LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

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\82U050621W.M
Title : SW846 8260

Signal : TIC: VU043616.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.787	127	139	150	rVB5	6246	10832	1.29%	0.158%
2	2.761	432	442	461	rBV	6769	13196	1.57%	0.192%
3	5.298	1212	1231	1244	rBV	135525	281658	33.45%	4.099%
4	5.382	1244	1257	1277	rVB	187231	379997	45.13%	5.530%
5	5.710	1345	1359	1382	rBV	161702	324201	38.50%	4.718%
6	6.256	1513	1529	1547	rBV	294968	552193	65.58%	8.036%
7	7.906	2025	2042	2059	rBV	492801	842015	100.00%	12.254%
8	8.260	2141	2152	2166	rVV8	4319	11142	1.32%	0.162%
9	9.423	2496	2514	2535	rBV	424575	694047	82.43%	10.100%
10	10.491	2837	2846	2854	rBV2	10904	16168	1.92%	0.235%
11	10.639	2880	2892	2908	rBV	351847	562880	66.85%	8.192%
12	11.427	3128	3137	3147	rBV2	7834	11759	1.40%	0.171%
13	11.616	3185	3196	3201	rBV3	5936	9937	1.18%	0.145%
14	11.648	3201	3206	3216	rVB2	12445	17859	2.12%	0.260%
15	11.819	3247	3259	3274	rBV	421917	667454	79.27%	9.713%
16	12.015	3308	3320	3329	rVB2	12434	18634	2.21%	0.271%
17	12.105	3329	3348	3359	rVV3	59460	93517	11.11%	1.361%
18	12.189	3362	3374	3391	rVB9	4708	10885	1.29%	0.158%
19	12.308	3399	3411	3422	rBV2	25060	39457	4.69%	0.574%
20	12.578	3485	3495	3501	rBV	35241	52034	6.18%	0.757%
21	12.619	3501	3508	3520	rVV2	32228	52488	6.23%	0.764%
22	12.690	3520	3530	3547	rVV	121609	207457	24.64%	3.019%
23	12.767	3547	3554	3563	rVB4	6228	9029	1.07%	0.131%
24	12.873	3579	3587	3601	rVB2	21241	33853	4.02%	0.493%
25	12.980	3601	3620	3632	rBV	83893	144672	17.18%	2.105%
26	13.121	3653	3664	3676	rVB6	10353	18258	2.17%	0.266%
27	13.259	3698	3707	3715	rVV3	23259	37816	4.49%	0.550%
28	13.301	3715	3720	3726	rVV2	11285	16352	1.94%	0.238%
29	13.336	3726	3731	3740	rVB5	6600	9582	1.14%	0.139%
30	13.462	3759	3770	3786	rVB3	156600	266249	31.62%	3.875%
31	13.536	3786	3793	3799	rBV3	7181	9646	1.15%	0.140%
32	13.632	3817	3823	3832	rVB6	8495	13579	1.61%	0.198%
33	13.790	3863	3872	3880	rBV4	10793	18079	2.15%	0.263%
34	13.844	3880	3889	3897	rVV4	39059	62339	7.40%	0.907%
35	13.893	3897	3904	3905	rVV	13714	15847	1.88%	0.231%
36	13.918	3906	3912	3915	rVV	27018	39180	4.65%	0.570%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

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\82U050621W.M
Title : SW846 8260

37	13.947	3916	3921	3944	rVB4	51545	107218	12.73%	1.560%
38	14.066	3944	3958	3966	rBV4	7918	13959	1.66%	0.203%
39	14.124	3967	3976	3986	rVB6	12698	20293	2.41%	0.295%
40	14.198	3987	3999	4012	rBV	90880	142142	16.88%	2.069%
41	14.295	4012	4029	4040	rBV	80187	142100	16.88%	2.068%
42	14.353	4040	4047	4058	rVB5	13213	19138	2.27%	0.279%
43	14.491	4079	4090	4098	rBV2	24786	41316	4.91%	0.601%
44	14.674	4137	4147	4172	rVB4	29965	67338	8.00%	0.980%
45	14.816	4182	4191	4201	rBV2	39319	60278	7.16%	0.877%
46	14.886	4204	4213	4226	rVV3	48690	85732	10.18%	1.248%
47	15.025	4245	4256	4267	rBV2	79413	130738	15.53%	1.903%
48	15.208	4306	4313	4315	rBV6	9856	11371	1.35%	0.165%
49	15.282	4328	4336	4346	rVB6	15357	25945	3.08%	0.378%
50	15.433	4372	4383	4391	rVV4	16632	35107	4.17%	0.511%
51	15.478	4391	4397	4402	rVV7	10997	18184	2.16%	0.265%
52	15.513	4402	4408	4418	rVB7	14018	23033	2.74%	0.335%
53	15.603	4421	4436	4445	rVB8	8529	18089	2.15%	0.263%
54	15.931	4526	4538	4555	rBV5	21156	48278	5.73%	0.703%
55	16.021	4555	4566	4580	rVB7	13113	24476	2.91%	0.356%
56	16.156	4601	4608	4613	rBV7	7591	11847	1.41%	0.172%
57	16.192	4613	4619	4628	rVB	15814	23939	2.84%	0.348%
58	16.269	4628	4643	4652	rBV2	23377	42799	5.08%	0.623%
59	16.417	4679	4689	4698	rVB3	28108	44264	5.26%	0.644%
60	16.536	4715	4726	4738	rBV6	7891	14961	1.78%	0.218%
61	16.623	4743	4753	4756	rVV4	16757	25784	3.06%	0.375%
62	16.655	4757	4763	4773	rVB3	26432	42463	5.04%	0.618%
63	16.793	4797	4806	4816	rBV3	24835	40565	4.82%	0.590%
64	17.105	4892	4903	4916	rBV3	8939	17325	2.06%	0.252%
65	17.552	5035	5042	5050	rVB5	5811	8449	1.00%	0.123%

