LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

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\82X061919W.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	1.276	16	20	36	rBV	301821	524712	41.18%	4.558%
2	1.404	38	41	44	rVV	34680	36437	2.86%	0.317%
3	1.788	100	104	114	rVB	182529	271583	21.31%	2.359%
4	2.001	134	139	150	rVB2	68694	119871	9.41%	1.041%
5	2.269	177	183	191	rVB	122060	201772	15.84%	1.753%
6	2.397	198	204	208	rBV	13838	27455	2.15%	0.238%
7	2.440	208	211	219	rVB	7927	13994	1.10%	0.122%
8	2.605	232	238	246	rVB	7165	13142	1.03%	0.114%
9	2.842	266	277	281	rBV2	37556	121702	9.55%	1.057%
10	2.891	281	285	288	rVV	47166	93310	7.32%	0.811%
11	2.934	288	292	305	rVB2	65391	145717	11.44%	1.266%
12	3.178	323	332	340	rBV2	42715	106525	8.36%	0.925%
13	3.403	360	369	376	rBV8	6068	16571	1.30%	0.144%
14	3.531	382	390	399	rBV2	16240	38849	3.05%	0.337%
15	3.818	427	437	445	rBV	29226	72086	5.66%	0.626%
16	3.928	448	455	462	rVV3	17564	46549	3.65%	0.404%
17	4.196	488	499	509	rBV5	20209	59416	4.66%	0.516%
18	4.312	511	518	523	rVV4	6979	20005	1.57%	0.174%
19	4.409	525	534	545	rVB	52133	151248	11.87%	1.314%
20	4.531	546	554	564	rVB3	10786	32012	2.51%	0.278%
21	5.305	670	681	692	rVB3	27051	95266	7.48%	0.828%
22	5.501	701	713	721	rBV	129971	374955	29.43%	3.257%
23	5.580	721	726	731	rVV4	24783	75286	5.91%	0.654%
24	5.659	731	739	765	rVB	213977	694001	54.47%	6.028%
25	5.933	772	784	796	rBV6	11619	40687	3.19%	0.353%
26	6.061	796	805	811	rVV	122990	326994	25.66%	2.840%
27	6.141	811	818	832	rVV	479602	1274195	100.00%	11.068%
28	6.305	832	845	846	rVV5	25093	73785	5.79%	0.641%
29	6.348	848	852	862	rVV3	36107	100169	7.86%	0.870%
30	6.458	862	870	882	rVB5	16881	49237	3.86%	0.428%
31	6.708	902	911	922	rVB5	4992	15533	1.22%	0.135%
32	6.860	925	936	945	rBV	326728	784795	61.59%	6.817%
33	6.933	945	948	960	rVV7	16693	46224	3.63%	0.402%
34	7.262	993	1002	1009	rVB2	18788	42621	3.34%	0.370%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

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\82X061919W.M
Title : SW846 8260

35	7.458	1023	1034	1044	rVB4	31339	91129	7.15%	0.792%
36	7.573	1044	1053	1058	rBV9	5596	12789	1.00%	0.111%
37	7.957	1113	1116	1123	rVV6	11973	27615	2.17%	0.240%
38	8.055	1123	1132	1140	rVB2	19563	49782	3.91%	0.432%
39	8.177	1145	1152	1155	rBV2	27412	58711	4.61%	0.510%
40	8.213	1155	1158	1163	rVV	29975	48446	3.80%	0.421%
41	8.378	1177	1185	1192	rBV10	5347	12937	1.02%	0.112%
42	8.567	1211	1216	1225	rVB3	29774	58983	4.63%	0.512%
43	8.713	1227	1240	1247	rBV	701688	1108631	87.01%	9.630%
44	8.780	1247	1251	1257	rVB	44839	68952	5.41%	0.599%
45	8.902	1264	1271	1280	rBV5	9633	22759	1.79%	0.198%
46	8.988	1280	1285	1289	rVV	8395	14103	1.11%	0.123%
47	9.165	1308	1314	1319	rVB2	10722	17372	1.36%	0.151%
48	9.219	1319	1323	1334	rVB	26550	44299	3.48%	0.385%
49	9.518	1369	1372	1376	rVV2	9194	14869	1.17%	0.129%
50	9.561	1376	1379	1385	rVB7	7668	15256	1.20%	0.133%
51	9.817	1414	1421	1427	rBV6	8742	16963	1.33%	0.147%
52	10.109	1464	1469	1479	rBV	700533	932713	73.20%	8.102%
53	10.250	1488	1492	1497	rVB	33608	42323	3.32%	0.368%
54	10.353	1503	1509	1516	rVB	126247	178873	14.04%	1.554%
55	10.640	1548	1556	1562	rBV10	5635	16917	1.33%	0.147%
56	11.018	1612	1618	1623	rBV	99040	129837	10.19%	1.128%
57	11.134	1632	1637	1643	rBV	605563	759581	59.61%	6.598%
58	11.359	1669	1674	1681	rVB	67367	83646	6.56%	0.727%
59	11.457	1684	1690	1694	rBV	10641	16199	1.27%	0.141%
60	11.664	1719	1724	1734	rVB	54421	75415	5.92%	0.655%
61	11.920	1761	1766	1768	rBV	9603	13734	1.08%	0.119%
62	12.073	1786	1791	1799	rBV	707282	916026	71.89%	7.957%
63	12.140	1799	1802	1806	rVB	14976	17912	1.41%	0.156%
64	12.298	1820	1828	1835	rBV	166641	215139	16.88%	1.869%
65	12.365	1835	1839	1847	rVB4	12795	24099	1.89%	0.209%
66	12.518	1859	1864	1869	rVB	15820	19840	1.56%	0.172%
67	12.664	1883	1888	1892	rBV	41155	51123	4.01%	0.444%
68	12.755	1899	1903	1910	rVB2	34250	50728	3.98%	0.441%
69	12.975	1932	1939	1943	rBV	28765	37782	2.97%	0.328%
70	13.225	1976	1980	1986	rVB	22526	32812	2.58%	0.285%
71	13.347	1995	2000	2006	rVB3	51964	72843	5.72%	0.633%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

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\82X061919W.M
Title : SW846 8260

72	14.694	2212	2221	2227	rBV	11066	17043	1.34%	0.148%
73	14.834	2240	2244	2252	rVB	10756	17291	1.36%	0.150%

