

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100120\
 Data File : VU040438.D
 Acq On : 01 Oct 2020 22:37
 Operator : SY/MD
 Sample : L4240-08
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG3S5

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.1

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\SOMUTR092920WMA.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	3.112	516	551	594	rBV4	67738288	189679024	78.62%	28.424%
2	3.404	614	642	684	rBV3	80263735	241263938	100.00%	36.154%
3	4.324	909	928	949	rBV	3860184	11002112	4.56%	1.649%
4	4.465	957	972	985	rBV	1954502	5158019	2.14%	0.773%
5	4.578	985	1007	1023	rVB3	59684837	161309112	66.86%	24.173%
6	5.407	1249	1265	1281	rVB2	3842265	8602420	3.57%	1.289%
7	10.902	2957	2974	2989	rBV	3213805	4786404	1.98%	0.717%
8	10.996	2989	3003	3009	rBV	6667794	11187412	4.64%	1.676%
9	11.086	3021	3031	3056	rVB	4710082	7070632	2.93%	1.060%
10	11.465	3136	3149	3165	rVB	7313619	11087344	4.60%	1.661%
11	12.182	3361	3372	3396	rVB3	2007952	3763116	1.56%	0.564%
12	12.565	3479	3491	3500	rBV	1962920	2950964	1.22%	0.442%
13	13.018	3623	3632	3646	rVB	2009558	2995984	1.24%	0.449%
14	13.442	3754	3764	3789	rVB4	1456848	2661270	1.10%	0.399%
15	13.828	3872	3884	3902	rVB2	2297818	3799300	1.57%	0.569%

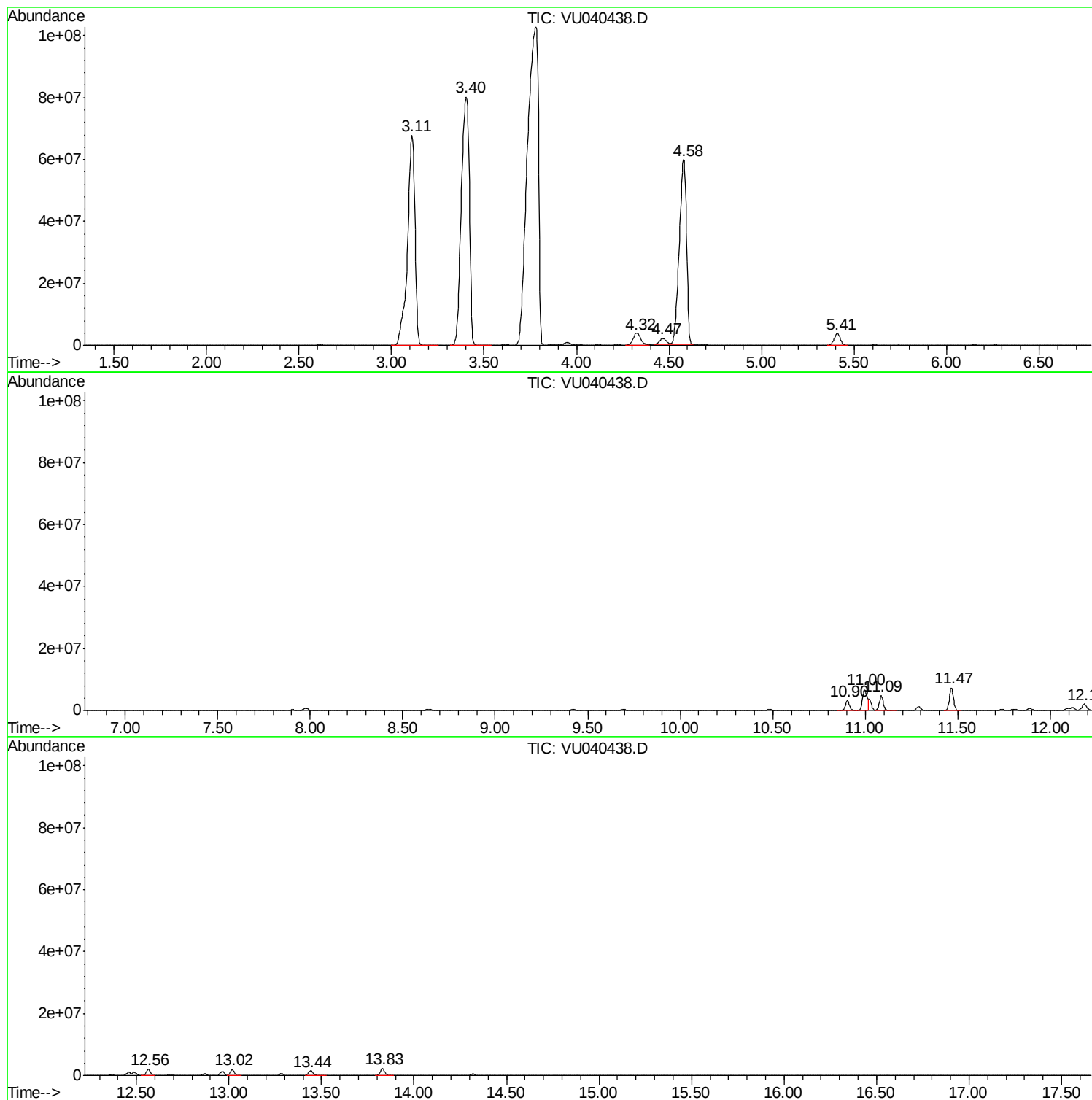
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
Quant Title : TRACE VOA SOM01.0

