LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-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: VU043612.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.298	1214	1231	1244	rBV	132846	281070	34.04%	3.680%
2	5.382	1244	1257	1275	rVB	179972	364925	44.19%	4.777%
3	5.710	1342	1359	1382	rBV	158794	324735	39.32%	4.251%
4	6.256	1512	1529	1548	rBV	288494	537899	65.14%	7.042%
5	7.906	2024	2042	2059	rBV	485035	825776	100.00%	10.810%
6	9.423	2501	2514	2535	rBV	421325	675485	81.80%	8.843%
7	10.491	2833	2846	2856	rBV	22797	34046	4.12%	0.446%
8	10.639	2879	2892	2906	rBV	349727	544674	65.96%	7.130%
9	10.912	2968	2977	2985	rBV	6743	9887	1.20%	0.129%
10	11.427	3129	3137	3146	rVB3	5825	8379	1.01%	0.110%
11	11.619	3185	3197	3199	rBV2	7583	10676	1.29%	0.140%
12	11.648	3199	3206	3218	rVB	29021	44228	5.36%	0.579%
13	11.819	3246	3259	3274	rBV	400527	634176	76.80%	8.302%
14	12.015	3310	3320	3329	rBV2	11093	15453	1.87%	0.202%
15	12.105	3338	3348	3359	rVB2	51447	79949	9.68%	1.047%
16	12.214	3364	3382	3390	rBV7	7795	18713	2.27%	0.245%
17	12.308	3398	3411	3424	rVB2	26602	39801	4.82%	0.521%
18	12.578	3476	3495	3501	rBV2	36406	58220	7.05%	0.762%
19	12.619	3501	3508	3519	rVV2	41106	65399	7.92%	0.856%
20	12.690	3519	3530	3549	rVV2	124368	237719	28.79%	3.112%
21	12.767	3549	3554	3564	rVV4	5974	8351	1.01%	0.109%
22	12.870	3579	3586	3602	rVB2	16071	26234	3.18%	0.343%
23	12.980	3602	3620	3634	rBV	94647	157912	19.12%	2.067%
24	13.115	3650	3662	3674	rVB5	17281	28830	3.49%	0.377%
25	13.205	3682	3690	3698	rBV3	6590	9992	1.21%	0.131%
26	13.259	3698	3707	3713	rBV3	30578	48187	5.84%	0.631%
27	13.301	3713	3720	3729	rVB	43522	61835	7.49%	0.810%
28	13.423	3745	3758	3760	rBV6	6625	10099	1.22%	0.132%
29	13.462	3760	3770	3785	rVB3	120946	206074	24.96%	2.698%
30	13.539	3785	3794	3798	rBV4	8701	12424	1.50%	0.163%
31	13.632	3817	3823	3835	rVB6	10683	17733	2.15%	0.232%
32	13.745	3848	3858	3863	rBV2	9472	13969	1.69%	0.183%
33	13.793	3863	3873	3880	rVV2	38196	66258	8.02%	0.867%
34	13.844	3880	3889	3898	rVV4	57501	94145	11.40%	1.232%
35	13.918	3898	3912	3915	rVV4	36395	75006	9.08%	0.982%
36	13.947	3916	3921	3937	rVB2	67083	119063	14.42%	1.559%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-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	14.066	3944	3958	3967	rBV5	12247	23417	2.84%	0.307%
38	14.124	3967	3976	3988	rVV3	12618	24234	2.93%	0.317%
39	14.198	3988	3999	4011	rVB	123471	192995	23.37%	2.527%
40	14.295	4011	4029	4040	rBV	121531	213067	25.80%	2.789%
41	14.356	4040	4048	4057	rVV4	21593	37536	4.55%	0.491%
42	14.491	4078	4090	4099	rBV	54754	89110	10.79%	1.167%
43	14.677	4136	4148	4171	rBV2	65388	149329	18.08%	1.955%
44	14.815	4182	4191	4202	rBV2	44270	68310	8.27%	0.894%
45	14.883	4203	4212	4224	rVB2	65812	109970	13.32%	1.440%
46	14.947	4224	4232	4243	rVB5	10265	16554	2.00%	0.217%
47	15.024	4244	4256	4269	rBV3	120876	207811	25.17%	2.721%
48	15.089	4271	4276	4290	rVB4	8323	13668	1.66%	0.179%
49	15.205	4305	4312	4317	rBV3	30625	46296	5.61%	0.606%
50	15.282	4328	4336	4348	rVB4	31338	54287	6.57%	0.711%
51	15.356	4351	4359	4367	rBV6	10576	16413	1.99%	0.215%
52	15.433	4372	4383	4389	rVV4	14919	30099	3.64%	0.394%
53	15.475	4389	4396	4403	rVV4	20492	37642	4.56%	0.493%
54	15.510	4403	4407	4418	rVB7	17616	25316	3.07%	0.331%
55	15.600	4422	4435	4446	rBV6	12892	29257	3.54%	0.383%
56	15.774	4475	4489	4497	rVB7	7184	13715	1.66%	0.180%
57	15.931	4525	4538	4555	rVB5	38311	86381	10.46%	1.131%
58	16.021	4555	4566	4579	rBV8	13152	23263	2.82%	0.305%
59	16.156	4600	4608	4613	rBV7	9376	13512	1.64%	0.177%
60	16.192	4613	4619	4628	rVB3	17207	24549	2.97%	0.321%
61	16.272	4629	4644	4653	rBV2	37848	68512	8.30%	0.897%
62	16.320	4653	4659	4671	rVB7	10926	20194	2.45%	0.264%
63	16.417	4679	4689	4699	rVB3	50892	78877	9.55%	1.033%
64	16.455	4699	4701	4717	rVB9	5272	9095	1.10%	0.119%
65	16.539	4717	4727	4735	rBV10	7343	14392	1.74%	0.188%
66	16.622	4740	4753	4756	rBV2	18701	29705	3.60%	0.389%
67	16.651	4756	4762	4772	rVB3	30287	49815	6.03%	0.652%
68	16.793	4797	4806	4816	rBV3	23302	39189	4.75%	0.513%
69	17.105	4890	4903	4917	rBV3	7150	14832	1.80%	0.194%

