LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : F4L18

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\SFAMXLM111121WMA.M
Title : VOC Analysis

Signal : TIC: VX025240.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.246	22	26	42	rBV	83292	188121	33.33%	3.206%
2	1.368	43	46	53	rVB	93339	90333	16.01%	1.540%
3	1.666	91	95	104	rVB	65082	86248	15.28%	1.470%
4	2.307	194	200	210	rBV	122208	197190	34.94%	3.361%
5	2.441	220	222	228	rBV2	250	460	0.08%	0.008%
6	2.947	296	305	312	rBV2	2079	5279	0.94%	0.090%
7	4.459	544	553	566	rBV	44017	121529	21.53%	2.071%
8	5.062	640	652	671	rVV	69678	216557	38.37%	3.691%
9	5.190	671	673	677	rVB2	188	252	0.04%	0.004%
10	5.233	677	680	684	rBV2	200	309	0.05%	0.005%
11	5.550	725	732	734	rBV3	324	611	0.11%	0.010%
12	5.977	789	802	822	rBV2	185224	564361	100.00%	9.618%
13	6.385	862	869	877	rVB2	1223	2887	0.51%	0.049%
14	6.629	908	909	913	rBV2	258	230	0.04%	0.004%
15	6.763	921	931	947	rBV	154483	360129	63.81%	6.138%
16	7.111	983	988	994	rBV3	934	1909	0.34%	0.033%
17	7.312	1011	1021	1035	rVV	112658	248286	43.99%	4.232%
18	7.428	1037	1040	1045	rVB2	189	312	0.06%	0.005%
19	7.482	1045	1049	1054	rBV3	318	499	0.09%	0.009%
20	7.806	1098	1102	1104	rBV2	374	492	0.09%	0.008%
21	7.927	1115	1122	1128	rVV5	768	1734	0.31%	0.030%
22	8.324	1181	1187	1200	rBV	81544	133268	23.61%	2.271%
23	8.476	1207	1212	1221	rVB3	728	1721	0.30%	0.029%
24	8.653	1234	1241	1248	rBV	291319	465013	82.40%	7.925%
25	8.720	1248	1252	1262	rVV	11740	18653	3.31%	0.318%
26	8.952	1284	1290	1299	rBV	59497	85925	15.23%	1.464%
27	9.281	1336	1344	1348	rBV2	1561	2081	0.37%	0.035%
28	9.385	1355	1361	1375	rBV	199909	277025	49.09%	4.721%
29	10.055	1465	1471	1489	rVB	316616	430979	76.37%	7.345%
30	10.195	1489	1494	1499	rBV	8140	10228	1.81%	0.174%
31	10.305	1505	1512	1517	rBV	5540	7607	1.35%	0.130%
32	10.464	1533	1538	1539	rVB2	251	312	0.06%	0.005%
33	10.647	1563	1568	1573	rBV	9056	12417	2.20%	0.212%
34	10.836	1597	1599	1602	rVB3	229	245	0.04%	0.004%
35	10.970	1616	1621	1625	rBV2	1113	1462	0.26%	0.025%
36	11.122	1642	1646	1651	rVB	2035	2407	0.43%	0.041%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

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\SFAMXLM111121WMA.M
Title : VOC Analysis

37	11.195	1652	1658	1665	rBV	217421	274053	48.56%	4.671%
38	11.323	1674	1679	1683	rVB4	2112	3994	0.71%	0.068%
39	11.378	1683	1688	1690	rBV2	679	946	0.17%	0.016%
40	11.402	1690	1692	1696	rVB	1264	1312	0.23%	0.022%
41	11.451	1696	1700	1703	rBV	3570	4443	0.79%	0.076%
42	11.616	1721	1727	1732	rBB	3785	4658	0.83%	0.079%
43	11.750	1745	1749	1754	rVB2	8488	11117	1.97%	0.189%
44	11.811	1756	1759	1762	rVB3	494	427	0.08%	0.007%
45	11.848	1762	1765	1770	rBV4	340	533	0.09%	0.009%
46	11.884	1770	1771	1775	rVB2	469	459	0.08%	0.008%
47	11.933	1775	1779	1781	rBV3	916	1180	0.21%	0.020%
48	12.024	1788	1794	1801	rBV	357289	433258	76.77%	7.384%
49	12.091	1801	1805	1808	rVV	5054	6284	1.11%	0.107%
50	12.244	1825	1830	1837	rVV	138373	165759	29.37%	2.825%
51	12.323	1837	1843	1851	rVB	336089	429771	76.15%	7.325%
52	12.414	1853	1858	1864	rVV	52948	64624	11.45%	1.101%
53	12.530	1873	1877	1879	rBV4	986	1423	0.25%	0.024%
54	12.616	1888	1891	1894	rBV	1304	1517	0.27%	0.026%
55	12.646	1894	1896	1897	rVV2	485	427	0.08%	0.007%
56	12.701	1901	1905	1914	rVB	7385	9278	1.64%	0.158%
57	12.799	1917	1921	1924	rBV	3520	4370	0.77%	0.074%
58	12.890	1932	1936	1944	rVB3	29373	53081	9.41%	0.905%
59	12.969	1945	1949	1952	rVV2	18067	28652	5.08%	0.488%
60	13.030	1956	1959	1964	rVB	9201	11154	1.98%	0.190%
61	13.134	1973	1976	1978	rBV3	523	582	0.10%	0.010%
62	13.170	1978	1982	1990	rVV2	4619	5873	1.04%	0.100%
63	13.256	1994	1996	1998	rBV2	362	331	0.06%	0.006%
64	13.292	1998	2002	2006	rVV2	10819	14878	2.64%	0.254%
65	13.341	2006	2010	2015	rVB2	13858	17778	3.15%	0.303%
66	13.427	2019	2024	2029	rBV	16050	20622	3.65%	0.351%
67	13.506	2032	2037	2041	rBV4	4453	7019	1.24%	0.120%
68	13.542	2041	2043	2046	rVV3	1461	1795	0.32%	0.031%
69	13.579	2046	2049	2054	rVB3	2995	3684	0.65%	0.063%
70	13.622	2054	2056	2057	rBV	512	542	0.10%	0.009%
71	13.664	2061	2063	2068	rVB3	1564	1988	0.35%	0.034%
72	13.780	2074	2082	2089	rBV	128692	181909	32.23%	3.100%
73	13.872	2093	2097	2103	rVV	31995	44608	7.90%	0.760%
74	13.926	2103	2106	2109	rVB3	2552	2922	0.52%	0.050%
75	13.963	2109	2112	2116	rBV2	4336	5249	0.93%	0.089%
76	14.000	2116	2118	2123	rVB4	2268	2741	0.49%	0.047%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : F4L18

