LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

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\82U070918W.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	1.088	136	155	178	rBV	705417	830396	73.04%	6.765%
2	1.404	246	253	263	rBV	11675	12850	1.13%	0.105%
3	1.757	354	363	376	rBV	30150	39747	3.50%	0.324%
4	1.947	415	422	436	rVB3	9780	14501	1.28%	0.118%
5	2.622	620	632	645	rBV	6199	11510	1.01%	0.094%
6	2.709	645	659	666	rBV3	30985	66702	5.87%	0.543%
7	2.796	679	686	702	rVB2	14136	26622	2.34%	0.217%
8	2.998	735	749	766	rBV	38372	81467	7.17%	0.664%
9	4.002	1043	1061	1074	rBV4	6951	19666	1.73%	0.160%
10	4.098	1074	1091	1112	rVV	57447	152236	13.39%	1.240%
11	4.230	1115	1132	1156	rVB2	94222	230170	20.24%	1.875%
12	4.886	1321	1336	1353	rBV2	169642	383436	33.72%	3.124%
13	4.989	1353	1368	1387	rVV3	303144	720387	63.36%	5.869%
14	5.085	1388	1398	1419	rVB4	16652	41380	3.64%	0.337%
15	5.227	1428	1442	1443	rBV5	8666	12335	1.08%	0.100%
16	5.313	1457	1469	1486	rVB	178942	375486	33.03%	3.059%
17	5.397	1486	1495	1508	rVB3	7428	15878	1.40%	0.129%
18	5.487	1508	1523	1530	rBV3	28929	65837	5.79%	0.536%
19	5.548	1531	1542	1557	rVV5	58678	158745	13.96%	1.293%
20	5.628	1557	1567	1582	rVB4	23376	50163	4.41%	0.409%
21	5.783	1594	1615	1616	rBV2	16884	48654	4.28%	0.396%
22	5.889	1636	1648	1667	rBV	350335	690133	60.70%	5.623%
23	6.188	1729	1741	1763	rVB2	96160	188815	16.61%	1.538%
24	6.416	1794	1812	1827	rBV3	57747	139591	12.28%	1.137%
25	6.927	1954	1971	1986	rBV5	27288	65834	5.79%	0.536%
26	7.053	1992	2010	2024	rBV	72487	146853	12.92%	1.196%
27	7.567	2147	2170	2190	rBV	627836	1136965	100.00%	9.263%
28	7.921	2267	2280	2291	rBV4	12626	22065	1.94%	0.180%
29	8.230	2364	2376	2396	rBV	130810	224706	19.76%	1.831%
30	8.551	2466	2476	2486	rVB3	10999	17978	1.58%	0.146%
31	9.095	2632	2645	2650	rBV	497395	876060	77.05%	7.137%
32	9.365	2718	2729	2736	rBV7	11207	22611	1.99%	0.184%
33	9.574	2784	2794	2804	rBV4	8422	13662	1.20%	0.111%
34	9.747	2826	2848	2859	rBV8	21134	61919	5.45%	0.504%

LSC Area Percent Report

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Sample : J3922-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

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	9.927	2886	2904	2906	rBV4	10267	21129	1.86%	0.172%
36	9.956	2907	2913	2924	rVB2	24243	38567	3.39%	0.314%
37	10.049	2929	2942	2959	rBV8	21293	45780	4.03%	0.373%
38	10.168	2970	2979	2988	rBV	56823	82102	7.22%	0.669%
39	10.226	2988	2997	3008	rVB	9077	18713	1.65%	0.152%
40	10.310	3011	3023	3036	rBV	484701	767495	67.50%	6.253%
41	10.377	3036	3044	3053	rVB4	21989	38380	3.38%	0.313%
42	10.500	3073	3082	3091	rVB5	24998	39987	3.52%	0.326%
43	10.590	3099	3110	3121	rBV	67591	107693	9.47%	0.877%
44	10.699	3136	3144	3156	rVB4	12586	25259	2.22%	0.206%
45	10.773	3156	3167	3174	rBV4	6817	11657	1.03%	0.095%
46	10.869	3187	3197	3207	rVB8	6900	14469	1.27%	0.118%
47	10.982	3219	3232	3248	rBV4	40567	77672	6.83%	0.633%
48	11.104	3261	3270	3278	rVB	25492	38838	3.42%	0.316%
49	11.152	3278	3285	3297	rBV10	8567	17425	1.53%	0.142%
50	11.294	3312	3329	3334	rBV3	37954	78412	6.90%	0.639%
51	11.326	3334	3339	3354	rVB2	42645	64822	5.70%	0.528%
52	11.410	3354	3365	3377	rVB2	6855	16470	1.45%	0.134%
53	11.487	3377	3389	3405	rBV	560964	929968	81.79%	7.577%
54	11.615	3420	3429	3441	rVB6	8447	17786	1.56%	0.145%
55	11.693	3442	3453	3466	rVB	106913	170516	15.00%	1.389%
56	11.776	3466	3479	3494	rBV3	124755	257823	22.68%	2.101%
57	11.869	3494	3508	3530	rVB8	54556	141698	12.46%	1.154%
58	11.985	3530	3544	3555	rBV2	30850	53306	4.69%	0.434%
59	12.255	3617	3628	3633	rBV	141609	212670	18.71%	1.733%
60	12.294	3633	3640	3650	rVV3	111088	199461	17.54%	1.625%
61	12.358	3650	3660	3679	rVV3	196784	492869	43.35%	4.016%
62	12.438	3679	3685	3694	rVB	21979	32836	2.89%	0.268%
63	12.500	3694	3704	3709	rBV	36128	59031	5.19%	0.481%
64	12.541	3709	3717	3733	rVB	112635	179495	15.79%	1.462%
65	12.651	3733	3751	3761	rBV2	106367	216633	19.05%	1.765%
66	12.709	3761	3769	3782	rVB2	27055	55042	4.84%	0.448%
67	12.792	3784	3795	3808	rBV3	40817	64992	5.72%	0.530%
68	12.882	3812	3823	3829	rBV	17591	25162	2.21%	0.205%
69	12.927	3829	3837	3843	rBV	23748	34333	3.02%	0.280%
70	12.969	3844	3850	3855	rVV	21422	30324	2.67%	0.247%
71	13.001	3855	3860	3871	rVB	23743	33228	2.92%	0.271%

