LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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: VV021887.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.201	45	50	68	rVB	57211	60208	0.19%	0.081%
2	1.304	76	82	96	rVB	214982	217604	0.70%	0.294%
3	1.561	155	162	174	rVB	162342	189919	0.61%	0.256%
4	2.105	321	331	344	rBV	340365	496732	1.61%	0.670%
5	2.349	401	407	409	rBV4	481	522	0.00%	0.001%
6	2.378	414	416	418	rBV	468	254	0.00%	0.000%
7	2.474	443	446	448	rBV	420	250	0.00%	0.000%
8	2.507	448	456	470	rVB2	5405	8839	0.03%	0.012%
9	2.770	528	538	549	rBV3	8595	17839	0.06%	0.024%
10	2.912	580	582	589	rVB2	405	337	0.00%	0.000%
11	2.950	589	594	596	rBV2	506	390	0.00%	0.001%
12	3.044	617	623	626	rVB4	433	411	0.00%	0.001%
13	3.101	638	641	645	rBV	297	283	0.00%	0.000%
14	3.204	671	673	675	rBV	501	214	0.00%	0.000%
15	3.269	690	693	694	rBV2	537	318	0.00%	0.000%
16	3.487	757	761	767	rVV2	324	340	0.00%	0.000%
17	3.539	772	777	781	rVB2	481	454	0.00%	0.001%
18	3.680	818	821	826	rBV2	365	276	0.00%	0.000%
19	3.745	837	841	844	rBV3	510	555	0.00%	0.001%
20	3.876	871	882	915	rBV	125636	338815	1.10%	0.457%
21	4.352	1014	1030	1060	rBV	239477	589528	1.91%	0.796%
22	4.844	1179	1183	1186	rVB2	326	284	0.00%	0.000%
23	4.918	1201	1206	1209	rBV2	411	371	0.00%	0.001%
24	4.934	1209	1211	1219	rVB2	481	415	0.00%	0.001%
25	4.982	1219	1226	1228	rBV2	271	277	0.00%	0.000%
26	5.050	1230	1247	1287	rBV2	595835	1583663	5.13%	2.138%
27	5.436	1365	1367	1371	rVB2	303	213	0.00%	0.000%
28	5.619	1411	1424	1466	rBV	498943	1009814	3.27%	1.363%
29	5.912	1504	1515	1533	rBV3	12533	27878	0.09%	0.038%
30	6.072	1550	1565	1587	rBV	326284	671421	2.17%	0.906%
31	6.297	1633	1635	1637	rBV2	472	249	0.00%	0.000%
32	6.410	1668	1670	1674	rVB2	606	301	0.00%	0.000%
33	6.593	1723	1727	1731	rBV2	314	264	0.00%	0.000%
34	6.648	1739	1744	1748	rBV4	879	1060	0.00%	0.001%
35	6.699	1758	1760	1763	rBV2	333	243	0.00%	0.000%
36	6.989	1839	1850	1875	rBV	190423	344453	1.12%	0.465%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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	7.320	1941	1953	1966	rBV	735612	1273928	4.12%	1.719%
38	7.391	1968	1975	1995	rVB	210857	369271	1.20%	0.498%
39	7.625	2038	2048	2069	rBV2	126626	217525	0.70%	0.294%
40	7.883	2124	2128	2132	rVB3	415	208	0.00%	0.000%
41	7.944	2140	2147	2153	rVV2	791	1053	0.00%	0.001%
42	7.986	2157	2160	2164	rVB4	742	428	0.00%	0.001%
43	8.092	2183	2193	2226	rBV	357435	645082	2.09%	0.871%
44	8.387	2281	2285	2289	rBV3	579	526	0.00%	0.001%
45	8.555	2331	2337	2340	rBV3	288	289	0.00%	0.000%
46	8.606	2351	2353	2355	rBV2	382	227	0.00%	0.000%
47	8.696	2375	2381	2384	rBV3	311	386	0.00%	0.001%
48	8.754	2394	2399	2402	rVB3	303	273	0.00%	0.000%
49	8.854	2419	2430	2456	rBV	745480	1214384	3.93%	1.639%
50	9.014	2469	2480	2505	rBV	899958	1417624	4.59%	1.913%
51	9.140	2508	2519	2543	rVB	427872	702326	2.27%	0.948%
52	9.281	2561	2563	2565	rBV3	471	214	0.00%	0.000%
53	9.320	2568	2575	2585	rVB7	2309	4119	0.01%	0.006%
54	9.445	2611	2614	2618	rVV2	340	252	0.00%	0.000%
55	9.551	2635	2647	2677	rVB3	309053	690186	2.23%	0.932%
56	9.738	2702	2705	2707	rVB	456	277	0.00%	0.000%
57	9.783	2713	2719	2727	rVB6	1541	2254	0.01%	0.003%
58	9.834	2733	2735	2737	rVB2	512	232	0.00%	0.000%
59	9.937	2758	2767	2779	rVB2	8404	13428	0.04%	0.018%
60	10.005	2784	2788	2792	rVB	253	237	0.00%	0.000%
61	10.034	2792	2797	2799	rBV2	876	884	0.00%	0.001%
62	10.130	2819	2827	2839	rVB3	4544	6332	0.02%	0.009%
63	10.217	2842	2854	2870	rBV	470901	753556	2.44%	1.017%
64	10.358	2892	2898	2913	rVB	14001	22268	0.07%	0.030%
65	10.452	2916	2927	2945	rBV	78503	165804	0.54%	0.224%
66	10.542	2946	2955	2973	rVB	76571	120583	0.39%	0.163%
67	10.654	2987	2990	2991	rBV3	865	414	0.00%	0.001%
68	10.741	3008	3017	3025	rBV	38323	60133	0.19%	0.081%
69	10.918	3057	3072	3085	rBV	236965	378581	1.23%	0.511%
70	10.989	3086	3094	3103	rBV	57518	84819	0.27%	0.114%
71	11.172	3138	3151	3165	rBV	599658	946671	3.07%	1.278%
72	11.249	3165	3175	3194	rVV	843628	1331791	4.31%	1.798%
73	11.339	3195	3203	3216	rVB	106502	165680	0.54%	0.224%
74	11.529	3251	3262	3282	rBV	884129	1362543	4.41%	1.839%
75	11.625	3282	3292	3315	rVB	804542	1275669	4.13%	1.722%
76	11.747	3317	3330	3349	rBV	7105497	10997510	35.61%	14.844%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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	12.021	3407	3415	3419	rBV2	25535	37856	0.12%	0.051%
78	12.117	3435	3445	3464	rBV	85699	135894	0.44%	0.183%
79	12.255	3480	3488	3496	rBV10	5225	8172	0.03%	0.011%
80	12.362	3513	3521	3532	rBV	122290	185551	0.60%	0.250%
81	12.464	3544	3553	3563	rBV2	351299	629763	2.04%	0.850%
82	12.728	3625	3635	3646	rBV	147873	219859	0.71%	0.297%
83	12.886	3668	3684	3695	rBV2	255621	393171	1.27%	0.531%
84	12.950	3697	3704	3716	rVB2	66050	102820	0.33%	0.139%
85	13.046	3723	3734	3750	rVB2	784364	1237545	4.01%	1.670%
86	13.156	3761	3768	3779	rVB5	19597	34085	0.11%	0.046%
87	13.506	3862	3877	3902	rBV	19158227	30883874	100.00%	41.685%
88	13.622	3903	3913	3933	rVB	975737	1518485	4.92%	2.050%
89	13.911	3996	4003	4016	rBV2	18018	29155	0.09%	0.039%
90	14.098	4051	4061	4062	rBV3	21201	28473	0.09%	0.038%
91	14.435	4160	4166	4177	rVB7	10502	18177	0.06%	0.025%
92	14.580	4203	4211	4217	rBV	76908	111026	0.36%	0.150%
93	14.632	4217	4227	4253	rVB	2839359	4282891	13.87%	5.781%
94	14.747	4255	4263	4273	rBV	79663	124565	0.40%	0.168%
95	14.818	4273	4285	4301	rBV	1573439	2494872	8.08%	3.367%
96	15.365	4446	4455	4468	rVB	251384	375051	1.21%	0.506%
97	15.667	4540	4549	4573	rVB	95230	183296	0.59%	0.247%
98	15.812	4585	4594	4601	rBV	139182	213362	0.69%	0.288%
99	16.487	4793	4804	4828	rBV	595151	938390	3.04%	1.267%
100	16.786	4890	4897	4916	rVB	67463	109065	0.35%	0.147%

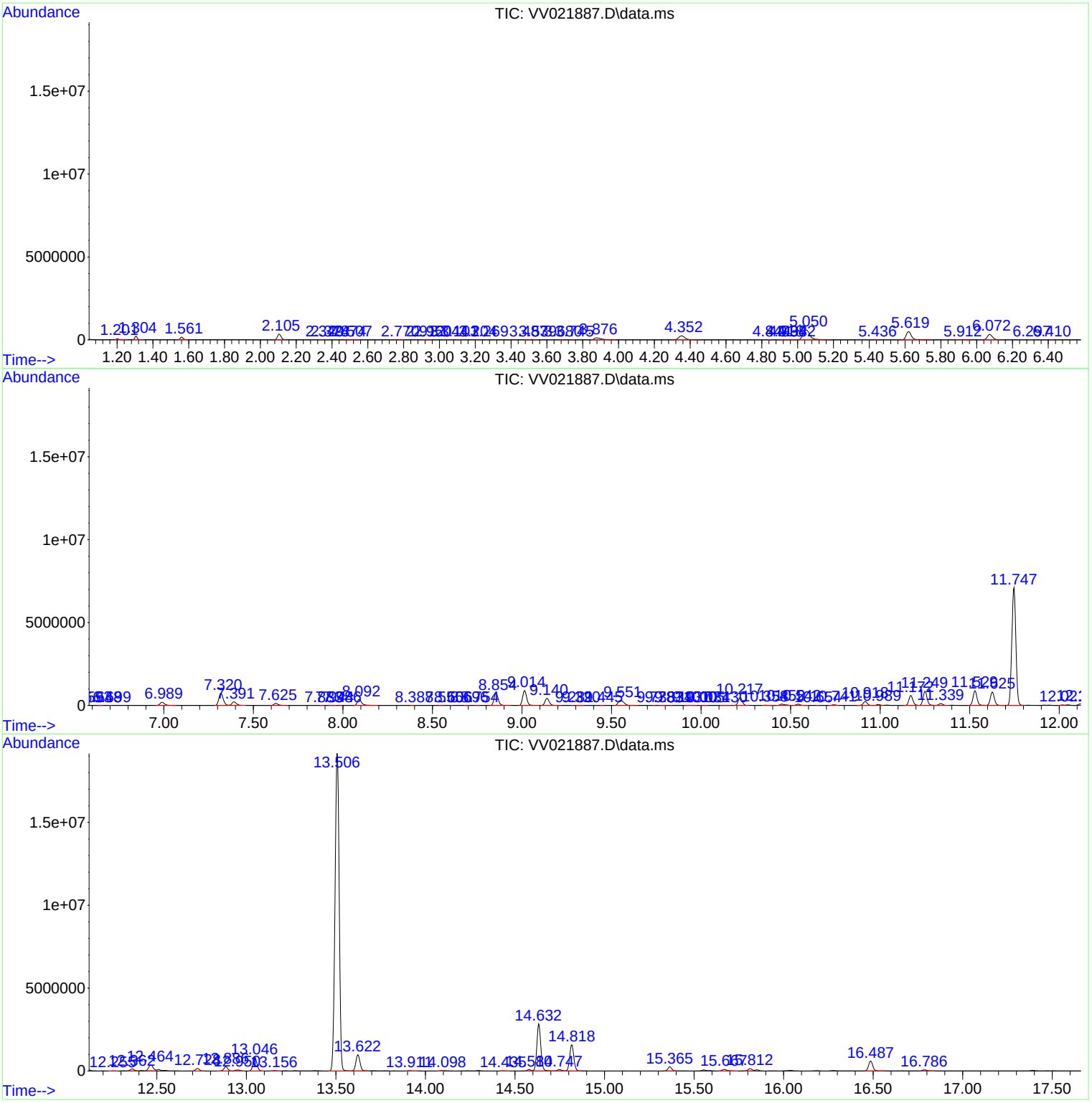
Sum of corrected areas: 74088359

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

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 : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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

