LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

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\SFAMXLM112221WMA.M
Title : VOC Analysis

Signal : TIC: VX025741.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	23	26	32	rBV2	5602	9470	0.04%	0.015%
2	1.374	43	47	58	rBV	1886222	1857349	7.12%	2.939%
3	1.685	91	98	110	rVB2	103088	178944	0.69%	0.283%
4	2.319	194	202	216	rBV3	144213	304610	1.17%	0.482%
5	2.520	233	235	245	rVB	1195	1848	0.01%	0.003%
6	2.794	272	280	287	rBV	10589	20621	0.08%	0.033%
7	2.965	300	308	314	rVB5	1404	3769	0.01%	0.006%
8	3.099	322	330	350	rVV	215533	470122	1.80%	0.744%
9	3.623	404	416	441	rBV	4589737	11077546	42.49%	17.526%
10	4.178	504	507	510	rBV	322	398	0.00%	0.001%
11	4.300	522	527	528	rBV	845	1107	0.00%	0.002%
12	4.507	546	561	597	rBV	9996901	26071182	100.00%	41.248%
13	5.068	643	653	669	rVB	43183	143025	0.55%	0.226%
14	5.391	692	706	720	rBV	295040	917823	3.52%	1.452%
15	5.836	776	779	782	rBV2	466	671	0.00%	0.001%
16	5.976	791	802	810	rBV2	109925	334660	1.28%	0.529%
17	6.245	842	846	847	rBV2	816	961	0.00%	0.002%
18	6.397	868	871	874	rVB2	322	392	0.00%	0.001%
19	6.434	874	877	878	rBV	357	345	0.00%	0.001%
20	6.769	923	932	945	rBV	83815	198875	0.76%	0.315%
21	7.129	982	991	1010	rBV	654377	1421222	5.45%	2.249%
22	7.312	1013	1021	1029	rBV	66494	148934	0.57%	0.236%
23	7.482	1044	1049	1055	rVB3	621	1338	0.01%	0.002%
24	7.562	1059	1062	1064	rBV	244	387	0.00%	0.001%
25	7.665	1074	1079	1085	rVB3	1745	3301	0.01%	0.005%
26	7.909	1113	1119	1125	rVB3	1786	3801	0.01%	0.006%
27	7.976	1125	1130	1134	rVB3	1040	1482	0.01%	0.002%
28	8.330	1181	1188	1196	rBV	49384	81870	0.31%	0.130%
29	8.580	1224	1229	1235	rBV3	2035	3750	0.01%	0.006%
30	8.653	1235	1241	1247	rVV	173303	268188	1.03%	0.424%
31	8.720	1247	1252	1259	rVV	149266	232103	0.89%	0.367%
32	8.805	1262	1266	1275	rVB2	3486	6182	0.02%	0.010%
33	8.952	1284	1290	1298	rBV	39664	58604	0.22%	0.093%
34	9.153	1318	1323	1331	rBV	45051	66827	0.26%	0.106%
35	9.275	1334	1343	1354	rVB	148568	218146	0.84%	0.345%
36	9.384	1356	1361	1378	rBV	110370	163498	0.63%	0.259%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

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\SFAMXLM112221WMA.M
Title : VOC Analysis

37	9.561	1386	1390	1395	rVB2	2124	3072	0.01%	0.005%
38	9.616	1397	1399	1404	rVB3	843	1180	0.00%	0.002%
39	9.726	1416	1417	1423	rVB3	585	784	0.00%	0.001%
40	10.012	1460	1464	1466	rBV	400	490	0.00%	0.001%
41	10.055	1466	1471	1479	rBV	181494	246753	0.95%	0.390%
42	10.195	1488	1494	1505	rBV	492117	636007	2.44%	1.006%
43	10.305	1506	1512	1524	rVB	898880	1239428	4.75%	1.961%
44	10.463	1535	1538	1542	rBV3	541	609	0.00%	0.001%
45	10.591	1555	1559	1562	rBV3	1003	1660	0.01%	0.003%
46	10.646	1562	1568	1581	rVB	1224505	1558521	5.98%	2.466%
47	10.780	1586	1590	1597	rVB3	4674	6199	0.02%	0.010%
48	10.854	1597	1602	1608	rBV5	1629	3057	0.01%	0.005%
49	10.963	1615	1620	1628	rBV	186801	232525	0.89%	0.368%
50	11.079	1634	1639	1640	rBV2	1575	2688	0.01%	0.004%
51	11.128	1645	1647	1652	rVV3	3493	4243	0.02%	0.007%
52	11.195	1652	1658	1668	rVV	139989	181897	0.70%	0.288%
53	11.305	1671	1676	1683	rVV	358916	447065	1.71%	0.707%
54	11.384	1683	1689	1696	rVV	1269715	2053480	7.88%	3.249%
55	11.457	1696	1701	1716	rVB	707508	894870	3.43%	1.416%
56	11.573	1716	1720	1722	rBV2	6331	8162	0.03%	0.013%
57	11.616	1722	1727	1736	rVB	983919	1199655	4.60%	1.898%
58	11.756	1745	1750	1755	rBV	3532978	4217255	16.18%	6.672%
59	11.866	1764	1768	1770	rBV	15080	19192	0.07%	0.030%
60	11.890	1770	1772	1778	rVB	34078	40063	0.15%	0.063%
61	11.975	1781	1786	1789	rBV	80072	98132	0.38%	0.155%
62	12.024	1789	1794	1799	rVB2	228127	300633	1.15%	0.476%
63	12.091	1799	1805	1810	rBV	1565120	1893583	7.26%	2.996%
64	12.183	1816	1820	1824	rVB	12378	16653	0.06%	0.026%
65	12.244	1825	1830	1834	rVV	562208	781163	3.00%	1.236%
66	12.323	1838	1843	1851	rVB3	397678	681084	2.61%	1.078%
67	12.408	1852	1857	1862	rBV2	23852	30309	0.12%	0.048%
68	12.463	1862	1866	1872	rVB	67567	82188	0.32%	0.130%
69	12.548	1873	1880	1887	rBV3	345922	607656	2.33%	0.961%
70	12.615	1887	1891	1894	rVV	153647	178862	0.69%	0.283%
71	12.646	1894	1896	1901	rVB	28311	31011	0.12%	0.049%
72	12.701	1901	1905	1914	rBV2	67269	123948	0.48%	0.196%
73	12.798	1916	1921	1925	rVV3	9253	17160	0.07%	0.027%
74	12.853	1925	1930	1935	rVB2	55612	71890	0.28%	0.114%
75	12.926	1935	1942	1945	rBV	71777	94330	0.36%	0.149%
76	12.963	1945	1948	1954	rVB	115121	137363	0.53%	0.217%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

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\SFAMXLM112221WMA.M
Title : VOC Analysis

77	13.030	1954	1959	1964	rBV4	6091	13258	0.05%	0.021%
78	13.140	1974	1977	1978	rBV	4831	4774	0.02%	0.008%
79	13.170	1978	1982	1986	rVB	26737	33793	0.13%	0.053%
80	13.256	1992	1996	1997	rBV	5274	5964	0.02%	0.009%
81	13.292	1997	2002	2007	rVV2	137791	196033	0.75%	0.310%
82	13.347	2007	2011	2019	rVB5	11087	18732	0.07%	0.030%
83	13.426	2019	2024	2033	rVB	41723	53361	0.20%	0.084%
84	13.512	2033	2038	2041	rBV4	5587	9056	0.03%	0.014%
85	13.585	2046	2050	2057	rVB5	7493	14470	0.06%	0.023%
86	13.719	2068	2072	2076	rBV2	13878	17893	0.07%	0.028%
87	13.780	2076	2082	2086	rVV	243376	305317	1.17%	0.483%
88	13.823	2086	2089	2099	rVB	16278	25378	0.10%	0.040%
89	13.932	2104	2107	2111	rVB3	2504	3315	0.01%	0.005%
90	13.969	2111	2113	2116	rBV3	1167	1205	0.00%	0.002%
91	14.005	2116	2119	2124	rVB2	2210	3323	0.01%	0.005%
92	14.079	2128	2131	2134	rBV2	1348	1814	0.01%	0.003%
93	14.115	2135	2137	2145	rVB3	2162	4251	0.02%	0.007%
94	14.231	2150	2156	2159	rVB4	3352	6015	0.02%	0.010%
95	14.316	2168	2170	2174	rVB2	1700	1819	0.01%	0.003%
96	14.377	2176	2180	2183	rVV2	1423	1554	0.01%	0.002%
97	14.481	2192	2197	2201	rBV2	2590	3335	0.01%	0.005%
98	14.640	2218	2223	2229	rBV	30480	36300	0.14%	0.057%
99	14.780	2242	2246	2253	rVB	20475	26592	0.10%	0.042%
100	15.011	2282	2284	2286	rBV	356	379	0.00%	0.001%

