

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
 Data File : VU043613.D
 Acq On : 07 May 2021 15:31
 Operator : SY/MD
 Sample : M2304-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 OWL-C4-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: VU043613.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.629	386	401	411	rBV	12300	21116	1.08%	0.120%
2	5.057	1141	1156	1181	rBV	110687	244388	12.49%	1.394%
3	5.298	1215	1231	1244	rBV2	139541	294415	15.05%	1.679%
4	5.382	1244	1257	1274	rVB	196409	401208	20.51%	2.288%
5	5.710	1345	1359	1379	rBV	169154	334943	17.12%	1.910%
6	6.256	1515	1529	1547	rBV	313384	577610	29.53%	3.294%
7	7.906	2023	2042	2060	rBV	515768	883208	45.15%	5.037%
8	9.423	2501	2514	2534	rBV	452695	731105	37.37%	4.170%
9	10.491	2834	2846	2858	rBV	109262	169304	8.66%	0.966%
10	10.639	2879	2892	2907	rBV	381939	594601	30.40%	3.391%
11	10.912	2966	2977	2992	rBV	55083	84479	4.32%	0.482%
12	11.616	3184	3196	3199	rBV	14995	20806	1.06%	0.119%
13	11.648	3199	3206	3220	rVB	82303	127966	6.54%	0.730%
14	11.819	3247	3259	3275	rBV	449919	701497	35.86%	4.001%
15	12.015	3310	3320	3330	rVB2	28312	42817	2.19%	0.244%
16	12.105	3330	3348	3359	rBV2	167548	258247	13.20%	1.473%
17	12.214	3363	3382	3393	rVV	28878	56218	2.87%	0.321%
18	12.307	3396	3411	3427	rBV	64910	99320	5.08%	0.566%
19	12.577	3477	3495	3502	rBV	336023	512898	26.22%	2.925%
20	12.619	3502	3508	3518	rVV	158440	241003	12.32%	1.375%
21	12.690	3518	3530	3548	rVV2	454108	864508	44.19%	4.931%
22	12.870	3567	3586	3596	rVV	33310	58810	3.01%	0.335%
23	12.979	3600	3620	3634	rBV	258596	417984	21.37%	2.384%
24	13.114	3652	3662	3676	rBV4	45506	81805	4.18%	0.467%
25	13.208	3681	3691	3698	rBV	30779	46451	2.37%	0.265%
26	13.259	3698	3707	3713	rVV3	61744	100319	5.13%	0.572%
27	13.301	3713	3720	3745	rVB	206964	341662	17.47%	1.949%
28	13.462	3745	3770	3786	rBV3	1124749	1956140	100.00%	11.157%
29	13.536	3786	3793	3797	rBV3	16049	22753	1.16%	0.130%
30	13.565	3797	3802	3810	rVB	28763	38217	1.95%	0.218%
31	13.632	3816	3823	3839	rVB4	29431	51292	2.62%	0.293%
32	13.741	3847	3857	3864	rBV2	42994	67040	3.43%	0.382%
33	13.793	3864	3873	3881	rVV	119648	209995	10.74%	1.198%
34	13.848	3881	3890	3898	rVV3	180612	308550	15.77%	1.760%
35	13.918	3898	3912	3915	rVV3	109373	208809	10.67%	1.191%
36	13.947	3916	3921	3943	rVB2	215547	419728	21.46%	2.394%

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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

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37	14.066	3943	3958	3967	rBV2	26709	47890	2.45%	0.273%
38	14.124	3967	3976	3987	rVV2	46609	75942	3.88%	0.433%
39	14.198	3988	3999	4011	rVB	251146	392714	20.08%	2.240%
40	14.294	4011	4029	4039	rBV2	317465	547410	27.98%	3.122%
41	14.352	4039	4047	4057	rVV2	99215	166570	8.52%	0.950%
42	14.401	4057	4062	4070	rVB3	15214	21175	1.08%	0.121%
43	14.491	4079	4090	4099	rBV	174809	286249	14.63%	1.633%
44	14.690	4135	4152	4172	rBV3	273826	698708	35.72%	3.985%
45	14.815	4182	4191	4202	rBV	80598	122633	6.27%	0.699%
46	14.883	4202	4212	4224	rVB2	183534	311363	15.92%	1.776%
47	14.947	4224	4232	4243	rVB3	21883	35373	1.81%	0.202%
48	15.024	4243	4256	4268	rBV2	449553	741878	37.93%	4.231%
49	15.089	4271	4276	4290	rVB4	17564	29692	1.52%	0.169%
50	15.204	4290	4312	4327	rBV5	94652	242481	12.40%	1.383%
51	15.282	4328	4336	4350	rVB3	66779	117825	6.02%	0.672%
52	15.356	4350	4359	4369	rBV3	24324	36758	1.88%	0.210%
53	15.433	4373	4383	4388	rVV3	22490	41571	2.13%	0.237%
54	15.474	4389	4396	4402	rVV3	48141	82793	4.23%	0.472%
55	15.510	4403	4407	4418	rVB4	33188	47668	2.44%	0.272%
56	15.587	4419	4431	4446	rVB3	35997	89128	4.56%	0.508%
57	15.774	4473	4489	4499	rBV3	19982	34822	1.78%	0.199%
58	15.934	4522	4539	4556	rBV4	99758	221829	11.34%	1.265%
59	16.021	4556	4566	4582	rVB6	23692	45021	2.30%	0.257%
60	16.156	4598	4608	4613	rBV	70696	108683	5.56%	0.620%
61	16.191	4613	4619	4628	rVB	60282	91110	4.66%	0.520%
62	16.269	4629	4643	4653	rBV	238869	413499	21.14%	2.358%
63	16.417	4672	4689	4700	rBV	277875	443583	22.68%	2.530%
64	16.539	4718	4727	4741	rBV5	14322	27711	1.42%	0.158%
65	16.648	4742	4761	4781	rVV2	96996	257999	13.19%	1.472%
66	16.793	4796	4806	4826	rVB2	60995	103224	5.28%	0.589%
67	17.105	4895	4903	4916	rVB4	17347	34387	1.76%	0.196%
68	17.548	5033	5041	5053	rVB2	14196	23994	1.23%	0.137%