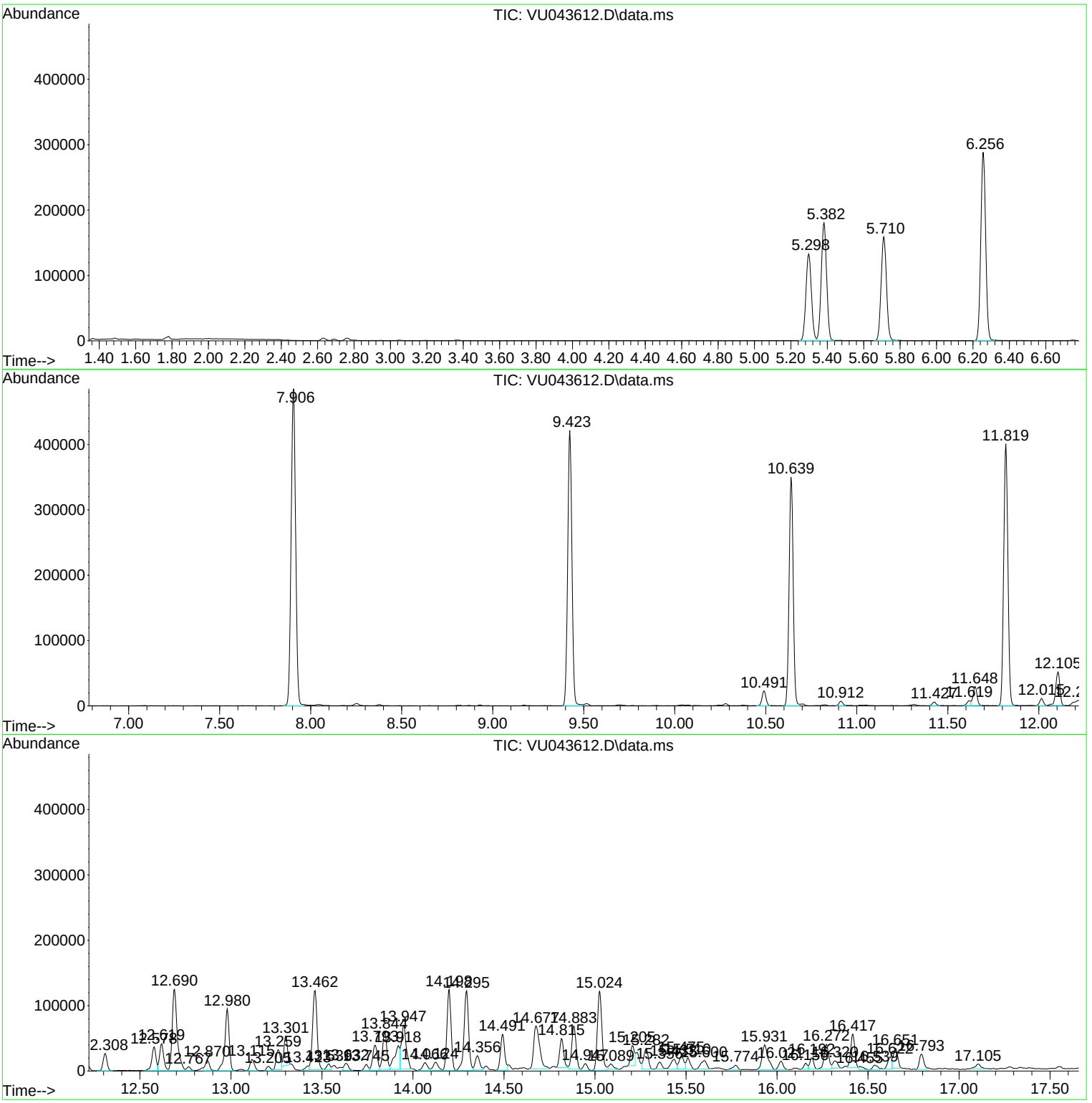
Sum of corrected areas: 7638664

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C2-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 : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

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

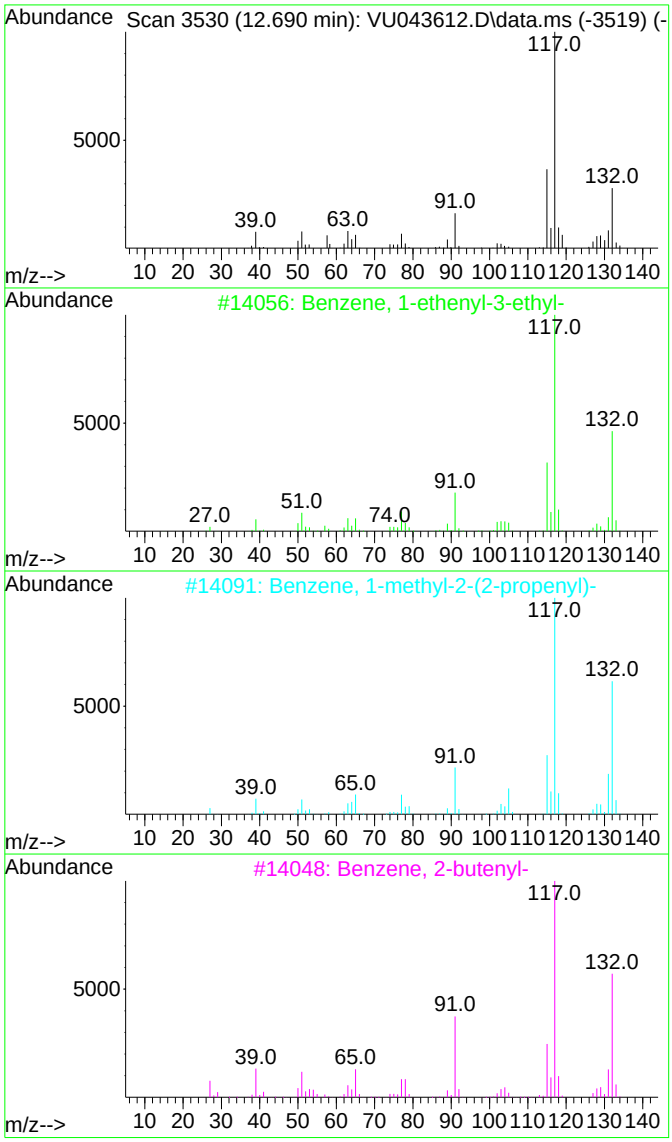
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

