LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

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\82X092618W.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	1.404	37	41	47	rBV	55433	52670	3.74%	0.251%
2	1.611	64	75	76	rBV6	18019	42802	3.04%	0.204%
3	1.770	95	101	110	rBV	290142	423421	30.03%	2.019%
4	1.989	131	137	148	rVB	145382	229579	16.28%	1.094%
5	2.105	151	156	161	rBV	22472	32167	2.28%	0.153%
6	2.257	177	181	189	rVB2	29724	44632	3.17%	0.213%
7	2.836	269	276	278	rVV2	55632	105071	7.45%	0.501%
8	2.885	278	284	295	rVB	167590	378325	26.83%	1.804%
9	3.019	301	306	315	rVB	22544	52082	3.69%	0.248%
10	3.166	321	330	345	rVB2	110768	261760	18.56%	1.248%
11	3.391	359	367	373	rBV2	14941	35805	2.54%	0.171%
12	3.513	379	387	402	rBV	78563	183714	13.03%	0.876%
13	3.653	403	410	417	rVV3	11534	31791	2.25%	0.152%
14	3.806	429	435	445	rVB	23424	56639	4.02%	0.270%
15	3.916	445	453	460	rBV3	13597	32870	2.33%	0.157%
16	4.184	482	497	505	rVV5	18040	52801	3.74%	0.252%
17	4.300	505	516	523	rVV2	40616	125812	8.92%	0.600%
18	4.391	523	531	543	rVV2	75847	225634	16.00%	1.076%
19	5.299	672	680	688	rVB2	16846	52737	3.74%	0.251%
20	5.507	698	714	722	rBV2	182772	549495	38.97%	2.620%
21	5.598	722	729	732	rBV3	46318	110423	7.83%	0.526%
22	5.665	733	740	746	rBV	176082	456373	32.37%	2.176%
23	5.921	772	782	796	rVB2	77580	227340	16.12%	1.084%
24	6.068	796	806	821	rVB	201784	523713	37.14%	2.497%
25	6.257	821	837	841	rBV2	28494	85189	6.04%	0.406%
26	6.342	842	851	862	rVV2	172984	562756	39.91%	2.683%
27	6.452	862	869	880	rVB3	29578	89239	6.33%	0.425%
28	6.702	901	910	918	rBV3	43746	115164	8.17%	0.549%
29	6.860	927	936	944	rBV	363383	901203	63.91%	4.296%
30	7.256	992	1001	1008	rBV4	15986	40043	2.84%	0.191%
31	7.458	1023	1034	1045	rBV3	76395	209945	14.89%	1.001%
32	7.573	1046	1053	1058	rBV	27021	56705	4.02%	0.270%
33	7.634	1058	1063	1074	rVV	34154	79151	5.61%	0.377%
34	7.732	1074	1079	1090	rVB	22776	46552	3.30%	0.222%

LSC Area Percent Report

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Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

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\82X092618W.M
Title : SW846 8260

35	7.848	1090	1098	1106	rVB3	16915	35786	2.54%	0.171%
36	8.055	1122	1132	1144	rBV	98906	217408	15.42%	1.036%
37	8.177	1144	1152	1161	rBV2	104361	257114	18.23%	1.226%
38	8.256	1161	1165	1174	rVB2	40109	79139	5.61%	0.377%
39	8.342	1174	1179	1182	rBV	44378	72229	5.12%	0.344%
40	8.378	1182	1185	1191	rVB2	29012	46809	3.32%	0.223%
41	8.500	1200	1205	1213	rBV	50494	98936	7.02%	0.472%
42	8.573	1213	1217	1226	rVB4	19428	43156	3.06%	0.206%
43	8.677	1226	1234	1235	rBV3	50997	86512	6.14%	0.412%
44	8.713	1235	1240	1248	rVV	831295	1410010	100.00%	6.722%
45	8.787	1248	1252	1257	rVB	64385	102632	7.28%	0.489%
46	8.976	1277	1283	1290	rBV2	45110	87618	6.21%	0.418%
47	9.043	1290	1294	1305	rVB5	18990	45447	3.22%	0.217%
48	9.165	1305	1314	1319	rBV4	18839	41212	2.92%	0.196%
49	9.219	1319	1323	1327	rVV	18766	31877	2.26%	0.152%
50	9.561	1375	1379	1385	rBV4	20466	45644	3.24%	0.218%
51	9.957	1441	1444	1451	rVB	34596	49256	3.49%	0.235%
52	10.067	1457	1462	1465	rBV	32367	44136	3.13%	0.210%
53	10.116	1465	1470	1480	rVV	785913	1139234	80.80%	5.431%
54	10.256	1487	1493	1501	rBV	258667	378947	26.88%	1.807%
55	10.359	1504	1510	1516	rVV	543253	765877	54.32%	3.651%
56	10.414	1516	1519	1531	rVB4	46752	91427	6.48%	0.436%
57	10.701	1560	1566	1580	rVB	112828	186748	13.24%	0.890%
58	10.835	1580	1588	1592	rBV4	18702	34065	2.42%	0.162%
59	11.018	1613	1618	1626	rBV	53241	77511	5.50%	0.370%
60	11.140	1632	1638	1650	rBV	870047	1177626	83.52%	5.614%
61	11.359	1667	1674	1681	rBV	203754	261100	18.52%	1.245%
62	11.439	1681	1687	1688	rVV	260917	355067	25.18%	1.693%
63	11.457	1688	1690	1694	rVV	296892	321593	22.81%	1.533%
64	11.506	1694	1698	1710	rVB	230244	337417	23.93%	1.609%
65	11.670	1720	1725	1732	rBV	224251	297373	21.09%	1.418%
66	11.810	1743	1748	1755	rBV	896468	1113543	78.97%	5.308%
67	11.920	1759	1766	1768	rBV2	23230	34552	2.45%	0.165%
68	11.945	1768	1770	1774	rVB	29381	34092	2.42%	0.163%
69	12.024	1780	1783	1787	rVB	33135	38069	2.70%	0.181%
70	12.079	1787	1792	1798	rBV	1087504	1368342	97.04%	6.523%
71	12.146	1798	1803	1807	rVB	225687	284804	20.20%	1.358%

LSC Area Percent Report

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Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

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\82X092618W.M
Title : SW846 8260