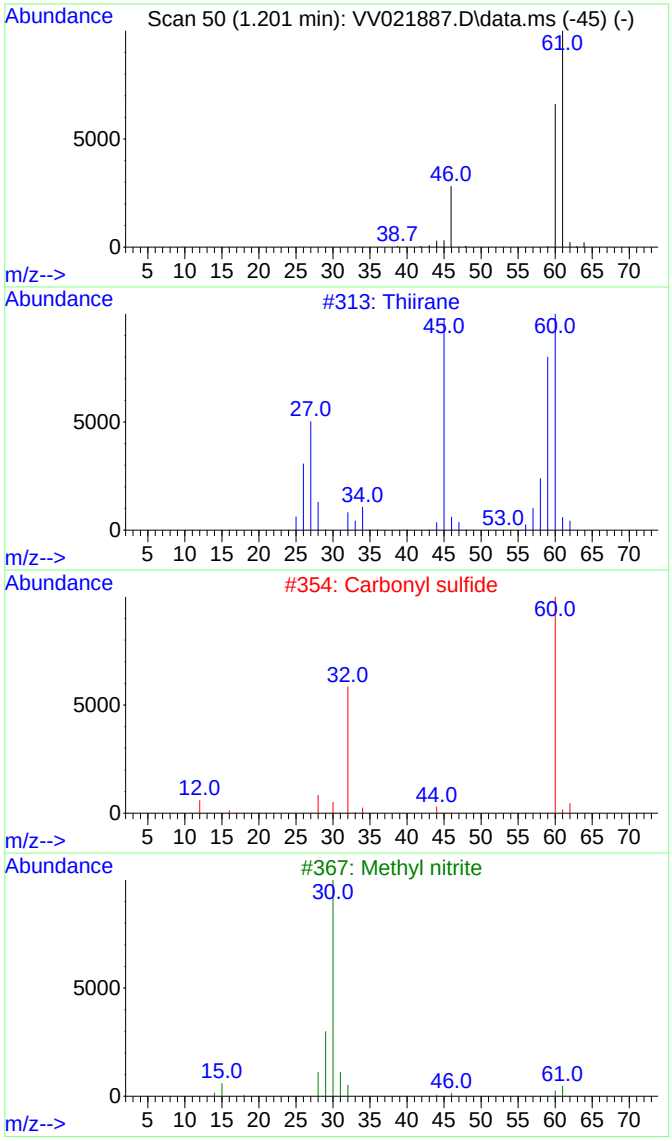
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

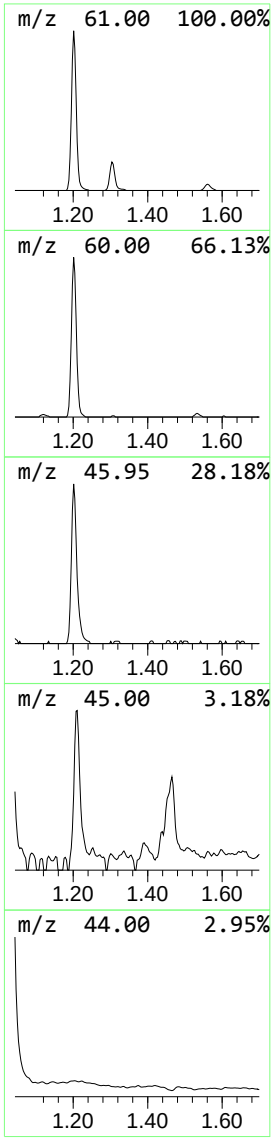
****

Peak Number 1 unknown-01 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.201	2.98 ug/L	60208	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiirane	60	C2H4S	000420-12-2	5
2	Carbonyl sulfide	60	COS	000463-58-1	5
3	Methyl nitrite	61	CH3NO2	000624-91-9	5
4	o-Ethylhydroxylamine	61	C2H7NO	000624-86-2	4
5	Methyl formate	60	C2H4O2	000107-31-3	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

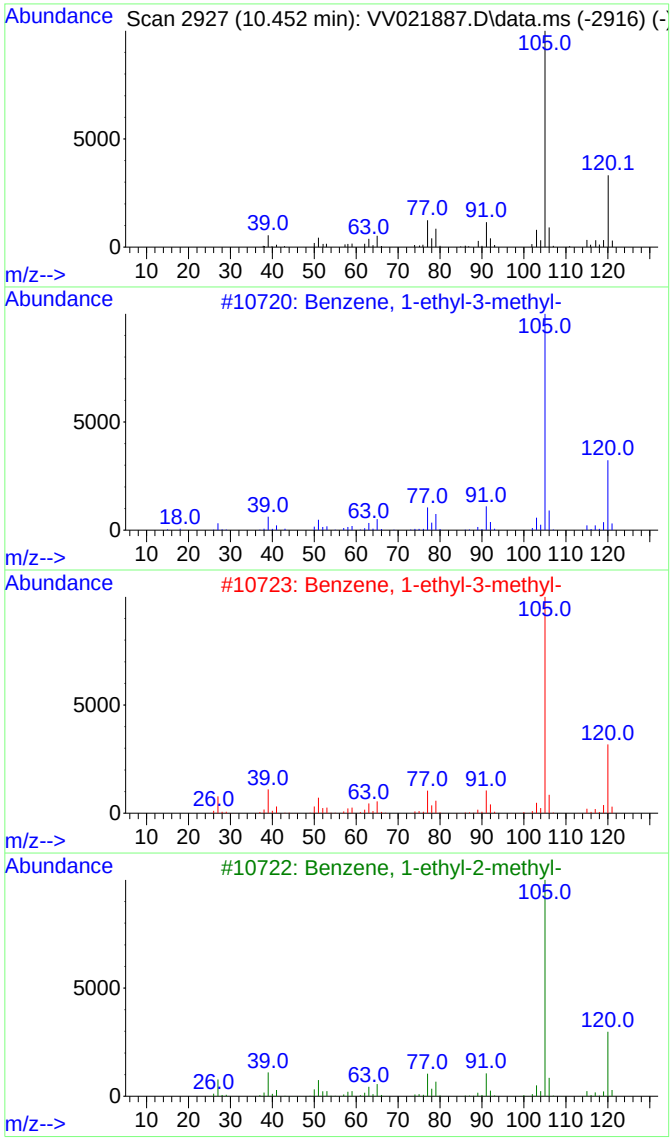
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

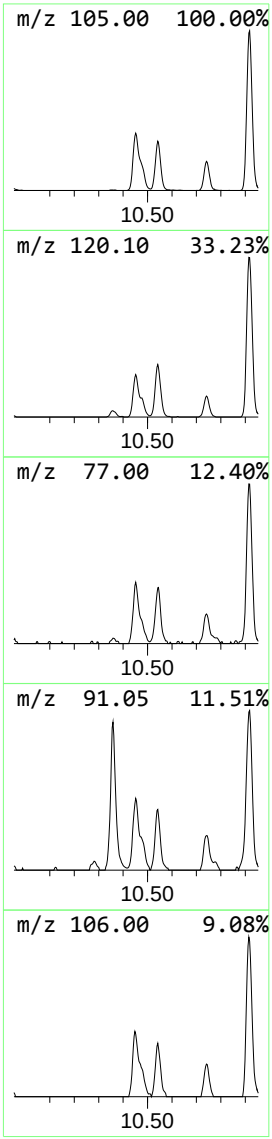
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.452	6.22 ug/L	165804	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

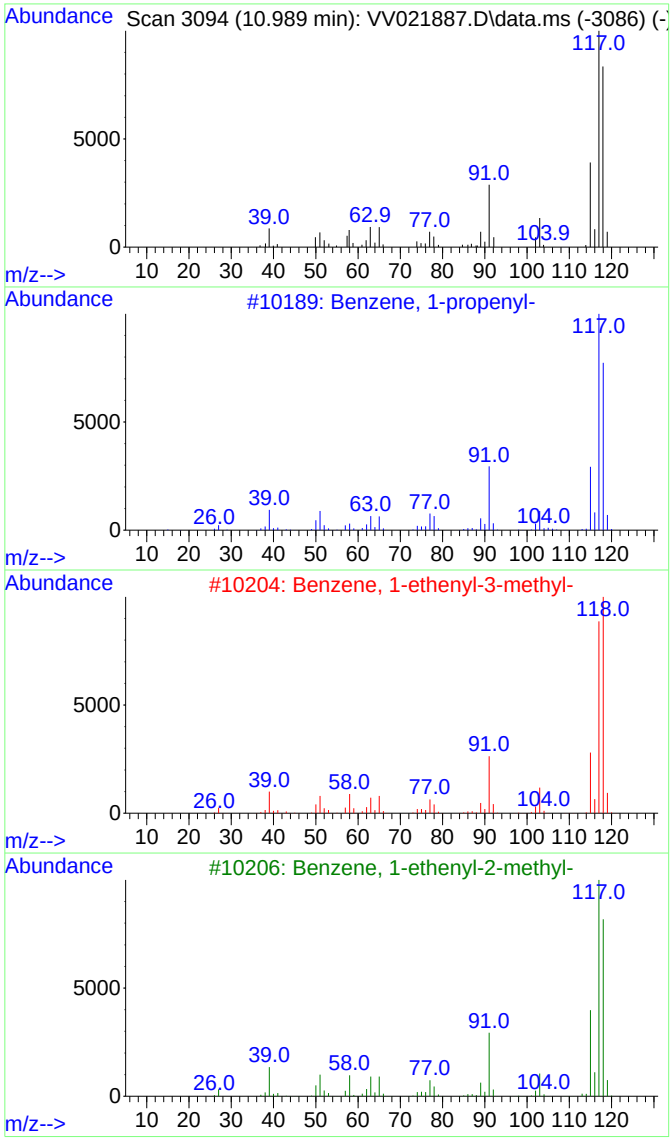
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

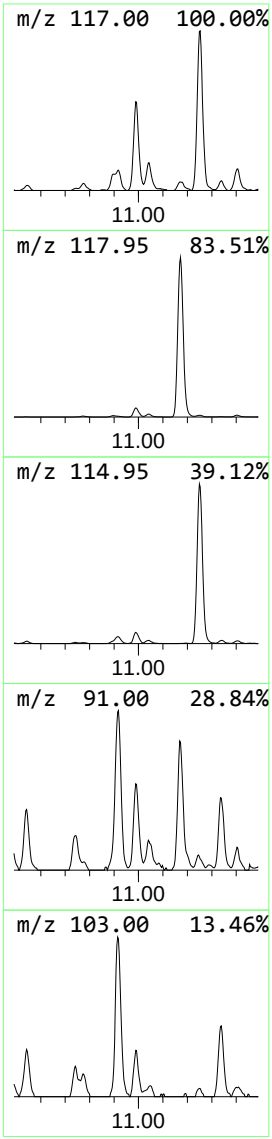
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-propenyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.989	3.18 ug/L	84819	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	95
2	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	95
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
5	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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

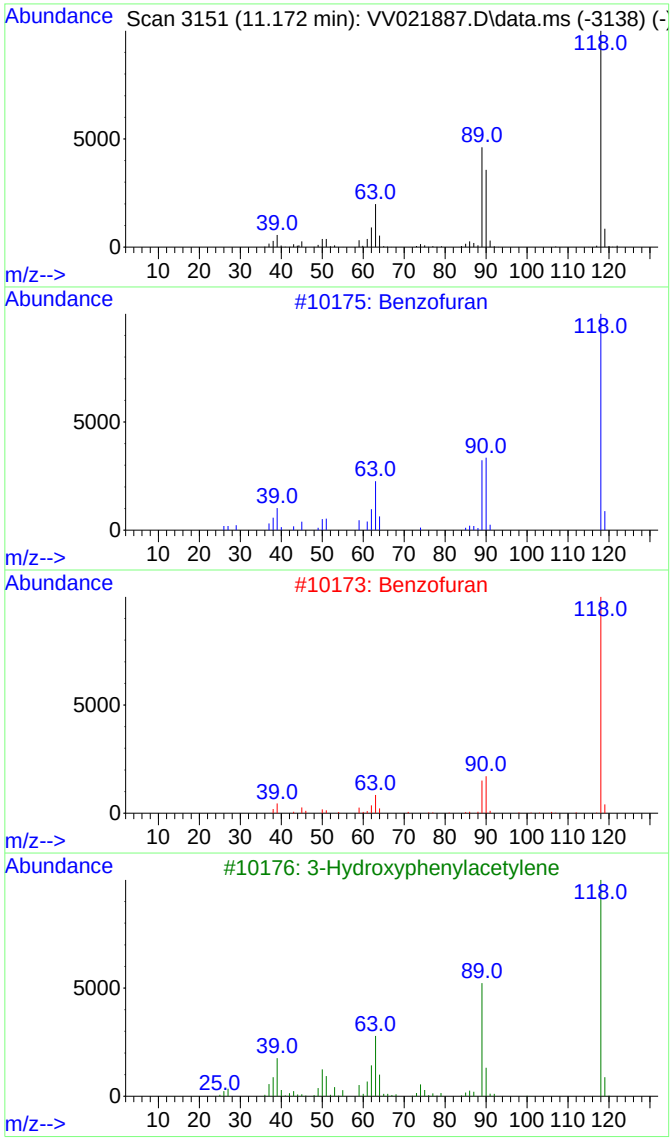
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