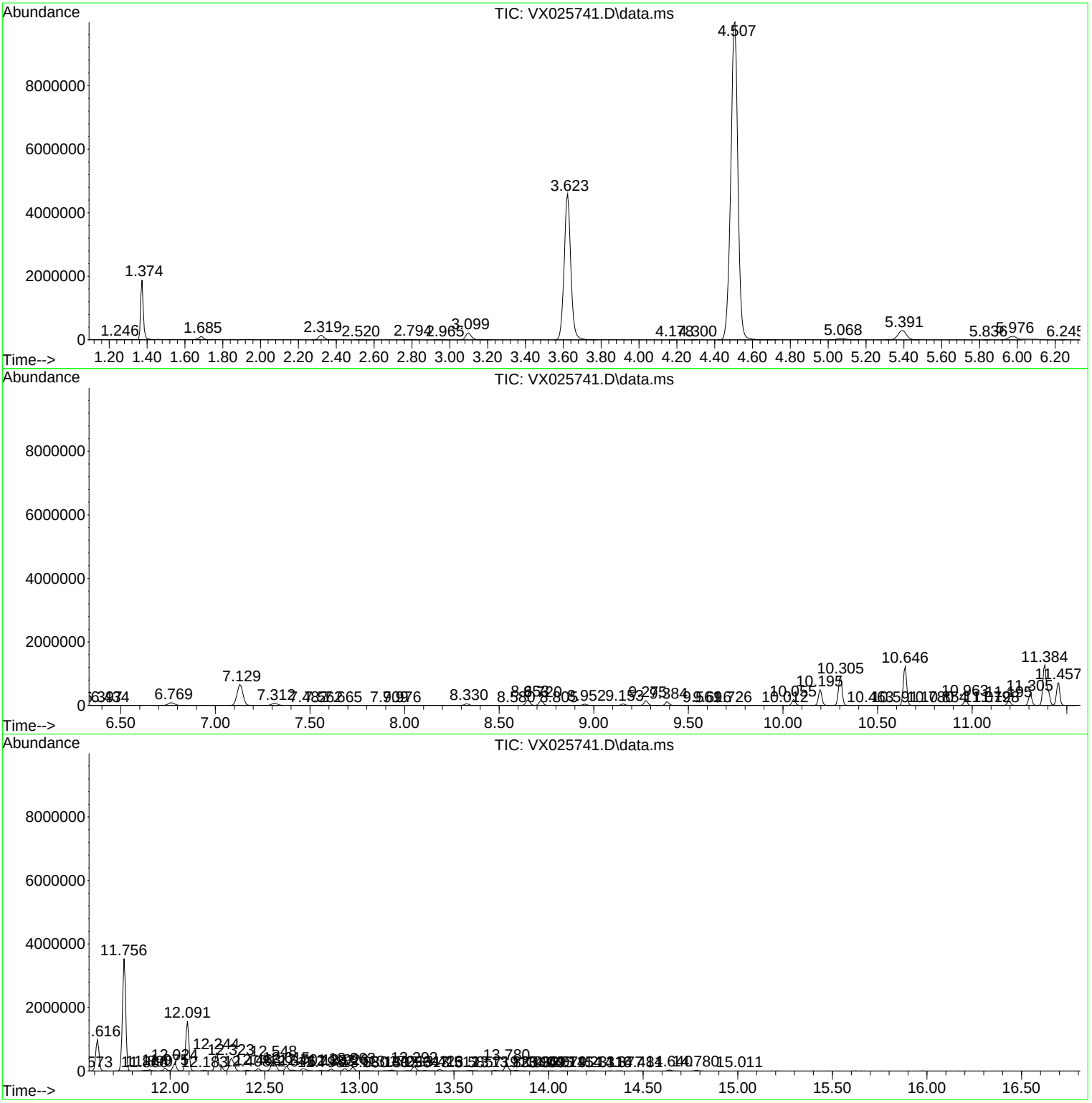
Sum of corrected areas: 63206102

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

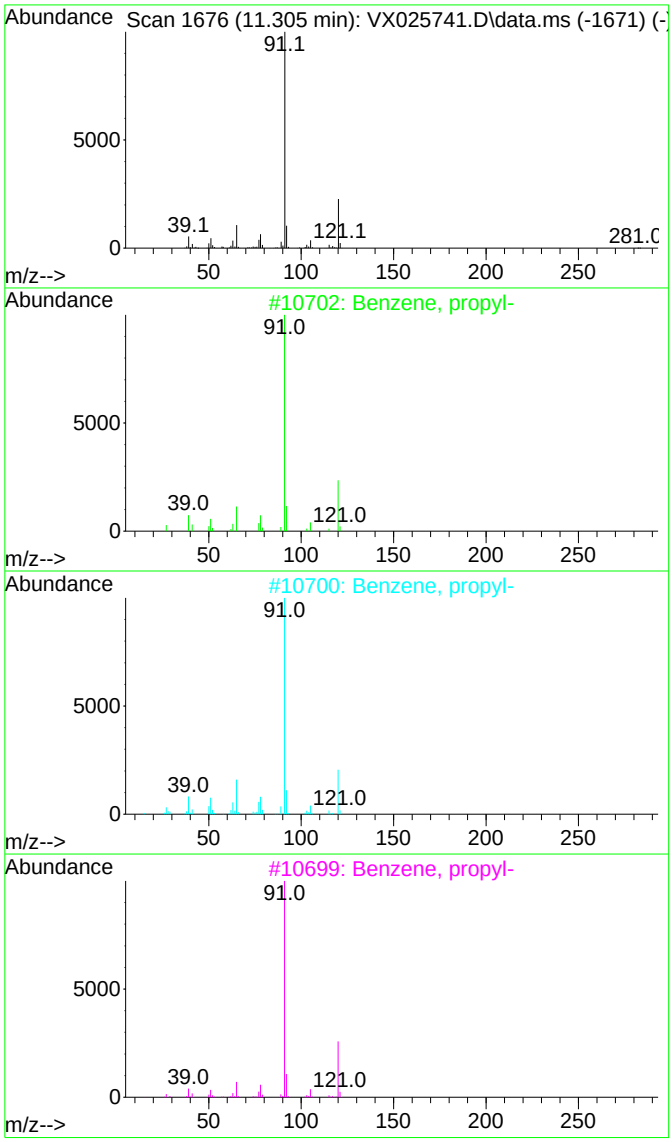
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

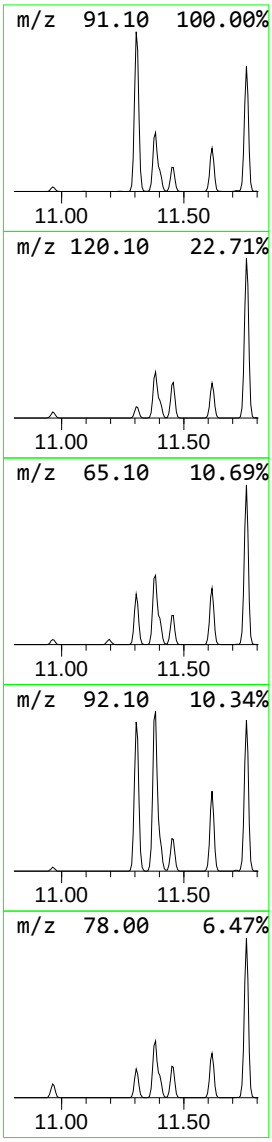
****

Peak Number 2 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.305	74.35 ug/L	447065	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	94
2	Benzene, propyl-	120	C9H12	000103-65-1	91
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	Benzene, propyl-	120	C9H12	000103-65-1	90
5	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

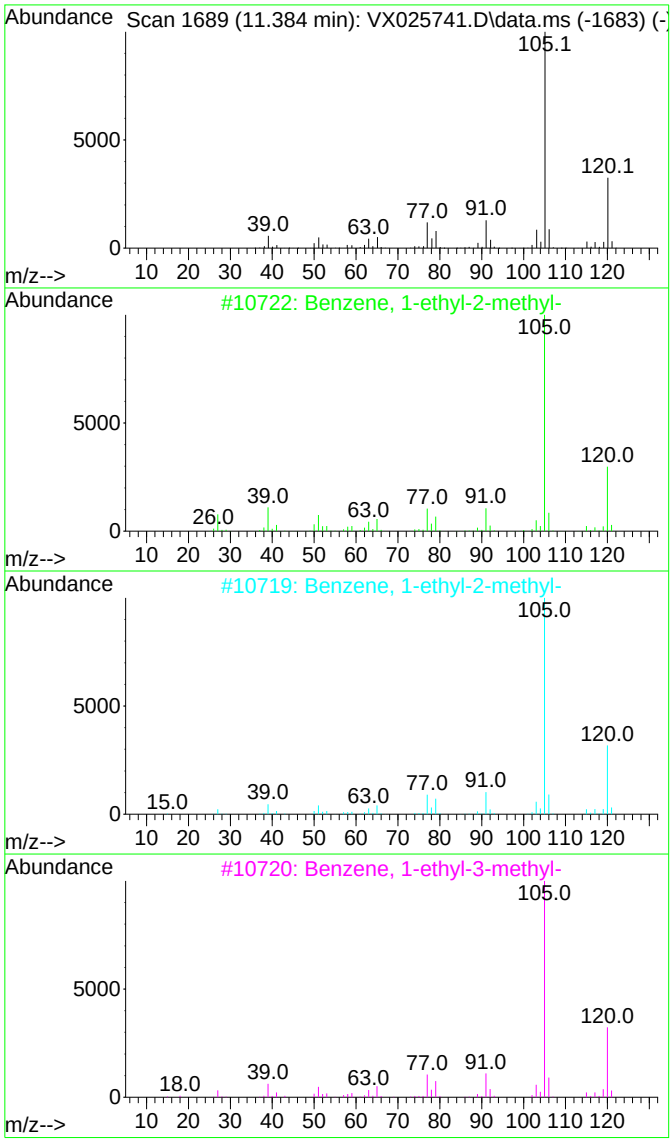
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

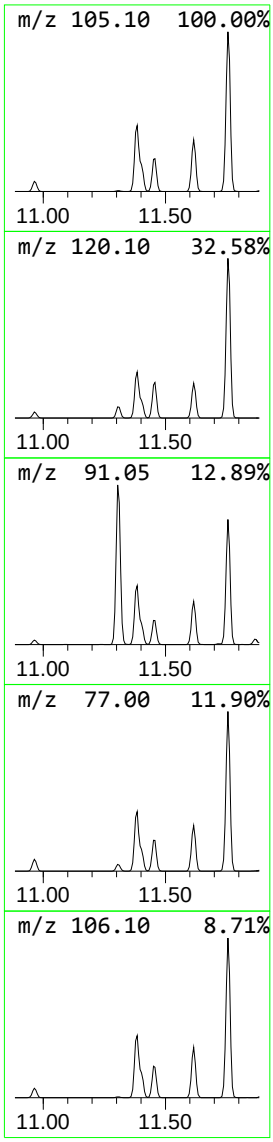
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	341.53 ug/L	2053480	1,4-Dichlorobenzene-d4	12.024

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	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

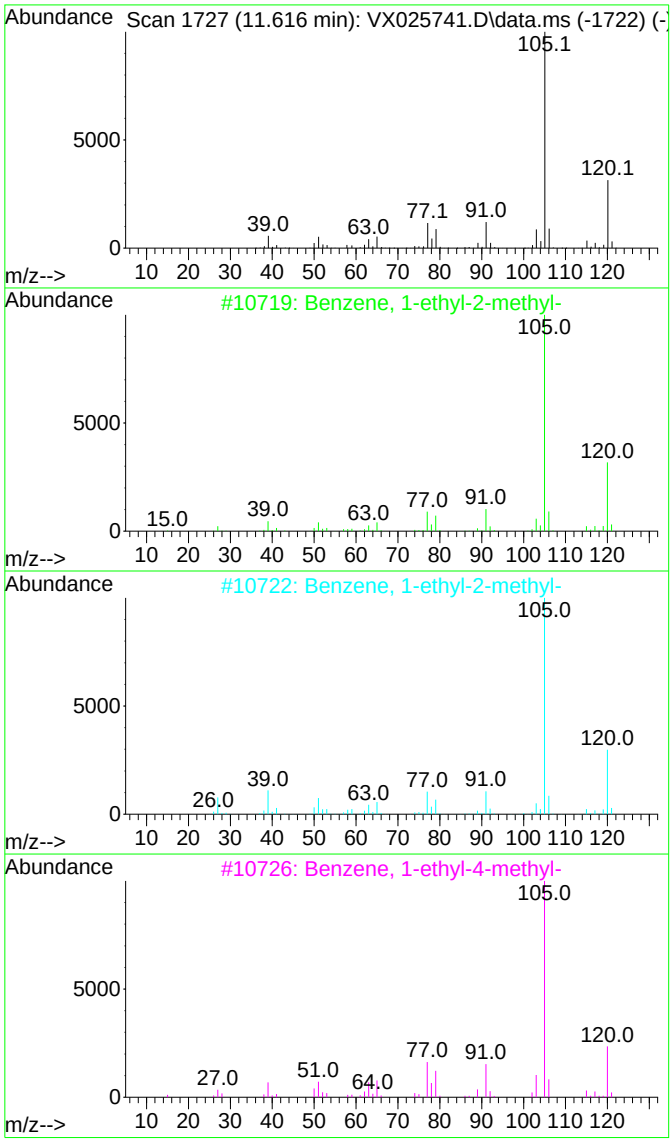
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

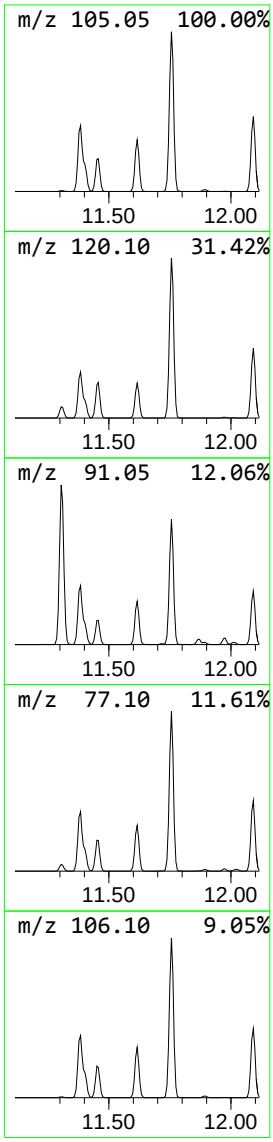
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	199.52 ug/L	1199660	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Mesitylene	120	C9H12	000108-67-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

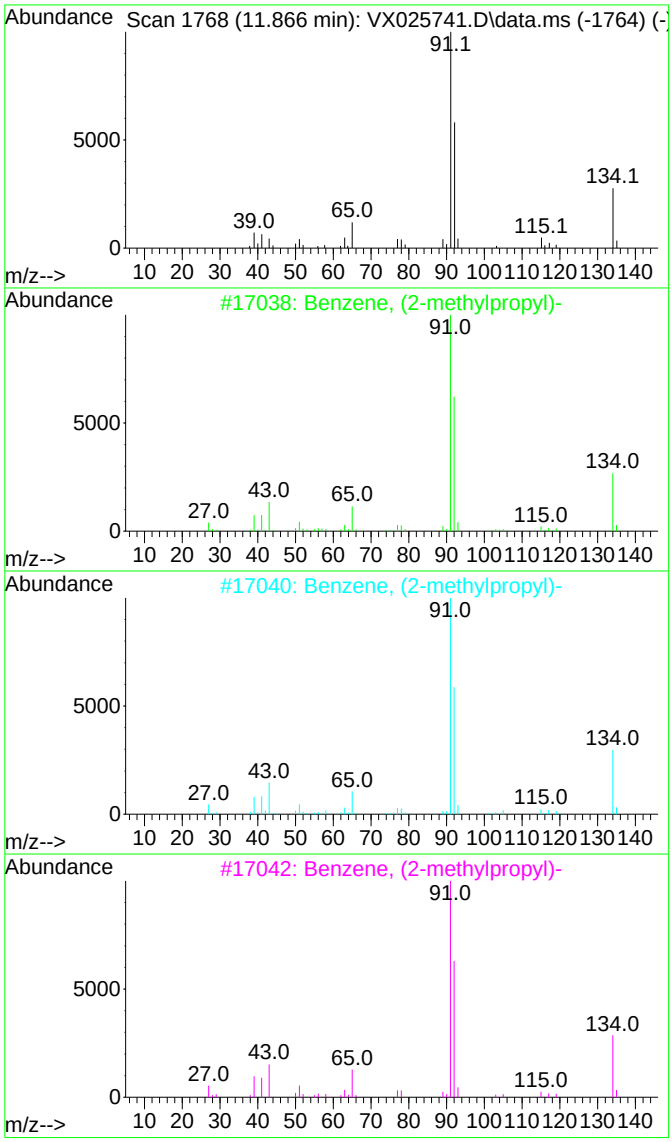
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

