LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

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_U\METHOD\82U092218W.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	1.402	63	71	79	rBV	44579	47408	5.32%	0.678%
2	1.759	174	182	196	rVB	52048	70230	7.89%	1.005%
3	1.948	234	241	253	rVB2	20402	29750	3.34%	0.426%
4	2.051	266	273	283	rVB	17025	22630	2.54%	0.324%
5	2.138	294	300	308	rBV2	5989	9284	1.04%	0.133%
6	2.186	308	315	328	rVB2	13171	20154	2.26%	0.288%
7	2.308	341	353	369	rVB3	13165	32282	3.63%	0.462%
8	2.443	383	395	418	rBV2	19986	41556	4.67%	0.594%
9	2.707	464	477	497	rBV	34789	80774	9.07%	1.155%
10	3.000	555	568	582	rVB3	9416	20158	2.26%	0.288%
11	4.000	861	879	896	rBV3	14852	39837	4.47%	0.570%
12	4.099	896	910	925	rVB4	5096	13514	1.52%	0.193%
13	4.202	925	942	960	rBV3	5079	15751	1.77%	0.225%
14	4.887	1138	1155	1172	rBV	119334	268669	30.18%	3.843%
15	4.983	1172	1185	1202	rVV	173745	381710	42.87%	5.460%
16	5.083	1202	1216	1235	rVB2	50476	123498	13.87%	1.767%
17	5.311	1272	1287	1305	rBV	144079	301670	33.88%	4.315%
18	5.488	1327	1342	1343	rBV7	5932	9649	1.08%	0.138%
19	5.540	1345	1358	1375	rVB2	54046	133361	14.98%	1.908%
20	5.887	1451	1466	1484	rBV	258974	510173	57.30%	7.298%
21	6.398	1611	1625	1640	rBV5	7801	20042	2.25%	0.287%
22	6.479	1640	1650	1657	rVV3	5903	10699	1.20%	0.153%
23	6.536	1657	1668	1679	rVV3	9098	19224	2.16%	0.275%
24	6.742	1714	1732	1744	rVB5	9618	19612	2.20%	0.281%
25	6.926	1772	1789	1804	rBV	38573	85212	9.57%	1.219%
26	7.051	1811	1828	1838	rBV2	39167	78871	8.86%	1.128%
27	7.106	1839	1845	1855	rVB	8777	14319	1.61%	0.205%
28	7.565	1964	1988	2002	rBV	492227	890344	100.00%	12.736%
29	7.639	2003	2011	2023	rVB	38660	71281	8.01%	1.020%
30	7.919	2087	2098	2111	rBV4	18168	32454	3.65%	0.464%
31	8.048	2127	2138	2147	rBV3	8869	15275	1.72%	0.219%
32	8.491	2255	2276	2284	rBV9	7925	23730	2.67%	0.339%
33	8.607	2302	2312	2323	rBV3	15224	29402	3.30%	0.421%
34	8.755	2345	2358	2369	rBV8	4557	9734	1.09%	0.139%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

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_U\METHOD\82U092218W.M
Title : SW846 8260

35	8.848	2372	2387	2401	rBV9	7261	17433	1.96%	0.249%
36	9.089	2450	2462	2484	rVV	368512	620201	69.66%	8.872%
37	9.253	2502	2513	2528	rBV	42453	75020	8.43%	1.073%
38	9.376	2536	2551	2568	rBV	99311	192352	21.60%	2.752%
39	9.450	2570	2574	2596	rVB	5605	13469	1.51%	0.193%
40	9.720	2642	2658	2668	rBV4	14400	31609	3.55%	0.452%
41	9.781	2668	2677	2703	rVB	27946	57617	6.47%	0.824%
42	9.954	2712	2731	2740	rBV4	13100	29968	3.37%	0.429%
43	10.048	2748	2760	2779	rVB10	10502	24185	2.72%	0.346%
44	10.308	2829	2841	2856	rBV	337538	541347	60.80%	7.744%
45	10.376	2857	2862	2871	rVB7	6280	9665	1.09%	0.138%
46	10.491	2891	2898	2908	rVB7	6985	11240	1.26%	0.161%
47	10.588	2919	2928	2939	rBV	15926	26463	2.97%	0.379%
48	10.684	2949	2958	2963	rBV	18159	30362	3.41%	0.434%
49	10.774	2978	2986	2993	rVB2	7693	11002	1.24%	0.157%
50	10.974	3039	3048	3061	rBV2	12252	21139	2.37%	0.302%
51	11.150	3093	3103	3117	rVB	37506	59222	6.65%	0.847%
52	11.292	3132	3147	3151	rBV3	6532	11252	1.26%	0.161%
53	11.324	3152	3157	3166	rVV	8602	14272	1.60%	0.204%
54	11.485	3195	3207	3222	rBV	409443	640286	71.91%	9.159%
55	11.572	3226	3234	3243	rVB2	9864	14838	1.67%	0.212%
56	11.781	3282	3299	3316	rBV3	36597	78166	8.78%	1.118%
57	11.864	3316	3325	3344	rVB3	27005	58789	6.60%	0.841%
58	11.983	3351	3362	3374	rBV4	6606	10761	1.21%	0.154%
59	12.060	3374	3386	3398	rVB3	27399	45766	5.14%	0.655%
60	12.253	3433	3446	3455	rBV	50775	85215	9.57%	1.219%
61	12.356	3468	3478	3501	rVB3	22609	55299	6.21%	0.791%
62	12.488	3507	3519	3529	rVB8	4527	9874	1.11%	0.141%
63	12.568	3539	3544	3552	rVB2	8369	11664	1.31%	0.167%
64	12.652	3552	3570	3580	rVV	34250	59342	6.67%	0.849%
65	12.703	3581	3586	3600	rVB6	6042	10415	1.17%	0.149%
66	12.787	3603	3612	3624	rBV2	16339	26421	2.97%	0.378%
67	12.912	3642	3651	3660	rBV5	7190	15001	1.68%	0.215%
68	12.964	3660	3667	3683	rVB2	22273	34373	3.86%	0.492%
69	13.125	3699	3717	3726	rBV3	46480	86107	9.67%	1.232%
70	13.182	3732	3735	3744	rVB4	7524	9859	1.11%	0.141%
71	13.240	3744	3753	3761	rBV	15985	23444	2.63%	0.335%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

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_U\METHOD\82U092218W.M
Title : SW846 8260