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



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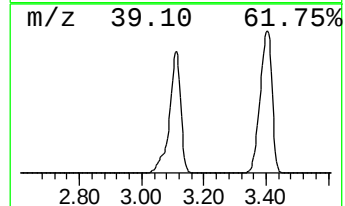
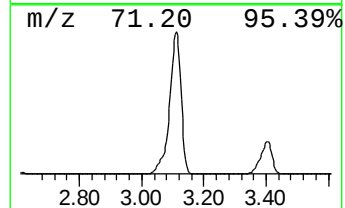
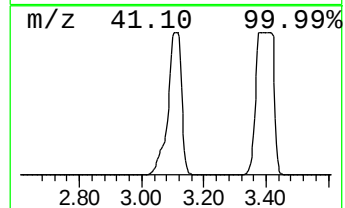
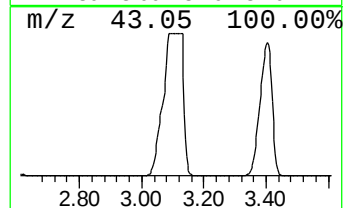
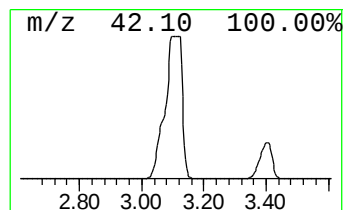
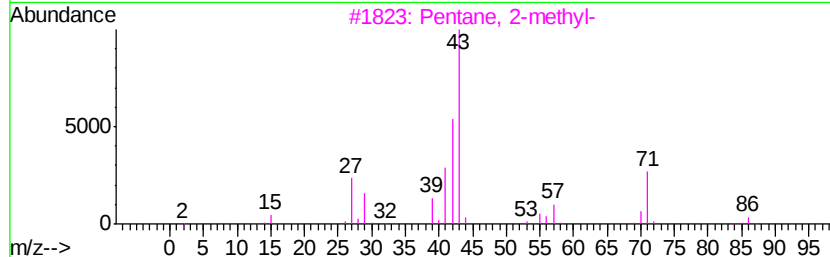
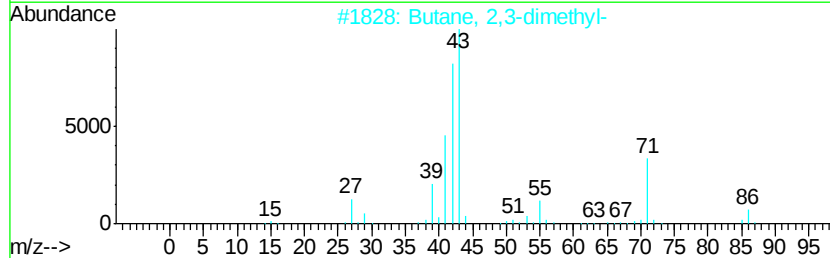
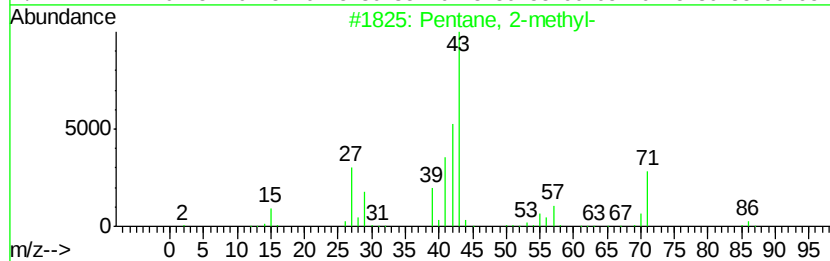
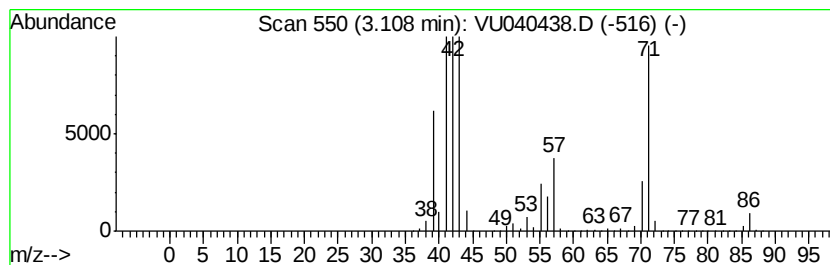
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.11	110.25 ug/L	189679000	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	49
2		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	38
3		Pentane, 2-methyl-	86	C6H14	000107-83-5	38
4		3-Penten-2-ol	86	C5H10O	001569-50-2	37
5		3-Penten-2-ol	86	C5H10O	001569-50-2	37



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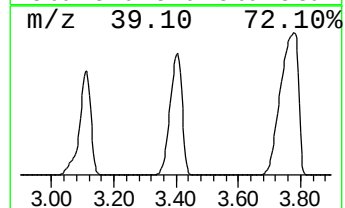
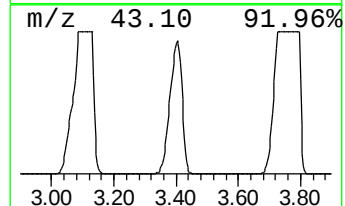
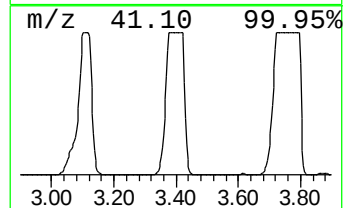
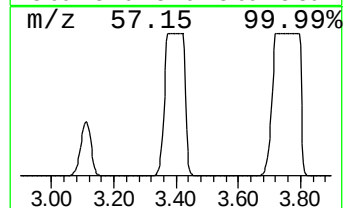
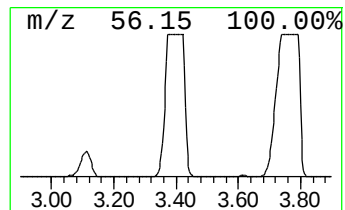
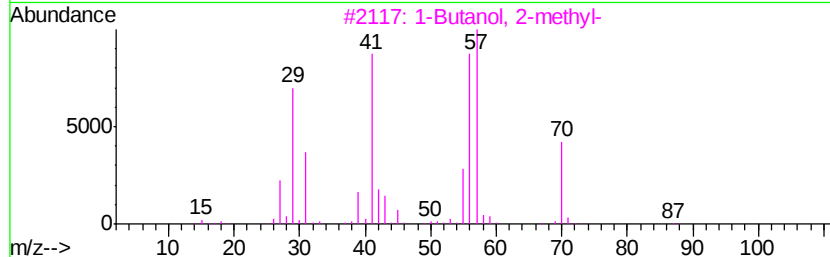
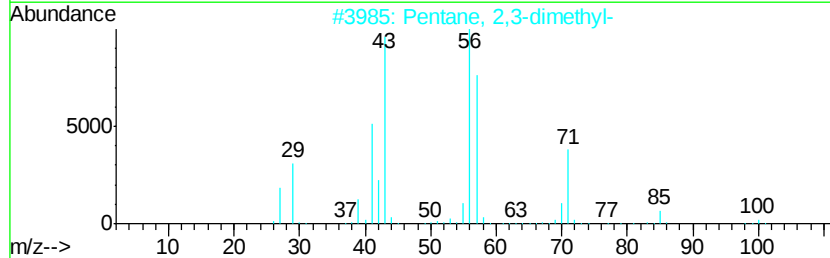
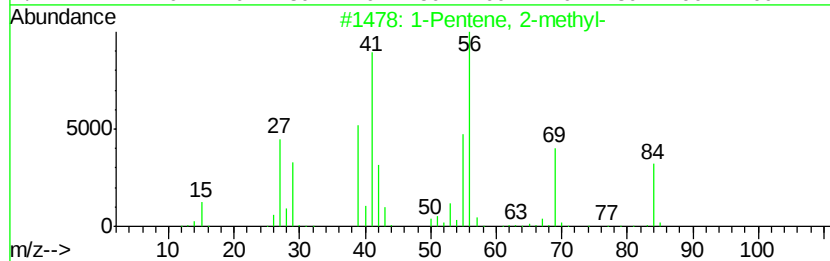
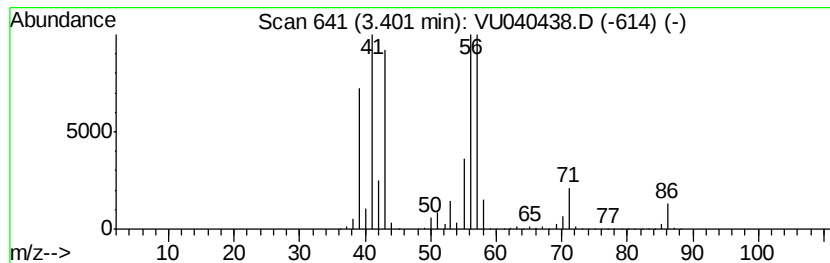
Instrument :
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ClientSampleId :
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TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.40	140.23 ug/L	241264000	1,4-Difluorobenzene	6.26	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	53
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	53
3	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	43
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	42
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	38



