

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100720\
 Data File : VX018821.D
 Acq On : 07 Oct 2020 19:22
 Operator : JC/SP
 Sample : L4240-11DL 4X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OK-01-100620

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

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.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	2.825	274	285	287	rBV	1267765	2409994	2.65%	1.030%
2	2.873	287	293	319	rVB	10924062	22516363	24.73%	9.622%
3	3.154	327	339	363	rBV	17914955	41273609	45.33%	17.637%
4	3.520	385	399	414	rBV	38238322	91048573	100.00%	38.906%
5	4.117	486	497	512	rBV	1162189	3716526	4.08%	1.588%
6	4.282	513	524	530	rVV	803815	2477306	2.72%	1.059%
7	4.379	530	540	554	rVV	11535254	33431574	36.72%	14.286%
8	5.550	717	732	747	rBV	996617	3103456	3.41%	1.326%
9	11.341	1677	1682	1689	rBV	1619853	1937141	2.13%	0.828%
10	11.421	1689	1695	1702	rVV	4918056	8564995	9.41%	3.660%
11	11.494	1702	1707	1715	rVB	2523531	3210712	3.53%	1.372%
12	11.652	1728	1733	1741	rVB	1499330	1761924	1.94%	0.753%
13	11.793	1751	1756	1761	rBV	7398836	8468950	9.30%	3.619%
14	12.128	1806	1811	1816	rVB	1162315	1403926	1.54%	0.600%
15	12.305	1830	1840	1844	rBV2	593387	1066141	1.17%	0.456%
16	12.353	1844	1848	1856	rVB2	1219899	1847582	2.03%	0.789%
17	12.652	1892	1897	1901	rBV	1107506	1294549	1.42%	0.553%
18	12.963	1940	1948	1951	rBV	732086	925512	1.02%	0.395%
19	13.000	1951	1954	1960	rVB	1082030	1269359	1.39%	0.542%
20	13.329	2004	2008	2017	rVB2	843913	1260583	1.38%	0.539%
21	13.628	2051	2057	2064	rVB2	592405	1032103	1.13%	0.441%

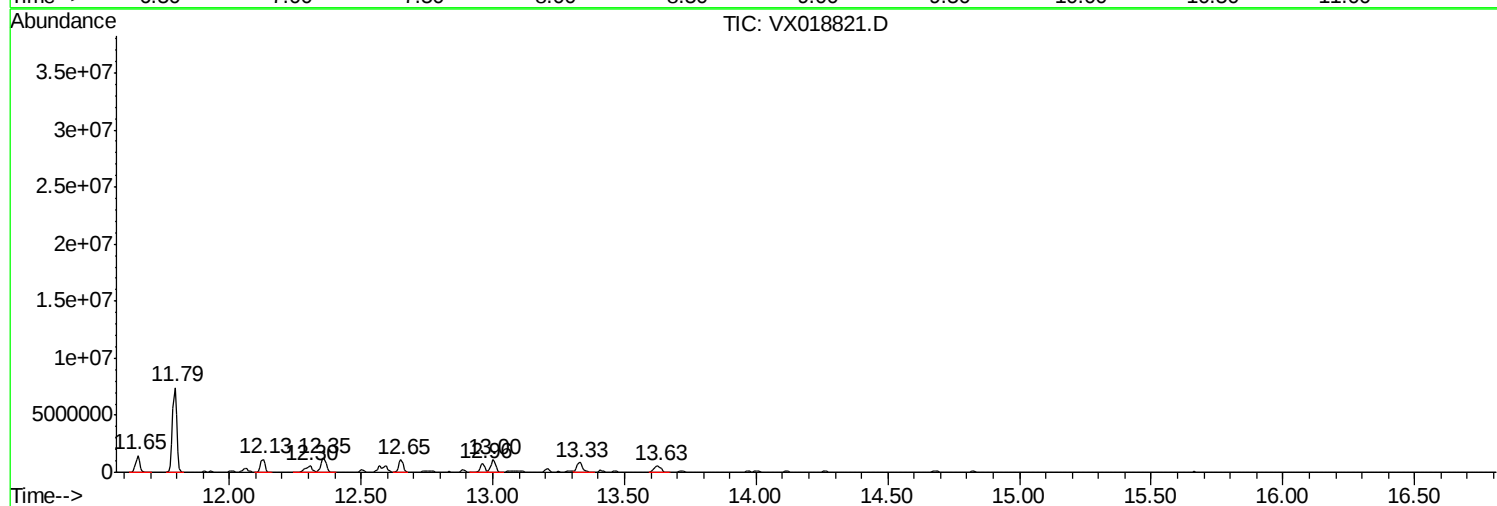
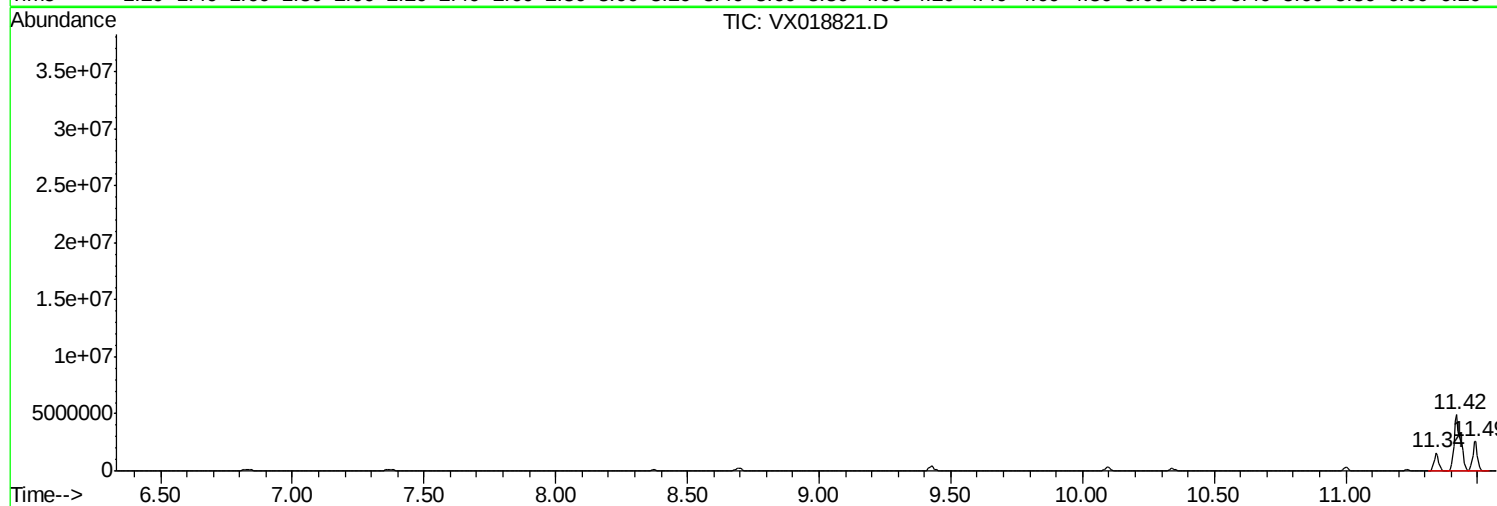
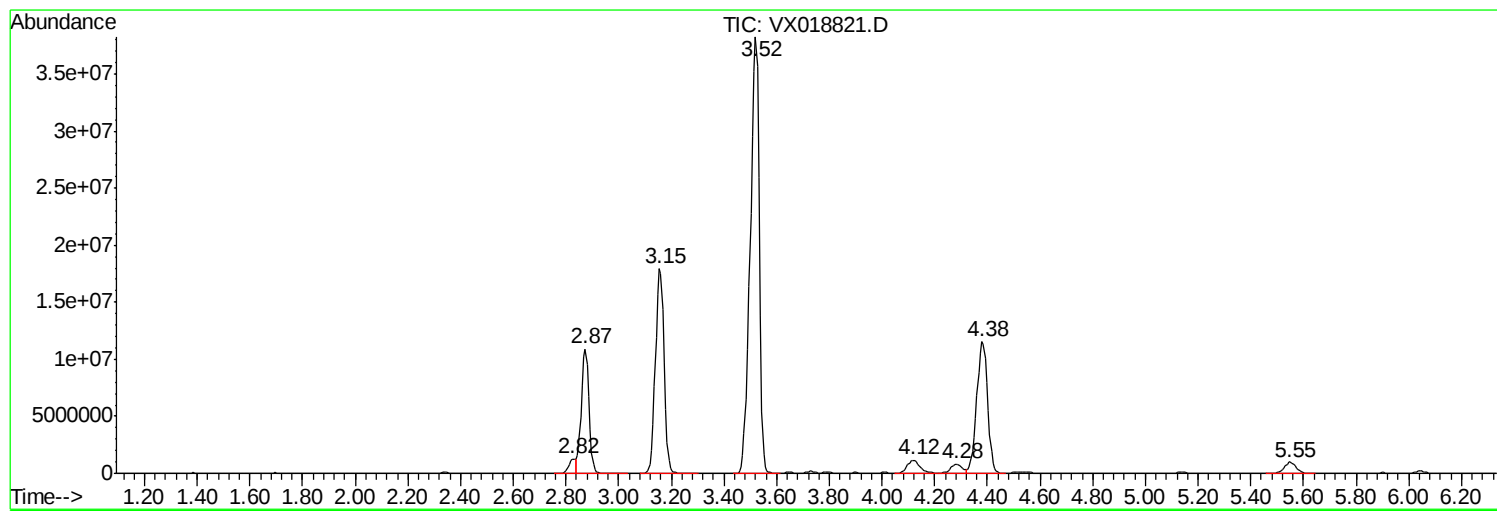
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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



