LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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	61	71	79	rBV	31085	33377	2.17%	0.181%
2	1.759	171	182	203	rBV	465519	608538	39.62%	3.296%
3	1.948	232	241	256	rVB	100799	139335	9.07%	0.755%
4	2.305	338	352	375	rVB	122311	233205	15.18%	1.263%
5	2.444	385	395	415	rBV2	17822	33122	2.16%	0.179%
6	2.707	463	477	483	rBV	143581	285920	18.61%	1.549%
7	2.794	495	504	526	rVB	402825	775868	50.51%	4.202%
8	2.997	551	567	593	rVB	353330	756440	49.25%	4.097%
9	3.855	808	834	858	rBV3	11524	40771	2.65%	0.221%
10	3.997	861	878	890	rBV2	11919	32028	2.09%	0.173%
11	4.096	890	909	930	rBV	466693	1233414	80.30%	6.680%
12	4.736	1089	1108	1134	rBV3	23704	65642	4.27%	0.356%
13	4.884	1139	1154	1171	rBV	110662	253630	16.51%	1.374%
14	4.996	1171	1189	1205	rBV2	612819	1535996	100.00%	8.319%
15	5.083	1205	1216	1235	rVB3	69455	169133	11.01%	0.916%
16	5.260	1251	1271	1278	rBV3	50108	131027	8.53%	0.710%
17	5.312	1278	1287	1305	rVB	138800	288733	18.80%	1.564%
18	5.485	1324	1341	1352	rBV3	94114	216563	14.10%	1.173%
19	5.556	1353	1363	1374	rVV2	76553	171403	11.16%	0.928%
20	5.627	1374	1385	1405	rVB2	114799	252855	16.46%	1.369%
21	5.887	1452	1466	1482	rBV	246280	482229	31.40%	2.612%
22	6.418	1610	1631	1646	rBV	692357	1518928	98.89%	8.227%
23	6.643	1689	1701	1716	rVV	57846	109143	7.11%	0.591%
24	6.739	1716	1731	1745	rVB4	27715	58750	3.82%	0.318%
25	6.906	1771	1783	1804	rVB5	25433	55856	3.64%	0.303%
26	7.067	1817	1833	1840	rBV5	8348	18303	1.19%	0.099%
27	7.260	1875	1893	1902	rBV3	24610	50085	3.26%	0.271%
28	7.443	1939	1950	1964	rVB6	7567	17123	1.11%	0.093%
29	7.566	1964	1988	2020	rBV	505685	1062109	69.15%	5.753%
30	7.710	2021	2033	2046	rVV2	55603	121017	7.88%	0.655%
31	7.774	2047	2053	2066	rVB5	22717	41707	2.72%	0.226%
32	7.919	2086	2098	2122	rVB	108614	197821	12.88%	1.071%
33	8.048	2124	2138	2149	rBV2	46542	82974	5.40%	0.449%
34	8.492	2260	2276	2284	rBV2	29952	56570	3.68%	0.306%

LSC Area Percent Report

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Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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	8.546	2284	2293	2303	rVV	67810	111415	7.25%	0.603%
36	8.607	2303	2312	2322	rVV2	40296	67610	4.40%	0.366%
37	9.090	2451	2462	2482	rBV	348286	587043	38.22%	3.179%
38	9.189	2482	2493	2504	rVV	53477	96363	6.27%	0.522%
39	9.250	2504	2512	2523	rVB4	9976	17759	1.16%	0.096%
40	9.366	2536	2548	2568	rVB6	33263	85314	5.55%	0.462%
41	9.569	2598	2611	2620	rBV3	9225	16286	1.06%	0.088%
42	9.733	2640	2662	2663	rBV7	10627	24728	1.61%	0.134%
43	9.954	2709	2731	2743	rBV6	23159	48077	3.13%	0.260%
44	10.048	2743	2760	2772	rBV10	11145	27549	1.79%	0.149%
45	10.167	2787	2797	2808	rBV	43131	66951	4.36%	0.363%
46	10.308	2829	2841	2855	rBV	325002	520180	33.87%	2.817%
47	10.376	2856	2862	2879	rVB10	7880	15584	1.01%	0.084%
48	10.456	2879	2887	2894	rBV4	12283	18912	1.23%	0.102%
49	10.588	2918	2928	2944	rBV	21940	38180	2.49%	0.207%
50	10.977	3039	3049	3069	rBV3	31433	75895	4.94%	0.411%
51	11.077	3069	3080	3083	rBV4	11908	18884	1.23%	0.102%
52	11.151	3095	3103	3112	rVB	18248	26533	1.73%	0.144%
53	11.324	3150	3157	3174	rVB	42116	74216	4.83%	0.402%
54	11.485	3195	3207	3222	rBV	394521	614448	40.00%	3.328%
55	11.572	3223	3234	3246	rVB	363611	551061	35.88%	2.985%
56	11.688	3258	3270	3280	rBV	31890	51563	3.36%	0.279%
57	11.768	3281	3295	3310	rBV2	483413	859225	55.94%	4.654%
58	11.864	3317	3325	3345	rVB2	37016	76985	5.01%	0.417%
59	11.983	3351	3362	3374	rBV	53663	85472	5.56%	0.463%
60	12.057	3374	3385	3400	rVV	102177	161813	10.53%	0.876%
61	12.250	3428	3445	3451	rBV	147670	230994	15.04%	1.251%
62	12.286	3451	3456	3467	rVV	90983	141373	9.20%	0.766%
63	12.356	3467	3478	3498	rVV2	182351	372397	24.24%	2.017%
64	12.549	3526	3538	3550	rVB3	75141	139364	9.07%	0.755%
65	12.649	3551	3569	3580	rVV	105035	179869	11.71%	0.974%
66	12.787	3599	3612	3623	rBV3	18712	30838	2.01%	0.167%
67	12.929	3647	3656	3660	rBV3	19342	30100	1.96%	0.163%
68	12.964	3660	3667	3676	rVB	56688	82521	5.37%	0.447%
69	13.122	3694	3716	3729	rBV2	380472	627533	40.86%	3.399%
70	13.289	3757	3768	3785	rVB	97355	164872	10.73%	0.893%
71	13.408	3791	3805	3812	rBV6	11300	20691	1.35%	0.112%