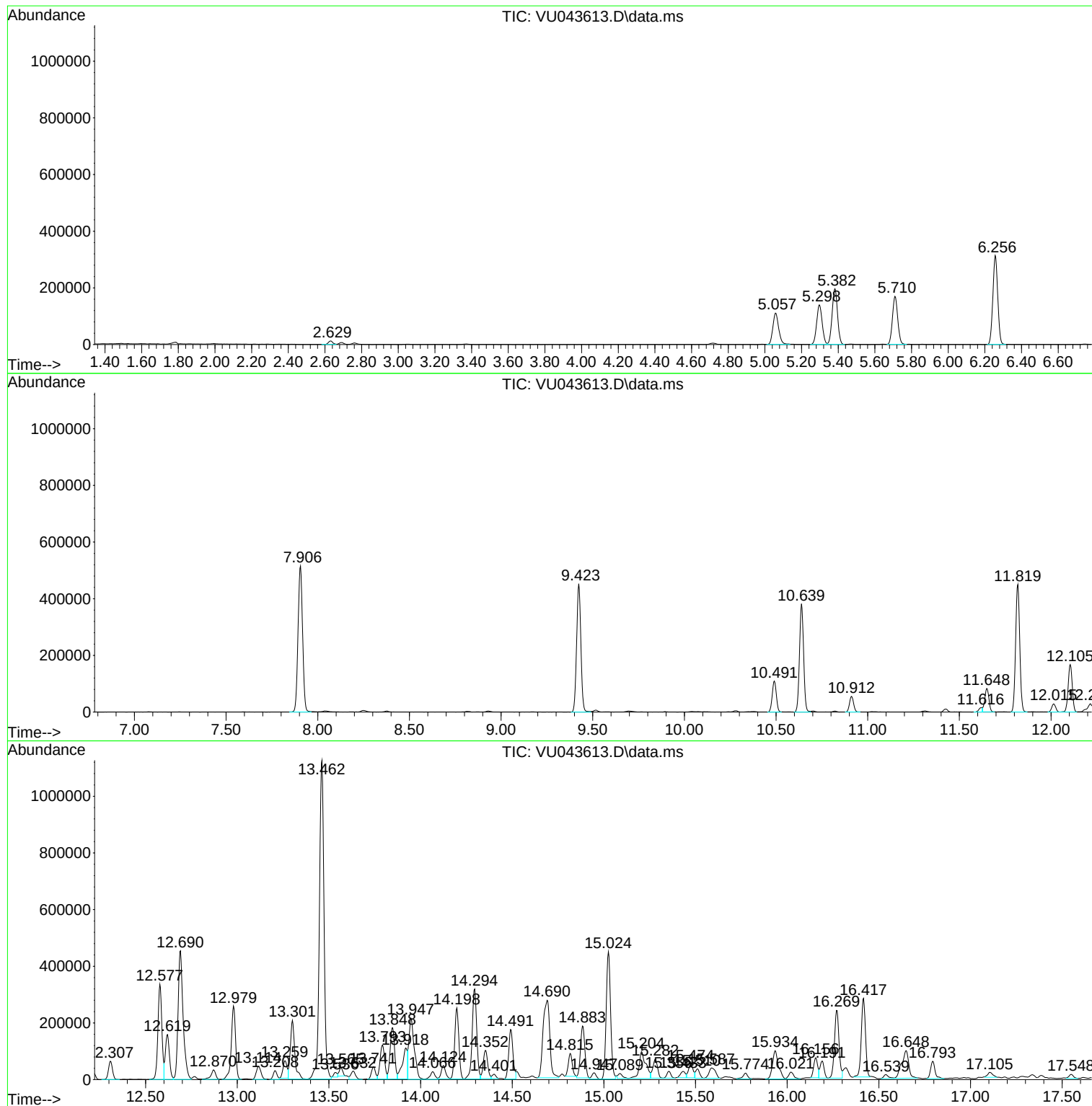
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ClientSampleId :
OWL-C4-GW

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



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Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

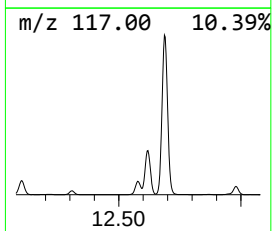
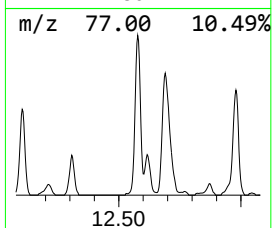
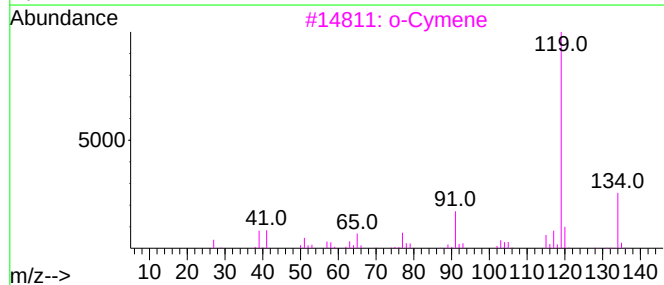
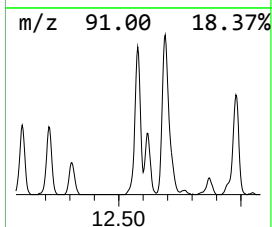
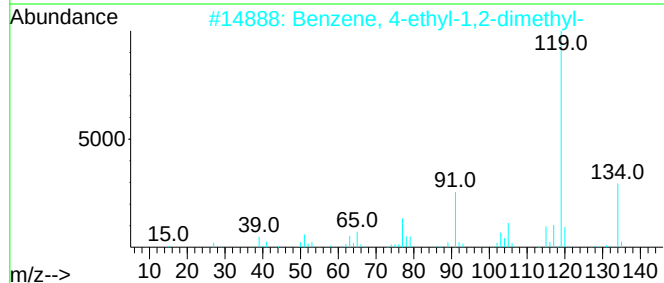
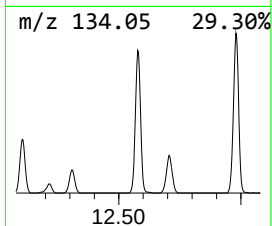
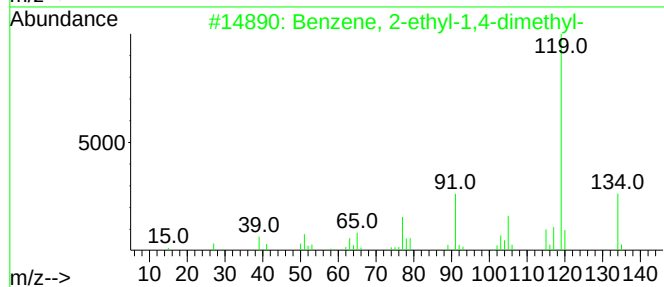
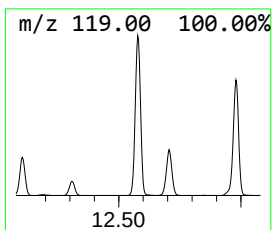
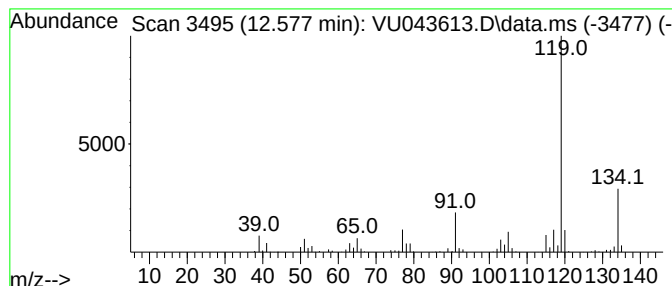
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.578	36.56 ug/l	512898	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
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



