LSC Area Percent Report

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.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	1.075	76	80	90	rBV	467350	537986	4.28%	0.410%
2	1.404	130	134	139	rBV	254724	253021	2.01%	0.193%
3	1.794	192	198	212	rVB	5034931	6781783	53.92%	5.166%
4	2.002	227	232	246	rVB	1192961	1742256	13.85%	1.327%
5	2.124	246	252	257	rBV	167321	245039	1.95%	0.187%
6	2.270	272	276	286	rVB	263706	412592	3.28%	0.314%
7	2.398	286	297	303	rBV	214570	423789	3.37%	0.323%
8	2.788	350	361	364	rBV4	99775	246574	1.96%	0.188%
9	2.849	364	371	374	rVV	594463	1323960	10.53%	1.008%
10	2.892	374	378	383	rVV	1394407	3008998	23.93%	2.292%
11	2.934	383	385	403	rVB2	670791	1484604	11.80%	1.131%
12	3.172	415	424	441	rVB	1069223	2418330	19.23%	1.842%
13	3.398	453	461	473	rBV2	118544	300469	2.39%	0.229%
14	3.526	473	482	496	rVB	193299	426653	3.39%	0.325%
15	3.666	496	505	512	rBV3	67551	163991	1.30%	0.125%
16	3.751	512	519	524	rVV2	59456	144832	1.15%	0.110%
17	3.818	524	530	540	rVB	239651	556863	4.43%	0.424%
18	3.928	540	548	554	rBV2	164280	416029	3.31%	0.317%
19	3.995	554	559	570	rVV2	101797	344506	2.74%	0.262%
20	4.202	584	593	602	rVB	175113	450290	3.58%	0.343%
21	4.410	616	627	639	rVB	1354177	3851306	30.62%	2.933%
22	4.538	639	648	661	rVB4	72131	228012	1.81%	0.174%
23	5.312	749	775	792	rVB4	142141	573792	4.56%	0.437%
24	5.513	792	808	811	rBV2	95367	273123	2.17%	0.208%
25	5.586	811	820	829	rVV2	729034	2398301	19.07%	1.827%
26	5.678	830	835	839	rVV	154500	426141	3.39%	0.325%
27	5.727	840	843	864	rVV6	153433	503817	4.01%	0.384%
28	5.934	864	877	891	rVV5	163511	544265	4.33%	0.415%
29	6.080	891	901	912	rVB	94605	245116	1.95%	0.187%
30	6.269	912	932	941	rBV4	166634	754442	6.00%	0.575%
31	6.361	942	947	956	rVV2	226169	687020	5.46%	0.523%
32	6.464	956	964	975	rVB2	160247	454357	3.61%	0.346%
33	6.708	991	1004	1011	rBV3	52484	133017	1.06%	0.101%
34	6.873	1022	1031	1037	rBV2	234836	637504	5.07%	0.486%

LSC Area Percent Report

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Sample : J2132-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Title : SW846 8260

35	6.946	1041	1043	1058	rVB3	143655	299074	2.38%	0.228%
36	7.263	1087	1095	1104	rBV2	115498	260228	2.07%	0.198%
37	7.470	1115	1129	1141	rBV3	667577	1668807	13.27%	1.271%
38	7.738	1167	1173	1181	rVB	90810	176940	1.41%	0.135%
39	7.964	1198	1210	1217	rBV5	113980	392776	3.12%	0.299%
40	8.061	1217	1226	1238	rVB3	79725	230752	1.83%	0.176%
41	8.183	1238	1246	1249	rBV	147836	308624	2.45%	0.235%
42	8.220	1249	1252	1257	rVB	162464	244484	1.94%	0.186%
43	8.519	1295	1301	1305	rBV2	87825	163606	1.30%	0.125%
44	8.586	1305	1312	1320	rVB3	258062	584027	4.64%	0.445%
45	8.677	1320	1327	1329	rBV4	85909	157167	1.25%	0.120%
46	8.726	1329	1335	1341	rVB	480596	802934	6.38%	0.612%
47	8.793	1341	1346	1351	rBV	234481	344392	2.74%	0.262%
48	8.921	1363	1367	1373	rVB	166412	260476	2.07%	0.198%
49	9.171	1401	1408	1413	rVB	73267	125858	1.00%	0.096%
50	9.226	1413	1417	1429	rVB	97915	175287	1.39%	0.134%
51	10.122	1553	1564	1569	rBV	529337	780462	6.21%	0.594%
52	10.262	1582	1587	1593	rBV	10020624	12576595	100.00%	9.579%
53	10.372	1599	1605	1611	rVB	6693917	9044180	71.91%	6.889%
54	10.707	1654	1660	1670	rVB	2367832	3006322	23.90%	2.290%
55	11.024	1706	1712	1719	rVB	644254	827883	6.58%	0.631%
56	11.146	1727	1732	1737	rVB	542708	664926	5.29%	0.506%
57	11.366	1759	1768	1775	rBV	1430206	1734212	13.79%	1.321%
58	11.445	1775	1781	1782	rVV	1697286	2212873	17.60%	1.686%
59	11.463	1782	1784	1788	rVV	1762698	1865281	14.83%	1.421%
60	11.512	1788	1792	1799	rVB	1408161	1721782	13.69%	1.311%
61	11.677	1813	1819	1828	rBV	2720859	3297711	26.22%	2.512%
62	11.817	1837	1842	1847	rVB	7174332	8090077	64.33%	6.162%
63	11.951	1862	1864	1871	rVB	113191	142920	1.14%	0.109%
64	12.030	1873	1877	1881	rBV	101171	128933	1.03%	0.098%
65	12.085	1881	1886	1892	rVB	832703	1084767	8.63%	0.826%
66	12.152	1892	1897	1902	rBV	1759882	2091295	16.63%	1.593%
67	12.311	1916	1923	1930	rBV	8715258	11162859	88.76%	8.503%
68	12.378	1930	1934	1944	rVB4	860734	1269972	10.10%	0.967%
69	12.475	1944	1950	1954	rBV2	267043	381580	3.03%	0.291%
70	12.524	1954	1958	1963	rVB	392841	473450	3.76%	0.361%
71	12.597	1964	1970	1972	rBV	768563	1043223	8.29%	0.795%

LSC Area Percent Report

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Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Title : SW846 8260