Sum of corrected areas: 11512176

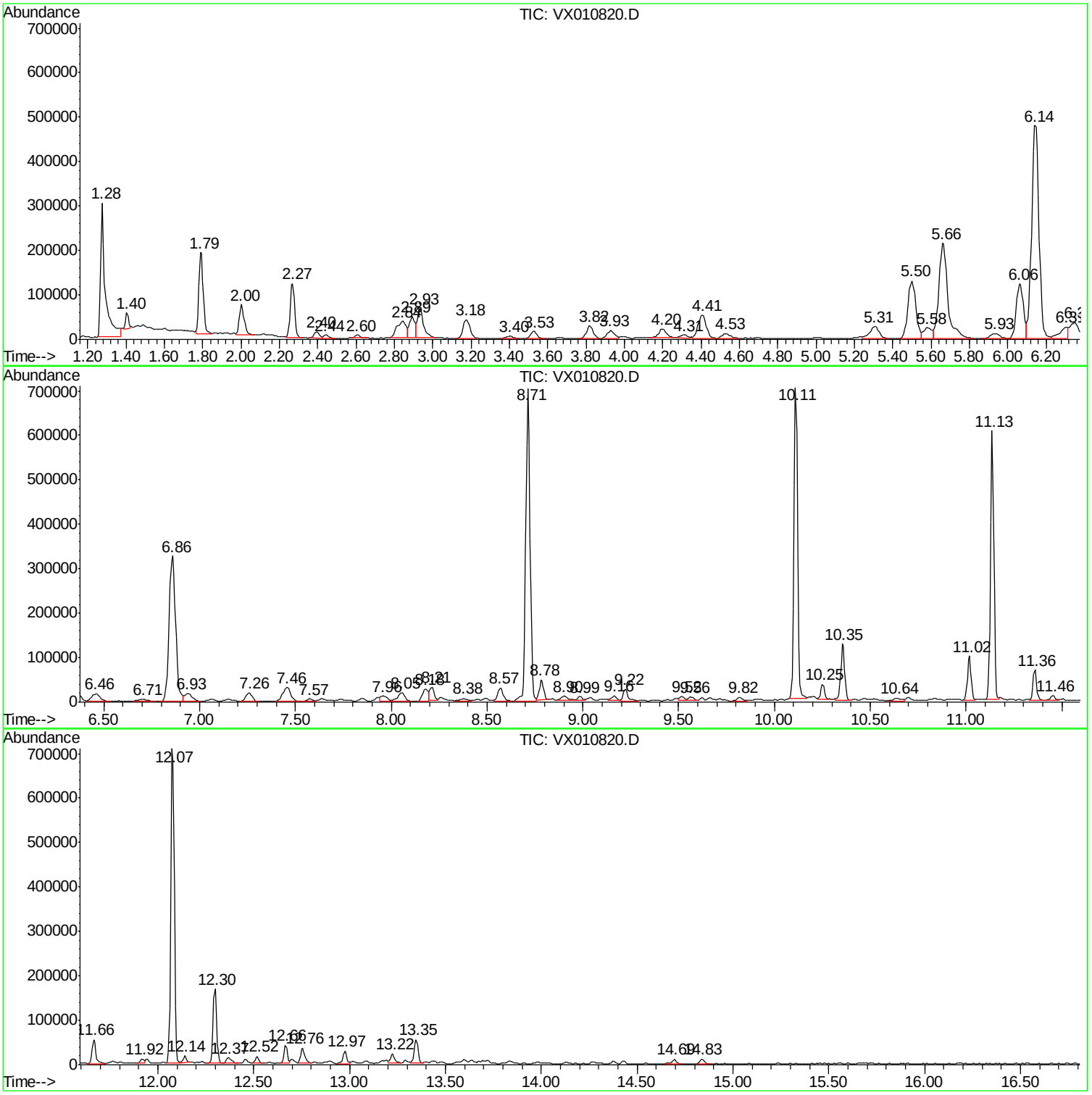
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

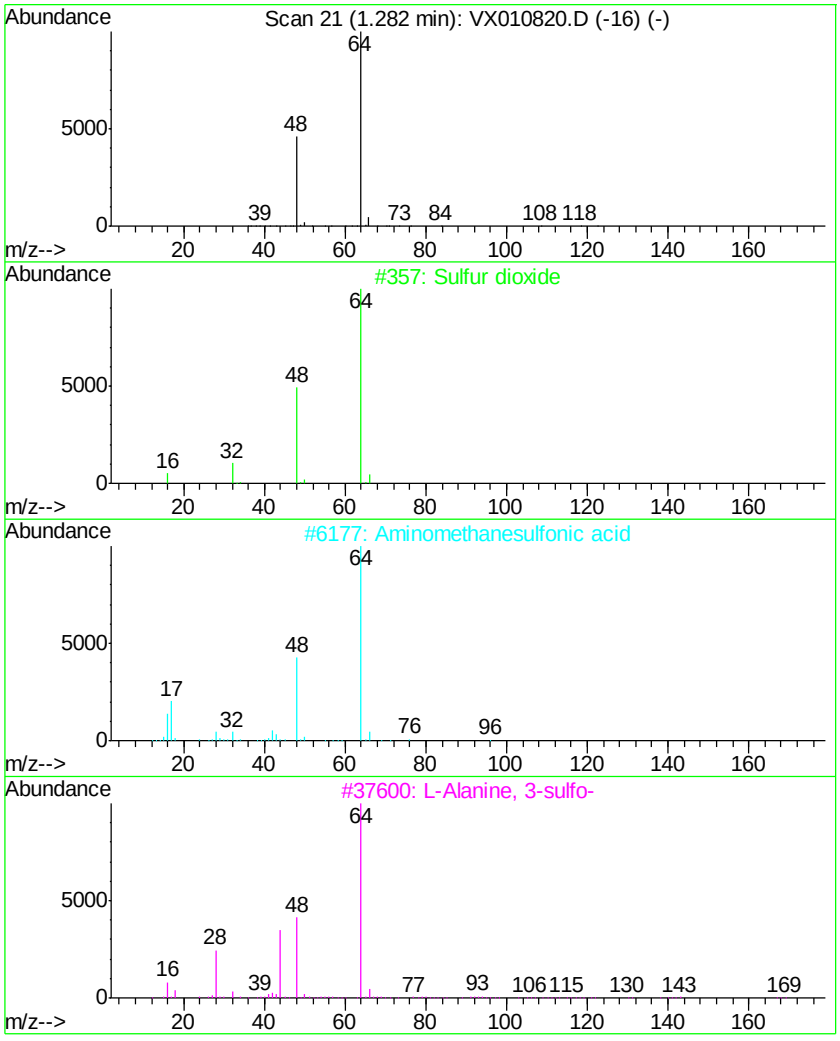
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