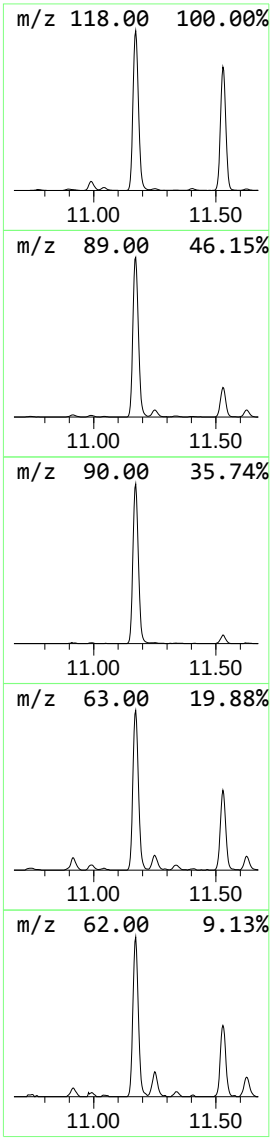
****

Peak Number 5 Benzofuran Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.172	35.54 ug/L	946671	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	95
2	Benzofuran	118	C8H6O	000271-89-6	91
3	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	91
4	Benzofuran	118	C8H6O	000271-89-6	90
5	Benzofuran	118	C8H6O	000271-89-6	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

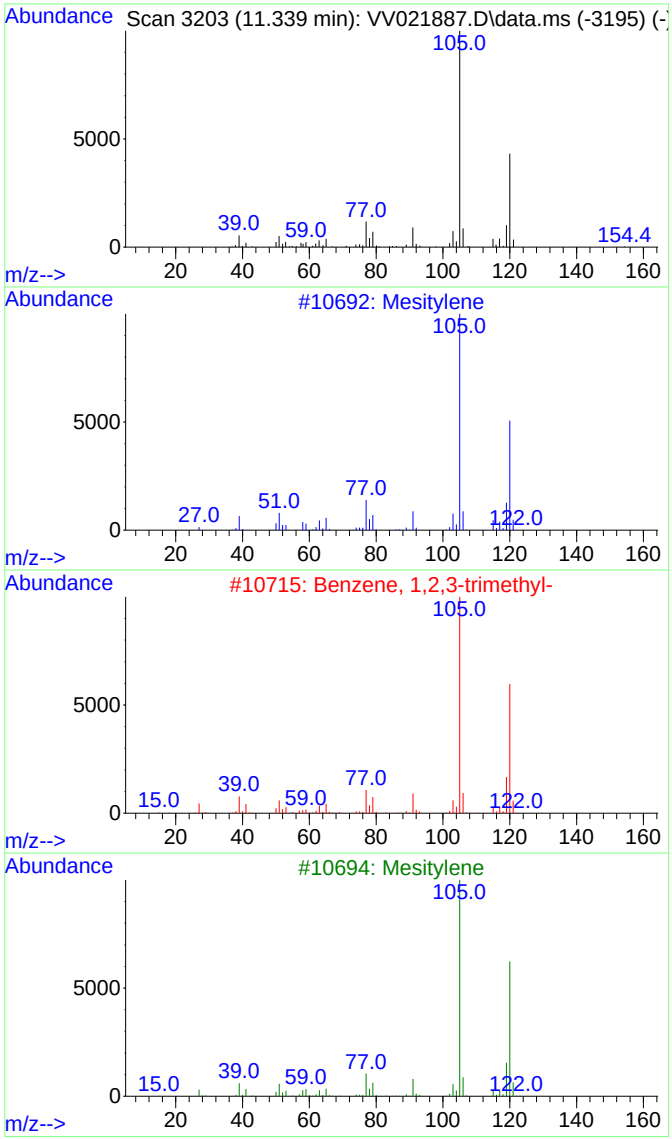
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

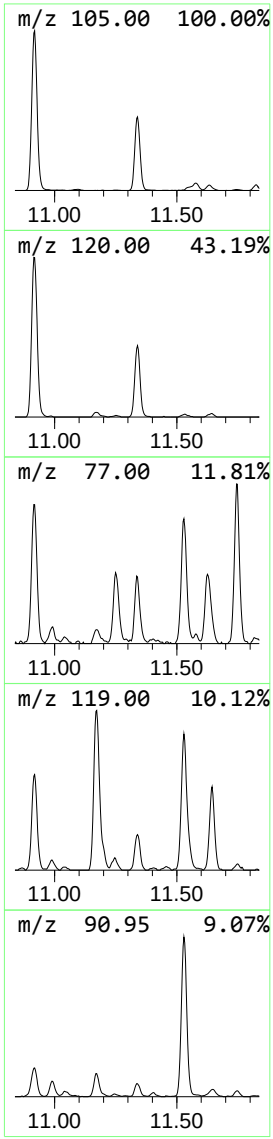
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.339	6.22 ug/L	165680	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	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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

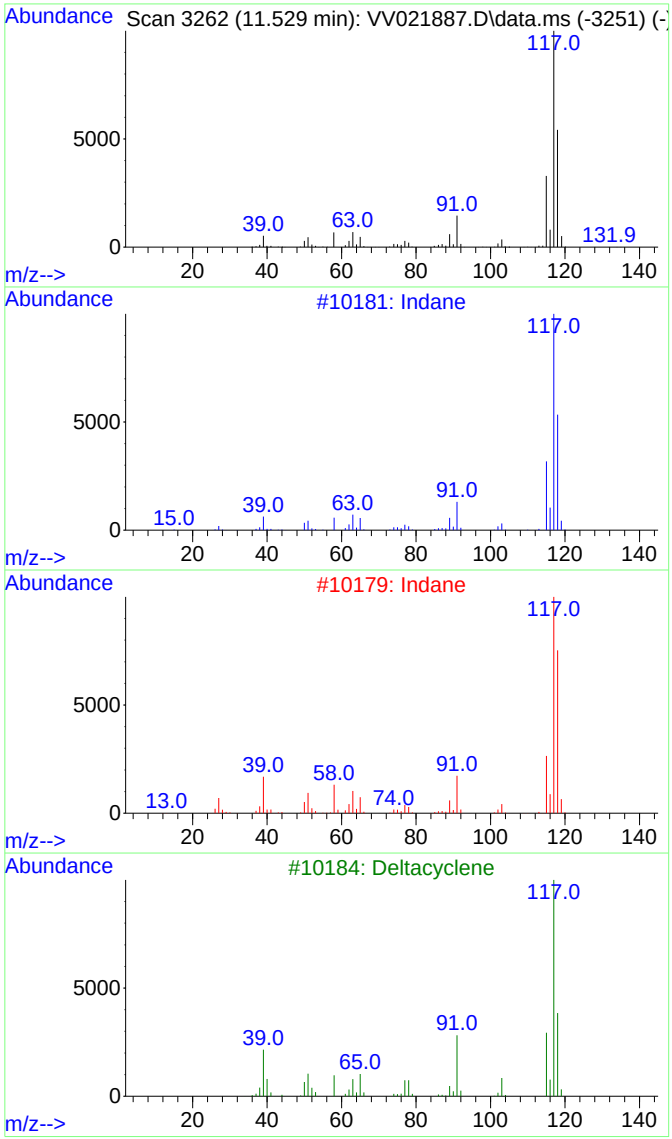
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

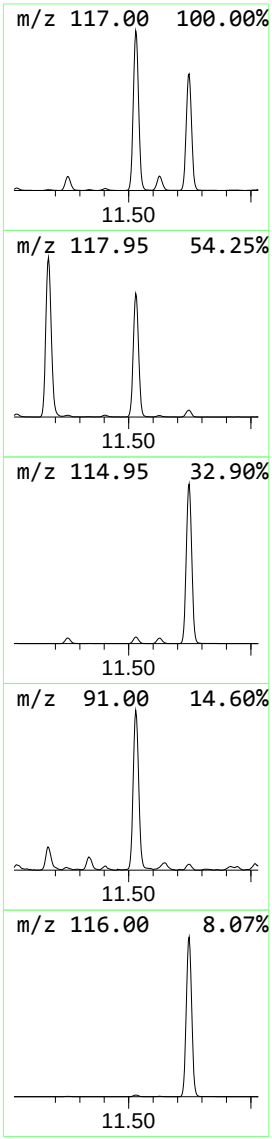
****

Peak Number 7 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	51.15 ug/L	1362540	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	Benzene, 1-propenyl-	118	C9H10	000637-50-3	72
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

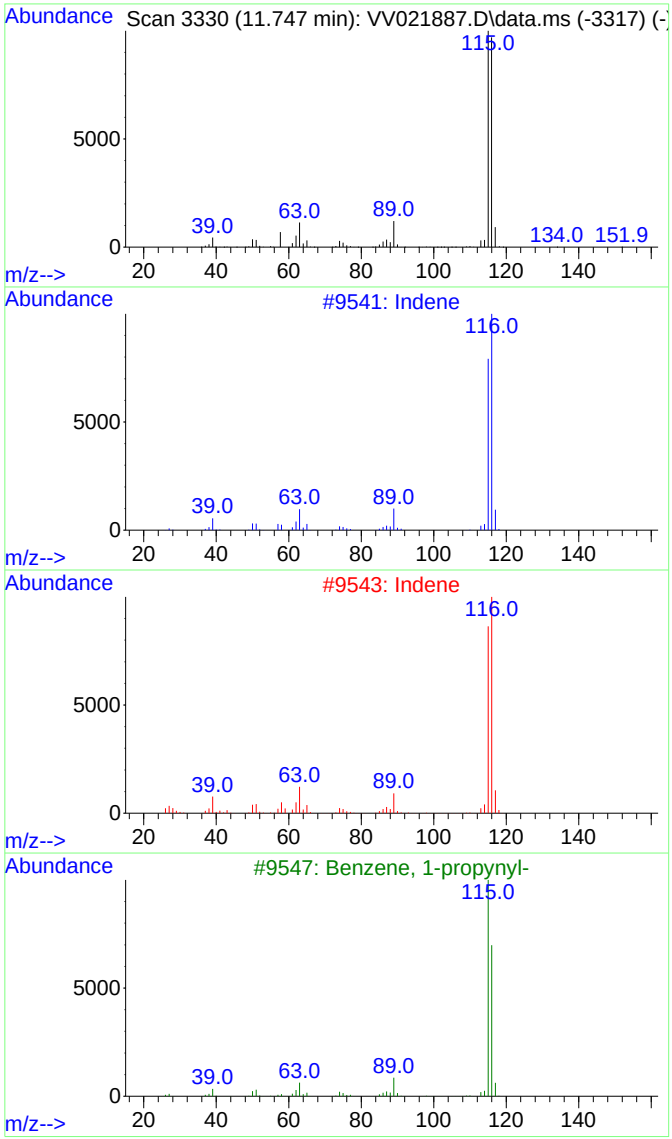
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

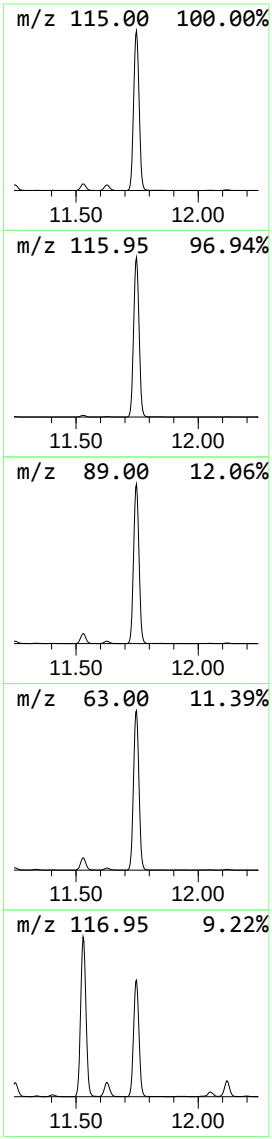
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.747	412.88 ug/L	10997500	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-ethynyl-4-methyl-	116	C9H8	000766-97-2	91





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

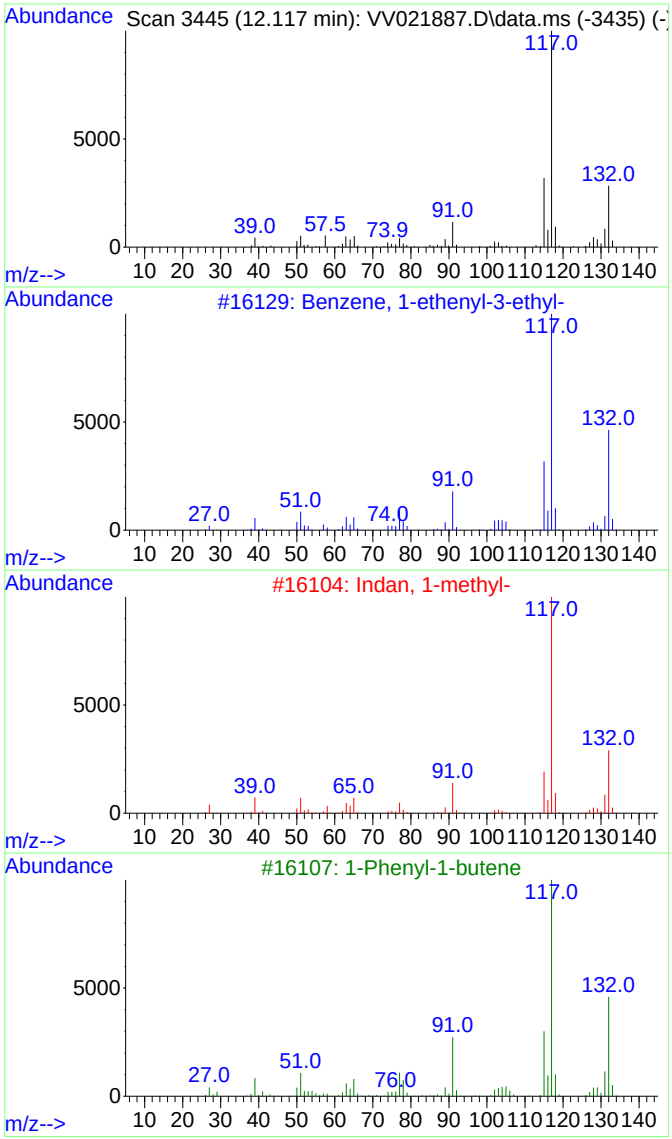
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

