

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
 Data File : VU045006.D
 Acq On : 29 Sep 2021 20:23
 Operator : SY/MD
 Sample : M3926-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EP-1-MW-4-0921

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\82U092821W.M
 Title : SW846 8260

Signal : TIC: VU045006.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.295	1213	1230	1243	rBV	179942	380391	1.71%	0.525%
2	5.378	1243	1256	1275	rVB	316439	648927	2.92%	0.896%
3	5.706	1341	1358	1382	rBV	218232	442215	1.99%	0.611%
4	6.253	1512	1528	1548	rBV	475018	886181	3.99%	1.224%
5	7.902	2017	2041	2058	rBV	732096	1267371	5.70%	1.750%
6	9.420	2500	2513	2529	rBV	675534	1092214	4.91%	1.508%
7	9.571	2550	2560	2573	rVB	318589	507471	2.28%	0.701%
8	9.697	2574	2599	2614	rBV3	238713	526339	2.37%	0.727%
9	10.275	2754	2779	2788	rBV4	142840	421601	1.90%	0.582%
10	10.488	2827	2845	2855	rBV	1246976	1934765	8.70%	2.672%
11	10.635	2879	2891	2902	rVB	555816	871049	3.92%	1.203%
12	10.816	2929	2947	2964	rVB5	173084	363733	1.64%	0.502%
13	10.909	2964	2976	2987	rBV	1927774	2930039	13.18%	4.046%
14	11.002	2995	3005	3010	rBV	370420	630207	2.84%	0.870%
15	11.089	3020	3032	3041	rBV	303328	507914	2.29%	0.701%
16	11.295	3085	3096	3115	rBV	536427	1007198	4.53%	1.391%
17	11.420	3127	3135	3141	rBV2	280434	421394	1.90%	0.582%
18	11.475	3141	3152	3175	rVB	14899264	22226281	100.00%	30.692%
19	11.613	3176	3195	3199	rBV	386417	658862	2.96%	0.910%
20	11.645	3199	3205	3217	rVB	732383	1149334	5.17%	1.587%
21	11.748	3229	3237	3246	rVB	170987	262999	1.18%	0.363%
22	11.815	3246	3258	3271	rVV	693963	1146888	5.16%	1.584%
23	11.899	3271	3284	3306	rVV	4928651	7530905	33.88%	10.399%
24	12.012	3308	3319	3334	rVB	373221	636184	2.86%	0.879%
25	12.098	3334	3346	3363	rBV2	2725125	4398579	19.79%	6.074%
26	12.188	3364	3374	3392	rVB3	558167	1159376	5.22%	1.601%
27	12.304	3400	3410	3423	rBV	388415	626768	2.82%	0.866%
28	12.381	3425	3434	3449	rVB	240876	424390	1.91%	0.586%
29	12.468	3450	3461	3466	rBV	825351	1313541	5.91%	1.814%
30	12.500	3466	3471	3480	rVV	868363	1241730	5.59%	1.715%
31	12.574	3484	3494	3501	rVV	1511349	2275188	10.24%	3.142%
32	12.616	3501	3507	3518	rVV	530222	828716	3.73%	1.144%
33	12.687	3518	3529	3548	rVV3	1347551	2941263	13.23%	4.062%
34	12.877	3578	3588	3599	rVB5	511529	967639	4.35%	1.336%
35	12.976	3600	3619	3627	rBV	811880	1418397	6.38%	1.959%
36	13.031	3627	3636	3651	rVB	1008675	1625748	7.31%	2.245%

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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\82U092821W.M
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37	13.111	3652	3661	3676	rVB4	171821	312850	1.41%	0.432%
38	13.455	3758	3768	3784	rVB3	1257351	2183169	9.82%	3.015%
39	13.626	3811	3821	3839	rVB	567000	915258	4.12%	1.264%
40	13.841	3879	3888	3899	rBV5	143581	254109	1.14%	0.351%
41	13.944	3916	3920	3943	rVB2	114798	231721	1.04%	0.320%
42	14.089	3949	3965	3987	rVB	486234	847568	3.81%	1.170%