Sum of corrected areas: 6871422

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
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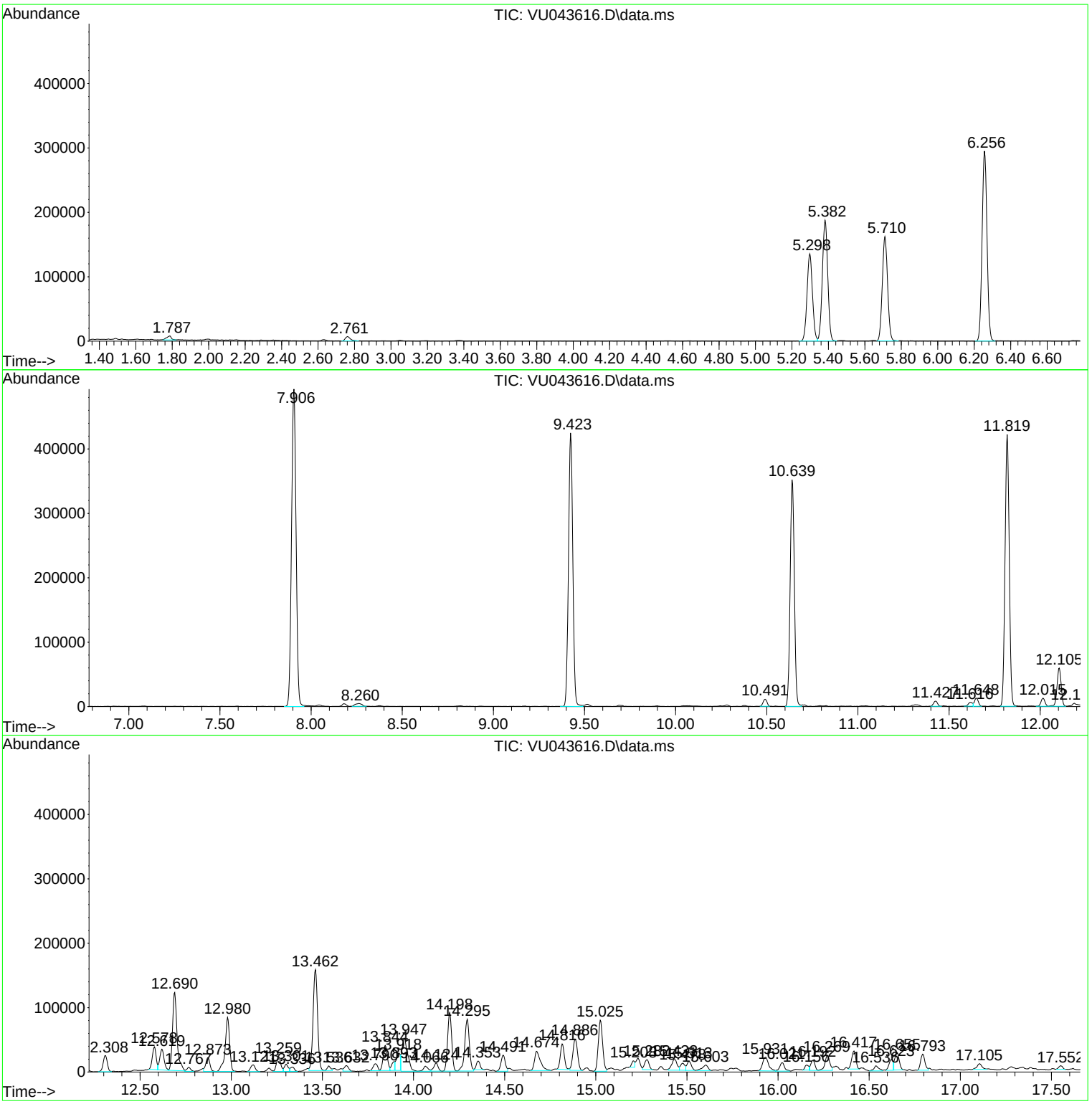
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C7-GW

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

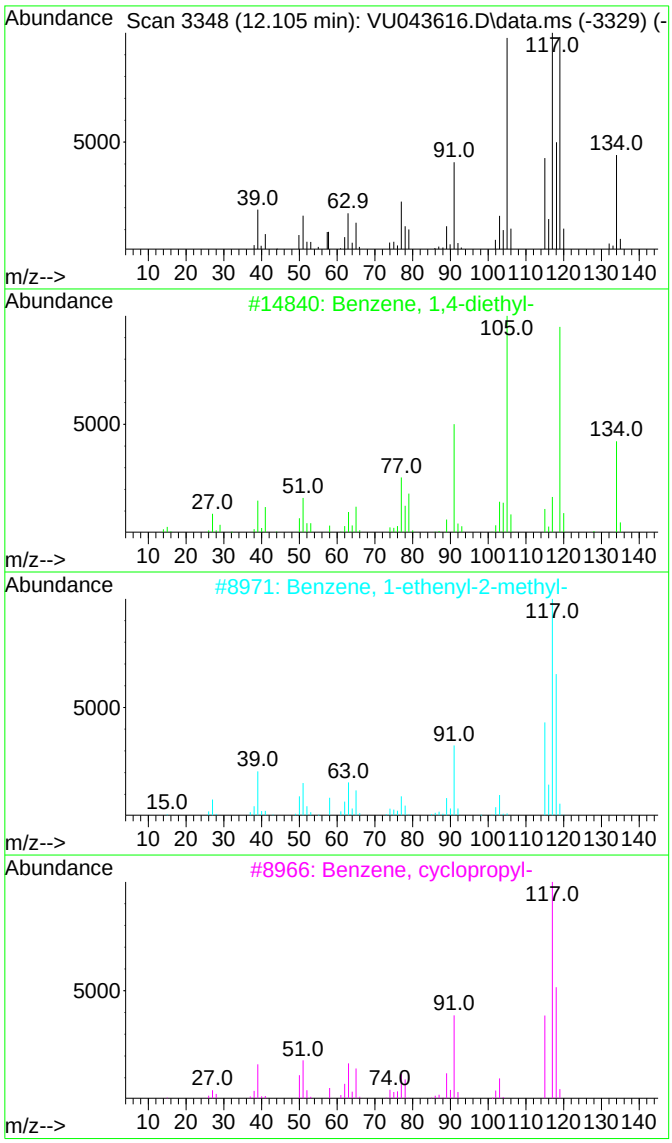
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

