LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

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: VU043687.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.761	431	442	460	rBV	6884	13392	1.51%	0.132%
2	5.298	1215	1231	1244	rBV	138283	295101	33.17%	2.917%
3	5.382	1244	1257	1275	rVB	191220	389554	43.79%	3.851%
4	5.710	1344	1359	1377	rBV	170581	339078	38.12%	3.352%
5	6.256	1515	1529	1550	rBV	310965	579231	65.11%	5.726%
6	7.906	2017	2042	2060	rBV	520608	889600	100.00%	8.794%
7	8.250	2136	2149	2159	rVB4	8523	15172	1.71%	0.150%
8	9.423	2499	2514	2532	rBV	453012	729530	82.01%	7.212%
9	9.517	2535	2543	2553	rVB3	8290	13954	1.57%	0.138%
10	9.697	2589	2599	2611	rVB9	4560	9060	1.02%	0.090%
11	10.279	2774	2780	2791	rVB2	6157	9426	1.06%	0.093%
12	10.491	2835	2846	2856	rBV2	12355	18079	2.03%	0.179%
13	10.639	2879	2892	2905	rBV	377981	583438	65.58%	5.767%
14	11.427	3128	3137	3144	rBV	6668	9291	1.04%	0.092%
15	11.616	3186	3196	3200	rBV	19645	29141	3.28%	0.288%
16	11.648	3200	3206	3218	rVB	47743	73297	8.24%	0.725%
17	11.819	3246	3259	3276	rBV	447811	709853	79.79%	7.017%
18	12.015	3311	3320	3330	rBV2	14665	20766	2.33%	0.205%
19	12.105	3338	3348	3358	rVB2	45090	69466	7.81%	0.687%
20	12.189	3364	3374	3389	rVB	57692	91221	10.25%	0.902%
21	12.308	3399	3411	3423	rVB2	18411	27514	3.09%	0.272%
22	12.574	3475	3494	3499	rBV6	7879	22902	2.57%	0.226%
23	12.619	3499	3508	3519	rVV	80876	131968	14.83%	1.305%
24	12.687	3519	3529	3548	rVV	142220	268535	30.19%	2.655%
25	12.767	3548	3554	3569	rVV2	12465	20326	2.28%	0.201%
26	12.870	3569	3586	3596	rVB3	43954	78626	8.84%	0.777%
27	12.947	3601	3610	3613	rVV3	16939	23508	2.64%	0.232%
28	12.980	3614	3620	3635	rVB2	40597	65737	7.39%	0.650%
29	13.121	3652	3664	3674	rBV2	31435	52083	5.85%	0.515%
30	13.208	3681	3691	3693	rBV3	7675	11041	1.24%	0.109%
31	13.256	3695	3706	3713	rVV	59949	105667	11.88%	1.045%
32	13.301	3713	3720	3726	rVV	64624	102296	11.50%	1.011%
33	13.333	3727	3730	3744	rVV2	31808	40469	4.55%	0.400%
34	13.417	3744	3756	3760	rVV5	8712	16760	1.88%	0.166%
35	13.459	3761	3769	3782	rVV4	56318	99640	11.20%	0.985%
36	13.536	3783	3793	3801	rVV4	26915	46191	5.19%	0.457%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

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	13.597	3805	3812	3817	rVV3	9894	16764	1.88%	0.166%
38	13.632	3817	3823	3838	rVB4	17143	30277	3.40%	0.299%
39	13.748	3848	3859	3864	rBV6	15984	28953	3.25%	0.286%
40	13.793	3865	3873	3881	rVV2	42919	76284	8.58%	0.754%
41	13.844	3881	3889	3896	rVV3	57361	93893	10.55%	0.928%
42	13.893	3896	3904	3905	rVV2	46521	57804	6.50%	0.571%
43	13.918	3907	3912	3916	rVV	77655	115514	12.98%	1.142%
44	13.947	3916	3921	3945	rVB2	100853	207789	23.36%	2.054%
45	14.066	3945	3958	3966	rBV3	20970	39263	4.41%	0.388%
46	14.124	3968	3976	3988	rVV6	19527	41151	4.63%	0.407%
47	14.198	3988	3999	4011	rVB	219198	343988	38.67%	3.400%
48	14.291	4011	4028	4040	rBV	261815	448203	50.38%	4.431%
49	14.352	4040	4047	4057	rVV4	26564	48734	5.48%	0.482%
50	14.401	4058	4062	4070	rVB3	8377	11278	1.27%	0.111%
51	14.491	4079	4090	4097	rBV	78487	130778	14.70%	1.293%
52	14.674	4137	4147	4172	rBV3	96560	213373	23.99%	2.109%
53	14.815	4182	4191	4201	rBV2	67754	103061	11.59%	1.019%
54	14.883	4202	4212	4224	rVB2	114787	196141	22.05%	1.939%
55	14.947	4224	4232	4243	rVB4	17851	30799	3.46%	0.304%
56	15.024	4243	4256	4269	rBV2	137715	254906	28.65%	2.520%
57	15.086	4270	4275	4289	rVB3	18671	31222	3.51%	0.309%
58	15.176	4291	4303	4304	rBV3	26492	34663	3.90%	0.343%
59	15.205	4306	4312	4317	rBV2	35682	47507	5.34%	0.470%
60	15.278	4327	4335	4349	rVB4	74051	131021	14.73%	1.295%
61	15.356	4350	4359	4366	rVV5	15043	24316	2.73%	0.240%
62	15.417	4367	4378	4389	rVV2	237559	415665	46.72%	4.109%
63	15.475	4390	4396	4401	rVV5	37909	62565	7.03%	0.618%
64	15.510	4402	4407	4418	rVV4	34250	55246	6.21%	0.546%
65	15.600	4418	4435	4445	rVB7	26139	63182	7.10%	0.625%
66	15.738	4470	4478	4483	rBV4	17967	27310	3.07%	0.270%
67	15.931	4519	4538	4555	rBV5	65313	163137	18.34%	1.613%
68	16.021	4557	4566	4580	rVB3	34939	59652	6.71%	0.590%
69	16.156	4600	4608	4612	rBV5	11500	16285	1.83%	0.161%
70	16.192	4612	4619	4628	rVB	25328	39144	4.40%	0.387%
71	16.272	4636	4644	4649	rBV2	14421	21331	2.40%	0.211%
72	16.317	4651	4658	4671	rVB7	27441	50817	5.71%	0.502%
73	16.417	4679	4689	4698	rBV	77441	120486	13.54%	1.191%
74	16.532	4717	4725	4738	rBV3	9817	21258	2.39%	0.210%
75	16.616	4743	4751	4756	rVV2	32195	53437	6.01%	0.528%
76	16.651	4757	4762	4777	rVB2	42304	68984	7.75%	0.682%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