72	12.304	1823	1829	1831	rBV2	226799	330396	23.43%	1.575%
73	12.371	1836	1840	1847	rVB3	246749	397114	28.16%	1.893%
74	12.475	1847	1857	1860	rBV3	36234	65432	4.64%	0.312%
75	12.524	1860	1865	1870	rVB2	90779	142340	10.09%	0.679%
76	12.585	1871	1875	1877	rBV	123141	152026	10.78%	0.725%
77	12.609	1877	1879	1884	rVV	102126	119703	8.49%	0.571%
78	12.670	1884	1889	1892	rVV	227714	267241	18.95%	1.274%
79	12.701	1892	1894	1899	rVB	24994	29696	2.11%	0.142%
80	12.755	1899	1903	1914	rVB4	71795	155938	11.06%	0.743%
81	12.908	1924	1928	1932	rVB	43768	57971	4.11%	0.276%
82	12.975	1932	1939	1943	rBV	101509	152183	10.79%	0.725%
83	13.018	1943	1946	1952	rVB	145511	175560	12.45%	0.837%
84	13.079	1952	1956	1958	rBV	29417	41212	2.92%	0.196%
85	13.103	1958	1960	1962	rVV	27577	30679	2.18%	0.146%
86	13.225	1972	1980	1984	rVB2	94980	146468	10.39%	0.698%
87	13.316	1990	1995	1997	rBV4	34918	57344	4.07%	0.273%
88	13.353	1997	2001	2006	rVB2	150190	209409	14.85%	0.998%
89	13.426	2011	2013	2018	rVB	32168	36343	2.58%	0.173%
90	13.481	2018	2022	2031	rVB2	33296	53147	3.77%	0.253%
91	13.566	2031	2036	2038	rBV2	18492	27845	1.97%	0.133%
92	13.603	2038	2042	2044	rBV	26833	31247	2.22%	0.149%
93	13.639	2044	2048	2054	rVB3	42526	75916	5.38%	0.362%
94	13.725	2059	2062	2068	rVB2	28324	49329	3.50%	0.235%
95	13.834	2075	2080	2086	rVB	152403	189989	13.47%	0.906%
96	13.987	2098	2105	2109	rBV4	20967	38625	2.74%	0.184%
97	14.133	2125	2129	2135	rBV	31289	38300	2.72%	0.183%
98	14.273	2148	2152	2162	rVB	29197	49667	3.52%	0.237%
99	14.694	2213	2221	2227	rBV	139539	190174	13.49%	0.907%
100	14.840	2239	2245	2254	rVB	71881	99604	7.06%	0.475%

Sum of corrected areas: 20976809

LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
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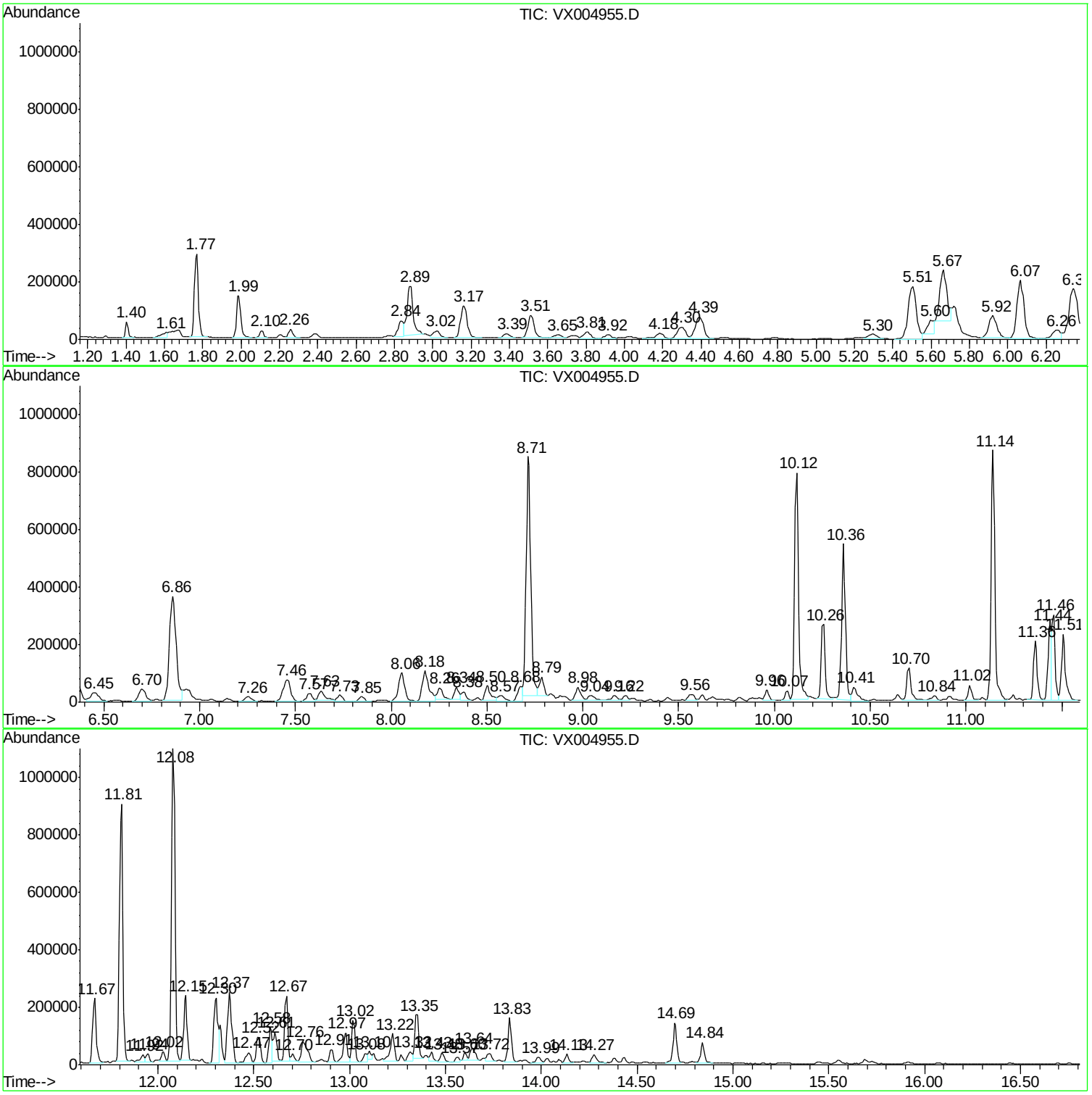
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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\VX100318\
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Quant Title : SW846 8260

