

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100120\
 Data File : VU040440.D
 Acq On : 01 Oct 2020 23:24
 Operator : SY/MD
 Sample : L4240-11
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG3T0

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	3.102	516	548	570	rBV	24271877	54030134	26.48%	12.127%
2	3.385	614	636	665	rBV2	36001286	82120619	40.25%	18.431%
3	3.748	720	749	770	rBV5	74076793	204035828	100.00%	45.794%
4	4.321	907	927	949	rBV	4077682	11180693	5.48%	2.509%
5	4.459	952	970	984	rVV	3171121	8180429	4.01%	1.836%
6	4.558	984	1001	1019	rVV	17587630	41897822	20.53%	9.404%
7	5.404	1243	1264	1279	rBV2	1696868	4158214	2.04%	0.933%
8	10.902	2959	2974	2990	rBV	1643309	2519480	1.23%	0.565%
9	10.996	2990	3003	3021	rVV	4898164	10608783	5.20%	2.381%
10	11.086	3021	3031	3052	rVB	2787204	4250944	2.08%	0.954%
11	11.288	3082	3094	3114	rVB	1437985	2216412	1.09%	0.497%
12	11.465	3135	3149	3171	rVB	7011996	10604362	5.20%	2.380%
13	12.182	3361	3372	3392	rVB7	1551530	3018708	1.48%	0.678%
14	12.565	3479	3491	3501	rBV	1492582	2220527	1.09%	0.498%
15	13.018	3623	3632	3646	rVB	1503272	2276916	1.12%	0.511%
16	13.442	3755	3764	3789	rVB3	1089252	2227876	1.09%	0.500%

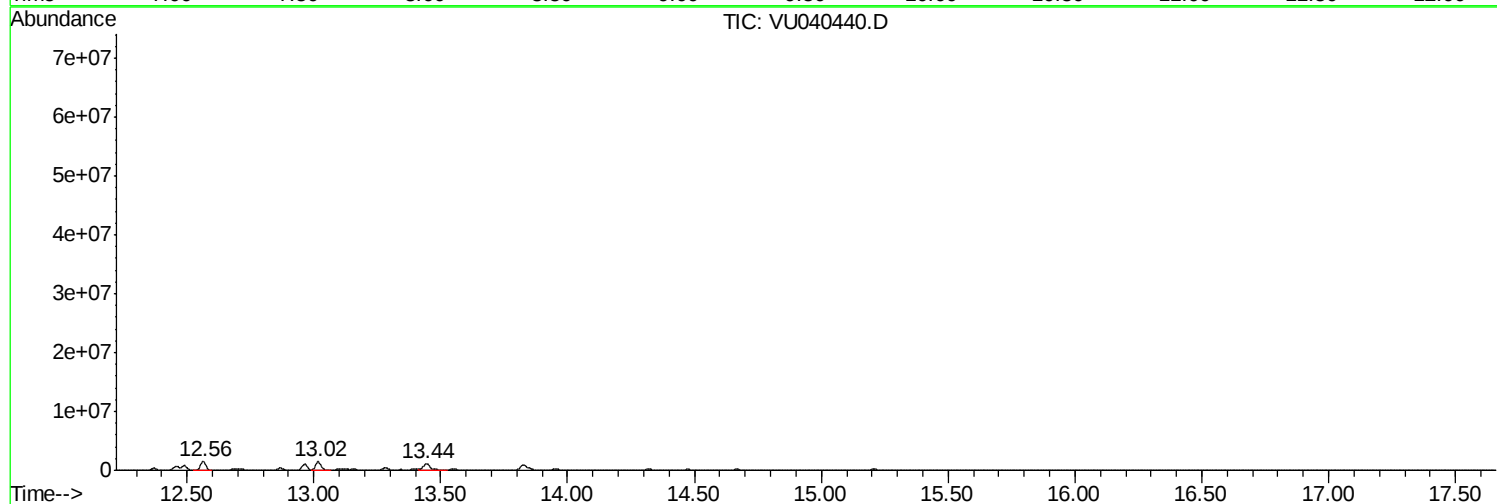
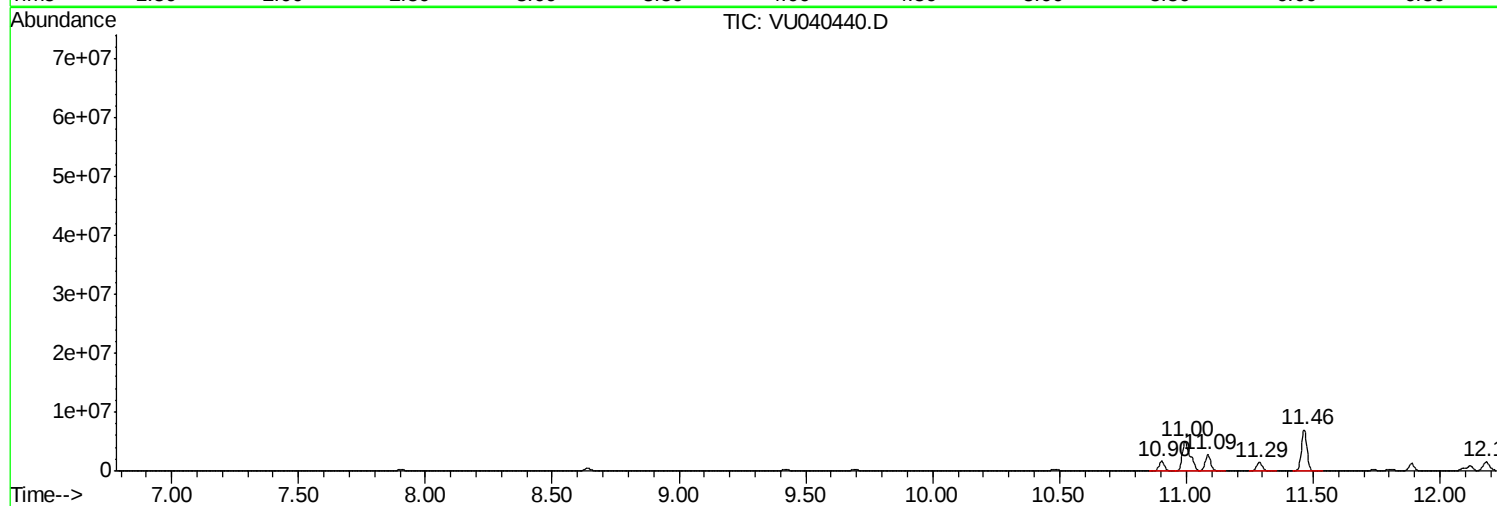
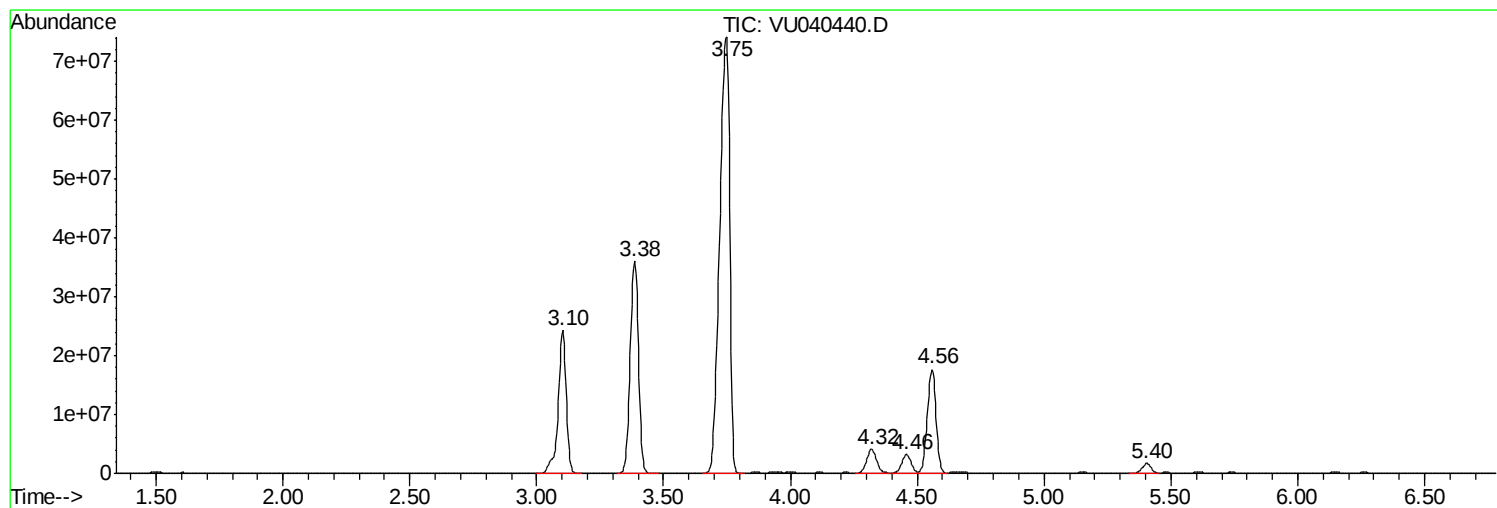
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
Quant Title : TRACE VOA SOM01.0

