LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027275.D
Acq On : 28 Sep 2018 16:38
Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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	204	rVB	390852	510136	5.55%	1.055%
2	2.305	336	352	372	rVB	69094	135439	1.47%	0.280%
3	2.707	459	477	497	rBV	419080	871150	9.48%	1.801%
4	3.000	550	568	592	rBV	915711	1940647	21.13%	4.013%
5	3.996	857	878	899	rBV2	113972	320232	3.49%	0.662%
6	4.736	1090	1108	1129	rBV3	35753	98730	1.07%	0.204%
7	4.884	1138	1154	1170	rBV	118032	268346	2.92%	0.555%
8	5.000	1170	1190	1205	rBV	2068823	5025653	54.71%	10.392%
9	5.083	1205	1216	1253	rVB	416357	1020123	11.11%	2.109%
10	5.260	1256	1271	1281	rBV	272348	655025	7.13%	1.354%
11	5.482	1324	1340	1351	rBV2	630788	1431180	15.58%	2.959%
12	5.556	1352	1363	1374	rVV3	629349	1500660	16.34%	3.103%
13	5.627	1374	1385	1415	rVB	776552	1734376	18.88%	3.586%
14	5.887	1451	1466	1484	rBV	257599	509614	5.55%	1.054%
15	6.421	1607	1632	1646	rBV	4253054	9185924	100.00%	18.994%
16	6.739	1715	1731	1748	rBV	168689	346346	3.77%	0.716%
17	6.910	1768	1784	1804	rBV3	229364	626296	6.82%	1.295%
18	7.051	1814	1828	1839	rBV	170547	357308	3.89%	0.739%
19	7.260	1878	1893	1902	rBV	113190	231207	2.52%	0.478%
20	7.530	1963	1977	1982	rBV2	608841	1091360	11.88%	2.257%
21	7.565	1983	1988	2019	rVB2	792567	1373359	14.95%	2.840%
22	7.710	2019	2033	2046	rBV2	259259	582757	6.34%	1.205%
23	7.778	2047	2054	2068	rVB	148636	261782	2.85%	0.541%
24	7.919	2086	2098	2122	rVB	537190	988947	10.77%	2.045%
25	8.048	2123	2138	2149	rBV	240825	438768	4.78%	0.907%
26	8.491	2257	2276	2284	rBV3	124880	244423	2.66%	0.505%
27	8.546	2284	2293	2303	rVV	492900	798946	8.70%	1.652%
28	8.607	2303	2312	2323	rVB	267420	447378	4.87%	0.925%
29	8.848	2369	2387	2413	rVB3	54821	136170	1.48%	0.282%
30	9.089	2451	2462	2479	rBV	373800	612814	6.67%	1.267%
31	9.189	2479	2493	2503	rBV	164610	282951	3.08%	0.585%
32	9.363	2526	2547	2553	rBV3	96673	212775	2.32%	0.440%
33	9.726	2643	2660	2678	rBV5	68388	226209	2.46%	0.468%
34	9.954	2704	2731	2748	rBV2	113042	251769	2.74%	0.521%

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Sample : J5041-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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

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Title : SW846 8260

35	10.048	2748	2760	2777	rBV5	74540	153187	1.67%	0.317%
36	10.167	2787	2797	2808	rBV	397100	616813	6.71%	1.275%
37	10.308	2830	2841	2853	rBV	353675	558739	6.08%	1.155%
38	10.495	2891	2899	2916	rVB4	53997	94082	1.02%	0.195%
39	10.588	2916	2928	2949	rBV	399984	625263	6.81%	1.293%
40	11.292	3128	3147	3151	rBV	83463	148778	1.62%	0.308%
41	11.324	3151	3157	3174	rVB	204215	324997	3.54%	0.672%
42	11.485	3196	3207	3221	rBV	426685	666164	7.25%	1.377%
43	11.572	3221	3234	3247	rBV	621430	939680	10.23%	1.943%
44	11.691	3260	3271	3283	rBV	99543	158487	1.73%	0.328%
45	11.774	3283	3297	3310	rBV3	376298	774382	8.43%	1.601%
46	11.864	3310	3325	3331	rBV2	73206	134438	1.46%	0.278%
47	11.983	3350	3362	3375	rBV	125078	201564	2.19%	0.417%
48	12.253	3421	3446	3452	rBV	387123	638369	6.95%	1.320%
49	12.286	3452	3456	3468	rVV	213945	314594	3.42%	0.650%
50	12.356	3468	3478	3497	rVV3	302681	755813	8.23%	1.563%
51	12.549	3526	3538	3550	rVB4	208670	404922	4.41%	0.837%
52	12.652	3550	3570	3582	rBV	287642	491411	5.35%	1.016%
53	12.925	3647	3655	3662	rVV3	64754	113936	1.24%	0.236%
54	13.121	3693	3716	3726	rBV2	878655	1421295	15.47%	2.939%
55	13.289	3759	3768	3785	rVB2	150164	261788	2.85%	0.541%
56	13.511	3827	3837	3845	rBV3	149662	233385	2.54%	0.483%
57	13.578	3845	3858	3862	rBV4	75015	137758	1.50%	0.285%
58	13.610	3863	3868	3875	rVB	75168	94935	1.03%	0.196%
59	13.742	3893	3909	3918	rBV2	247888	440949	4.80%	0.912%
60	13.787	3918	3923	3936	rVV	65398	101569	1.11%	0.210%
61	13.961	3967	3977	3986	rBV5	54099	111088	1.21%	0.230%
62	14.154	4026	4037	4043	rBV	53890	98374	1.07%	0.203%
63	14.334	4083	4093	4109	rBV2	100929	225513	2.45%	0.466%
64	14.539	4148	4157	4170	rVB4	67609	126653	1.38%	0.262%
65	14.681	4190	4201	4217	rBV4	72220	154969	1.69%	0.320%
66	14.877	4251	4262	4272	rVV	459973	739242	8.05%	1.529%
67	15.064	4309	4320	4334	rBV	455520	764211	8.32%	1.580%
68	15.916	4573	4585	4607	rBV	127610	250141	2.72%	0.517%
69	16.060	4620	4630	4637	rBV	155277	245776	2.68%	0.508%
70	16.099	4637	4642	4656	rVB	80779	120383	1.31%	0.249%

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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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