LSC Area Percent Report

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Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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	13.456	3812	3820	3828	rVV2	22494	38280	2.49%	0.207%
73	13.507	3828	3836	3845	rVV4	44422	75064	4.89%	0.407%
74	13.578	3845	3858	3862	rVV5	27160	58653	3.82%	0.318%
75	13.610	3863	3868	3887	rVB3	40241	86820	5.65%	0.470%
76	13.742	3896	3909	3917	rBV2	50554	91466	5.95%	0.495%
77	13.787	3918	3923	3935	rVB2	16218	24685	1.61%	0.134%
78	13.855	3935	3944	3954	rBV4	19781	33058	2.15%	0.179%
79	13.954	3961	3975	3986	rBV3	30004	57783	3.76%	0.313%
80	14.012	3986	3993	4004	rVB2	13127	20243	1.32%	0.110%
81	14.154	4027	4037	4041	rBV3	13220	20704	1.35%	0.112%
82	14.347	4083	4097	4109	rBV5	26873	63098	4.11%	0.342%
83	14.540	4149	4157	4171	rVB5	12548	22982	1.50%	0.124%
84	14.678	4190	4200	4218	rBV2	35881	63403	4.13%	0.343%
85	14.877	4250	4262	4273	rBV	68951	116689	7.60%	0.632%
86	15.064	4308	4320	4335	rBV	73295	127503	8.30%	0.691%
87	15.916	4572	4585	4597	rBV3	13251	25263	1.64%	0.137%
88	16.060	4621	4630	4637	rBV3	18184	30140	1.96%	0.163%
89	16.099	4637	4642	4658	rVB2	12401	19332	1.26%	0.105%

Sum of corrected areas: 18463379

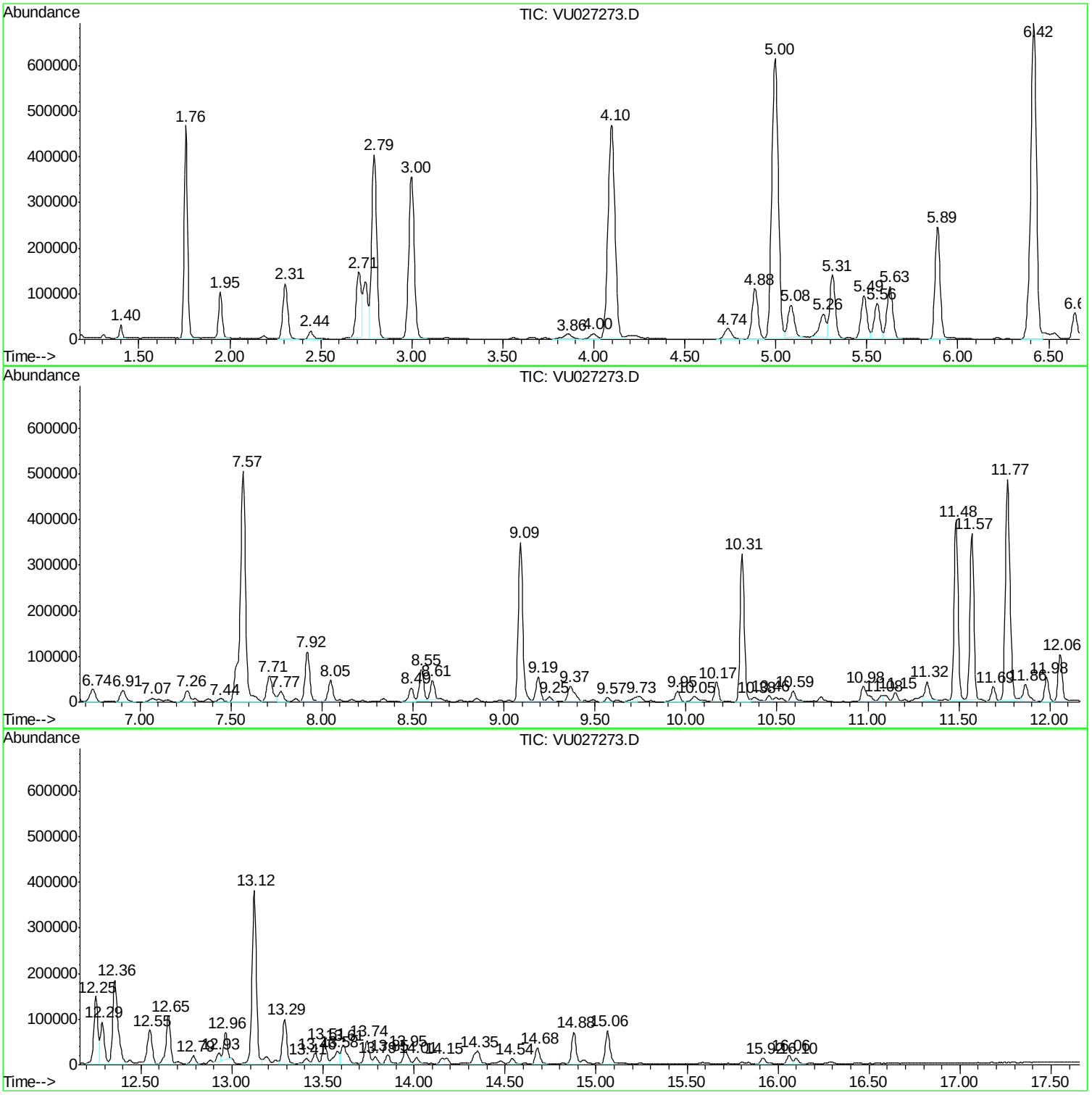
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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 : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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

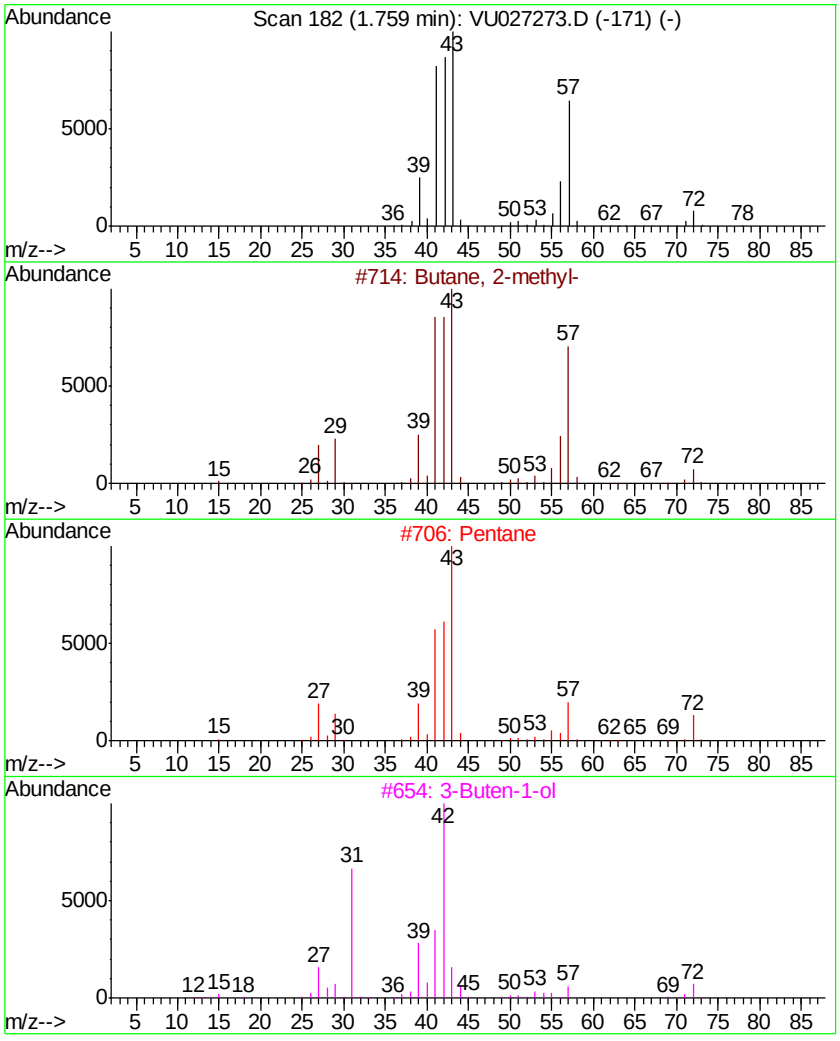
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

