LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010699.D
Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

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\82X061919W.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	1.788	100	104	110	rVB	31256	45640	1.16%	0.189%
2	2.849	270	278	286	rVB2	21909	48601	1.24%	0.202%
3	2.934	286	292	300	rVB	26331	54430	1.39%	0.226%
4	3.172	323	331	341	rBV	60830	141355	3.60%	0.586%
5	4.403	524	533	545	rVB	55452	159033	4.05%	0.659%
6	5.501	701	713	720	rBV2	127867	371366	9.46%	1.540%
7	5.580	720	726	732	rVV3	44848	115200	2.93%	0.478%
8	5.659	732	739	748	rVV	197978	520730	13.26%	2.159%
9	5.952	775	787	796	rBV3	17545	55759	1.42%	0.231%
10	6.061	796	805	814	rVV	126756	332528	8.47%	1.379%
11	6.141	814	818	827	rVV2	14626	39396	1.00%	0.163%
12	6.263	829	838	847	rVV4	31915	113125	2.88%	0.469%
13	6.360	847	854	861	rVV2	32362	90289	2.30%	0.374%
14	6.458	862	870	883	rVB2	32148	91212	2.32%	0.378%
15	6.860	924	936	950	rBV	328241	778235	19.82%	3.227%
16	7.458	1023	1034	1048	rBV2	115623	293061	7.46%	1.215%
17	7.848	1090	1098	1107	rVB2	24713	54661	1.39%	0.227%
18	8.025	1119	1127	1139	rBV2	31172	75356	1.92%	0.312%
19	8.390	1179	1187	1193	rBV2	23119	47126	1.20%	0.195%
20	8.665	1224	1232	1235	rBV2	51598	95463	2.43%	0.396%
21	8.713	1235	1240	1249	rVB	697615	1106542	28.18%	4.588%
22	8.841	1254	1261	1268	rBV	51175	92874	2.37%	0.385%
23	9.037	1288	1293	1301	rVB	106992	173885	4.43%	0.721%
24	9.165	1306	1314	1320	rBV2	45402	77166	1.97%	0.320%
25	9.567	1375	1380	1385	rBV2	29427	45710	1.16%	0.190%
26	9.677	1393	1398	1403	rBV	78571	118947	3.03%	0.493%
27	10.110	1464	1469	1476	rBV	692032	925236	23.56%	3.836%
28	10.183	1476	1481	1486	rVB	34924	53508	1.36%	0.222%
29	10.250	1486	1492	1501	rBV	263406	355741	9.06%	1.475%
30	10.359	1501	1510	1516	rVV	290989	403454	10.27%	1.673%
31	10.646	1547	1557	1562	rBV4	24667	70284	1.79%	0.291%
32	10.695	1562	1565	1575	rVB2	41150	65741	1.67%	0.273%
33	10.835	1584	1588	1593	rVB	29754	41806	1.06%	0.173%
34	11.018	1613	1618	1622	rBV	180977	222767	5.67%	0.924%

LSC Area Percent Report

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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

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\82X061919W.M
Title : SW846 8260

35	11.134	1631	1637	1642	rBV	607947	762939	19.43%	3.163%
36	11.359	1669	1674	1683	rVB	221089	277803	7.07%	1.152%
37	11.451	1683	1689	1694	rBV	83044	113358	2.89%	0.470%
38	11.664	1716	1724	1734	rBV	146702	212364	5.41%	0.880%
39	11.804	1742	1747	1753	rVB	3335968	3926935	100.00%	16.281%
40	11.945	1767	1770	1778	rVB	65715	86521	2.20%	0.359%
41	12.079	1786	1792	1797	rBV	731299	955678	24.34%	3.962%
42	12.140	1797	1802	1812	rVB	1602678	1918436	48.85%	7.954%
43	12.231	1812	1817	1821	rBV	58935	73558	1.87%	0.305%
44	12.298	1822	1828	1833	rBV	474046	580223	14.78%	2.406%
45	12.365	1833	1839	1849	rVB2	115390	187680	4.78%	0.778%
46	12.457	1849	1854	1859	rBV	142277	177412	4.52%	0.736%
47	12.518	1859	1864	1870	rBV	299860	364998	9.29%	1.513%
48	12.585	1870	1875	1877	rBV	326584	421695	10.74%	1.748%
49	12.609	1877	1879	1883	rVV	354373	356960	9.09%	1.480%
50	12.664	1883	1888	1892	rVV	366782	462740	11.78%	1.919%
51	12.701	1892	1894	1899	rVV	94877	105815	2.69%	0.439%
52	12.755	1899	1903	1910	rVB2	288704	465591	11.86%	1.930%
53	12.902	1922	1927	1932	rVB2	232777	318391	8.11%	1.320%
54	12.975	1932	1939	1942	rBV	269440	352270	8.97%	1.461%
55	13.018	1942	1946	1952	rVB	461123	559768	14.25%	2.321%
56	13.079	1952	1956	1963	rVB3	66580	93208	2.37%	0.386%
57	13.146	1963	1967	1971	rBV	40993	59109	1.51%	0.245%
58	13.194	1971	1975	1977	rBV	72964	82111	2.09%	0.340%
59	13.225	1977	1980	1983	rVB	63763	64285	1.64%	0.267%
60	13.261	1983	1986	1989	rVB2	34765	44175	1.12%	0.183%
61	13.304	1989	1993	1995	rBV	69352	87545	2.23%	0.363%
62	13.341	1995	1999	2005	rVB2	815721	1114565	28.38%	4.621%
63	13.475	2017	2021	2027	rVB	368990	485382	12.36%	2.012%
64	13.560	2027	2035	2038	rBV3	33793	62356	1.59%	0.259%
65	13.597	2038	2041	2044	rVV	43449	65408	1.67%	0.271%
66	13.639	2044	2048	2053	rVV3	105779	227494	5.79%	0.943%
67	13.694	2053	2057	2059	rVV3	57456	98737	2.51%	0.409%
68	13.719	2059	2061	2069	rVB2	90415	140815	3.59%	0.584%
69	13.834	2069	2080	2087	rBV	430877	658864	16.78%	2.732%
70	13.908	2087	2092	2097	rVB3	44981	68412	1.74%	0.284%
71	13.987	2097	2105	2108	rBV2	79218	119969	3.06%	0.497%

