

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092620\
 Data File : VU040316.D
 Acq On : 25 Sep 2020 21:43
 Operator : SY/MD
 Sample : L4139-04
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG494

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\SOMUTR083120WMA.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	313	326	343	rBV	2255629	3450568	1.79%	0.643%
2	3.716	720	739	764	rVB	881910	1938255	1.00%	0.361%
3	4.015	812	832	852	rBV	3999813	10325927	5.35%	1.925%
4	4.552	980	999	1019	rBV	1343499	3400634	1.76%	0.634%
5	5.809	1351	1390	1411	rBV2	51899156	125137688	64.80%	23.325%
6	8.009	2051	2074	2126	rVB3	72836659	193113752	100.00%	35.995%
7	9.452	2503	2523	2547	rBV	7347144	12563824	6.51%	2.342%
8	9.578	2547	2562	2585	rVV2	22768901	37212719	19.27%	6.936%
9	9.710	2585	2603	2635	rVB2	45093201	81635921	42.27%	15.216%
10	10.105	2710	2726	2755	rVB	17988709	28987131	15.01%	5.403%
11	10.484	2828	2844	2856	rBV	1596427	2494968	1.29%	0.465%
12	10.899	2956	2973	2990	rBV	1362037	2483295	1.29%	0.463%
13	10.996	2990	3003	3009	rBV	3034842	5210782	2.70%	0.971%
14	11.086	3022	3031	3044	rVB2	1732428	2648817	1.37%	0.494%
15	11.288	3084	3094	3116	rVB	1655192	2638205	1.37%	0.492%
16	11.465	3137	3149	3166	rVB	7053429	10672957	5.53%	1.989%
