LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : C0AY3

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_X\Method\SFAMXLM072624WMA.M
Title : VOC Analysis

Signal : TIC: VX042764.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.239	22	25	39	rBV	18562	24910	5.39%	0.556%
2	1.368	43	46	56	rVB	31847	39359	8.52%	0.878%
3	1.648	89	92	98	rVB	25488	34774	7.52%	0.776%
4	2.294	192	198	210	rBV	76233	120684	26.11%	2.692%
5	2.386	210	213	220	rVB2	872	1383	0.30%	0.031%
6	2.508	229	233	237	rBB2	225	441	0.10%	0.010%
7	2.538	237	238	241	rBV	140	154	0.03%	0.003%
8	2.575	241	244	248	rVB	334	454	0.10%	0.010%
9	2.794	275	280	283	rBB2	608	945	0.20%	0.021%
10	2.843	286	288	290	rBV	147	160	0.03%	0.004%
11	2.892	293	296	298	rBB	91	113	0.02%	0.003%
12	2.940	298	304	311	rBB3	1211	2553	0.55%	0.057%
13	3.824	446	449	452	rBB	207	179	0.04%	0.004%
14	4.245	517	518	522	rBV	156	185	0.04%	0.004%
15	4.471	547	555	574	rBV2	24315	71766	15.53%	1.601%
16	4.635	580	582	584	rBV2	167	145	0.03%	0.003%
17	4.684	588	590	591	rBV2	169	114	0.02%	0.003%
18	4.757	600	602	604	rBV2	205	246	0.05%	0.005%
19	4.806	609	610	613	rBB2	132	112	0.02%	0.002%
20	5.056	640	651	664	rBV	49686	151923	32.87%	3.388%
21	5.385	703	705	710	rBB	251	338	0.07%	0.008%
22	5.678	749	753	755	rBB	125	155	0.03%	0.003%
23	5.726	759	761	764	rVB	209	199	0.04%	0.004%
24	5.964	790	800	821	rBV2	131846	382353	82.72%	8.528%
25	6.348	861	863	865	rVB	232	179	0.04%	0.004%
26	6.763	921	931	958	rBV	104972	254249	55.01%	5.671%
27	6.970	963	965	967	rVB	235	193	0.04%	0.004%
28	7.062	978	980	983	rBB	133	130	0.03%	0.003%
29	7.116	987	989	992	rBV2	406	324	0.07%	0.007%
30	7.196	1000	1002	1005	rBV2	178	240	0.05%	0.005%
31	7.245	1008	1010	1012	rBB2	252	137	0.03%	0.003%
32	7.305	1012	1020	1037	rBV	77514	169565	36.69%	3.782%
33	7.494	1048	1051	1055	rBB2	174	303	0.07%	0.007%
34	7.598	1066	1068	1070	rVB	195	136	0.03%	0.003%
35	7.641	1071	1075	1076	rBV	153	175	0.04%	0.004%
36	7.879	1110	1114	1115	rBB	178	164	0.04%	0.004%

LSC Area Percent Report

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Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : C0AY3

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_X\Method\SFAMXLM072624WMA.M
Title : VOC Analysis

37	8.007	1132	1135	1137	rBB	142	159	0.03%	0.004%
38	8.098	1147	1150	1153	rBB	191	180	0.04%	0.004%
39	8.324	1182	1187	1196	rBV	49122	83810	18.13%	1.869%
40	8.488	1211	1214	1216	rBB	104	116	0.03%	0.003%
41	8.537	1220	1222	1225	rBB	112	121	0.03%	0.003%
42	8.574	1225	1228	1230	rBB	184	148	0.03%	0.003%
43	8.647	1234	1240	1250	rBV	200178	319303	69.08%	7.122%
44	8.915	1282	1284	1285	rBV	172	142	0.03%	0.003%
45	8.952	1285	1290	1306	rVB	36863	56130	12.14%	1.252%
46	9.067	1306	1309	1311	rBB	236	224	0.05%	0.005%
47	9.128	1316	1319	1322	rBB2	201	244	0.05%	0.005%
48	9.275	1339	1343	1349	rVB2	931	1441	0.31%	0.032%
49	9.384	1356	1361	1373	rBV	111165	173584	37.56%	3.872%
50	9.665	1405	1407	1408	rBV	213	157	0.03%	0.004%
51	9.714	1411	1415	1417	rBV	154	254	0.05%	0.006%
52	9.756	1421	1422	1425	rBB	142	133	0.03%	0.003%
53	9.988	1457	1460	1461	rBB	135	113	0.02%	0.003%
54	10.055	1465	1471	1488	rBV	223064	308565	66.76%	6.882%
55	10.201	1491	1495	1501	rVB	2025	2915	0.63%	0.065%
56	10.305	1508	1512	1517	rBV2	1590	2255	0.49%	0.050%
57	10.646	1564	1568	1572	rVV2	821	1120	0.24%	0.025%
58	10.732	1581	1582	1585	rBV2	236	230	0.05%	0.005%
59	10.896	1608	1609	1611	rBB2	257	155	0.03%	0.003%
60	10.927	1612	1614	1616	rBV	284	333	0.07%	0.007%
61	11.006	1626	1627	1628	rBV	232	132	0.03%	0.003%
62	11.195	1652	1658	1669	rVB	152779	208186	45.04%	4.643%
63	11.311	1675	1677	1685	rVB3	1143	1490	0.32%	0.033%
64	11.384	1686	1689	1690	rBV	362	279	0.06%	0.006%
65	11.421	1693	1695	1698	rVB	134	115	0.02%	0.003%
66	11.451	1698	1700	1704	rBB	423	580	0.13%	0.013%
67	11.512	1708	1710	1712	rBB	189	145	0.03%	0.003%
68	11.603	1722	1725	1729	rBB	243	245	0.05%	0.005%
69	11.646	1731	1732	1735	rBV	192	134	0.03%	0.003%
70	11.695	1738	1740	1741	rBV	302	222	0.05%	0.005%
71	11.756	1745	1750	1753	rBB	1212	1340	0.29%	0.030%
72	11.963	1780	1784	1788	rBB3	424	660	0.14%	0.015%
73	12.024	1788	1794	1802	rBV	224148	293678	63.54%	6.550%
74	12.170	1816	1818	1819	rBV	349	261	0.06%	0.006%
75	12.244	1822	1830	1837	rVV2	13155	22158	4.79%	0.494%
76	12.317	1837	1842	1853	rVB	230802	301894	65.32%	6.733%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : C0AY3