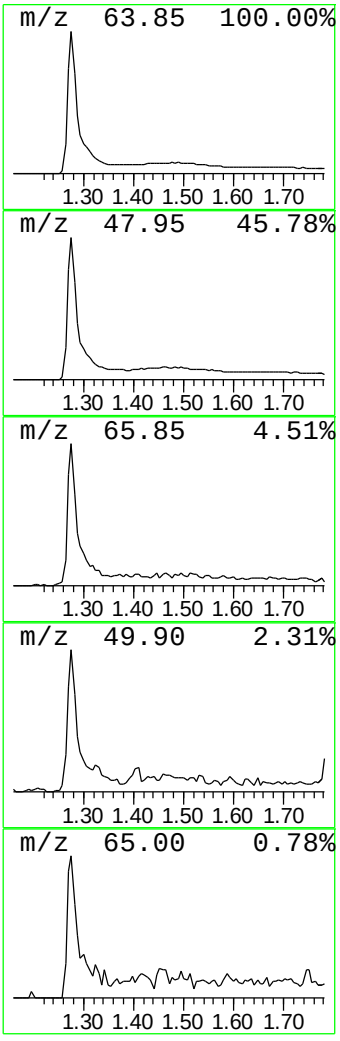
****

Peak Number 1 Sulfur dioxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.28	37.80 ug/l	524712	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4
5	Ethyl Chloride	64	C2H5Cl	000075-00-3	3





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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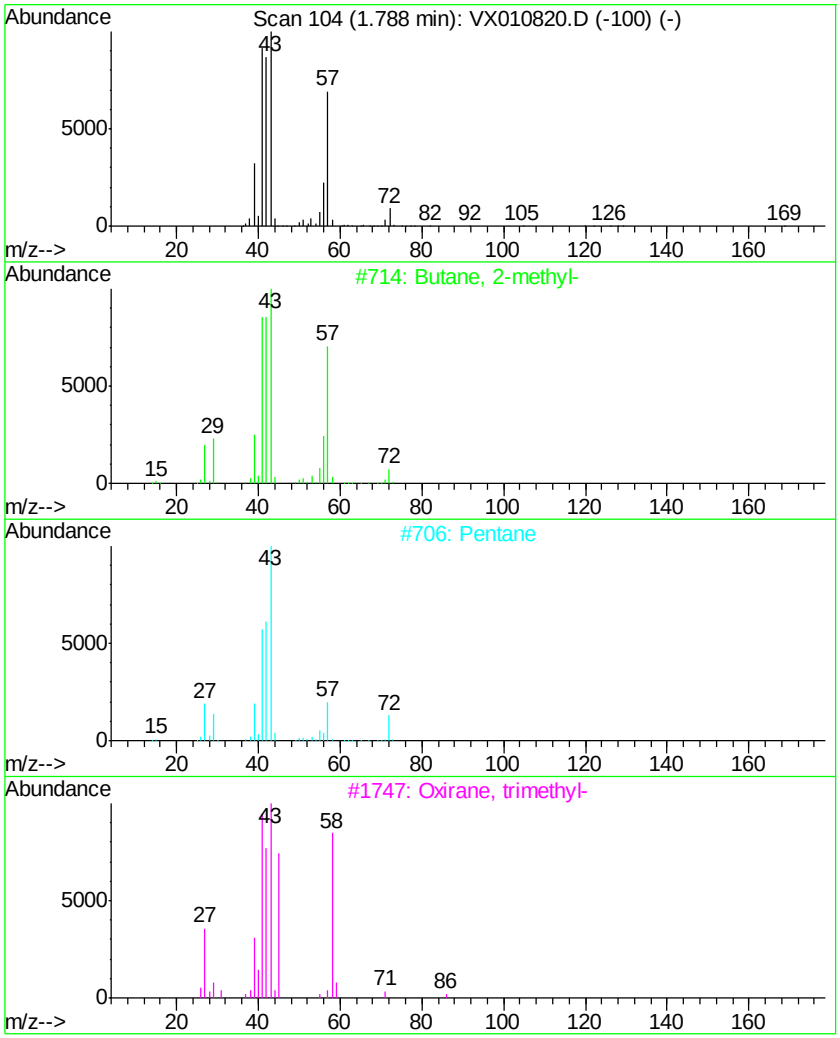
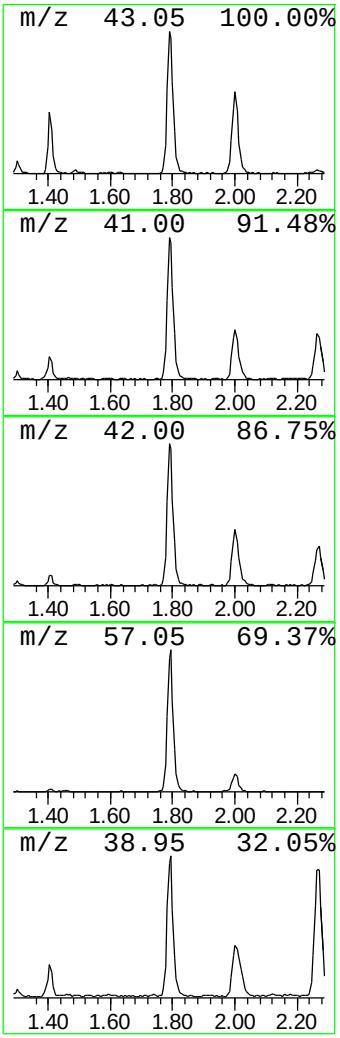
Peak Number 2 Butane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.79	19.57 ug/l	271583	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Pentane	72	C5H12	000109-66-0	37
3	Oxirane, trimethyl-	86	C5H10O	005076-19-7	36
4	3-Buten-1-ol	72	C4H8O	000627-27-0	28
5	2,5-Furandione, dihydro-3-methyl-	114	C5H6O3	004100-80-5	23





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

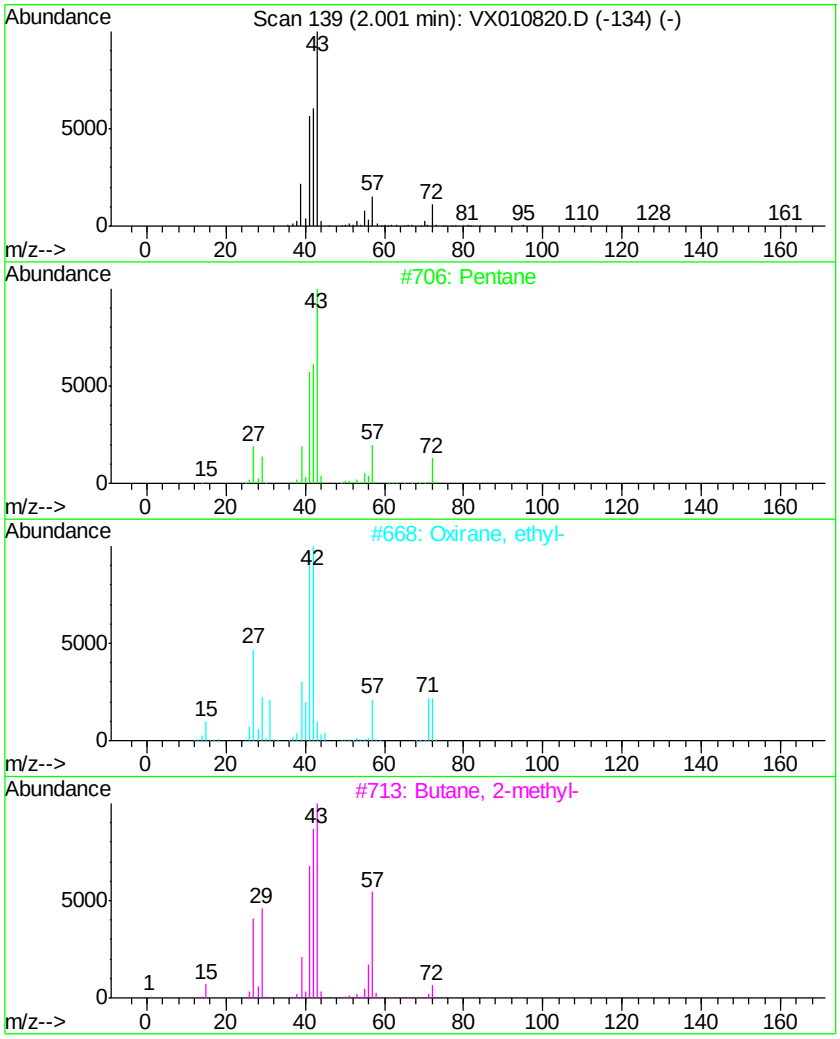
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

