LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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	1.759	172	182	195	rBV	96181	125877	9.49%	0.871%
2	1.948	232	241	254	rVB	86206	115104	8.68%	0.797%
3	2.305	339	352	370	rVB4	37420	75862	5.72%	0.525%
4	2.443	383	395	416	rBV	16879	31228	2.35%	0.216%
5	2.710	465	478	481	rBV2	63345	110896	8.36%	0.767%
6	2.746	482	489	497	rVV	119430	234350	17.66%	1.622%
7	2.791	497	503	524	rVB	83080	158485	11.95%	1.097%
8	3.000	545	568	584	rBV	175346	372472	28.08%	2.578%
9	3.305	649	663	683	rVB2	30307	66554	5.02%	0.461%
10	3.861	811	836	853	rBV4	8788	27769	2.09%	0.192%
11	4.000	863	879	890	rBV4	8543	23213	1.75%	0.161%
12	4.096	890	909	930	rVV	276583	742154	55.94%	5.136%
13	4.202	931	942	964	rVB5	5375	21212	1.60%	0.147%
14	4.736	1087	1108	1133	rBV3	21813	61977	4.67%	0.429%
15	4.884	1139	1154	1171	rBV	100765	227667	17.16%	1.576%
16	4.993	1171	1188	1204	rVV4	377270	947924	71.45%	6.560%
17	5.083	1204	1216	1234	rVB3	63906	160403	12.09%	1.110%
18	5.222	1244	1259	1267	rBV	65032	153644	11.58%	1.063%
19	5.312	1278	1287	1301	rVB2	126894	258073	19.45%	1.786%
20	5.392	1302	1312	1326	rVB2	32550	67325	5.07%	0.466%
21	5.485	1327	1341	1352	rBV2	73625	164894	12.43%	1.141%
22	5.556	1353	1363	1373	rVV	54196	116645	8.79%	0.807%
23	5.627	1373	1385	1405	rVB2	127944	285557	21.52%	1.976%
24	5.887	1452	1466	1485	rBV	220258	434514	32.75%	3.007%
25	6.418	1609	1631	1647	rBV	598048	1326699	100.00%	9.182%
26	6.643	1689	1701	1715	rVB2	47310	88669	6.68%	0.614%
27	6.739	1715	1731	1745	rVB5	33442	74682	5.63%	0.517%
28	6.906	1770	1783	1804	rBV4	42719	93371	7.04%	0.646%
29	7.260	1886	1893	1901	rVB4	19302	31030	2.34%	0.215%
30	7.344	1912	1919	1924	rBV3	14114	21391	1.61%	0.148%
31	7.443	1939	1950	1963	rVB6	9975	22766	1.72%	0.158%
32	7.533	1963	1978	1980	rBV3	134385	203460	15.34%	1.408%
33	7.565	1981	1988	2020	rVB	476330	875543	65.99%	6.060%
34	7.710	2020	2033	2045	rBV2	64011	139407	10.51%	0.965%

LSC Area Percent Report

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Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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

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Peak separation: 5

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

35	7.778	2046	2054	2069	rVB2	44495	81227	6.12%	0.562%
36	7.919	2086	2098	2123	rVB2	143598	265378	20.00%	1.837%
37	8.048	2123	2138	2149	rBV	62195	116511	8.78%	0.806%
38	8.337	2217	2228	2238	rBV6	14961	27444	2.07%	0.190%
39	8.492	2259	2276	2284	rBV4	47894	98258	7.41%	0.680%
40	8.546	2284	2293	2302	rVB	89527	141968	10.70%	0.983%
41	8.607	2302	2312	2322	rBV	128466	216284	16.30%	1.497%
42	8.758	2349	2359	2368	rBV2	13822	23985	1.81%	0.166%
43	8.848	2380	2387	2400	rVV3	29763	59856	4.51%	0.414%
44	9.090	2451	2462	2479	rVV	310435	520097	39.20%	3.600%
45	9.189	2481	2493	2503	rVV	52654	94653	7.13%	0.655%
46	9.250	2503	2512	2527	rVB4	48072	87199	6.57%	0.603%
47	9.366	2536	2548	2550	rBV6	39842	62764	4.73%	0.434%
48	9.572	2602	2612	2624	rVB2	23515	40486	3.05%	0.280%
49	9.723	2642	2659	2676	rBV6	40792	129486	9.76%	0.896%
50	9.954	2702	2731	2744	rBV4	56104	122453	9.23%	0.847%
51	10.048	2750	2760	2785	rBV8	27419	72509	5.47%	0.502%
52	10.218	2804	2813	2826	rBV8	8244	19450	1.47%	0.135%
53	10.308	2830	2841	2854	rBV	286624	455200	34.31%	3.150%
54	10.379	2854	2863	2872	rVB3	21432	33285	2.51%	0.230%
55	10.495	2890	2899	2911	rVB3	38513	64939	4.89%	0.449%
56	10.707	2955	2965	2972	rBV7	13290	22747	1.71%	0.157%
57	10.974	3037	3048	3066	rBV	107530	186327	14.04%	1.290%
58	11.151	3093	3103	3114	rVB	122370	186338	14.05%	1.290%
59	11.273	3130	3141	3150	rBV9	11004	23677	1.78%	0.164%
60	11.334	3152	3160	3172	rVV8	9355	20136	1.52%	0.139%
61	11.485	3196	3207	3219	rBV	357143	559298	42.16%	3.871%
62	11.572	3220	3234	3245	rBV	138572	213759	16.11%	1.479%
63	11.688	3262	3270	3281	rVB2	20519	34702	2.62%	0.240%
64	11.774	3282	3297	3314	rVV4	68334	158238	11.93%	1.095%
65	11.868	3315	3326	3342	rVB6	54353	126756	9.55%	0.877%
66	11.980	3351	3361	3373	rBV	43370	70219	5.29%	0.486%
67	12.057	3374	3385	3398	rBV	97712	152811	11.52%	1.058%
68	12.147	3399	3413	3417	rBV	33859	55374	4.17%	0.383%
69	12.176	3417	3422	3433	rVB	49611	71689	5.40%	0.496%
70	12.250	3436	3445	3453	rBV	77920	117014	8.82%	0.810%
71	12.292	3454	3458	3470	rVB8	21485	34633	2.61%	0.240%

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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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