72	13.408	3797	3805	3812	rBV3	7525	11326	1.27%	0.162%
73	13.456	3812	3820	3829	rVV2	15793	25094	2.82%	0.359%
74	13.510	3829	3837	3845	rBV3	27334	39613	4.45%	0.567%
75	13.610	3863	3868	3873	rBV	7728	9616	1.08%	0.138%
76	13.639	3874	3877	3890	rVB	10976	13497	1.52%	0.193%
77	13.742	3891	3909	3920	rBV	60495	101848	11.44%	1.457%
78	13.957	3965	3976	3987	rBV7	9428	19908	2.24%	0.285%
79	14.015	3987	3994	4001	rVB4	7819	11172	1.25%	0.160%
80	14.153	4026	4037	4049	rBV3	14093	21303	2.39%	0.305%
81	14.337	4084	4094	4106	rBV4	14021	28604	3.21%	0.409%
82	14.478	4131	4138	4148	rBV2	8669	13222	1.49%	0.189%
83	14.543	4149	4158	4170	rVB4	10166	18287	2.05%	0.262%
84	14.880	4252	4263	4277	rBV	36308	64284	7.22%	0.920%
85	15.063	4310	4320	4332	rBV	27484	47120	5.29%	0.674%

Sum of corrected areas: 6990589

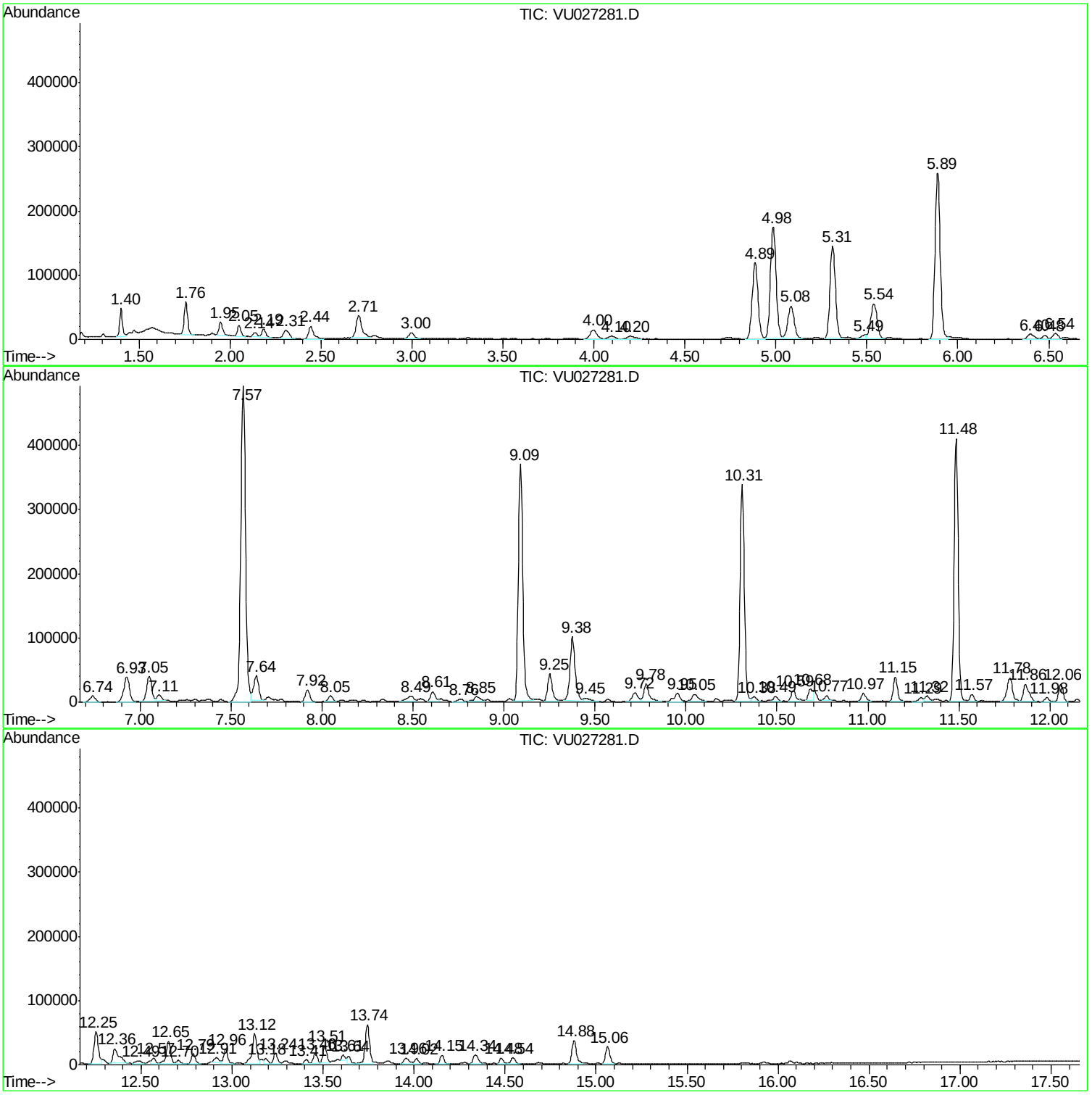
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.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 U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

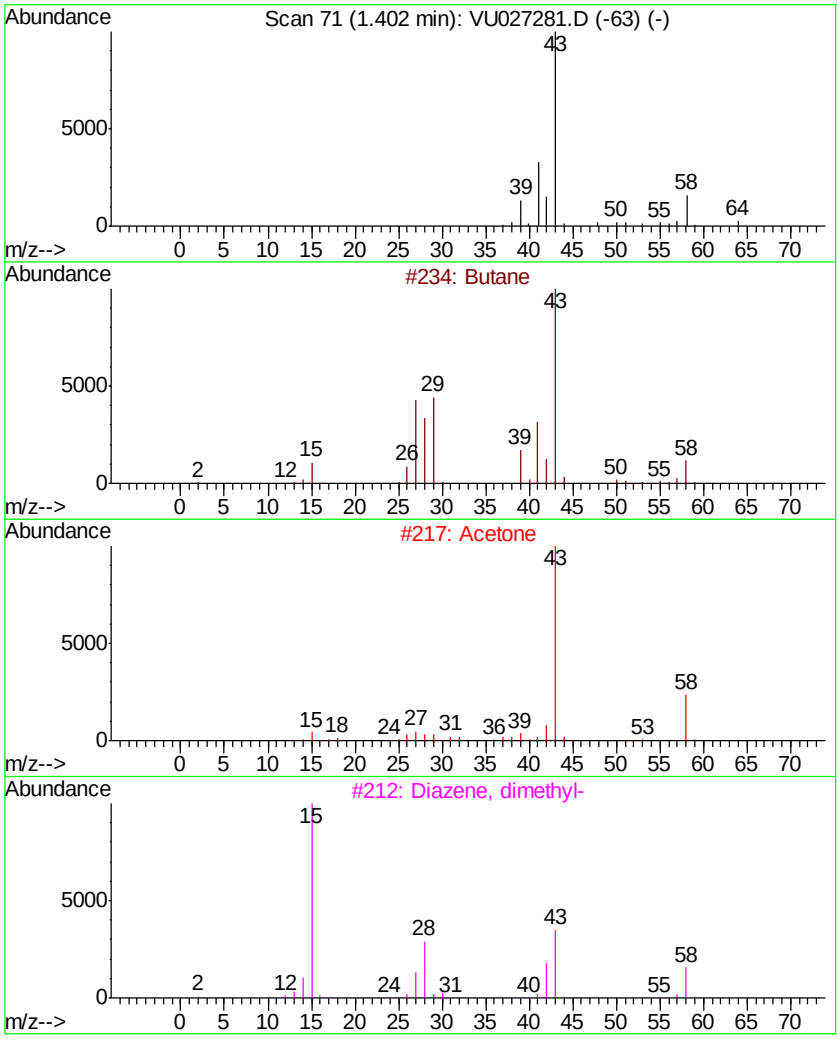
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