LSC Area Percent Report

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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

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\82X061919W.M
Title : SW846 8260

72	14.286	2146	2154	2158	rVB2	77772	159143	4.05%	0.660%
73	14.426	2174	2177	2181	rVB2	31684	40542	1.03%	0.168%
74	14.536	2190	2195	2203	rBV2	101609	142291	3.62%	0.590%
75	14.694	2214	2221	2224	rBV	78034	110676	2.82%	0.459%
76	14.834	2239	2244	2249	rBV	174396	236888	6.03%	0.982%

Sum of corrected areas: 24119337

LSC Report - Integrated Chromatogram

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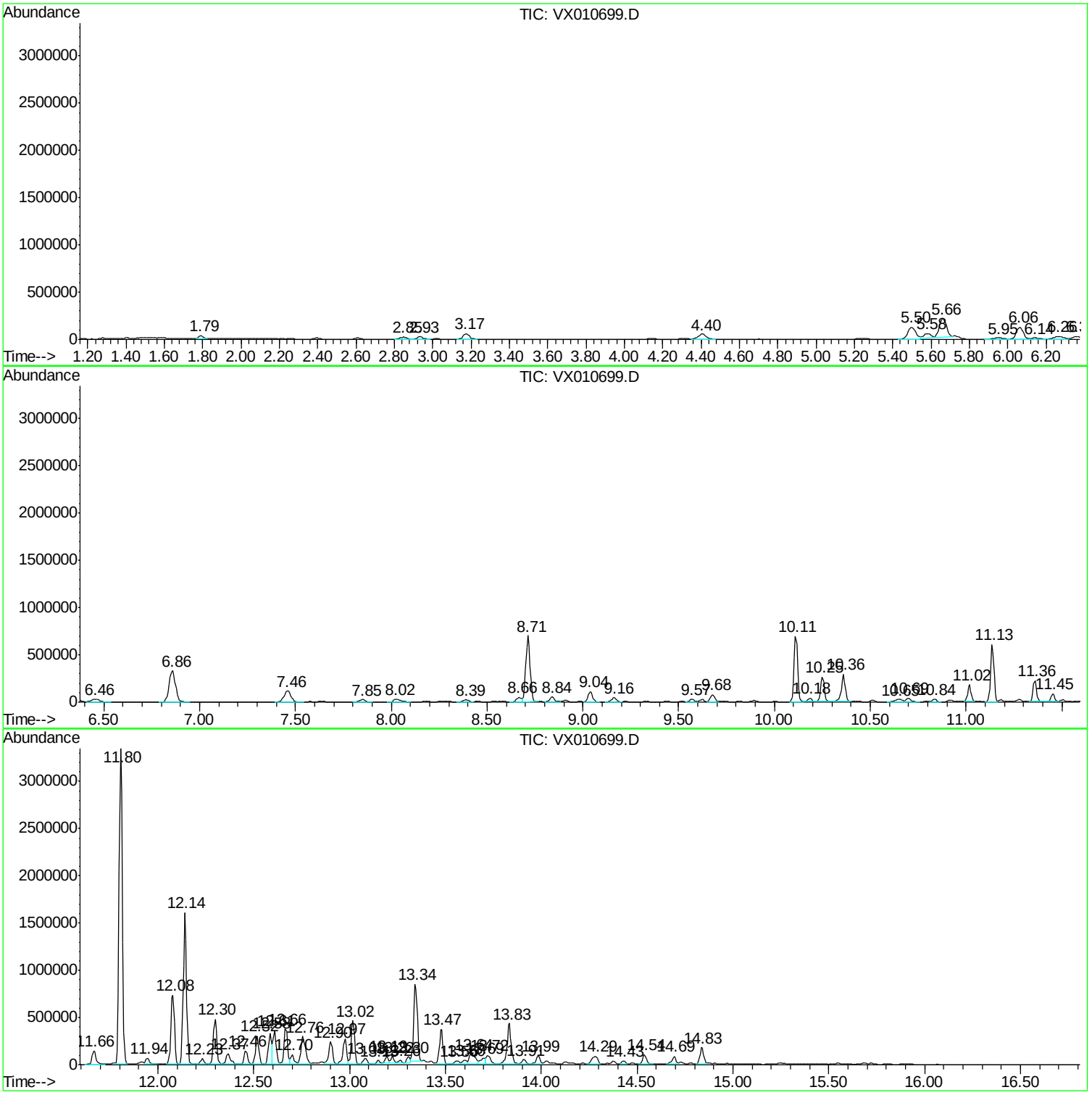
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



Library Search Compound Report

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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

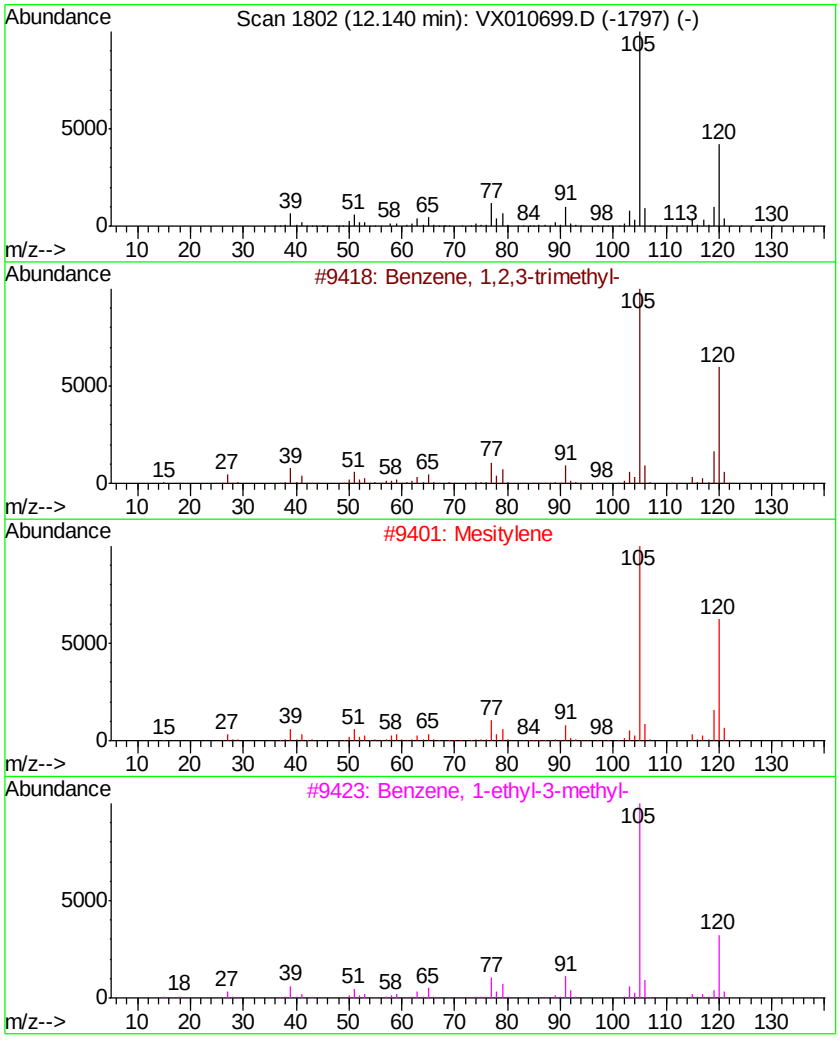
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