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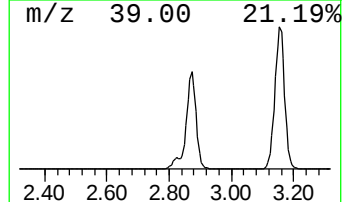
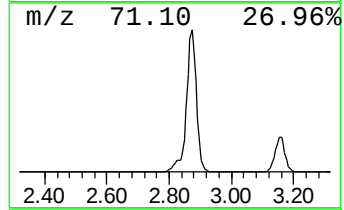
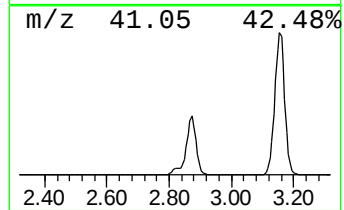
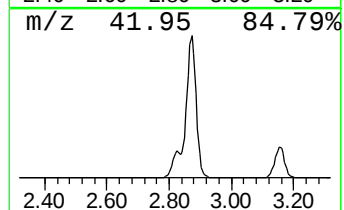
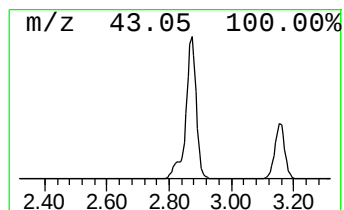
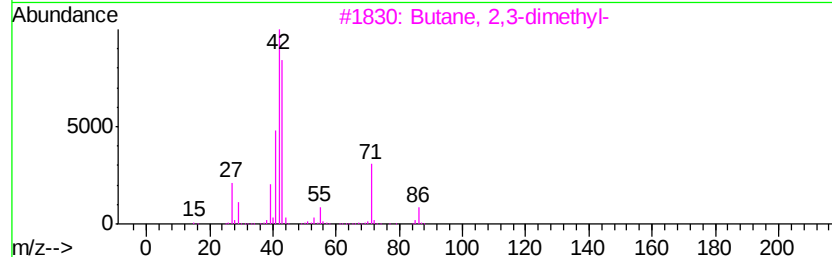
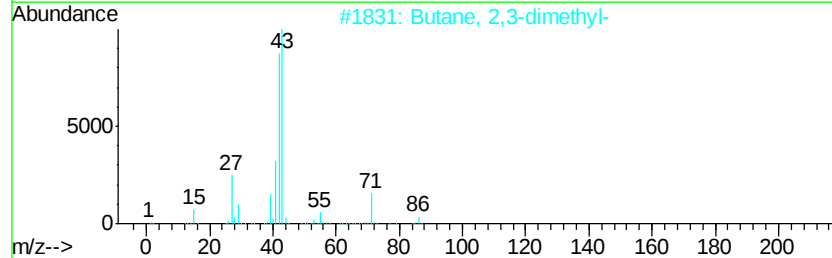
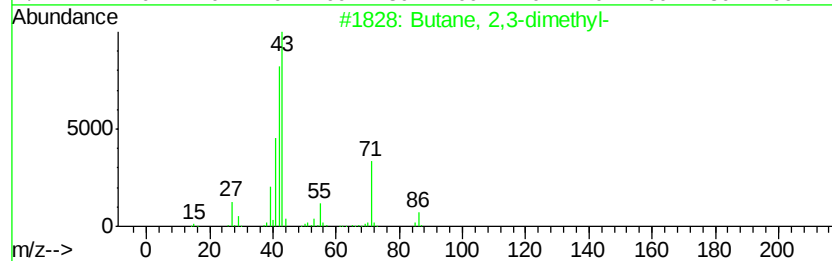
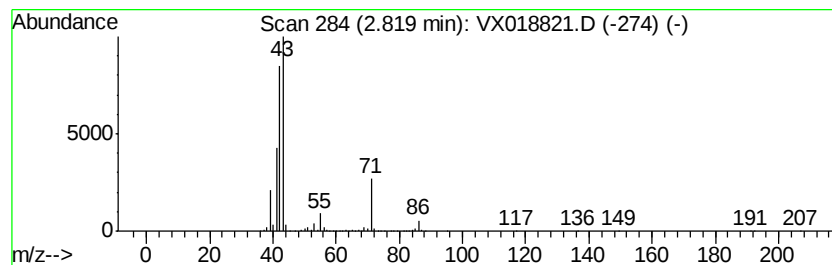
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.82	3.88 ug/L	2409990	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	91
2		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
4		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
5		Pentane, 2-methyl-	86	C6H14	000107-83-5	50



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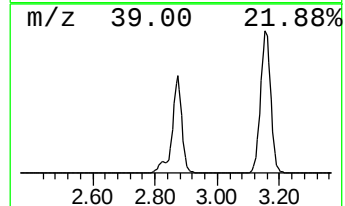
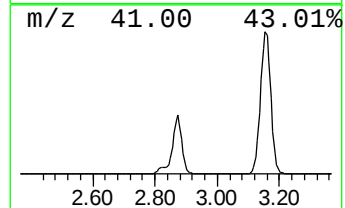
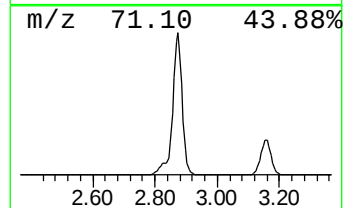
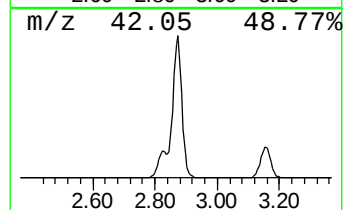
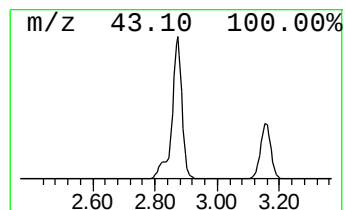
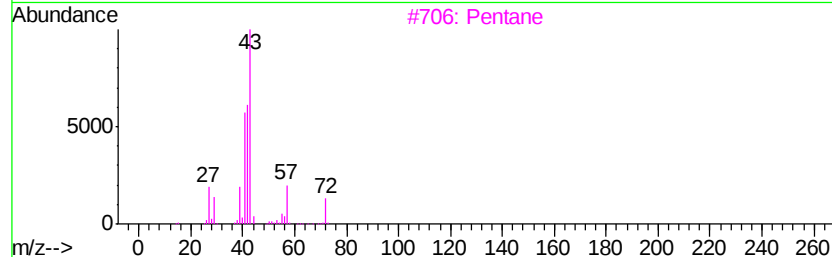
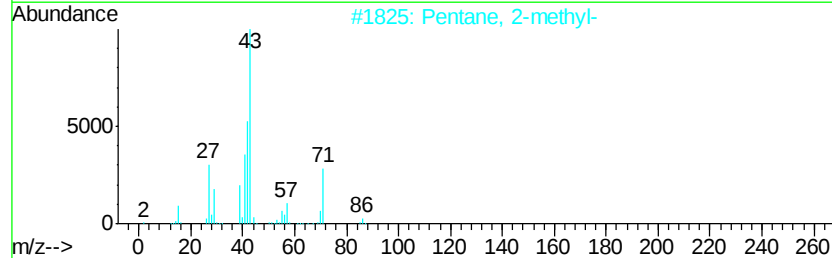
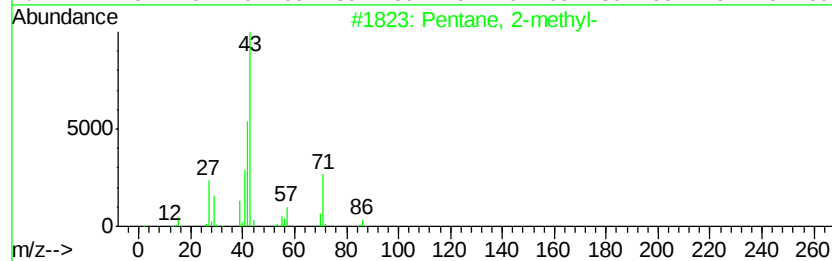
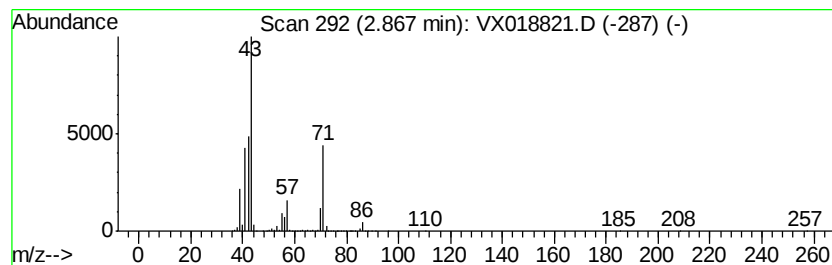
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	36.28 ug/L	22516400	1,4-Difluorobenzene	6.83