Sum of corrected areas: 72416472

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EP-1-MW-4-0921

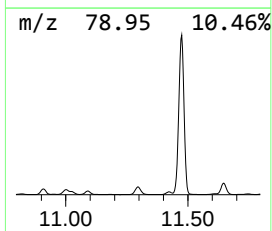
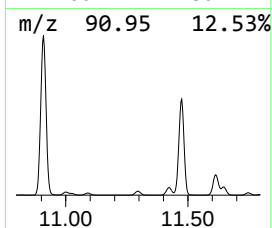
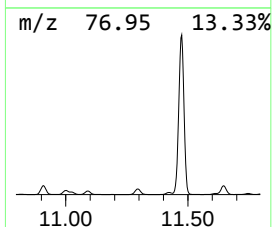
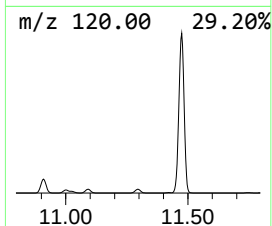
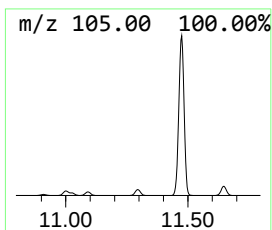
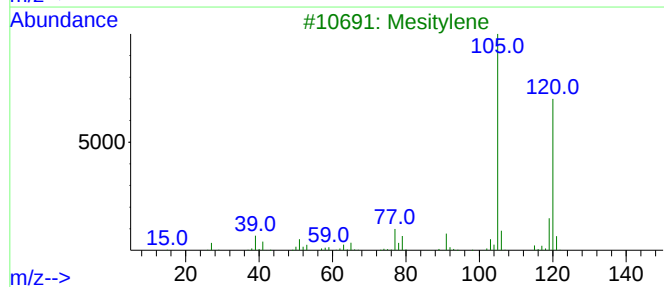
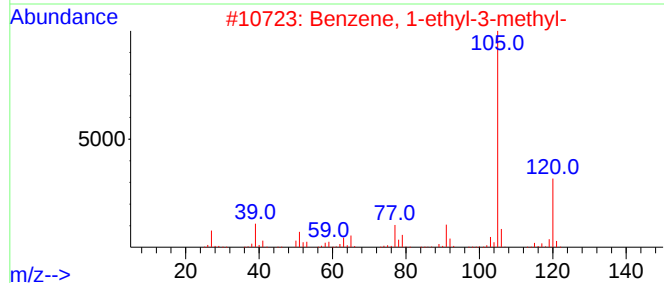
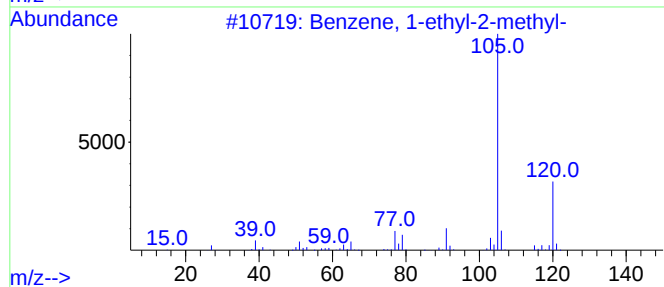
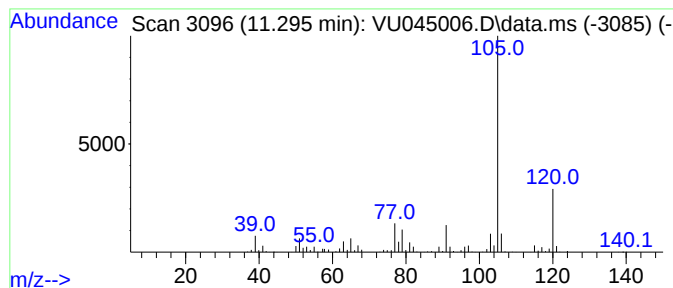
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.295	43.91 ug/l	1007200	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

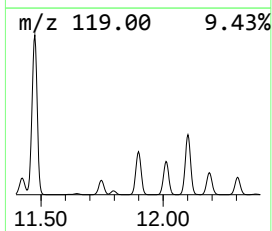
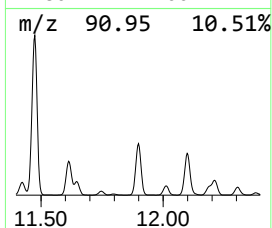
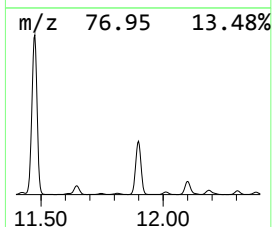
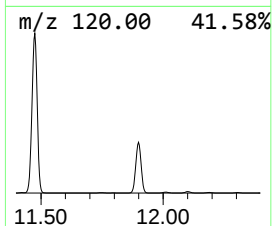
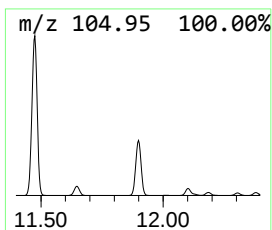
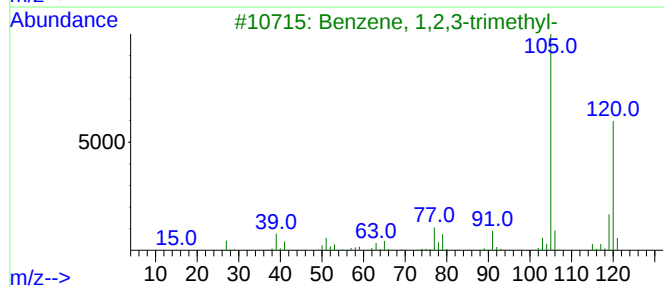
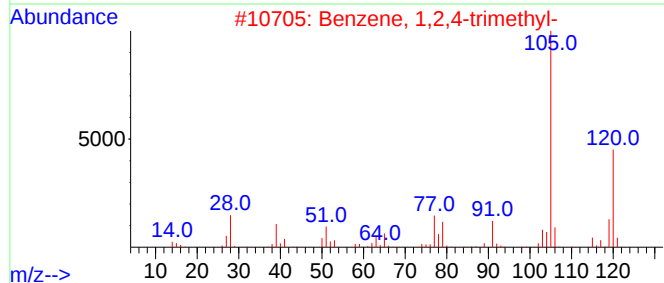
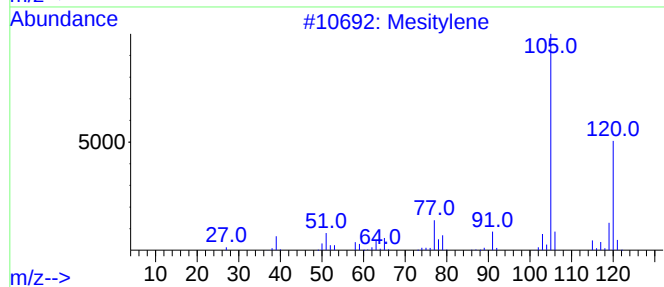
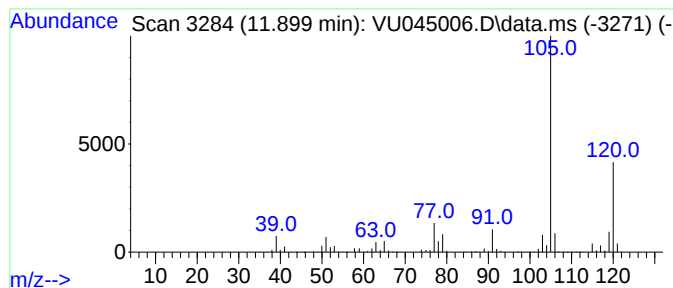
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.899	328.32 ug/l	7530910	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