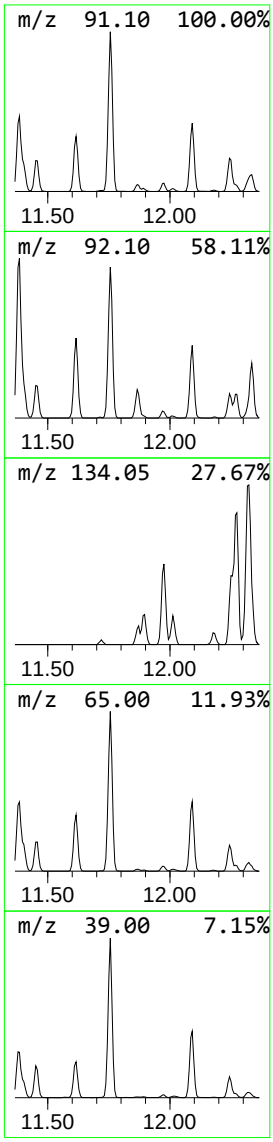
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, (2-methylpropyl)- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.866	3.19 ug/L	19192	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
4	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
5	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

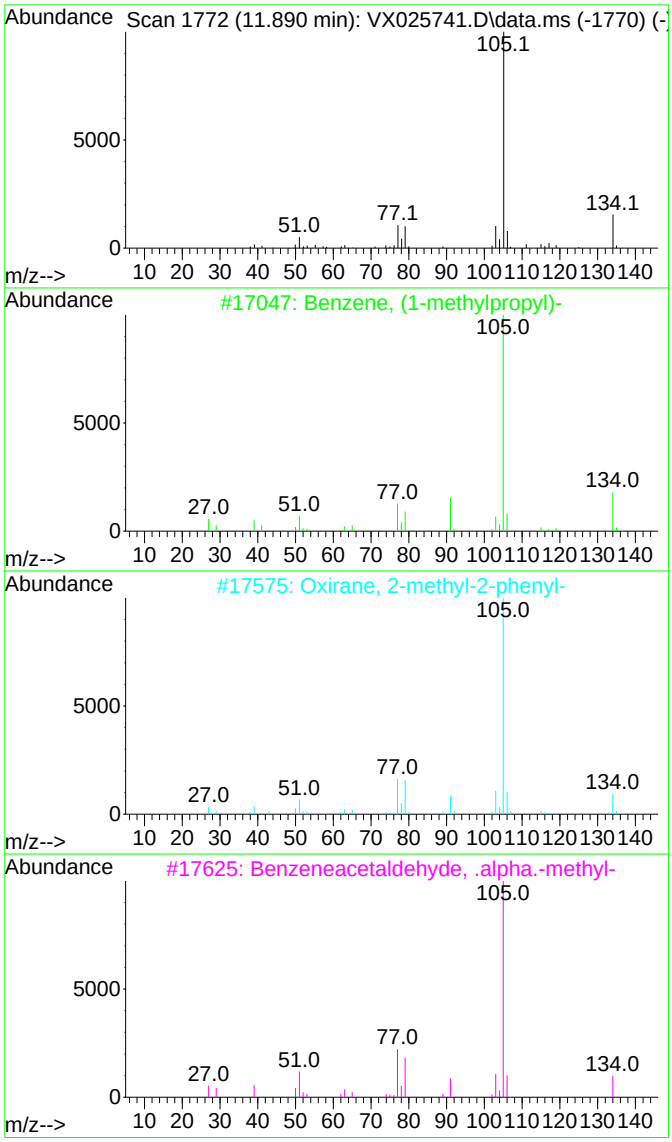
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

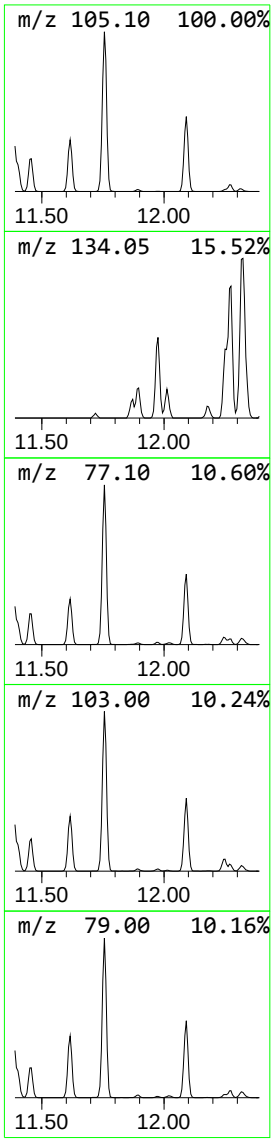
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (1-methylpropyl)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.890	6.66 ug/L	40063	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
2	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	87
3	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	87
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	83
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	78





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

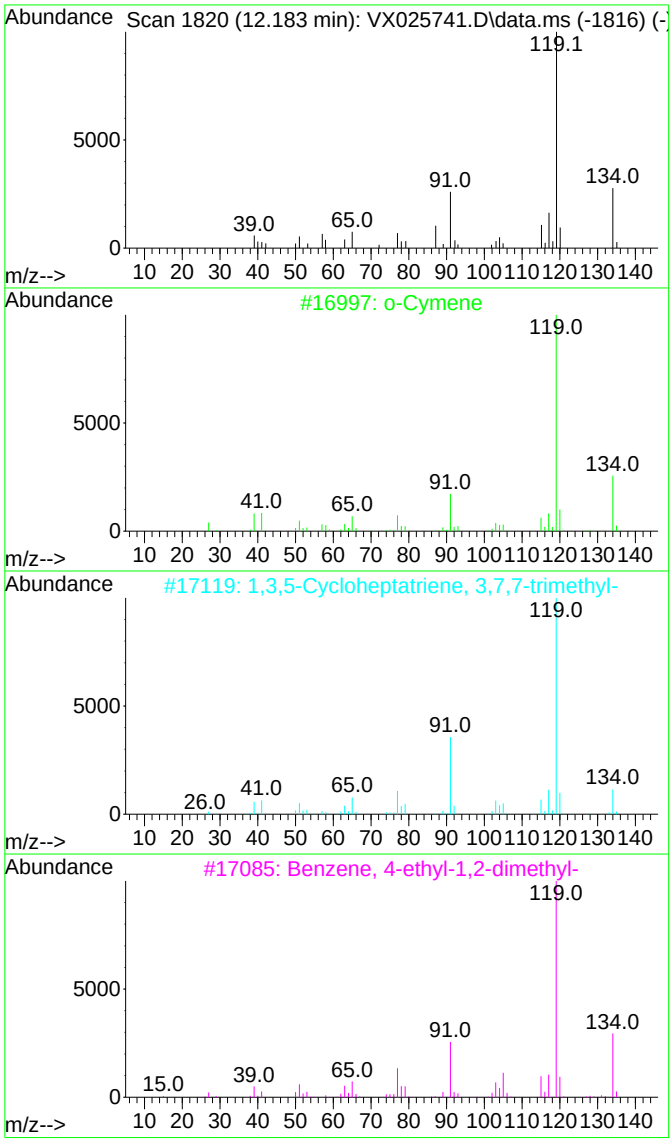
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

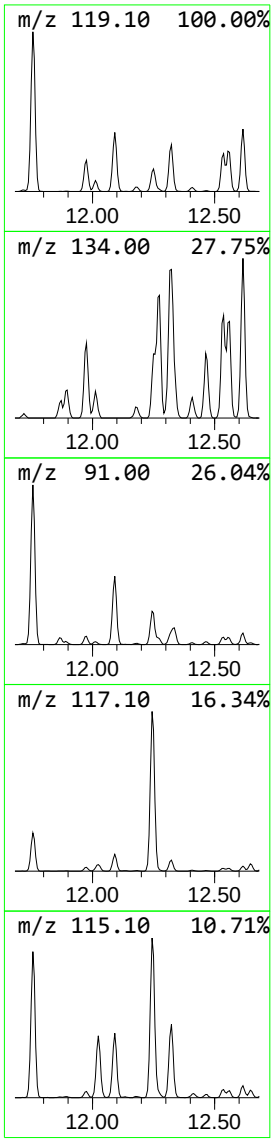
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 o-Cymene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.183	2.77 ug/L	16653	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	90
2	1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	90
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90
4	o-Cymene	134	C10H14	000527-84-4	90
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

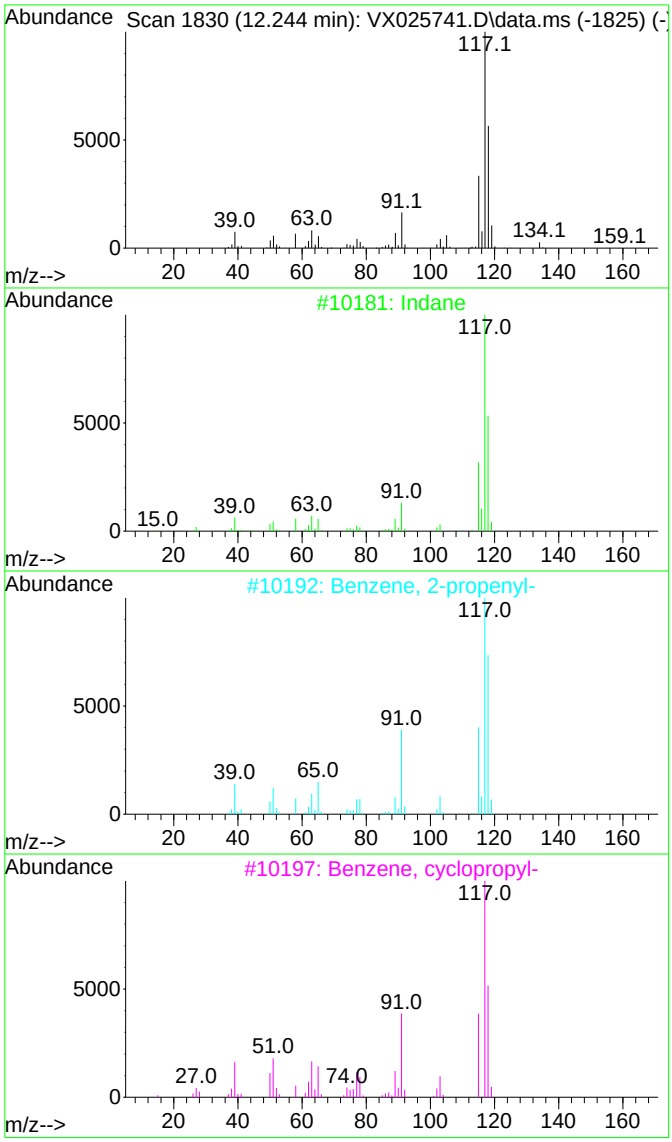
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

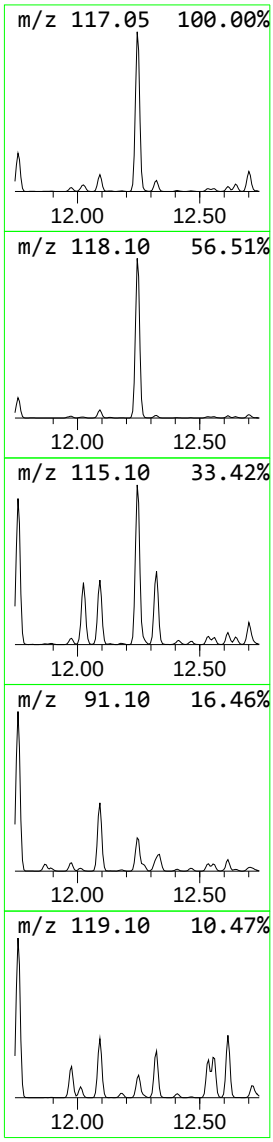
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	129.92 ug/L	781163	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	90
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4	Indane	118	C9H10	000496-11-7	81
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