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



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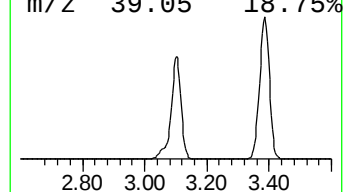
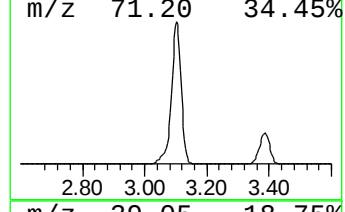
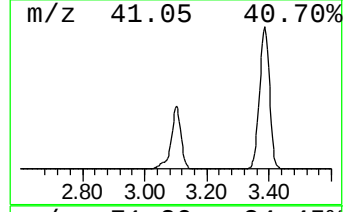
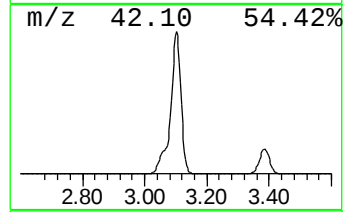
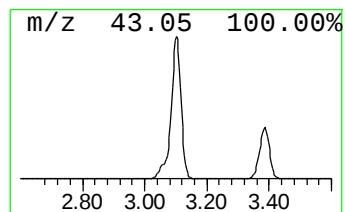
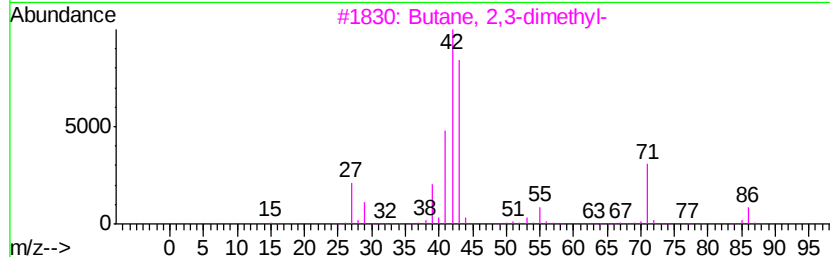
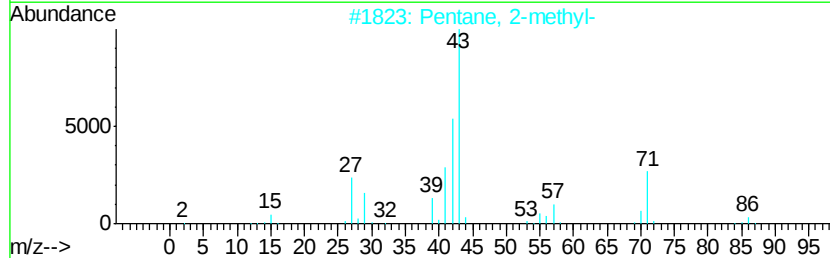
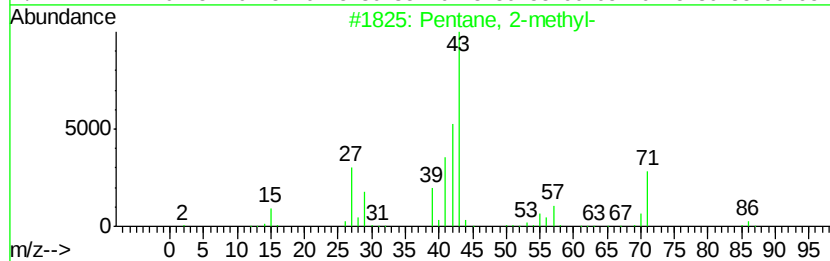
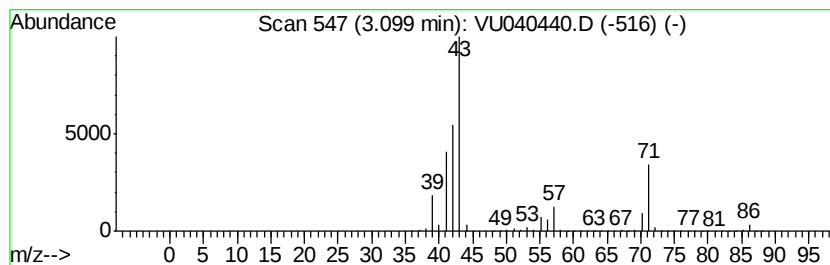
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	64.97 ug/L	54030100	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2		Pentane, 2-methyl-	86	C6H14	000107-83-5	91
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	58
4		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	52
5		Pentane	72	C5H12	000109-66-0	40



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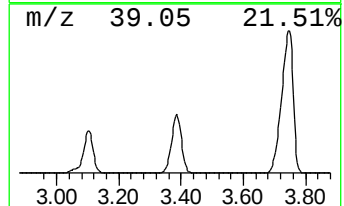
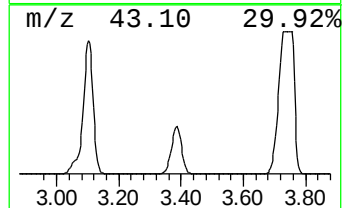
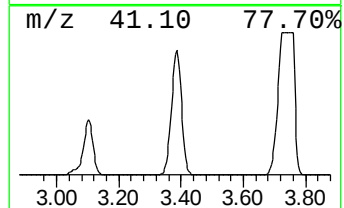
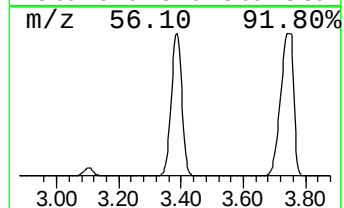
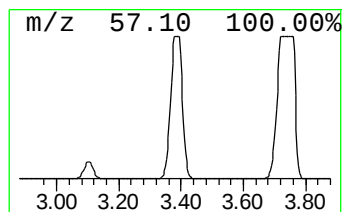
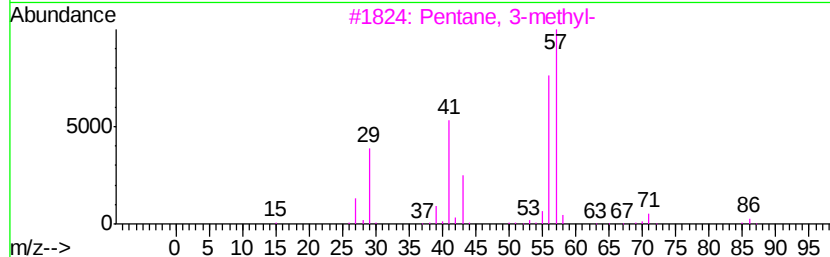
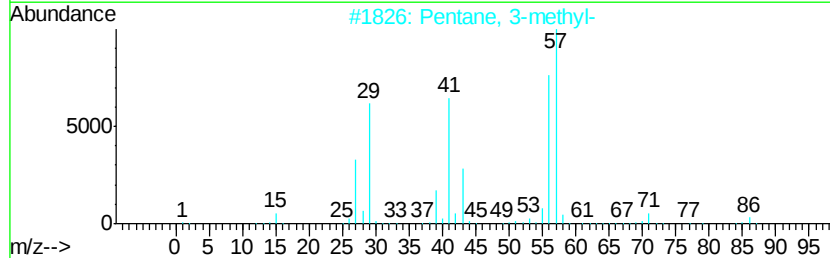
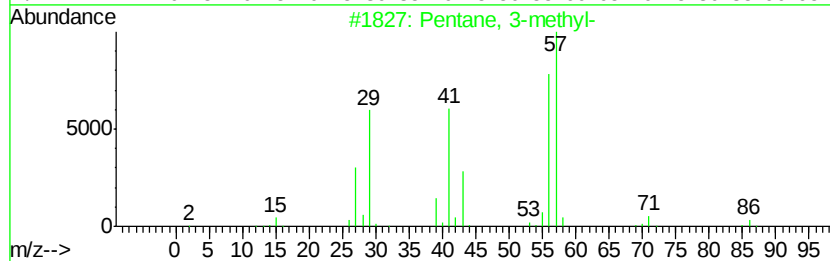
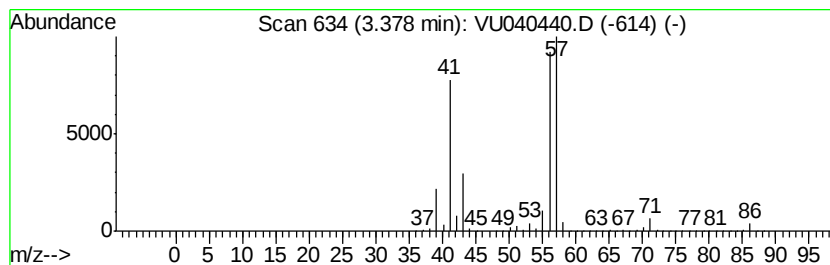
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.38	98.75 ug/L	82120600	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	90
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	72
4		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
5		1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40