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		Pentane	72	C5H12	000109-66-0	49
4		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	49
5		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	42



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ClientSampleId :
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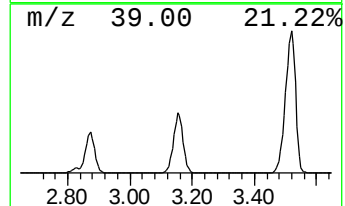
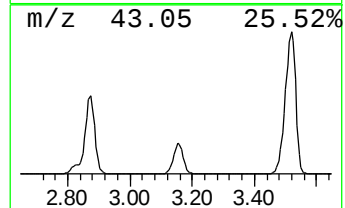
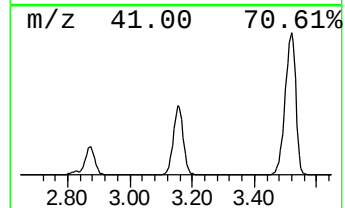
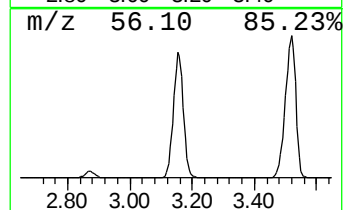
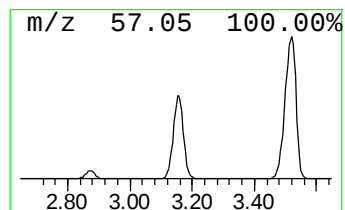
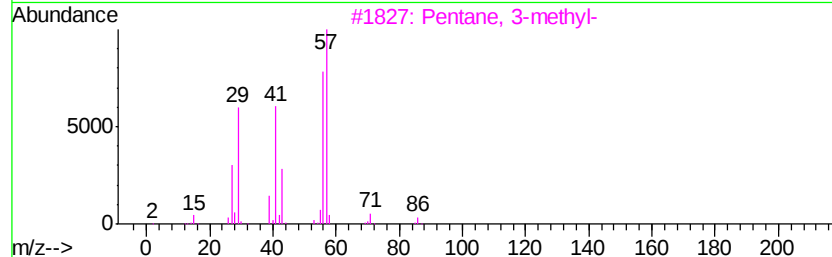
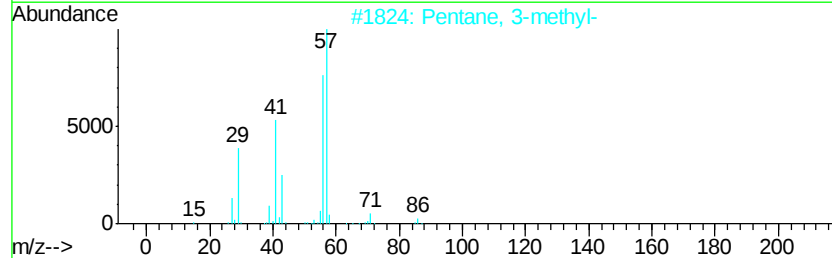
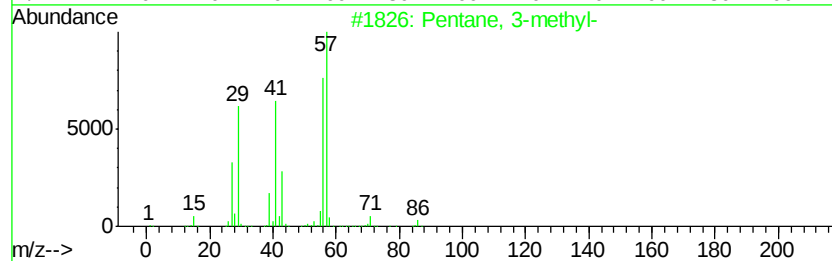
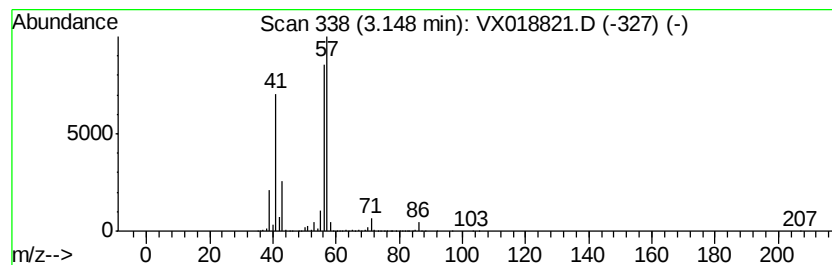
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	66.50 ug/L	41273600	1,4-Difluorobenzene	6.83

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	90
4		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
5		n-Hexane	86	C6H14	000110-54-3	40



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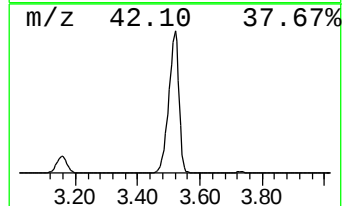
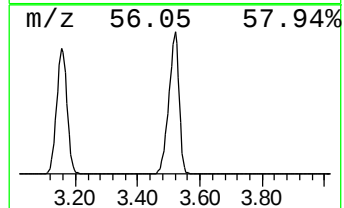
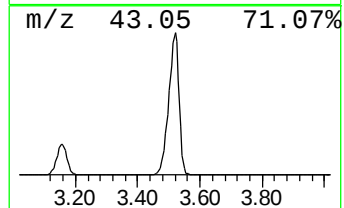
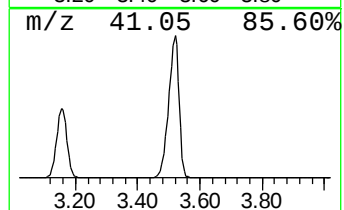
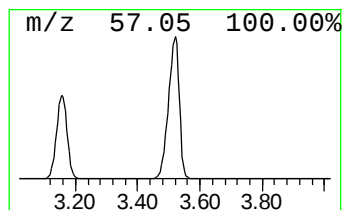
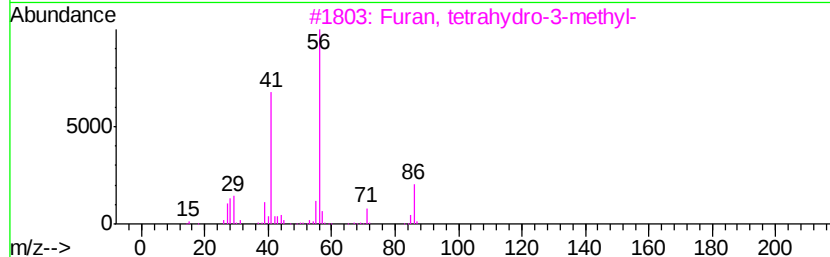
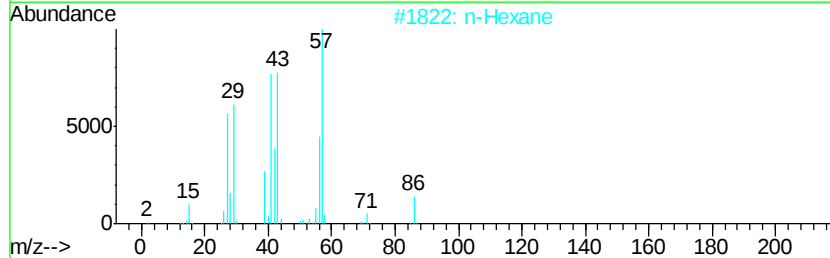
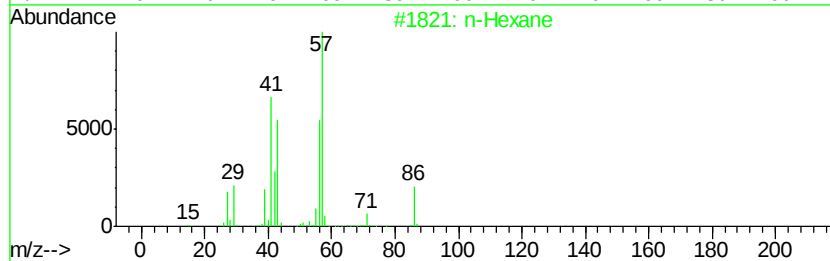
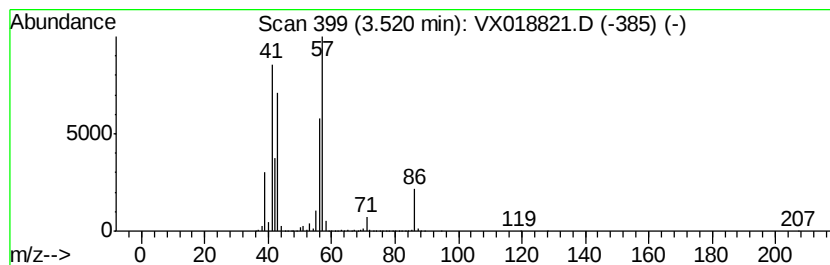
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	146.69 ug/L	91048600	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	94
2		n-Hexane	86	C6H14	000110-54-3	64
3		Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	53
4		n-Hexane	86	C6H14	000110-54-3	50
5		Butanal, 2-methyl-	86	C5H10O	000096-17-3	43



