

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081420\
 Data File : VX017897.D
 Acq On : 14 Aug 2020 14:04
 Operator : JC/SP
 Sample : L3697-02
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG0M9

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_X\METHOD\SOMXTR080720WMA.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	2.166	172	177	186	rBV	2436451	3485651	2.41%	0.712%
2	3.825	437	449	469	rBV	3349079	8897302	6.16%	1.817%
3	4.373	528	539	554	rBV2	902404	2726649	1.89%	0.557%
4	4.562	555	570	582	rBV2	1334027	3771569	2.61%	0.770%
5	6.129	803	827	841	rBV2	39177412	113638844	78.67%	23.204%
6	8.787	1253	1263	1268	rBV3	71479897	144447886	100.00%	29.495%
7	10.122	1473	1482	1490	rBV	19493892	27086834	18.75%	5.531%
8	10.238	1495	1501	1506	rBV	24295859	31808476	22.02%	6.495%
9	10.348	1512	1519	1527	rBV	51474009	74290740	51.43%	15.169%
10	10.683	1569	1574	1586	rVB	18819306	24215333	16.76%	4.944%
11	11.000	1621	1626	1633	rBV	4742021	6149233	4.26%	1.256%
12	11.342	1675	1682	1689	rBV	2491986	3346597	2.32%	0.683%
13	11.421	1689	1695	1697	rBV2	2638312	3981576	2.76%	0.813%
14	11.494	1703	1707	1712	rVB	1958946	2406820	1.67%	0.491%
15	11.652	1729	1733	1742	rVB	2654080	3232867	2.24%	0.660%
16	11.793	1751	1756	1763	rVB	8556109	10323515	7.15%	2.108%
17	12.079	1795	1803	1807	rBV3	1352201	2730164	1.89%	0.557%
18	12.128	1807	1811	1816	rVB	3774829	4595148	3.18%	0.938%
19	12.286	1831	1837	1844	rBV	5001796	6711856	4.65%	1.370%
20	12.372	1844	1851	1859	rVB4	2018418	4104416	2.84%	0.838%
21	12.585	1879	1886	1893	rBV3	782885	1903535	1.32%	0.389%
22	12.872	1930	1933	1940	rVB2	1973933	2527926	1.75%	0.516%
23	13.817	2080	2088	2096	rBV	2644137	3360601	2.33%	0.686%