72	12.356	3470	3478	3482	rBV	46268	70630	5.32%	0.489%
73	12.498	3512	3522	3528	rBV7	24297	44777	3.38%	0.310%
74	12.549	3529	3538	3549	rVB3	74778	141636	10.68%	0.980%
75	12.623	3553	3561	3563	rBV2	22382	28150	2.12%	0.195%
76	12.649	3564	3569	3577	rVV	45227	63948	4.82%	0.443%
77	12.704	3577	3586	3600	rVB	93468	160051	12.06%	1.108%
78	12.790	3603	3613	3627	rBV3	40835	73585	5.55%	0.509%
79	12.932	3648	3657	3664	rBV5	32450	55332	4.17%	0.383%
80	13.080	3695	3703	3707	rBV3	19018	30064	2.27%	0.208%
81	13.122	3707	3716	3730	rVB2	148790	242421	18.27%	1.678%
82	13.289	3760	3768	3776	rVB3	46701	76125	5.74%	0.527%
83	13.408	3797	3805	3813	rBV3	15685	28893	2.18%	0.200%
84	13.459	3814	3821	3827	rVB	14971	19725	1.49%	0.137%
85	13.507	3827	3836	3843	rBV4	59226	100241	7.56%	0.694%
86	13.732	3893	3906	3916	rBV5	34609	68338	5.15%	0.473%
87	13.787	3916	3923	3935	rVB	33448	52472	3.96%	0.363%
88	13.858	3935	3945	3961	rBV4	24099	52030	3.92%	0.360%
89	13.958	3965	3976	3987	rBV5	34566	77097	5.81%	0.534%
90	14.015	3988	3994	4002	rVB4	13588	20823	1.57%	0.144%
91	14.150	4027	4036	4041	rBV5	11914	20047	1.51%	0.139%
92	14.343	4084	4096	4110	rBV8	37834	95305	7.18%	0.660%
93	14.478	4130	4138	4148	rBV3	17550	28985	2.18%	0.201%
94	14.546	4149	4159	4170	rVB5	22281	43633	3.29%	0.302%
95	14.610	4170	4179	4190	rVB4	15259	23403	1.76%	0.162%
96	14.681	4190	4201	4227	rBV4	39120	87752	6.61%	0.607%
97	14.867	4250	4259	4273	rVV10	12644	33946	2.56%	0.235%
98	14.938	4275	4281	4295	rVB4	14844	28729	2.17%	0.199%
99	15.070	4312	4322	4333	rVB6	12094	27792	2.09%	0.192%
100	15.588	4473	4483	4499	rVB6	12188	25135	1.89%	0.174%

Sum of corrected areas: 14449032

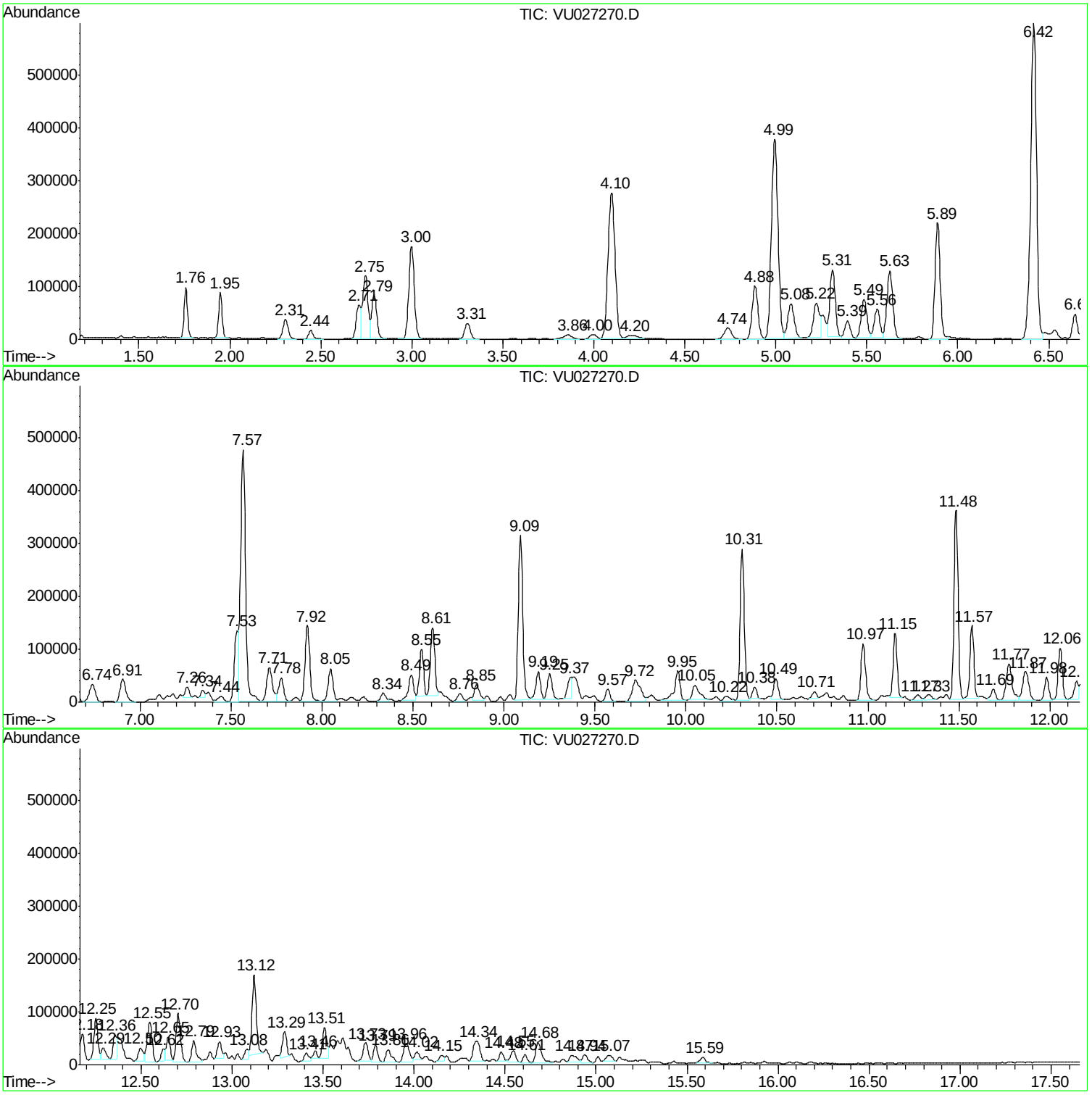
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
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Quant Title : SW846 8260