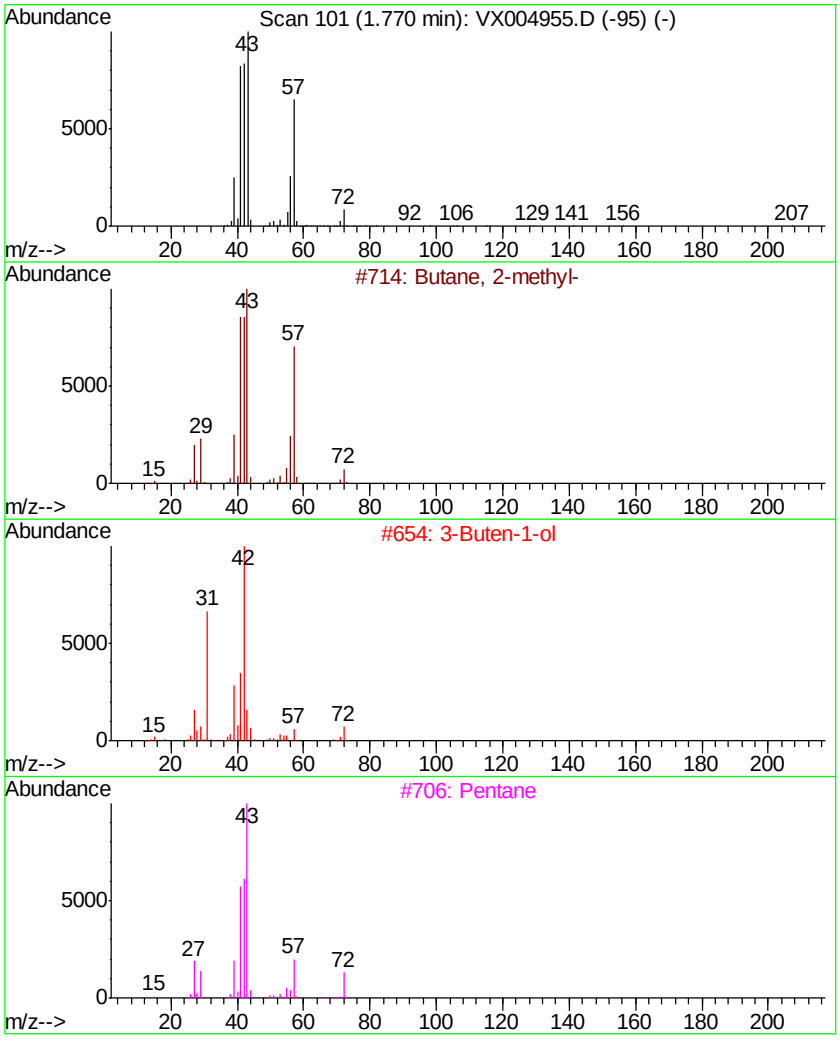
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

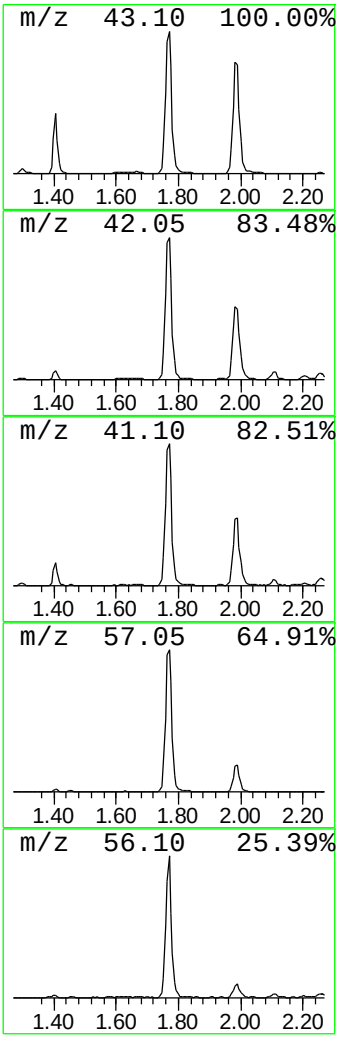
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.77	46.39 ug/l	423421	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	47
3	Pentane	72	C5H12	000109-66-0	45
4	1-Butene	56	C4H8	000106-98-9	9
5	Azetidine	57	C3H7N	000503-29-7	9





Library Search Compound Report

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ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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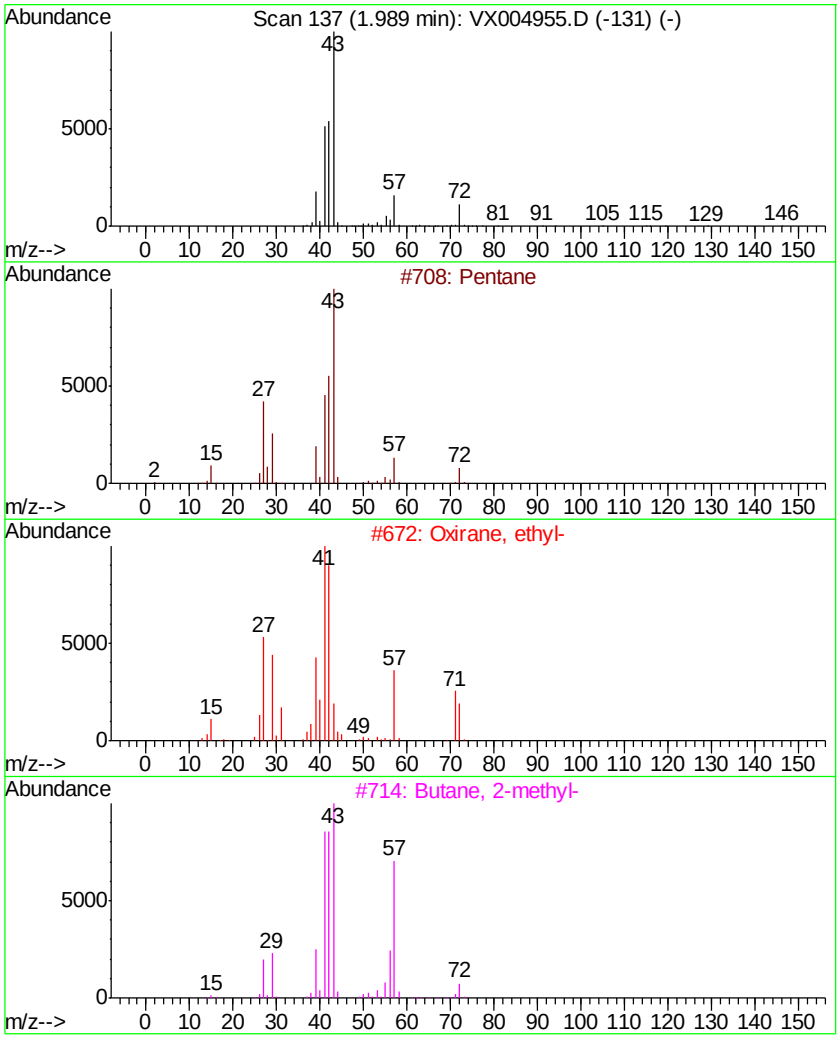
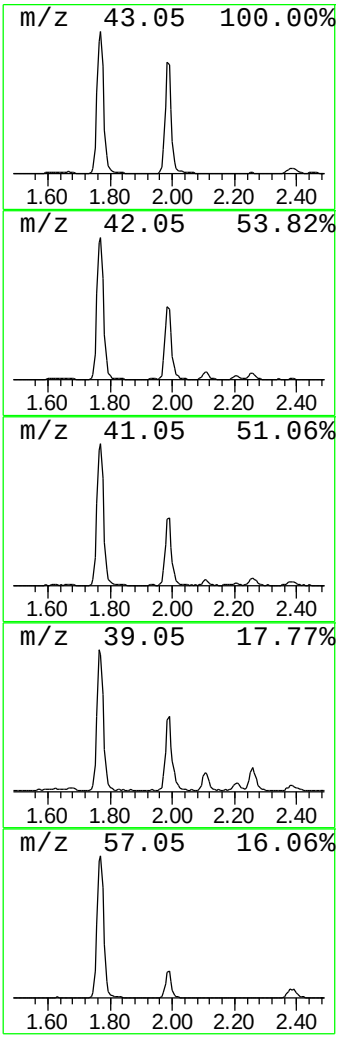
Peak Number 2 Pentane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.99	25.15 ug/l	229579	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	90
2	Oxirane, ethyl-	72	C4H8O	000106-88-7	50
3	Butane, 2-methyl-	72	C5H12	000078-78-4	42
4	Tetrahydrofuran	72	C4H8O	000109-99-9	9
5	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	9