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

77	16.793	4796	4806	4813	rBV2	32280	49231	5.53%	0.487%
78	17.102	4888	4902	4914	rVB5	12580	27390	3.08%	0.271%
79	17.278	4949	4957	4969	rVV9	7726	15697	1.76%	0.155%
80	17.336	4969	4975	4984	rVV7	5456	9039	1.02%	0.089%
81	17.548	5034	5041	5051	rVB3	11313	17593	1.98%	0.174%

Sum of corrected areas: 10116044

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
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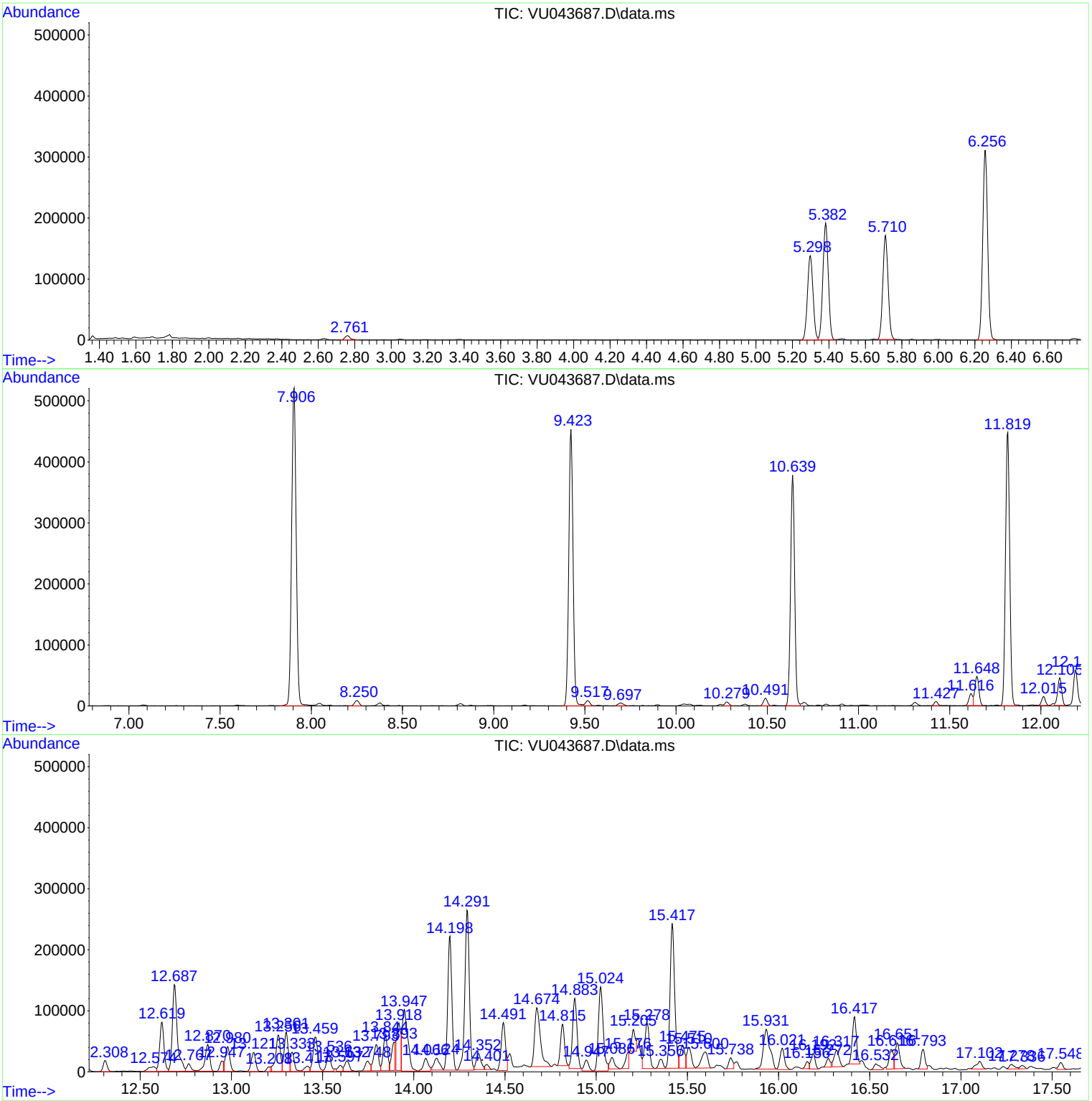
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

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\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