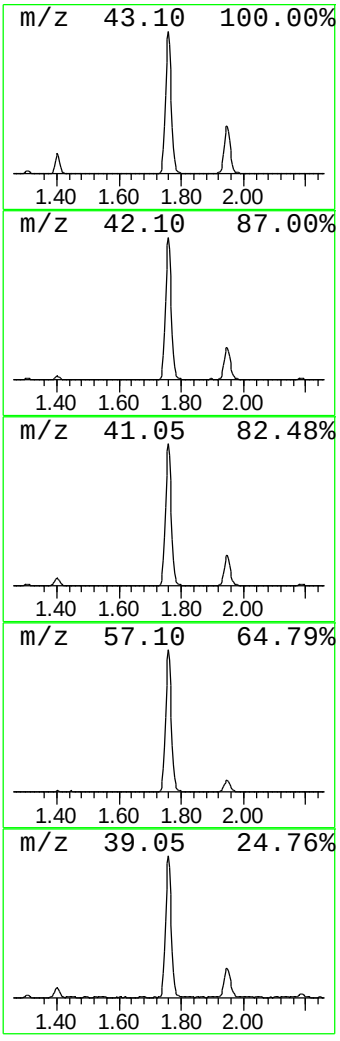
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	19.81 ug/l	608538	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2	Pentane	72	C5H12	000109-66-0	33
3	3-Buten-1-ol	72	C4H8O	000627-27-0	25
4	1-Butene	56	C4H8	000106-98-9	10
5	Methane, isocyanato-	57	C2H3NO	000624-83-9	9





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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

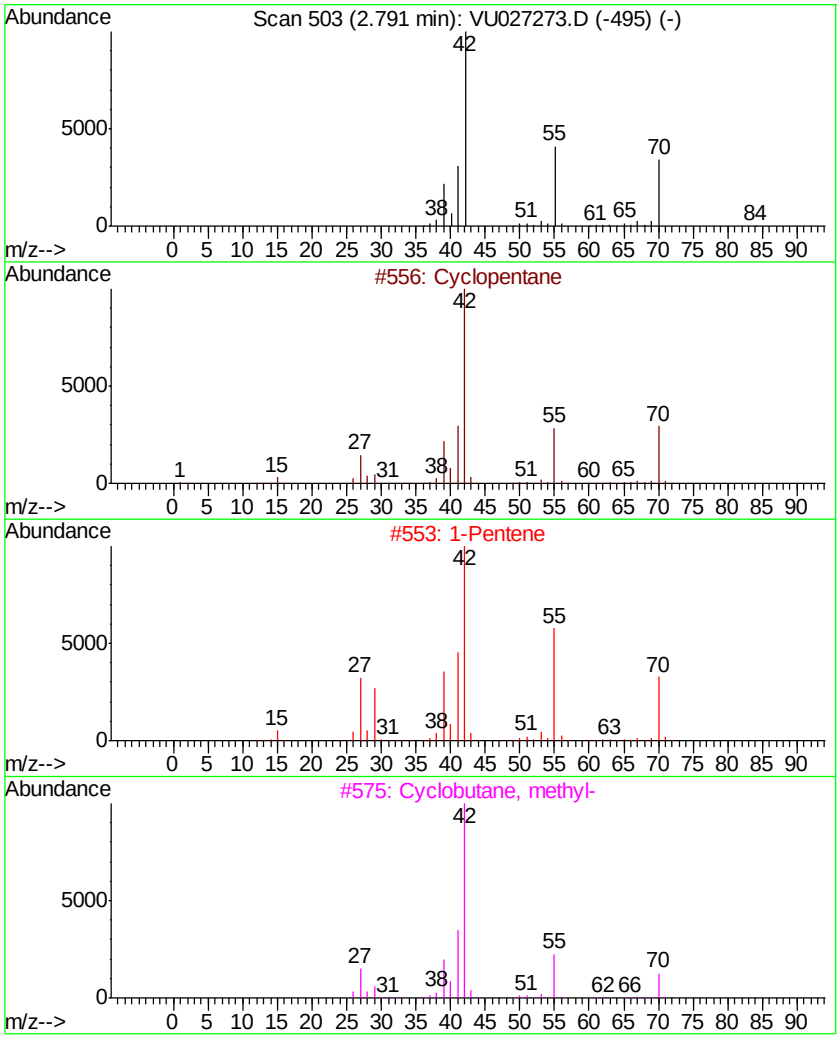
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

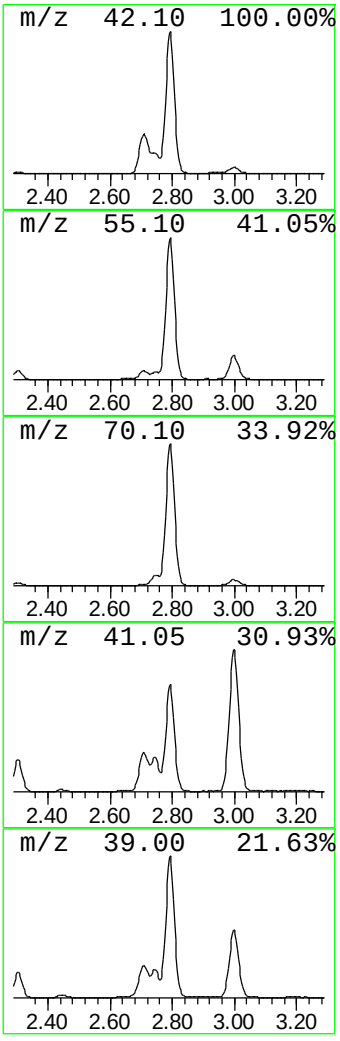
****

Peak Number 2 Cyclopentane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	25.26 ug/l	775868	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane	70	C5H10	000287-92-3	86
2	1-Pentene	70	C5H10	000109-67-1	78
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	9
5	Methyl propargyl ether	70	C4H6O	000627-41-8	7





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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