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_X\Method\SFAMXLM072624WMA.M
Title : VOC Analysis

77	12.500	1870	1872	1874	rVB	183	115	0.02%	0.003%
78	12.542	1874	1879	1880	rBV	443	563	0.12%	0.013%
79	12.670	1898	1900	1901	rVB	228	115	0.02%	0.003%
80	12.811	1901	1923	1943	rBV3	87626	462203	100.00%	10.309%
81	13.774	2076	2081	2092	rVB	235499	301408	65.21%	6.723%
82	13.871	2093	2097	2107	rVB	10088	15243	3.30%	0.340%
83	13.999	2114	2118	2121	rBV3	1026	1375	0.30%	0.031%
84	14.079	2129	2131	2132	rBV	513	297	0.06%	0.007%
85	14.243	2156	2158	2159	rBV	229	201	0.04%	0.004%
86	14.323	2168	2171	2173	rBV	1045	1011	0.22%	0.023%
87	14.414	2184	2186	2187	rBV	191	164	0.04%	0.004%
88	14.463	2191	2194	2196	rBV	173	155	0.03%	0.003%
89	14.487	2196	2198	2200	rBV	377	226	0.05%	0.005%
90	14.597	2211	2216	2219	rBV3	1905	3609	0.78%	0.080%
91	14.640	2219	2223	2233	rVB	27212	38311	8.29%	0.854%
92	14.780	2241	2246	2254	rVB	59869	78842	17.06%	1.758%
93	15.054	2288	2291	2292	rBV	350	292	0.06%	0.007%
94	15.206	2312	2316	2322	rVB2	16045	24191	5.23%	0.540%
95	15.383	2339	2345	2348	rBV5	3427	6554	1.42%	0.146%
96	15.481	2355	2361	2370	rVB3	12303	21977	4.75%	0.490%
97	15.615	2378	2383	2386	rBV2	13800	20739	4.49%	0.463%
98	16.090	2455	2461	2469	rVB3	7514	13168	2.85%	0.294%
99	16.322	2492	2499	2509	rBV	147626	259184	56.08%	5.781%
100	16.688	2552	2559	2567	rVB	96340	190819	41.28%	4.256%

Sum of corrected areas: 4483553

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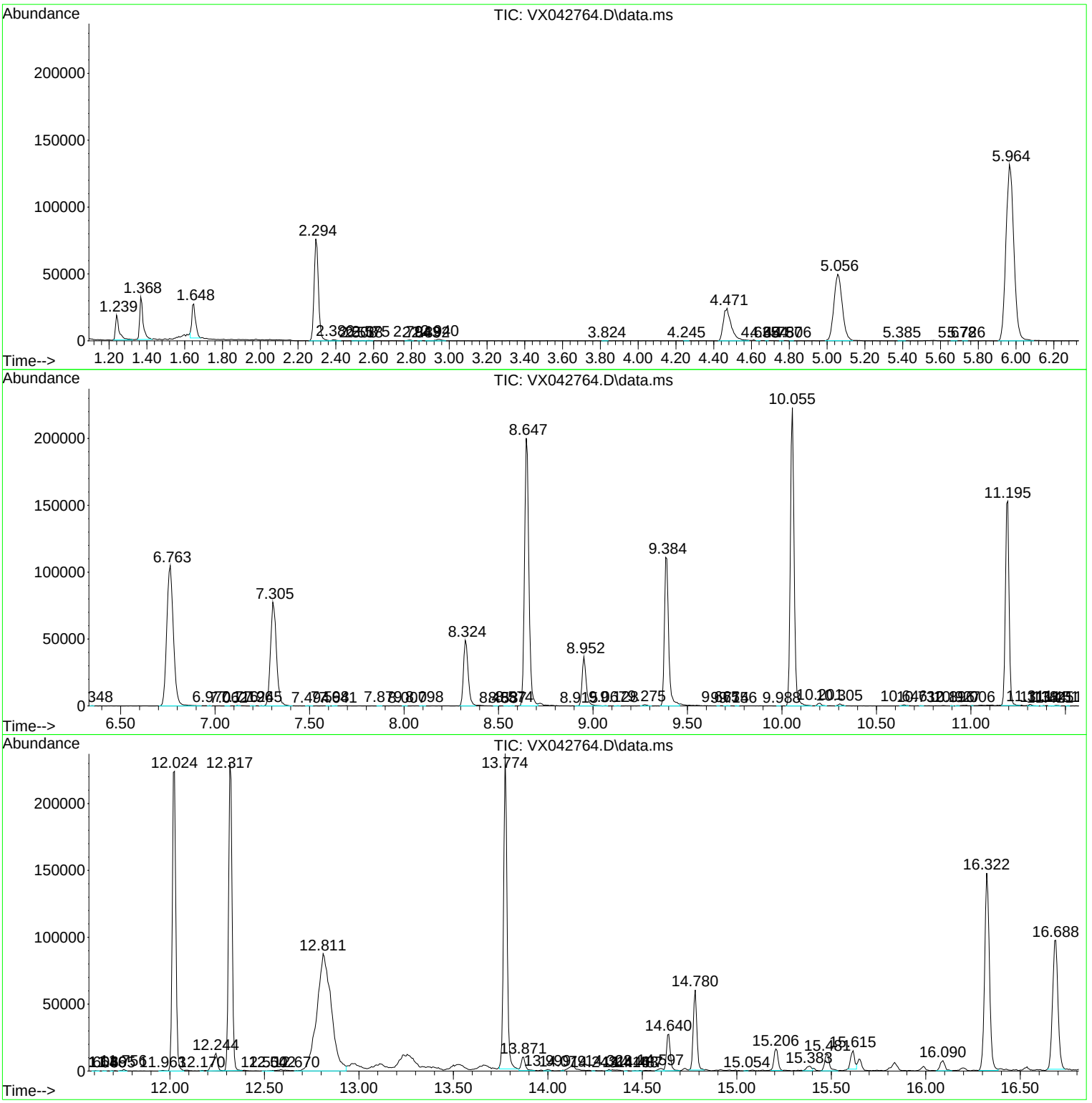
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
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Instrument :
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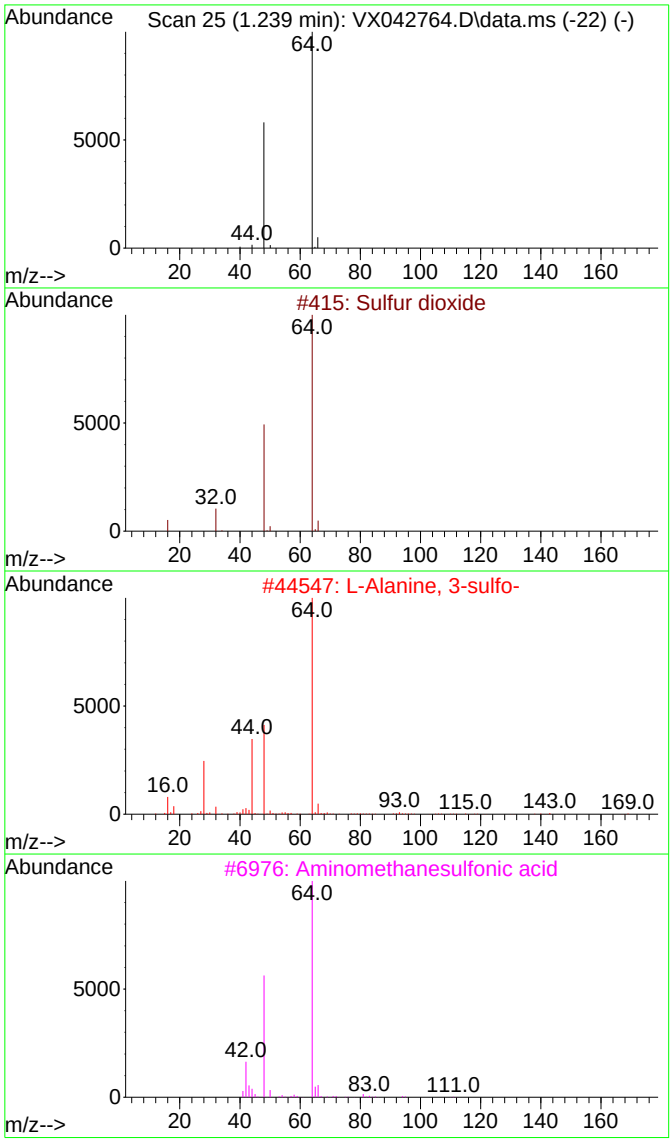
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