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ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

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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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Title : SW846 8260

72	13.075	3874	3883	3887	rBV4	11380	18656	1.64%	0.152%
73	13.123	3888	3898	3911	rVB2	180782	288227	25.35%	2.348%
74	13.200	3916	3922	3929	rVV5	10635	16000	1.41%	0.130%
75	13.262	3929	3941	3944	rVV4	13997	25435	2.24%	0.207%
76	13.294	3945	3951	3969	rVB3	26010	47618	4.19%	0.388%
77	13.413	3979	3988	3993	rBV2	7691	11888	1.05%	0.097%
78	13.458	3995	4002	4010	rVV3	14962	24767	2.18%	0.202%
79	13.512	4010	4019	4027	rVV3	33736	54423	4.79%	0.443%
80	13.580	4027	4040	4046	rVV4	23849	57071	5.02%	0.465%
81	13.612	4046	4050	4069	rVB3	21250	34101	3.00%	0.278%
82	13.734	4078	4088	4100	rBV4	9833	18520	1.63%	0.151%

Sum of corrected areas: 12274109

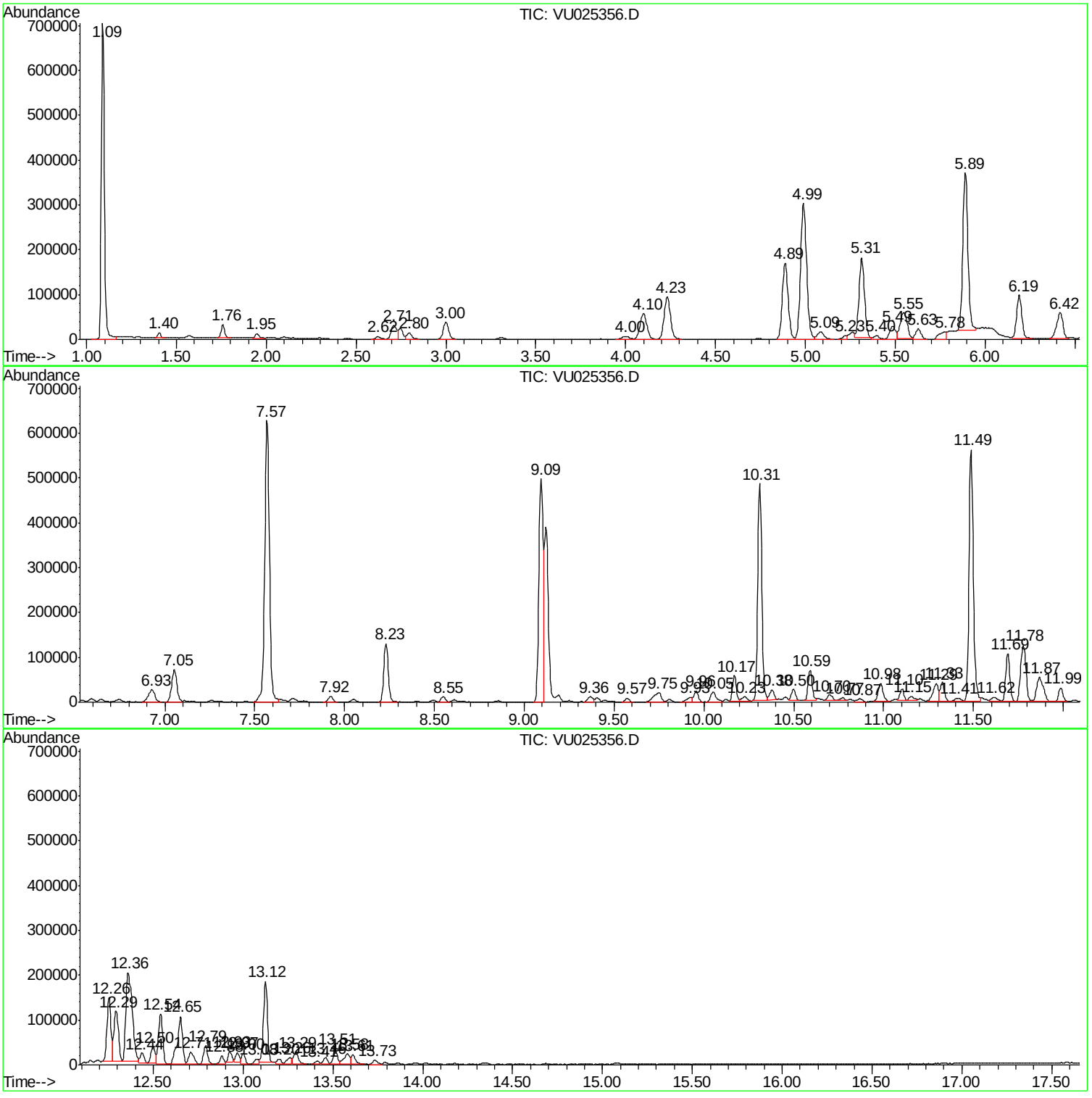
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
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ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