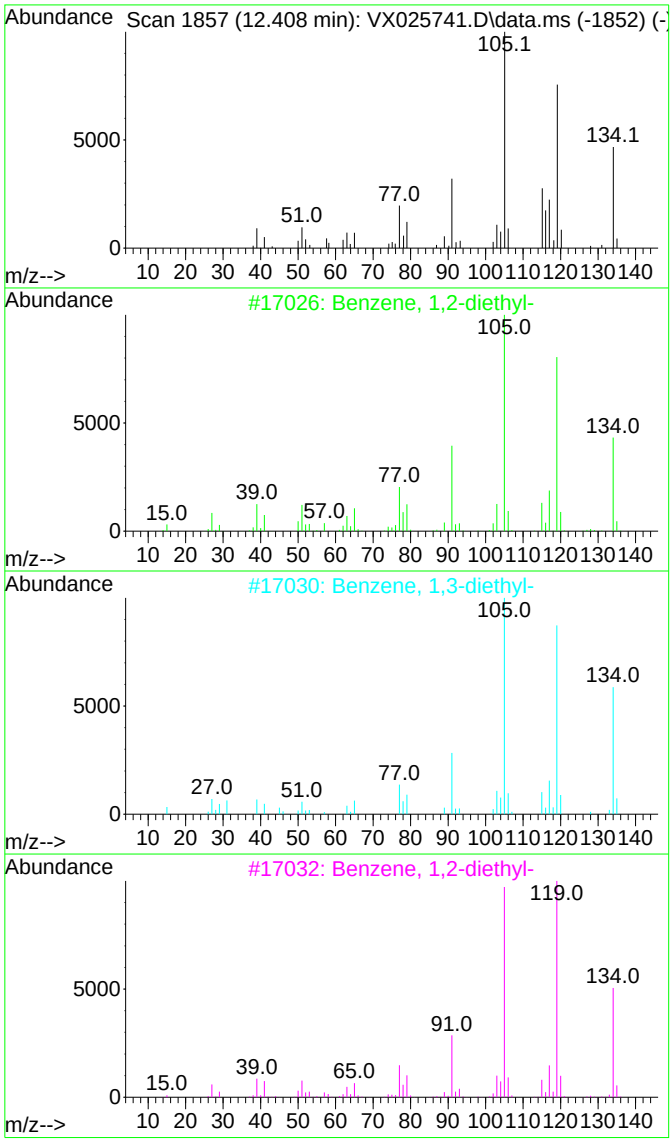
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

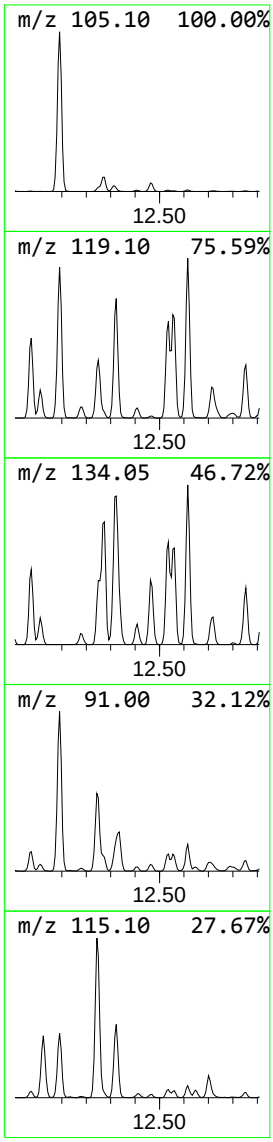
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1,2-diethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.408	5.04 ug/L	30309	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	81
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	81
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

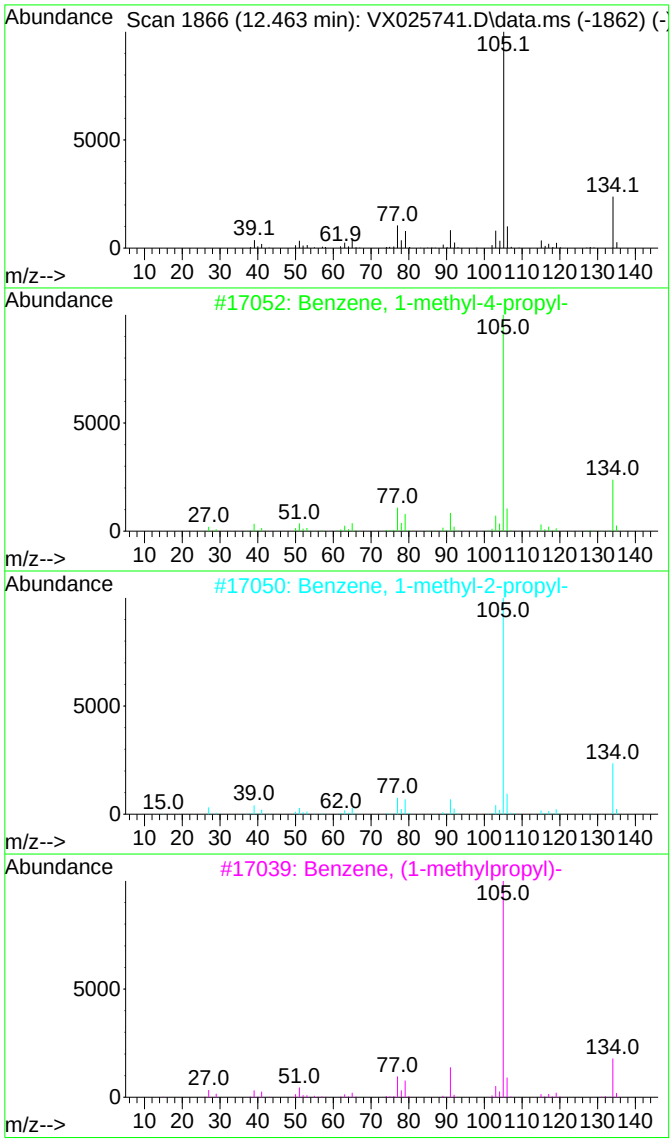
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

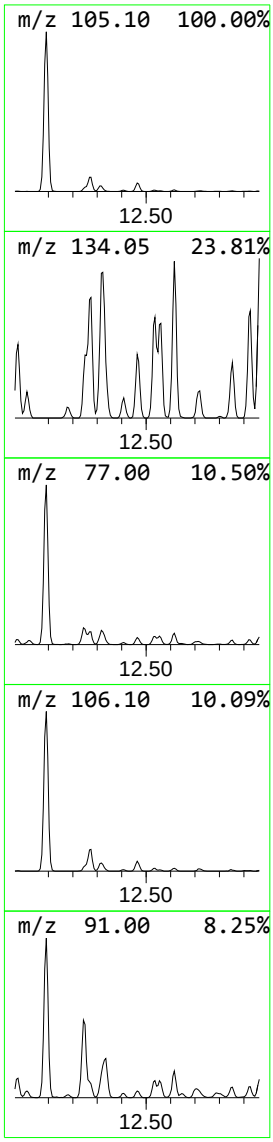
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-methyl-4-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.463	13.67 ug/L	82188	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

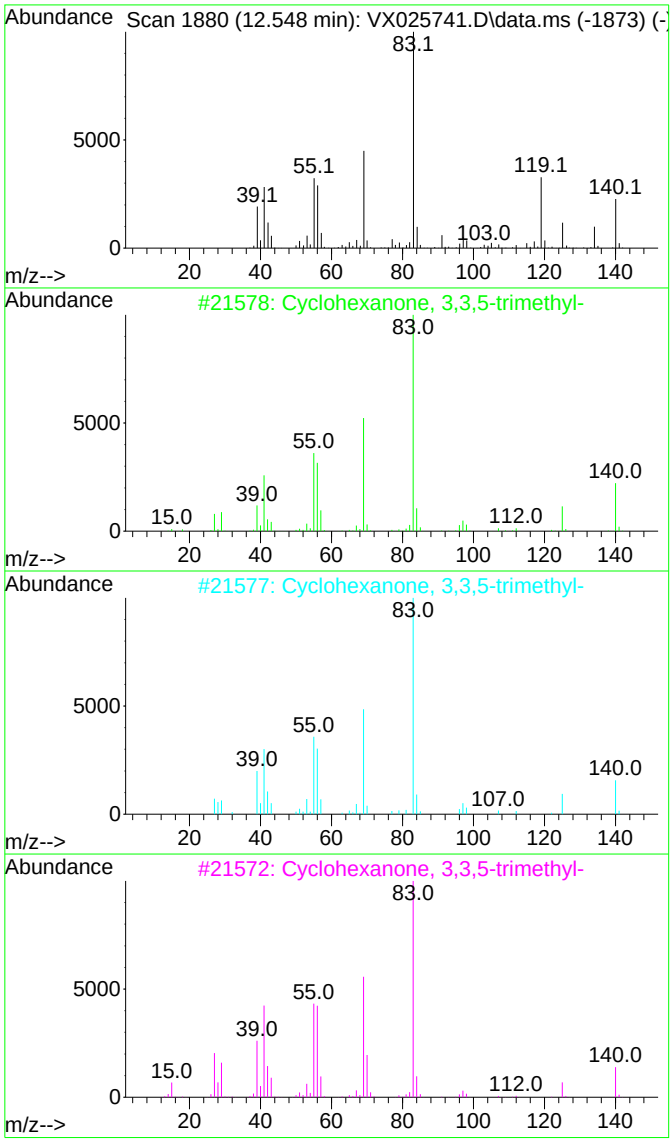
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

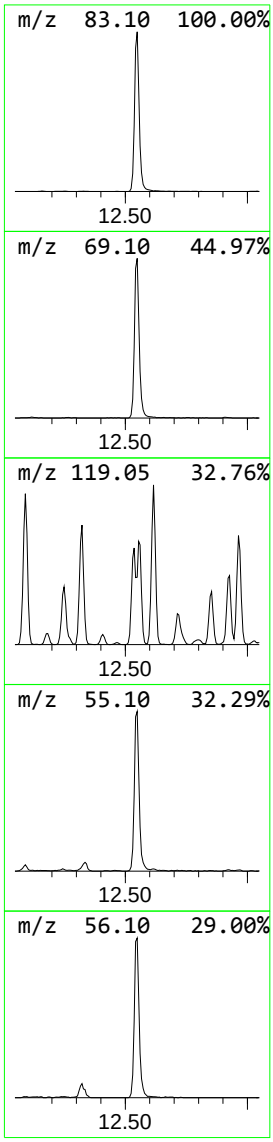
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Cyclohexanone, 3,3,5-trimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.548	101.06 ug/L	607656	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	81
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	81
5	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

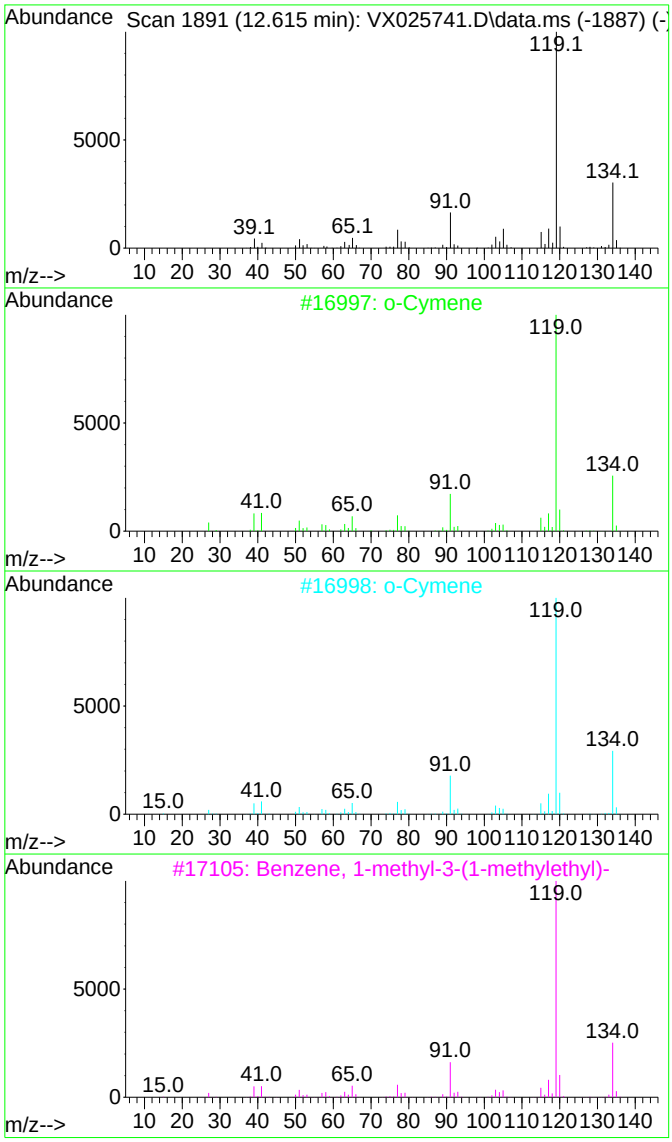
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