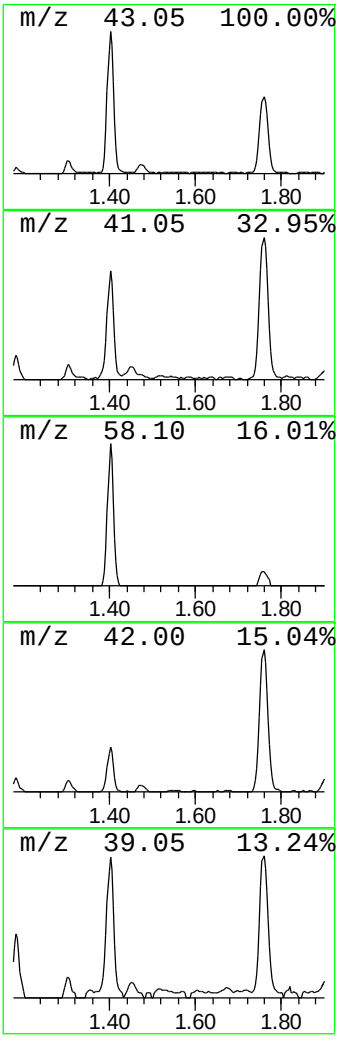
****

Peak Number 1 Butane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.40	6.21 ug/l	47408	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane	58	C4H10	000106-97-8	64
2	Acetone	58	C3H6O	000067-64-1	7
3	Diazene, dimethyl-	58	C2H6N2	000503-28-6	5
4	Isopropylsulfonyl chloride	142	C3H7ClO2S	010147-37-2	4
5	Allyl acetate	100	C5H8O2	000591-87-7	4





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

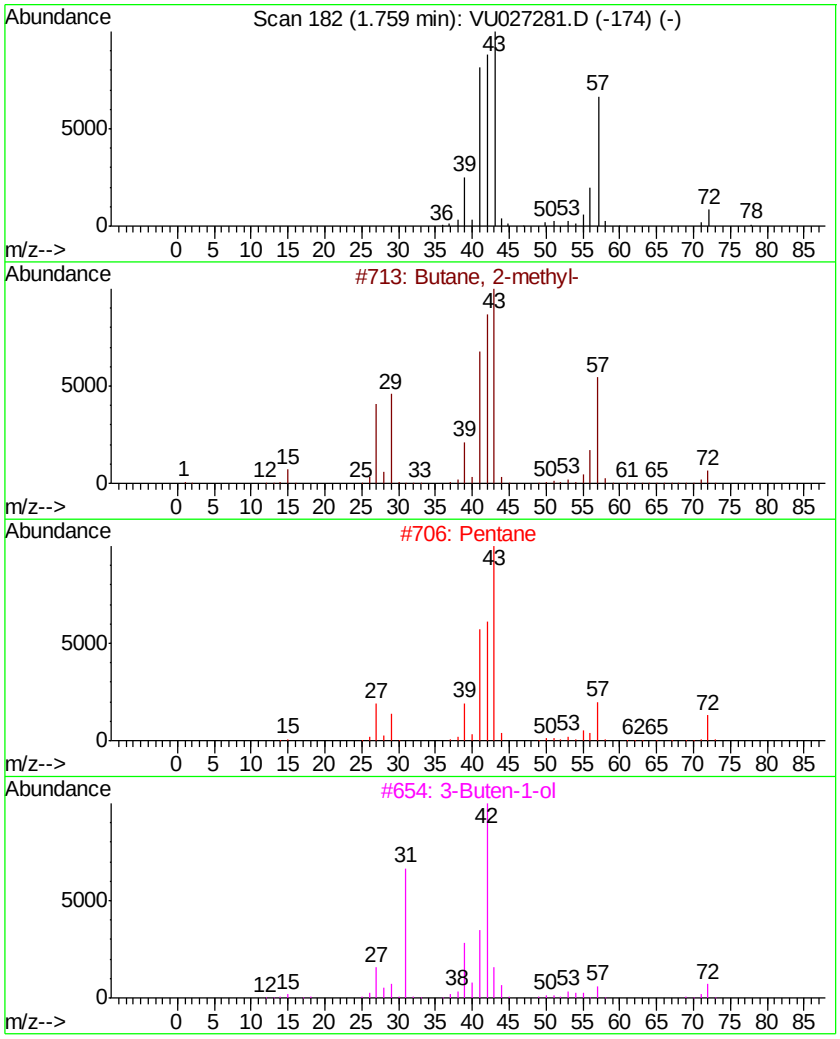
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

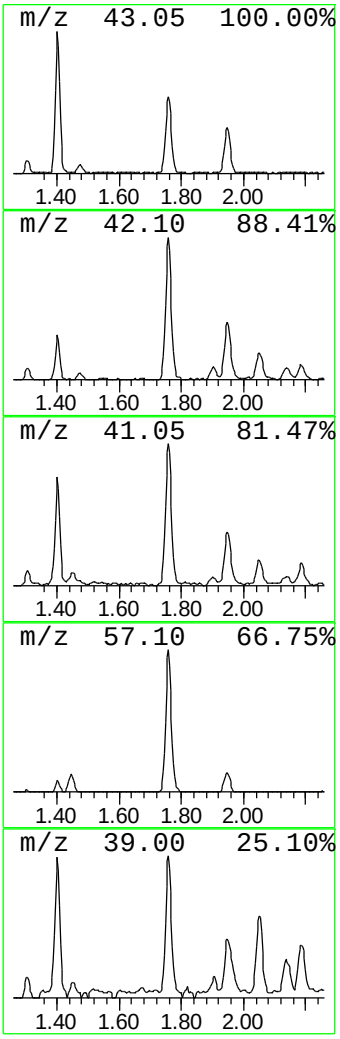
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	9.20 ug/l	70230	Pentafluorobenzene	4.99

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	50
3	3-Buten-1-ol	72	C4H8O	000627-27-0	47
4	Isobutylene epoxide	72	C4H8O	000558-30-5	42
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

