LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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	1.794	99	105	116	rVB	369200	529429	3.18%	0.321%
2	2.001	133	139	151	rVB	657890	918198	5.51%	0.557%
3	2.397	195	204	220	rVB	233338	480151	2.88%	0.291%
4	2.848	270	278	280	rBV	239563	471908	2.83%	0.286%
5	2.891	280	285	305	rVB	867715	2249413	13.50%	1.365%
6	3.178	321	332	347	rVB	751959	1749909	10.50%	1.062%
7	3.525	379	389	404	rBV	643882	1471968	8.83%	0.893%
8	4.403	523	533	546	rVB	761460	2243543	13.46%	1.361%
9	5.500	700	713	718	rBV	145947	424626	2.55%	0.258%
10	5.586	718	727	735	rVV4	516755	1919208	11.52%	1.164%
11	5.659	736	739	746	rVV	258536	661367	3.97%	0.401%
12	5.726	746	750	769	rVB	113861	337410	2.02%	0.205%
13	5.927	771	783	796	rBV2	286723	874617	5.25%	0.531%
14	6.061	796	805	818	rVV	144247	377426	2.27%	0.229%
15	6.262	818	838	848	rVV3	82255	328140	1.97%	0.199%
16	6.360	848	854	861	rVV2	74957	202343	1.21%	0.123%
17	6.458	861	870	881	rVB2	131873	370940	2.23%	0.225%
18	6.708	898	911	920	rBV	234809	557844	3.35%	0.338%
19	6.860	927	936	944	rBV	360159	860398	5.16%	0.522%
20	7.463	1021	1035	1047	rBV	1203942	2717966	16.31%	1.649%
21	7.732	1071	1079	1090	rVB	154998	308461	1.85%	0.187%
22	8.500	1200	1205	1210	rBV	107908	184800	1.11%	0.112%
23	8.664	1226	1232	1235	rBV	269273	466220	2.80%	0.283%
24	8.713	1235	1240	1251	rVV	837338	1408383	8.45%	0.855%
25	8.835	1254	1260	1265	rVV	119894	209949	1.26%	0.127%
26	8.969	1278	1282	1288	rVV	140808	223620	1.34%	0.136%
27	9.036	1288	1293	1300	rVB	228757	378515	2.27%	0.230%
28	9.164	1305	1314	1320	rBV	123796	202332	1.21%	0.123%
29	9.573	1374	1381	1385	rBV3	119432	204804	1.23%	0.124%
30	9.622	1385	1389	1394	rVB	327962	445366	2.67%	0.270%
31	9.677	1394	1398	1402	rBV	146102	206365	1.24%	0.125%
32	10.109	1465	1469	1475	rVB	784573	1053403	6.32%	0.639%
33	10.183	1475	1481	1486	rVV	117006	179703	1.08%	0.109%
34	10.250	1486	1492	1498	rVV	7924657	10474417	62.86%	6.355%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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	10.359	1501	1510	1516	rVV	10016720	13590627	81.56%	8.246%
36	10.701	1560	1566	1576	rVB	9050059	11717712	70.32%	7.110%
37	10.835	1579	1588	1593	rVB2	117784	184029	1.10%	0.112%
38	11.018	1612	1618	1623	rBV	1424103	1782162	10.70%	1.081%
39	11.134	1631	1637	1642	rBV	757165	962700	5.78%	0.584%
40	11.359	1668	1674	1681	rBV	1859976	2277617	13.67%	1.382%
41	11.432	1681	1686	1694	rVV	8508567	16065520	96.41%	9.748%
42	11.505	1694	1698	1707	rVB	4847144	5846408	35.09%	3.547%
43	11.670	1719	1725	1733	rBV	7620282	9993114	59.97%	6.063%
44	11.810	1742	1748	1753	rVB	13127664	16663399	100.00%	10.110%
45	11.944	1767	1770	1775	rVB	331211	400496	2.40%	0.243%
46	12.024	1775	1783	1786	rBV	819645	993825	5.96%	0.603%
47	12.072	1786	1791	1797	rVB2	1097826	1704170	10.23%	1.034%
48	12.139	1797	1802	1808	rBV	7659497	9729691	58.39%	5.903%
49	12.231	1812	1817	1822	rVB	191705	233802	1.40%	0.142%
50	12.298	1822	1828	1831	rBV	2527688	3560894	21.37%	2.161%
51	12.322	1831	1832	1835	rVV	1117388	738194	4.43%	0.448%
52	12.371	1835	1840	1849	rVB2	1824133	2606733	15.64%	1.582%
53	12.456	1849	1854	1859	rBV	560672	757228	4.54%	0.459%
54	12.517	1859	1864	1869	rVB	1432515	1718114	10.31%	1.042%
55	12.585	1870	1875	1877	rBV	1270011	1684549	10.11%	1.022%
56	12.609	1877	1879	1883	rVB	1516993	1514494	9.09%	0.919%
57	12.664	1883	1888	1892	rBV	1505198	1860877	11.17%	1.129%
58	12.700	1892	1894	1898	rVB	357828	365750	2.19%	0.222%
59	12.755	1898	1903	1911	rBV3	1063011	1912475	11.48%	1.160%
60	12.847	1915	1918	1922	rBV2	144460	168562	1.01%	0.102%
61	12.902	1922	1927	1932	rVB2	1105348	1354083	8.13%	0.822%
62	12.975	1932	1939	1942	rBV	846049	1099994	6.60%	0.667%
63	13.017	1942	1946	1951	rVB	1651931	1972705	11.84%	1.197%
64	13.078	1951	1956	1961	rBV4	204233	435055	2.61%	0.264%
65	13.194	1971	1975	1977	rBV2	266101	344884	2.07%	0.209%
66	13.219	1977	1979	1983	rVB	380113	439870	2.64%	0.267%
67	13.261	1983	1986	1990	rVB3	162984	207486	1.25%	0.126%
68	13.304	1990	1993	1995	rBV	275872	365496	2.19%	0.222%
69	13.340	1995	1999	2005	rVB2	2713261	3733243	22.40%	2.265%
70	13.475	2017	2021	2029	rVB	1582958	2082400	12.50%	1.263%
71	13.560	2030	2035	2038	rBV3	158940	242548	1.46%	0.147%

LSC Area Percent Report

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Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

72	13.597	2038	2041	2044	rBV	156124	193840	1.16%	0.118%
73	13.639	2044	2048	2053	rVV3	332095	614300	3.69%	0.373%
74	13.718	2058	2061	2067	rVB2	354243	557358	3.34%	0.338%
75	13.834	2072	2080	2086	rVB	1611766	2317498	13.91%	1.406%
76	13.907	2087	2092	2097	rVB	268910	386524	2.32%	0.235%
77	13.981	2097	2104	2108	rBV2	391398	585498	3.51%	0.355%
78	14.285	2146	2154	2159	rVB	548143	899776	5.40%	0.546%
79	14.535	2190	2195	2205	rBV2	523960	696154	4.18%	0.422%
80	14.694	2214	2221	2225	rBV	626328	876593	5.26%	0.532%
81	14.834	2238	2244	2249	rBV	515255	684714	4.11%	0.415%