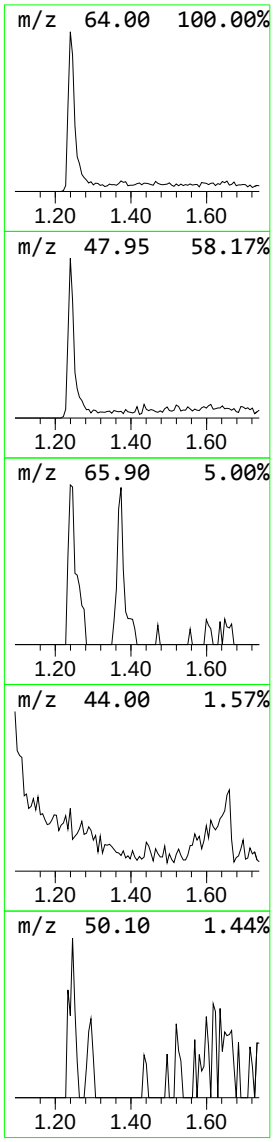
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Sulfur dioxide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.239	4.90 ug/L	24910	1,4-Difluorobenzene	6.763

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

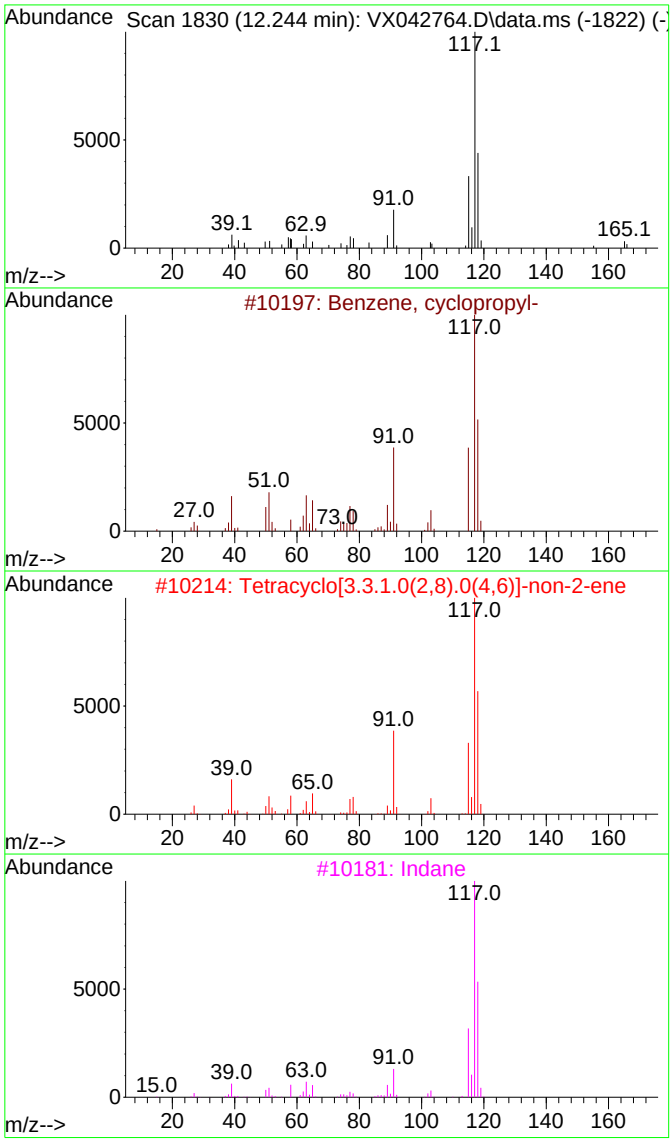
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