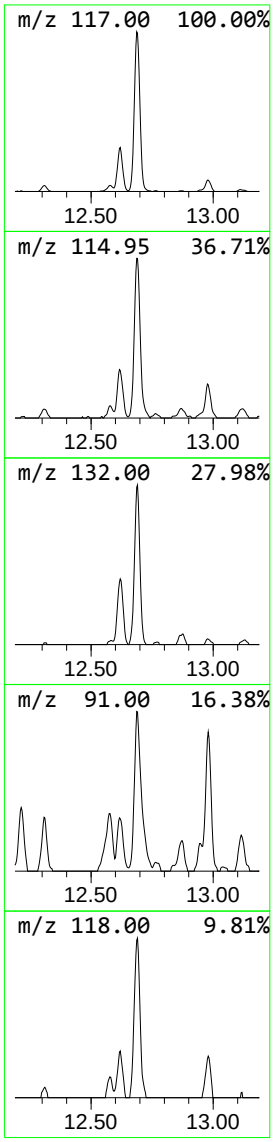
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.690	18.74 ug/l	237719	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

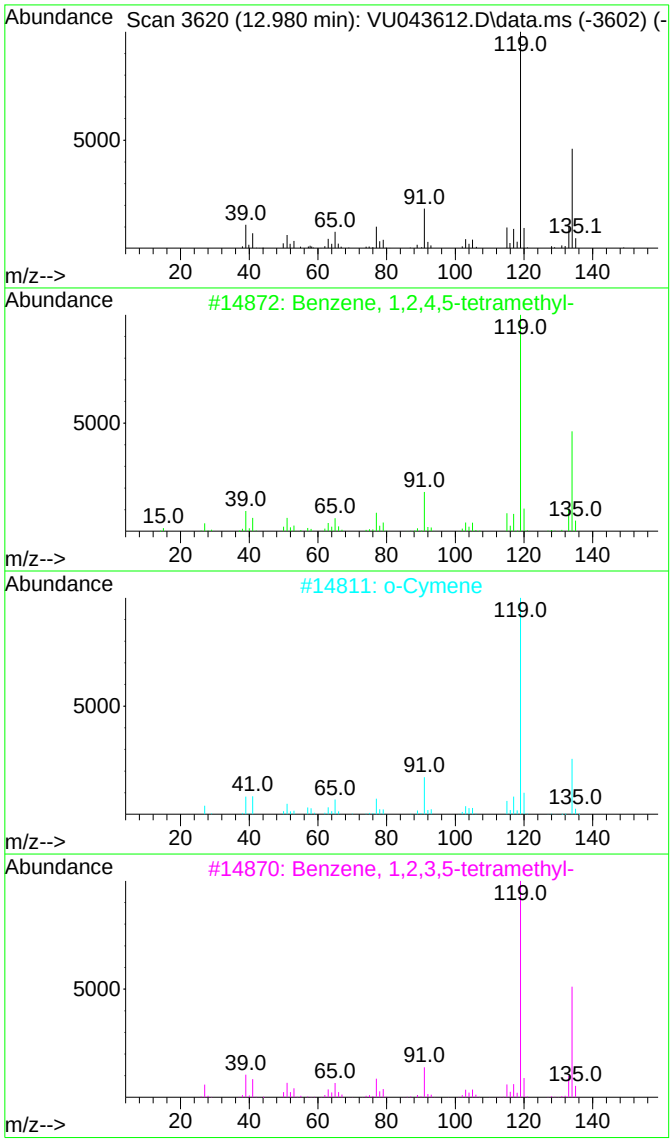
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

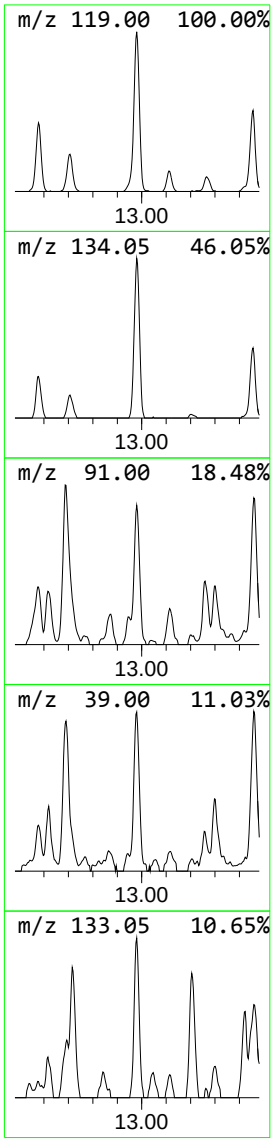
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	12.45 ug/l	157912	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	97
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	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 : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