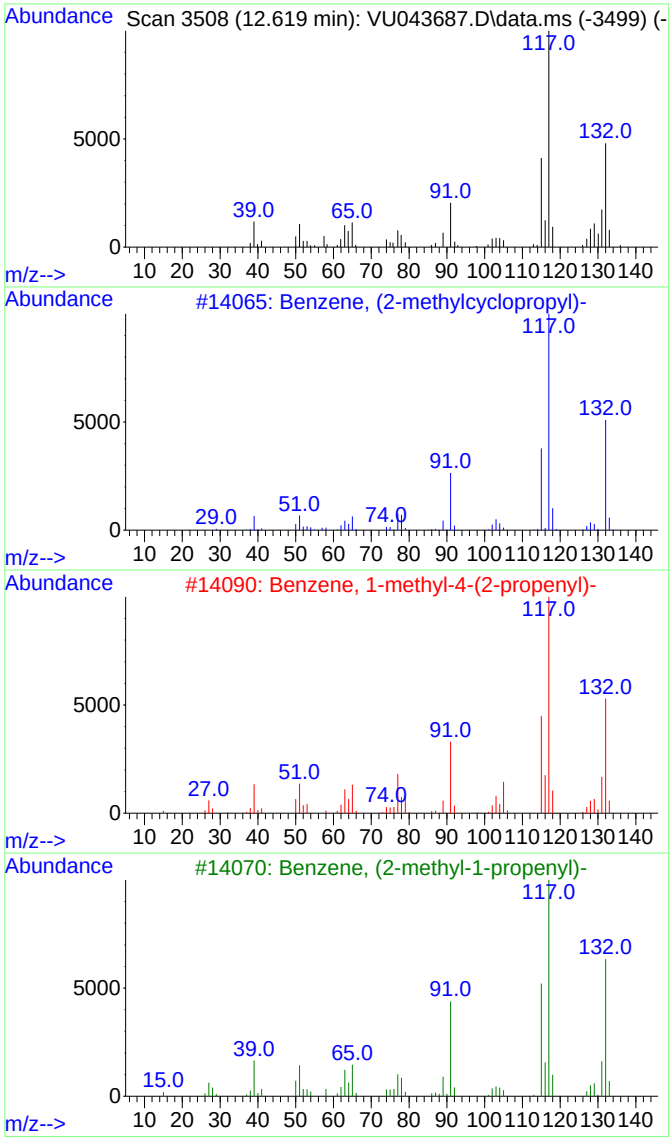
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

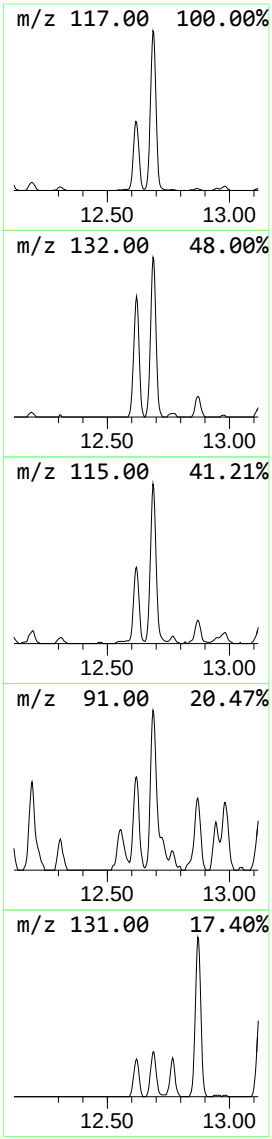
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, (2-methylcycloprop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.619	9.30 ug/l	131968	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	94
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	93
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	93
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	93
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

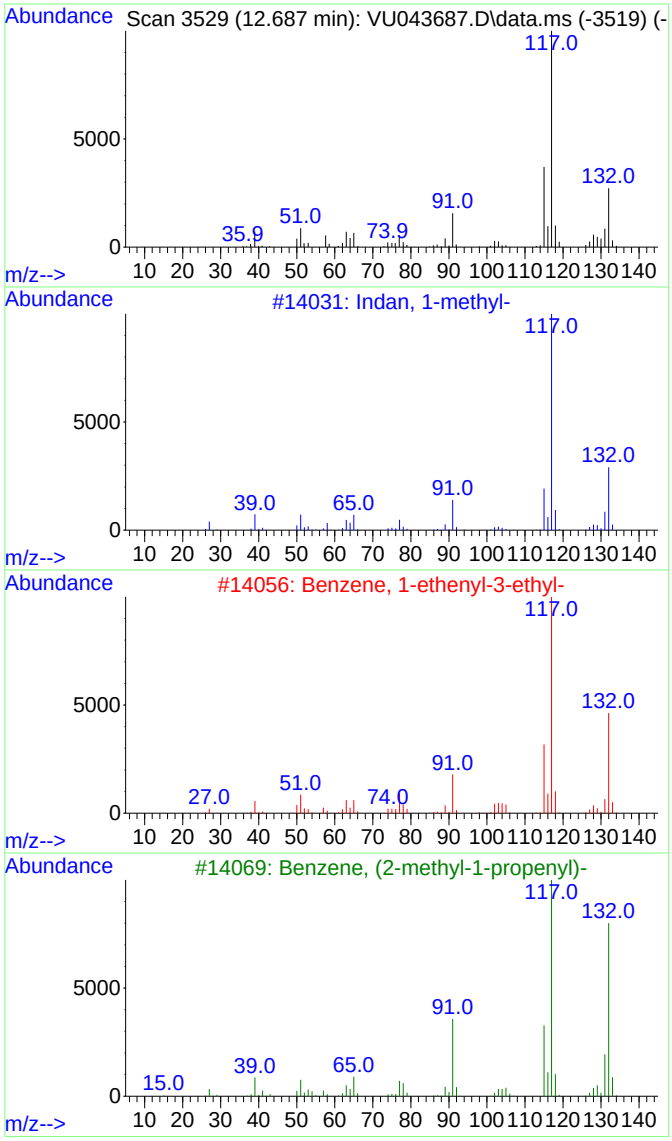
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

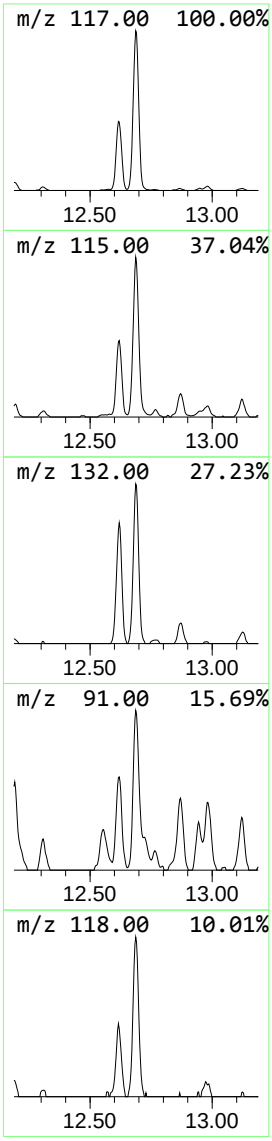
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indan, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	18.91 ug/l	268535	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	87
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