Sum of corrected areas: 164816303

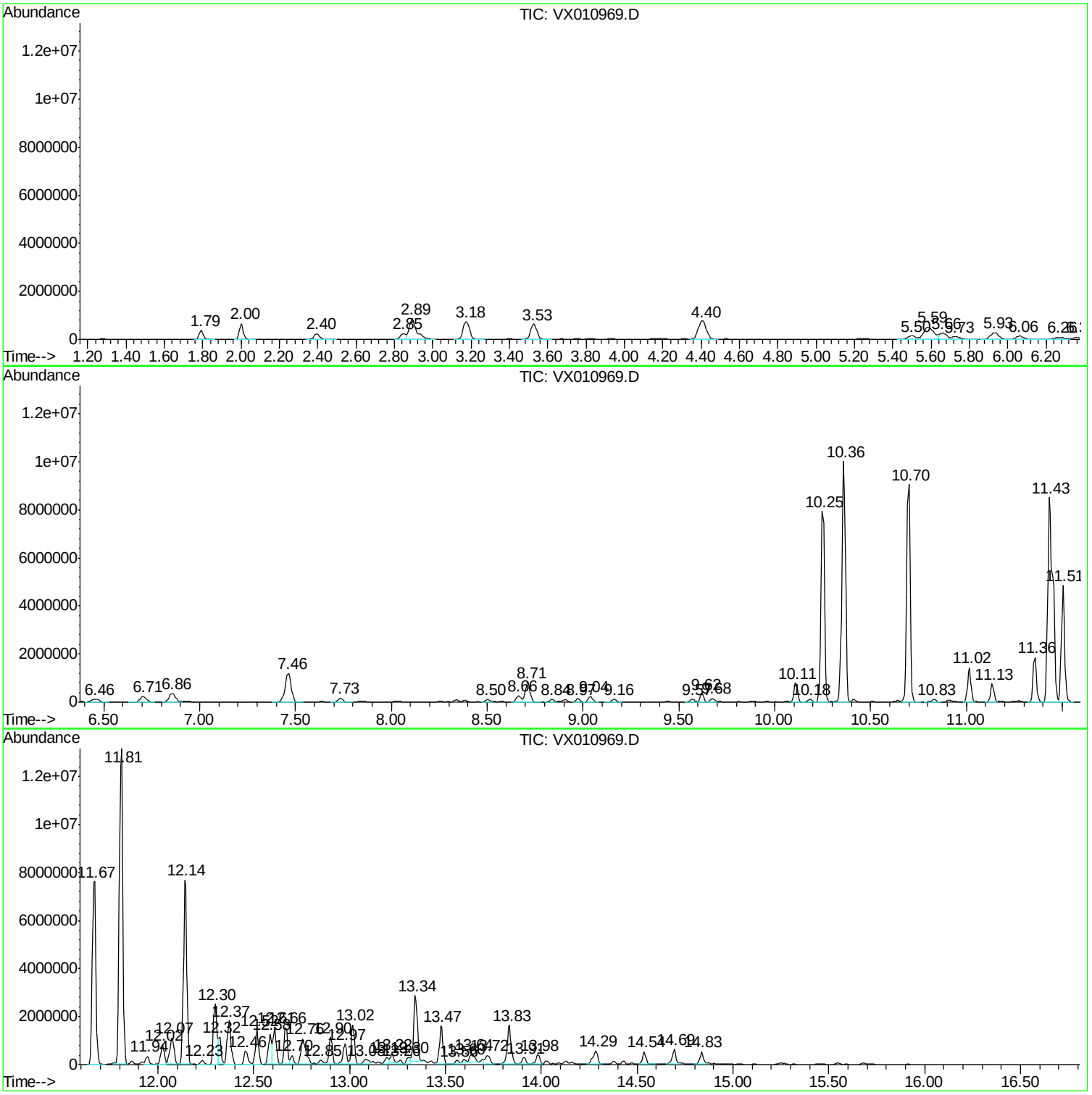
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

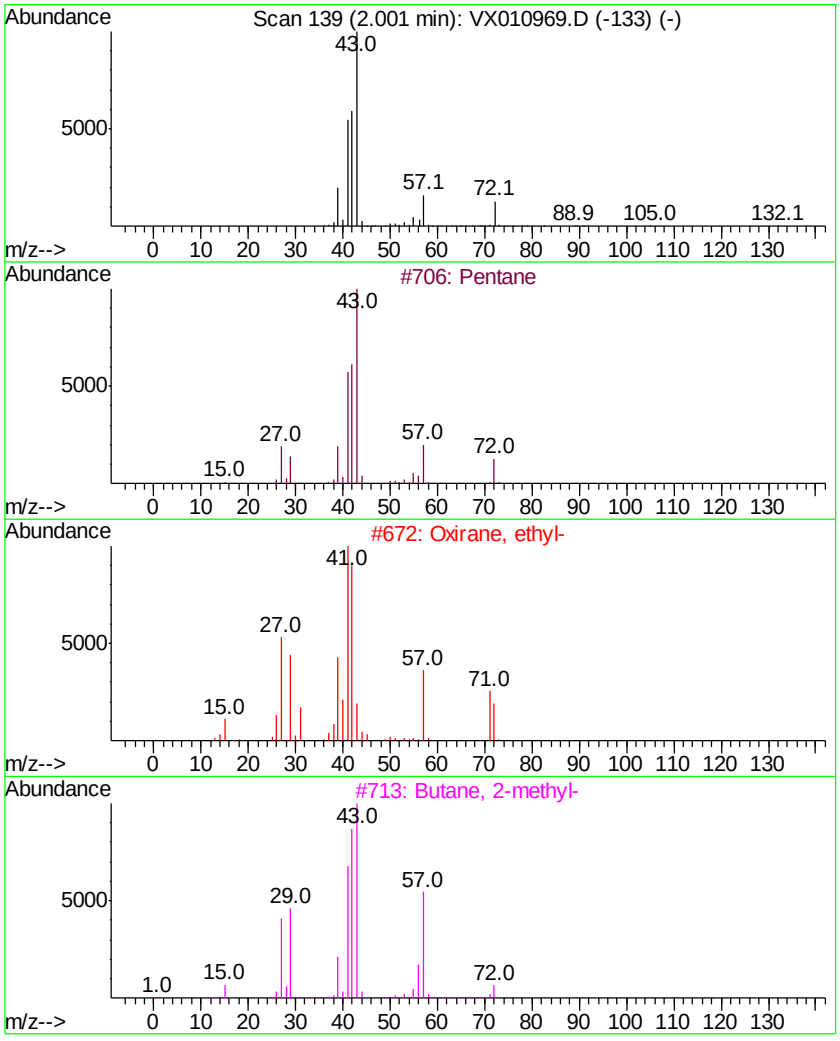
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