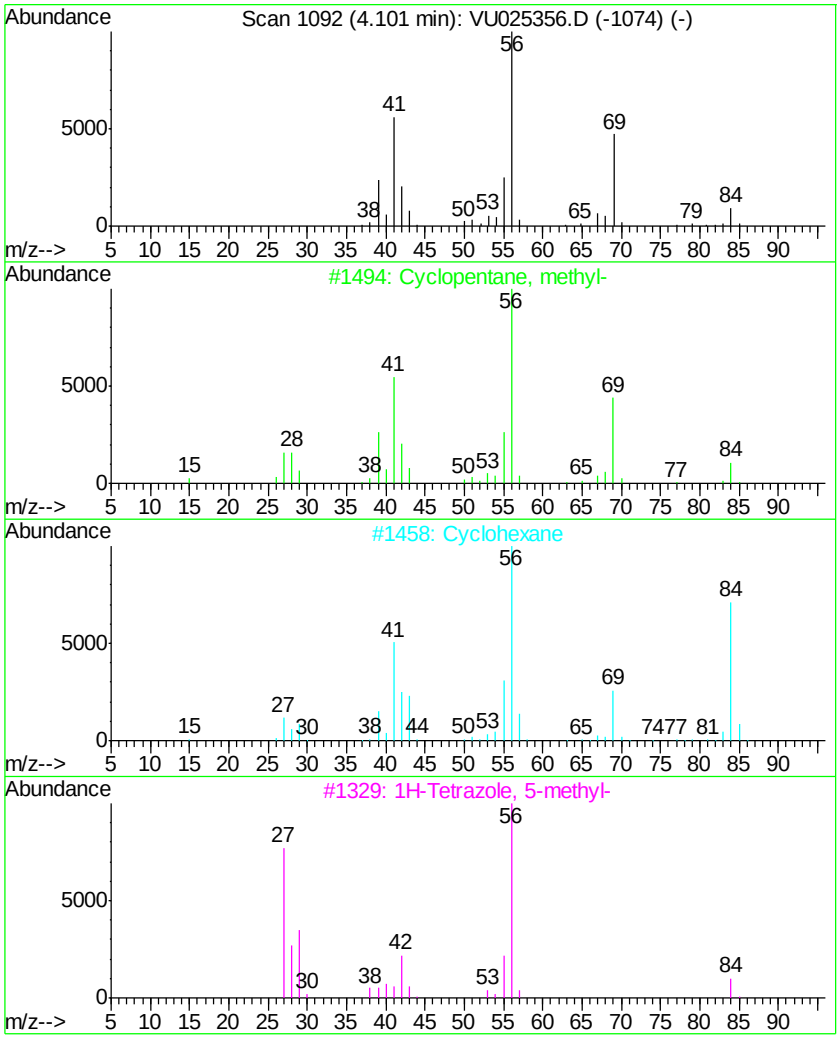
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

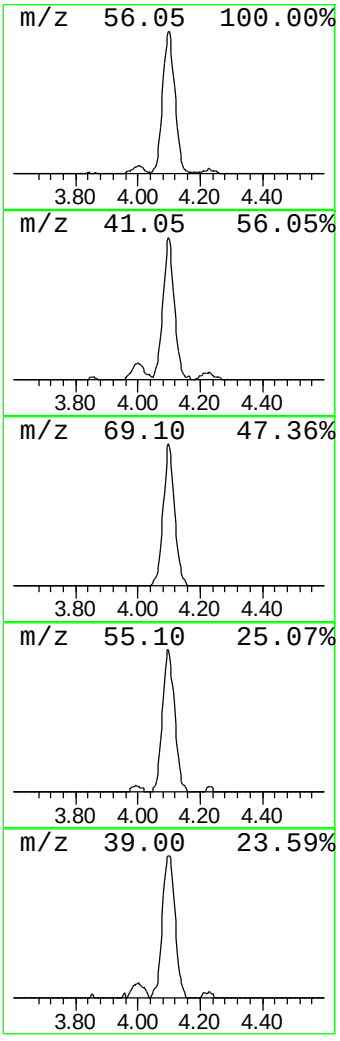
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.10	10.57 ug/l	152236	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	83
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

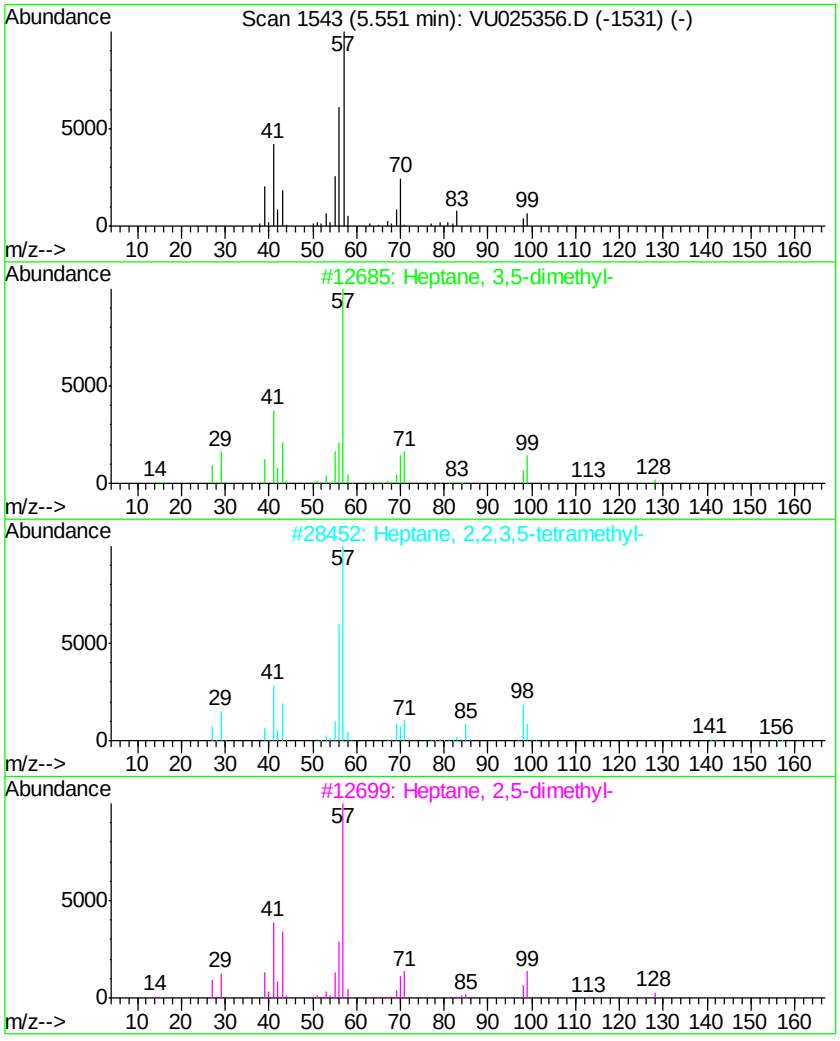
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

