LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

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\82X091118W.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	2.446	207	212	222	rVB2	7178	14832	1.24%	0.118%
2	2.617	234	240	249	rVB	34116	57545	4.80%	0.459%
3	2.836	271	276	287	rBV4	5806	16458	1.37%	0.131%
4	4.300	507	516	525	rBV3	10673	31195	2.60%	0.249%
5	5.507	700	714	729	rBV2	142108	392630	32.76%	3.134%
6	5.659	729	739	761	rVB	190202	614519	51.27%	4.906%
7	6.068	795	806	822	rBV	134732	349806	29.18%	2.792%
8	6.342	839	851	864	rVV	333584	971129	81.02%	7.752%
9	6.860	926	936	952	rBV2	296802	690030	57.57%	5.508%
10	7.458	1024	1034	1044	rVV4	9335	23804	1.99%	0.190%
11	7.573	1047	1053	1057	rVV2	9415	19357	1.61%	0.155%
12	7.634	1057	1063	1065	rVV4	15853	32144	2.68%	0.257%
13	7.677	1065	1070	1086	rVB	43520	100307	8.37%	0.801%
14	8.055	1115	1132	1141	rBV	382749	723320	60.35%	5.774%
15	8.177	1143	1152	1160	rBV	622703	1198594	100.00%	9.568%
16	8.256	1160	1165	1170	rVB	19896	32745	2.73%	0.261%
17	8.713	1234	1240	1250	rVB	692721	1102619	91.99%	8.802%
18	9.043	1288	1294	1300	rVB2	15354	25534	2.13%	0.204%
19	9.165	1307	1314	1319	rVB2	7779	12212	1.02%	0.097%
20	9.677	1393	1398	1403	rBV2	7600	12048	1.01%	0.096%
21	10.116	1464	1470	1478	rBV	601332	796944	66.49%	6.362%
22	10.189	1478	1482	1486	rVV	11517	17267	1.44%	0.138%
23	10.250	1486	1492	1497	rVB	9653	15263	1.27%	0.122%
24	10.341	1502	1507	1516	rBV7	9036	25637	2.14%	0.205%
25	10.652	1547	1558	1564	rBV6	7900	22410	1.87%	0.179%
26	11.018	1611	1618	1622	rVV4	24593	50401	4.21%	0.402%
27	11.140	1632	1638	1643	rBV	569709	726532	60.62%	5.800%
28	11.280	1657	1661	1666	rVB3	9713	13848	1.16%	0.111%
29	11.359	1670	1674	1684	rBV	38522	51758	4.32%	0.413%
30	11.676	1720	1726	1733	rVB3	7423	12720	1.06%	0.102%
31	11.920	1759	1766	1768	rBV	19450	31006	2.59%	0.248%
32	11.945	1768	1770	1776	rVB	44110	55092	4.60%	0.440%
33	12.079	1787	1792	1798	rBV	706351	859971	71.75%	6.865%
34	12.231	1813	1817	1822	rBV	19697	27066	2.26%	0.216%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

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\82X091118W.M
Title : SW846 8260

35	12.304	1822	1829	1834	rVV	97737	118286	9.87%	0.944%
36	12.371	1835	1840	1850	rVB2	45560	95600	7.98%	0.763%
37	12.463	1850	1855	1860	rBV	50761	63561	5.30%	0.507%
38	12.670	1879	1889	1892	rBV	160940	200870	16.76%	1.604%
39	12.700	1892	1894	1899	rVV2	39054	57472	4.79%	0.459%
40	12.761	1899	1904	1911	rVV4	93662	181869	15.17%	1.452%
41	12.816	1911	1913	1919	rVV2	11523	22077	1.84%	0.176%
42	12.896	1919	1926	1931	rVB3	45822	75737	6.32%	0.605%
43	12.981	1931	1940	1944	rBV	132245	192547	16.06%	1.537%
44	13.030	1944	1948	1953	rVV3	12337	20506	1.71%	0.164%
45	13.085	1953	1957	1964	rVB2	30023	44980	3.75%	0.359%
46	13.152	1964	1968	1972	rBV2	25675	39095	3.26%	0.312%
47	13.194	1972	1975	1978	rVV2	43629	55767	4.65%	0.445%
48	13.249	1982	1984	1989	rVB	14181	16692	1.39%	0.133%
49	13.347	1995	2000	2007	rVB2	289264	399188	33.30%	3.187%
50	13.481	2016	2022	2028	rVB	118900	153587	12.81%	1.226%
51	13.566	2032	2036	2039	rBV2	15699	25953	2.17%	0.207%
52	13.603	2039	2042	2045	rVB	21999	24360	2.03%	0.194%
53	13.639	2045	2048	2052	rBV2	60975	85023	7.09%	0.679%
54	13.700	2052	2058	2059	rBV2	41483	59426	4.96%	0.474%
55	13.725	2059	2062	2070	rVB2	67841	108309	9.04%	0.865%
56	13.810	2070	2076	2078	rBV	25131	34838	2.91%	0.278%
57	13.841	2078	2081	2082	rBV	19026	23916	2.00%	0.191%
58	13.914	2088	2093	2098	rVB	54483	66677	5.56%	0.532%
59	13.987	2098	2105	2109	rBV2	87793	125912	10.50%	1.005%
60	14.030	2109	2112	2116	rVV2	26263	36995	3.09%	0.295%
61	14.133	2124	2129	2131	rBV2	15224	20687	1.73%	0.165%
62	14.164	2131	2134	2139	rVB	19601	22972	1.92%	0.183%
63	14.286	2146	2154	2161	rVB2	59960	127613	10.65%	1.019%
64	14.377	2166	2169	2173	rVB	26325	31175	2.60%	0.249%
65	14.432	2173	2178	2182	rVB2	26080	34725	2.90%	0.277%
66	14.475	2182	2185	2190	rVB	11960	15651	1.31%	0.125%
67	14.542	2190	2196	2205	rBV2	124787	170807	14.25%	1.364%
68	14.700	2214	2222	2225	rBV	106182	154186	12.86%	1.231%
69	14.731	2225	2227	2231	rVV2	22885	26889	2.24%	0.215%
70	14.840	2240	2245	2250	rVV	186978	246477	20.56%	1.968%
71	14.907	2254	2256	2261	rVV5	9676	12123	1.01%	0.097%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