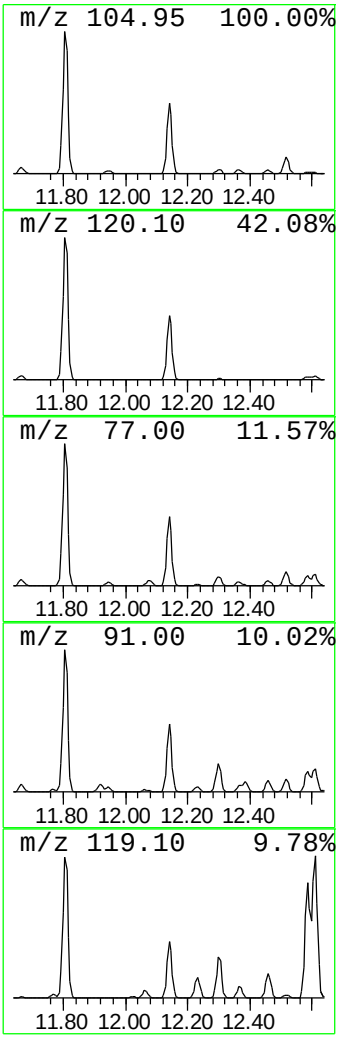
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	100.37 ug/l	1918440	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-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





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Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

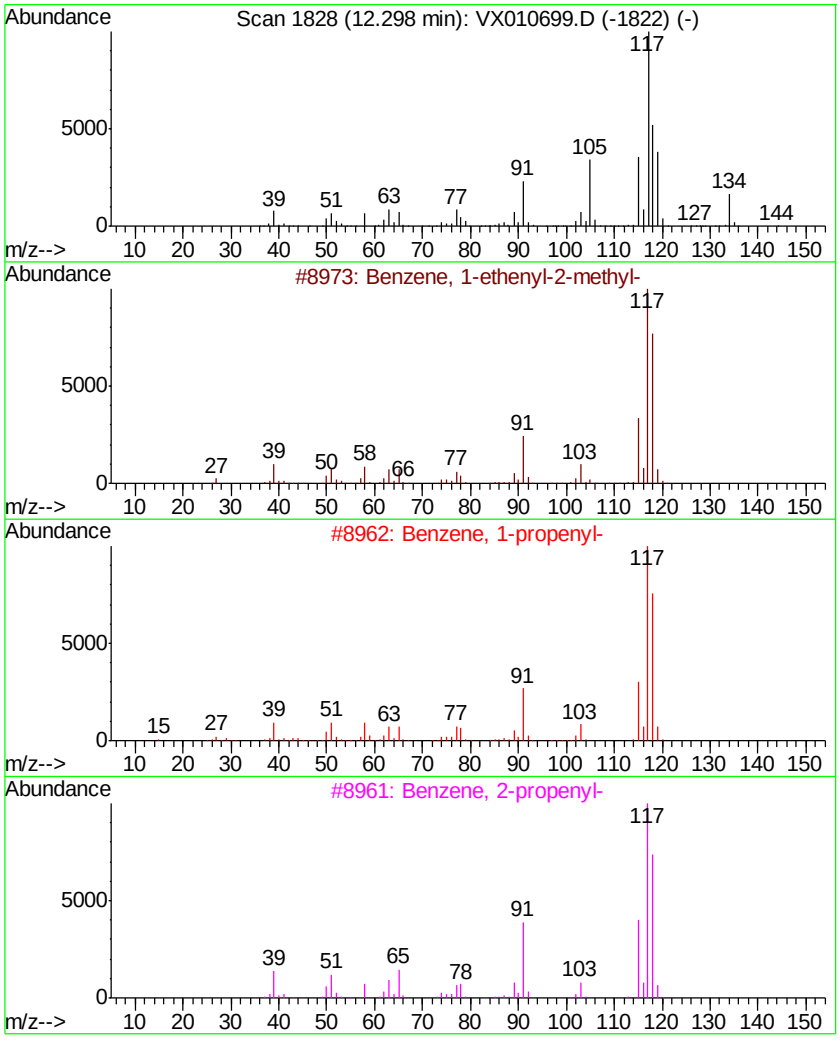
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

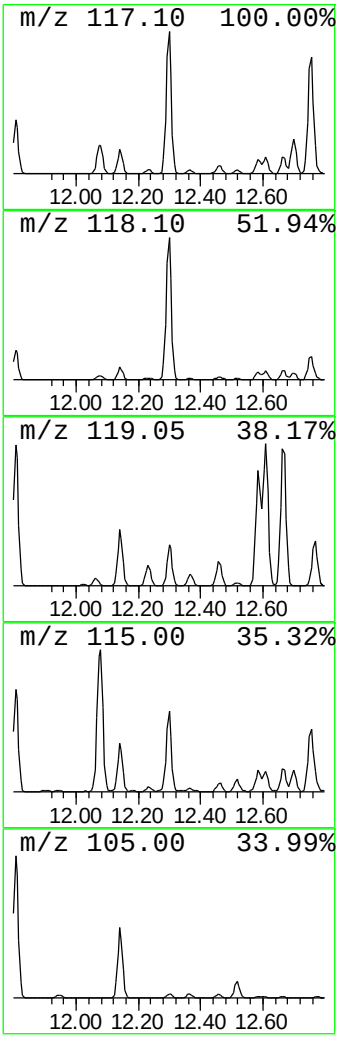
****

Peak Number 2 Benzene, 1-ethenyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	30.36 ug/l	580223	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	87
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	87
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
5	Indane	118	C9H10	000496-11-7	55





