LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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	2.321	349	357	373	rVB	12970	21748	3.20%	0.310%
2	2.443	385	395	414	rBV2	15600	27848	4.10%	0.397%
3	2.829	498	515	528	rBV2	5157	10895	1.60%	0.155%
4	3.003	550	569	591	rBV	90023	197678	29.08%	2.821%
5	4.202	926	942	951	rBV2	3314	8304	1.22%	0.118%
6	4.884	1135	1154	1171	rBV	96909	218990	32.22%	3.125%
7	4.983	1171	1185	1208	rVB	127757	285920	42.06%	4.080%
8	5.311	1272	1287	1303	rBV	123882	260374	38.30%	3.715%
9	5.392	1303	1312	1328	rVB3	9632	21290	3.13%	0.304%
10	5.887	1452	1466	1485	rBV	199922	390513	57.45%	5.572%
11	6.417	1614	1631	1642	rVB5	4953	10525	1.55%	0.150%
12	7.565	1969	1988	2025	rBV	376681	679763	100.00%	9.700%
13	8.176	2166	2178	2187	rBV4	3359	7119	1.05%	0.102%
14	9.089	2450	2462	2485	rBV	281593	477721	70.28%	6.817%
15	9.369	2539	2549	2567	rVB9	3322	8622	1.27%	0.123%
16	10.166	2786	2797	2808	rBV	40694	62748	9.23%	0.895%
17	10.308	2830	2841	2858	rBV	253419	406653	59.82%	5.803%
18	10.491	2891	2898	2915	rVB5	4140	7271	1.07%	0.104%
19	10.588	2915	2928	2950	rBV	51174	84807	12.48%	1.210%
20	11.099	3075	3087	3095	rBV4	7169	12194	1.79%	0.174%
21	11.292	3134	3147	3150	rBV3	8108	12223	1.80%	0.174%
22	11.324	3150	3157	3173	rVB	26610	44061	6.48%	0.629%
23	11.485	3195	3207	3222	rBV	317215	489916	72.07%	6.991%
24	11.572	3225	3234	3242	rVB2	7458	11419	1.68%	0.163%
25	11.687	3262	3270	3281	rVB2	15486	23251	3.42%	0.332%
26	11.768	3282	3295	3314	rVV4	179642	337730	49.68%	4.819%
27	11.867	3316	3326	3345	rVB8	19232	48245	7.10%	0.688%
28	11.983	3352	3362	3373	rBV2	30916	47407	6.97%	0.676%
29	12.250	3432	3445	3452	rBV	102713	163327	24.03%	2.331%
30	12.285	3452	3456	3468	rVV2	45992	67517	9.93%	0.963%
31	12.356	3468	3478	3498	rVV2	121932	256864	37.79%	3.665%
32	12.539	3526	3535	3549	rVB2	24921	41323	6.08%	0.590%
33	12.652	3549	3570	3581	rBV	94440	159283	23.43%	2.273%
34	12.713	3581	3589	3599	rVV7	5099	11076	1.63%	0.158%

LSC Area Percent Report

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Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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	12.787	3603	3612	3621	rVB4	13556	21208	3.12%	0.303%
36	12.880	3633	3641	3648	rBV2	10452	19550	2.88%	0.279%
37	12.928	3649	3656	3661	rVV4	12916	21937	3.23%	0.313%
38	12.964	3661	3667	3674	rVB	19090	27477	4.04%	0.392%
39	13.121	3694	3716	3734	rBV2	342980	561298	82.57%	8.009%
40	13.240	3746	3753	3757	rBV3	5082	7527	1.11%	0.107%
41	13.289	3758	3768	3787	rVB	73002	123529	18.17%	1.763%
42	13.401	3793	3803	3812	rBV10	11242	21466	3.16%	0.306%
43	13.456	3812	3820	3828	rVV	15087	24379	3.59%	0.348%
44	13.510	3828	3837	3845	rVV	42503	68107	10.02%	0.972%
45	13.578	3845	3858	3862	rVV4	28579	57611	8.48%	0.822%
46	13.613	3863	3869	3889	rVB3	43631	96719	14.23%	1.380%
47	13.742	3896	3909	3918	rBV	79849	137798	20.27%	1.966%
48	13.787	3919	3923	3932	rVB	14418	19875	2.92%	0.284%
49	13.854	3934	3944	3958	rBV2	28318	46044	6.77%	0.657%
50	13.951	3958	3974	3986	rBV3	36414	73006	10.74%	1.042%
51	14.012	3987	3993	4003	rVB3	14135	21689	3.19%	0.309%
52	14.153	4026	4037	4041	rBV3	14514	23916	3.52%	0.341%
53	14.350	4082	4098	4110	rBV4	39667	94234	13.86%	1.345%
54	14.478	4129	4138	4148	rVB2	11946	18827	2.77%	0.269%
55	14.539	4148	4157	4170	rVB4	17921	32716	4.81%	0.467%
56	14.678	4189	4200	4214	rBV2	64590	110456	16.25%	1.576%
57	14.877	4250	4262	4273	rBV	44691	80117	11.79%	1.143%
58	14.935	4273	4280	4295	rVB10	9137	20927	3.08%	0.299%
59	15.063	4309	4320	4334	rBV	85699	149929	22.06%	2.139%
60	15.588	4470	4483	4498	rVB6	6908	13959	2.05%	0.199%
61	15.803	4542	4550	4555	rBV5	7294	11583	1.70%	0.165%
62	15.835	4556	4560	4570	rVB3	6591	8932	1.31%	0.127%
63	15.915	4574	4585	4596	rBV3	19454	35354	5.20%	0.504%
64	16.060	4620	4630	4638	rBV	43725	73297	10.78%	1.046%
65	16.099	4638	4642	4654	rVB4	16321	22460	3.30%	0.320%
66	16.263	4682	4693	4695	rBV4	7573	9412	1.38%	0.134%
67	16.292	4697	4702	4720	rVB4	11035	19331	2.84%	0.276%
68	16.433	4734	4746	4760	rBV3	6612	10784	1.59%	0.154%
69	16.742	4833	4842	4852	rBV3	5048	7871	1.16%	0.112%

