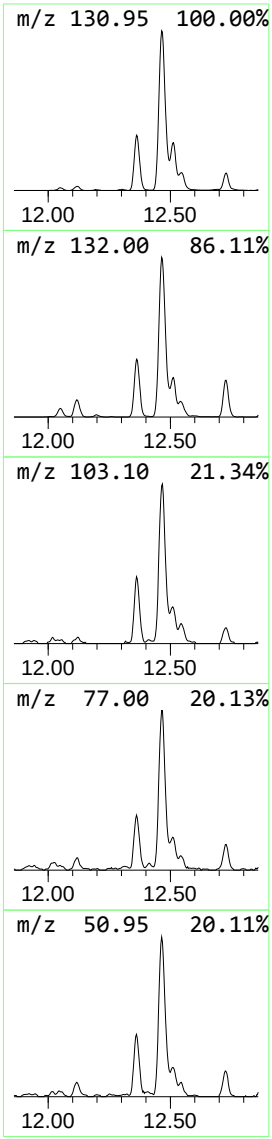
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzofuran, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.361	6.80 ug/L	180930	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, 2-methyl-	132	C9H8O	004265-25-2	93
3	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	1H-Indenol	132	C9H8O	056631-57-3	91





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

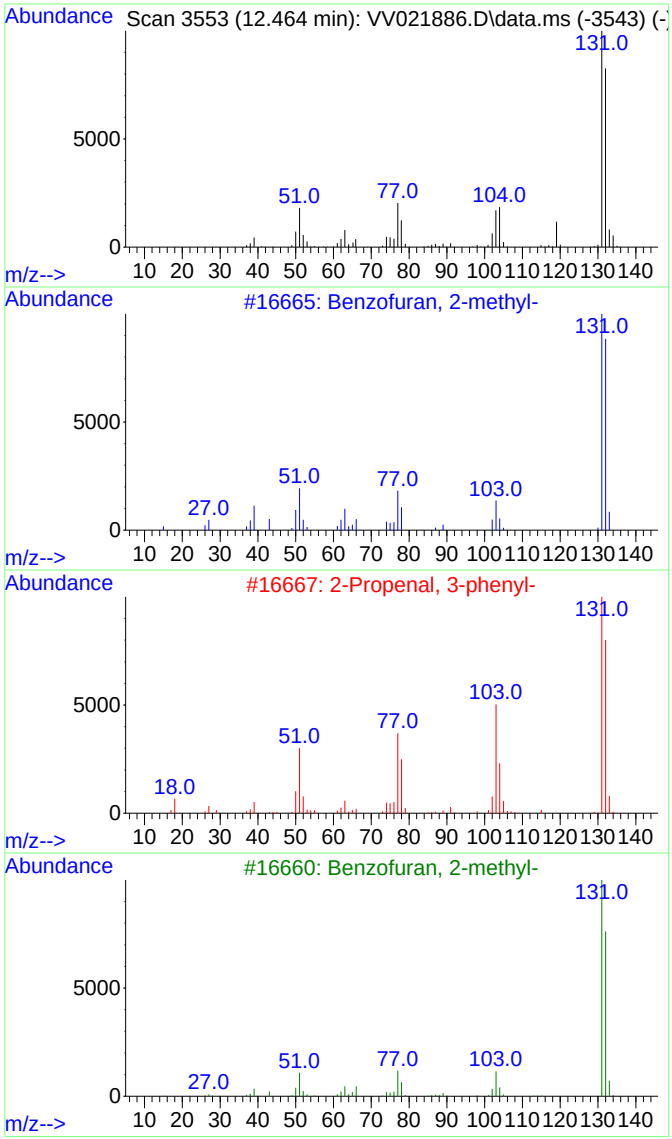
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

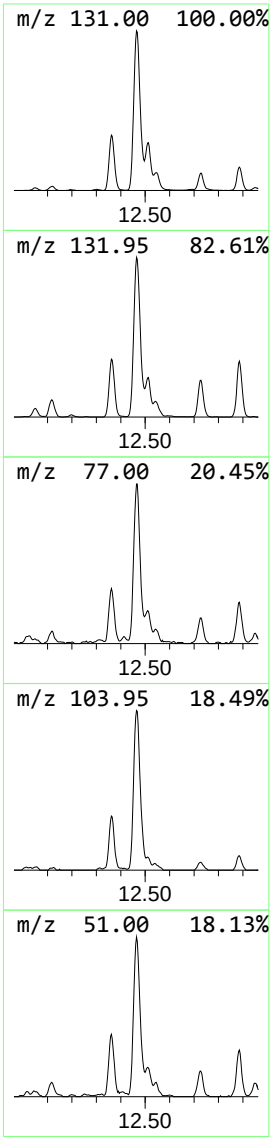
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 2-Propenal, 3-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	23.23 ug/L	617867	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	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

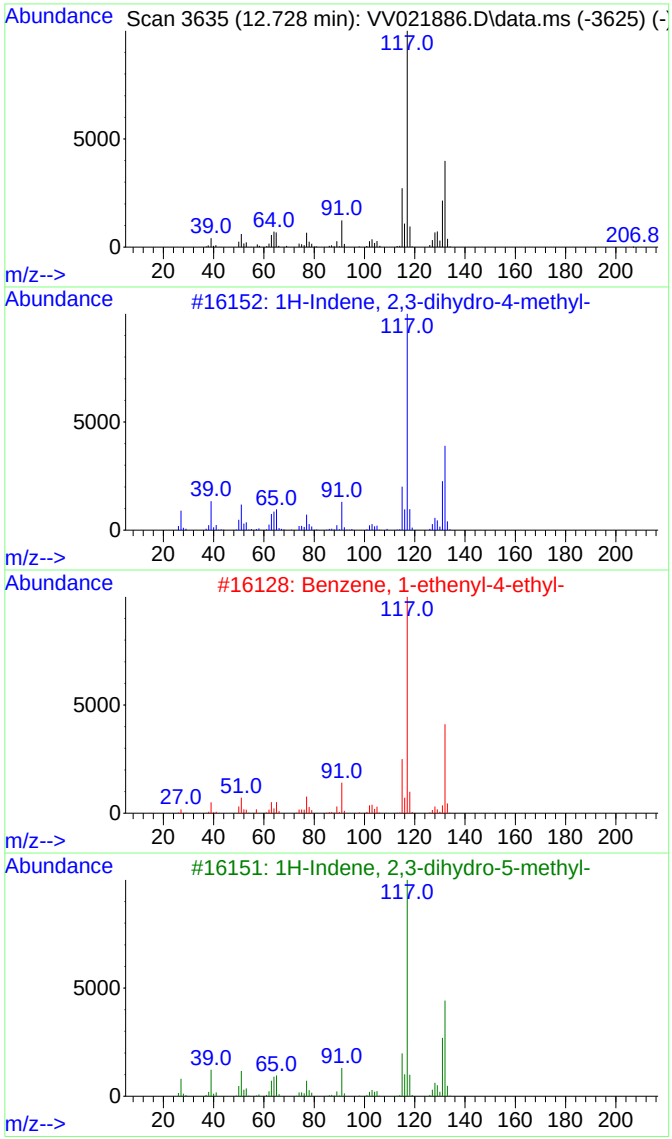
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