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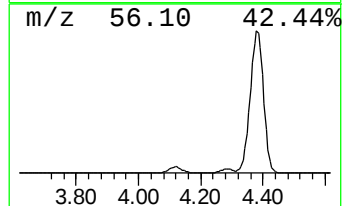
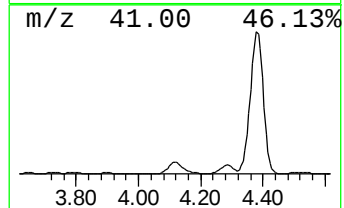
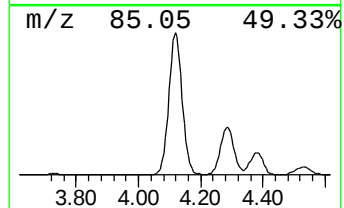
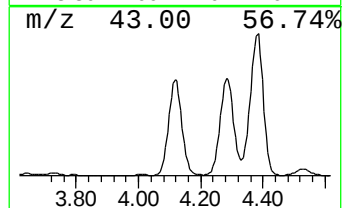
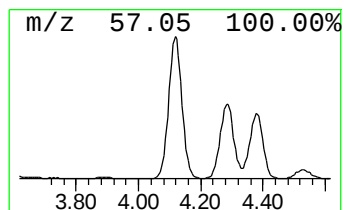
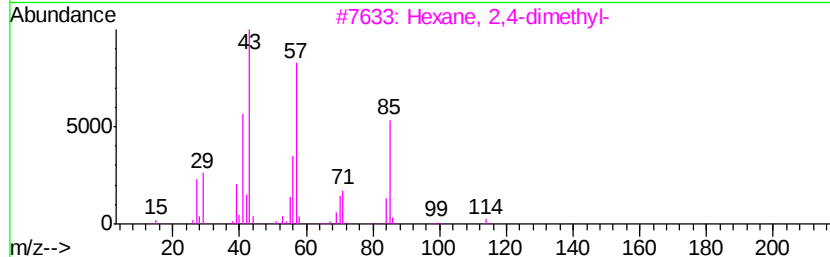
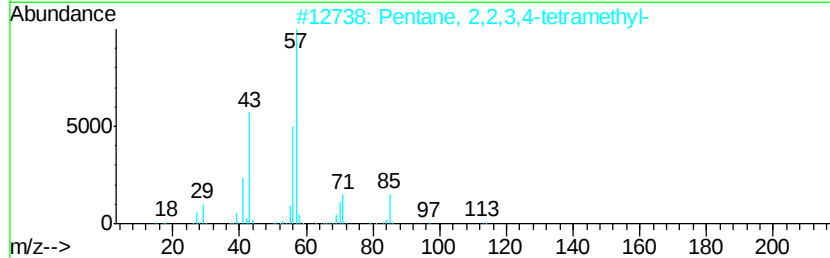
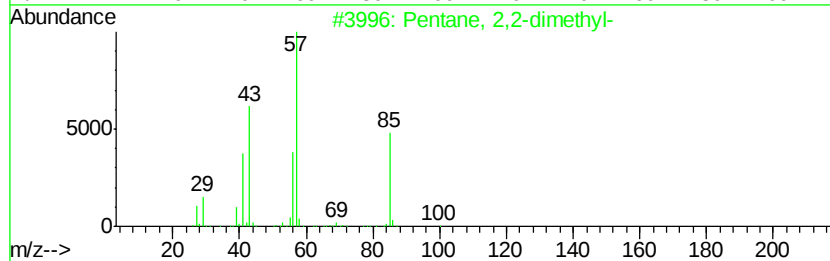
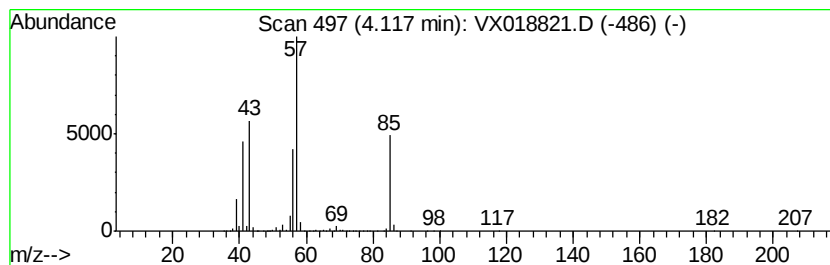
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.12	5.99 ug/L	3716530	1,4-Difluorobenzene	6.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	90	
2	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	78	
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72	
4	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64	
5	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64	



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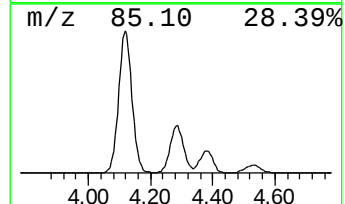
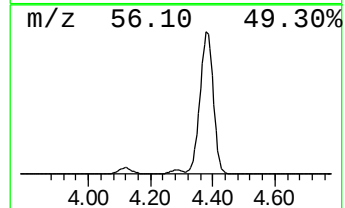
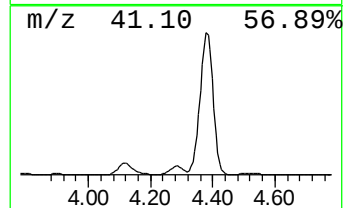
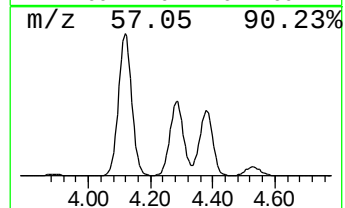
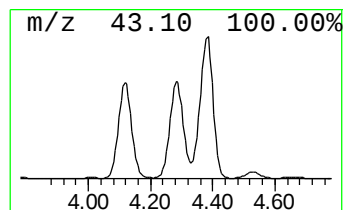
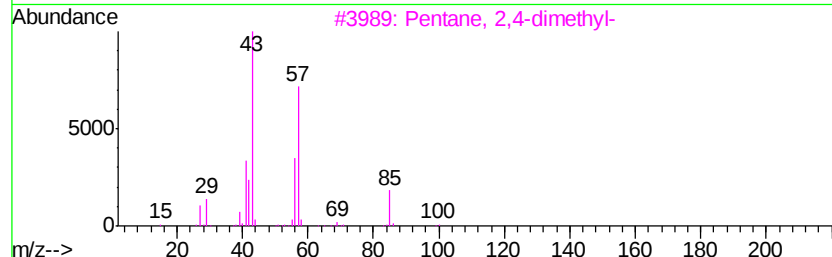
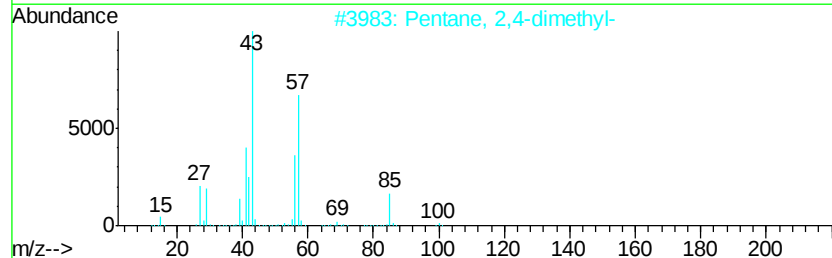
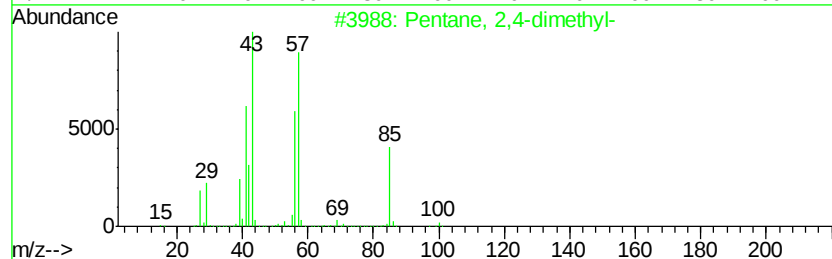
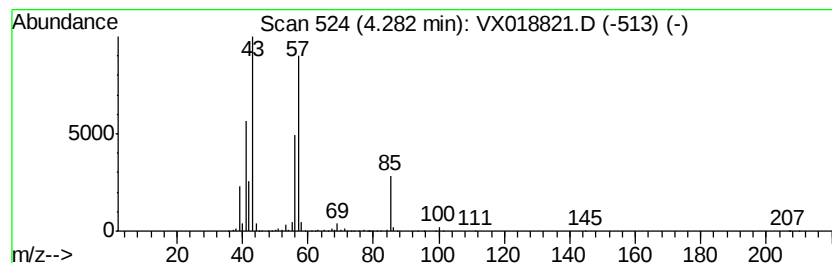
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	3.99 ug/L	2477310	1,4-Difluorobenzene	6.83

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	91
3		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	74
4		Hexane, 2-methyl-	100	C7H16	000591-76-4	64
5		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