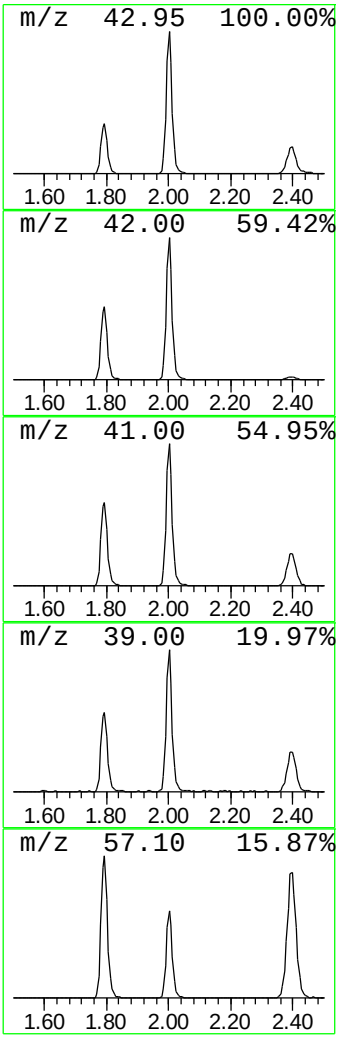
****

Peak Number 1 Pentane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	69.42 ug/l	918198	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	50
3	Butane, 2-methyl-	72	C5H12	000078-78-4	45
4	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	35
5	Butanenitrile, 3-methyl-	83	C5H9N	000625-28-5	9





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

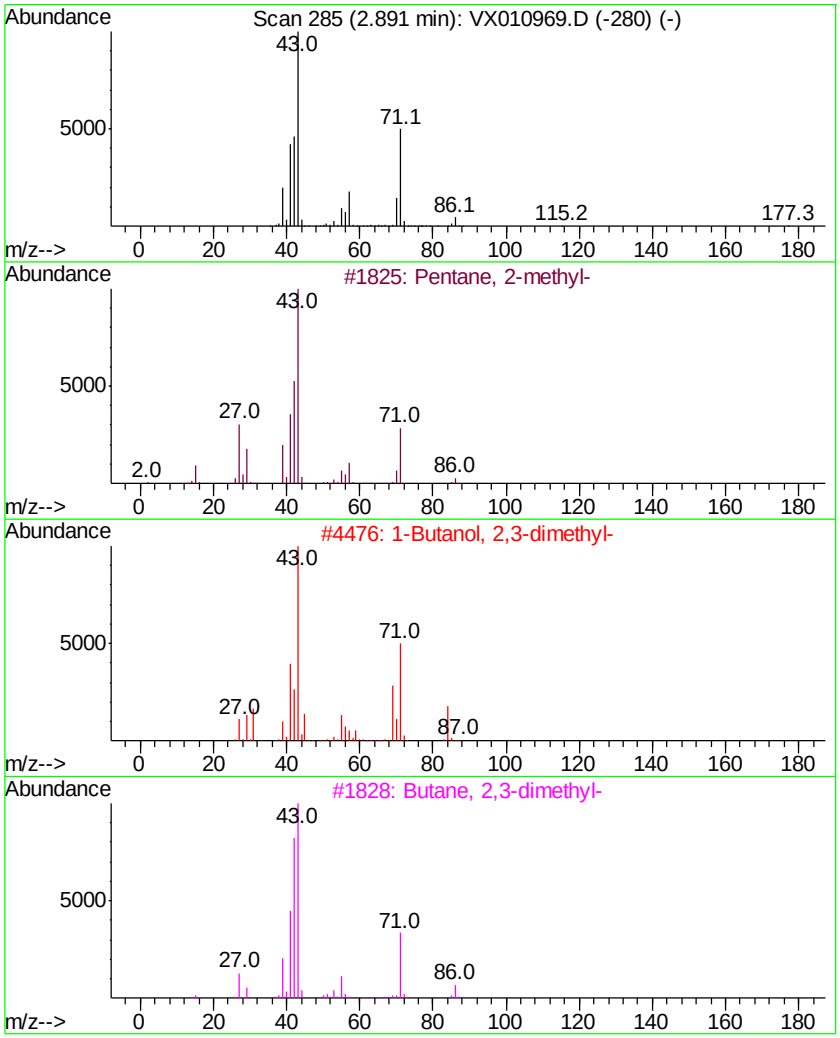
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

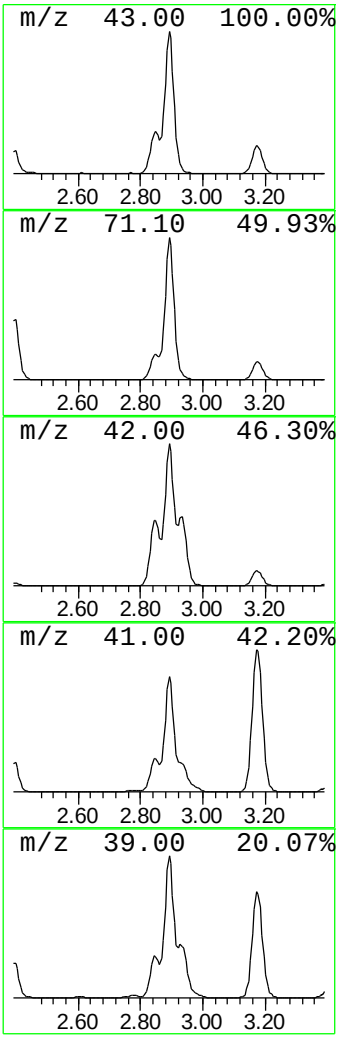
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	170.06 ug/l	2249410	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
4	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
5	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