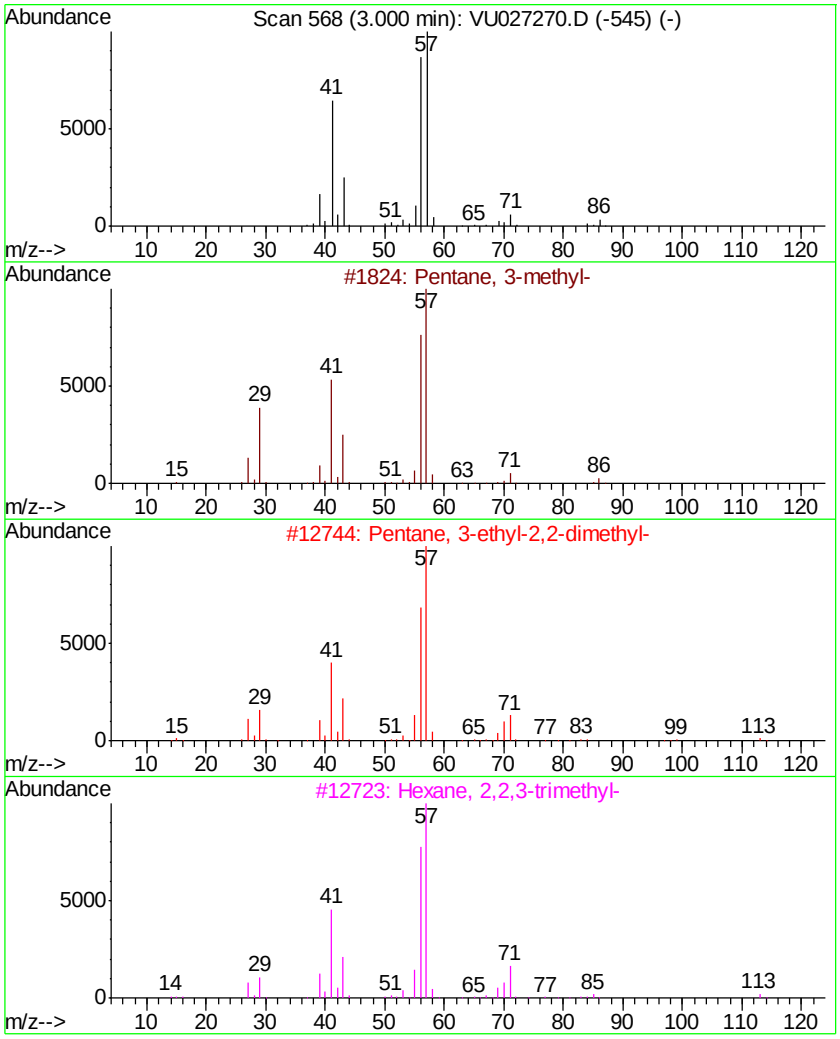
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

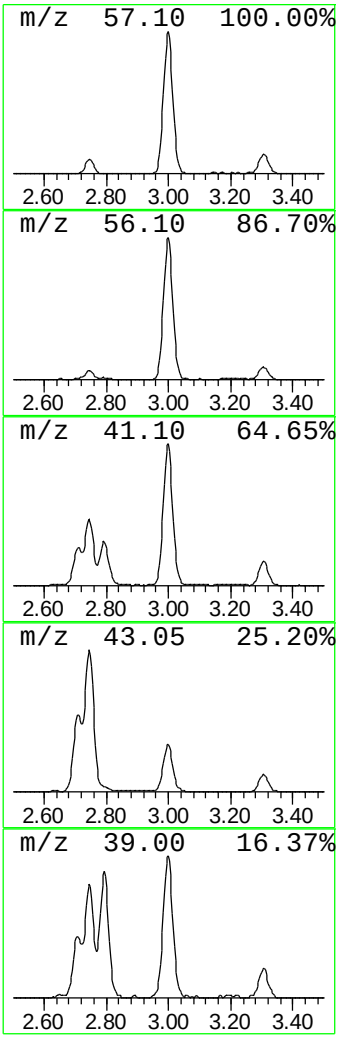
****

Peak Number 1 Pentane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	19.65 ug/l	372472	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	53
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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

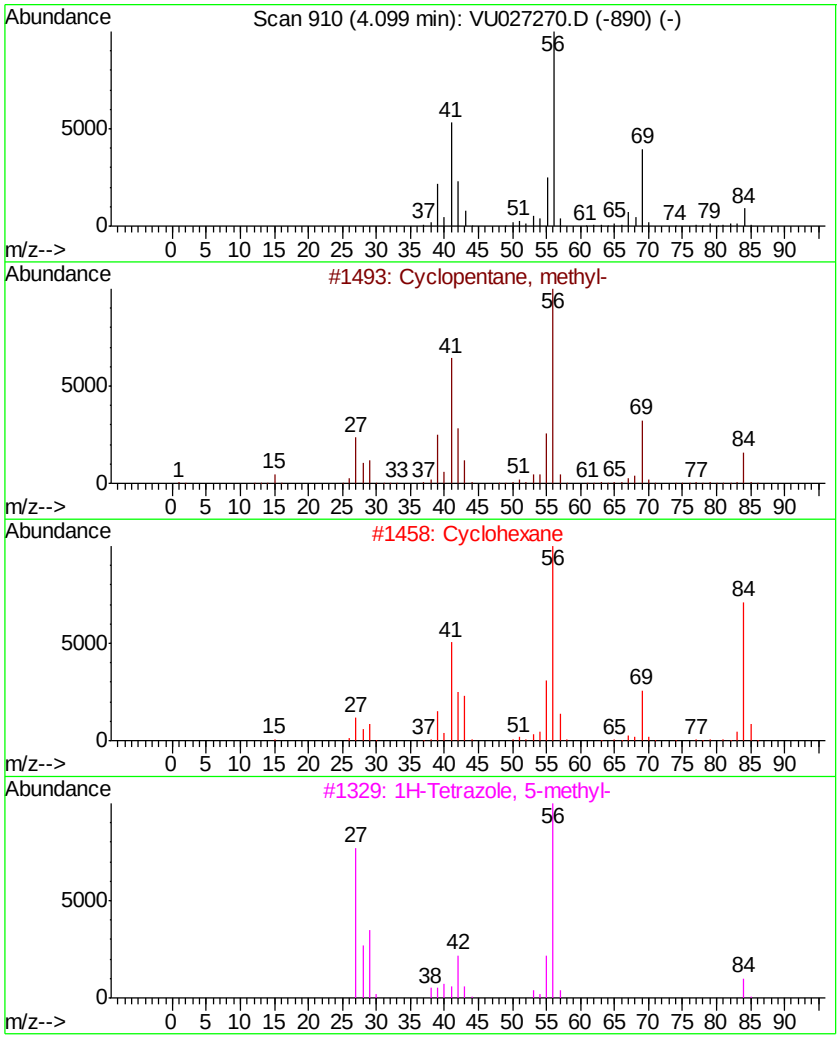
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

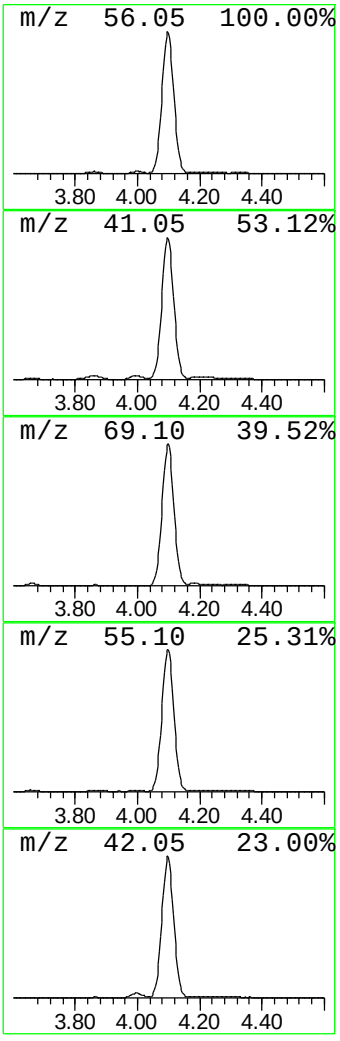
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.10	39.15 ug/l	742154	Pentafluorobenzene	4.99

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	72
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Cyclobutane, butyl-	112	C8H16	013152-44-8	50





