LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

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\82X112619W.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	1.070	26	30	39	rBV	2521834	2938240	7.36%	0.920%
2	1.783	141	147	159	rVB	1637979	2240700	5.61%	0.702%
3	2.838	311	320	323	rBV	480303	1029321	2.58%	0.322%
4	2.881	323	327	332	rVV	696533	1515377	3.80%	0.474%
5	2.923	332	334	347	rVB2	317236	650872	1.63%	0.204%
6	3.167	364	374	387	rBV2	1255943	3148283	7.89%	0.986%
7	3.807	467	479	489	rBV	268184	649939	1.63%	0.203%
8	3.917	489	497	505	rBV	190984	488618	1.22%	0.153%
9	4.191	527	542	549	rBV2	190874	551072	1.38%	0.173%
10	4.399	566	576	587	rVB	825468	2391104	5.99%	0.749%
11	5.295	701	723	734	rBV5	216138	939174	2.35%	0.294%
12	5.490	743	755	762	rBV	416047	1242975	3.11%	0.389%
13	5.581	762	770	775	rVV4	439312	1544479	3.87%	0.484%
14	5.654	776	782	788	rVV	667312	2019559	5.06%	0.632%
15	5.715	789	792	810	rVB4	365635	1136447	2.85%	0.356%
16	5.923	814	826	837	rBV2	375890	1151877	2.89%	0.361%
17	6.057	837	848	853	rVV	400244	1045106	2.62%	0.327%
18	6.136	853	861	872	rVV	3102563	8153666	20.42%	2.553%
19	6.252	872	880	889	rVV3	383640	1653045	4.14%	0.518%
20	6.349	890	896	902	rVV2	580523	1749625	4.38%	0.548%
21	6.423	902	908	924	rVB5	753931	2765196	6.93%	0.866%
22	6.849	968	978	984	rBV	1009831	2568171	6.43%	0.804%
23	6.935	986	992	1004	rVB3	618296	1826643	4.58%	0.572%
24	7.252	1036	1044	1052	rVB	488346	1027875	2.57%	0.322%
25	7.453	1064	1077	1088	rBV4	808145	2357524	5.91%	0.738%
26	7.953	1147	1159	1166	rBV4	268501	917092	2.30%	0.287%
27	8.050	1166	1175	1187	rVB2	283214	815613	2.04%	0.255%
28	8.172	1187	1195	1197	rBV3	431420	821826	2.06%	0.257%
29	8.209	1197	1201	1205	rVV	588274	1031871	2.58%	0.323%
30	8.569	1254	1260	1269	rVB3	758387	1532796	3.84%	0.480%
31	8.666	1269	1276	1278	rBV2	313960	554322	1.39%	0.174%
32	8.709	1278	1283	1289	rVB	2165239	3462552	8.67%	1.084%
33	8.782	1289	1295	1300	rBV	689736	1112802	2.79%	0.348%
34	9.160	1350	1357	1362	rVB2	277774	468003	1.17%	0.147%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

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\82X112619W.M
Title : SW846 8260

35	9.215	1362	1366	1372	rBV	535408	797174	2.00%	0.250%
36	9.556	1418	1422	1428	rVB3	215976	455991	1.14%	0.143%
37	10.111	1508	1513	1517	rBV	1966225	2629287	6.59%	0.823%
38	10.245	1530	1535	1541	rVB	13068449	17651237	44.21%	5.527%
39	10.355	1546	1553	1559	rVB	29033209	39922088	100.00%	12.500%
40	10.696	1603	1609	1615	rVB	2368809	3068394	7.69%	0.961%
41	11.013	1656	1661	1669	rVB	2210023	2694228	6.75%	0.844%
42	11.135	1676	1681	1685	rBV	1683551	2129723	5.33%	0.667%
43	11.355	1712	1717	1723	rVB	3679059	4345037	10.88%	1.360%
44	11.428	1723	1729	1731	rVV	3290430	4570440	11.45%	1.431%
45	11.452	1731	1733	1737	rVV	4178706	4267267	10.69%	1.336%
46	11.501	1737	1741	1747	rVB	3797821	4511758	11.30%	1.413%
47	11.666	1762	1768	1777	rBV	6279914	7973547	19.97%	2.497%
48	11.806	1785	1791	1796	rVB	28713674	34746817	87.04%	10.879%
49	12.074	1830	1835	1839	rBV	2176538	2783874	6.97%	0.872%
50	12.141	1839	1846	1851	rVB	13091395	16230458	40.66%	5.082%
51	12.294	1866	1871	1878	rBV	20385192	25718166	64.42%	8.052%
52	12.367	1878	1883	1890	rVB3	1832207	2821121	7.07%	0.883%
53	12.464	1893	1899	1903	rBV2	1520001	2092944	5.24%	0.655%
54	12.513	1903	1907	1914	rVB2	1424294	1931353	4.84%	0.605%
55	12.580	1914	1918	1920	rBV	2058089	2443282	6.12%	0.765%
56	12.604	1920	1922	1927	rVB	1477575	1620674	4.06%	0.507%
57	12.665	1927	1932	1935	rBV	5116468	6190497	15.51%	1.938%
58	12.696	1935	1937	1941	rVB	1243829	1337304	3.35%	0.419%
59	12.751	1941	1946	1954	rBV	5987402	7804961	19.55%	2.444%
60	12.897	1965	1970	1975	rVB2	1188010	1525536	3.82%	0.478%
61	12.970	1975	1982	1986	rBV	4276966	5420234	13.58%	1.697%
62	13.013	1986	1989	1995	rVB	3896307	4511452	11.30%	1.413%
63	13.074	1995	1999	2006	rBV3	402147	816844	2.05%	0.256%
64	13.190	2009	2018	2019	rBV4	332852	491524	1.23%	0.154%
65	13.220	2019	2023	2028	rVB	3647598	4127208	10.34%	1.292%
66	13.312	2033	2038	2039	rVV	399257	581553	1.46%	0.182%
67	13.342	2039	2043	2048	rVV2	7858005	10930319	27.38%	3.422%
68	13.397	2048	2052	2054	rVV3	706733	1025008	2.57%	0.321%
69	13.421	2054	2056	2061	rVV	410655	477557	1.20%	0.150%
70	13.476	2061	2065	2073	rVB	1640463	2042996	5.12%	0.640%
71	13.556	2073	2078	2081	rBV2	456535	641478	1.61%	0.201%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