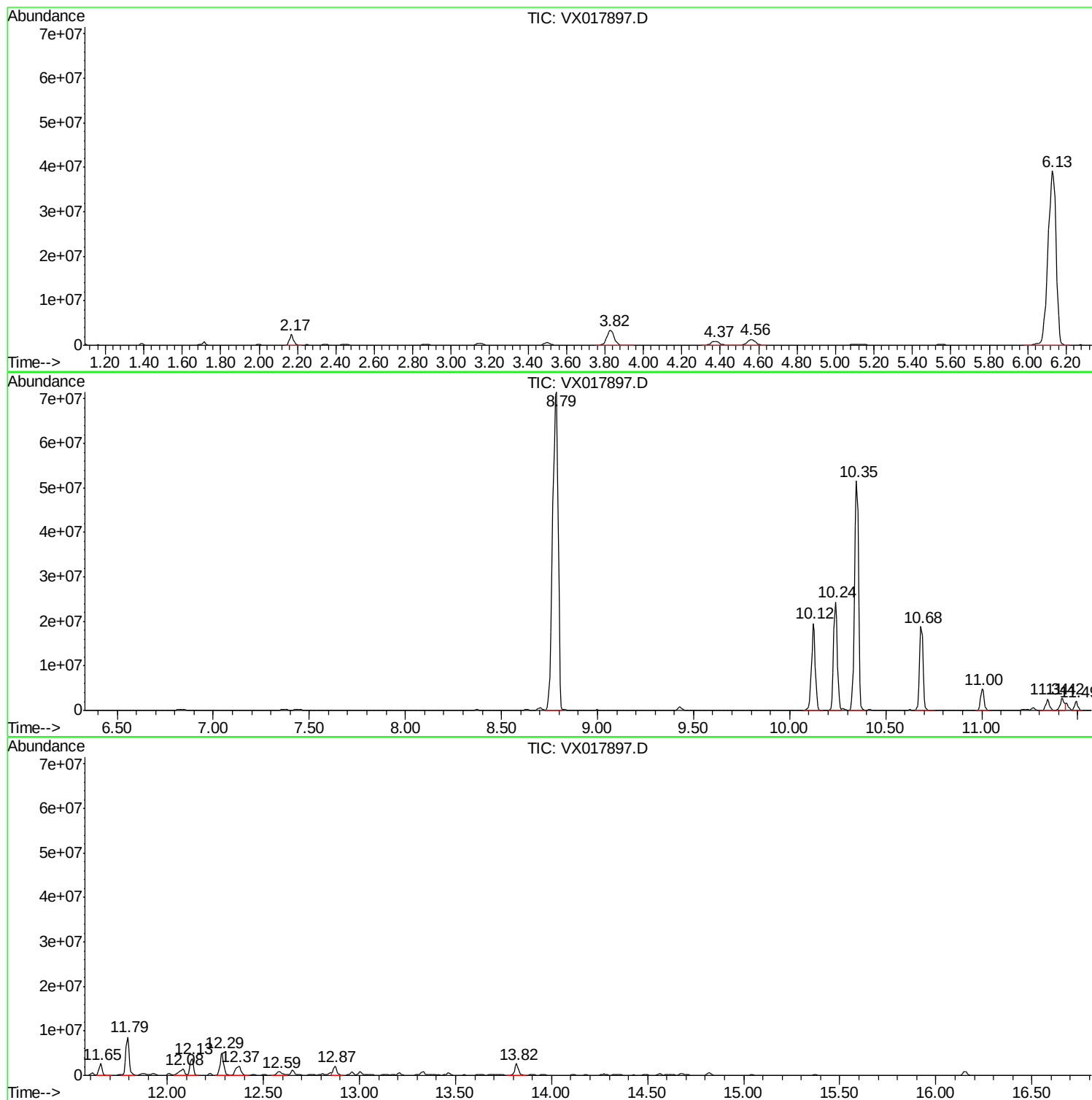
Sum of corrected areas: 489743538

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG0M9

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

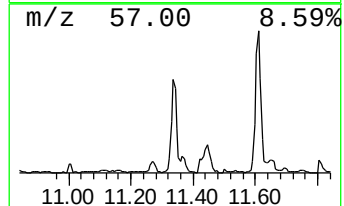
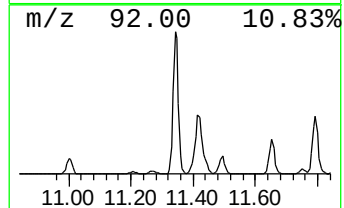
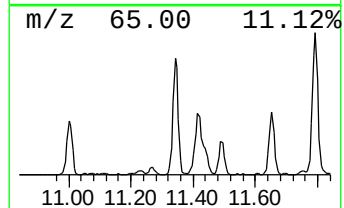
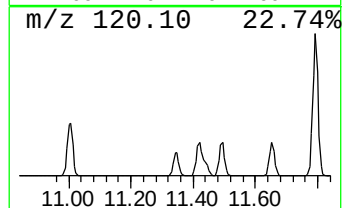
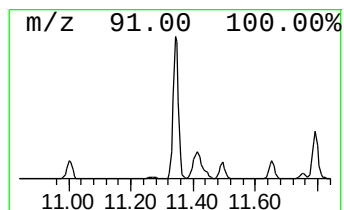
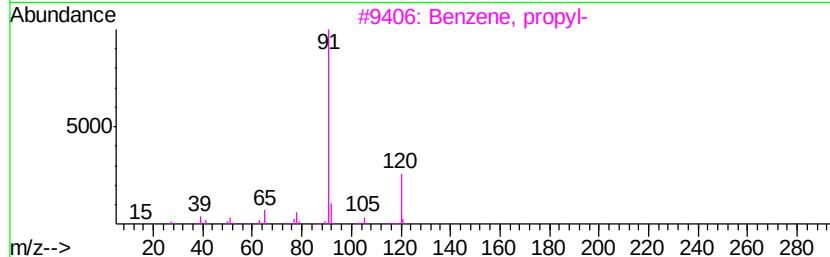
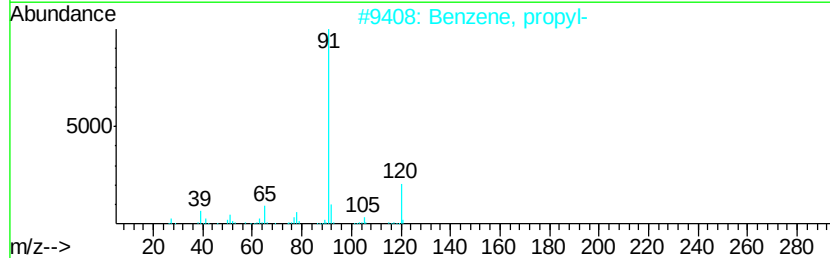
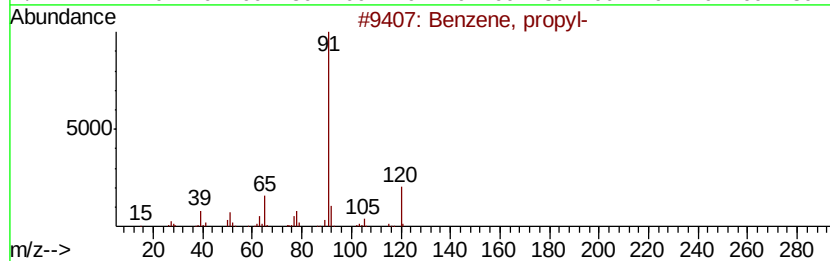
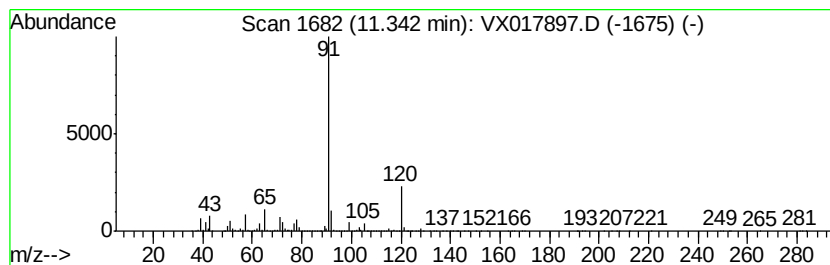
Instrument :
MSVOA_X
ClientSampled :
BG0M9

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.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 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.34	6.13 ug/L	3346600	1,4-Dichlorobenzene-d4	12.06	
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	81
3	Benzene, propyl-	120	C9H12	000103-65-1	74
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
5	1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG0M9