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ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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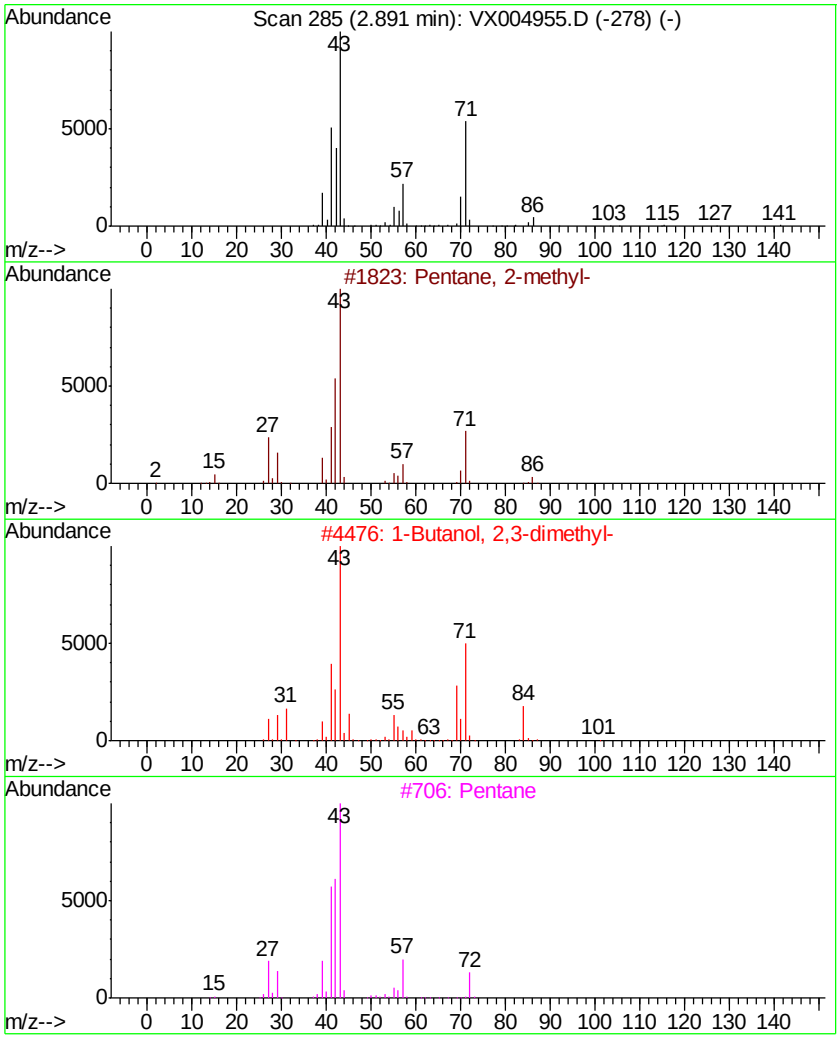
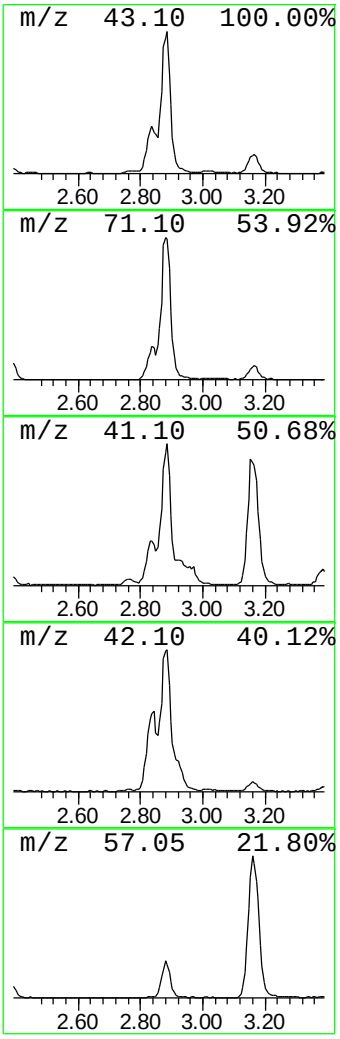
Peak Number 3 Pentane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	41.45 ug/l	378325	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	Pentane	72	C5H12	000109-66-0	38
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28
5	Isobutane	58	C4H10	000075-28-5	25





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ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

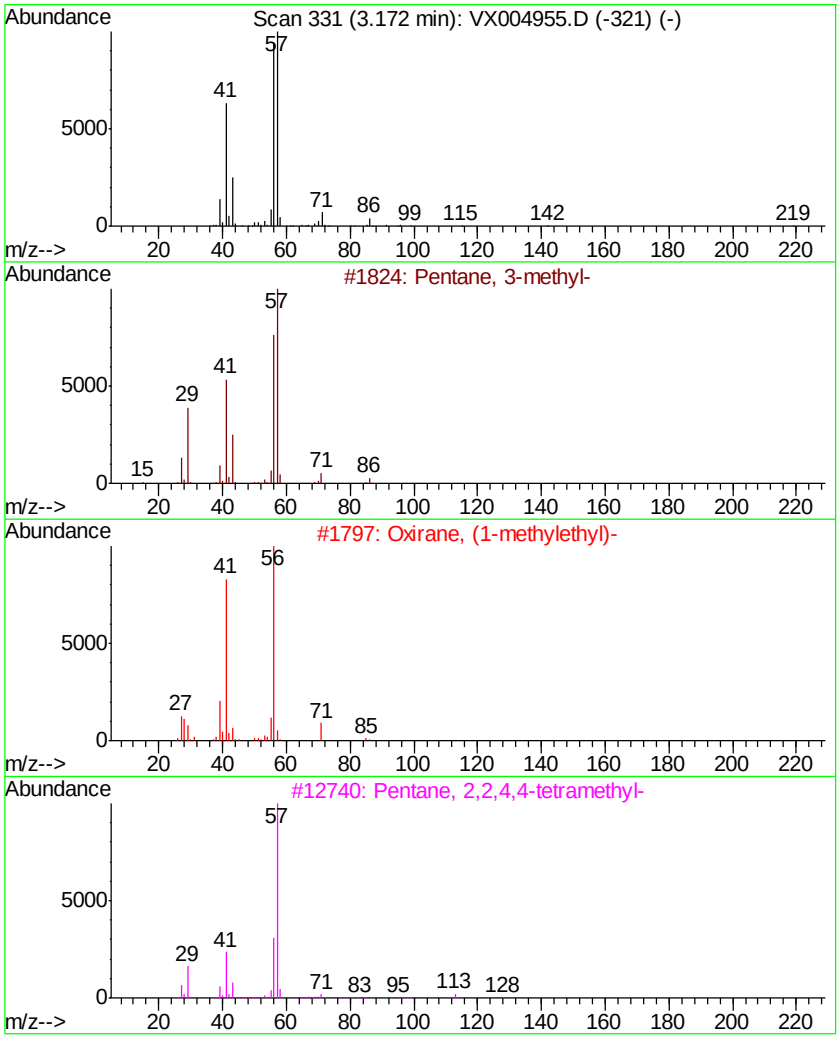
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