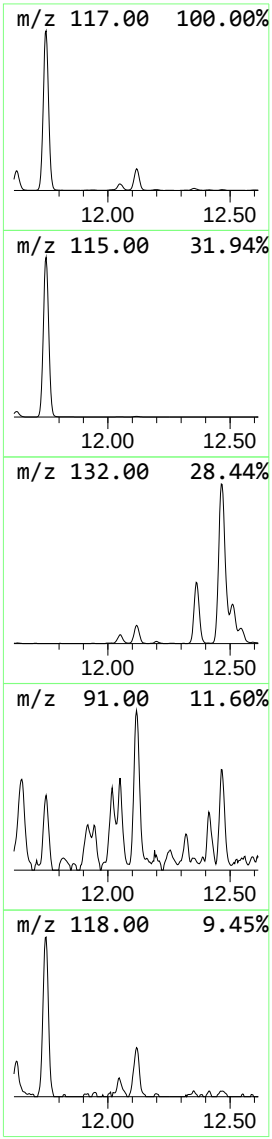
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	5.10 ug/L	135894	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	90
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

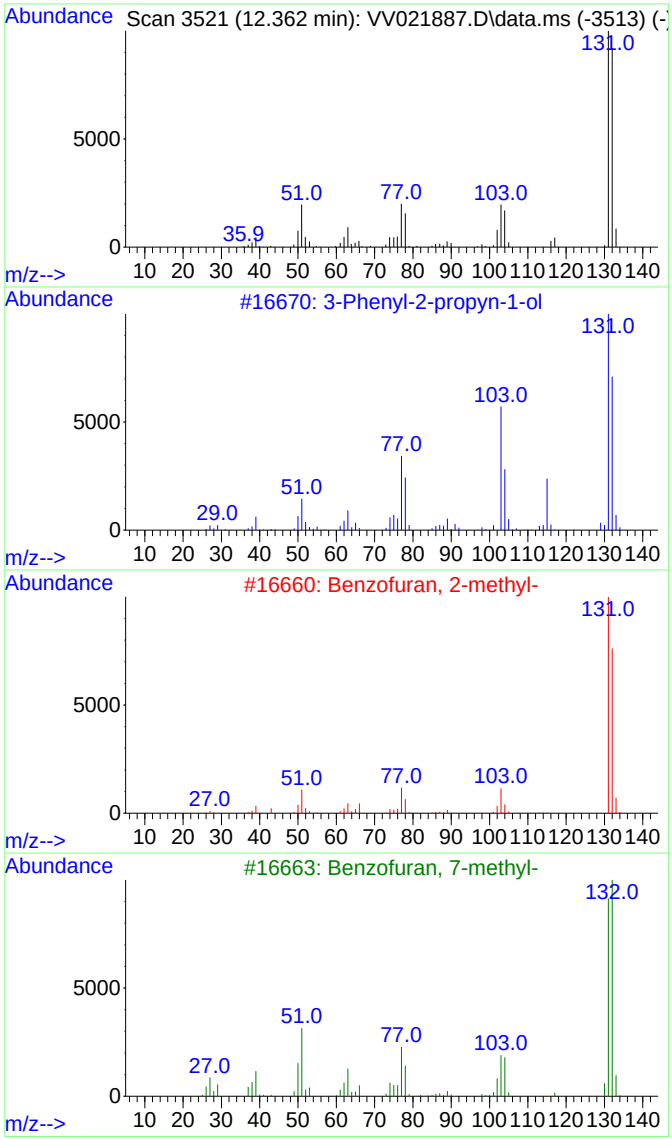
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

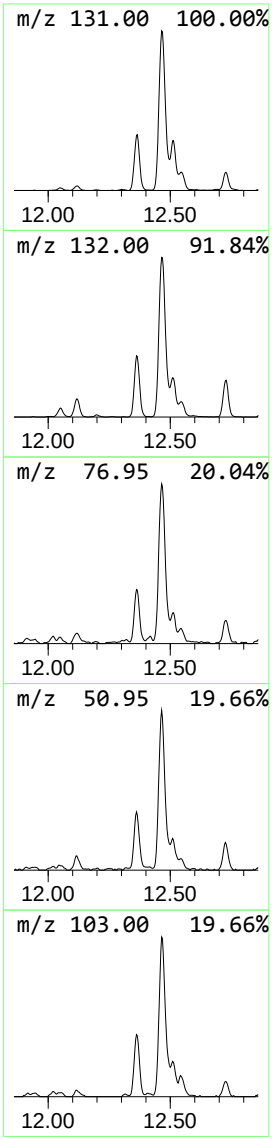
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 3-Phenyl-2-propyn-1-ol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.361	6.97 ug/L	185551	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
4	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

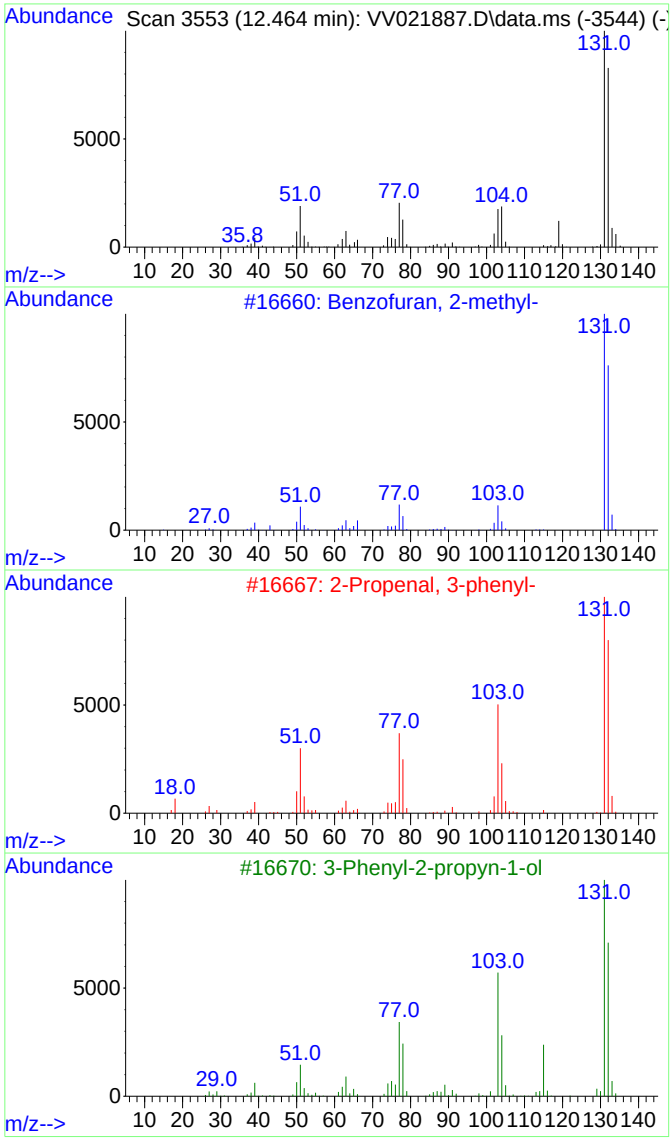
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

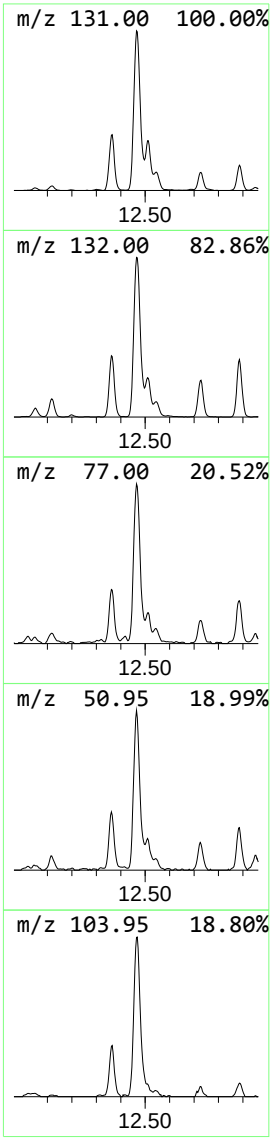
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzofuran, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	23.64 ug/L	629763	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	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	89





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

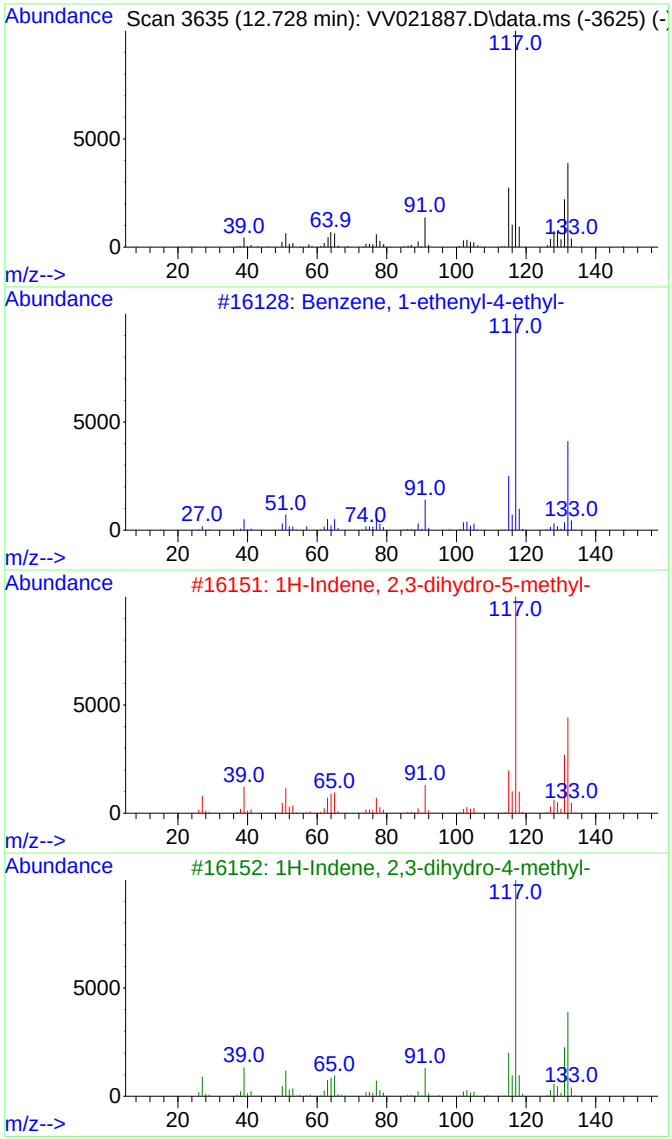
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

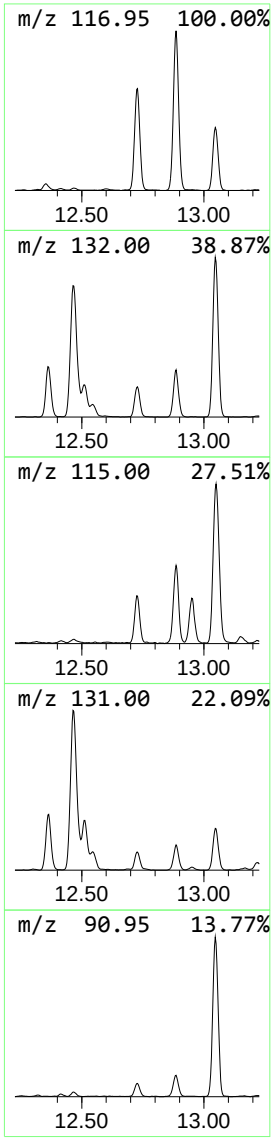
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.728	8.25 ug/L	219859	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	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