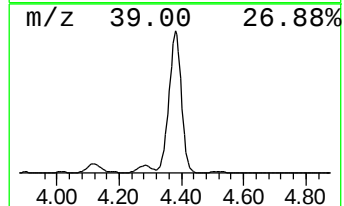
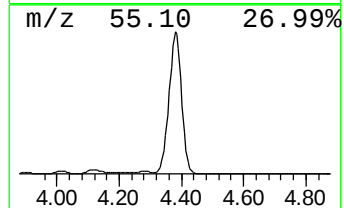
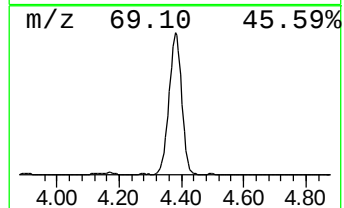
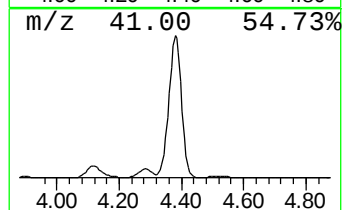
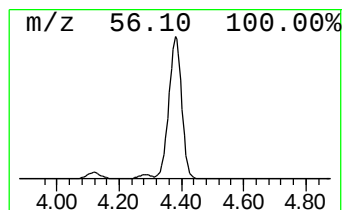
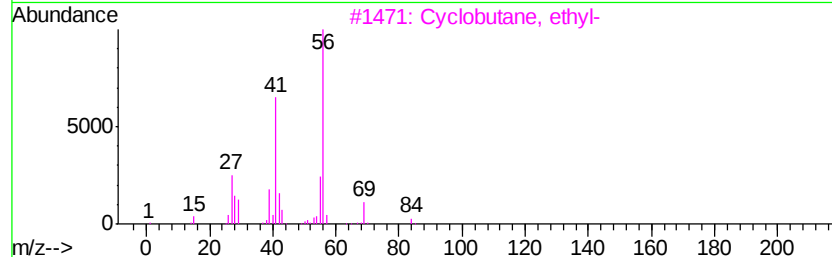
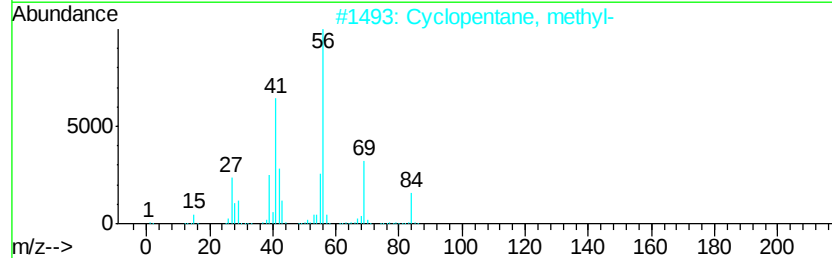
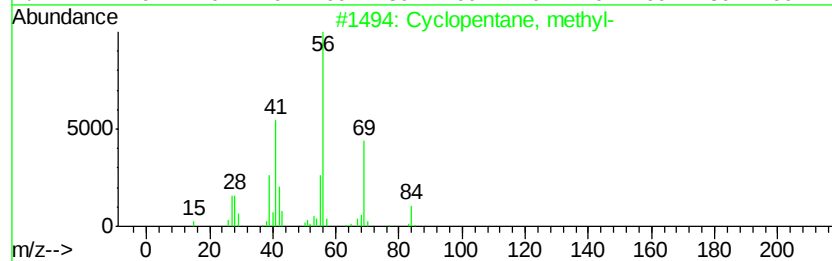
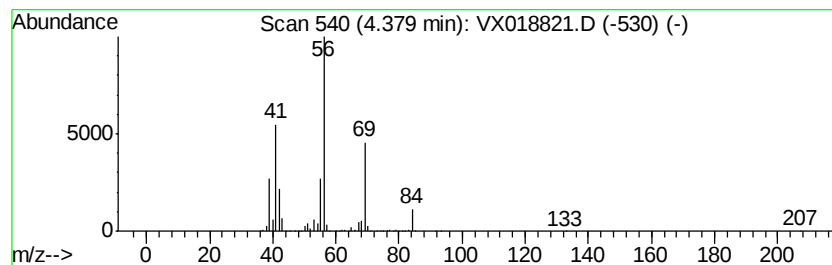
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 (DEL) Alkane: Cyclic4.38 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	53.86 ug/L	33431600	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	86
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4		Cyclohexane	84	C6H12	000110-82-7	64
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

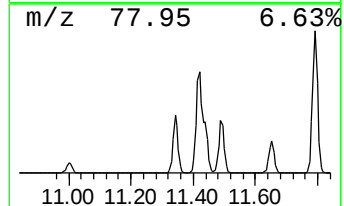
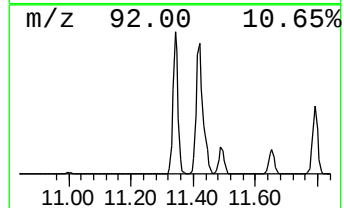
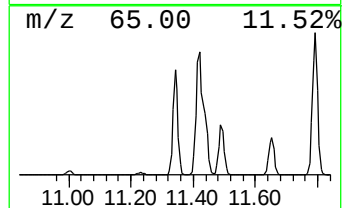
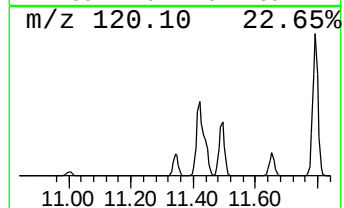
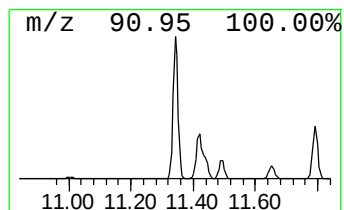
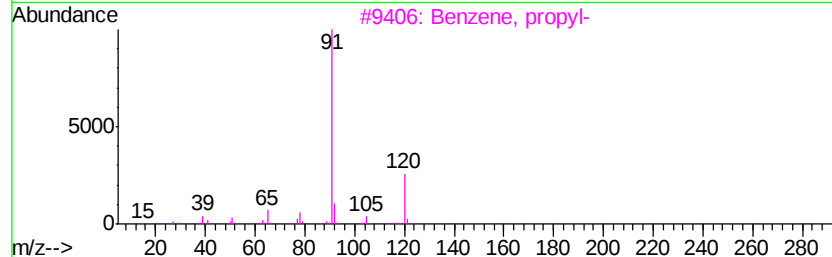
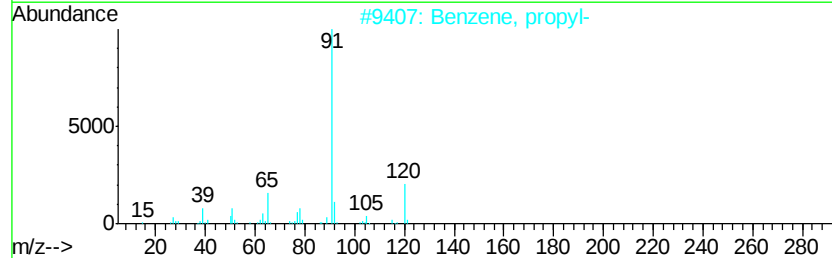
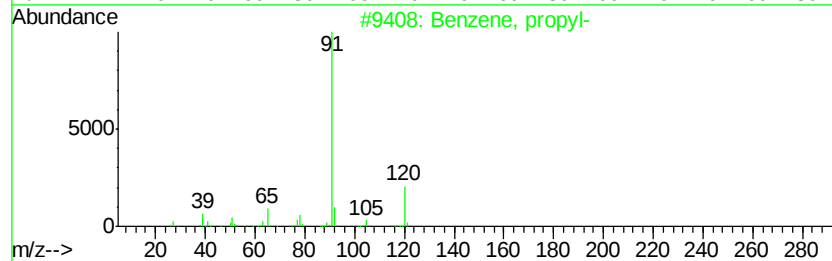
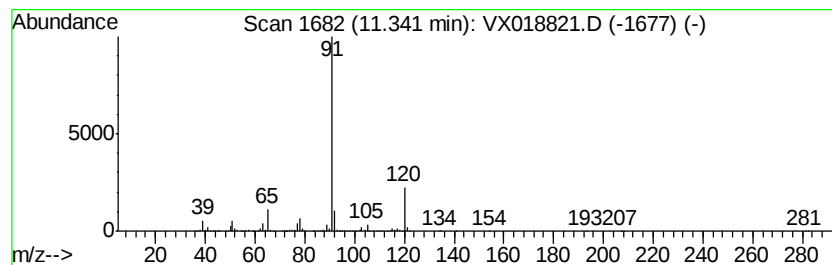
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 Benzene, propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	6.90 ug/L	1937140	1,4-Dichlorobenzene-d4	12.06

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	87
4		Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	78
5		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