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Title : SW846 8260

Sum of corrected areas: 48362448

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
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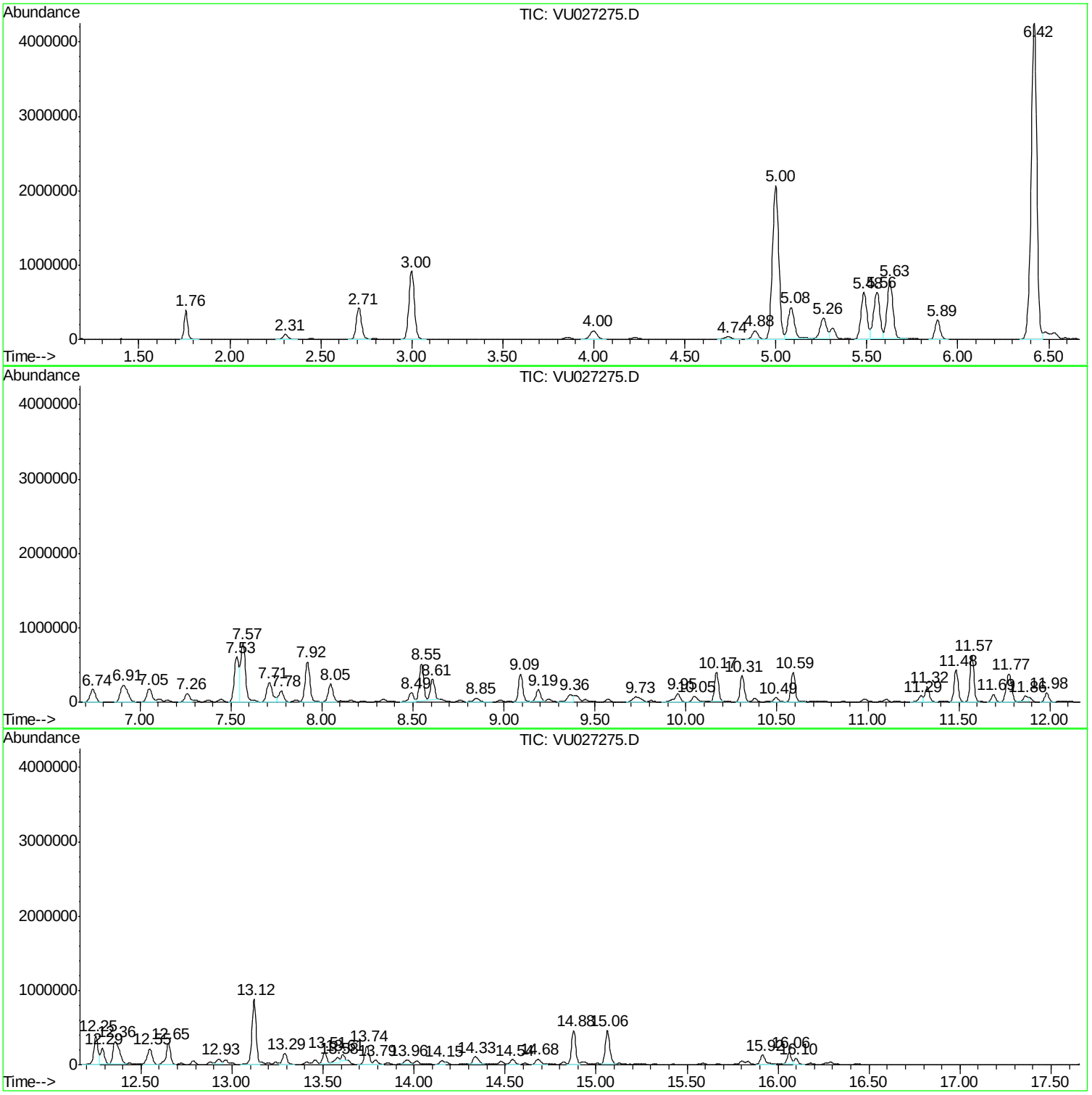
MSVOA_U
ClientSampleId :
S13-0246

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

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

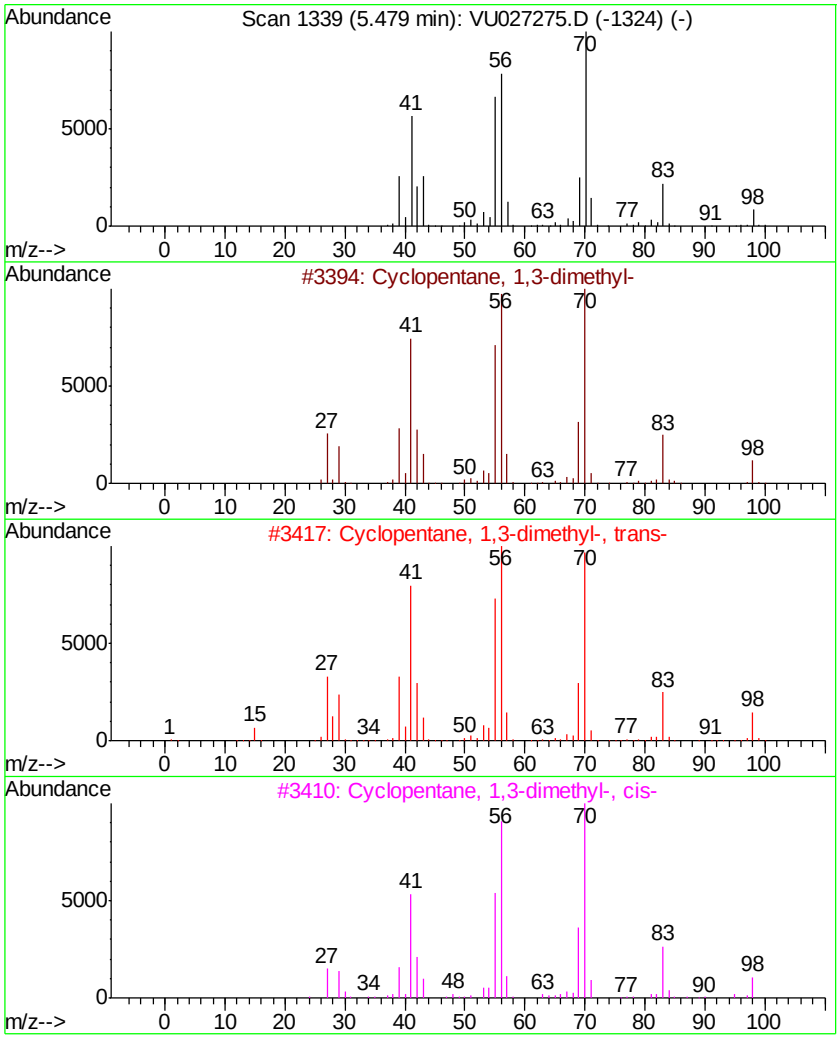
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