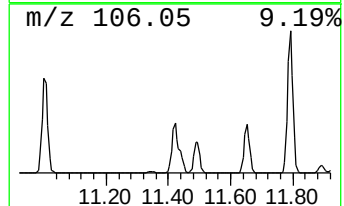
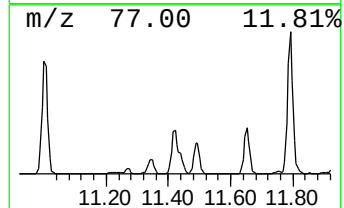
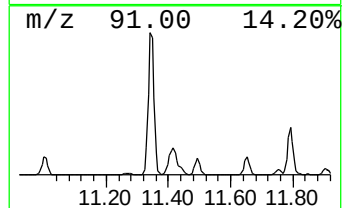
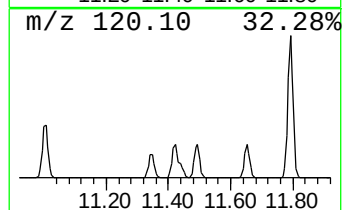
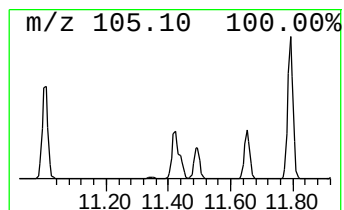
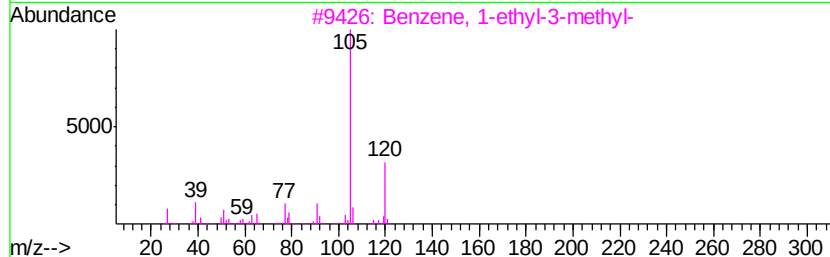
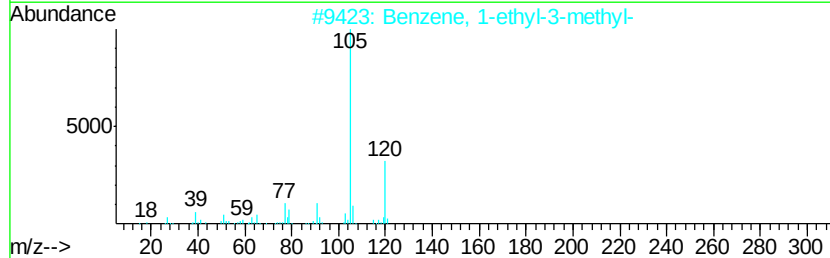
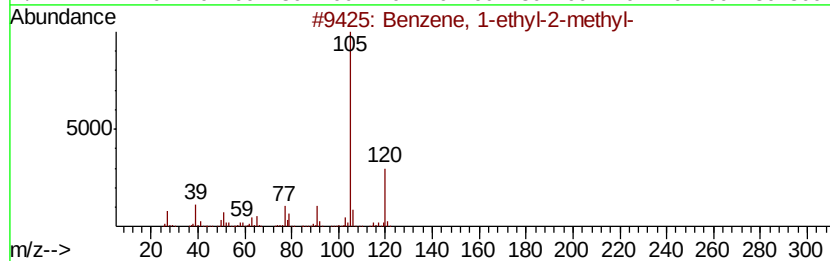
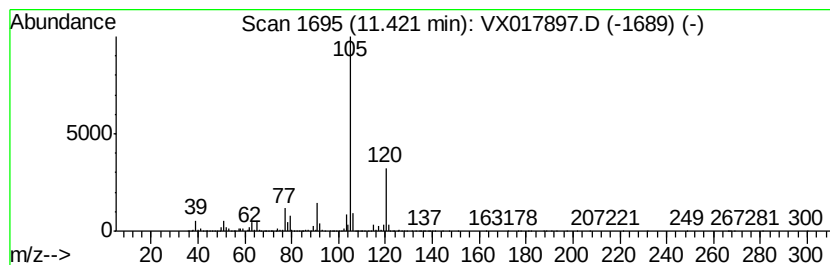
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	7.29 ug/L	3981580	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

