LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\82X071619W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.276	14	20	23	rBV	13515	20171	1.92%	0.167%
2	1.318	23	27	36	rVV2	12695	37399	3.56%	0.309%
3	2.397	197	204	208	rBV4	5295	10509	1.00%	0.087%
4	2.611	230	239	245	rBV	10958	21686	2.06%	0.179%
5	2.848	272	278	284	rBV6	4922	12572	1.20%	0.104%
6	3.172	323	331	338	rBV2	14621	32715	3.11%	0.270%
7	5.500	700	713	724	rBV	125654	357359	34.01%	2.953%
8	5.659	729	739	760	rVB	210226	611975	58.23%	5.057%
9	6.061	795	805	821	rBV	123961	325249	30.95%	2.688%
10	6.860	926	936	946	rBV	315572	747271	71.11%	6.175%
11	7.841	1088	1097	1106	rBV3	5559	13792	1.31%	0.114%
12	8.384	1178	1186	1192	rBV5	6864	14376	1.37%	0.119%
13	8.713	1228	1240	1248	rBV	684087	1050884	100.00%	8.684%
14	8.841	1256	1261	1265	rBV4	7474	11182	1.06%	0.092%
15	9.036	1288	1293	1300	rVB	21292	37073	3.53%	0.306%
16	9.164	1306	1314	1319	rBV2	8624	15935	1.52%	0.132%
17	9.445	1356	1360	1367	rVB3	8240	14871	1.42%	0.123%
18	9.567	1372	1380	1387	rBV4	13332	31028	2.95%	0.256%
19	9.676	1392	1398	1403	rBV	53313	81980	7.80%	0.677%
20	9.817	1416	1421	1426	rVB3	7477	12825	1.22%	0.106%
21	9.890	1426	1433	1437	rBV2	15968	32131	3.06%	0.266%
22	10.061	1455	1461	1464	rBV3	6135	11058	1.05%	0.091%
23	10.109	1464	1469	1475	rVV	669181	913678	86.94%	7.551%
24	10.182	1476	1481	1486	rVV3	11938	24317	2.31%	0.201%
25	10.250	1486	1492	1497	rVV	78343	113668	10.82%	0.939%
26	10.359	1498	1510	1516	rVV2	66084	148487	14.13%	1.227%
27	10.408	1516	1518	1526	rVB3	11405	19309	1.84%	0.160%
28	10.512	1531	1535	1541	rVB3	11880	18174	1.73%	0.150%
29	10.640	1545	1556	1564	rBV3	31362	73668	7.01%	0.609%
30	10.713	1564	1568	1579	rVB7	8730	19912	1.89%	0.165%
31	10.835	1579	1588	1592	rBV4	35454	59988	5.71%	0.496%
32	10.908	1592	1600	1604	rBV3	32414	53064	5.05%	0.439%
33	10.945	1604	1606	1612	rVB2	14648	18641	1.77%	0.154%
34	11.018	1612	1618	1621	rBV	45938	60941	5.80%	0.504%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\82X071619W.M
Title : SW846 8260

35	11.054	1621	1624	1632	rVB7	8461	16930	1.61%	0.140%
36	11.134	1632	1637	1642	rBV	592048	744490	70.84%	6.152%
37	11.182	1642	1645	1648	rVB2	14077	16747	1.59%	0.138%
38	11.274	1648	1660	1669	rVB4	37798	65252	6.21%	0.539%
39	11.359	1669	1674	1681	rBV	79492	113757	10.82%	0.940%
40	11.432	1681	1686	1688	rBV	68784	96668	9.20%	0.799%
41	11.505	1693	1698	1701	rVB	33673	41806	3.98%	0.345%
42	11.579	1706	1710	1715	rVB6	8643	12959	1.23%	0.107%
43	11.664	1715	1724	1732	rBV	309181	406894	38.72%	3.363%
44	11.768	1738	1741	1743	rVV2	15531	17903	1.70%	0.148%
45	11.804	1743	1747	1758	rVB	459062	560749	53.36%	4.634%
46	11.914	1758	1765	1767	rBV3	17591	30972	2.95%	0.256%
47	11.944	1767	1770	1775	rVB	49808	61834	5.88%	0.511%
48	12.024	1775	1783	1786	rBV	82646	104088	9.90%	0.860%
49	12.078	1786	1792	1797	rVB	692818	934737	88.95%	7.725%
50	12.139	1797	1802	1807	rBV	154249	184840	17.59%	1.528%
51	12.231	1811	1817	1823	rVB	24761	40940	3.90%	0.338%
52	12.298	1823	1828	1830	rBV2	105221	138480	13.18%	1.144%
53	12.322	1830	1832	1835	rVV	96971	98439	9.37%	0.813%
54	12.365	1835	1839	1847	rVB2	87227	129947	12.37%	1.074%
55	12.456	1849	1854	1859	rBV	65000	86095	8.19%	0.711%
56	12.517	1859	1864	1869	rVB	188584	230795	21.96%	1.907%
57	12.585	1869	1875	1877	rBV	105712	140800	13.40%	1.164%
58	12.609	1877	1879	1883	rVV	115559	121411	11.55%	1.003%
59	12.664	1885	1888	1891	rVV	104706	124715	11.87%	1.031%
60	12.700	1891	1894	1897	rVV	26618	31634	3.01%	0.261%
61	12.755	1898	1903	1911	rVV4	84323	207682	19.76%	1.716%
62	12.847	1915	1918	1921	rVV4	24566	33433	3.18%	0.276%
63	12.902	1923	1927	1931	rVB2	83642	111228	10.58%	0.919%
64	12.975	1931	1939	1942	rBV2	82307	131258	12.49%	1.085%
65	13.017	1942	1946	1952	rVB	146776	190503	18.13%	1.574%
66	13.078	1952	1956	1963	rVB3	40902	64877	6.17%	0.536%
67	13.145	1963	1967	1971	rBV	21494	33948	3.23%	0.281%
68	13.194	1971	1975	1982	rVV3	57567	88487	8.42%	0.731%
69	13.261	1982	1986	1989	rVB2	23355	33206	3.16%	0.274%
70	13.304	1989	1993	1995	rBV	45079	54496	5.19%	0.450%
71	13.340	1995	1999	2004	rVV2	255625	370554	35.26%	3.062%

LSC Area Percent Report

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Operator : JC/SP
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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Title : SW846 8260