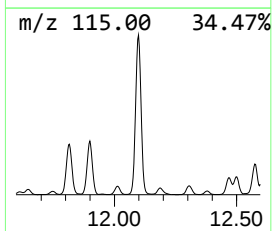
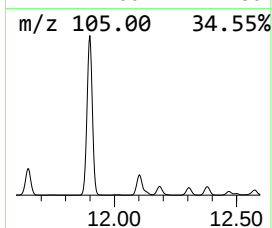
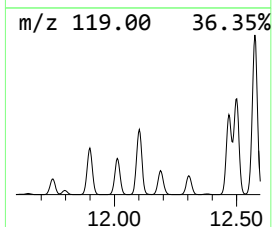
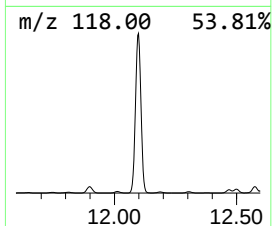
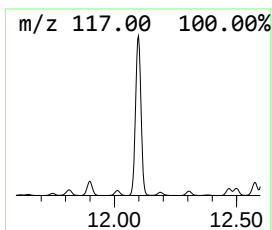
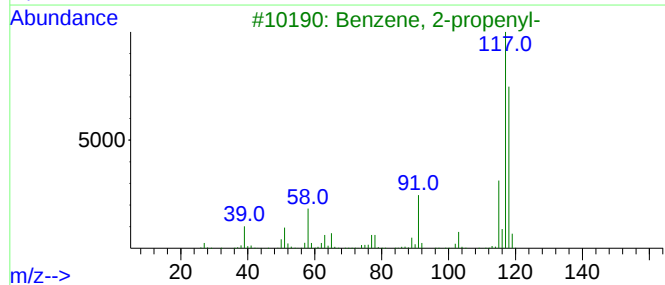
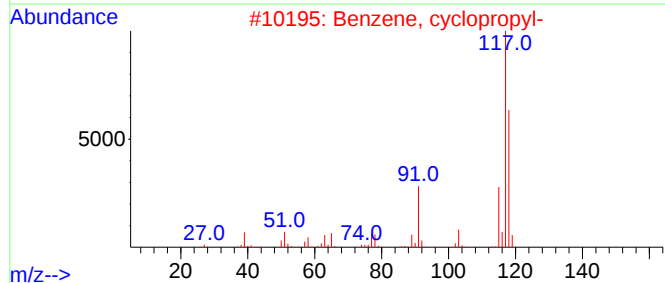
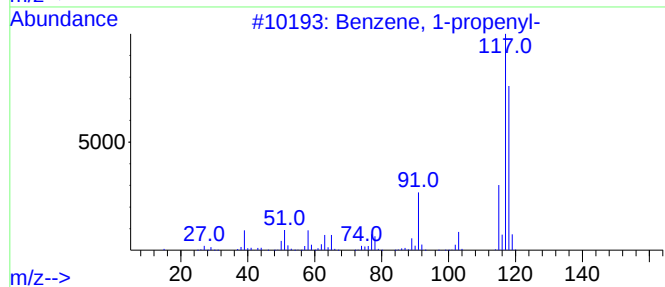
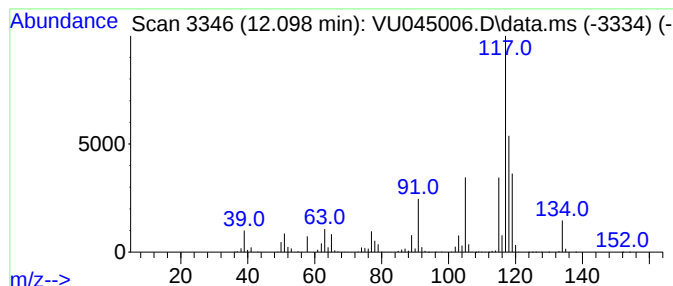
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-propenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.098	191.76 ug/l	4398580	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	92
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
5			Indane	118	C9H10	000496-11-7	60



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Instrument :
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ClientSampleId :
EP-1-MW-4-0921