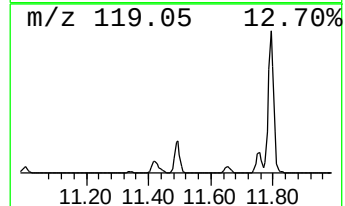
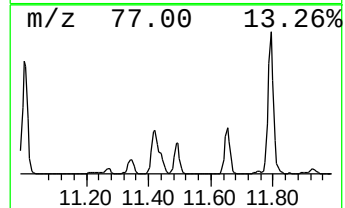
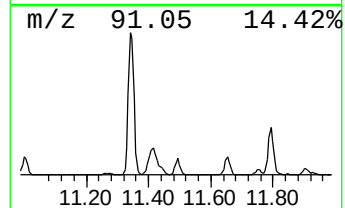
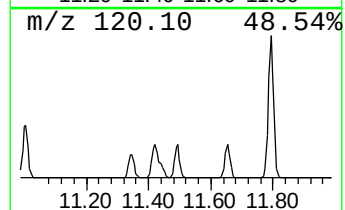
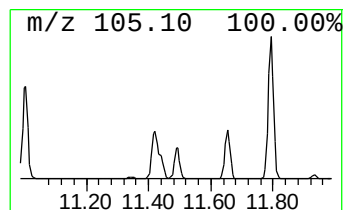
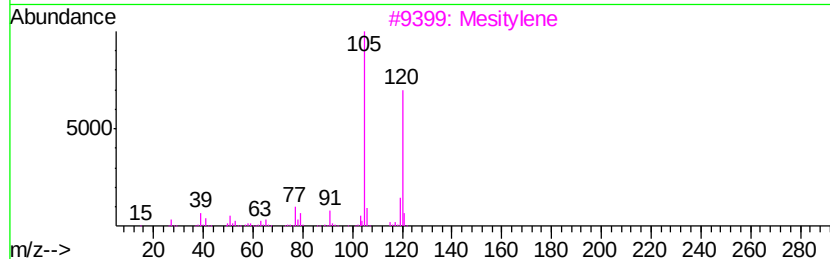
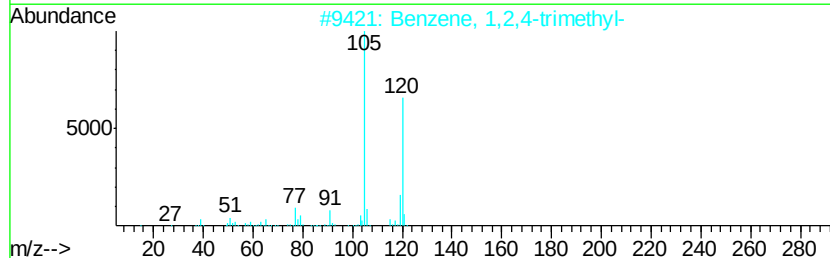
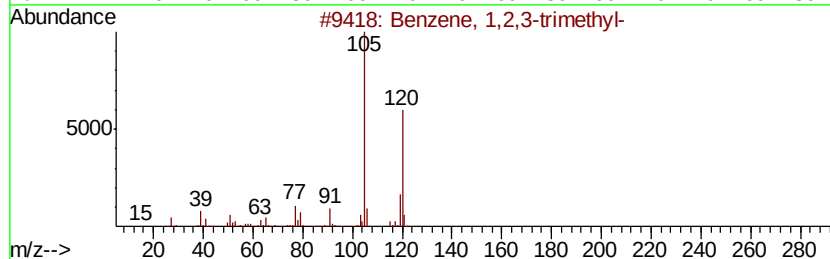
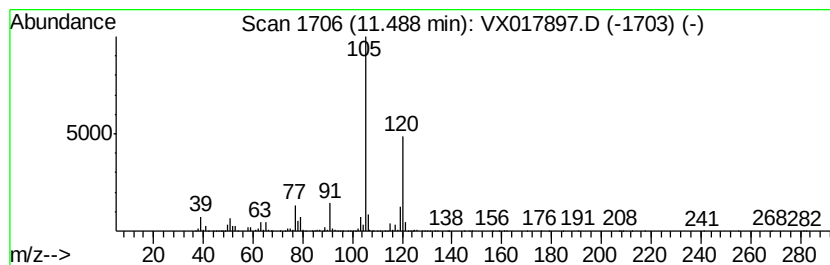
Instrument :
MSVOA_X
ClientSampleId :
BG0M9

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	4.41 ug/L	2406820	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 94
2		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 94
3		Mesitylene	120 C9H12	000108-67-8 94
4		Mesitylene	120 C9H12	000108-67-8 94
5		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG0M9