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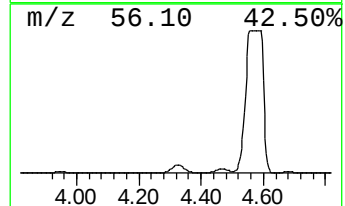
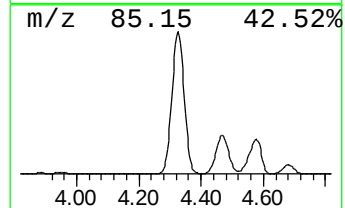
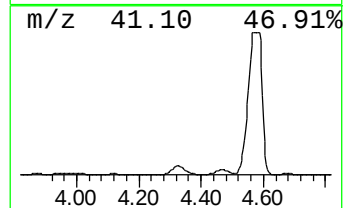
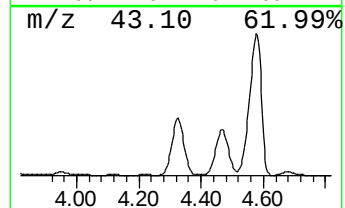
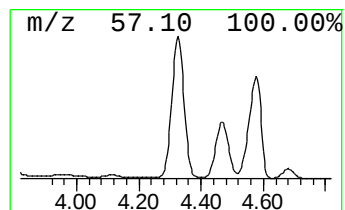
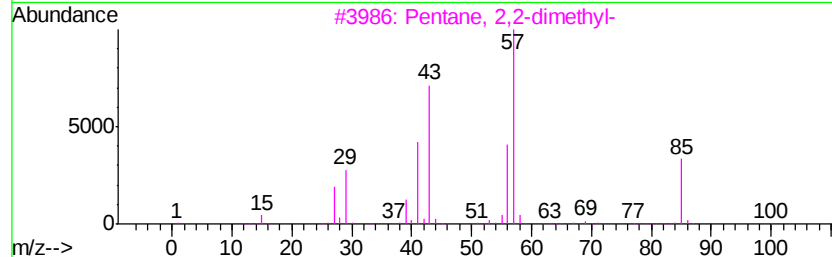
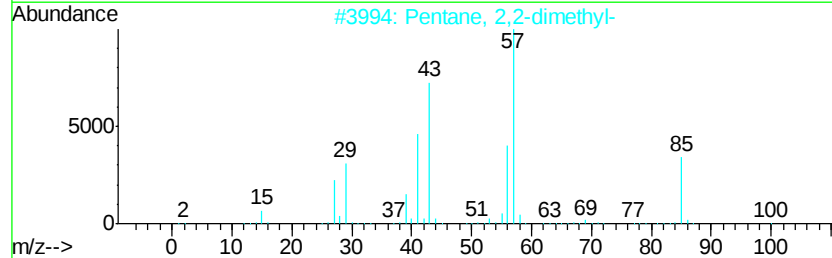
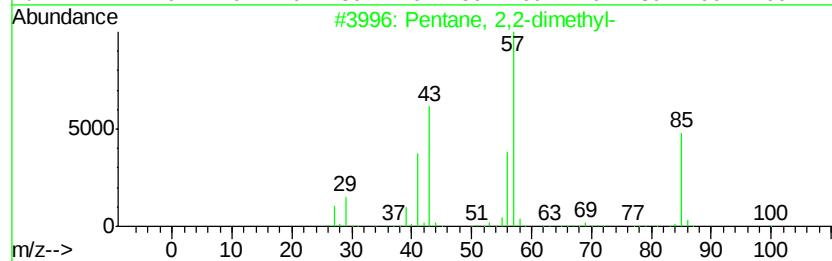
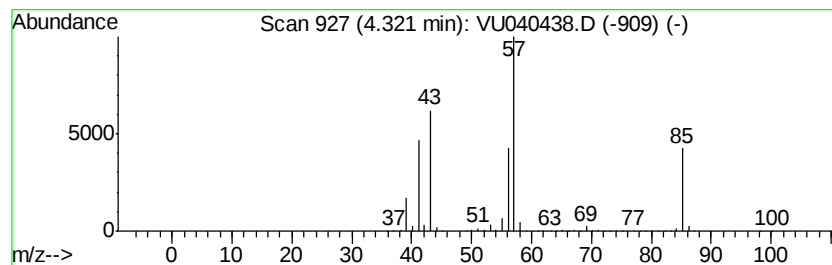
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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.32	6.39 ug/L	11002100	1,4-Difluorobenzene	6.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	90	
2	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83	
3	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83	
4	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83	
5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	64	



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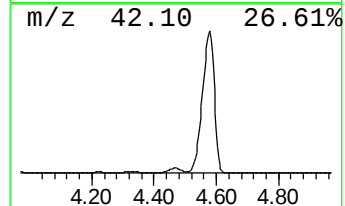
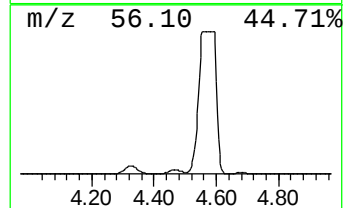
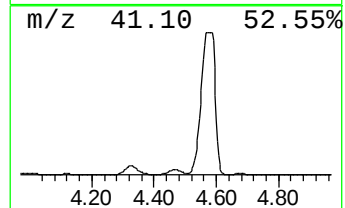
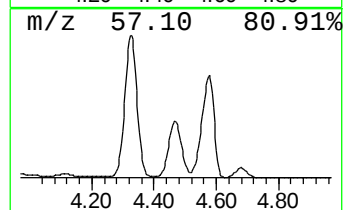
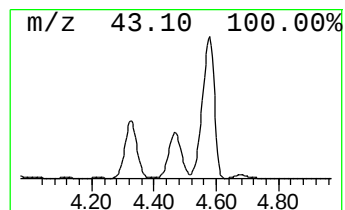
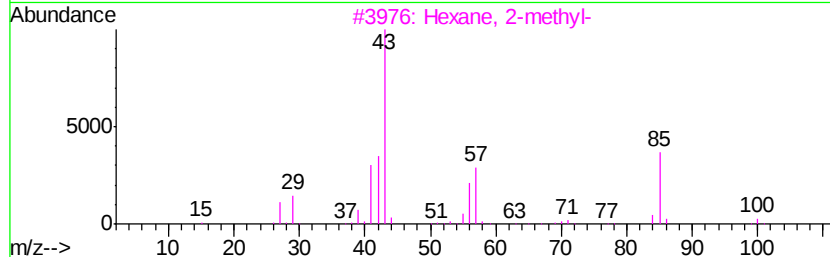
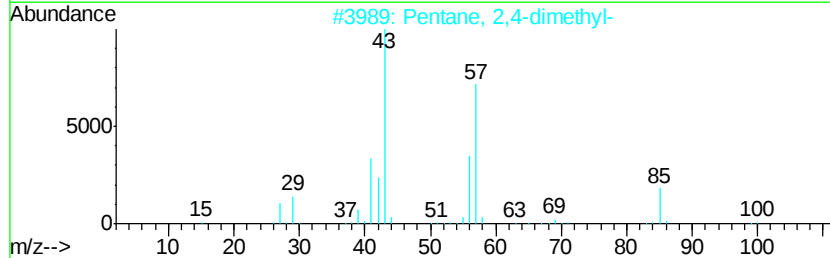
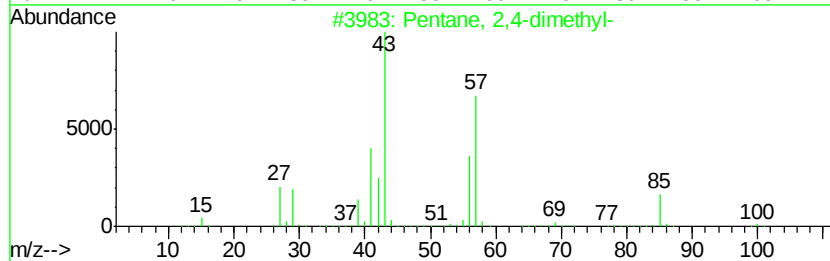
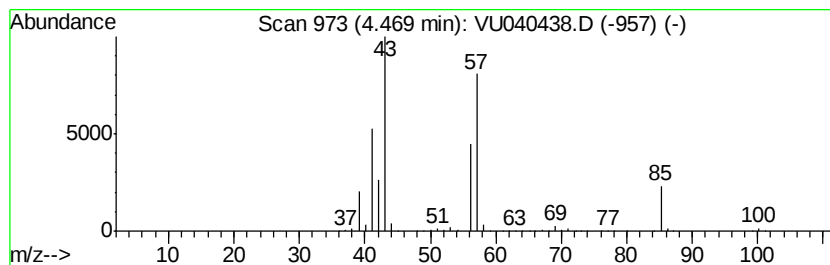
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.47	3.00 ug/L	5158020	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	90
3		Hexane, 2-methyl-	100	C7H16	000591-76-4	72
4		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
5		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64