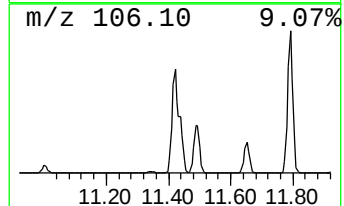
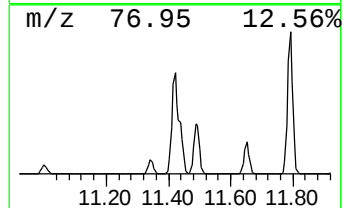
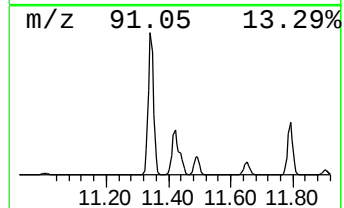
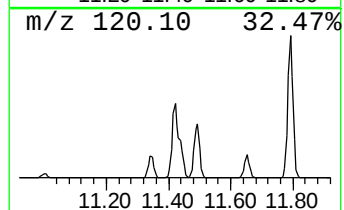
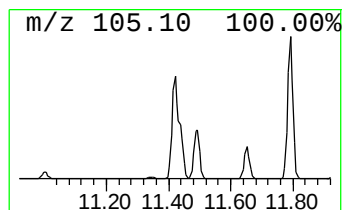
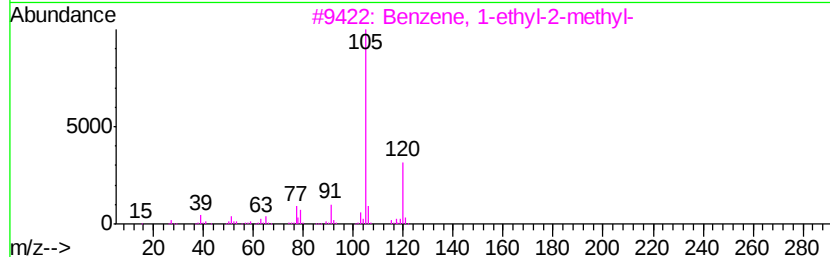
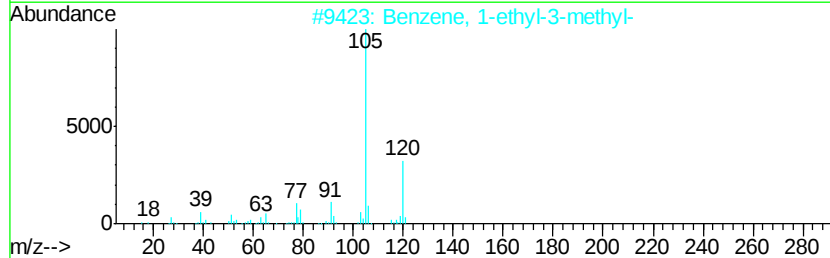
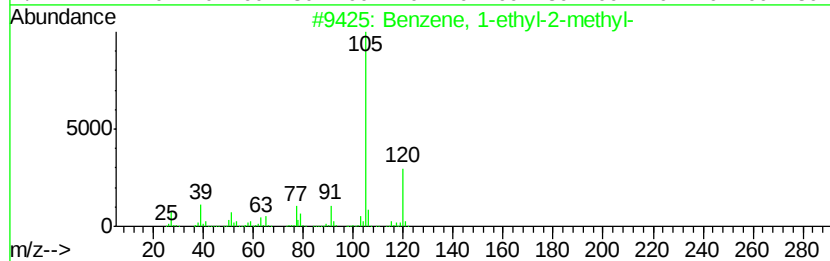
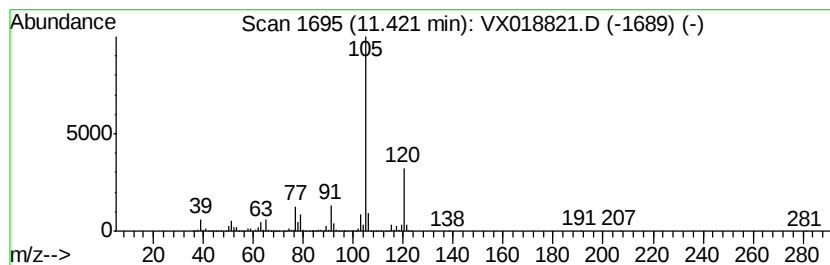
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	30.50 ug/L	8565000	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	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	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

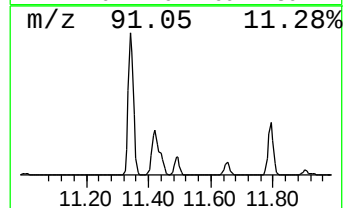
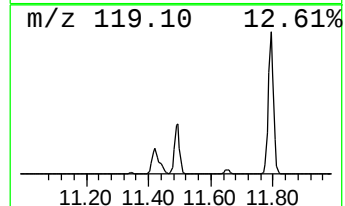
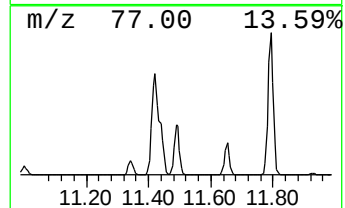
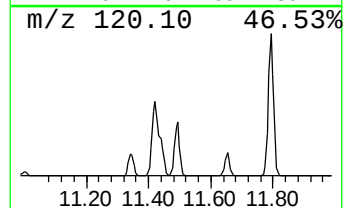
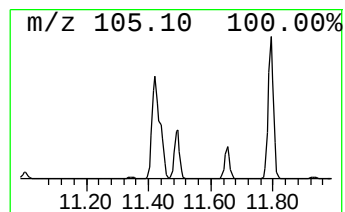
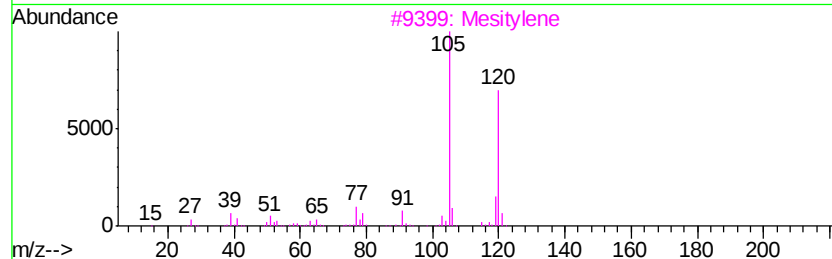
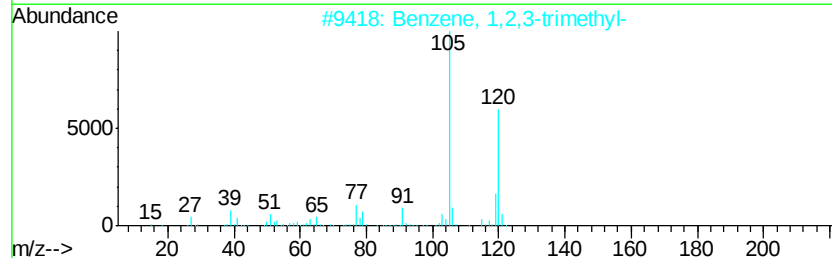
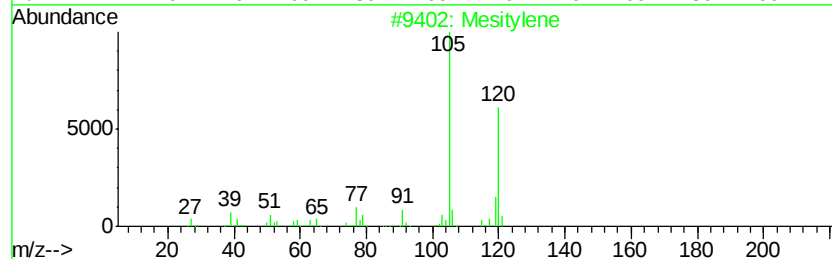
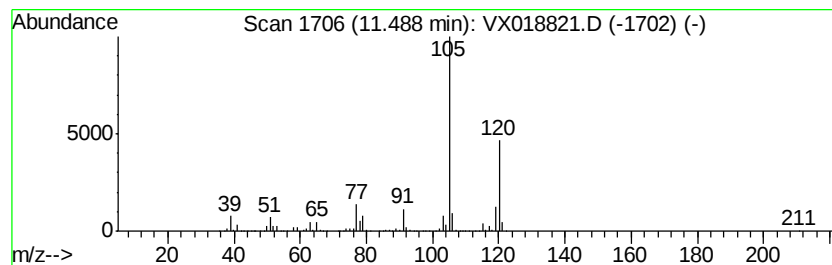
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	11.43 ug/L	3210710	1,4-Dichlorobenzene-d4	12.06