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010699.D
Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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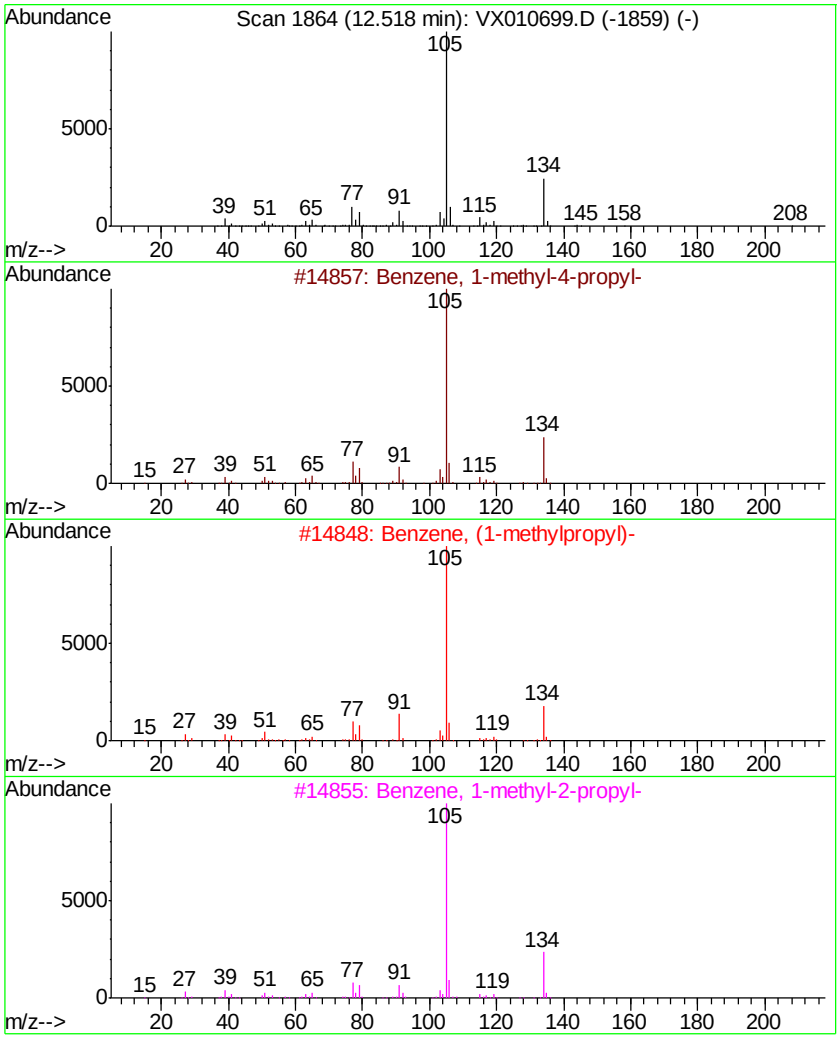
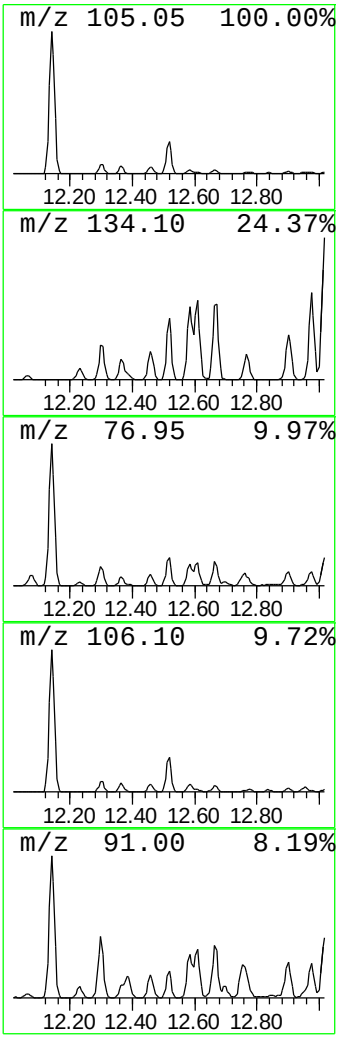
****

Peak Number 3 Benzene, 1-methyl-4-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	19.10 ug/l	364998	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	94
2	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	87
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	80