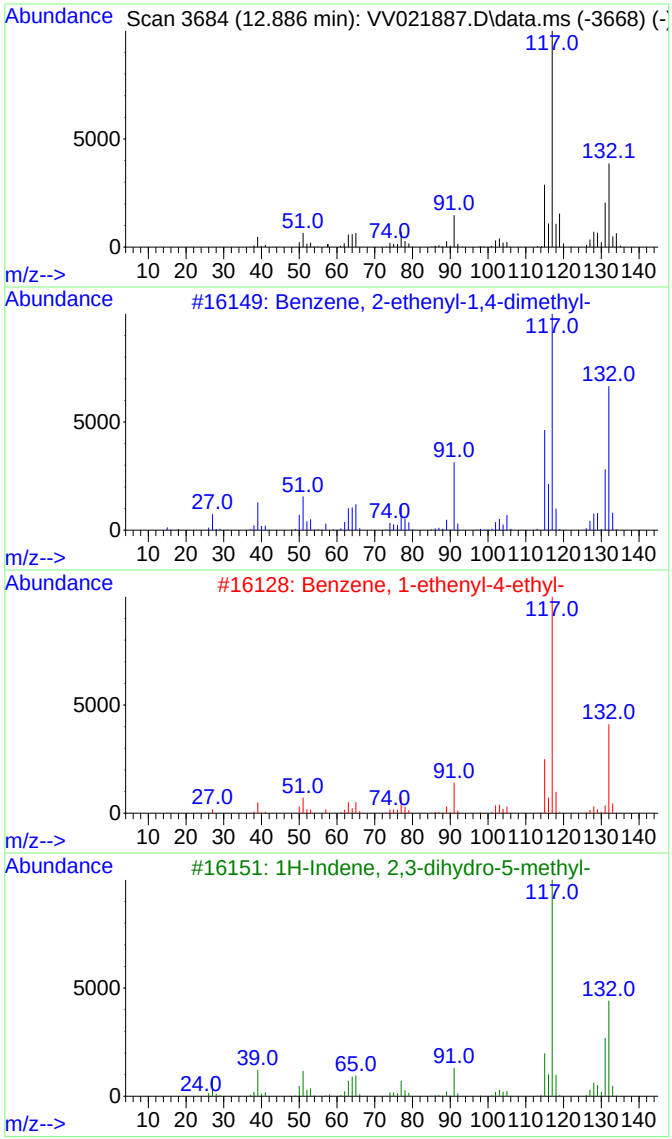
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

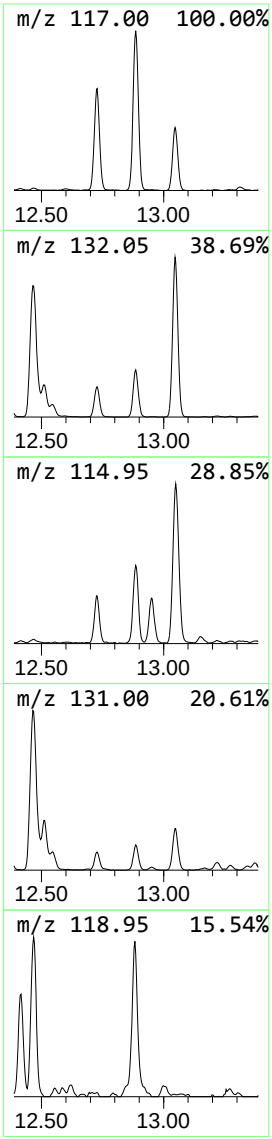
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	14.76 ug/L	393171	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	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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

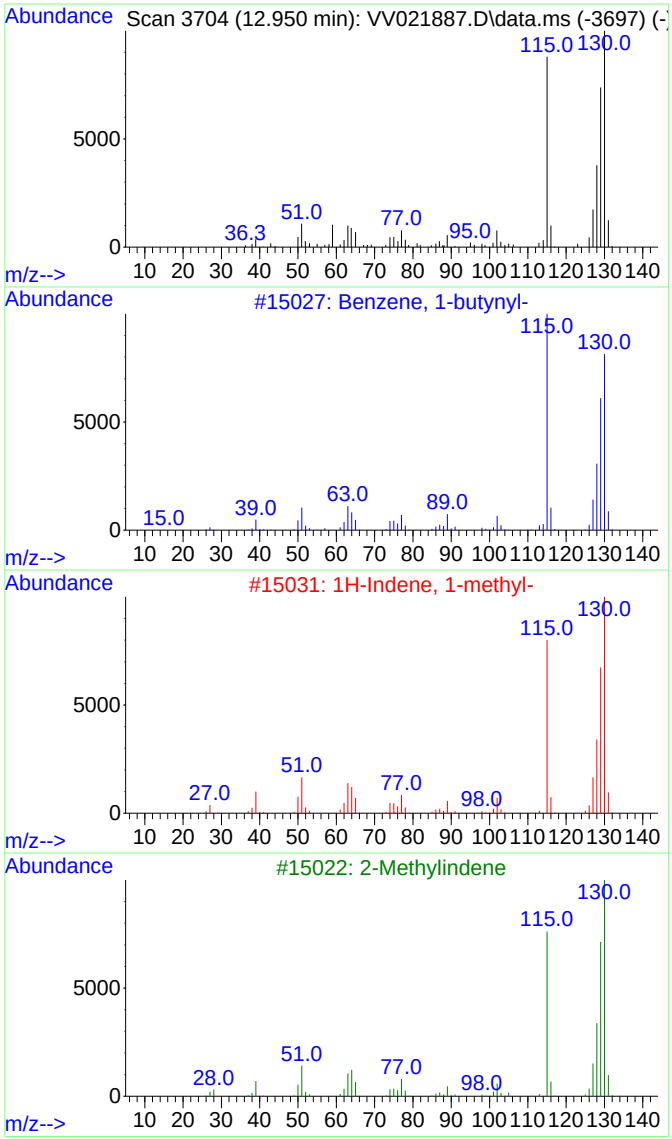
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

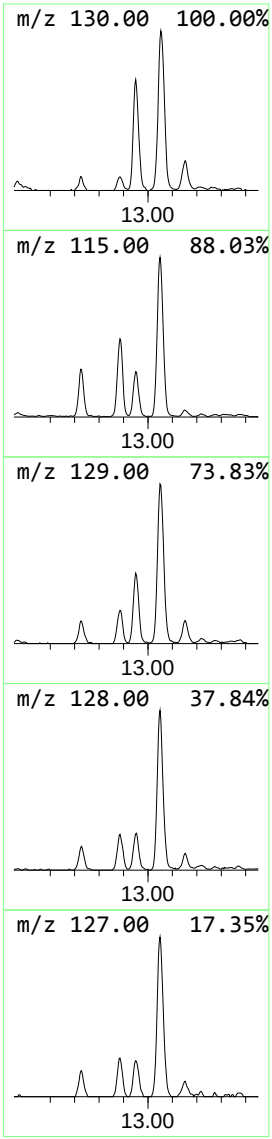
****

Peak Number 14 Benzene, 1-butynyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.950	3.86 ug/L	102820	1,4-Dichlorobenzene-d4	11.252

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-butynyl-	130	C10H10	000622-76-4	96
2	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
3	2-Methylindene	130	C10H10	002177-47-1	94
4	Benzene, 1-methyl-4-(1-propynyl)-	130	C10H10	002749-93-1	94
5	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

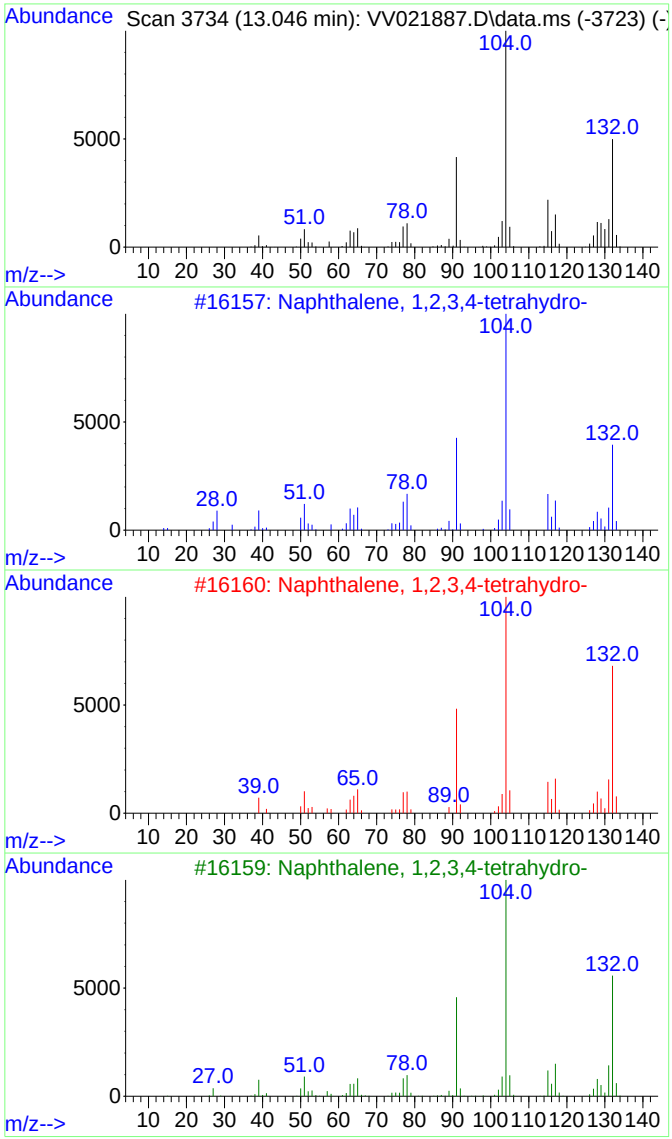
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

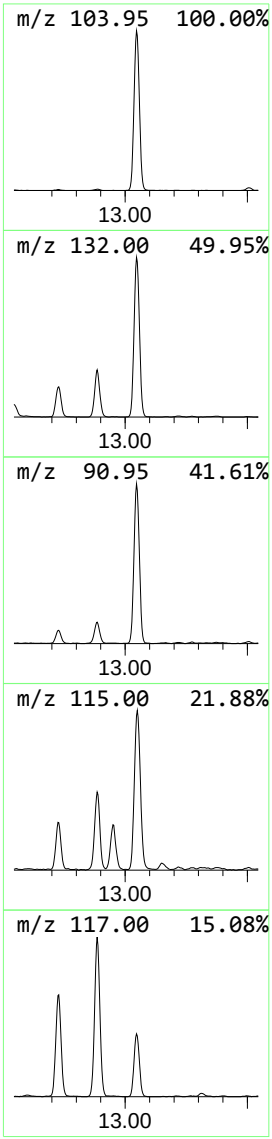
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.046	46.46 ug/L	1237550	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	96
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	94





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Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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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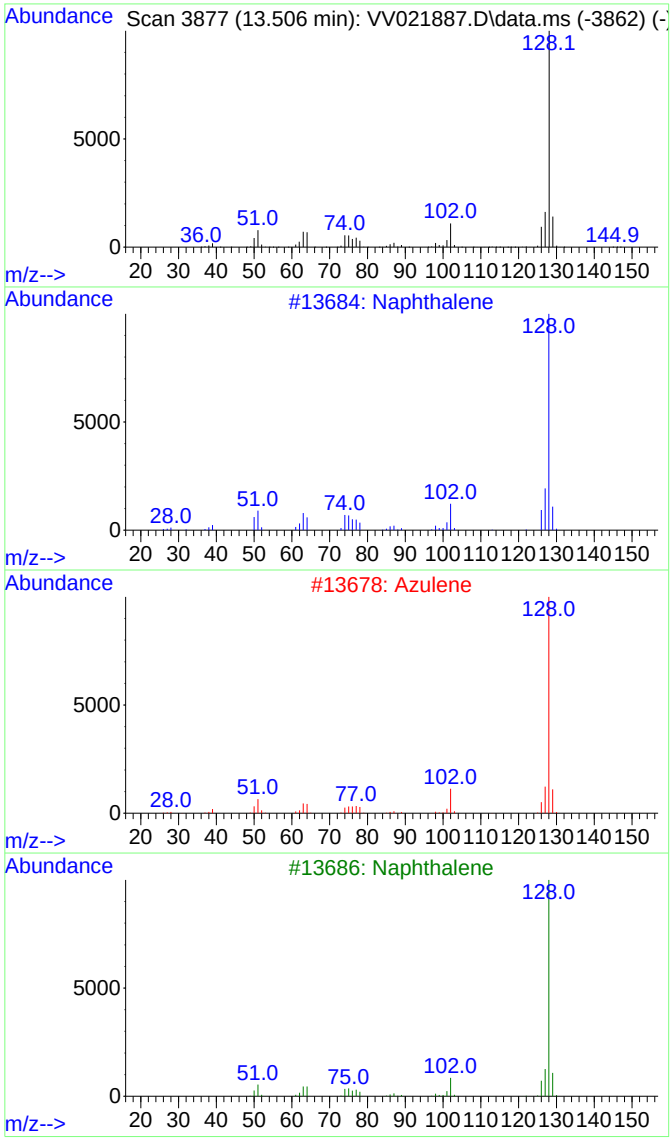
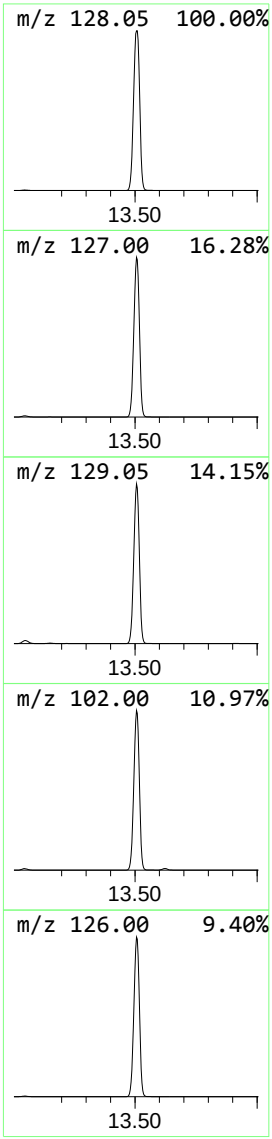
Peak Number 16 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.506	1159.49 ug/L	30883900	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	Azulene	128	C10H8	000275-51-4	91
5	Naphthalene	128	C10H8	000091-20-3	91