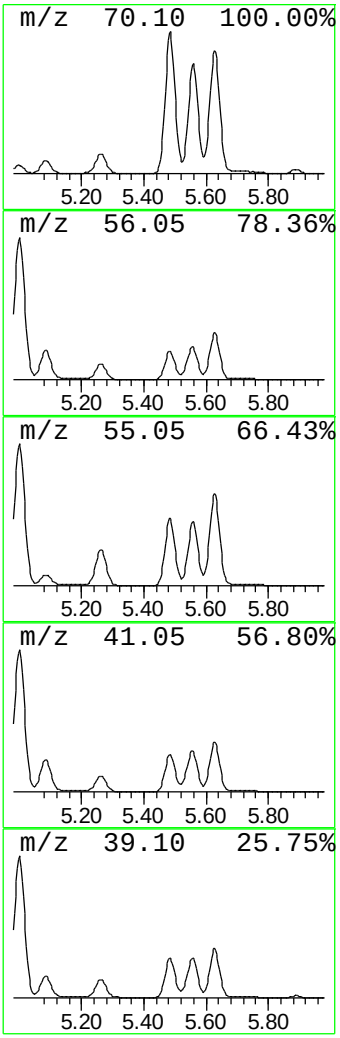
****

Peak Number 1 Cyclopentane, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.48	140.42 ug/l	1431180	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	95
2	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	87
5	Cycloheptane	98	C7H14	000291-64-5	50





Library Search Compound Report

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Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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

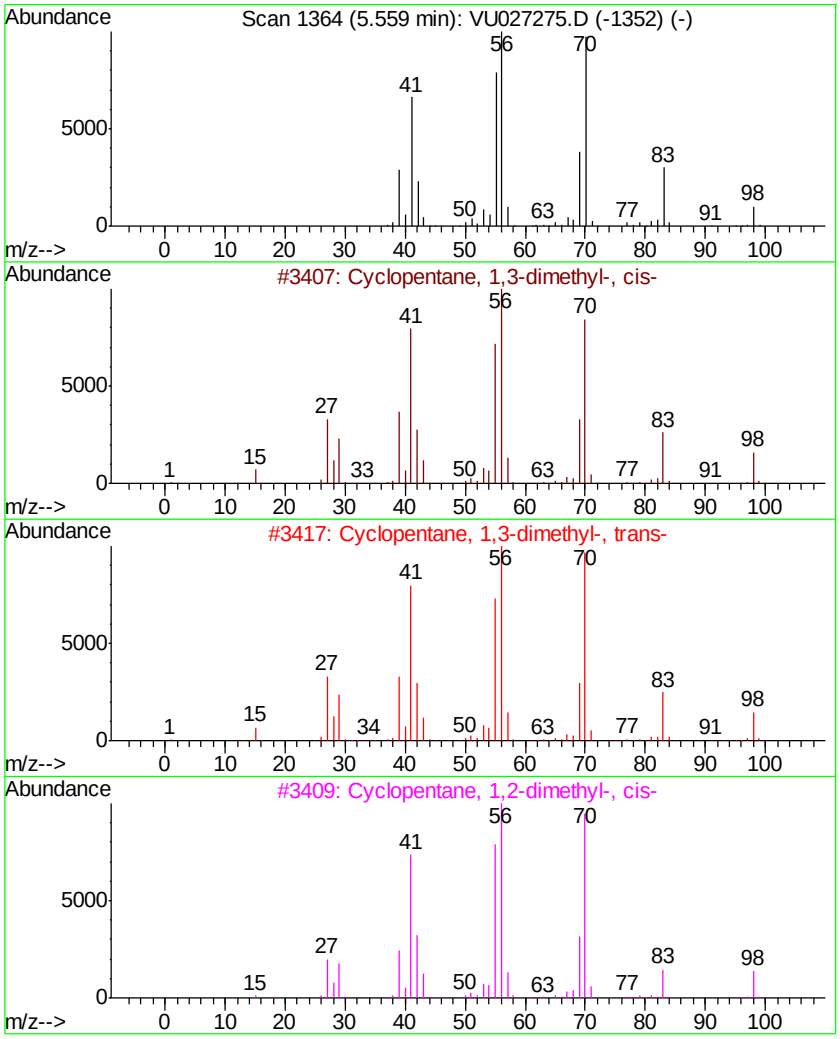
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

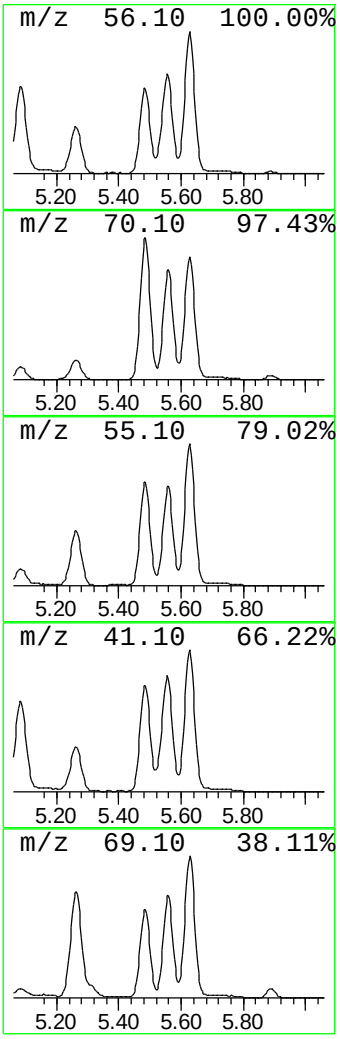
****

Peak Number 2 Cyclopentane, 1,3-dimethyl-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.56	147.24 ug/l	1500660	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	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	91
3	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	86
4	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	78
5	Isopropylcyclobutane	98	C7H14	000872-56-0	53





Library Search Compound Report

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Data File : VU027275.D
Acq On : 28 Sep 2018 16:38
Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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

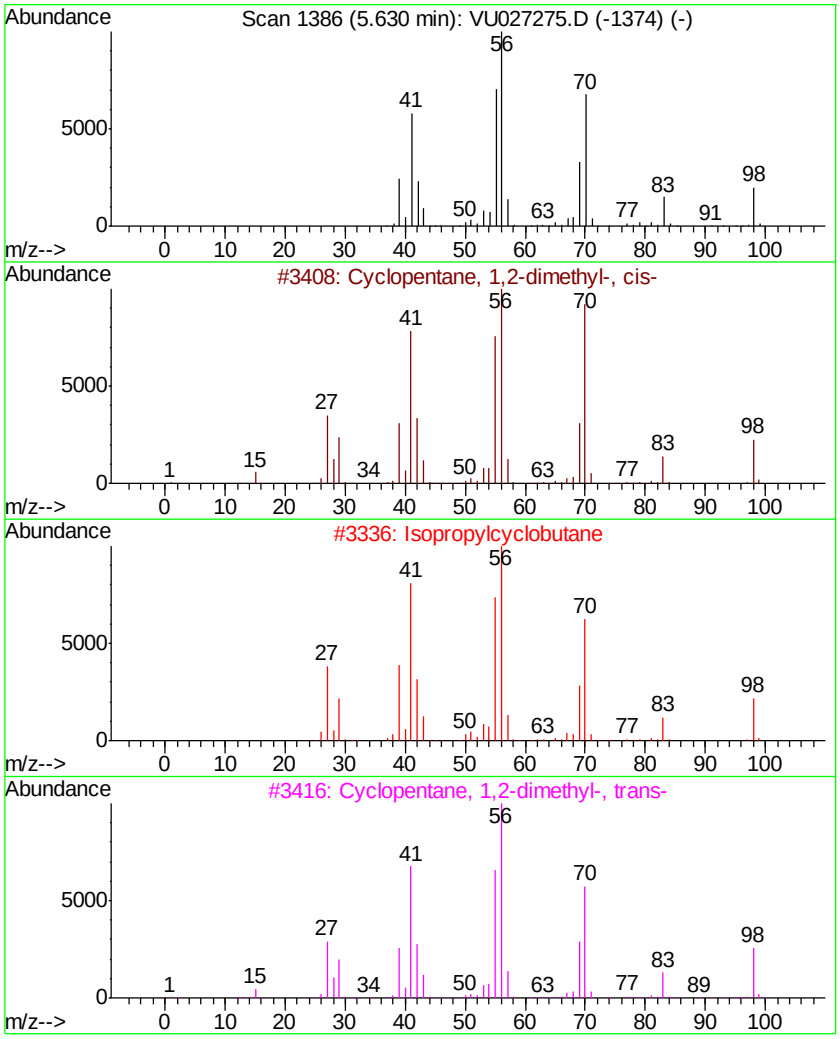
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