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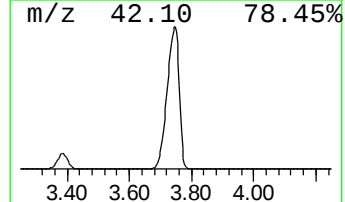
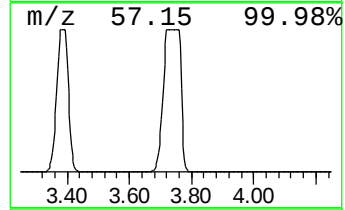
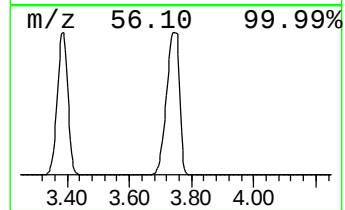
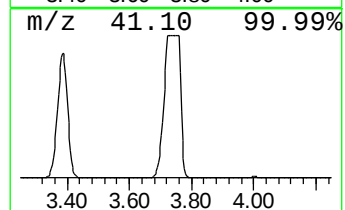
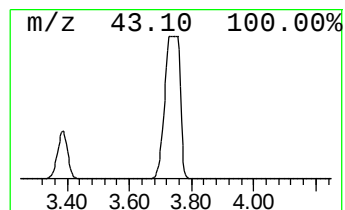
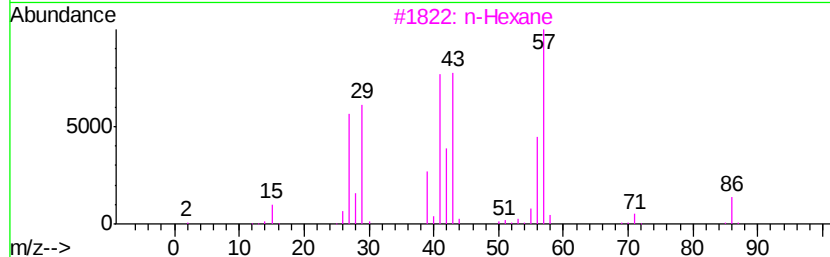
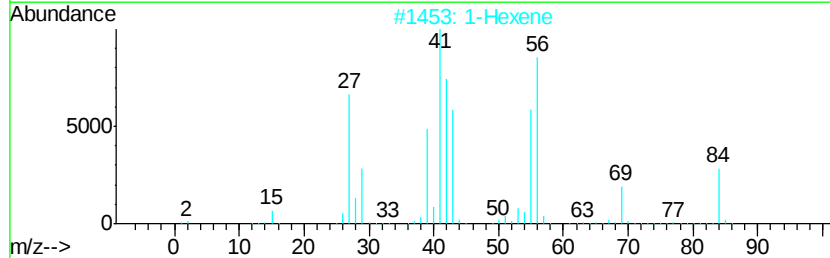
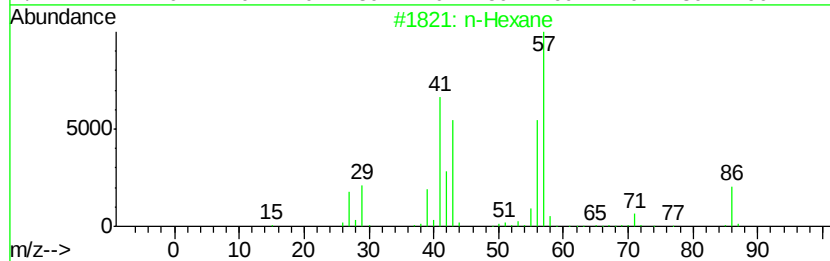
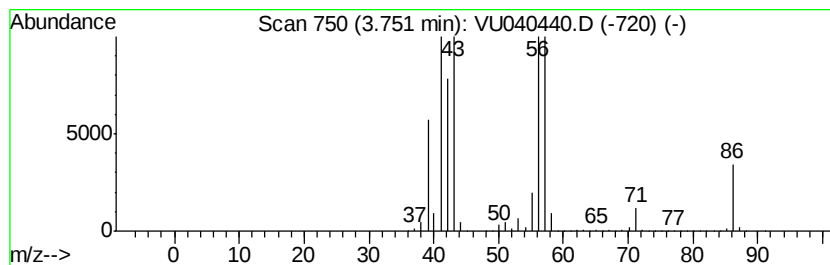
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.75	245.34 ug/L	204036000	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	64
2		1-Hexene	84	C6H12	000592-41-6	59
3		n-Hexane	86	C6H14	000110-54-3	50
4		n-Hexane	86	C6H14	000110-54-3	50
5		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	33



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Instrument :
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ClientSampleId :
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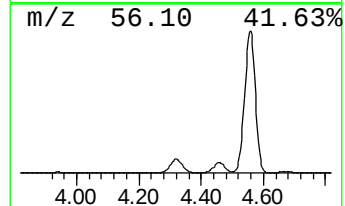
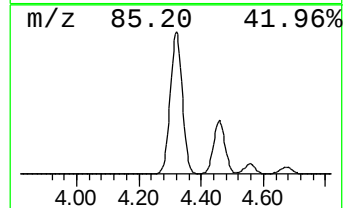
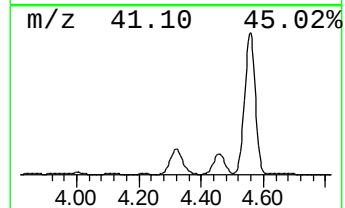
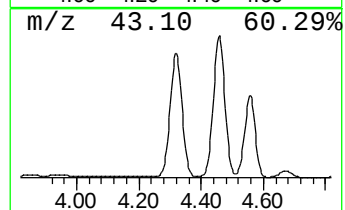
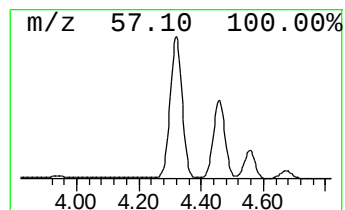
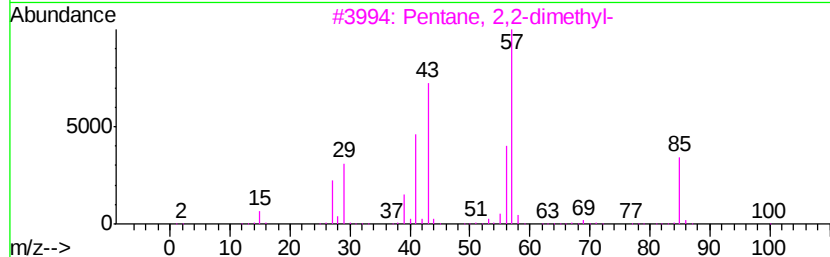
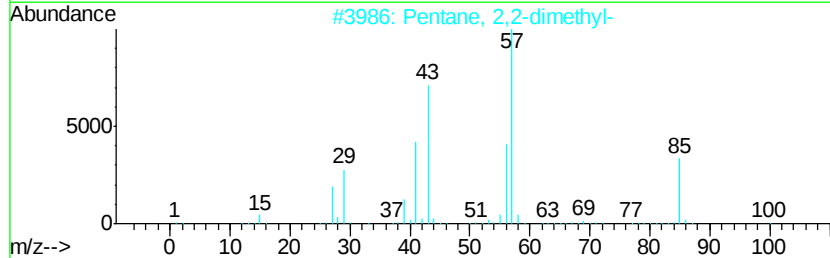
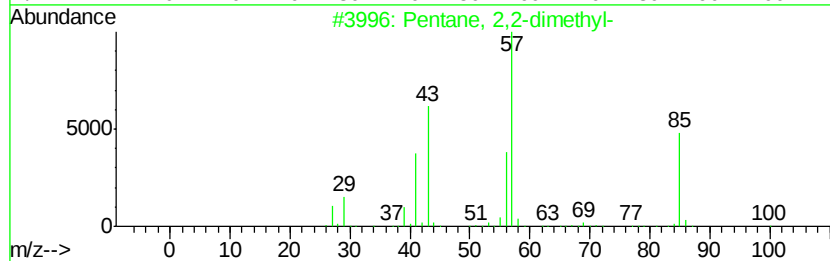
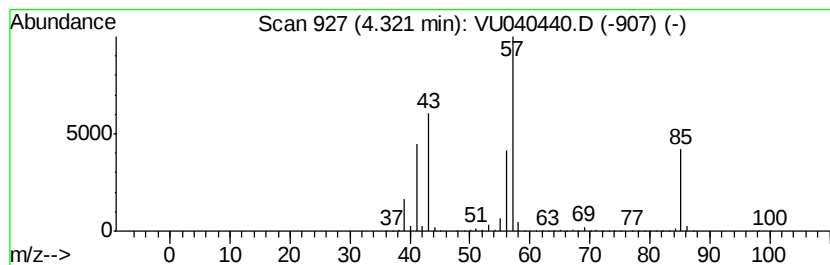
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.32	13.44 ug/L	11180700	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			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72
5			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64