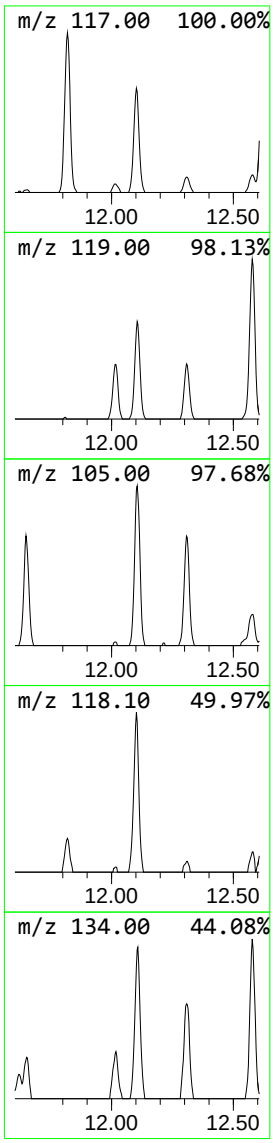
****

Peak Number 1 Benzene, 1,4-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.105	7.01 ug/l	93517	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	87
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	83
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	55
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

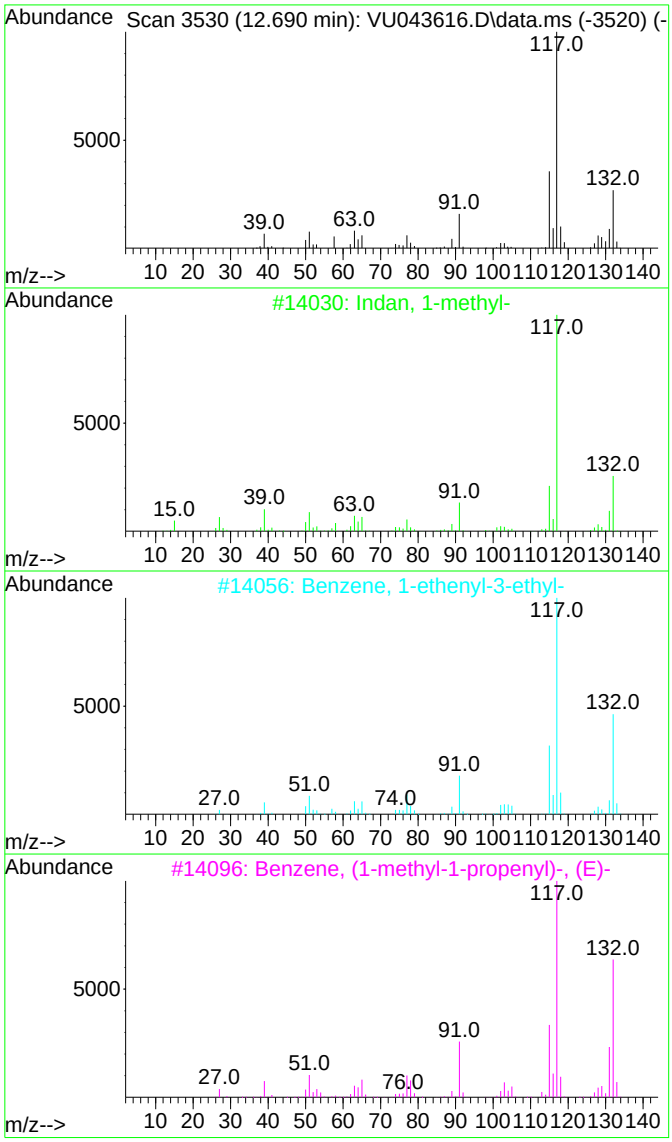
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

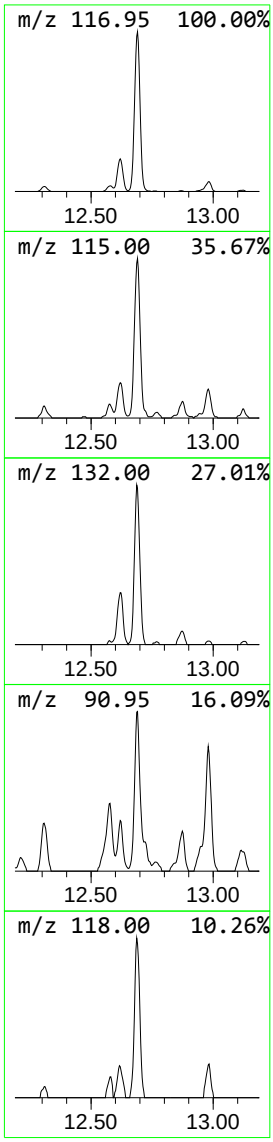
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.690	15.54 ug/l	207457	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	80
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