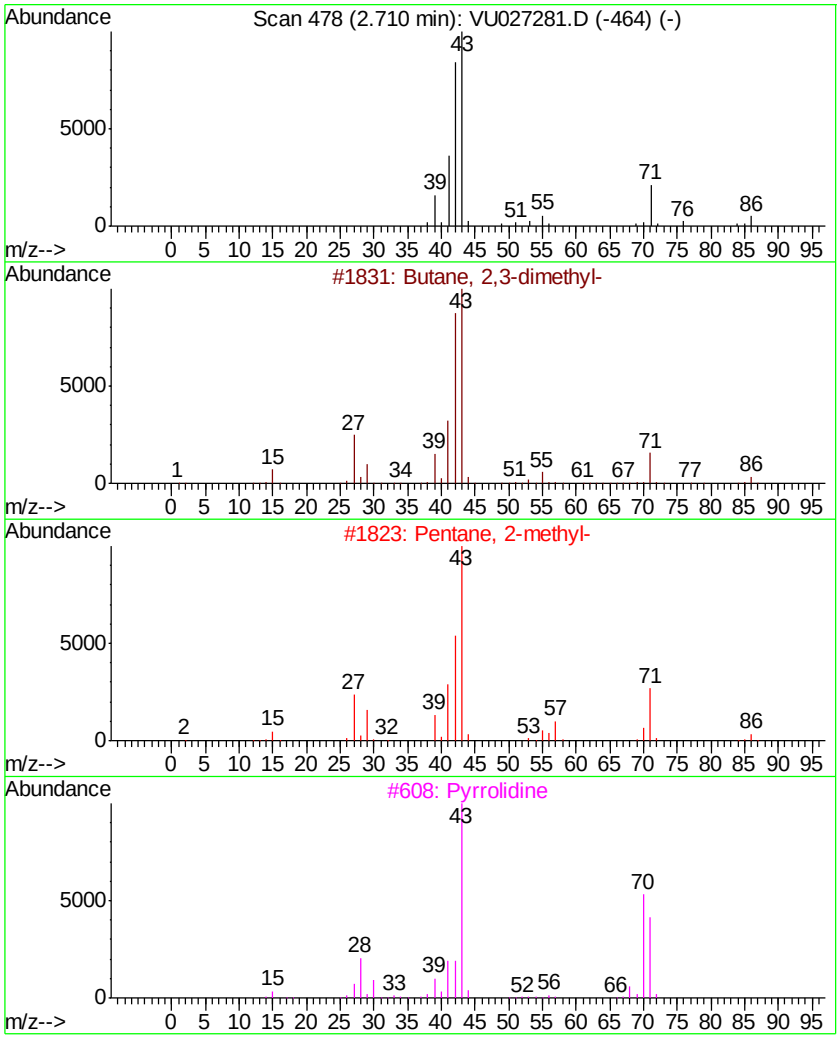
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

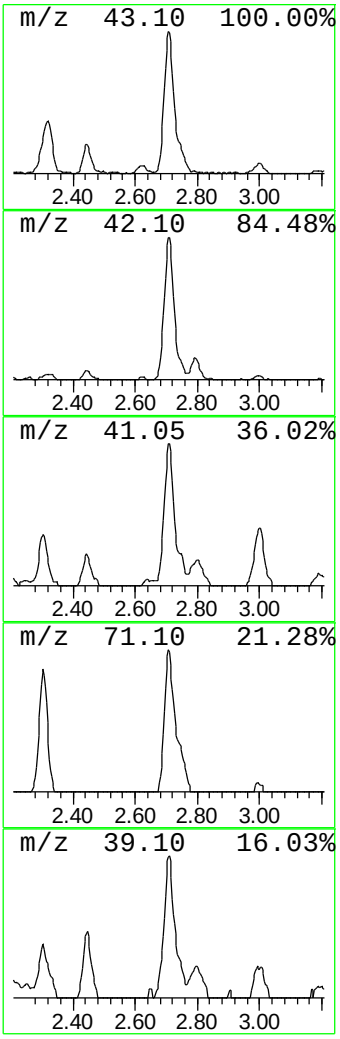
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.71	10.58 ug/l	80774	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	90
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	33
3	Pyrrolidine	71	C4H9N	000123-75-1	10
4	Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	10
5	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

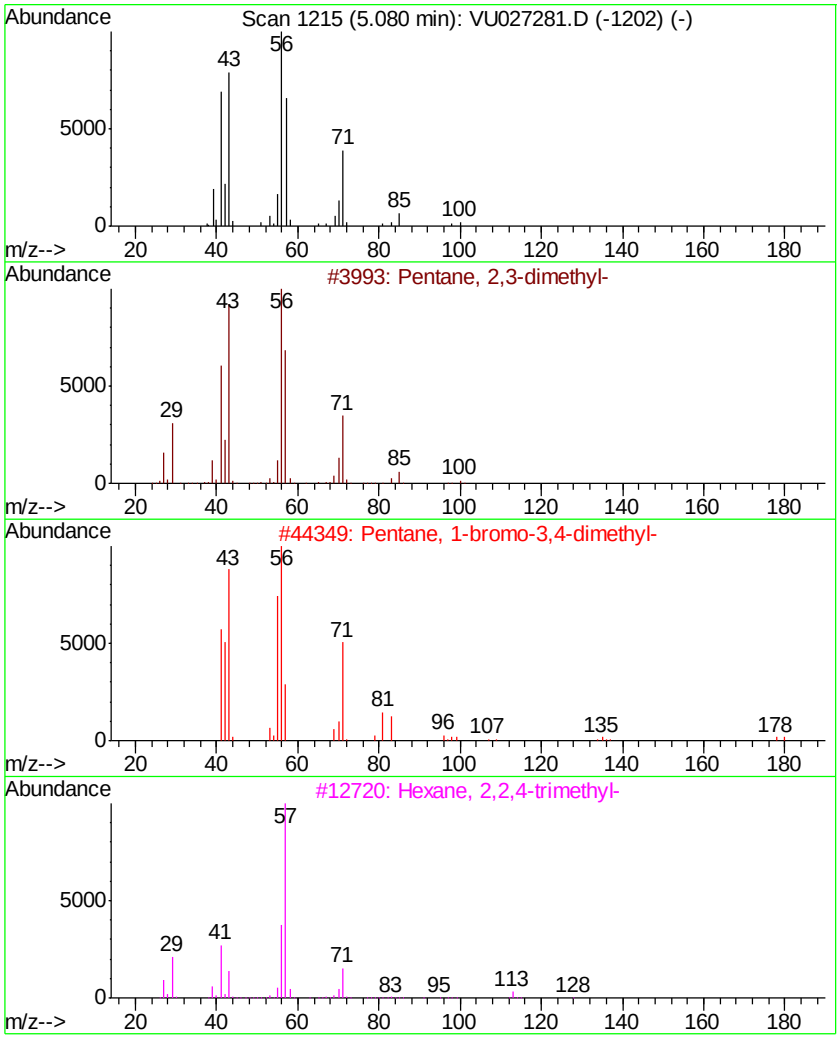
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