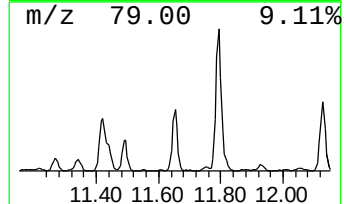
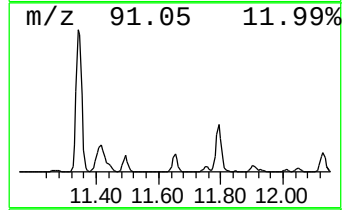
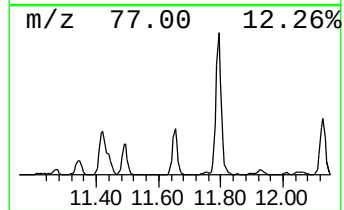
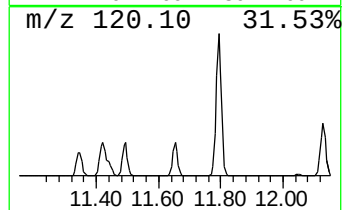
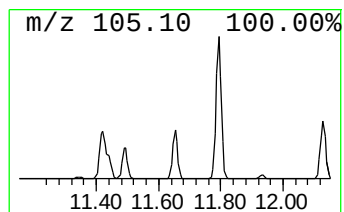
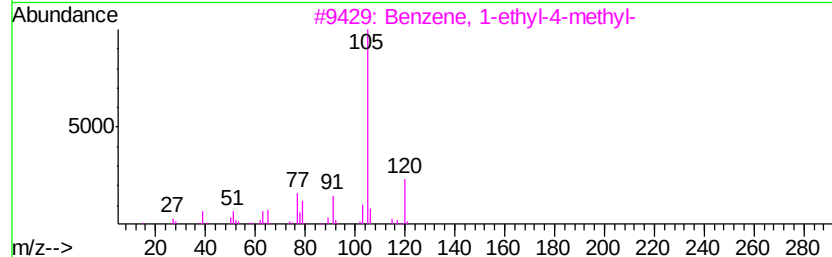
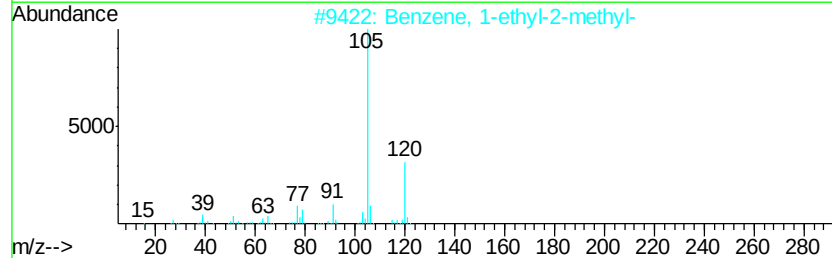
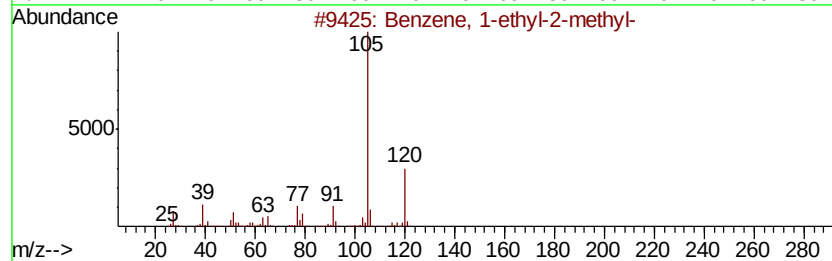
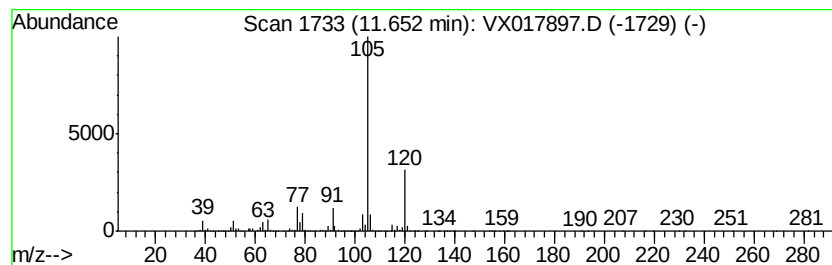
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	5.92 ug/L	3232870	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG0M9

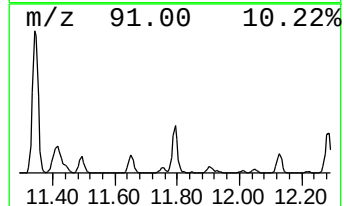
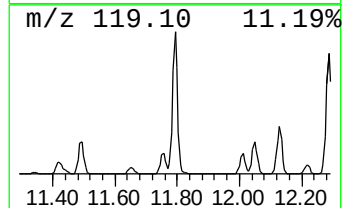
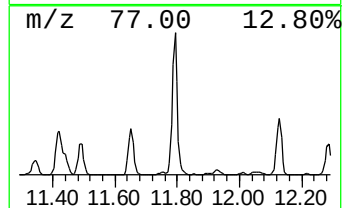
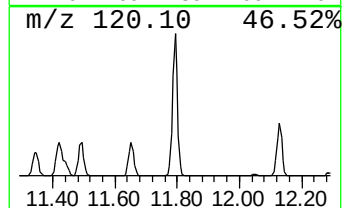
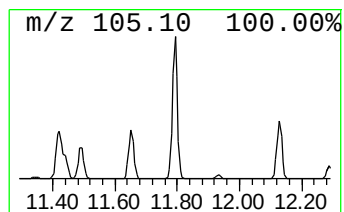
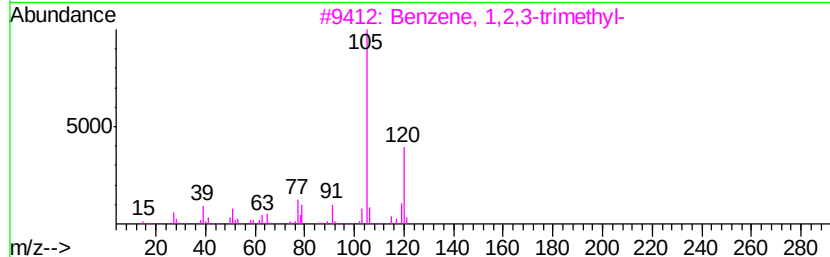
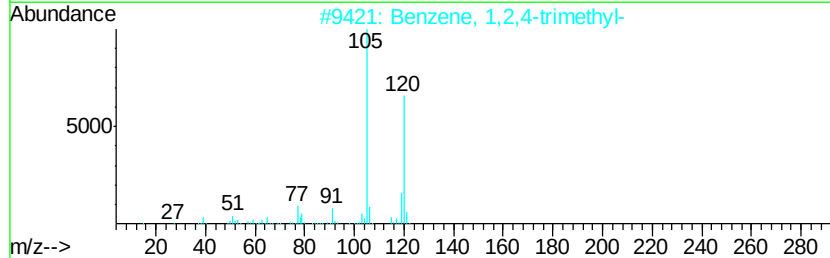
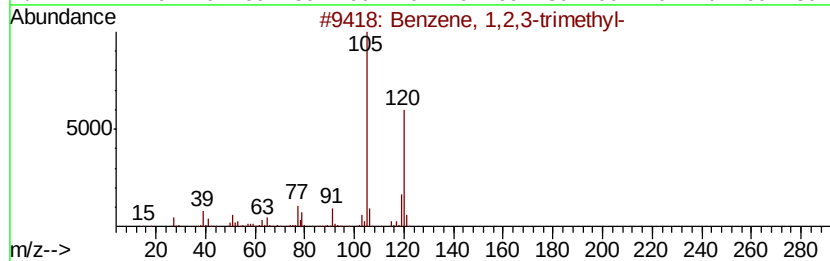
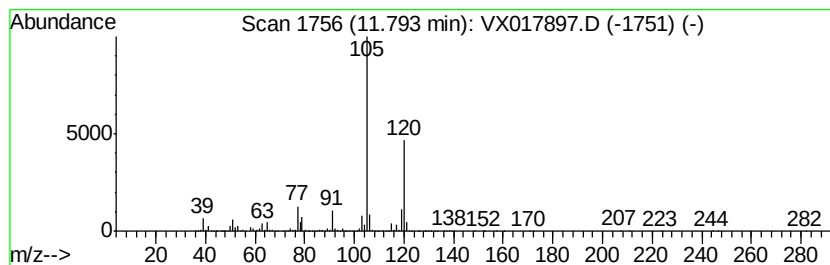
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	18.91 ug/L	10323500	1,4-Dichlorobenzene-d4	12.06