Sum of corrected areas: 7007950

LSC Report - Integrated Chromatogram

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Acq On : 28 Sep 2018 00:29
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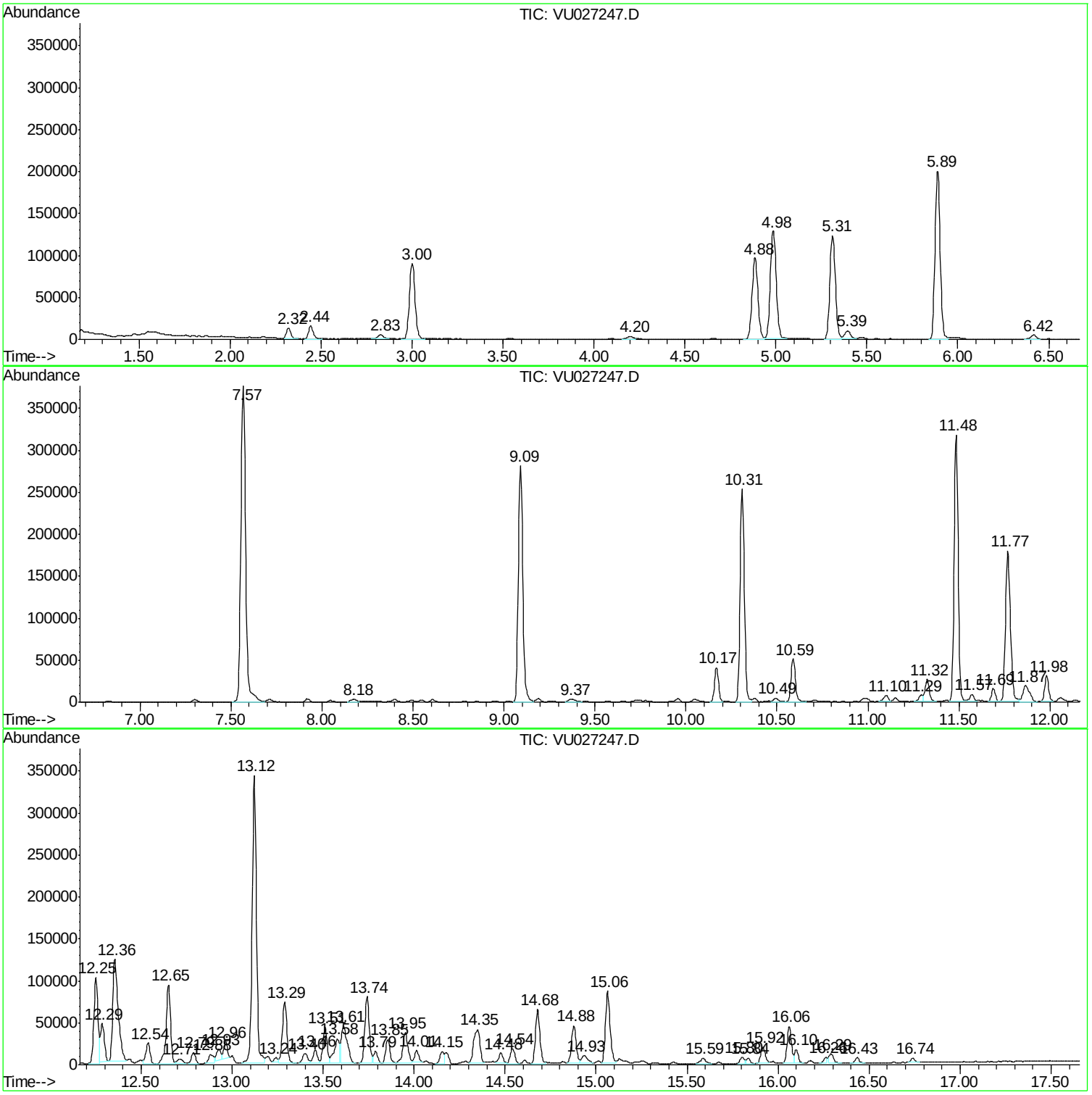
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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 : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

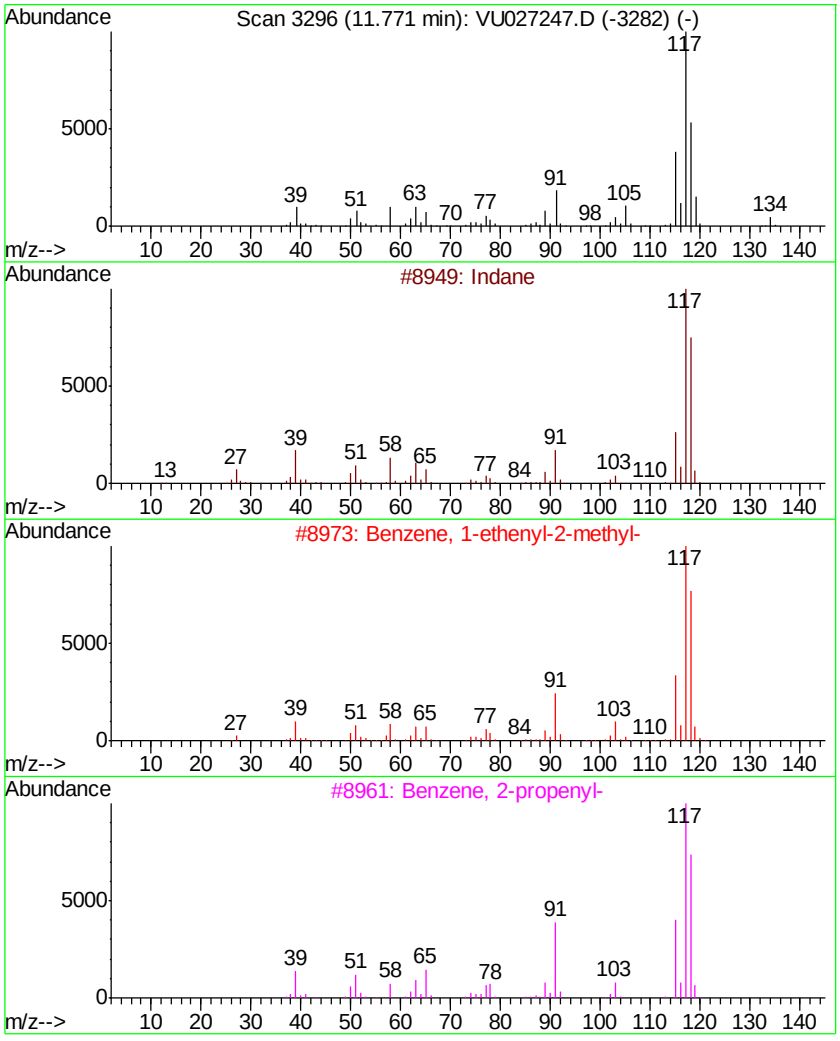
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

