LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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	1.758	172	182	199	rVB	281570	367985	3.40%	0.421%
2	2.305	339	352	369	rVB	87436	169597	1.57%	0.194%
3	2.710	464	478	482	rBV	193315	364428	3.37%	0.417%
4	2.746	483	489	496	rVV	317711	590378	5.45%	0.676%
5	2.794	496	504	523	rVB	553777	1042927	9.63%	1.194%
6	3.000	550	568	593	rBV	650053	1382786	12.77%	1.583%
7	4.096	890	909	932	rVV	1599206	4268654	39.42%	4.886%
8	4.736	1087	1108	1128	rBV2	45962	127664	1.18%	0.146%
9	4.887	1138	1155	1170	rBV	108929	248614	2.30%	0.285%
10	5.000	1170	1190	1206	rBV2	1986206	4849184	44.78%	5.551%
11	5.083	1206	1216	1236	rVB2	184909	446240	4.12%	0.511%
12	5.221	1245	1259	1267	rBV	193134	465153	4.30%	0.532%
13	5.311	1280	1287	1300	rVB	134172	263009	2.43%	0.301%
14	5.392	1301	1312	1326	rVB	94374	191974	1.77%	0.220%
15	5.485	1326	1341	1352	rBV2	310187	697397	6.44%	0.798%
16	5.556	1353	1363	1374	rVV	261790	577561	5.33%	0.661%
17	5.626	1374	1385	1413	rVB	449135	992684	9.17%	1.136%
18	5.887	1452	1466	1482	rBV	242255	472942	4.37%	0.541%
19	6.417	1606	1631	1646	rBV	2875911	6132069	56.63%	7.020%
20	6.643	1688	1701	1715	rVV	243566	456487	4.22%	0.523%
21	6.739	1716	1731	1747	rVB2	87192	191499	1.77%	0.219%
22	6.906	1770	1783	1805	rVB2	91134	199890	1.85%	0.229%
23	7.530	1965	1977	1981	rVV2	363722	600984	5.55%	0.688%
24	7.565	1982	1988	2019	rVB	629915	1129346	10.43%	1.293%
25	7.710	2019	2033	2046	rBV4	151943	382987	3.54%	0.438%
26	7.774	2048	2053	2066	rVB2	99864	171913	1.59%	0.197%
27	7.919	2086	2098	2122	rVB	302832	553132	5.11%	0.633%
28	8.048	2123	2138	2149	rBV	148971	270115	2.49%	0.309%
29	8.491	2251	2276	2283	rBV2	123723	273546	2.53%	0.313%
30	8.546	2283	2293	2303	rVV	377776	628520	5.80%	0.719%
31	8.607	2303	2312	2322	rVB	197409	337410	3.12%	0.386%
32	9.089	2453	2462	2479	rVB	346629	554391	5.12%	0.635%
33	9.189	2481	2493	2501	rBV	143571	241869	2.23%	0.277%
34	9.250	2501	2512	2533	rVV	3655228	5881259	54.31%	6.732%

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ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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	9.372	2536	2550	2567	rVB7	127029	392493	3.62%	0.449%
36	9.491	2580	2587	2601	rVB	95685	167139	1.54%	0.191%
37	9.723	2640	2659	2677	rBV6	68778	225243	2.08%	0.258%
38	9.954	2711	2731	2743	rBV4	142855	291658	2.69%	0.334%
39	10.044	2749	2759	2769	rBV4	110240	206139	1.90%	0.236%
40	10.166	2786	2797	2809	rBV	862788	1343035	12.40%	1.537%
41	10.308	2831	2841	2854	rBV	338711	560700	5.18%	0.642%
42	10.588	2917	2928	2941	rBV	946783	1464755	13.53%	1.677%
43	10.707	2955	2965	2972	rBV	97759	160906	1.49%	0.184%
44	10.774	2972	2986	3007	rVB	1194805	1887225	17.43%	2.160%
45	10.974	3026	3048	3069	rBV	817019	1403160	12.96%	1.606%
46	11.150	3091	3103	3117	rVB	7213038	10828370	100.00%	12.396%
47	11.295	3128	3148	3151	rBV	91073	171046	1.58%	0.196%
48	11.324	3151	3157	3173	rVB	207093	349774	3.23%	0.400%
49	11.482	3195	3206	3218	rBV2	529931	840119	7.76%	0.962%
50	11.572	3222	3234	3258	rVB	4073975	6261251	57.82%	7.167%
51	11.691	3259	3271	3283	rBV	87786	150303	1.39%	0.172%
52	11.768	3283	3295	3312	rBV2	1155095	2108728	19.47%	2.414%
53	11.874	3314	3328	3347	rVB5	503279	1148935	10.61%	1.315%
54	11.983	3351	3362	3374	rBV	170649	276503	2.55%	0.317%
55	12.057	3374	3385	3401	rVB	426642	653421	6.03%	0.748%
56	12.147	3402	3413	3417	rBV	476819	689214	6.36%	0.789%
57	12.176	3418	3422	3432	rVB	517501	690696	6.38%	0.791%
58	12.253	3434	3446	3452	rVV	701794	1051826	9.71%	1.204%
59	12.285	3453	3456	3468	rVV	206800	267178	2.47%	0.306%
60	12.356	3468	3478	3499	rVV3	459287	1079525	9.97%	1.236%
61	12.494	3510	3521	3528	rBV3	75824	137950	1.27%	0.158%
62	12.552	3528	3539	3550	rVB3	410801	717516	6.63%	0.821%
63	12.649	3550	3569	3577	rBV	487547	845417	7.81%	0.968%
64	12.703	3577	3586	3602	rVB	743826	1145166	10.58%	1.311%
65	12.790	3602	3613	3619	rBV2	113755	186795	1.73%	0.214%
66	12.964	3661	3667	3676	rVB	234328	328690	3.04%	0.376%
67	13.086	3695	3705	3706	rBV3	91049	113431	1.05%	0.130%
68	13.121	3706	3716	3731	rVB2	1537592	2485172	22.95%	2.845%
69	13.289	3758	3768	3795	rVB	773099	1262163	11.66%	1.445%
70	13.408	3796	3805	3812	rBV2	75181	123342	1.14%	0.141%
71	13.456	3813	3820	3827	rBV	83947	117038	1.08%	0.134%

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Instrument : MSVOA_U
ClientSampleId : S13-0251