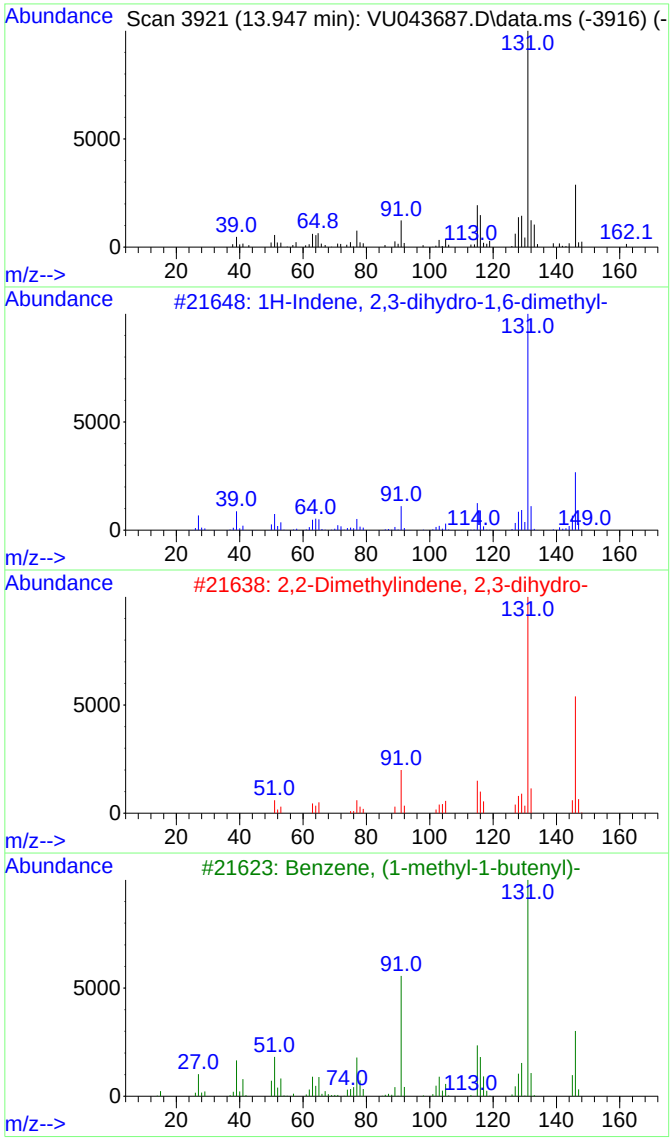
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

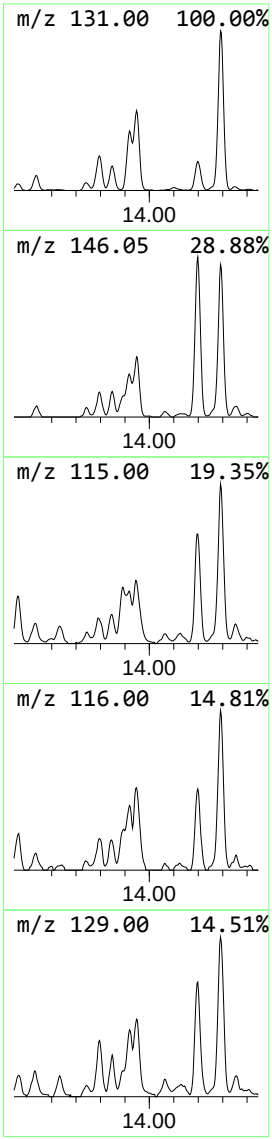
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.947	14.64 ug/l	207789	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	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

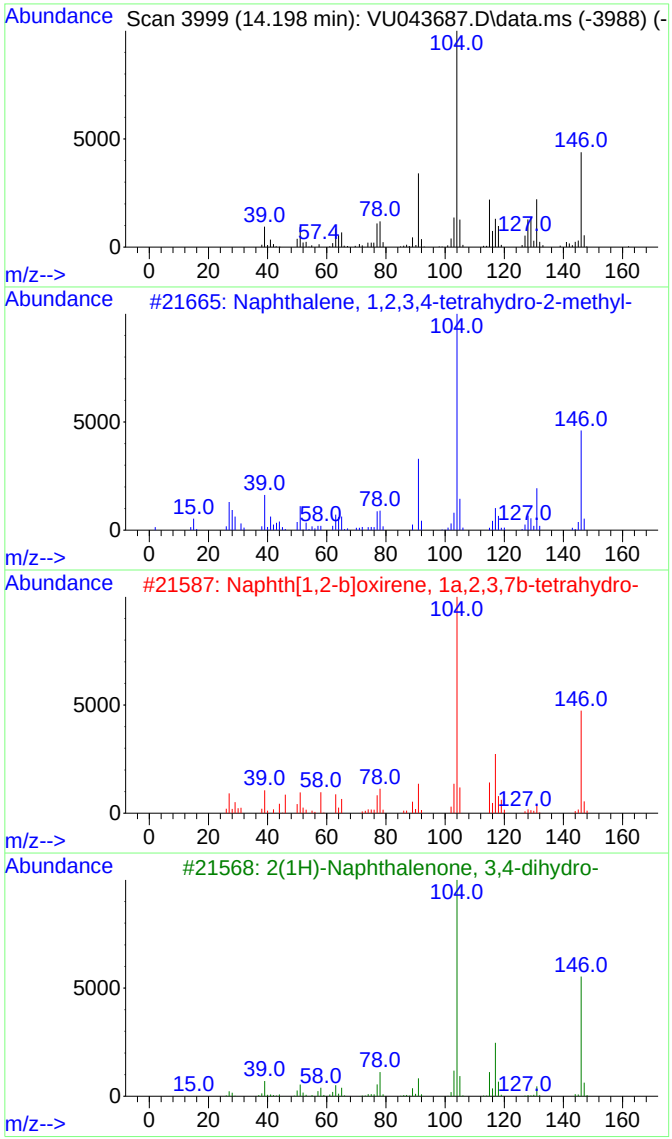
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