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ClientSampleId : S13-0258

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

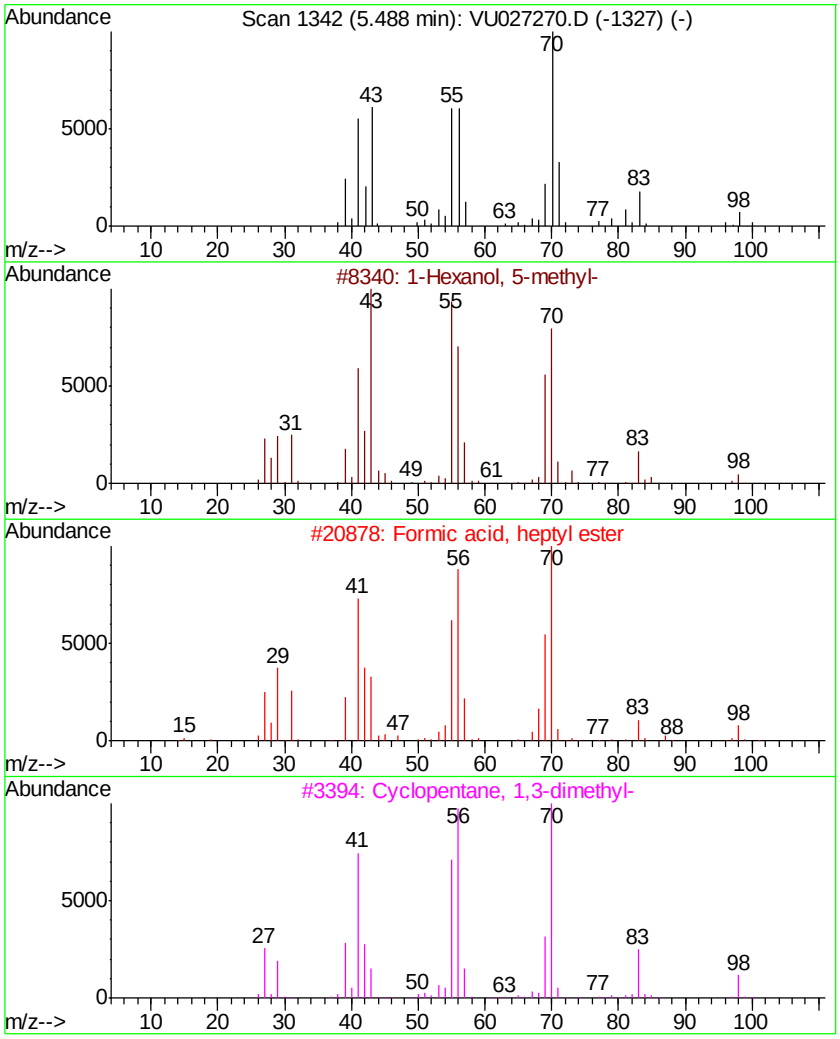
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

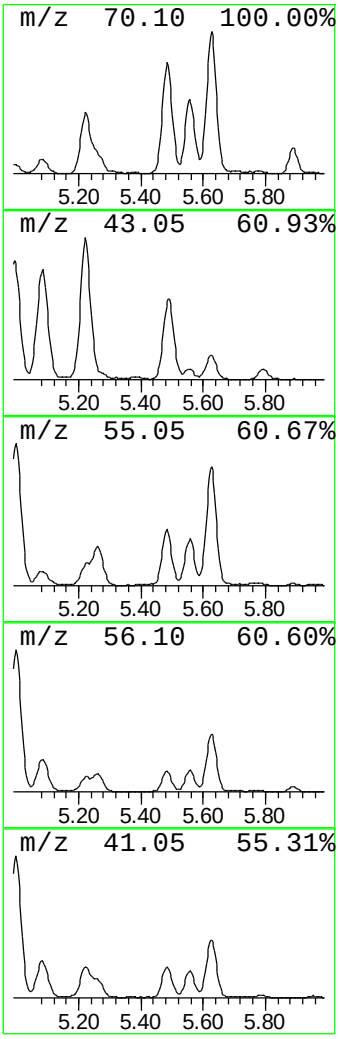
****

Peak Number 3 1-Hexanol, 5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.49	18.97 ug/l	164894	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 5-methyl-	116	C7H16O	000627-98-5	64
2	Formic acid, heptyl ester	144	C8H16O2	000112-23-2	64
3	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	64
4	2,3-Dimethyl-aziridine	71	C4H9N	002549-68-0	64
5	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	60





Library Search Compound Report

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Acq On : 28 Sep 2018 14:42
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Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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

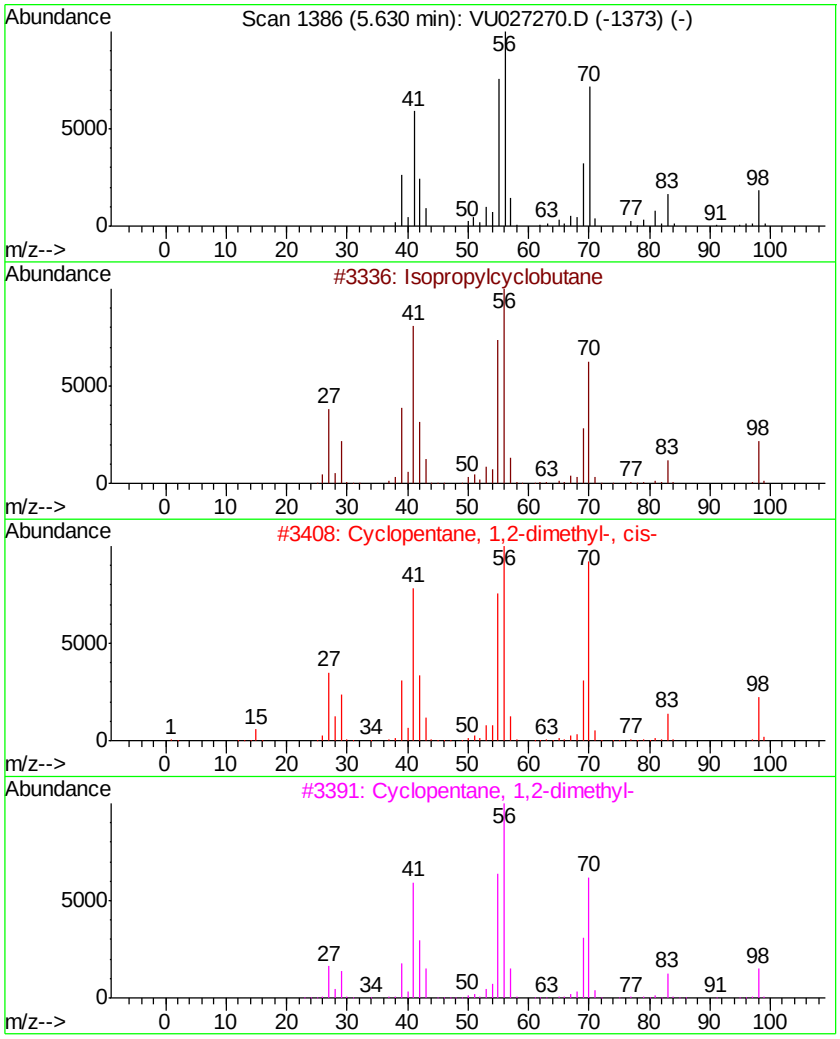
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