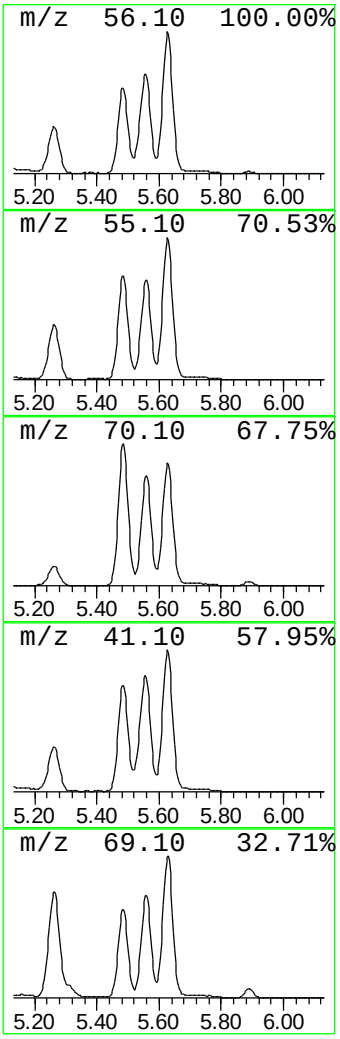
****

Peak Number 3 Cyclopentane, 1,2-dimethyl-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	170.17 ug/l	1734380	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	94
2	Isopropylcyclobutane	98	C7H14	000872-56-0	94
3	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
4	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	74
5	1-Heptene	98	C7H14	000592-76-7	72





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ClientSampleId : S13-0246

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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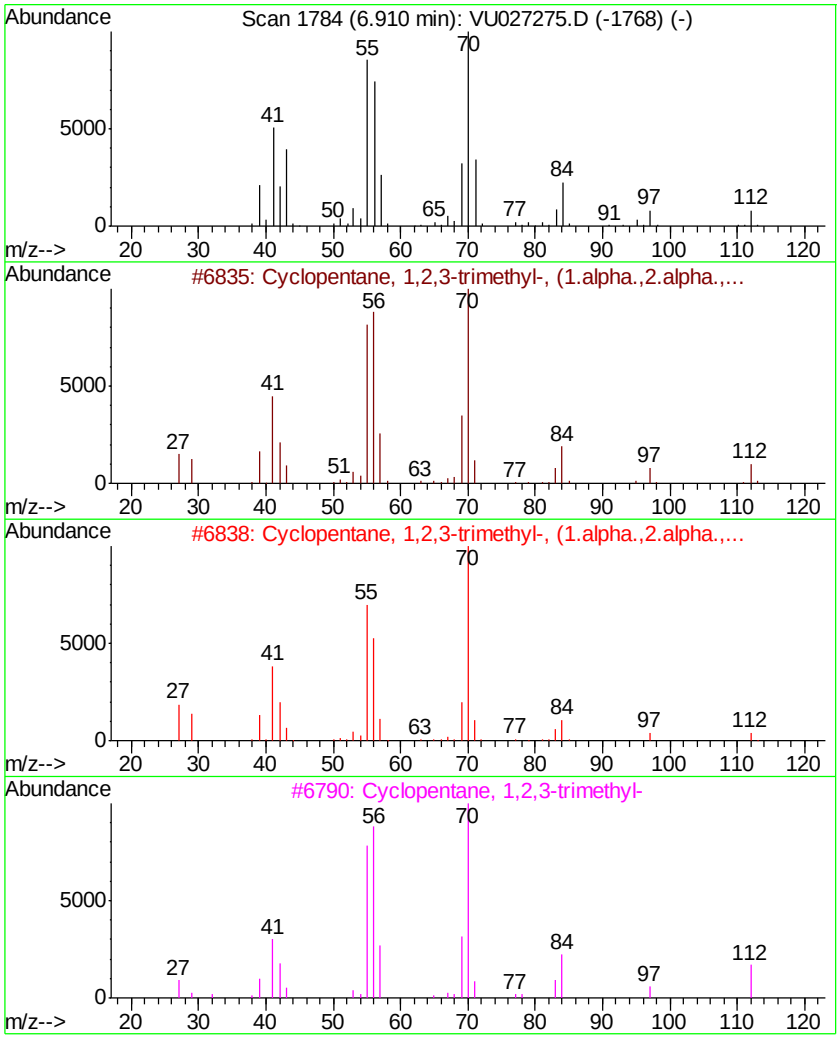
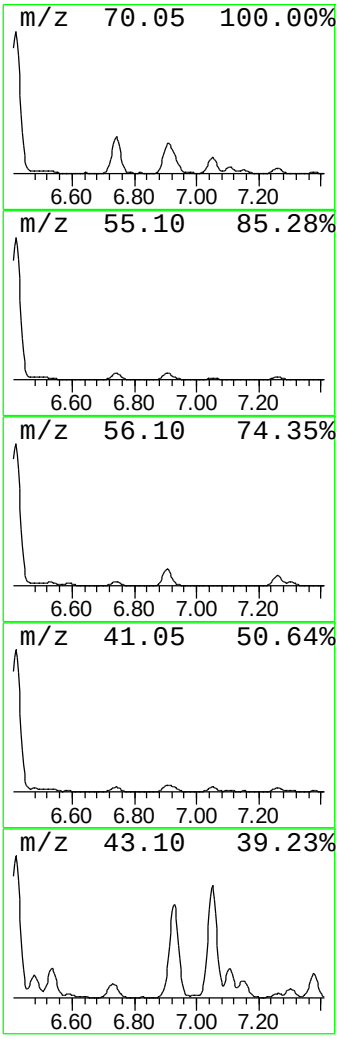
Peak Number 4 Cyclopentane, 1,2,3-trimeth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.91	61.45 ug/l	626296	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	015890-40-1	90
2	Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	87
3	Cyclopentane, 1,2,3-trimethyl-	112	C8H16	002815-57-8	80
4	cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	64
5	trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	64