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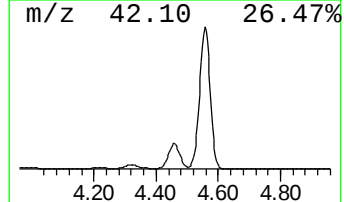
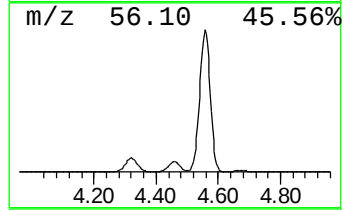
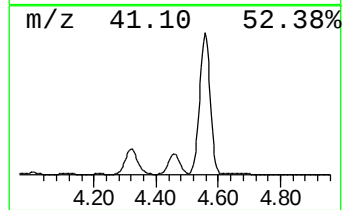
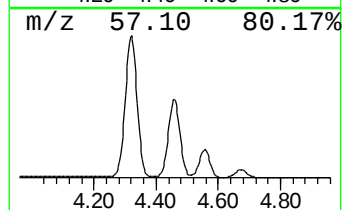
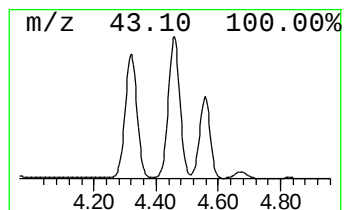
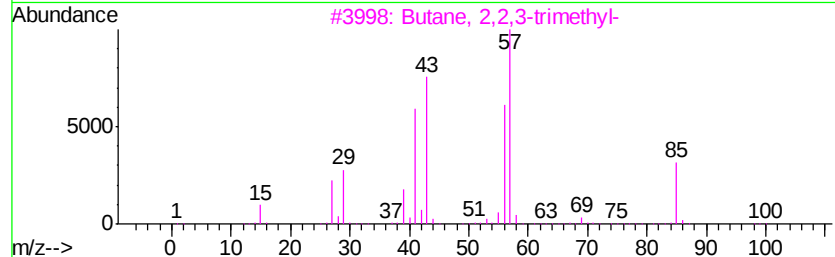
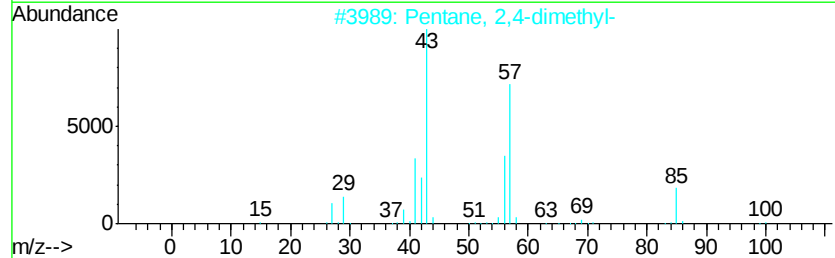
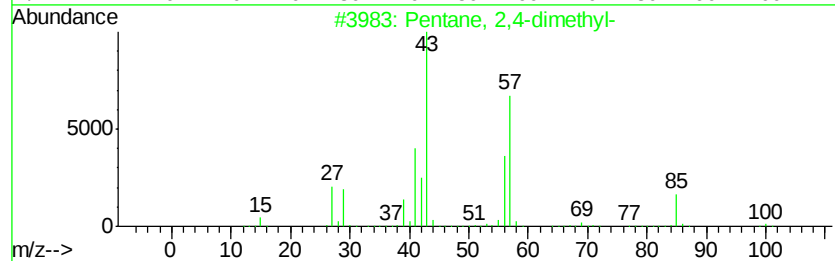
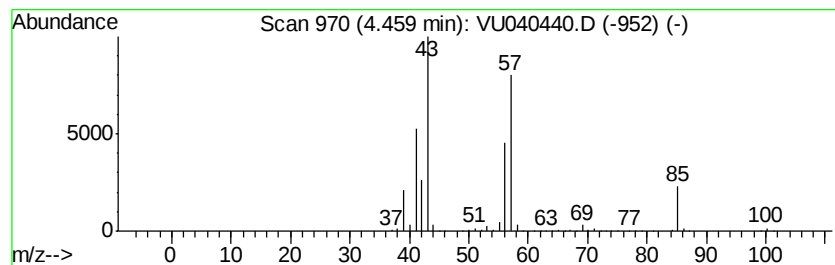
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	9.84 ug/L	8180430	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	83
3	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
5	n-Hexane	86	C6H14	000110-54-3	50



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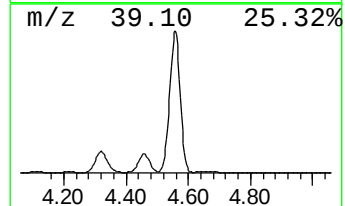
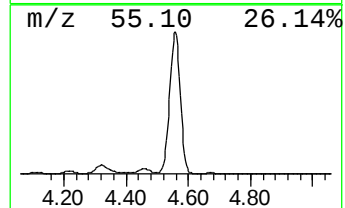
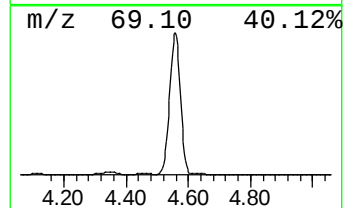
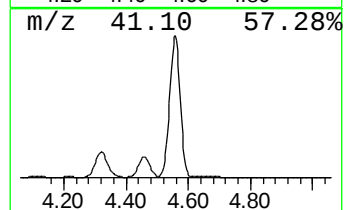
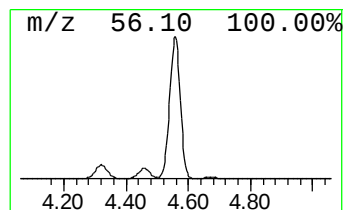
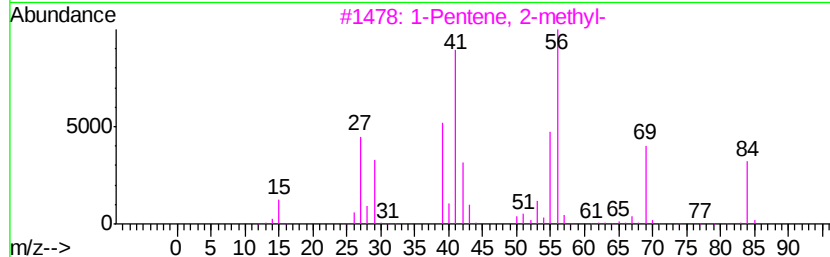
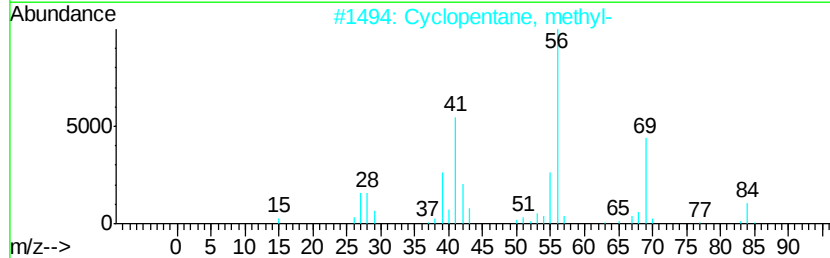
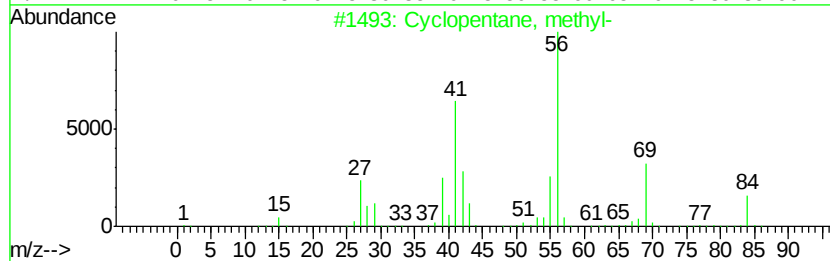
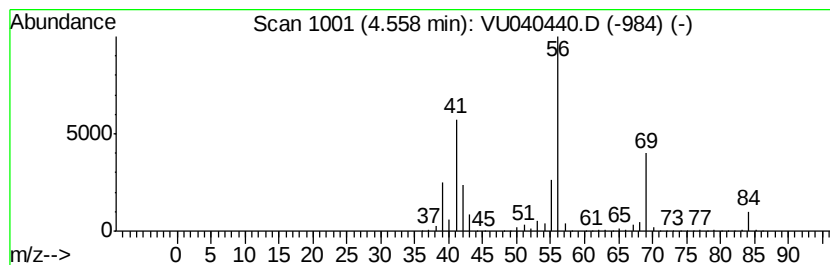
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 (DEL) Alkane: Cyclic4.56 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	50.38 ug/L	41897800	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	90
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	78
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	78
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