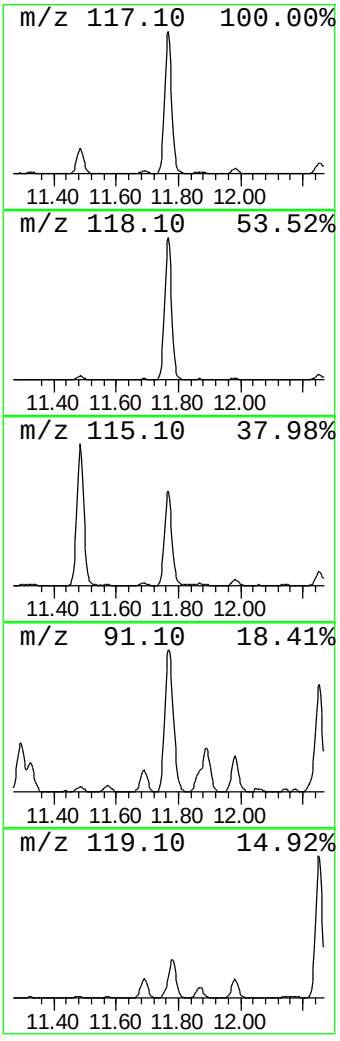
****

Peak Number 1 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	34.47 ug/l	337730	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	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
4	7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	64
5	Deltacyclene	118	C9H10	007785-10-6	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

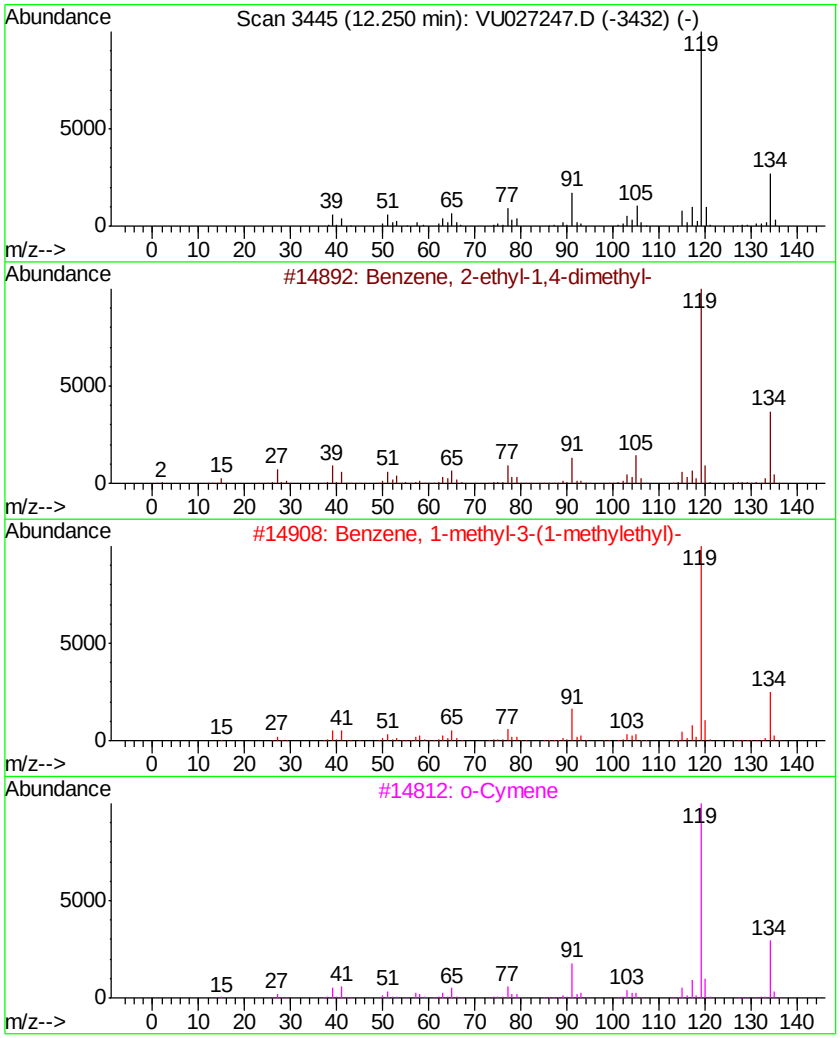
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260

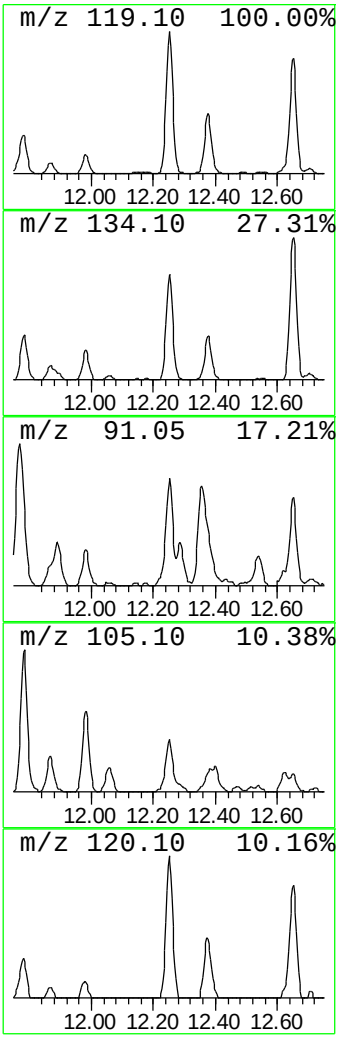
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	16.67 ug/l	163327	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

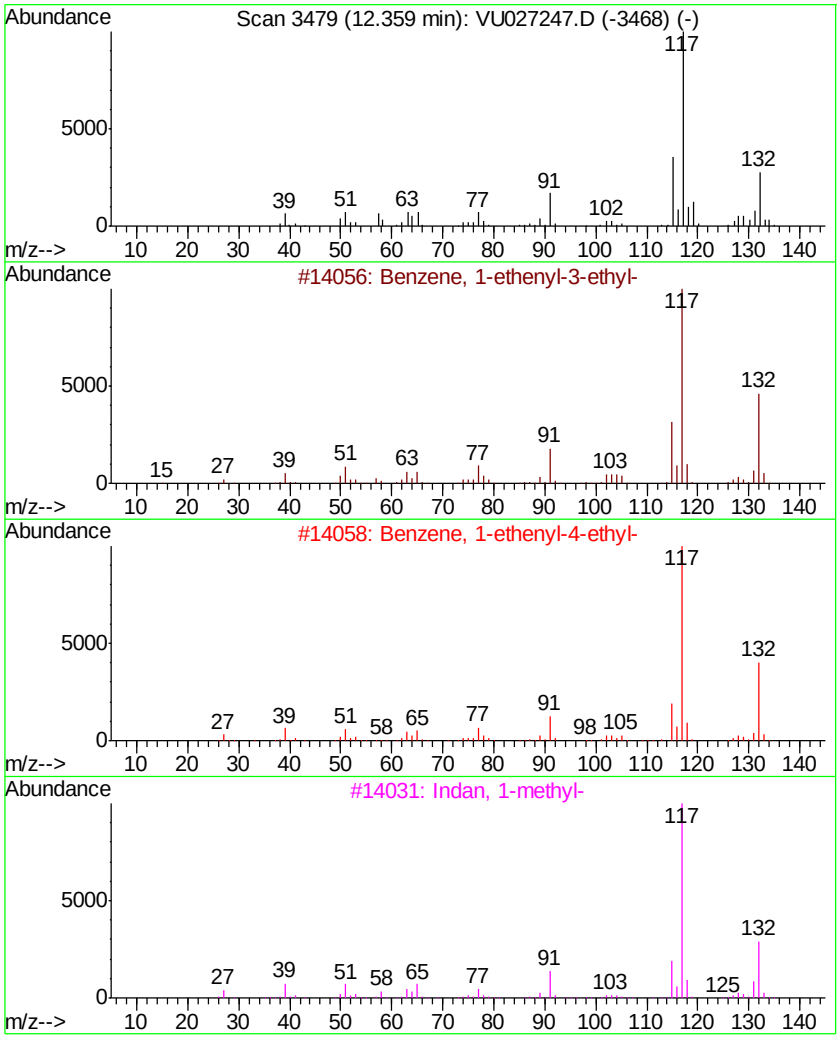
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