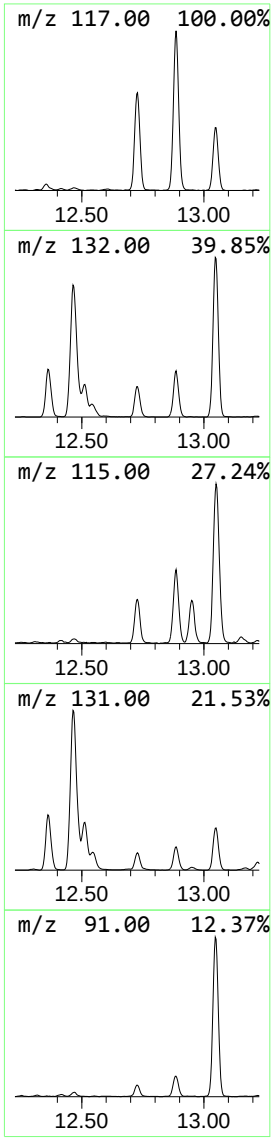
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.728	8.17 ug/L	217192	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

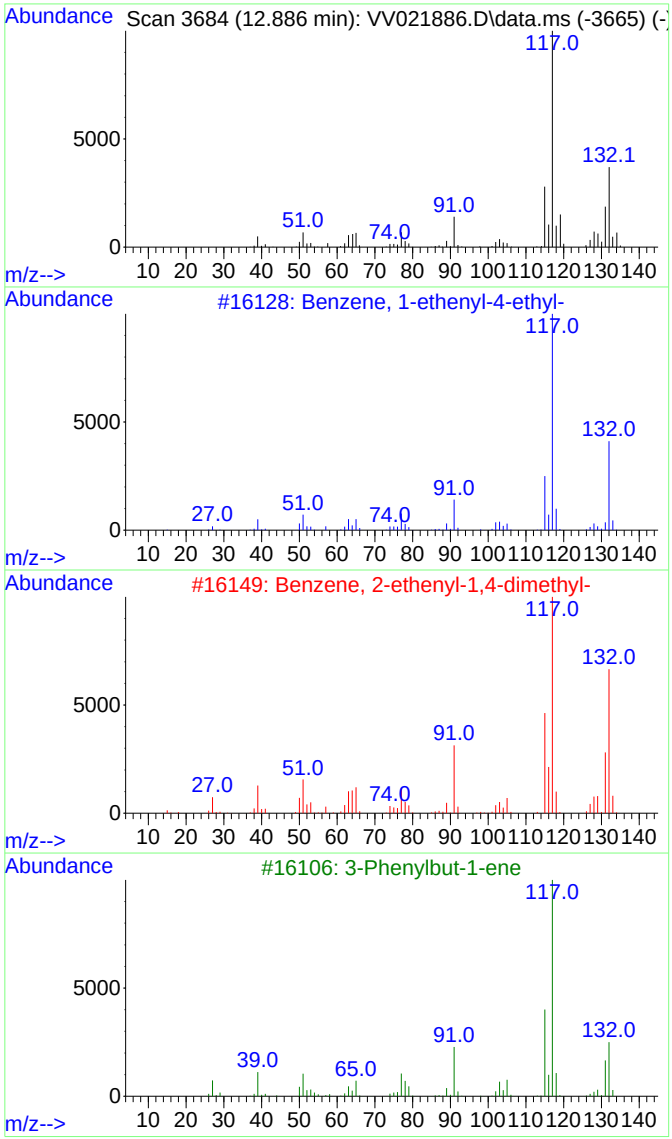
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

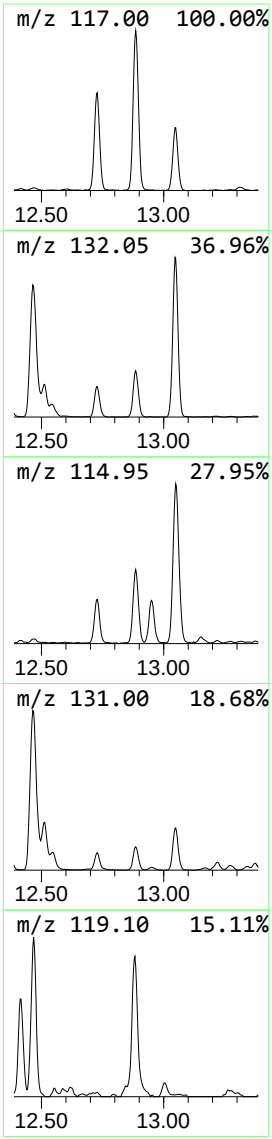
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	14.43 ug/L	383827	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

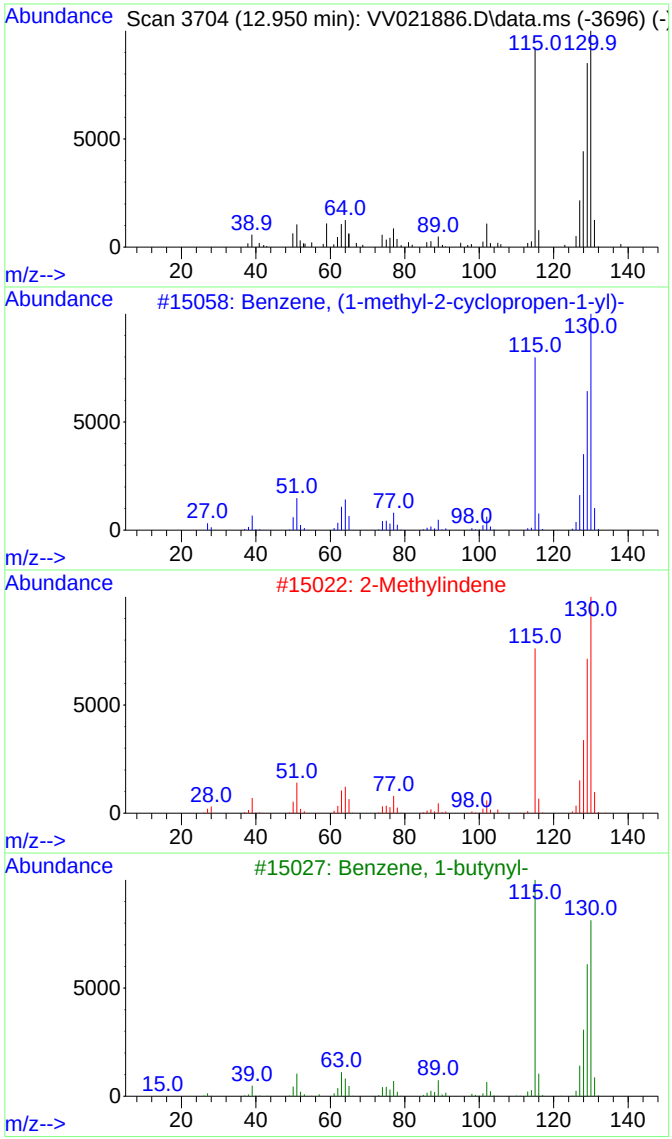
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

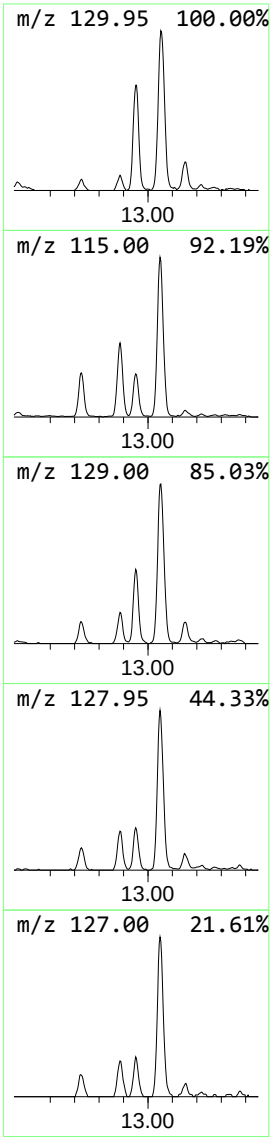
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, (1-methyl-2-cyclop... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.950	3.81 ug/L	101217	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	94
2	2-Methylindene	130	C10H10	002177-47-1	93
3	Benzene, 1-butynyl-	130	C10H10	000622-76-4	93
4	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	92
5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	91