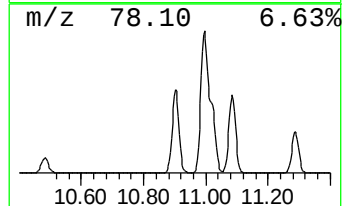
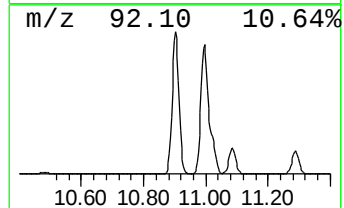
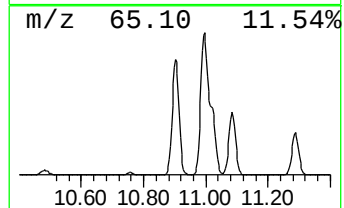
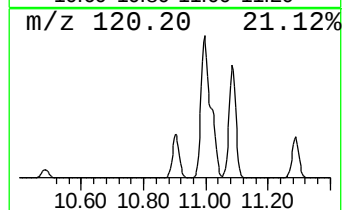
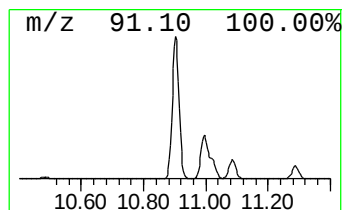
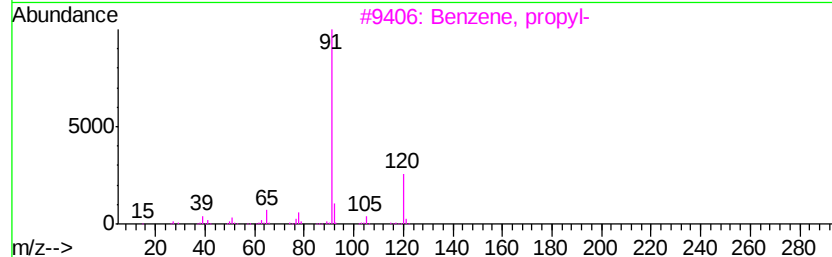
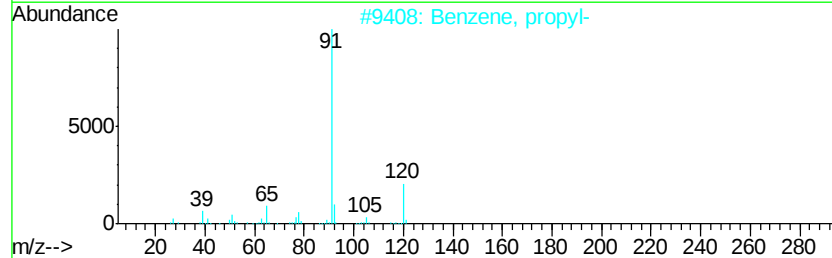
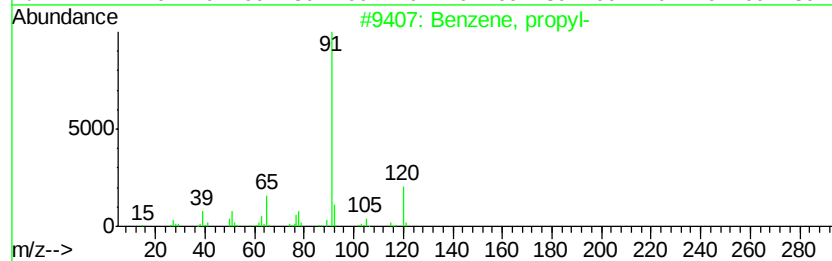
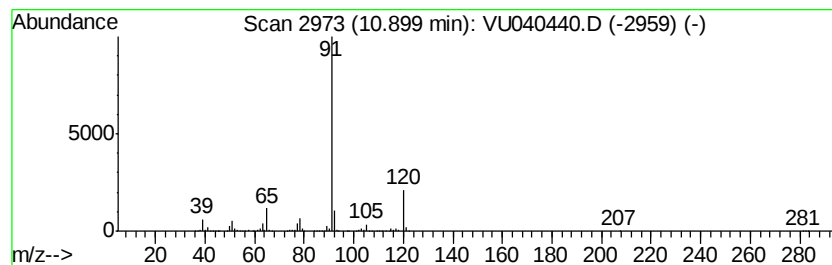
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 Benzene, propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	1.19 ug/L	2519480	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	90
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 : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

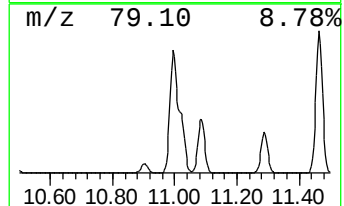
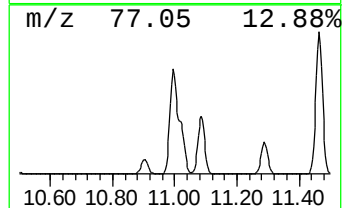
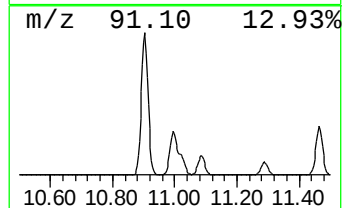
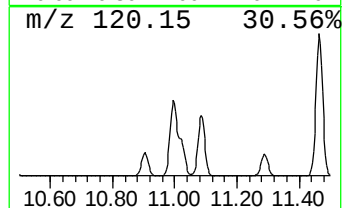
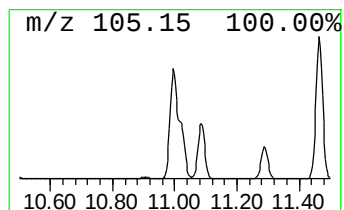
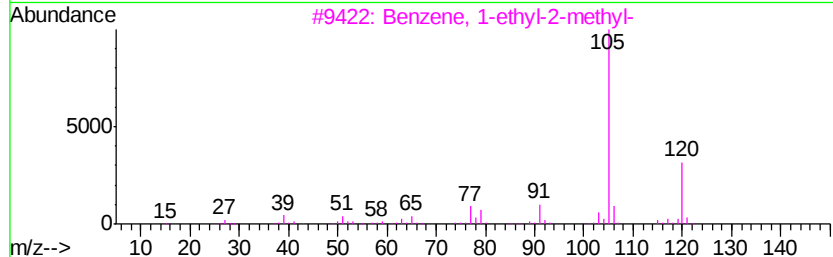
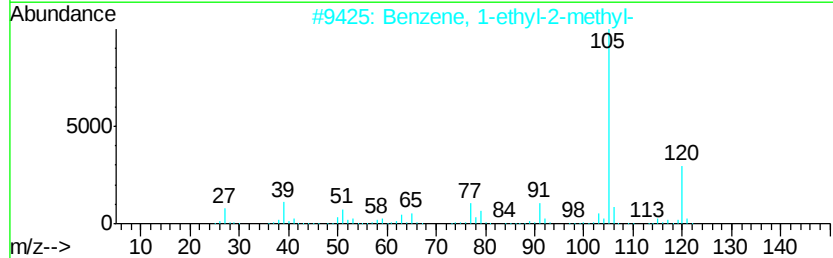
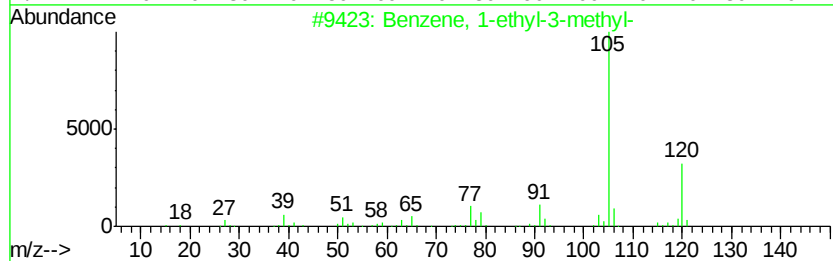
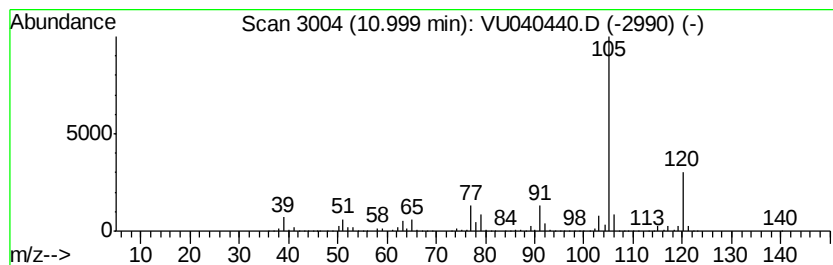
Instrument :
MSVOA_U
ClientSampleId :
BG3T0

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 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	5.00 ug/L	10608800	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	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:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