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

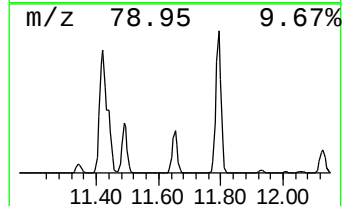
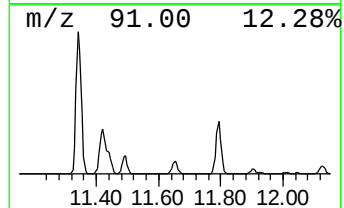
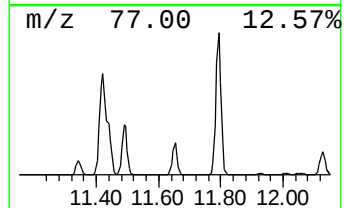
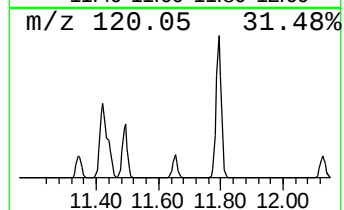
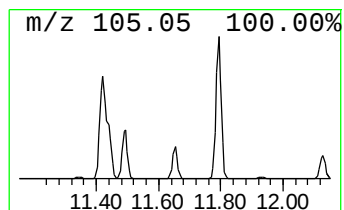
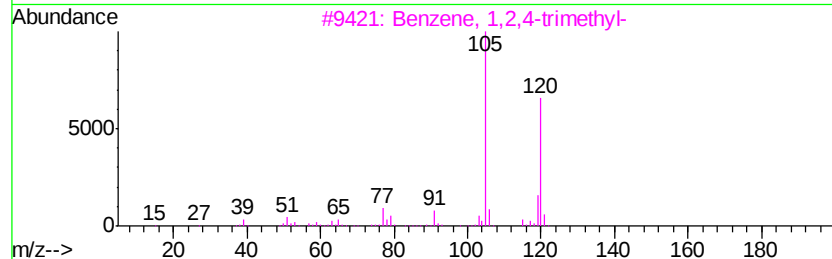
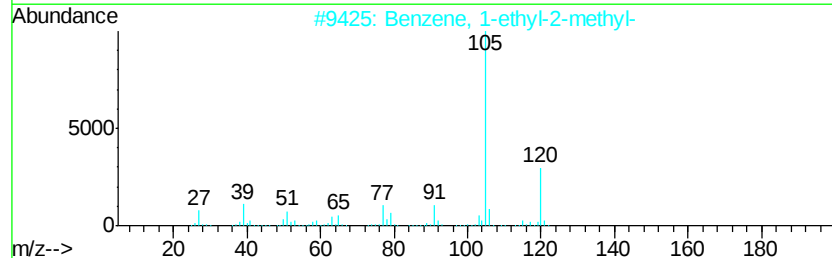
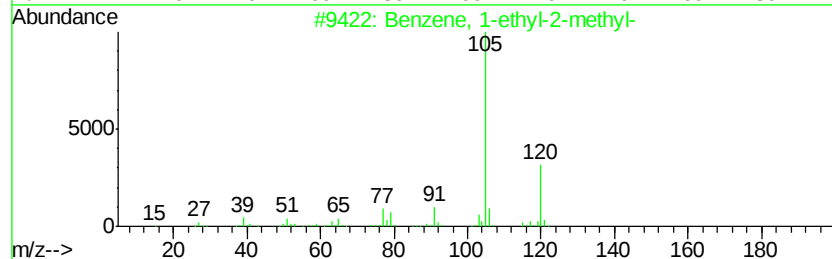
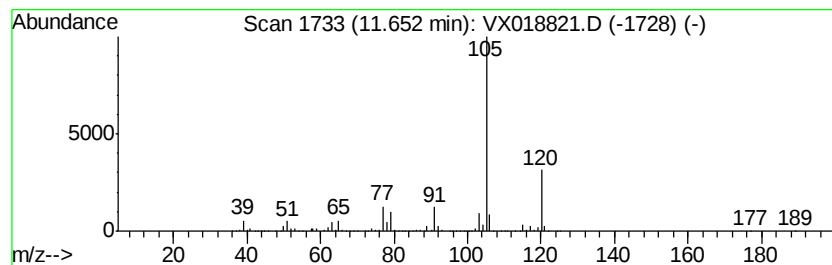
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Benzene, 1,2,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	6.27 ug/L	1761920	1,4-Dichlorobenzene-d4	12.06

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	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

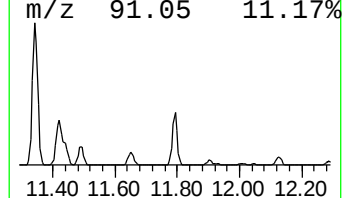
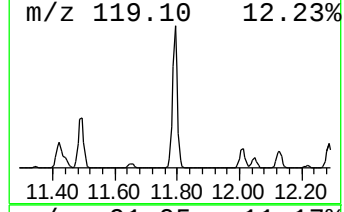
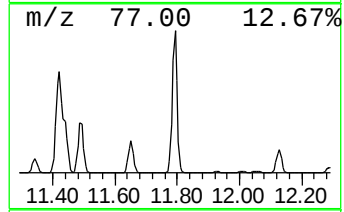
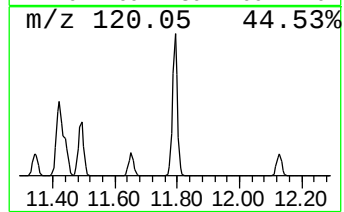
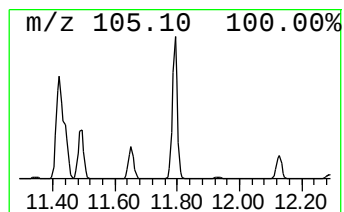
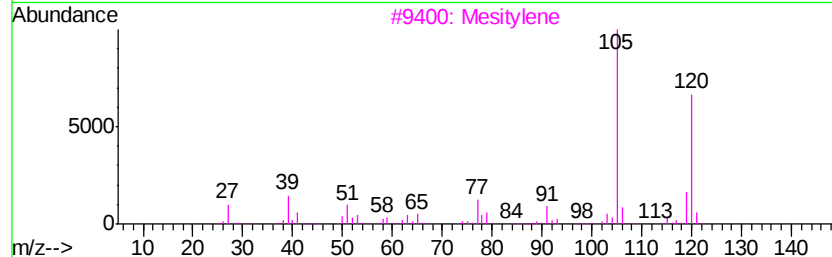
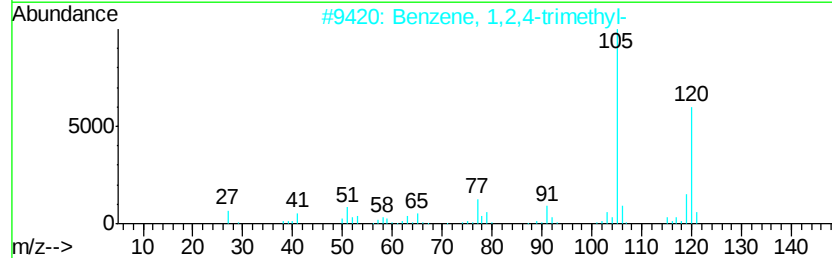
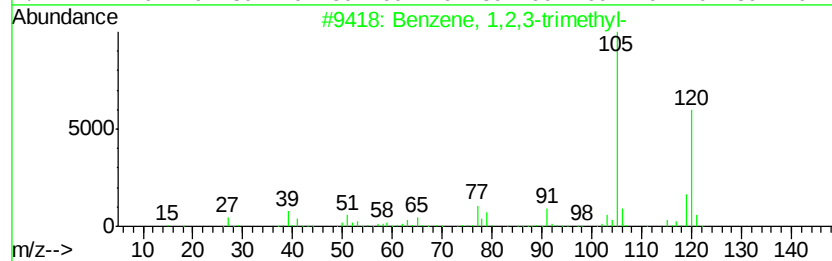
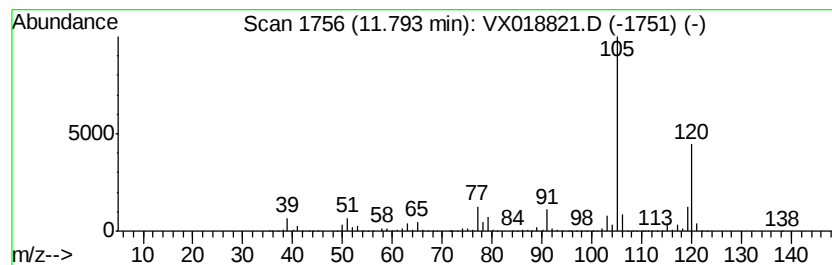
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	30.16 ug/L	8468950	1,4-Dichlorobenzene-d4	12.06