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Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

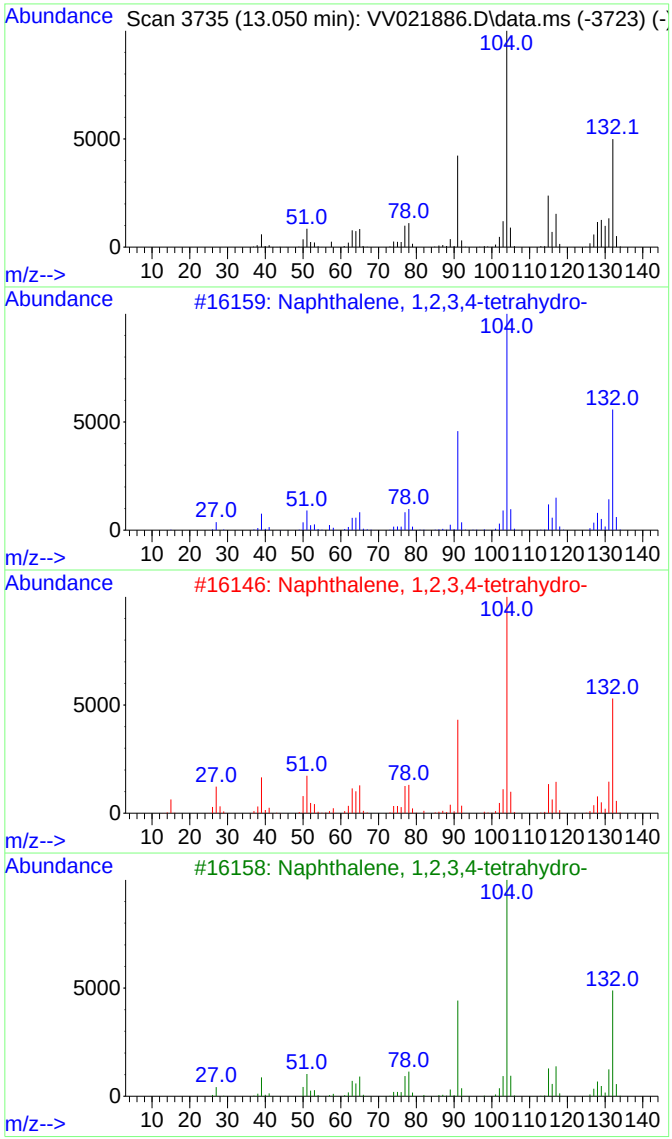
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

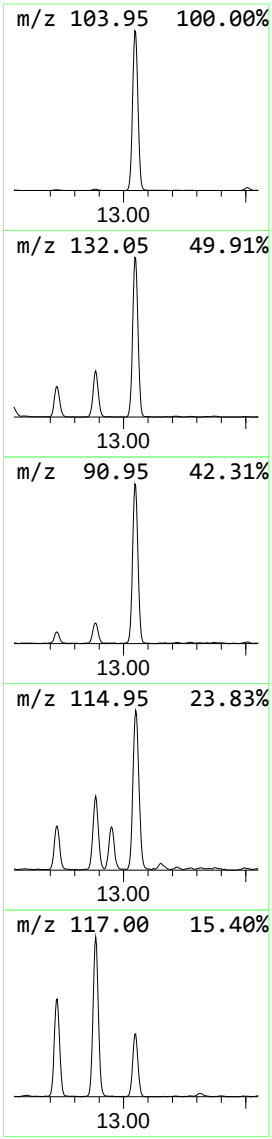
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.050	45.53 ug/L	1210920	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





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Data File : VV021886.D
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Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

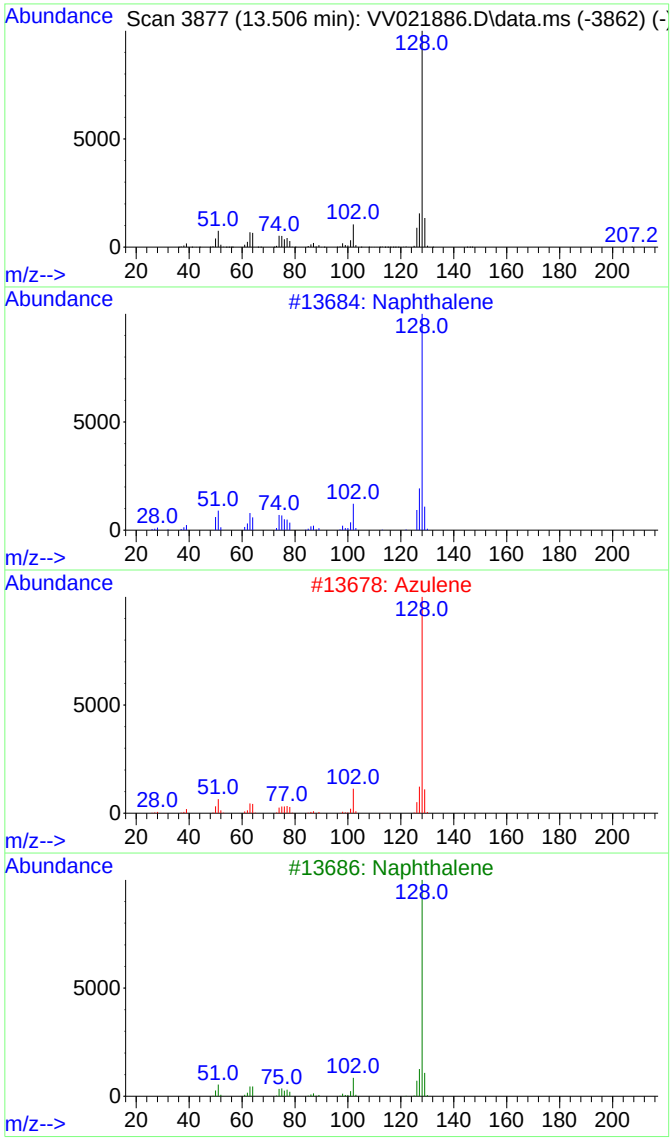
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

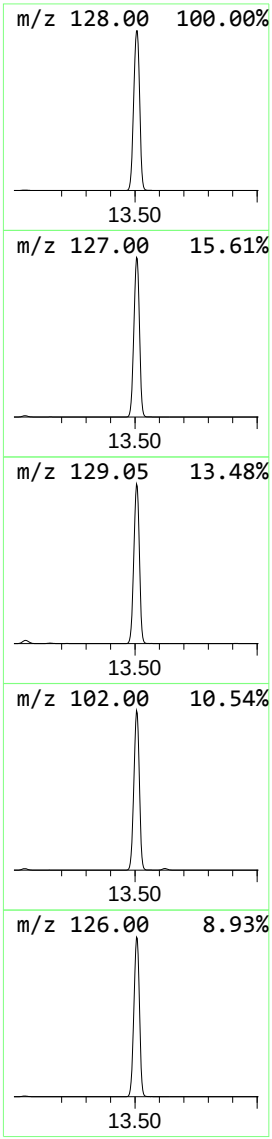
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.506	1121.97 ug/L	29838700	1,4-Dichlorobenzene-d4	11.252