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\SFAMXLM111121WMA.M
Title : VOC Analysis

77	14.073	2127	2130	2134	rBV2	1188	2069	0.37%	0.035%
78	14.219	2150	2154	2157	rBV3	3052	4085	0.72%	0.070%
79	14.323	2166	2171	2174	rBV3	2214	2938	0.52%	0.050%
80	14.384	2176	2181	2186	rVB2	12660	16671	2.95%	0.284%
81	14.451	2189	2192	2193	rBV2	310	312	0.06%	0.005%
82	14.597	2212	2216	2220	rBV	14205	18023	3.19%	0.307%
83	14.640	2220	2223	2232	rVB	8181	11453	2.03%	0.195%
84	14.725	2234	2237	2241	rBV3	1028	1158	0.21%	0.020%
85	14.780	2241	2246	2257	rBV2	69882	93563	16.58%	1.595%
86	14.999	2280	2282	2287	rVB2	409	452	0.08%	0.008%
87	15.121	2300	2302	2305	rBV3	225	322	0.06%	0.005%
88	15.207	2312	2316	2321	rVB2	11303	16518	2.93%	0.282%
89	15.262	2323	2325	2329	rVB2	405	357	0.06%	0.006%
90	15.341	2335	2338	2340	rBV3	357	438	0.08%	0.007%
91	15.408	2346	2349	2353	rVB3	2237	3191	0.57%	0.054%
92	15.481	2356	2361	2365	rBV2	4341	6907	1.22%	0.118%
93	15.615	2374	2383	2388	rBV2	9062	14577	2.58%	0.248%
94	15.731	2400	2402	2406	rBV2	433	754	0.13%	0.013%
95	15.816	2411	2416	2417	rBV3	3034	4588	0.81%	0.078%
96	15.993	2440	2445	2451	rVB2	2813	4333	0.77%	0.074%
97	16.091	2457	2461	2468	rVB2	1917	3248	0.58%	0.055%
98	16.328	2493	2500	2512	rBV	158097	275142	48.75%	4.689%
99	16.536	2530	2534	2540	rVB2	1322	2345	0.42%	0.040%
100	16.688	2553	2559	2568	rVB	11848	21719	3.85%	0.370%