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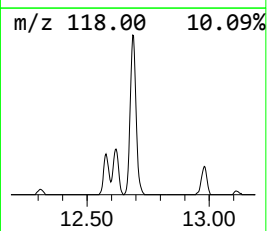
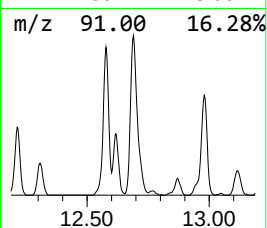
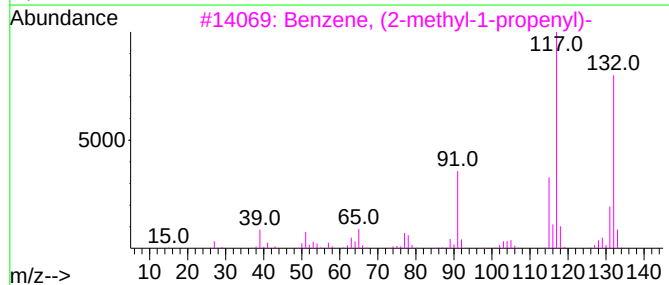
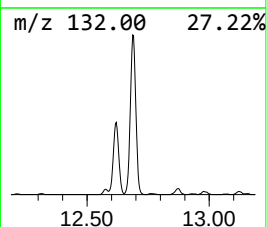
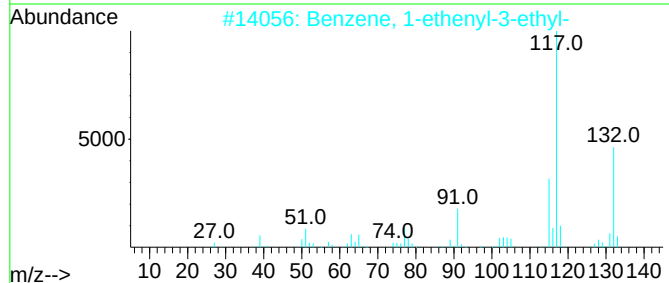
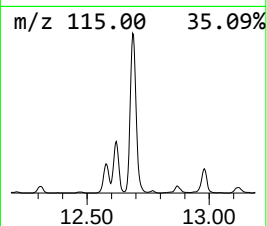
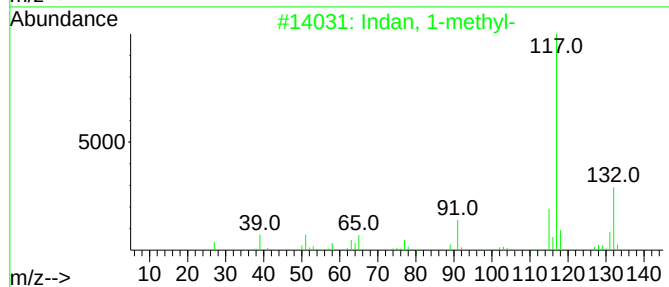
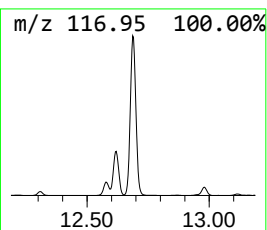
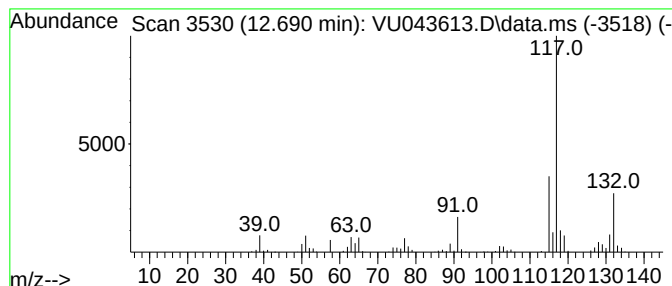
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

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	61.62 ug/l	864508	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