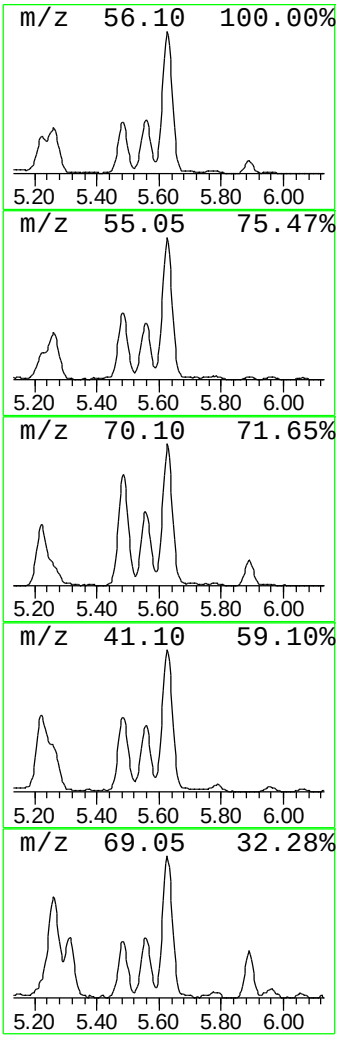
****

Peak Number 4 Isopropylcyclobutane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	32.86 ug/l	285557	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-, cis-	98	C7H14	001192-18-3	91
3	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
4	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
5	5-Methyl-2-hexene,c&t	98	C7H14	003404-62-4	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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

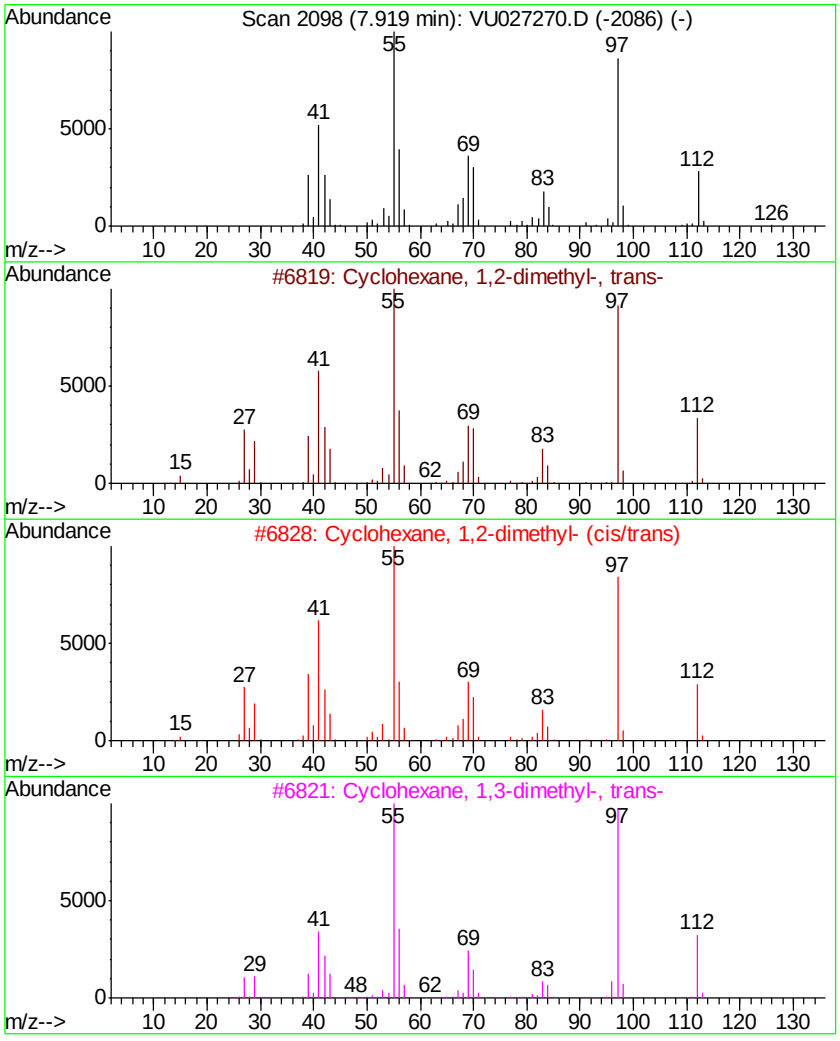
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