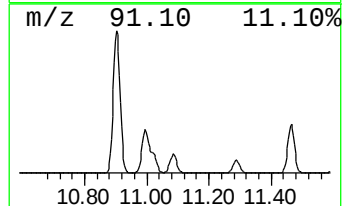
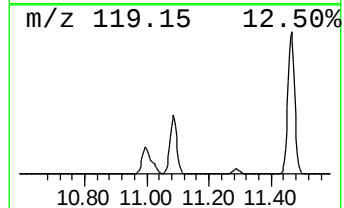
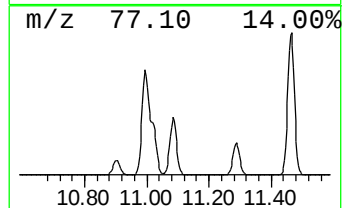
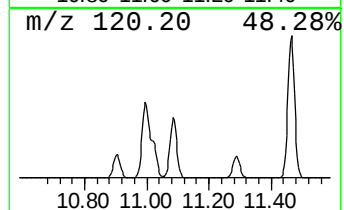
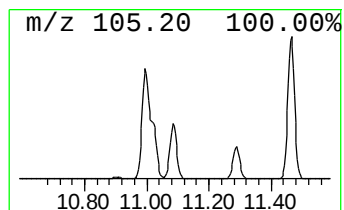
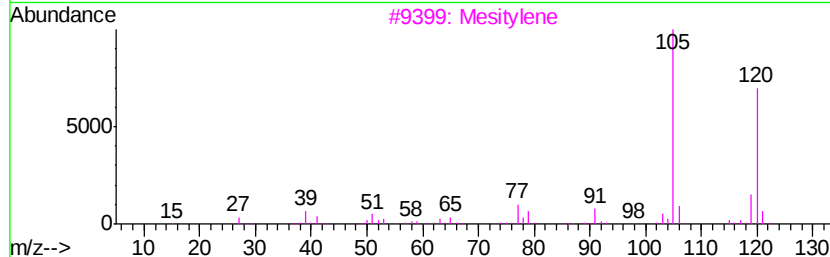
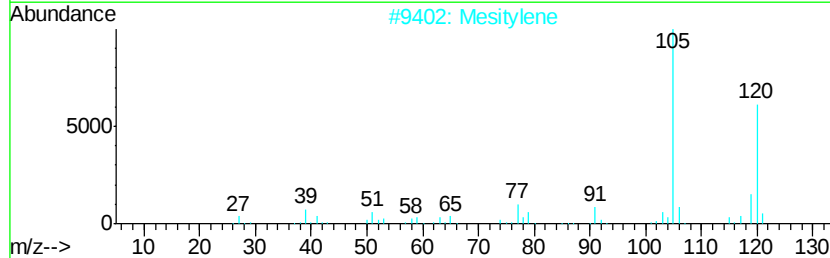
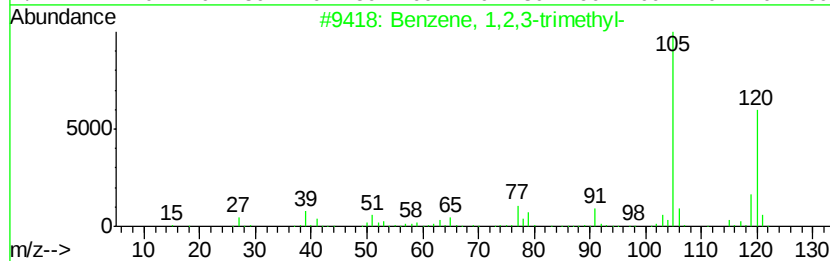
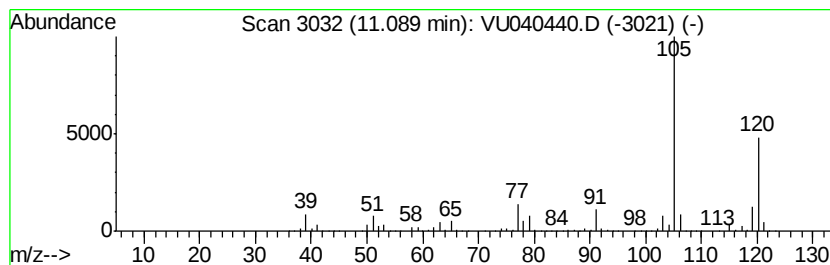
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	2.00 ug/L	4250940	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 : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

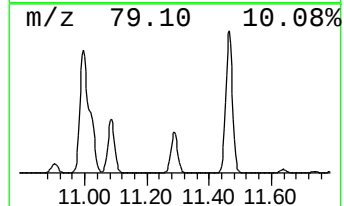
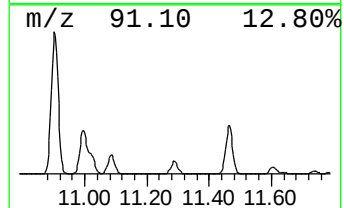
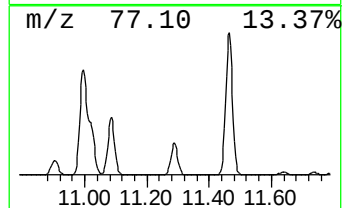
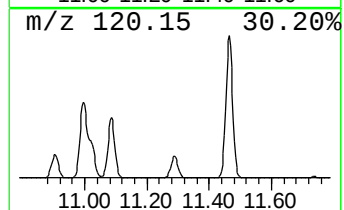
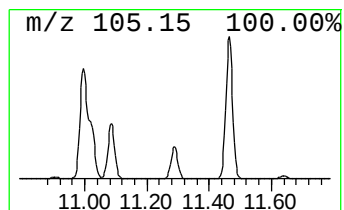
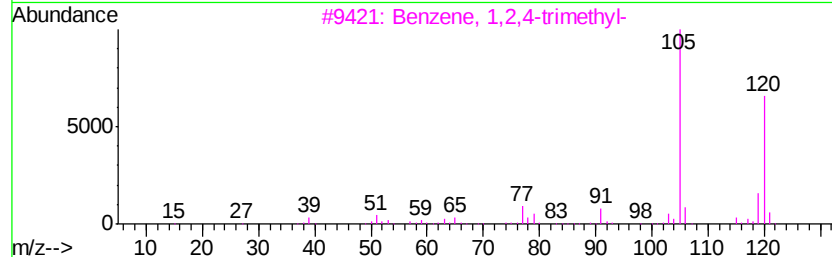
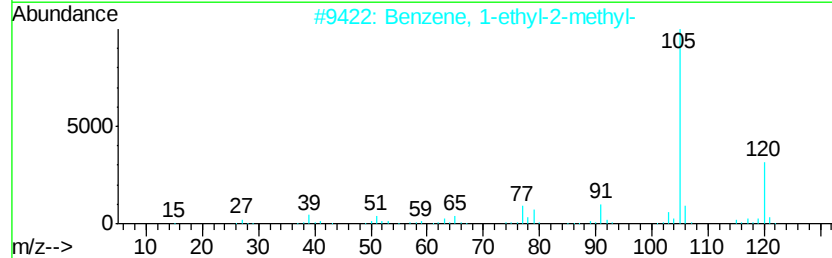
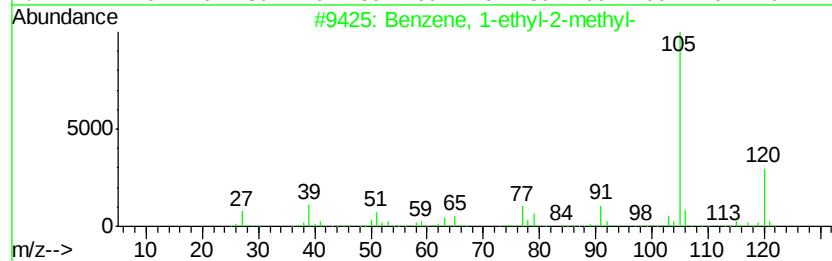
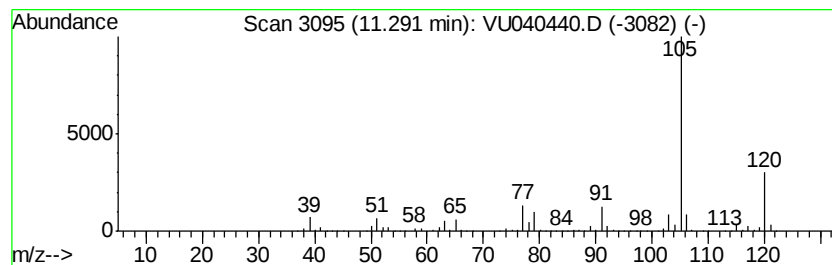
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	1.05 ug/L	2216410	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-2-methyl-	120	C9H12	000611-14-3	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