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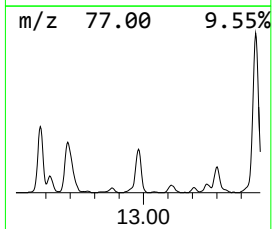
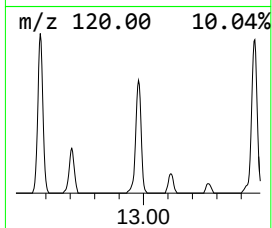
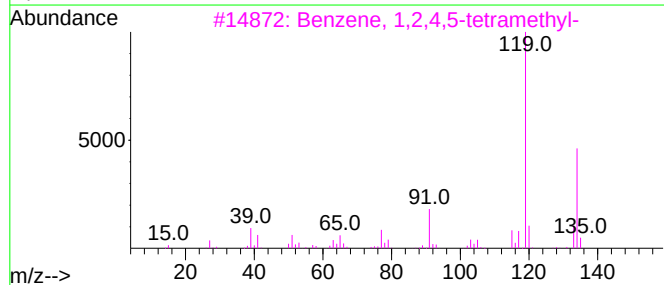
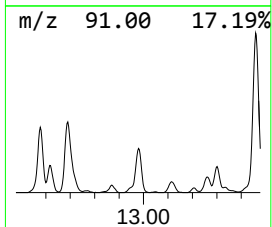
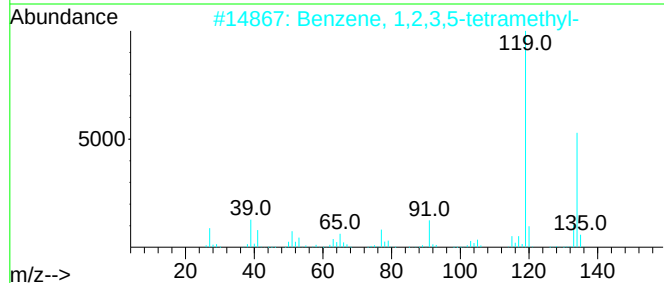
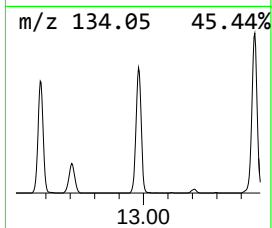
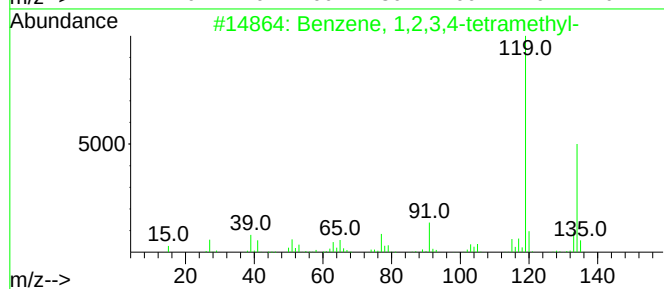
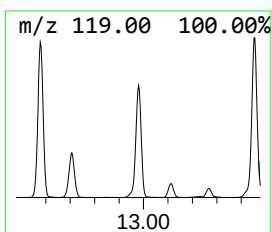
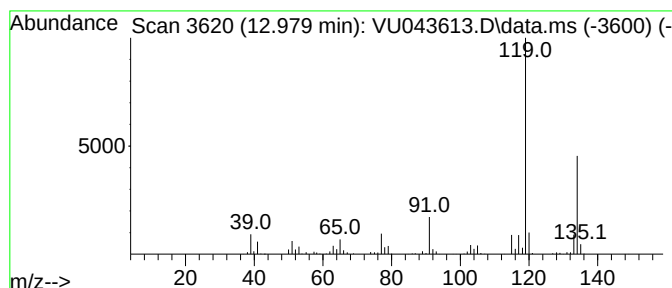
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.979	29.79 ug/l	417984	1,4-Dichlorobenzene-d4	11.819

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



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Misc : 5.0mL/MSVOA_U/WATER
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Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

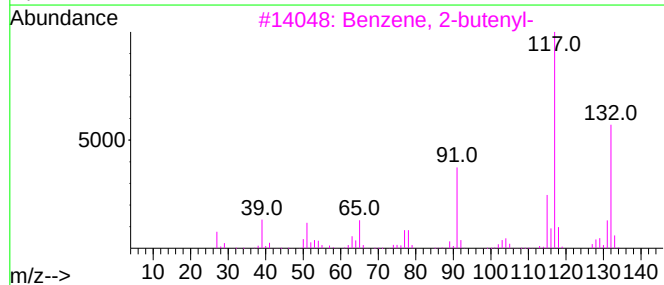
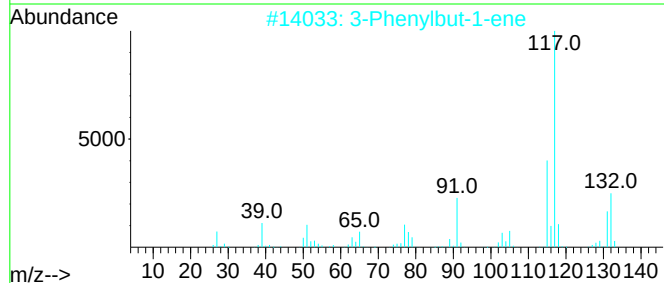
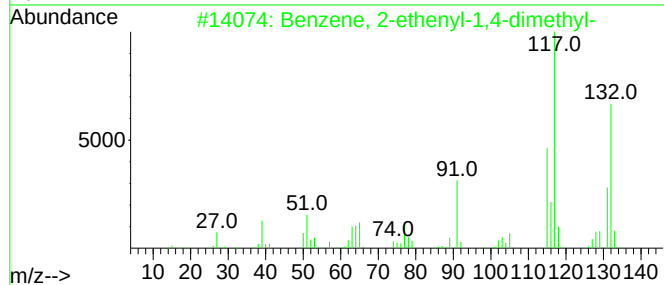
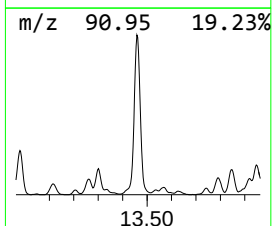
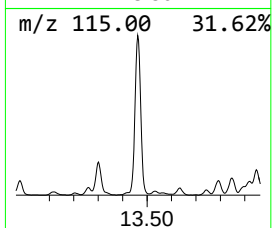
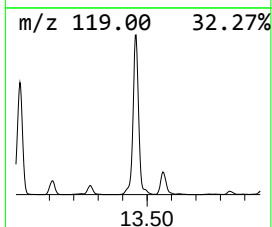
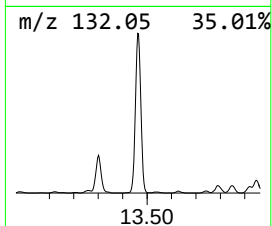
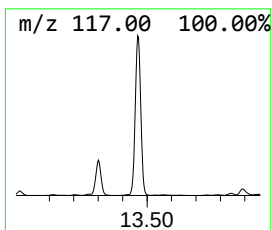
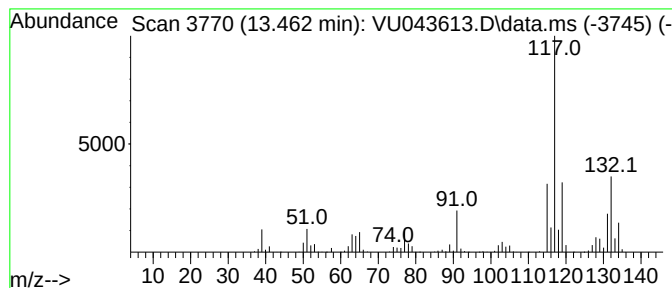
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