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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

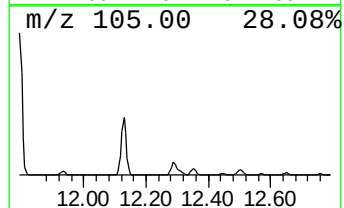
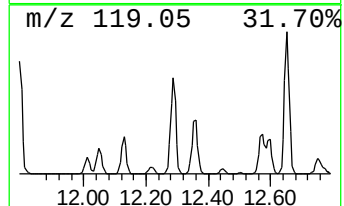
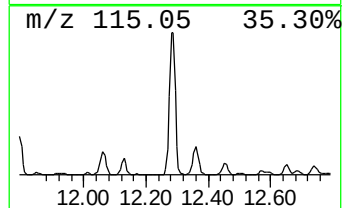
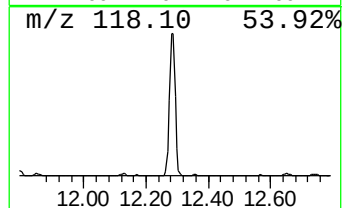
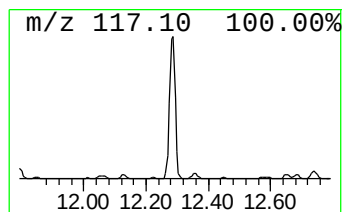
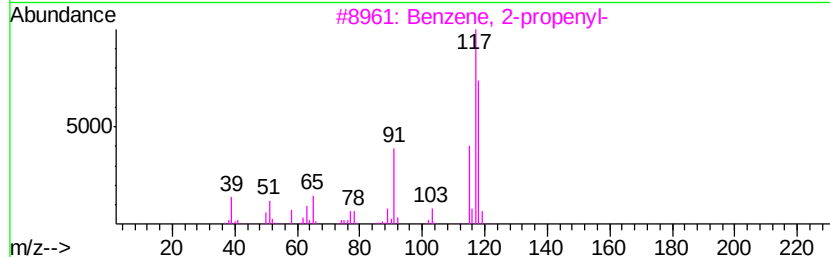
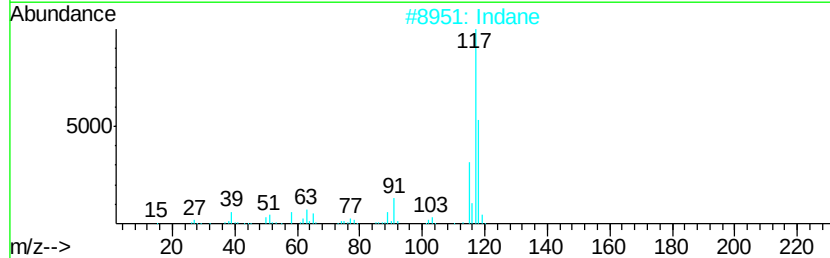
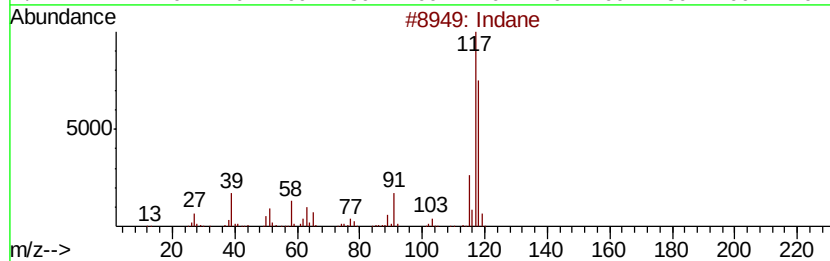
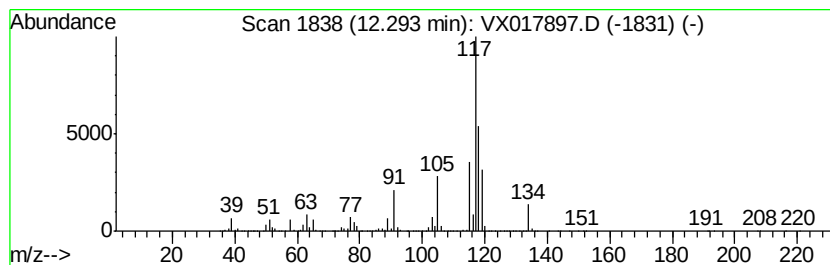
Instrument :
MSVOA_X
ClientSampleId :
BG0M9