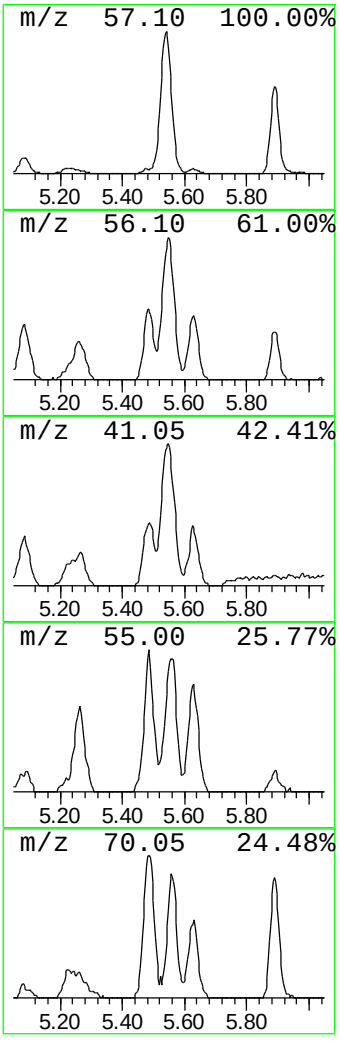
****

Peak Number 3 Heptane, 3,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.55	11.50 ug/l	158745	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	59
2	Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	56
3	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	53
4	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
5	1-Hexene, 5-methyl-	98	C7H14	003524-73-0	50





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ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

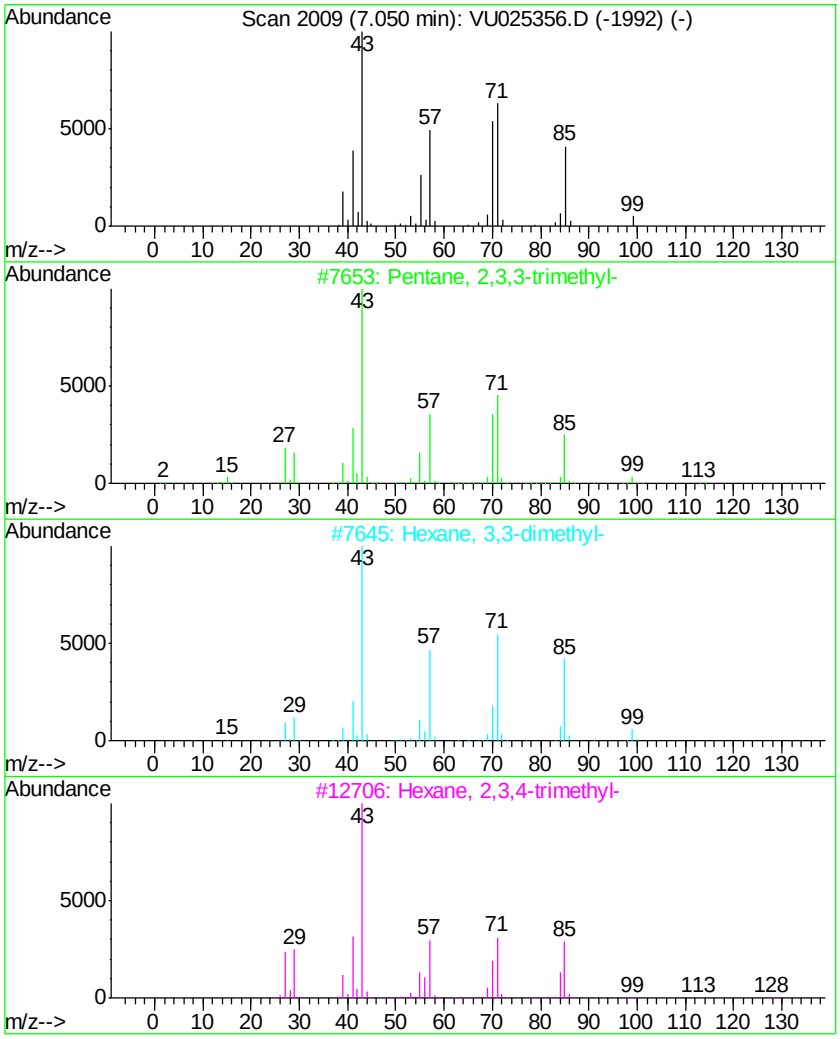
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

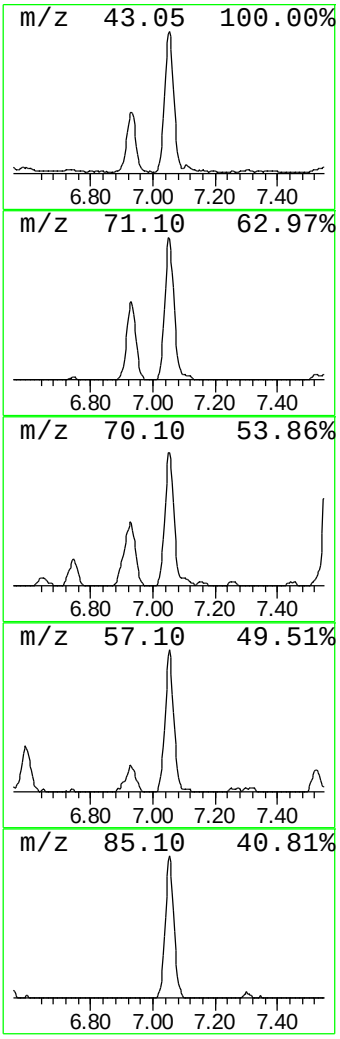
****

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	10.64 ug/l	146853	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	90
2	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	50
5	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