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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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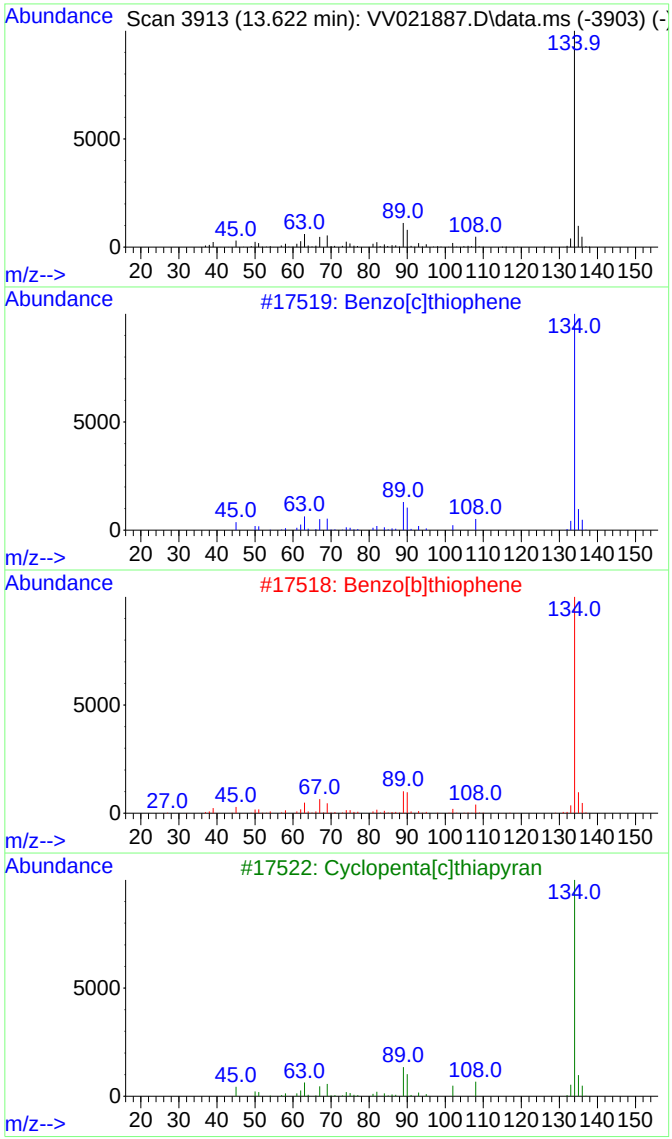
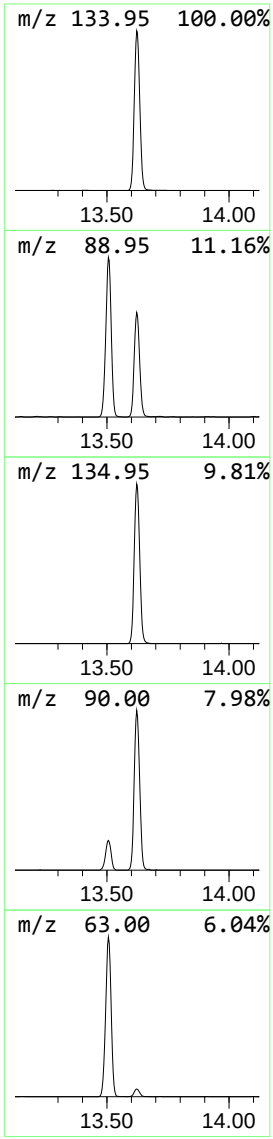
Peak Number 17 Benzo[c]thiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	57.01 ug/L	1518490	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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

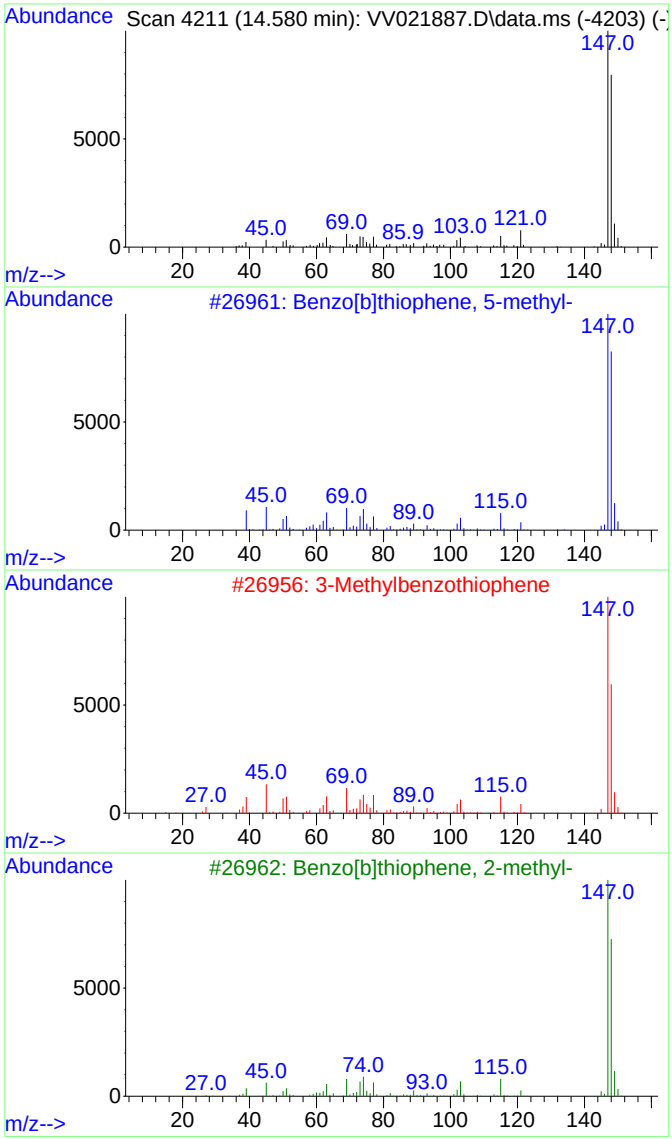
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

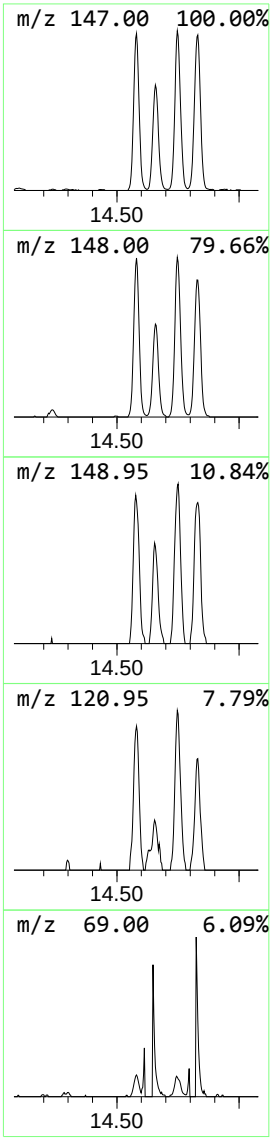
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzo[b]thiophene, 5-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.580	4.17 ug/L	111026	1,4-Dichlorobenzene-d4	11.252

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

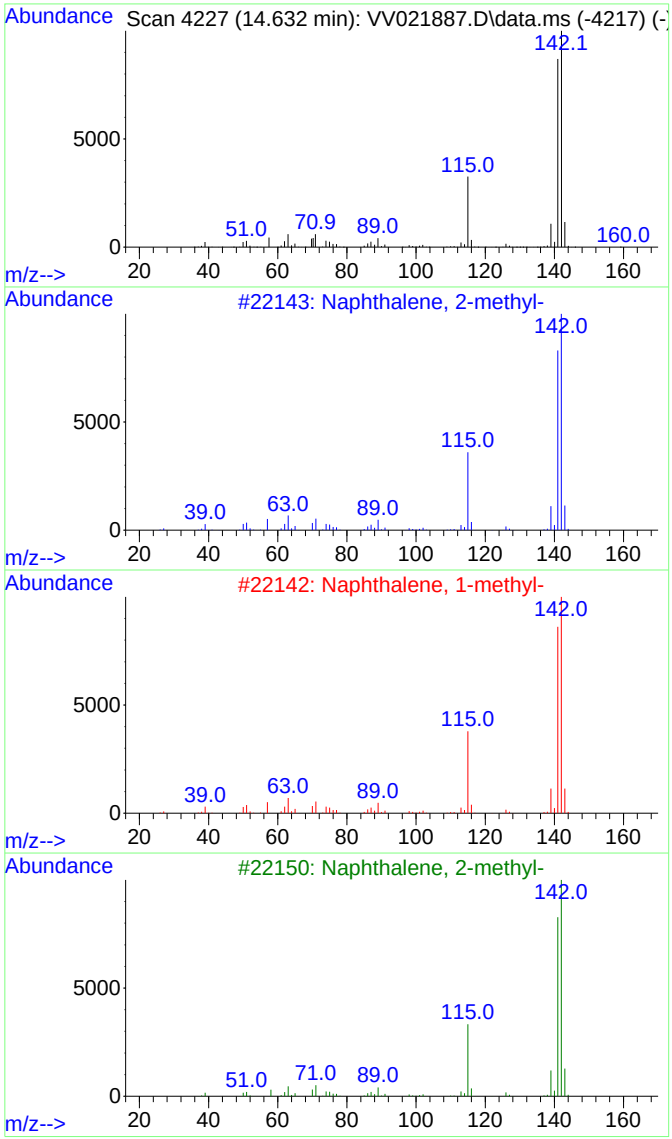
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

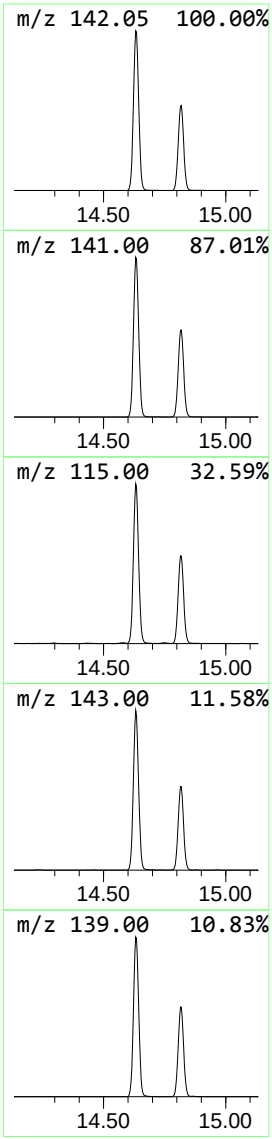
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.632	160.79 ug/L	4282890	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, 1-methyl-	142	C11H10	000090-12-0	94
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

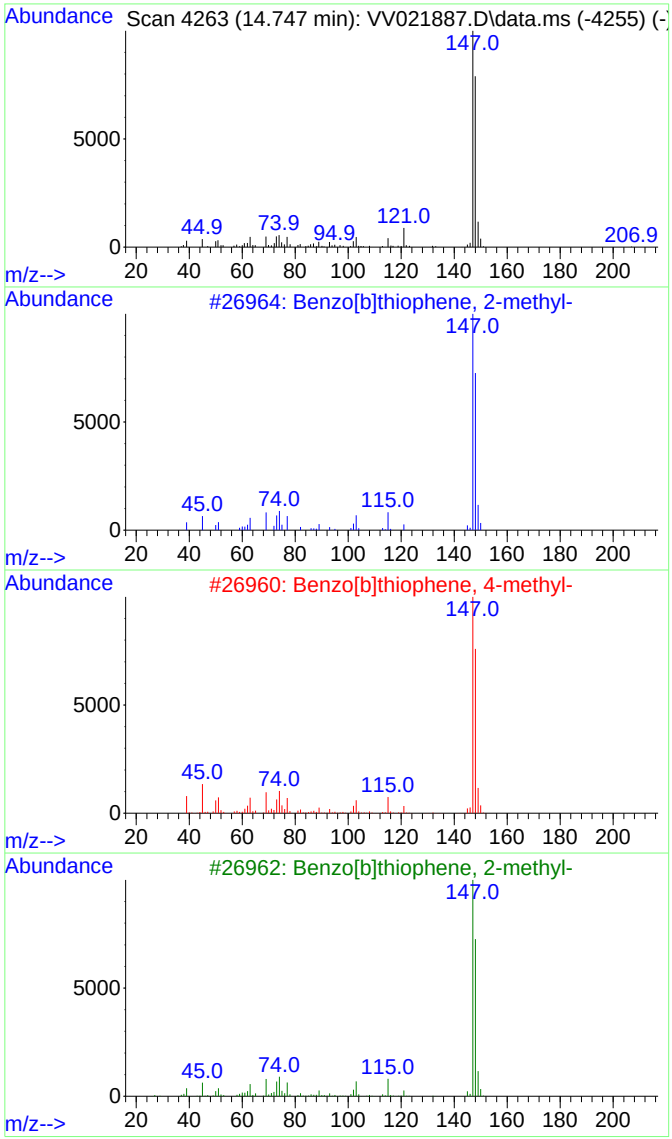
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