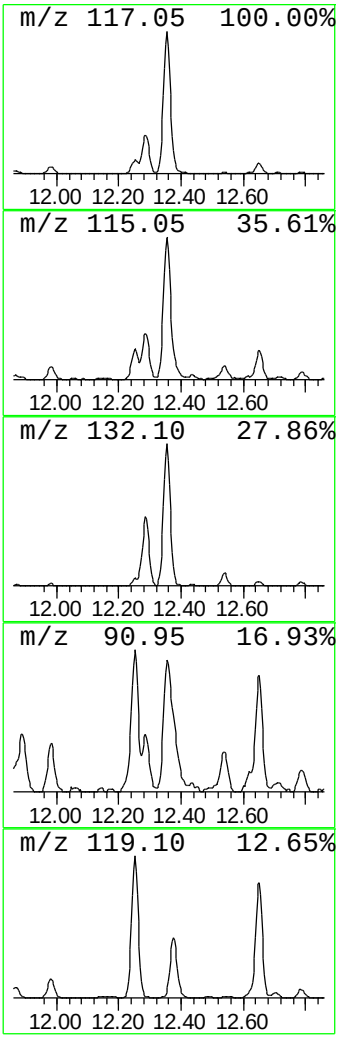
****

Peak Number 3 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	26.22 ug/l	256864	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3	Indan, 1-methyl-	132	C10H12	000767-58-8	83
4	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	80
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

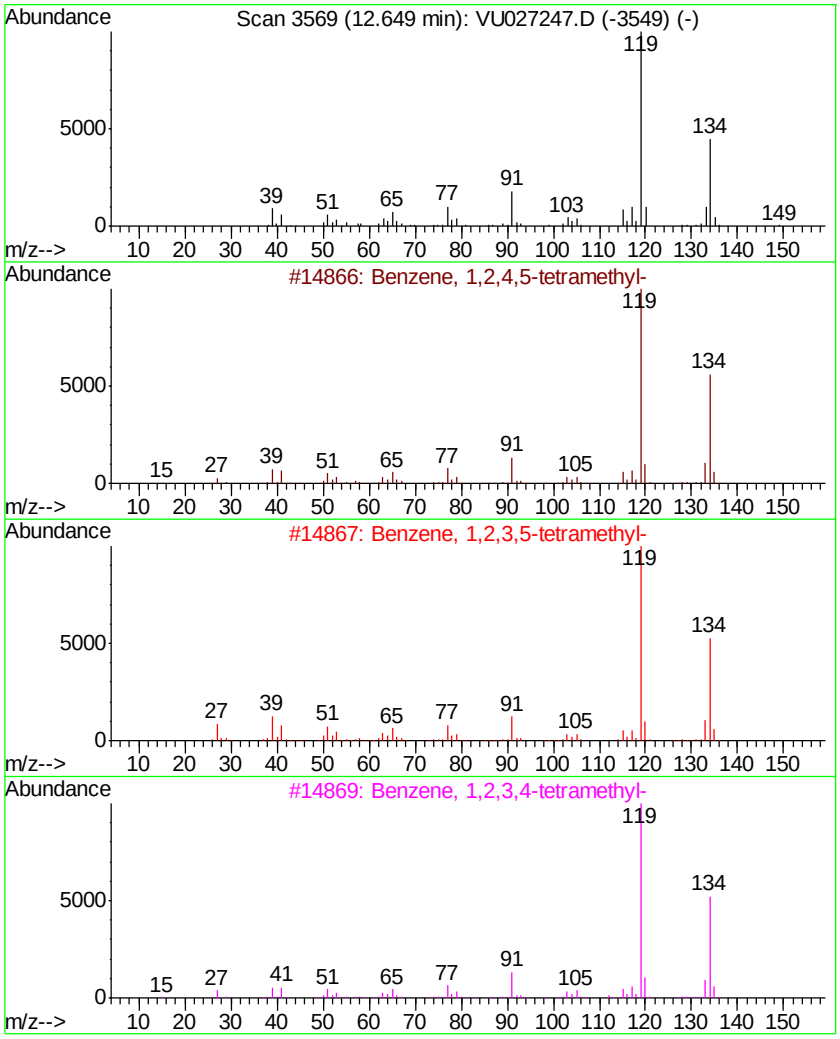
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

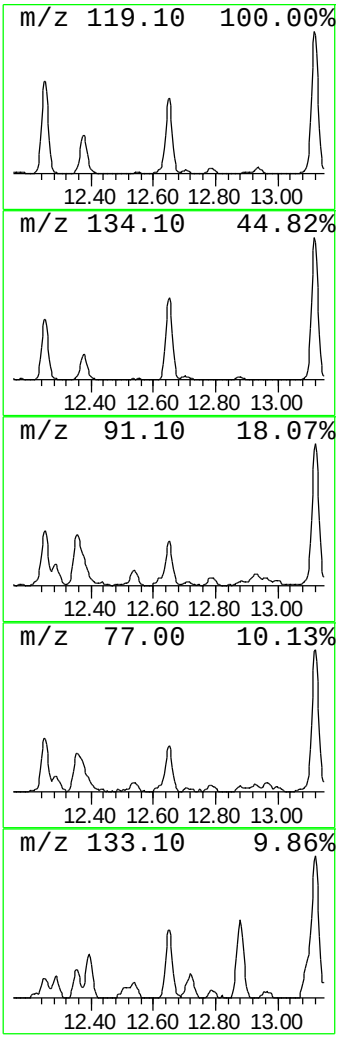
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	16.26 ug/l	159283	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	96
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	o-Cymene	134	C10H14	000527-84-4	93