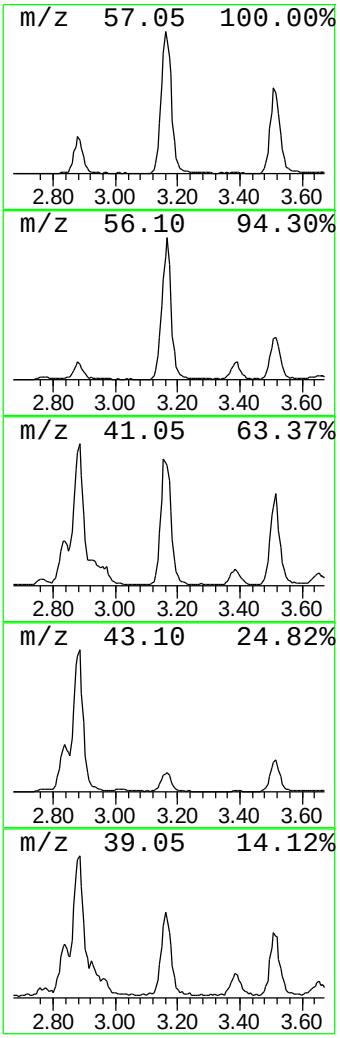
****

Peak Number 4 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	28.68 ug/l	261760	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	59
3	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	42
4	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40
5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	39





Library Search Compound Report

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Operator : JC/MD
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Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

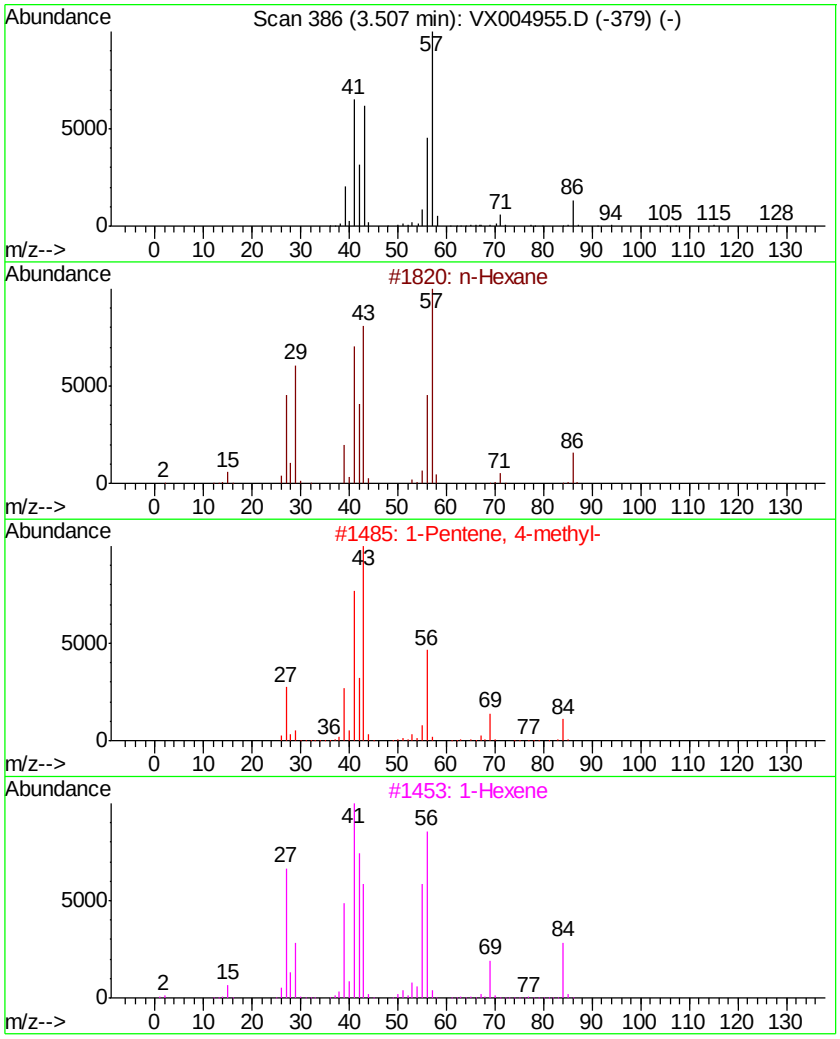
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

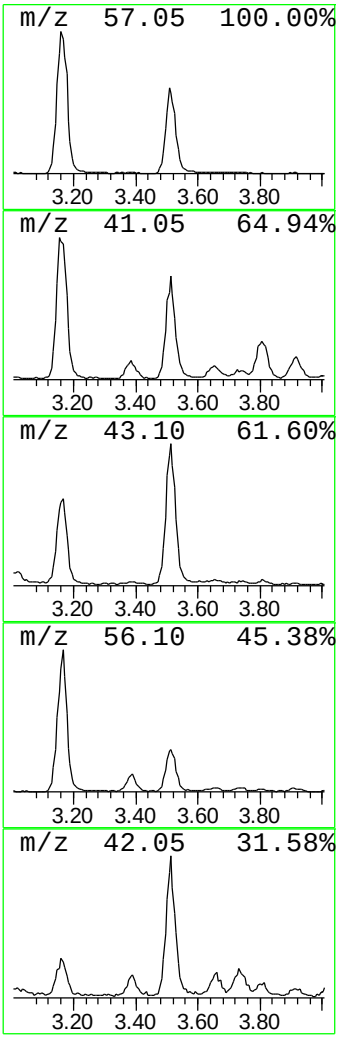
****

Peak Number 5 n-Hexane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.51	20.13 ug/l	183714	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	91
2	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	43
3	1-Hexene	84	C6H12	000592-41-6	38
4	2(3H)-Furanone, dihydro-3-methyl-	100	C5H8O2	001679-47-6	38
5	Butanal, 2-methyl-	86	C5H10O	000096-17-3	35