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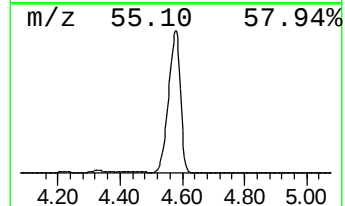
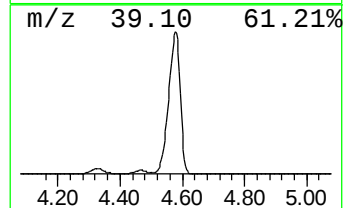
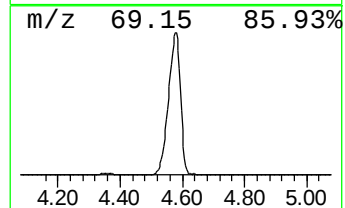
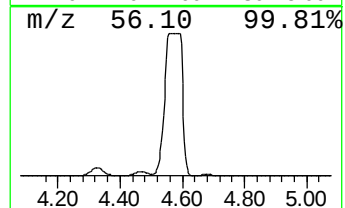
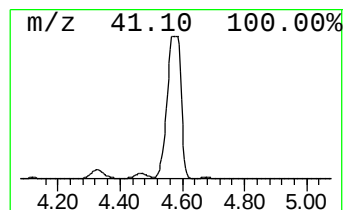
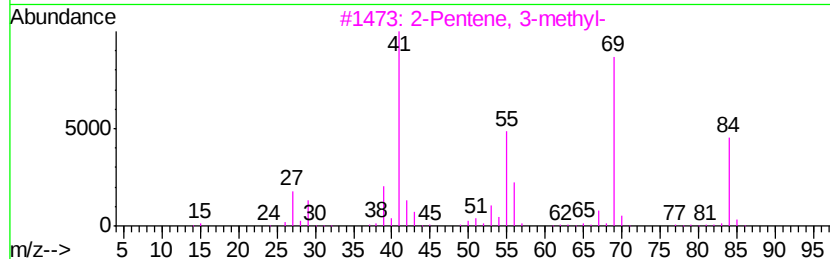
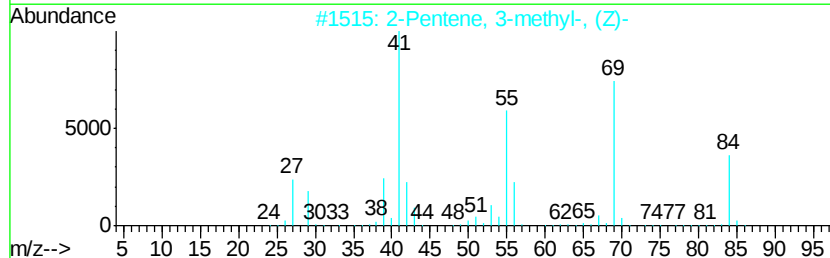
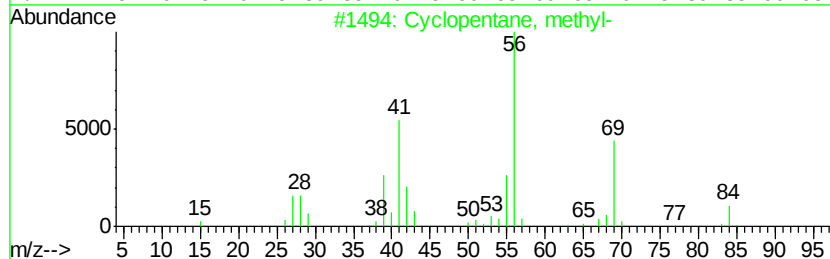
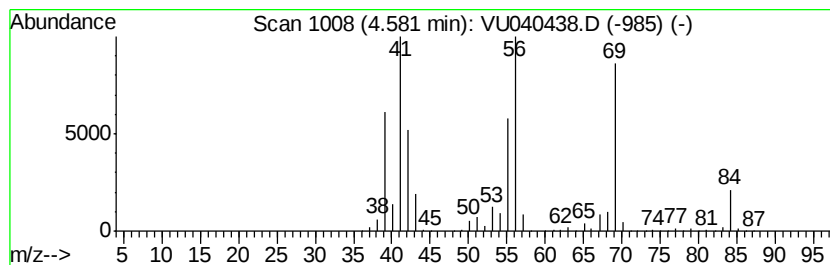
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic4.58 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.58	93.76 ug/L	161309000	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	60
2		2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	55
3		2-Pentene, 3-methyl-	84	C6H12	000922-61-2	49
4		Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	46
5		Cyclopropane, 1,2,3-trimethyl-	84	C6H12	042984-19-0	43



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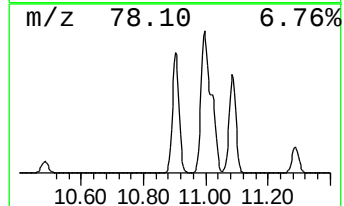
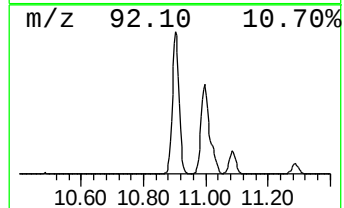
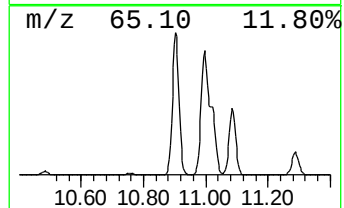
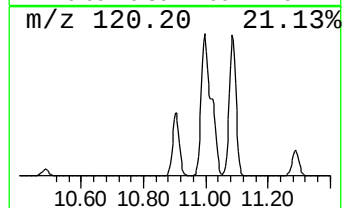
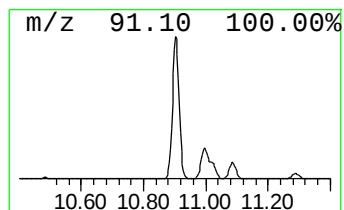
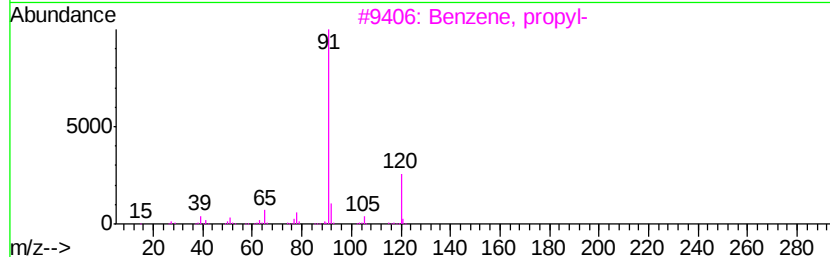
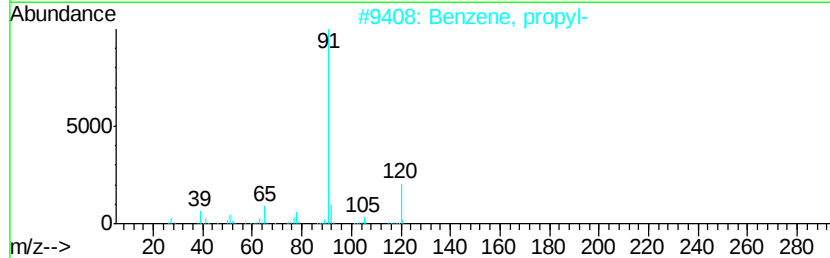
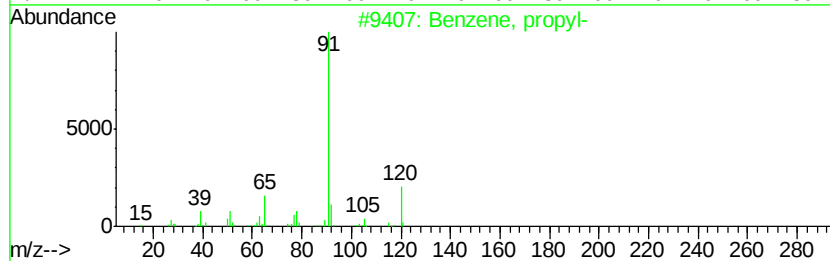
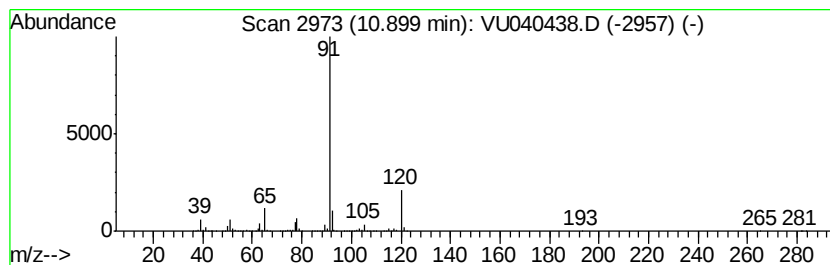
Instrument :
MSVOA_U
ClientSampleId :
BG3S5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 Benzene, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	2.16 ug/L	4786400	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, propyl-	120 C9H12	000103-65-1 94
2		Benzene, propyl-	120 C9H12	000103-65-1 90
3		Benzene, propyl-	120 C9H12	000103-65-1 86
4		N-Benzyl-2-phenethylamine	211 C15H17N	003647-71-0 83
5		N-Benzyl-2-phenethylamine	211 C15H17N	003647-71-0 78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040438.D
Acq On : 01 Oct 2020 22:37
Operator : SY/MD
Sample : L4240-08
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3S5