17	11.890	3271	3281	3308	rVB	2253045	3620643	1.87%	0.675%
18	12.089	3330	3343	3361	rBV3	1087214	2030399	1.05%	0.378%
19	12.201	3361	3378	3396	rVB5	978035	2484469	1.29%	0.463%
20	12.475	3445	3463	3482	rBV2	940210	2200900	1.14%	0.410%
21	12.851	3572	3580	3596	rVB2	1428448	2251494	1.17%	0.420%

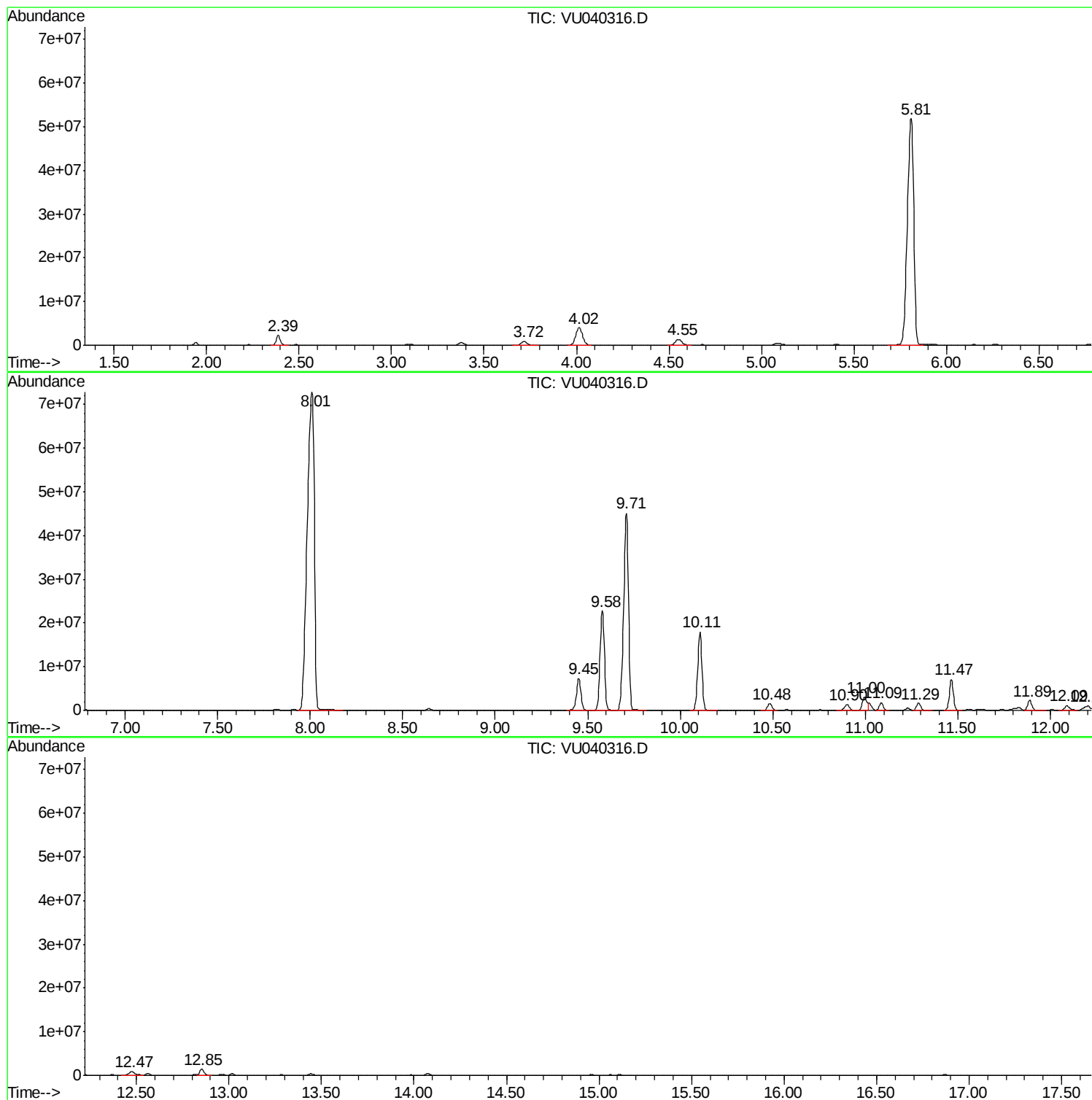
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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



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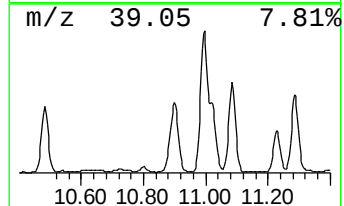
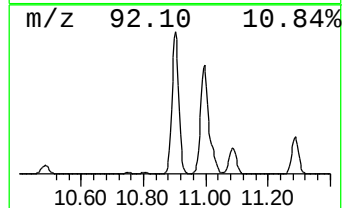
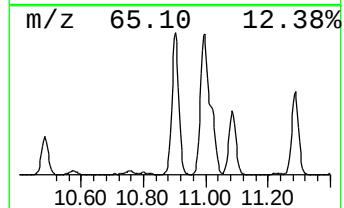
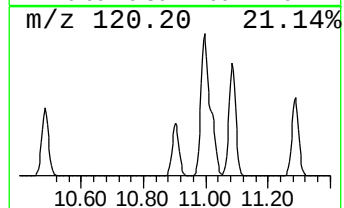
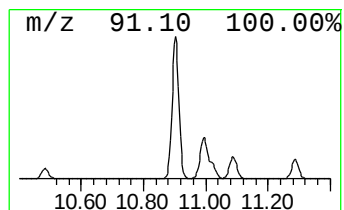
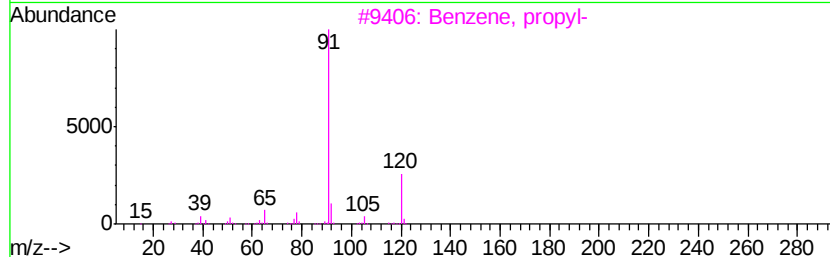
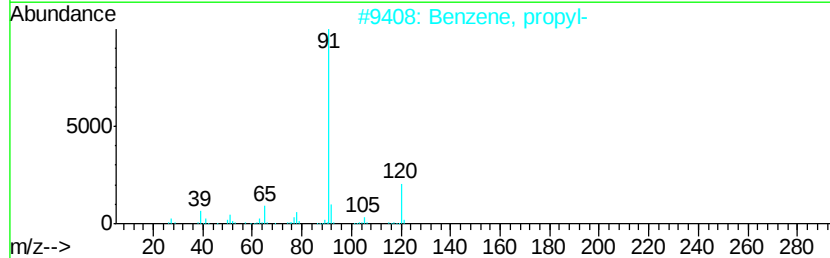
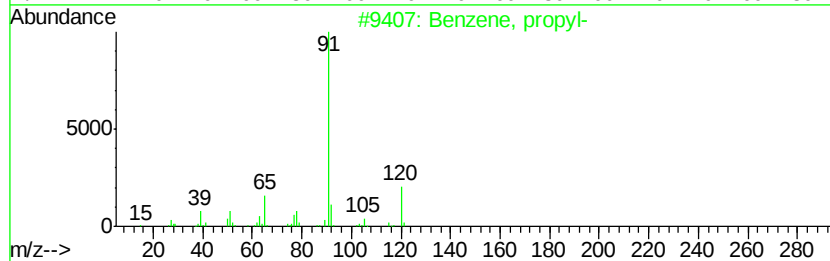
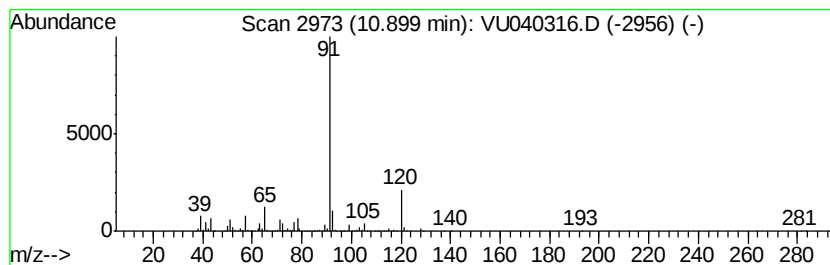
Instrument :
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ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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	3.43 ug/L	2483300	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, propyl-	120 C9H12	000103-65-1 93
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 64