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\82X112619W.M
Title : SW846 8260

72	13.598	2081	2085	2088	rVV	1317667	1875646	4.70%	0.587%
73	13.635	2088	2091	2097	rVV4	1273959	2303638	5.77%	0.721%
74	13.696	2097	2101	2102	rVV	649105	877384	2.20%	0.275%
75	13.714	2102	2104	2110	rVB2	1087382	1647374	4.13%	0.516%
76	13.830	2116	2123	2132	rBV	9884951	12023584	30.12%	3.765%
77	13.921	2132	2138	2142	rVB	608858	816473	2.05%	0.256%
78	13.976	2142	2147	2151	rBV2	461159	656706	1.64%	0.206%
79	14.025	2151	2155	2158	rVB	386407	439561	1.10%	0.138%
80	14.129	2167	2172	2176	rBV	815077	996702	2.50%	0.312%
81	14.269	2190	2195	2200	rBV	687866	1090910	2.73%	0.342%
82	14.372	2208	2212	2217	rBV	439195	596934	1.50%	0.187%
83	14.427	2217	2221	2226	rVB	536822	671477	1.68%	0.210%
84	14.689	2260	2264	2274	rVB	1291123	1767213	4.43%	0.553%
85	14.836	2282	2288	2298	rBV	2094494	2789000	6.99%	0.873%

Sum of corrected areas: 319383698

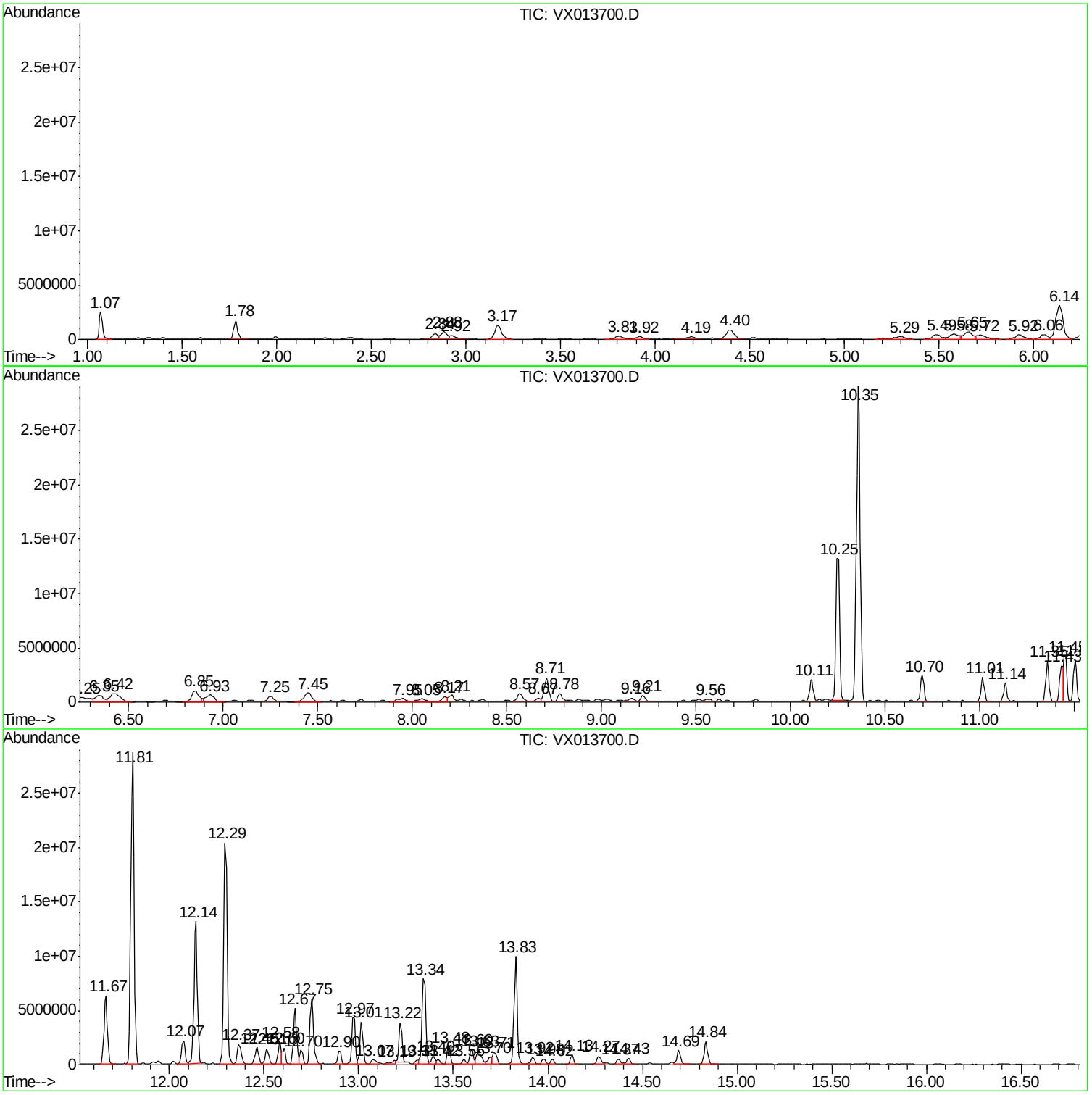
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.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\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