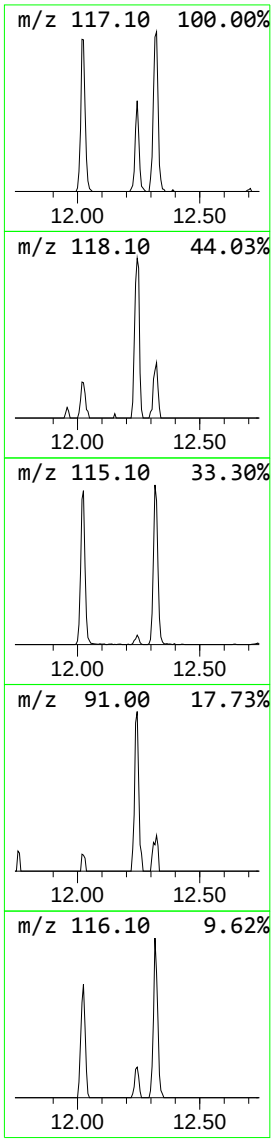
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, cyclopropyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	3.77 ug/L	22158	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
3	Indane	118	C9H10	000496-11-7	72
4	Deltacyclene	118	C9H10	007785-10-6	64
5	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

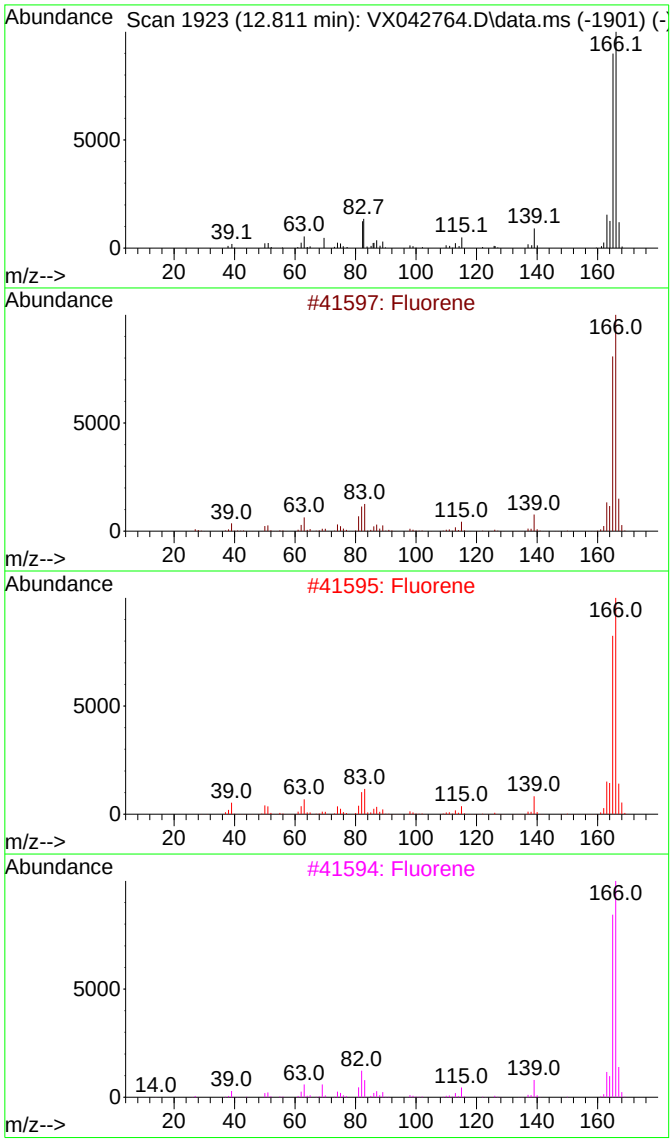
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

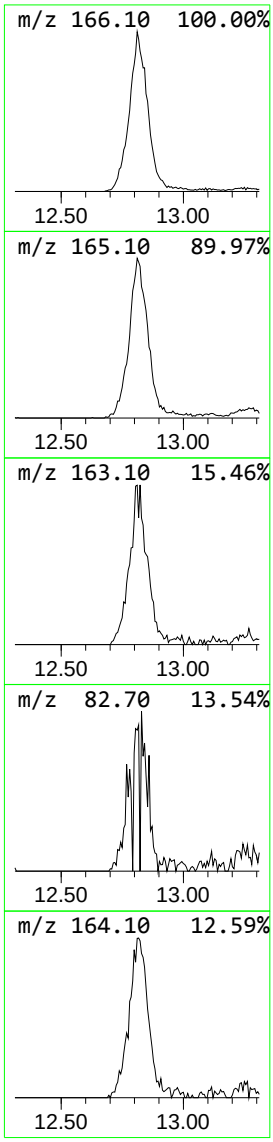
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1H-Phenalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.811	78.69 ug/L	462203	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluorene	166	C13H10	000086-73-7	96
2	Fluorene	166	C13H10	000086-73-7	94
3	Fluorene	166	C13H10	000086-73-7	91
4	Fluorene	166	C13H10	000086-73-7	90
5	1H-Phenalene	166	C13H10	000203-80-5	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

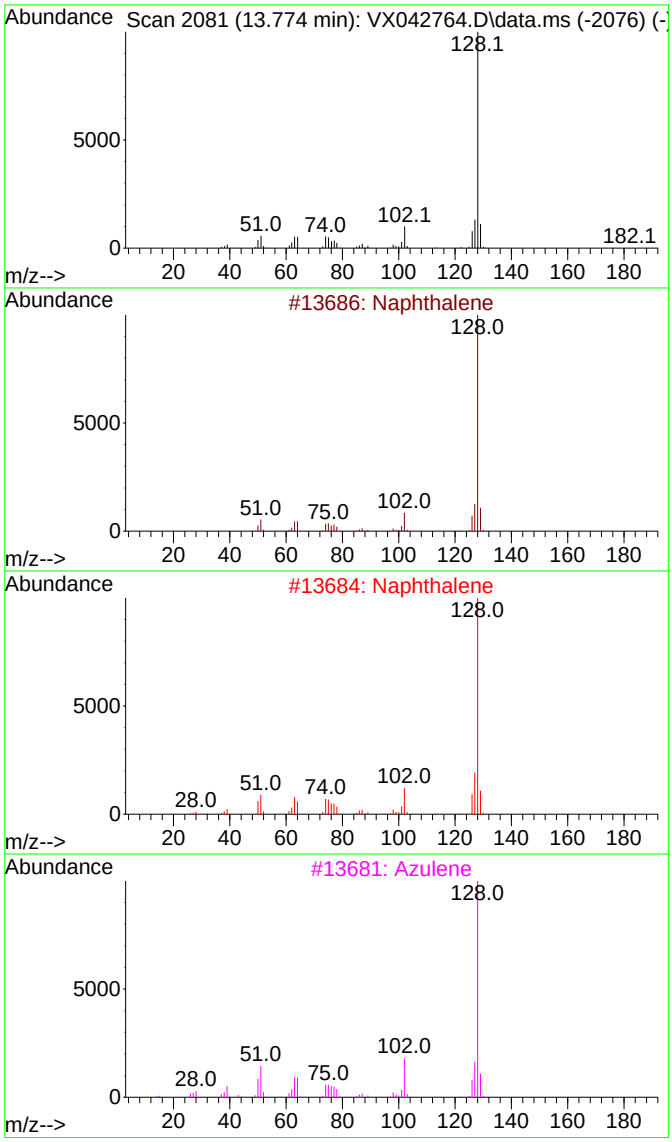
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