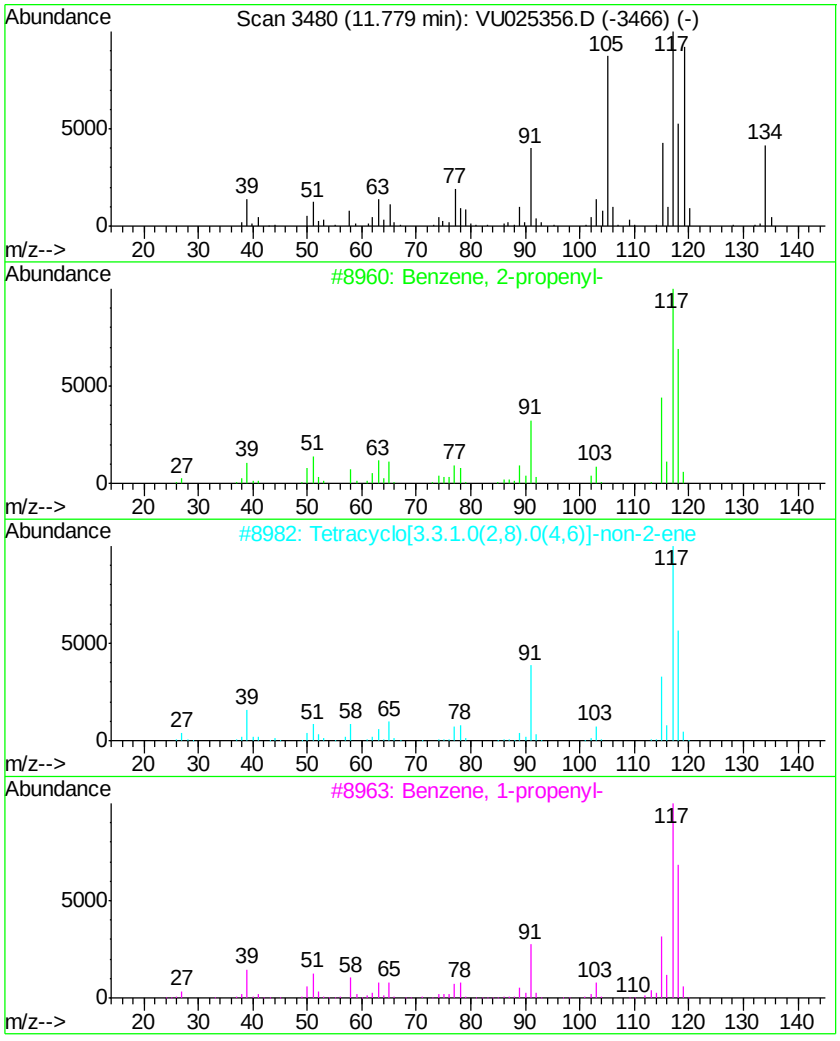
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

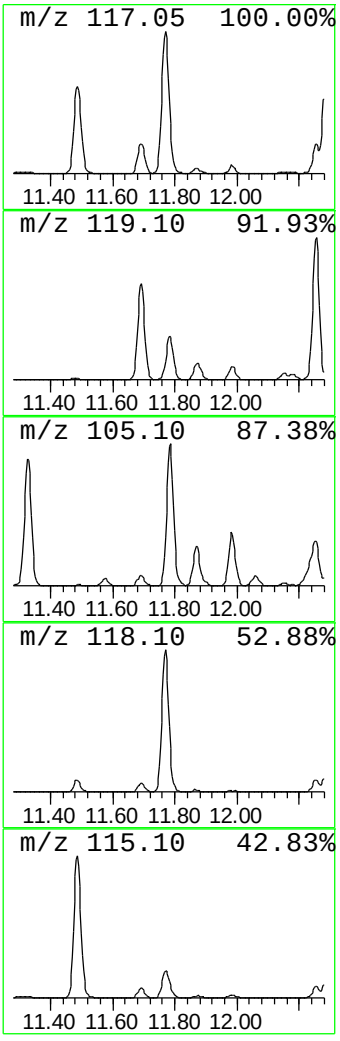
****

Peak Number 5 Benzene, 2-propenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	13.86 ug/l	257823	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	55
3	Benzene, 1-propenyl-	118	C9H10	000637-50-3	55
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	50





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ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

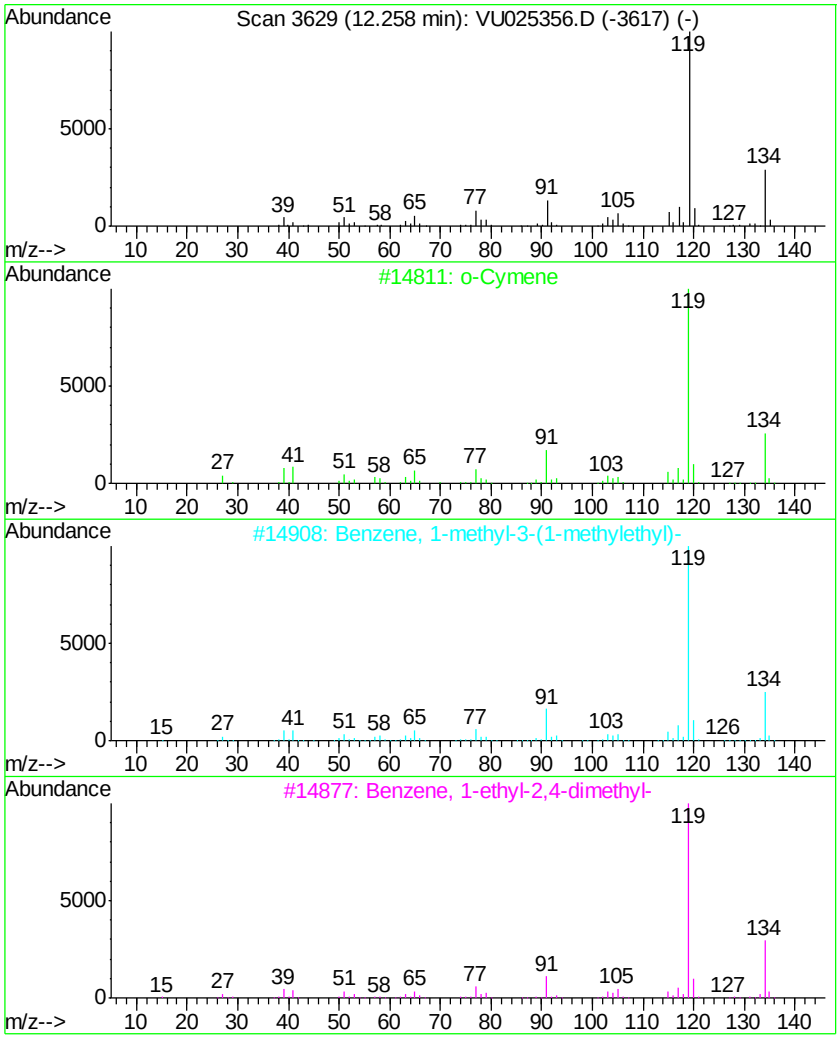
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