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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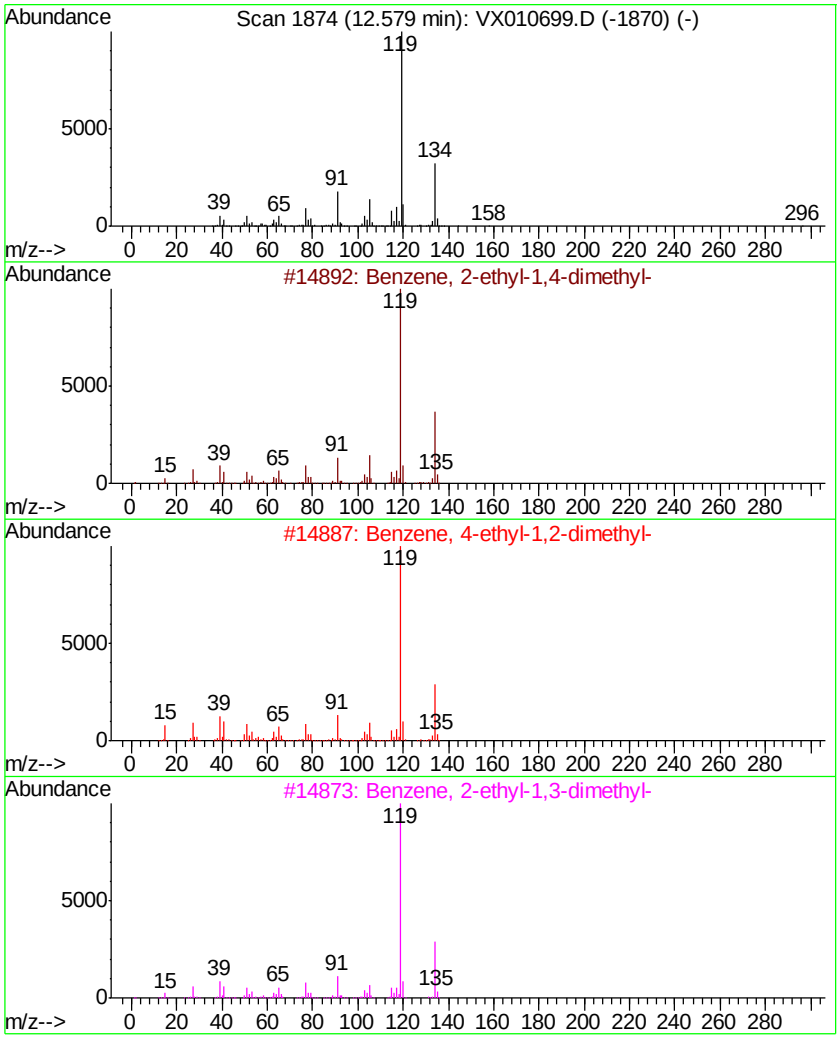
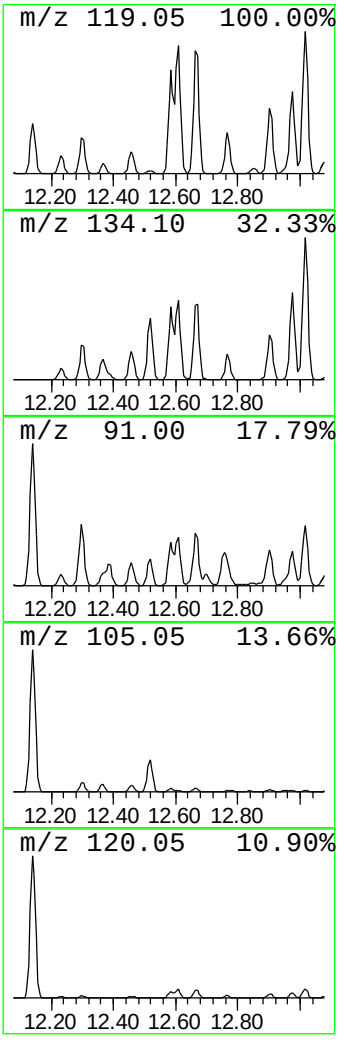
Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	22.06 ug/l	421695	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	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

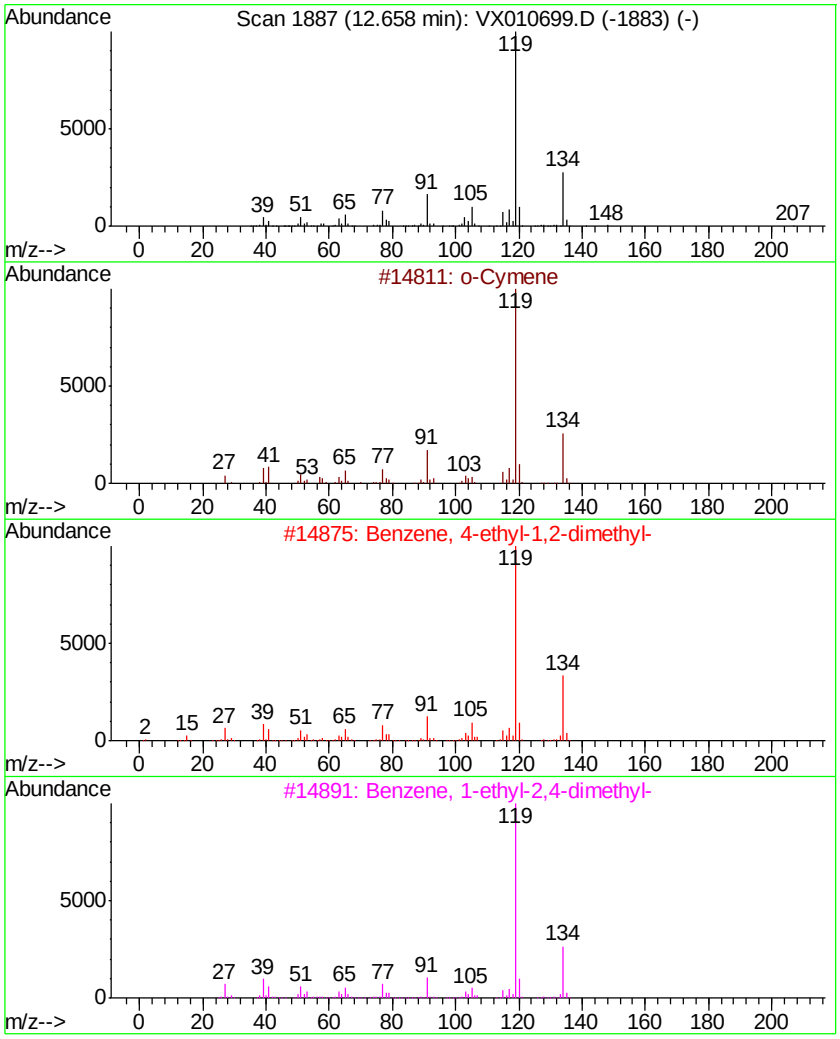
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