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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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

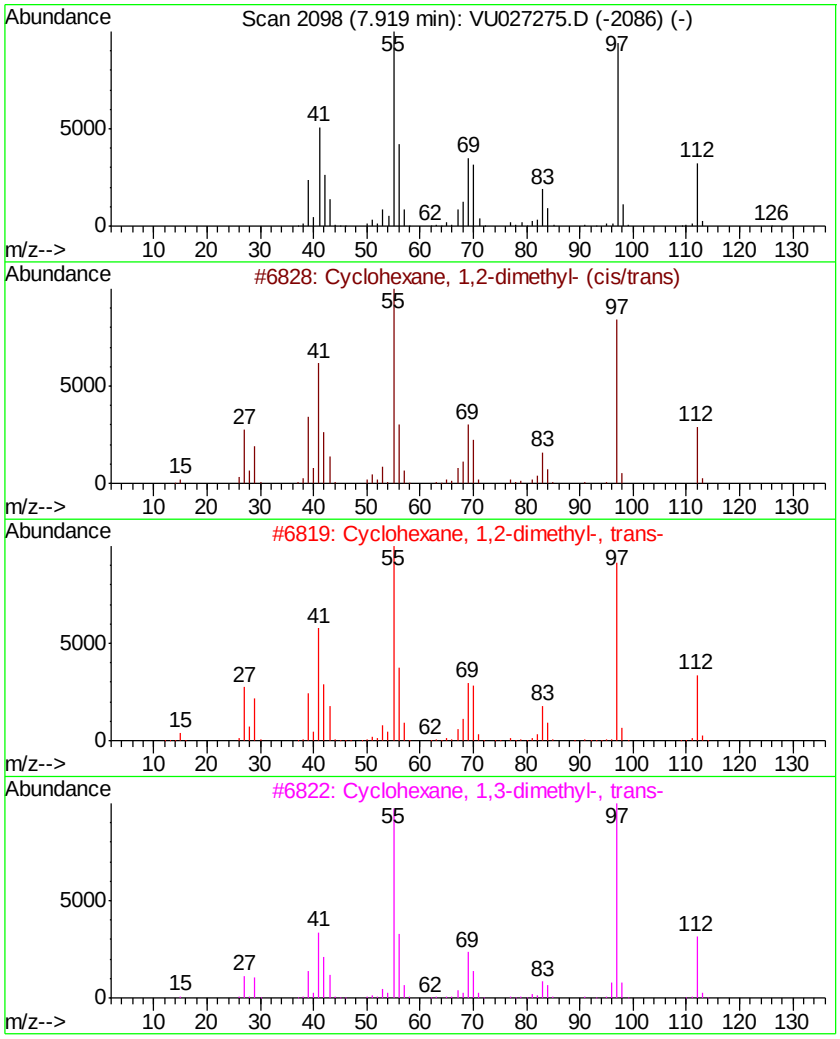
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

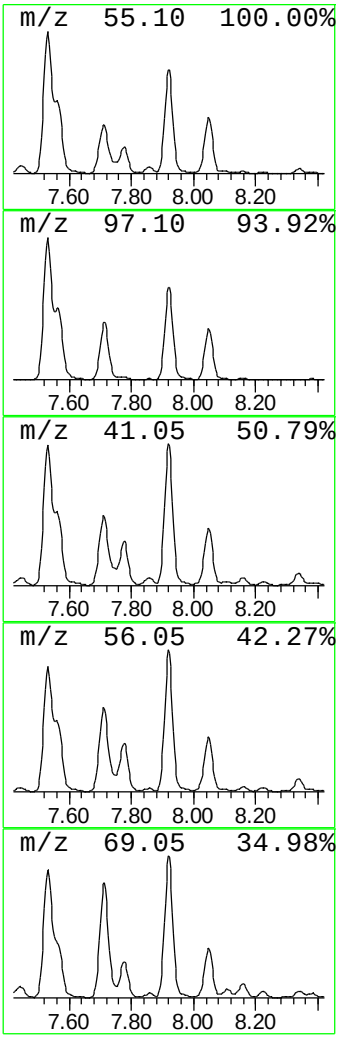
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.92	80.69 ug/l	988947	Chlorobenzene-d5	9.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	97
2	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	96
3	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
4	1H-Pyrazole, 4,5-dihydro-3,5,5-t...	112	C6H12N2	003975-85-7	76
5	Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	72





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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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

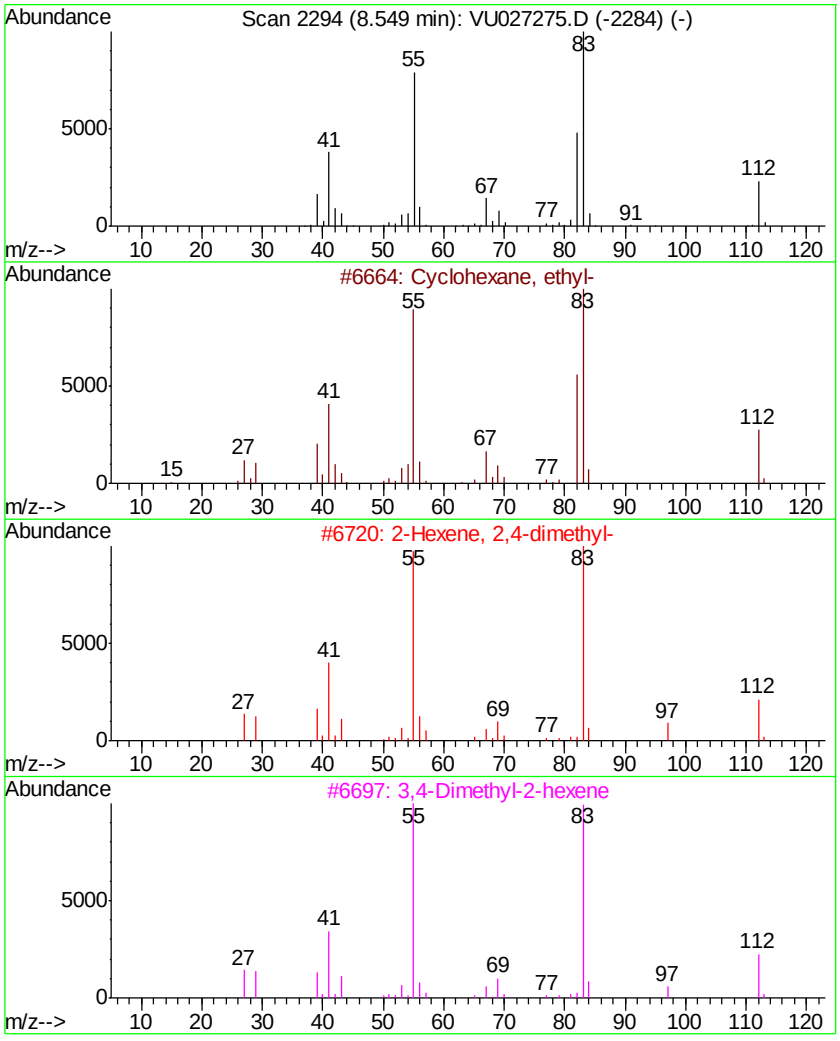
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