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\82X091118W.M
Title : SW846 8260

72	14.981	2261	2268	2274	rVB8	7109	19089	1.59%	0.152%
73	15.249	2307	2312	2320	rVB2	14436	26125	2.18%	0.209%
74	15.548	2357	2361	2366	rVV	18640	28029	2.34%	0.224%
75	15.688	2376	2384	2387	rBV2	26931	51018	4.26%	0.407%
76	15.730	2387	2391	2398	rVB	22702	36435	3.04%	0.291%
77	15.889	2413	2417	2419	rVV2	7680	12019	1.00%	0.096%
78	15.926	2419	2423	2430	rVB	10194	18753	1.56%	0.150%
79	16.175	2459	2464	2473	rVB5	8485	16025	1.34%	0.128%

Sum of corrected areas: 12526784

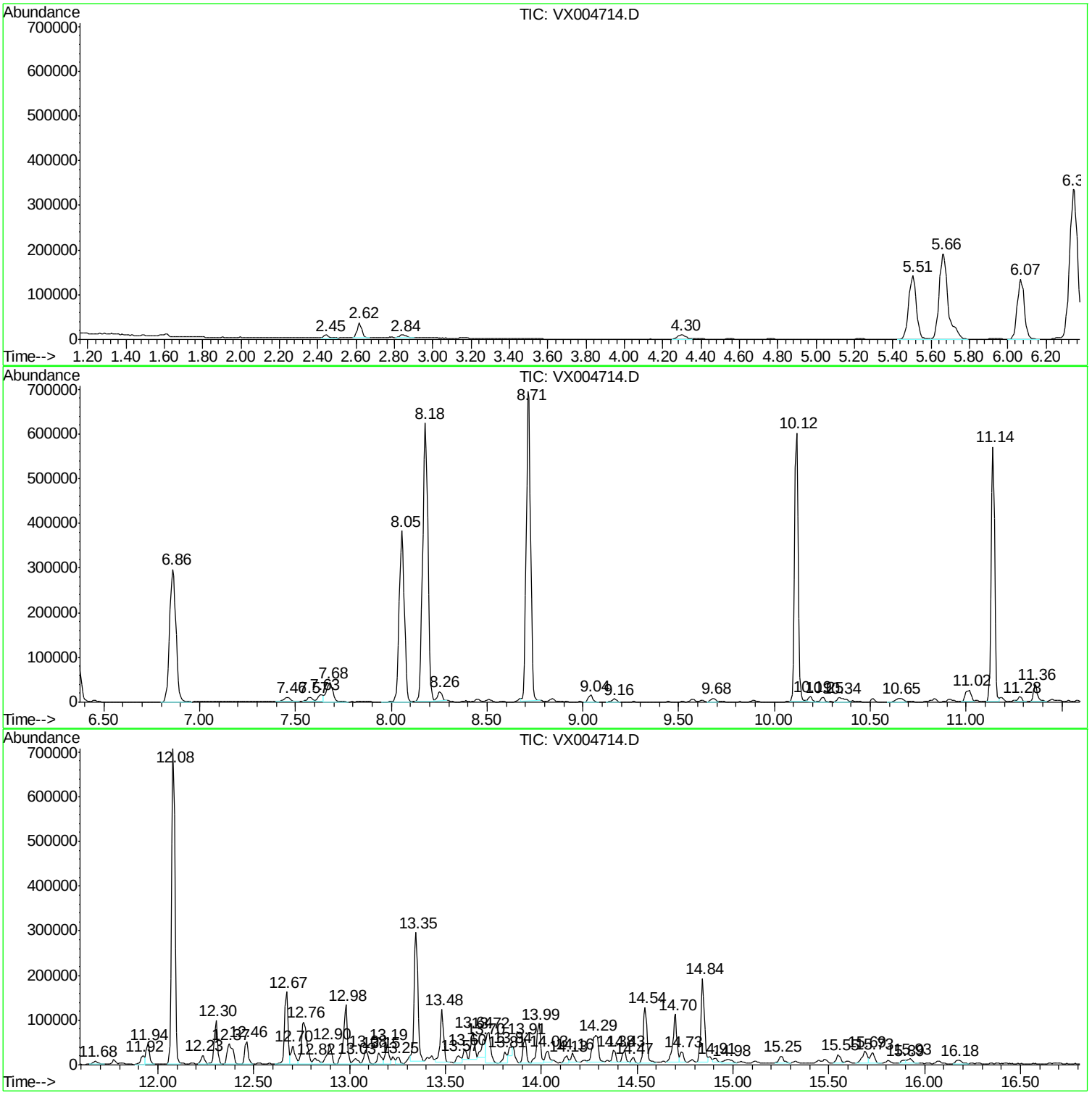
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