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

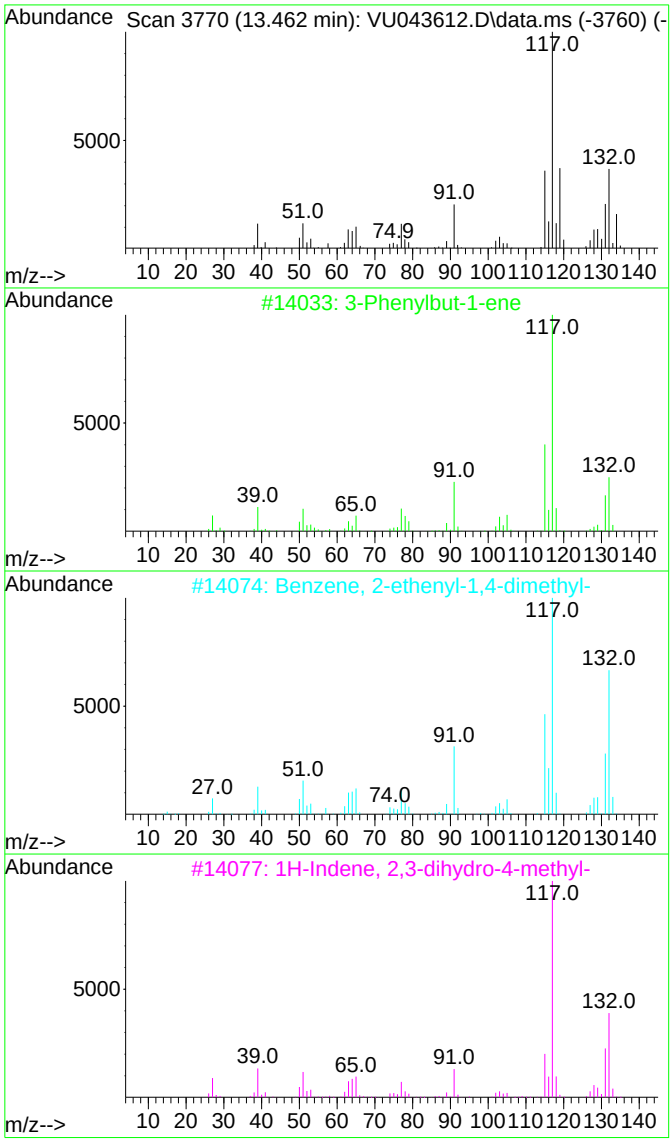
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

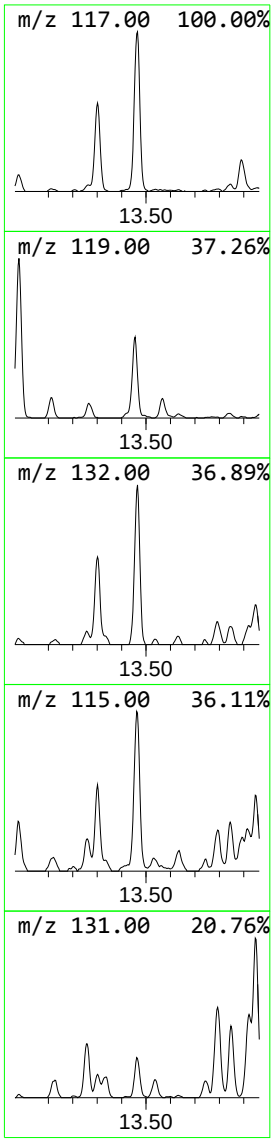
****

Peak Number 3 3-Phenylbut-1-ene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.462	16.25 ug/l	206074	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	74
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

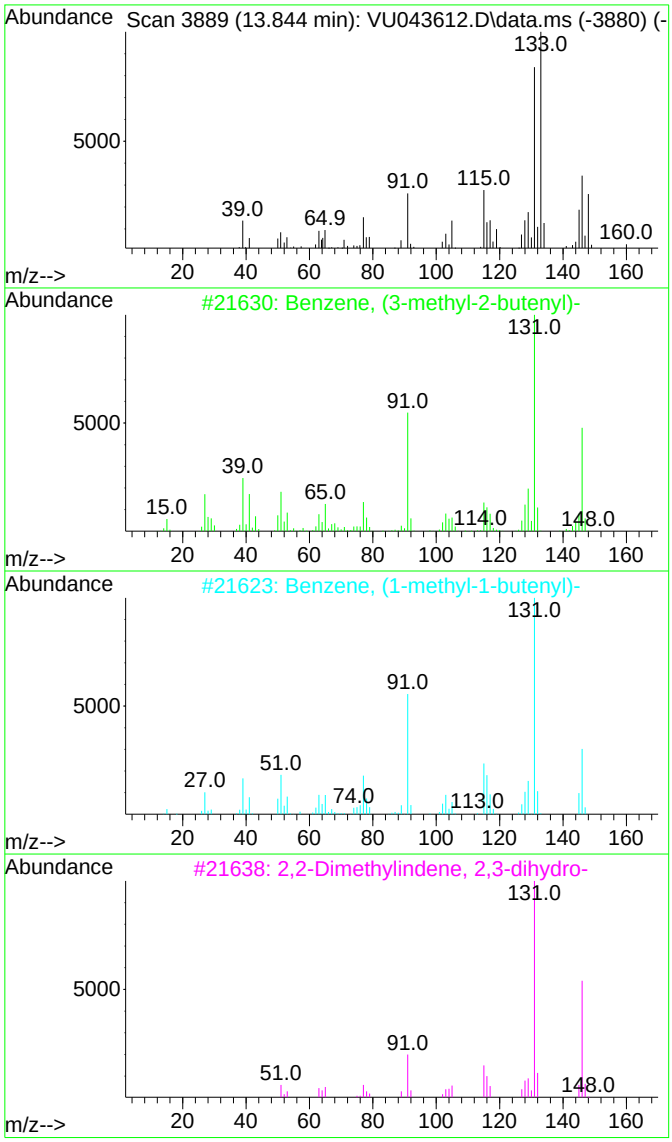
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