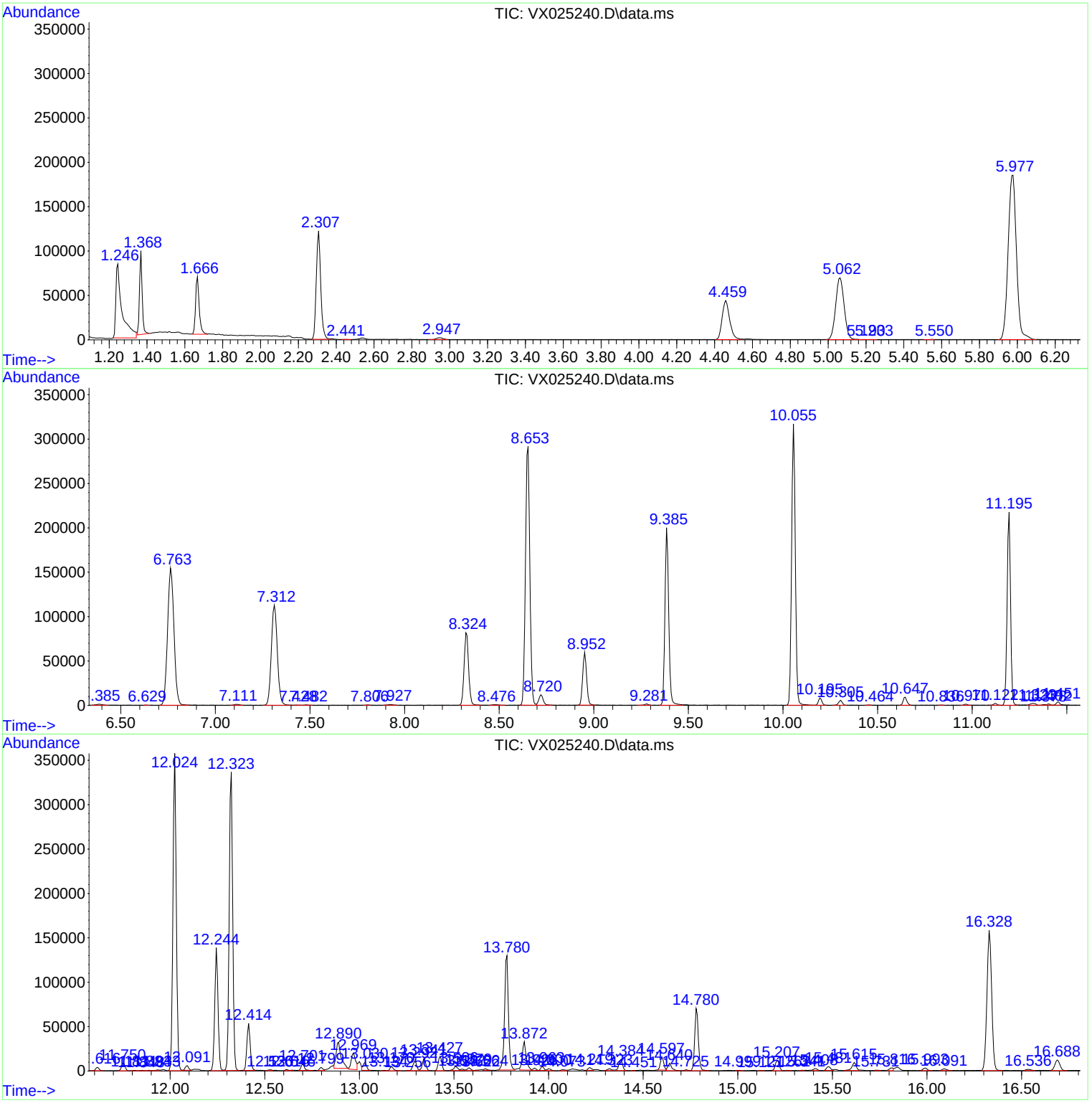
Sum of corrected areas: 5867475

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

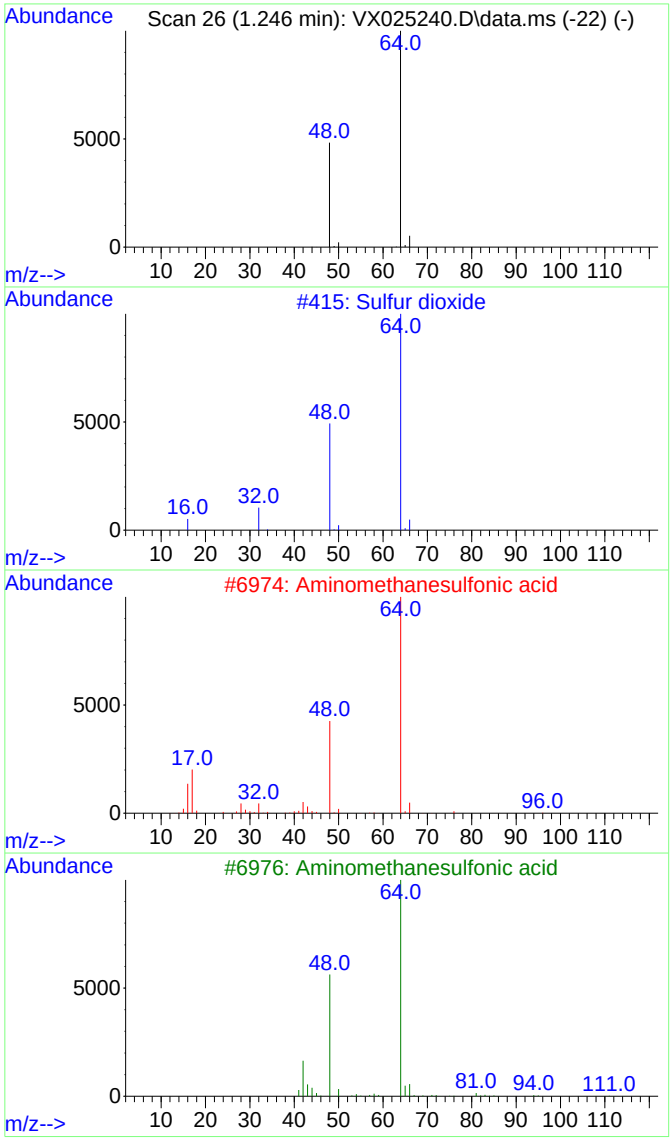
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

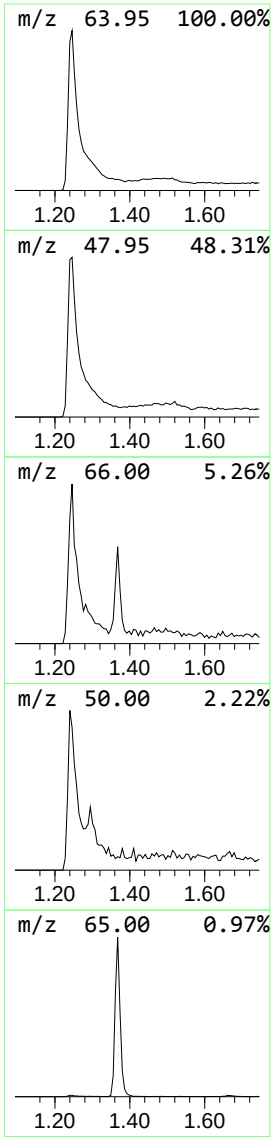
****

Peak Number 1 Sulfur dioxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.246	26.12 ug/L	188121	1,4-Difluorobenzene	6.763