72	12.616	1972	1973	1978	rVB	419100	378323	3.01%	0.288%
73	12.677	1978	1983	1986	rBV	1716923	2007405	15.96%	1.529%
74	12.707	1986	1988	1993	rVV	594399	693203	5.51%	0.528%
75	12.762	1993	1997	2005	rVB	2196173	2958710	23.53%	2.254%
76	12.908	2016	2021	2026	rVB3	232088	327026	2.60%	0.249%
77	12.987	2026	2034	2037	rBV	909864	1155028	9.18%	0.880%
78	13.024	2037	2040	2046	rVB	773047	926307	7.37%	0.706%
79	13.091	2046	2051	2056	rBV3	168321	363654	2.89%	0.277%
80	13.231	2070	2074	2079	rVB	1644289	1998719	15.89%	1.522%
81	13.317	2084	2088	2090	rBV2	151022	199169	1.58%	0.152%
82	13.359	2090	2095	2100	rVV2	3565139	4608311	36.64%	3.510%
83	13.408	2100	2103	2106	rVV4	268218	355401	2.83%	0.271%
84	13.487	2112	2116	2122	rVB	518622	672786	5.35%	0.512%
85	13.567	2124	2129	2132	rBV2	196564	277894	2.21%	0.212%
86	13.609	2132	2136	2139	rVV	518535	701929	5.58%	0.535%
87	13.652	2139	2143	2147	rVV2	430892	779178	6.20%	0.593%
88	13.731	2148	2156	2161	rVB2	525009	1118198	8.89%	0.852%
89	13.841	2167	2174	2184	rBV	4040469	5090329	40.47%	3.877%
90	13.932	2184	2189	2194	rVB	314297	448081	3.56%	0.341%
91	13.993	2194	2199	2203	rBV2	188152	265813	2.11%	0.202%
92	14.036	2203	2206	2210	rVB2	170320	194720	1.55%	0.148%
93	14.140	2218	2223	2228	rBV	324207	394163	3.13%	0.300%
94	14.280	2242	2246	2252	rBV	239118	378215	3.01%	0.288%
95	14.438	2268	2272	2277	rVB	185110	222177	1.77%	0.169%
96	14.664	2304	2309	2312	rBV	150756	206139	1.64%	0.157%
97	14.707	2312	2316	2326	rVB	433496	611737	4.86%	0.466%
98	14.847	2334	2339	2348	rVB2	550282	765590	6.09%	0.583%

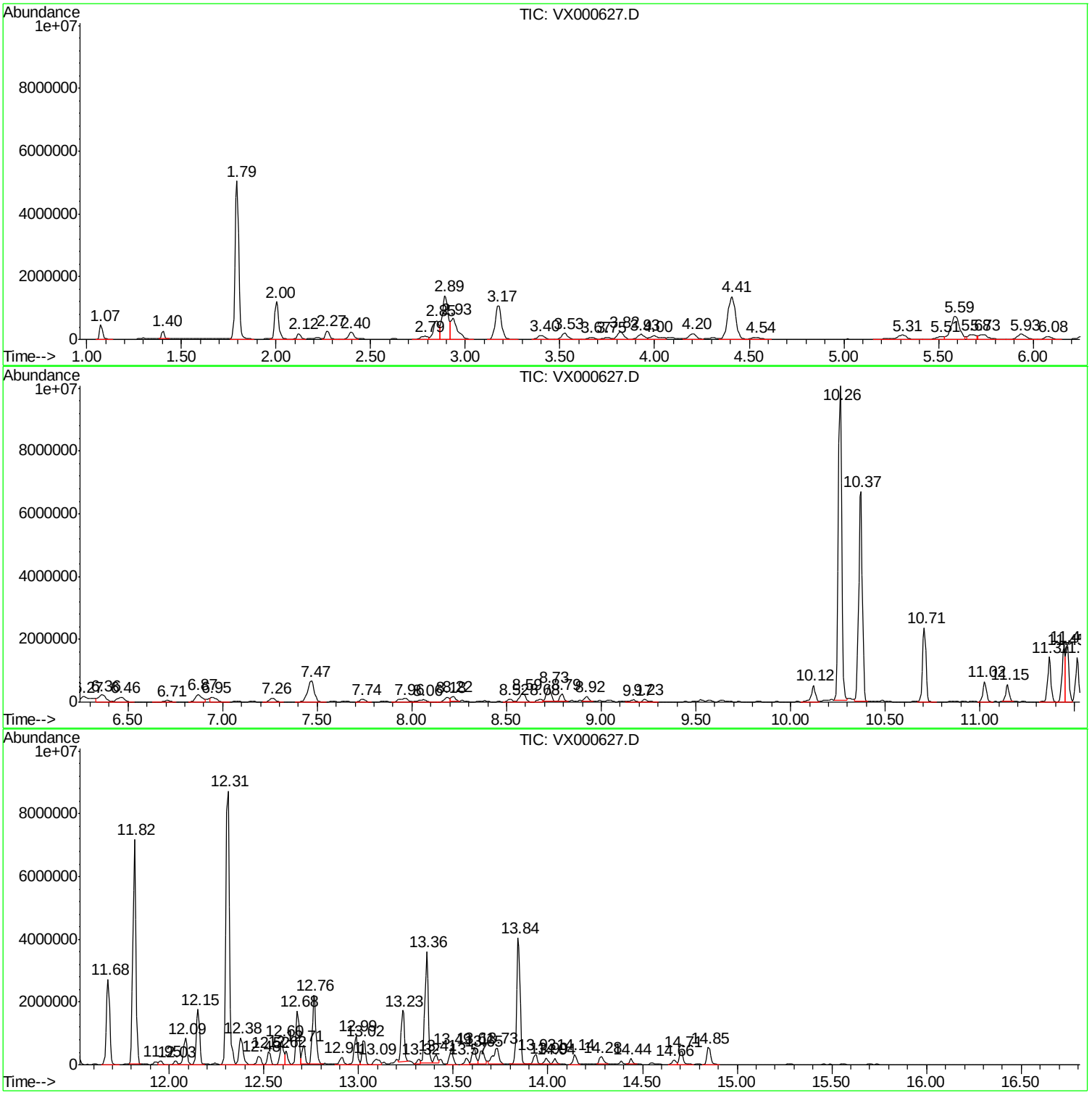
Sum of corrected areas: 131287718

LSC Report - Integrated Chromatogram

Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

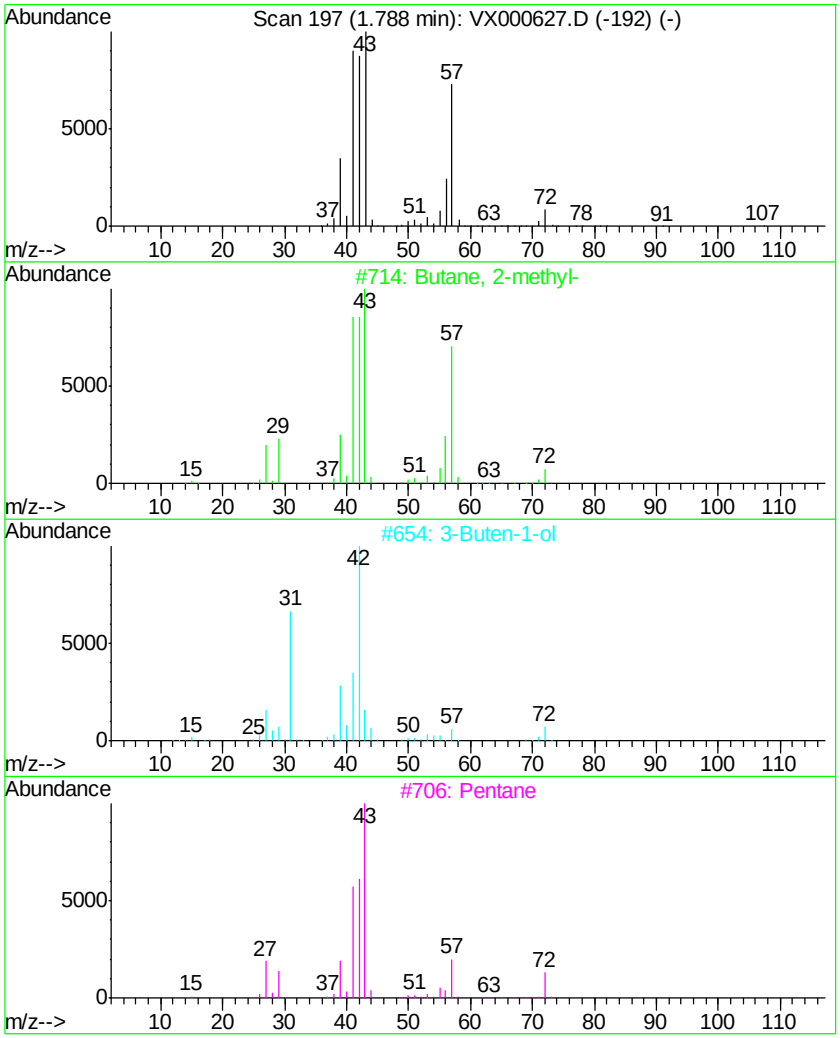
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