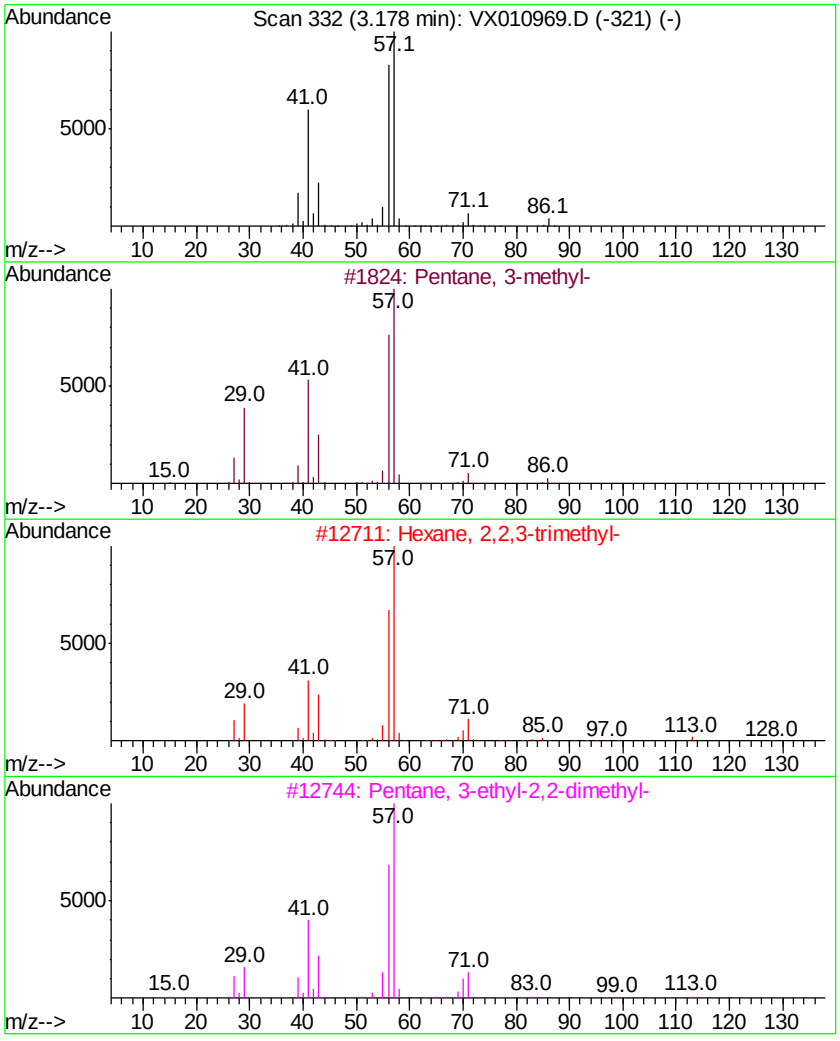
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

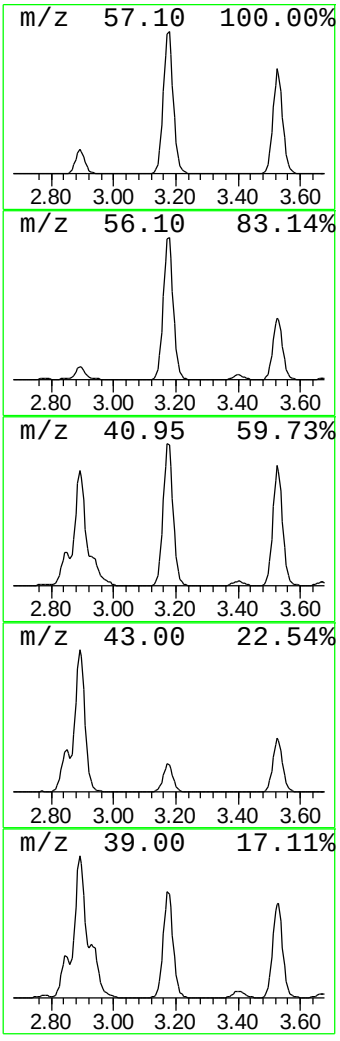
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	132.30 ug/l	1749910	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
4	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40
5	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
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Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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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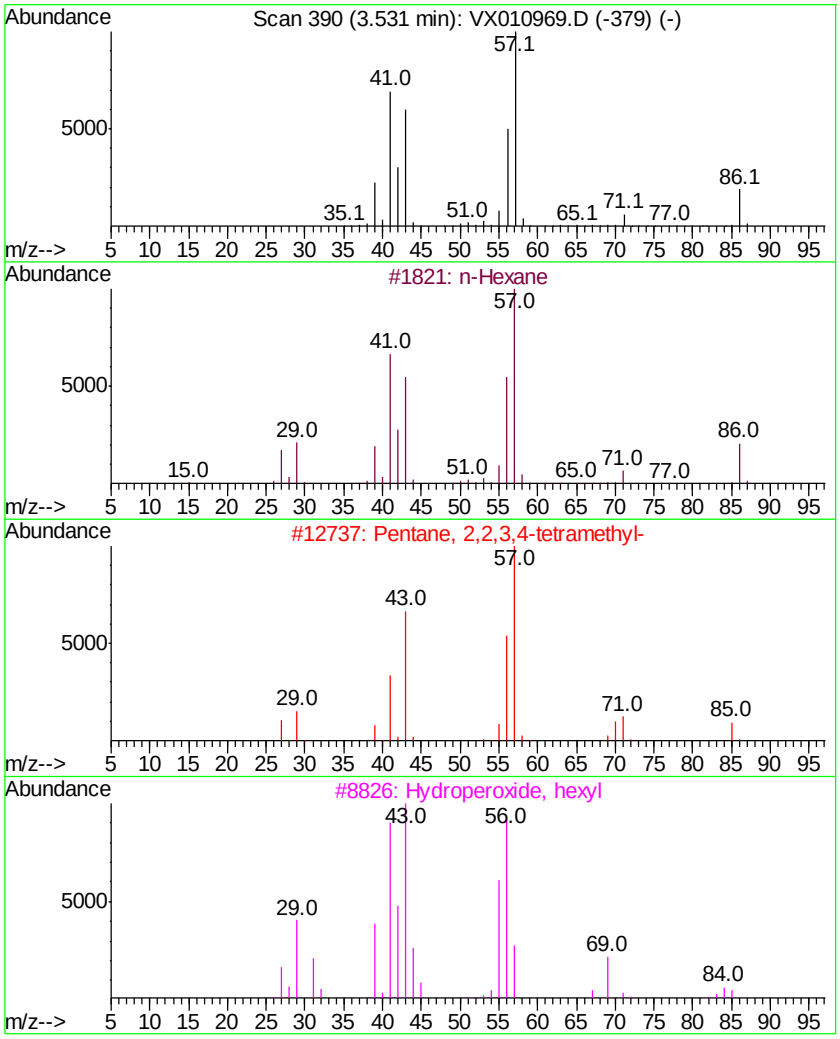
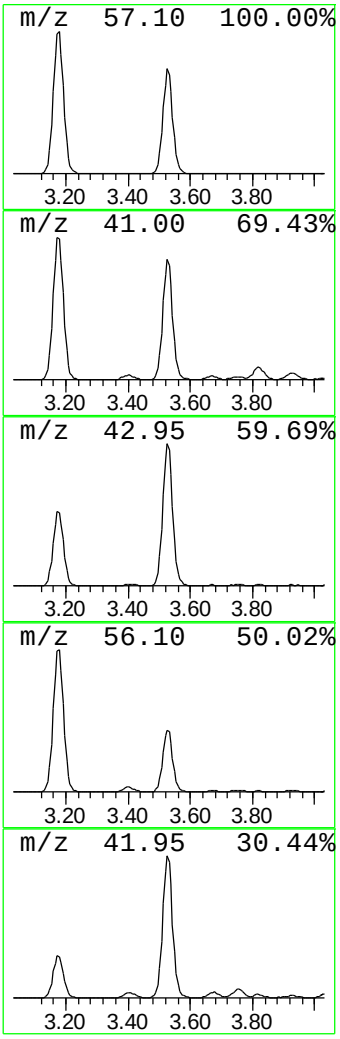
Peak Number 4 n-Hexane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.53	111.28 ug/l	1471970	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	94
2	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	42
3	Hydroperoxide, hexyl	118	C6H14O2	004312-76-9	40
4	Methane, isocyanato-	57	C2H3NO	000624-83-9	37
5	Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	35