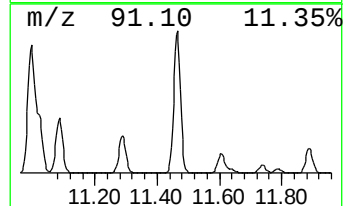
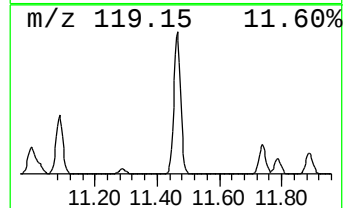
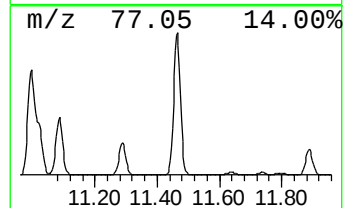
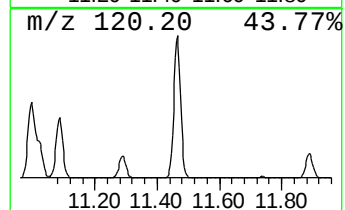
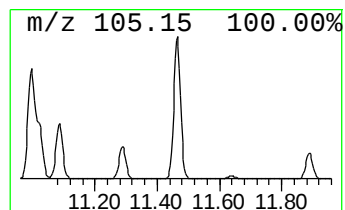
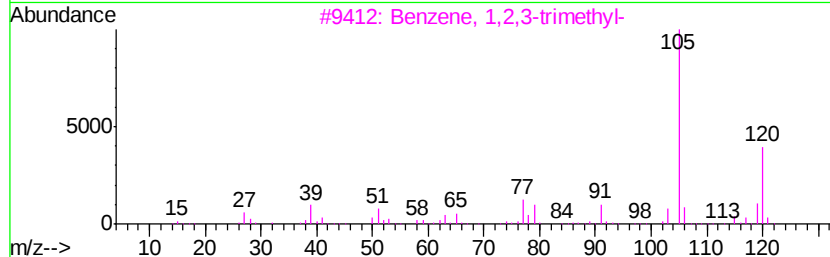
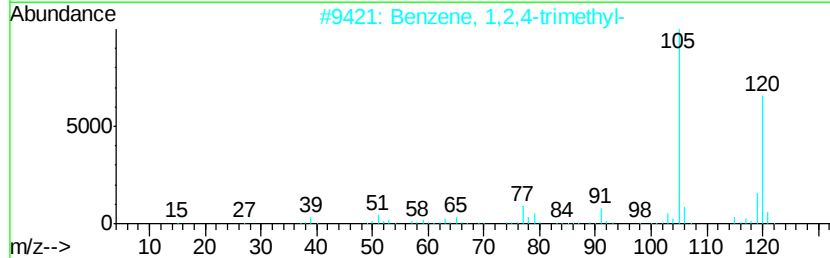
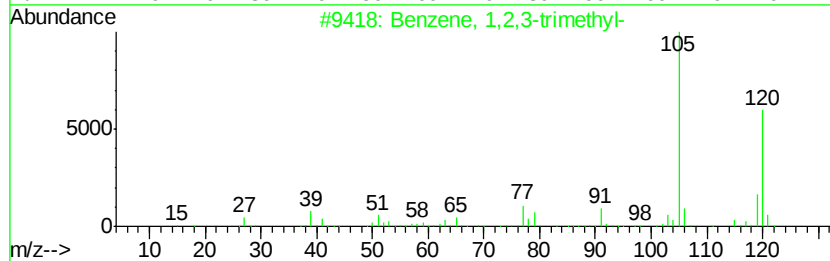
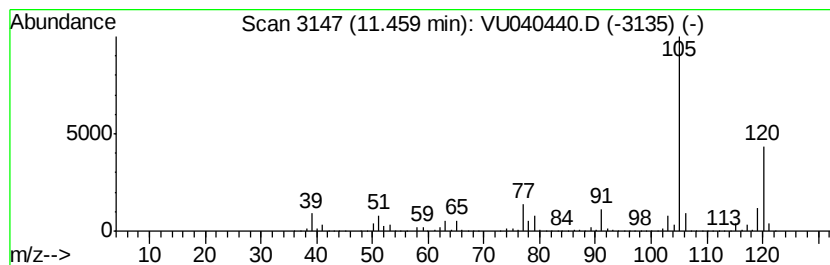
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Benzene, 1,2,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	5.00 ug/L	10604400	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	94
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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

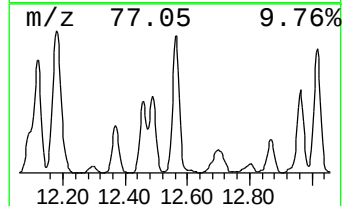
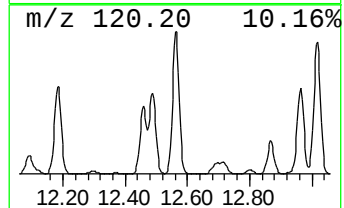
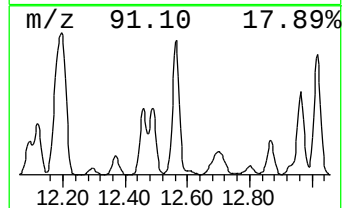
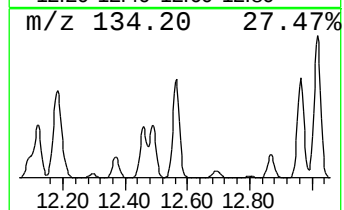
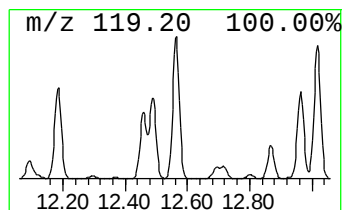
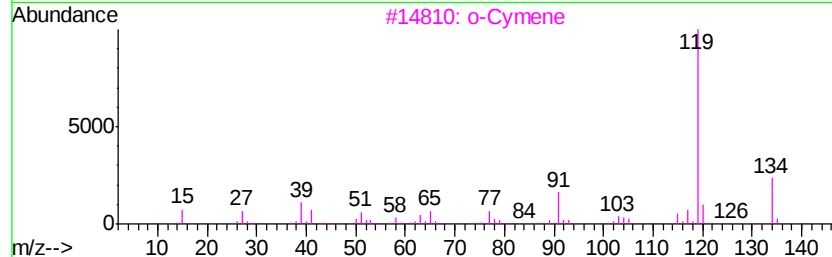
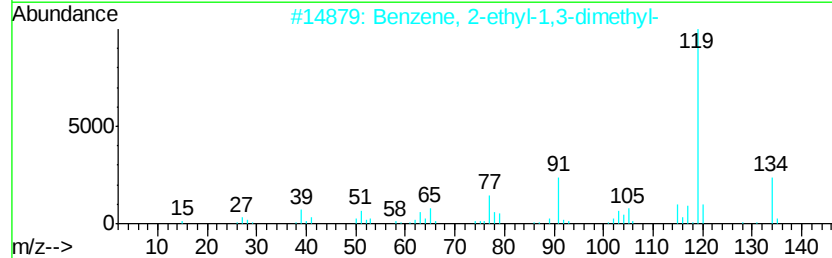
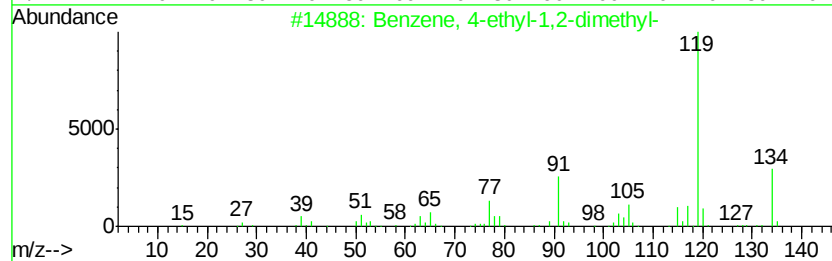
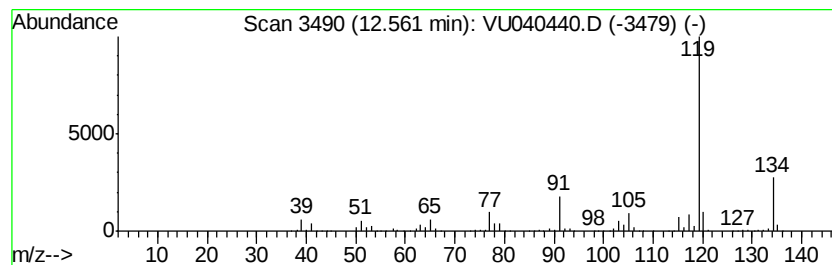
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	1.05 ug/L	2220530	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	97
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	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 : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