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	13.510	3827	3837	3844	rBV4	217981	365841	3.38%	0.419%
73	13.742	3896	3909	3919	rBV	2145449	3374558	31.16%	3.863%
74	13.787	3919	3923	3935	rVB	91228	128316	1.18%	0.147%
75	13.854	3935	3944	3959	rBV2	98907	186228	1.72%	0.213%
76	13.954	3960	3975	3986	rBV4	115763	250531	2.31%	0.287%
77	14.153	4026	4037	4042	rBV	70754	124550	1.15%	0.143%
78	14.350	4083	4098	4118	rVB4	242682	529374	4.89%	0.606%
79	14.539	4148	4157	4170	rVB6	73522	140620	1.30%	0.161%
80	14.678	4190	4200	4218	rBV2	183275	335485	3.10%	0.384%
81	14.877	4250	4262	4273	rBV	1305670	2033046	18.78%	2.327%
82	15.060	4308	4319	4334	rBV	881143	1428403	13.19%	1.635%
83	15.607	4472	4489	4500	rVB3	48356	118706	1.10%	0.136%
84	15.915	4573	4585	4610	rBV2	138420	274389	2.53%	0.314%
85	16.060	4620	4630	4636	rBV	153483	242655	2.24%	0.278%
86	16.099	4637	4642	4658	rVB	112672	167856	1.55%	0.192%

Sum of corrected areas: 87357174

LSC Report - Integrated Chromatogram

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Sample : J5041-03
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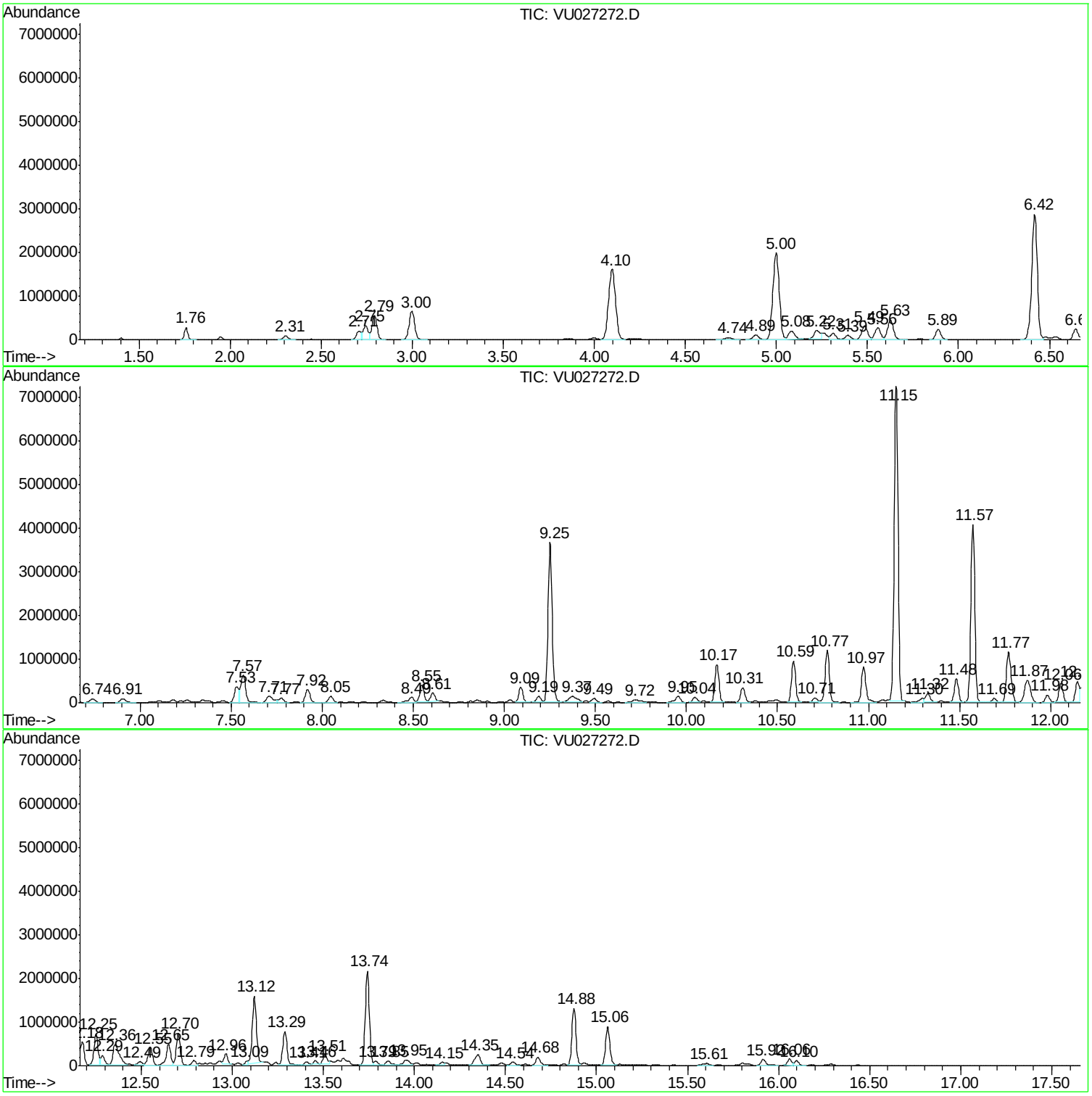
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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 : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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

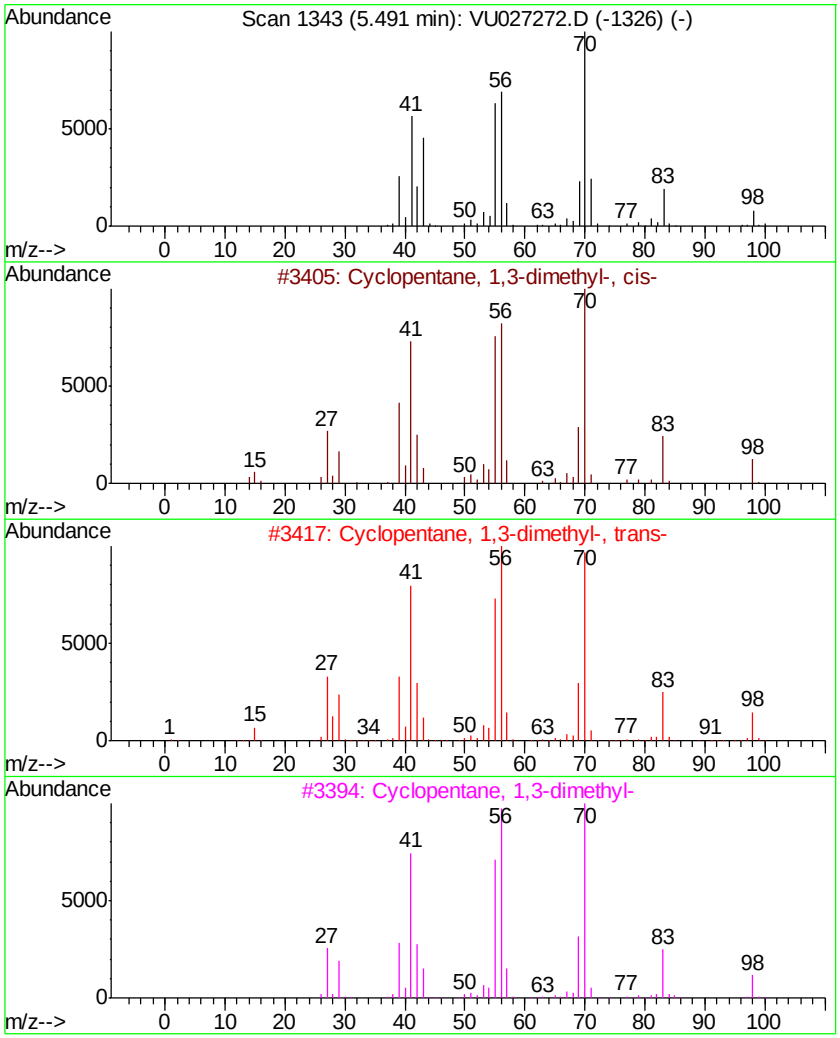
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