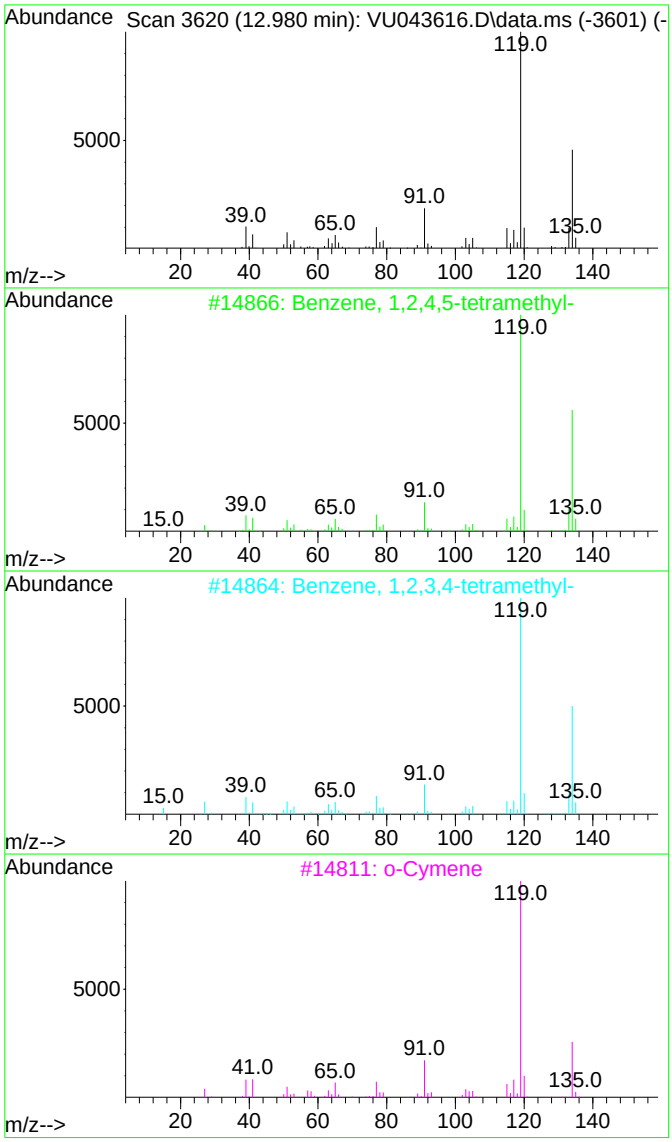
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

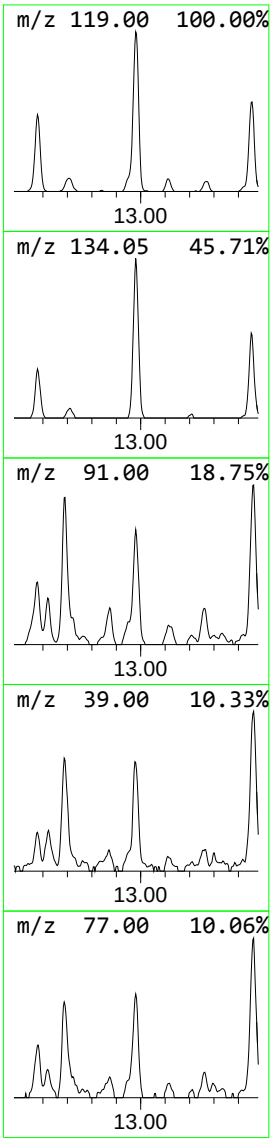
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	10.84 ug/l	144672	1,4-Dichlorobenzene-d4	11.819

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

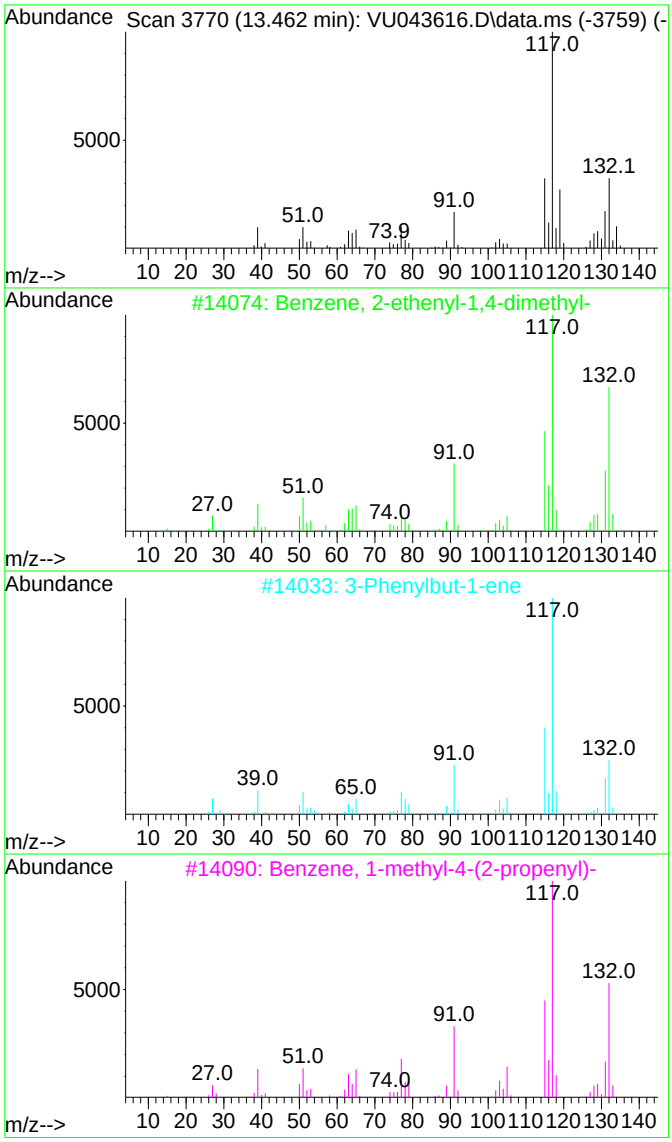
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