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

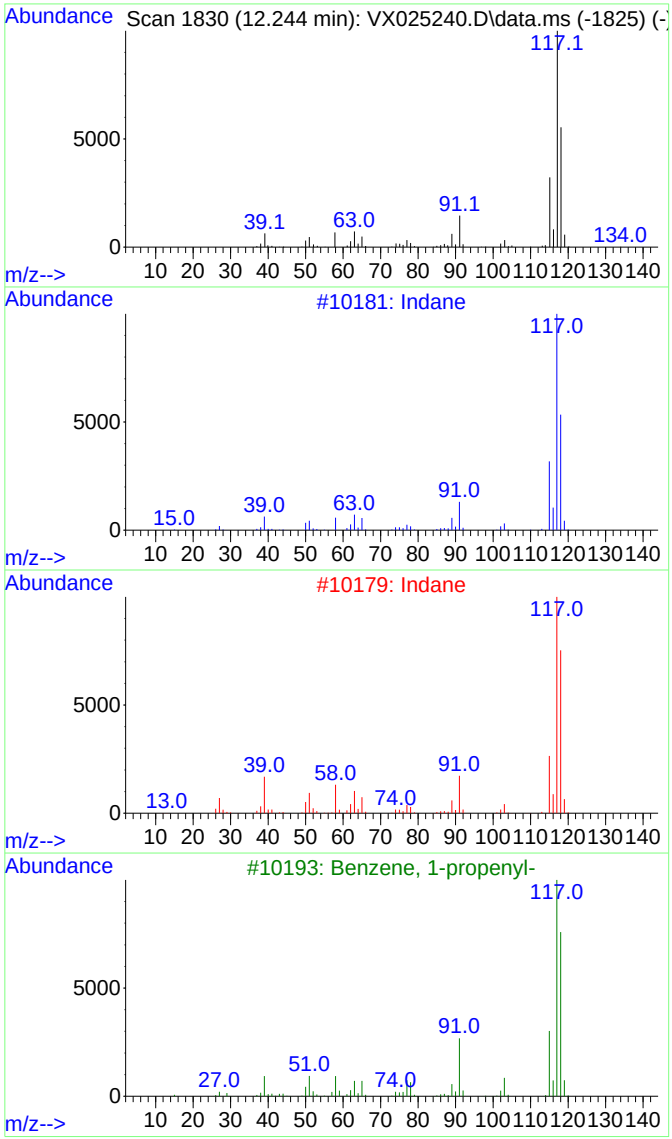
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

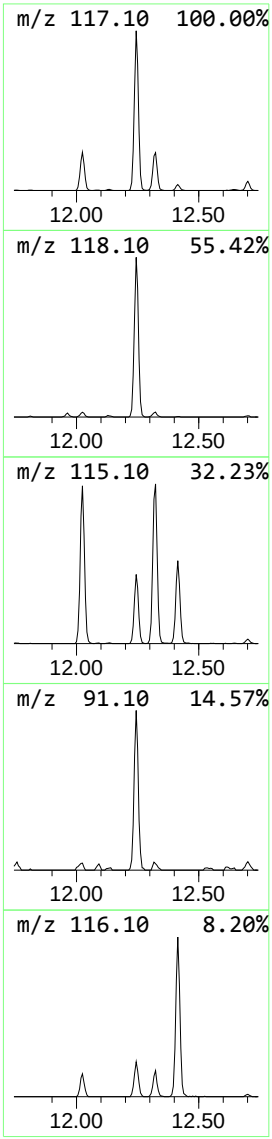
****

Peak Number 3 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	19.13 ug/L	165759	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Indane	118	C9H10	000496-11-7	91
3	Benzene, 1-propenyl-	118	C9H10	000637-50-3	83
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	83
5	Deltacyclene	118	C9H10	007785-10-6	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

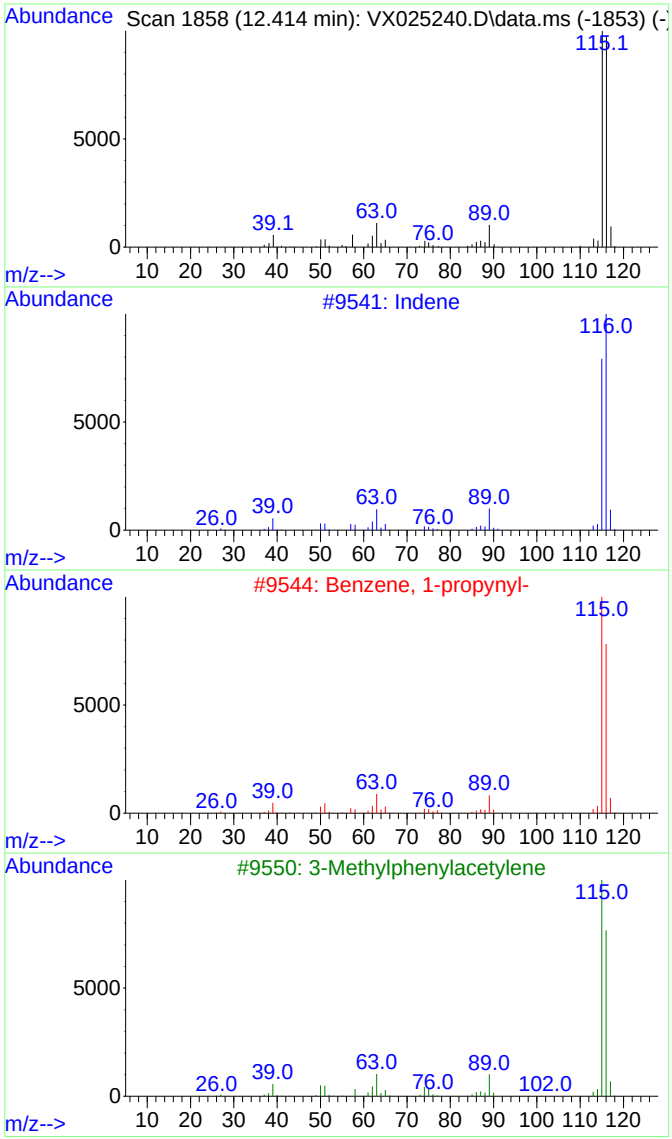
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

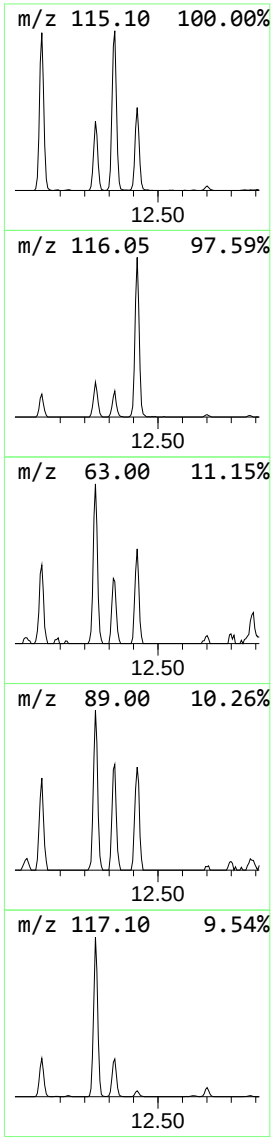
****

Peak Number 4 Indene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.415	7.46 ug/L	64624	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
3	3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
5	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

