

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070920\
 Data File : VU039419.D
 Acq On : 10 Jul 2020 01:29
 Operator : SY/MD
 Sample : L3261-07
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BE506

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
 Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.388	314	326	347	rVB	980834	1443423	1.25%	0.501%
2	4.015	812	832	854	rBV2	1202423	3111657	2.70%	1.081%
3	4.687	1020	1041	1075	rVB2	574113	1443967	1.25%	0.501%
4	5.800	1350	1387	1415	rBV2	31130458	60642547	52.53%	21.061%
5	7.996	2052	2070	2124	rBV3	59505834	115438918	100.00%	40.092%
6	9.452	2502	2523	2548	rBV	7677395	12514142	10.84%	4.346%
7	9.578	2548	2562	2585	rVB	10054578	15941852	13.81%	5.537%
8	9.703	2585	2601	2616	rBV	25809009	41496907	35.95%	14.412%
9	10.105	2711	2726	2753	rVB	7554839	11968194	10.37%	4.157%
10	10.484	2828	2844	2857	rBV	1873600	2926205	2.53%	1.016%
11	10.902	2957	2974	2990	rBV	848242	1626979	1.41%	0.565%
12	10.996	2990	3003	3009	rBV2	1033443	1782185	1.54%	0.619%
13	11.288	3084	3094	3116	rVB	905475	1474097	1.28%	0.512%
14	11.465	3138	3149	3165	rVB	3244571	4932326	4.27%	1.713%
15	11.832	3243	3263	3271	rBV3	533254	1423822	1.23%	0.494%
16	11.890	3271	3281	3293	rVB	1399156	2248058	1.95%	0.781%
17	12.092	3327	3344	3361	rVV2	2049036	3364970	2.91%	1.169%
18	12.205	3361	3379	3396	rVB5	900065	2330270	2.02%	0.809%
19	14.079	3945	3962	3986	rVB	1094156	1826329	1.58%	0.634%

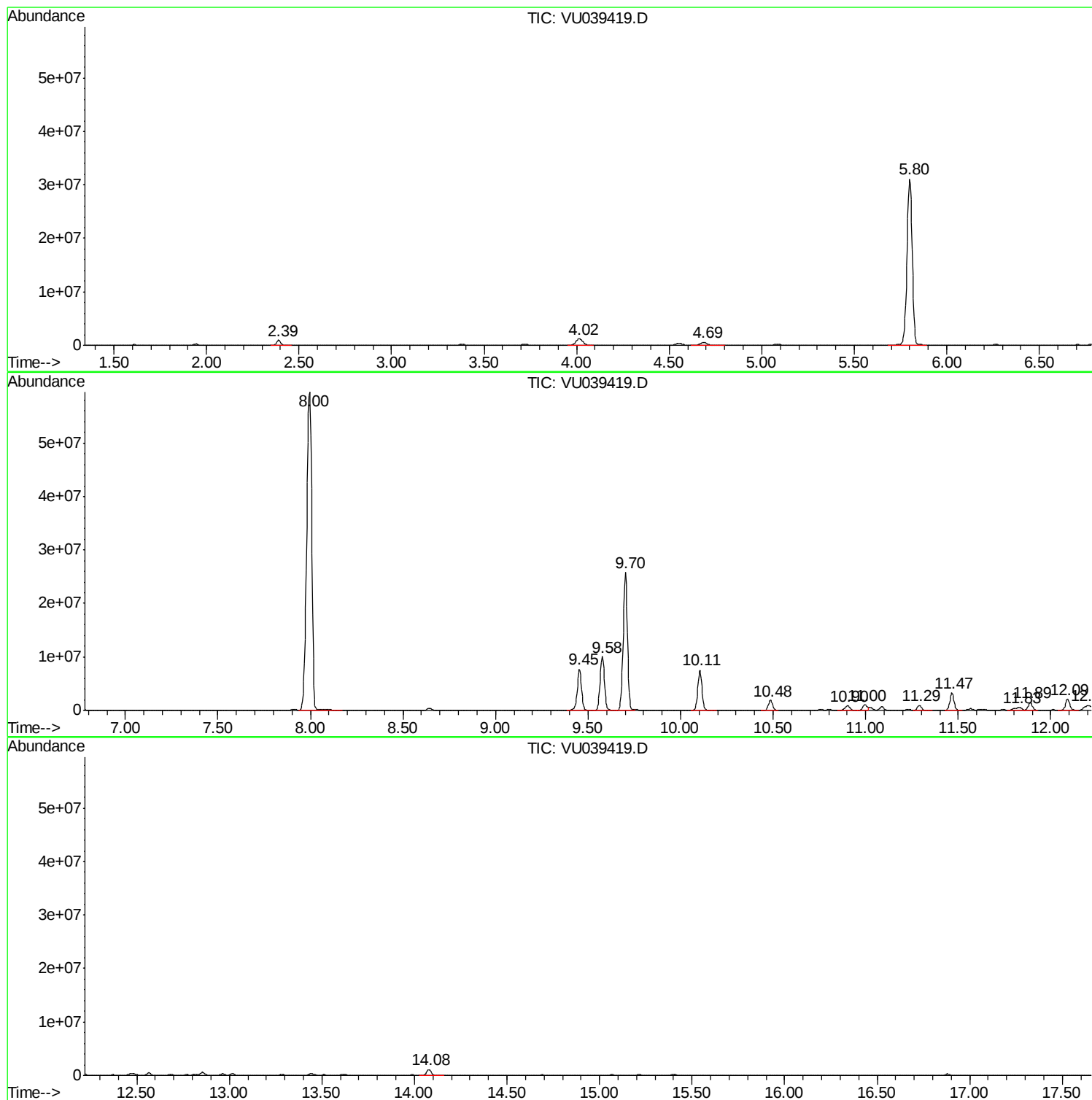
Sum of corrected areas: 287936848

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

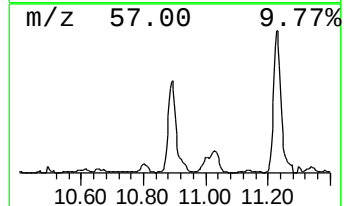