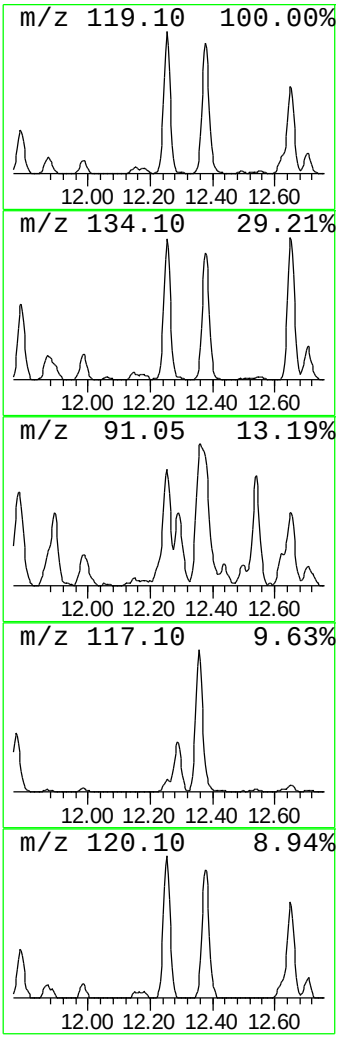
****

Peak Number 6 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	11.43 ug/l	212670	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95





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ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

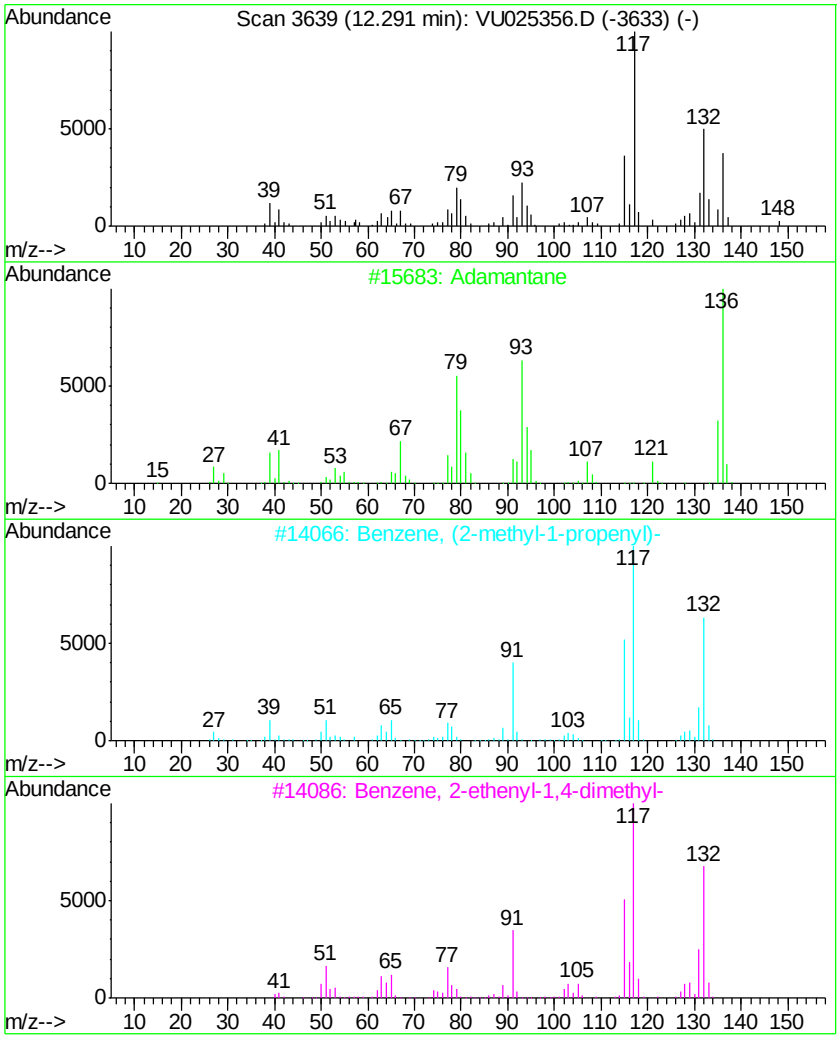
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

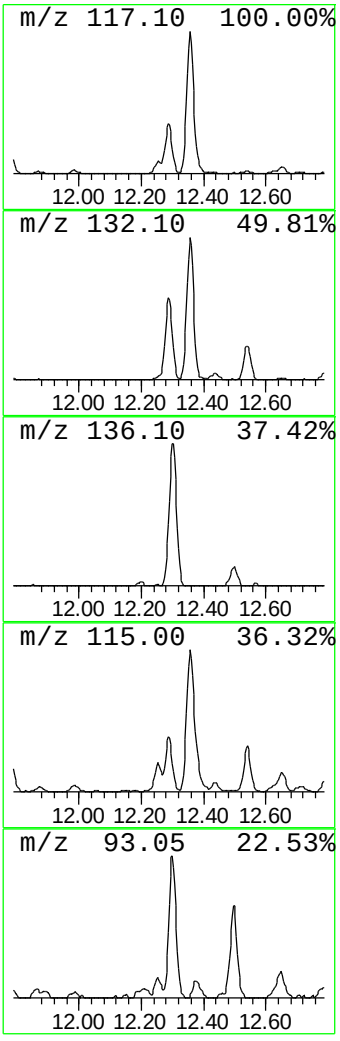
****

Peak Number 7 Adamantane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	10.72 ug/l	199461	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adamantane	136	C10H16	000281-23-2	91
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