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	139.43 ug/l	1956140	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	95
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

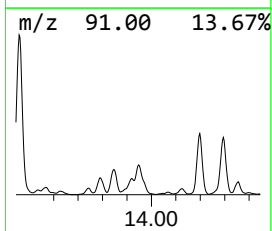
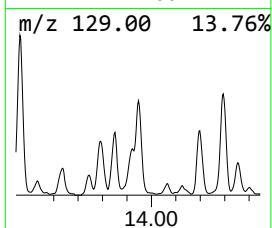
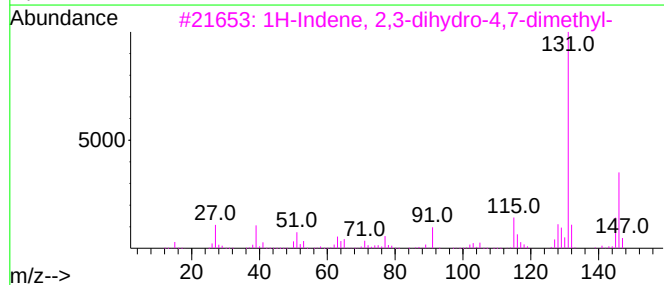
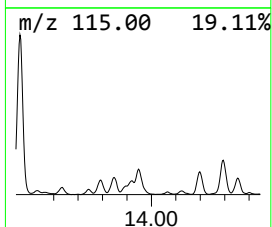
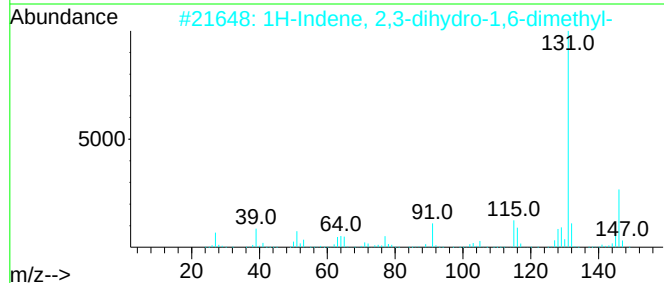
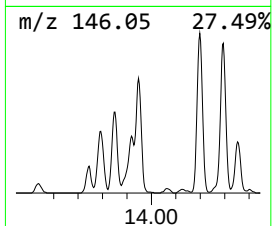
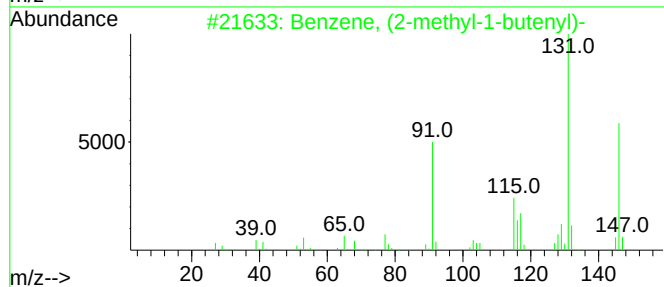
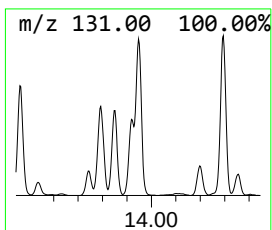
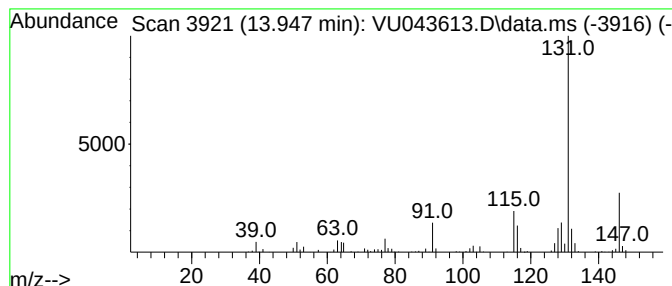
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.947	29.92 ug/l	419728	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