72	13.383	2004	2006	2010	rVV	31102	38777	3.69%	0.320%
73	13.426	2010	2013	2018	rVV3	17218	30466	2.90%	0.252%
74	13.475	2018	2021	2031	rVB	92830	137725	13.11%	1.138%
75	13.560	2031	2035	2038	rBV4	21305	34041	3.24%	0.281%
76	13.597	2038	2041	2044	rVV	17602	22782	2.17%	0.188%
77	13.633	2044	2047	2053	rVV3	59108	123241	11.73%	1.018%
78	13.694	2055	2057	2059	rVV3	26231	33855	3.22%	0.280%
79	13.718	2059	2061	2066	rVB3	46793	64862	6.17%	0.536%
80	13.804	2072	2075	2077	rBV	14716	16223	1.54%	0.134%
81	13.846	2077	2082	2087	rVB3	35748	74770	7.11%	0.618%
82	13.907	2087	2092	2097	rVB4	33453	60267	5.73%	0.498%
83	13.981	2097	2104	2109	rBV2	48994	86192	8.20%	0.712%
84	14.029	2109	2112	2115	rBV2	16484	22302	2.12%	0.184%
85	14.127	2123	2128	2131	rBV3	17304	26497	2.52%	0.219%
86	14.279	2147	2153	2159	rVB4	44669	95140	9.05%	0.786%
87	14.377	2166	2169	2174	rVB2	17817	23106	2.20%	0.191%
88	14.426	2174	2177	2181	rVB3	18132	23700	2.26%	0.196%
89	14.474	2181	2185	2191	rVB	14492	20780	1.98%	0.172%
90	14.535	2191	2195	2204	rBV2	49189	74467	7.09%	0.615%
91	14.669	2214	2217	2218	rBV3	17030	18529	1.76%	0.153%
92	14.694	2218	2221	2224	rVV	31537	38743	3.69%	0.320%
93	14.730	2224	2227	2231	rVB2	14013	18060	1.72%	0.149%
94	14.779	2231	2235	2239	rBV4	6562	10740	1.02%	0.089%
95	14.840	2240	2245	2249	rVV2	37730	56165	5.34%	0.464%
96	15.249	2307	2312	2320	rVB5	8915	20608	1.96%	0.170%
97	15.541	2356	2360	2366	rVB6	7700	12646	1.20%	0.105%
98	15.682	2377	2383	2387	rBV2	7902	14692	1.40%	0.121%

Sum of corrected areas: 12100818

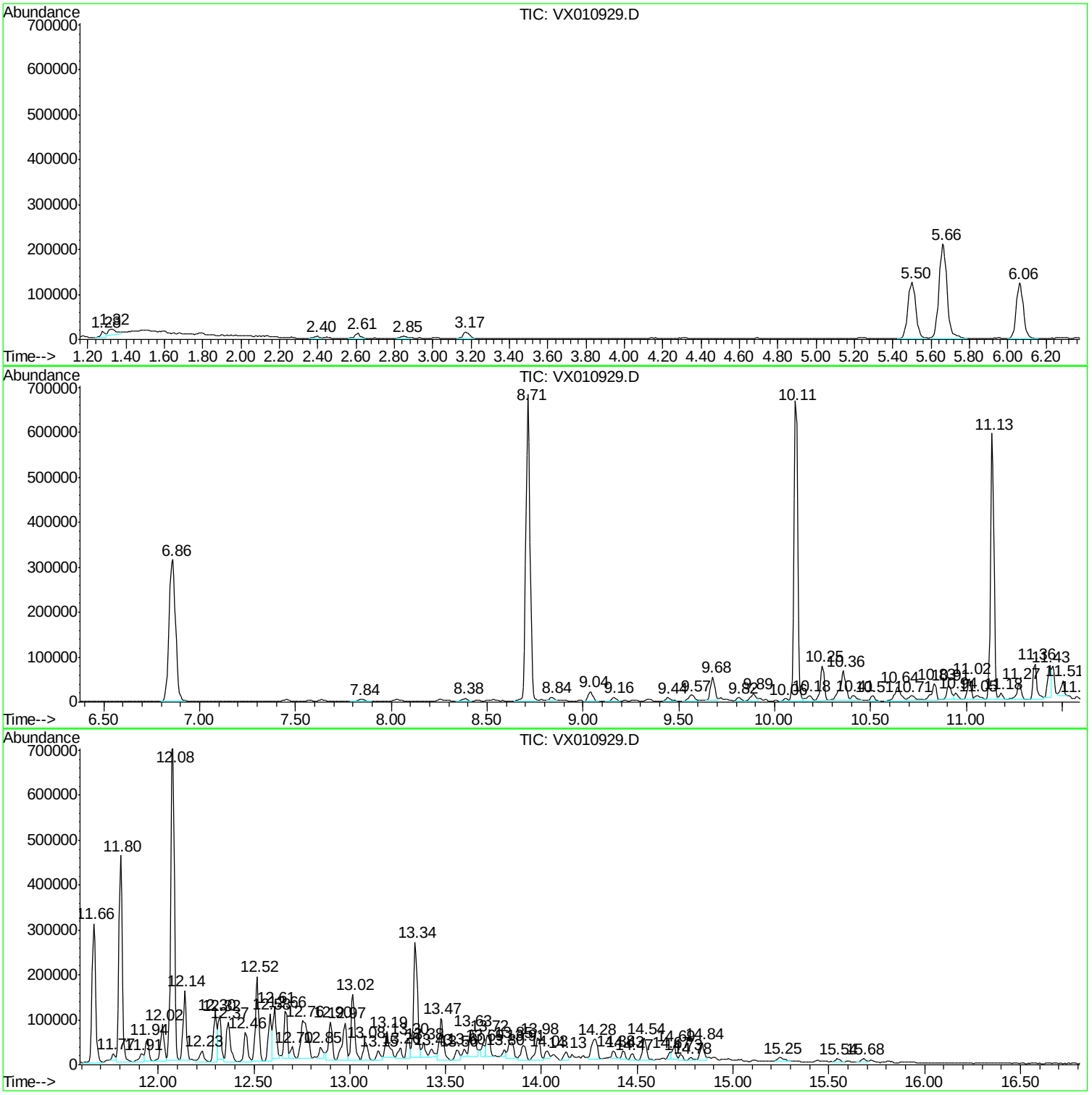
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