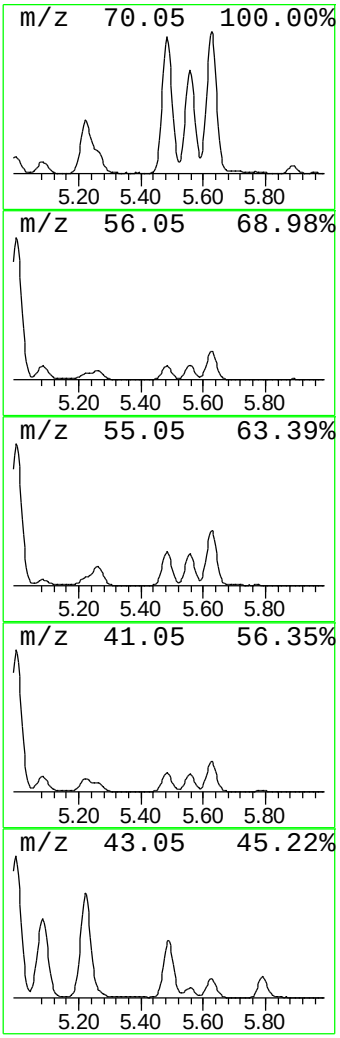
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.49	73.73 ug/l	697397	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	92
2	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	87
3	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	70
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	60
5	Octane, 3-methyl-6-methylene-	140	C10H20	074630-07-2	53





Library Search Compound Report

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Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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

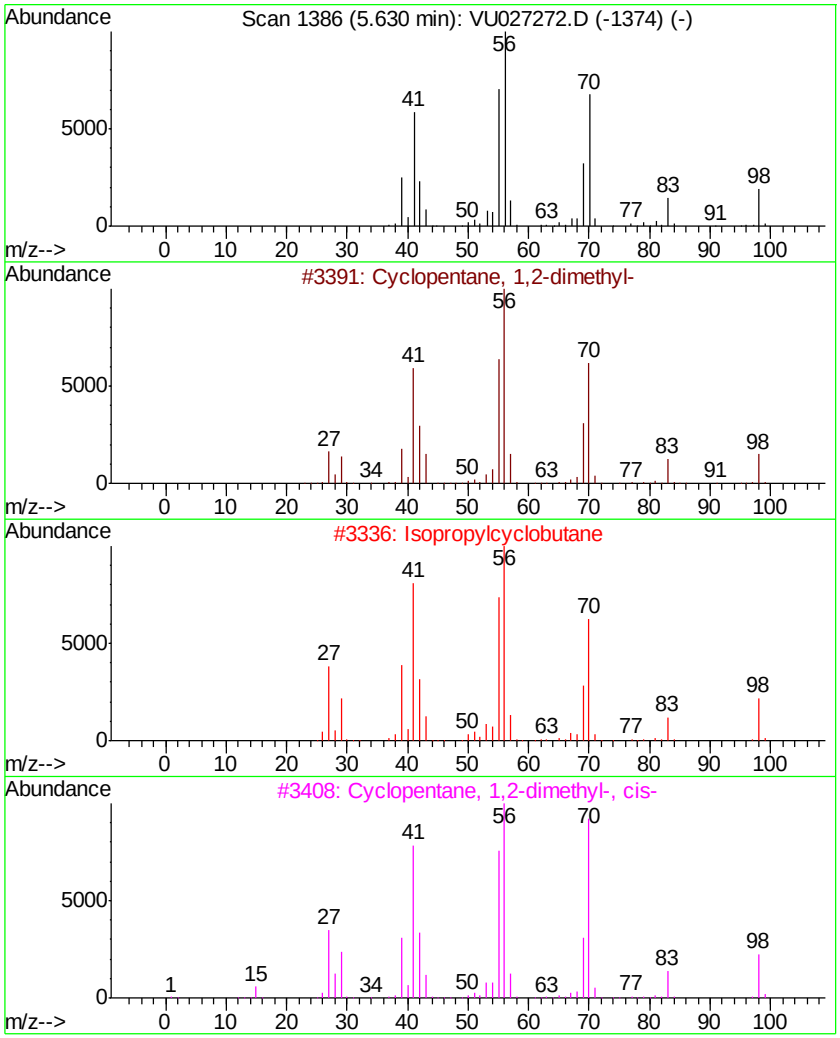
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

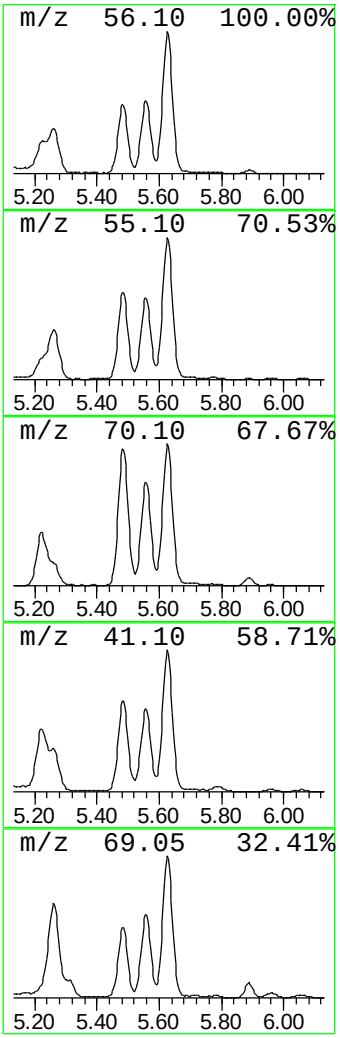
****

Peak Number 2 Cyclopentane, 1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	104.95 ug/l	992684	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	94
2	Isopropylcyclobutane	98	C7H14	000872-56-0	94
3	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	91
4	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
5	2,2-Dimethylglutaric anhydride	142	C7H10O3	002938-48-9	72





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ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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