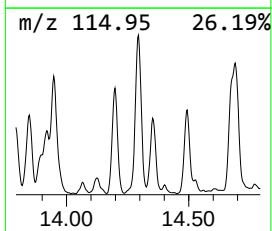
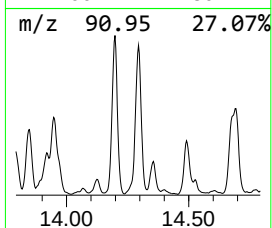
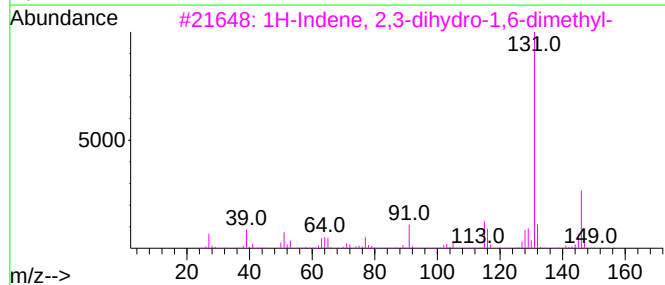
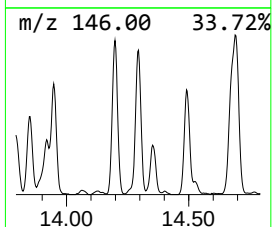
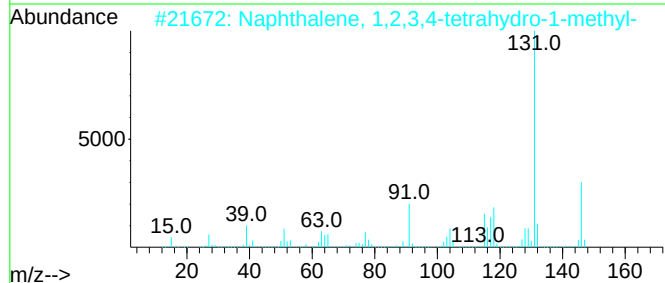
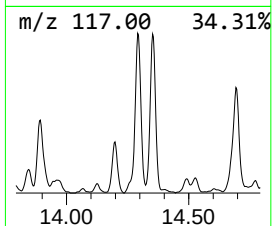
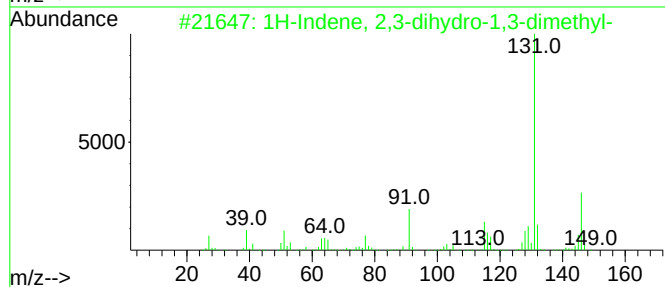
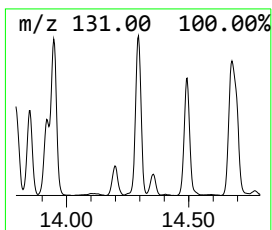
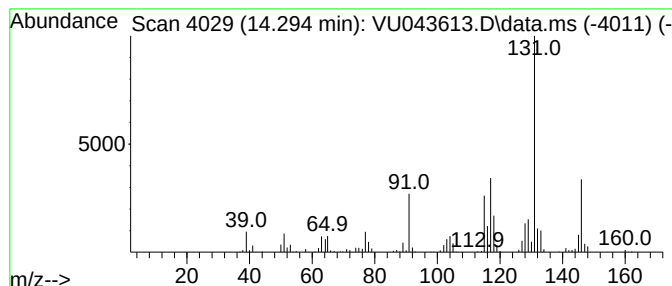
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.294	39.02 ug/l	547410	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	93
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	90
5			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

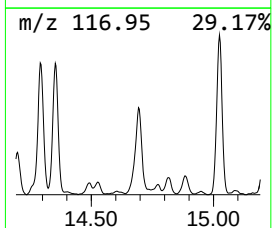
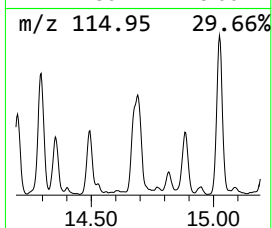
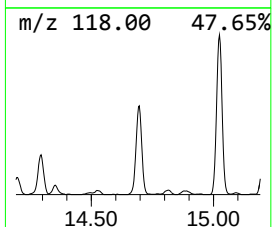
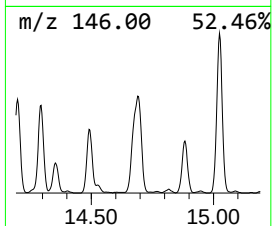
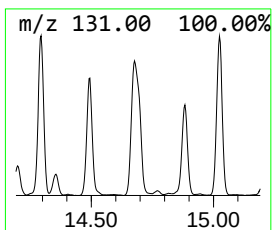
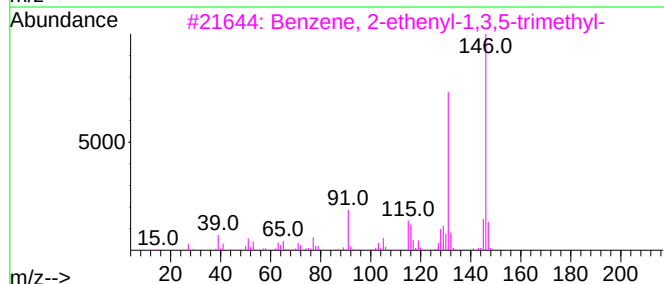
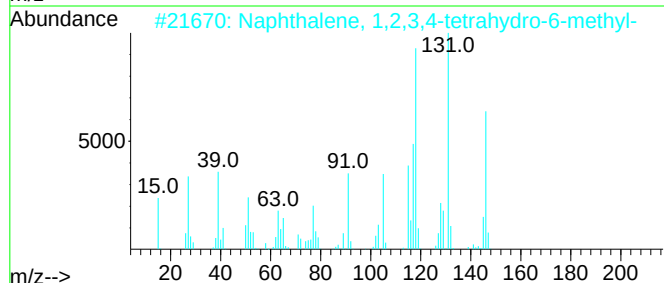
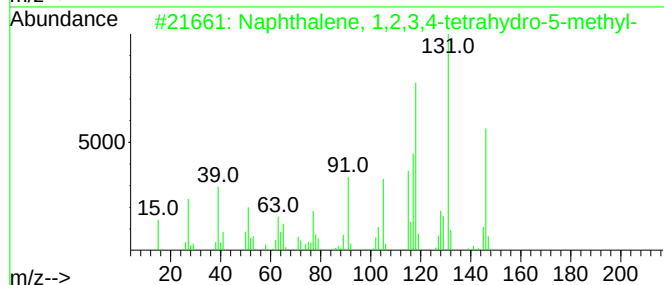
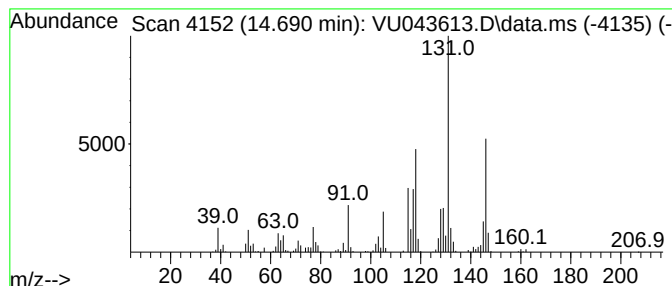
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.690	49.80 ug/l	698708	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	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	89
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	76
4			1H-Inden-1-one, 2,3-dihydro-3-me...	146	C10H10O	006072-57-7	76
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