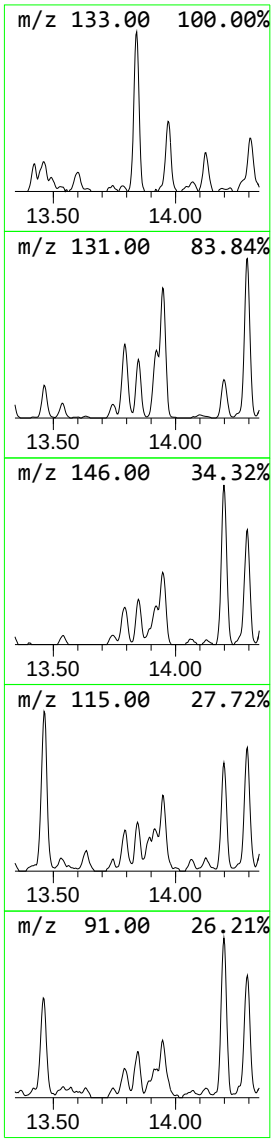
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, (3-methyl-2-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.844	7.42 ug/l	94145	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	64
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	50
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	50
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

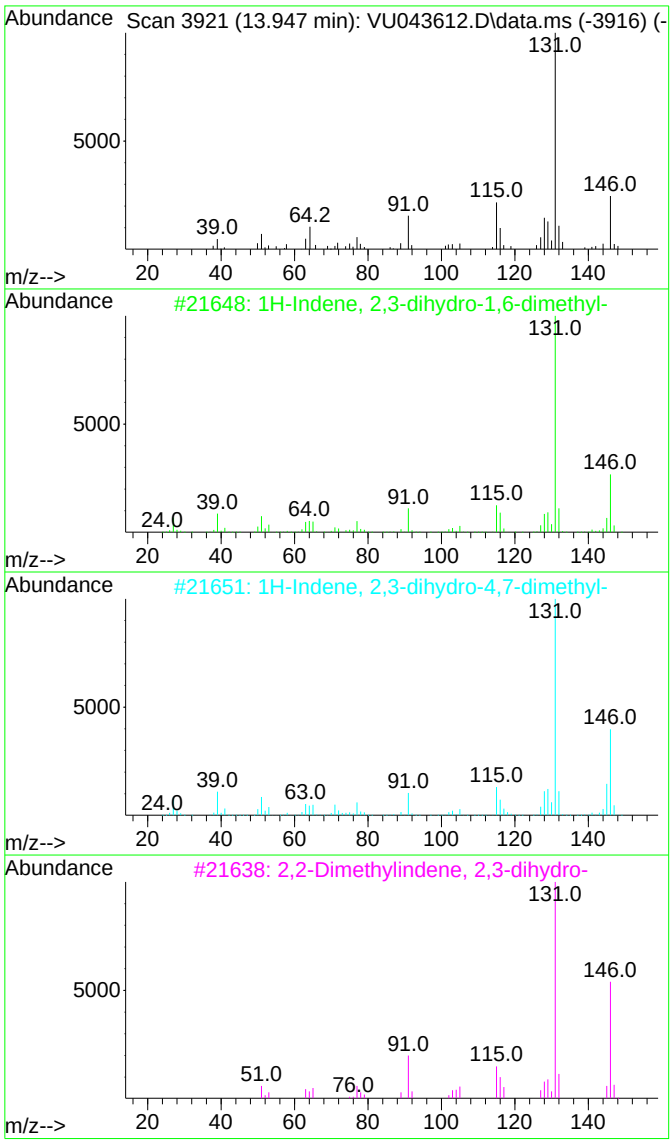
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

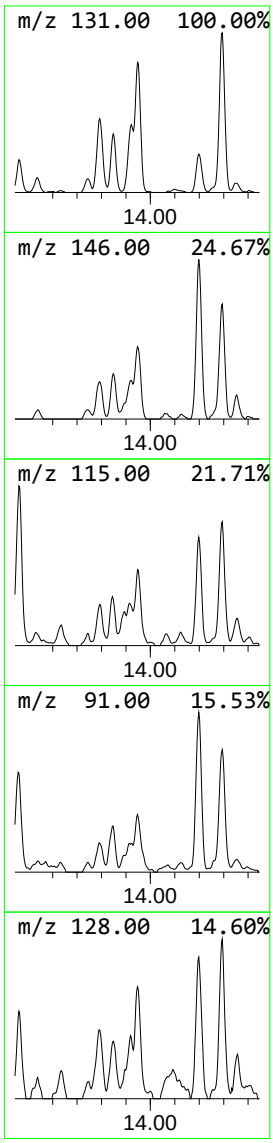
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.947	9.39 ug/l	119063	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	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

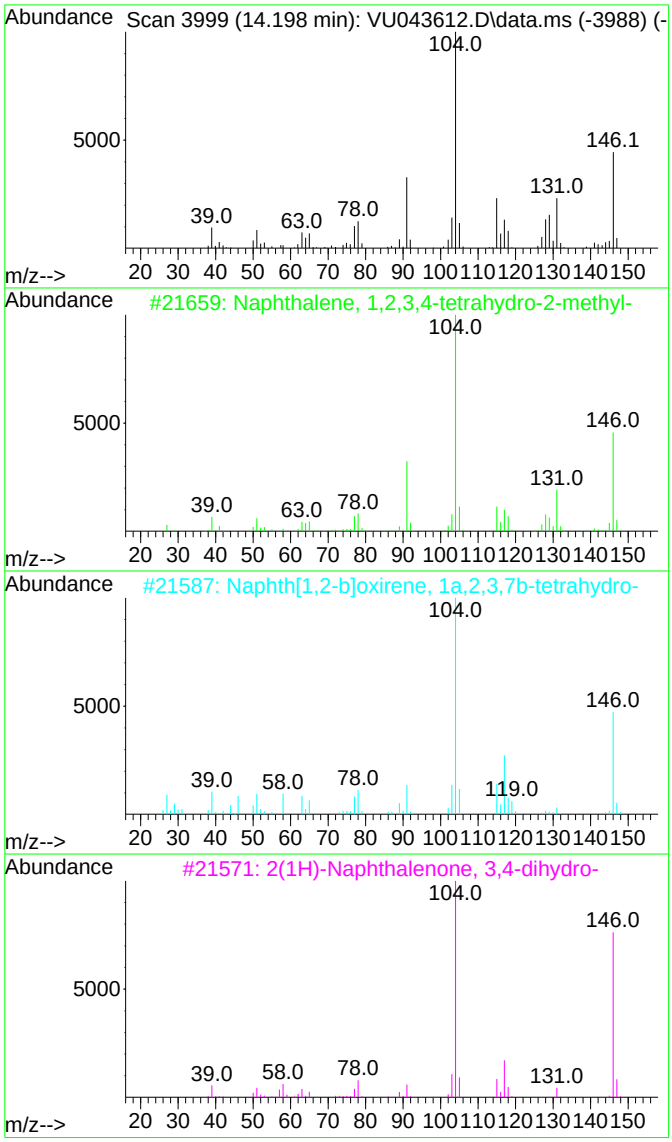
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