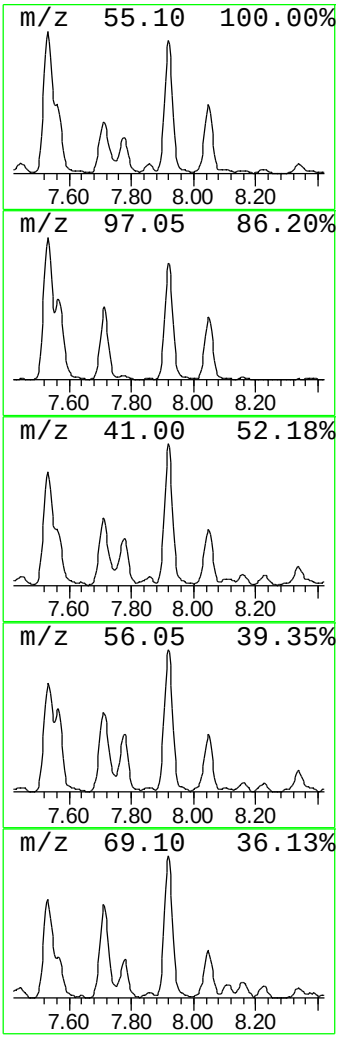
****

Peak Number 5 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.92	25.51 ug/l	265378	Chlorobenzene-d5	9.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	97
2	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	96
3	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
4	Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	76
5	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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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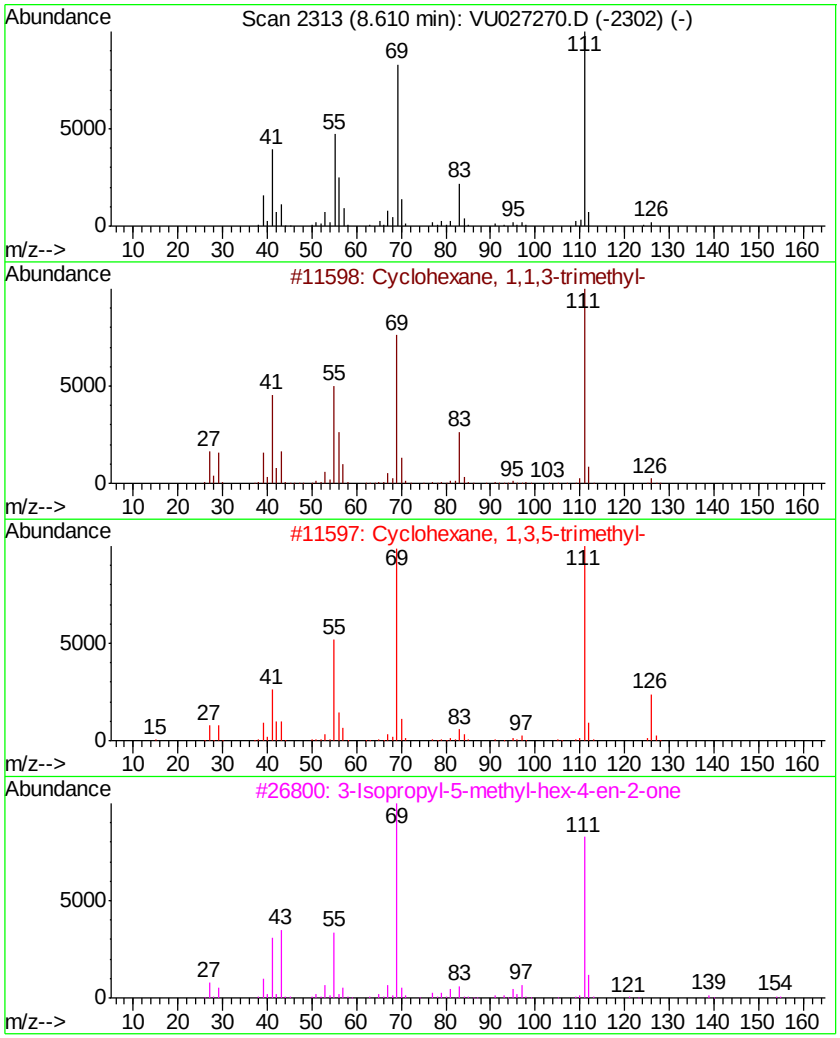
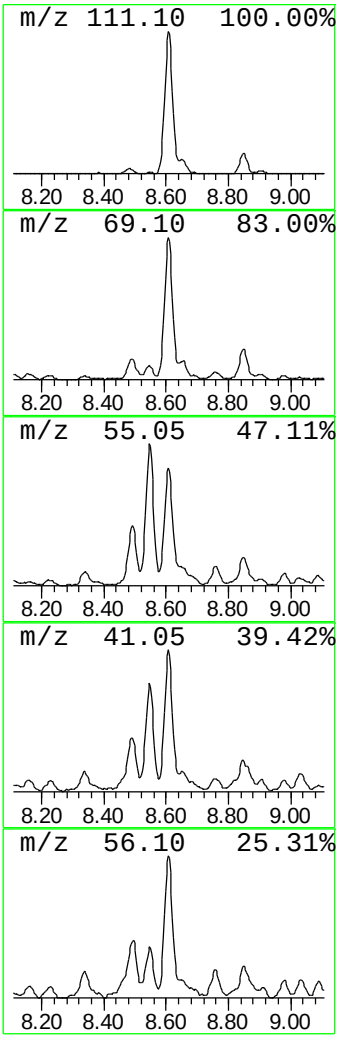
Peak Number 6 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	20.79 ug/l	216284	Chlorobenzene-d5	9.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	97
2	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
3	3-Isopropyl-5-methyl-hex-4-en-2-one	154	C10H18O	077142-85-9	59
4	Cyclopentane, 1,1,3-trimethyl-3-...	166	C12H22	074421-09-3	59
5	3-Heptene, 2-methyl-, (E)-	112	C8H16	000692-96-6	49