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





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

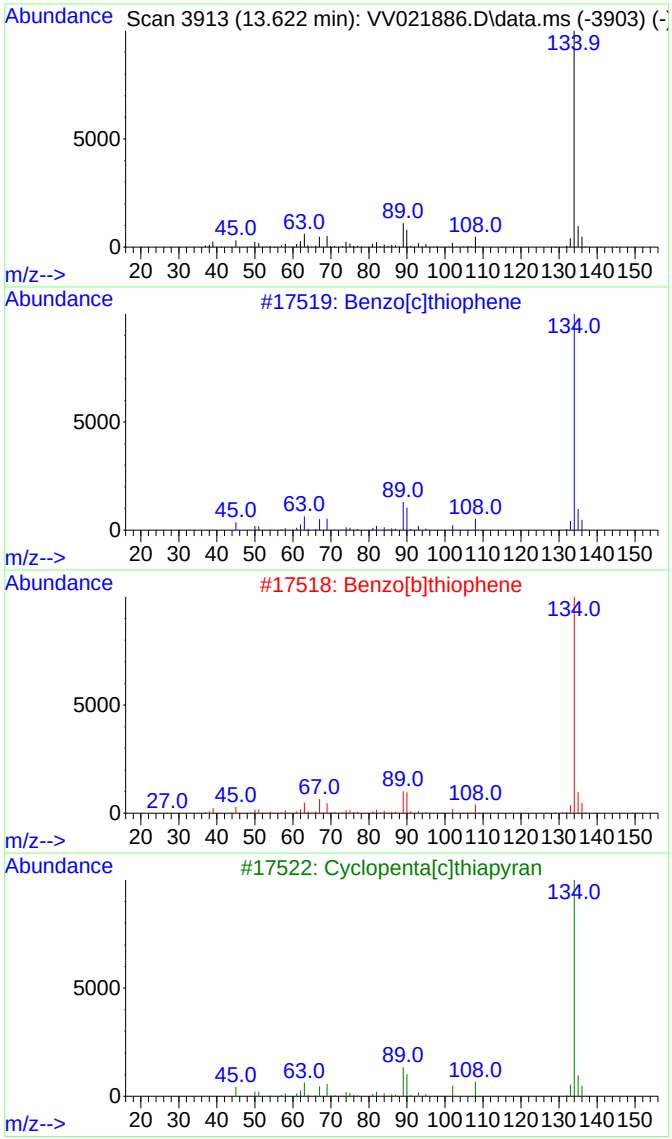
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

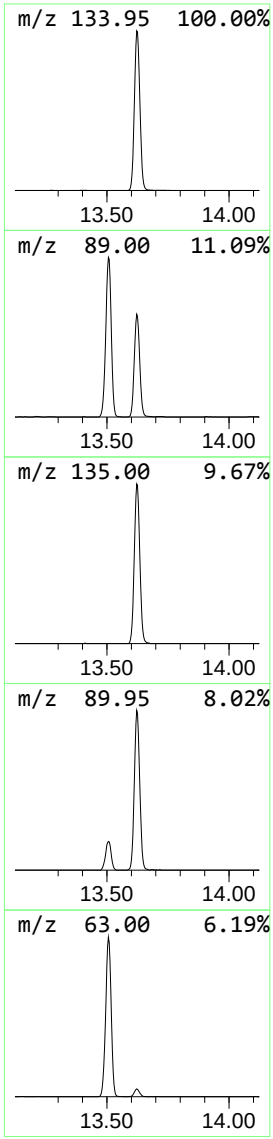
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzo[c]thiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	54.62 ug/L	1452490	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	91





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

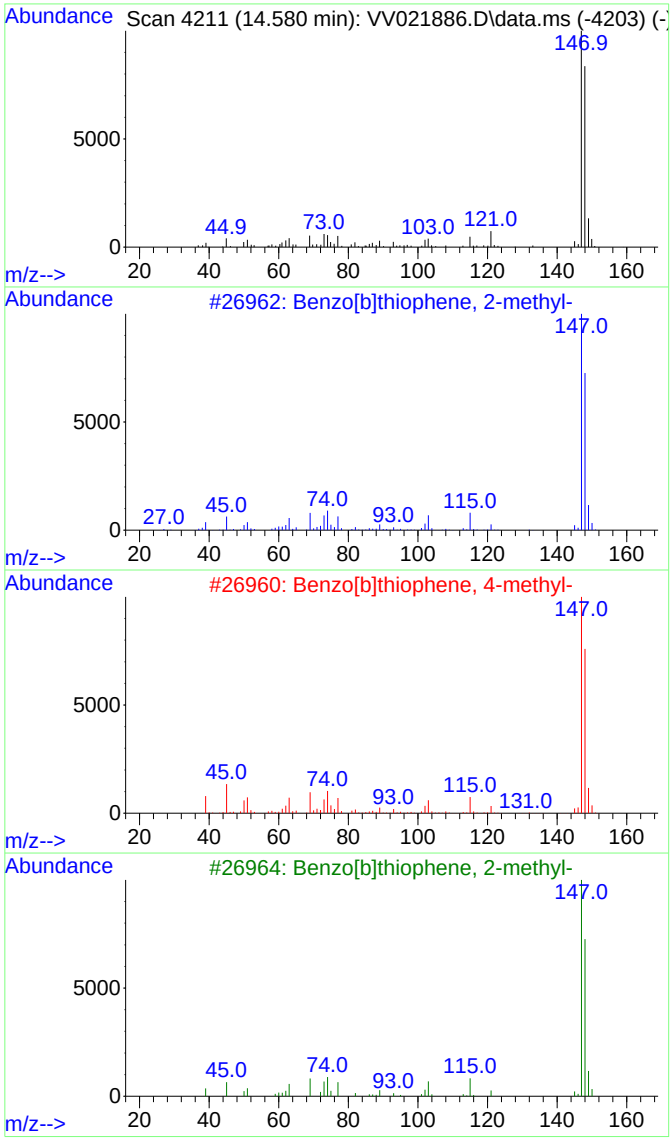
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

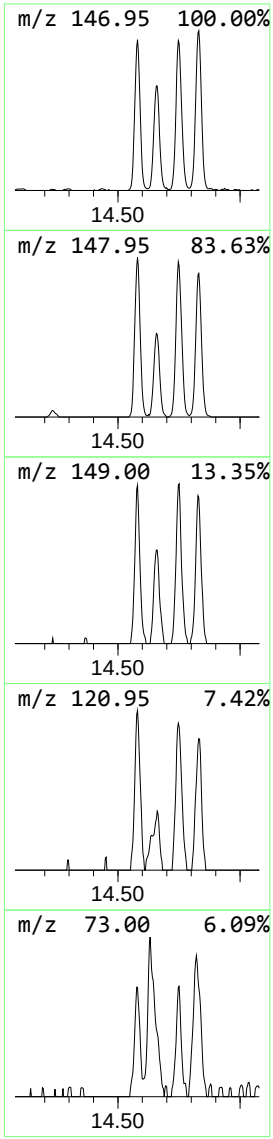
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzo[b]thiophene, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.580	3.94 ug/L	104779	1,4-Dichlorobenzene-d4	11.252