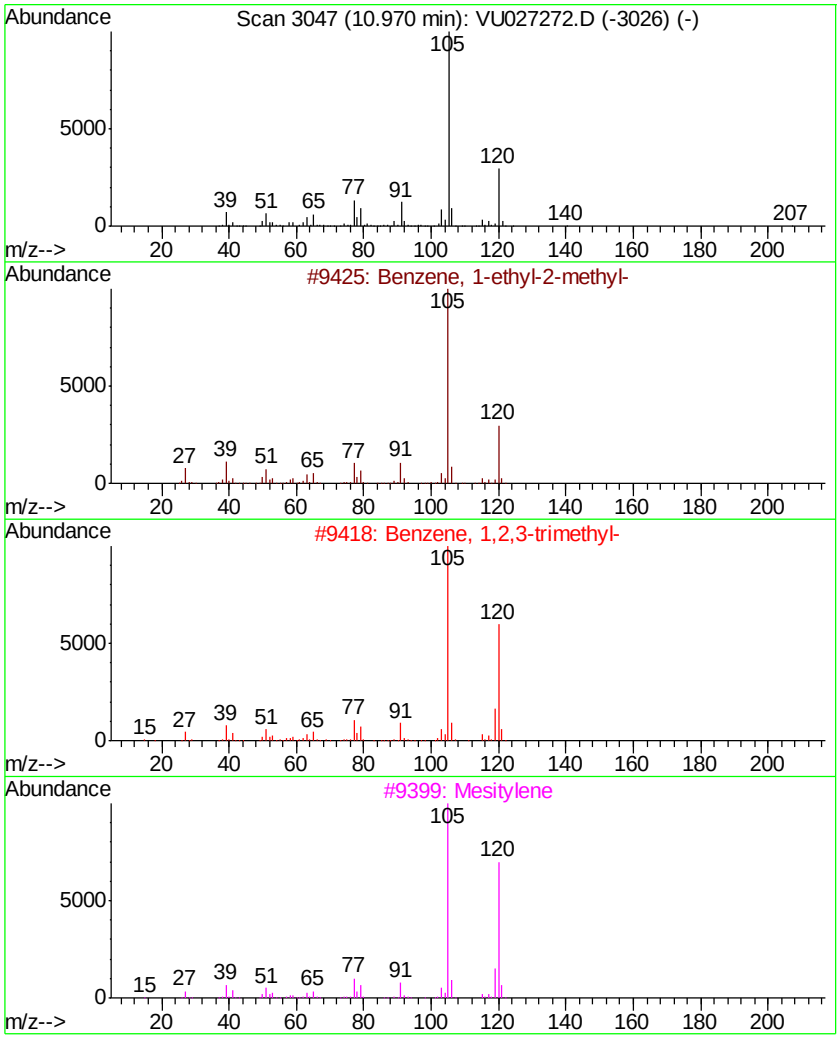
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

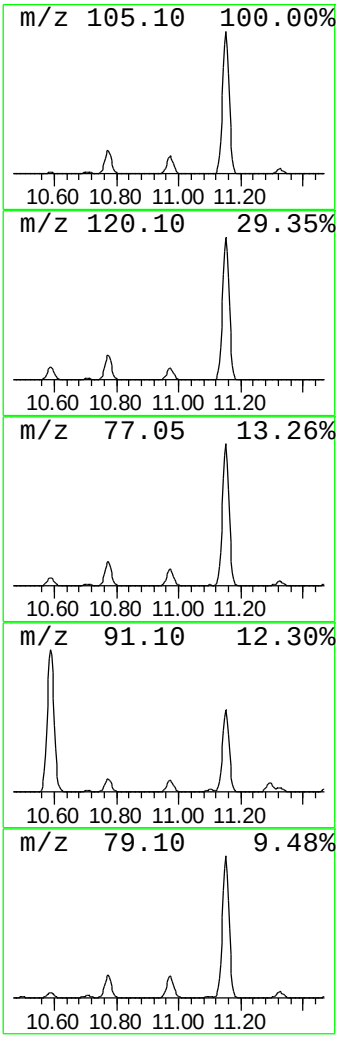
****

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	83.51 ug/l	1403160	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





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ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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

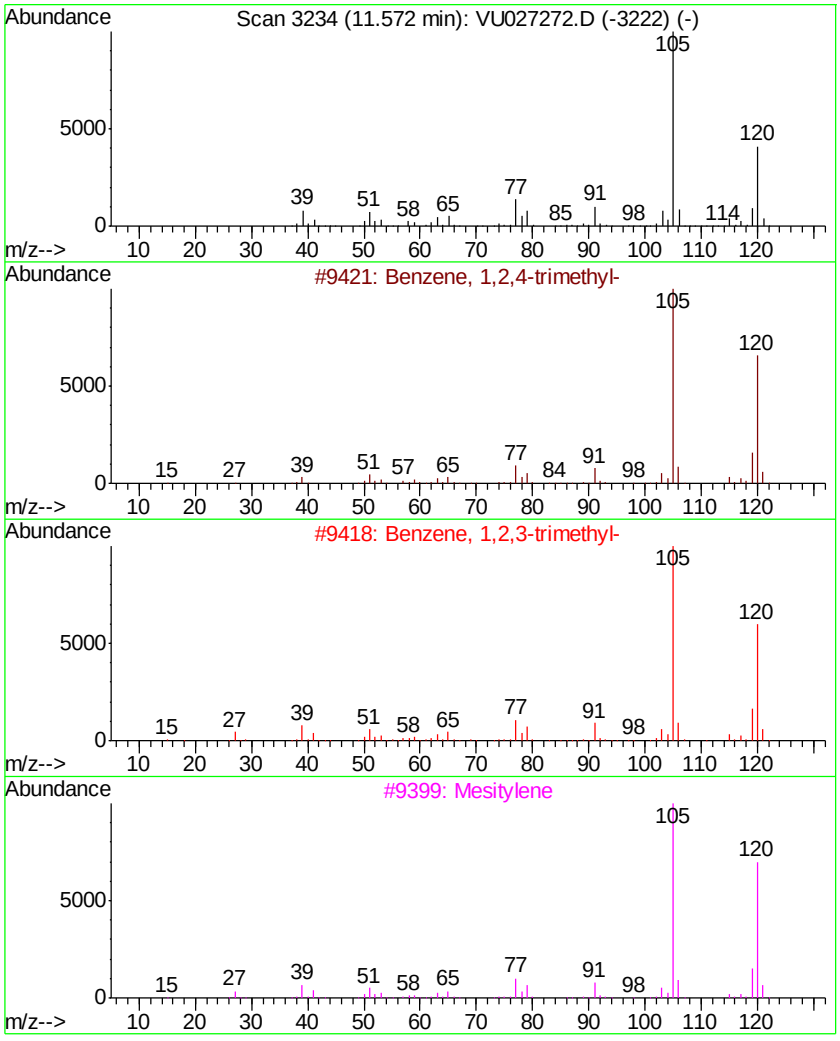
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