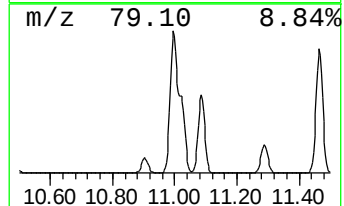
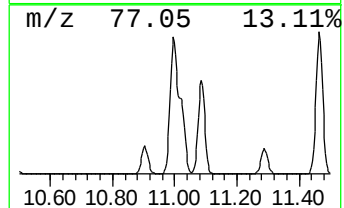
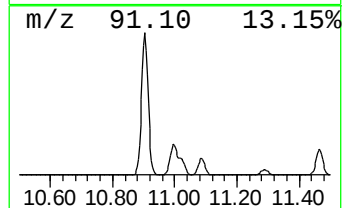
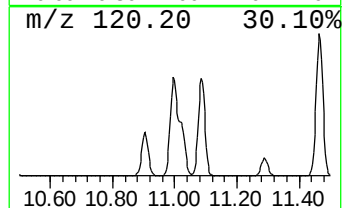
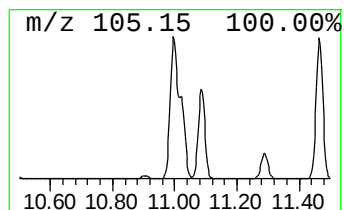
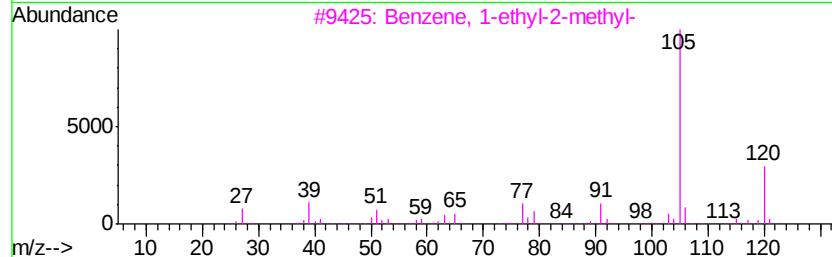
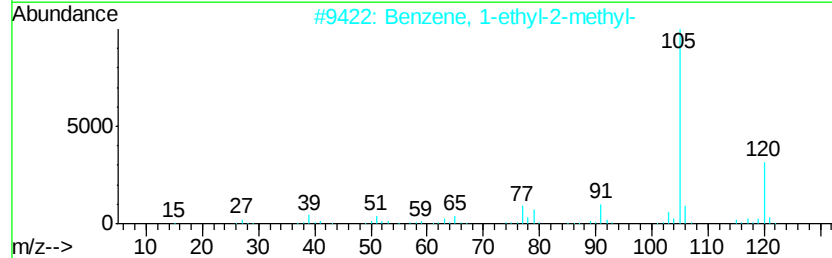
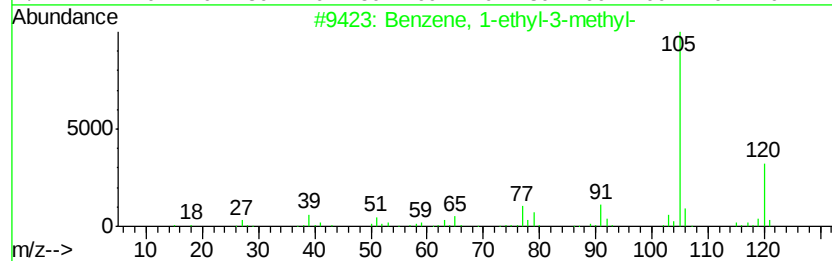
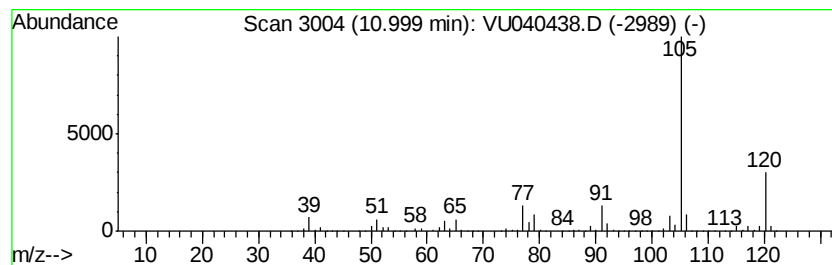
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	5.05 ug/L	11187400	1,4-Dichlorobenzene-d4	11.81

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040438.D
Acq On : 01 Oct 2020 22:37
Operator : SY/MD
Sample : L4240-08
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3S5