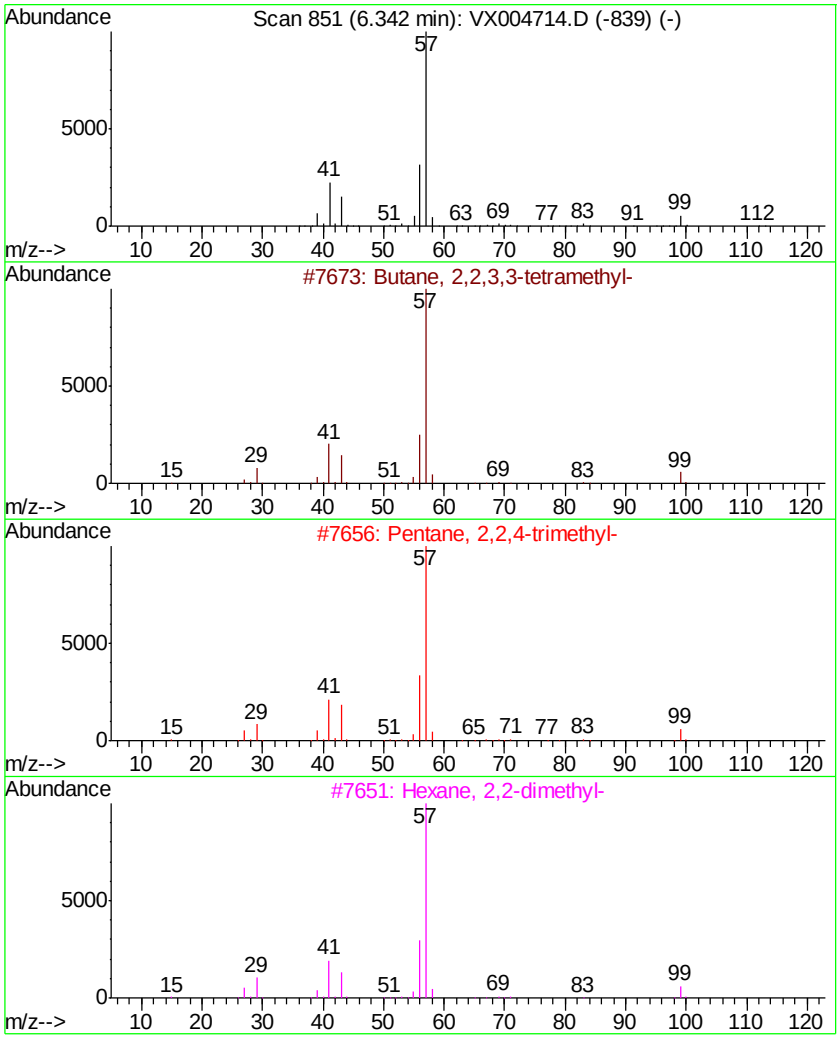
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

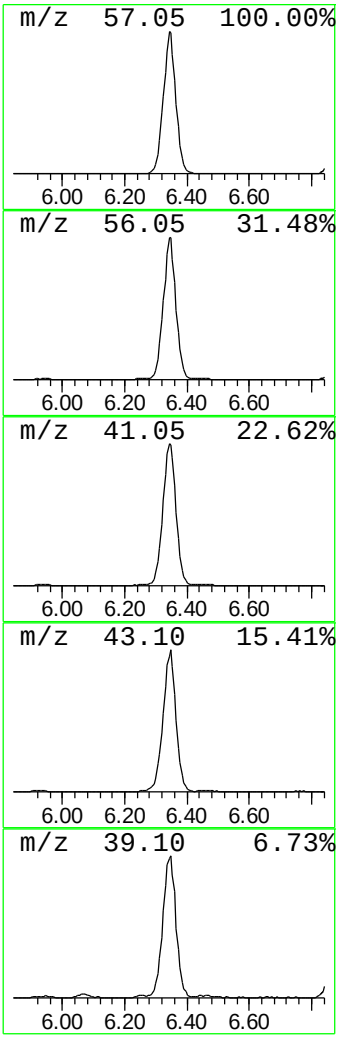
****

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	70.37 ug/l	971129	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
4	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

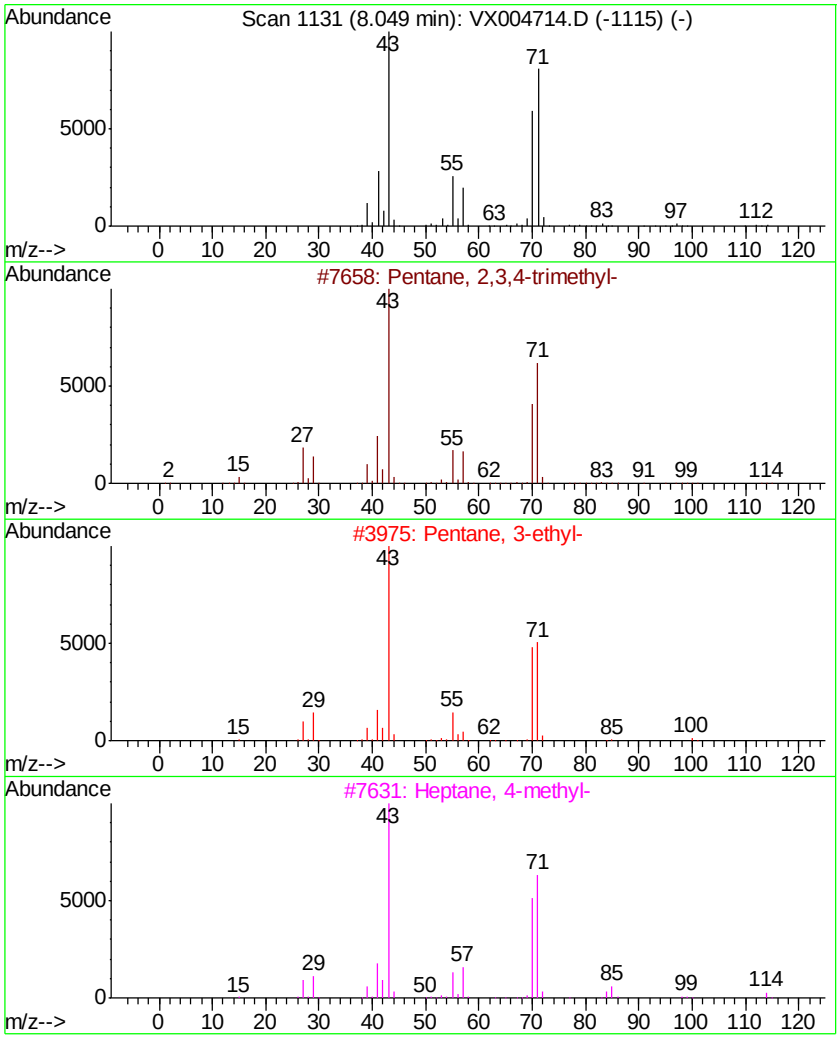
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