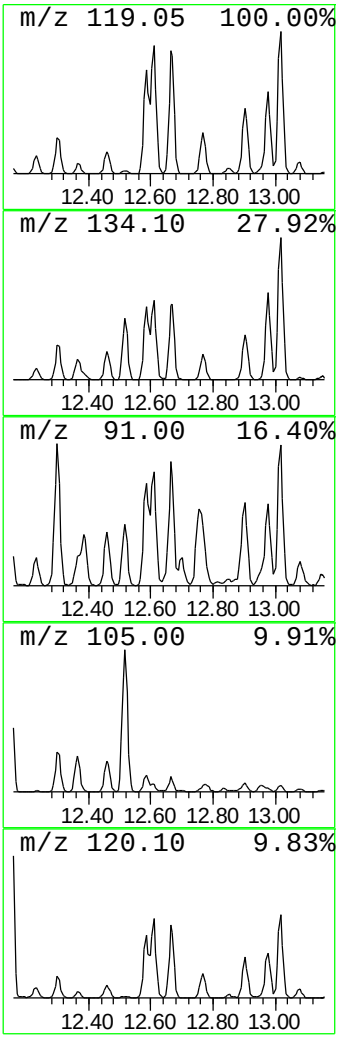
****

Peak Number 6 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	24.21 ug/l	462740	1,4-Dichlorobenzene-d4	12.08

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	94
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	p-Cymene	134	C10H14	000099-87-6	94





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Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

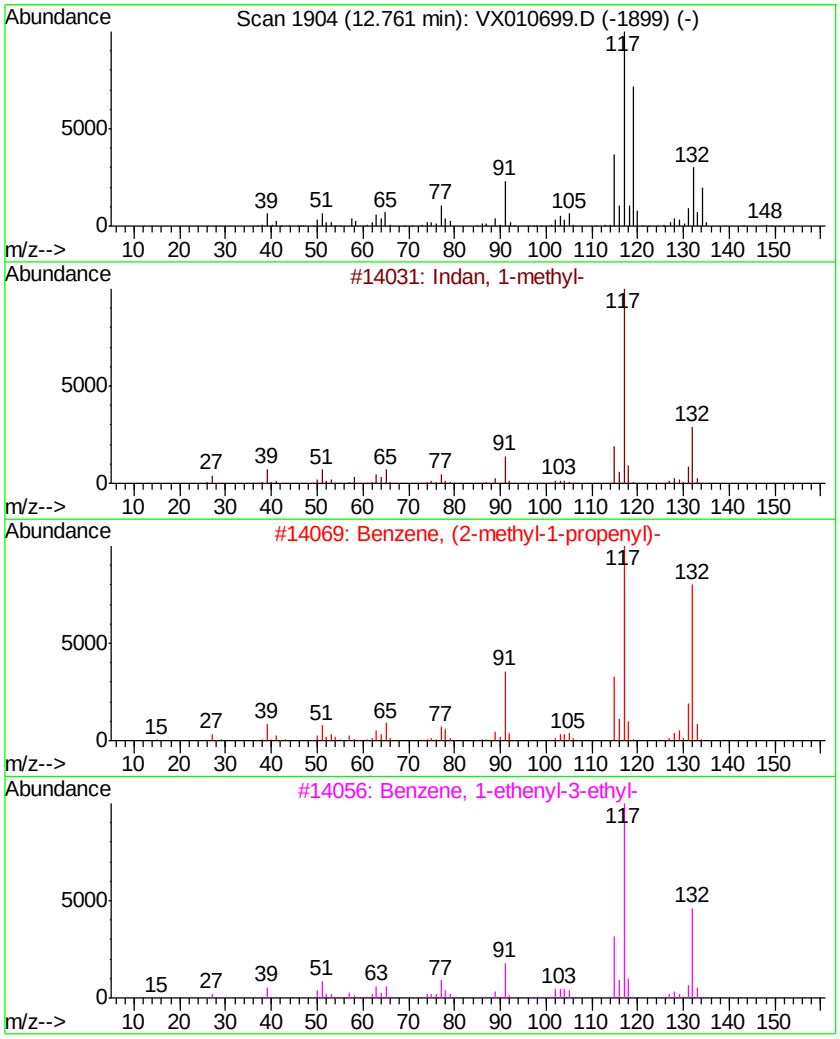
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

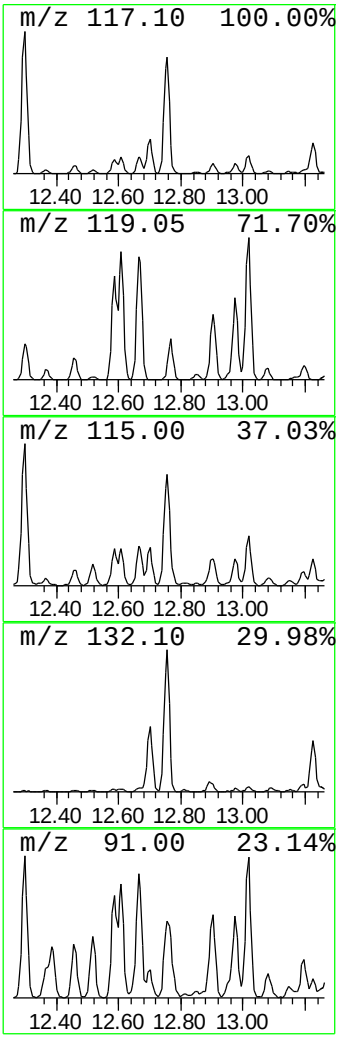
****

Peak Number 7 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	24.36 ug/l	465591	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	74