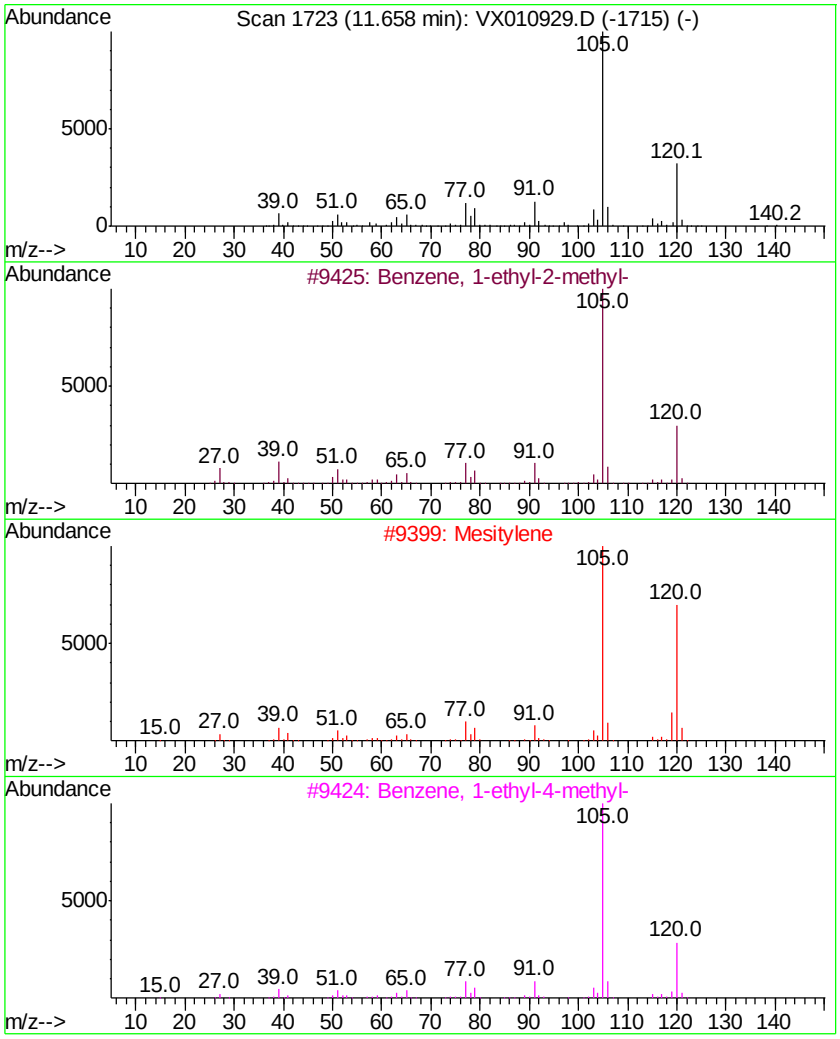
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

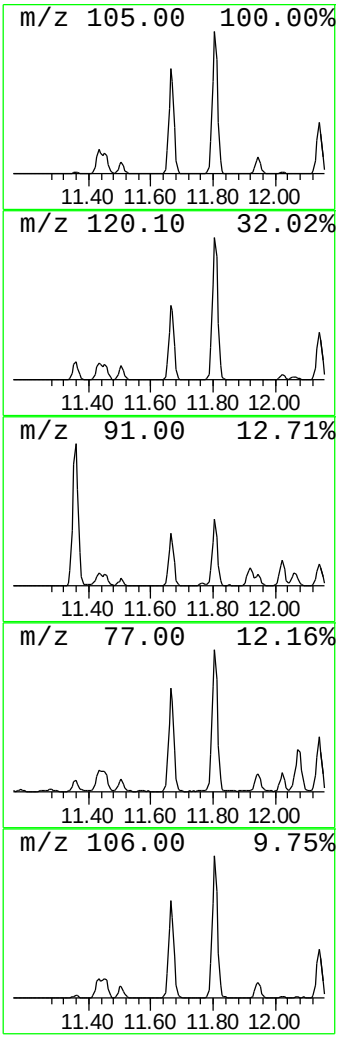
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	21.77 ug/l	406894	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Mesitylene	120	C9H12	000108-67-8	90
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

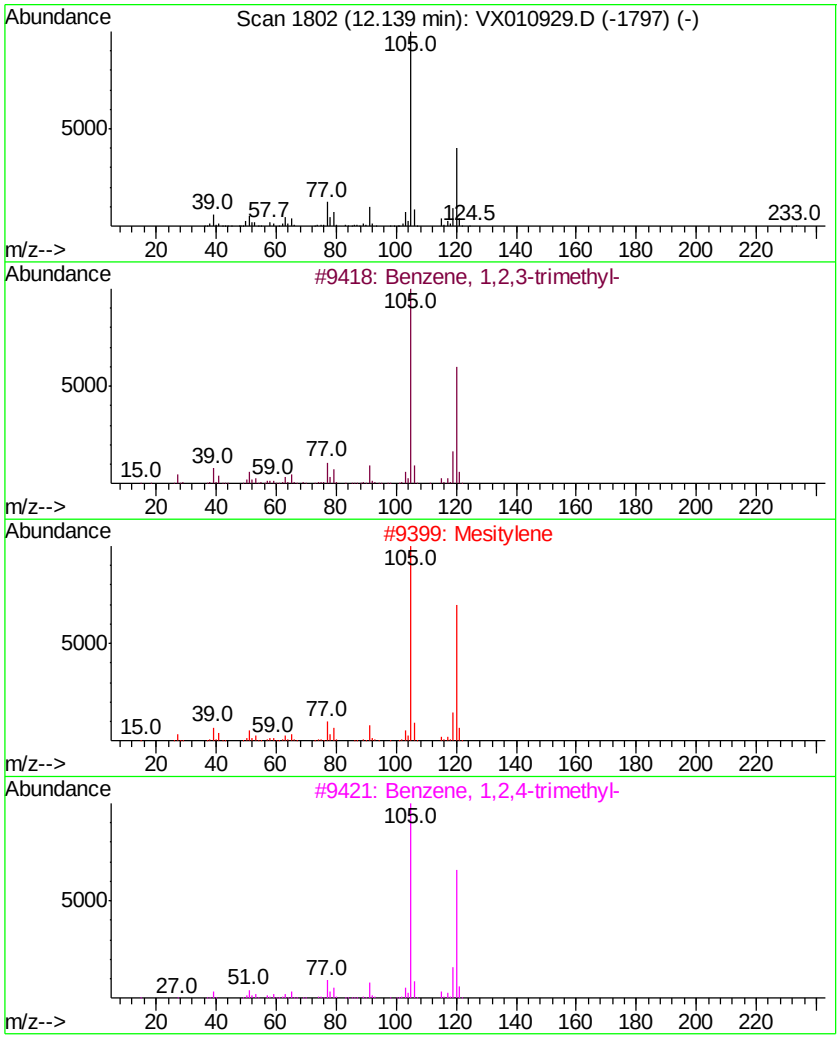
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