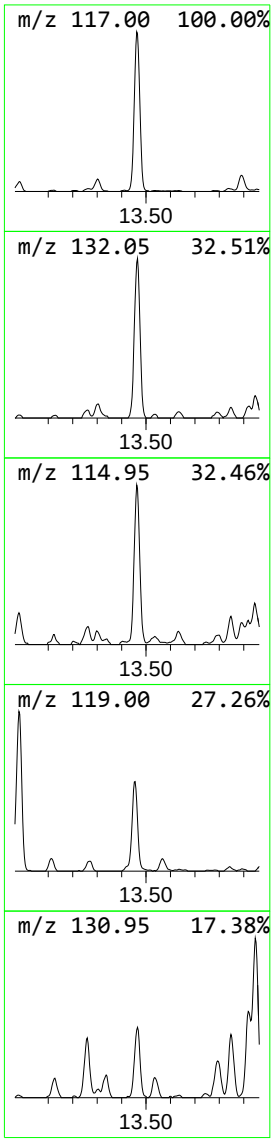
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.462	19.95 ug/l	266249	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	78
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

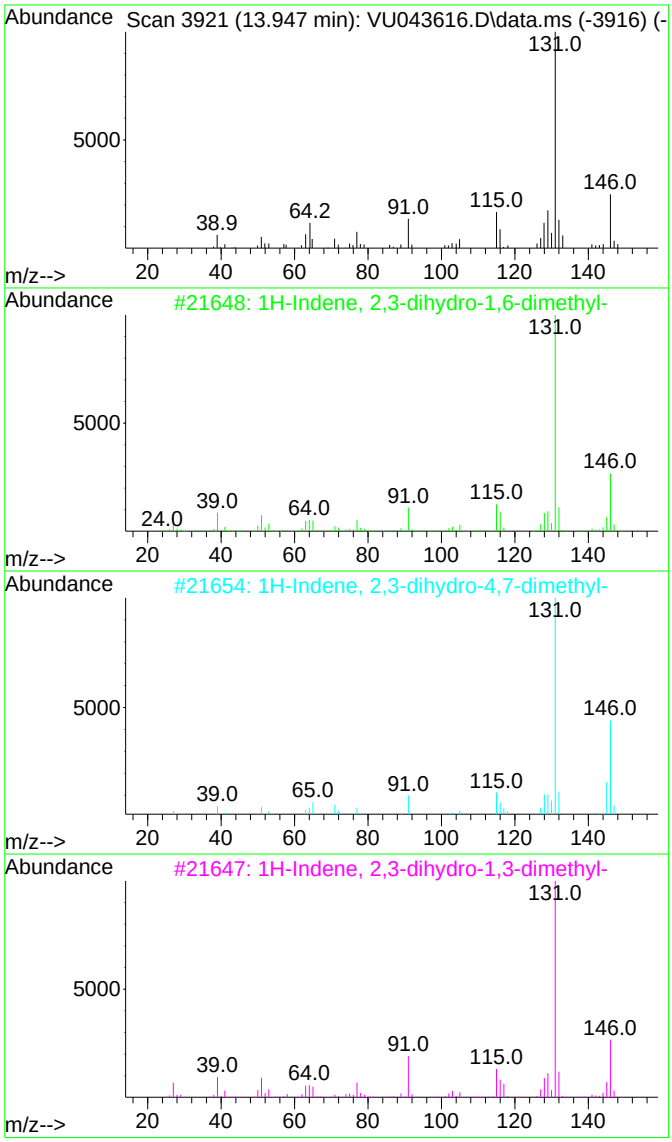
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

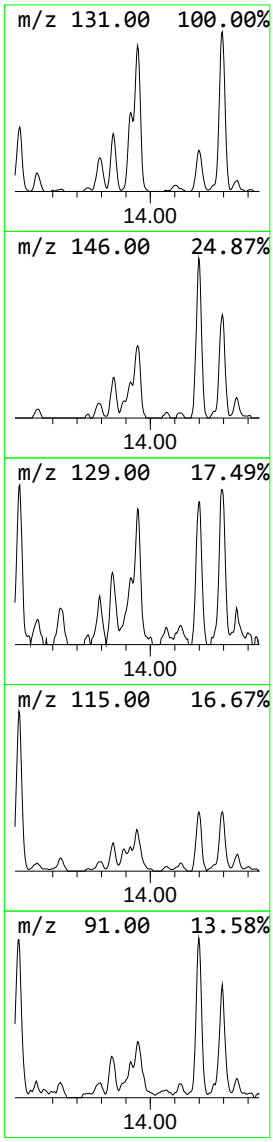
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.947	8.03 ug/l	107218	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90





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Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

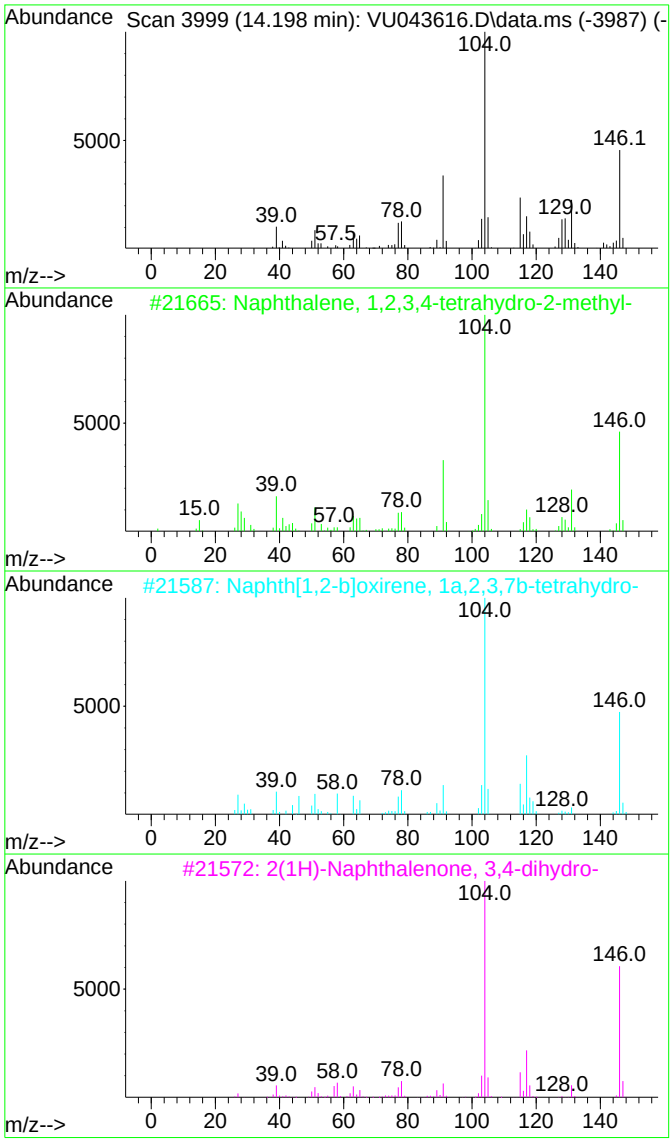
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