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Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

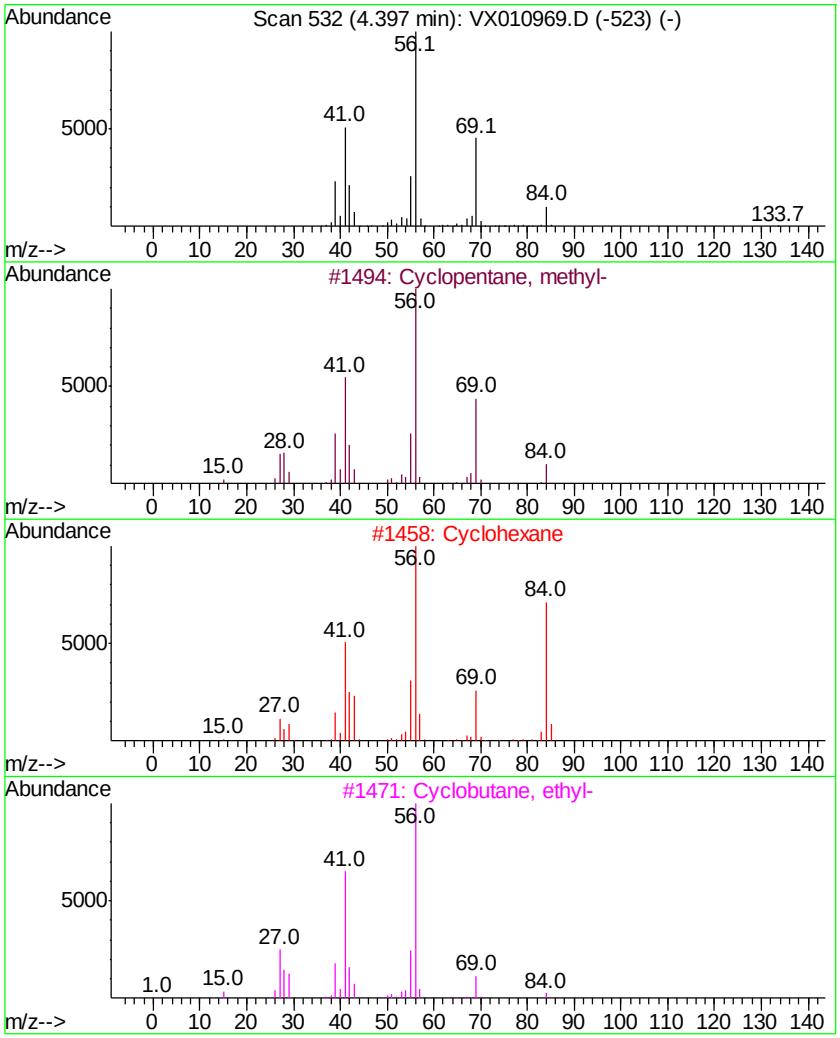
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

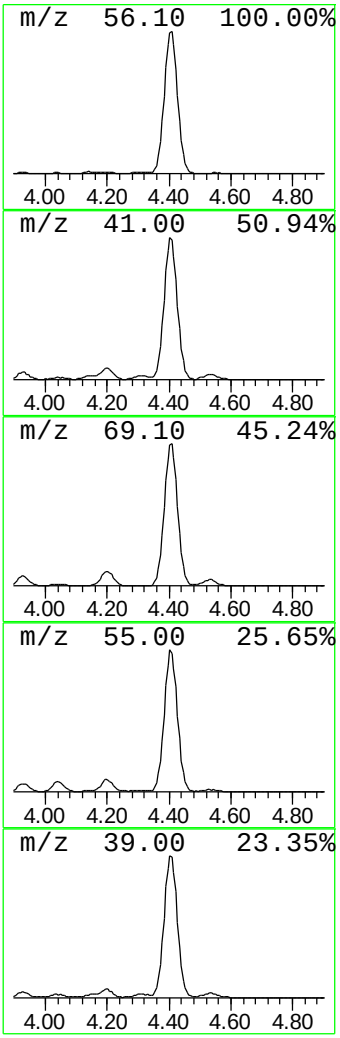
****

Peak Number 5 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	169.61 ug/l	2243540	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	78
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	42





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

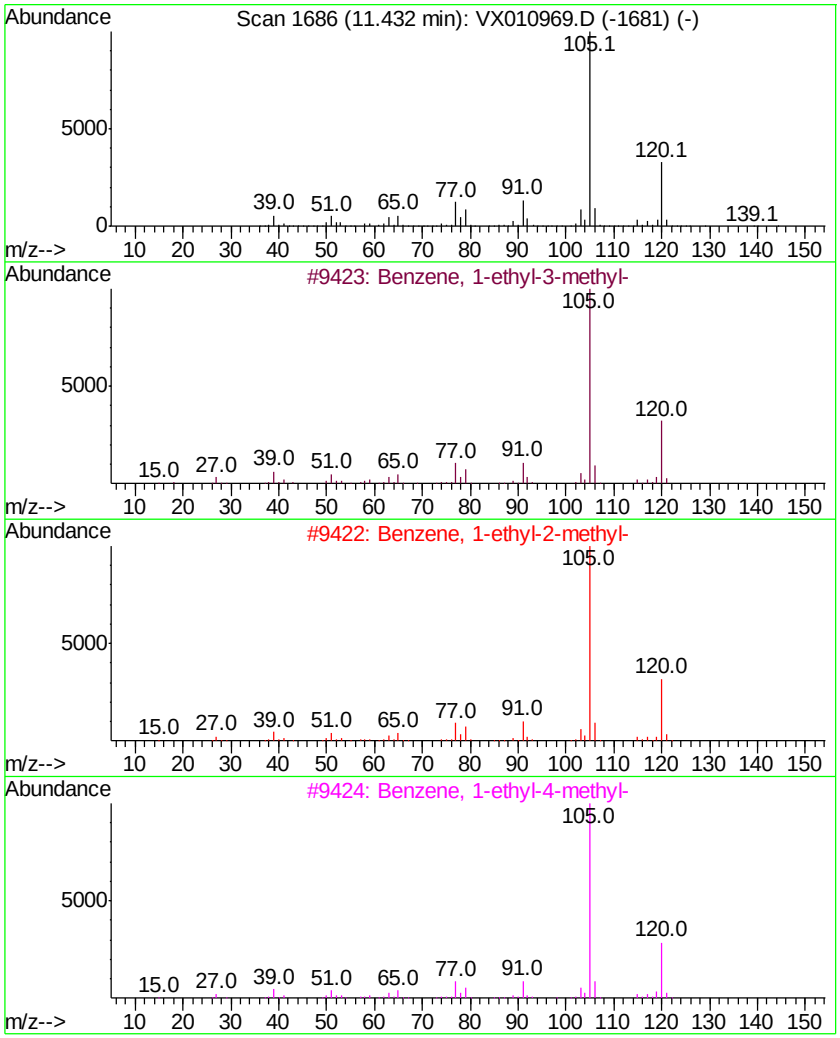
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