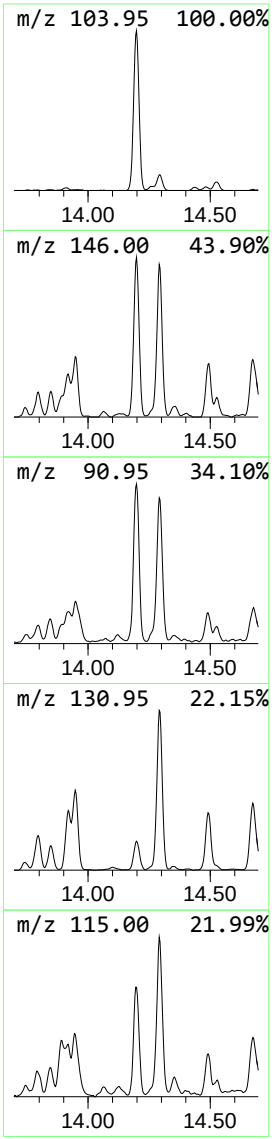
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.198	24.23 ug/l	343988	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	62
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5	Benzenemethanimine, .alpha.-(1,1...	161	C11H15N	033611-54-0	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

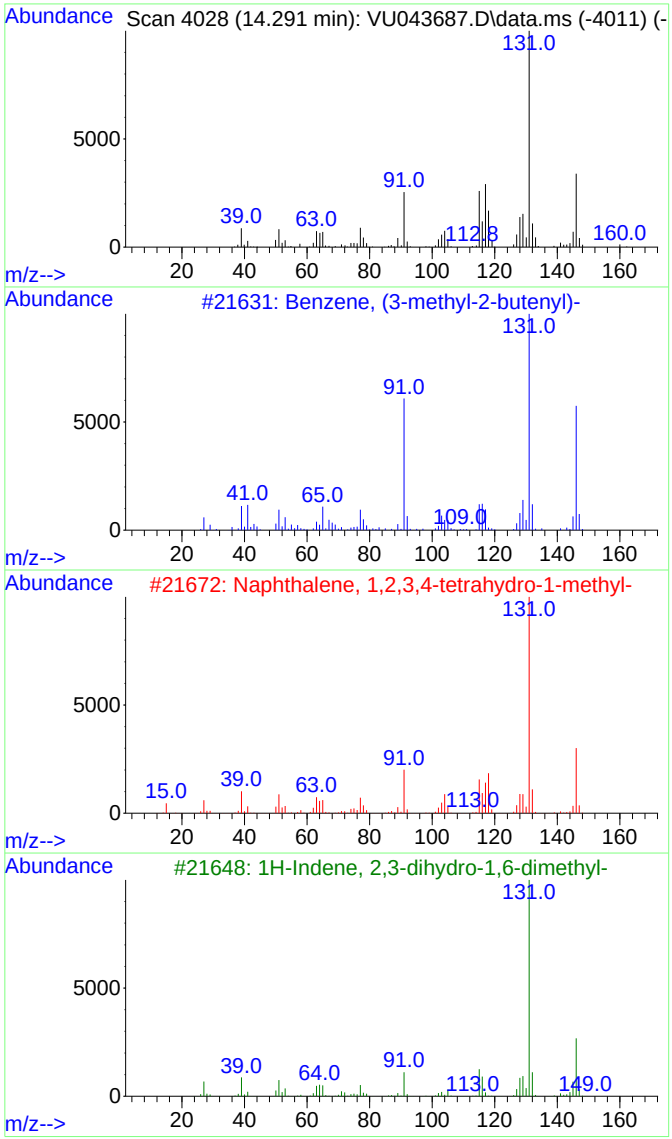
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

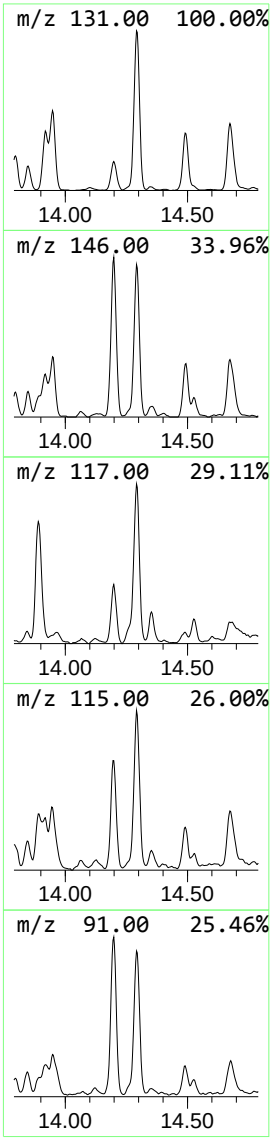
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, (3-methyl-2-butenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.291	31.57 ug/l	448203	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	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

