

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
 Data File : VU045007.D
 Acq On : 29 Sep 2021 20:47
 Operator : SY/MD
 Sample : M3926-04DL 10X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 24 Sample Multiplier: 1

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

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 1 % 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: VU045007.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.475	34	42	55	rBV	163127	162478	7.50%	1.243%
2	5.295	1214	1230	1242	rBV	177760	374607	17.29%	2.865%
3	5.379	1242	1256	1278	rVB	310526	639442	29.51%	4.891%
4	5.707	1343	1358	1381	rBV	214393	438772	20.25%	3.356%
5	6.253	1511	1528	1546	rBV	470004	879463	40.59%	6.727%
6	7.903	2022	2041	2062	rBV	724813	1233699	56.94%	9.436%
7	9.420	2498	2513	2529	rBV	655501	1068848	49.33%	8.175%
8	9.575	2550	2561	2576	rVB	30745	48559	2.24%	0.371%
9	9.697	2584	2599	2615	rBV2	24462	47104	2.17%	0.360%
10	10.488	2826	2845	2855	rBV	120454	186520	8.61%	1.427%
11	10.636	2877	2891	2904	rBV	523294	814514	37.59%	6.230%
12	10.816	2939	2947	2957	rVB6	12245	21730	1.00%	0.166%
13	10.909	2964	2976	2994	rBV	187631	289664	13.37%	2.216%
14	11.002	2995	3005	3010	rBV	38908	64873	2.99%	0.496%
15	11.092	3022	3033	3043	rVB	33408	51841	2.39%	0.397%
16	11.295	3085	3096	3114	rBV	55618	98337	4.54%	0.752%
17	11.420	3127	3135	3140	rBV3	24624	35978	1.66%	0.275%
18	11.472	3140	3151	3173	rVB	1408321	2166780	100.00%	16.573%
19	11.613	3185	3195	3199	rBV2	33972	52242	2.41%	0.400%
20	11.645	3199	3205	3218	rVB	71767	112068	5.17%	0.857%
21	11.748	3229	3237	3245	rVB	16727	24152	1.11%	0.185%
22	11.816	3245	3258	3271	rBV	656515	1048915	48.41%	8.023%
23	11.899	3272	3284	3305	rVB	517057	800892	36.96%	6.126%
24	12.012	3308	3319	3331	rVB	38051	61451	2.84%	0.470%
25	12.099	3333	3346	3363	rBV2	268829	441823	20.39%	3.379%
26	12.189	3363	3374	3394	rVB3	50128	108538	5.01%	0.830%
27	12.304	3396	3410	3425	rBV	39852	65878	3.04%	0.504%
28	12.381	3425	3434	3445	rVB	21058	33491	1.55%	0.256%
29	12.468	3449	3461	3466	rBV	79369	126377	5.83%	0.967%
30	12.497	3466	3470	3481	rVV	80168	114028	5.26%	0.872%
31	12.574	3484	3494	3501	rVV	139621	214800	9.91%	1.643%
32	12.616	3501	3507	3518	rVB	54276	84808	3.91%	0.649%
33	12.687	3518	3529	3548	rBV3	137199	298912	13.80%	2.286%
34	12.877	3577	3588	3599	rVB4	52325	97323	4.49%	0.744%
35	12.976	3601	3619	3627	rVV	79821	139086	6.42%	1.064%
36	13.028	3627	3635	3652	rVB	102067	159644	7.37%	1.221%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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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Title : SW846 8260

37	13.115	3652	3662	3676	rBV3	15457	27261	1.26%	0.209%
38	13.455	3757	3768	3784	rVB3	123836	219087	10.11%	1.676%
39	13.626	3809	3821	3831	rVB	55996	89509	4.13%	0.685%
40	13.844	3879	3889	3899	rBV6	14163	23999	1.11%	0.184%
41	14.089	3948	3965	3984	rBV	62458	106817	4.93%	0.817%

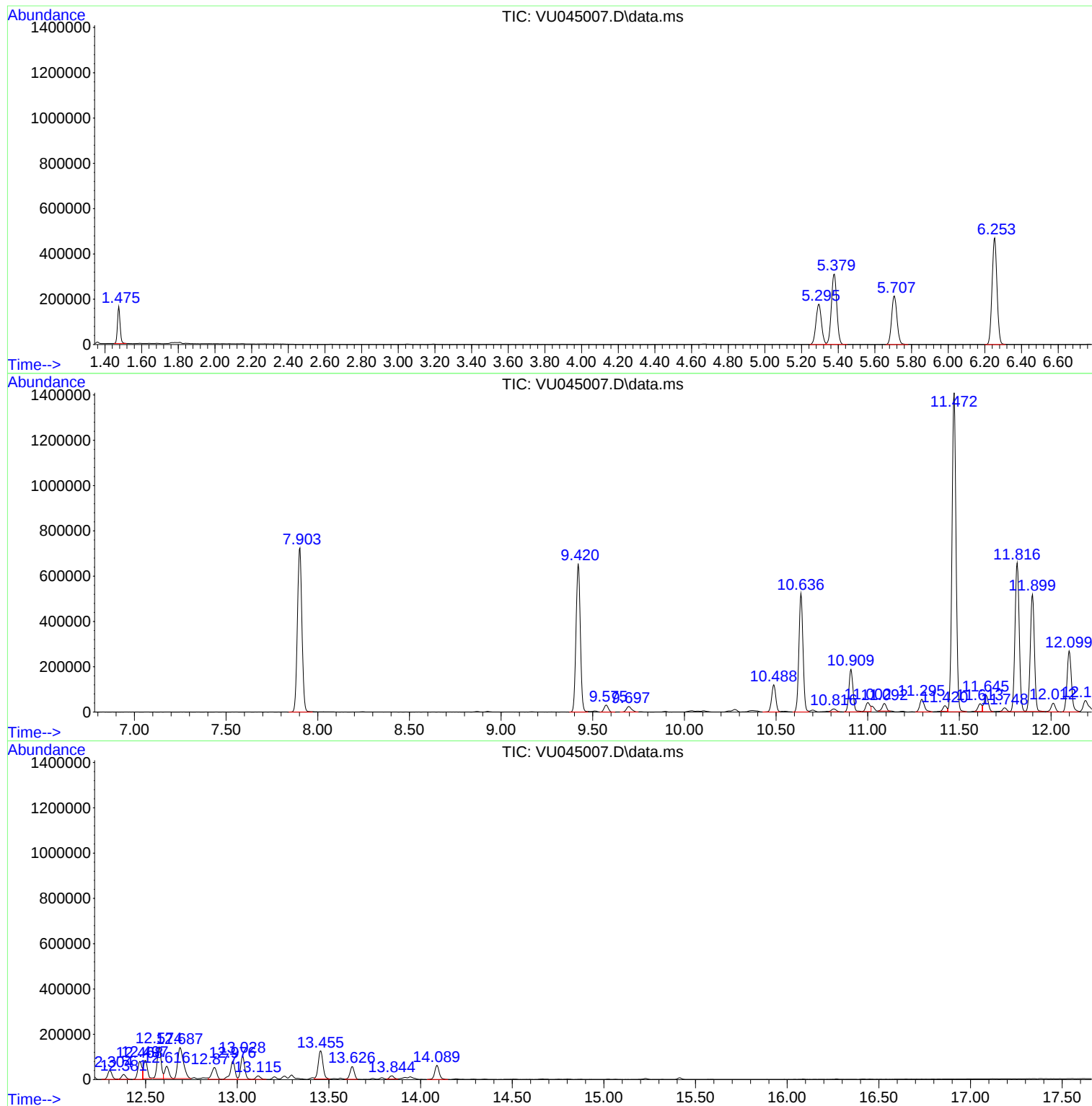
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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