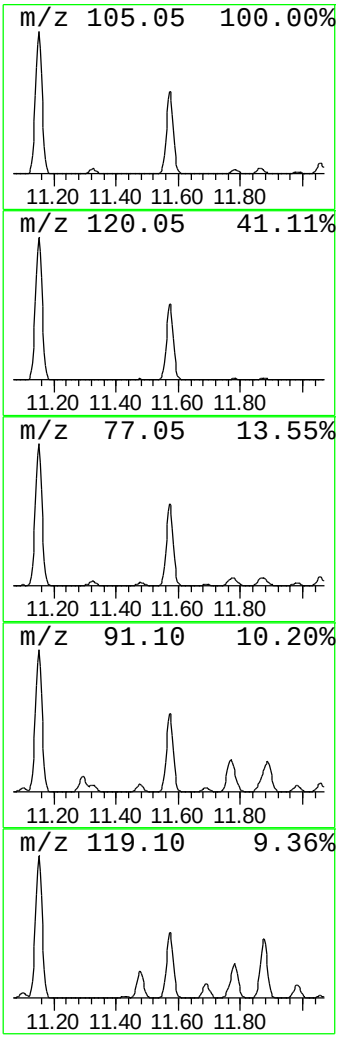
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	372.64 ug/l	6261250	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





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ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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

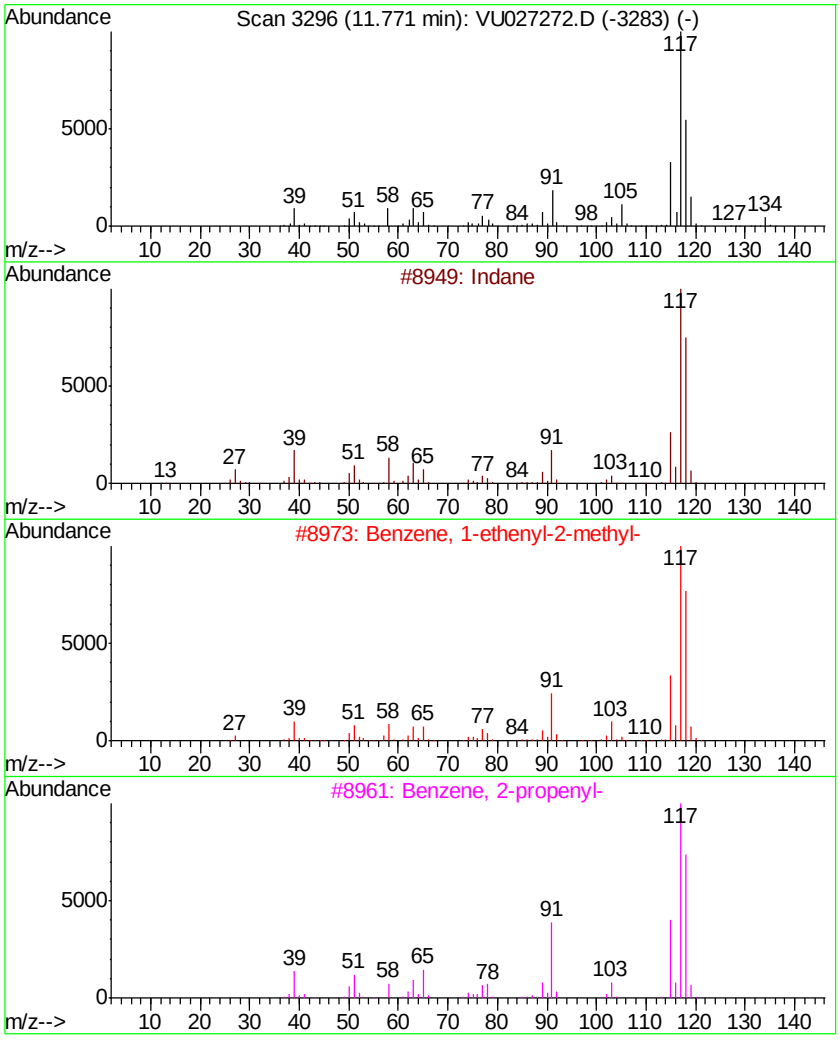
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

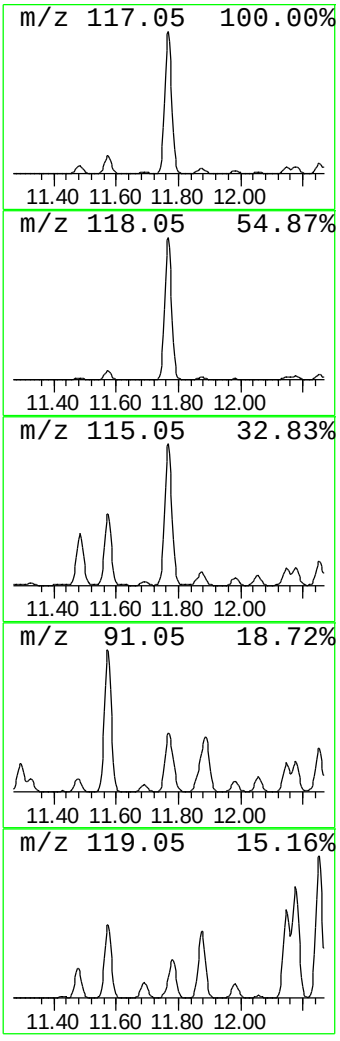
****

Peak Number 5 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	125.50 ug/l	2108730	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	89
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	87
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
5	Deltacyclene	118	C9H10	007785-10-6	58