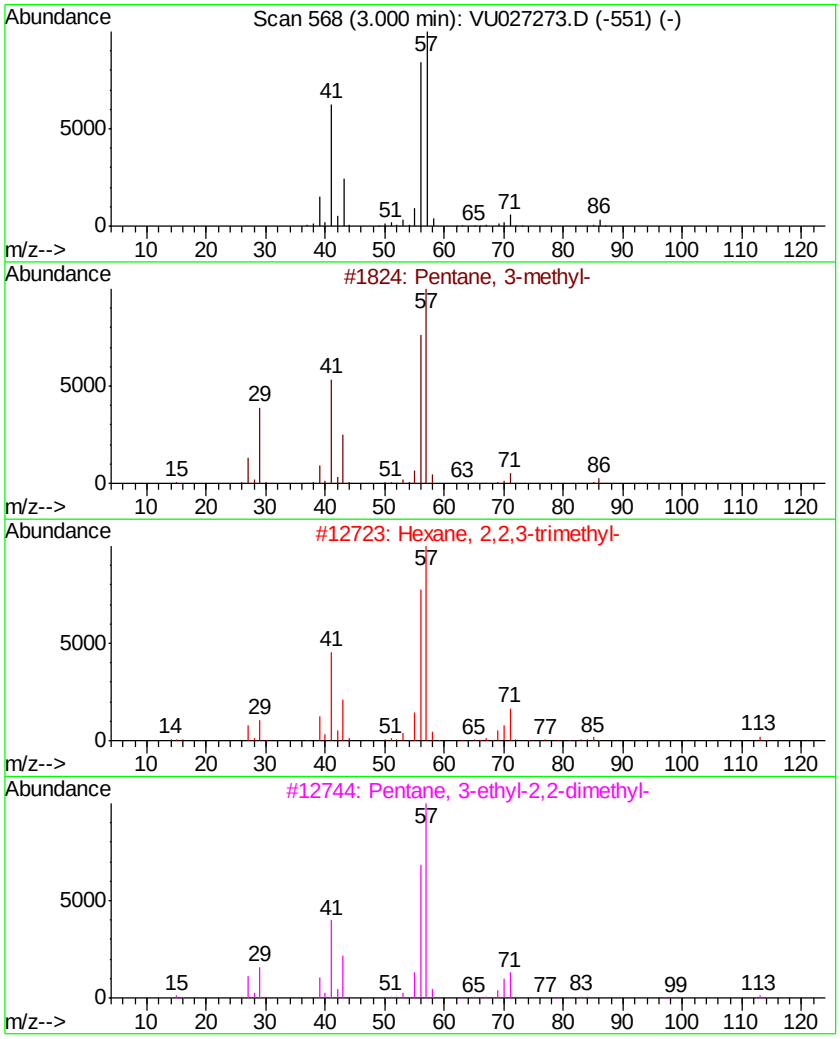
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

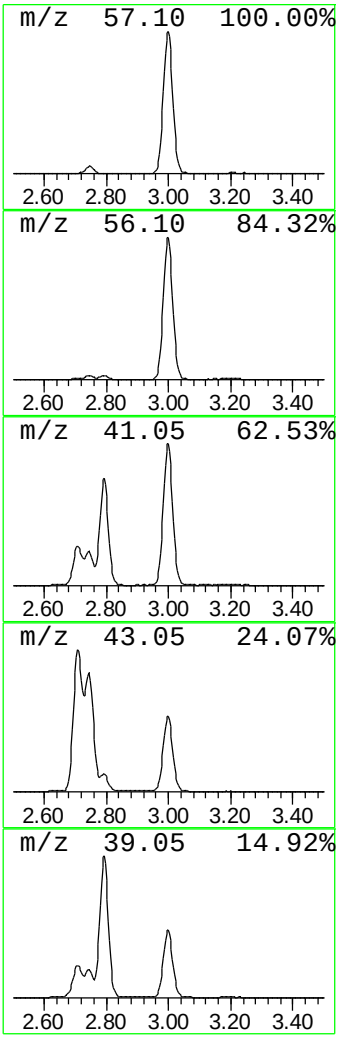
****

Peak Number 3 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	24.62 ug/l	756440	Pentafluorobenzene	4.99

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	78
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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

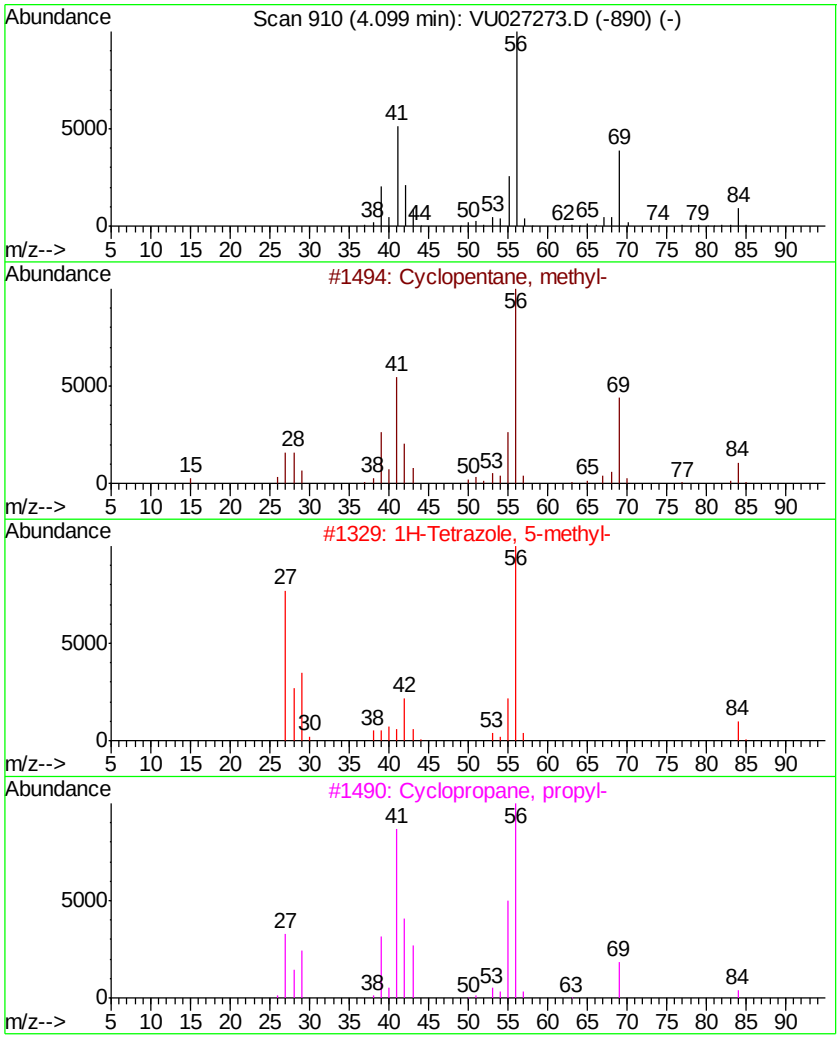
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