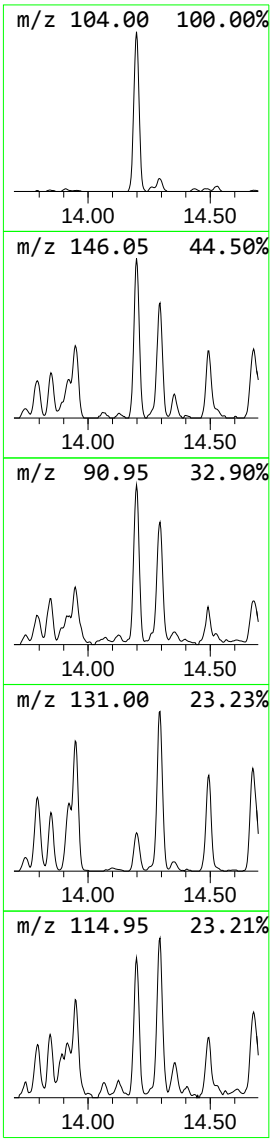
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.198	15.22 ug/l	192995	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	70
3	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	58
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	50
5	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

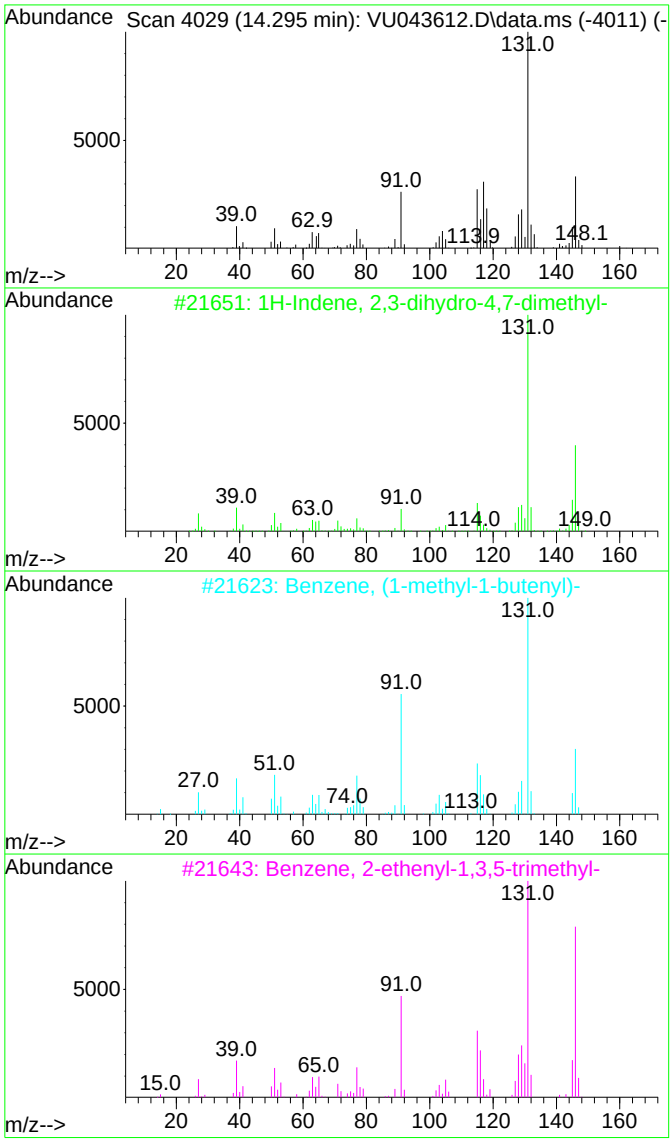
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

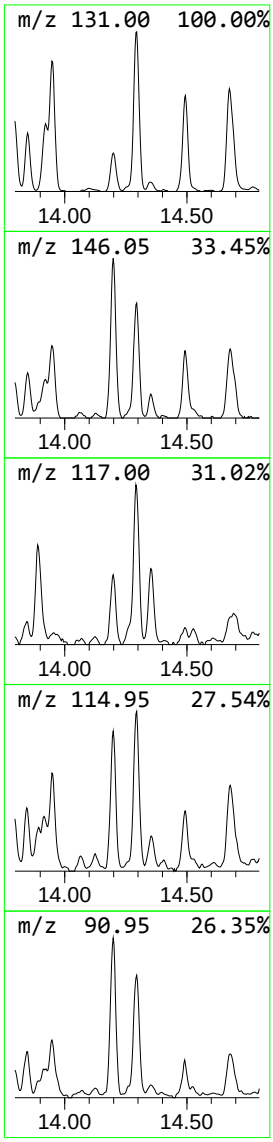
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.295	16.80 ug/l	213067	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	93
3	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	91
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	90
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