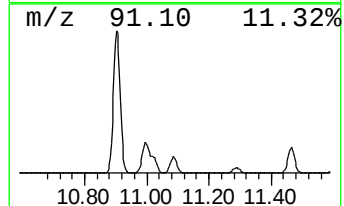
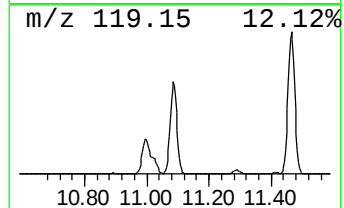
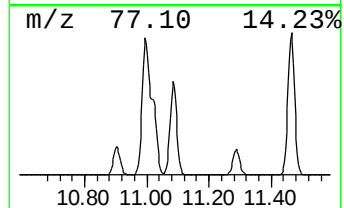
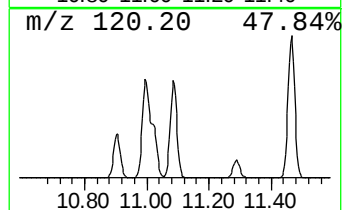
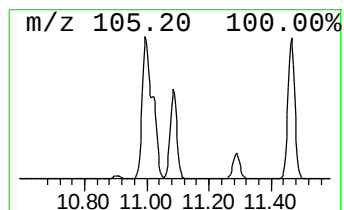
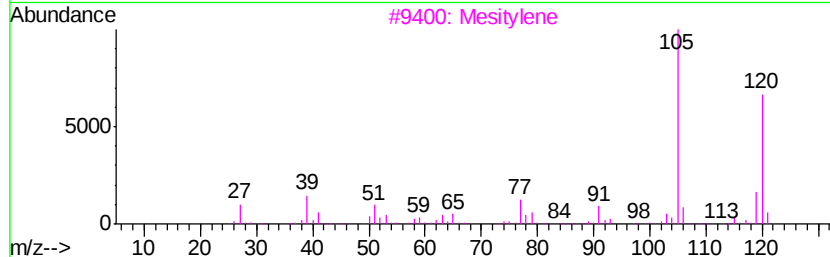
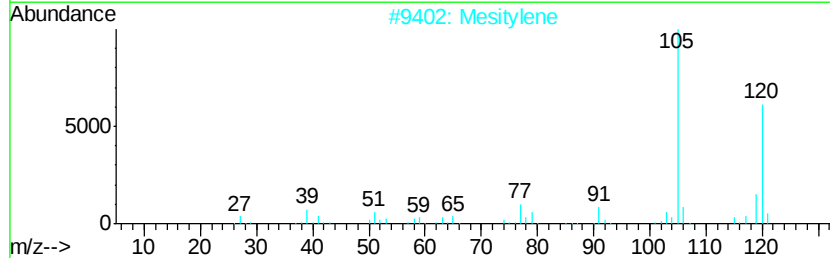
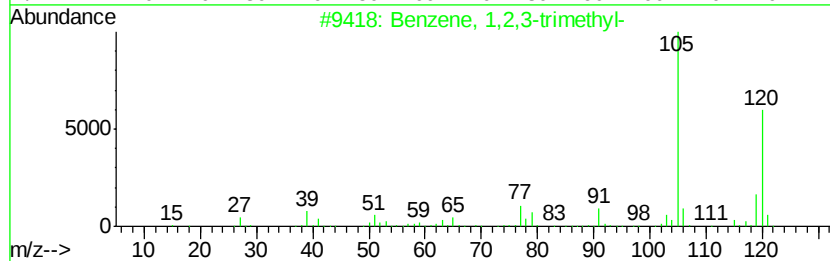
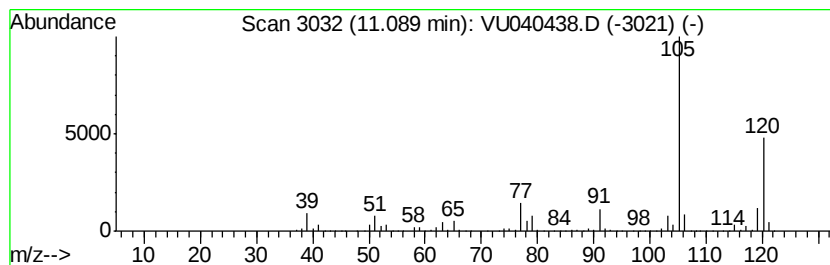
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	3.19 ug/L	7070630	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		Mesitylene	120	C9H12	000108-67-8	97
3		Mesitylene	120	C9H12	000108-67-8	95
4		Mesitylene	120	C9H12	000108-67-8	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040438.D
Acq On : 01 Oct 2020 22:37
Operator : SY/MD
Sample : L4240-08
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3S5

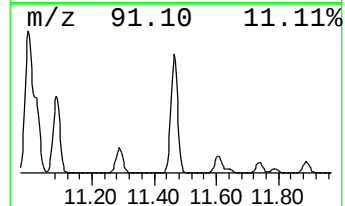
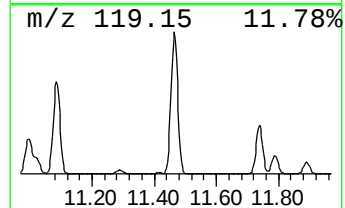
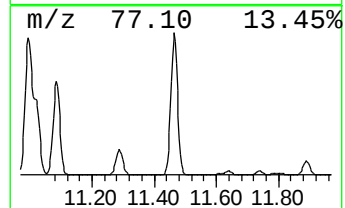
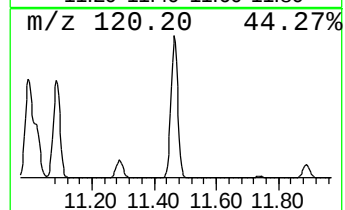
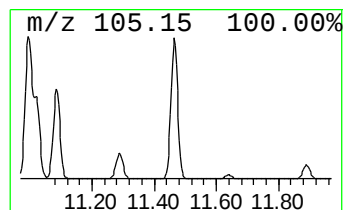
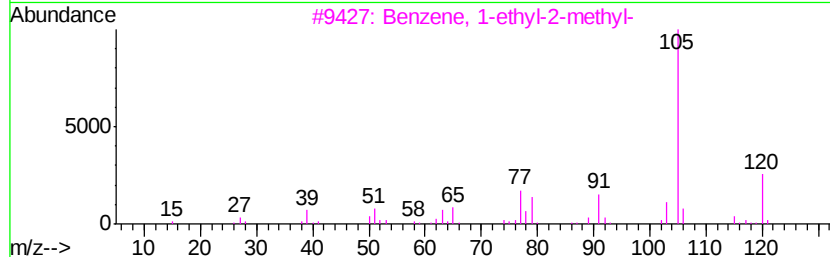
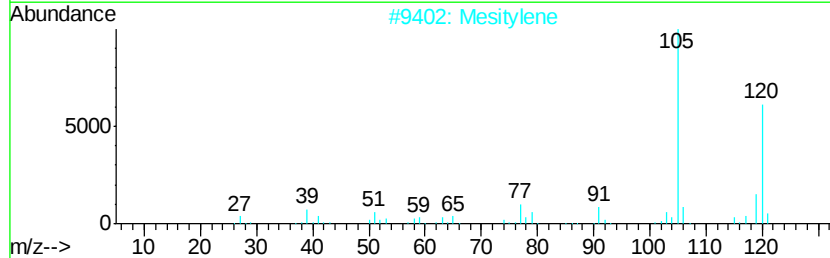
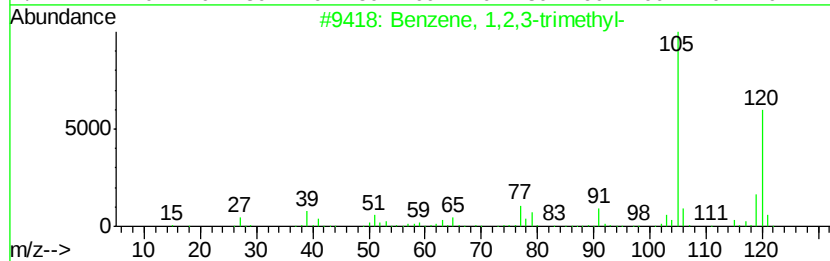
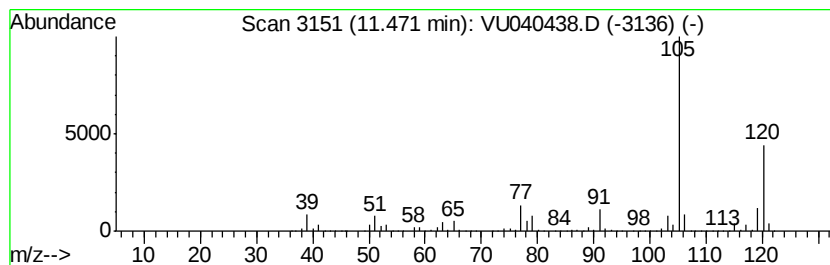
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	5.00 ug/L	11087300	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		Mesitylene	120	C9H12	000108-67-8	95
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040438.D
Acq On : 01 Oct 2020 22:37
Operator : SY/MD
Sample : L4240-08
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3S5