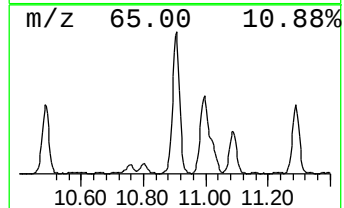
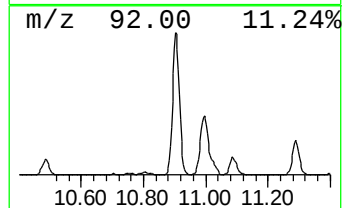
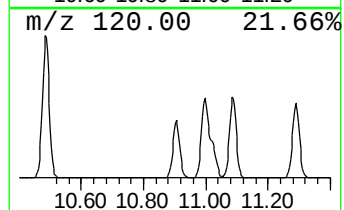
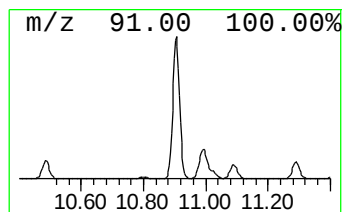
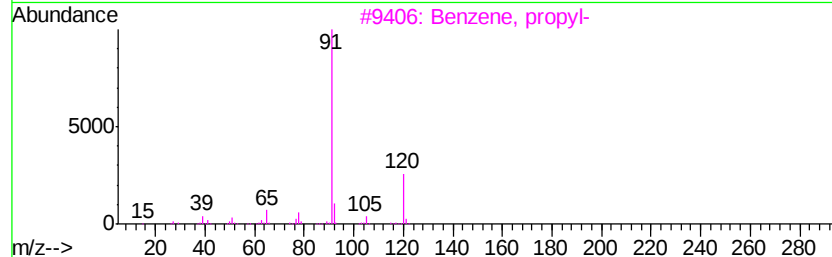
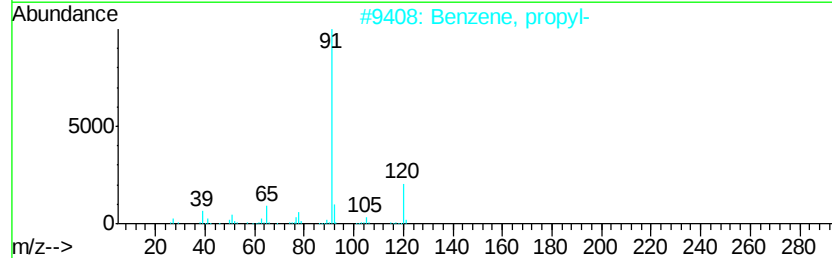
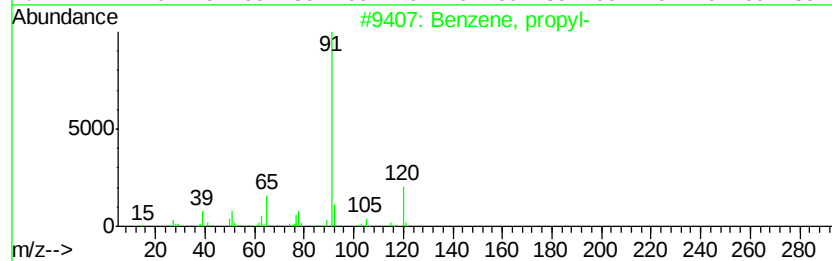
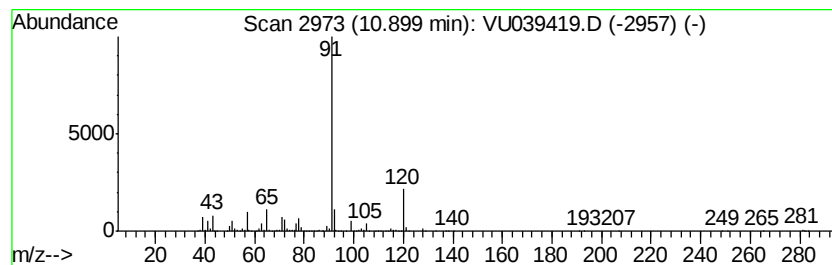
Instrument :
MSVOA_U
ClientSampleId :
BE506

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.90	5.71 ug/L	1626980	1,4-Dichlorobenzene-d4	11.81	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	87
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	80
4	1-Benzylamino-2-benzvloxvethane	241	C16H19NO	038336-06-0	59
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

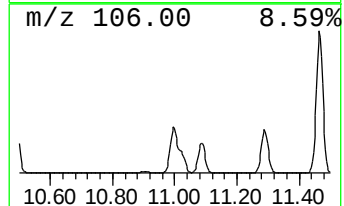
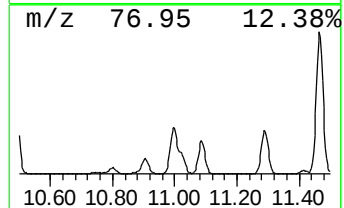