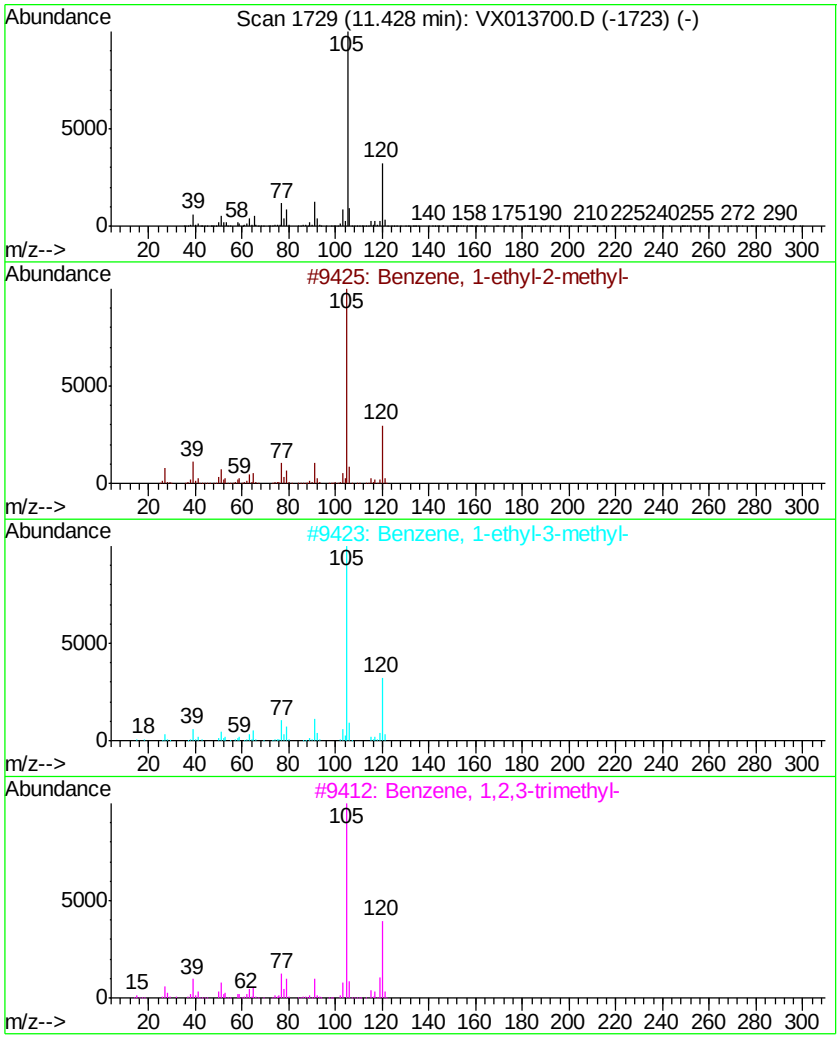
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

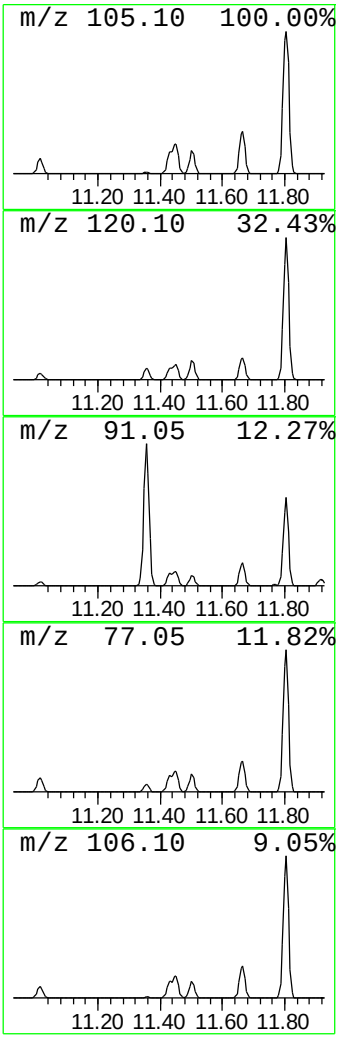
****

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	82.09 ug/l	4570440	1,4-Dichlorobenzene-d4	12.07

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-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

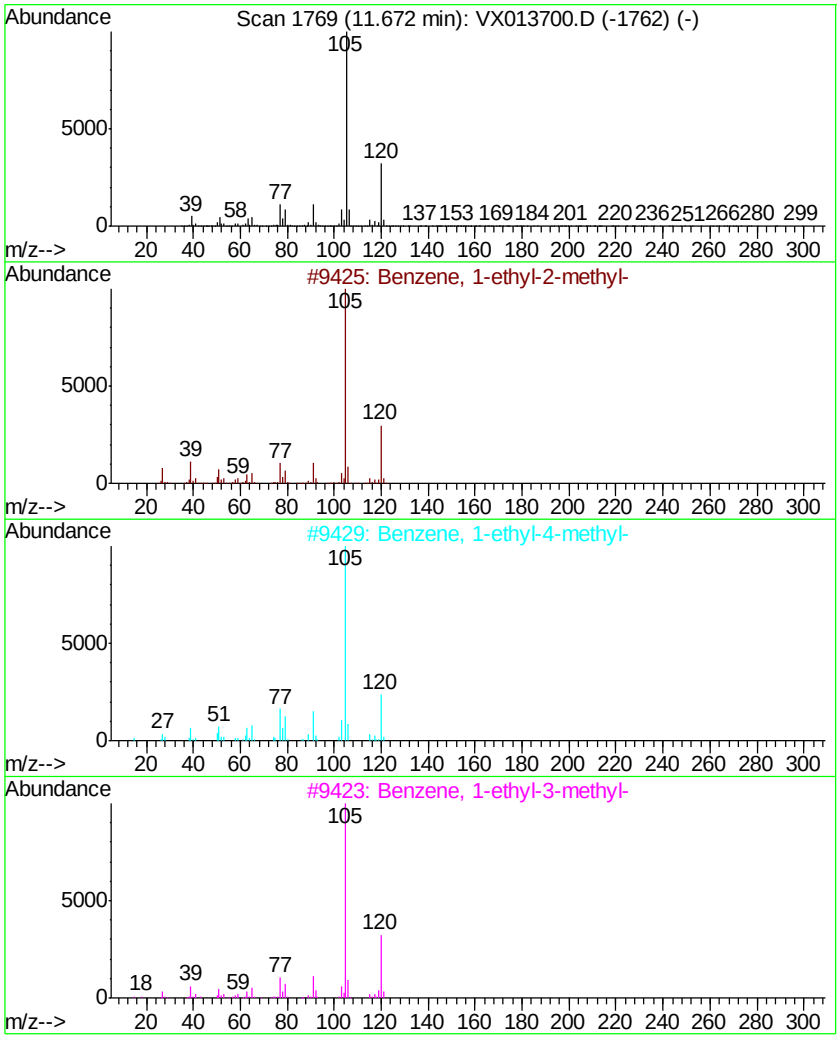
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