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ClientSampleId : S13-0251

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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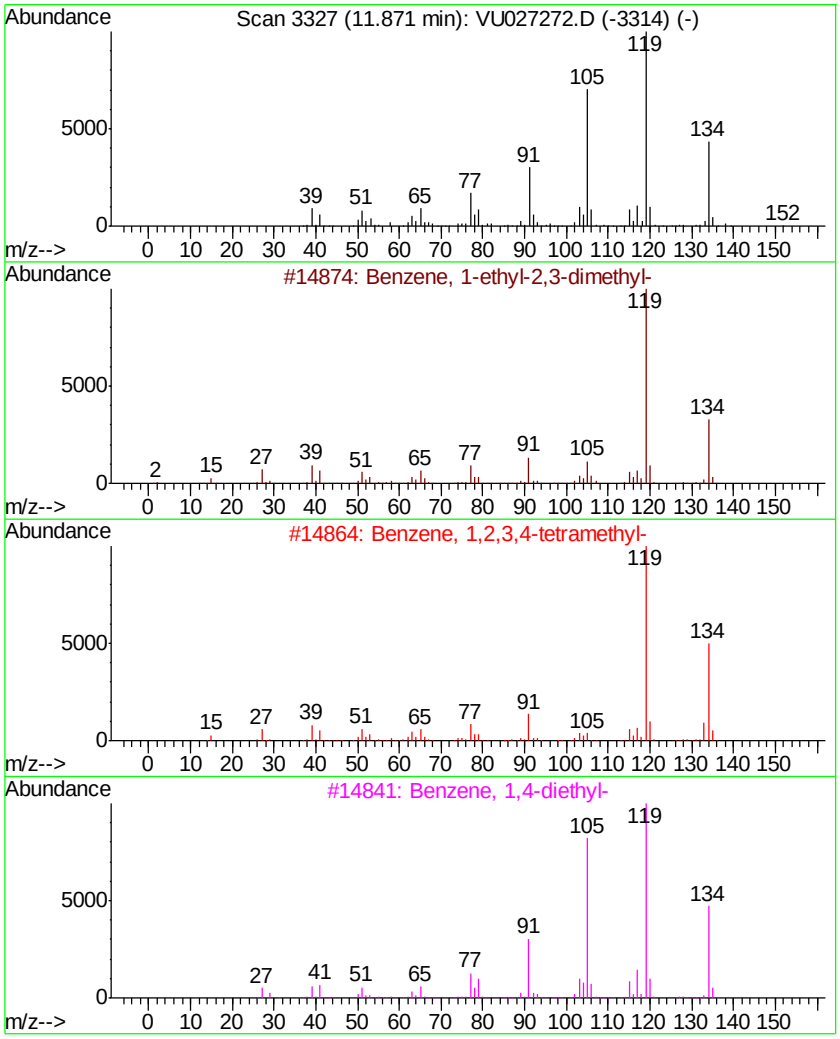
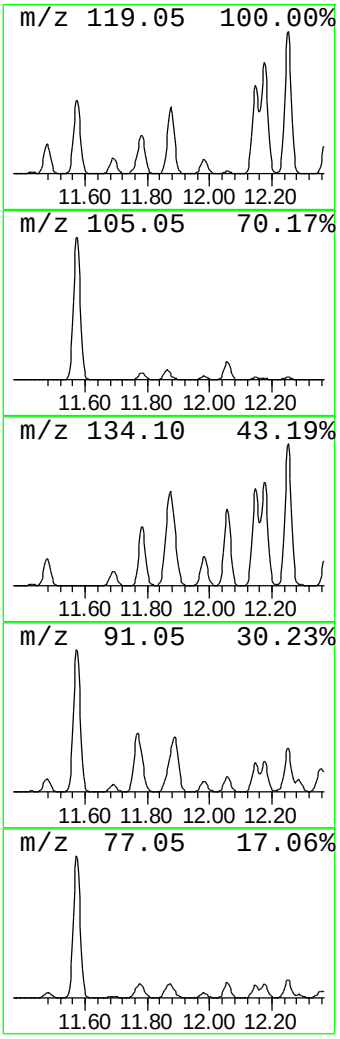
Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	68.38 ug/l	1148940	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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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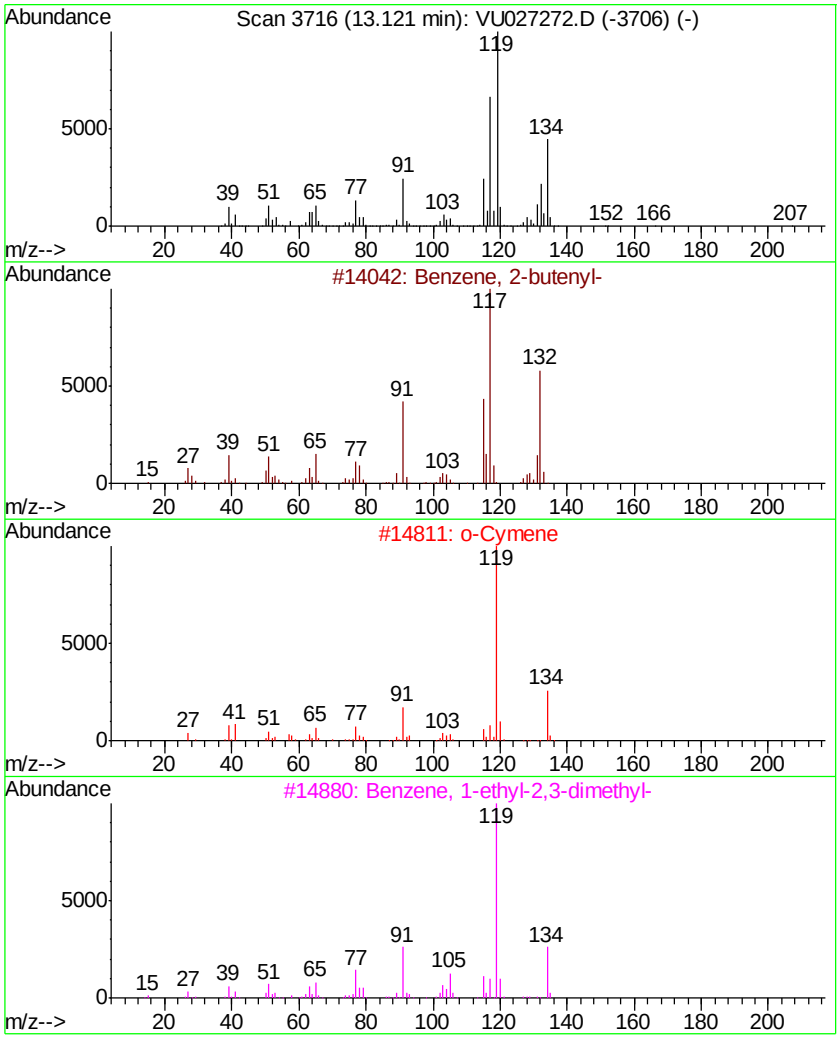
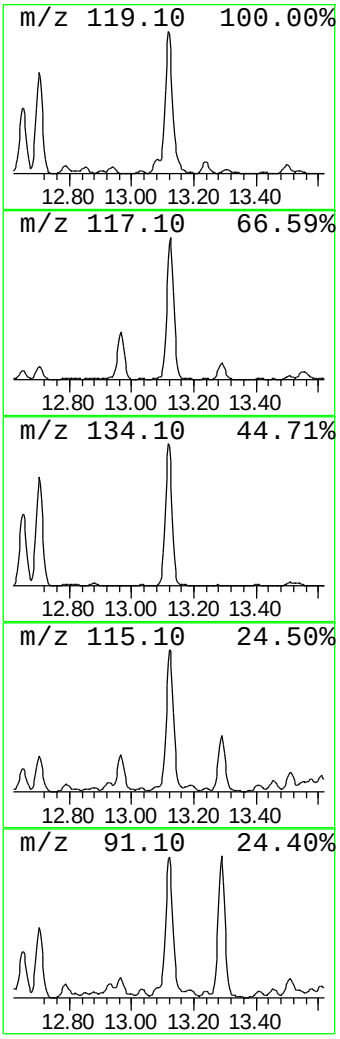
Peak Number 7 Benzene, 2-butenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	147.91 ug/l	2485170	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
2	o-Cymene	134	C10H14	000527-84-4	60
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	60
4	p-Cymene	134	C10H14	000099-87-6	60
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	60