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Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

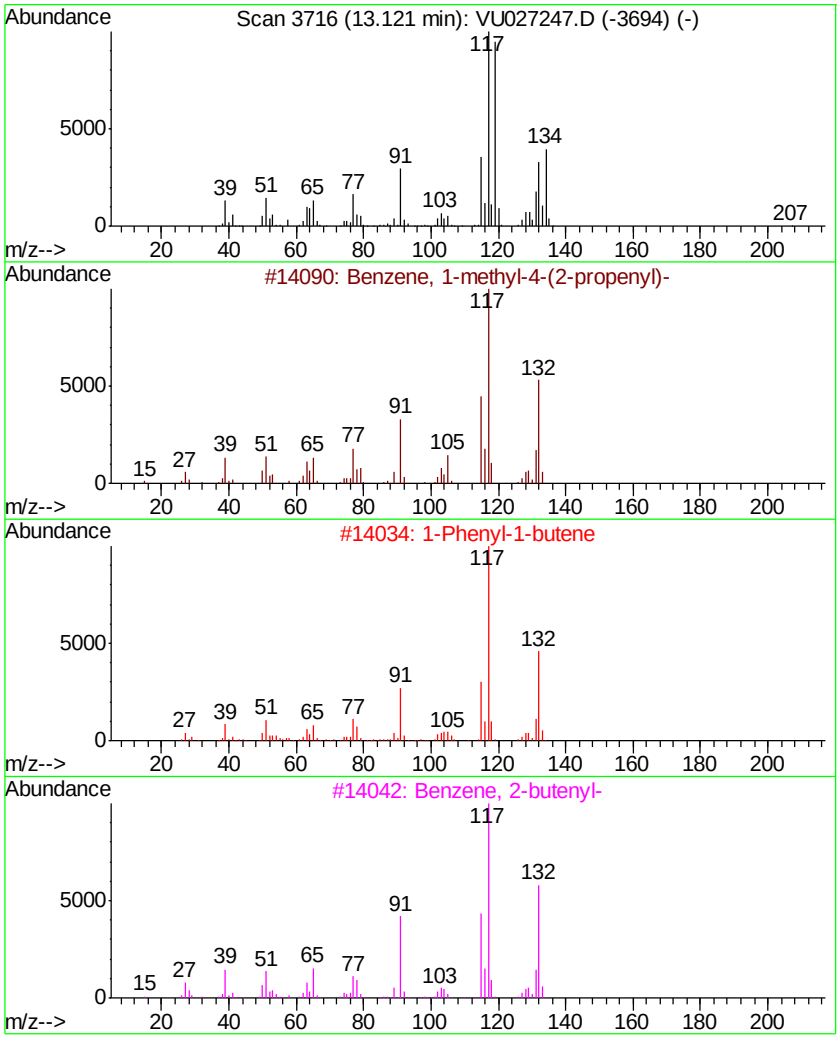
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

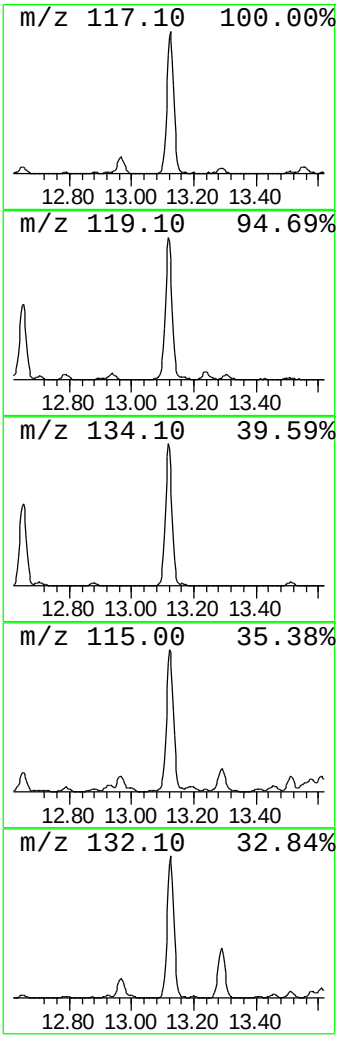
****

Peak Number 5 Benzene, 1-methyl-4-(2-prop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	57.29 ug/l	561298	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	1-Phenyl-1-butene	132	C10H12	000824-90-8	50
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	50
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	50
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	50





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Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

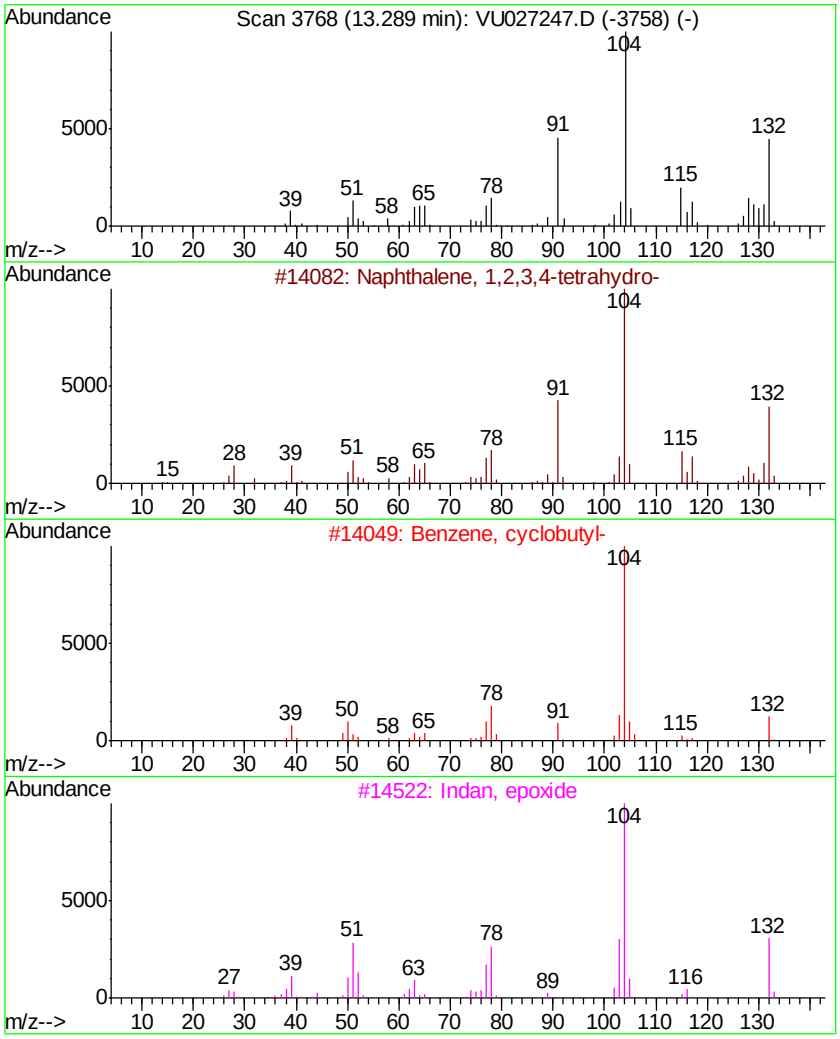
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