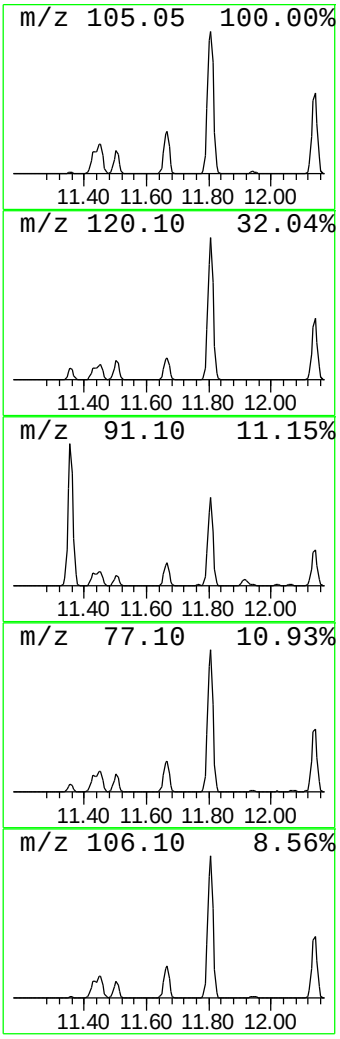
****

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	143.21 ug/l	7973550	1,4-Dichlorobenzene-d4	12.07

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-4-methyl-	120	C9H12	000622-96-8	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

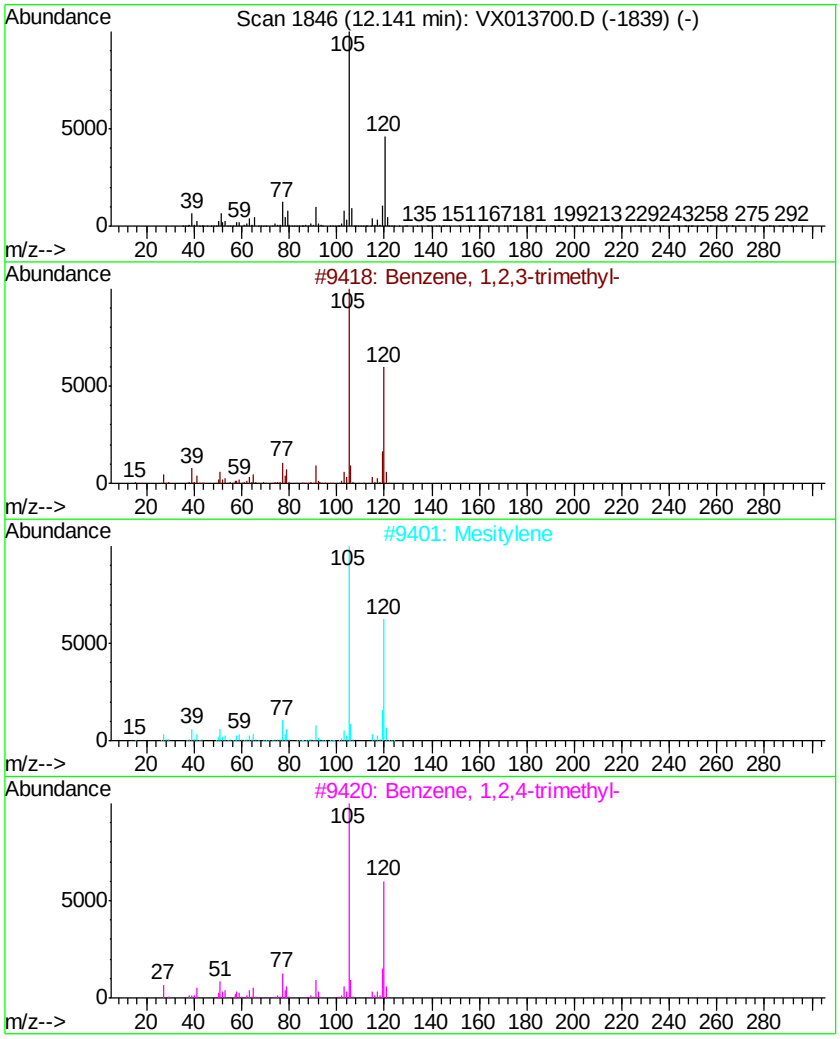
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

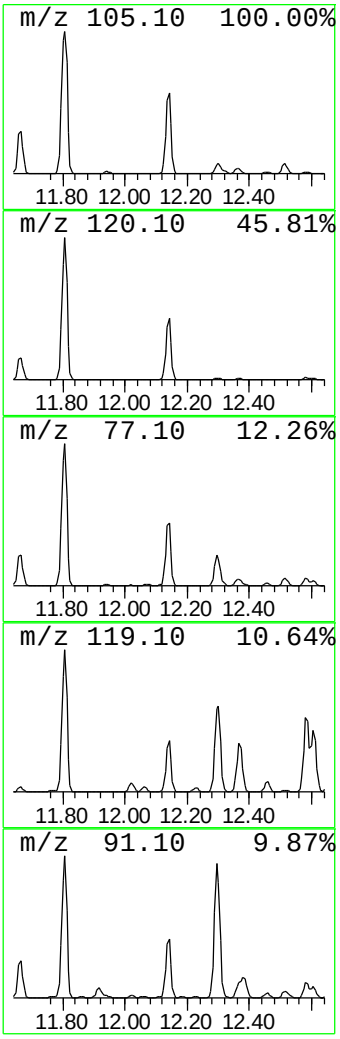
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	291.51 ug/l	16230500	1,4-Dichlorobenzene-d4	12.07