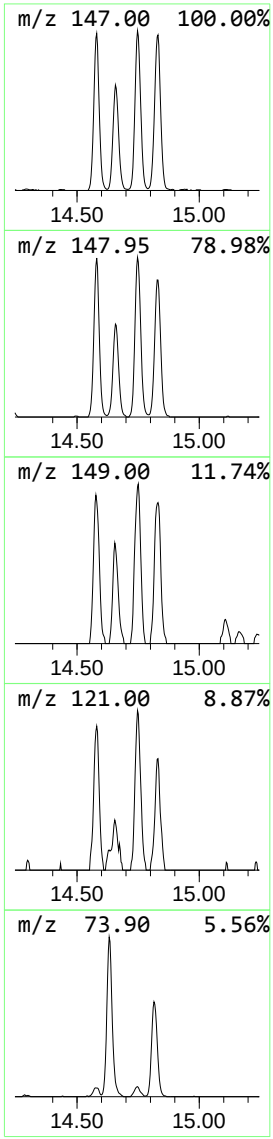
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzo[b]thiophene, 2-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.747	4.68 ug/L	124565	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	3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
5	3-Methylbenzothiophene	148	C9H8S	001455-18-1	86





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Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

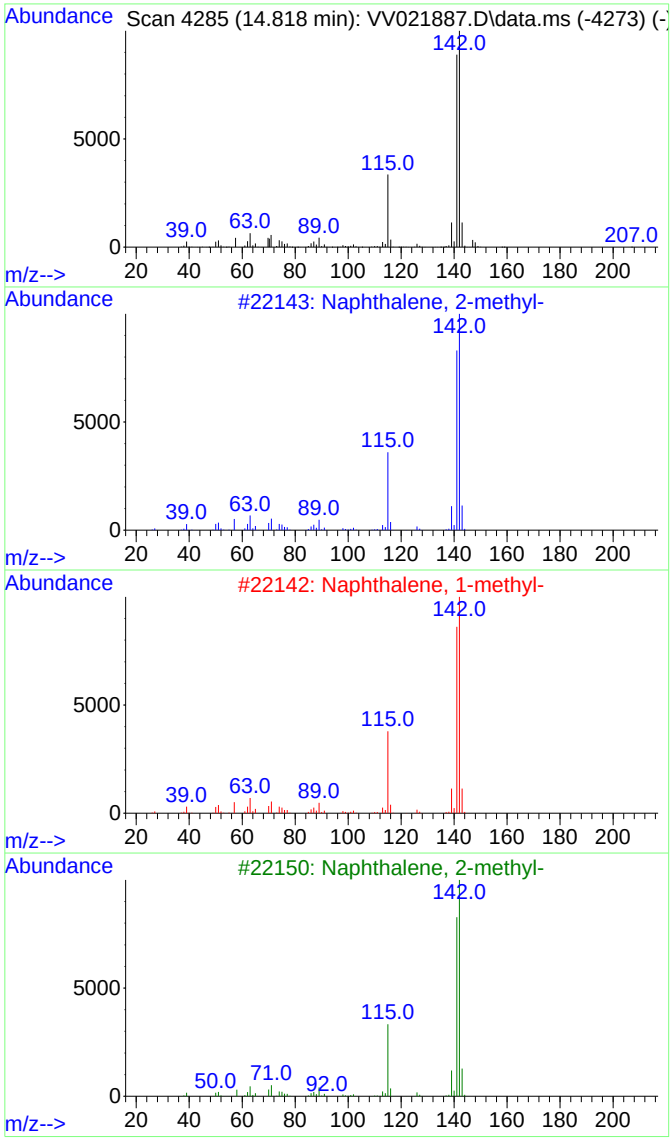
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

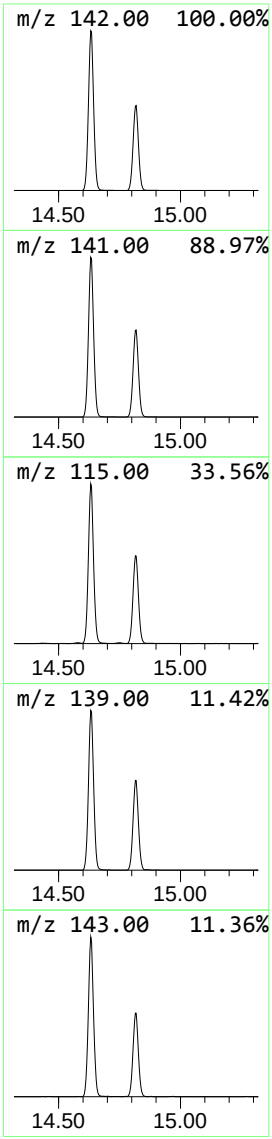
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.818	93.67 ug/L	2494870	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, 1-methyl-	142	C11H10	000090-12-0	95
5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

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

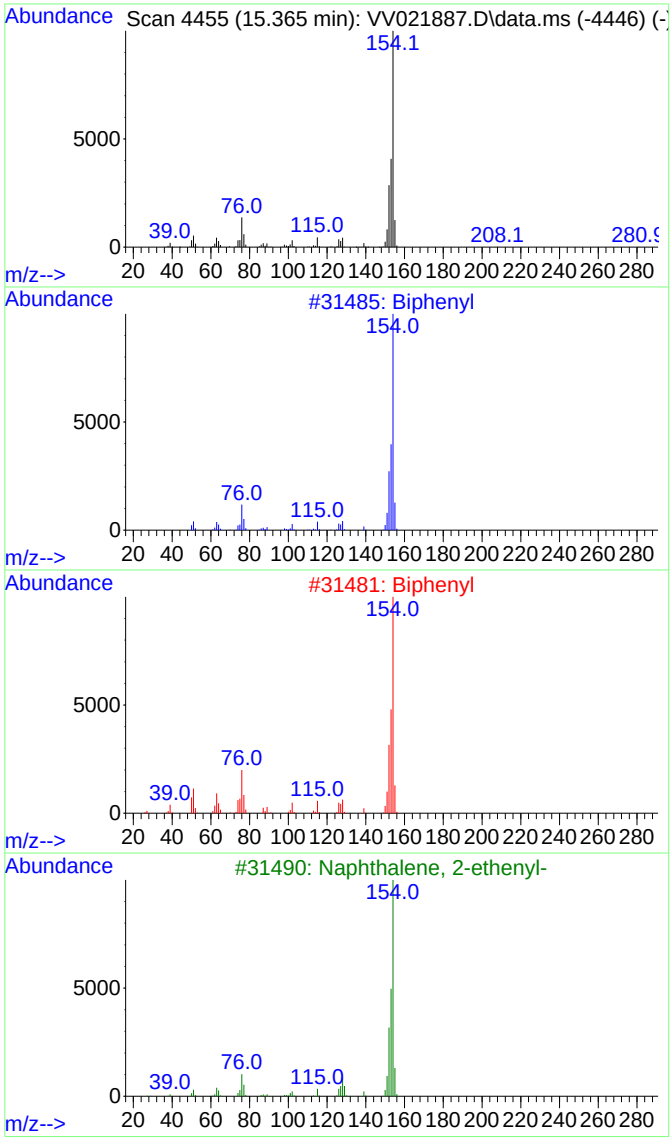
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

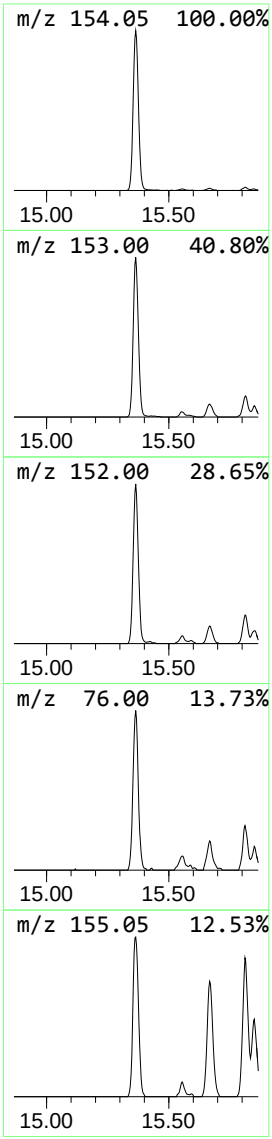
****

Peak Number 22 Biphenyl Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.365	14.08 ug/L	375051	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	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	91
4	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	83
5	Biphenyl	154	C12H10	000092-52-4	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

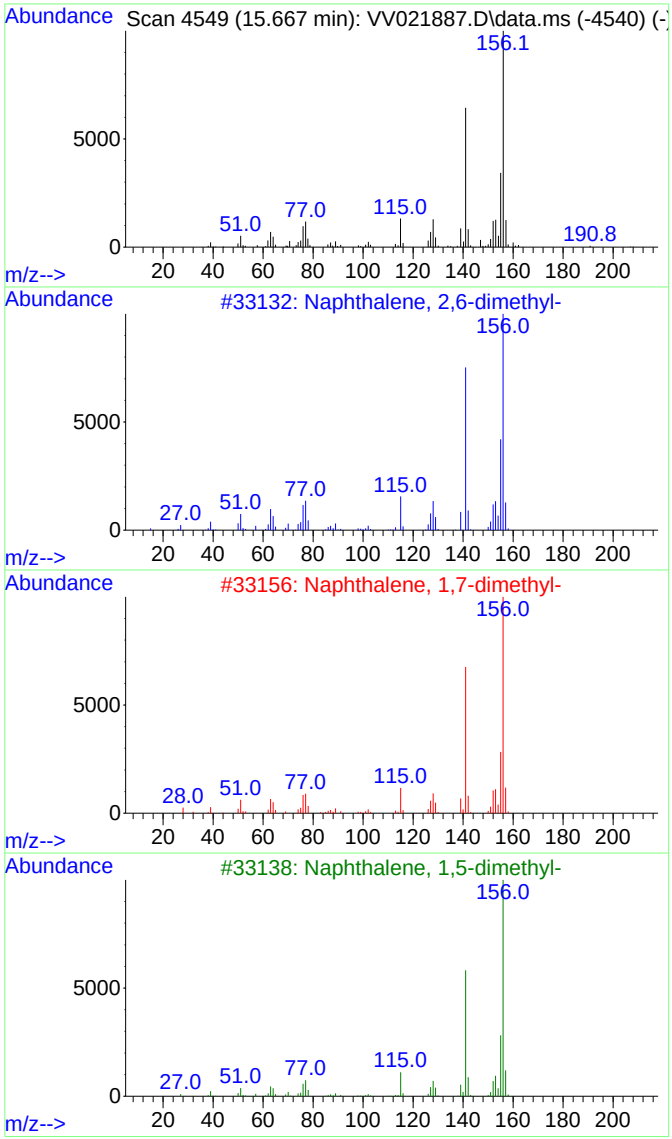
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

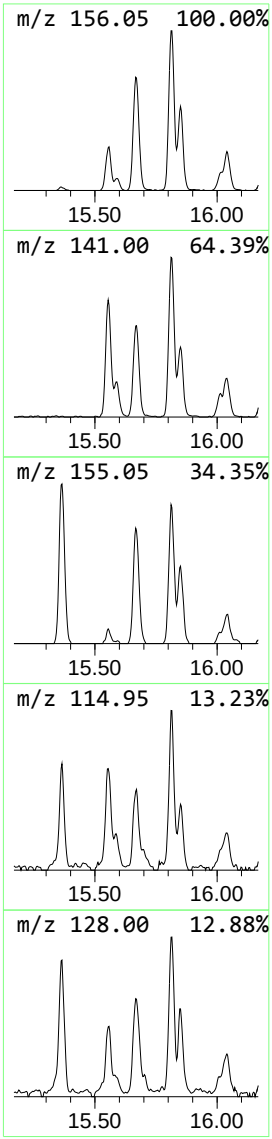
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Naphthalene, 2,6-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.667	6.88 ug/L	183296	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, 1,7-dimethyl-	156	C12H12	000575-37-1	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, 2,7-dimethyl-	156	C12H12	000582-16-1	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