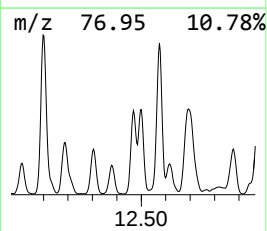
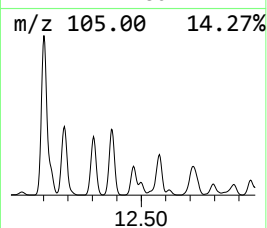
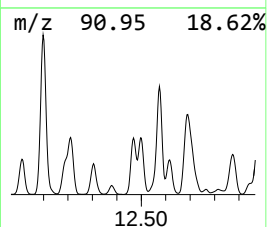
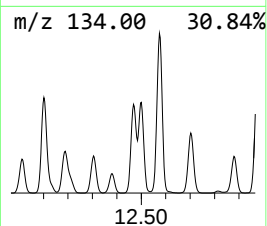
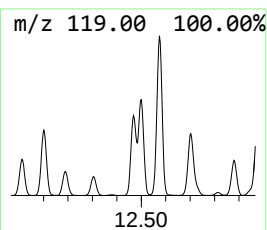
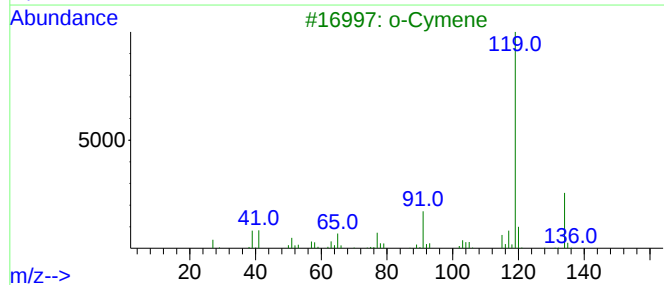
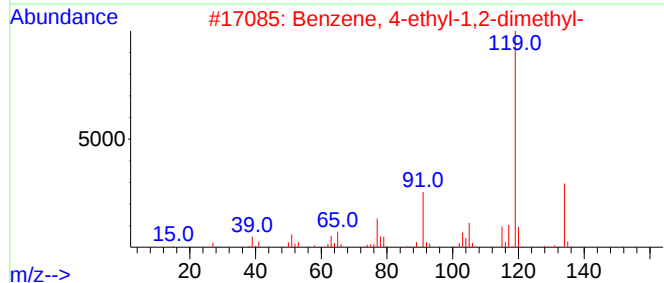
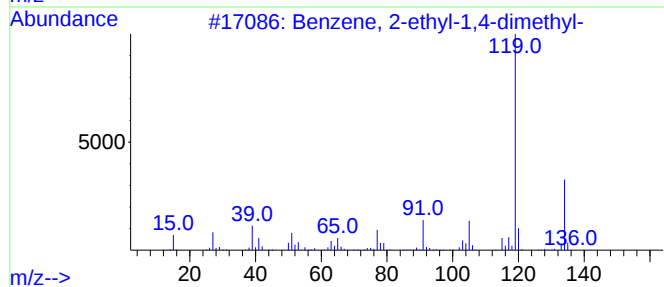
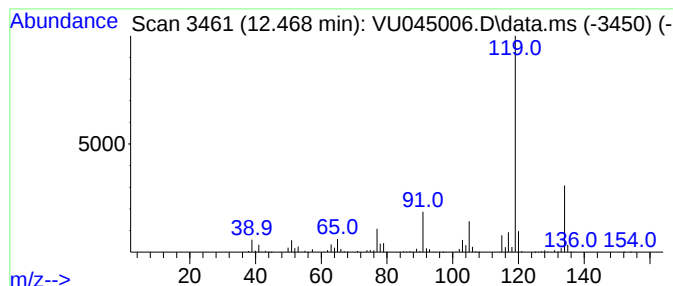
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.468	57.27 ug/l	1313540	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

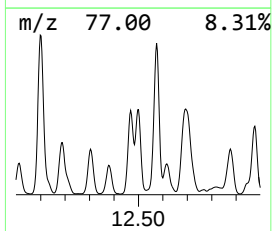
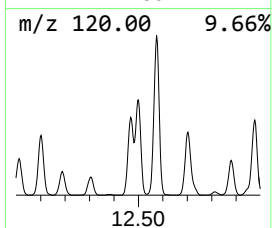
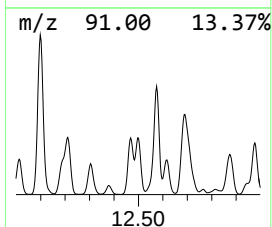
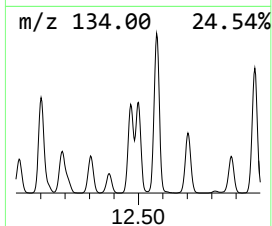
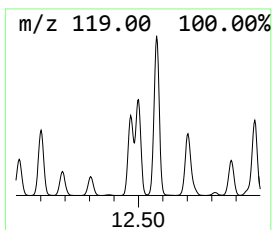
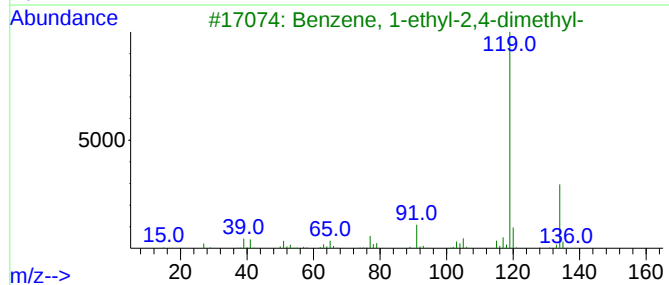
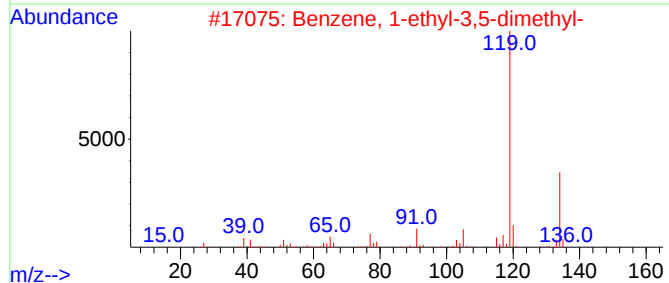
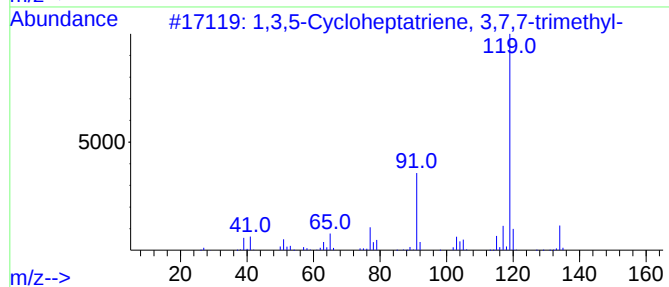
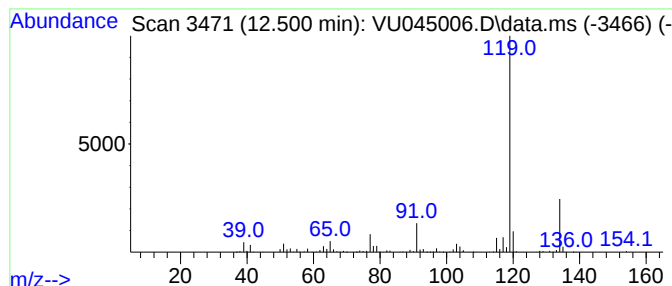
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,3,5-Cycloheptatriene, 3,7... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.500	54.13 ug/l	1241730	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	91
2			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045006.D
Acq On : 29 Sep 2021 20:23
Operator : SY/MD
Sample : M3926-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