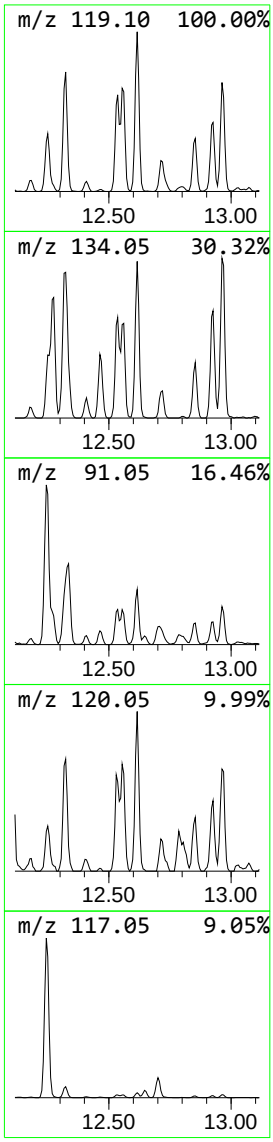
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1-methyl-3-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	29.75 ug/L	178862	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

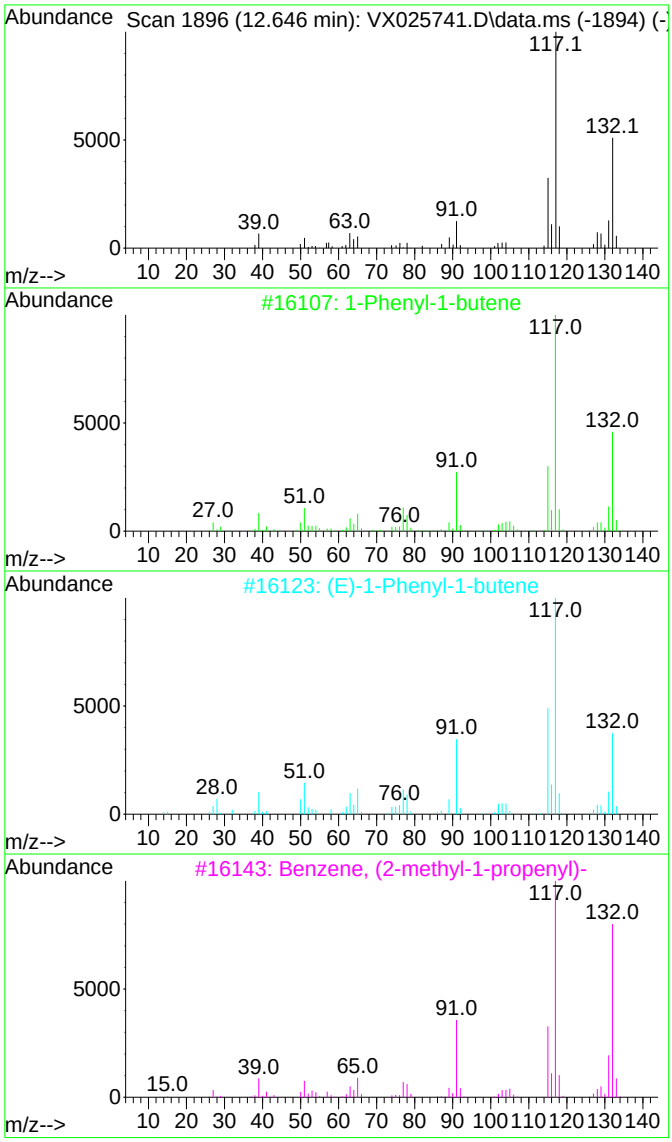
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

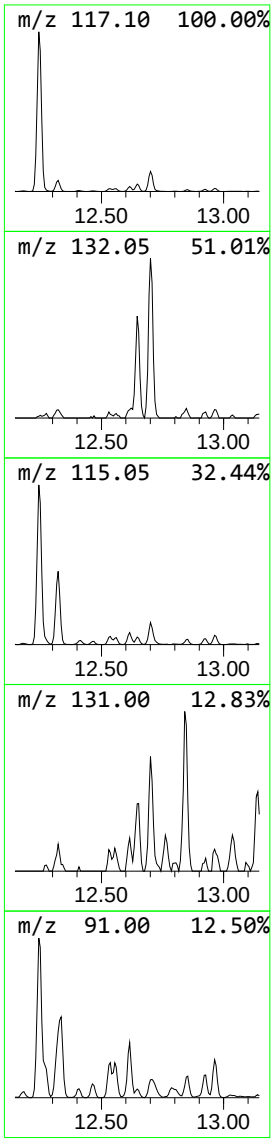
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1-Phenyl-1-butene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.646	5.16 ug/L	31011	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	94
2	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	91
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
4	Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	91
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

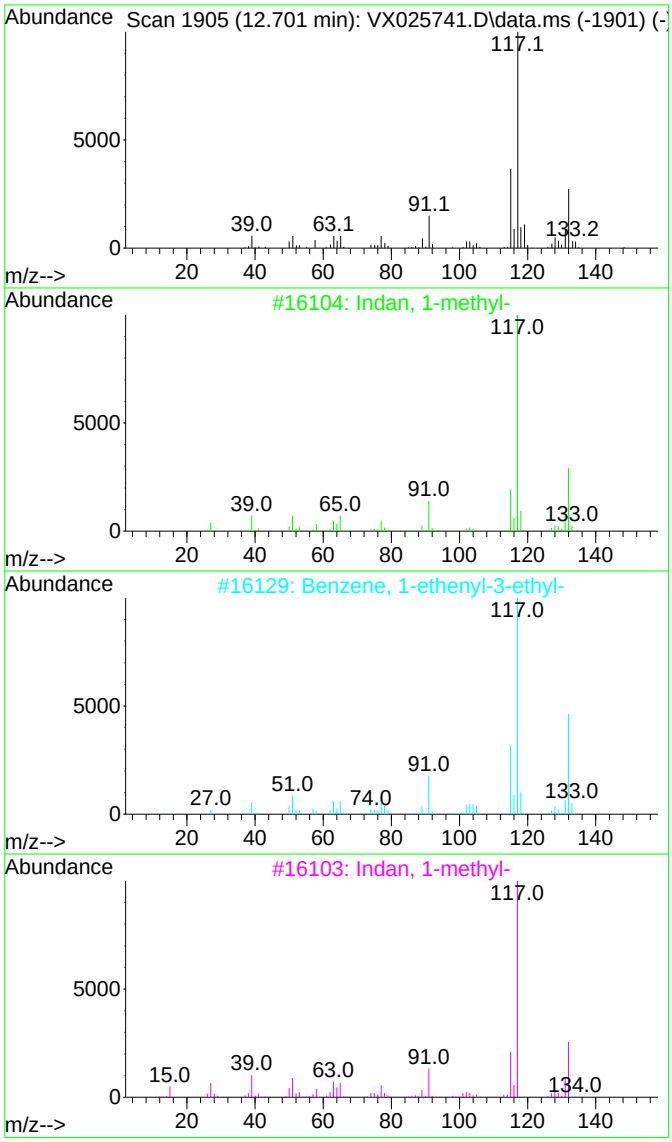
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

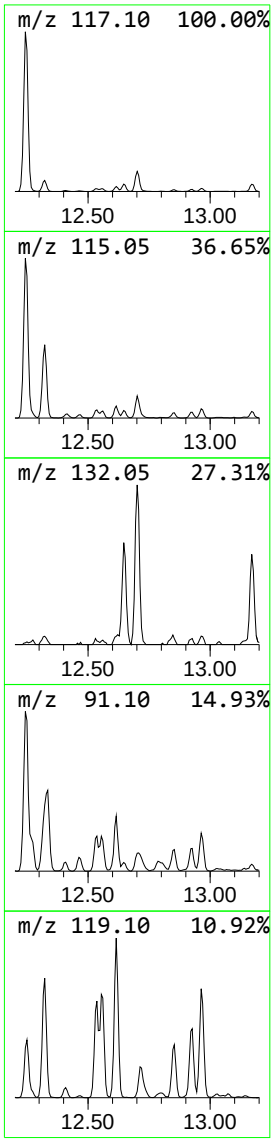
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Indan, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	20.61 ug/L	123948	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
3	Indan, 1-methyl-	132	C10H12	000767-58-8	83
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	81
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

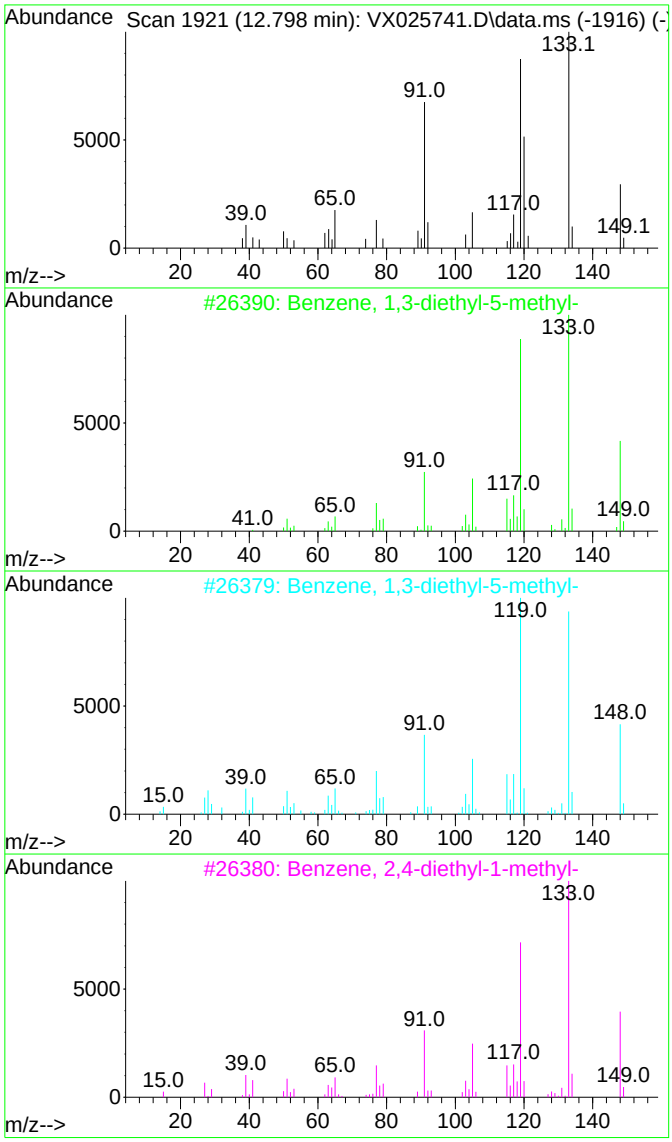
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

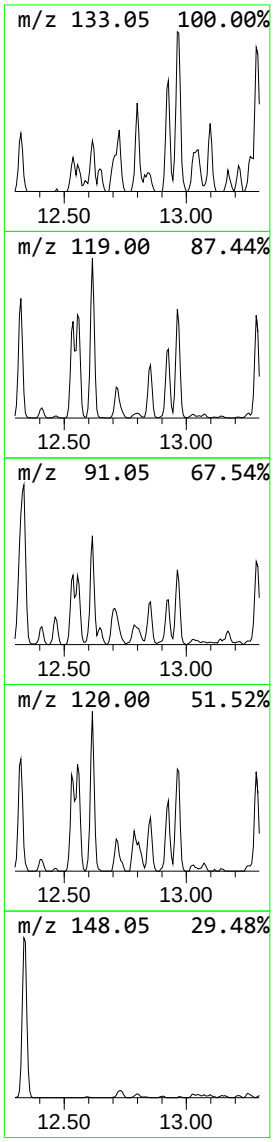
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.798	2.85 ug/L	17160	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	81
2	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	64
3	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	58
4	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	52
5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