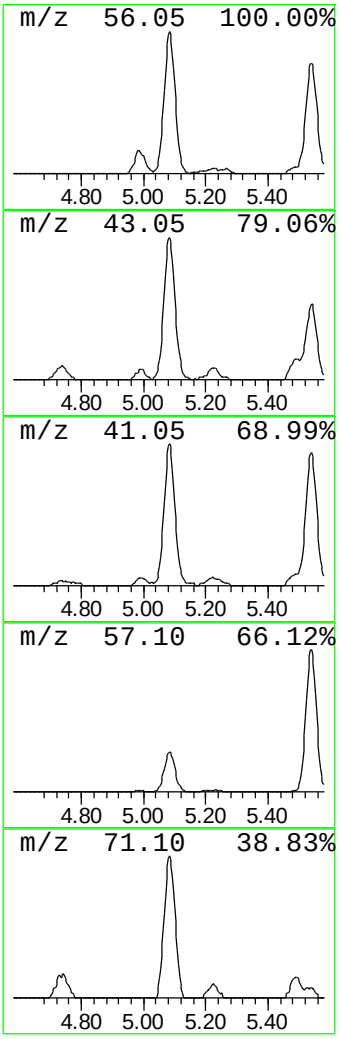
****

Peak Number 4 Pentane, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	16.18 ug/l	123498	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	94
2	Pentane, 1-bromo-3,4-dimethyl-	178	C7H15Br	006570-92-9	50
3	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45
4	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	43
5	Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

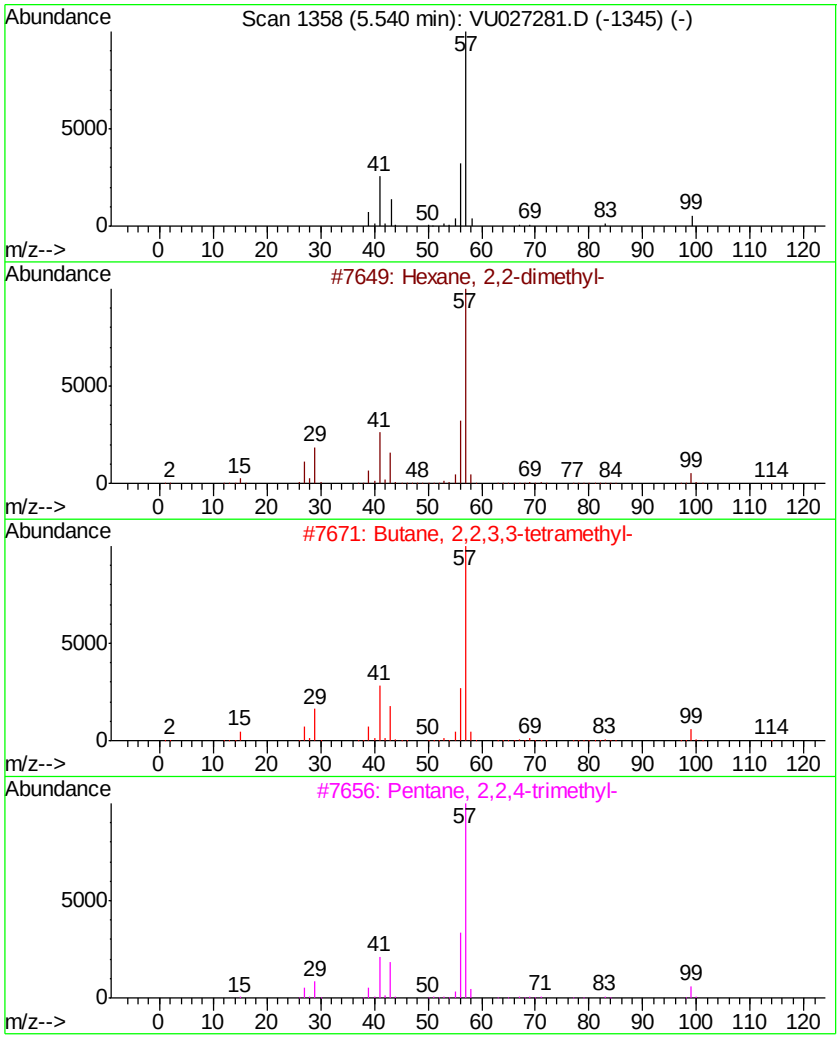
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

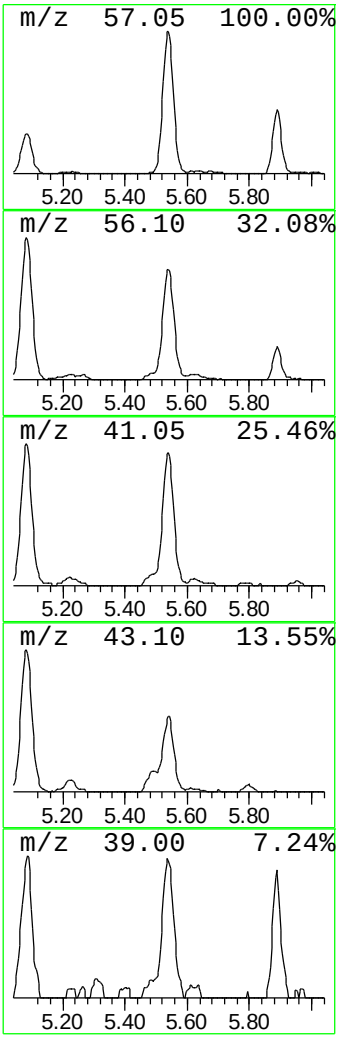
****

Peak Number 5 Hexane, 2,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.54	13.07 ug/l	133361	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	78
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

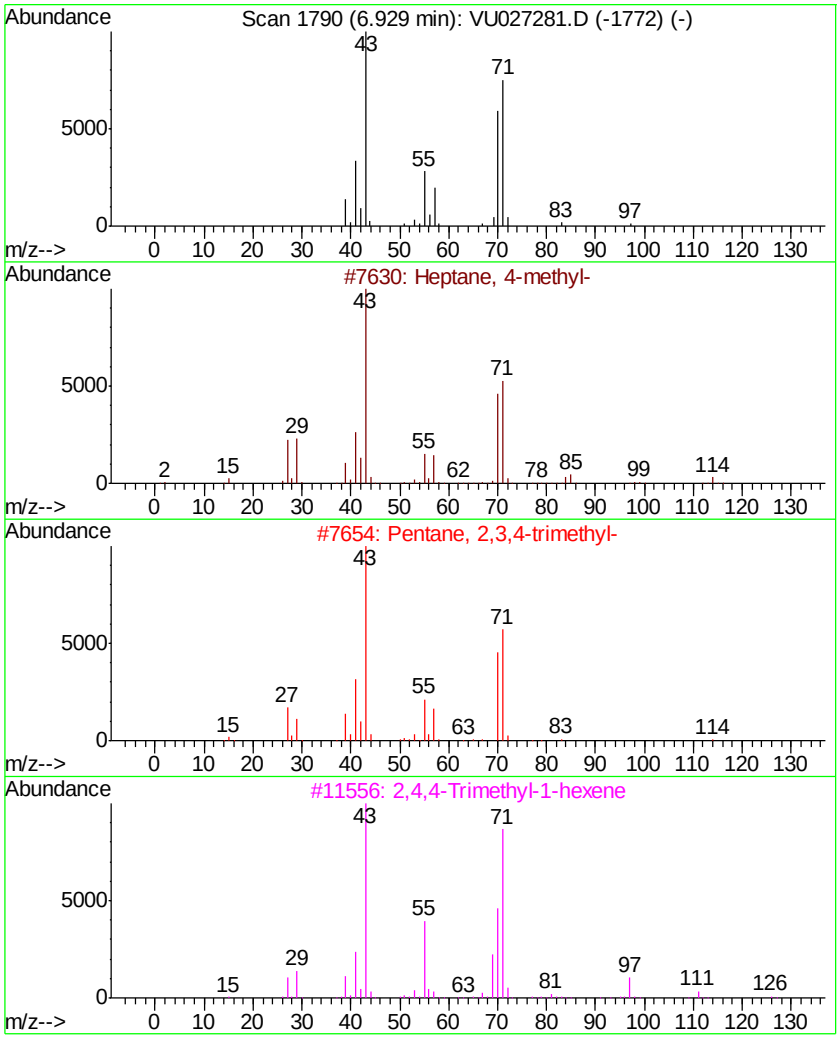
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