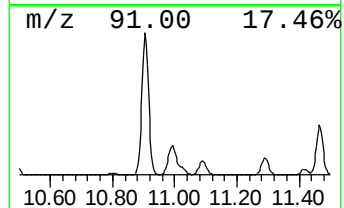
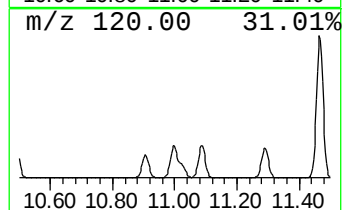
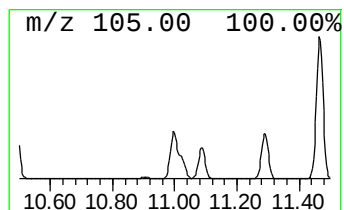
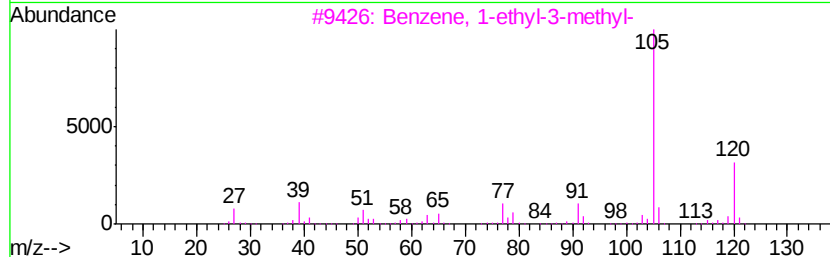
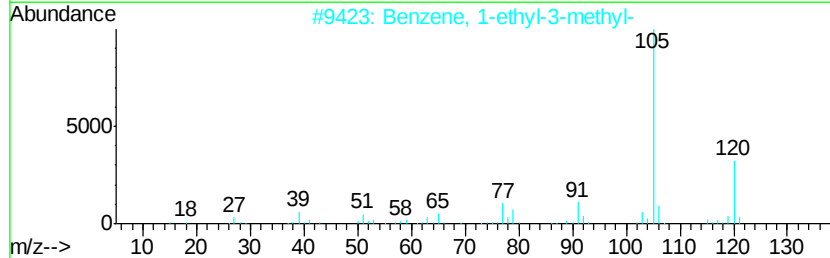
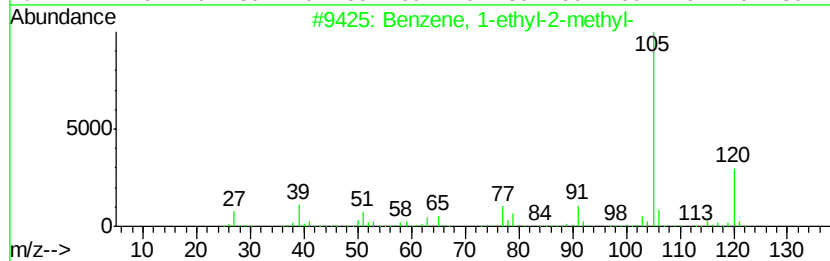
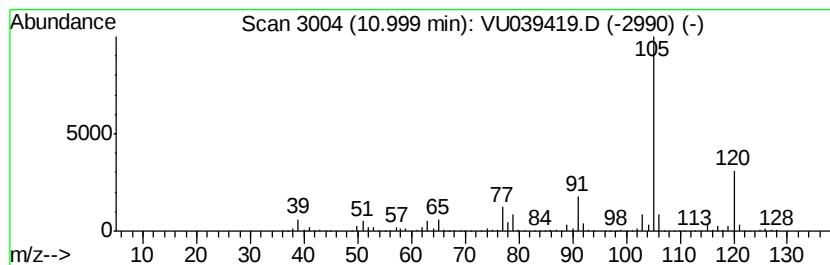
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	6.26 ug/L	1782190	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	1-ethyl-2-methyl-	120	C9H12	000611-14-3	90	
2	Benzene,	1-ethyl-3-methyl-	120	C9H12	000620-14-4	90	
3	Benzene,	1-ethyl-3-methyl-	120	C9H12	000620-14-4	90	
4	Benzene,	1-ethyl-2-methyl-	120	C9H12	000611-14-3	90	
5	Benzene,	1-ethyl-4-methyl-	120	C9H12	000622-96-8	90	



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

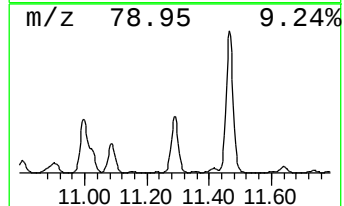