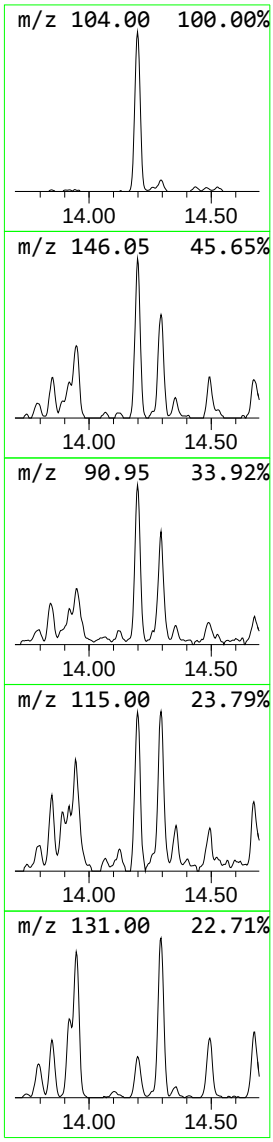
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.198	10.65 ug/l	142142	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	62
3	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	58
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	50
5	1,3-Propanediamine, 2-phenyl-	150	C9H14N2	055165-09-8	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C7-GW

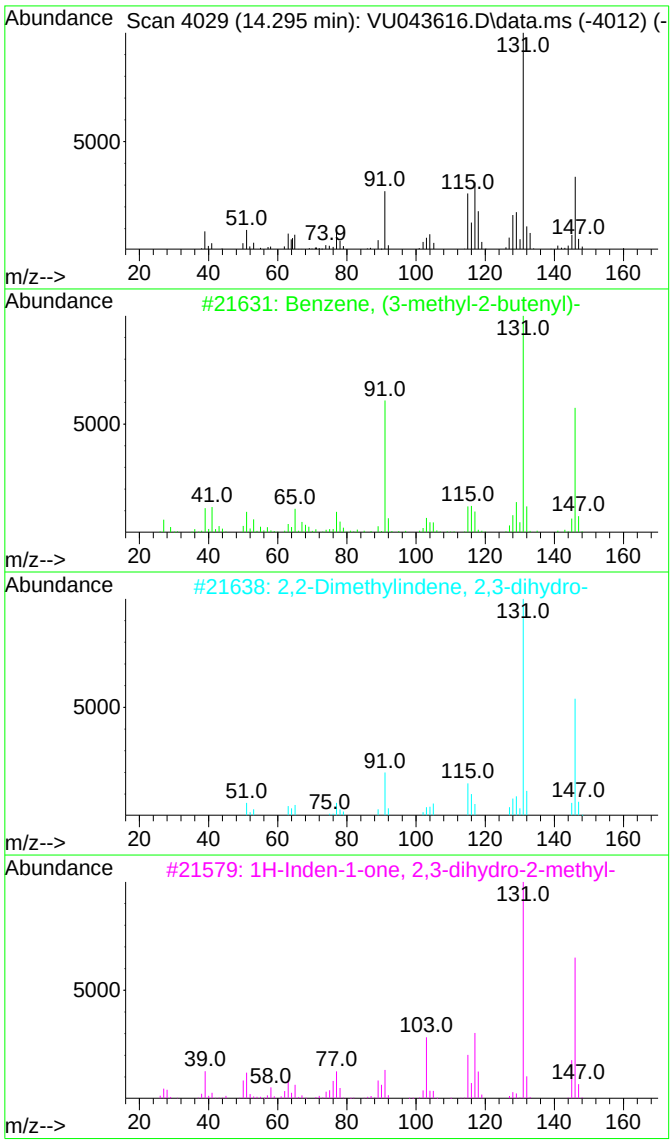
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

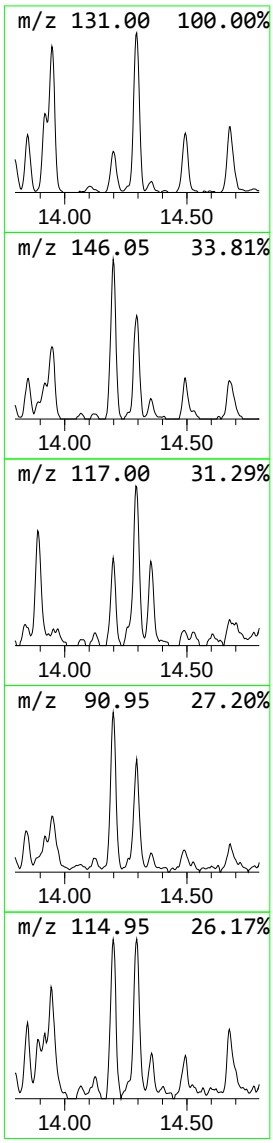
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, (3-methyl-2-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.295	10.64 ug/l	142100	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	87
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	87
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86





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Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