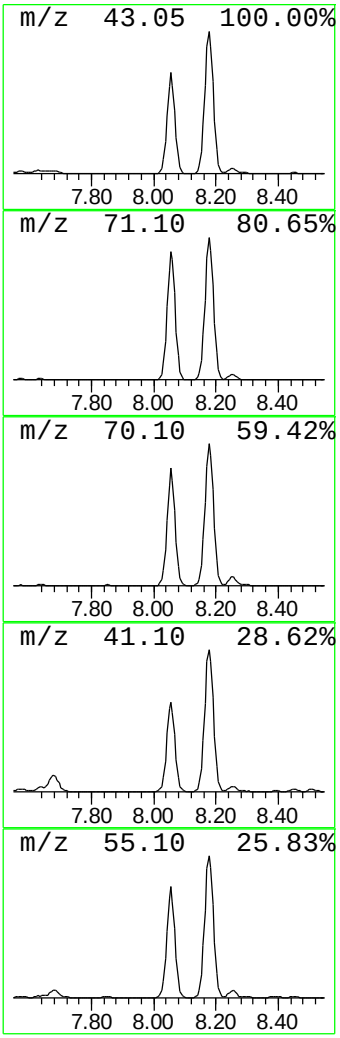
****

Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	52.41 ug/l	723320	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	78
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

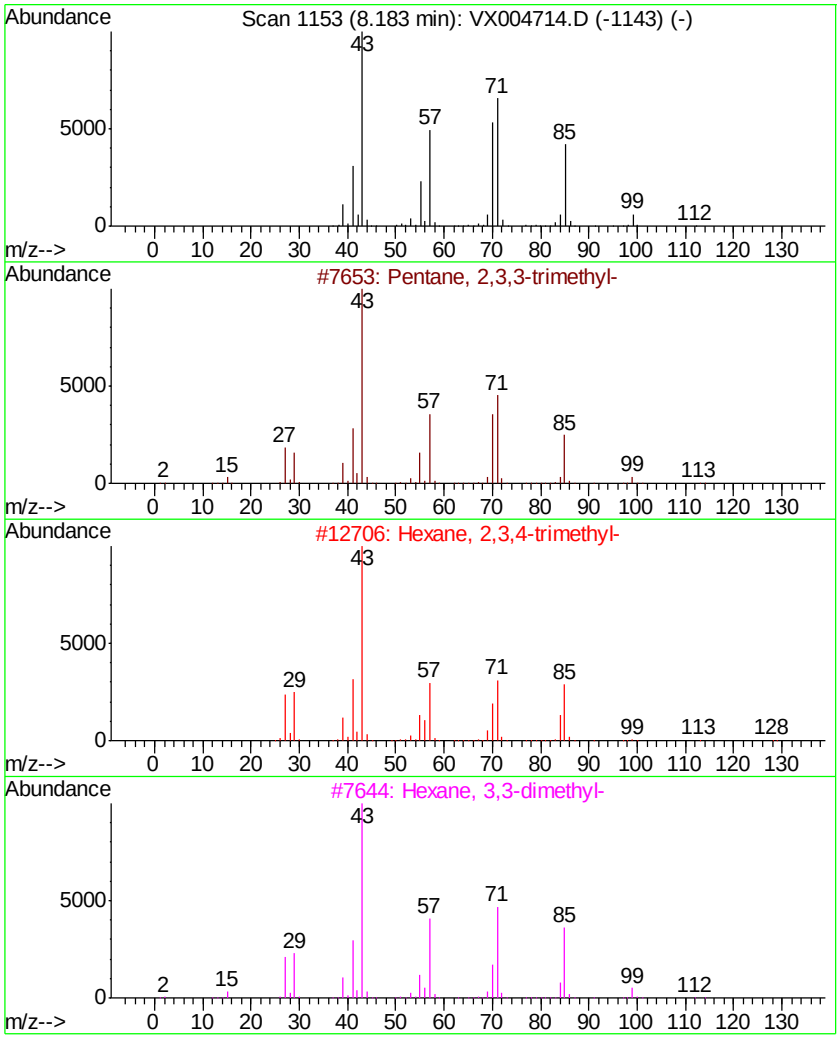
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

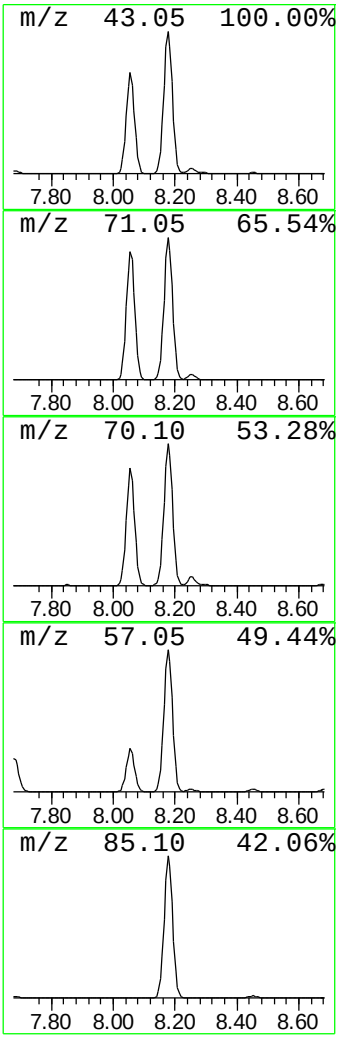
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	86.85 ug/l	1198590	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	50
5	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