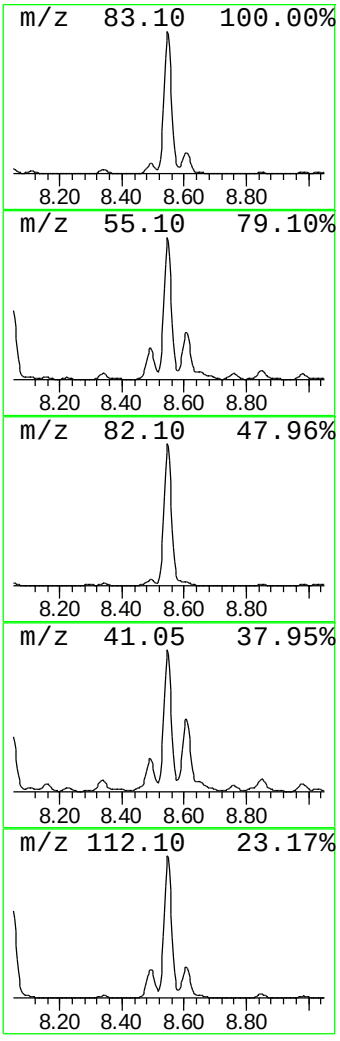
****

Peak Number 6 Cyclohexane, ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.55	65.19 ug/l	798946	Chlorobenzene-d5	9.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	93
2	2-Hexene, 2,4-dimethyl-	112	C8H16	014255-23-3	76
3	3,4-Dimethyl-2-hexene	112	C8H16	002213-37-8	76
4	2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	64
5	3-Heptene, 5-methyl-	112	C8H16	013172-91-3	64