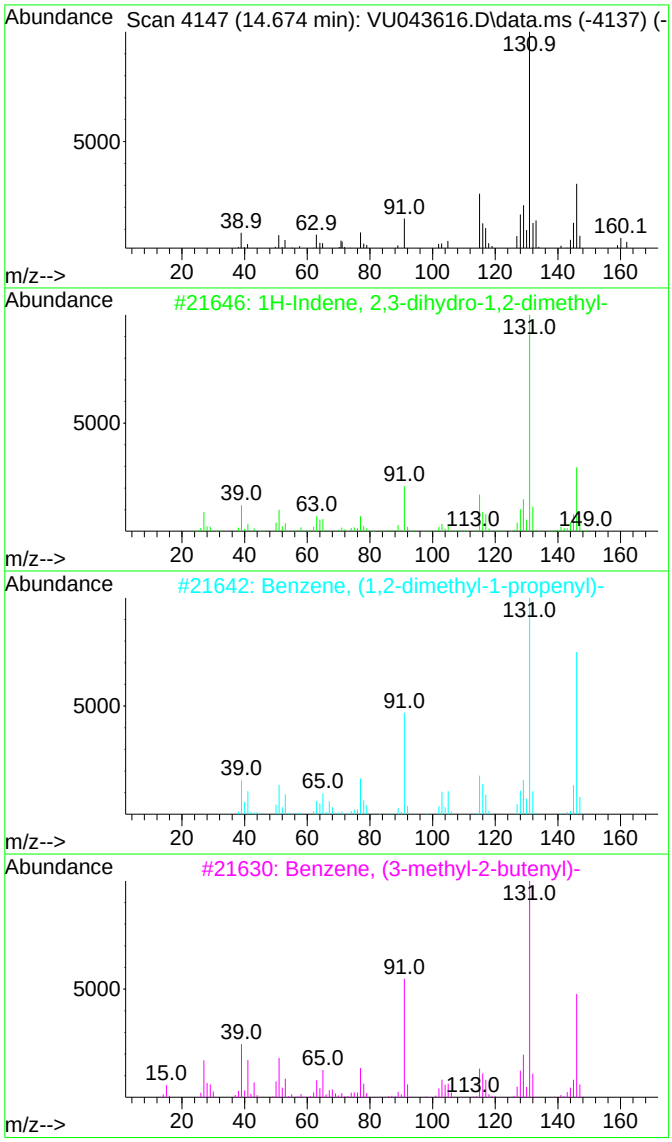
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

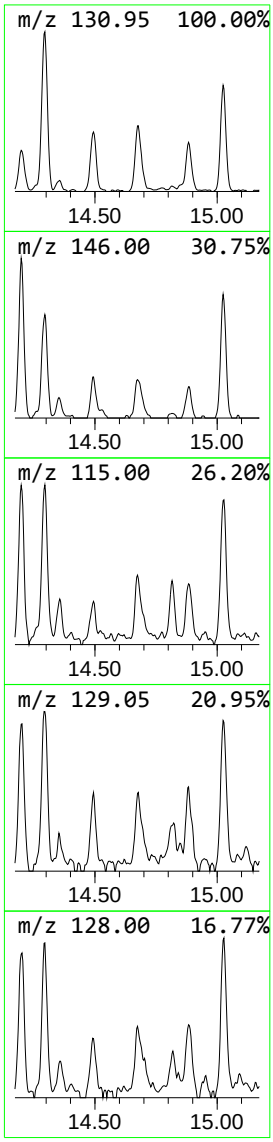
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.674	5.04 ug/l	67338	1,4-Dichlorobenzene-d4	11.819

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	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	87
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

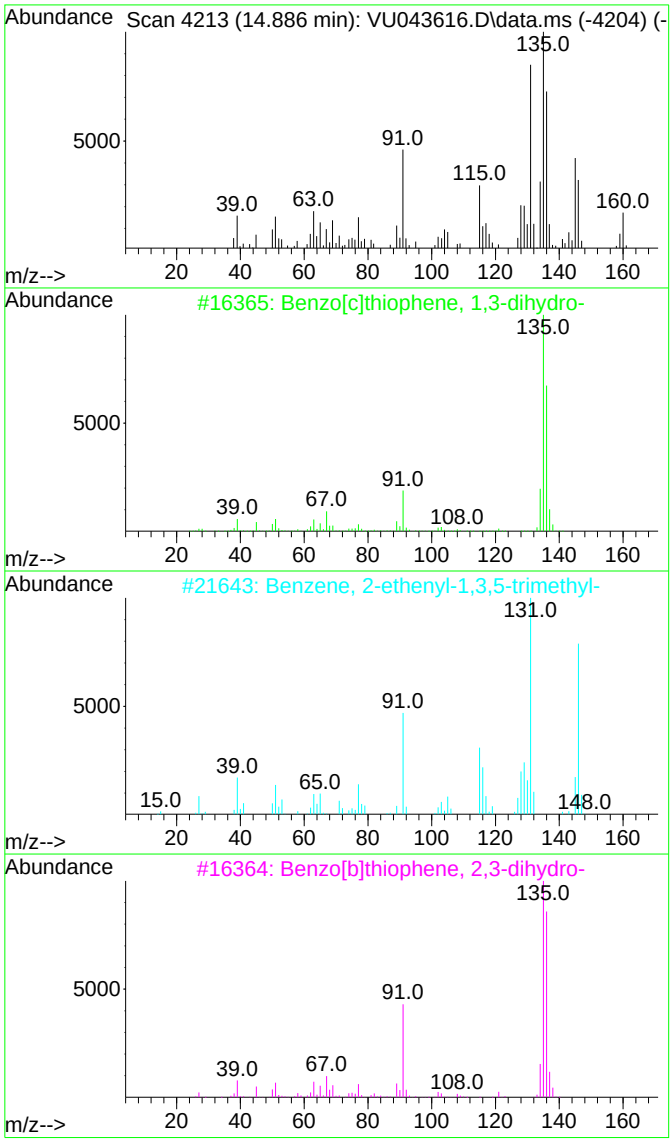
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