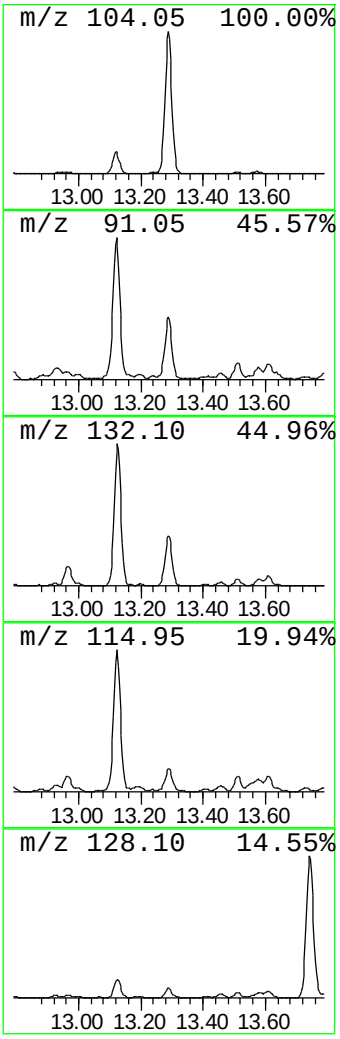
****

Peak Number 6 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	12.61 ug/l	123529	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	90
2	Benzene, cyclobutyl-	132	C10H12	004392-30-7	50
3	Indan, epoxide	132	C9H8O	000768-22-9	49
4	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	38
5	3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	27





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

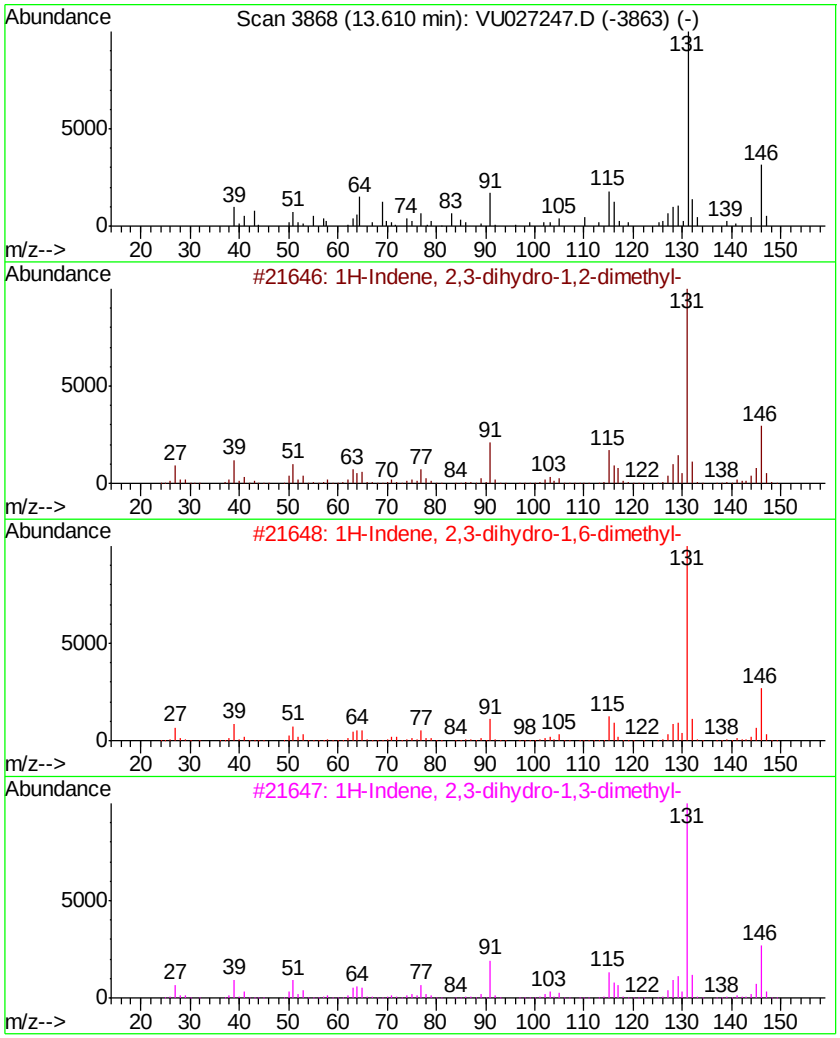
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