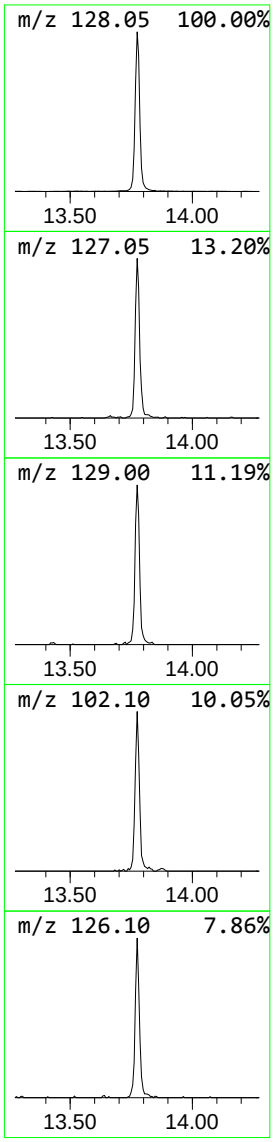
****

Peak Number 5 Azulene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.774	51.32 ug/L	301408	1,4-Dichlorobenzene-d4	12.024

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	91
4	Azulene	128	C10H8	000275-51-4	91
5	Azulene	128	C10H8	000275-51-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

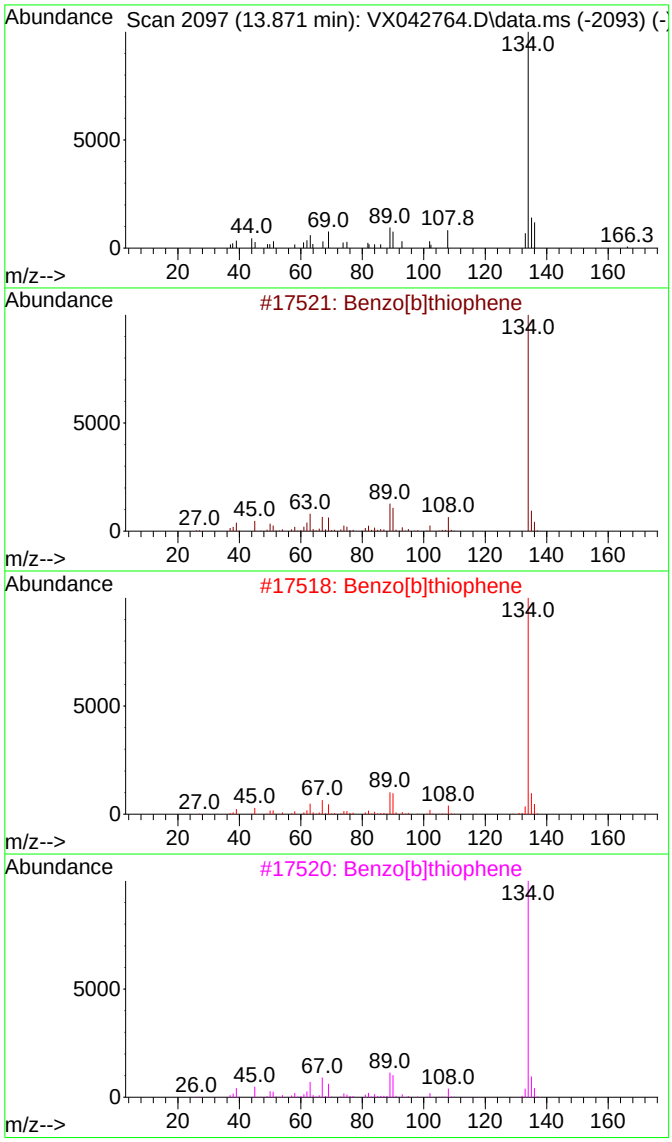
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

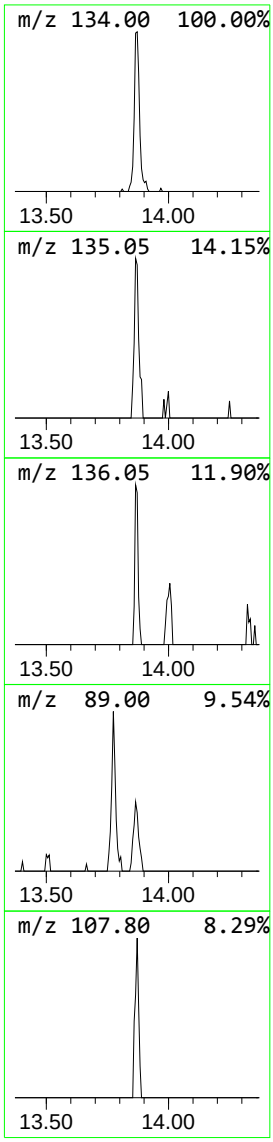
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzo[b]thiophene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.871	2.60 ug/L	15243	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene	134	C8H6S	000095-15-8	83
2	Benzo[b]thiophene	134	C8H6S	000095-15-8	80
3	Benzo[b]thiophene	134	C8H6S	000095-15-8	78
4	[1]Benzothieno[2',3':3,4]cyclobu...	246	C13H10O3S	034002-19-2	72
5	Benzo[c]thiophene	134	C8H6S	000270-82-6	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