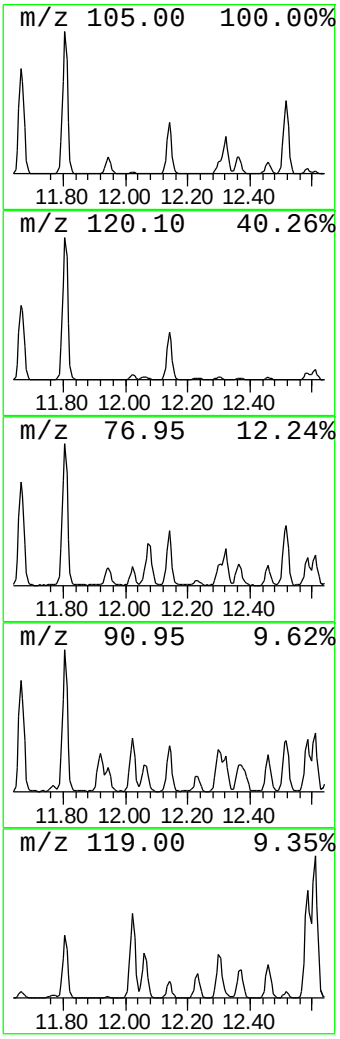
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	9.89 ug/l	184840	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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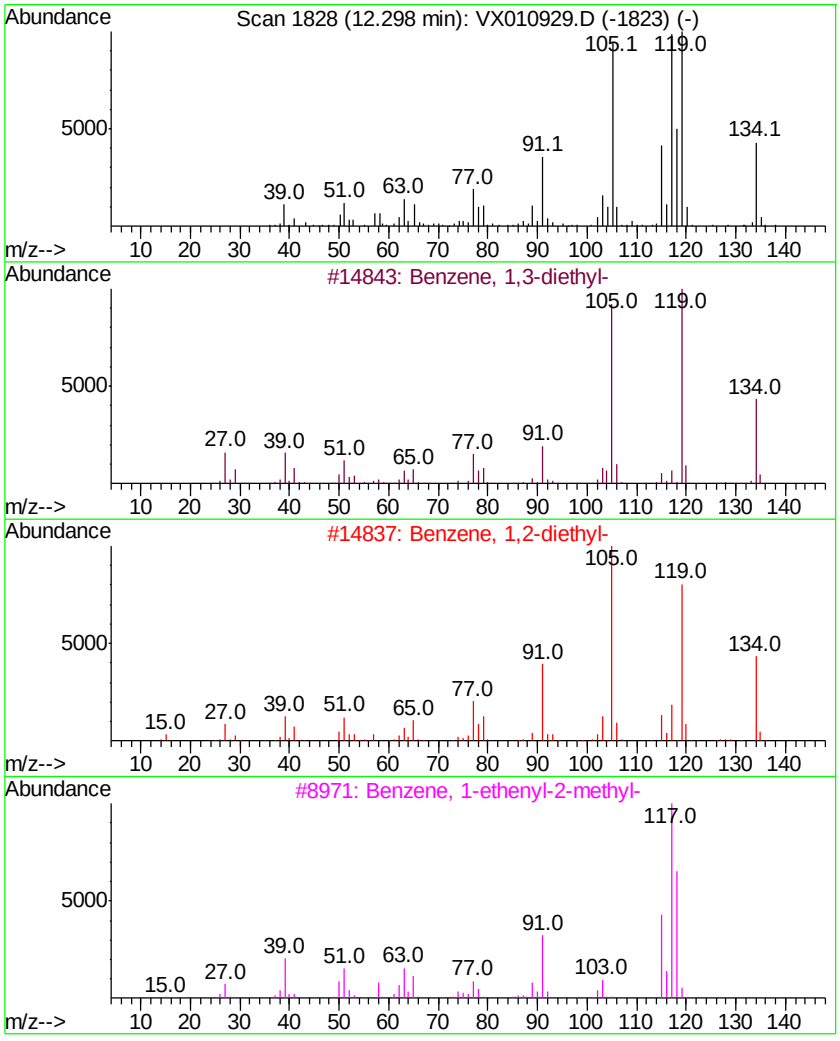
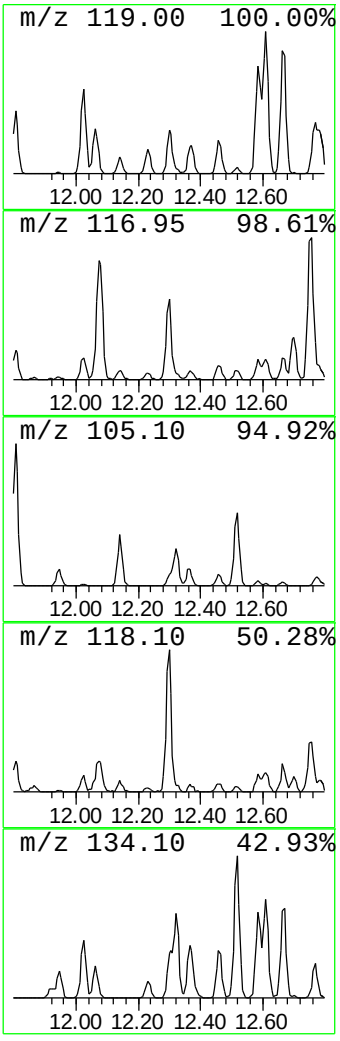
Peak Number 3 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	7.41 ug/l	138480	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	86
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	86
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	80
5	Indane	118	C9H10	000496-11-7	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