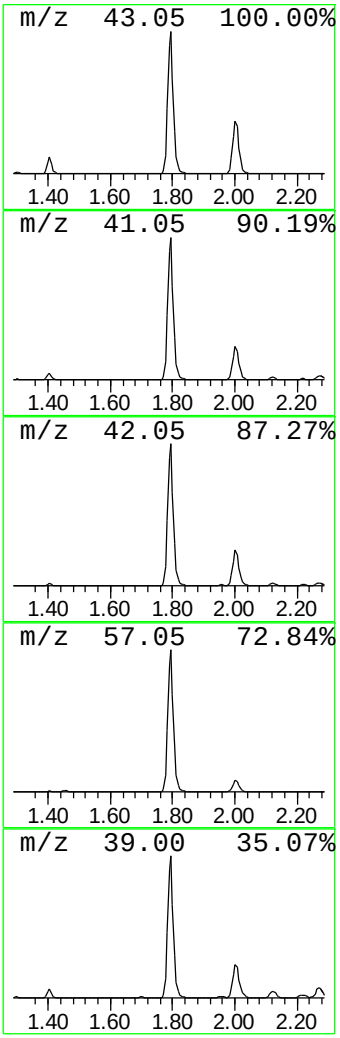
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.79	795.72 ug/l	6781780	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2	3-Buten-1-ol	72	C4H8O	000627-27-0	43
3	Pentane	72	C5H12	000109-66-0	28
4	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	27
5	Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	10





Library Search Compound Report

Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

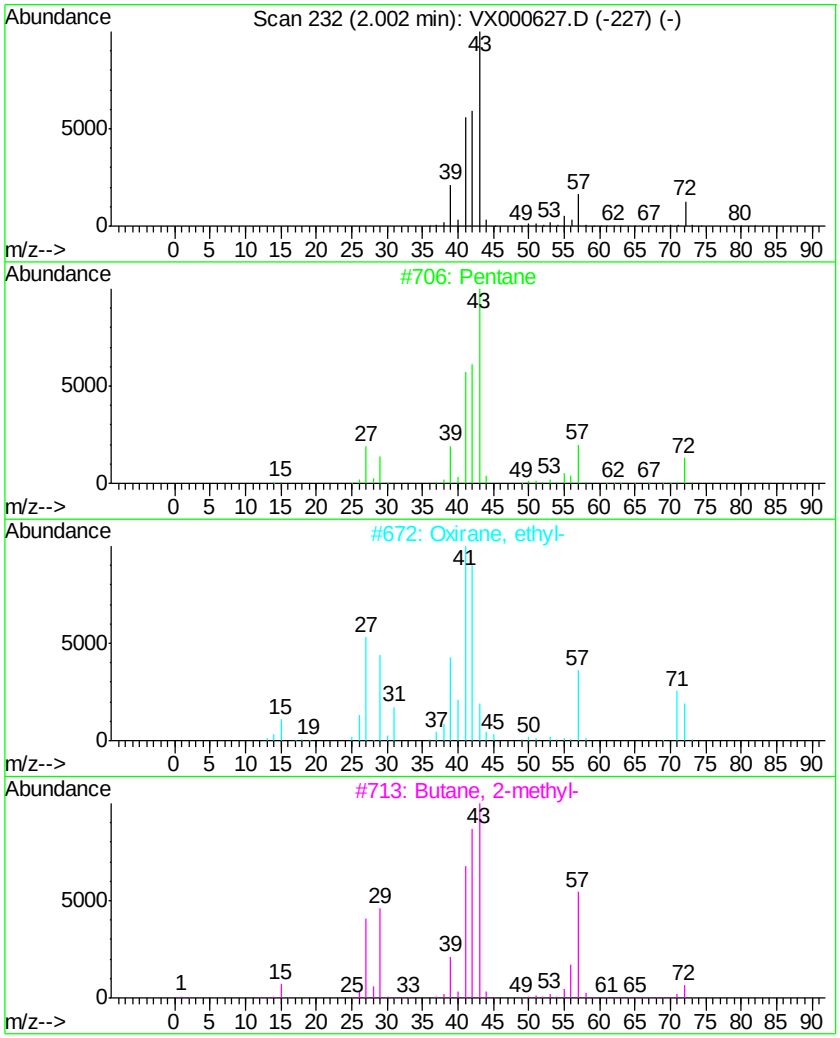
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

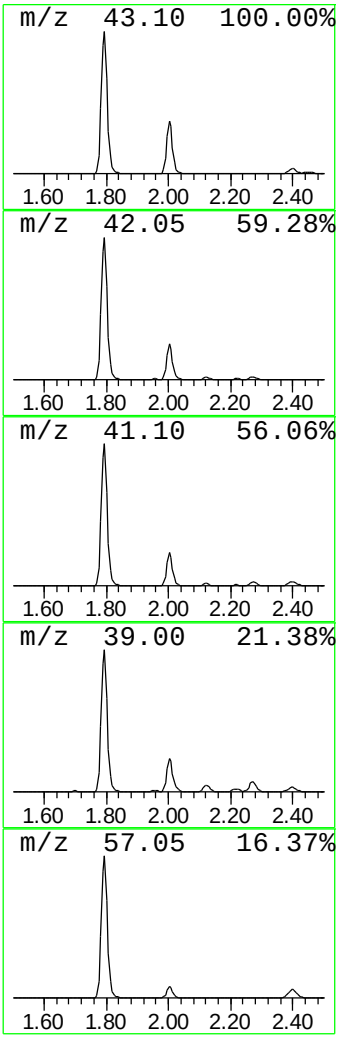
****

Peak Number 2 Pentane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	204.42 ug/l	1742260	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Oxirane, ethyl-	72	C4H8O	000106-88-7	53
3	Butane, 2-methyl-	72	C5H12	000078-78-4	39
4	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	38
5	Butanal, 2,2-dimethyl-	100	C6H12O	002094-75-9	9





Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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

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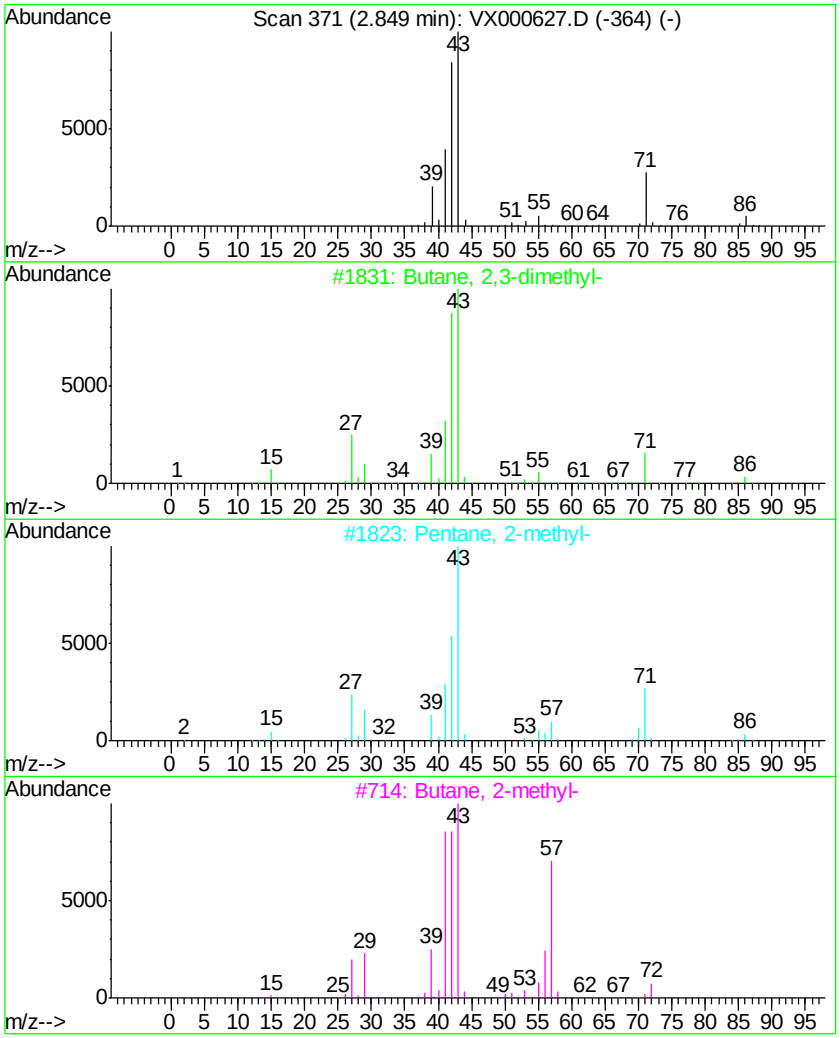
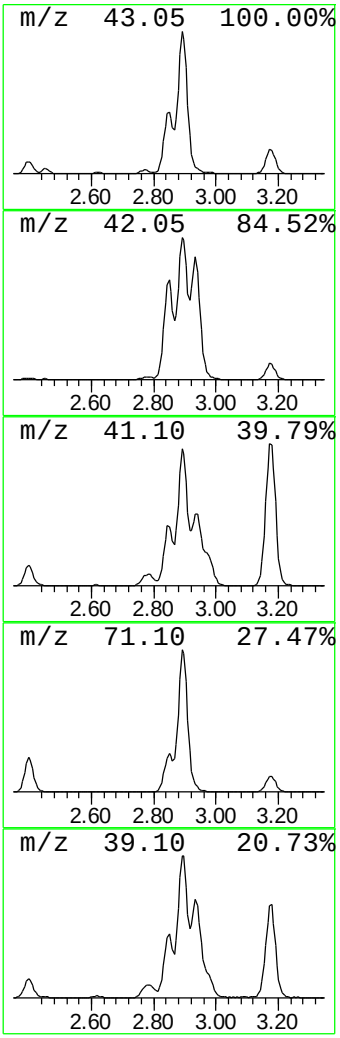
Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.85	155.34 ug/l	1323960	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	53
3	Butane, 2-methyl-	72	C5H12	000078-78-4	36
4	Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	12
5	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9





Library Search Compound Report

Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

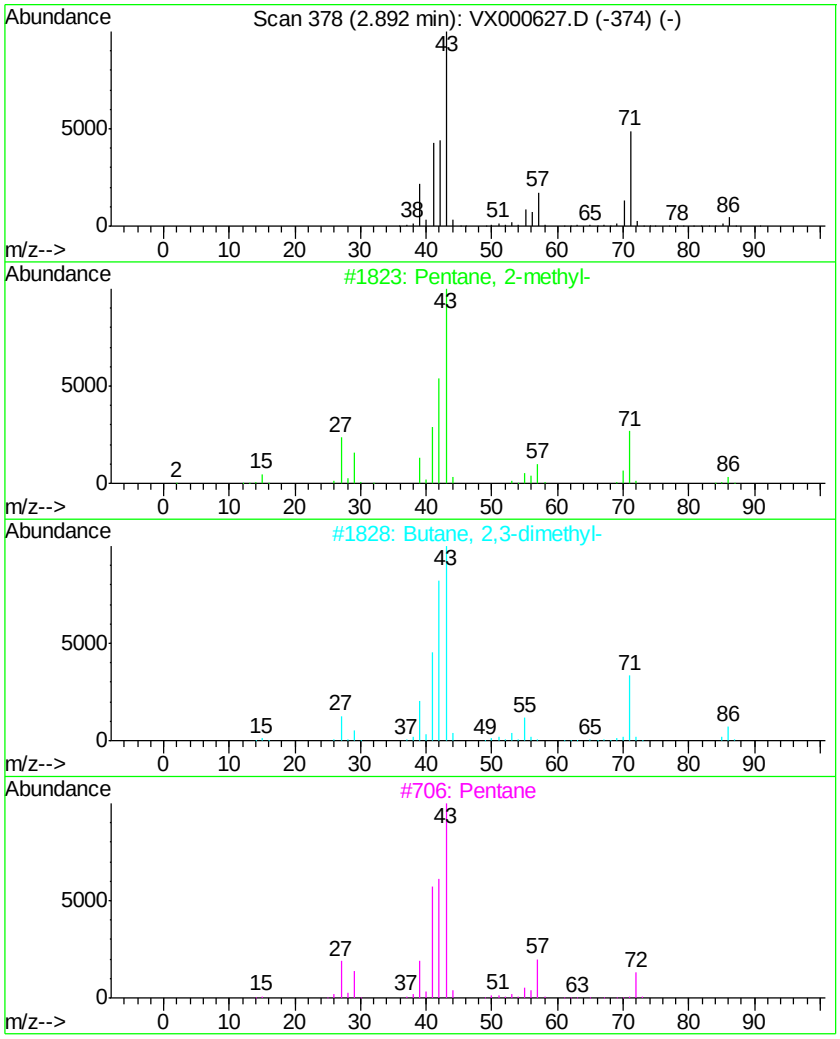
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