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010699.D
Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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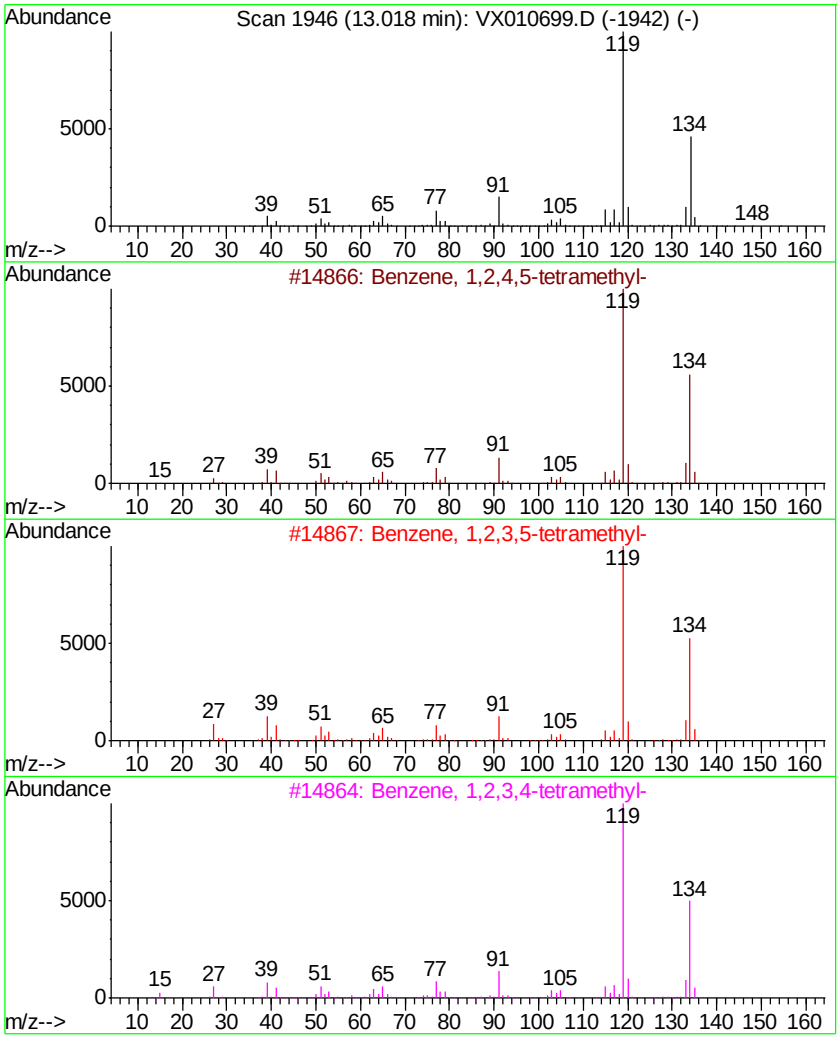
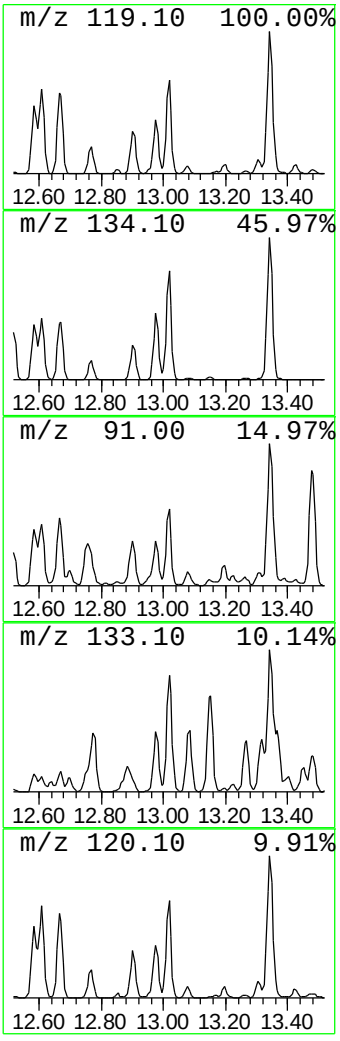
Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	29.29 ug/l	559768	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	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010699.D
Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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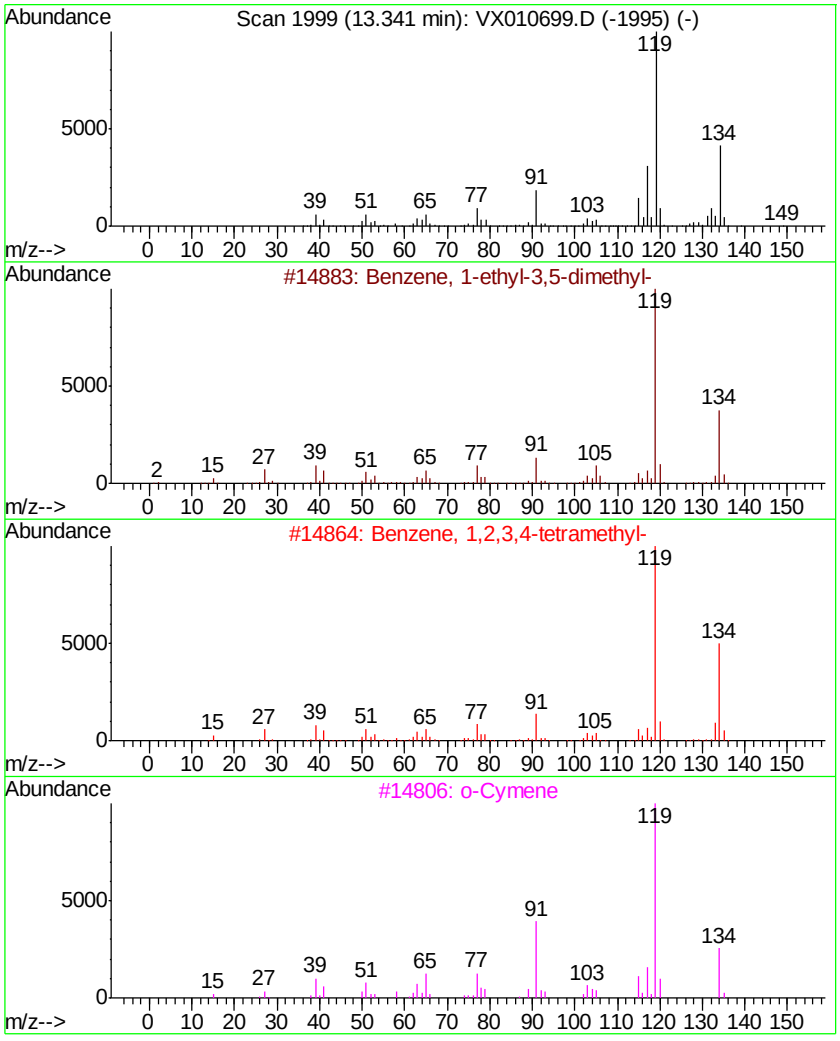
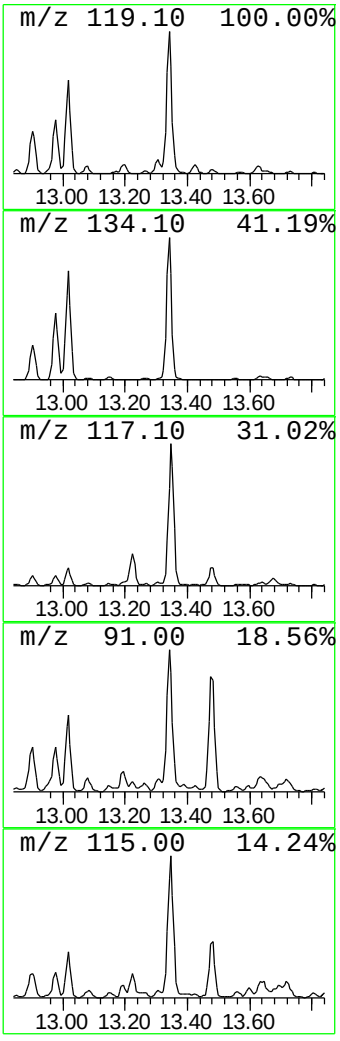
Peak Number 9 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	58.31 ug/l	1114570	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	81
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81
3	o-Cymene	134	C10H14	000527-84-4	81
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	76
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	76