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

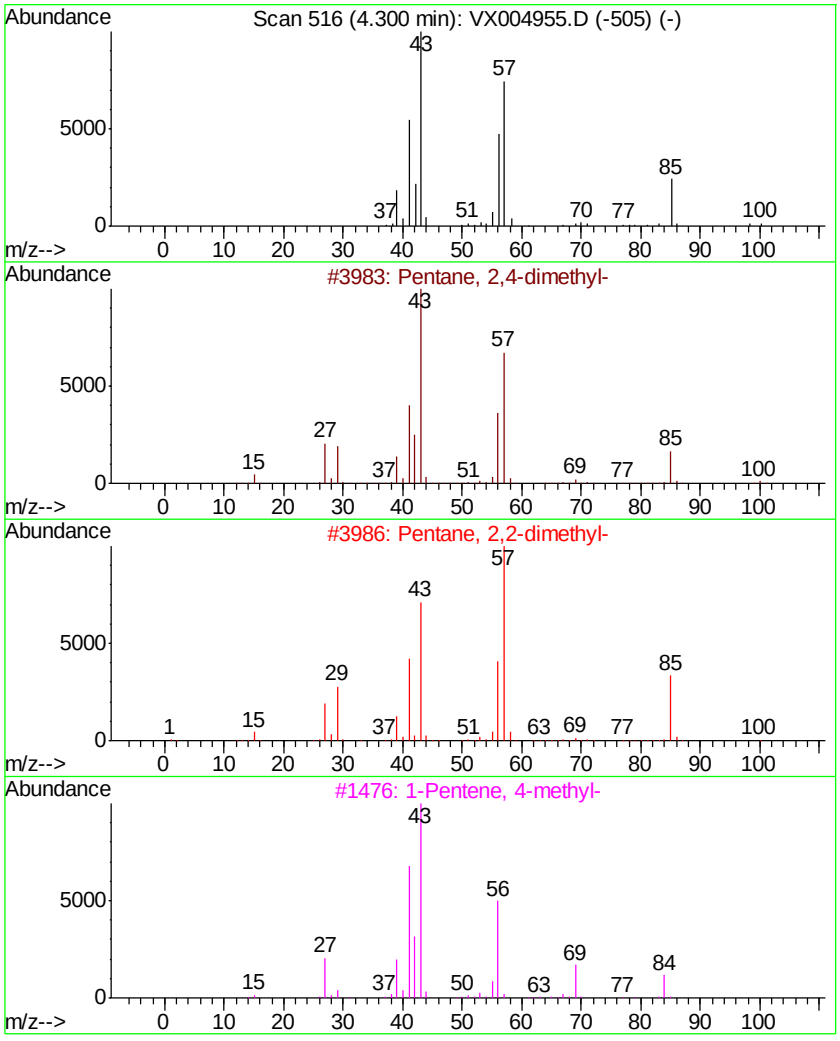
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

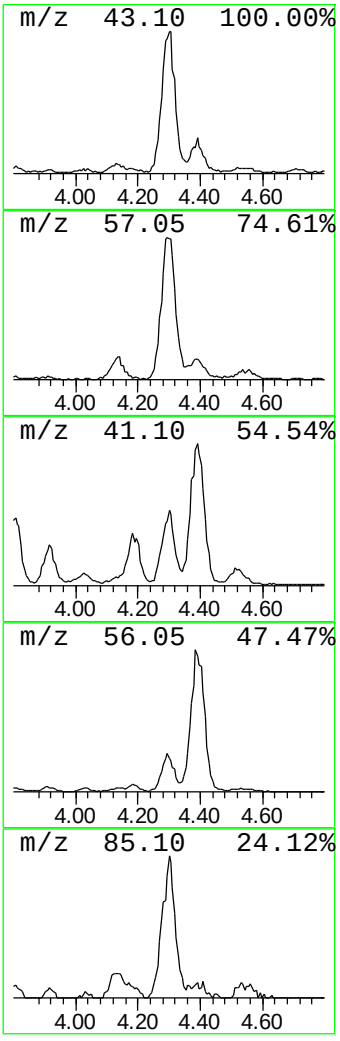
****

Peak Number 6 Pentane, 2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.30	13.78 ug/l	125812	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
3	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	47
4	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	42
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	42





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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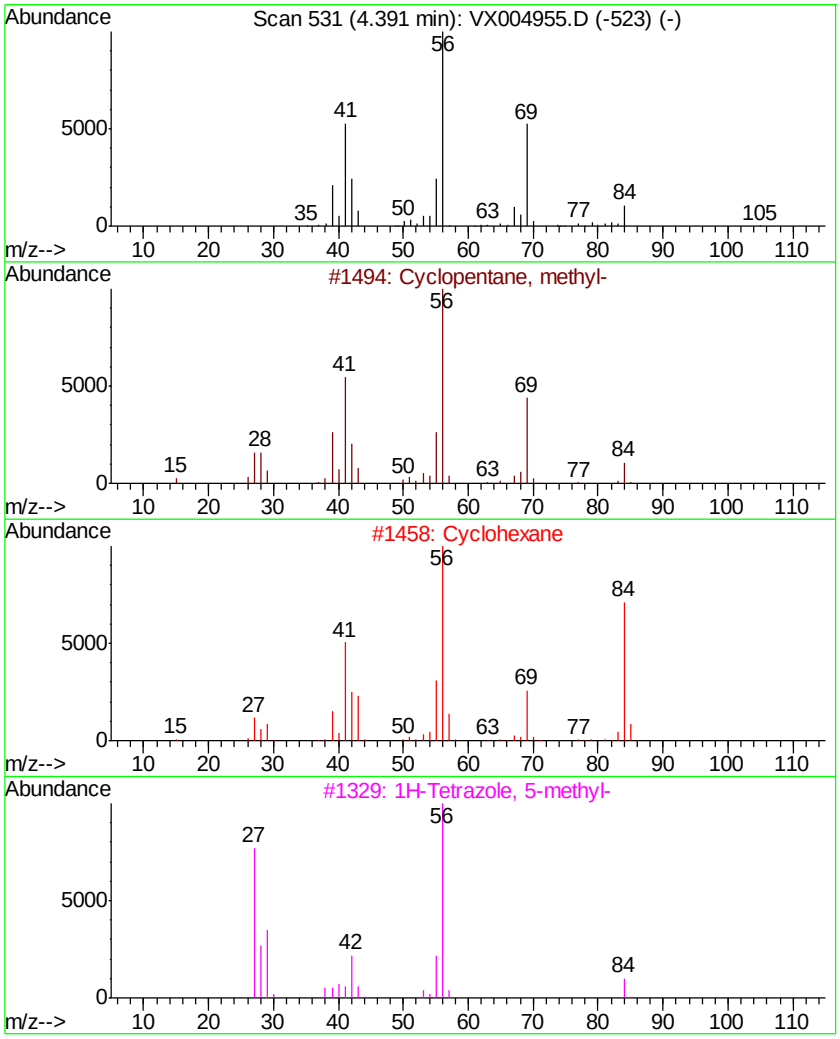
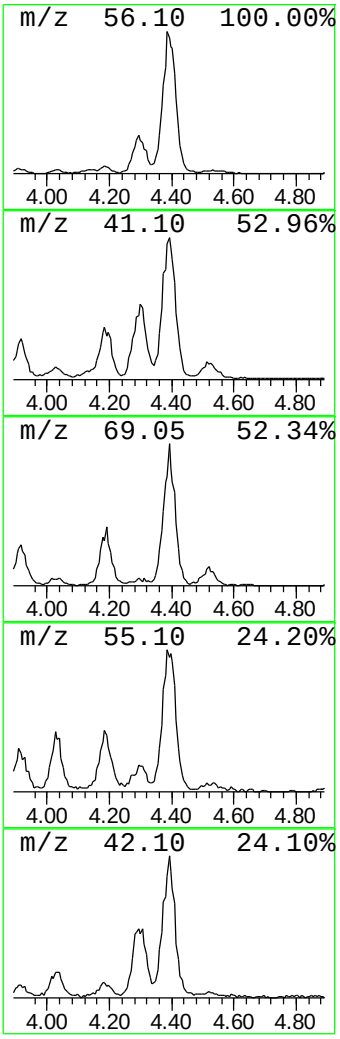
Peak Number 7 Cyclopentane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	24.72 ug/l	225634	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclohexane	84	C6H12	000110-82-7	78
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Tetramethylene diisocyanate	140	C6H8N2O2	004538-37-8	40