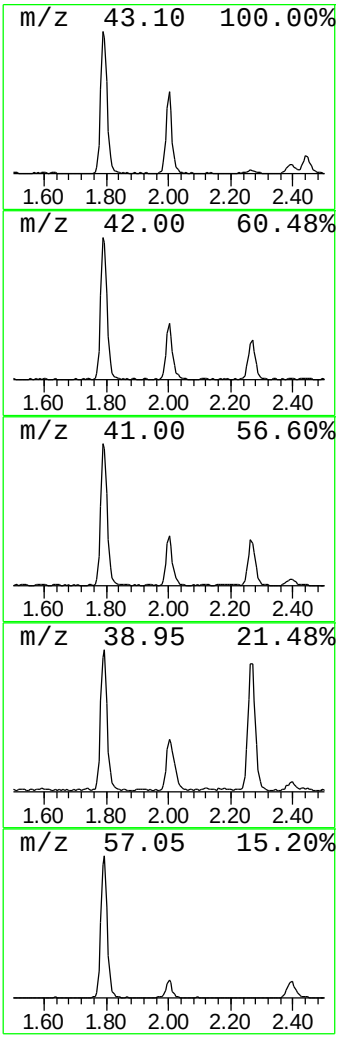
****

Peak Number 3 Pentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	8.64 ug/l	119871	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	87
2	Oxirane, ethyl-	72	C4H8O	000106-88-7	42
3	Butane, 2-methyl-	72	C5H12	000078-78-4	38
4	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	38
5	2-Hexanone, 3,4-dimethyl-	128	C8H16O	019550-10-8	25





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

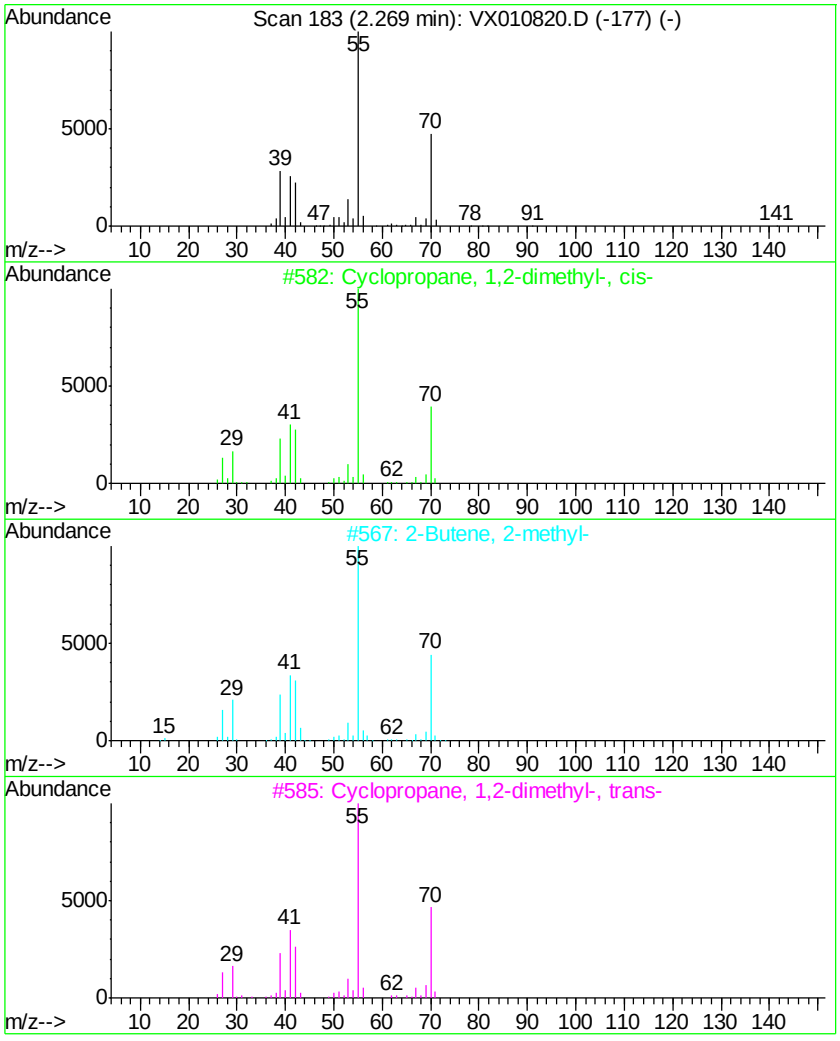
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

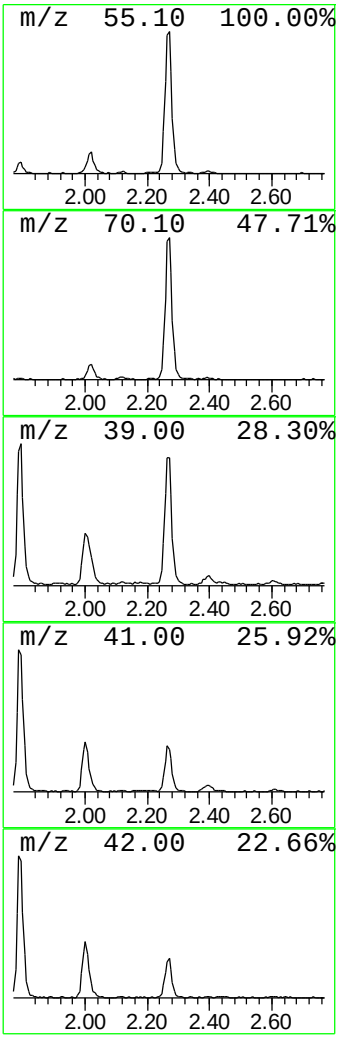
****

Peak Number 4 Cyclopropane, 1,2-dimethyl-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	14.54 ug/l	201772	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
2	2-Butene, 2-methyl-	70	C5H10	000513-35-9	91
3	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	86
4	2-Pentene, (E)-	70	C5H10	000646-04-8	86
5	1-Butene, 3-methyl-	70	C5H10	000563-45-1	78