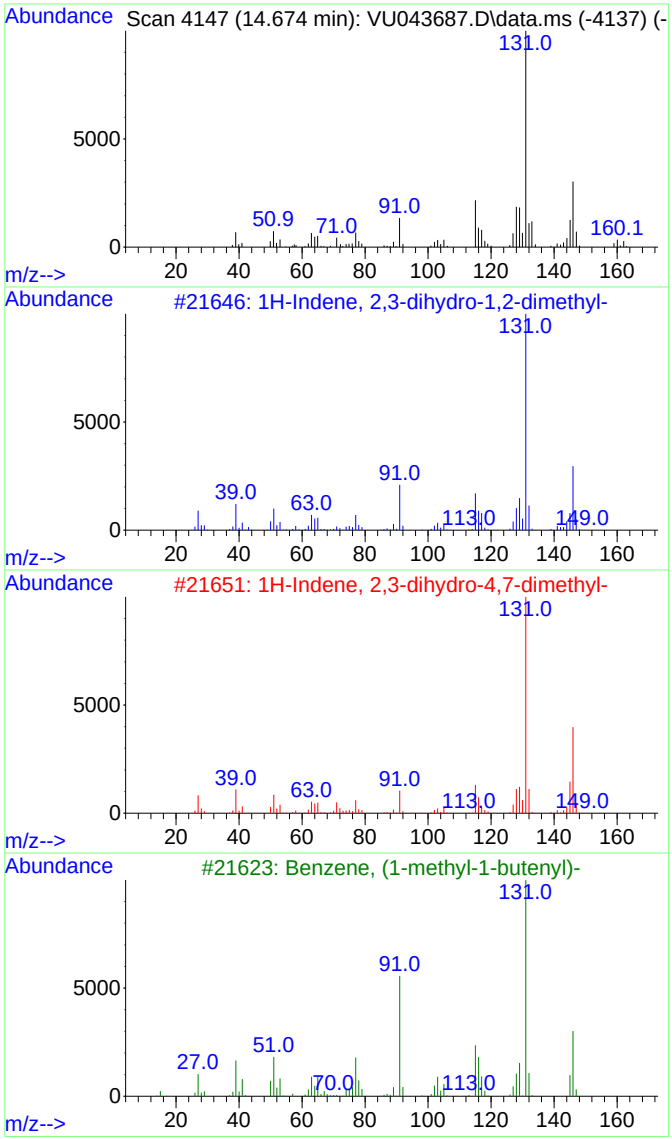
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

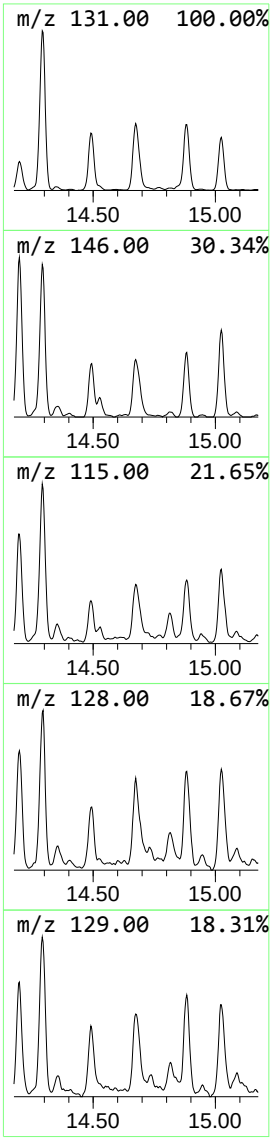
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.674	15.03 ug/l	213373	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	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	87
4	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	83
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81





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Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

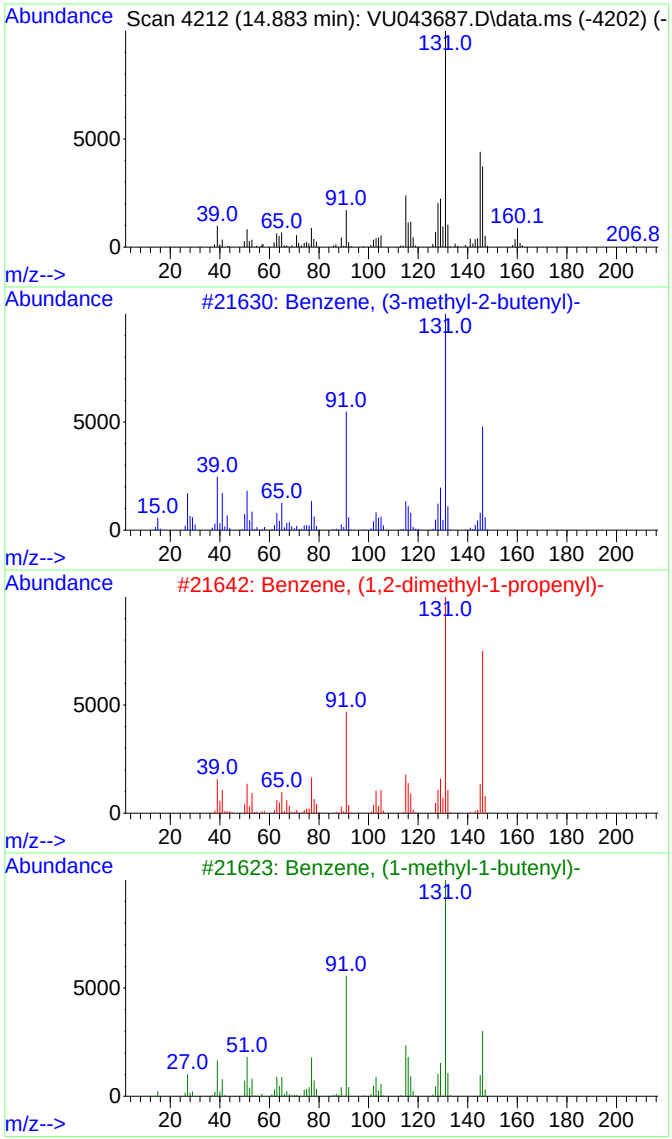
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

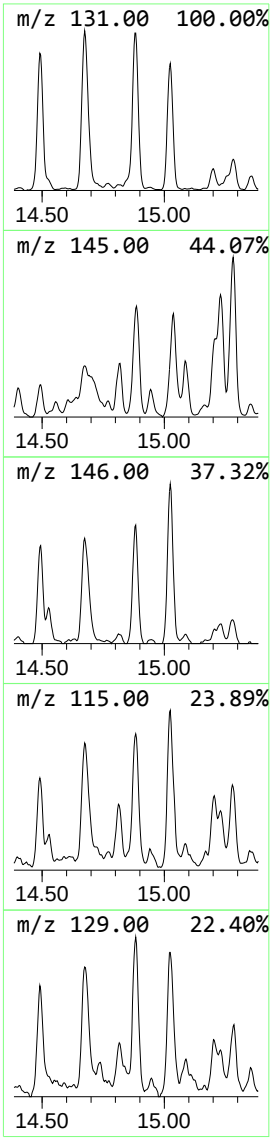
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.883	13.82 ug/l	196141	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	76
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	76
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	62
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