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	94
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\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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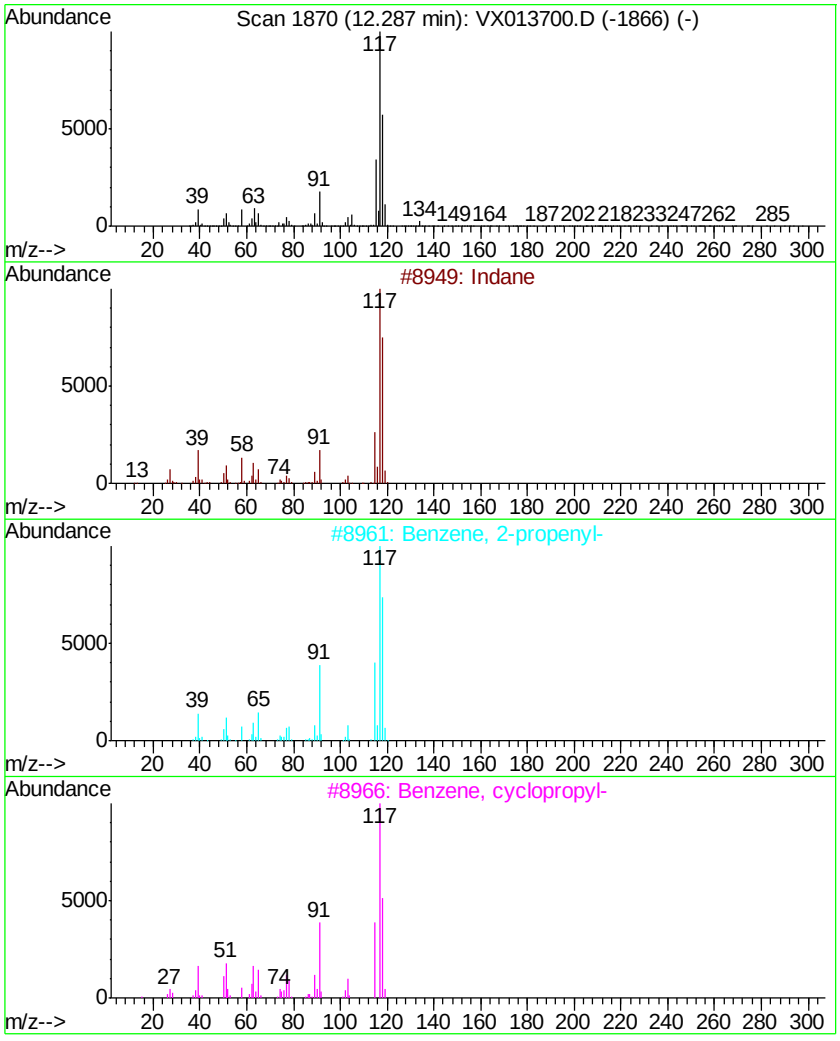
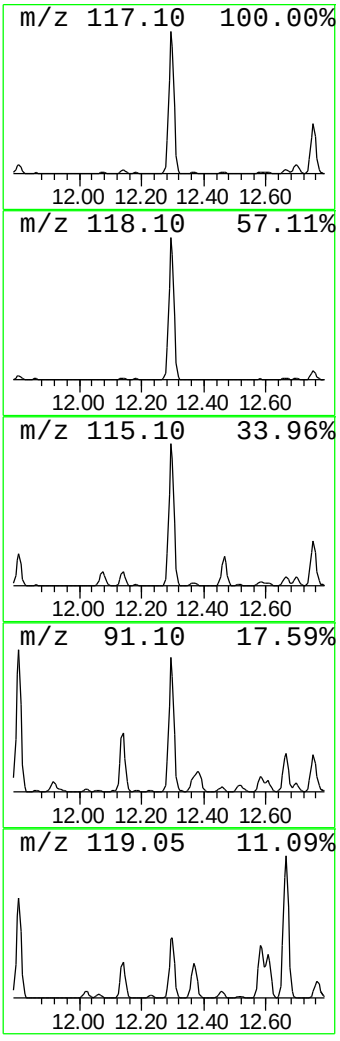
Peak Number 4 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	461.91 ug/l	25718200	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4	Deltacyclene	118	C9H10	007785-10-6	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	58





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

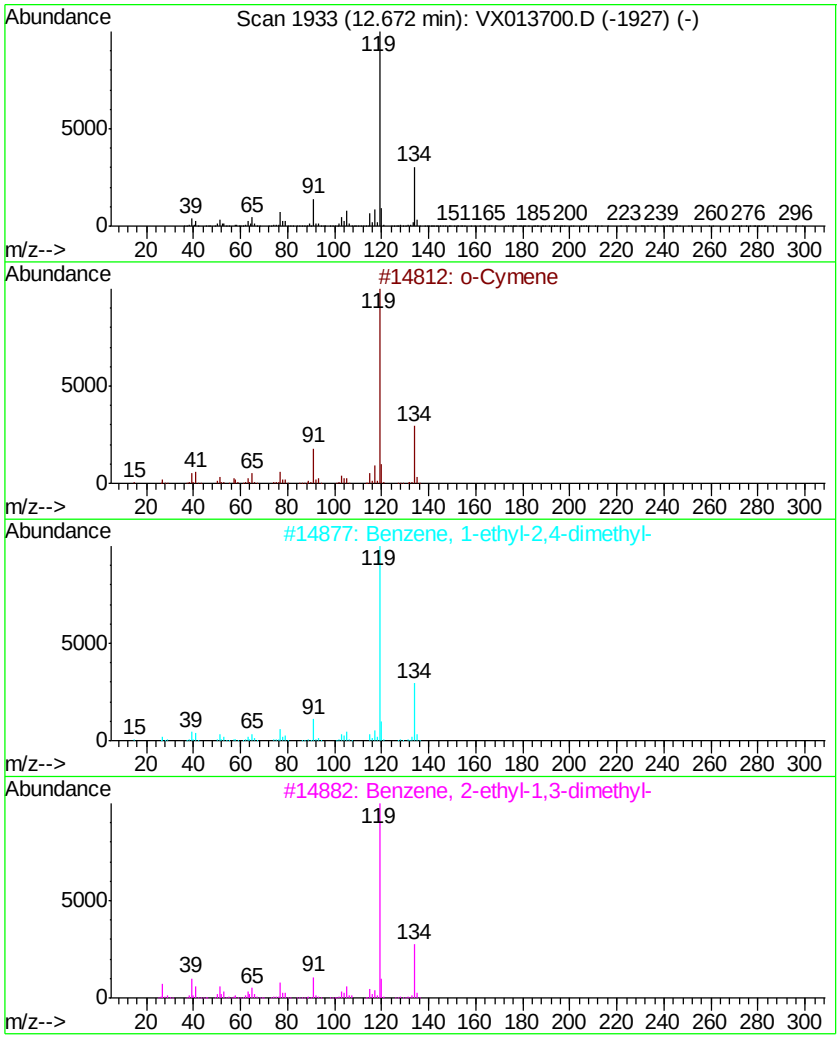
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