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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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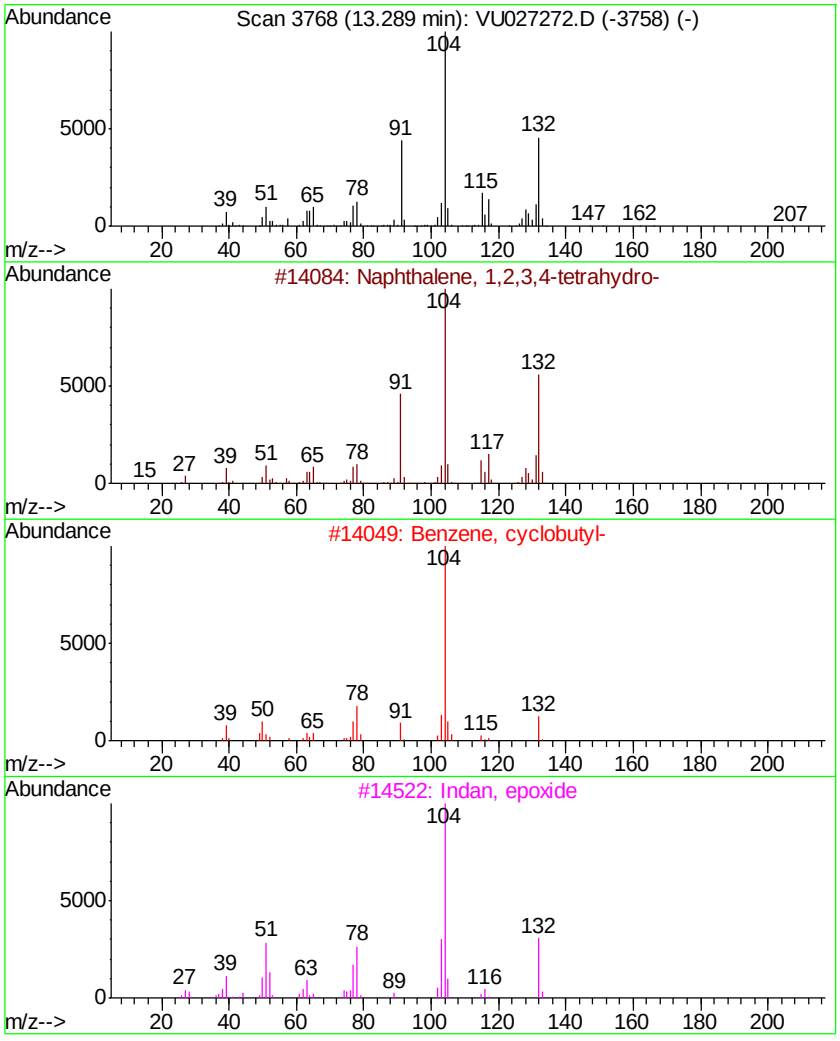
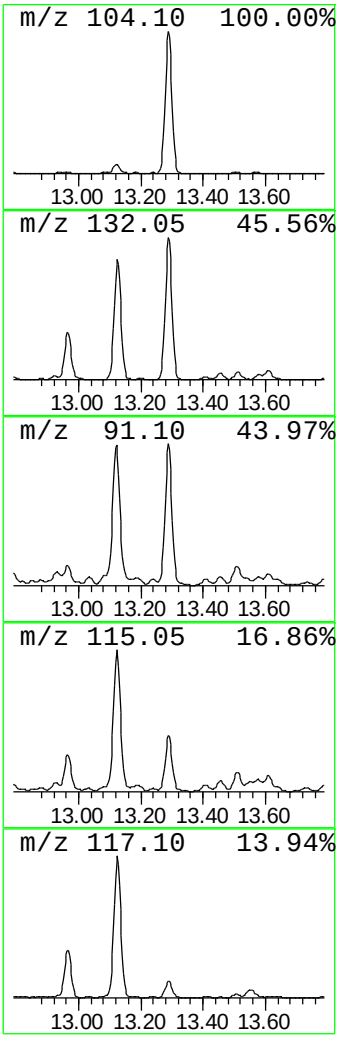
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	75.12 ug/l	1262160	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2	Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
3	Indan, epoxide	132	C9H8O	000768-22-9	58
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
5	N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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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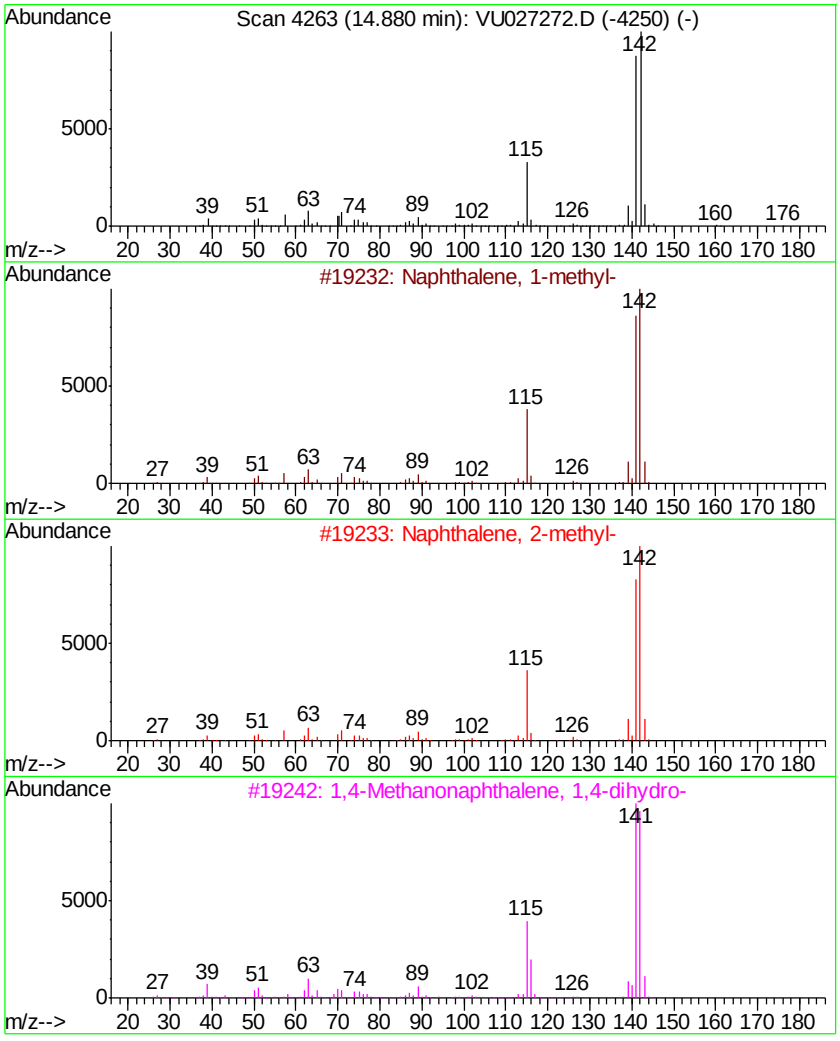
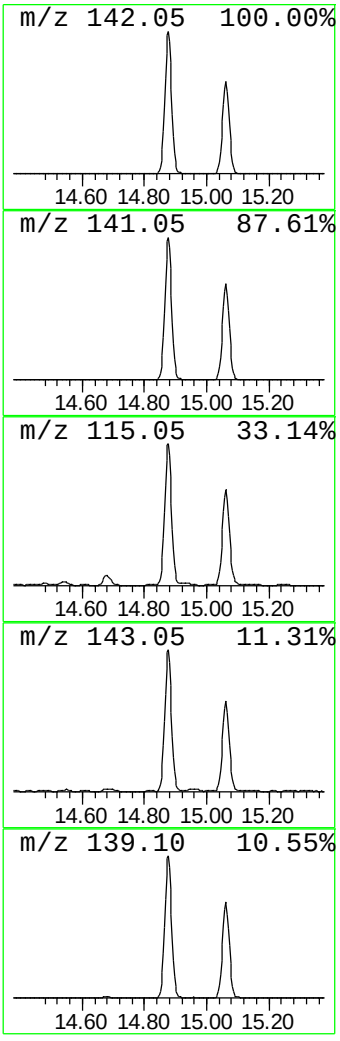
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	121.00 ug/l	2033050	1,4-Dichlorobenzene-d4	11.48