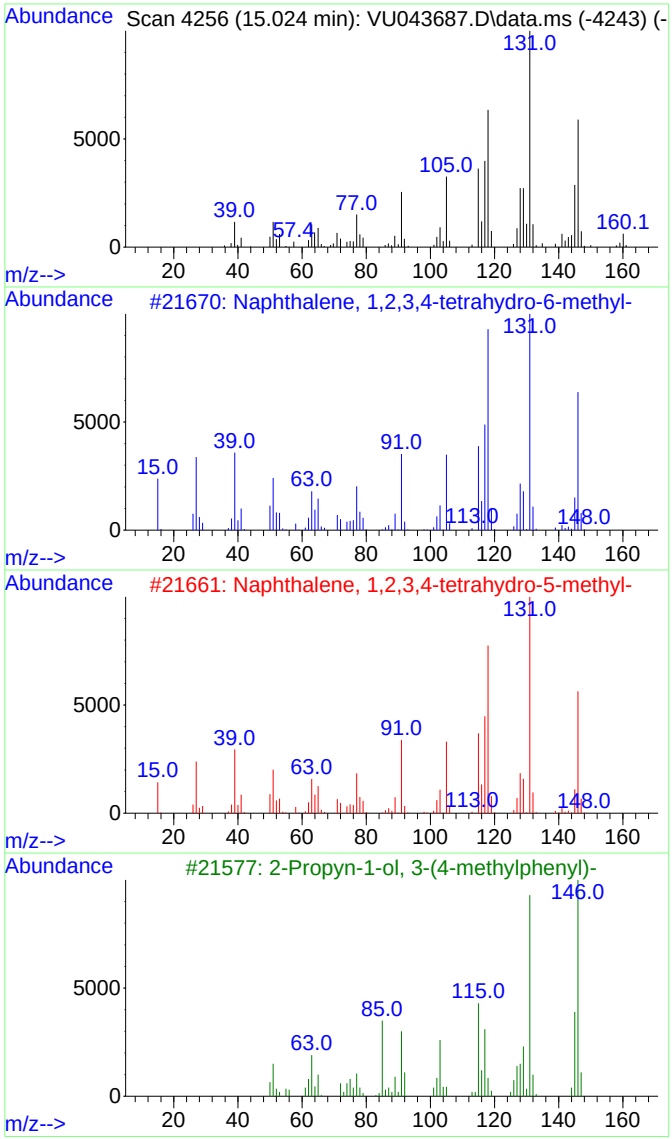
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

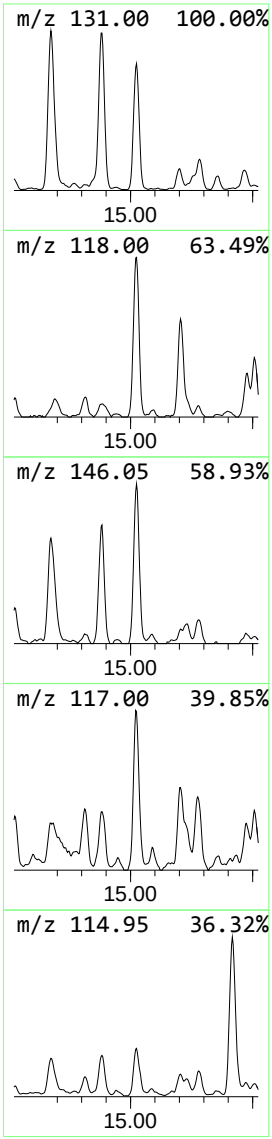
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.024	17.95 ug/l	254906	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	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
3	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	74
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

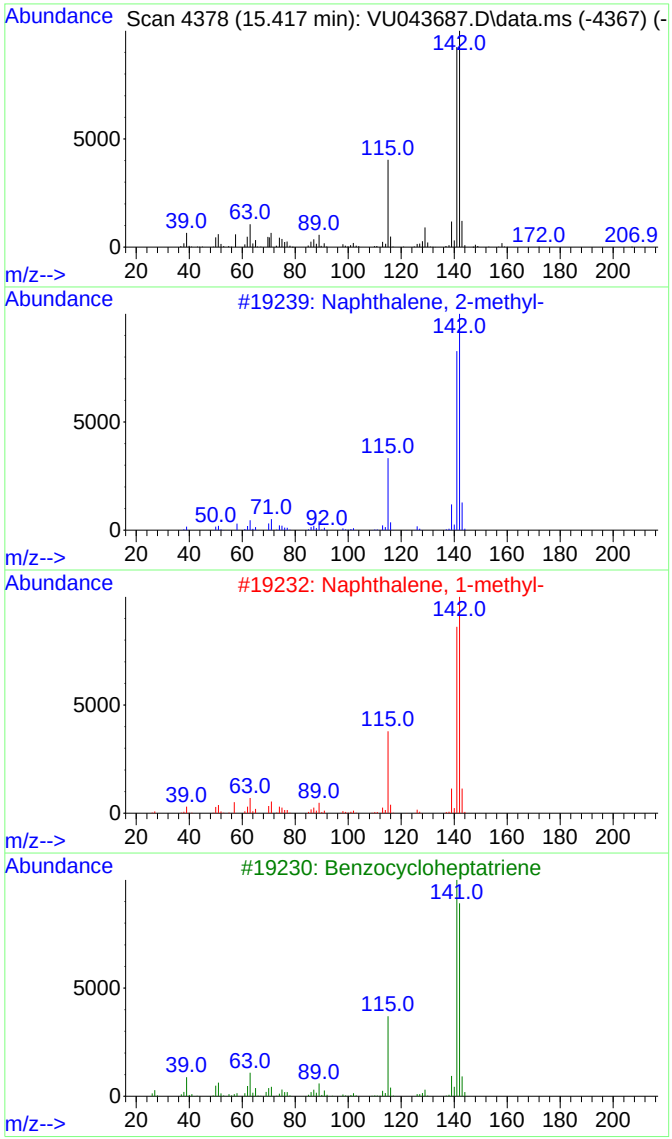
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