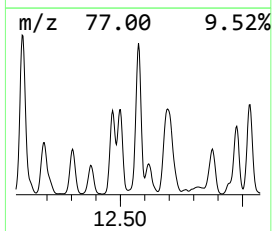
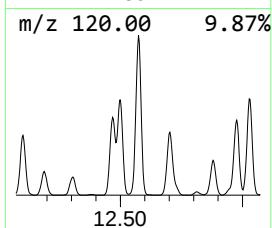
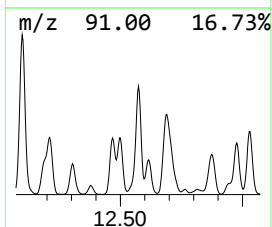
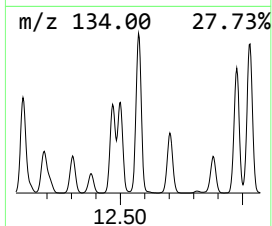
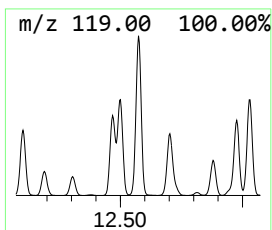
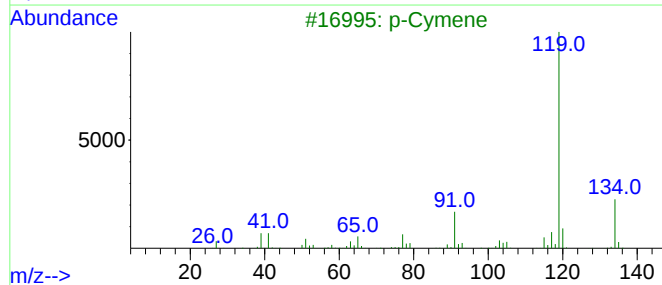
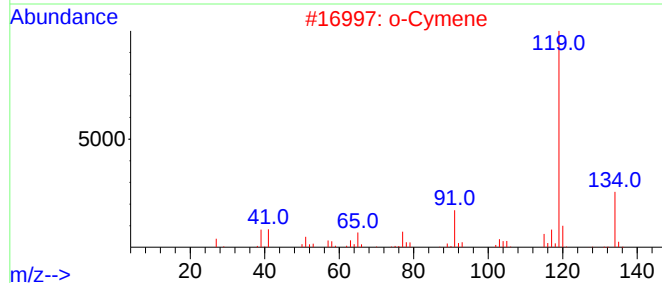
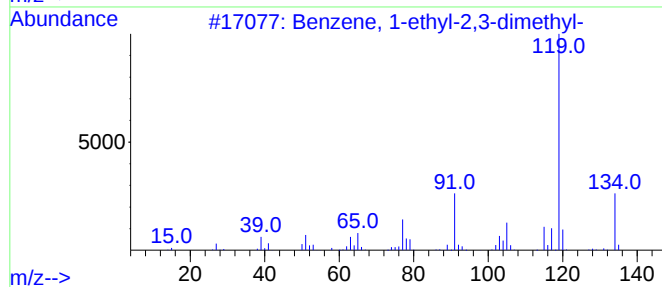
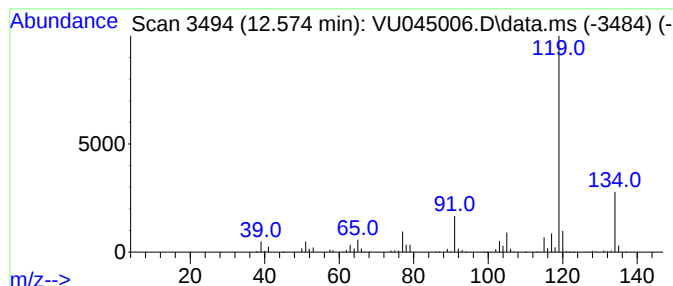
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.574	99.19 ug/l	2275190	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			o-Cymene	134	C10H14	000527-84-4	95
3			p-Cymene	134	C10H14	000099-87-6	94
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045006.D
Acq On : 29 Sep 2021 20:23
Operator : SY/MD
Sample : M3926-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