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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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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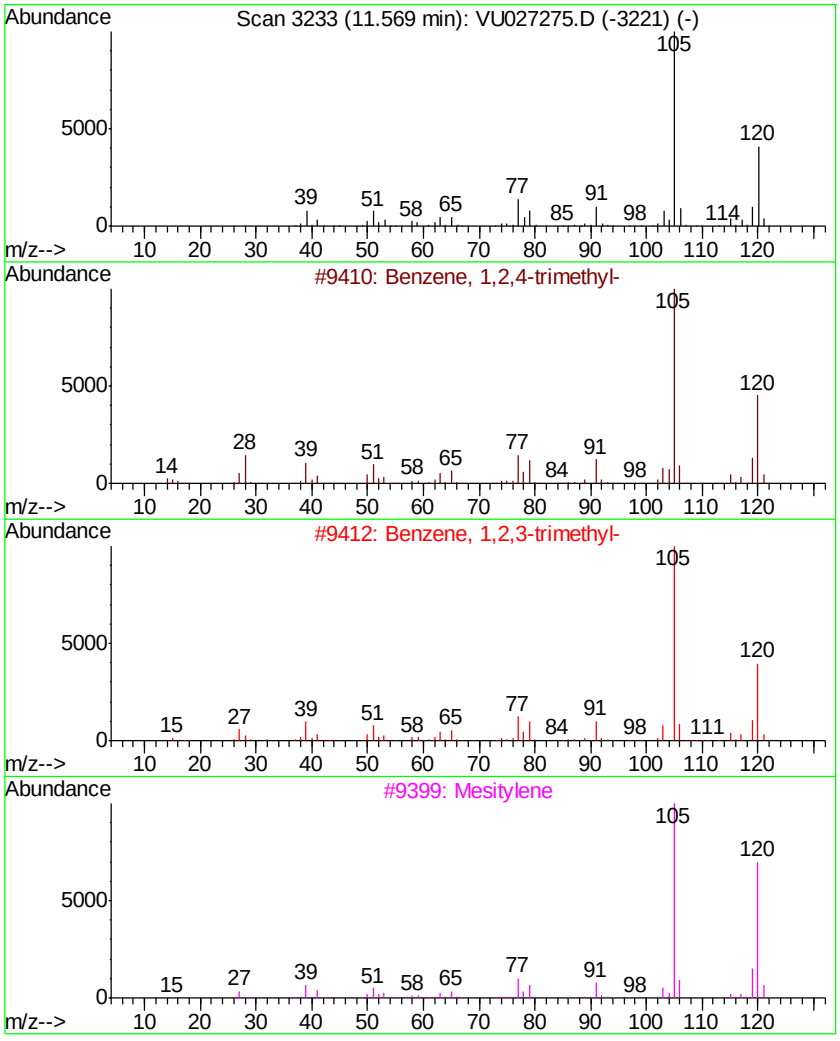
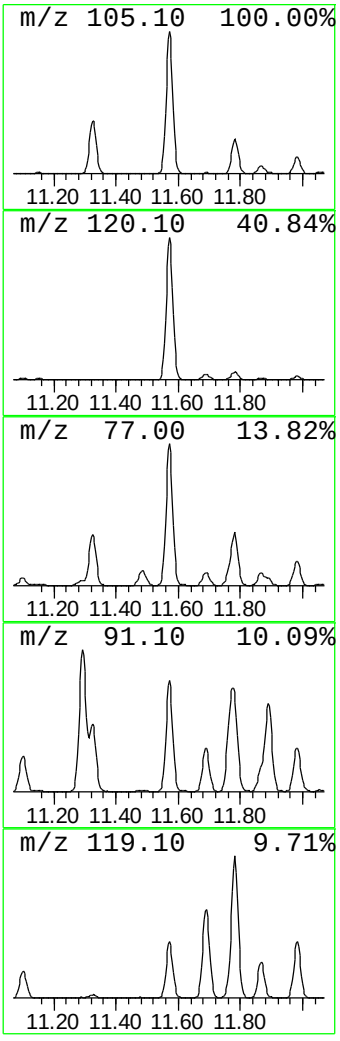
Peak Number 7 Benzene, 1,2,4-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	70.53 ug/l	939680	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	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027275.D
Acq On : 28 Sep 2018 16:38
Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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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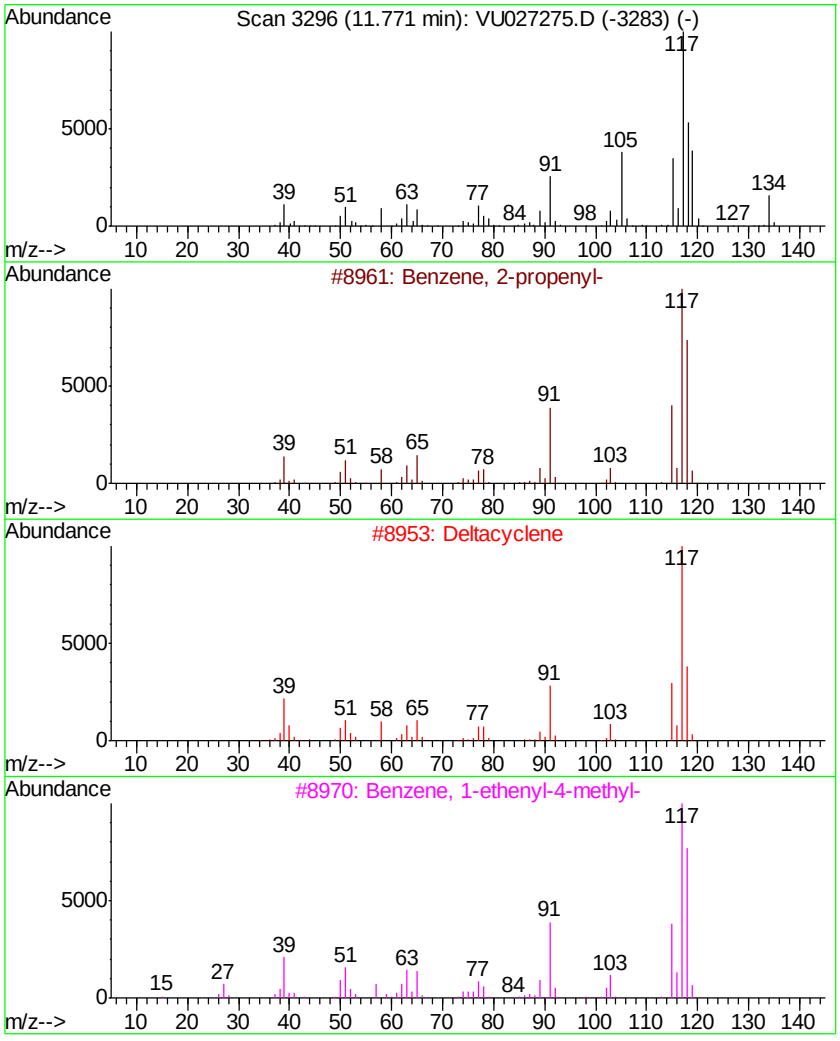
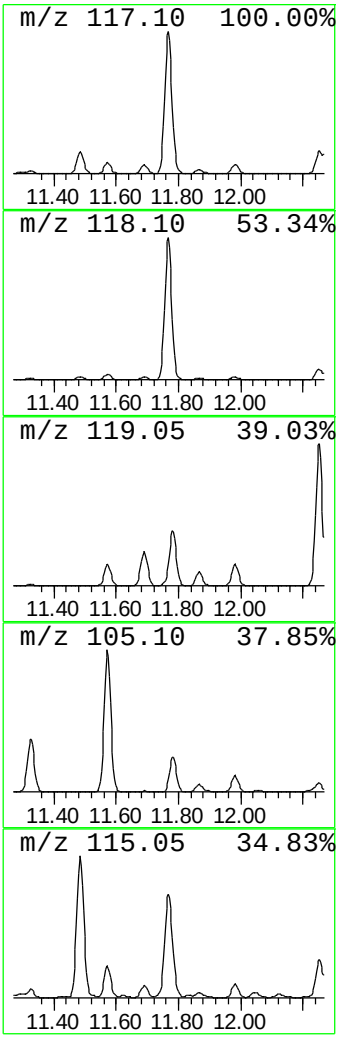
Peak Number 8 Benzene, 2-propenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	58.12 ug/l	774382	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	91
2	Deltacyclene	118	C9H10	007785-10-6	90
3	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	80
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027275.D
Acq On : 28 Sep 2018 16:38
Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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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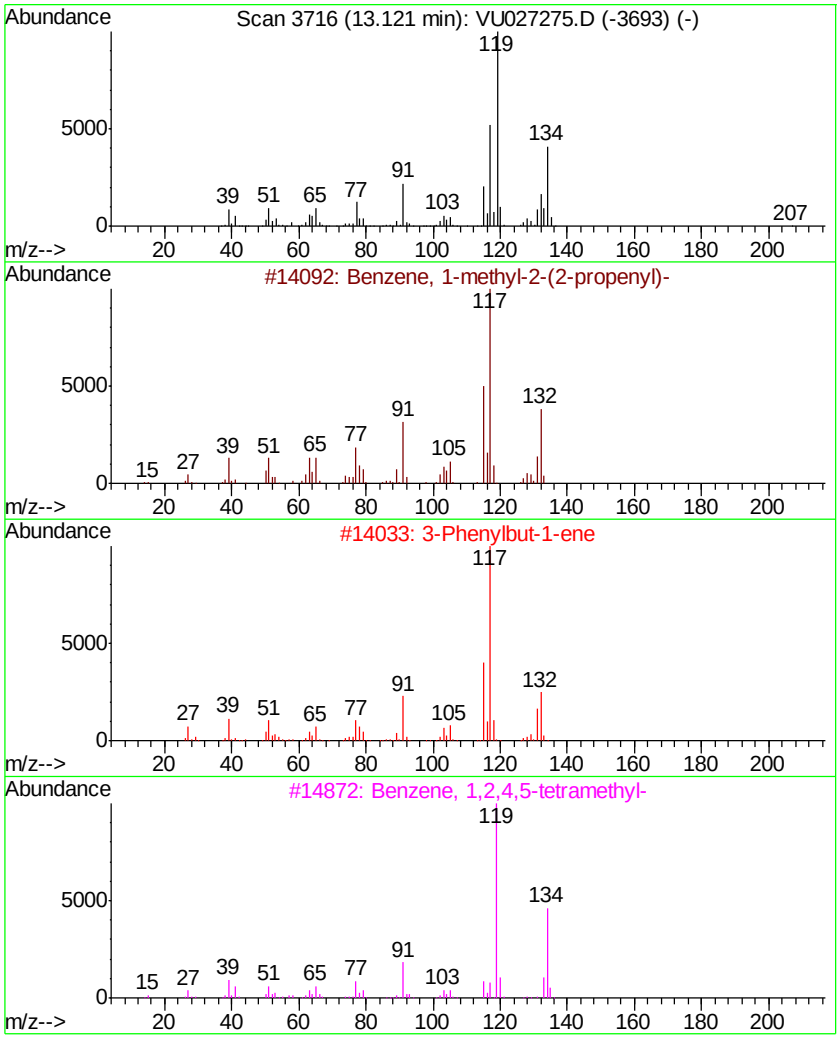
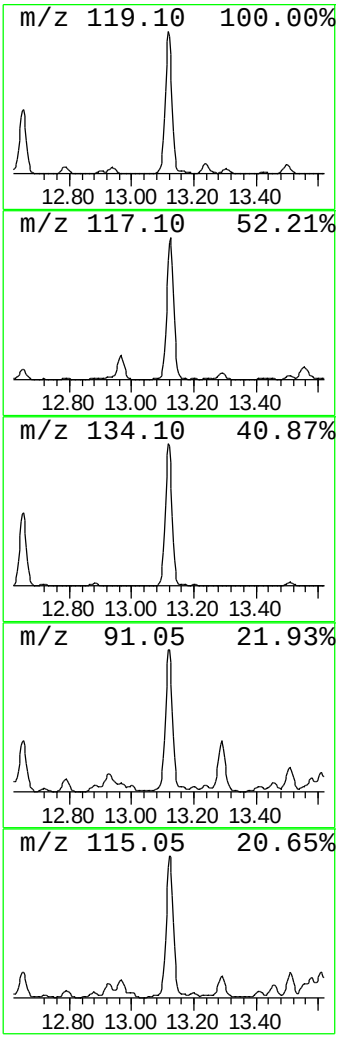