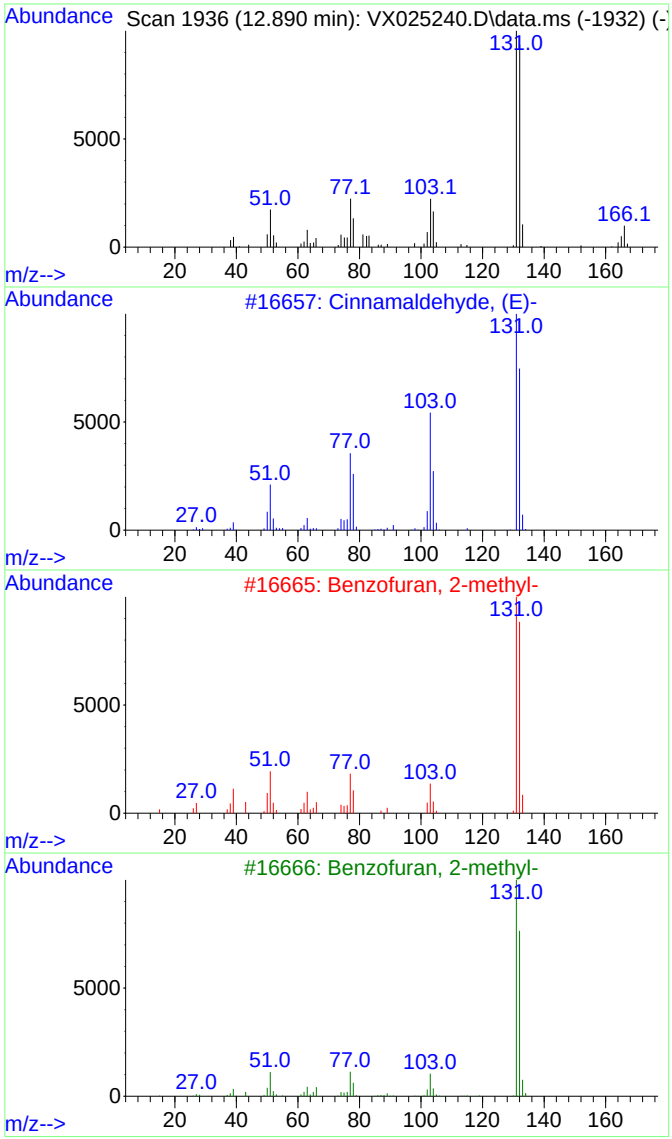
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

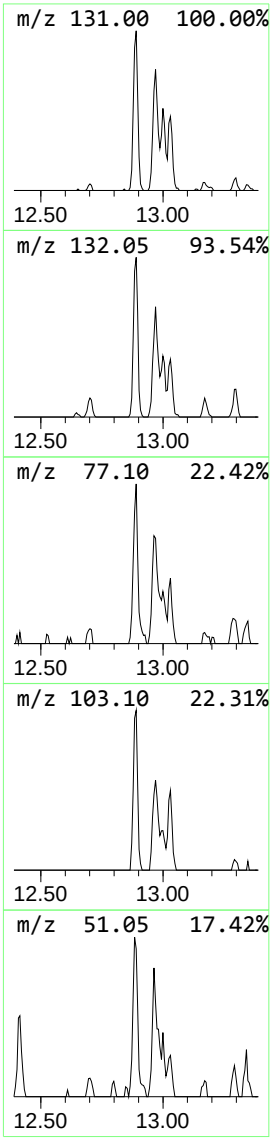
****

Peak Number 5 Cinnamaldehyde, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	6.13 ug/L	53081	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	90
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
4	5-Methylbenzimidazole	132	C8H8N2	000614-97-1	87
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

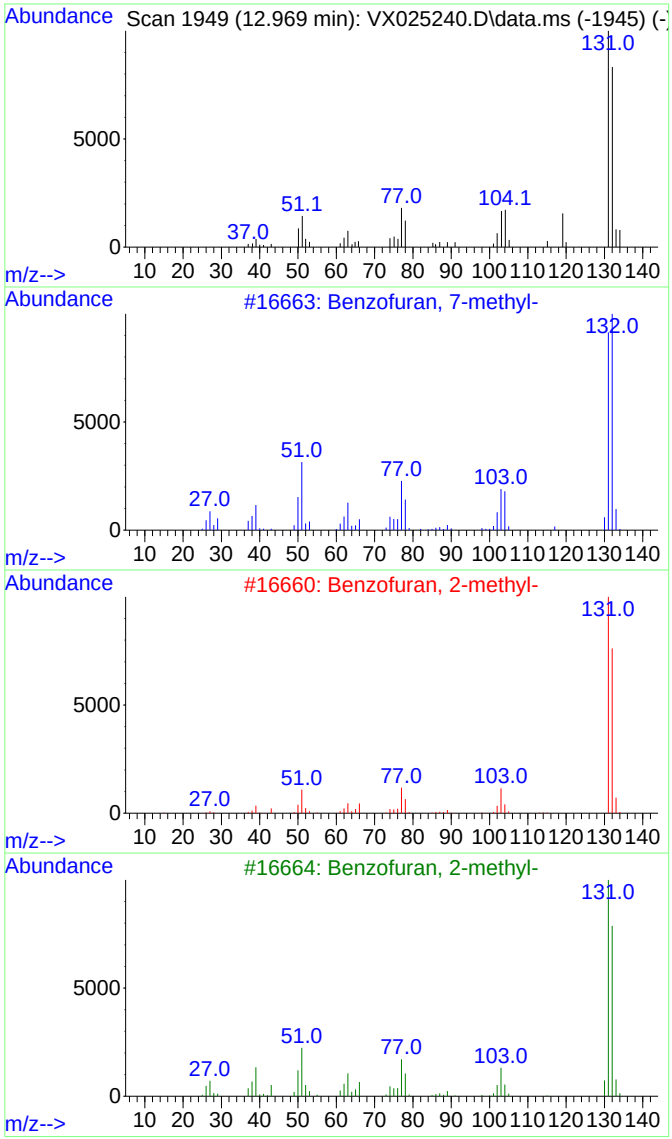
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