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.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.
12.29	12.29 ug/L	6711860	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 70
2	Indane		118 C9H10	000496-11-7 64
3	Benzene, 2-propenyl-		118 C9H10	000300-57-2 64
4	Benzene, cyclopropyl-		118 C9H10	000873-49-4 64
5	Benzene, cyclopropyl-		118 C9H10	000873-49-4 58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG0M9

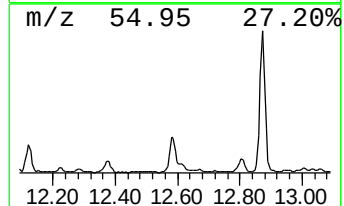
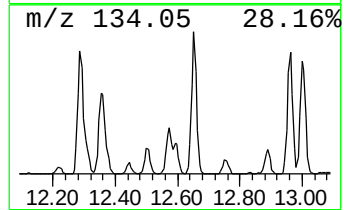
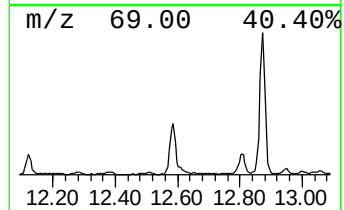
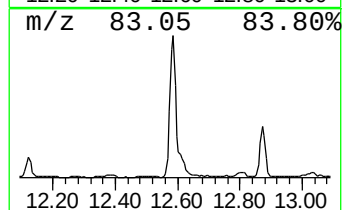
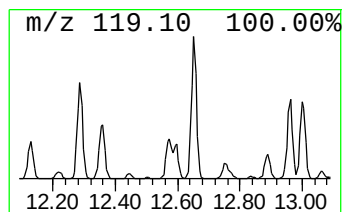
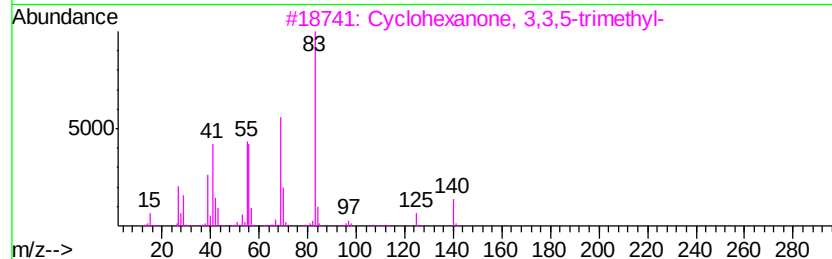
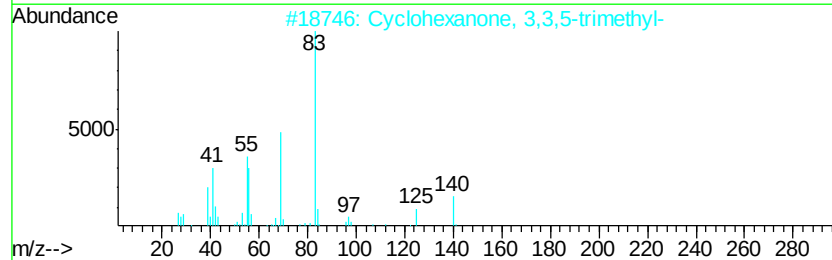
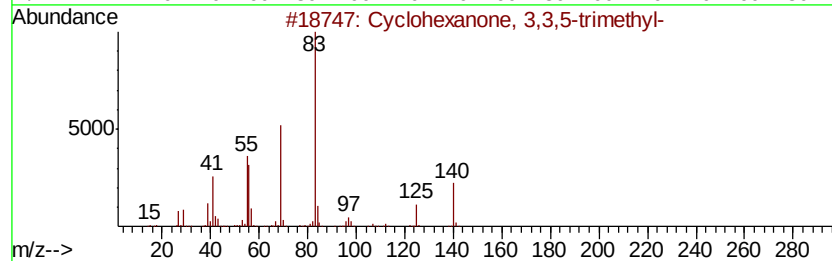
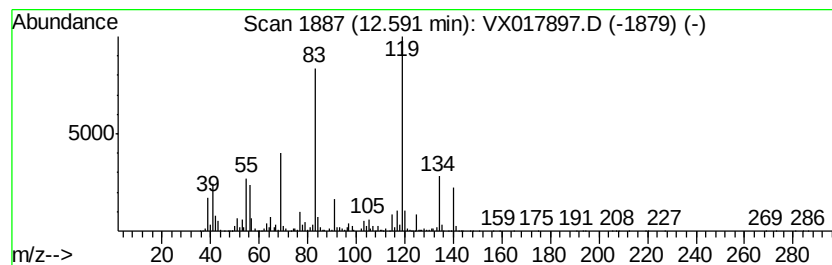
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 Cyclohexanone, 3,3,5-trimet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	3.49 ug/L	1903540	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
2		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
4		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	68
5		p-Cymene	134	C10H14	000099-87-6	68



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VO081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