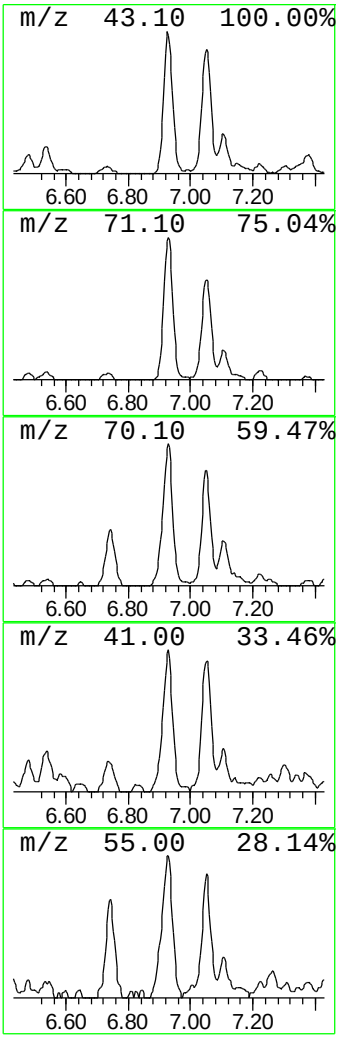
****

Peak Number 6 Heptane, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	8.35 ug/l	85212	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
2	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
3	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	64
4	Ethanedioic acid, bis(3-methylbu...	230	C12H22O4	002051-00-5	59
5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	56





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

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

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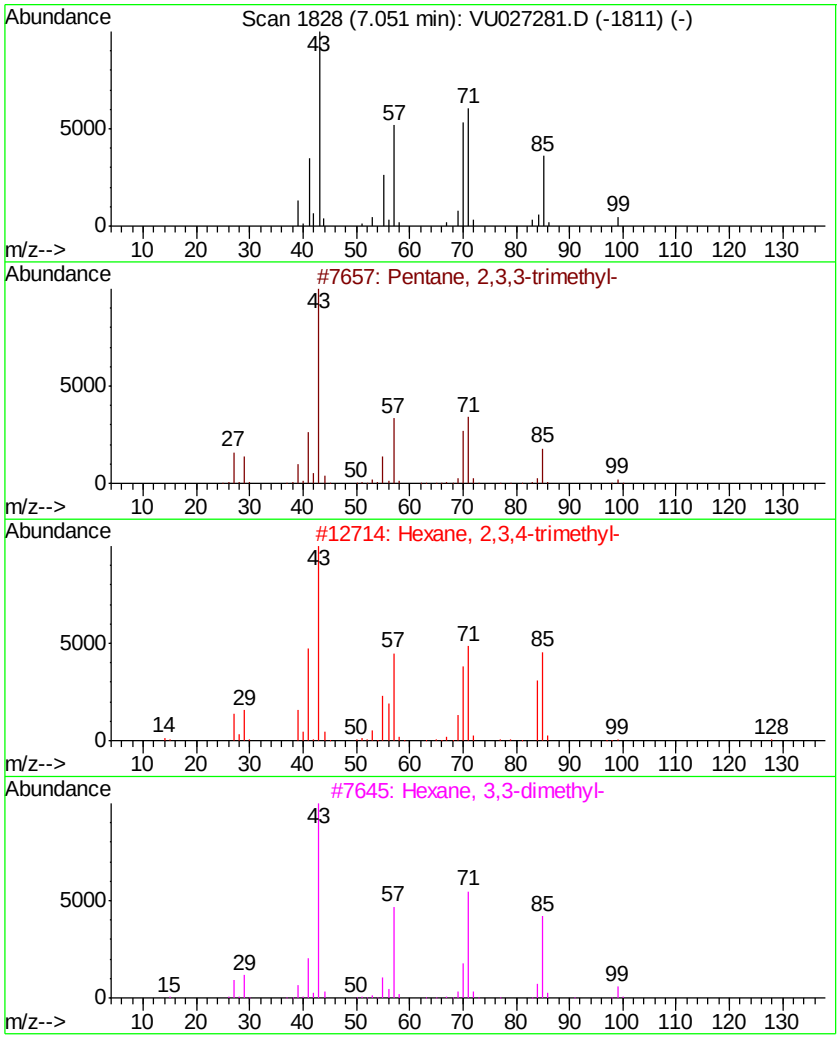
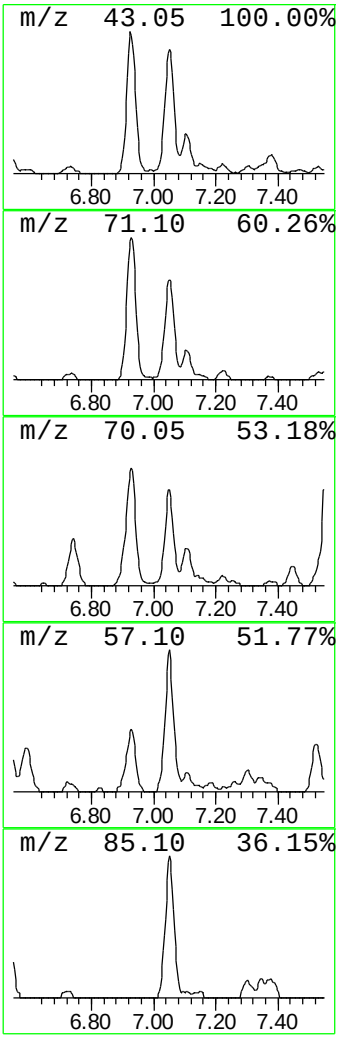
Peak Number 7 Pentane, 2,3,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	7.73 ug/l	78871	1,4-Difluorobenzene	5.89