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

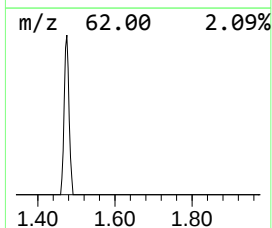
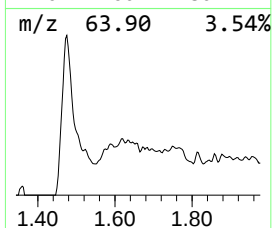
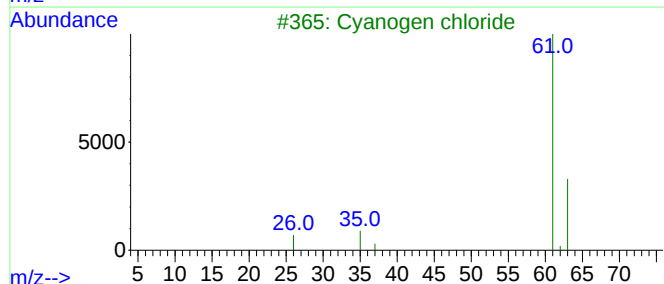
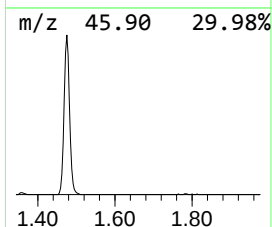
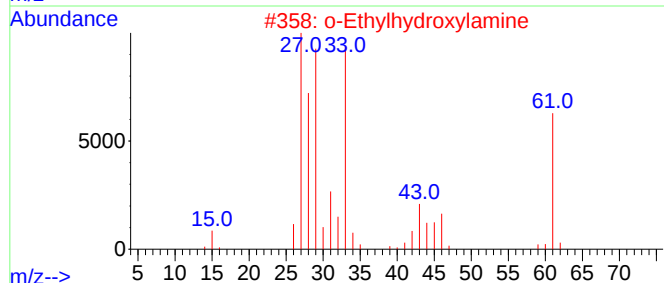
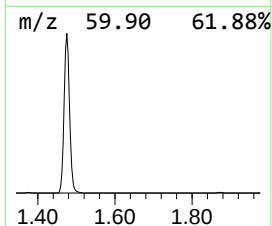
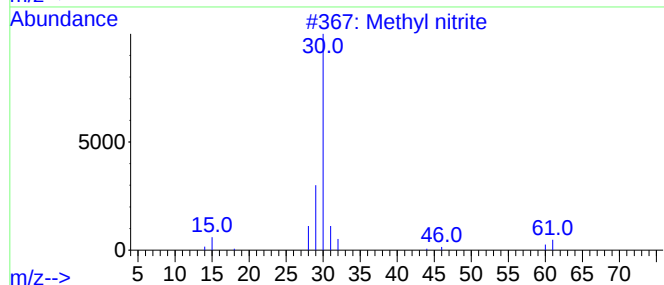
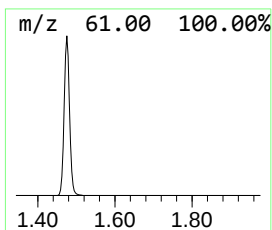
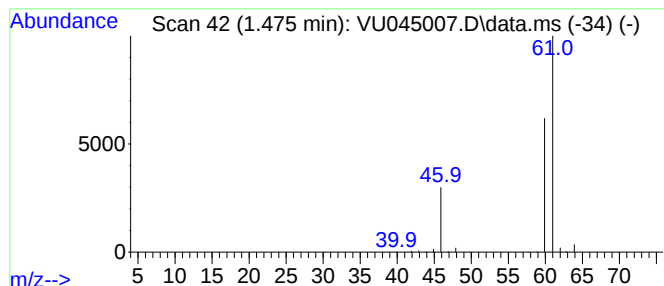
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 1 Methyl nitrite Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.475	12.70 ug/l	162478	Pentafluorobenzene	5.379

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl nitrite	61	CH3NO2	000624-91-9	5
2			o-Ethylhydroxylamine	61	C2H7NO	000624-86-2	4
3			Cyanogen chloride	61	CClN	000506-77-4	4
4			Thiirane	60	C2H4S	000420-12-2	3
5			Pentaborane(11)	66	B5H11	018433-84-6	3