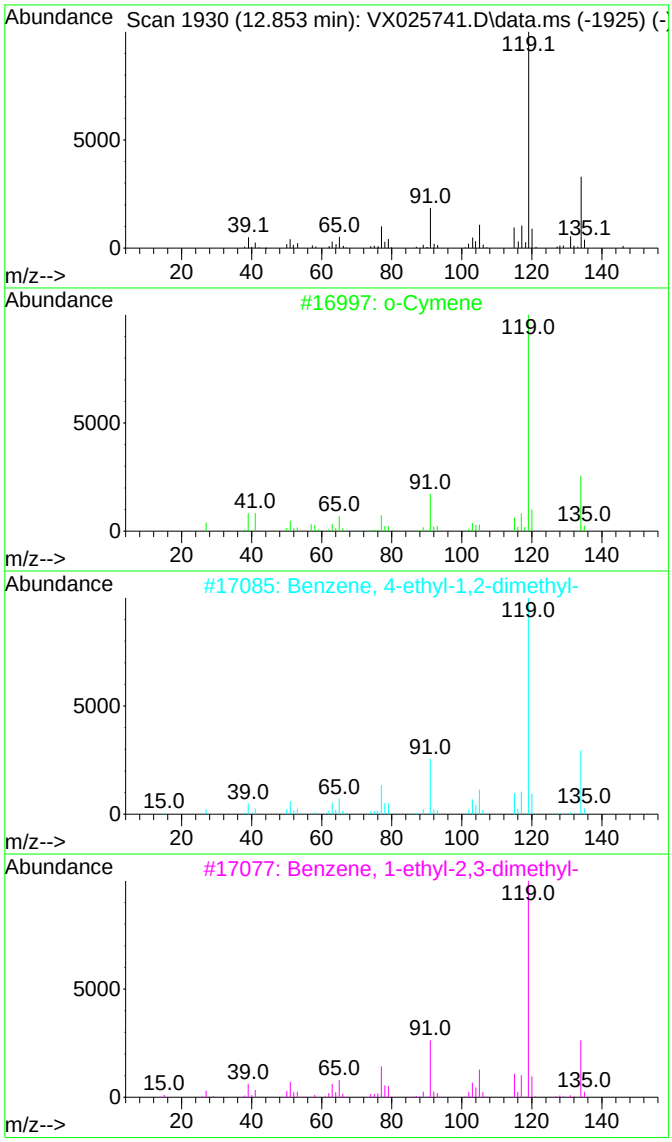
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

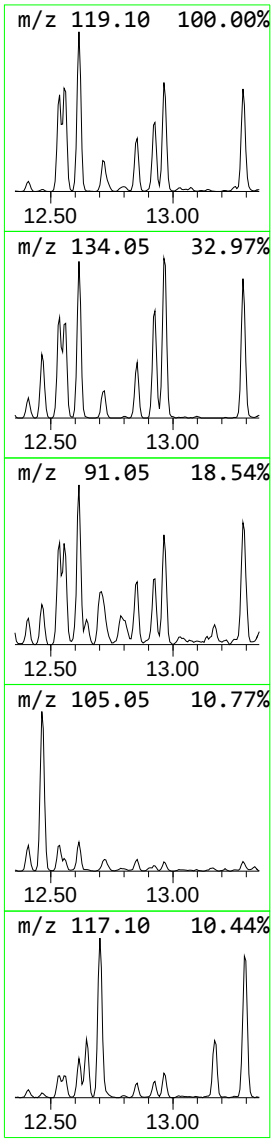
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.853	11.96 ug/L	71890	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	o-Cymene	134	C10H14	000527-84-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

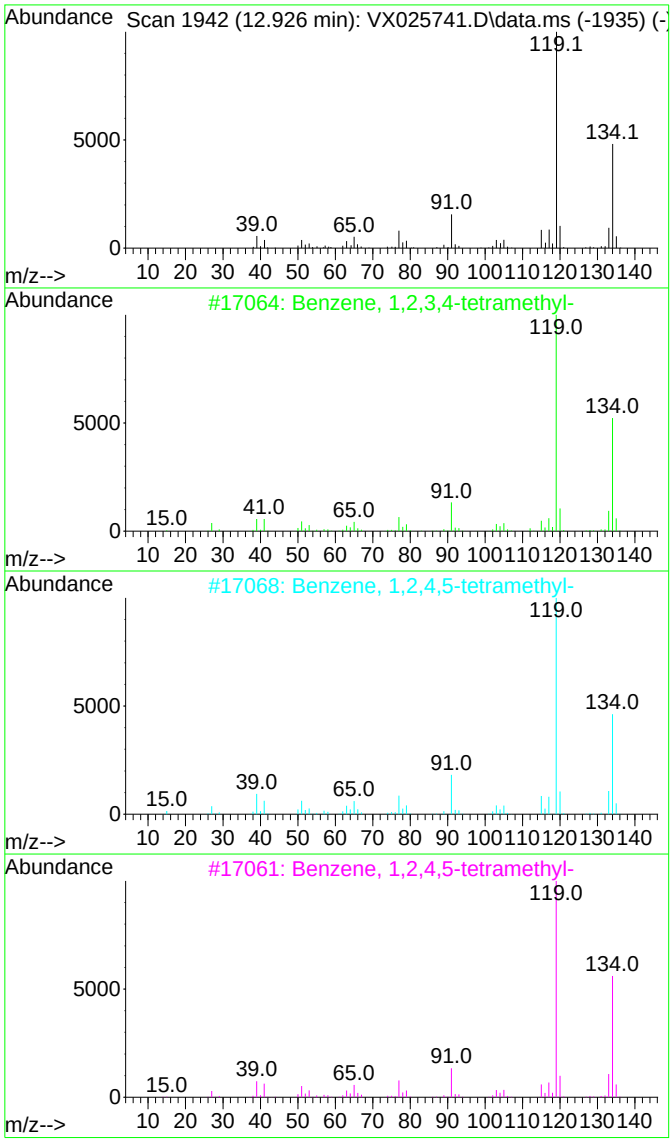
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

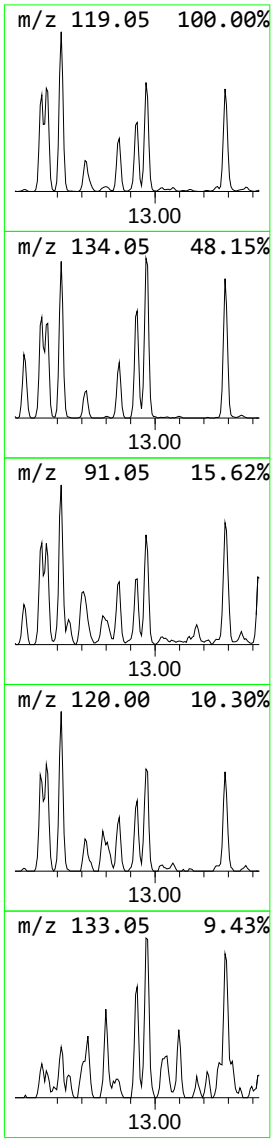
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.926	15.69 ug/L	94330	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

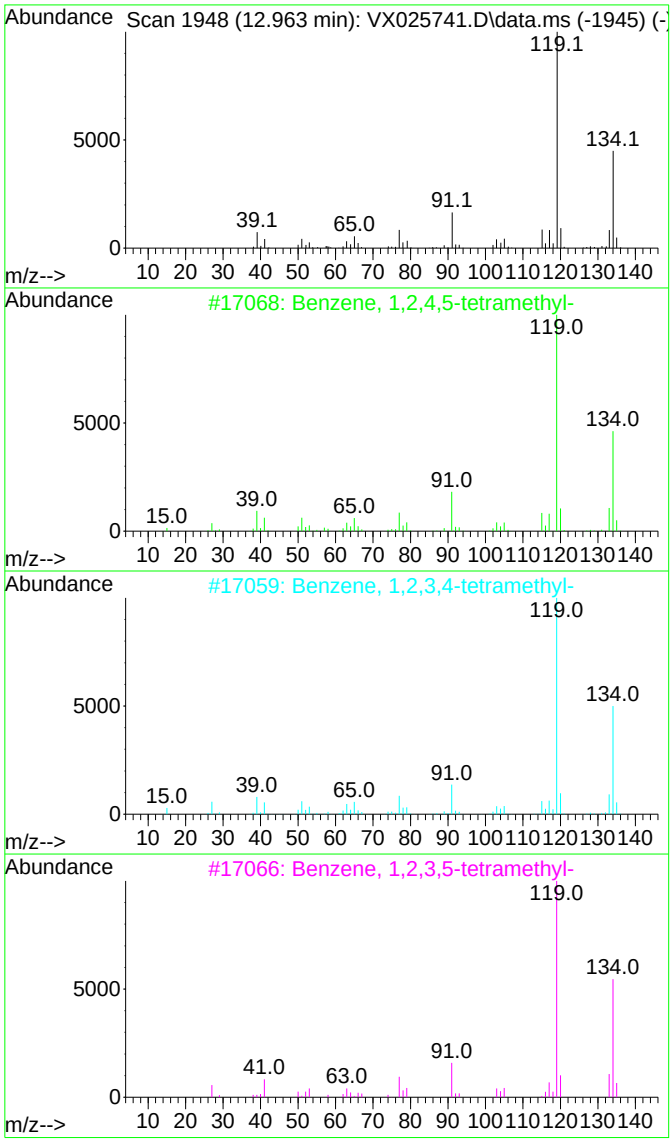
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

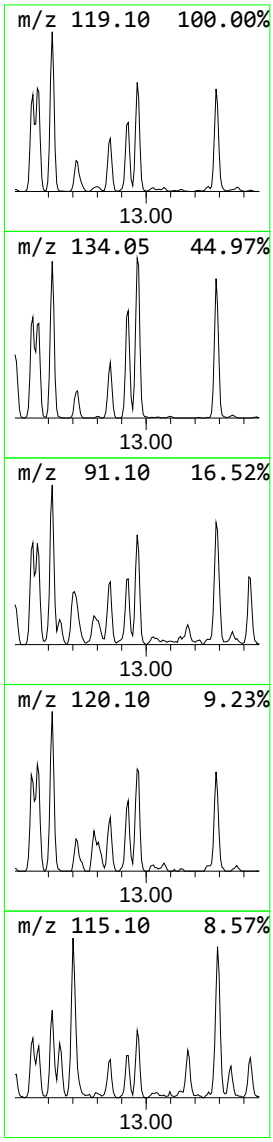
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	22.85 ug/L	137363	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

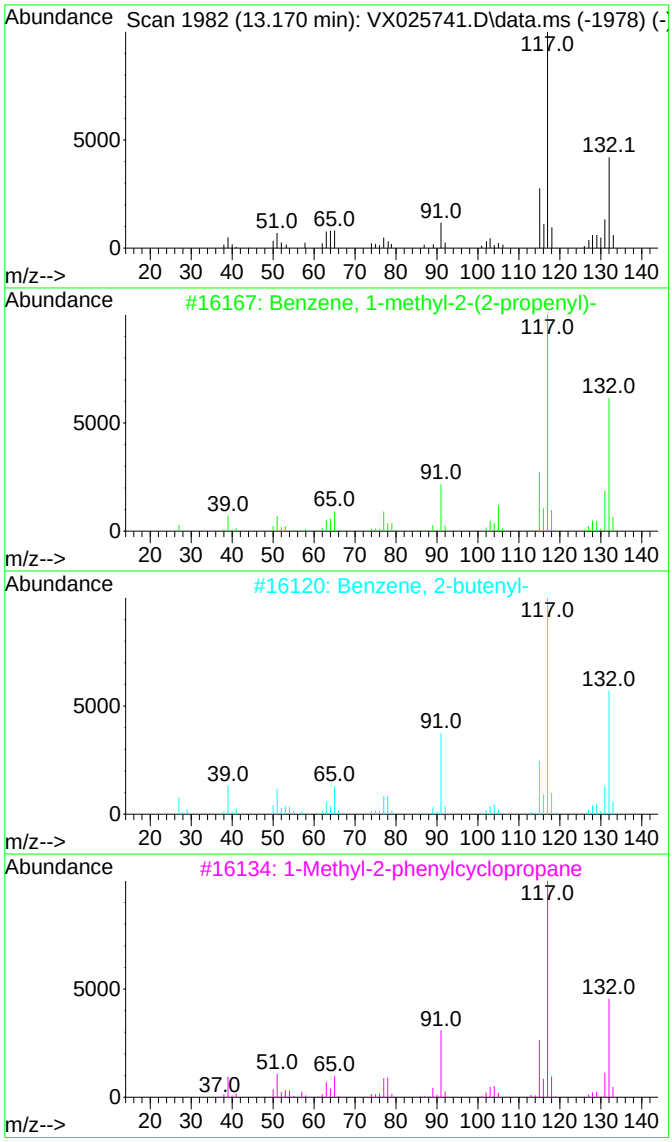
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

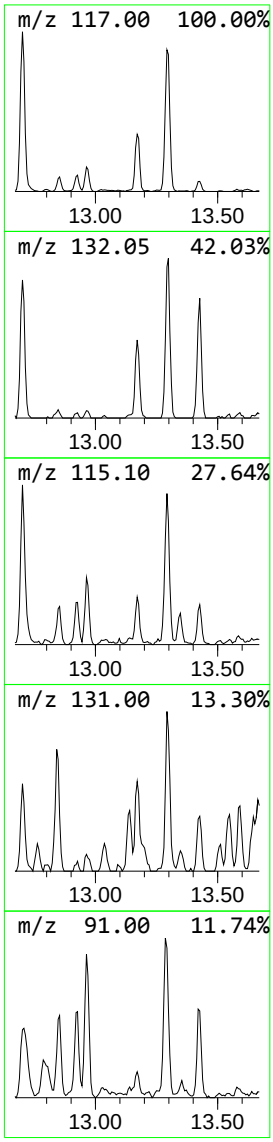
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1-methyl-2-(2-prop... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	5.62 ug/L	33793	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	94
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