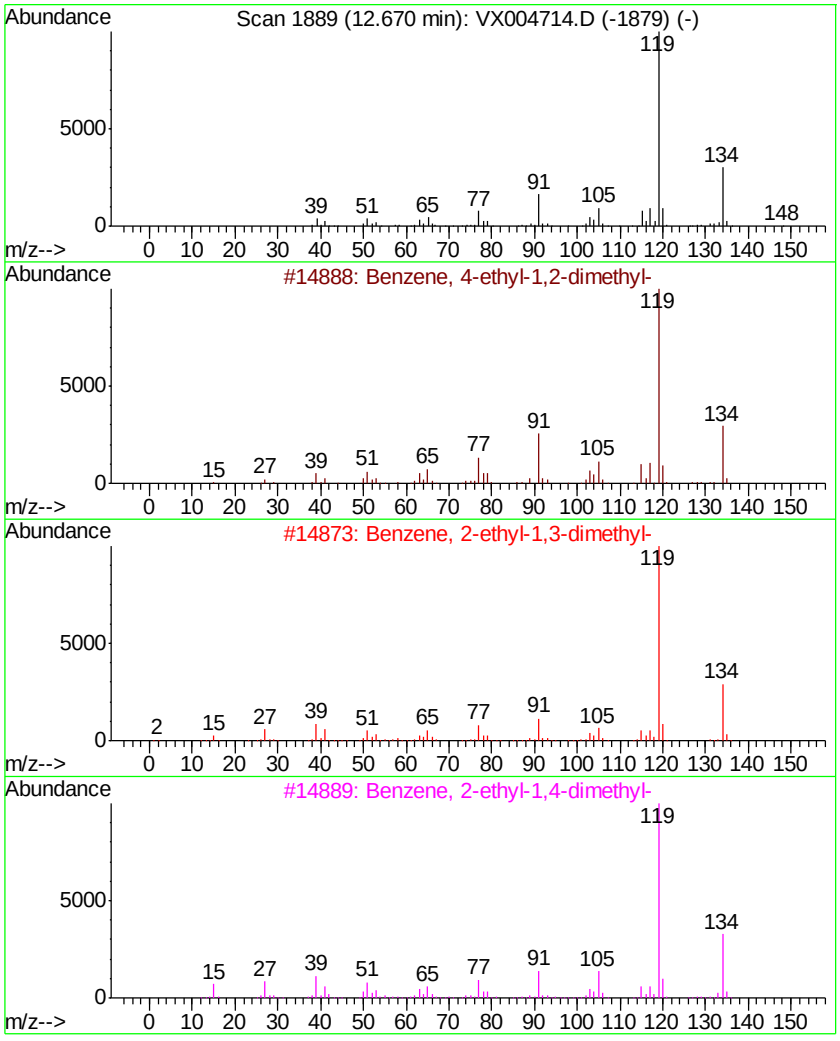
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

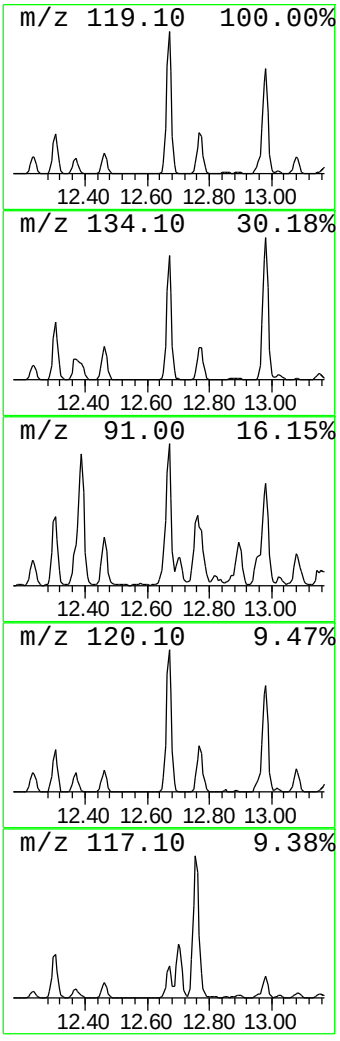
****

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	11.68 ug/l	200870	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

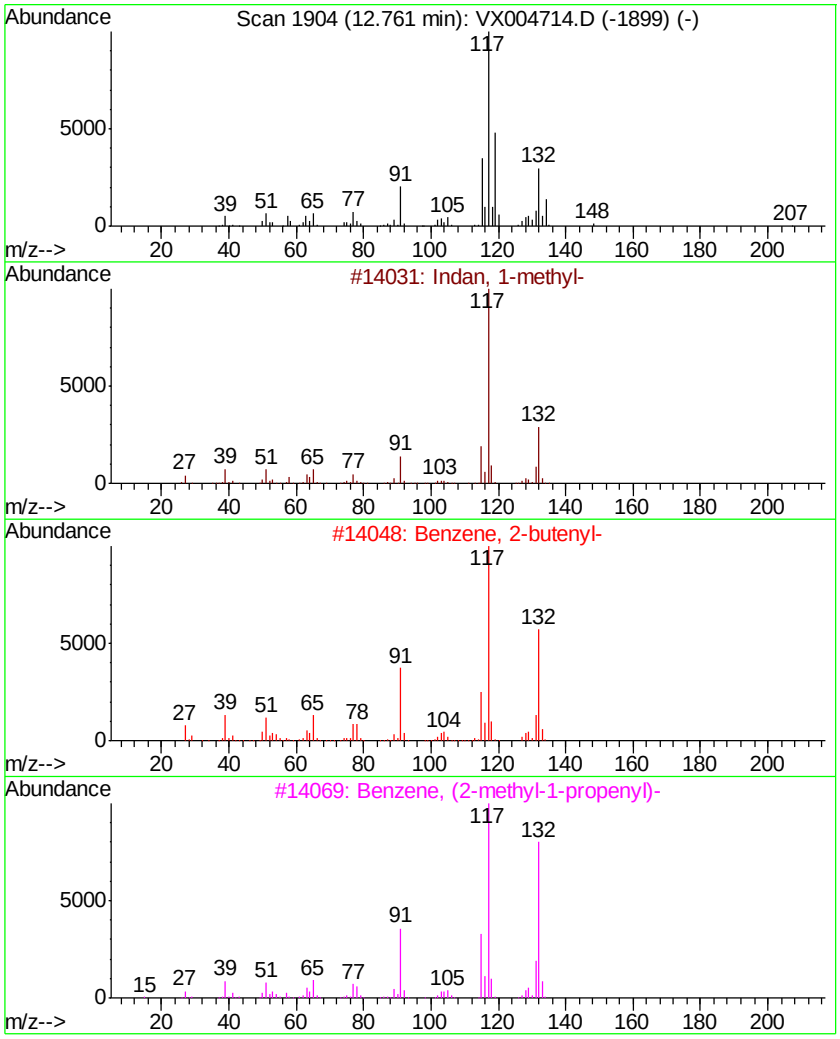
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