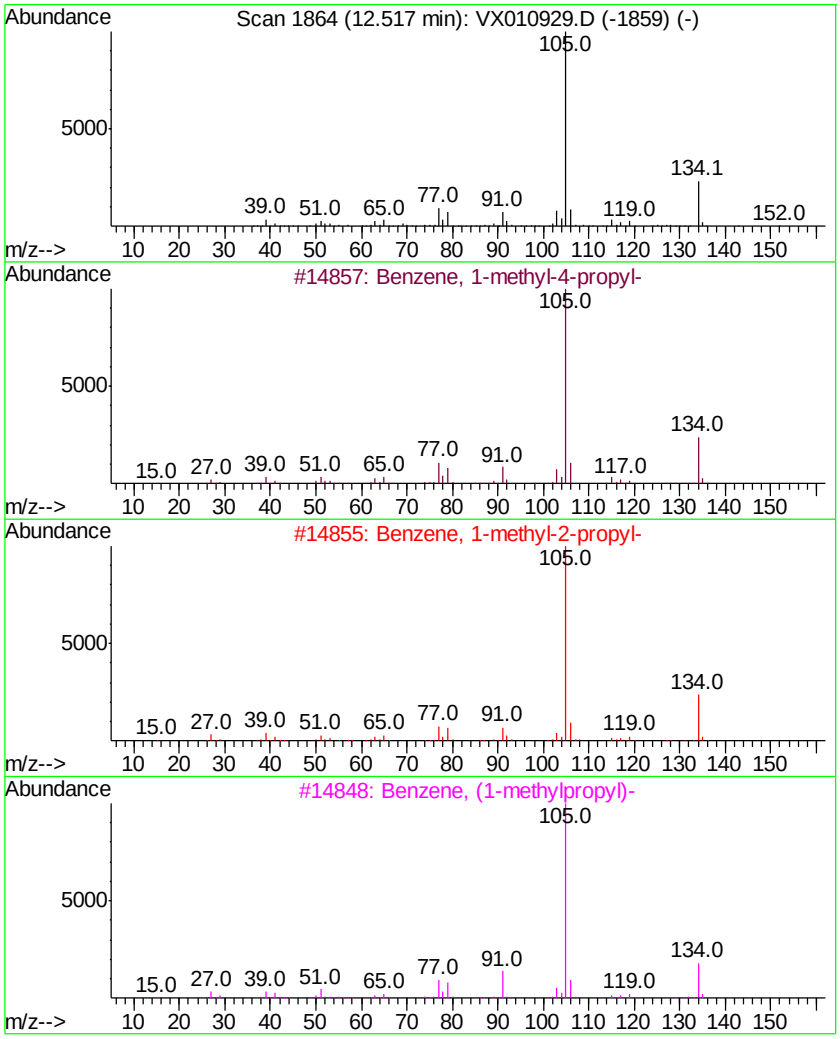
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

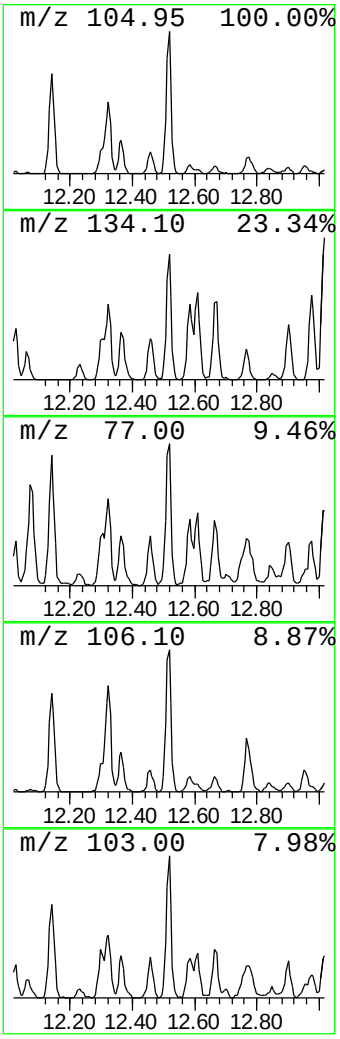
****

Peak Number 4 Benzene, 1-methyl-4-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	12.35 ug/l	230795	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	93
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	80
5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

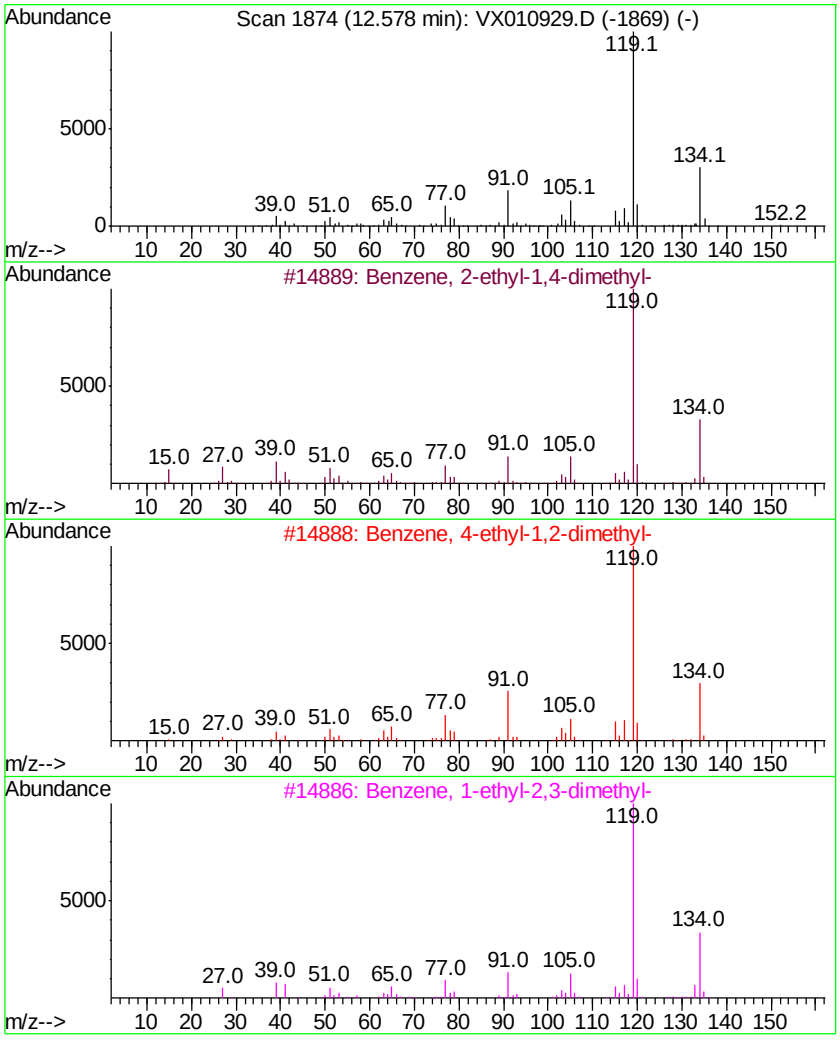
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