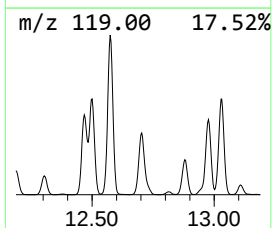
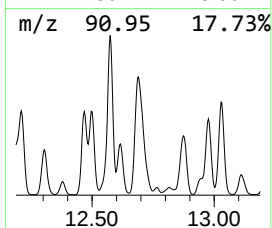
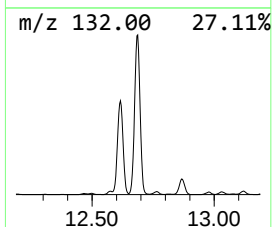
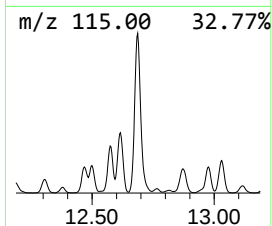
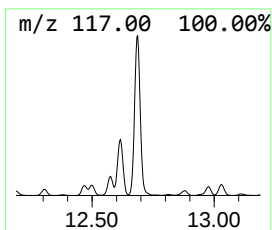
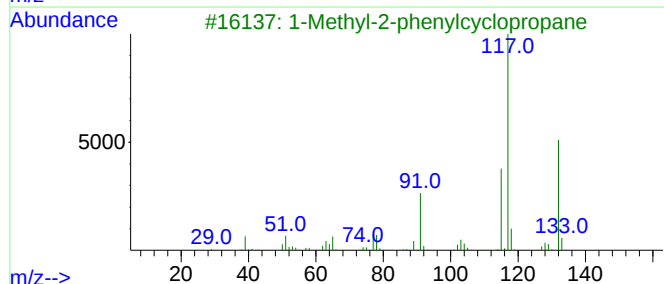
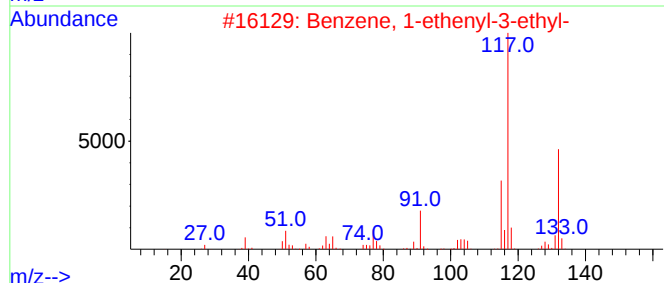
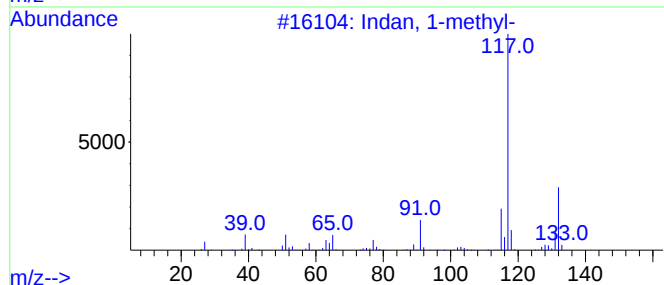
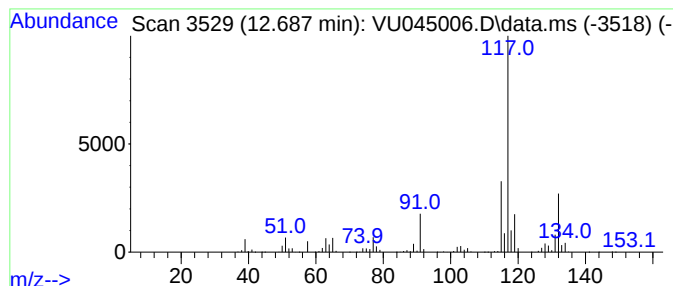
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	128.23 ug/l	2941260	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	94
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045006.D
Acq On : 29 Sep 2021 20:23
Operator : SY/MD
Sample : M3926-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