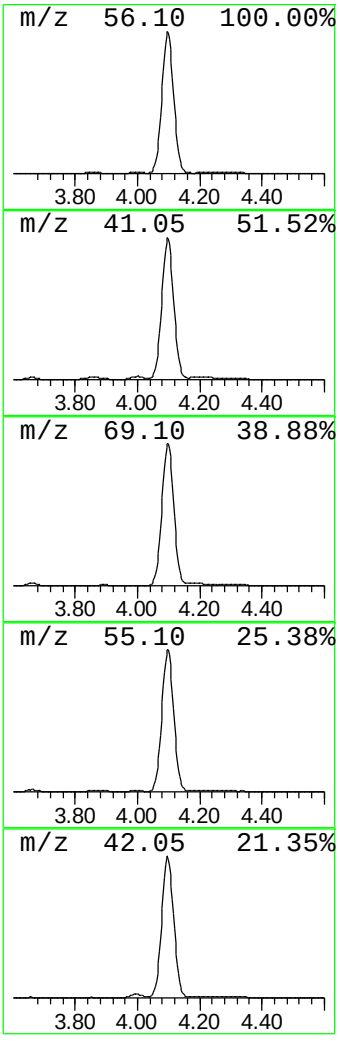
****

Peak Number 4 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.10	40.15 ug/l	1233410	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	78
4	Cyclohexane	84	C6H12	000110-82-7	78
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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

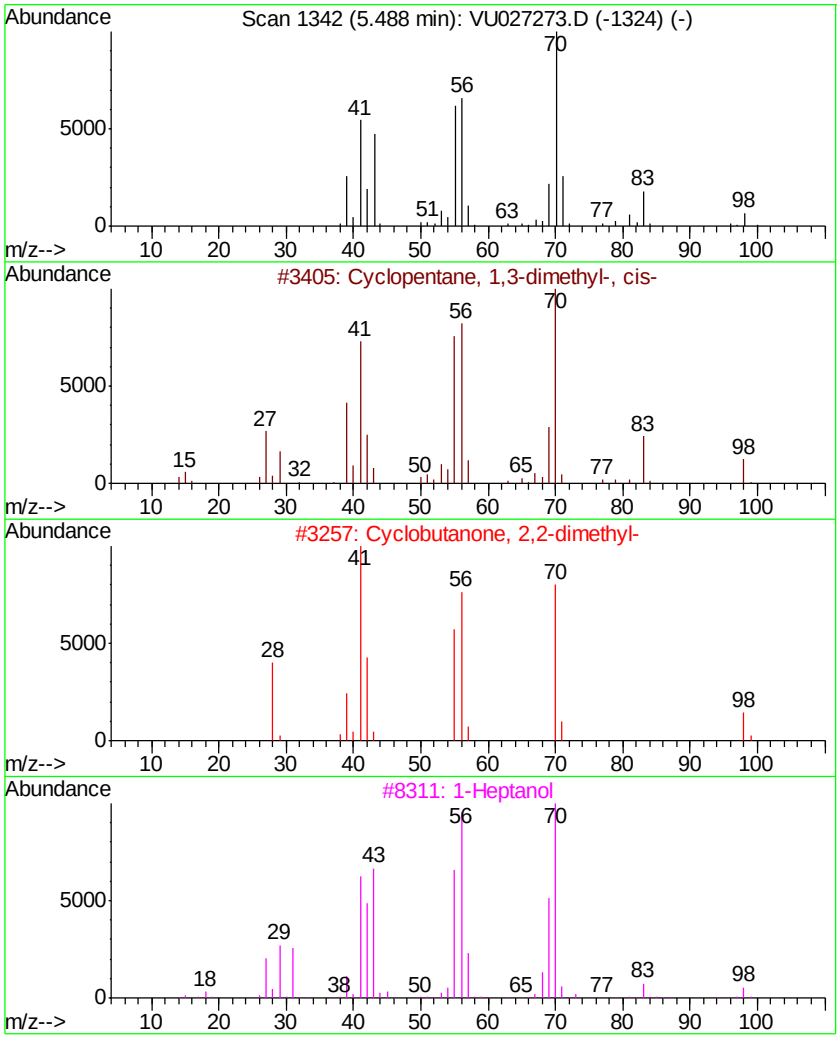
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