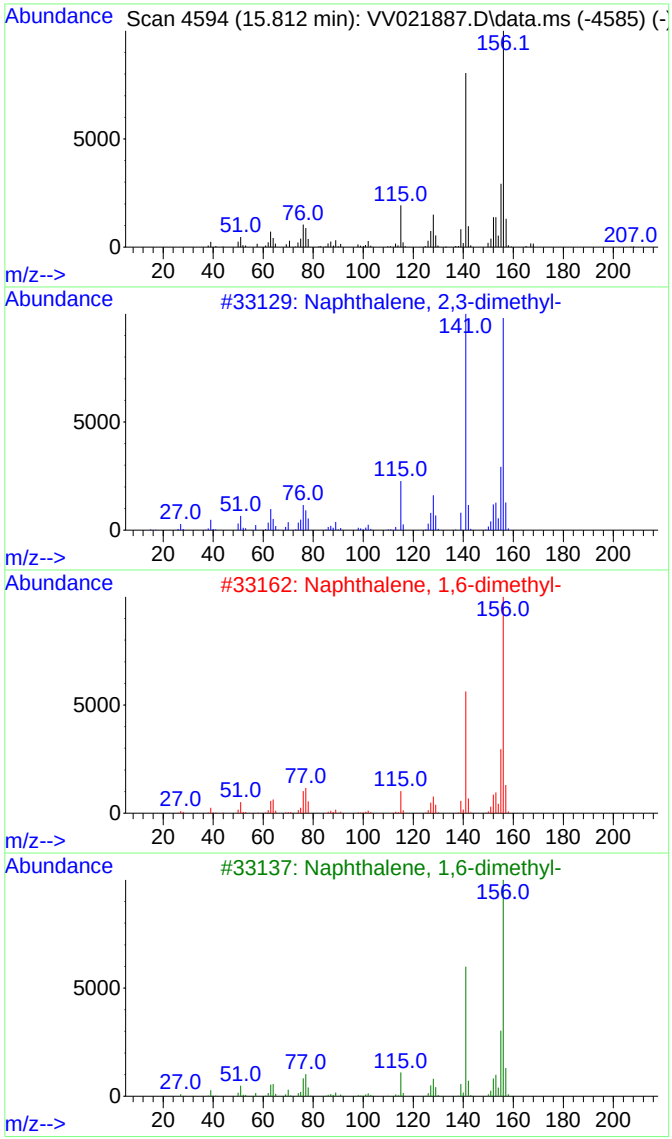
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

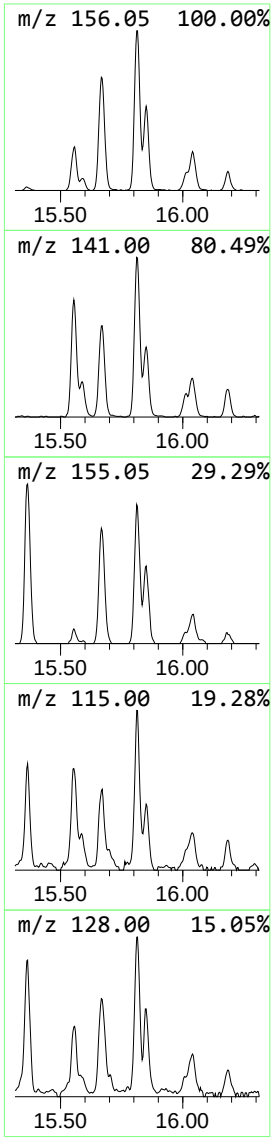
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Naphthalene, 2,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.812	8.01 ug/L	213362	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, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

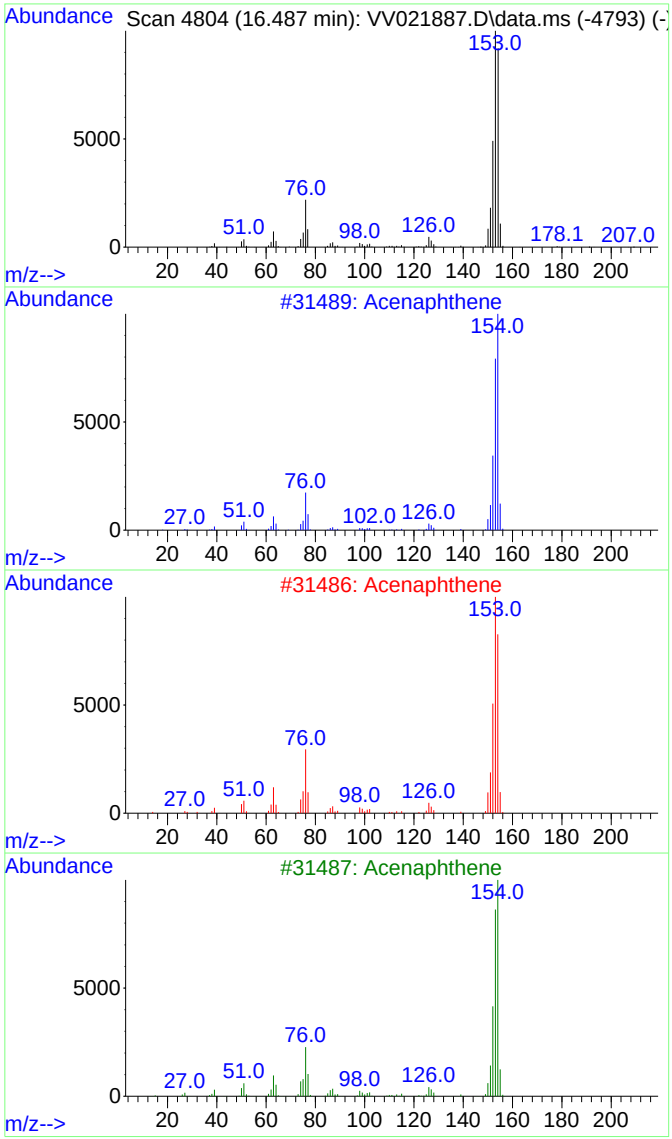
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

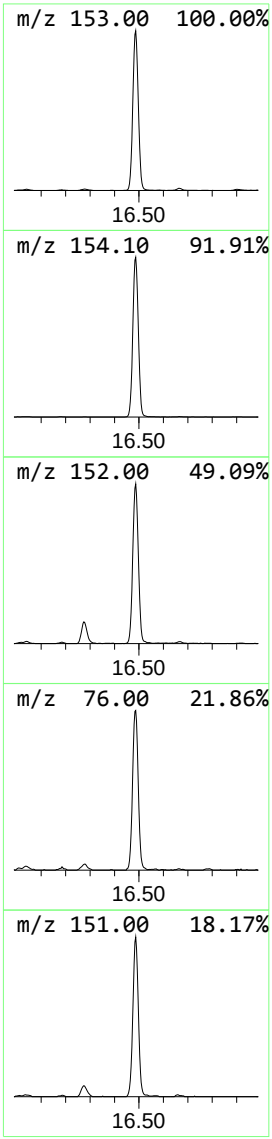
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Acenaphthene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.487	35.23 ug/L	938390	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	59
5	Acenaphthene	154	C12H10	000083-32-9	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKB0DL

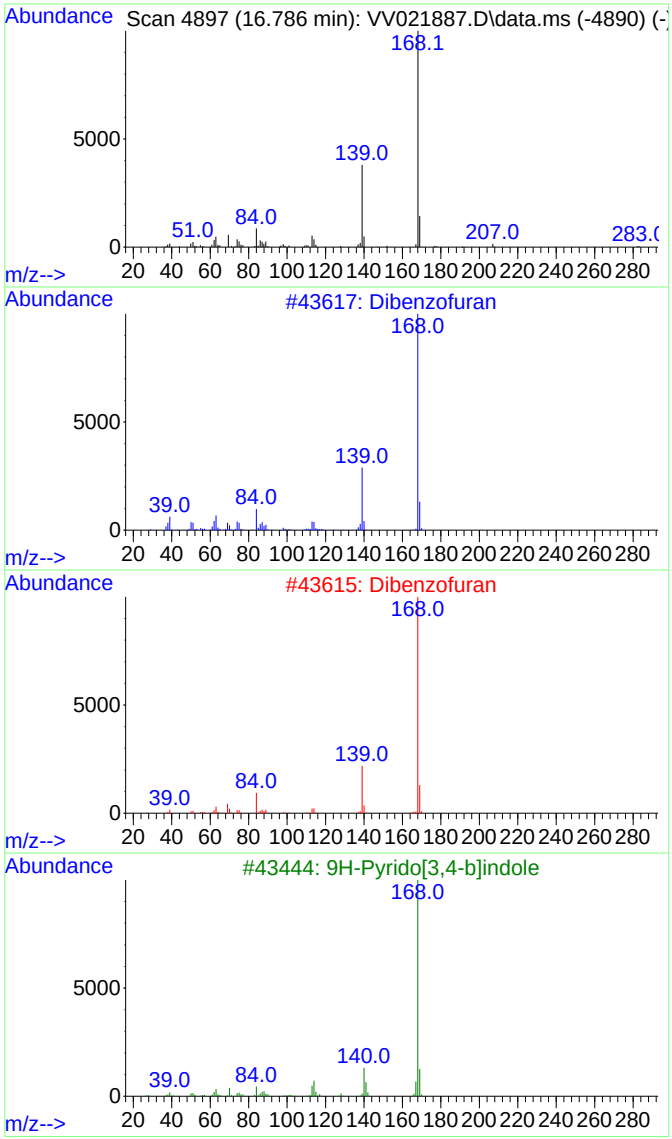
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

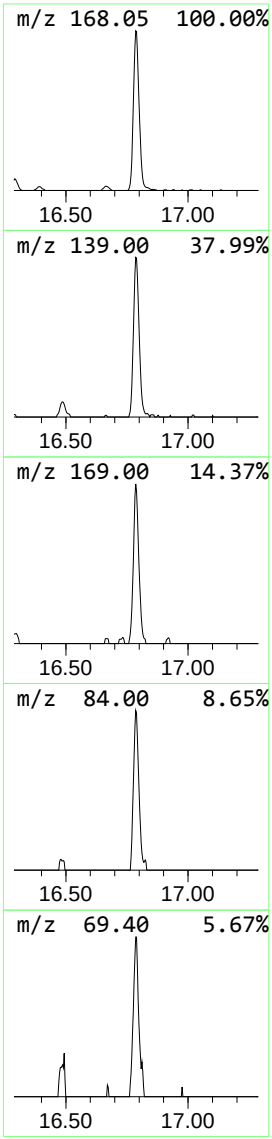
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Dibenzofuran Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.786	4.09 ug/L	109065	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	86
3	9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	59
4	9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	59
5	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081921\
Data File : VV021887.D
Acq On : 19 Aug 2021 13:35
Operator : SY/MD
Sample : M3422-09DL 5X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : DBKB0DL

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
unknown-01	1.201	3.0	ug/L	60208	1	5.619	1009810	50.0
Benzene, 1-ethy...	10.452	6.2	ug/L	165804	3	11.252	1331790	50.0
Benzene, 1-prop...	10.989	3.2	ug/L	84819	3	11.252	1331790	50.0
Benzofuran	11.172	35.5	ug/L	946671	3	11.252	1331790	50.0
Benzene, 1,2,3-...	11.339	6.2	ug/L	165680	3	11.252	1331790	50.0
Indane	11.529	51.1	ug/L	1362540	3	11.252	1331790	50.0
Indene	11.747	412.9	ug/L	10997500	3	11.252	1331790	50.0
Benzene, 1-ethe...	12.117	5.1	ug/L	135894	3	11.252	1331790	50.0
3-Phenyl-2-prop...	12.361	7.0	ug/L	185551	3	11.252	1331790	50.0
Benzofuran, 2-m...	12.464	23.6	ug/L	629763	3	11.252	1331790	50.0
Benzene, 1-ethe...	12.728	8.3	ug/L	219859	3	11.252	1331790	50.0
Benzene, 2-ethe...	12.886	14.8	ug/L	393171	3	11.252	1331790	50.0
Benzene, 1-buty...	12.950	3.9	ug/L	102820	3	11.252	1331790	50.0
Naphthalene, 1,...	13.046	46.5	ug/L	1237550	3	11.252	1331790	50.0
Azulene	13.506	1159.5	ug/L	30883900	3	11.252	1331790	50.0
Benzo[c]thiophene	13.622	57.0	ug/L	1518490	3	11.252	1331790	50.0
Benzo[b]thiophe...	14.580	4.2	ug/L	111026	3	11.252	1331790	50.0
Naphthalene, 2-...	14.632	160.8	ug/L	4282890	3	11.252	1331790	50.0
Benzo[b]thiophe...	14.747	4.7	ug/L	124565	3	11.252	1331790	50.0
Naphthalene, 1-...	14.818	93.7	ug/L	2494870	3	11.252	1331790	50.0
Biphenyl	15.365	14.1	ug/L	375051	3	11.252	1331790	50.0
Naphthalene, 2,...	15.667	6.9	ug/L	183296	3	11.252	1331790	50.0
Naphthalene, 2,...	15.812	8.0	ug/L	213362	3	11.252	1331790	50.0
Acenaphthene	16.487	35.2	ug/L	938390	3	11.252	1331790	50.0
Dibenzofuran	16.786	4.1	ug/L	109065	3	11.252	1331790	50.0