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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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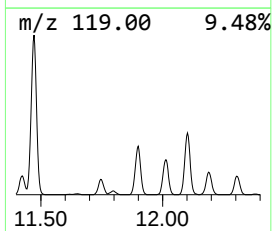
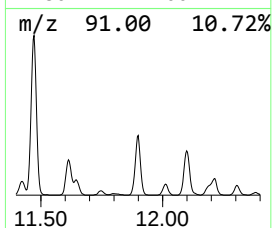
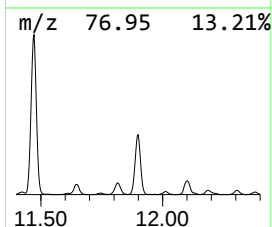
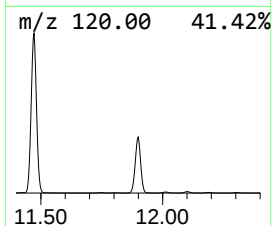
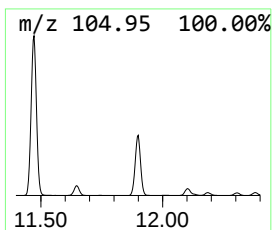
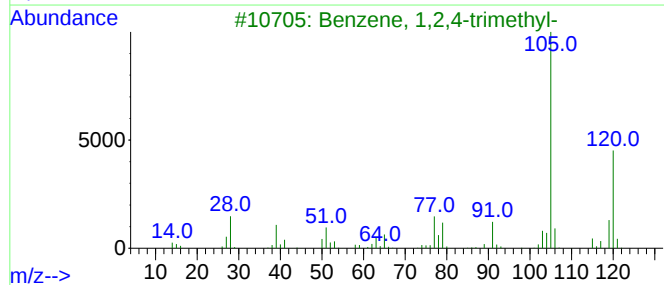
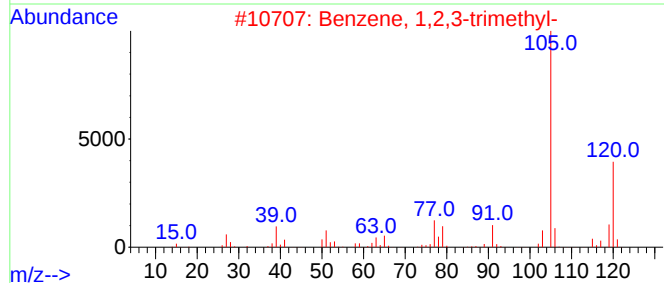
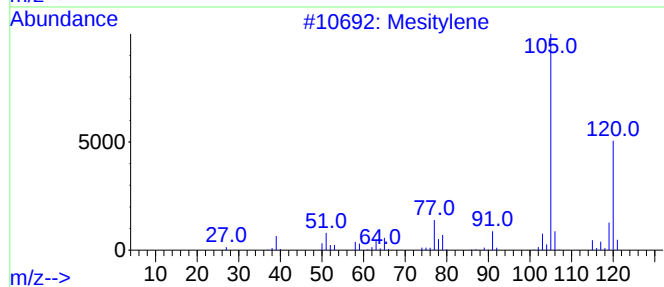
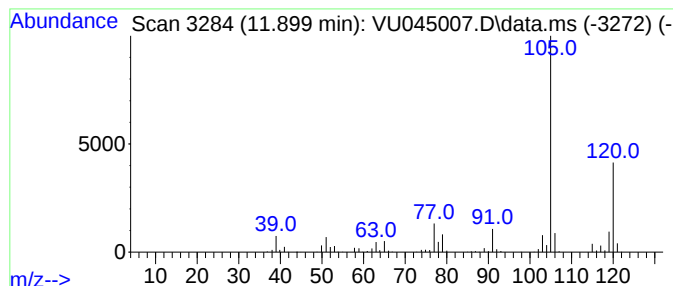
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Mesitylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.899	38.18 ug/l	800892	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



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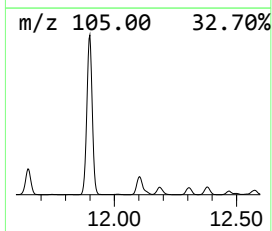
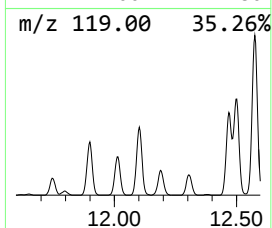
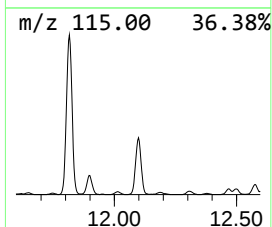
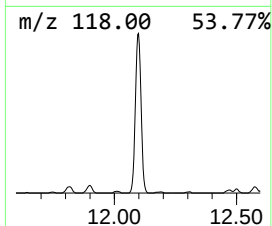
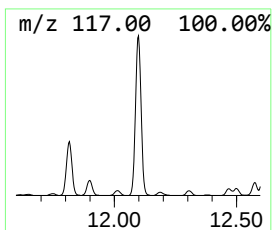
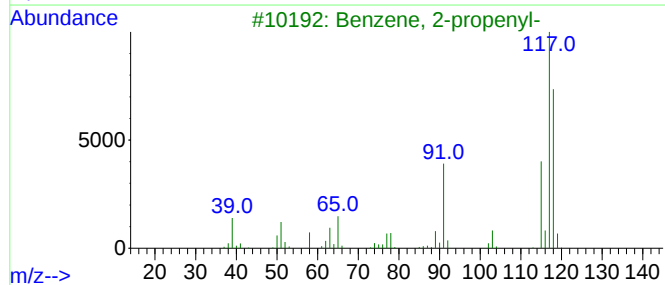
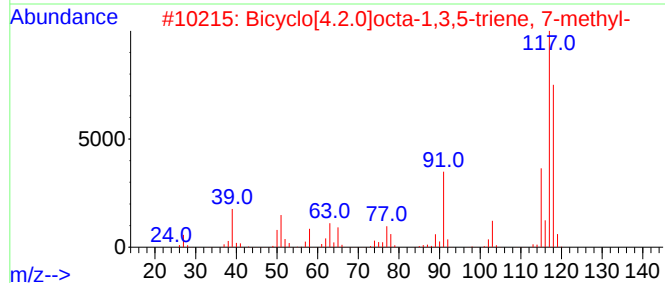
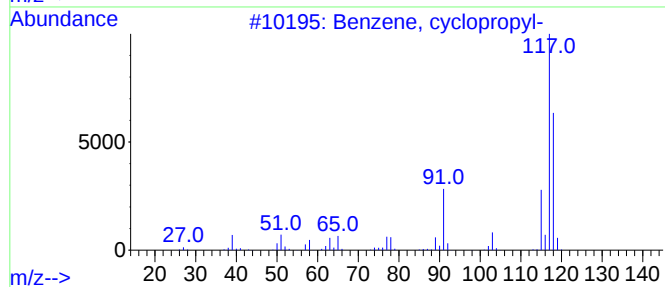
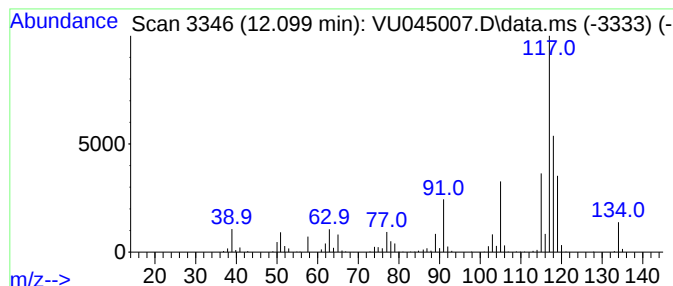
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.099	21.06 ug/l	441823	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, cyclopropyl-	118	C9H10	000873-49-4	91
2			Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	83
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64