5		1,2-Ethanediamine, N-(phenylmeth...	150 C9H14N2	004152-09-4 59



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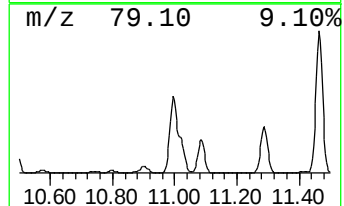
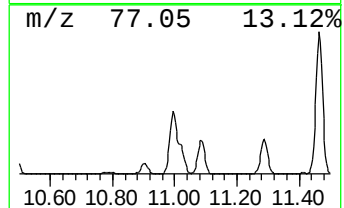
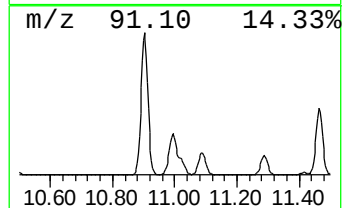
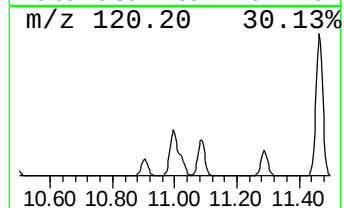
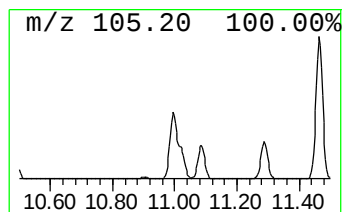
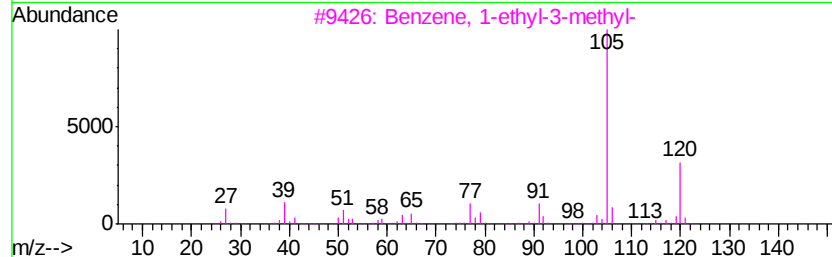
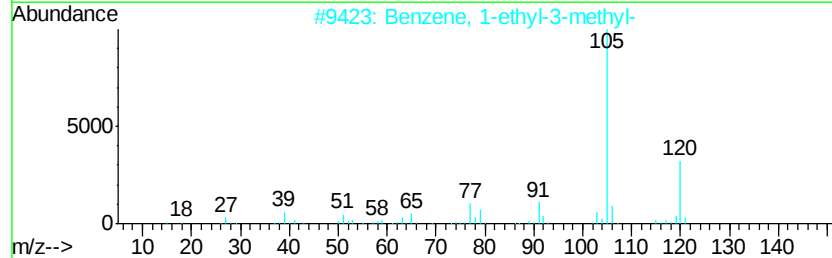
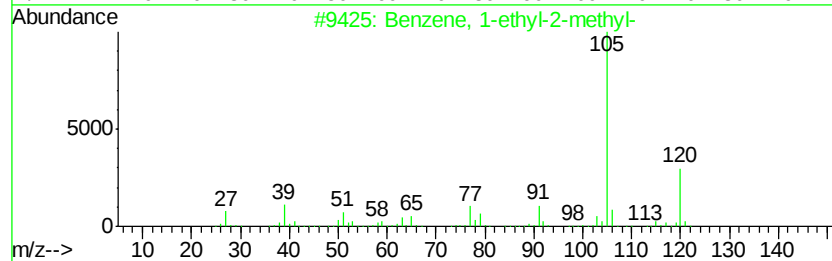
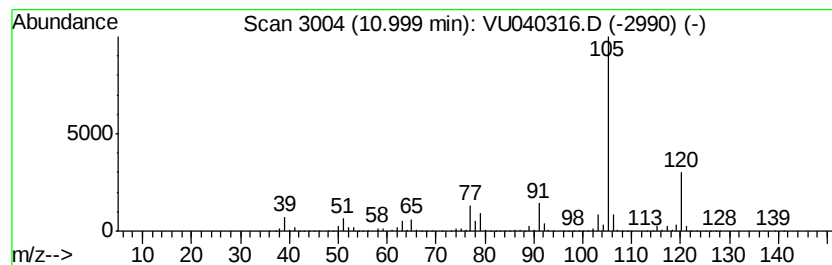
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	7.20 ug/L	5210780	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	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		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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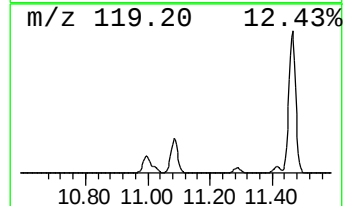
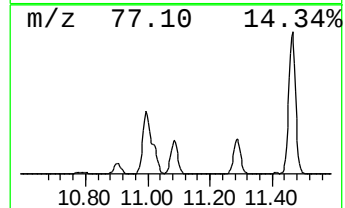
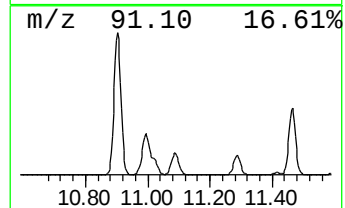
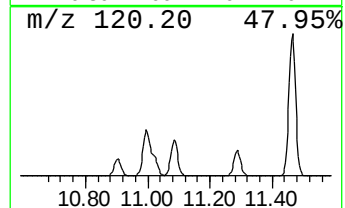
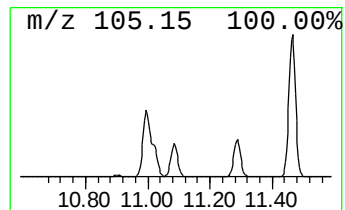
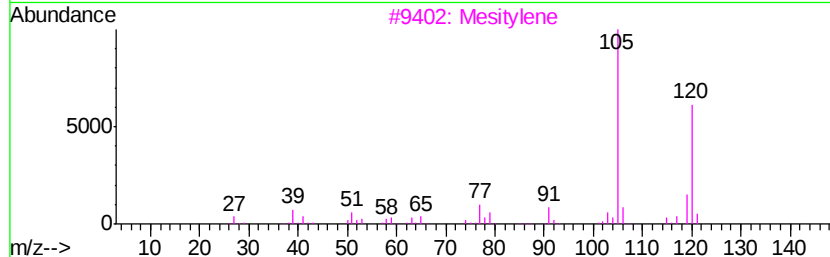
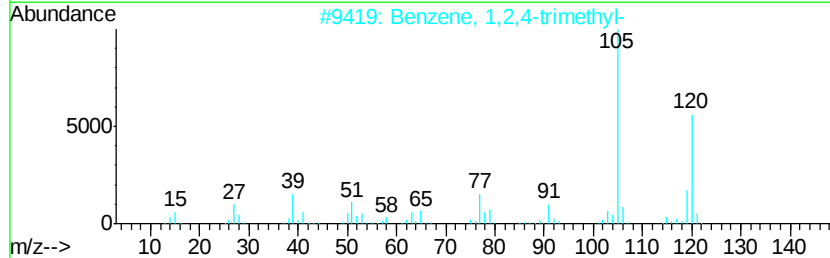
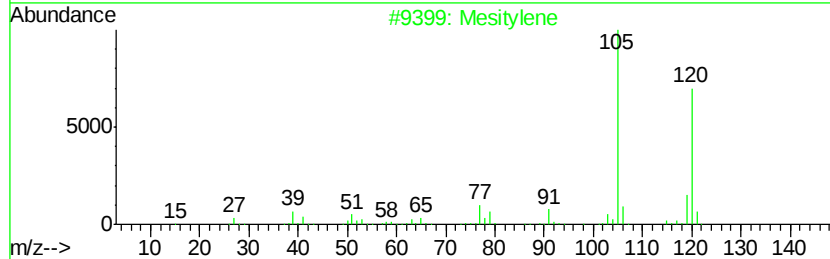
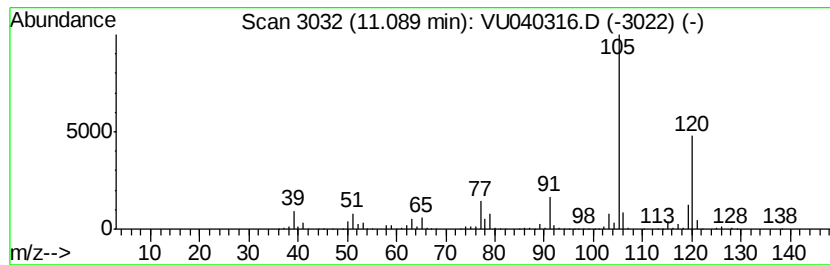