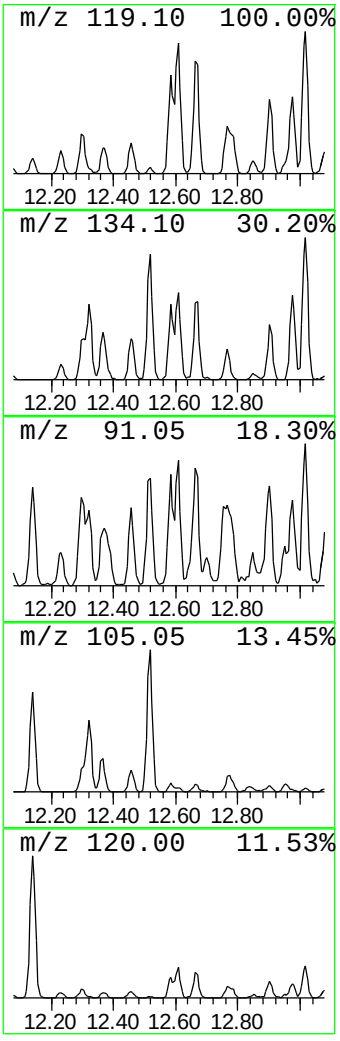
****

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	7.53 ug/l	140800	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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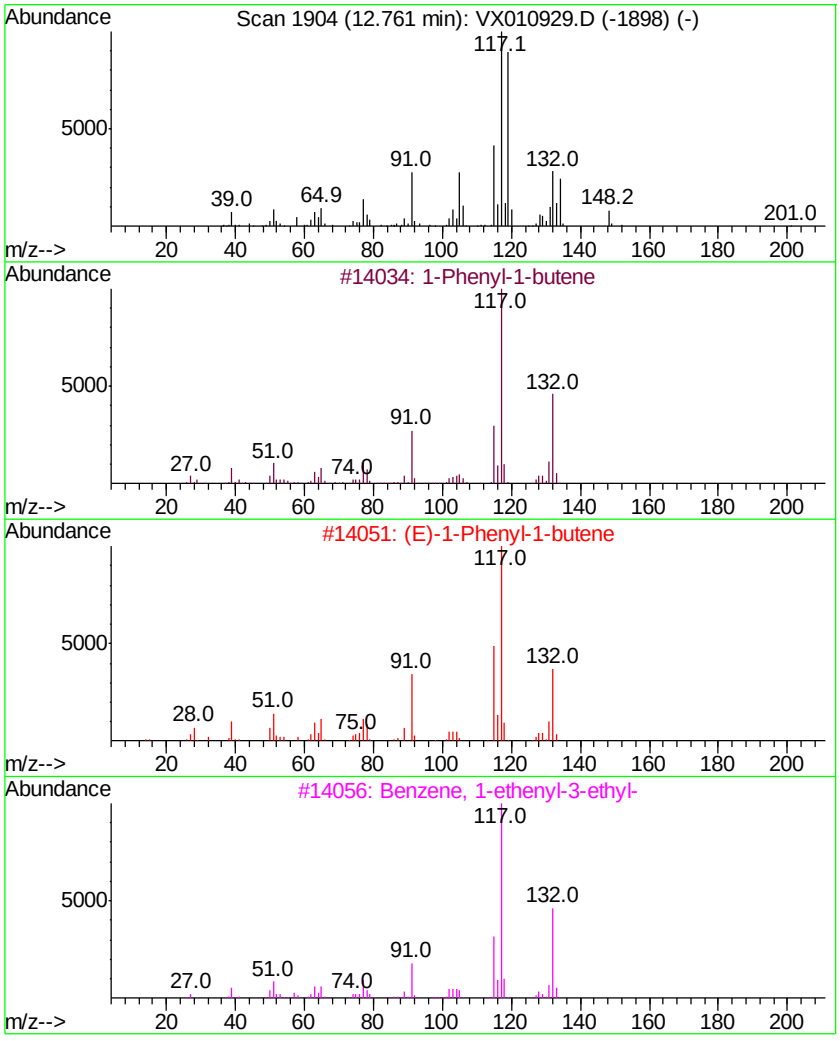
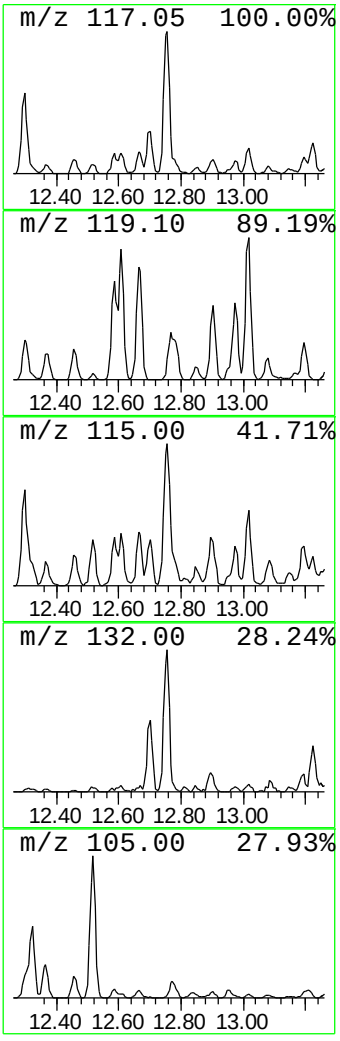
Peak Number 6 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	11.11 ug/l	207682	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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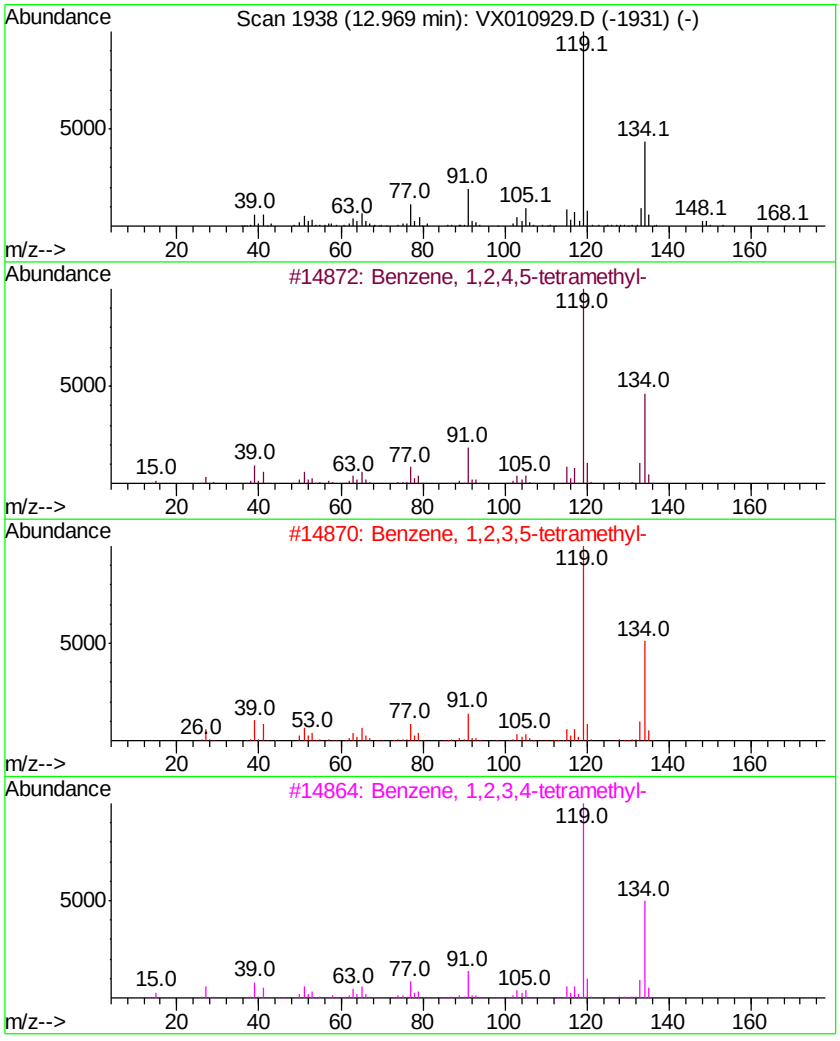