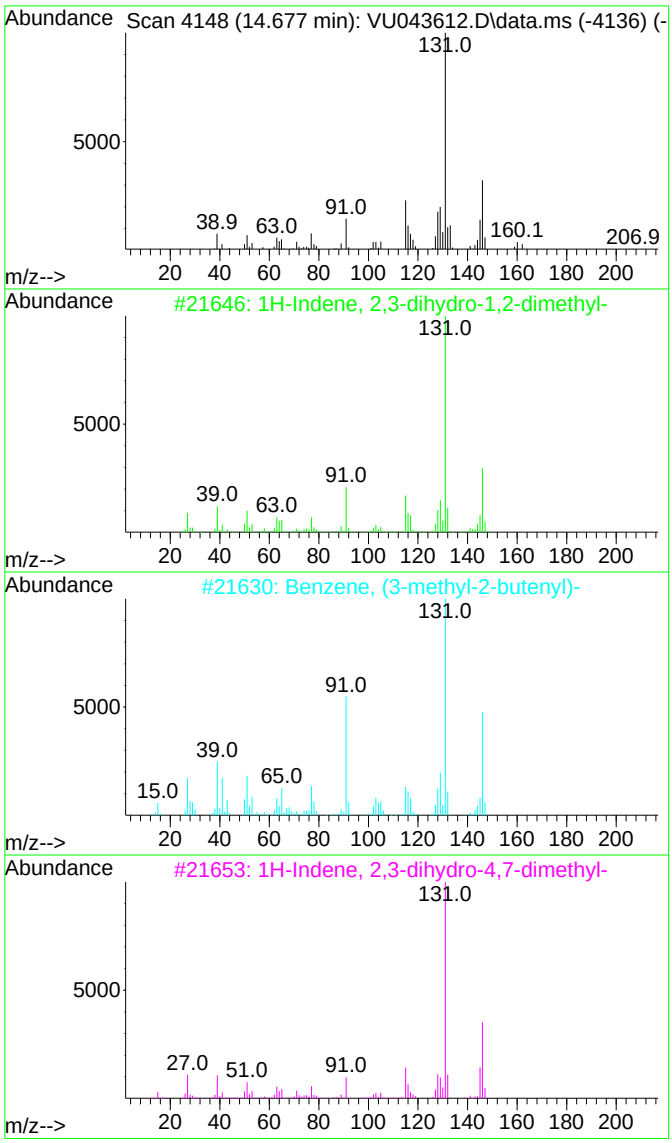
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

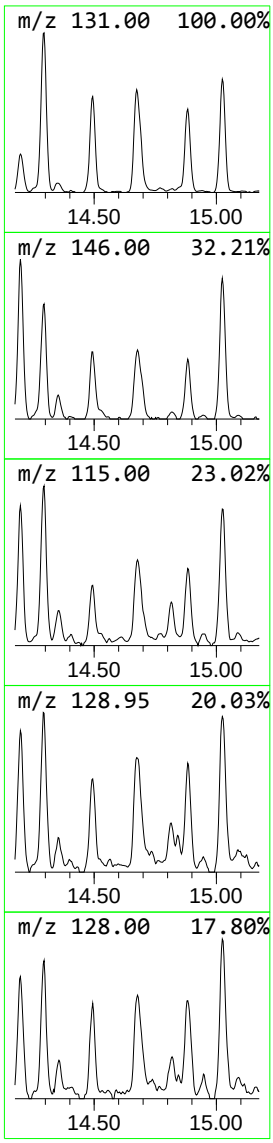
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.677	11.77 ug/l	149329	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	93
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	87
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	83
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

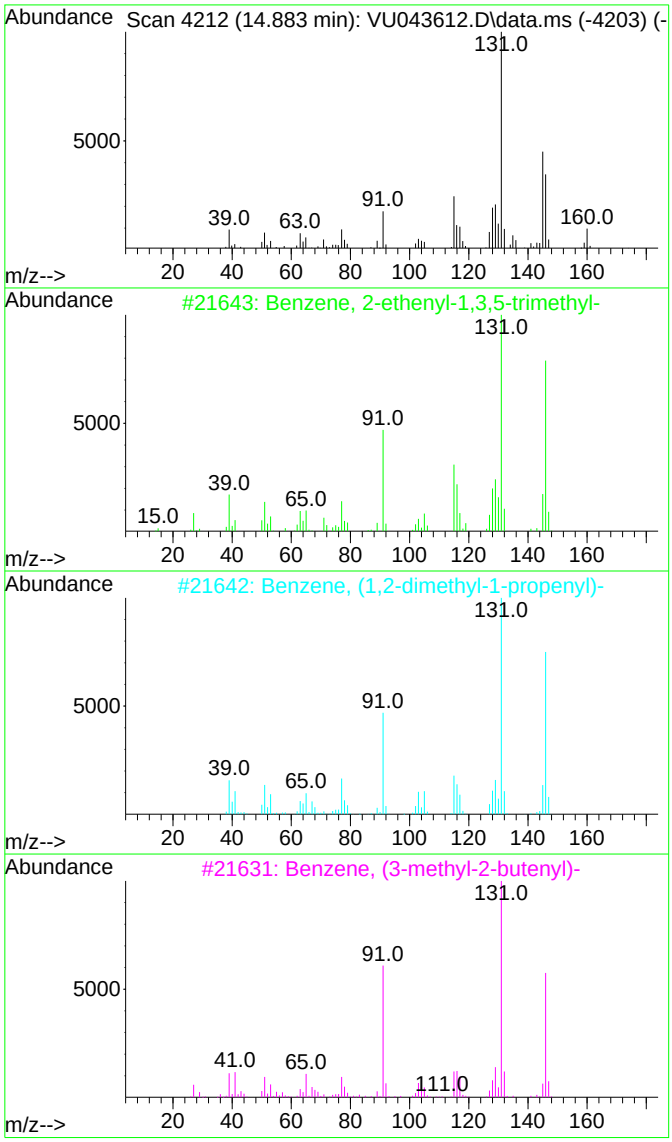
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