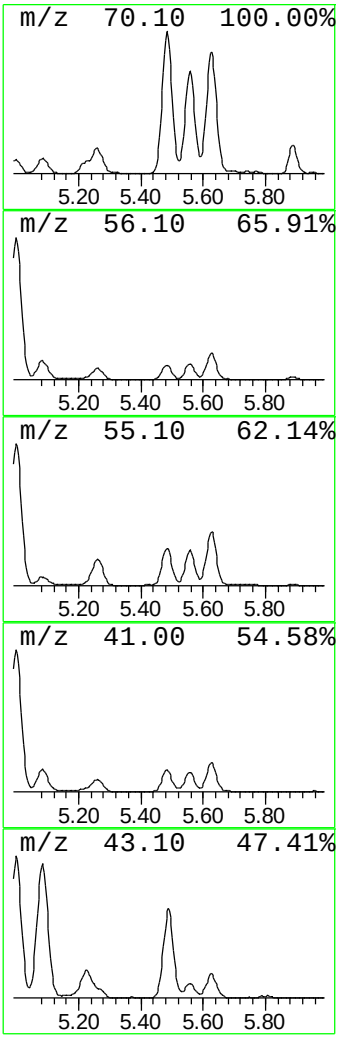
****

Peak Number 5 Cyclopentane, 1,3-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.49	22.45 ug/l	216563	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	91
2	Cyclobutanone, 2,2-dimethyl-	98	C6H10O	001192-14-9	72
3	1-Heptanol	116	C7H16O	000111-70-6	64
4	2,3-Dimethyl-aziridine	71	C4H9N	002549-68-0	64
5	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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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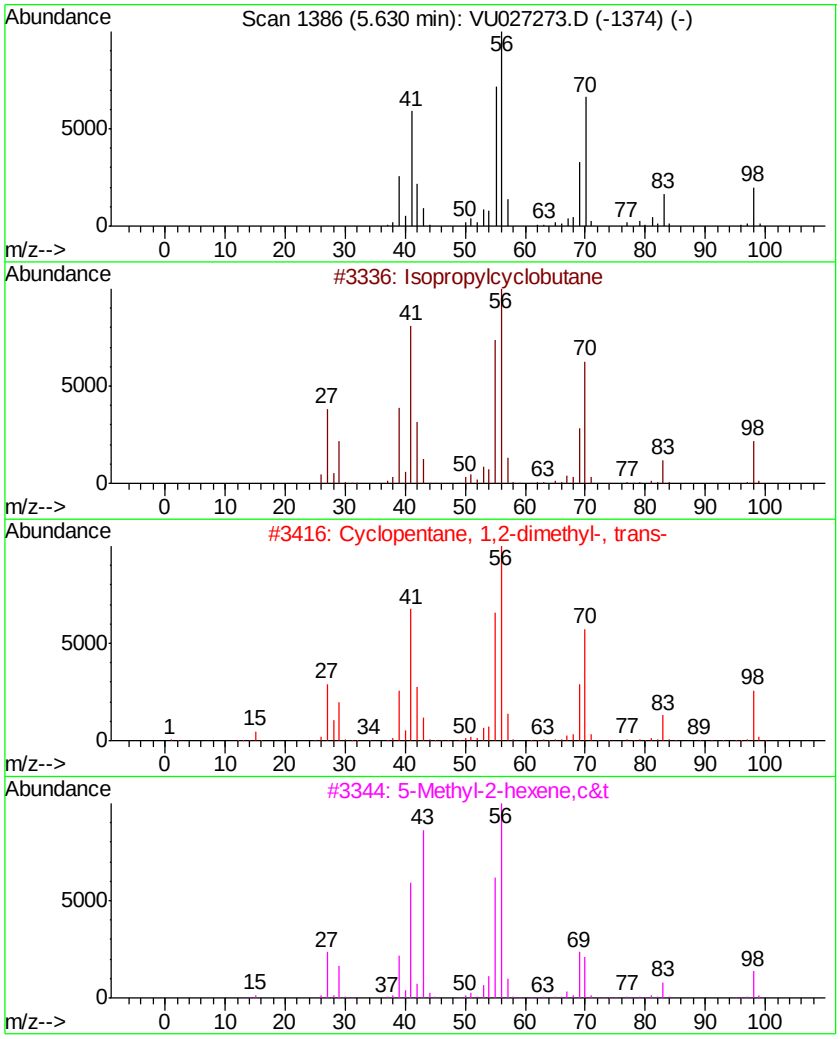
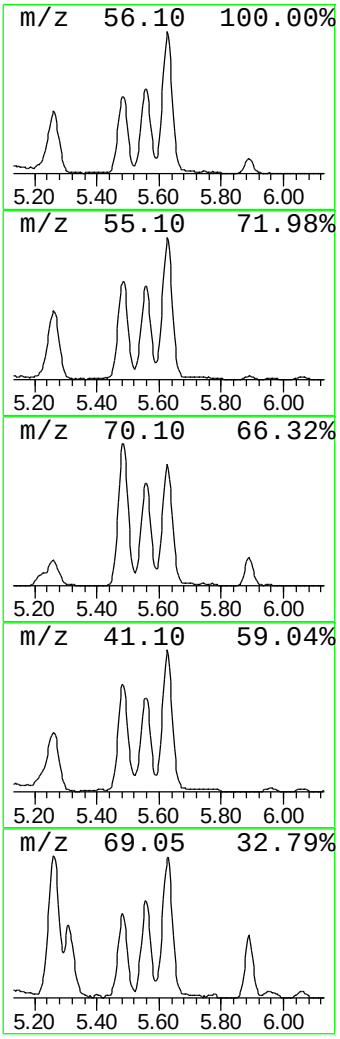
Peak Number 6 Isopropylcyclobutane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	26.22 ug/l	252855	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropylcyclobutane	98	C7H14	000872-56-0	93
2	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
3	5-Methyl-2-hexene,c&t	98	C7H14	003404-62-4	72
4	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	72