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





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

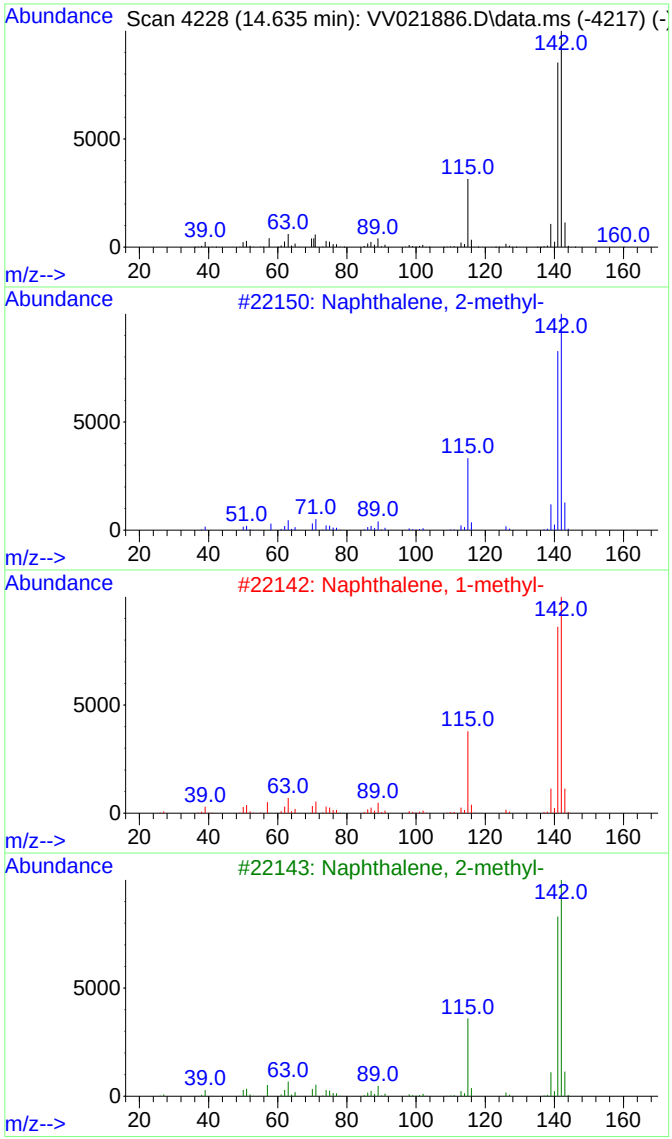
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

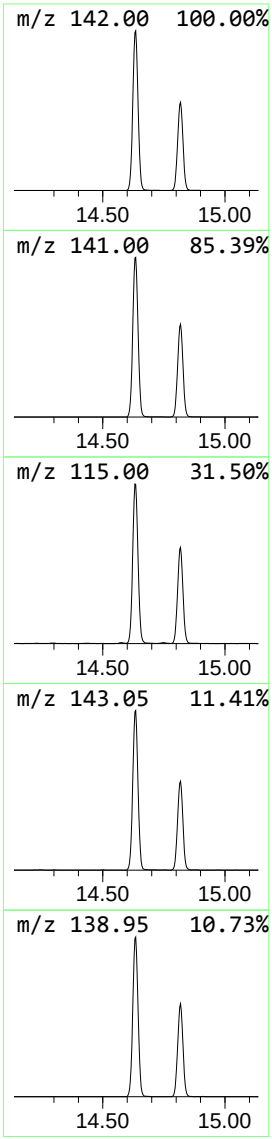
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Naphthalene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.635	147.46 ug/L	3921670	1,4-Dichlorobenzene-d4	11.252

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

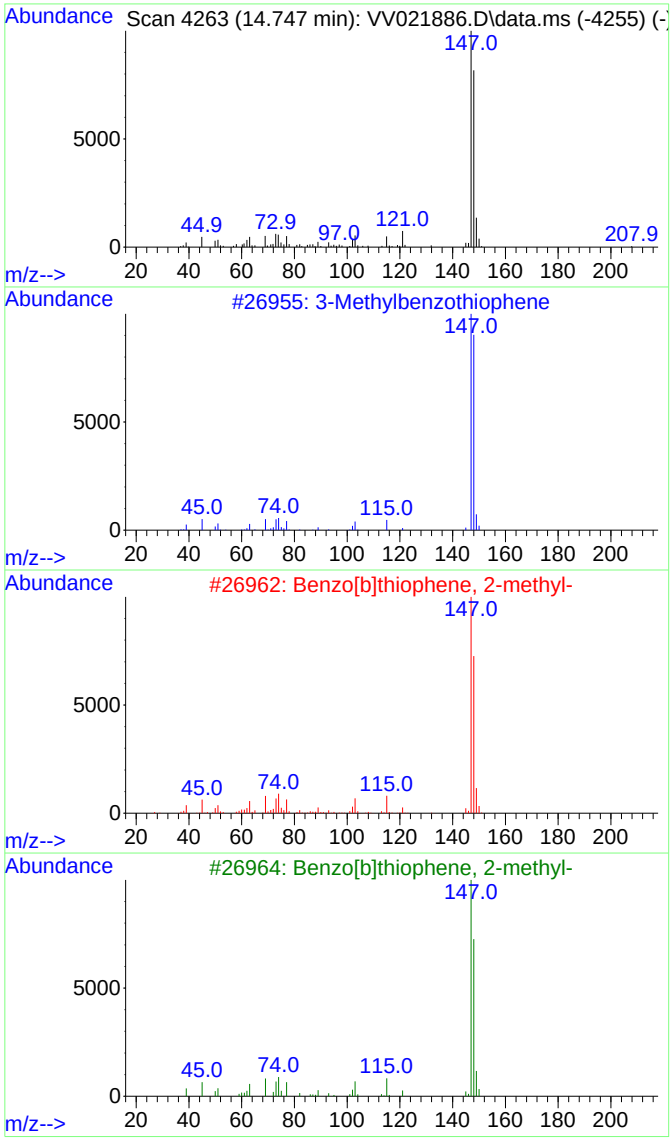
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

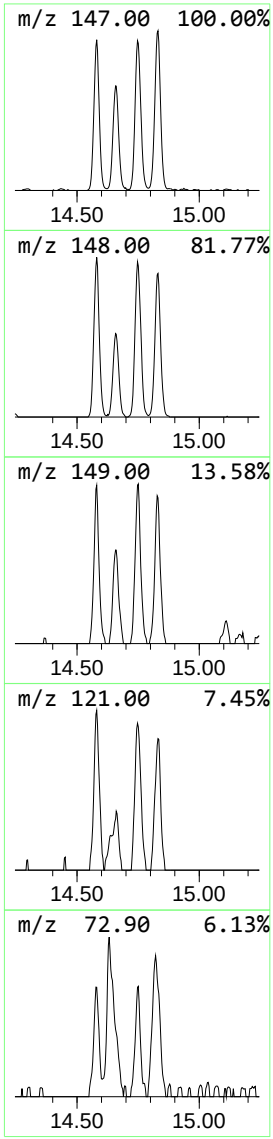
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 3-Methylbenzothiophene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.747	4.34 ug/L	115301	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylbenzothiophene	148	C9H8S	001455-18-1	95
2	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
3	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
4	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	91
5	3-Methylbenzothiophene	148	C9H8S	001455-18-1	91





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Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

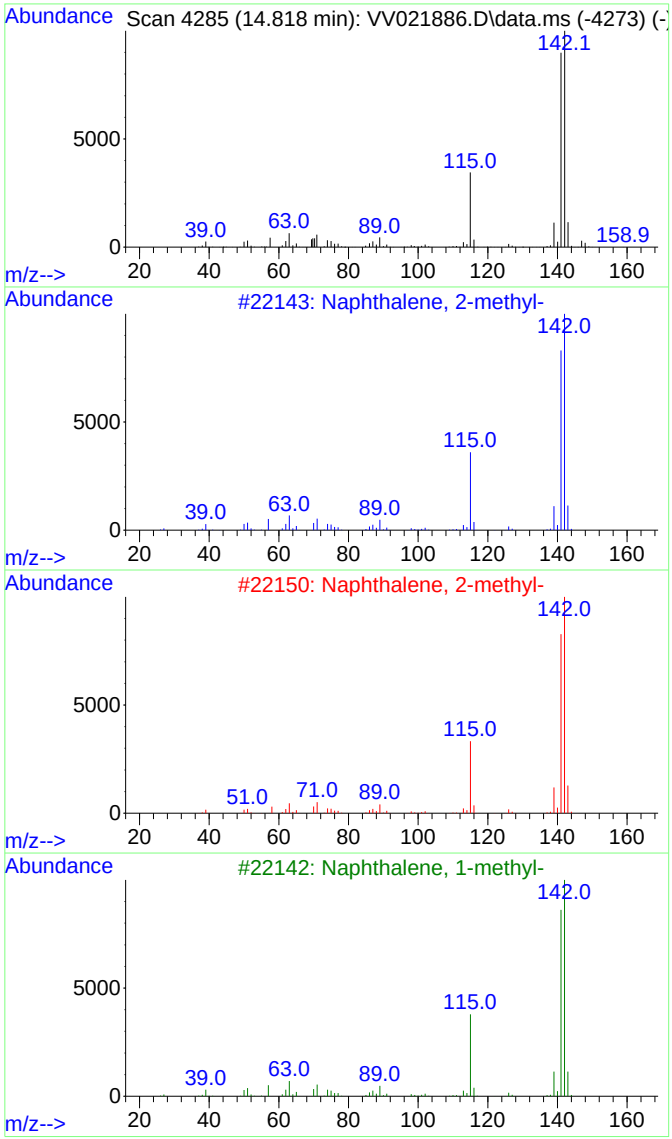
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