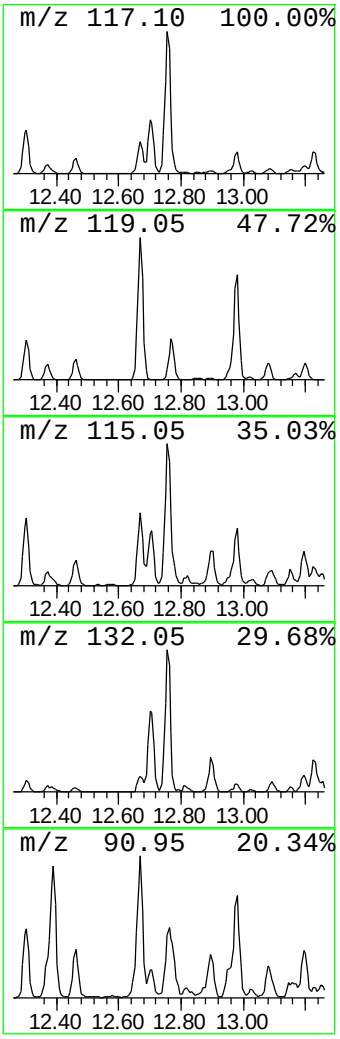
****

Peak Number 5 Indan, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	10.57 ug/l	181869	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	68
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	68
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

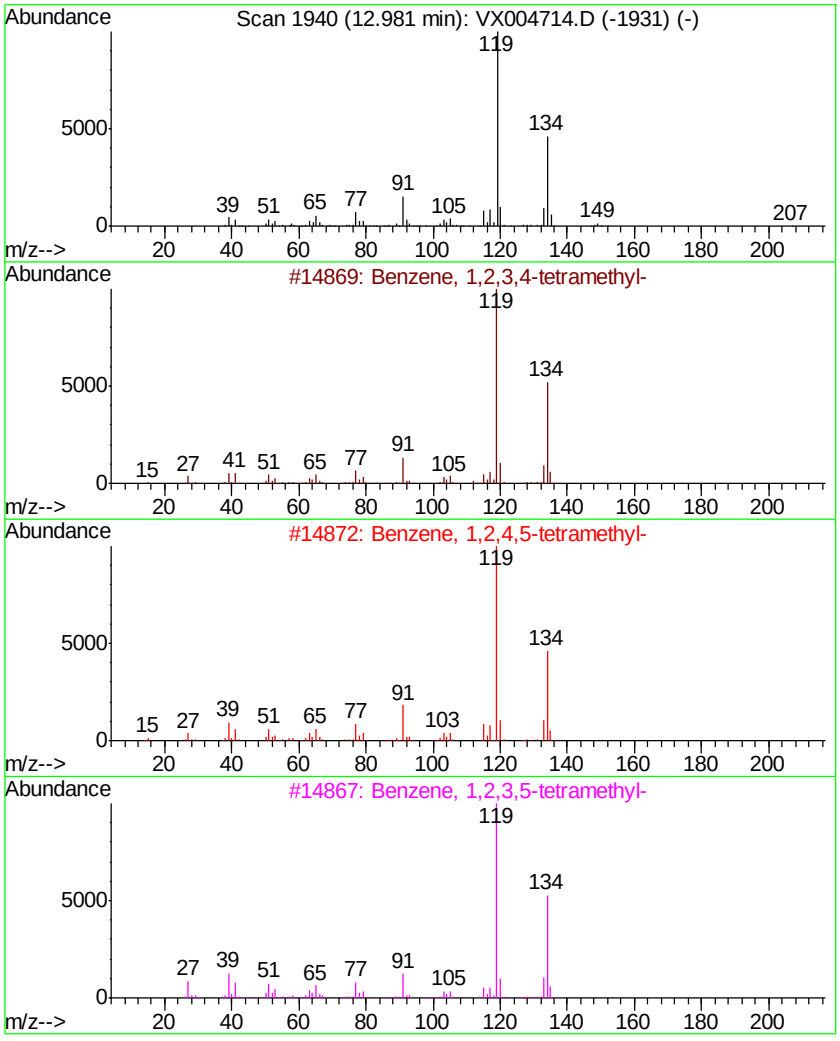
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

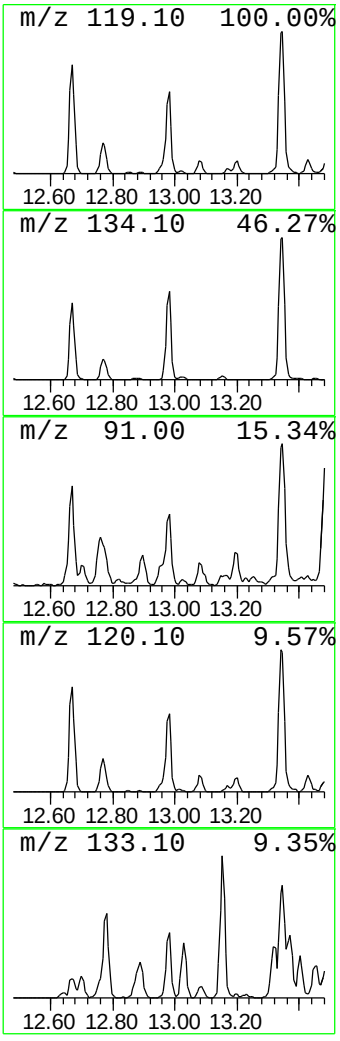
****

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	11.20 ug/l	192547	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	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