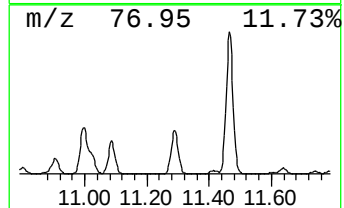
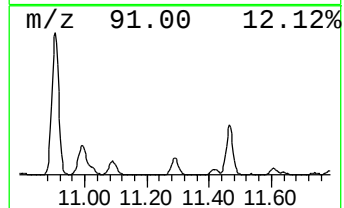
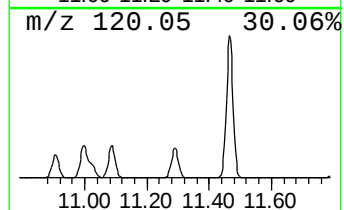
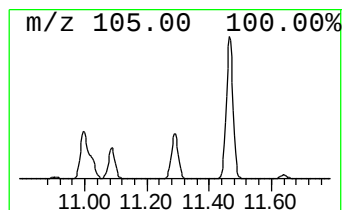
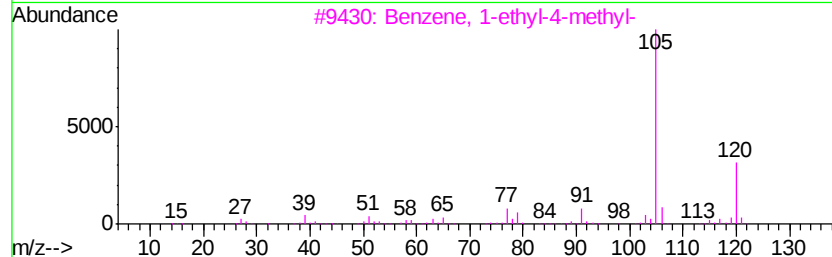
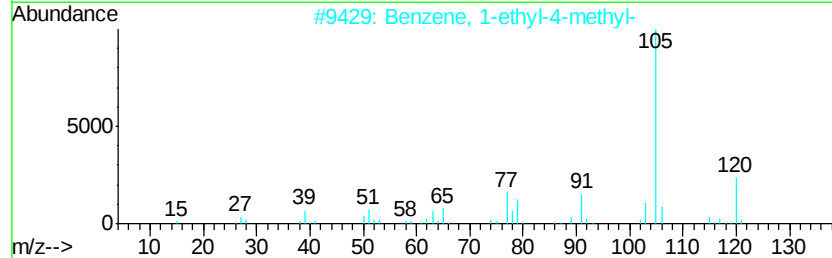
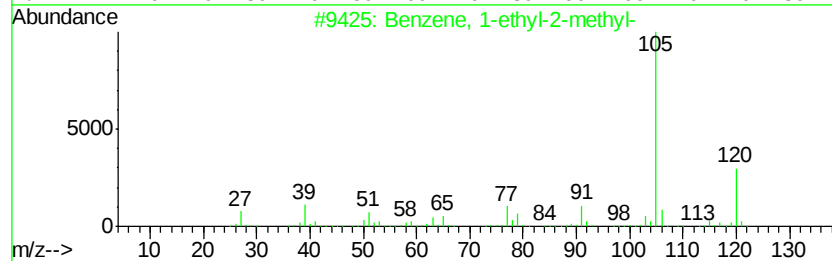
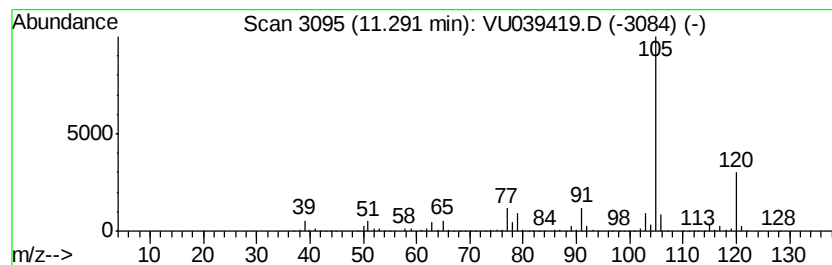
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	5.18 ug/L	1474100	1,4-Dichlorobenzene-d4	11.81

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

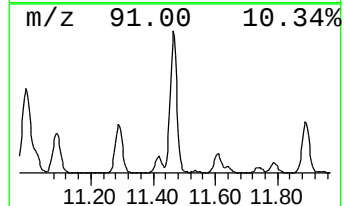
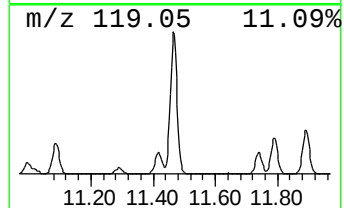
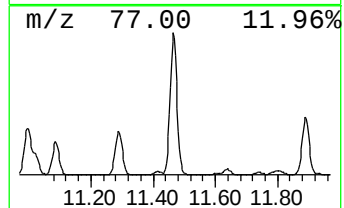
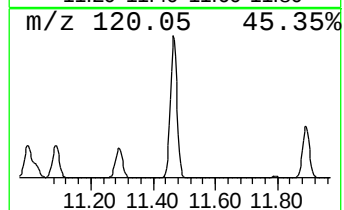
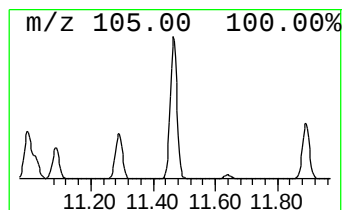
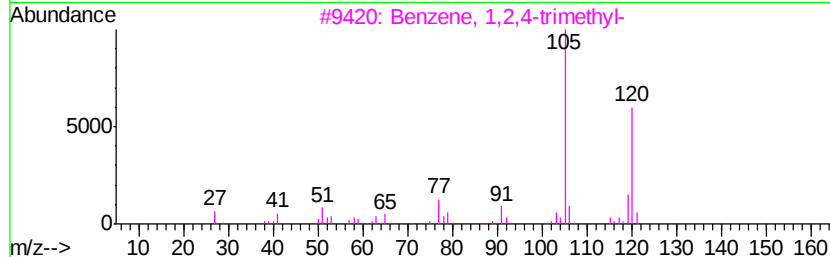
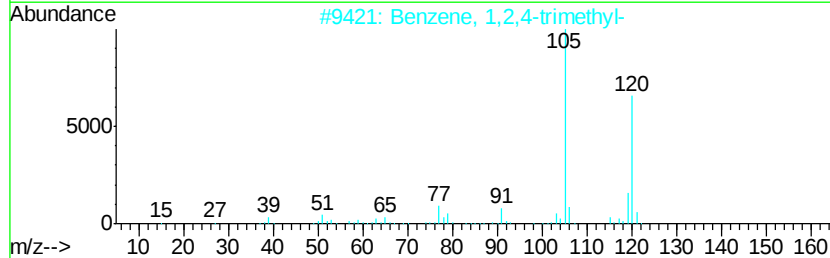
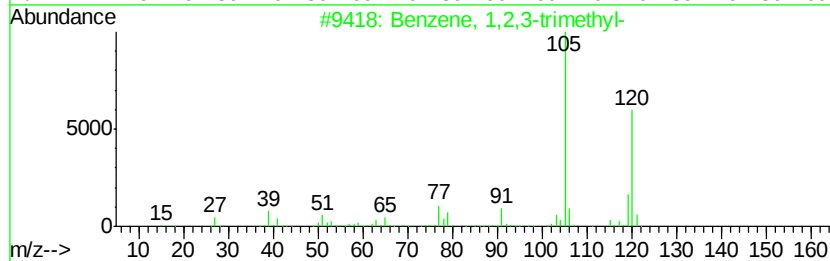