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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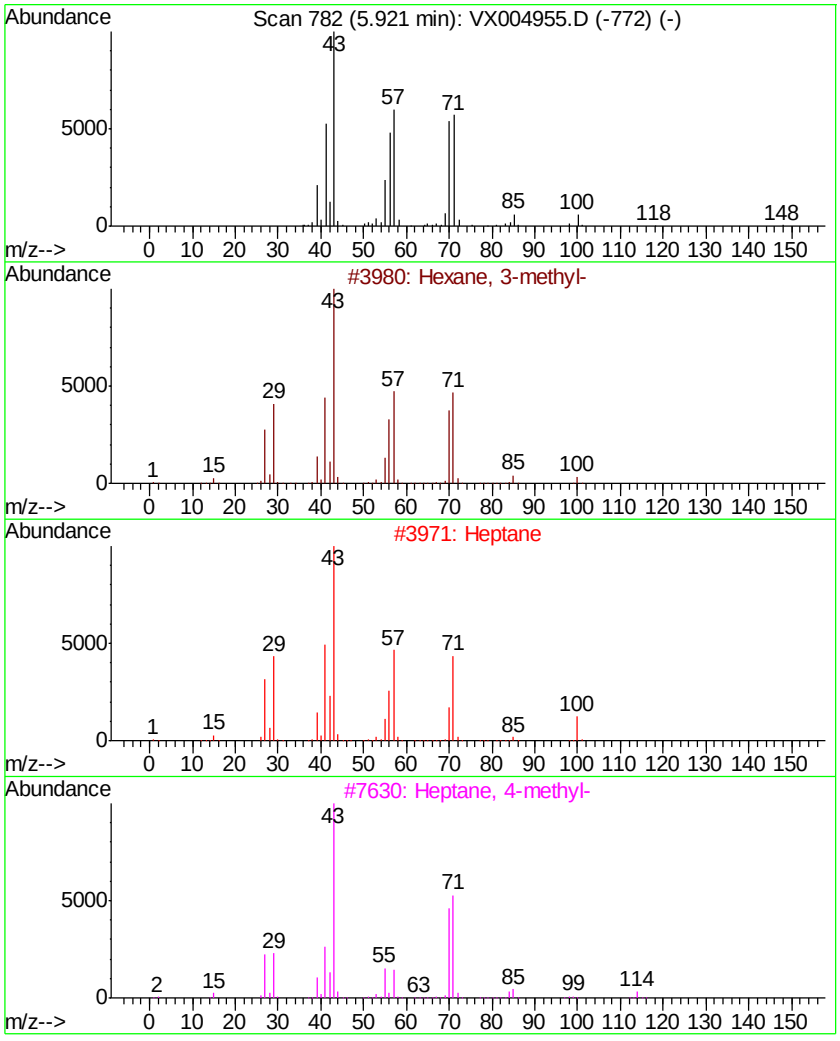
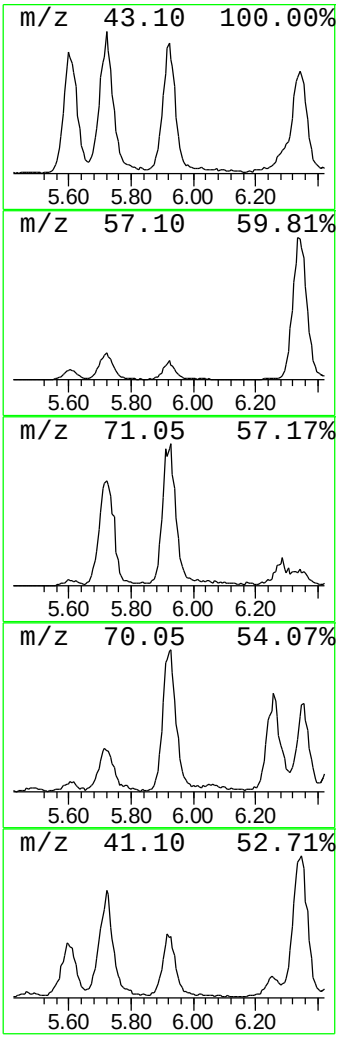
Peak Number 8 Hexane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	24.91 ug/l	227340	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2	Heptane	100	C7H16	000142-82-5	72
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	53
4	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	53
5	Dodecane, 4-methyl-	184	C13H28	006117-97-1	47