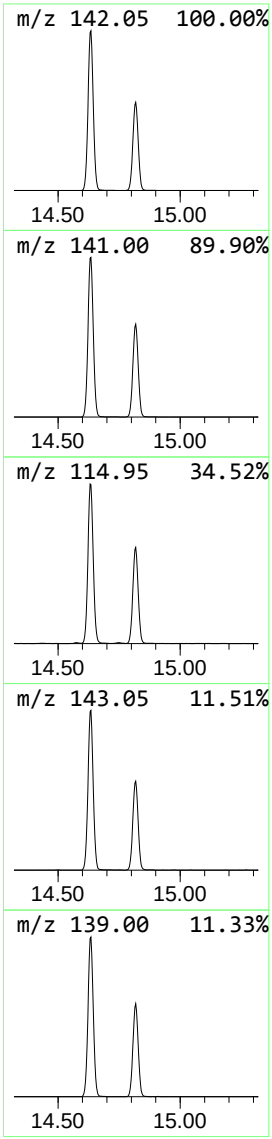
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.818	87.85 ug/L	2336430	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

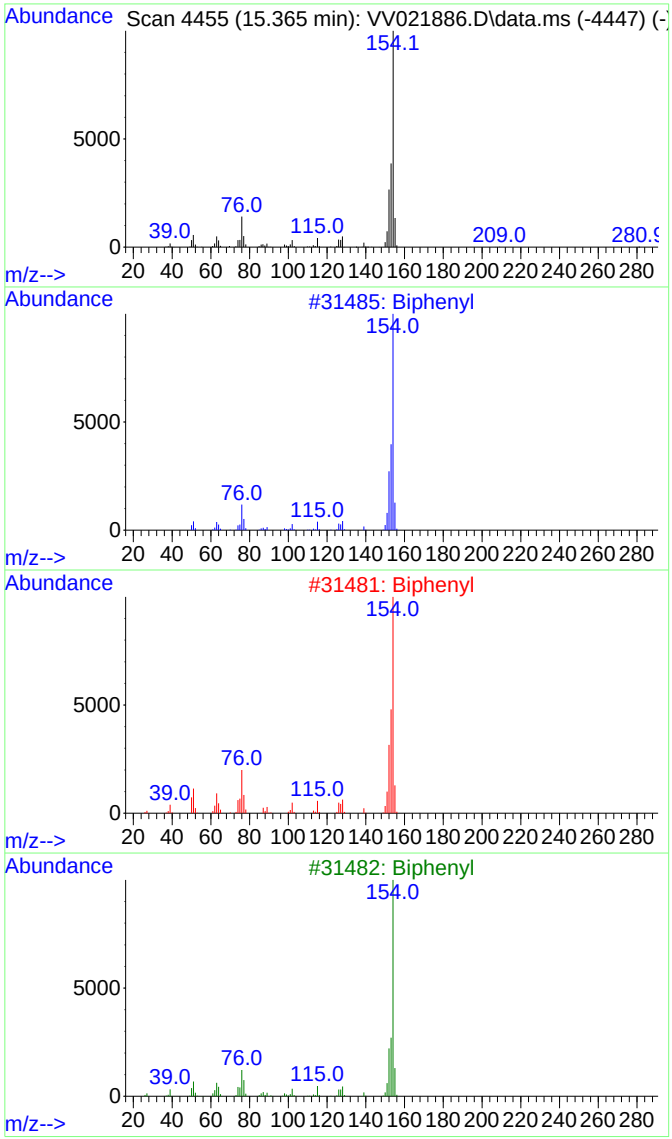
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

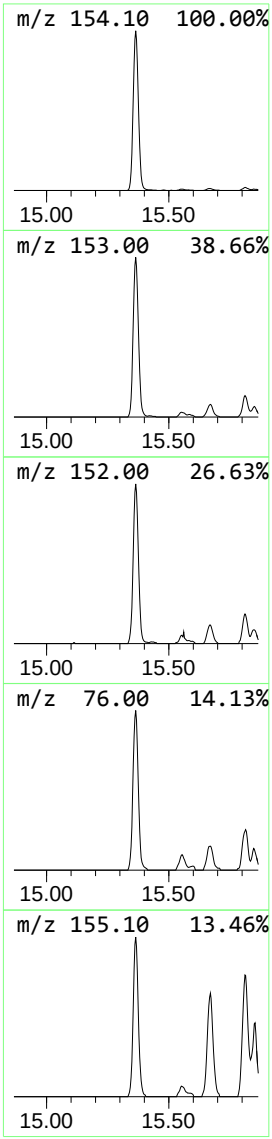
****

Peak Number 22 Biphenyl Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.365	11.75 ug/L	312367	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	93
3	Biphenyl	154	C12H10	000092-52-4	92
4	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	87
5	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

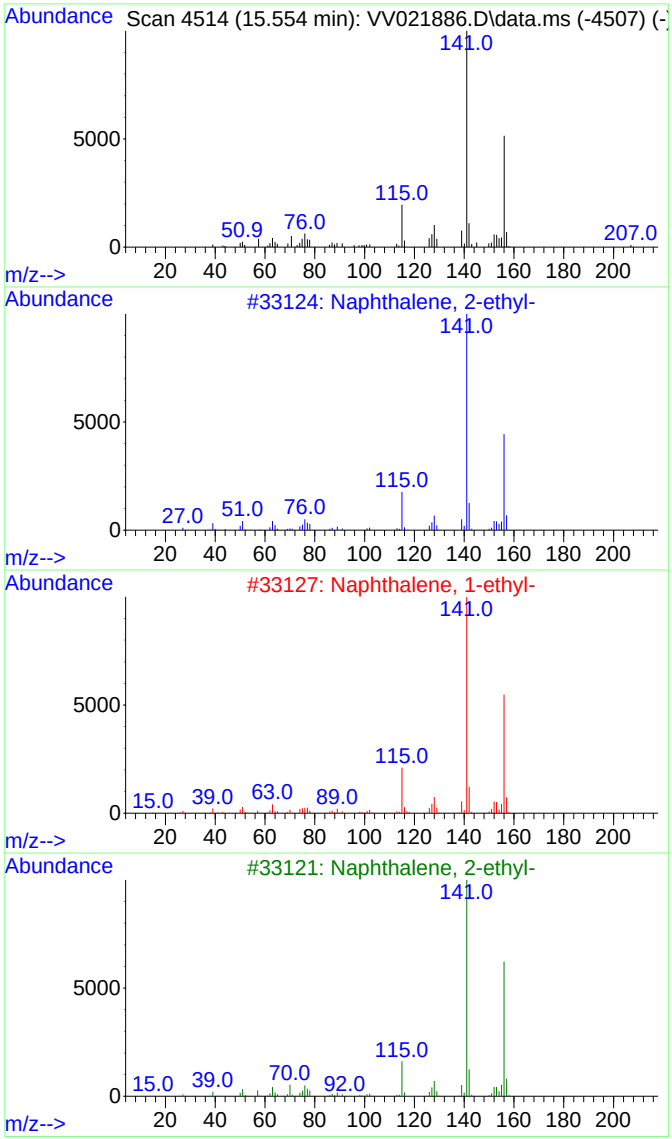
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