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Mesitylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	3.66 ug/L	2648820	1,4-Dichlorobenzene-d4	11.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene		120	C9H12	000108-67-8	95
2	Benzene, 1,2,4-trimethyl-		120	C9H12	000095-63-6	94
3	Mesitylene		120	C9H12	000108-67-8	94
4	Benzene, 1,2,3-trimethyl-		120	C9H12	000526-73-8	94
5	Mesitylene		120	C9H12	000108-67-8	91



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ALS Vial : 31 Sample Multiplier: 1

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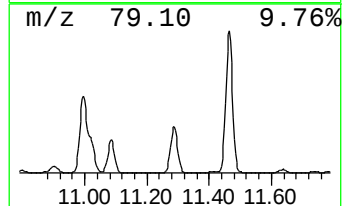
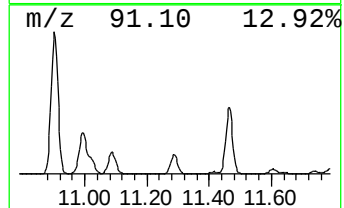
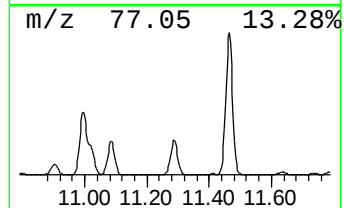
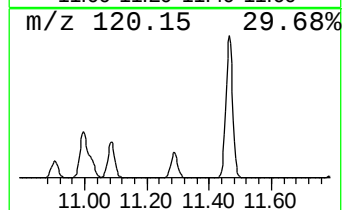
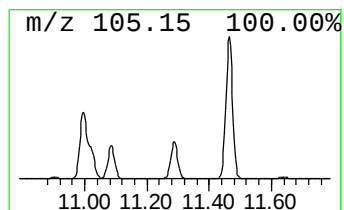
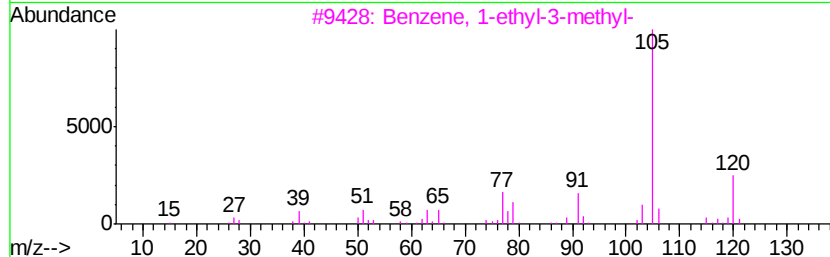
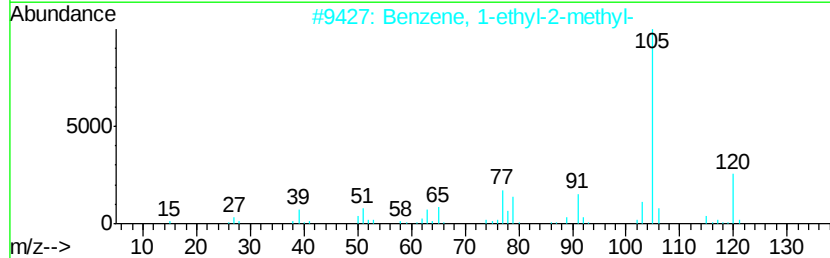
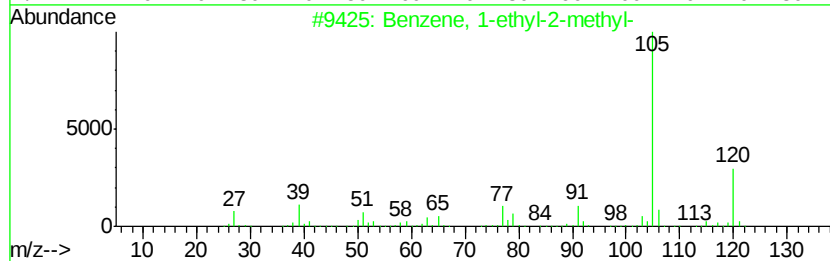
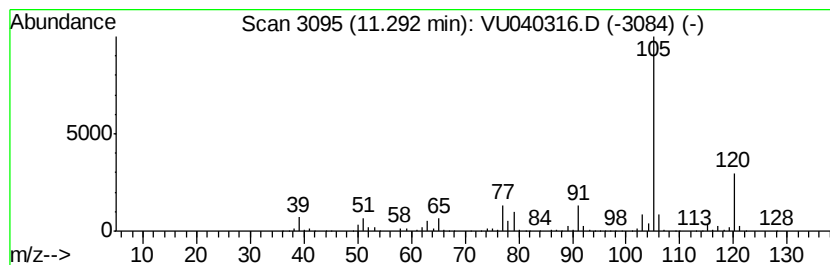
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	3.64 ug/L	2638210	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	94
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	93
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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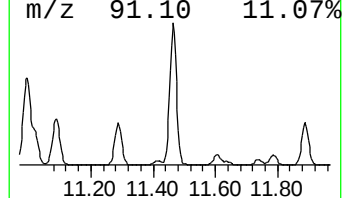
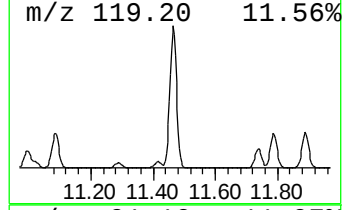