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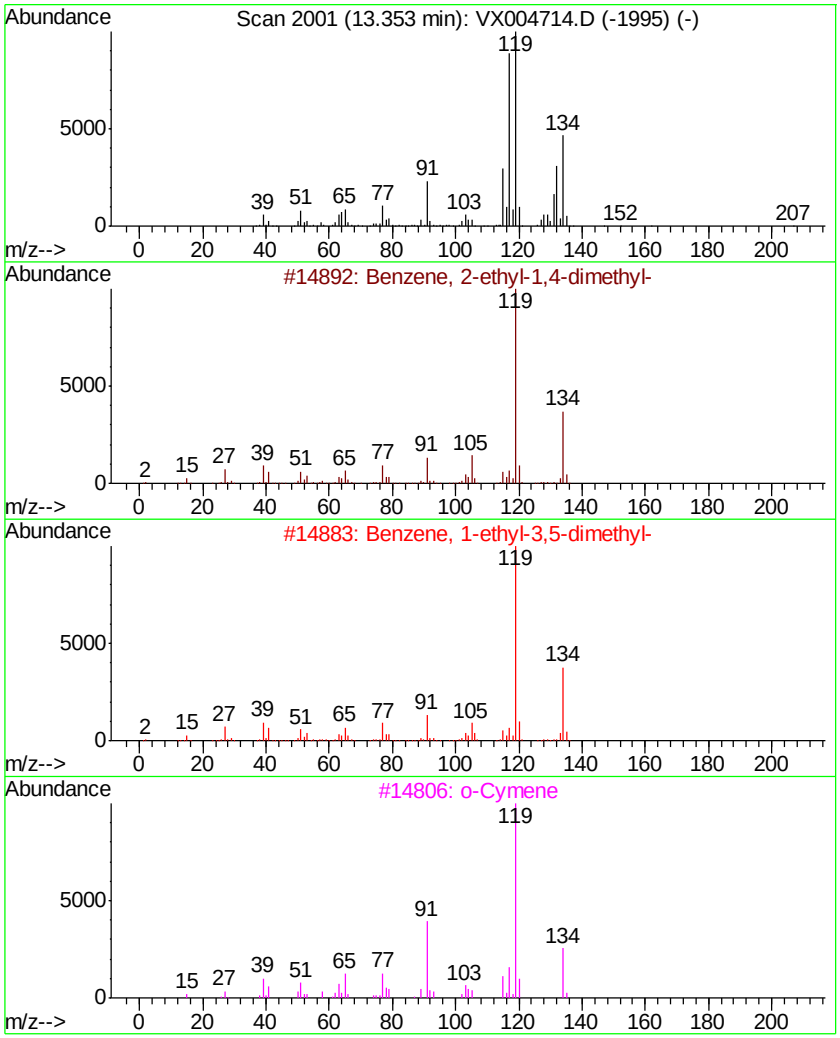
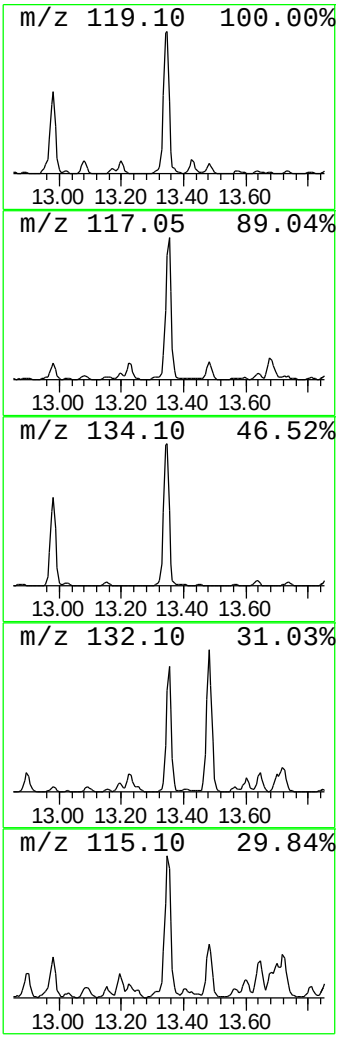
Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	23.21 ug/l	399188	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	76
2	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	76
3	o-Cymene	134	C10H14	000527-84-4	70
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

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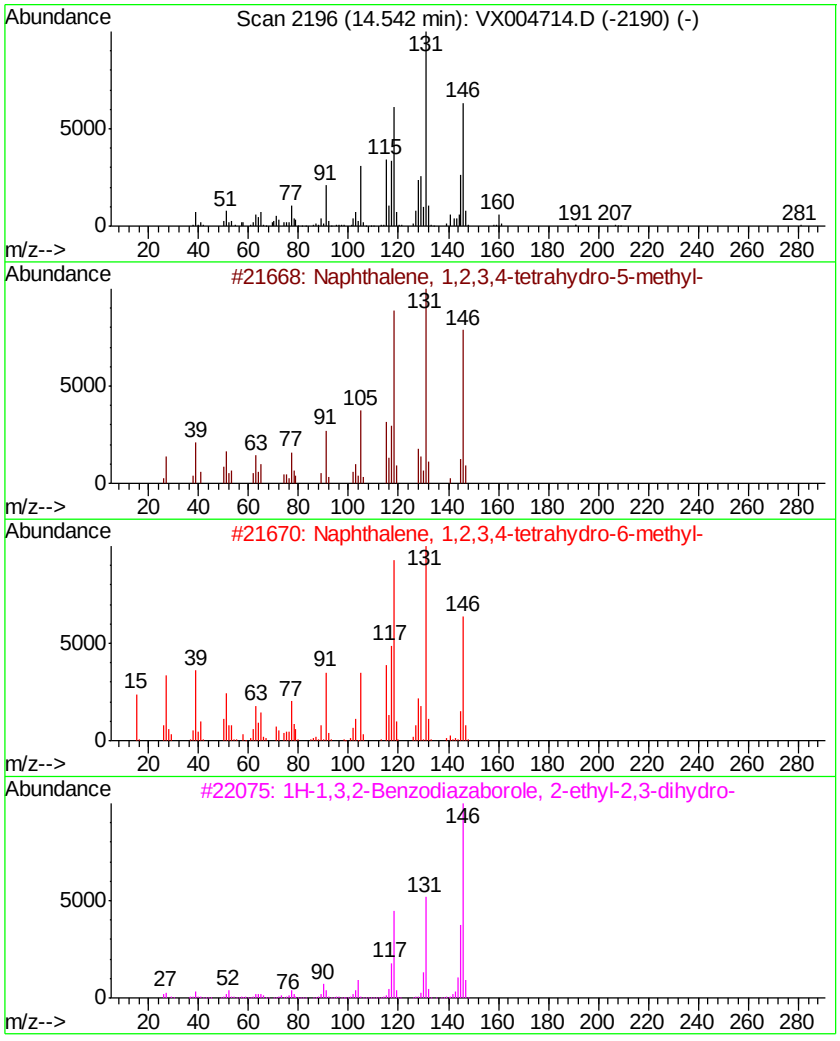
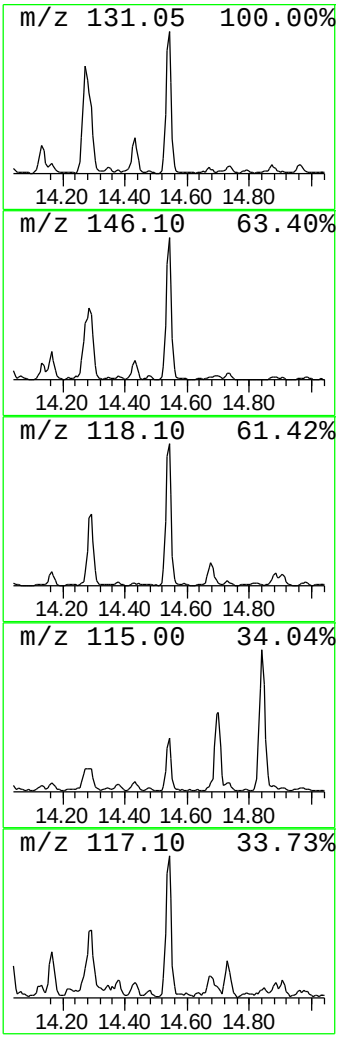
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	9.93 ug/l	170807	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	90
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	83
4	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	68
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