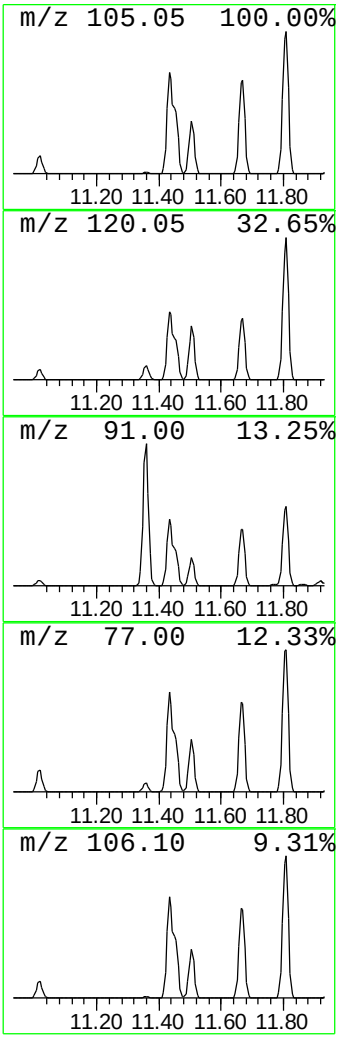
****

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	471.36 ug/l	16065500	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

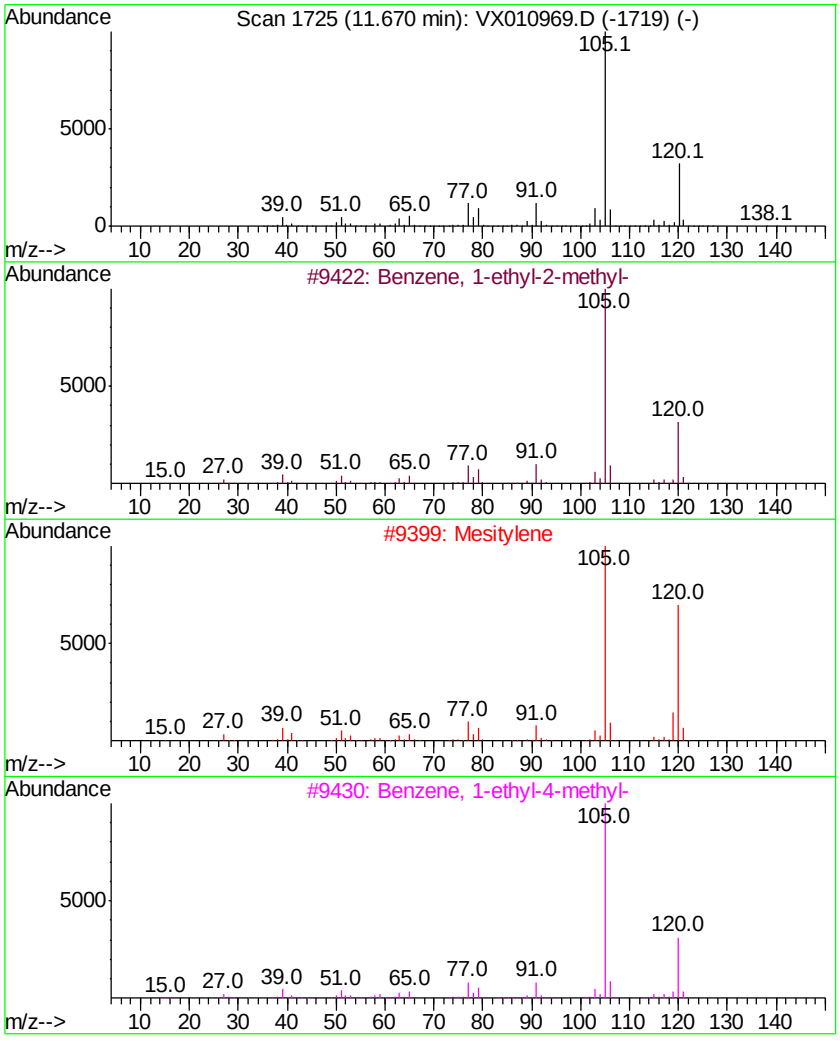
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