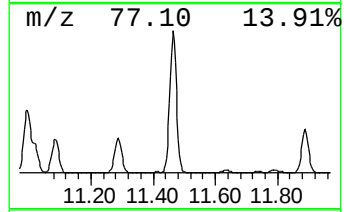
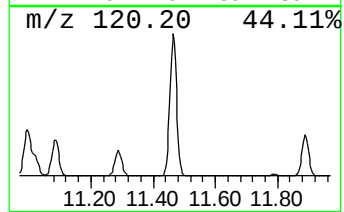
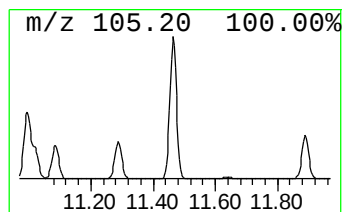
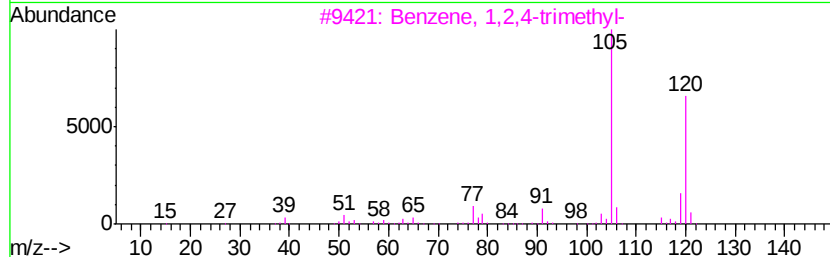
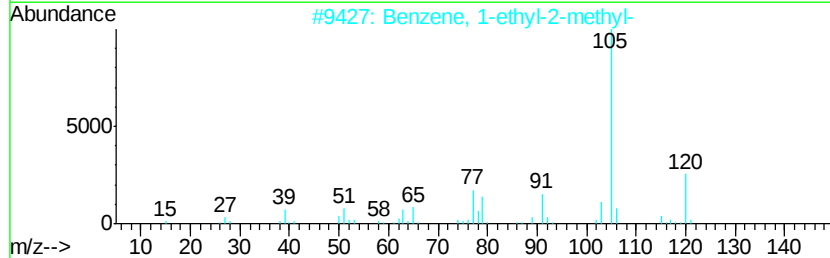
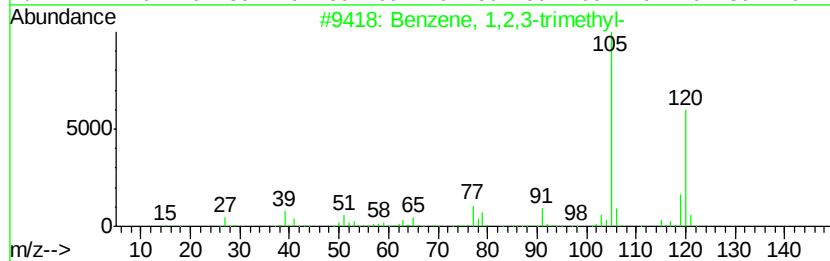
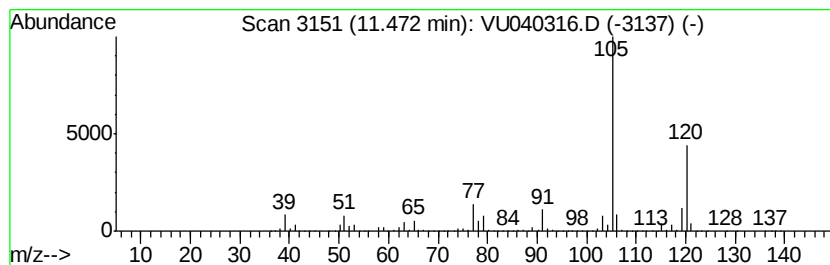
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	14.74 ug/L	10673000	1,4-Dichlorobenzene-d4	11.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Mesitylene	120	C9H12	000108-67-8	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



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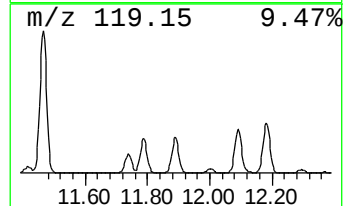
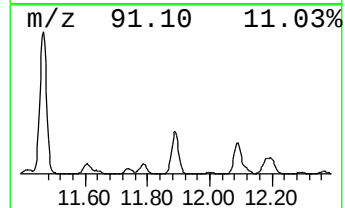
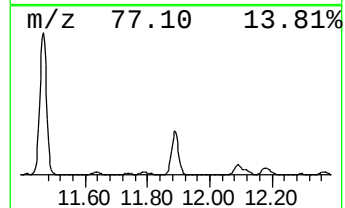
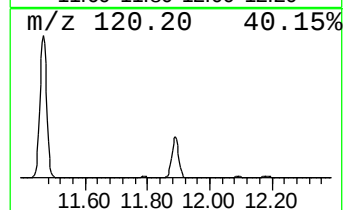
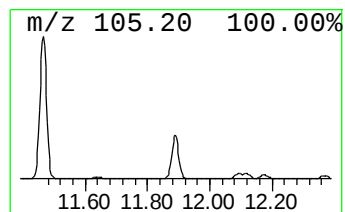
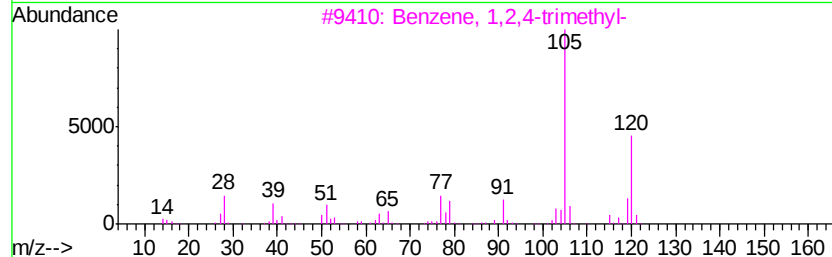
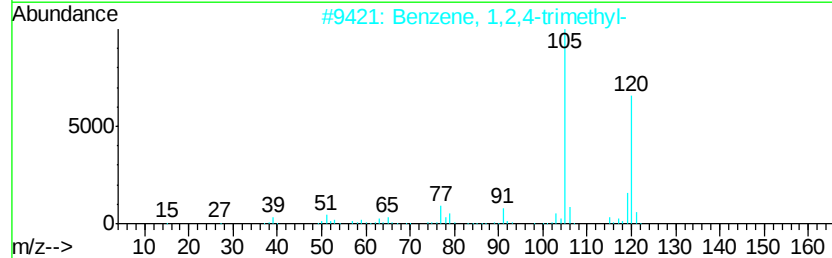
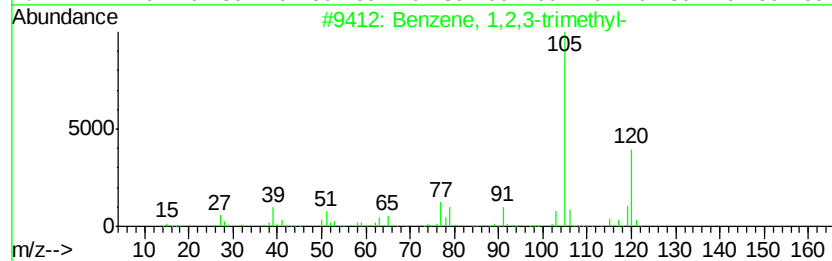
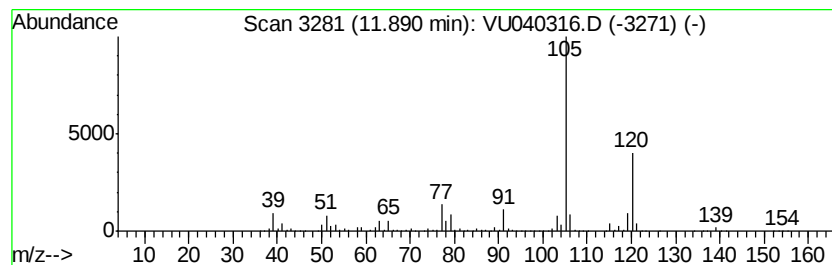
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ClientSampled :
BG494

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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	5.00 ug/L	3620640	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 97