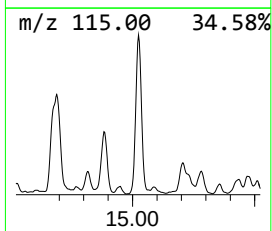
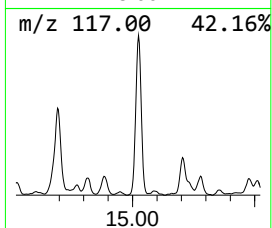
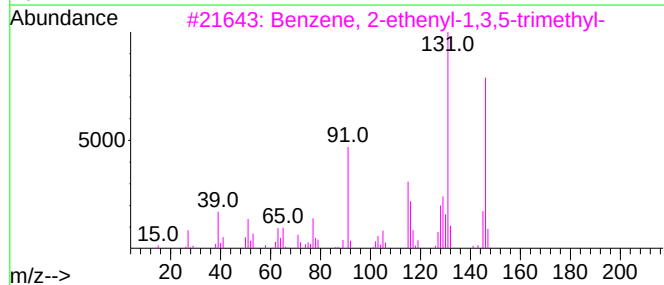
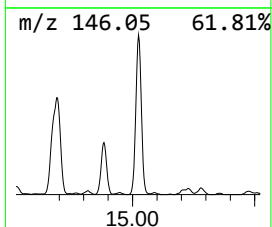
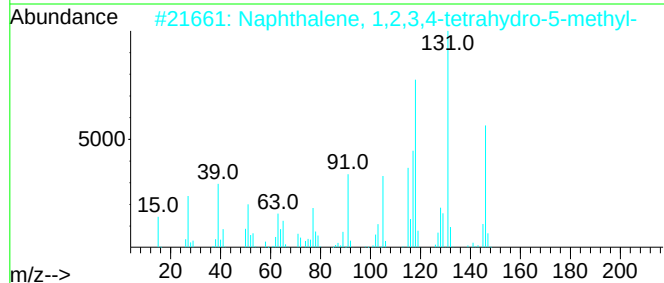
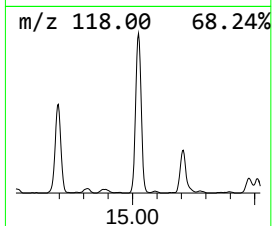
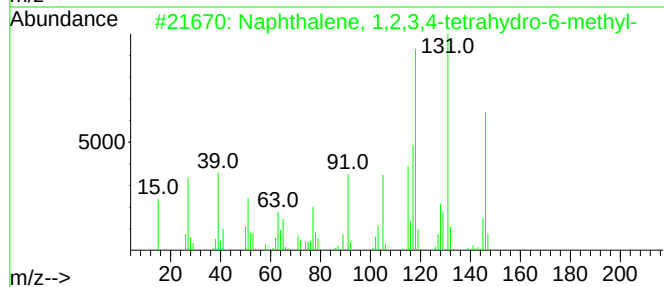
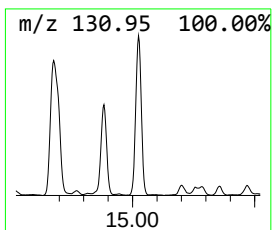
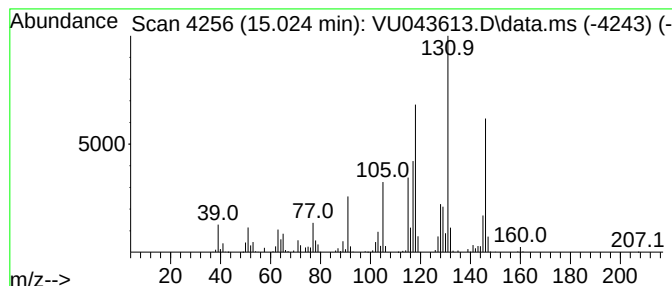
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.024	52.88 ug/l	741878	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	97
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	86
4			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	70
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

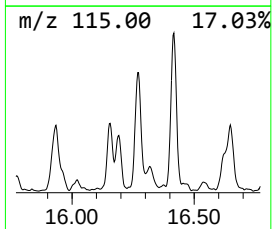
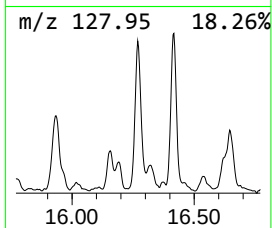
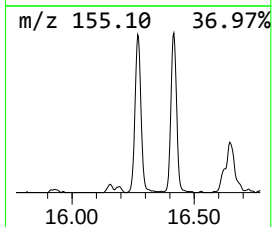
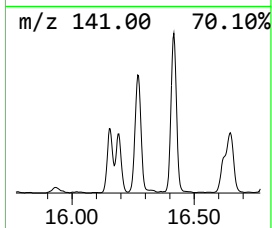
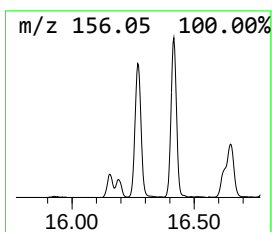
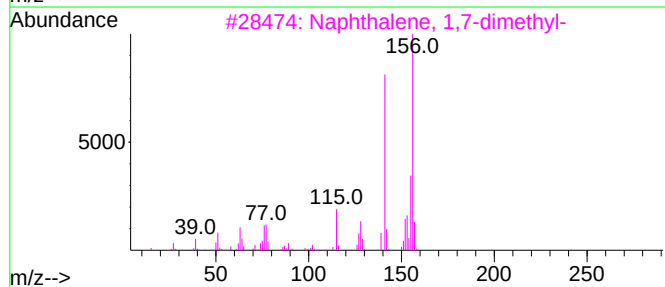
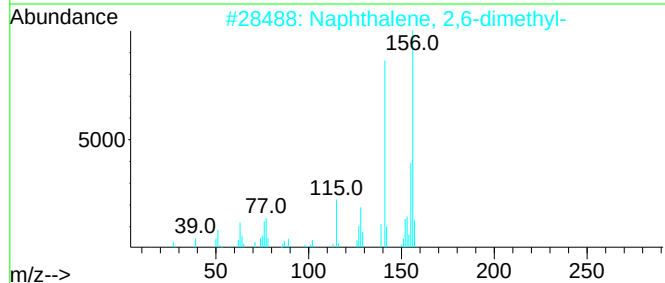
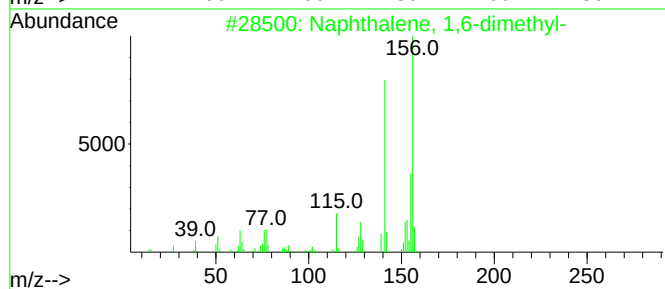
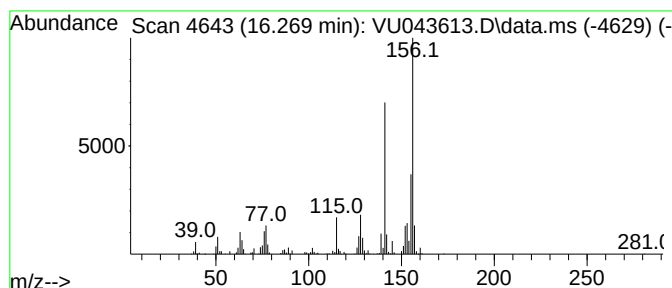
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.269	29.47 ug/l	413499	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-GW