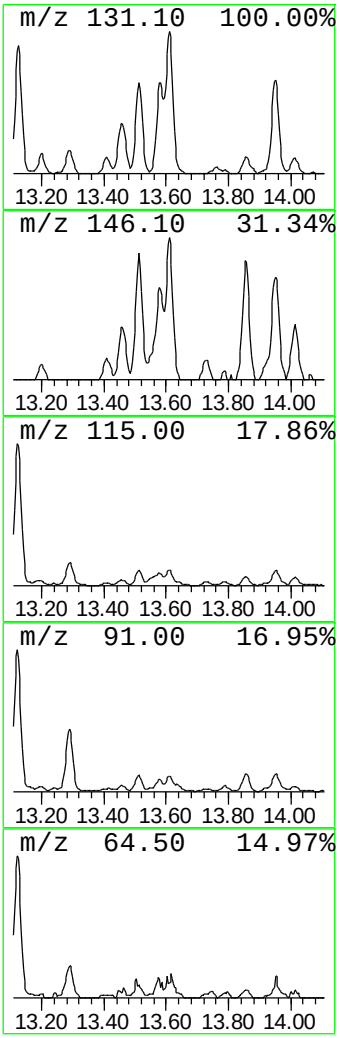
****

Peak Number 7 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.61	9.87 ug/l	96719	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	90
5	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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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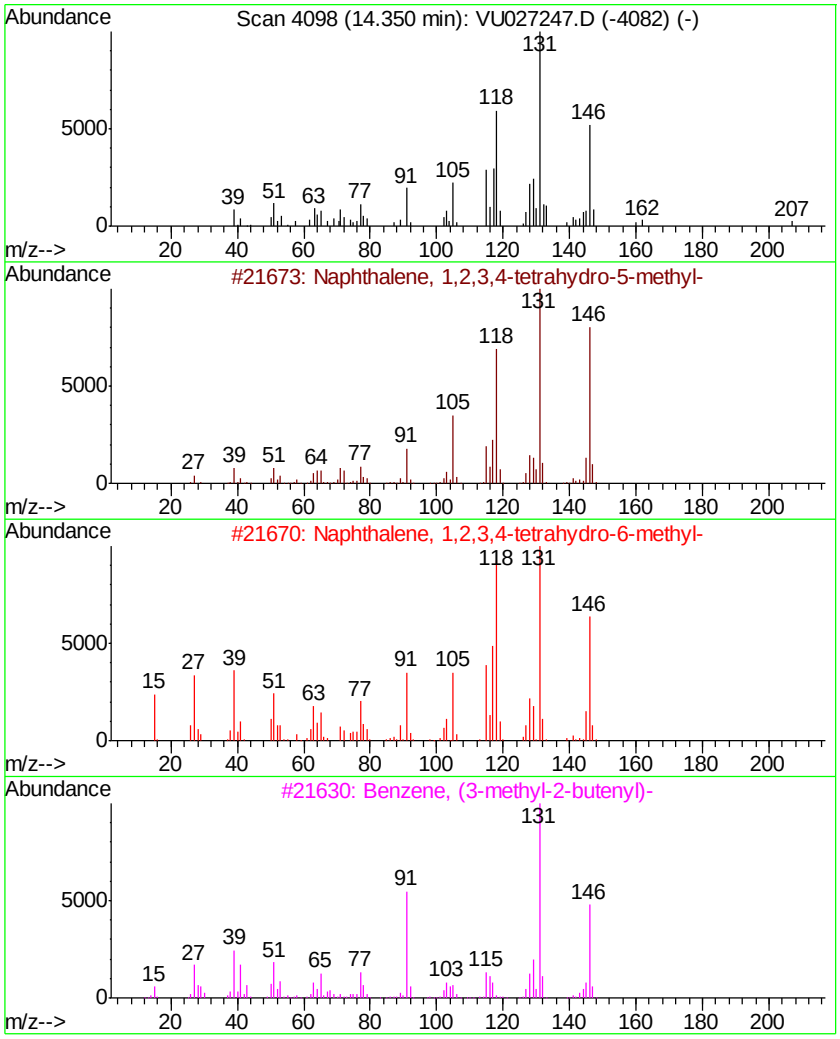
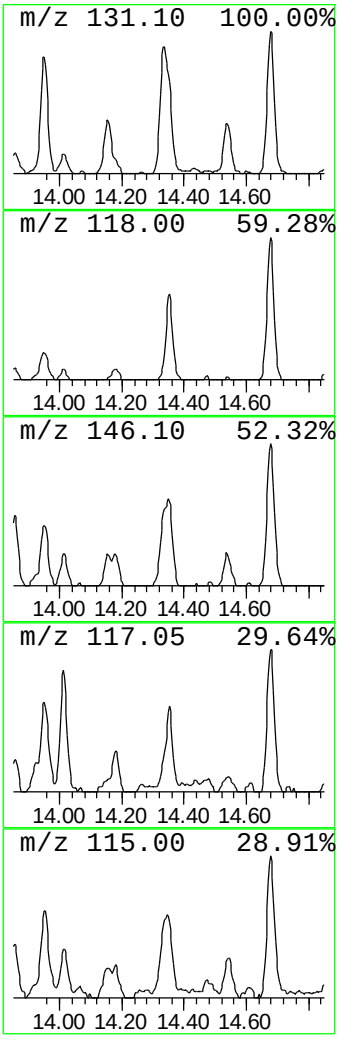
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.35	9.62 ug/l	94234	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
4	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	74
5	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	62





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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

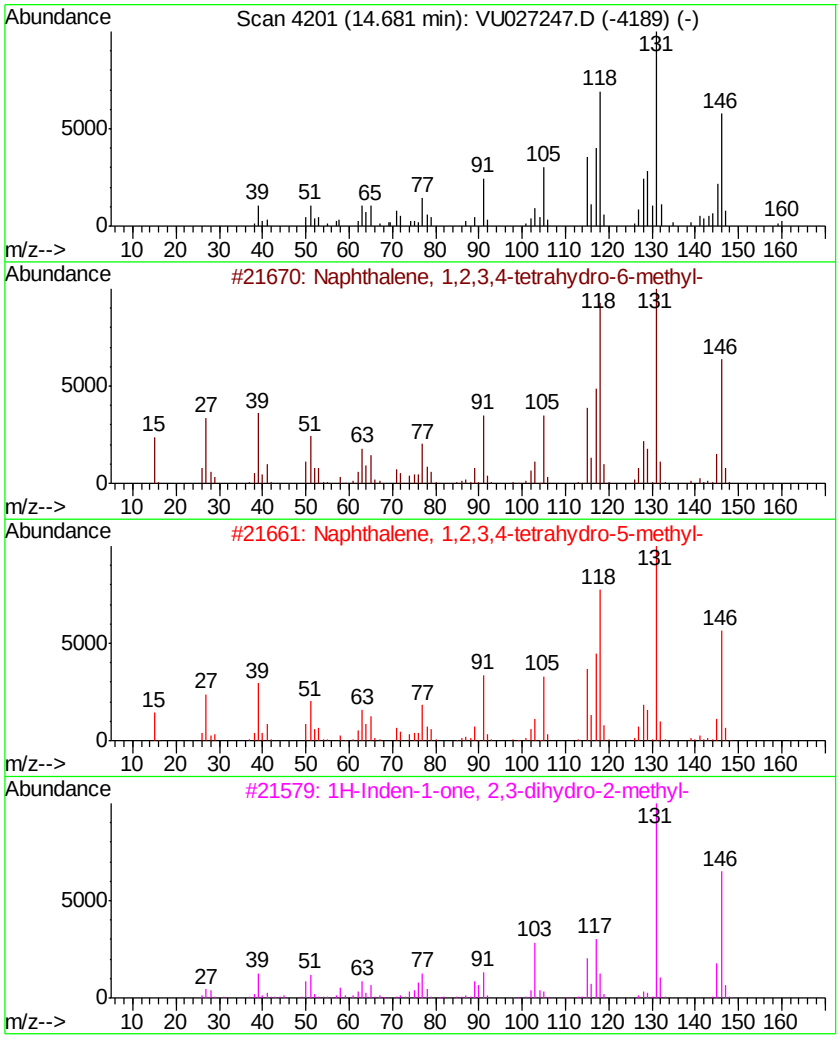
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