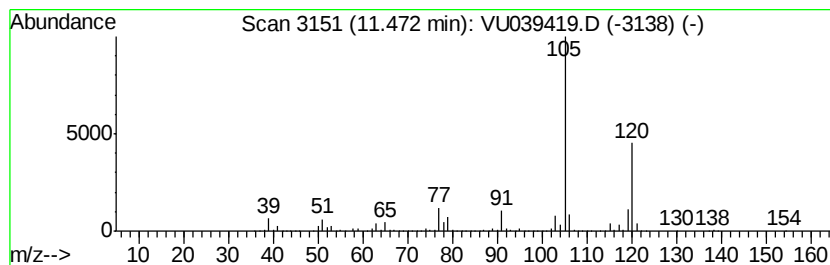
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	17.32 ug/L	4932330	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5		Mesitylene	120	C9H12	000108-67-8	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

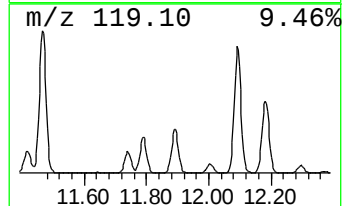
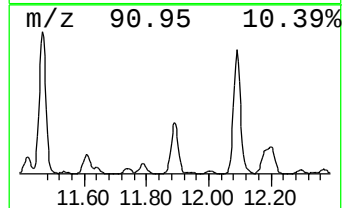
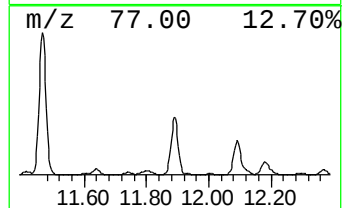
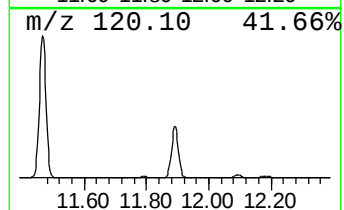
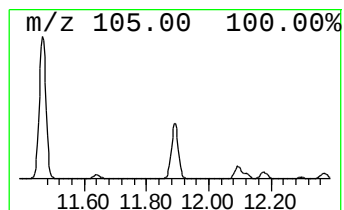
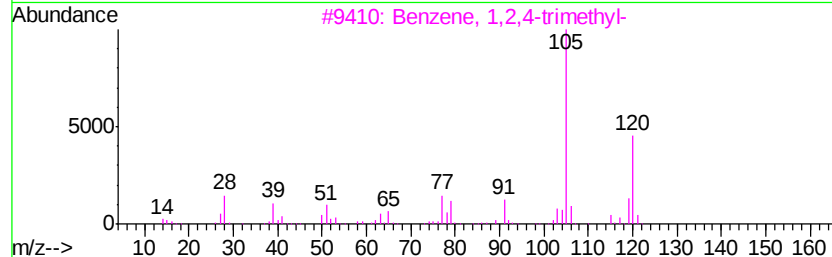
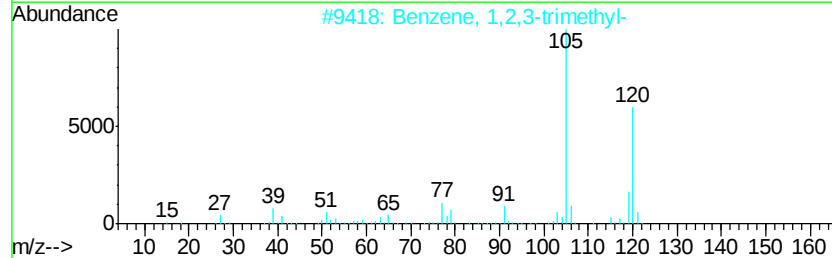
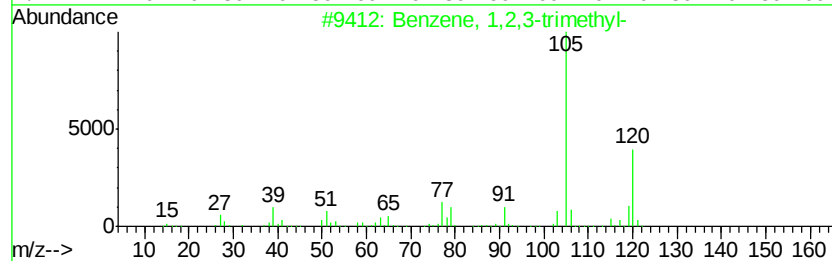
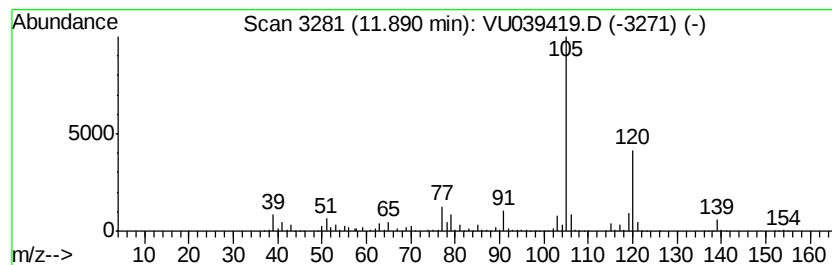
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 Benzene, 1,2,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	7.89 ug/L	2248060	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