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	64
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4	Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	59
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	53





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

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

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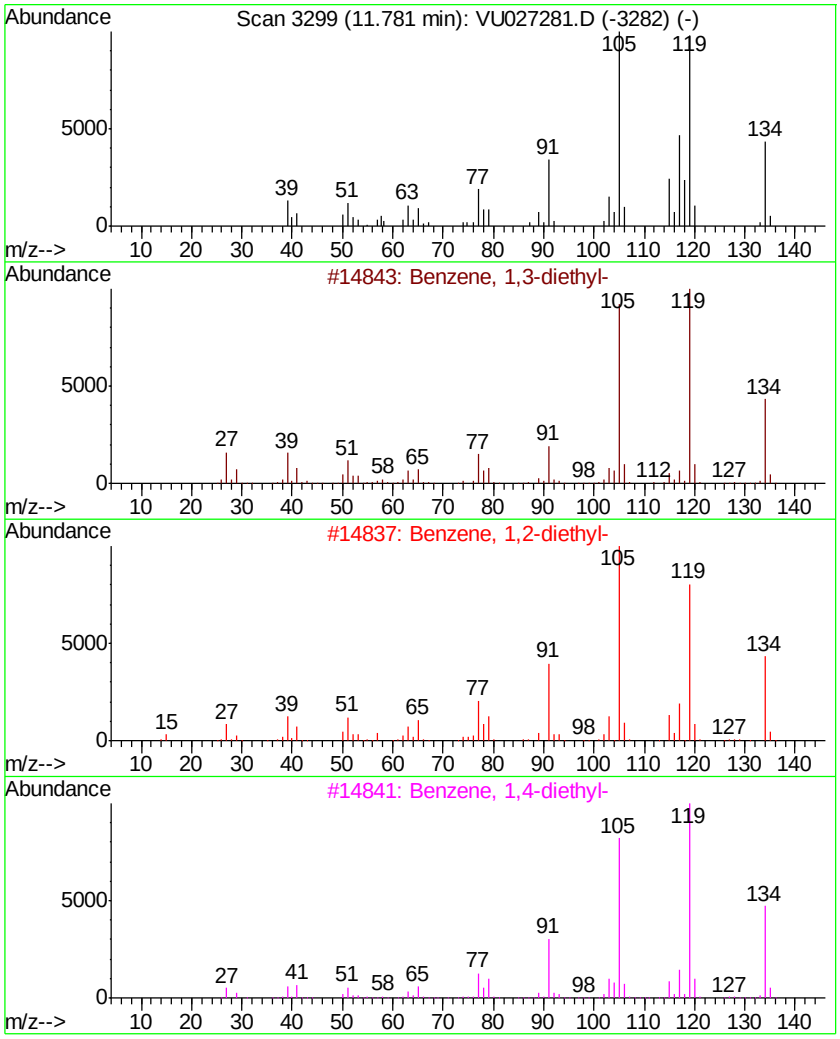
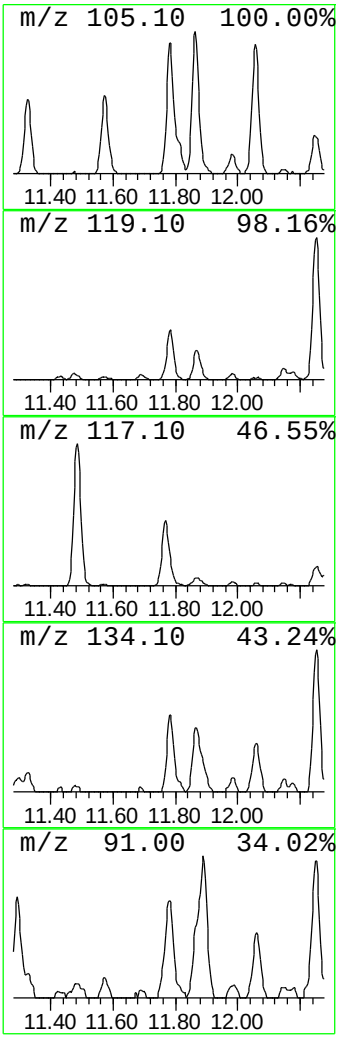
Peak Number 8 Benzene, 1,3-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	6.10 ug/l	78166	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	92
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	89
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	58
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	58





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

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

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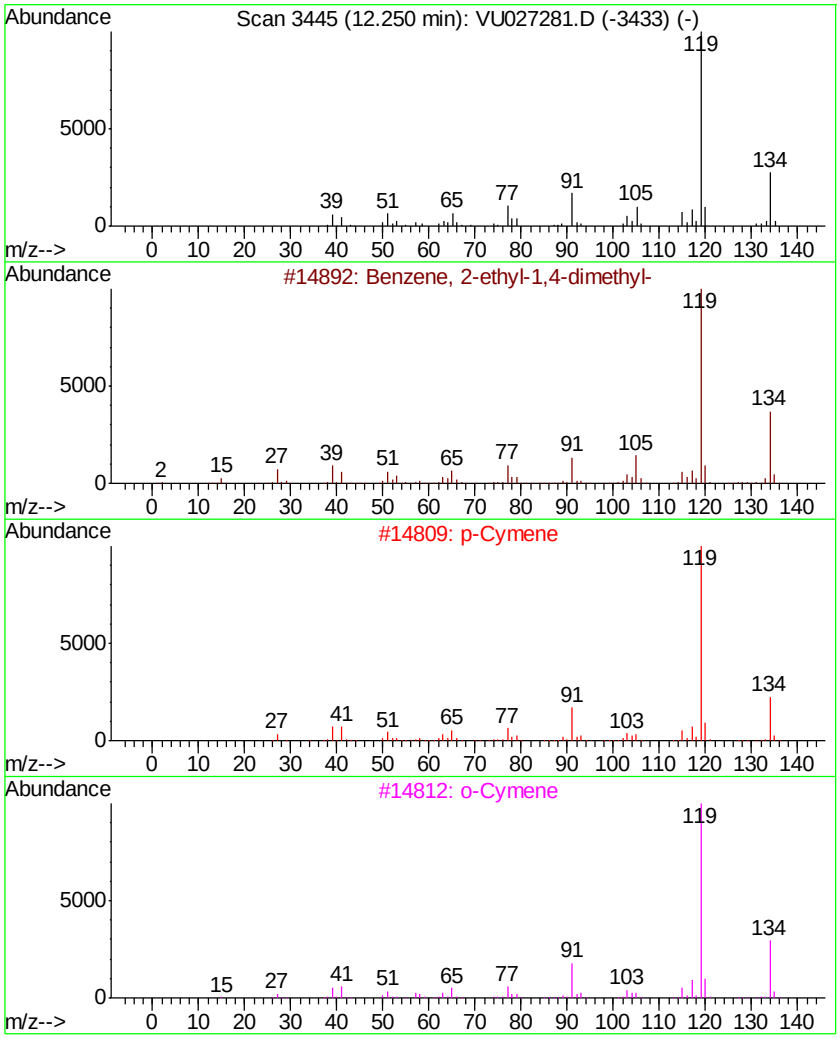
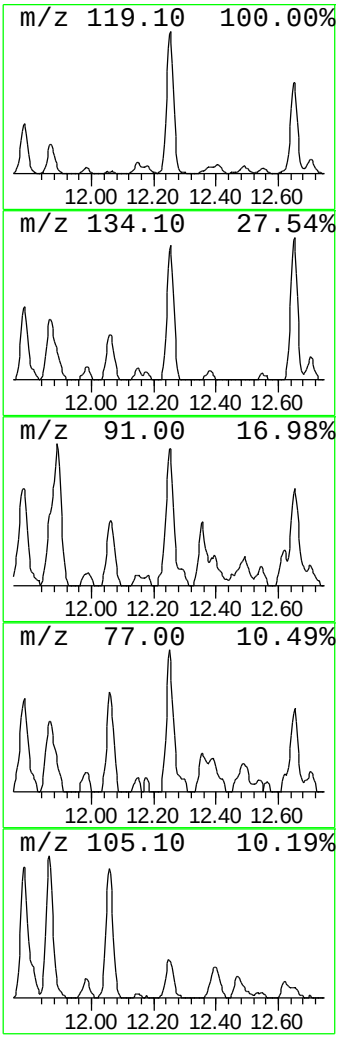
Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	6.65 ug/l	85215	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	p-Cymene	134	C10H14	000099-87-6	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