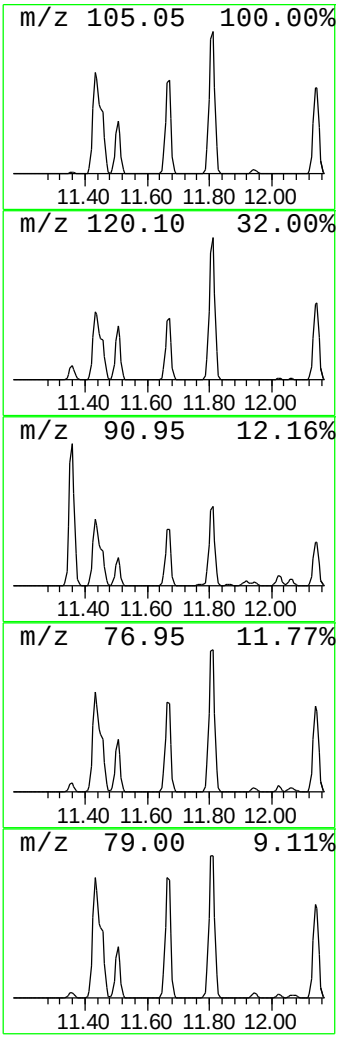
****

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	293.20 ug/l	9993110	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	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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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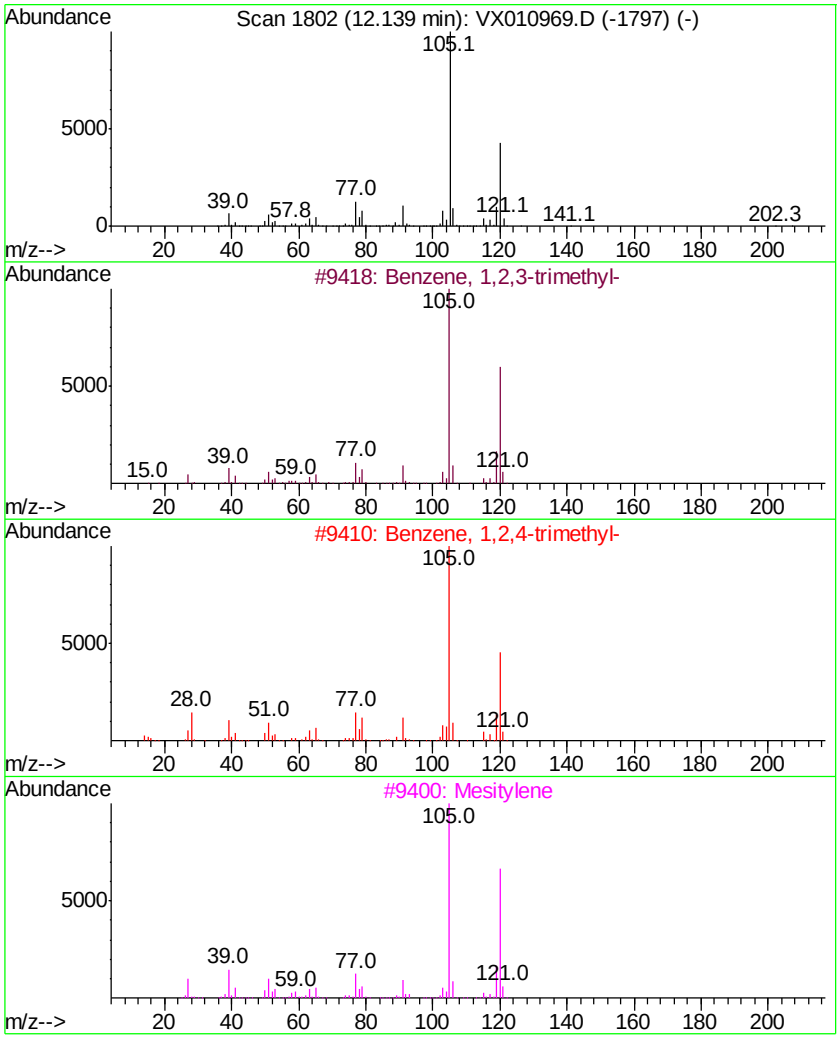
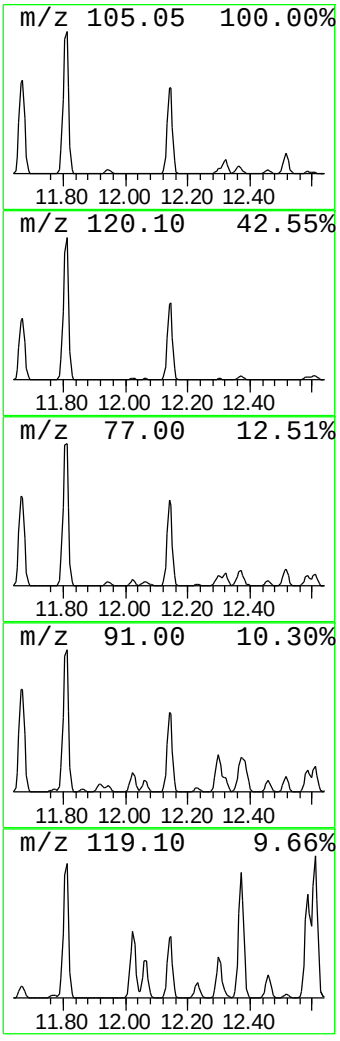
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	285.47 ug/l	9729690	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	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3	Mesitylene	120	C9H12	000108-67-8	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\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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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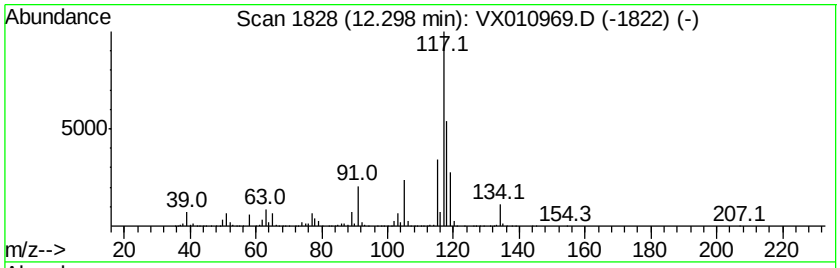
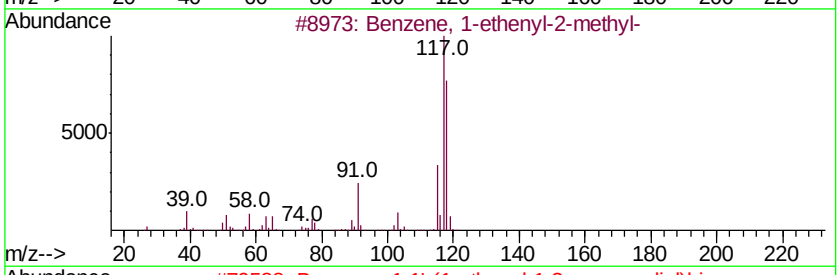
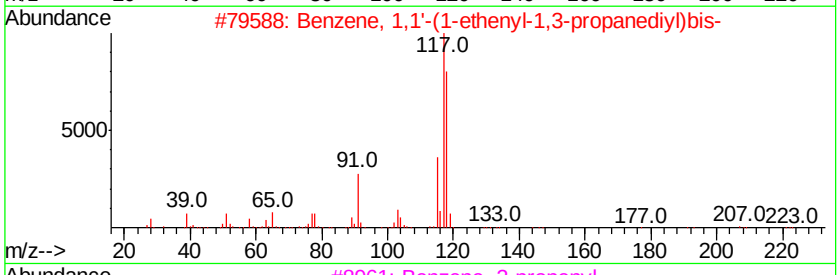
Peak Number 9 Benzene, 1-ethenyl-2-methyl- Concentration Rank 9

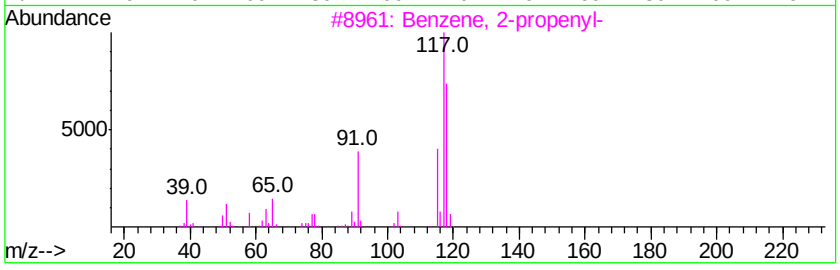
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	104.48 ug/l	3560890	1,4-Dichlorobenzene-d4	12.08