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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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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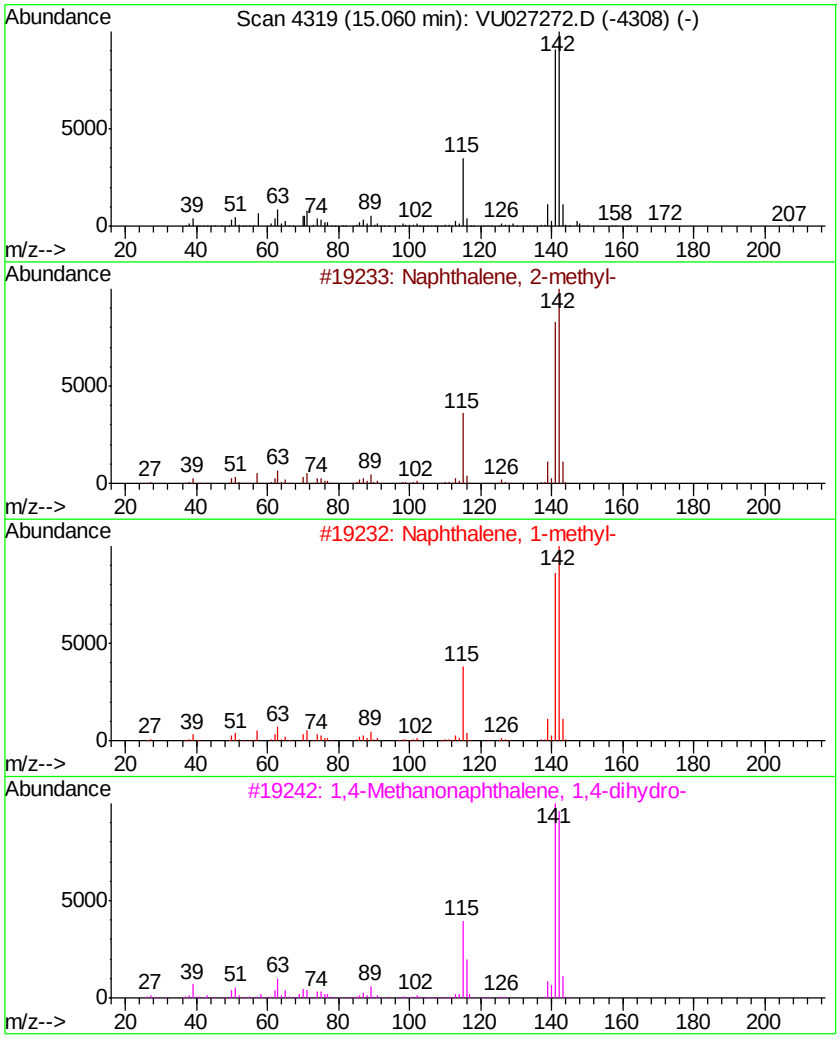
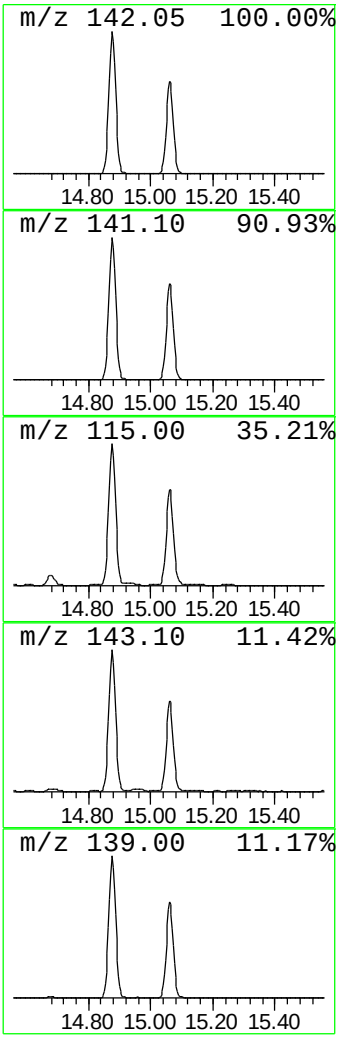
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	85.01 ug/l	1428400	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	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027272.D
Acq On : 28 Sep 2018 15:28
Operator : MD/SY
Sample : J5041-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : S13-0251

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...	5.49	73.7	ug/l	697397	2	5.89	472942	50.0
Cyclopentane, 1,2...	5.63	105.0	ug/l	992684	2	5.89	472942	50.0
Benzene, 1-ethyl-...	10.97	83.5	ug/l	1403160	4	11.48	840119	50.0
Benzene, 1,2,4-tr...	11.57	372.6	ug/l	6261250	4	11.48	840119	50.0
Indane	11.77	125.5	ug/l	2108730	4	11.48	840119	50.0
Benzene, 1-ethyl-...	11.87	68.4	ug/l	1148940	4	11.48	840119	50.0
Benzene, 2-butenyl-	13.12	147.9	ug/l	2485170	4	11.48	840119	50.0
Naphthalene, 1,2,...	13.29	75.1	ug/l	1262160	4	11.48	840119	50.0
Naphthalene, 1-me...	14.88	121.0	ug/l	2033050	4	11.48	840119	50.0
1,4-Methanonaphth...	15.06	85.0	ug/l	1428400	4	11.48	840119	50.0