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010699.D
Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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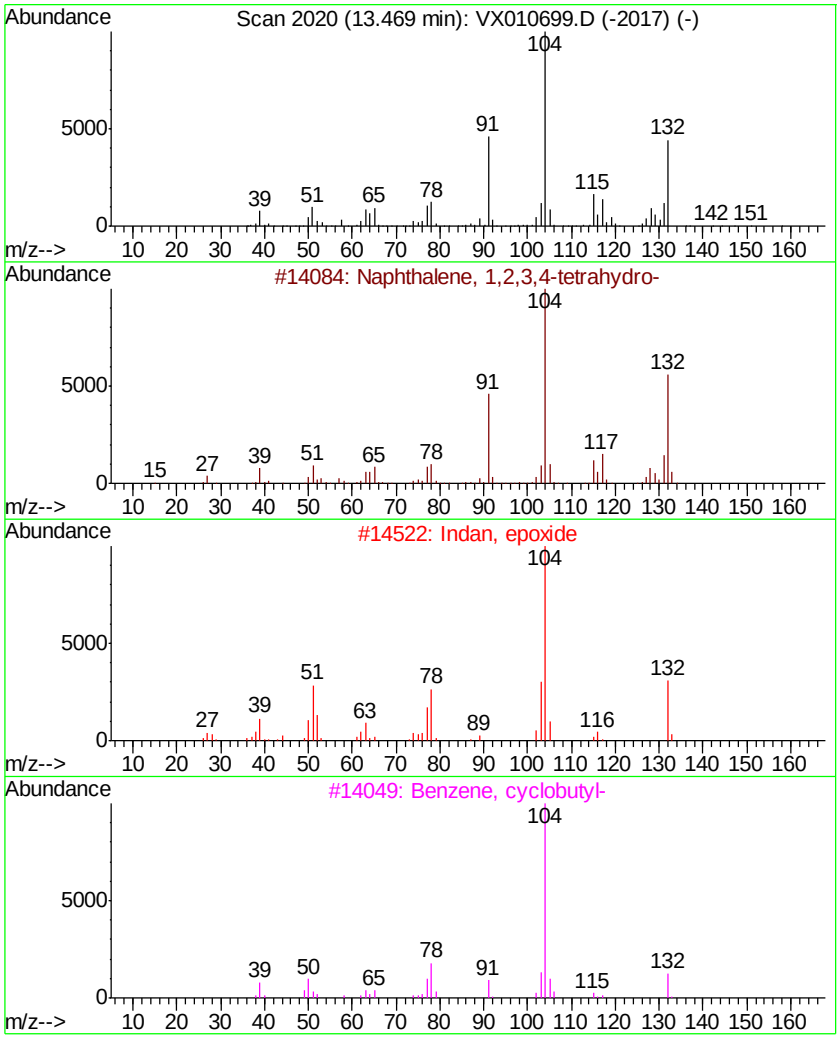
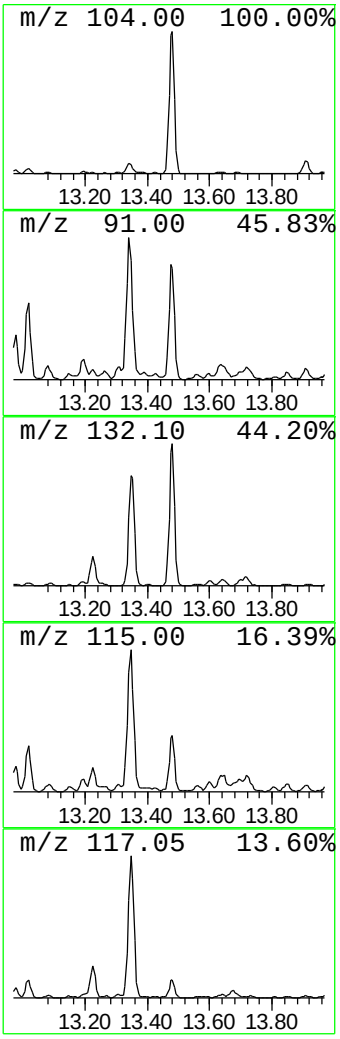
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	25.39 ug/l	485382	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	91
2	Indan, epoxide	132	C9H8O	000768-22-9	58
3	Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
5	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010699.D
Acq On : 08 Jul 2019 16:36
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0580

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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, 1,2,3-tr...	12.14	100.4	ug/l	1918440	4	12.08	955678	50.0
Benzene, 1-etheny...	12.30	30.4	ug/l	580223	4	12.08	955678	50.0
Benzene, 1-methyl...	12.52	19.1	ug/l	364998	4	12.08	955678	50.0
Benzene, 2-ethyl-...	12.58	22.1	ug/l	421695	4	12.08	955678	50.0
o-Cymene	12.66	24.2	ug/l	462740	4	12.08	955678	50.0
Indan, 1-methyl-	12.76	24.4	ug/l	465591	4	12.08	955678	50.0
Benzene, 1,2,4,5-...	13.02	29.3	ug/l	559768	4	12.08	955678	50.0
Benzene, 1-ethyl-...	13.34	58.3	ug/l	1114570	4	12.08	955678	50.0
Naphthalene, 1,2,...	13.47	25.4	ug/l	485382	4	12.08	955678	50.0