2		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 94
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		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092620\
Data File : VU040316.D
Acq On : 25 Sep 2020 21:43
Operator : SY/MD
Sample : L4139-04
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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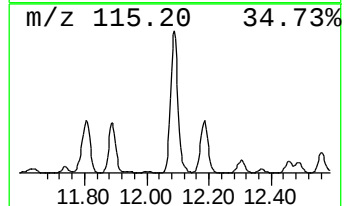
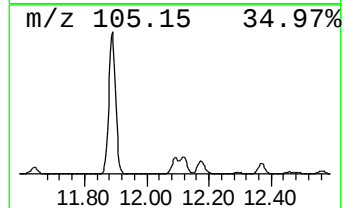
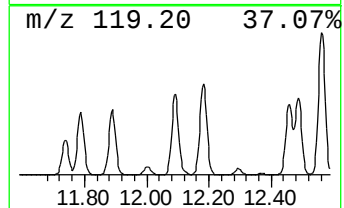
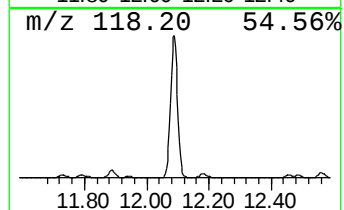
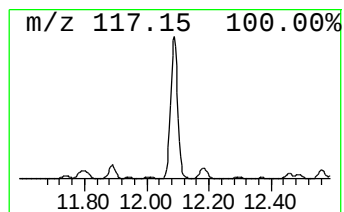
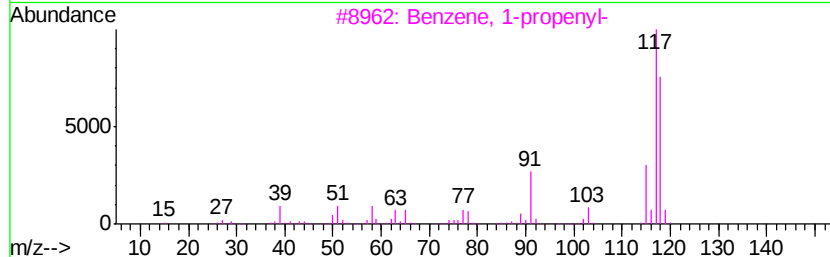
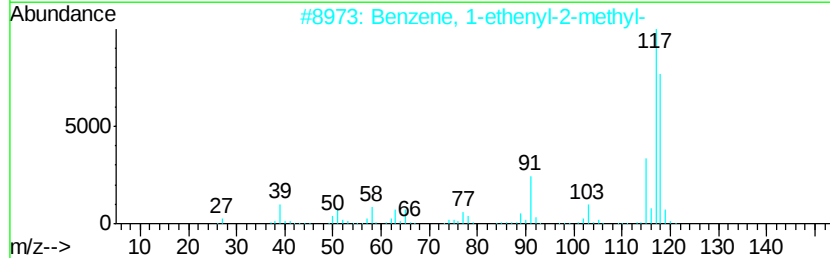
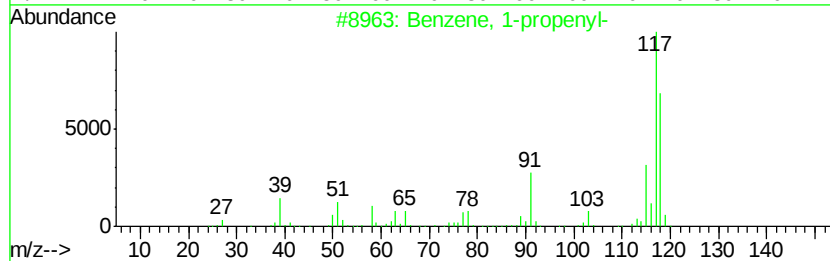
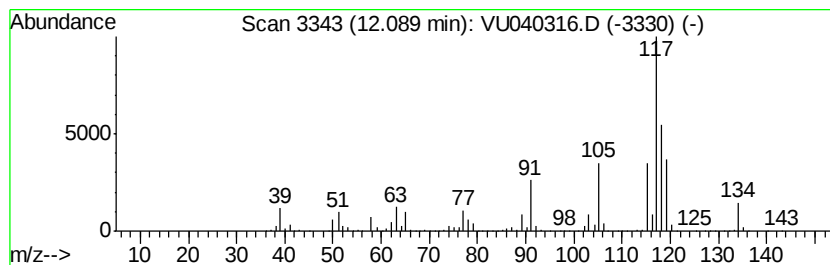
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 Benzene, 1-propenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	2.80 ug/L	2030400	1,4-Dichlorobenzene-d4	11.81

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092620\
Data File : VU040316.D
Acq On : 25 Sep 2020 21:43
Operator : SY/MD
Sample : L4139-04
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG494

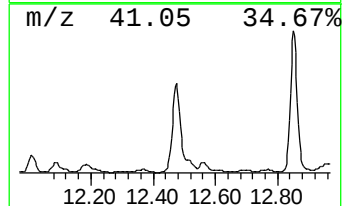
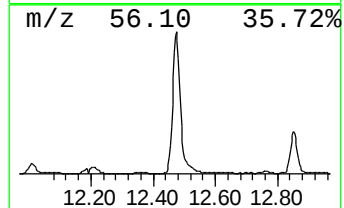
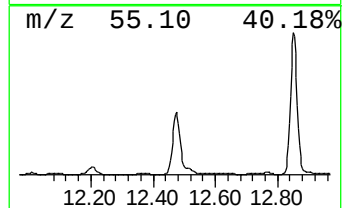
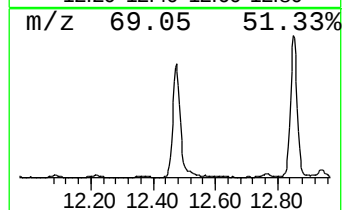