5	1-Heptene	98	C7H14	000592-76-7	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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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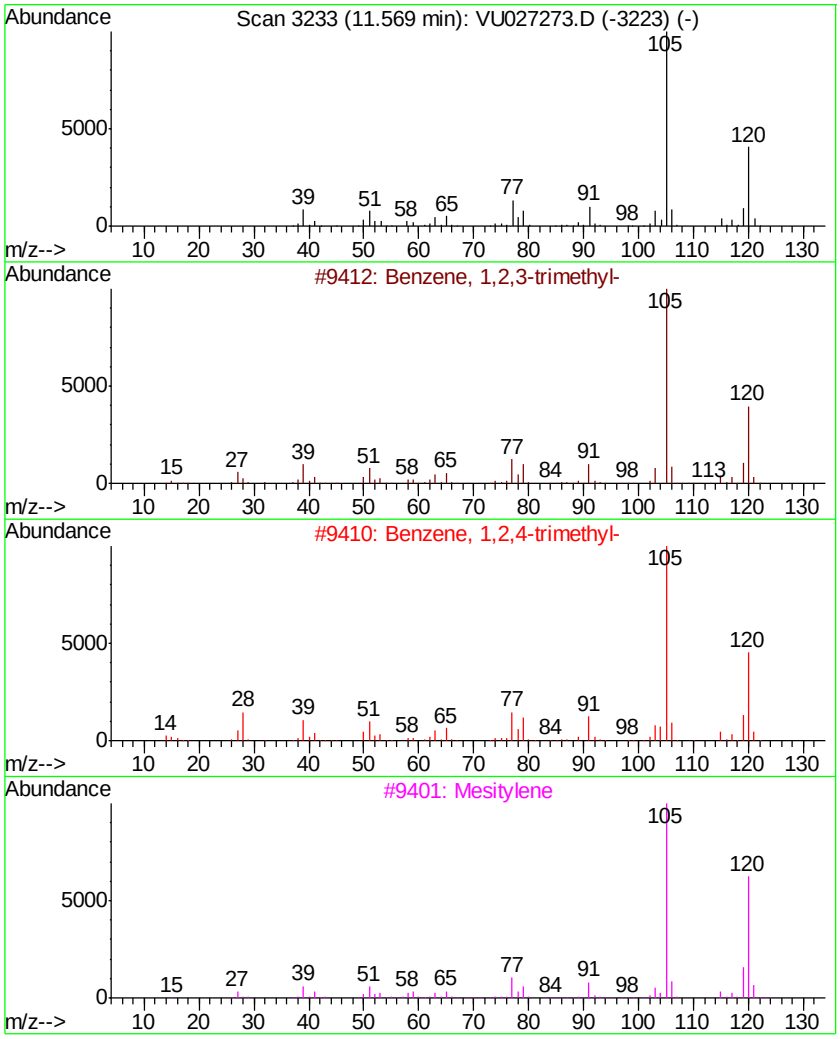
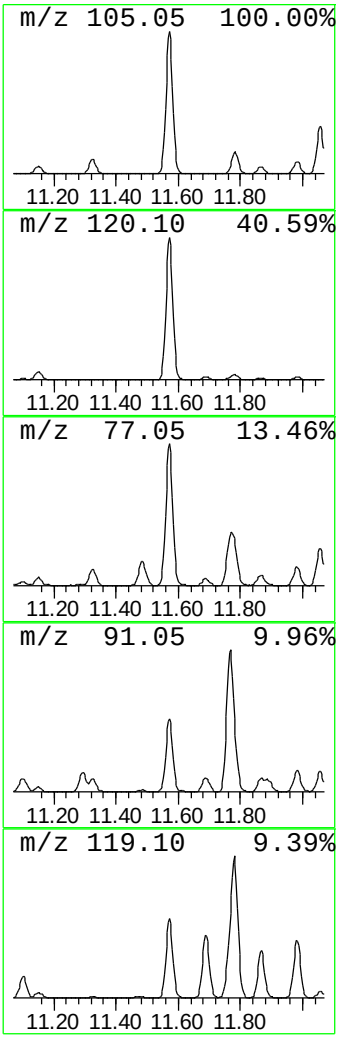
Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	44.84 ug/l	551061	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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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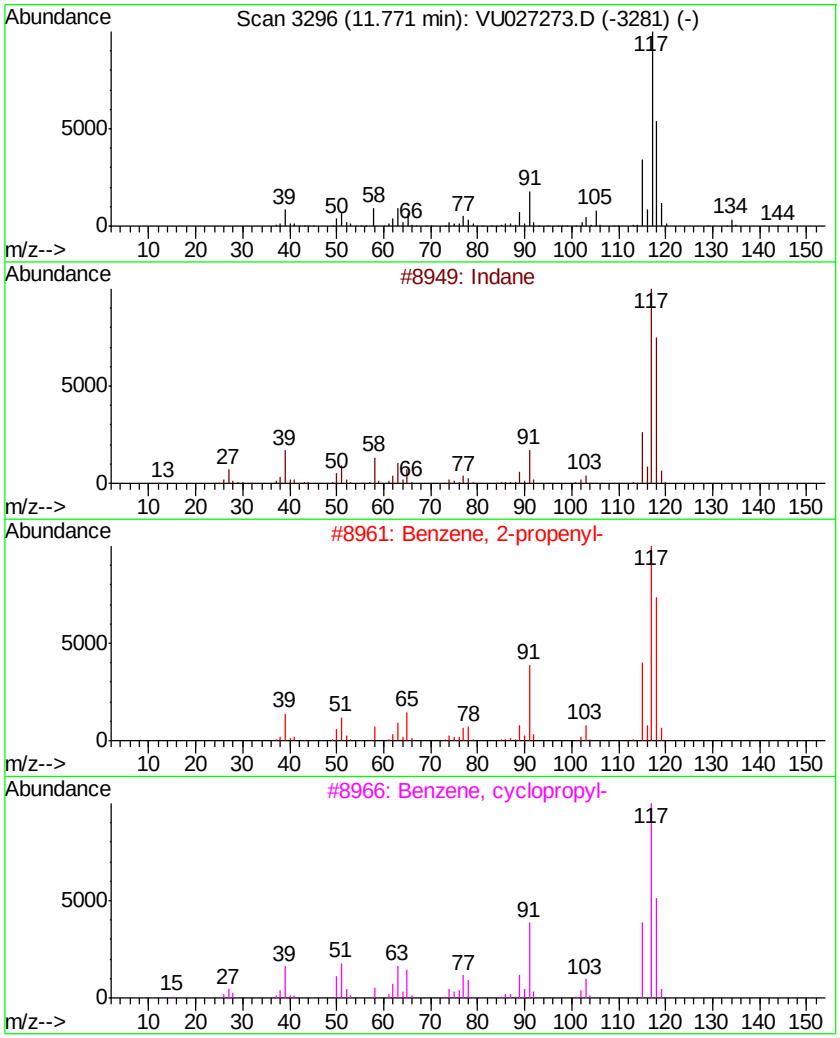
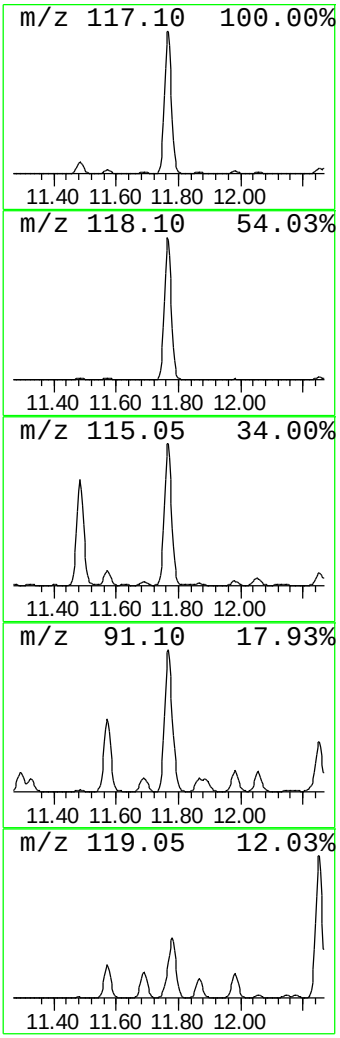
Peak Number 8 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	69.92 ug/l	859225	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	96
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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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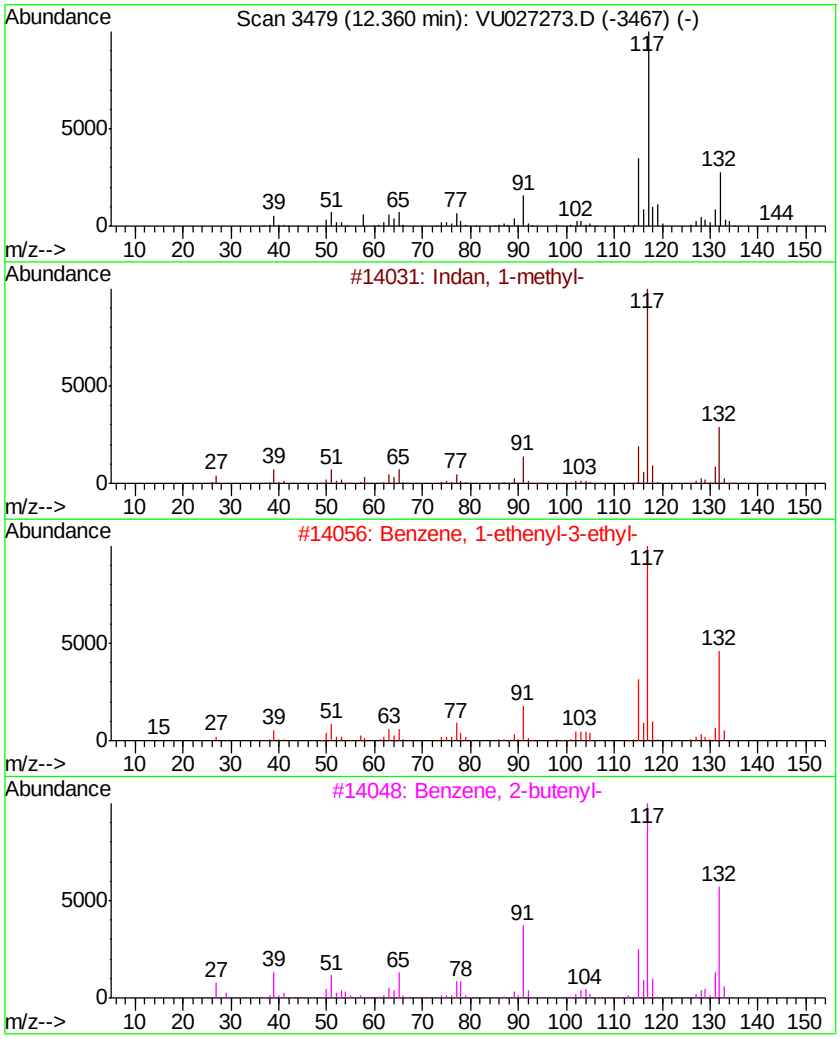
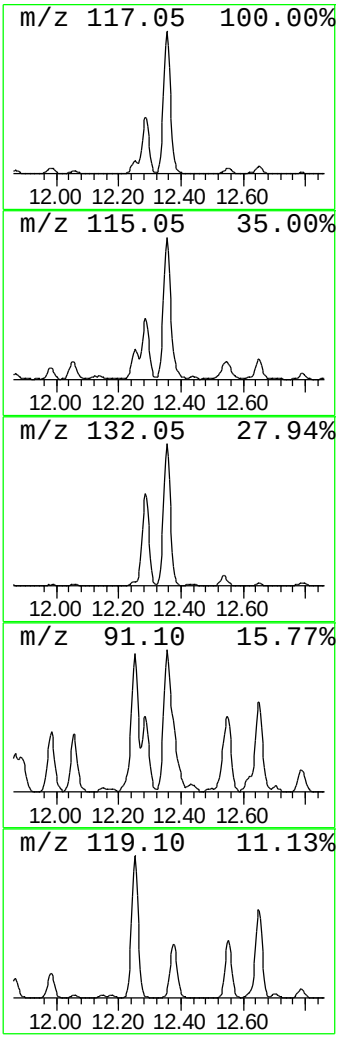
Peak Number 9 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	30.30 ug/l	372397	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	83