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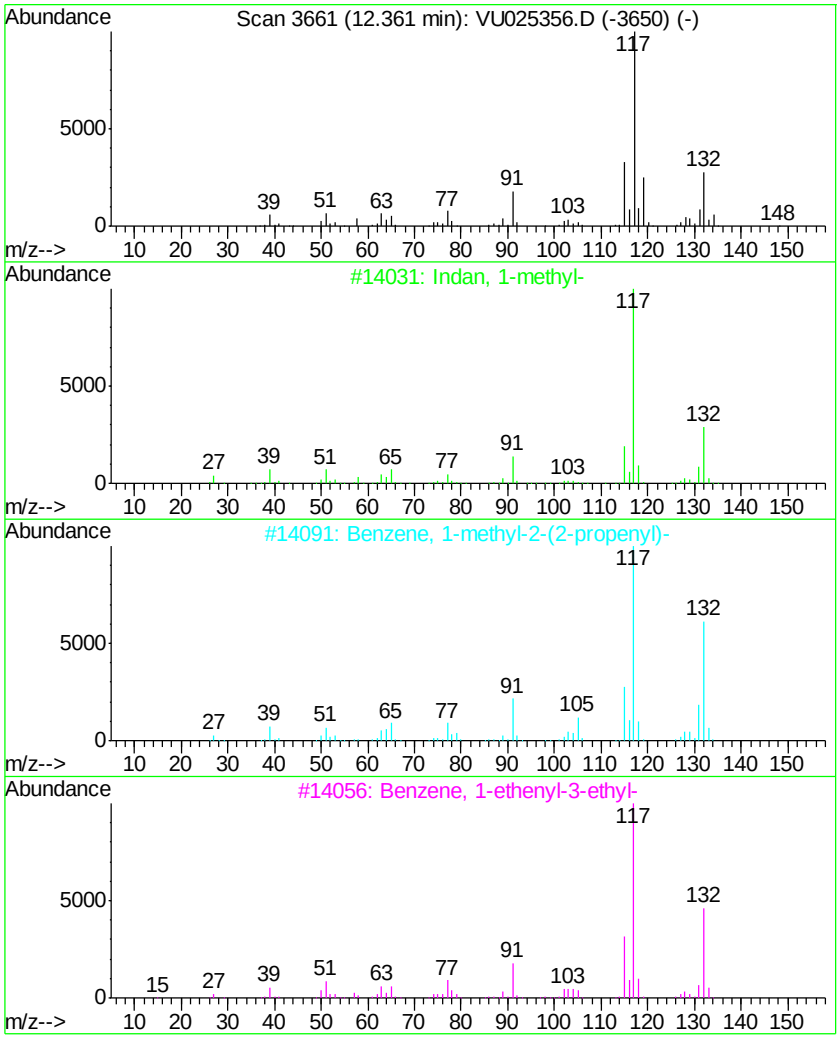
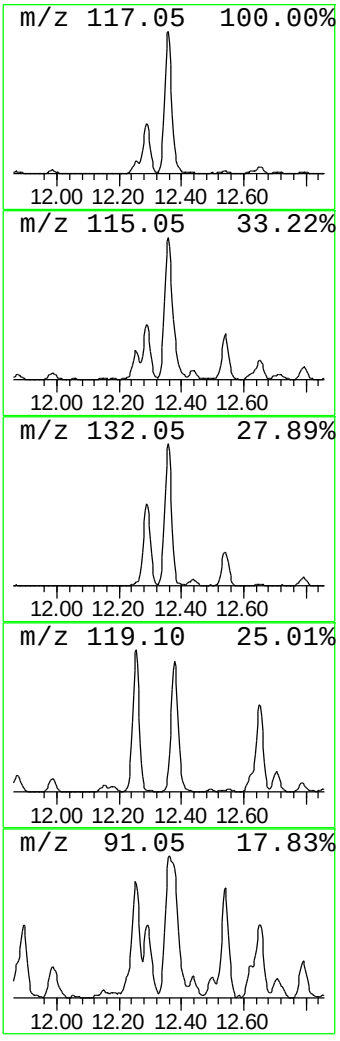
Peak Number 8 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	26.50 ug/l	492869	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	80





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

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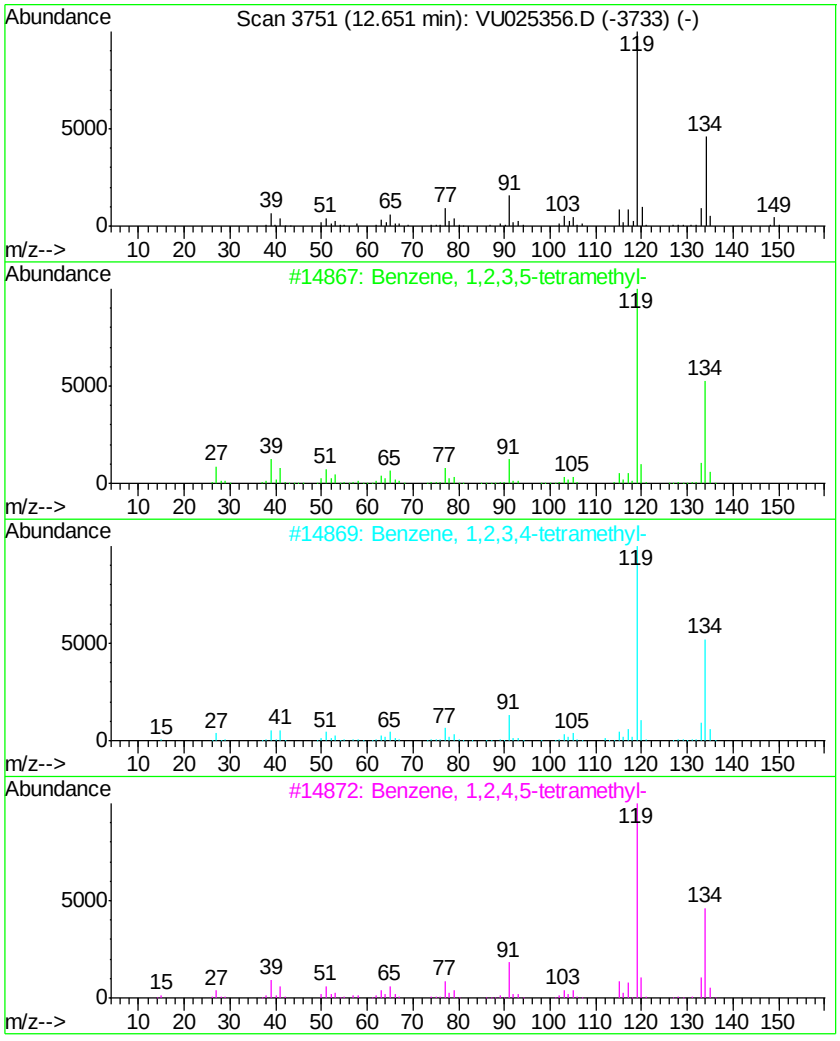
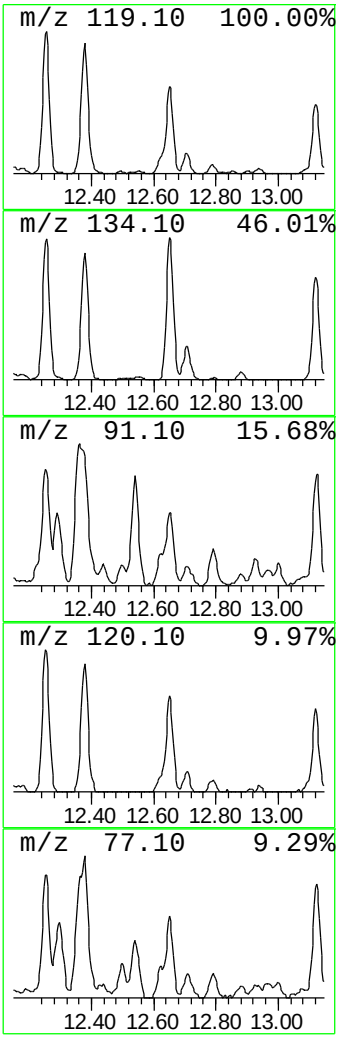
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	11.65 ug/l	216633	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5	o-Cymene	134	C10H14	000527-84-4	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