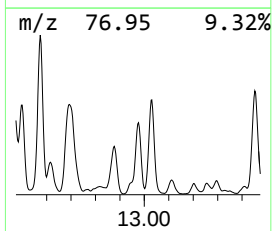
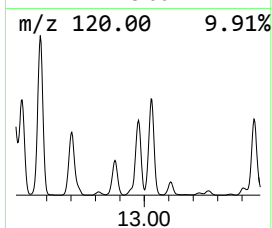
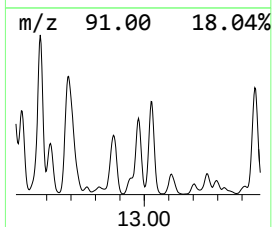
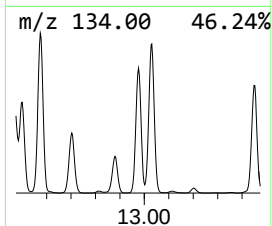
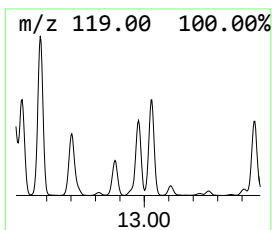
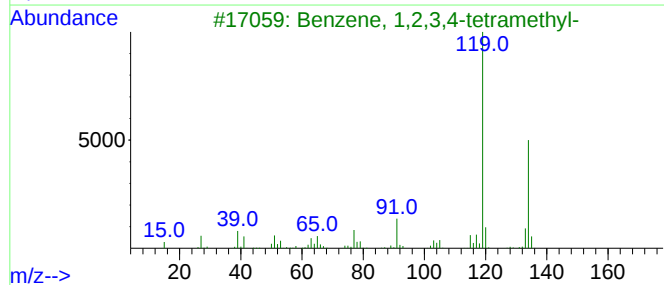
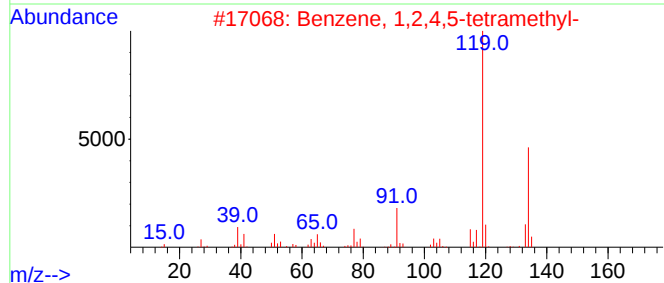
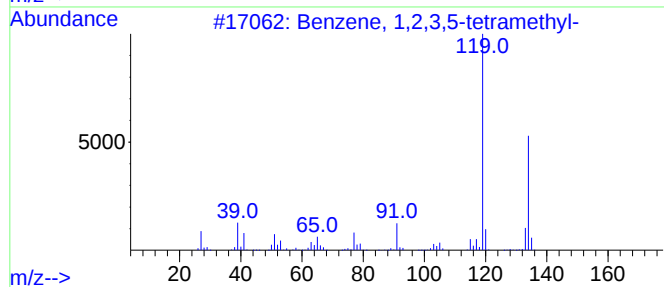
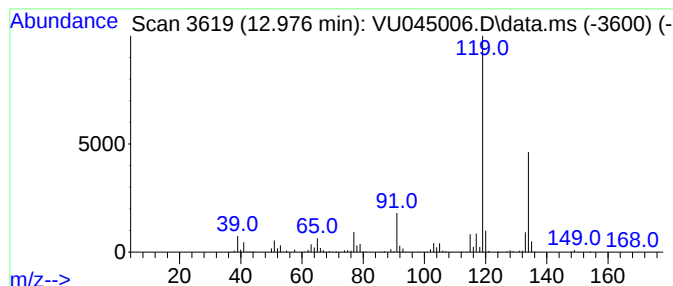
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.976	61.84 ug/l	1418400	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045006.D
Acq On : 29 Sep 2021 20:23
Operator : SY/MD
Sample : M3926-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

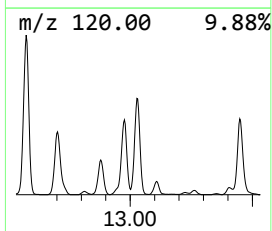
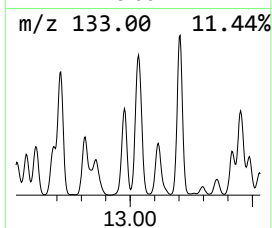
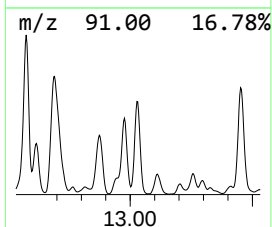
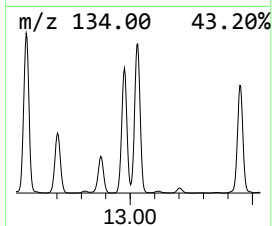
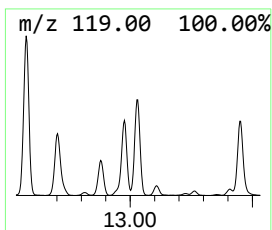
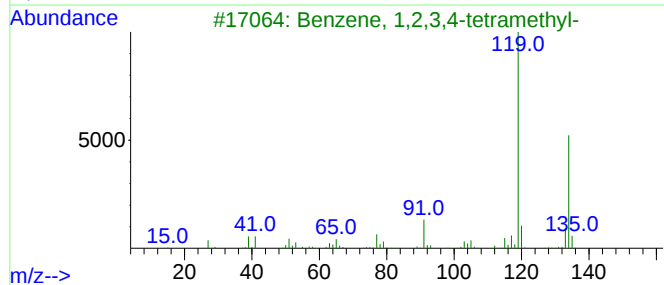
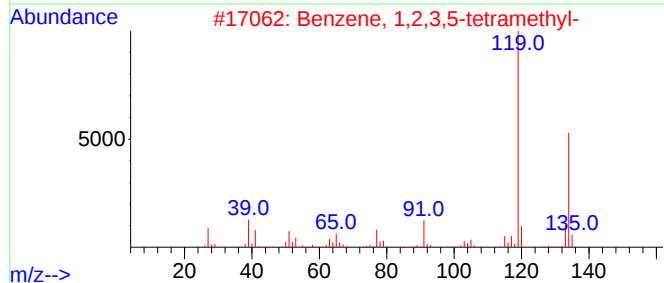
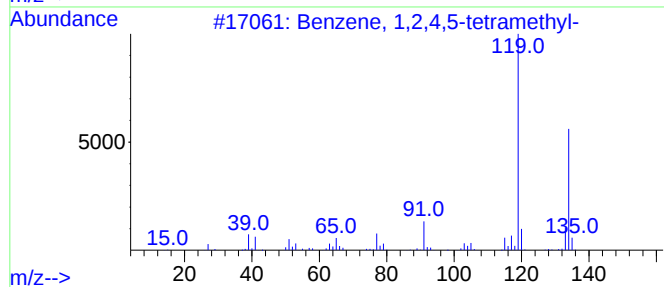
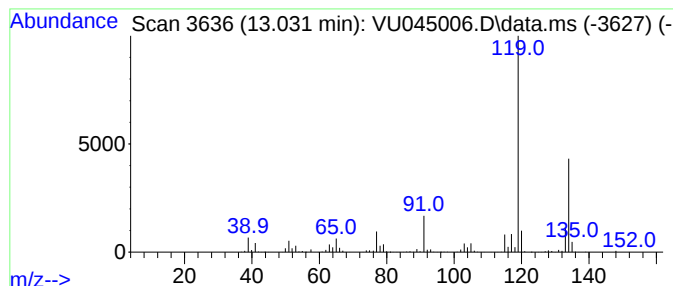
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.031	70.88 ug/l	1625750	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045006.D
Acq On : 29 Sep 2021 20:23
Operator : SY/MD
Sample : M3926-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