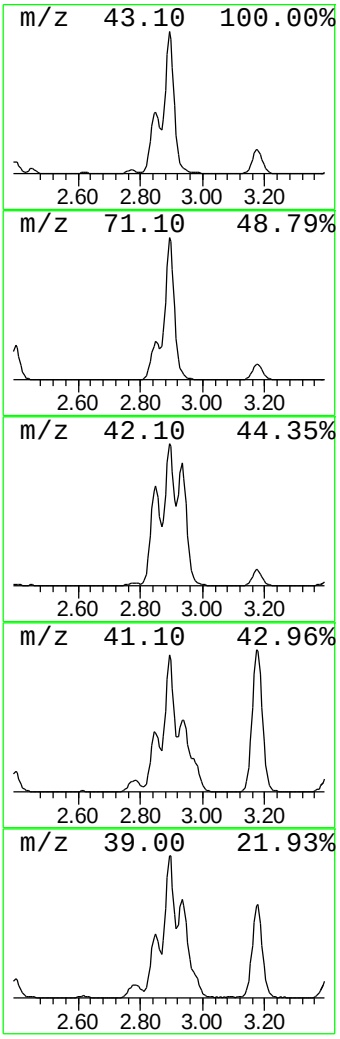
****

Peak Number 4 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	353.05 ug/l	3009000	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	50
3	Pentane	72	C5H12	000109-66-0	46
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38





Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

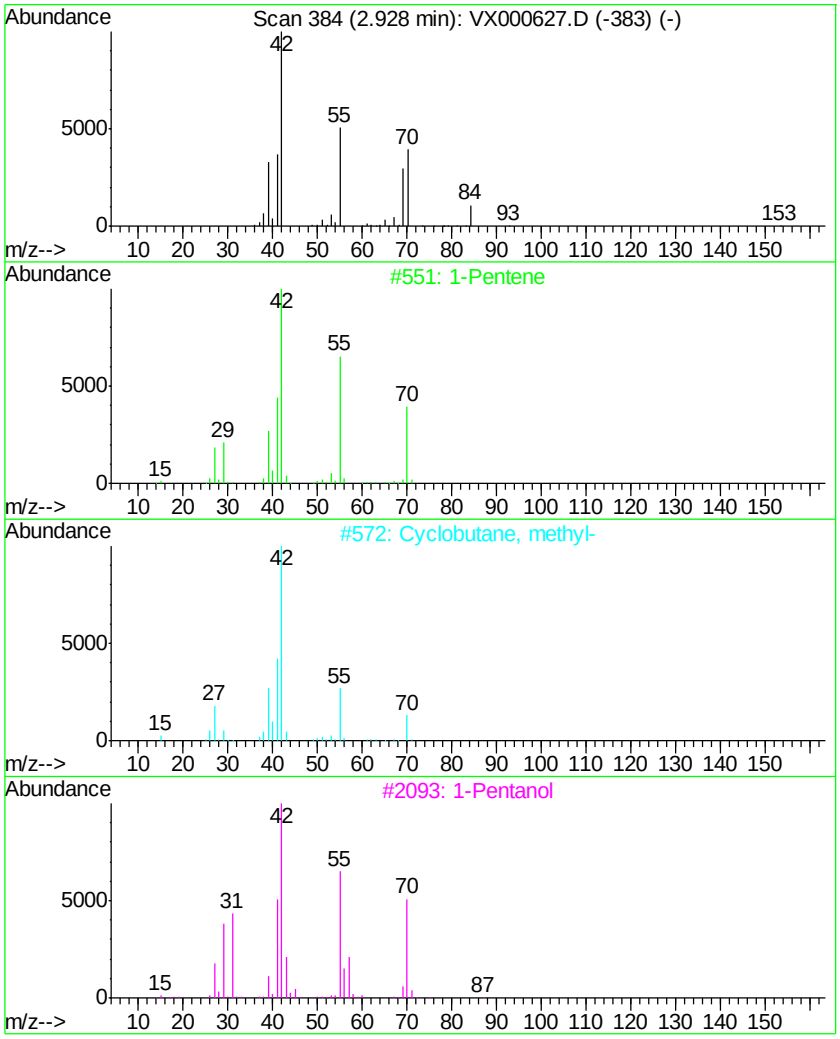
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

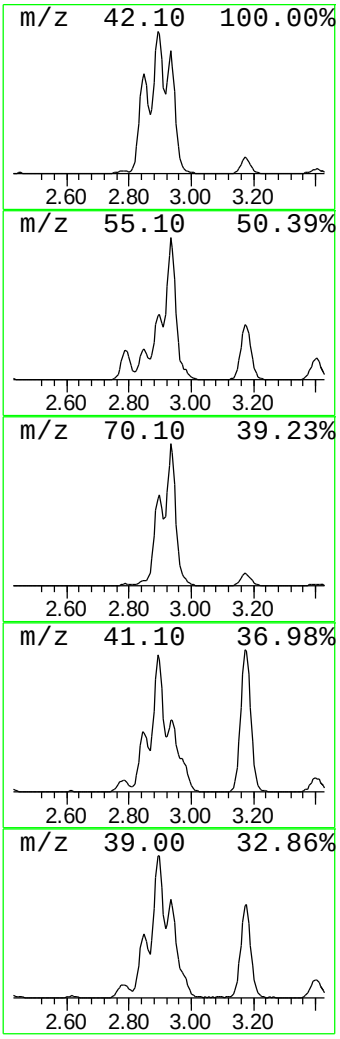
****

Peak Number 5 1-Pentene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	174.19 ug/l	1484600	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene	70	C5H10	000109-67-1	72
2	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
3	1-Pentanol	88	C5H12O	000071-41-0	50
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	42
5	3-Hexene, (Z)-	84	C6H12	007642-09-3	22





Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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