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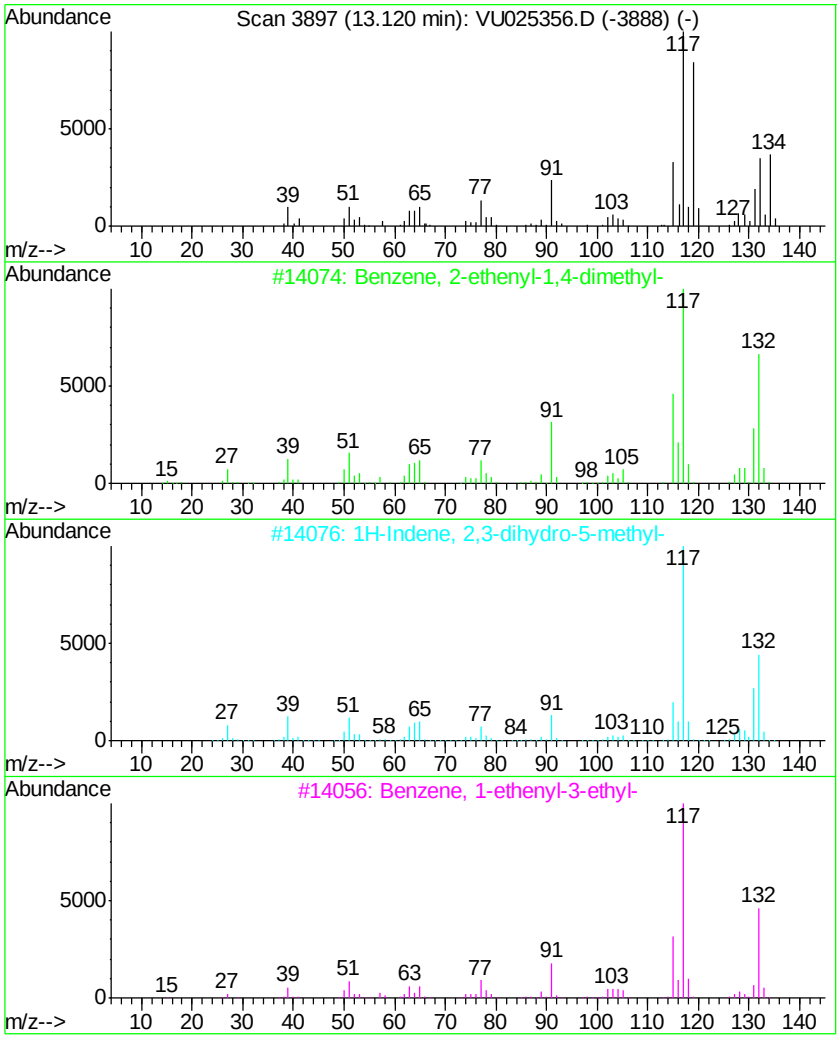
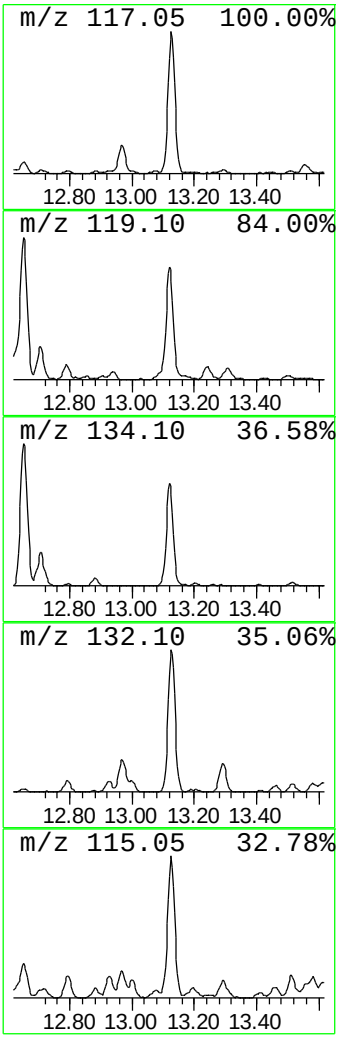
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	15.50 ug/l	288227	1,4-Dichlorobenzene-d4	11.49

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-5-methyl-	132	C10H12	000874-35-1	92
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
5	2,4-Dimethylstyrene	132	C10H12	002234-20-0	89





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071218\
Data File : VU025356.D
Acq On : 12 Jul 2018 21:59
Operator : MD/SY
Sample : J3922-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SS037MW266-180709

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.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, met...	4.10	10.6	ug/l	152236	1	4.99	720387	50.0
Heptane, 3,5-dime...	5.55	11.5	ug/l	158745	2	5.89	690133	50.0
Pentane, 2,3,3-tr...	7.05	10.6	ug/l	146853	2	5.89	690133	50.0
Benzene, 2-propenyl-	11.78	13.9	ug/l	257823	4	11.49	929968	50.0
o-Cymene	12.26	11.4	ug/l	212670	4	11.49	929968	50.0
Adamantane	12.29	10.7	ug/l	199461	4	11.49	929968	50.0
Indan, 1-methyl-	12.36	26.5	ug/l	492869	4	11.49	929968	50.0
Benzene, 1,2,3,5-...	12.65	11.7	ug/l	216633	4	11.49	929968	50.0
Benzene, 2-etheny...	13.12	15.5	ug/l	288227	4	11.49	929968	50.0