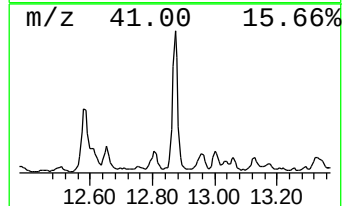
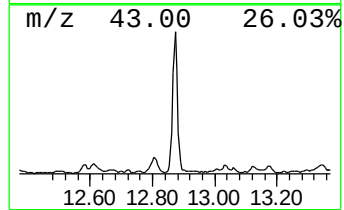
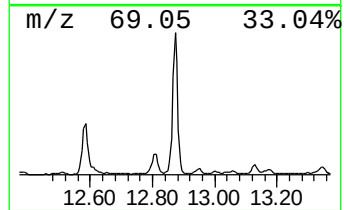
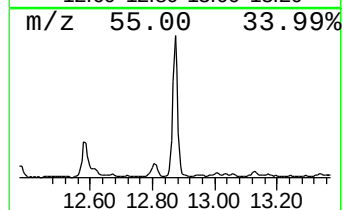
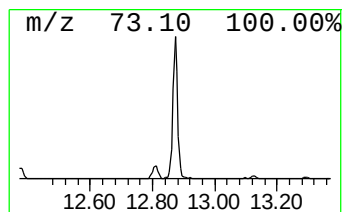
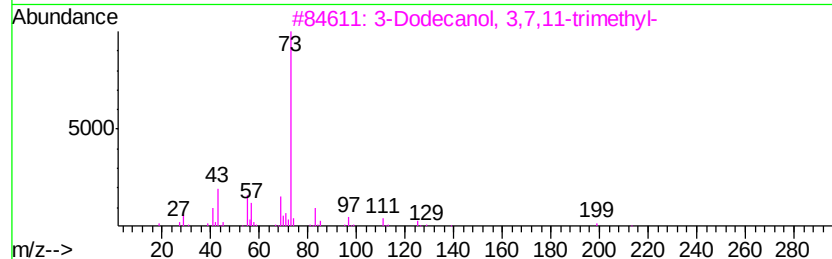
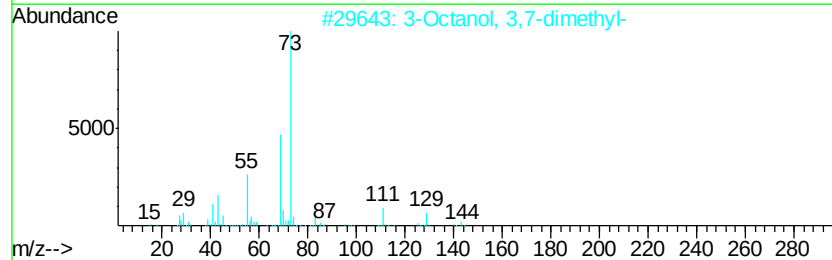
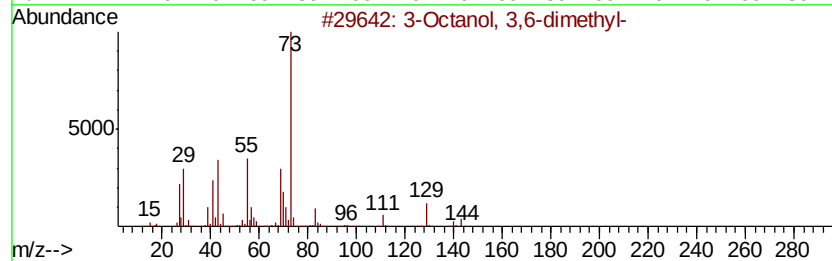
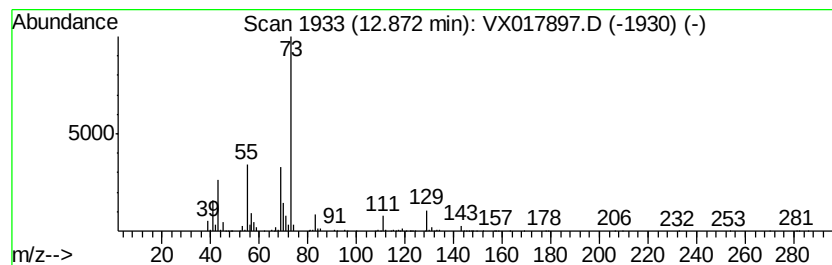
Instrument :
MSVOA_X
ClientSampleId :
BG0M9

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 3-Octanol, 3,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	4.63 ug/L	2527930	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	3-Octanol, 3,6-dimethyl-	158 C10H22O	000151-19-9	86
2	3-Octanol, 3,7-dimethyl-	158 C10H22O	000078-69-3	59
3	3-Dodecanol, 3,7,11-trimethyl-	228 C15H32O	007278-65-1	37
4	3-Octanol, 3,7-dimethyl-	158 C10H22O	000078-69-3	28
5	2,5-Dimethyl-4-hydroxy-3-hexanone	144 C8H16O2	000815-77-0	27



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX081420\
Data File : VX017897.D
Acq On : 14 Aug 2020 14:04
Operator : JC/SP
Sample : L3697-02
Misc : 25.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG0M9

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR080720WMA.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-	11.34	6.1	ug/L	3346600	3	12.06	2730160	5.0
Benzene, 1-ethyl-...	11.42	7.3	ug/L	3981580	3	12.06	2730160	5.0
Benzene, 1,2,3-tr...	11.49	4.4	ug/L	2406820	3	12.06	2730160	5.0
Benzene, 1-ethyl-...	11.65	5.9	ug/L	3232870	3	12.06	2730160	5.0
Benzene, 1,2,4-tr...	11.79	18.9	ug/L	10323500	3	12.06	2730160	5.0
Indane	12.29	12.3	ug/L	6711860	3	12.06	2730160	5.0
Cyclohexanone, 3,...	12.59	3.5	ug/L	1903540	3	12.06	2730160	5.0
3-Octanol, 3,6-di...	12.87	4.6	ug/L	2527930	3	12.06	2730160	5.0