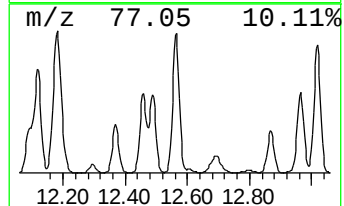
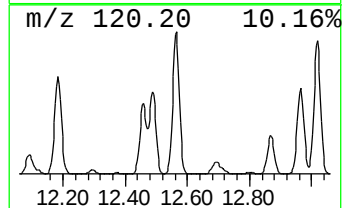
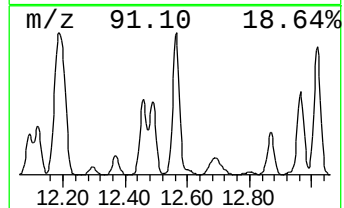
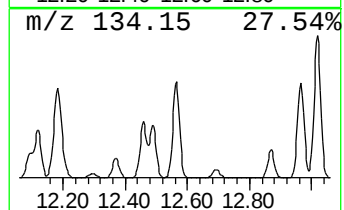
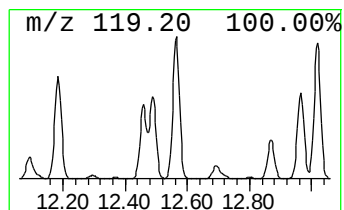
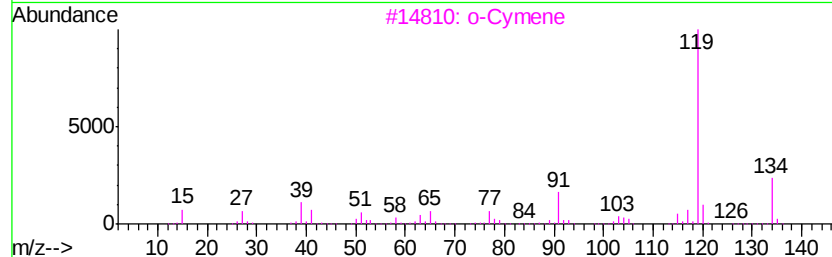
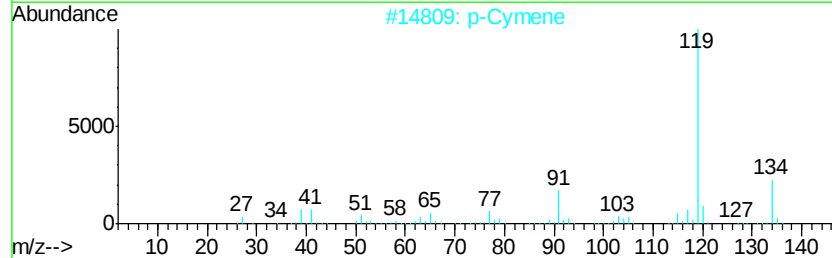
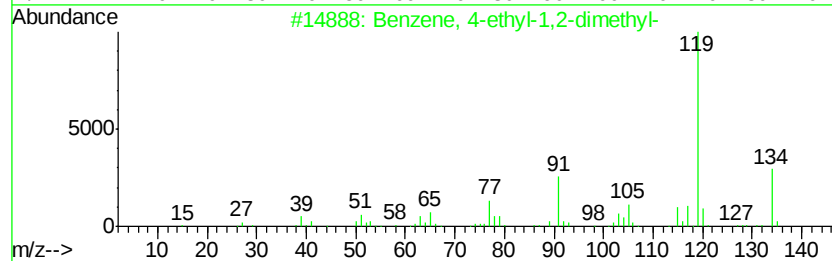
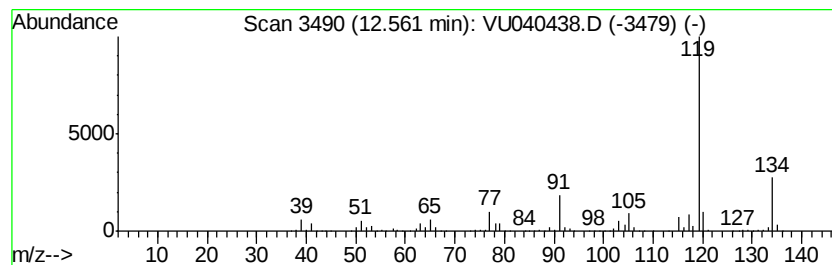
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	1.33 ug/L	2950960	1,4-Dichlorobenzene-d4	11.81

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040438.D
Acq On : 01 Oct 2020 22:37
Operator : SY/MD
Sample : L4240-08
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3S5

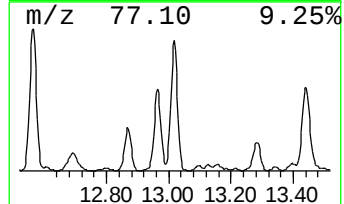
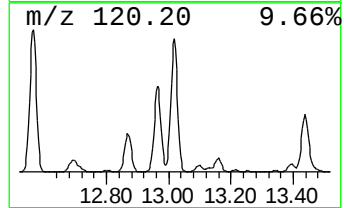
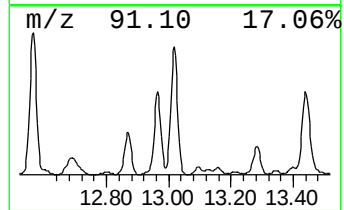
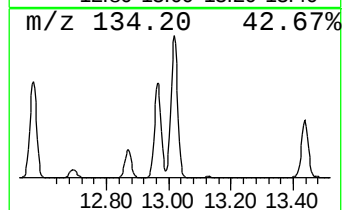
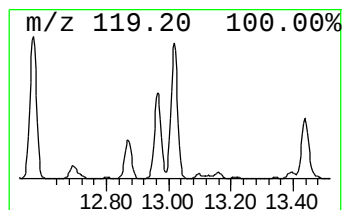
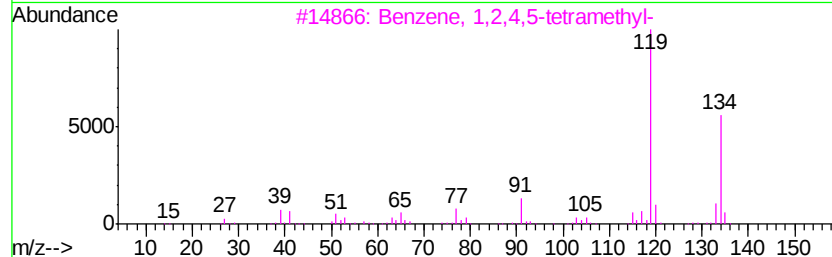
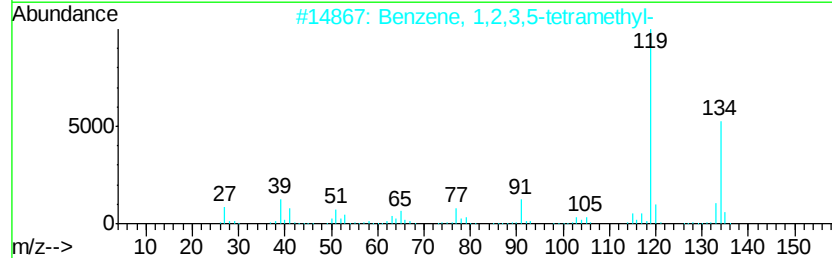
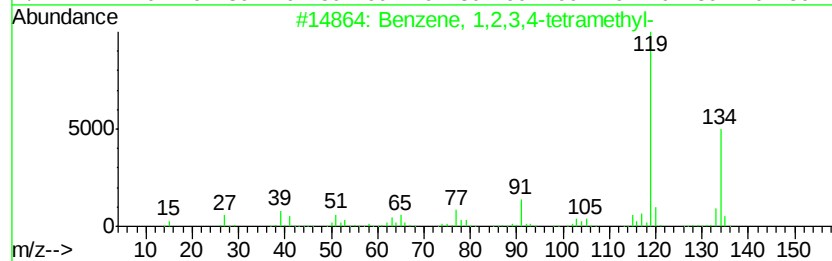
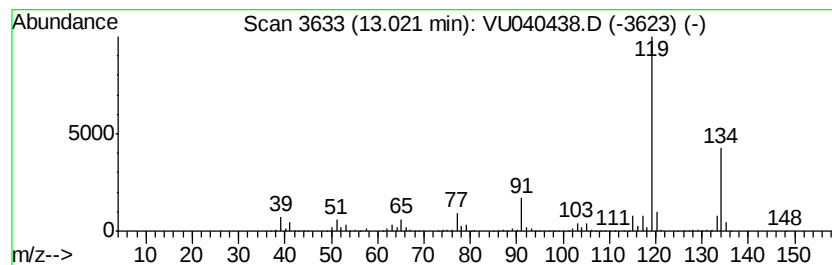
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	1.35 ug/L	2995980	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040438.D
Acq On : 01 Oct 2020 22:37
Operator : SY/MD
Sample : L4240-08
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3S5