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

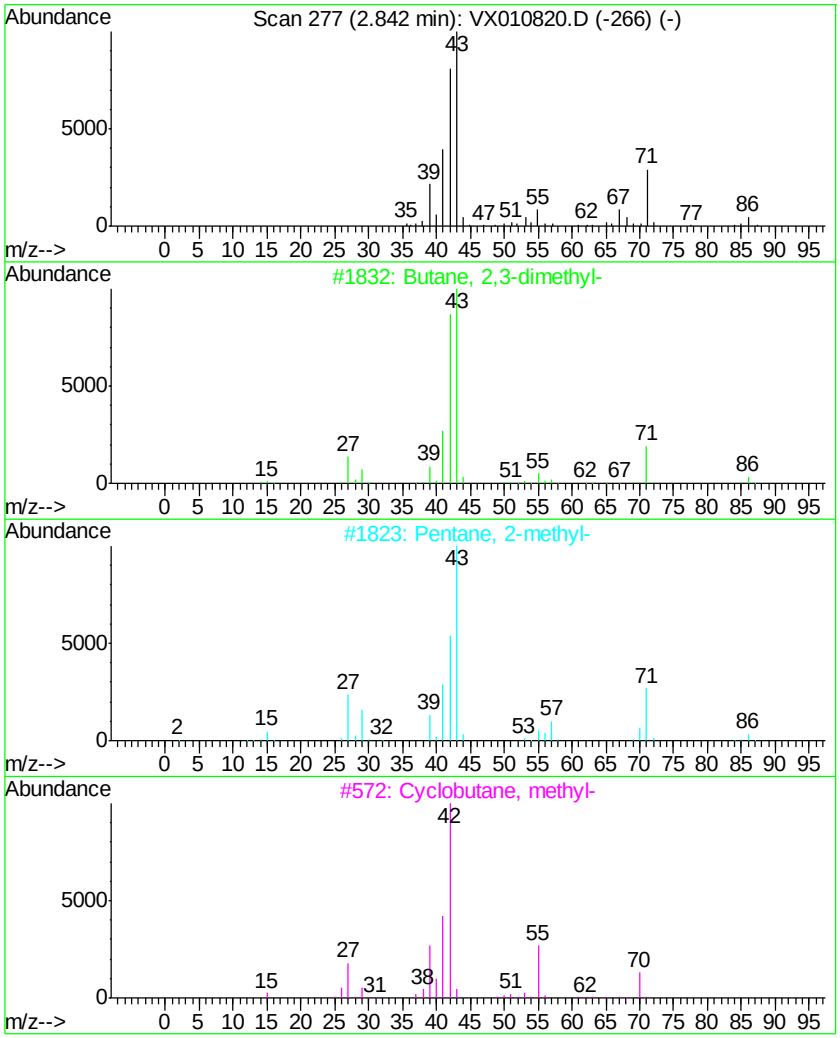
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

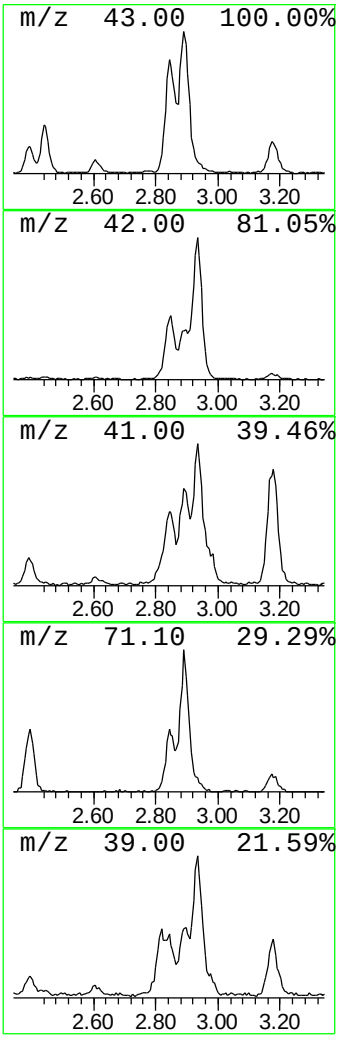
****

Peak Number 5 Butane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.84	8.77 ug/l	121702	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	64
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	47
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	46
4	Butane, 2-chloro-3-methyl-	106	C5H11Cl	000631-65-2	42
5	Tetrahydrofuran	72	C4H8O	000109-99-9	36





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

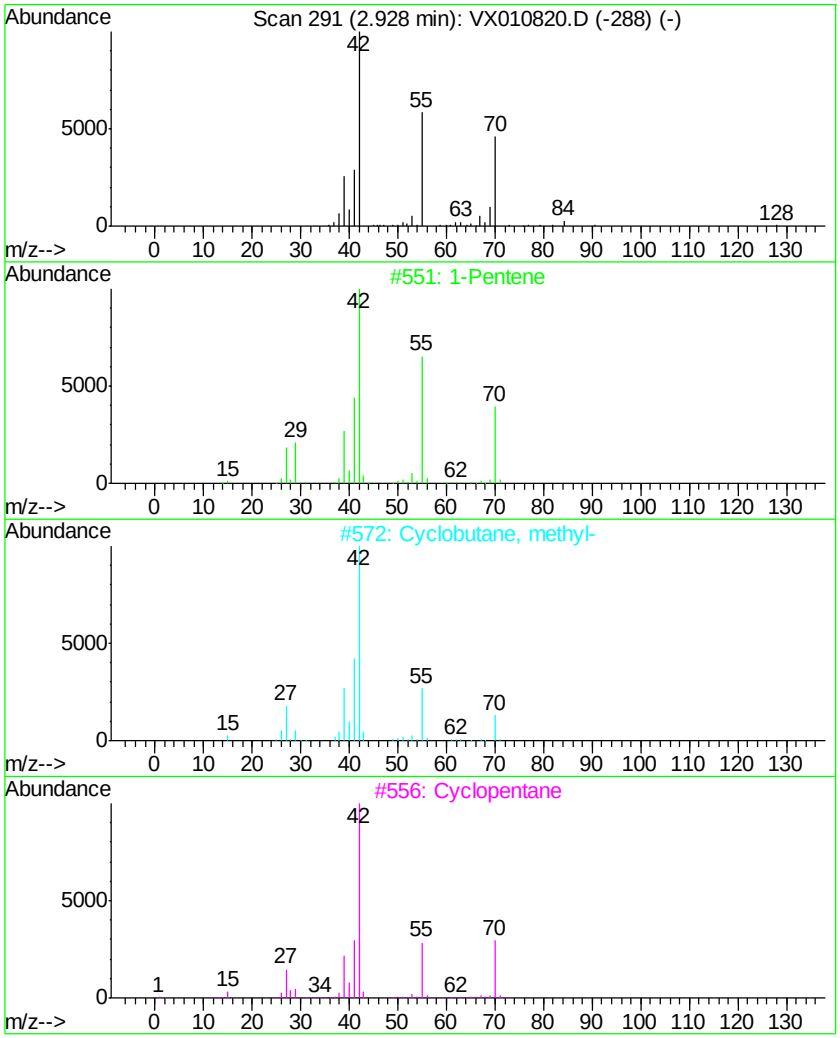
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