Peak Number 9 Benzene, 1-methyl-2-(2-prop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	106.68 ug/l	1421300	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4	o-Cymene	134	C10H14	000527-84-4	64
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027275.D
Acq On : 28 Sep 2018 16:38
Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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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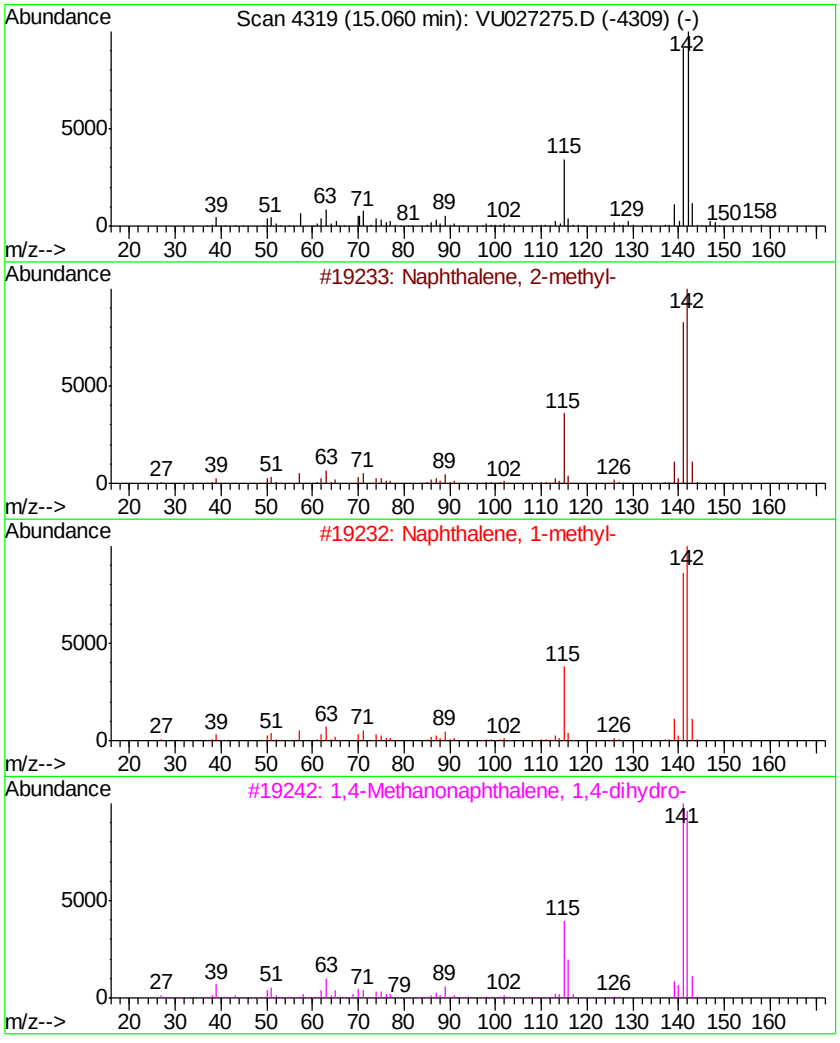
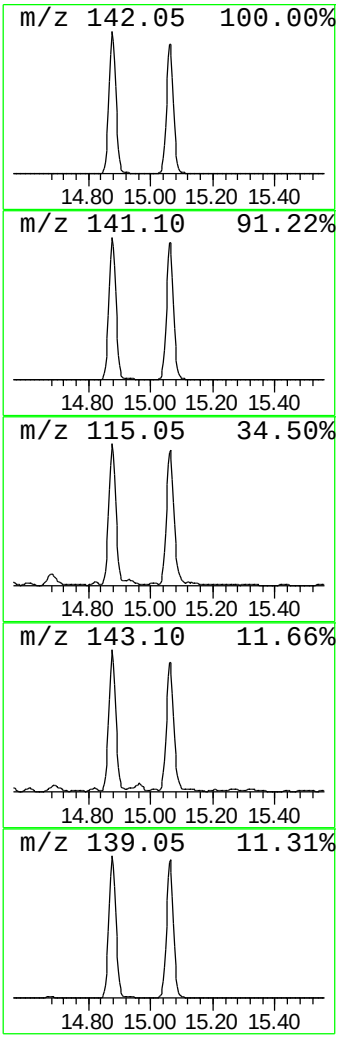
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	57.36 ug/l	764211	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027275.D
Acq On : 28 Sep 2018 16:38
Operator : MD/SY
Sample : J5041-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0246

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
Cyclopentane, 1,3...	5.48	140.4	ug/l	1431180	2	5.89	509614	50.0
Cyclopentane, 1,3...	5.56	147.2	ug/l	1500660	2	5.89	509614	50.0
Cyclopentane, 1,2...	5.63	170.2	ug/l	1734380	2	5.89	509614	50.0
Cyclopentane, 1,2...	6.91	61.5	ug/l	626296	2	5.89	509614	50.0
Cyclohexane, 1,2-...	7.92	80.7	ug/l	988947	3	9.09	612814	50.0
Cyclohexane, ethyl-	8.55	65.2	ug/l	798946	3	9.09	612814	50.0
Benzene, 1,2,4-tr...	11.57	70.5	ug/l	939680	4	11.48	666164	50.0
Benzene, 2-propenyl-	11.77	58.1	ug/l	774382	4	11.48	666164	50.0
Benzene, 1-methyl...	13.12	106.7	ug/l	1421300	4	11.48	666164	50.0
Naphthalene, 1-me...	15.06	57.4	ug/l	764211	4	11.48	666164	50.0