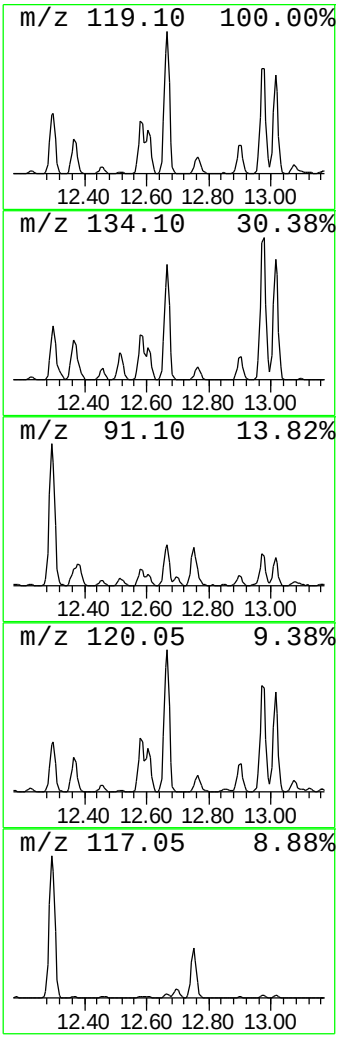
****

Peak Number 5 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	111.19 ug/l	6190500	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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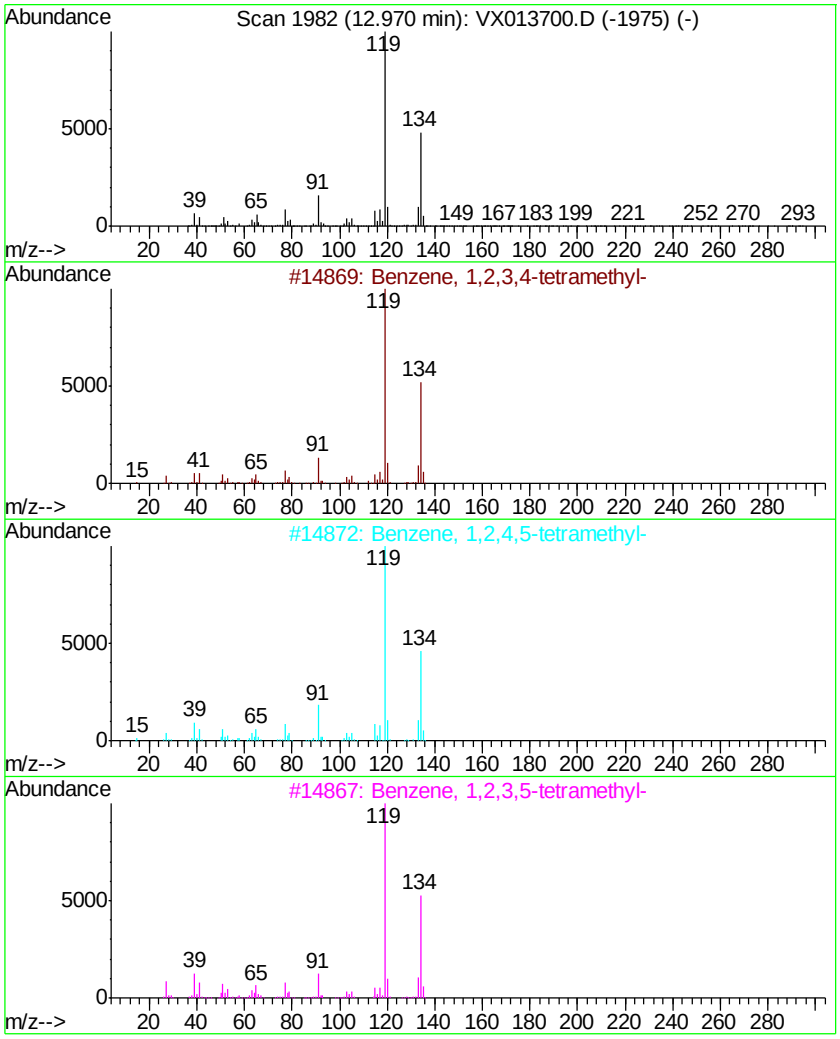
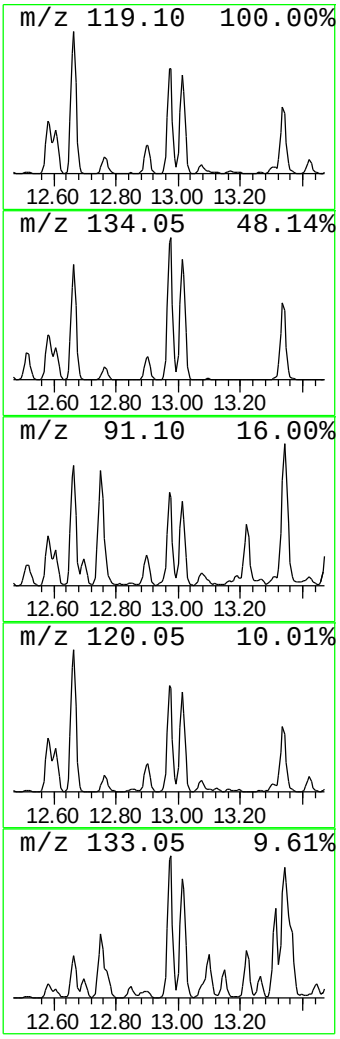
Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	97.35 ug/l	5420230	1,4-Dichlorobenzene-d4	12.07