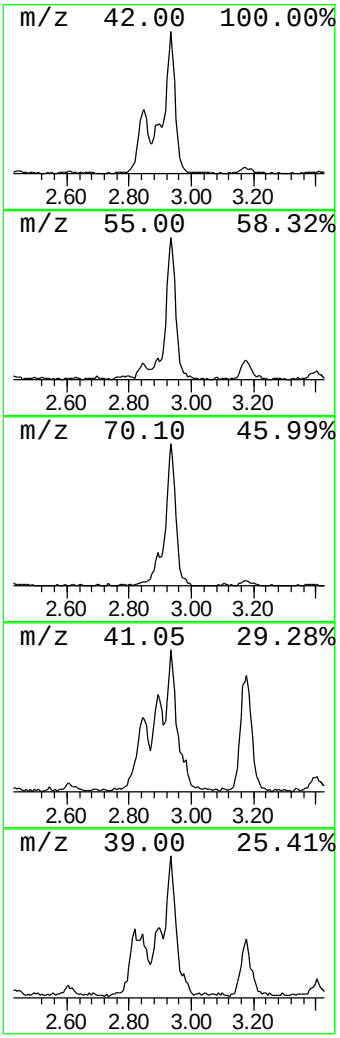
****

Peak Number 6 1-Pentene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	10.50 ug/l	145717	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene	70	C5H10	000109-67-1	80
2	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
3	Cyclopentane	70	C5H10	000287-92-3	64
4	2H-Tetrazole, 2,5-dimethyl-	98	C3H6N4	004135-93-7	39
5	2-Pentene, (E)-	70	C5H10	000646-04-8	32





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

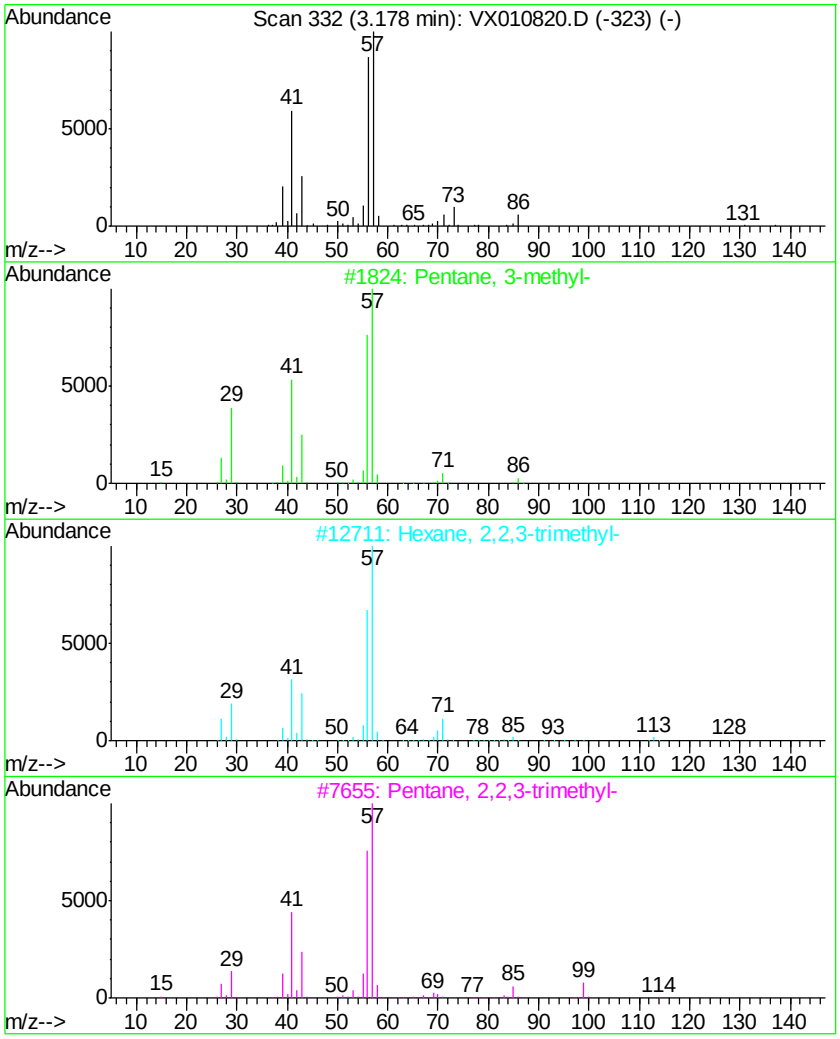
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

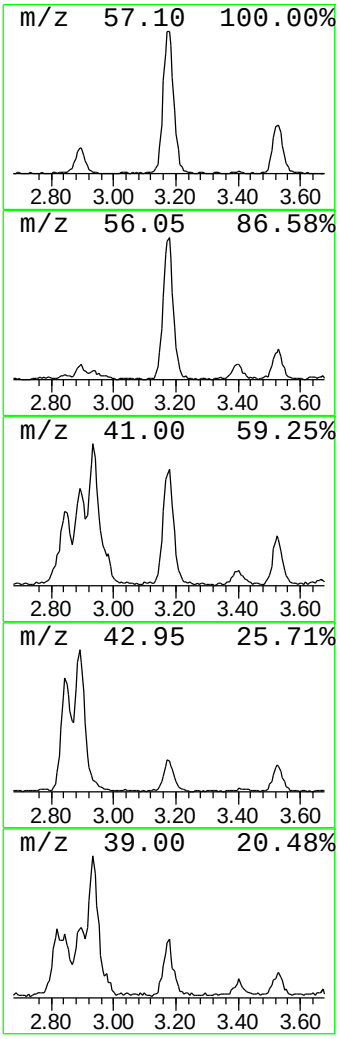
****

Peak Number 7 Pentane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	7.67 ug/l	106525	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	87
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
3	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	64
4	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	56