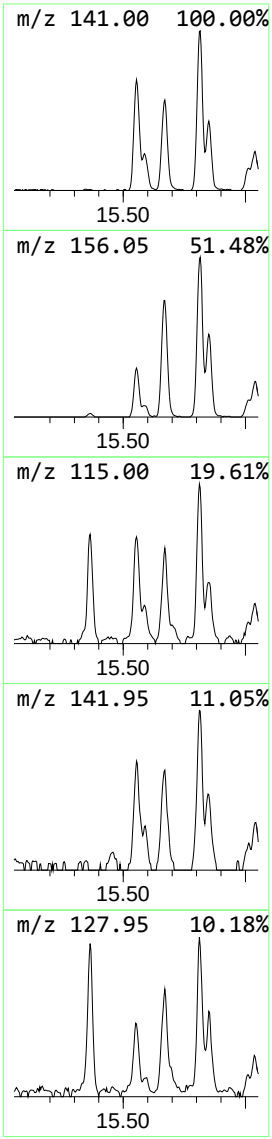
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Naphthalene, 2-ethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.554	2.53 ug/L	67323	1,4-Dichlorobenzene-d4	11.252

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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

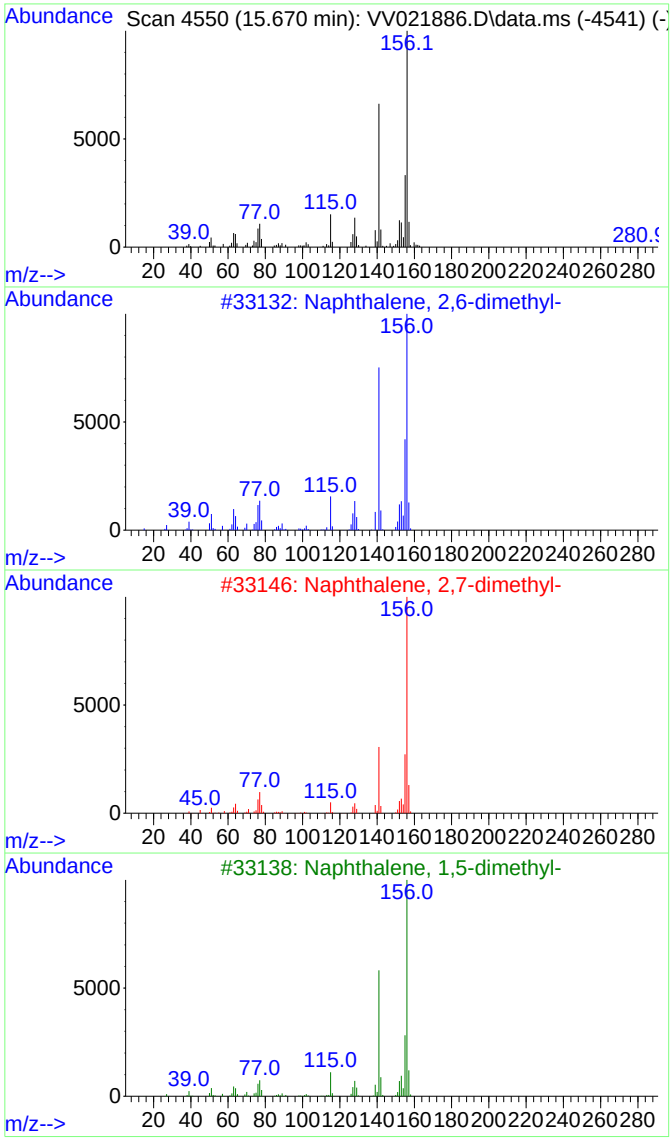
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

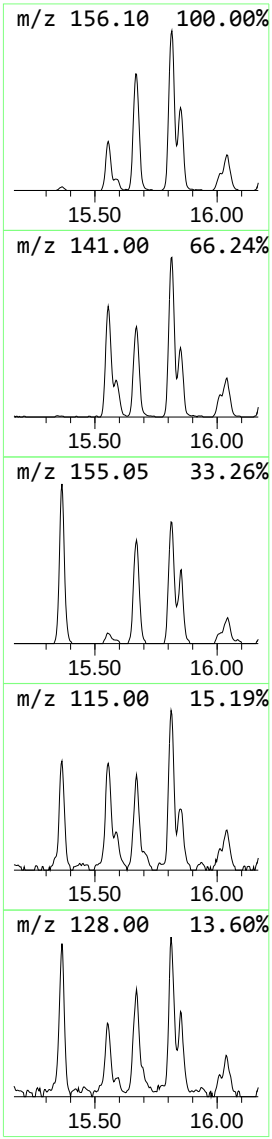
****

Peak Number 24 Naphthalene, 2,6-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.670	5.82 ug/L	154687	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, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

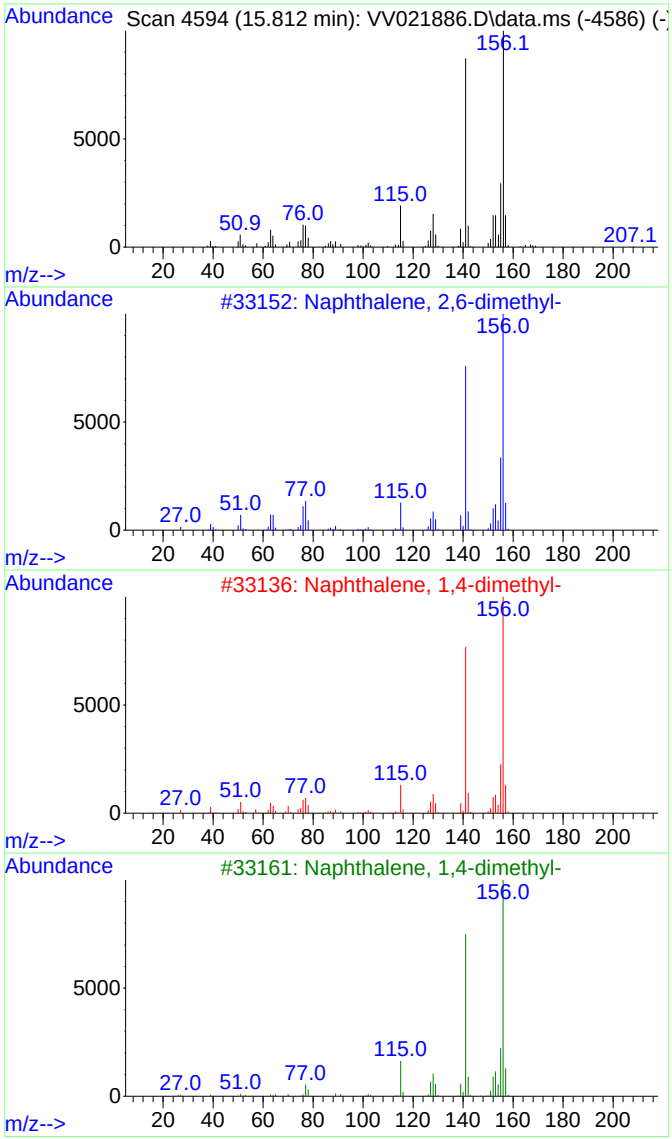
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