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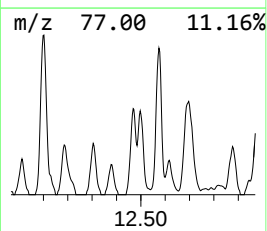
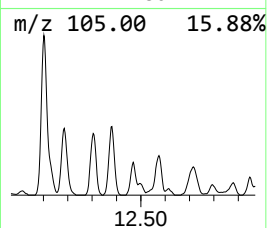
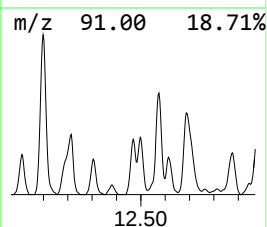
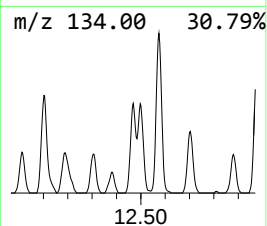
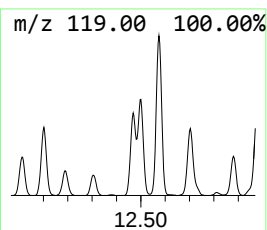
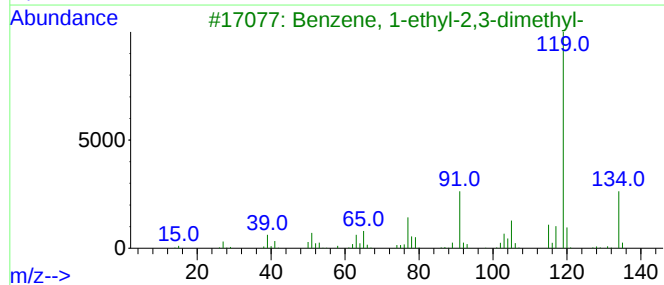
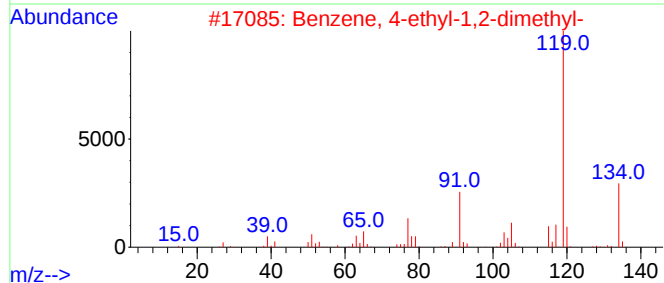
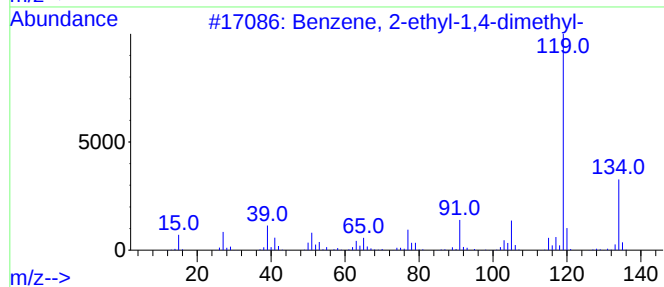
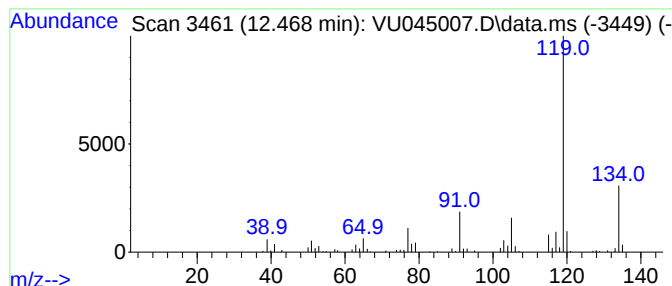
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.468	6.02 ug/l	126377	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	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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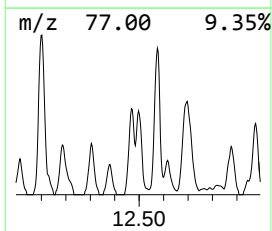
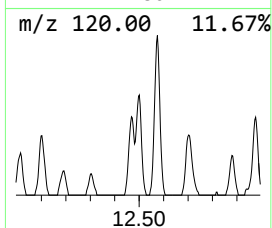
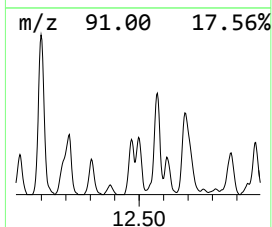
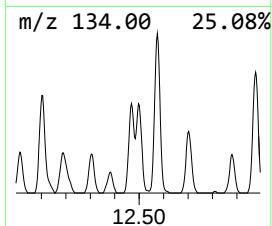
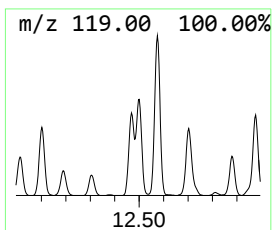
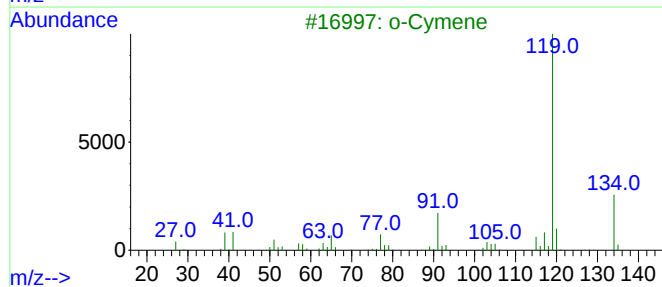
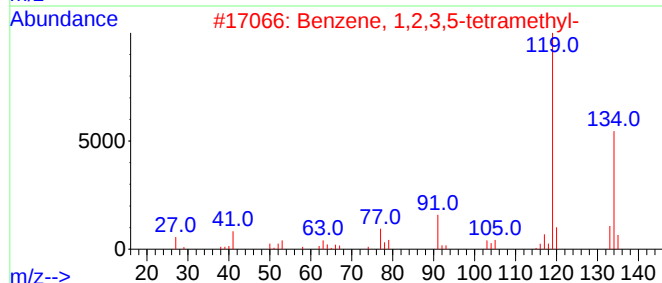
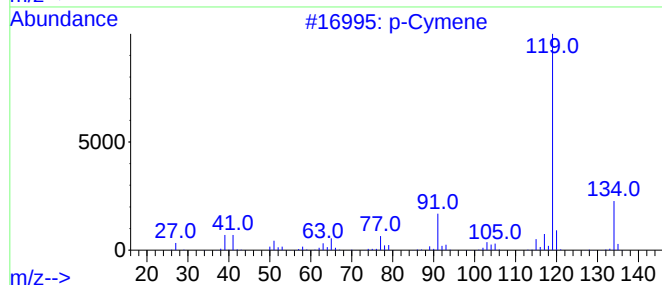
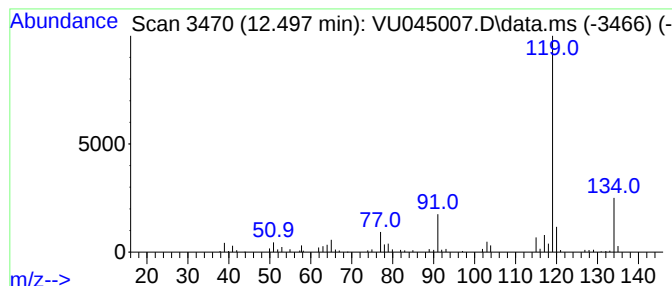
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 5 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.497	5.44 ug/l	114028	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	94
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
3			o-Cymene	134	C10H14	000527-84-4	91
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045007.D
Acq On : 29 Sep 2021 20:47
Operator : SY/MD
Sample : M3926-04DL 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