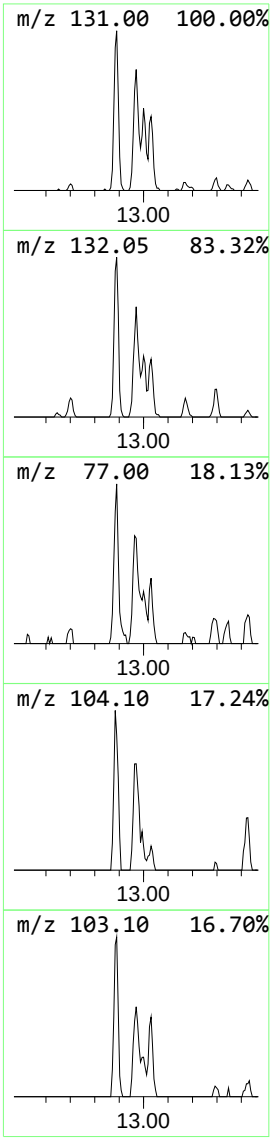
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzofuran, 7-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.969	3.31 ug/L	28652	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

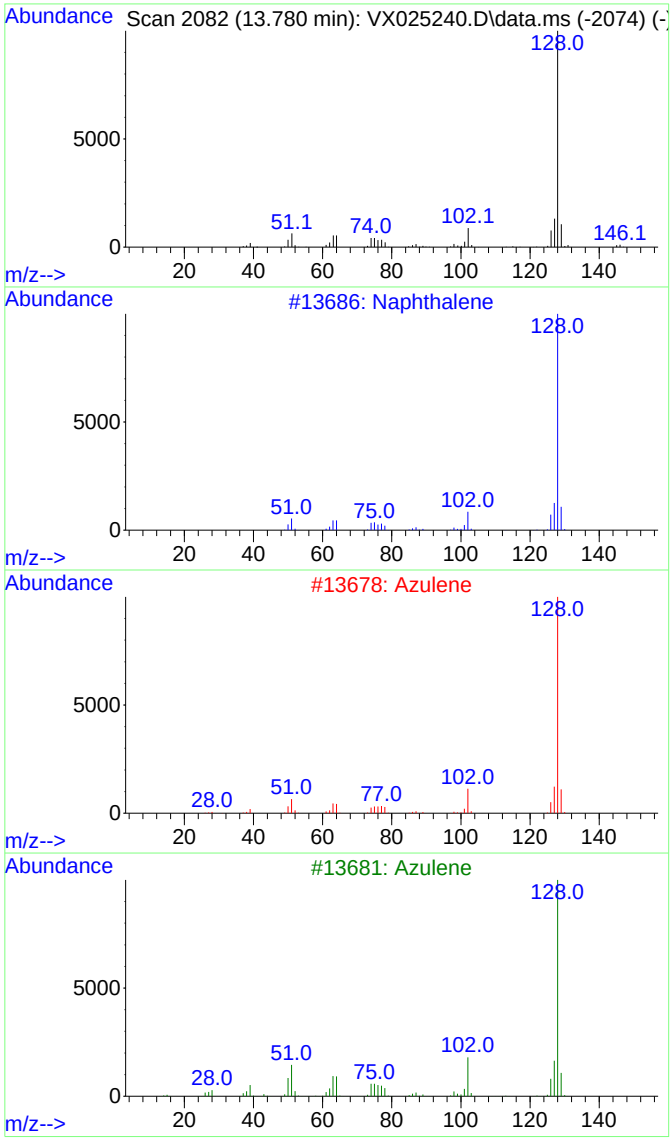
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

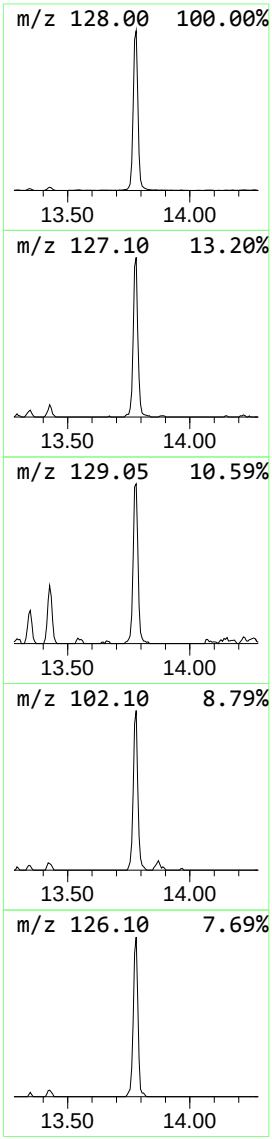
****

Peak Number 7 Azulene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	20.99 ug/L	181909	1,4-Dichlorobenzene-d4	12.024