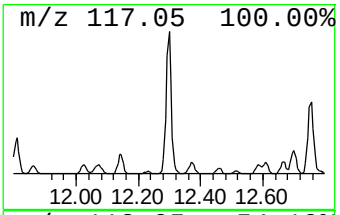
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
2	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	64
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4	Indane	118	C9H10	000496-11-7	64
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64

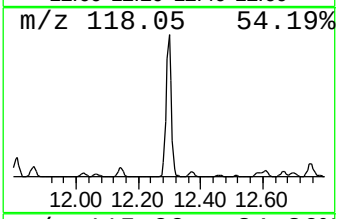


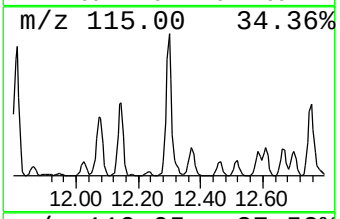


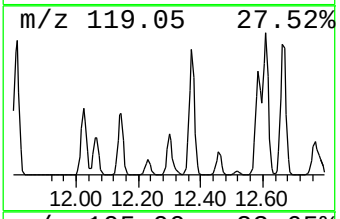


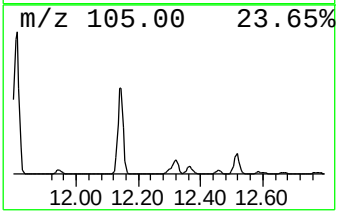












Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
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ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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

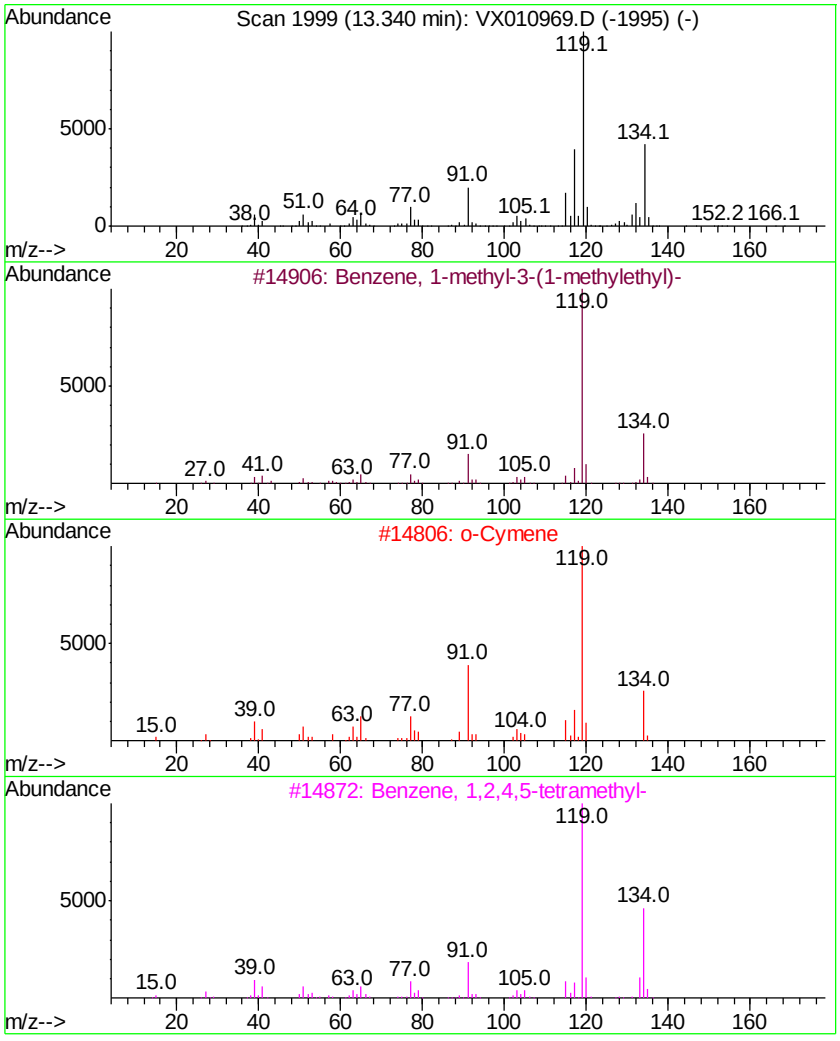
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

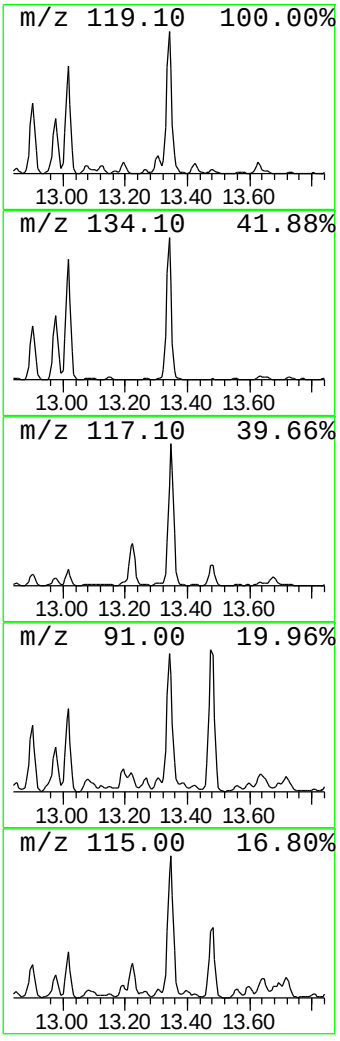
****

Peak Number 10 Benzene, 1-methyl-3-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	109.53 ug/l	3733240	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70
2	o-Cymene	134	C10H14	000527-84-4	70
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	68
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010969.D
Acq On : 18 Jul 2019 19:15
Operator : JC/SP
Sample : K3884-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : FL-0625

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
Pentane	2.00	69.4	ug/l	918198	1	5.67	661367	50.0
Pentane, 2-methyl-	2.89	170.1	ug/l	2249410	1	5.67	661367	50.0
Pentane, 3-methyl-	3.18	132.3	ug/l	1749910	1	5.67	661367	50.0
n-Hexane	3.53	111.3	ug/l	1471970	1	5.67	661367	50.0
Cyclopentane, met...	4.40	169.6	ug/l	2243540	1	5.67	661367	50.0
Benzene, 1-ethyl-...	11.43	471.4	ug/l	16065500	4	12.08	1704170	50.0
Benzene, 1-ethyl-...	11.67	293.2	ug/l	9993110	4	12.08	1704170	50.0
Benzene, 1,2,3-tr...	12.14	285.5	ug/l	9729690	4	12.08	1704170	50.0
Benzene, 1-etheny...	12.30	104.5	ug/l	3560890	4	12.08	1704170	50.0
Benzene, 1-methyl...	13.34	109.5	ug/l	3733240	4	12.08	1704170	50.0