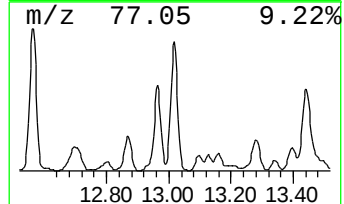
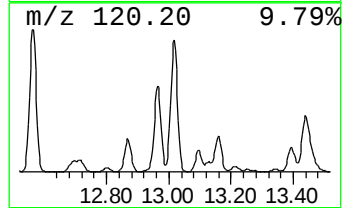
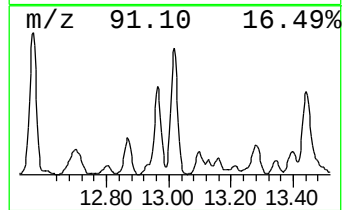
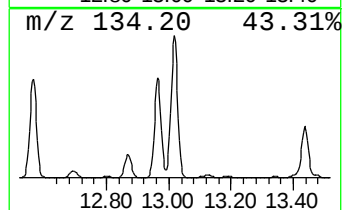
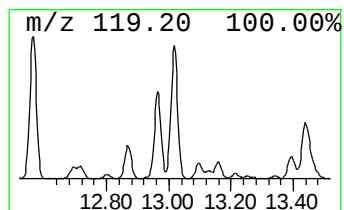
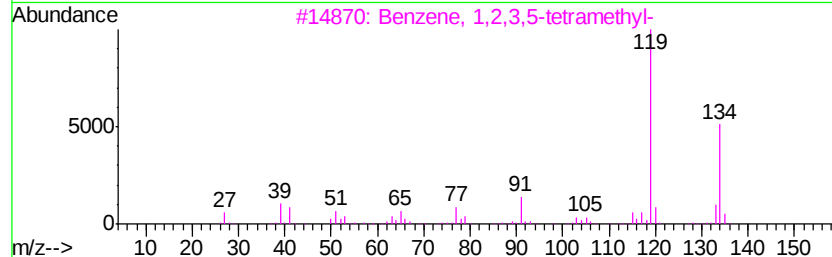
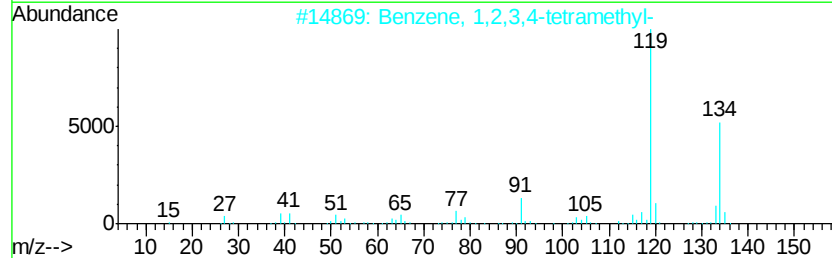
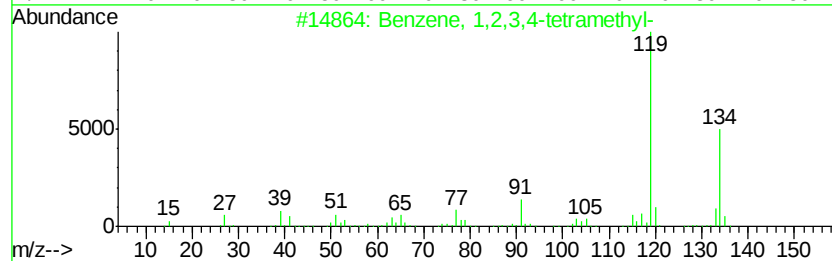
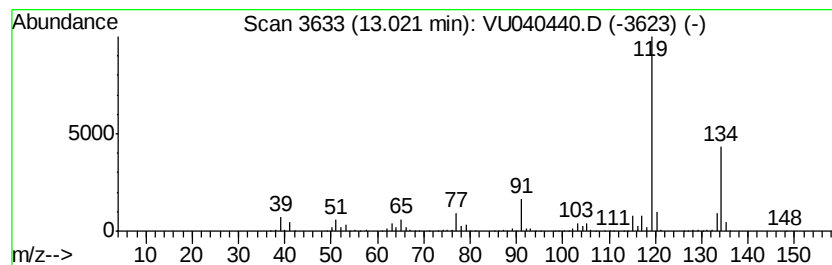
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	1.07 ug/L	2276920	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	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100120\
Data File : VU040440.D
Acq On : 01 Oct 2020 23:24
Operator : SY/MD
Sample : L4240-11
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0

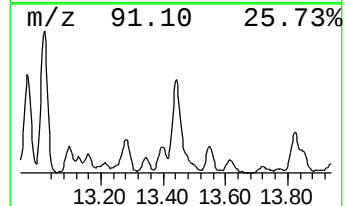
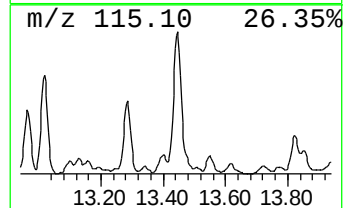
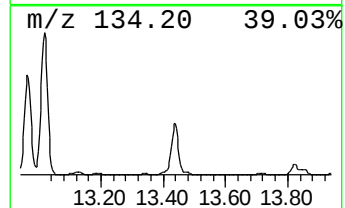
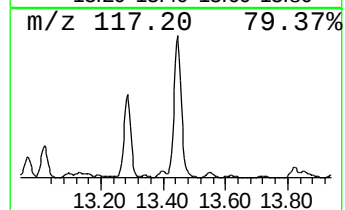
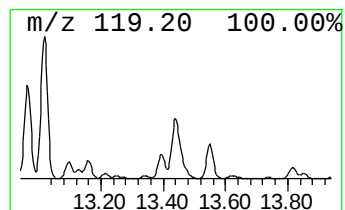
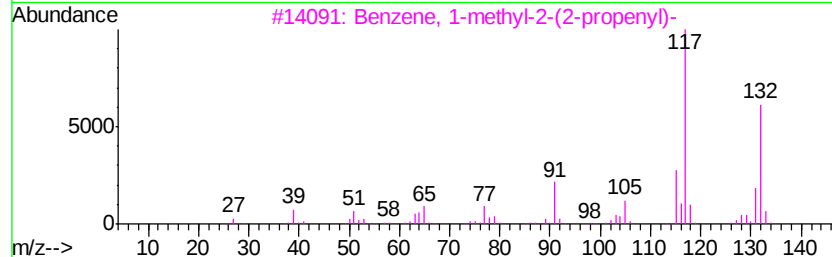
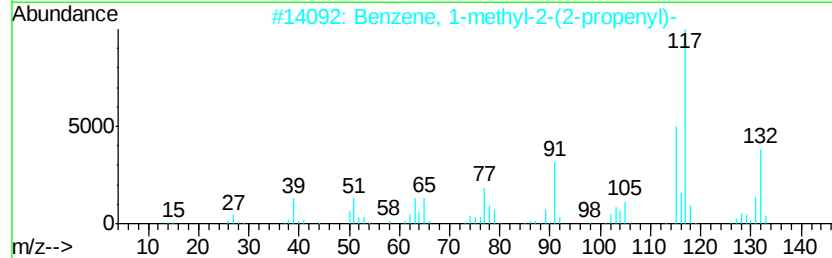
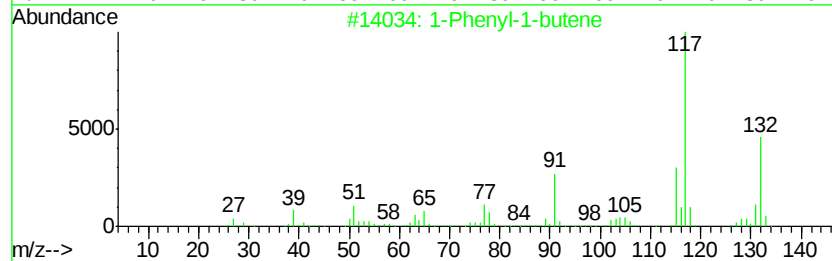
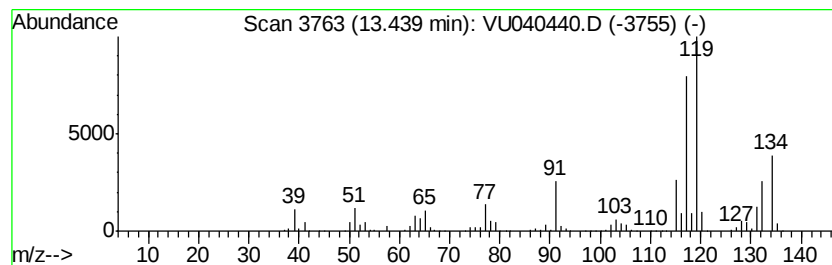
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 1-Phenyl-1-butene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	1.05 ug/L	2227880	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55
4			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	55
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100120\
 Data File : VU040440.D
 Acq On : 01 Oct 2020 23:24
 Operator : SY/MD
 Sample : L4240-11
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG3T0

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...	3.10	65.0	ug/L	54030100	1	6.26	4158210	5.0
(DEL) Alkane: Str...	3.38	98.8	ug/L	82120600	1	6.26	4158210	5.0
(DEL) Alkane: Str...	3.75	245.3	ug/L	204036000	1	6.26	4158210	5.0
(DEL) Alkane: Str...	4.32	13.4	ug/L	11180700	1	6.26	4158210	5.0
(DEL) Alkane: Str...	4.46	9.8	ug/L	8180430	1	6.26	4158210	5.0
(DEL) Alkane: Cyc...	4.56	50.4	ug/L	41897800	1	6.26	4158210	5.0
Benzene, propyl-	10.90	1.2	ug/L	2519480	3	11.81	10604400	5.0
Benzene, 1-ethyl-...	11.00	5.0	ug/L	10608800	3	11.81	10604400	5.0
Benzene, 1,2,3-tr...	11.09	2.0	ug/L	4250940	3	11.81	10604400	5.0
Benzene, 1-ethyl-...	11.29	1.1	ug/L	2216410	3	11.81	10604400	5.0
Benzene, 1,2,4-tr...	11.46	5.0	ug/L	10604400	3	11.81	10604400	5.0
Benzene, 4-ethyl-...	12.56	1.1	ug/L	2220530	3	11.81	10604400	5.0
Benzene, 1,2,3,4-...	13.02	1.1	ug/L	2276920	3	11.81	10604400	5.0
1-Phenyl-1-butene	13.44	1.1	ug/L	2227880	3	11.81	10604400	5.0