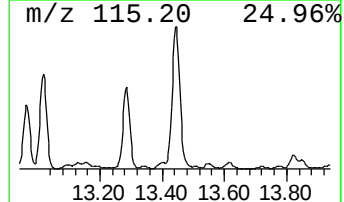
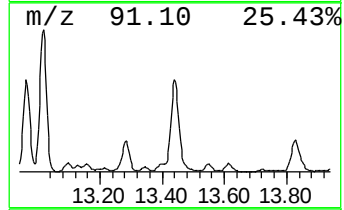
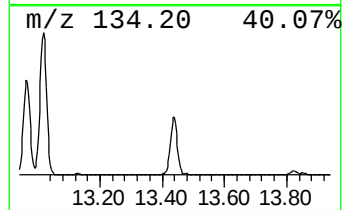
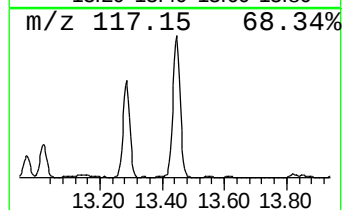
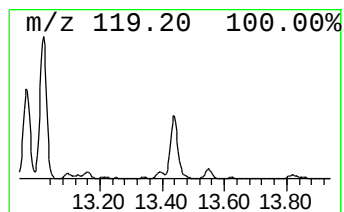
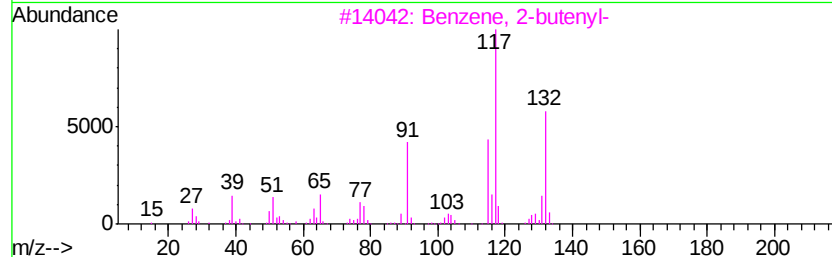
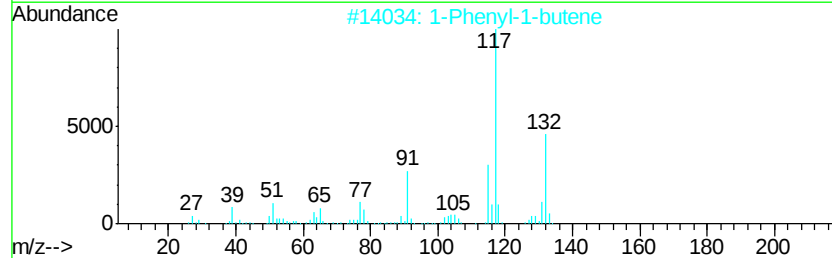
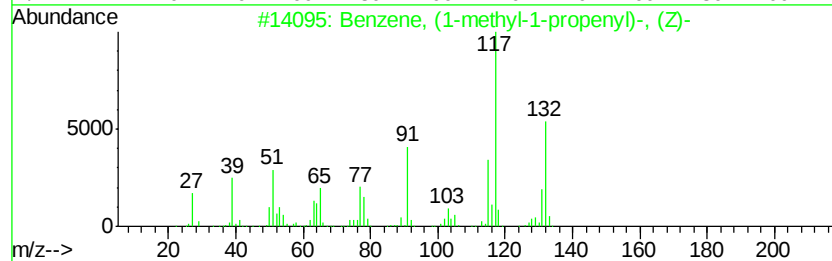
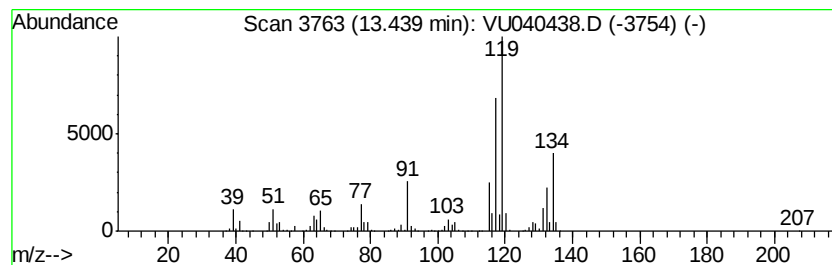
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Benzene, (1-methyl-1-propen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	1.20 ug/L	2661270	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000767-99-7	87
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	74
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100120\
 Data File : VU040438.D
 Acq On : 01 Oct 2020 22:37
 Operator : SY/MD
 Sample : L4240-08
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG3S5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.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
(DEL) Alkane: Str...	3.11	110.3	ug/L	189679000	1	6.26	8602420	5.0
(DEL) Alkane: Str...	3.40	140.2	ug/L	241264000	1	6.26	8602420	5.0
(DEL) Alkane: Str...	4.32	6.4	ug/L	11002100	1	6.26	8602420	5.0
(DEL) Alkane: Str...	4.47	3.0	ug/L	5158020	1	6.26	8602420	5.0
(DEL) Alkane: Cyc...	4.58	93.8	ug/L	161309000	1	6.26	8602420	5.0
Benzene, propyl-	10.90	2.2	ug/L	4786400	3	11.81	11087300	5.0
Benzene, 1-ethyl-...	11.00	5.0	ug/L	11187400	3	11.81	11087300	5.0
Benzene, 1,2,3-tr...	11.09	3.2	ug/L	7070630	3	11.81	11087300	5.0
Mesitylene	11.47	5.0	ug/L	11087300	3	11.81	11087300	5.0
Benzene, 4-ethyl-...	12.56	1.3	ug/L	2950960	3	11.81	11087300	5.0
Benzene, 1,2,3,4-...	13.02	1.4	ug/L	2995980	3	11.81	11087300	5.0
Benzene, (1-methy...	13.44	1.2	ug/L	2661270	3	11.81	11087300	5.0