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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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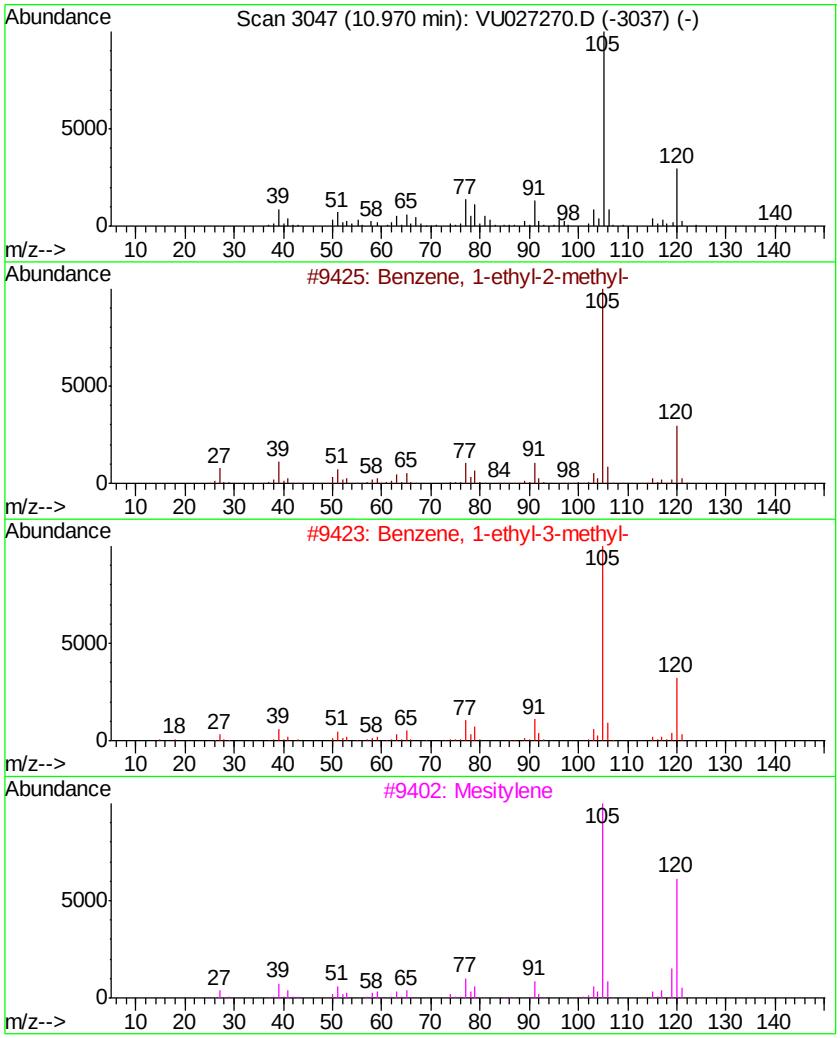
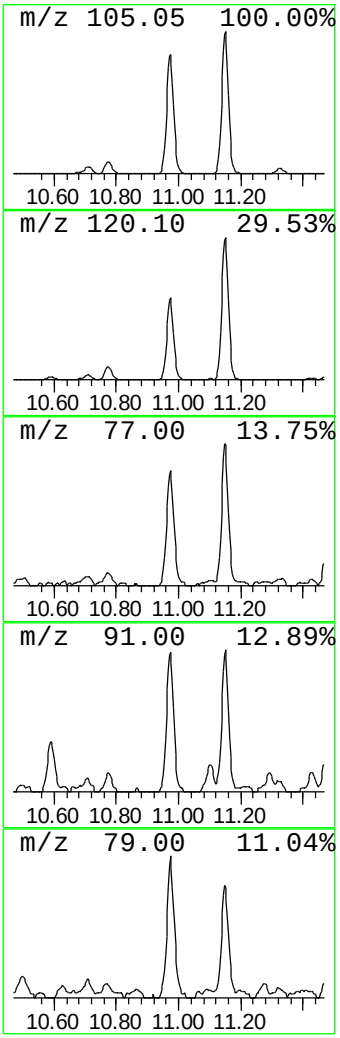
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	16.66 ug/l	186327	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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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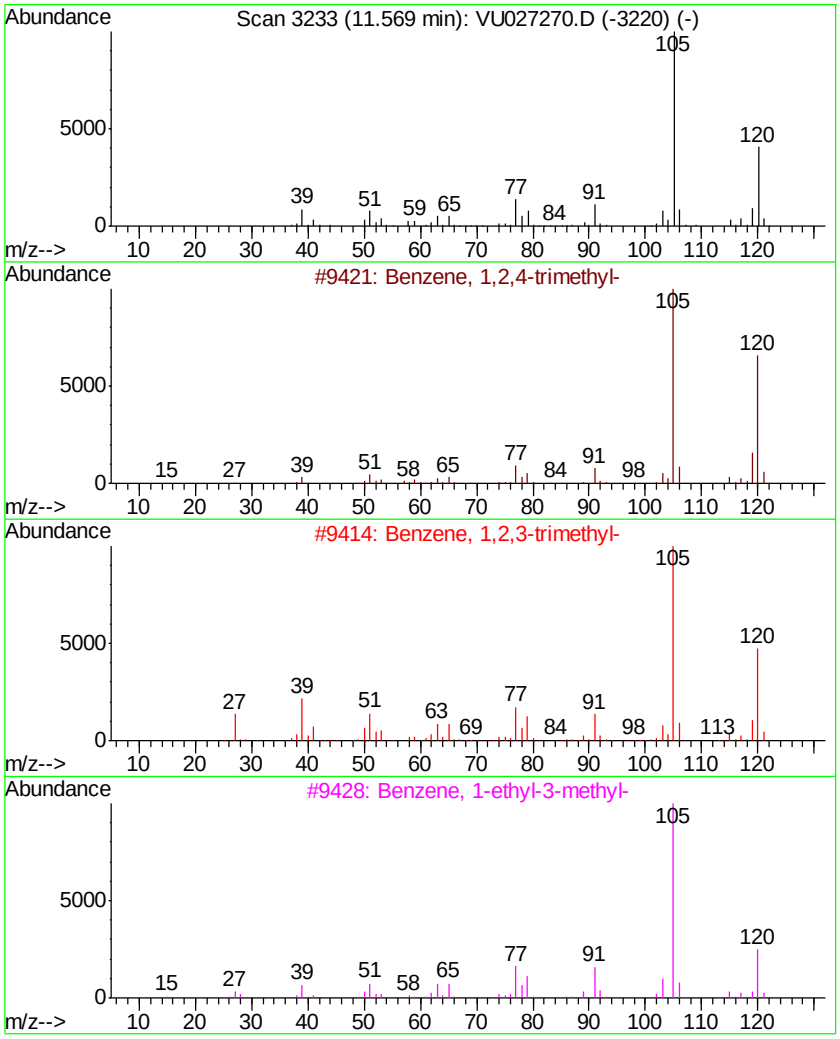
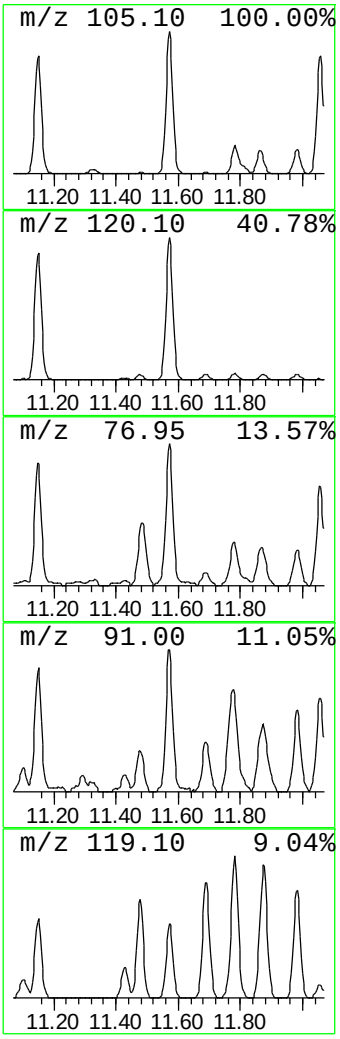
Peak Number 8 Benzene, 1,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	19.11 ug/l	213759	1,4-Dichlorobenzene-d4	11.48