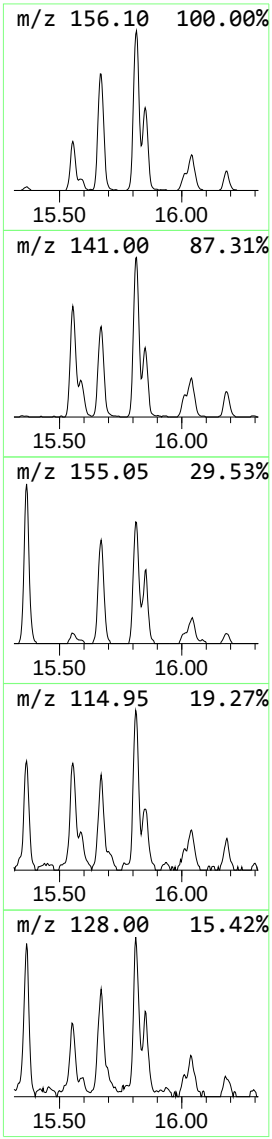
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Naphthalene, 1,4-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.812	6.57 ug/L	174695	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	97
2	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	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, 2,6-dimethyl-	156	C12H12	000581-42-0	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKA9DL

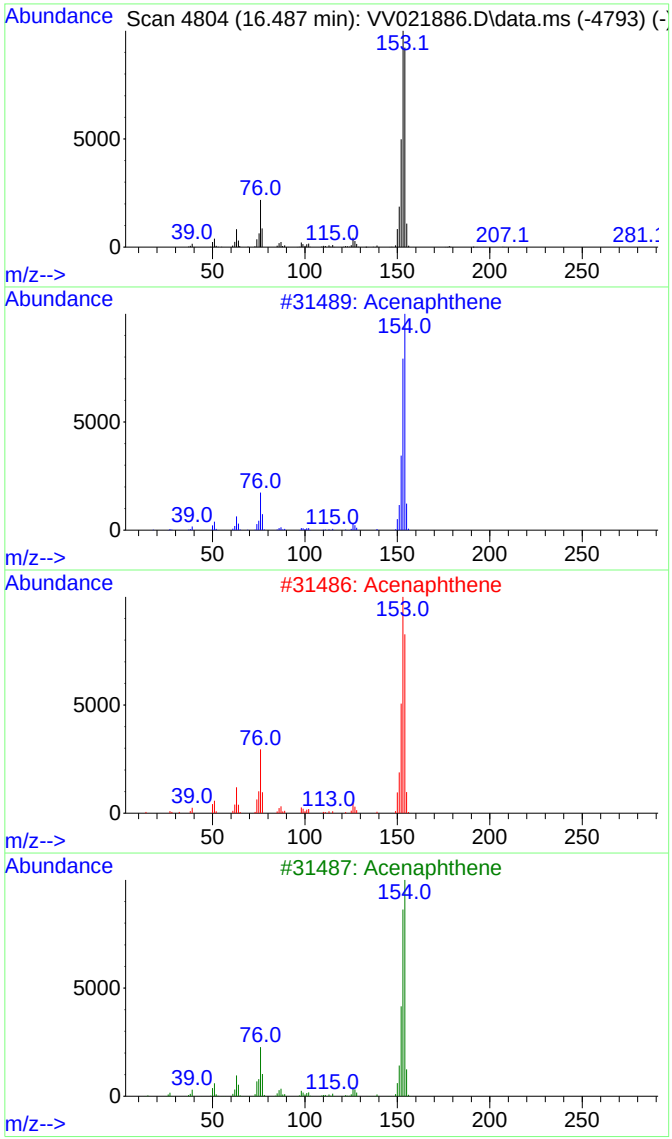
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

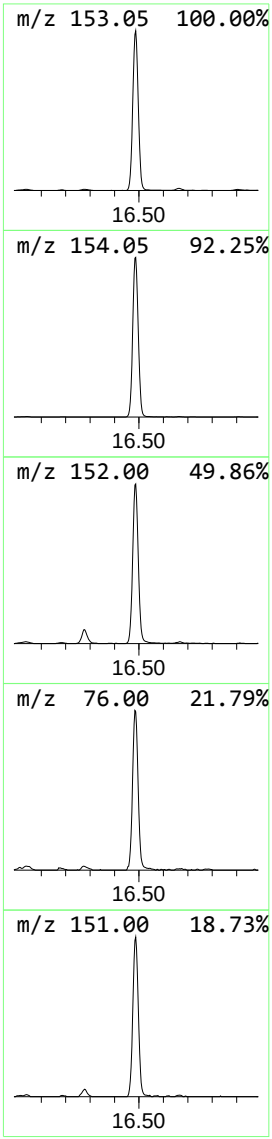
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Acenaphthene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.487	29.43 ug/L	782614	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	74
4	Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	59
5	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021886.D
Acq On : 19 Aug 2021 13:12
Operator : SY/MD
Sample : M3422-08DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKA9DL

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
Benzene, 1-ethy...	10.452	6.1	ug/L	161596	3	11.252	1329750	50.0
Benzene, 1-ethy...	10.741	2.7	ug/L	71562	3	11.252	1329750	50.0
Benzene, 1-ethe...	10.989	3.2	ug/L	83975	3	11.252	1329750	50.0
Benzofuran	11.172	34.7	ug/L	922517	3	11.252	1329750	50.0
Benzene, 1,2,3-...	11.339	6.0	ug/L	158440	3	11.252	1329750	50.0
Indane	11.529	50.2	ug/L	1334680	3	11.252	1329750	50.0
Indene	11.744	408.2	ug/L	10856300	3	11.252	1329750	50.0
1-Phenyl-1-butene	12.117	4.9	ug/L	129791	3	11.252	1329750	50.0
Benzofuran, 2-m...	12.361	6.8	ug/L	180930	3	11.252	1329750	50.0
2-Propenal, 3-p...	12.464	23.2	ug/L	617867	3	11.252	1329750	50.0
1H-Indene, 2,3-...	12.728	8.2	ug/L	217192	3	11.252	1329750	50.0
Benzene, 1-ethe...	12.886	14.4	ug/L	383827	3	11.252	1329750	50.0
Benzene, (1-met...	12.950	3.8	ug/L	101217	3	11.252	1329750	50.0
Naphthalene, 1,...	13.050	45.5	ug/L	1210920	3	11.252	1329750	50.0
Azulene	13.506	1122.0	ug/L	29838700	3	11.252	1329750	50.0
Benzo[c]thiophene	13.622	54.6	ug/L	1452490	3	11.252	1329750	50.0
Benzo[b]thiophe...	14.580	3.9	ug/L	104779	3	11.252	1329750	50.0
Naphthalene, 2-...	14.635	147.5	ug/L	3921670	3	11.252	1329750	50.0
3-Methylbenzoth...	14.747	4.3	ug/L	115301	3	11.252	1329750	50.0
Naphthalene, 1-...	14.818	87.8	ug/L	2336430	3	11.252	1329750	50.0
Biphenyl	15.365	11.8	ug/L	312367	3	11.252	1329750	50.0
Naphthalene, 2-...	15.554	2.5	ug/L	67323	3	11.252	1329750	50.0
Naphthalene, 2,...	15.670	5.8	ug/L	154687	3	11.252	1329750	50.0
Naphthalene, 1,...	15.812	6.6	ug/L	174695	3	11.252	1329750	50.0
Acenaphthene	16.487	29.4	ug/L	782614	3	11.252	1329750	50.0