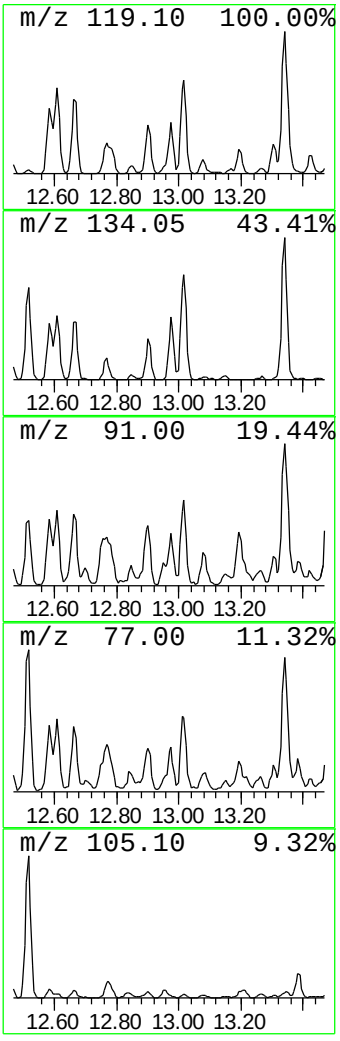
Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	7.02 ug/l	131258	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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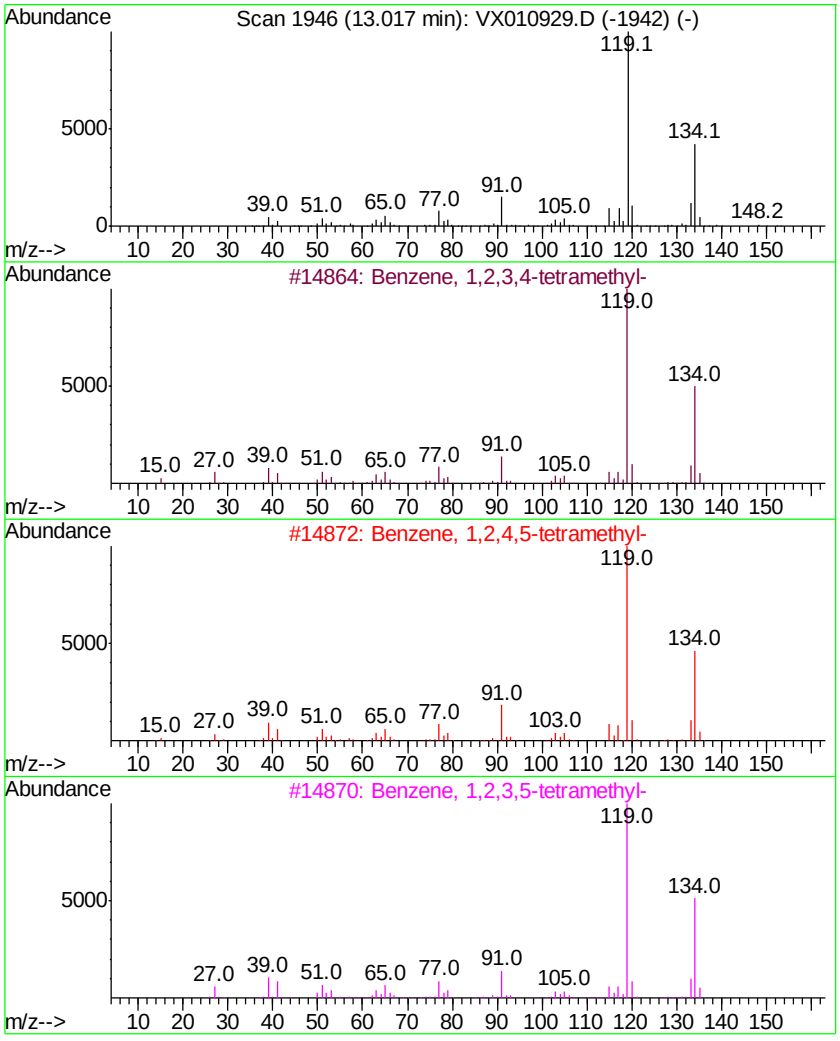
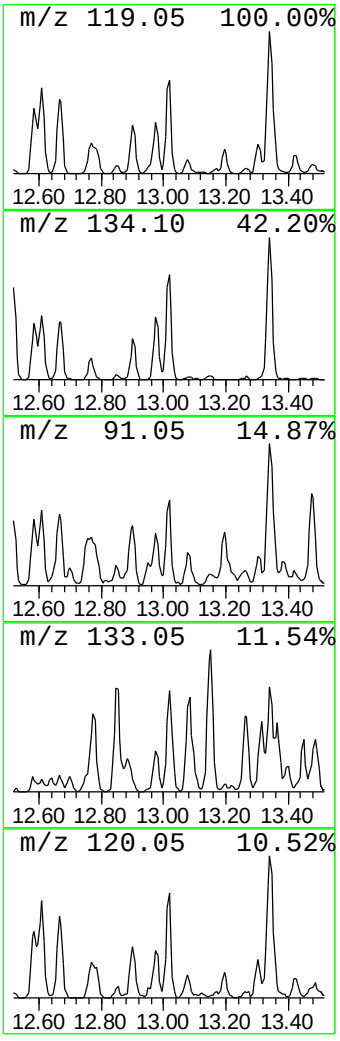
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	10.19 ug/l	190503	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94
5	o-Cymene	134	C10H14	000527-84-4	94