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

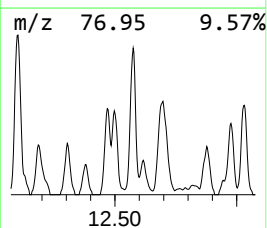
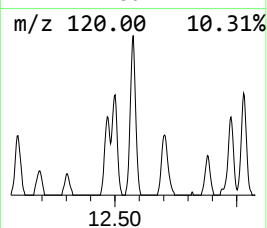
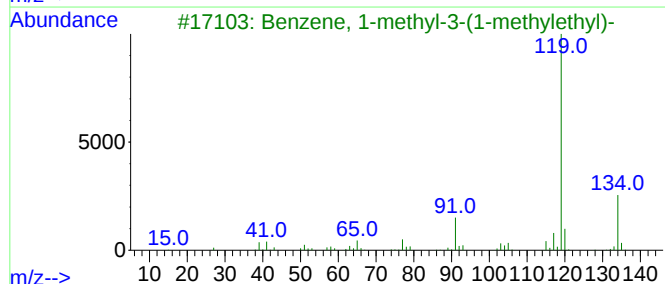
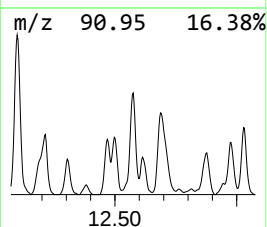
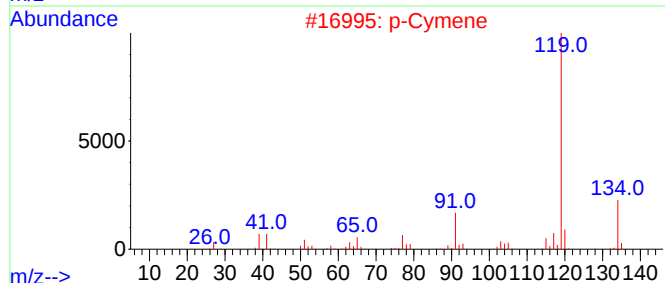
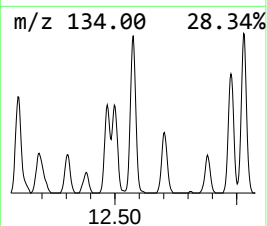
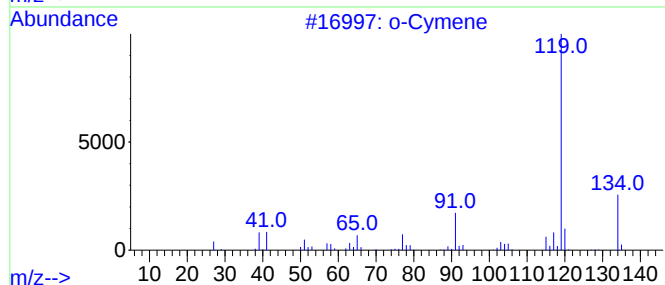
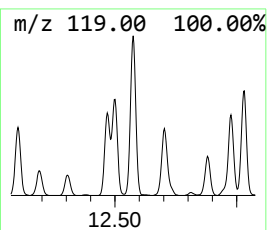
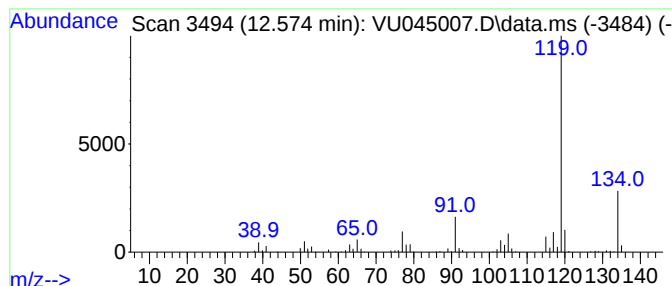
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.574	10.24 ug/l	214800	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			p-Cymene	134	C10H14	000099-87-6	97
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045007.D
Acq On : 29 Sep 2021 20:47
Operator : SY/MD
Sample : M3926-04DL 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