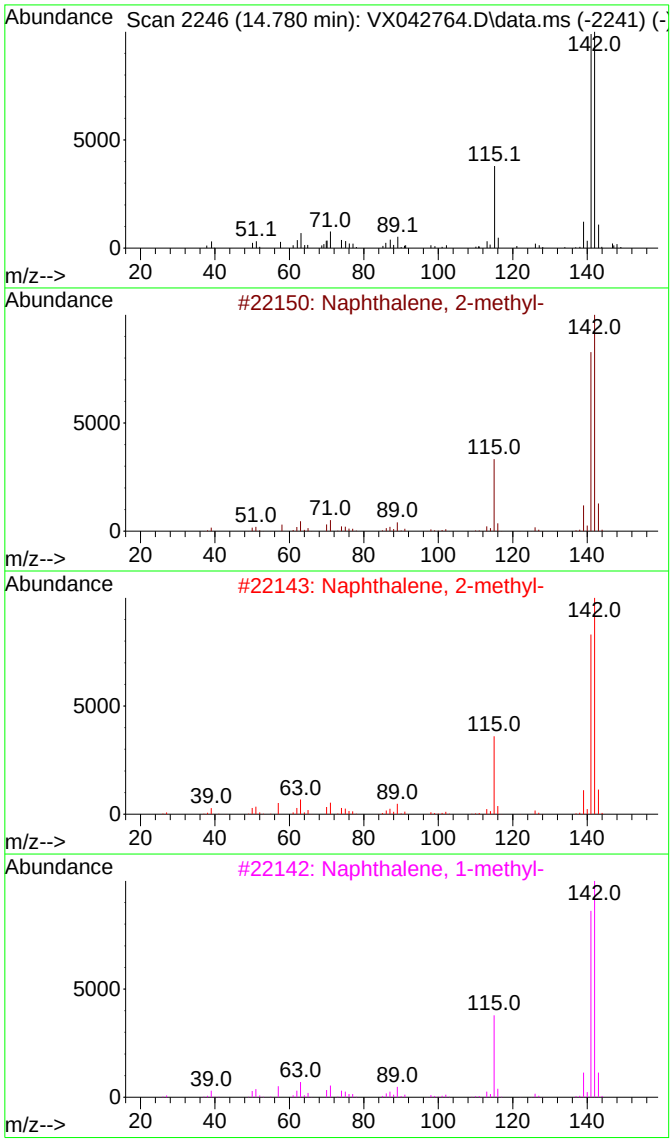
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

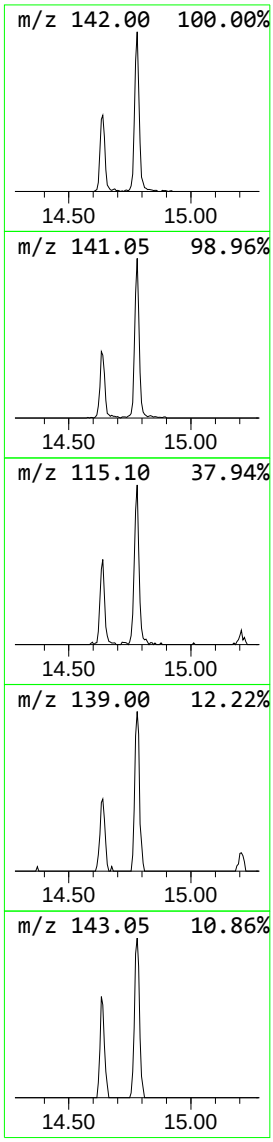
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 1-ethylidene- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	13.42 ug/L	78842	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	95
5	Benzocycloheptatriene	142	C11H10	000264-09-5	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

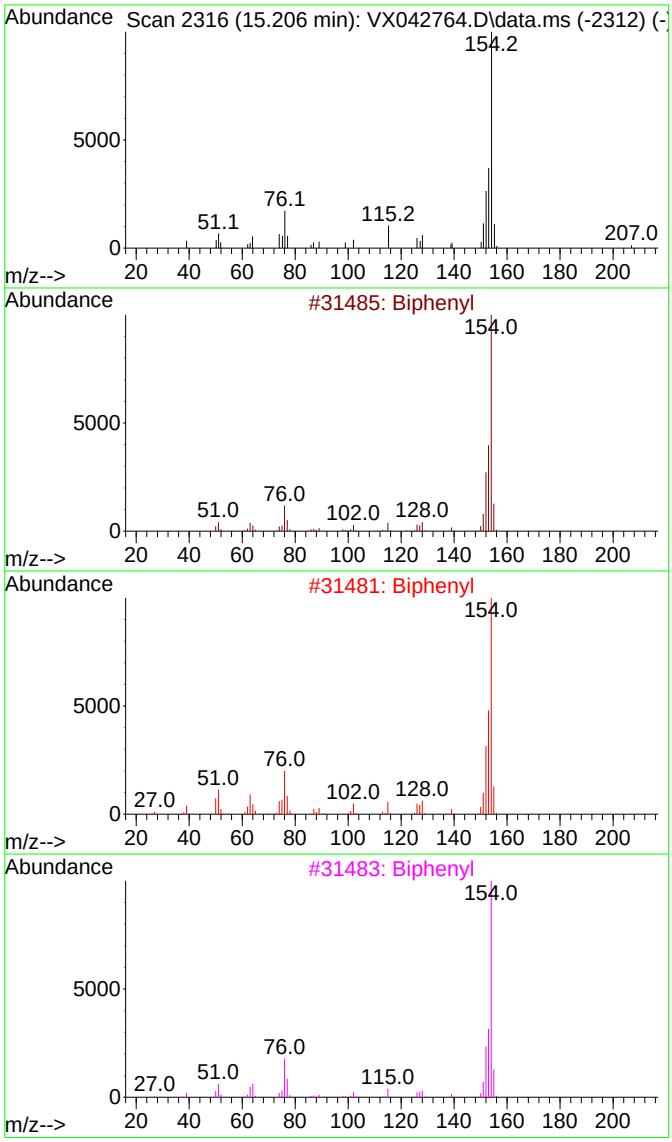
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

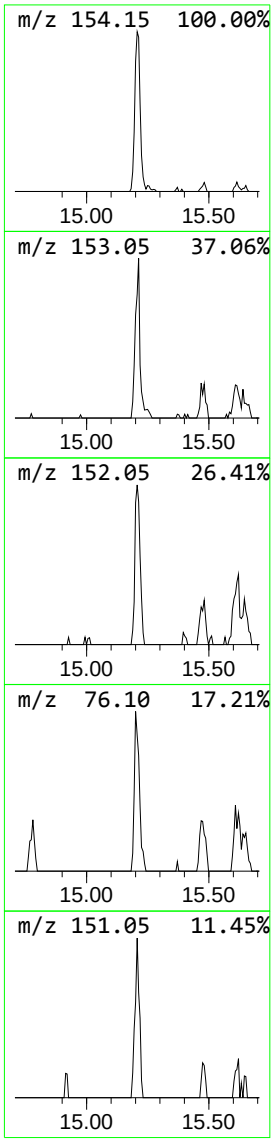
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 2-ethenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.207	4.12 ug/L	24191	1,4-Dichlorobenzene-d4	12.024