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Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

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

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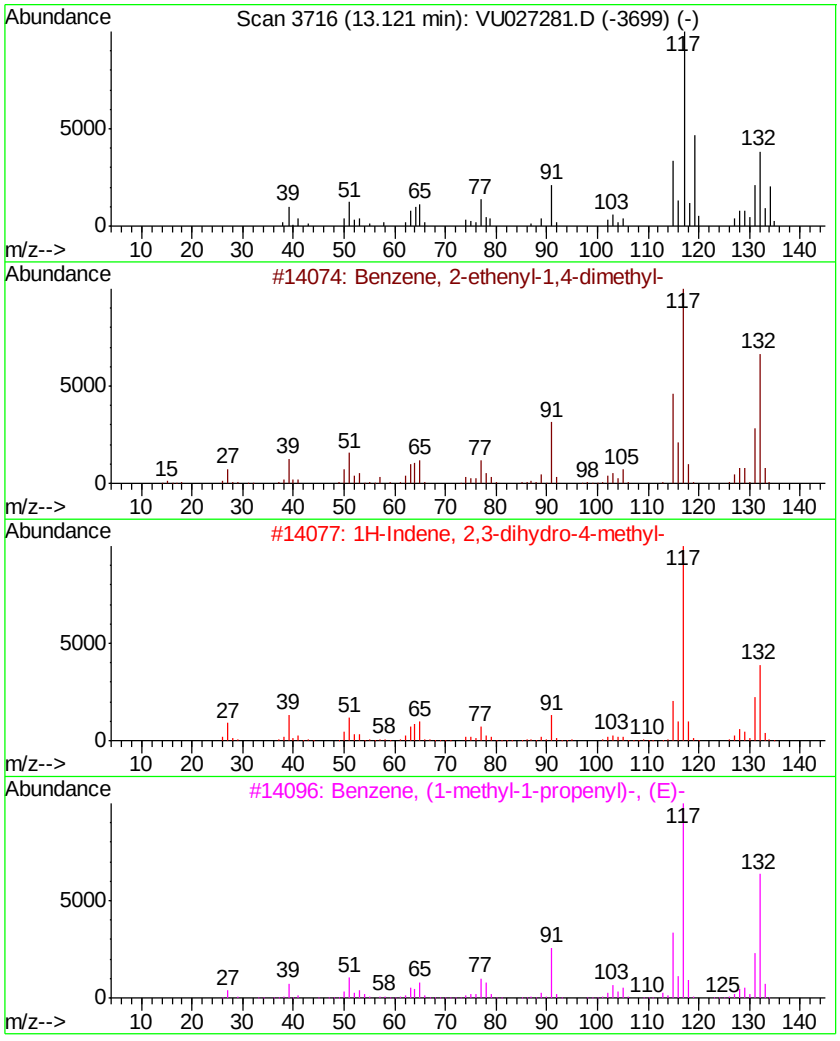
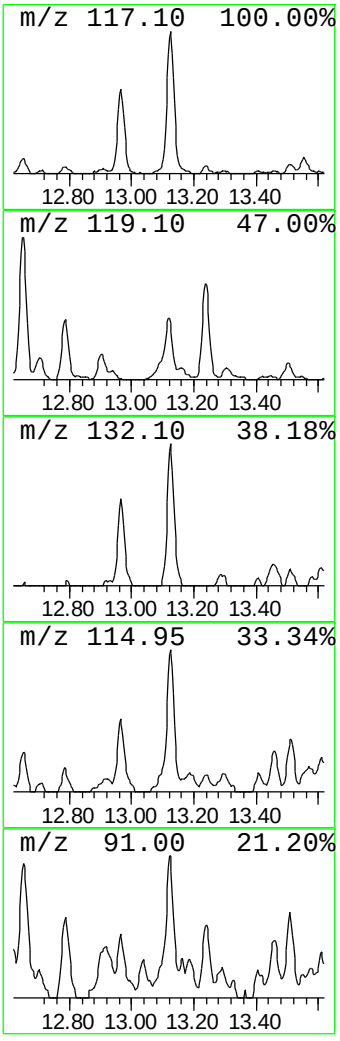
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	6.72 ug/l	86107	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027281.D
Acq On : 28 Sep 2018 18:58
Operator : MD/SY
Sample : J5084-24
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ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : B-11(TW-3)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.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	1.40	6.2	ug/l	47408	1	4.99	381710	50.0
Butane, 2-methyl-	1.76	9.2	ug/l	70230	1	4.99	381710	50.0
Butane, 2,3-dimet...	2.71	10.6	ug/l	80774	1	4.99	381710	50.0
Pentane, 2,3-dime...	5.08	16.2	ug/l	123498	1	4.99	381710	50.0
Hexane, 2,2-dimet...	5.54	13.1	ug/l	133361	2	5.89	510173	50.0
Heptane, 4-methyl-	6.93	8.3	ug/l	85212	2	5.89	510173	50.0
Pentane, 2,3,3-tr...	7.05	7.7	ug/l	78871	2	5.89	510173	50.0
Benzene, 1,3-diet...	11.78	6.1	ug/l	78166	4	11.48	640286	50.0
Benzene, 2-ethyl-...	12.25	6.7	ug/l	85215	4	11.48	640286	50.0
Benzene, 2-etheny...	13.12	6.7	ug/l	86107	4	11.48	640286	50.0