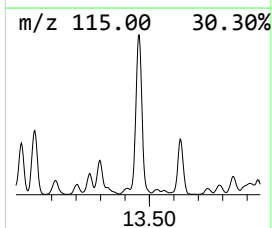
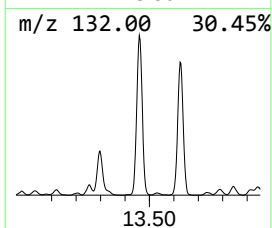
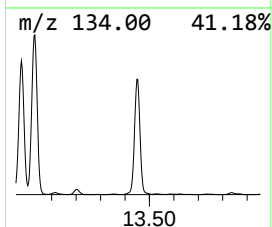
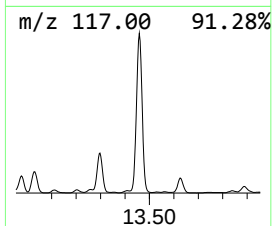
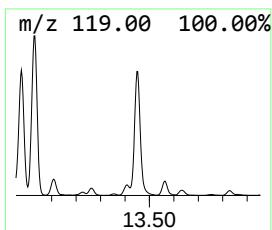
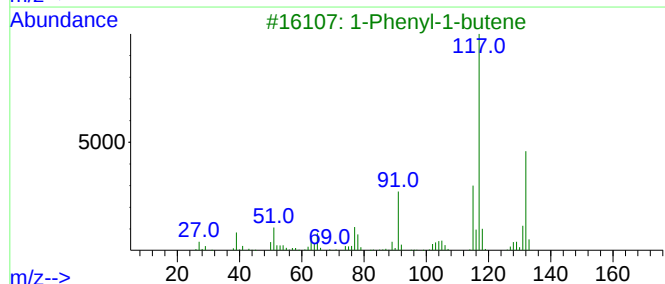
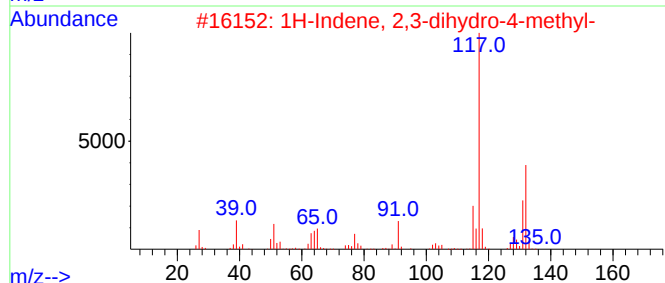
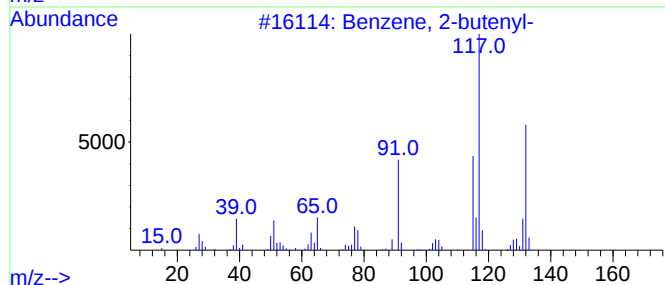
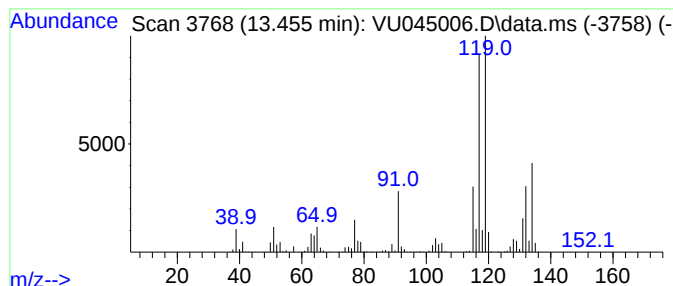
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 2-butenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.455	95.18 ug/l	2183170	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	86
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045006.D
Acq On : 29 Sep 2021 20:23
Operator : SY/MD
Sample : M3926-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP-1-MW-4-0921

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1,2,3-...	11.899	328.3	ug/l	7530910	4	11.816	1146890	50.0
Benzene, 1-prop...	12.098	191.8	ug/l	4398580	4	11.816	1146890	50.0
Benzene, 2-ethy...	12.468	57.3	ug/l	1313540	4	11.816	1146890	50.0
1,3,5-Cyclohept...	12.500	54.1	ug/l	1241730	4	11.816	1146890	50.0
Benzene, 1-ethy...	12.574	99.2	ug/l	2275190	4	11.816	1146890	50.0
Indan, 1-methyl-	12.687	128.2	ug/l	2941260	4	11.816	1146890	50.0
Benzene, 1,2,3,...	12.976	61.8	ug/l	1418400	4	11.816	1146890	50.0
Benzene, 1,2,4,...	13.031	70.9	ug/l	1625750	4	11.816	1146890	50.0
Benzene, 2-bute...	13.455	95.2	ug/l	2183170	4	11.816	1146890	50.0