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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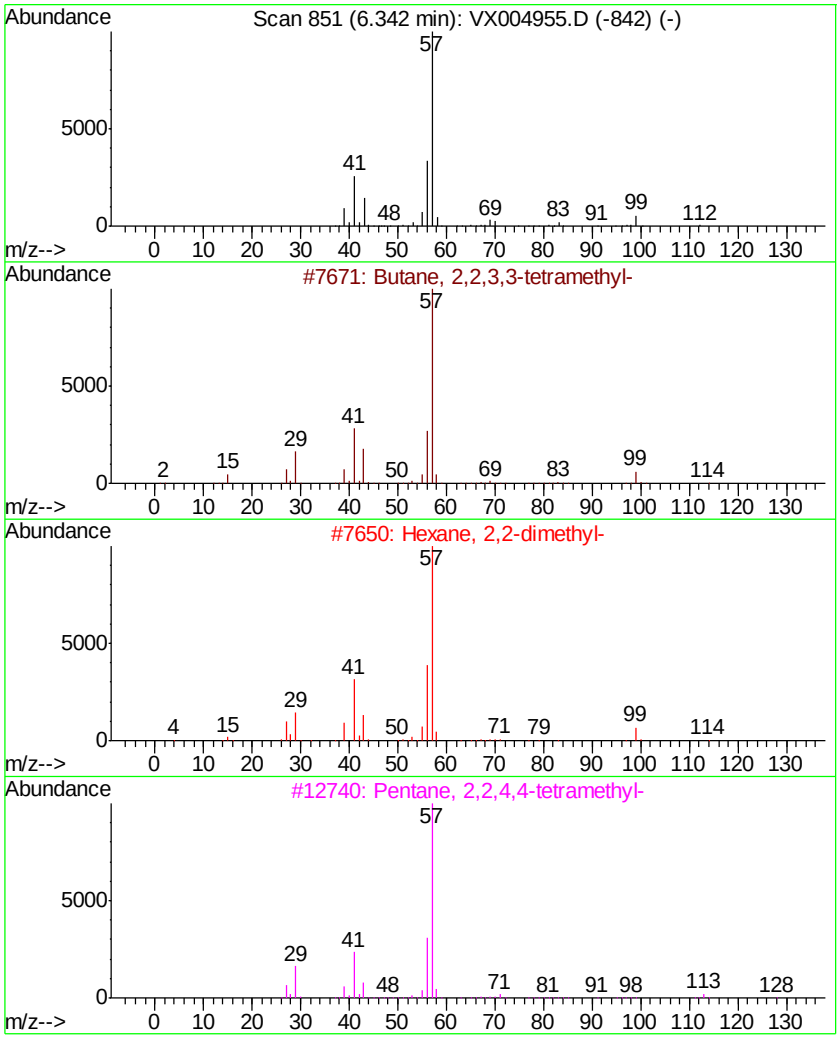
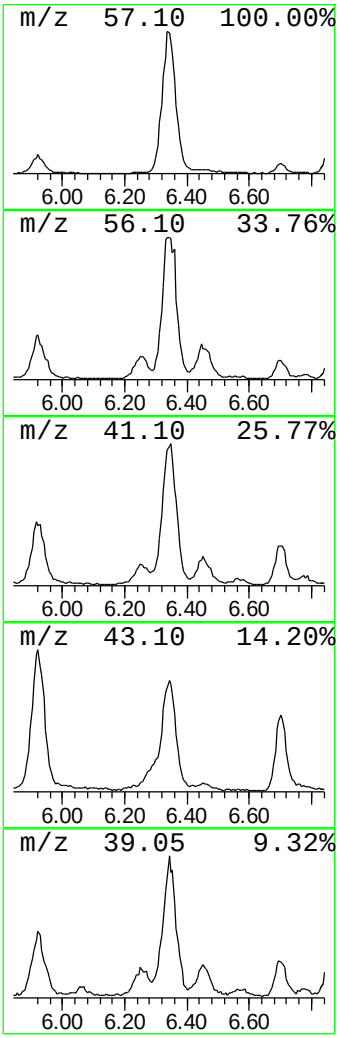
Peak Number 9 Butane, 2,2,3,3-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	31.22 ug/l	562756	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	64
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
3	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	50
5	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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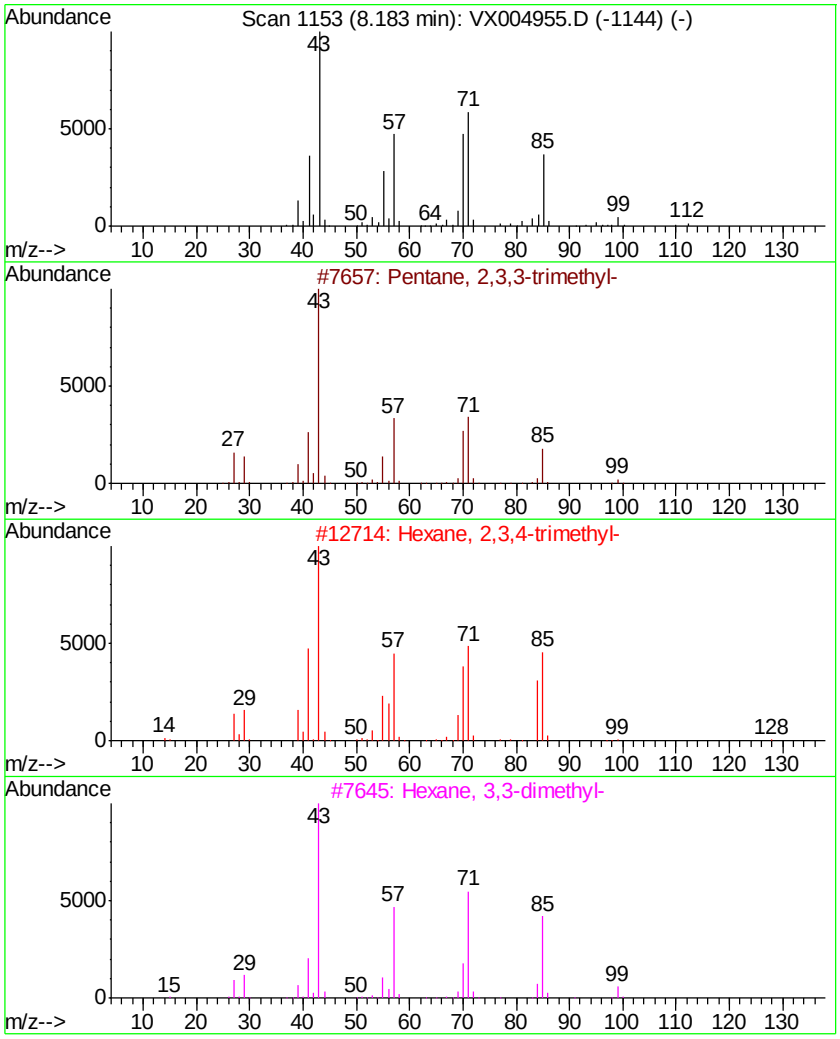
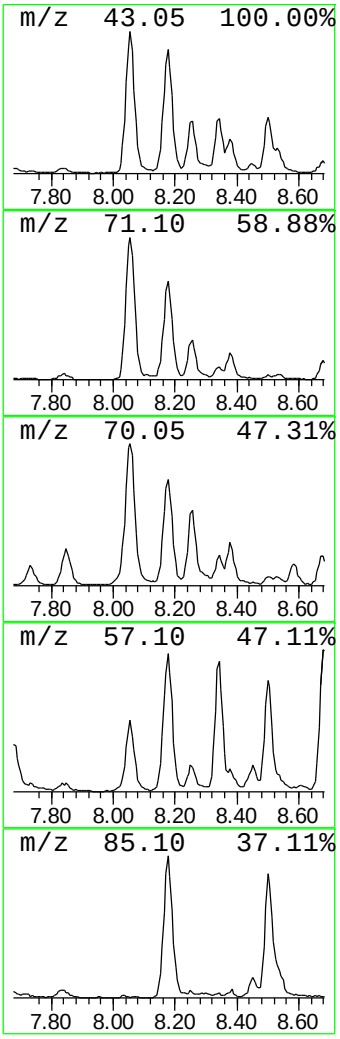
Peak Number 10 Pentane, 2,3,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	14.27 ug/l	257114	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	83
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004955.D
Acq On : 03 Oct 2018 08:53
Operator : JC/MD
Sample : J5084-03 10X
Misc : 5.82g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-9(11.5-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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-	1.77	46.4	ug/l	423421	1	5.67	456373	50.0
Pentane	1.99	25.1	ug/l	229579	1	5.67	456373	50.0
Pentane, 2-methyl-	2.89	41.5	ug/l	378325	1	5.67	456373	50.0
Pentane, 3-methyl-	3.17	28.7	ug/l	261760	1	5.67	456373	50.0
n-Hexane	3.51	20.1	ug/l	183714	1	5.67	456373	50.0
Pentane, 2,4-dime...	4.30	13.8	ug/l	125812	1	5.67	456373	50.0
Cyclopentane, met...	4.39	24.7	ug/l	225634	1	5.67	456373	50.0
Hexane, 3-methyl-	5.92	24.9	ug/l	227340	1	5.67	456373	50.0
Butane, 2,2,3,3-t...	6.34	31.2	ug/l	562756	2	6.86	901203	50.0
Pentane, 2,3,3-tr...	8.18	14.3	ug/l	257114	2	6.86	901203	50.0