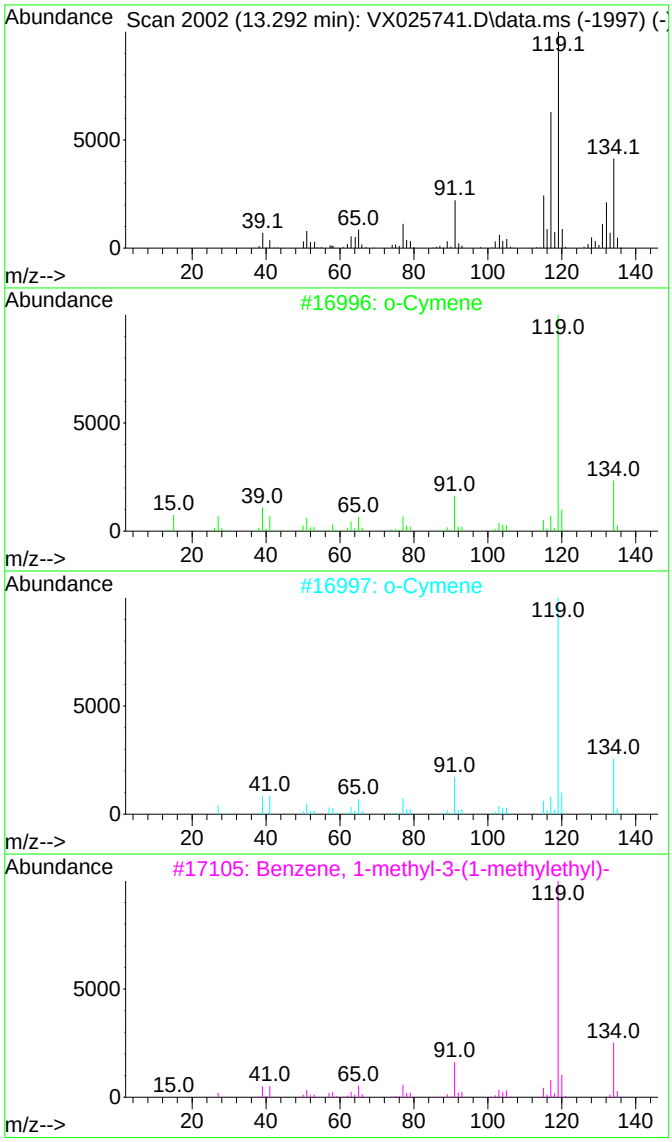
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

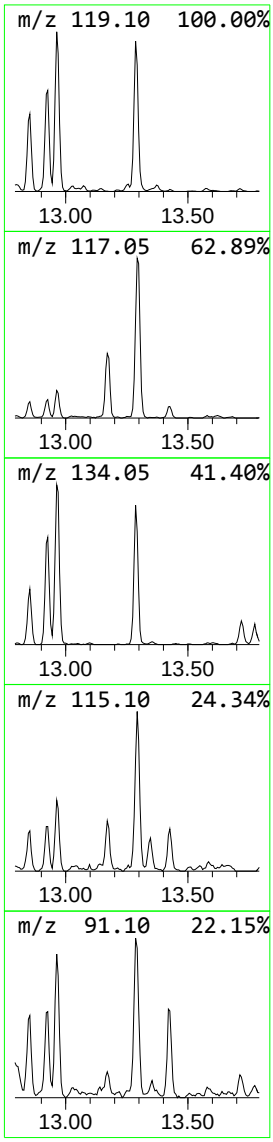
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	32.60 ug/L	196033	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	60
2	o-Cymene	134	C10H14	000527-84-4	60
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	60
5	p-Cymene	134	C10H14	000099-87-6	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

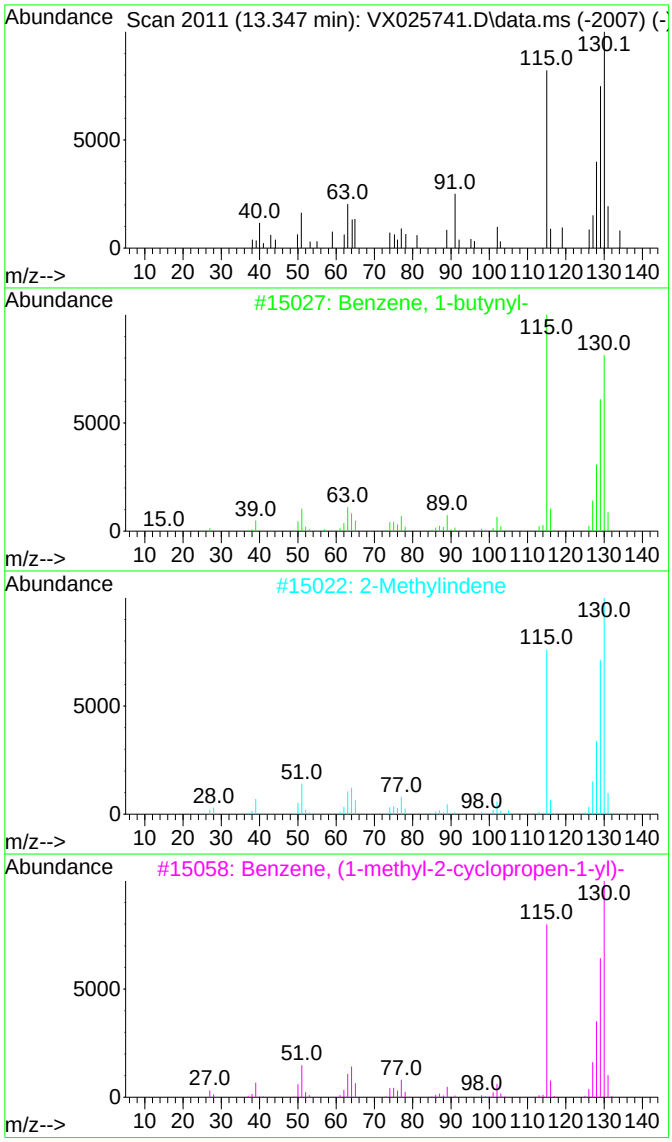
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

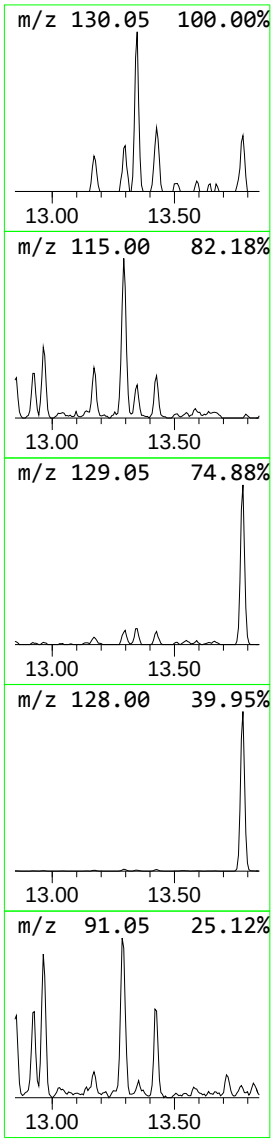
****

Peak Number 22 Benzene, 1-butynyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.347	3.12 ug/L	18732	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-butynyl-	130	C10H10	000622-76-4	81
2	2-Methylindene	130	C10H10	002177-47-1	81
3	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	81
4	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	81
5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

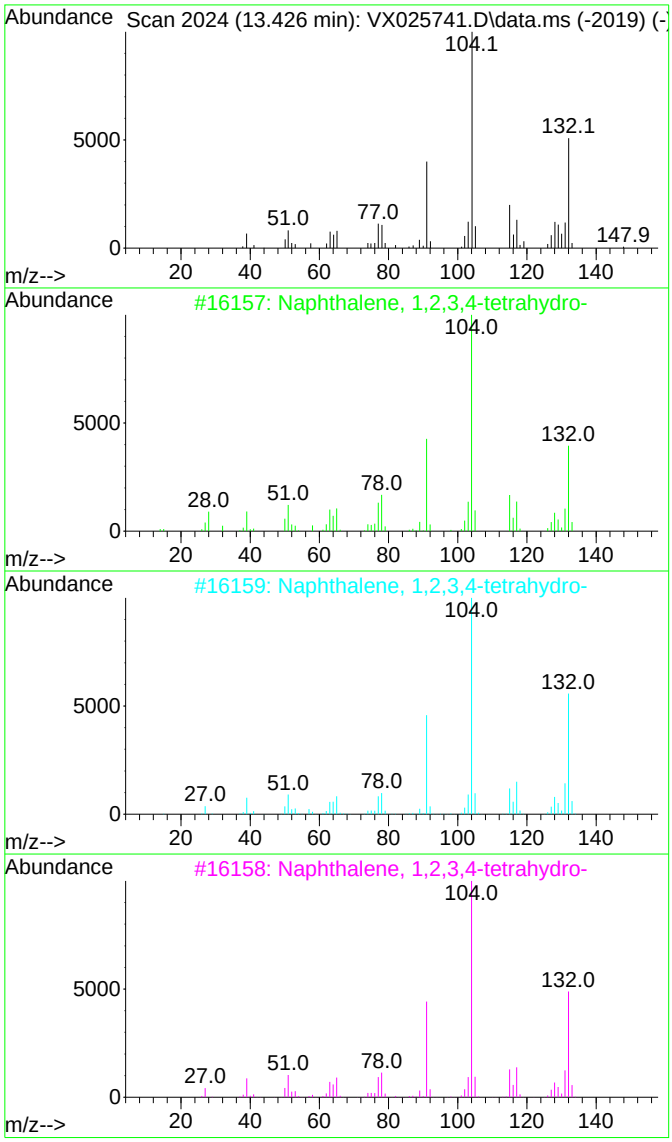
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

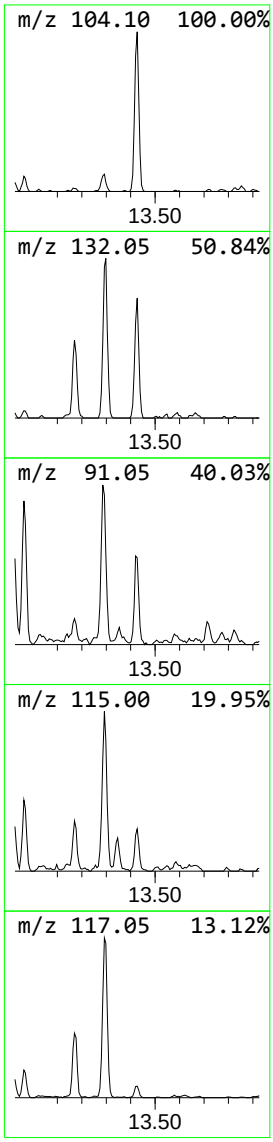
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.426	8.87 ug/L	53361	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
3	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
4	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

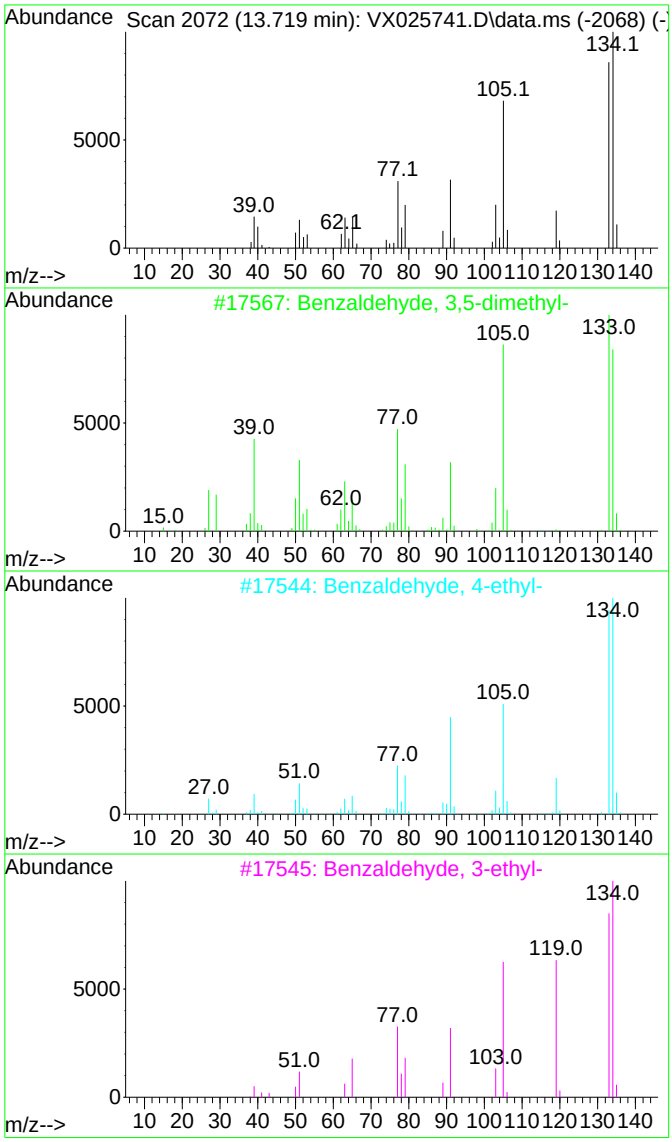
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