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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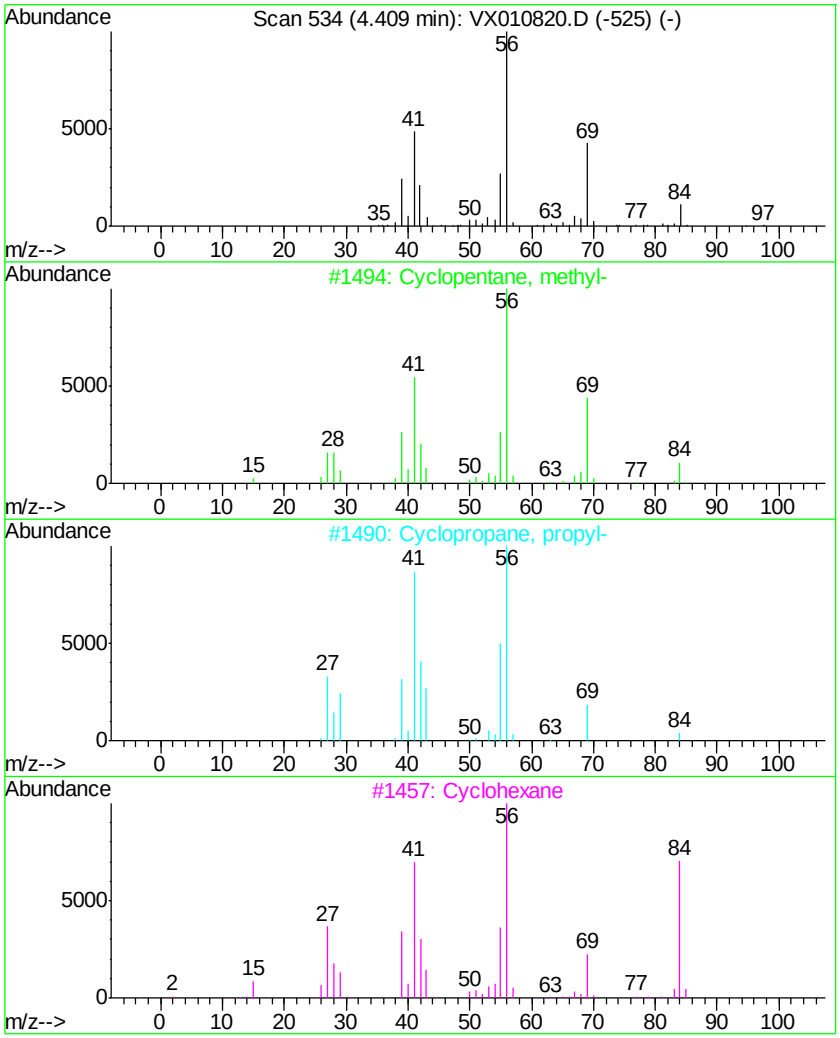
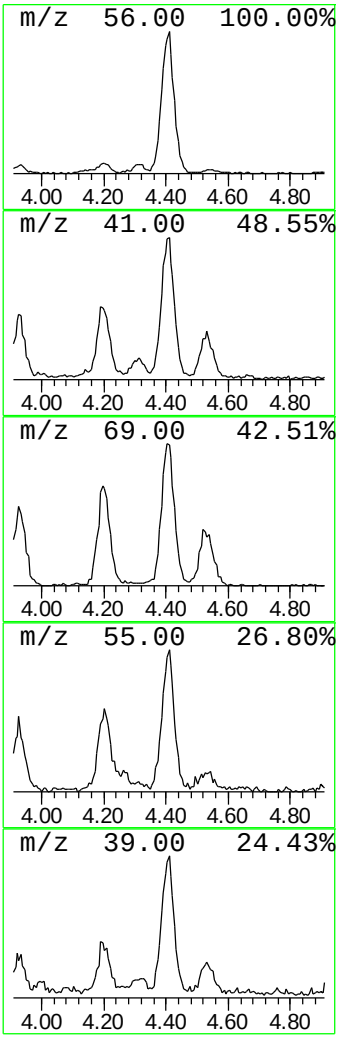
Peak Number 8 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	10.90 ug/l	151248	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclopropane, propyl-	84	C6H12	002415-72-7	72
3	Cyclohexane	84	C6H12	000110-82-7	72
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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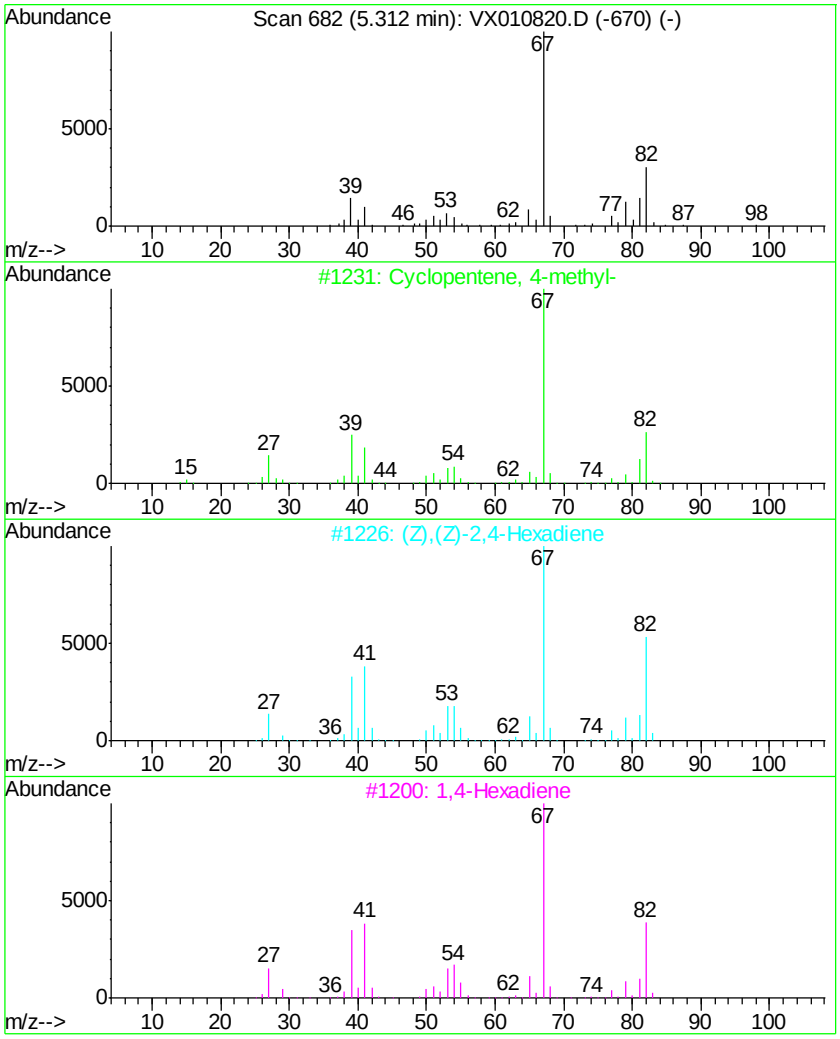
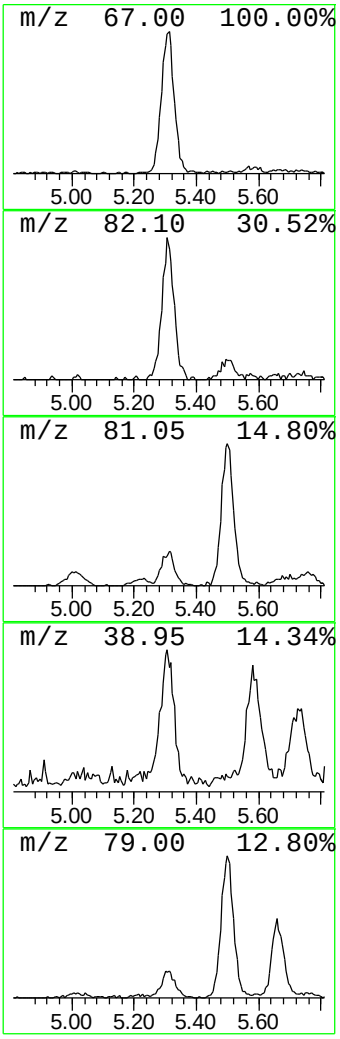
Peak Number 9 Cyclopentene, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.31	6.86 ug/l	95266	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 4-methyl-	82	C6H10	001759-81-5	86
2	(Z),(Z)-2,4-Hexadiene	82	C6H10	006108-61-8	83
3	1,4-Hexadiene	82	C6H10	000592-45-0	80
4	2,4-Hexadiene, (E,Z)-	82	C6H10	005194-50-3	80
5	1,3-Pentadiene, 3-methyl-, (Z)-	82	C6H10	002787-45-3	80