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Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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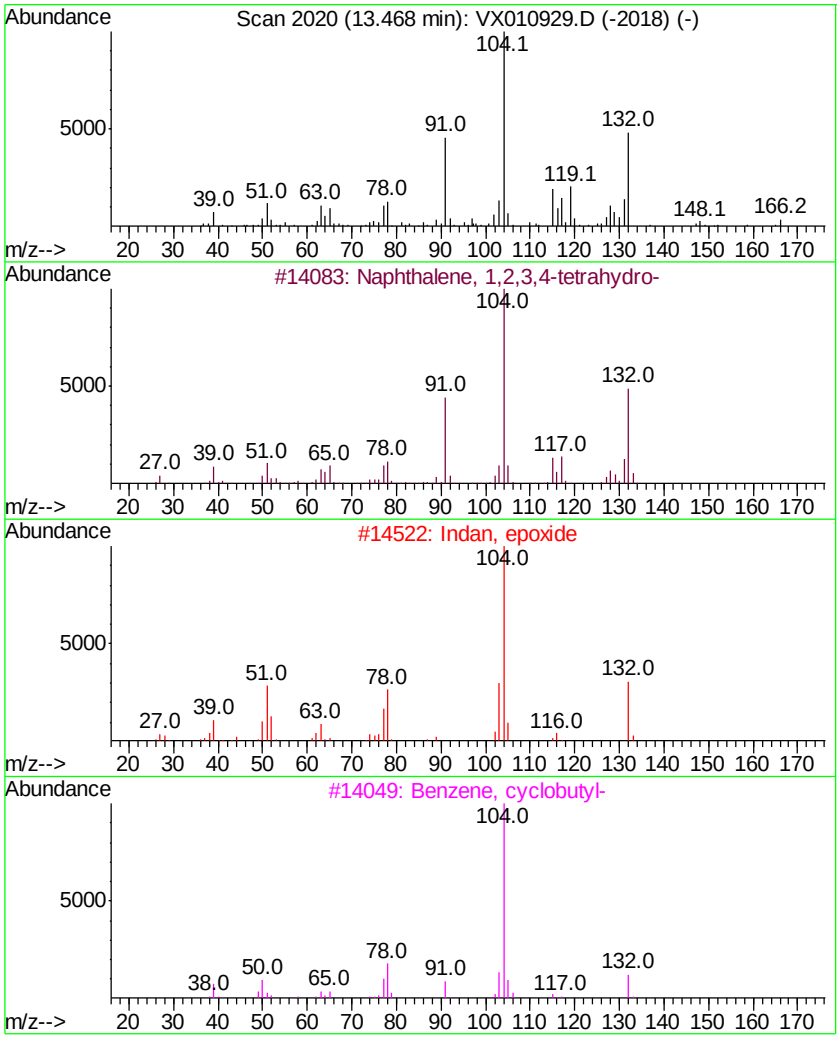
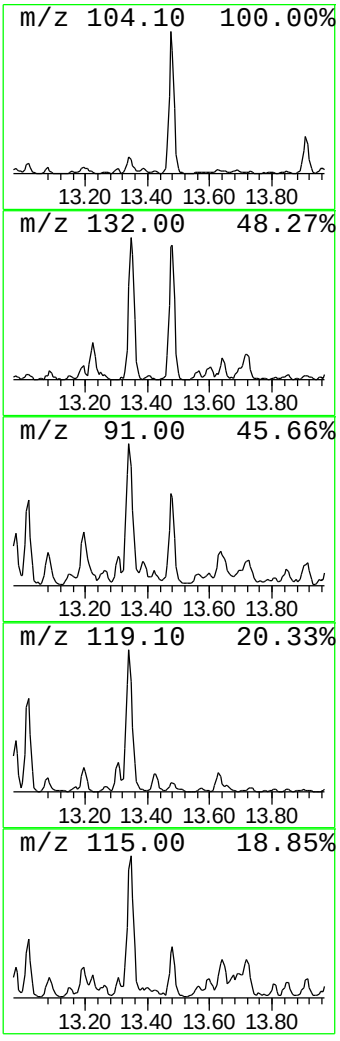
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	7.37 ug/l	137725	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Indan, epoxide	132	C9H8O	000768-22-9	60
3	Benzene, cyclobutyl-	132	C10H12	004392-30-7	50
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46
5	Pentafluoropropionic acid, 2-phe...	268	C11H9F5O2	081089-48-7	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071719\
Data File : VX010929.D
Acq On : 17 Jul 2019 12:36
Operator : JC/SP
Sample : K3791-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0628

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-ethyl-...	11.66	21.8	ug/l	406894	4	12.08	934737	50.0
Benzene, 1,2,3-tr...	12.14	9.9	ug/l	184840	4	12.08	934737	50.0
Benzene, 1,3-diet...	12.30	7.4	ug/l	138480	4	12.08	934737	50.0
Benzene, 1-methyl...	12.52	12.4	ug/l	230795	4	12.08	934737	50.0
Benzene, 2-ethyl-...	12.58	7.5	ug/l	140800	4	12.08	934737	50.0
1-Phenyl-1-butene	12.76	11.1	ug/l	207682	4	12.08	934737	50.0
Benzene, 1,2,4,5-...	12.97	7.0	ug/l	131258	4	12.08	934737	50.0
Benzene, 1,2,3,4-...	13.02	10.2	ug/l	190503	4	12.08	934737	50.0
Naphthalene, 1,2,...	13.47	7.4	ug/l	137725	4	12.08	934737	50.0