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

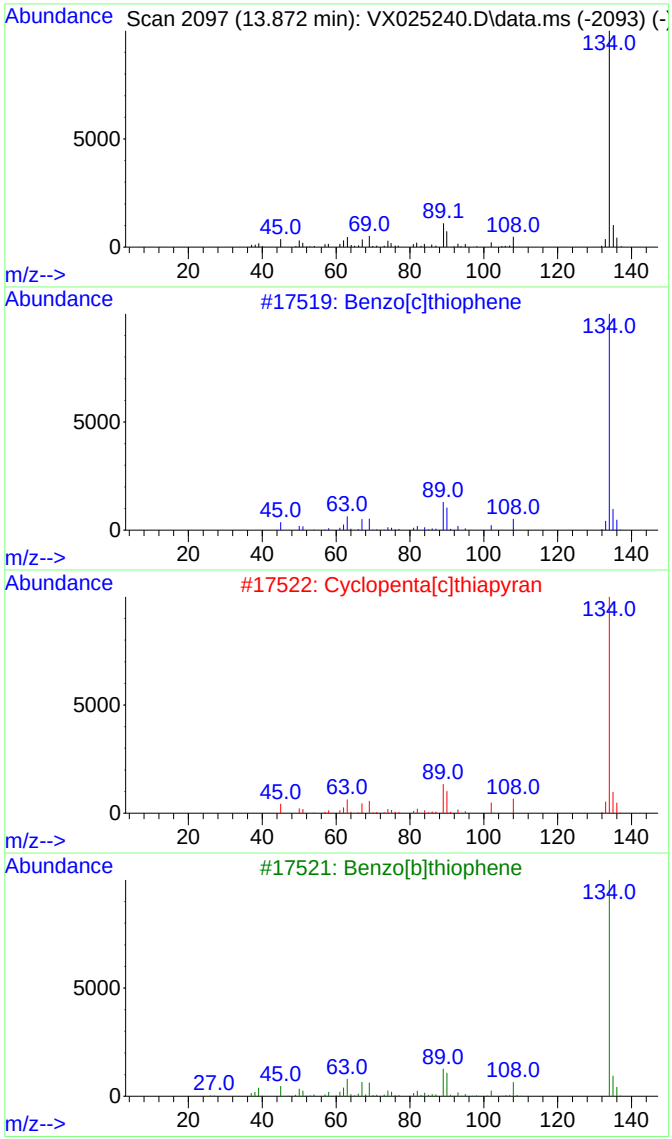
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

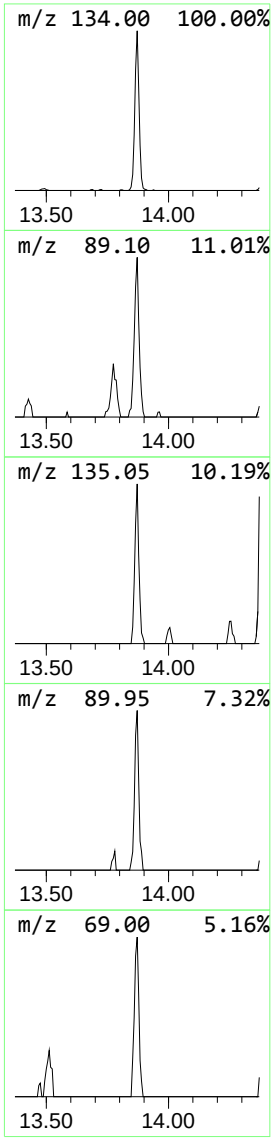
****

Peak Number 8 Benzo[c]thiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.872	5.15 ug/L	44608	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
3	Benzo[b]thiophene	134	C8H6S	000095-15-8	93
4	Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5	Benzo[b]thiophene	134	C8H6S	000095-15-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

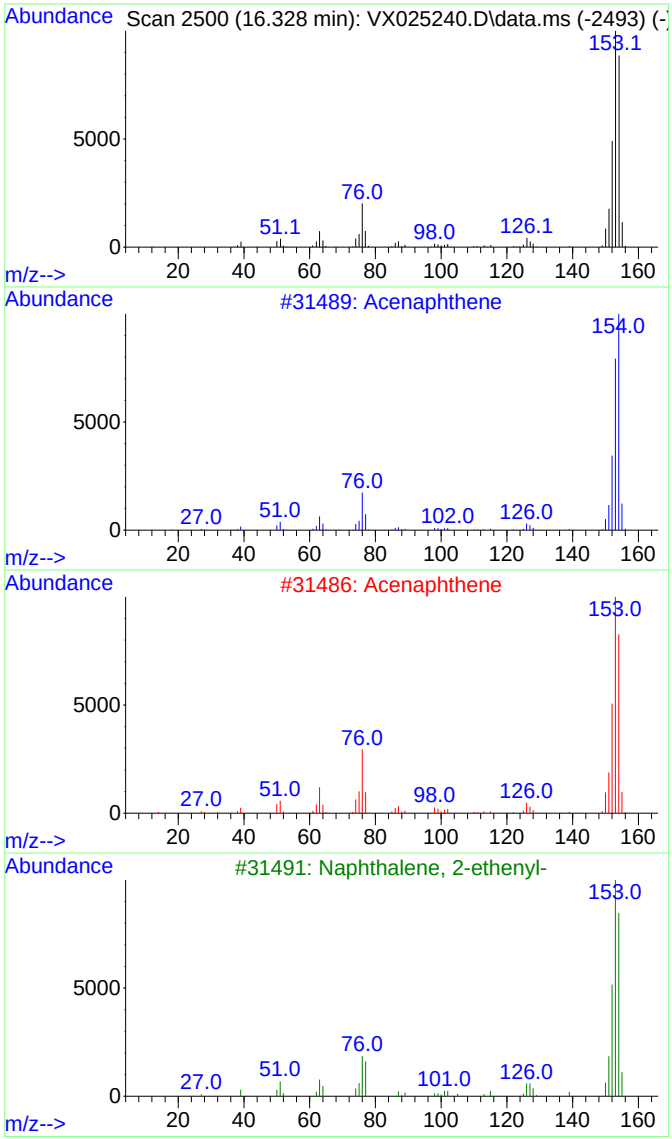
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