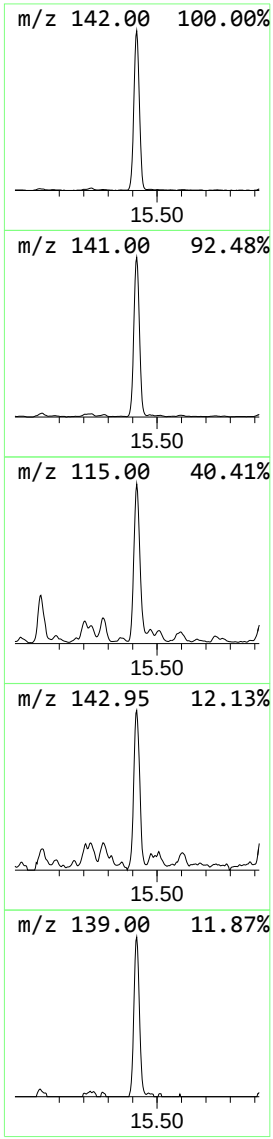
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.417	29.28 ug/l	415665	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

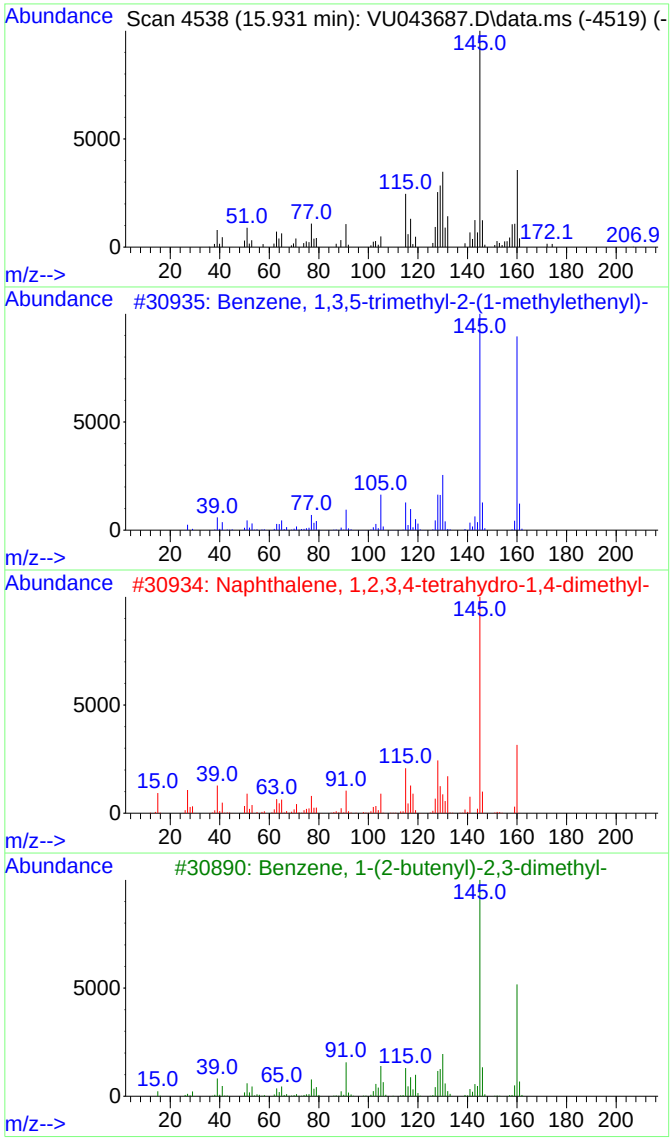
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

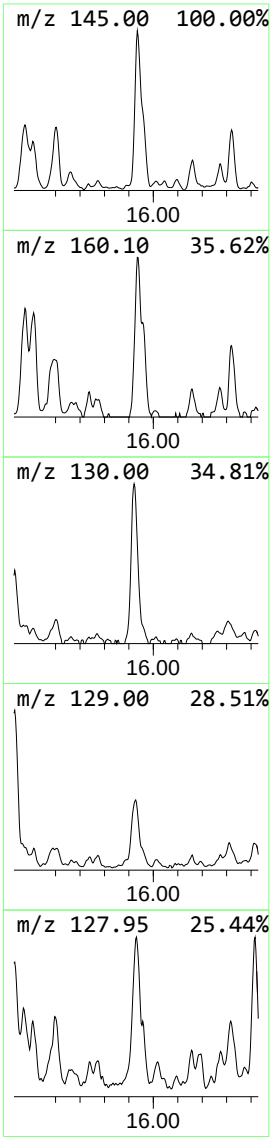
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.931	11.49 ug/l	163137	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	91
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	86
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	68
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	64
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051221\
Data File : VU043687.D
Acq On : 12 May 2021 18:57
Operator : SY/MD
Sample : M2365-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-3

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, (2-met...	12.619	9.3	ug/l	131968	4	11.819	709853	50.0
Indan, 1-methyl-	12.687	18.9	ug/l	268535	4	11.819	709853	50.0
1H-Indene, 2,3-...	13.947	14.6	ug/l	207789	4	11.819	709853	50.0
Naphthalene, 1,...	14.198	24.2	ug/l	343988	4	11.819	709853	50.0
Benzene, (3-met...	14.291	31.6	ug/l	448203	4	11.819	709853	50.0
1H-Indene, 2,3-...	14.674	15.0	ug/l	213373	4	11.819	709853	50.0
Benzene, (1,2-d...	14.883	13.8	ug/l	196141	4	11.819	709853	50.0
Naphthalene, 1,...	15.024	17.9	ug/l	254906	4	11.819	709853	50.0
Naphthalene, 1-...	15.417	29.3	ug/l	415665	4	11.819	709853	50.0
Benzene, 1,3,5-...	15.931	11.5	ug/l	163137	4	11.819	709853	50.0