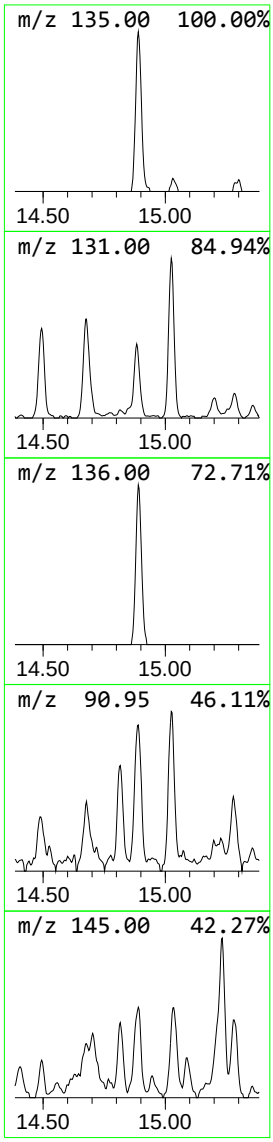
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzo[c]thiophene, 1,3-dihy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.886	6.42 ug/l	85732	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	50
2	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	42
3	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	42
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	42
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

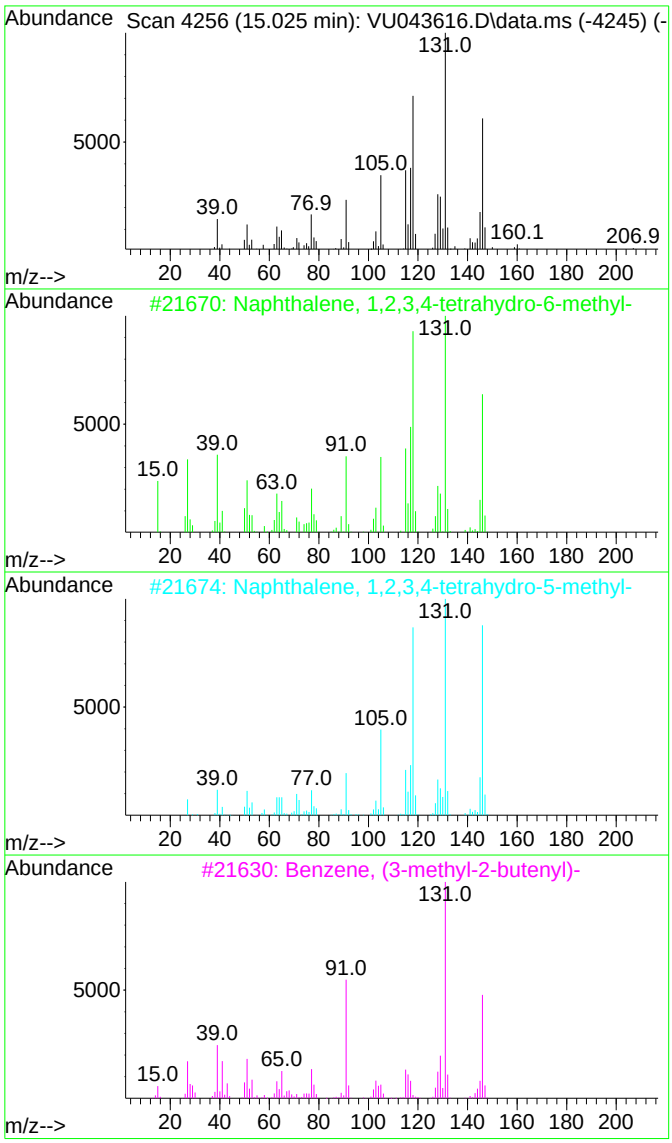
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

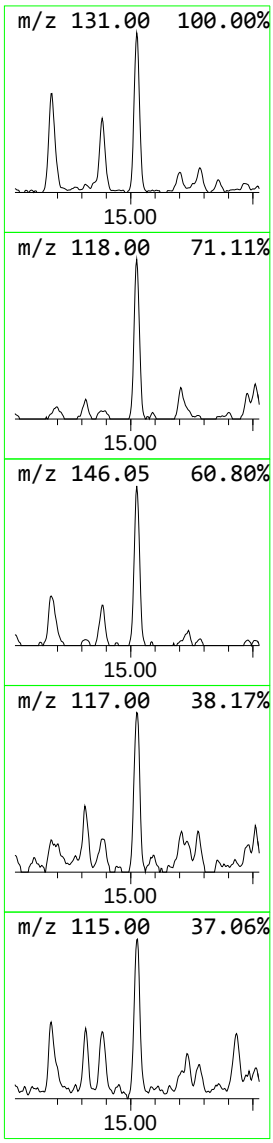
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.025	9.79 ug/l	130738	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
5	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043616.D
Acq On : 07 May 2021 16:40
Operator : SY/MD
Sample : M2304-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C7-GW

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.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
Benzene, 1,4-di...	12.105	7.0	ug/l	93517	4	11.819	667454	50.0
Indan, 1-methyl-	12.690	15.5	ug/l	207457	4	11.819	667454	50.0
Benzene, 1,2,4,...	12.980	10.8	ug/l	144672	4	11.819	667454	50.0
Benzene, 2-ethe...	13.462	19.9	ug/l	266249	4	11.819	667454	50.0
1H-Indene, 2,3-...	13.947	8.0	ug/l	107218	4	11.819	667454	50.0
Naphthalene, 1,...	14.198	10.7	ug/l	142142	4	11.819	667454	50.0
Benzene, (3-met...	14.295	10.6	ug/l	142100	4	11.819	667454	50.0
1H-Indene, 2,3-...	14.674	5.0	ug/l	67338	4	11.819	667454	50.0
Benzo[c]thiophe...	14.886	6.4	ug/l	85732	4	11.819	667454	50.0
Naphthalene, 1,...	15.025	9.8	ug/l	130738	4	11.819	667454	50.0