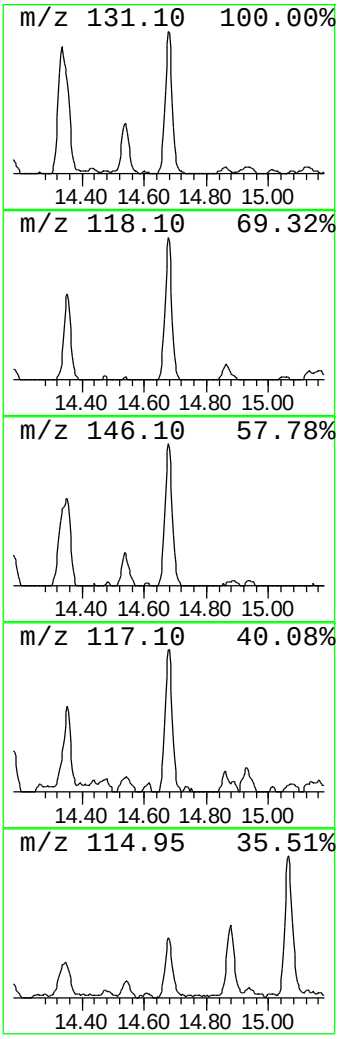
****

Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	11.27 ug/l	110456	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	90
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	70
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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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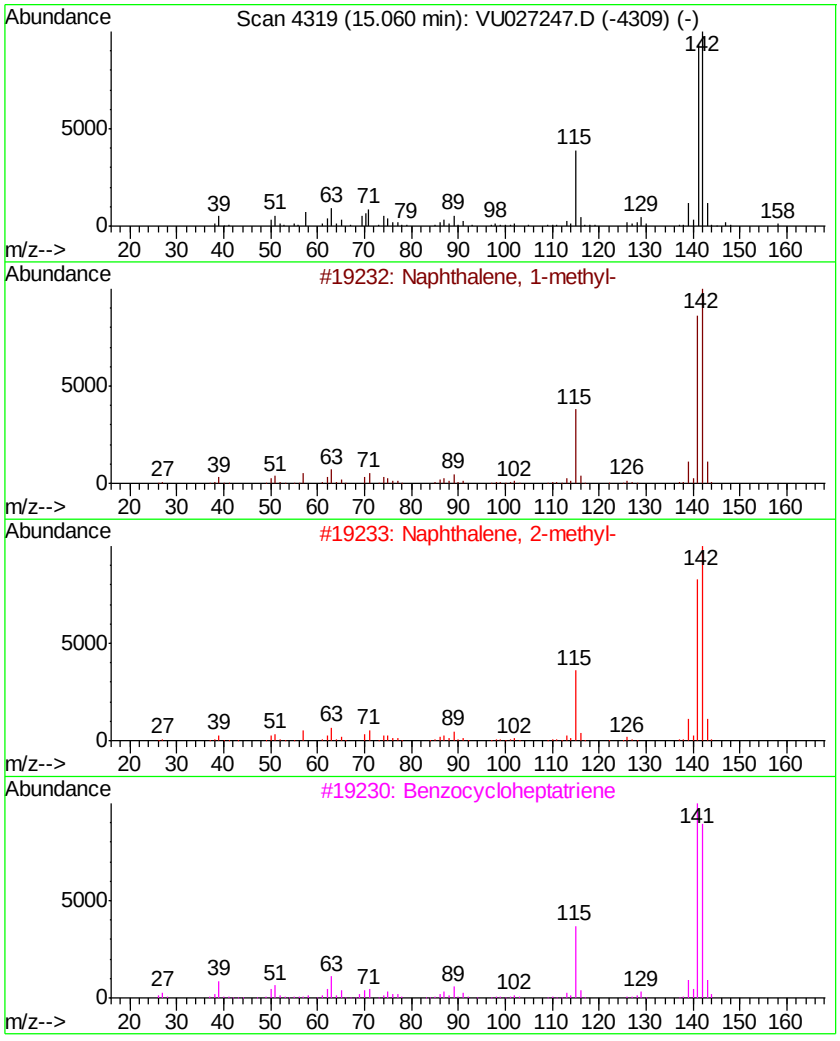
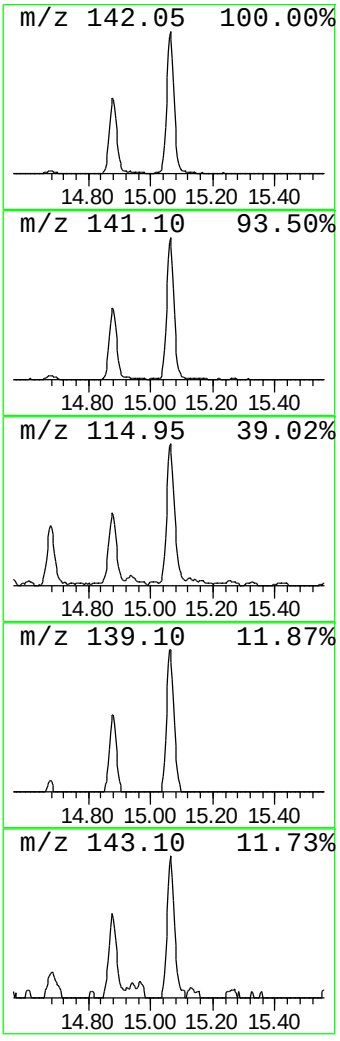
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	15.30 ug/l	149929	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	Benzocycloheptatriene	142	C11H10	000264-09-5	94
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
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 : VU027247.D
Acq On : 28 Sep 2018 00:29
Operator : MD/SY
Sample : J5029-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T11-MW-13-8.0-092018

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
Indane	11.77	34.5	ug/l	337730	4	11.48	489916	50.0
Benzene, 2-ethyl-...	12.25	16.7	ug/l	163327	4	11.48	489916	50.0
Benzene, 1-etheny...	12.36	26.2	ug/l	256864	4	11.48	489916	50.0
Benzene, 1,2,4,5-...	12.65	16.3	ug/l	159283	4	11.48	489916	50.0
Benzene, 1-methyl...	13.12	57.3	ug/l	561298	4	11.48	489916	50.0
Naphthalene, 1,2,...	13.29	12.6	ug/l	123529	4	11.48	489916	50.0
1H-Indene, 2,3-di...	13.61	9.9	ug/l	96719	4	11.48	489916	50.0
Naphthalene, 1,2,...	14.35	9.6	ug/l	94234	4	11.48	489916	50.0
Naphthalene, 1,2,...	14.68	11.3	ug/l	110456	4	11.48	489916	50.0
Naphthalene, 1-me...	15.06	15.3	ug/l	149929	4	11.48	489916	50.0