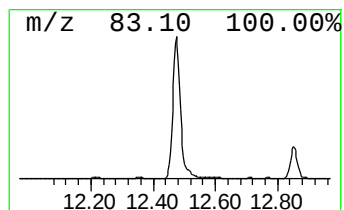
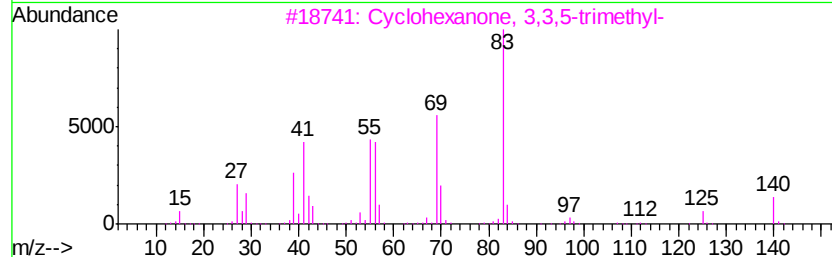
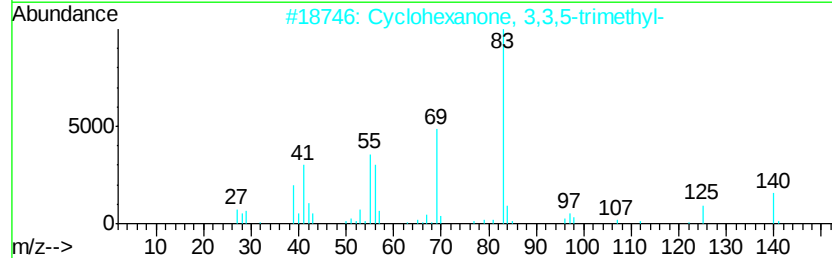
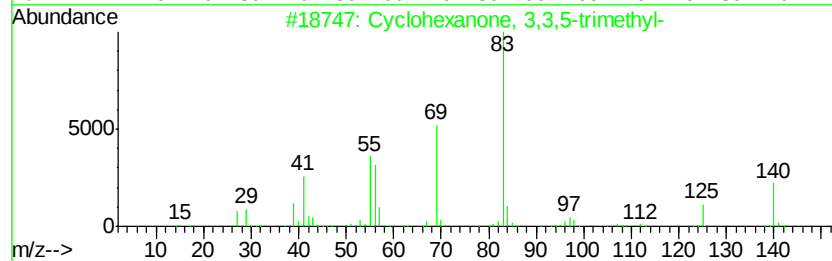
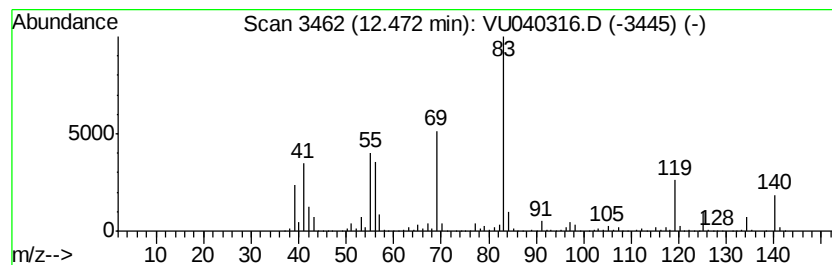
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 Cyclohexanone, 3,3,5-trimet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	3.04 ug/L	2200900	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4		1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	47
5		Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	47



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092620\
Data File : VU040316.D
Acq On : 25 Sep 2020 21:43
Operator : SY/MD
Sample : L4139-04
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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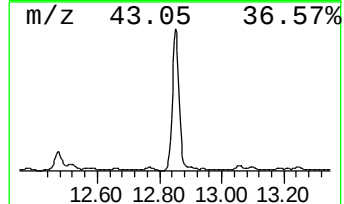
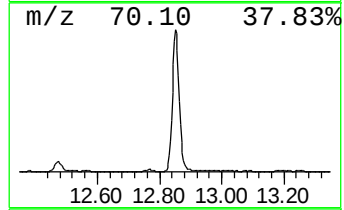
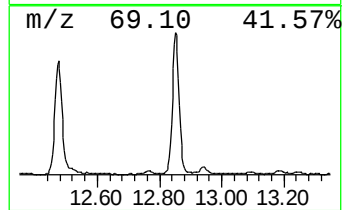
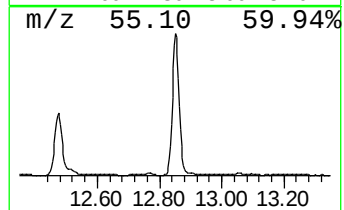
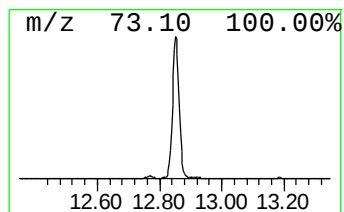
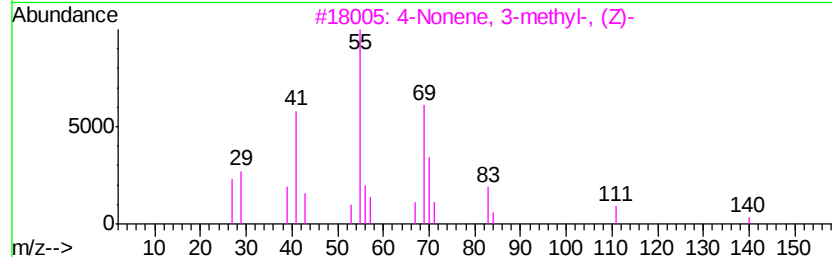
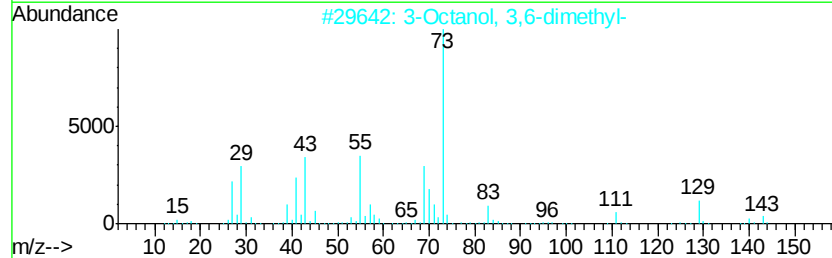
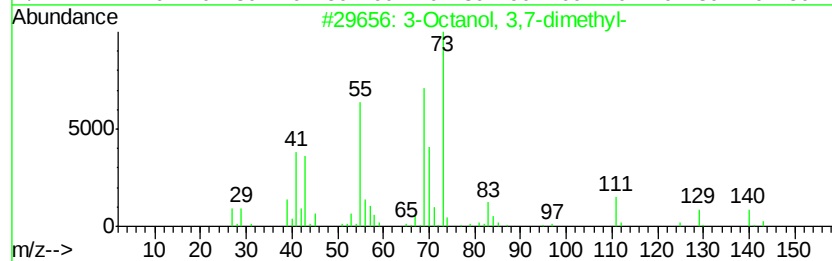
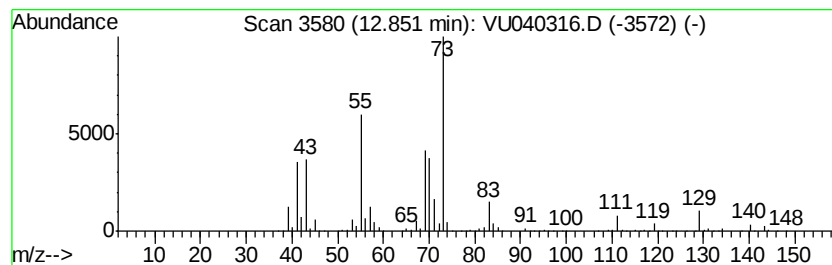