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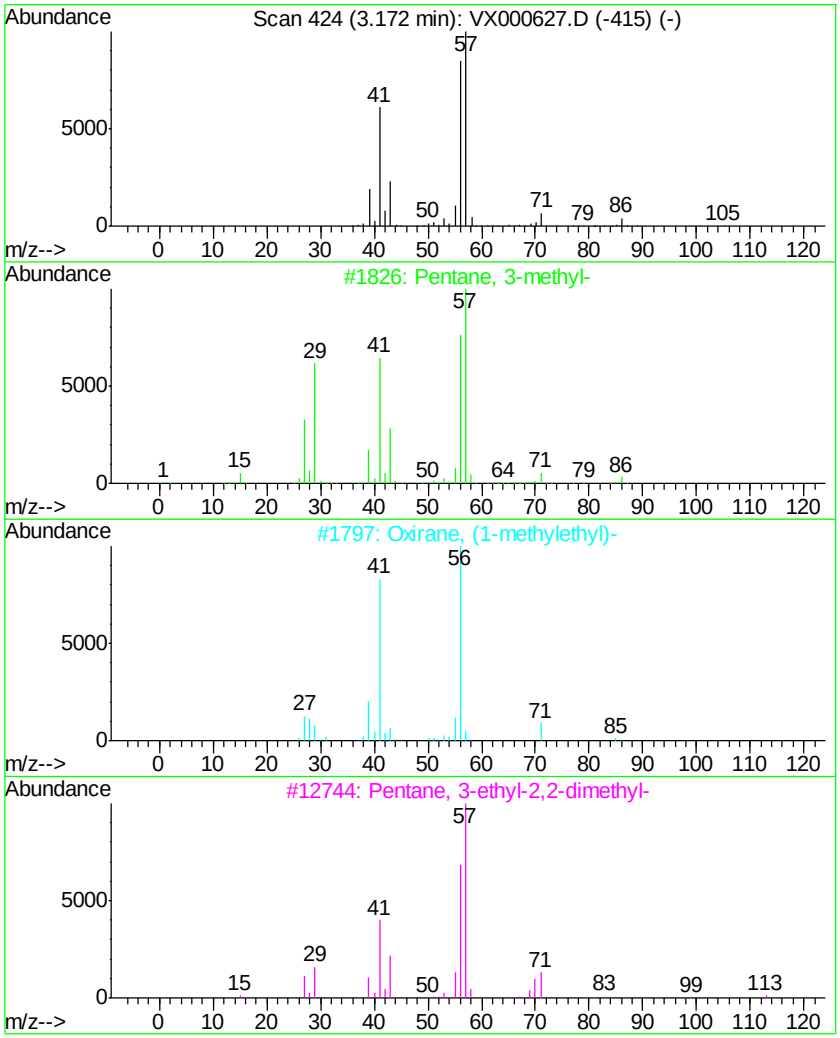
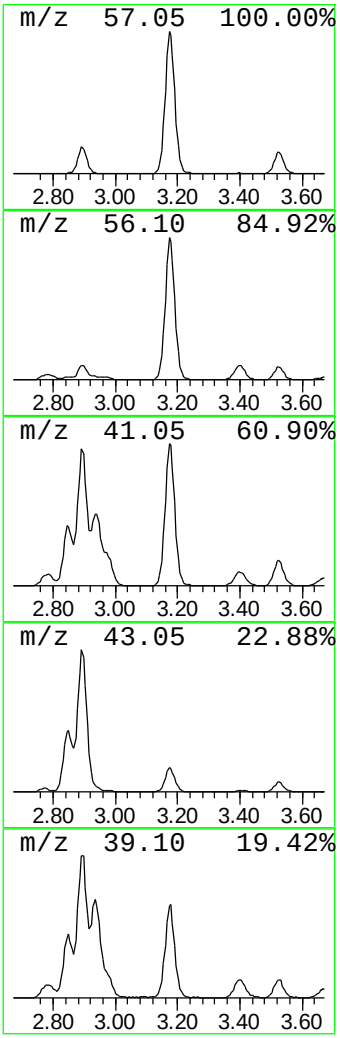
Peak Number 6 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	283.75 ug/l	2418330	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
4	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40
5	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40





Library Search Compound Report

Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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

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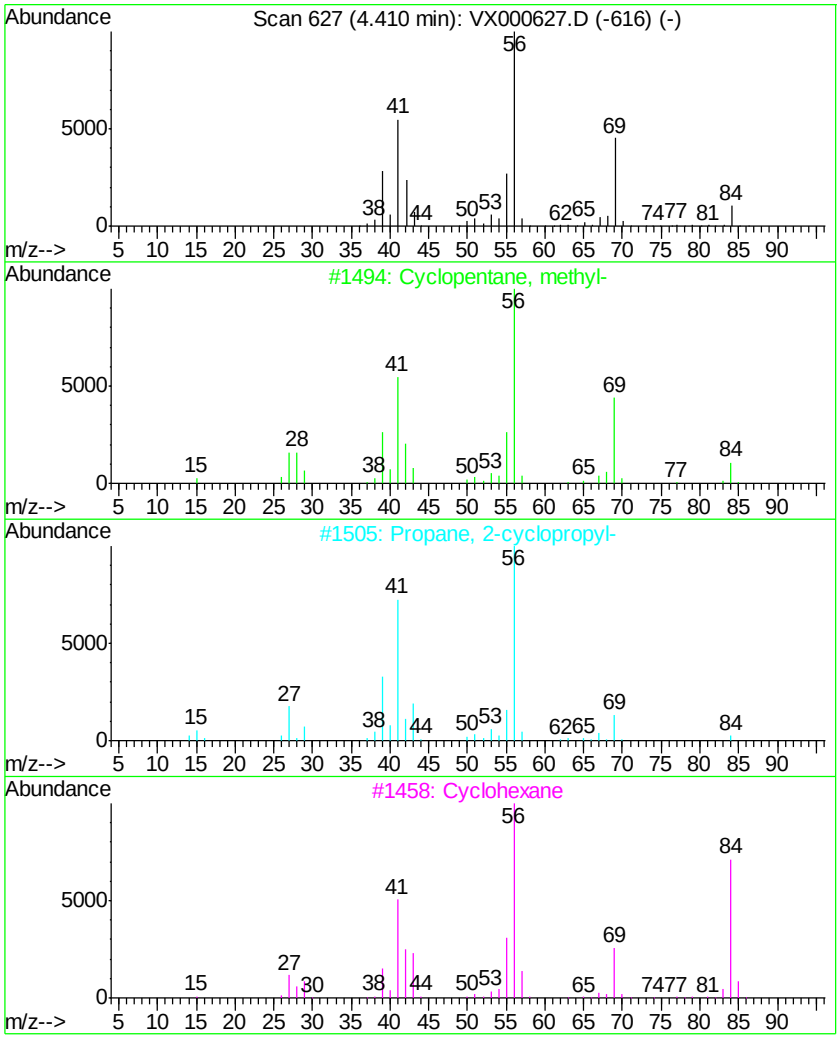
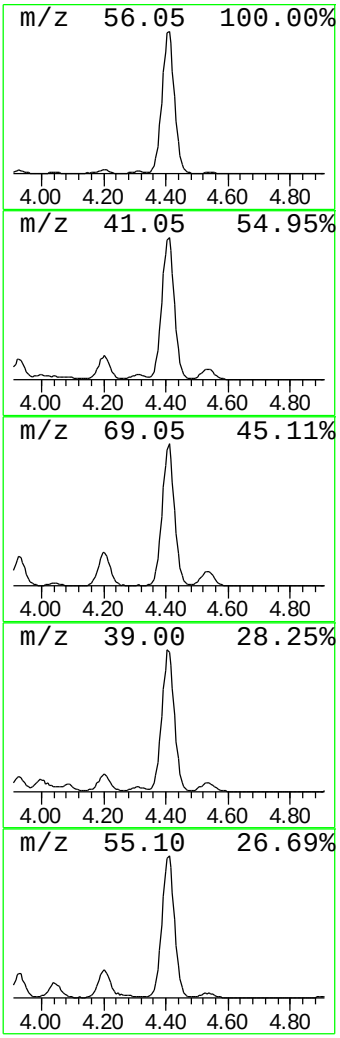
Peak Number 7 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	451.88 ug/l	3851310	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	80
3	Cyclohexane	84	C6H12	000110-82-7	78
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Cyclobutane	56	C4H8	000287-23-0	53