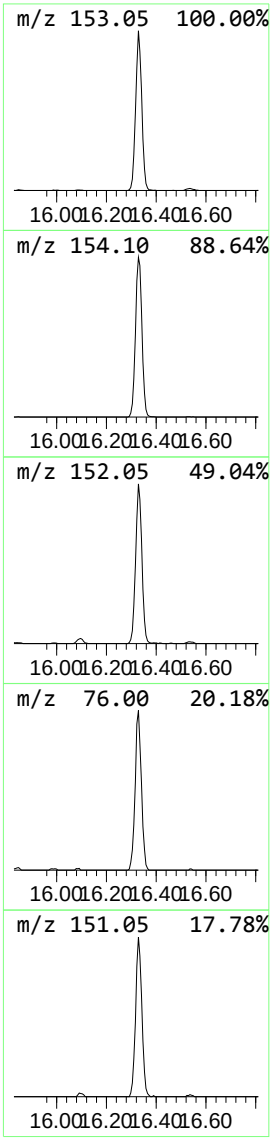
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 2-ethenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.328	31.75 ug/L	275142	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthene	154	C12H10	000083-32-9	93
2	Acenaphthene	154	C12H10	000083-32-9	89
3	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	64
4	Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	59
5	2,4,6-Trihydroxybenzaldehyde	154	C7H6O4	000487-70-7	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

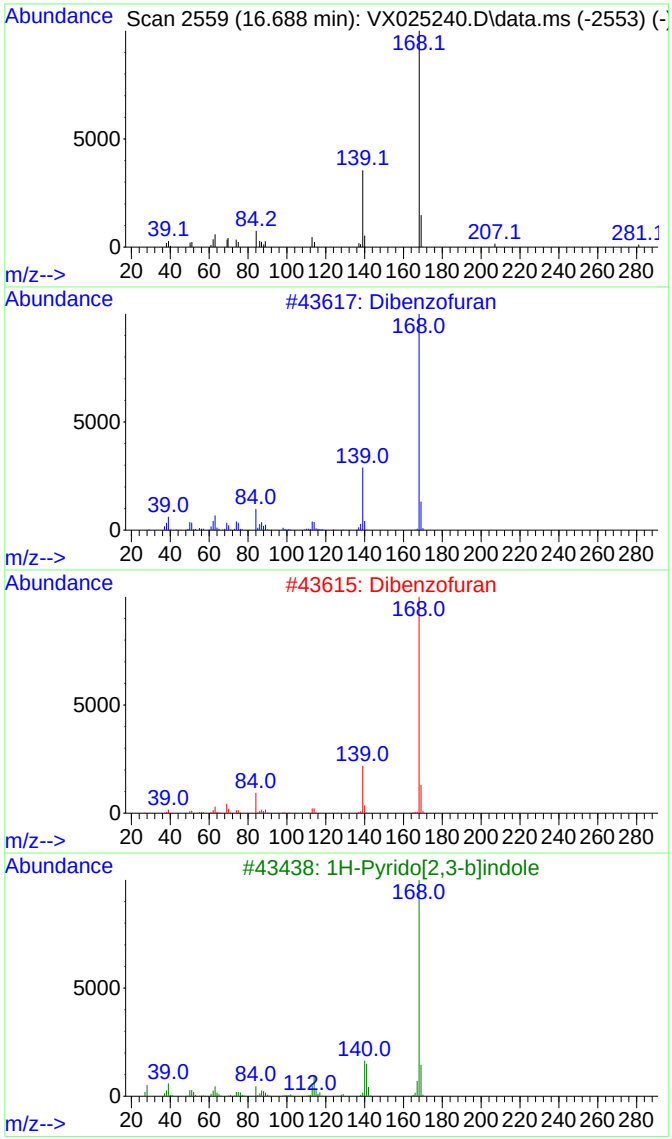
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M
Quant Title : VOC Analysis

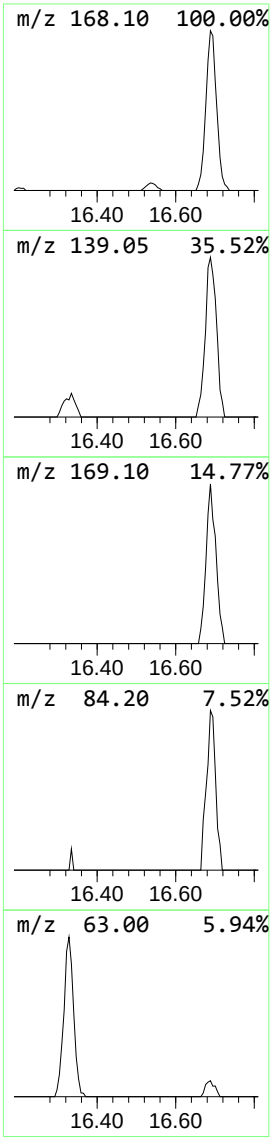
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1H-Pyrido[2,3-b]indole Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.688	2.51 ug/L	21719	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	90
2	Dibenzofuran	168	C12H8O	000132-64-9	80
3	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	59
4	1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	59
5	9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111921\
Data File : VX025240.D
Acq On : 19 Nov 2021 17:52
Operator : JC/MD
Sample : M4779-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
F4L18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.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	1.246	26.1	ug/L	188121	1	6.763	360129	50.0
Indane	12.244	19.1	ug/L	165759	3	12.024	433258	50.0
Indene	12.415	7.5	ug/L	64624	3	12.024	433258	50.0
Cinnamaldehyde,...	12.890	6.1	ug/L	53081	3	12.024	433258	50.0
Benzofuran, 7-m...	12.969	3.3	ug/L	28652	3	12.024	433258	50.0
Azulene	13.780	21.0	ug/L	181909	3	12.024	433258	50.0
Benzo[c]thiophene	13.872	5.2	ug/L	44608	3	12.024	433258	50.0
Naphthalene, 2-...	16.328	31.8	ug/L	275142	3	12.024	433258	50.0
1H-Pyrido[2,3-b...	16.688	2.5	ug/L	21719	3	12.024	433258	50.0