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 3-Octanol, 3,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.85	3.11 ug/L	2251490	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	53
2		3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	47
3		4-Nonene, 3-methyl-, (Z)-	140	C10H20	063830-69-3	46
4		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	43
5		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	43



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092620\
Data File : VU040316.D
Acq On : 25 Sep 2020 21:43
Operator : SY/MD
Sample : L4139-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG494

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.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	3.4	ug/L	2483300	3	11.81	3620640	5.0
Benzene, 1-ethyl-...	11.00	7.2	ug/L	5210780	3	11.81	3620640	5.0
Mesitylene	11.09	3.7	ug/L	2648820	3	11.81	3620640	5.0
Benzene, 1-ethyl-...	11.29	3.6	ug/L	2638210	3	11.81	3620640	5.0
Benzene, 1,2,3-tr...	11.47	14.7	ug/L	10673000	3	11.81	3620640	5.0
Benzene, 1,2,4-tr...	11.89	5.0	ug/L	3620640	3	11.81	3620640	5.0
Benzene, 1-propenyl-	12.09	2.8	ug/L	2030400	3	11.81	3620640	5.0
Cyclohexanone, 3,...	12.47	3.0	ug/L	2200900	3	11.81	3620640	5.0
3-Octanol, 3,7-di...	12.85	3.1	ug/L	2251490	3	11.81	3620640	5.0