Library Search Compound Report

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Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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

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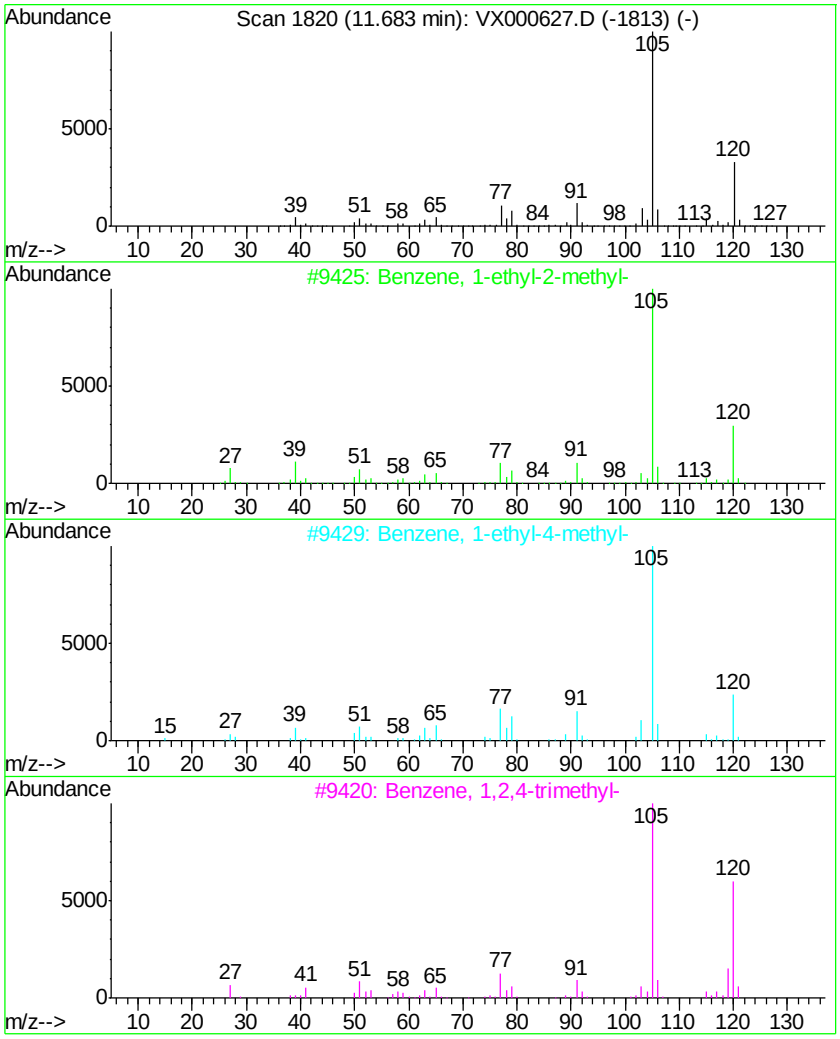
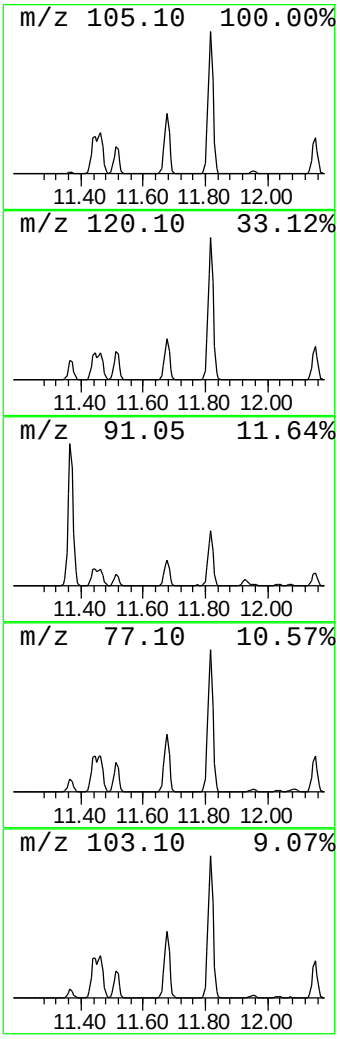
Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.68	152.00 ug/l	3297710	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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

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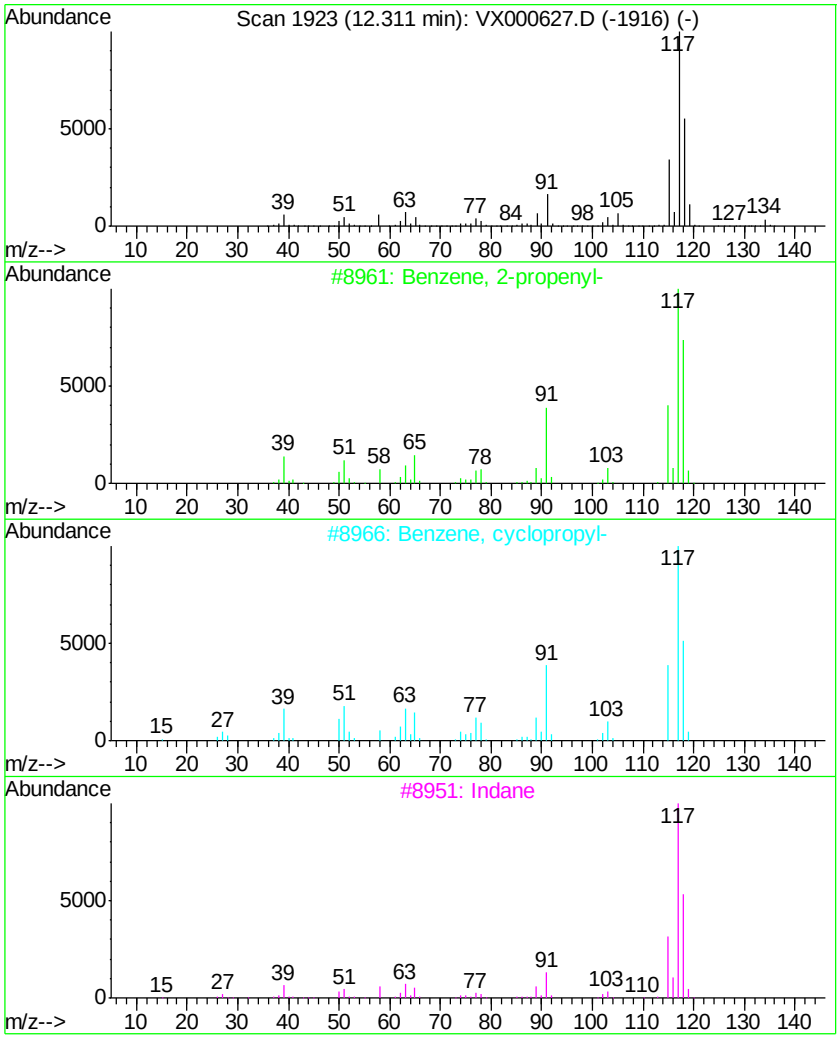
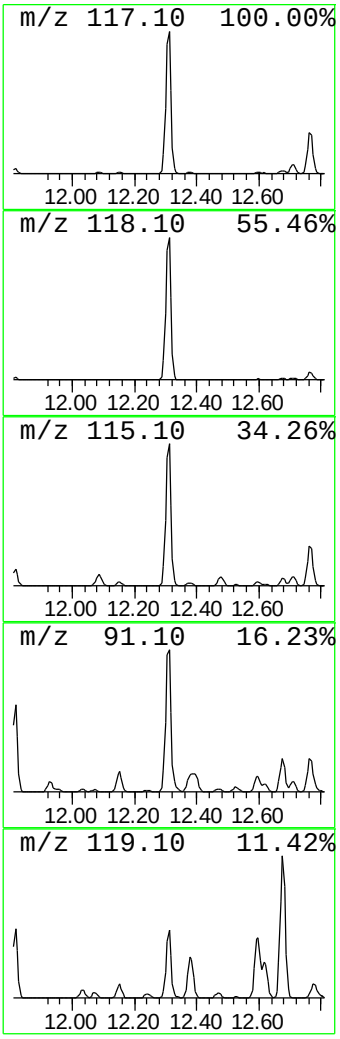
Peak Number 9 Benzene, 2-propenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.31	514.53 ug/l	11162900	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	87
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
3	Indane	118	C9H10	000496-11-7	81
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64





Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
Quant Title : SW846 8260

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

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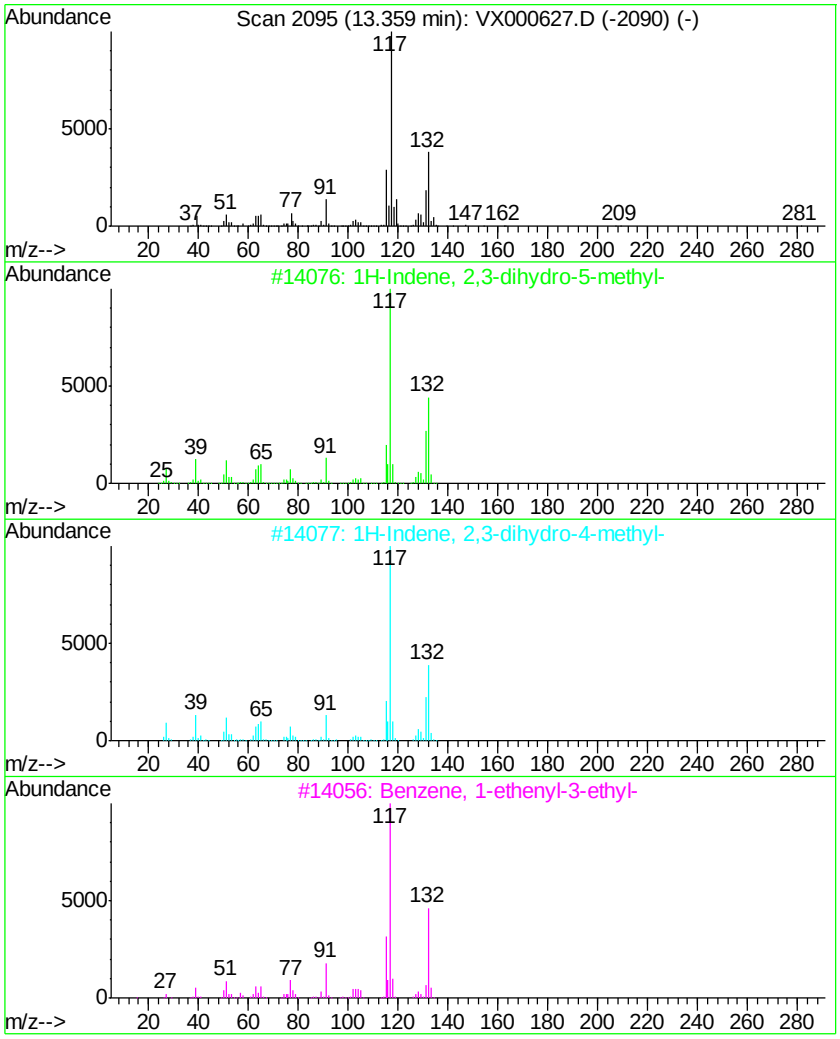
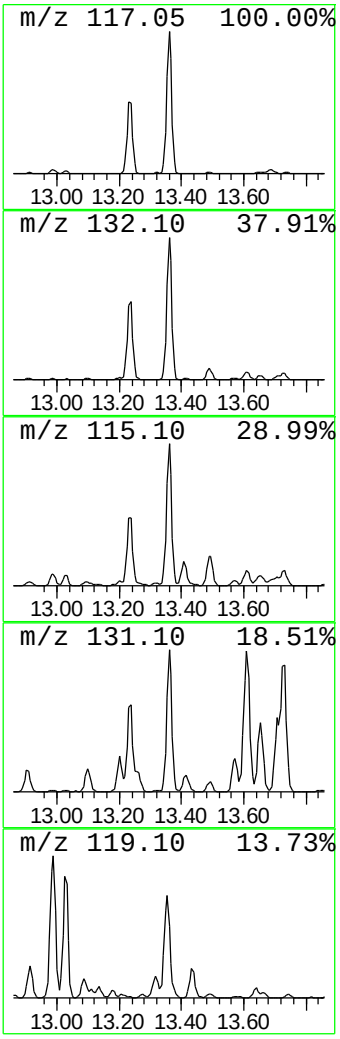
Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	212.41 ug/l	4608310	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	80
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	80
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	78





Tentatively Identified Compound (LSC) summary

Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX033118\
Data File : VX000627.D
Acq On : 01 Apr 2018 06:36
Operator : JC/MD
Sample : J2132-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.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-methyl-	1.79	795.7	ug/l	6781780	1	5.67	426141	50.0
Pentane	2.00	204.4	ug/l	1742260	1	5.67	426141	50.0
Butane, 2,3-dimet...	2.85	155.3	ug/l	1323960	1	5.67	426141	50.0
Pentane, 2-methyl-	2.89	353.1	ug/l	3009000	1	5.67	426141	50.0
1-Pentene	2.93	174.2	ug/l	1484600	1	5.67	426141	50.0
Pentane, 3-methyl-	3.17	283.8	ug/l	2418330	1	5.67	426141	50.0
Cyclopentane, met...	4.41	451.9	ug/l	3851310	1	5.67	426141	50.0
Benzene, 1-ethyl-...	11.68	152.0	ug/l	3297710	4	12.09	1084770	50.0
Benzene, 2-propenyl-	12.31	514.5	ug/l	11162900	4	12.09	1084770	50.0
1H-Indene, 2,3-di...	13.36	212.4	ug/l	4608310	4	12.09	1084770	50.0