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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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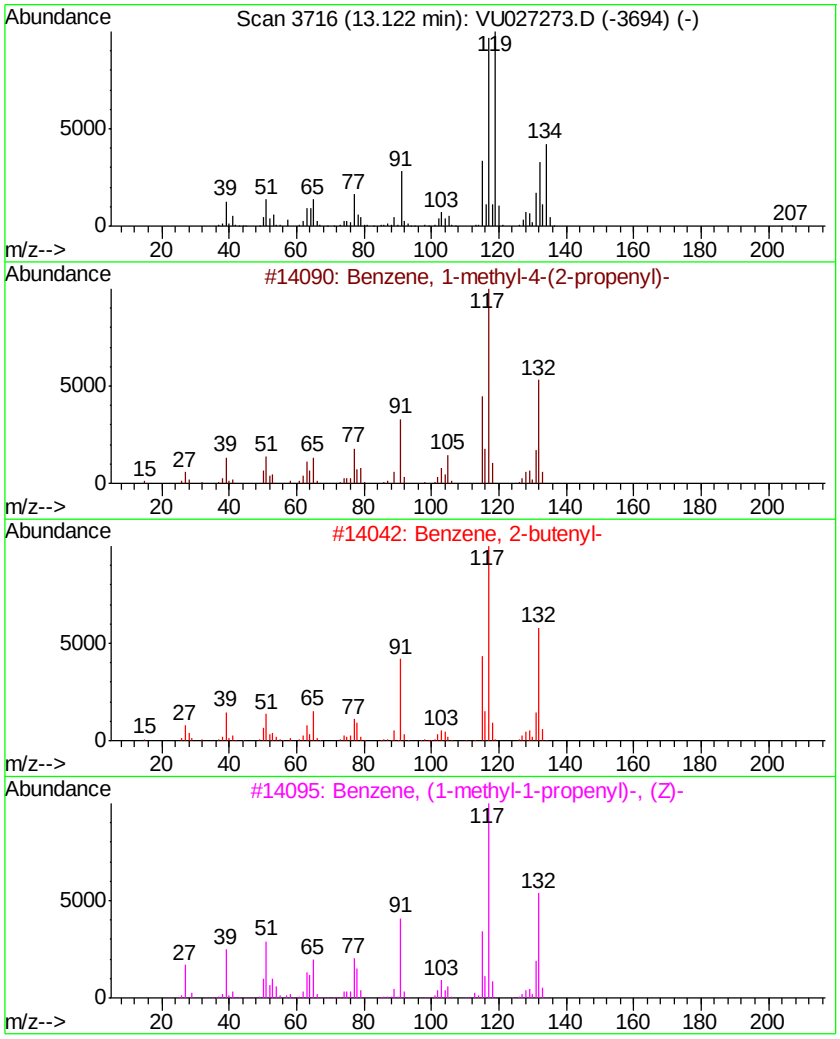
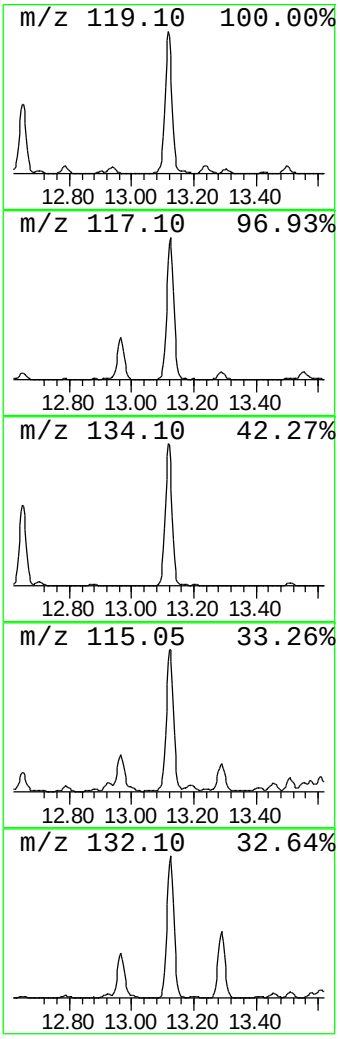
Peak Number 10 Benzene, 1-methyl-4-(2-prop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	51.06 ug/l	627533	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	50
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	50
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027273.D
Acq On : 28 Sep 2018 15:52
Operator : MD/SY
Sample : J5041-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0247

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, 2-methyl-	1.76	19.8	ug/l	608538	1	4.99	1536000	50.0
Cyclopentane	2.79	25.3	ug/l	775868	1	4.99	1536000	50.0
Pentane, 3-methyl-	3.00	24.6	ug/l	756440	1	4.99	1536000	50.0
Cyclopentane, met...	4.10	40.1	ug/l	1233410	1	4.99	1536000	50.0
Cyclopentane, 1,3...	5.49	22.4	ug/l	216563	2	5.89	482229	50.0
Isopropylcyclobutane	5.63	26.2	ug/l	252855	2	5.89	482229	50.0
Benzene, 1,2,3-tr...	11.57	44.8	ug/l	551061	4	11.48	614448	50.0
Indane	11.77	69.9	ug/l	859225	4	11.48	614448	50.0
Indan, 1-methyl-	12.36	30.3	ug/l	372397	4	11.48	614448	50.0
Benzene, 1-methyl...	13.12	51.1	ug/l	627533	4	11.48	614448	50.0