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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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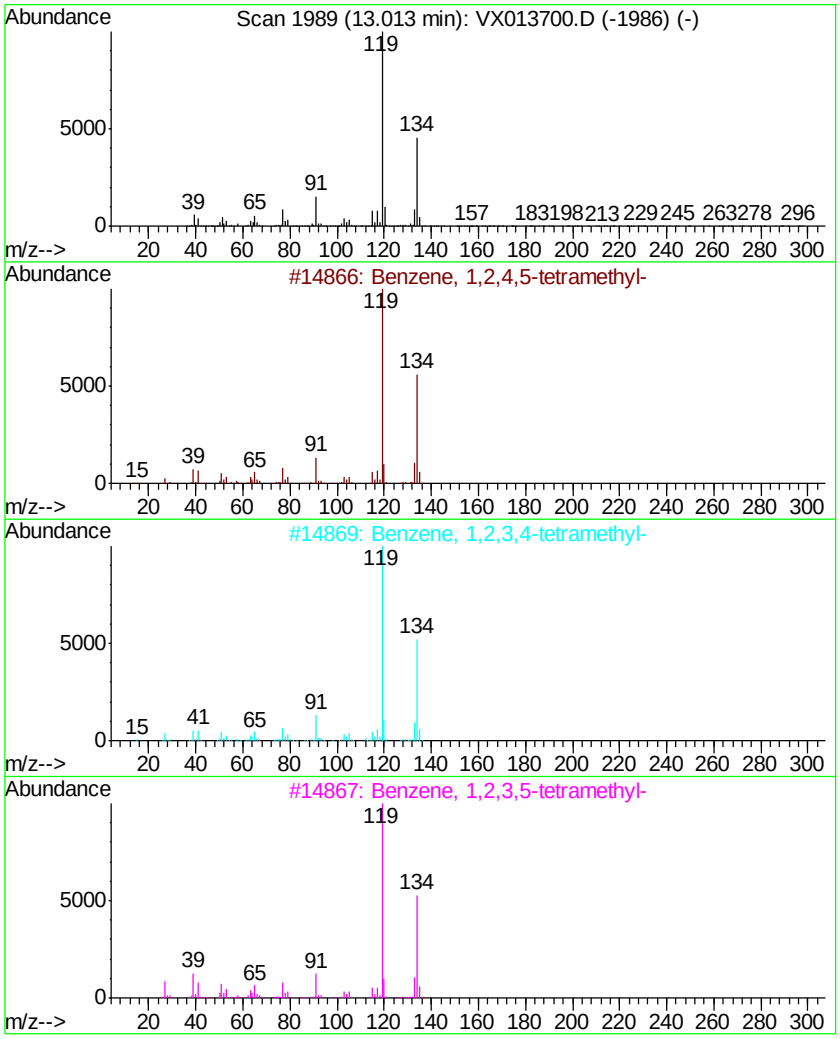
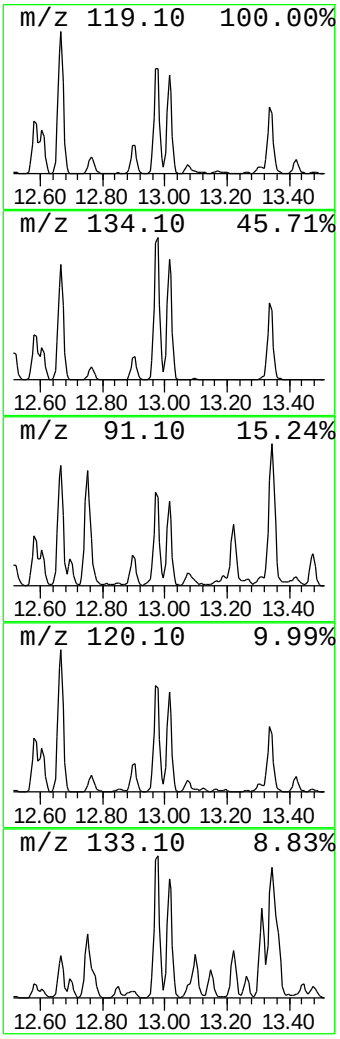
Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	81.03 ug/l	4511450	1,4-Dichlorobenzene-d4	12.07