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		Mesitylene	120	C9H12	000108-67-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

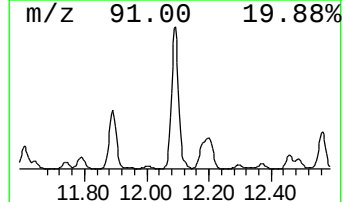
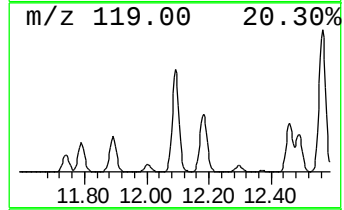
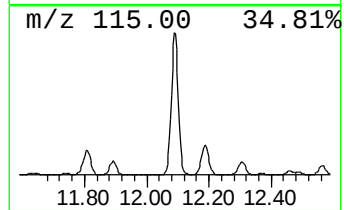
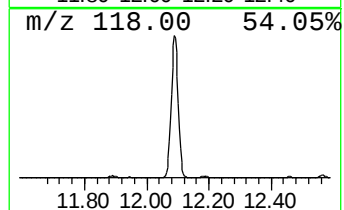
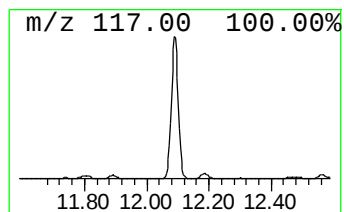
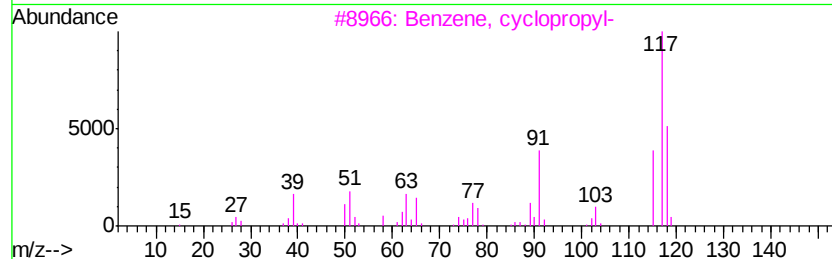
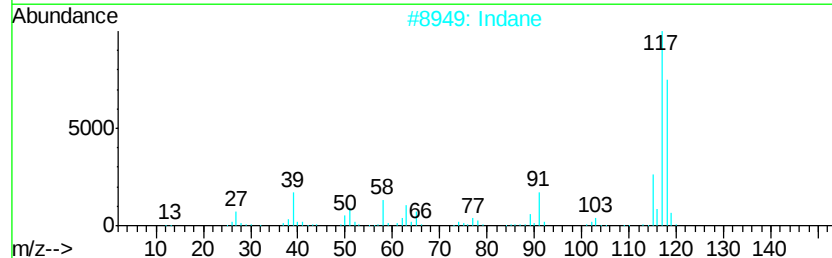
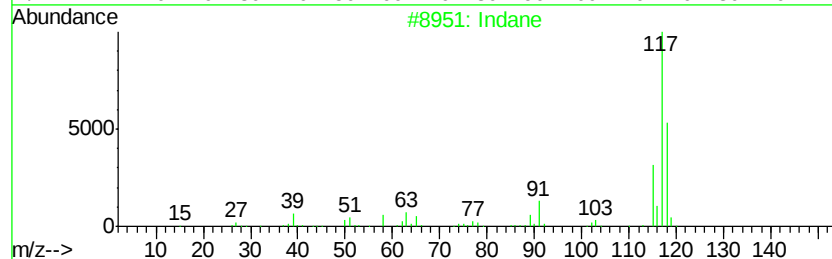
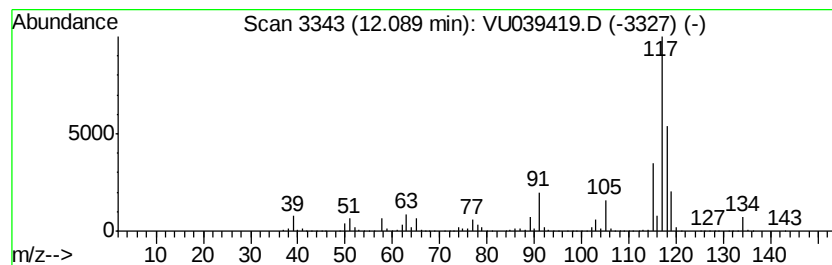
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	11.82 ug/L	3364970	1,4-Dichlorobenzene-d4	11.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	70
2	Indane		118	C9H10	000496-11-7	70
3	Benzene, cvclopopyl-		118	C9H10	000873-49-4	70
4	Benzene, cvclopopyl-		118	C9H10	000873-49-4	70
5	Benzene, 2-propenyl-		118	C9H10	000300-57-2	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU070920\
Data File : VU039419.D
Acq On : 10 Jul 2020 01:29
Operator : SY/MD
Sample : L3261-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE506

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR070720WMA.M
Quant Title : TRACE VOA SOM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, propyl-	10.90	5.7	ug/L	1626980	3	11.81	1423820	5.0
Benzene, 1-ethyl-...	11.00	6.3	ug/L	1782190	3	11.81	1423820	5.0
Benzene, 1-ethyl-...	11.29	5.2	ug/L	1474100	3	11.81	1423820	5.0
Benzene, 1,2,3-tr...	11.47	17.3	ug/L	4932330	3	11.81	1423820	5.0
Benzene, 1,2,4-tr...	11.89	7.9	ug/L	2248060	3	11.81	1423820	5.0
Indane	12.09	11.8	ug/L	3364970	3	11.81	1423820	5.0