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	Biphenyl	154	C12H10	000092-52-4	81
4	Biphenyl	154	C12H10	000092-52-4	72
5	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

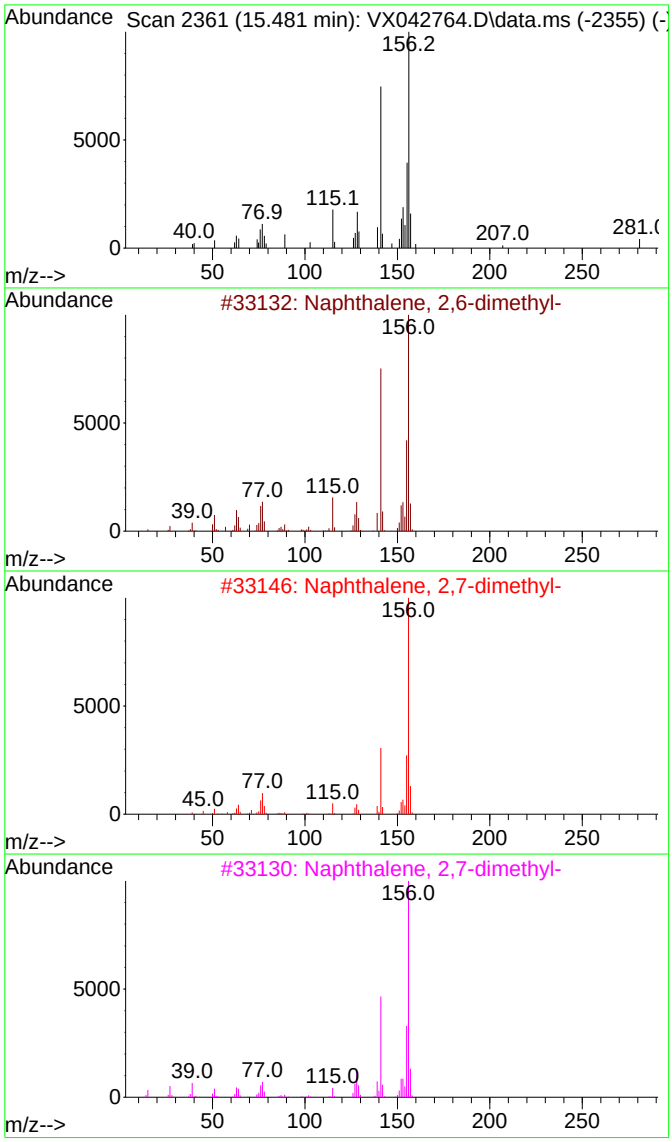
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

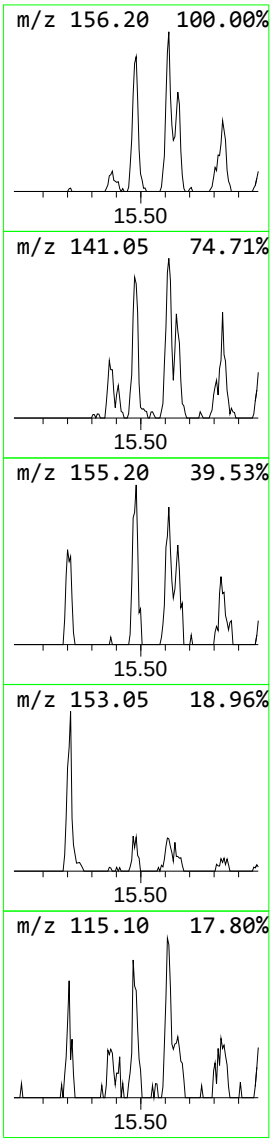
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 2,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.481	3.74 ug/L	21977	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

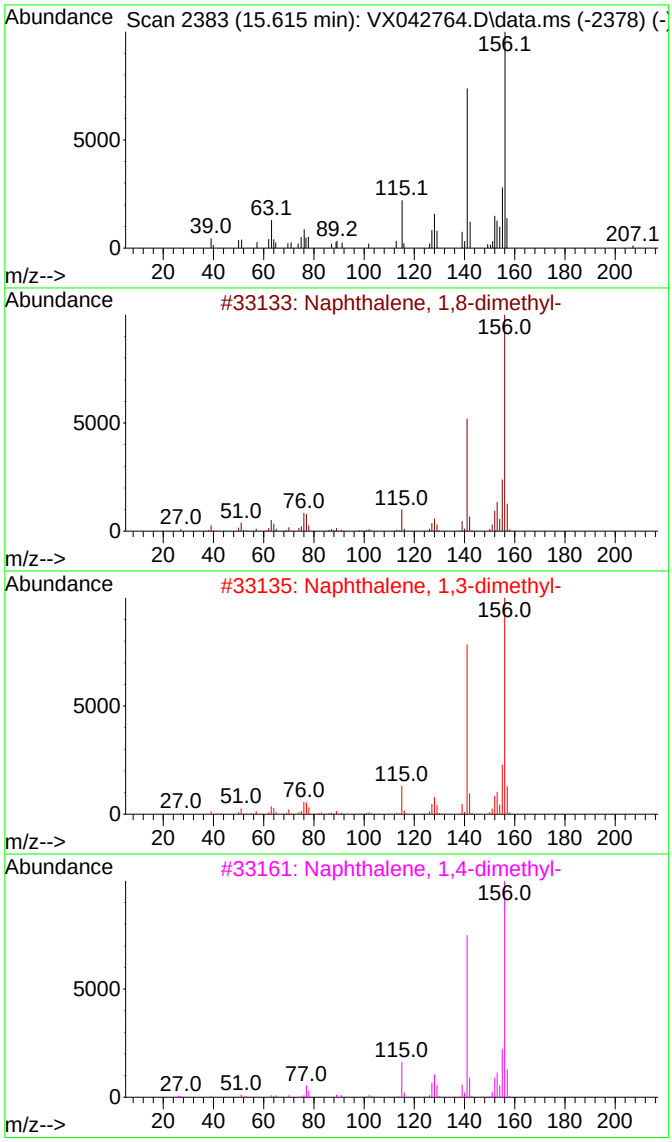
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