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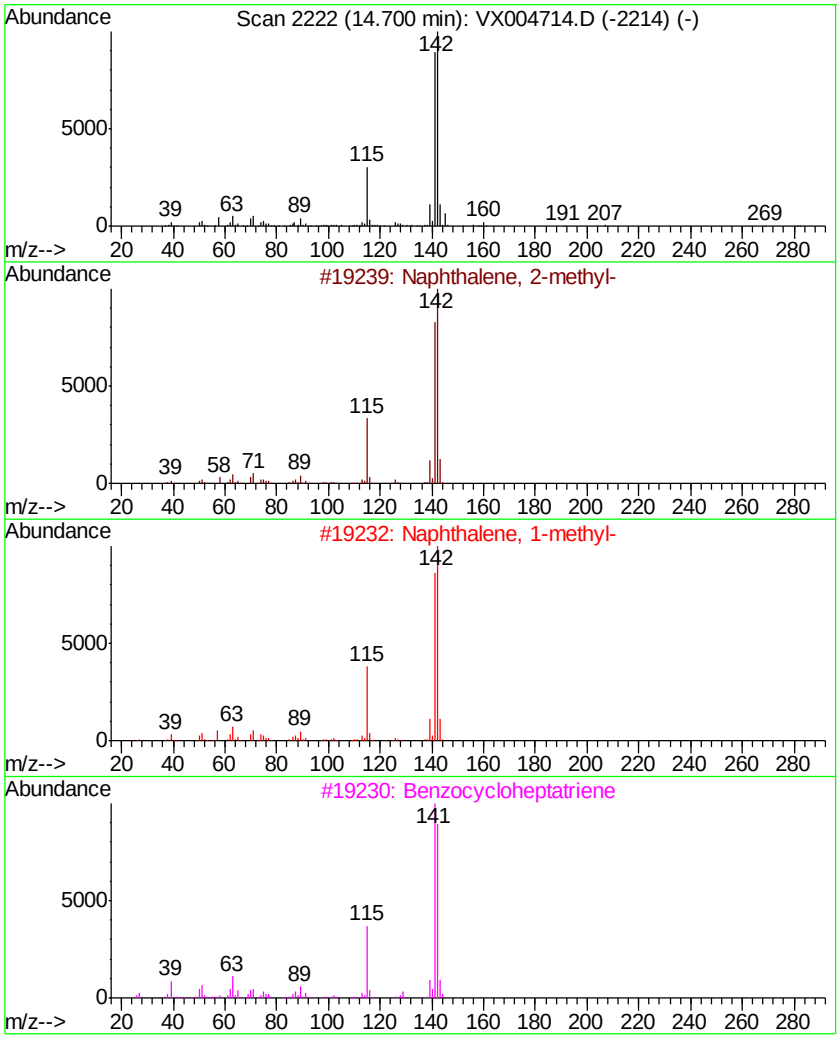
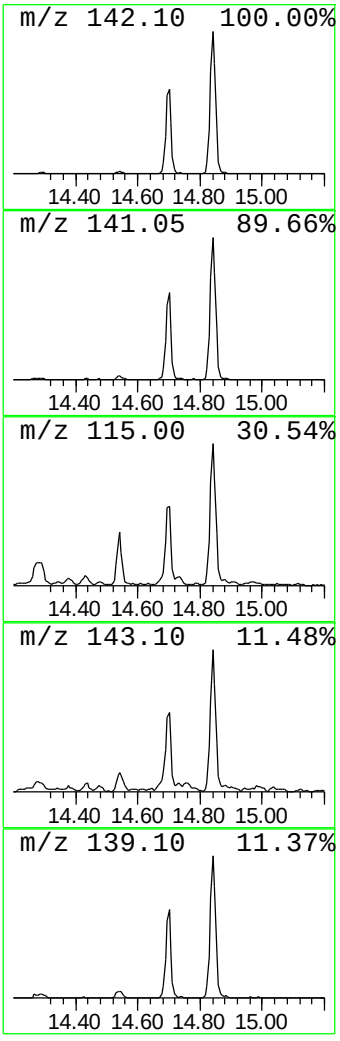
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	8.96 ug/l	154186	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091918\
Data File : VX004714.D
Acq On : 19 Sep 2018 18:59
Operator : JC/MD
Sample : J4916-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FC-MW02SR1-20180911

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.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
Butane, 2,2,3,3-t...	6.34	70.4	ug/l	971129	2	6.86	690030	50.0
Pentane, 2,3,4-tr...	8.05	52.4	ug/l	723320	2	6.86	690030	50.0
Pentane, 2,3,3-tr...	8.18	86.8	ug/l	1198590	2	6.86	690030	50.0
Benzene, 4-ethyl-...	12.67	11.7	ug/l	200870	4	12.08	859971	50.0
Indan, 1-methyl-	12.76	10.6	ug/l	181869	4	12.08	859971	50.0
Benzene, 1,2,3,4-...	12.98	11.2	ug/l	192547	4	12.08	859971	50.0
Benzene, 2-ethyl-...	13.35	23.2	ug/l	399188	4	12.08	859971	50.0
Naphthalene, 1,2,...	14.54	9.9	ug/l	170807	4	12.08	859971	50.0
Naphthalene, 1-me...	14.70	9.0	ug/l	154186	4	12.08	859971	50.0