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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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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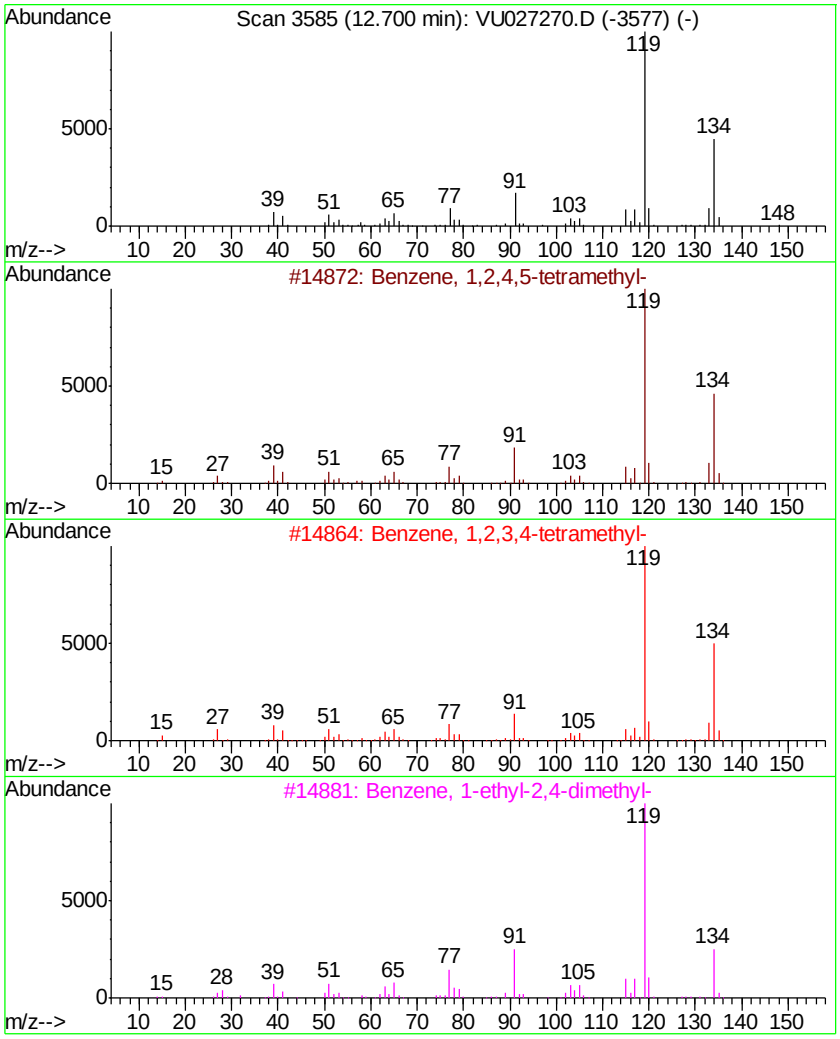
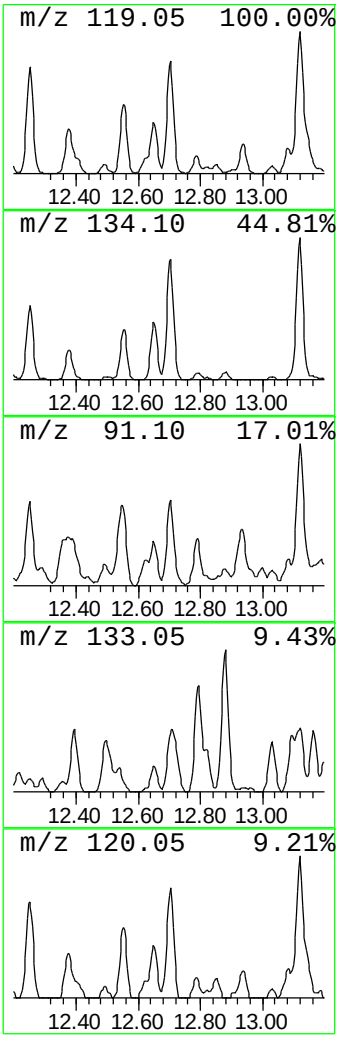
Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	14.31 ug/l	160051	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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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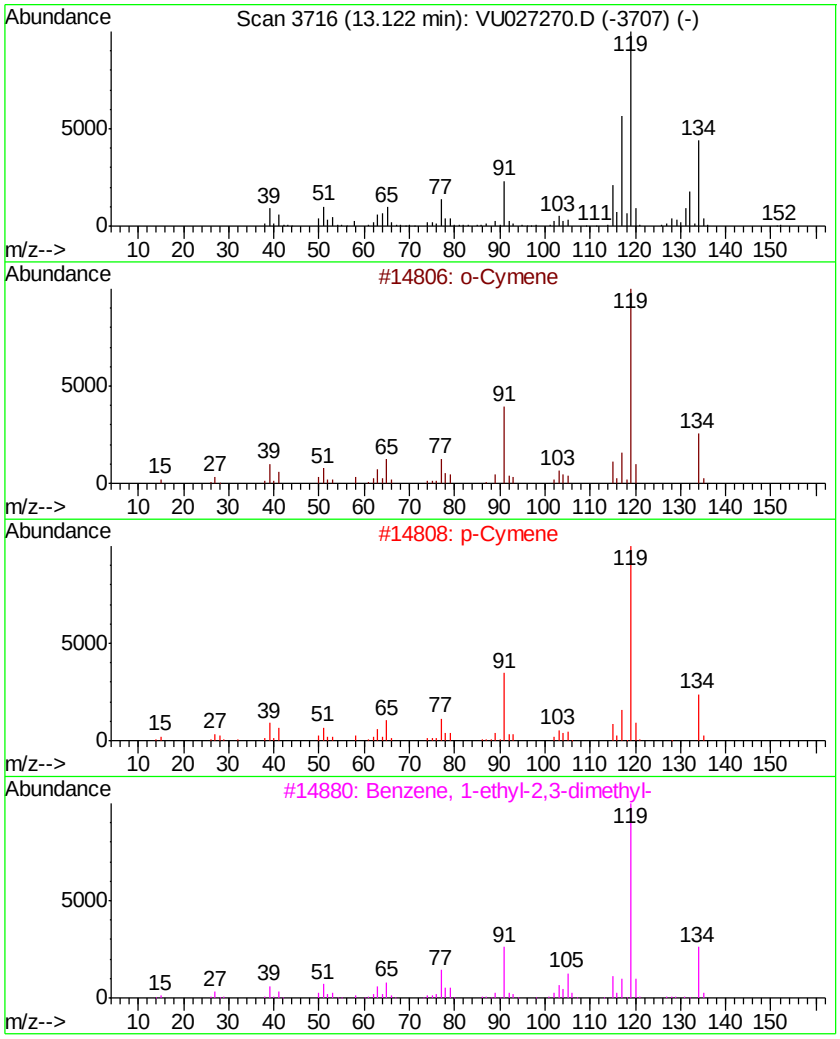
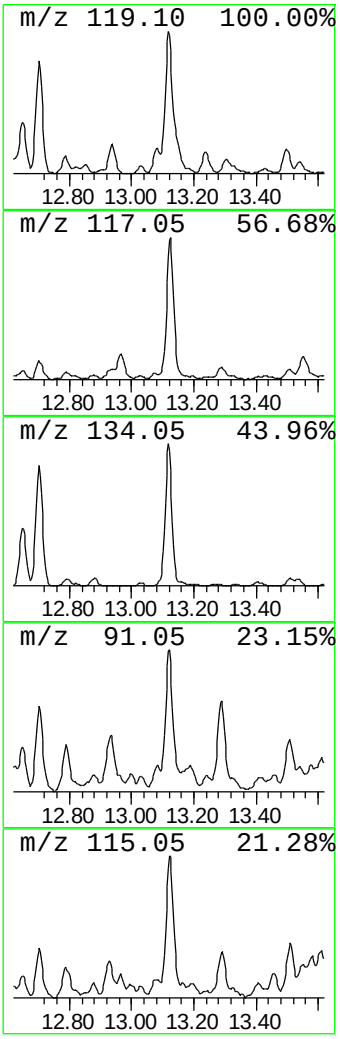
Peak Number 10 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	21.67 ug/l	242421	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	64
2	p-Cymene	134	C10H14	000099-87-6	64
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	64
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027270.D
Acq On : 28 Sep 2018 14:42
Operator : MD/SY
Sample : J5041-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0258

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
Pentane, 3-methyl-	3.00	19.6	ug/l	372472	1	4.99	947924	50.0
Cyclopentane, met...	4.10	39.1	ug/l	742154	1	4.99	947924	50.0
1-Hexanol, 5-methyl-	5.49	19.0	ug/l	164894	2	5.89	434514	50.0
Isopropylcyclobutane	5.63	32.9	ug/l	285557	2	5.89	434514	50.0
Cyclohexane, 1,2-...	7.92	25.5	ug/l	265378	3	9.09	520097	50.0
Cyclohexane, 1,1,...	8.61	20.8	ug/l	216284	3	9.09	520097	50.0
Benzene, 1-ethyl-...	10.97	16.7	ug/l	186327	4	11.48	559298	50.0
Benzene, 1,2,4-tr...	11.57	19.1	ug/l	213759	4	11.48	559298	50.0
Benzene, 1,2,4,5-...	12.70	14.3	ug/l	160051	4	11.48	559298	50.0
o-Cymene	13.12	21.7	ug/l	242421	4	11.48	559298	50.0