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	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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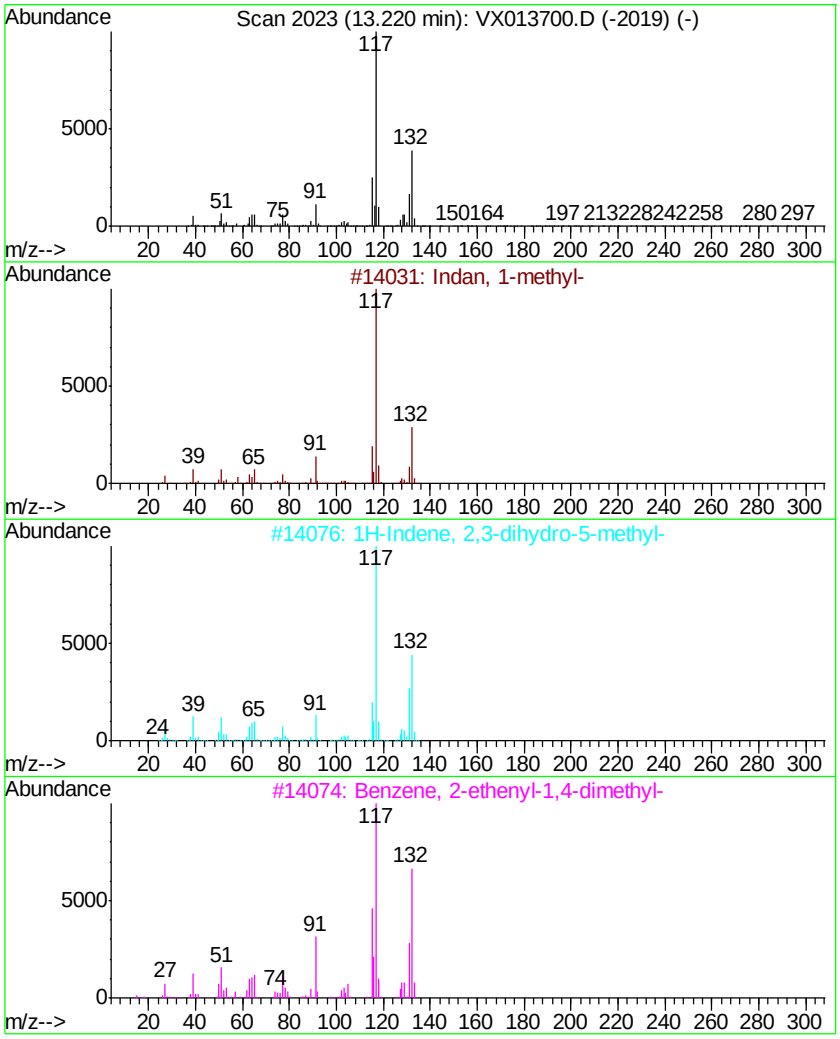
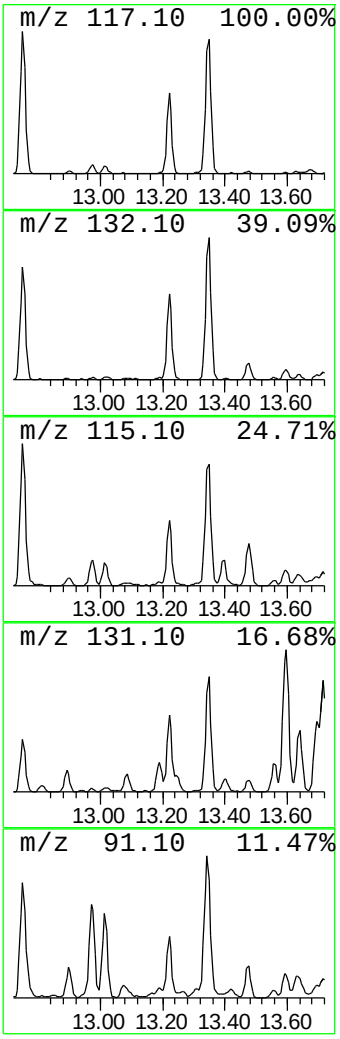
Peak Number 9 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	74.13 ug/l	4127210	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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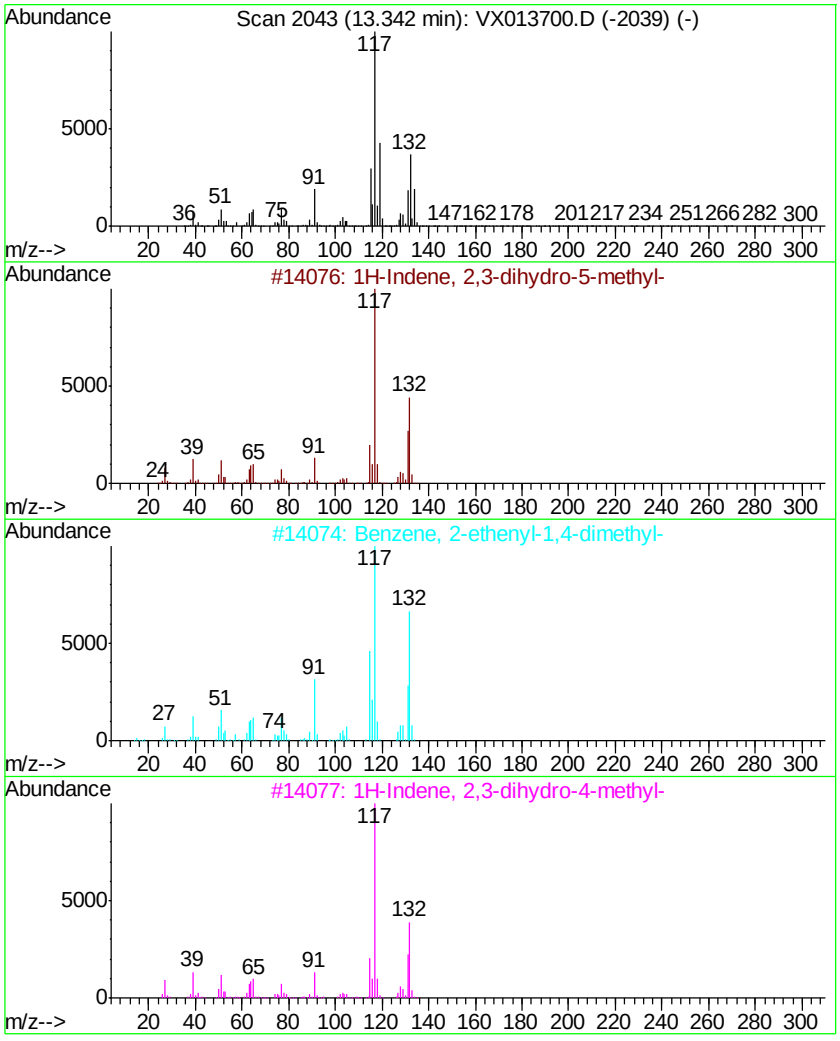
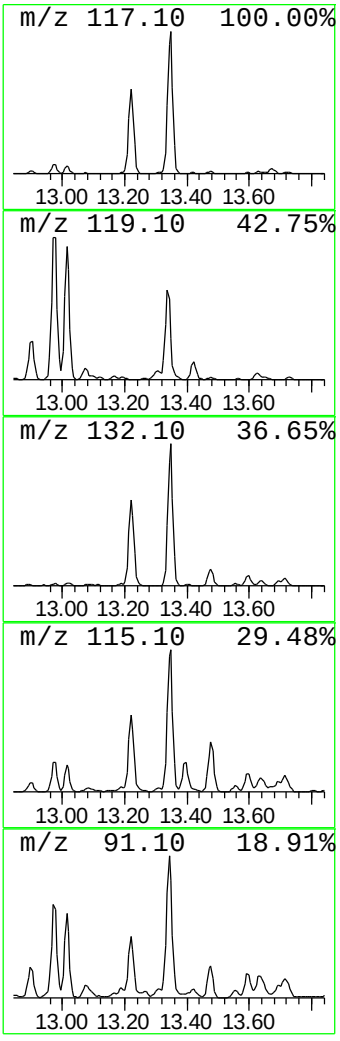
Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	196.32 ug/l	10930300	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	92
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013700.D
Acq On : 29 Nov 2019 14:50
Operator : JC/SP
Sample : K6027-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.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-...	11.43	82.1	ug/l	4570440	4	12.07	2783870	50.0
Benzene, 1-ethyl-...	11.67	143.2	ug/l	7973550	4	12.07	2783870	50.0
Benzene, 1,2,3-tr...	12.14	291.5	ug/l	16230500	4	12.07	2783870	50.0
Indane	12.29	461.9	ug/l	25718200	4	12.07	2783870	50.0
o-Cymene	12.67	111.2	ug/l	6190500	4	12.07	2783870	50.0
Benzene, 1,2,3,4-...	12.97	97.3	ug/l	5420230	4	12.07	2783870	50.0
Benzene, 1,2,4,5-...	13.01	81.0	ug/l	4511450	4	12.07	2783870	50.0
Indan, 1-methyl-	13.22	74.1	ug/l	4127210	4	12.07	2783870	50.0
1H-Indene, 2,3-di...	13.34	196.3	ug/l	10930300	4	12.07	2783870	50.0