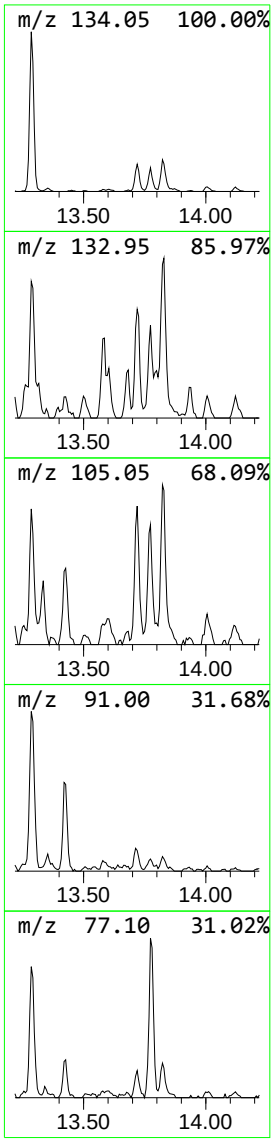
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzaldehyde, 3,5-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.719	2.98 ug/L	17893	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dimethyl-	134	C9H10O	005779-95-3	95
2	Benzaldehyde, 4-ethyl-	134	C9H10O	004748-78-1	94
3	Benzaldehyde, 3-ethyl-	134	C9H10O	034246-54-3	93
4	Benzaldehyde, 2,4-dimethyl-	134	C9H10O	015764-16-6	90
5	Benzaldehyde, 2,5-dimethyl-	134	C9H10O	005779-94-2	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

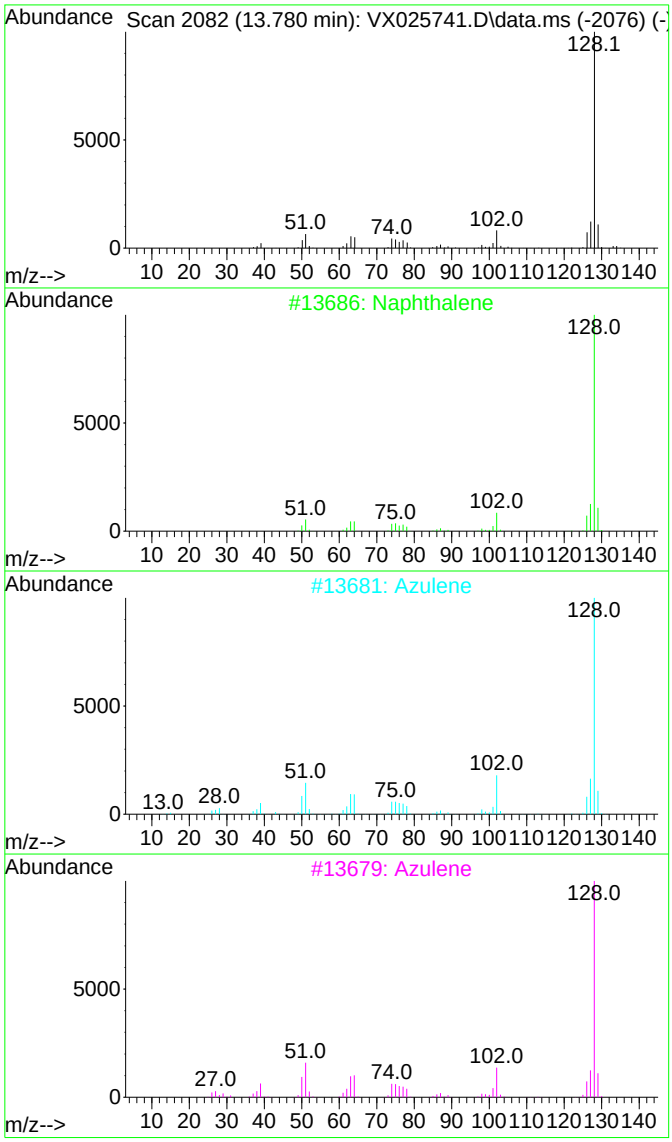
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

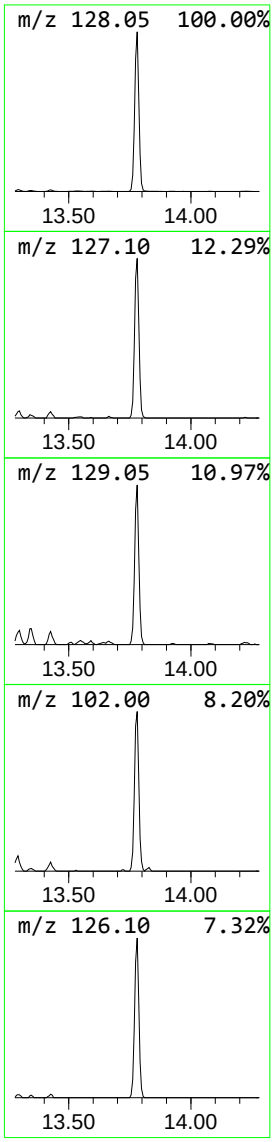
****

Peak Number 25 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	50.78 ug/L	305317	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	91
3	Azulene	128	C10H8	000275-51-4	90
4	Azulene	128	C10H8	000275-51-4	90
5	Naphthalene	128	C10H8	000091-20-3	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

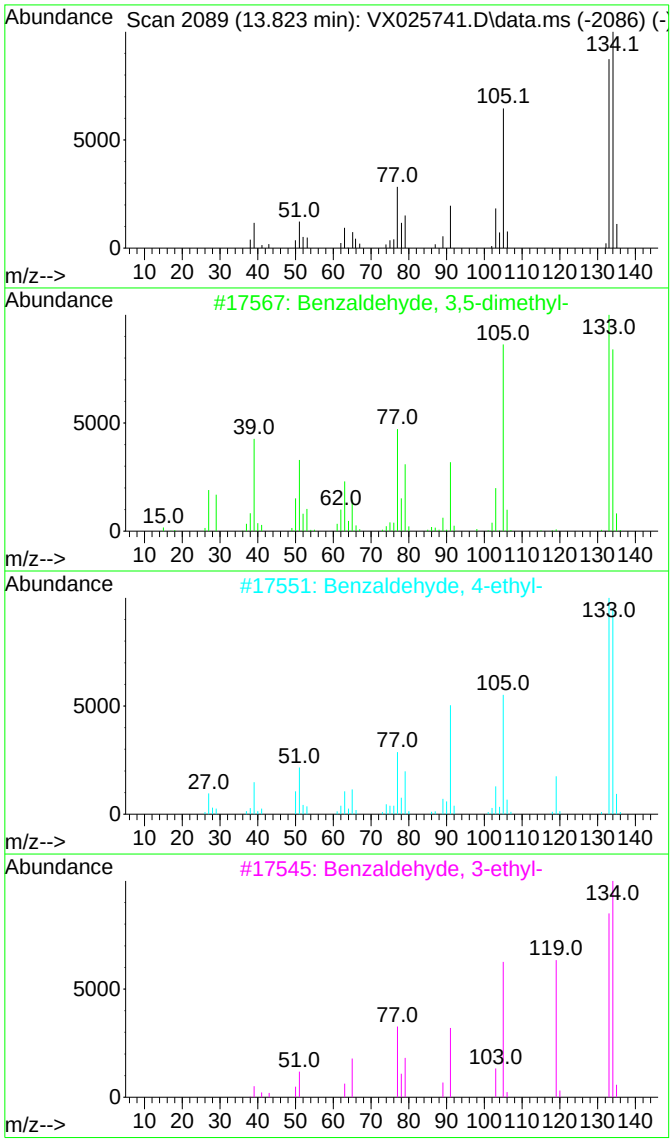
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.M
Quant Title : VOC Analysis

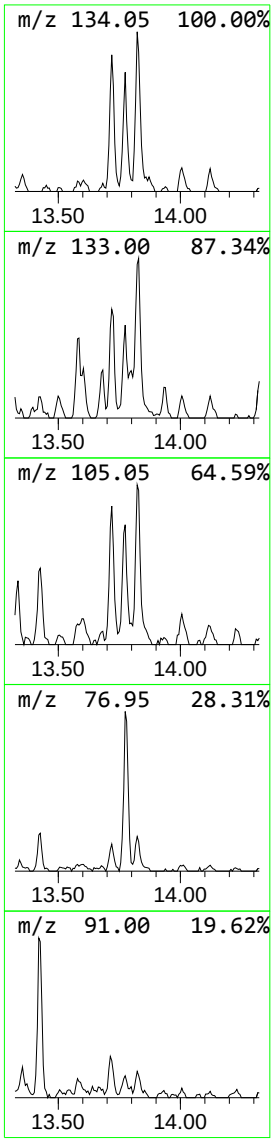
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Benzaldehyde, 4-ethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.823	4.22 ug/L	25378	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dimethyl-	134	C9H10O	005779-95-3	94
2	Benzaldehyde, 4-ethyl-	134	C9H10O	004748-78-1	91
3	Benzaldehyde, 3-ethyl-	134	C9H10O	034246-54-3	91
4	Benzaldehyde, 2,4-dimethyl-	134	C9H10O	015764-16-6	91
5	Benzaldehyde, 4-ethyl-	134	C9H10O	004748-78-1	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121121\
Data File : VX025741.D
Acq On : 11 Dec 2021 14:25
Operator : JC/MD
Sample : M4997-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW5R1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM112221WMA.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, propyl-	11.305	74.3	ug/L	447065	3	12.024	300633	50.0
Benzene, 1-ethy...	11.384	341.5	ug/L	2053480	3	12.024	300633	50.0
Benzene, 1-ethy...	11.616	199.5	ug/L	1199660	3	12.024	300633	50.0
Benzene, (2-met...	11.866	3.2	ug/L	19192	3	12.024	300633	50.0
Benzene, (1-met...	11.890	6.7	ug/L	40063	3	12.024	300633	50.0
o-Cymene	12.183	2.8	ug/L	16653	3	12.024	300633	50.0
Indane	12.244	129.9	ug/L	781163	3	12.024	300633	50.0
Benzene, 1,2-di...	12.408	5.0	ug/L	30309	3	12.024	300633	50.0
Benzene, 1-meth...	12.463	13.7	ug/L	82188	3	12.024	300633	50.0
Cyclohexanone, ...	12.548	101.1	ug/L	607656	3	12.024	300633	50.0
Benzene, 1-meth...	12.616	29.8	ug/L	178862	3	12.024	300633	50.0
1-Phenyl-1-butene	12.646	5.2	ug/L	31011	3	12.024	300633	50.0
Indan, 1-methyl-	12.701	20.6	ug/L	123948	3	12.024	300633	50.0
Benzene, 1,3-di...	12.798	2.9	ug/L	17160	3	12.024	300633	50.0
Benzene, 4-ethy...	12.853	12.0	ug/L	71890	3	12.024	300633	50.0
Benzene, 1,2,3,...	12.926	15.7	ug/L	94330	3	12.024	300633	50.0
Benzene, 1,2,4,...	12.963	22.9	ug/L	137363	3	12.024	300633	50.0
Benzene, 1-meth...	13.170	5.6	ug/L	33793	3	12.024	300633	50.0
Benzene, 1,2,3,...	13.292	32.6	ug/L	196033	3	12.024	300633	50.0
Benzene, 1-buty...	13.347	3.1	ug/L	18732	3	12.024	300633	50.0
Naphthalene, 1,...	13.426	8.9	ug/L	53361	3	12.024	300633	50.0
Benzaldehyde, 3...	13.719	3.0	ug/L	17893	3	12.024	300633	50.0
Azulene	13.780	50.8	ug/L	305317	3	12.024	300633	50.0
Benzaldehyde, 4...	13.823	4.2	ug/L	25378	3	12.024	300633	50.0