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

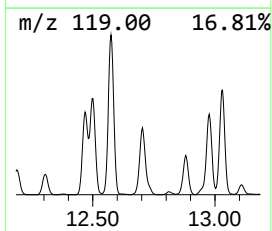
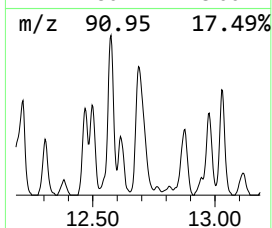
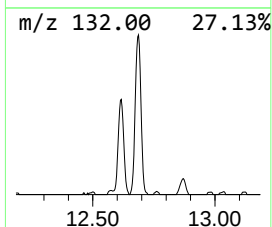
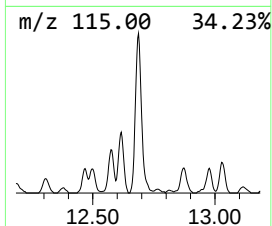
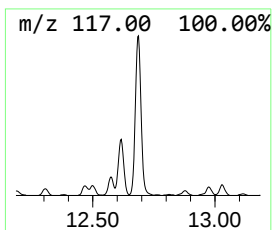
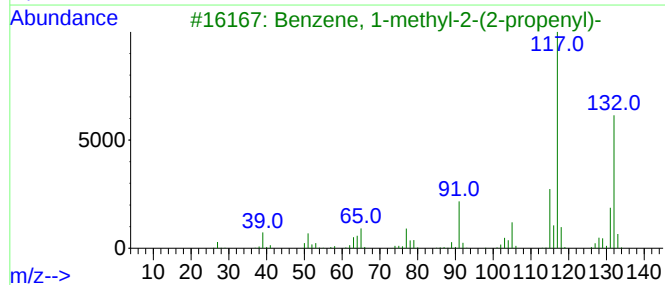
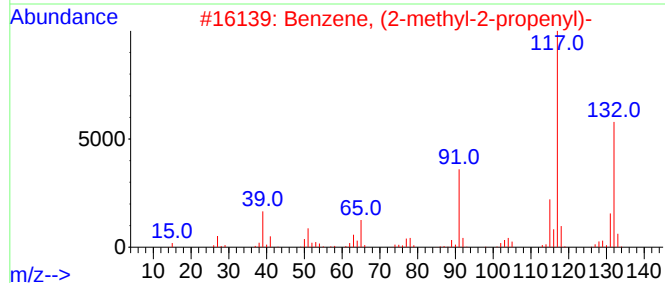
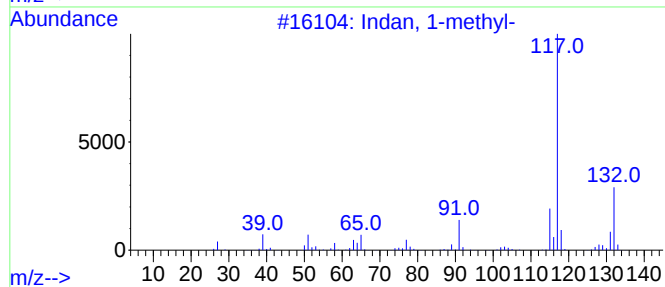
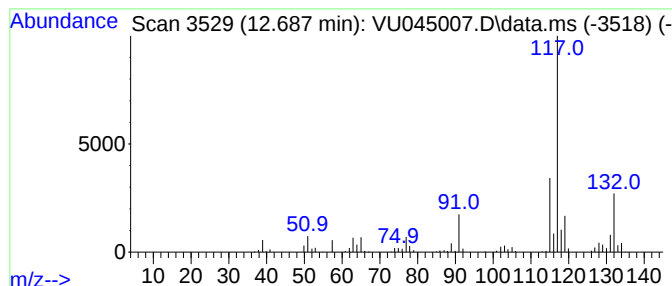
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Quant Title : SW846 8260

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	14.25 ug/l	298912	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045007.D
Acq On : 29 Sep 2021 20:47
Operator : SY/MD
Sample : M3926-04DL 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