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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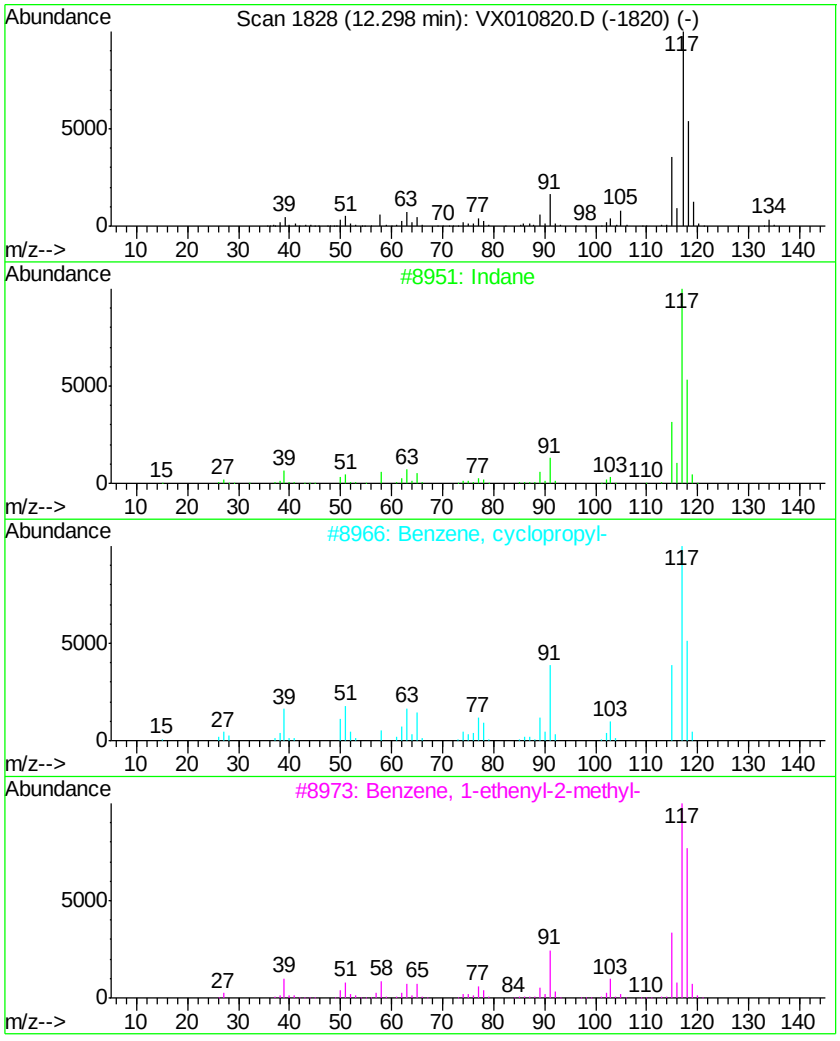
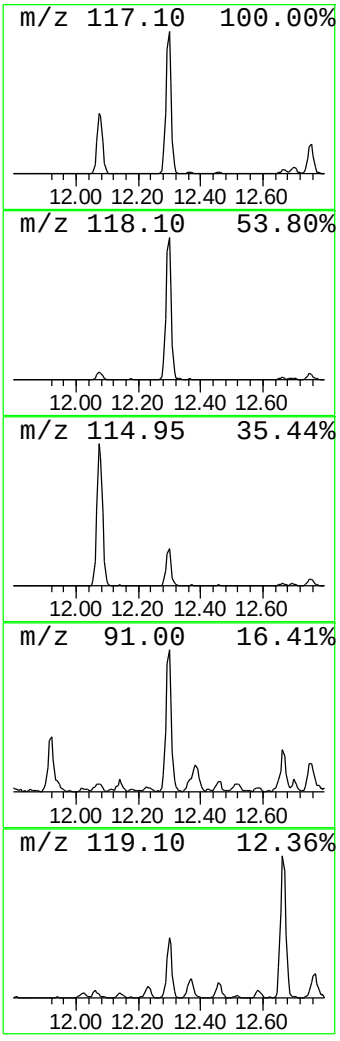
Peak Number 10 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	11.74 ug/l	215139	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	74
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071119\
Data File : VX010820.D
Acq On : 11 Jul 2019 20:17
Operator : JC/SP
Sample : K3749-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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
Sulfur dioxide	1.28	37.8	ug/l	524712	1	5.67	694001	50.0
Butane, 2-methyl-	1.79	19.6	ug/l	271583	1	5.67	694001	50.0
Pentane	2.00	8.6	ug/l	119871	1	5.67	694001	50.0
Cyclopropane, 1,2...	2.27	14.5	ug/l	201772	1	5.67	694001	50.0
Butane, 2,3-dimet...	2.84	8.8	ug/l	121702	1	5.67	694001	50.0
1-Pentene	2.93	10.5	ug/l	145717	1	5.67	694001	50.0
Pentane, 3-methyl-	3.18	7.7	ug/l	106525	1	5.67	694001	50.0
Cyclopentane, met...	4.41	10.9	ug/l	151248	1	5.67	694001	50.0
Cyclopentene, 4-m...	5.31	6.9	ug/l	95266	1	5.67	694001	50.0
Indane	12.30	11.7	ug/l	215139	4	12.08	916026	50.0