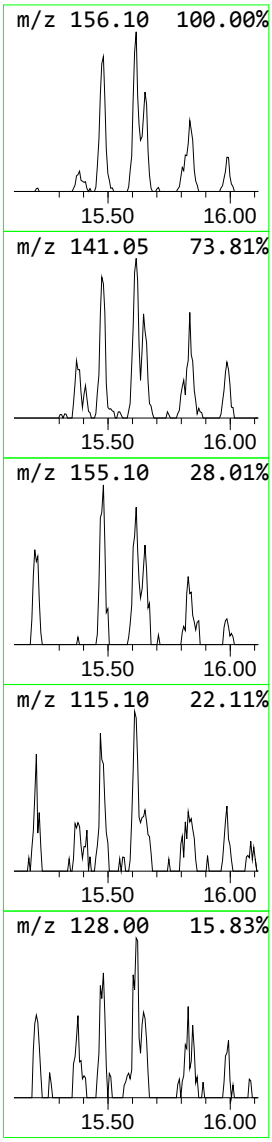
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Naphthalene, 1,8-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	3.53 ug/L	20739	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

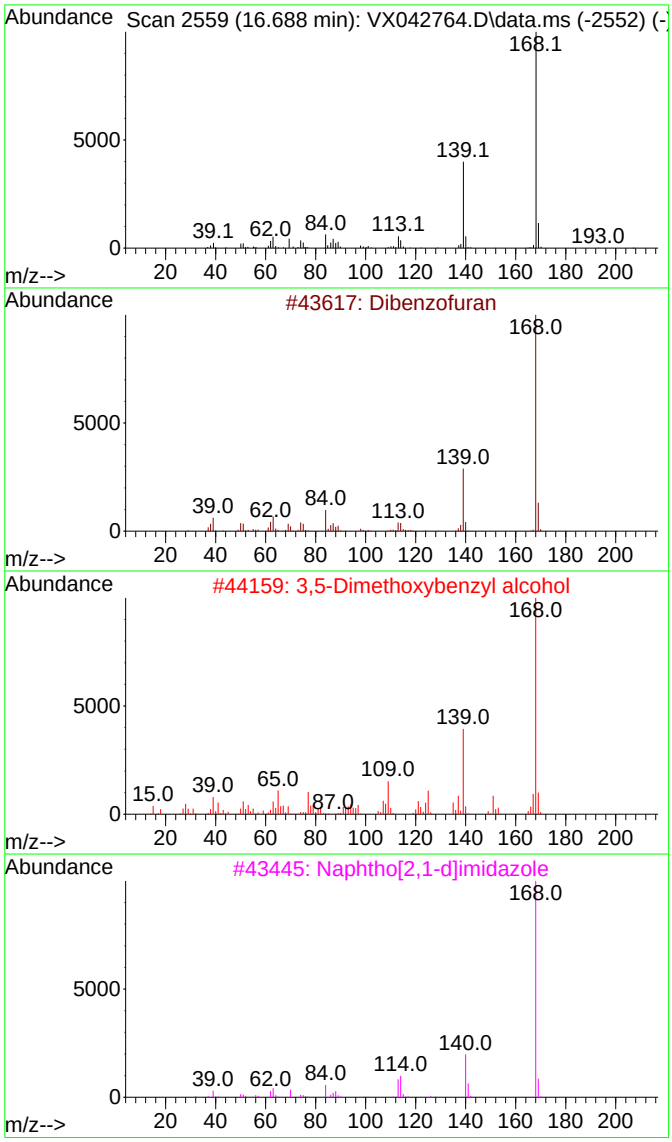
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.M
Quant Title : VOC Analysis

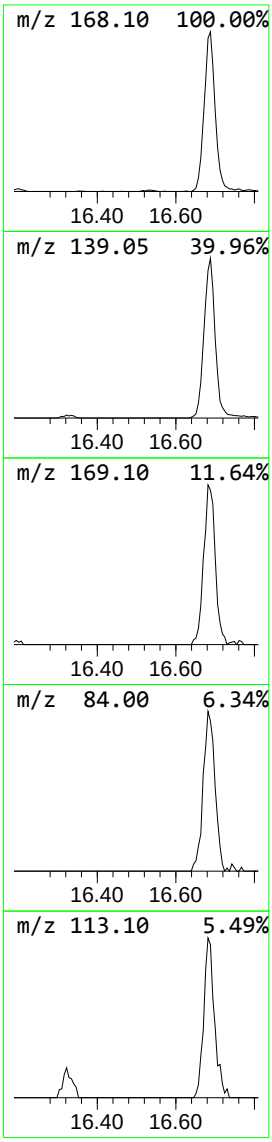
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 3,5-Dimethoxybenzyl alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.688	32.49 ug/L	190819	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	87
2	3,5-Dimethoxybenzyl alcohol	168	C9H12O3	000705-76-0	56
3	Naphtho[2,1-d]imidazole	168	C11H8N2	000233-53-4	52
4	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	50
5	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073124\
Data File : VX042764.D
Acq On : 31 Jul 2024 11:44
Operator : JC/MD
Sample : P3378-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
C0AY3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM072624WMA.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	1.239	4.9	ug/L	24910	1	6.763	254249	50.0
Benzene, cyclop...	12.244	3.8	ug/L	22158	3	12.024	293678	50.0
1H-Phenalene	12.811	78.7	ug/L	462203	3	12.024	293678	50.0
Azulene	13.774	51.3	ug/L	301408	3	12.024	293678	50.0
Benzo[b]thiophene	13.871	2.6	ug/L	15243	3	12.024	293678	50.0
1H-Indene, 1-et...	14.780	13.4	ug/L	78842	3	12.024	293678	50.0
Naphthalene, 2-...	15.207	4.1	ug/L	24191	3	12.024	293678	50.0
Naphthalene, 2,...	15.481	3.7	ug/L	21977	3	12.024	293678	50.0
Naphthalene, 1,...	15.615	3.5	ug/L	20739	3	12.024	293678	50.0
3,5-Dimethoxybe...	16.688	32.5	ug/L	190819	3	12.024	293678	50.0