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

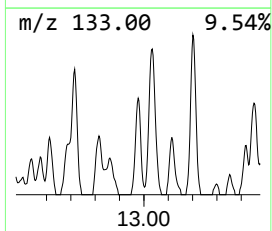
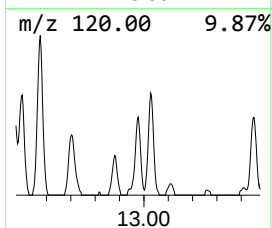
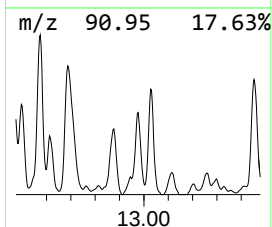
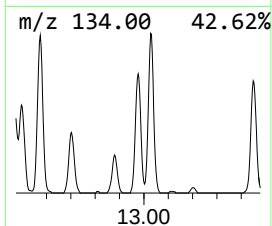
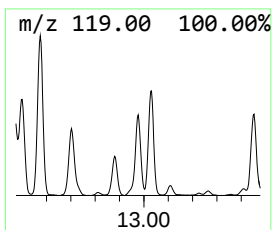
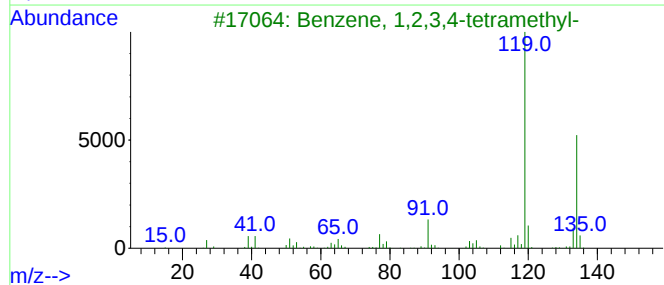
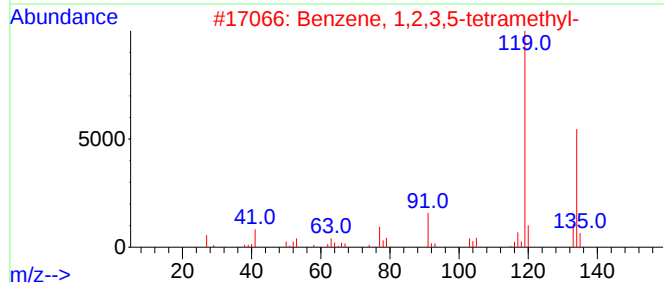
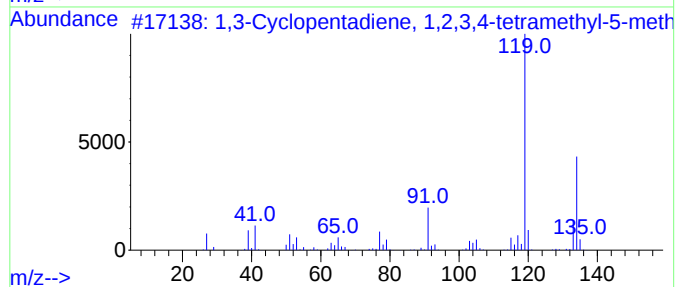
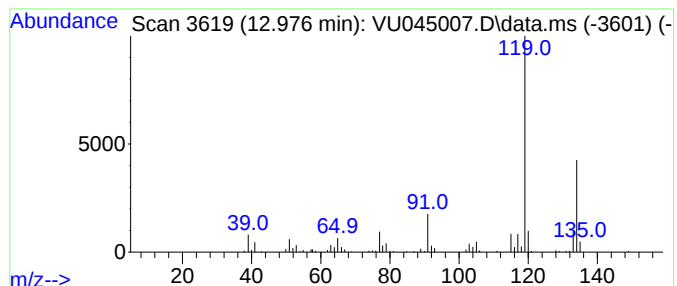
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 8 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.976	6.63 ug/l	139086	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045007.D
Acq On : 29 Sep 2021 20:47
Operator : SY/MD
Sample : M3926-04DL 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

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

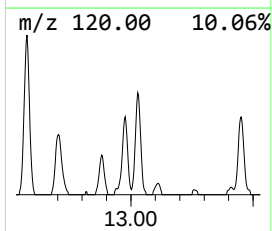
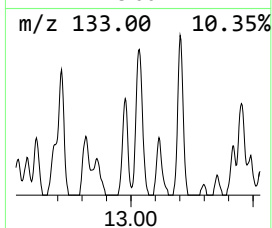
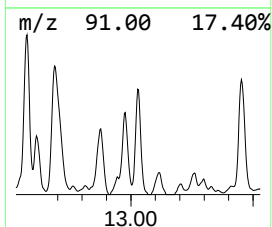
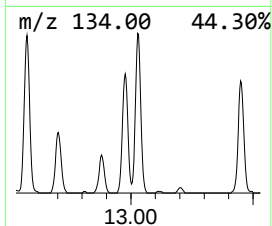
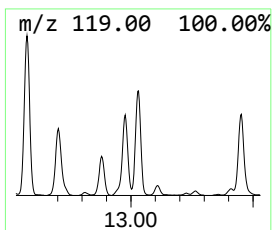
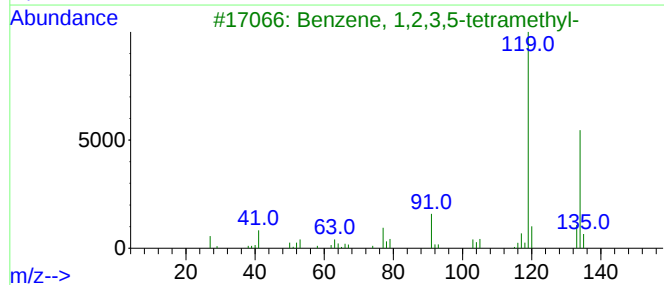
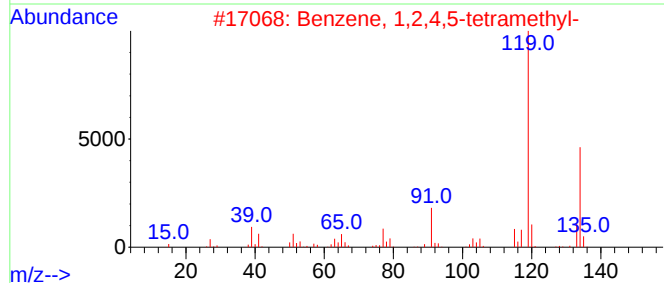
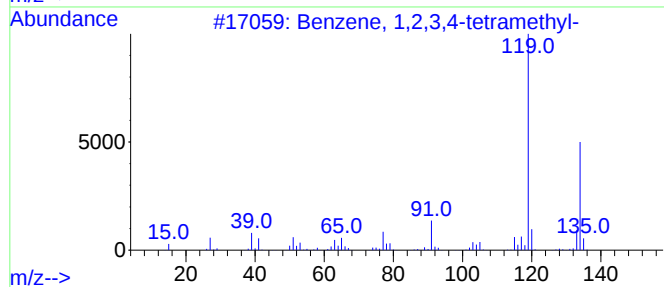
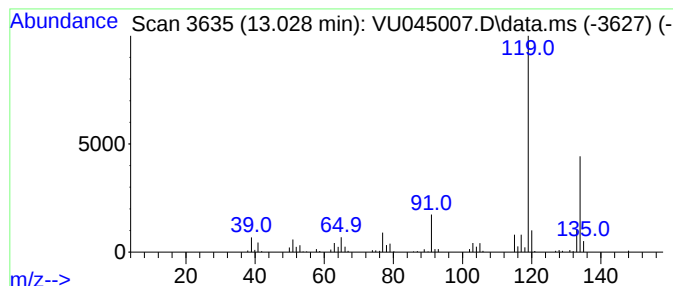
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	7.61 ug/l	159644	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			o-Cymene	134	C10H14	000527-84-4	94
5			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU045007.D
Acq On : 29 Sep 2021 20:47
Operator : SY/MD
Sample : M3926-04DL 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