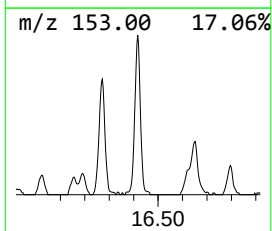
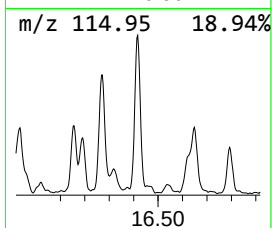
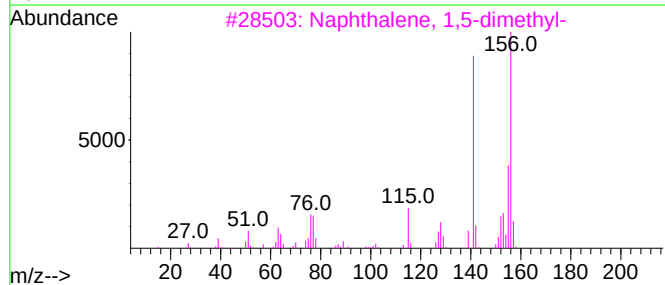
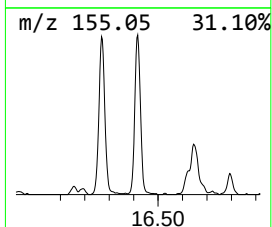
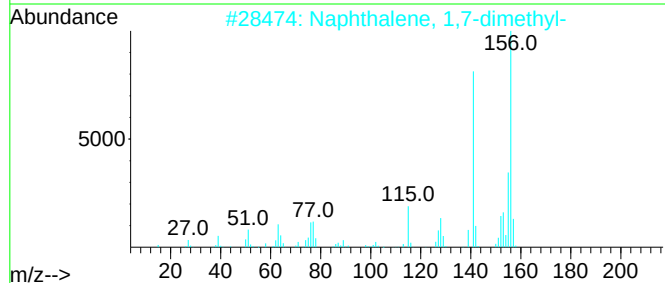
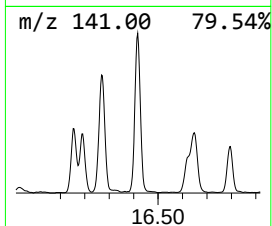
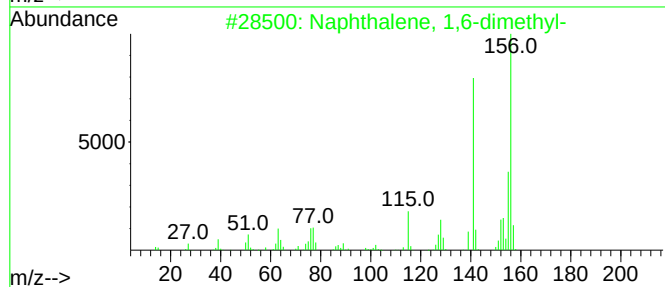
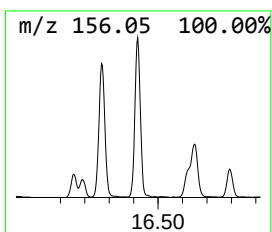
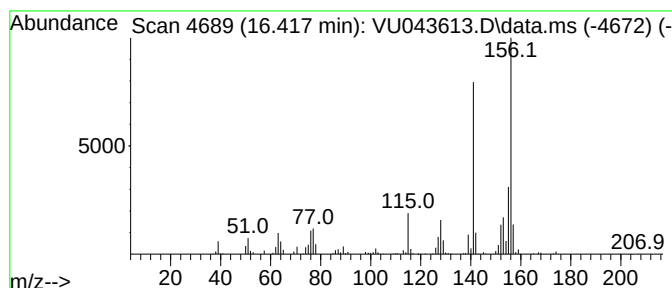
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.417	31.62 ug/l	443583	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	98
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050721\
Data File : VU043613.D
Acq On : 07 May 2021 15:31
Operator : SY/MD
Sample : M2304-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
OWL-C4-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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Indan, 1-methyl-	12.690	61.6	ug/l	864508	4	11.819	701497	50.0
Benzene, 1,2,3,...	12.979	29.8	ug/l	417984	4	11.819	701497	50.0
Benzene, 2-ethe...	13.462	139.4	ug/l	1956140	4	11.819	701497	50.0
Benzene, (2-met...	13.947	29.9	ug/l	419728	4	11.819	701497	50.0
1H-Indene, 2,3-...	14.294	39.0	ug/l	547410	4	11.819	701497	50.0
Naphthalene, 1,...	14.690	49.8	ug/l	698708	4	11.819	701497	50.0
Naphthalene, 1,...	15.024	52.9	ug/l	741878	4	11.819	701497	50.0
Naphthalene, 1,...	16.269	29.5	ug/l	413499	4	11.819	701497	50.0
Naphthalene, 1,...	16.417	31.6	ug/l	443583	4	11.819	701497	50.0