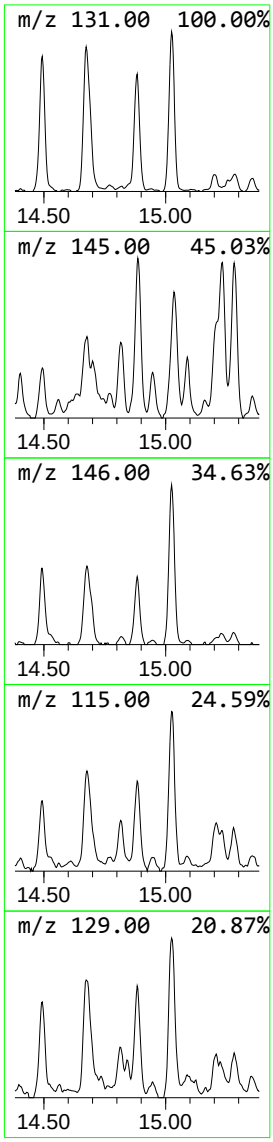
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.883	8.67 ug/l	109970	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	76
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	58
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	52





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-GW

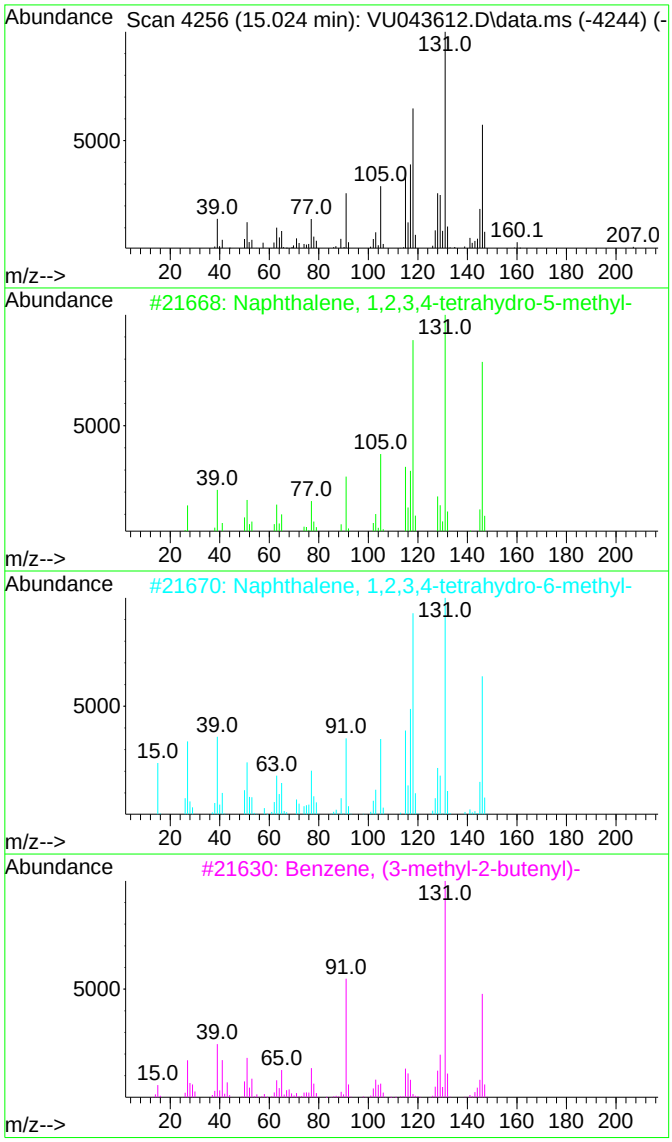
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

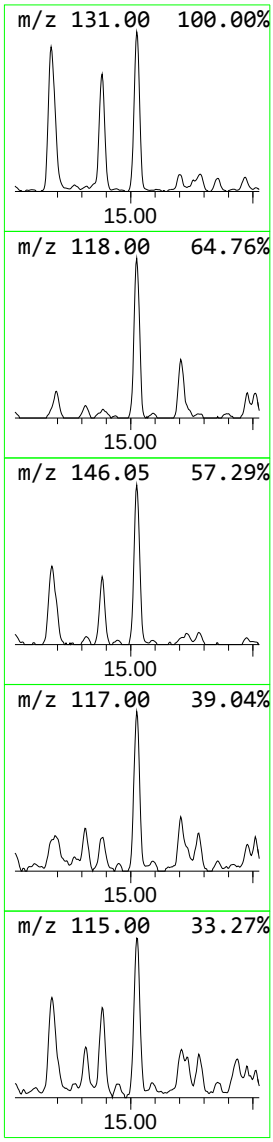
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.025	16.38 ug/l	207811	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
4	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	64
5	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	62





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043612.D
Acq On : 07 May 2021 15:07
Operator : SY/MD
Sample : M2304-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : OWL-C2-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-ethe...	12.690	18.7	ug/l	237719	4	11.819	634176	50.0
Benzene, 1,2,4,...	12.980	12.4	ug/l	157912	4	11.819	634176	50.0
3-Phenylbut-1-ene	13.462	16.3	ug/l	206074	4	11.819	634176	50.0
Benzene, (3-met...	13.844	7.4	ug/l	94145	4	11.819	634176	50.0
1H-Indene, 2,3-...	13.947	9.4	ug/l	119063	4	11.819	634176	50.0
Naphthalene, 1,...	14.198	15.2	ug/l	192995	4	11.819	634176	50.0
1H-Indene, 2,3-...	14.295	16.8	ug/l	213067	4	11.819	634176	50.0
1H-Indene, 2,3-...	14.677	11.8	ug/l	149329	4	11.819	634176	50.0
Benzene, 2-ethe...	14.883	8.7	ug/l	109970	4	11.819	634176	50.0
Naphthalene, 1,...	15.025	16.4	ug/l	207811	4	11.819	634176	50.0