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

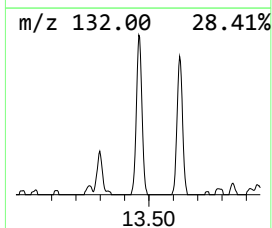
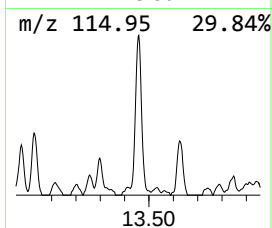
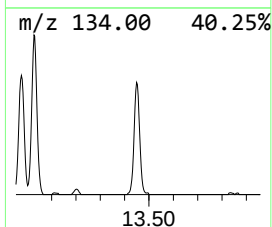
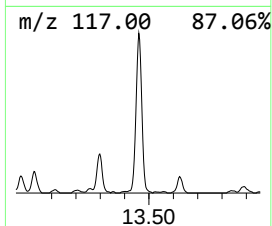
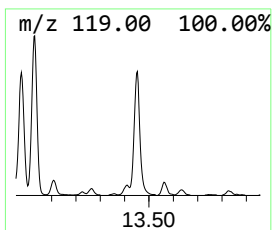
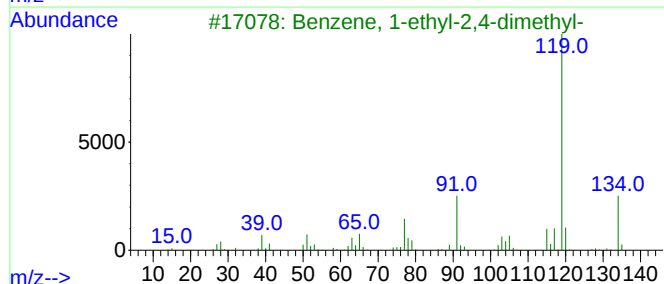
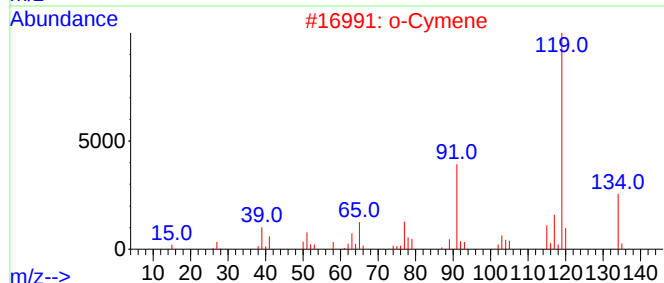
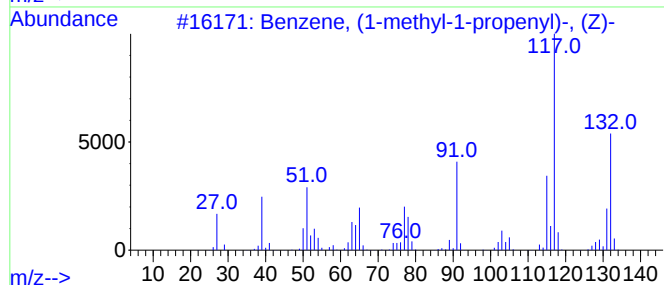
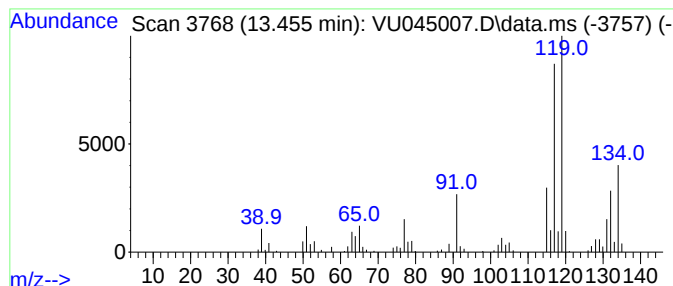
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1-methyl-1-propen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.455	10.44 ug/l	219087	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
2			o-Cymene	134	C10H14	000527-84-4	55
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	50
4			p-Cymene	134	C10H14	000099-87-6	50
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
 Data File : VU045007.D
 Acq On : 29 Sep 2021 20:47
 Operator : SY/MD
 Sample : M3926-04DL 10X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 24 Sample Multiplier: 1

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

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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Mesitylene	11.899	38.2	ug/l	800892	4	11.816	1048920	50.0
Benzene, cyclop...	12.099	21.1	ug/l	441823	4	11.816	1048920	50.0
Benzene, 2-ethy...	12.468	6.0	ug/l	126377	4	11.816	1048920	50.0
p-Cymene	12.497	5.4	ug/l	114028	4	11.816	1048920	50.0
o-Cymene	12.574	10.2	ug/l	214800	4	11.816	1048920	50.0
Indan, 1-methyl-	12.687	14.2	ug/l	298912	4	11.816	1048920	50.0
1,3-Cyclopentad...	12.976	6.6	ug/l	139086	4	11.816	1048920	50.0
Benzene, 1,2,3,...	13.028	7.6	ug/l	159644	4	11.816	1048920	50.0
Benzene, (1-met...	13.455	10.4	ug/l	219087	4	11.816	1048920	50.0