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

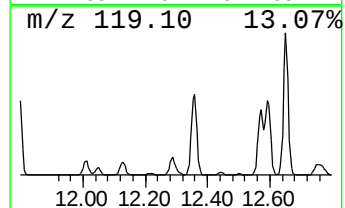
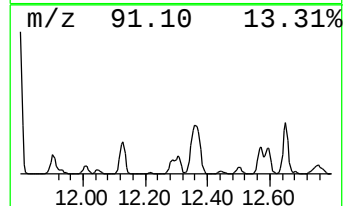
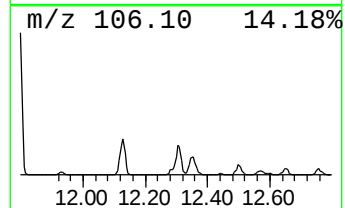
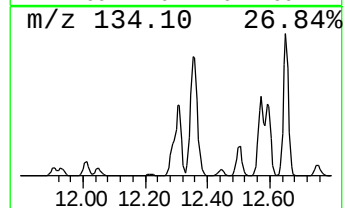
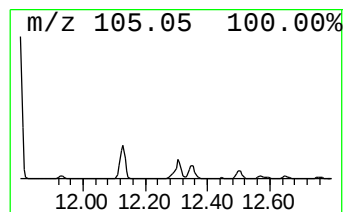
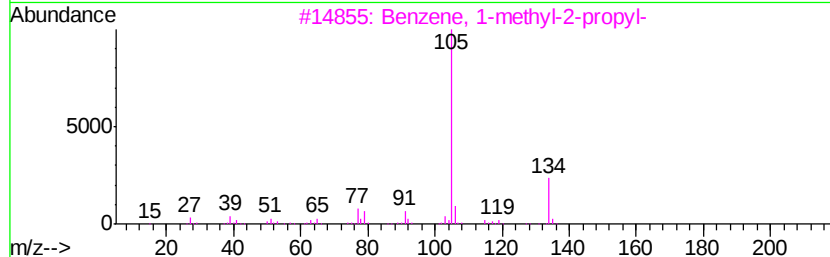
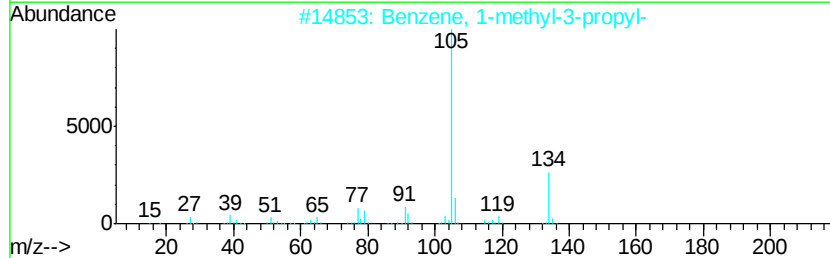
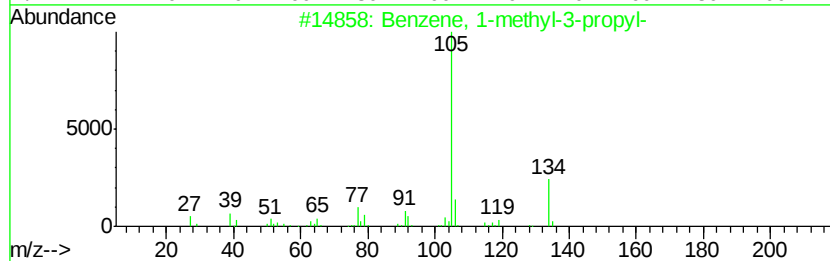
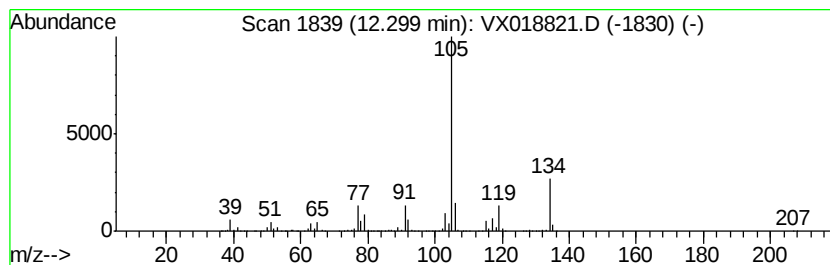
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 Benzene, 1-methyl-3-propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	3.80 ug/L	1066140	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	95
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	94
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

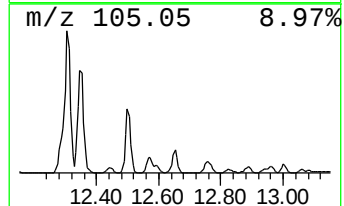
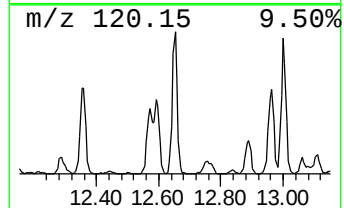
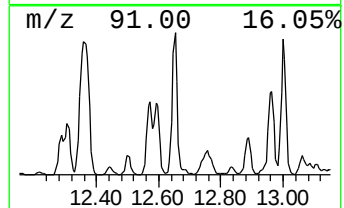
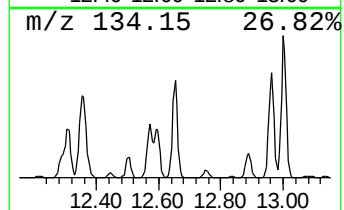
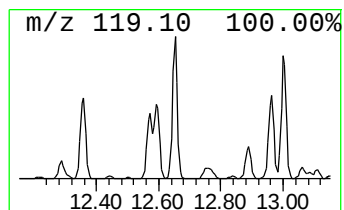
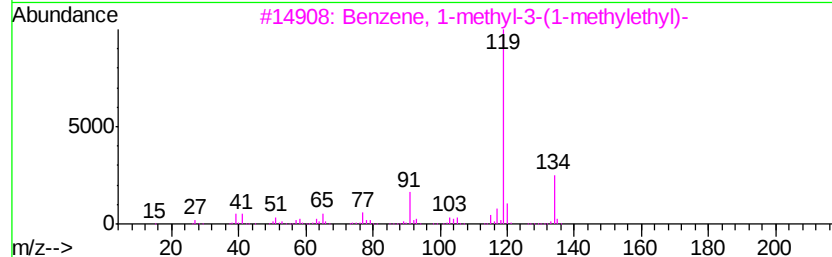
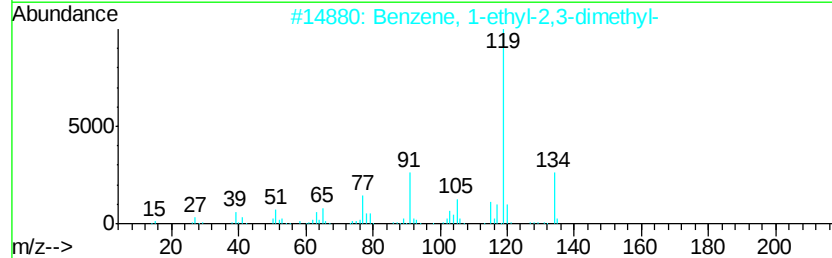
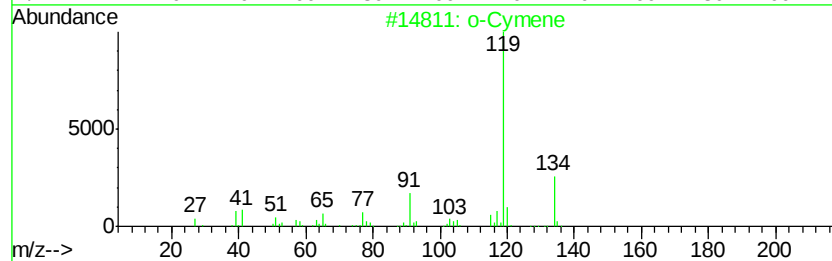
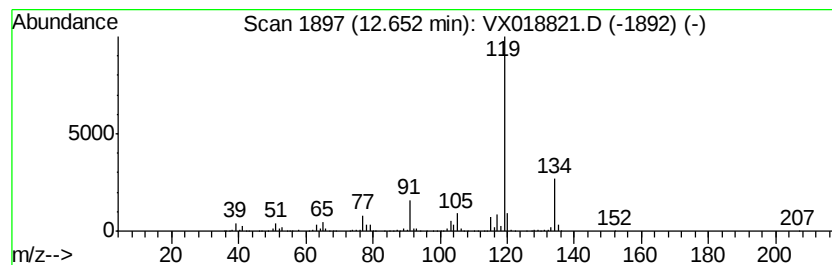
Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 o-Cymene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	4.61 ug/L	1294550	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 97
2		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 96
3		Benzene, 1-methyl-3-(1-methylethyl-)	134 C10H14	000535-77-3 95
4		o-Cymene	134 C10H14	000527-84-4 95
5		p-Cymene	134 C10H14	000099-87-6 95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-100620

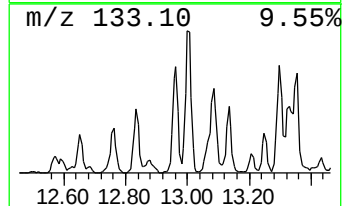
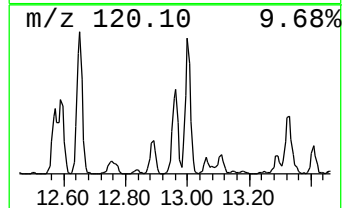
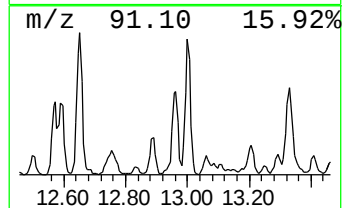
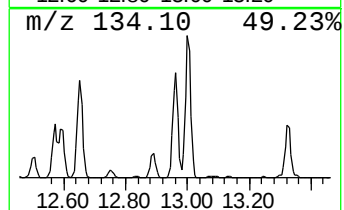
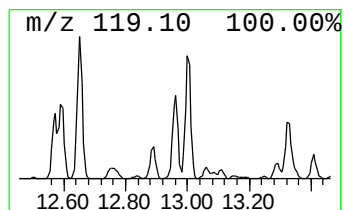
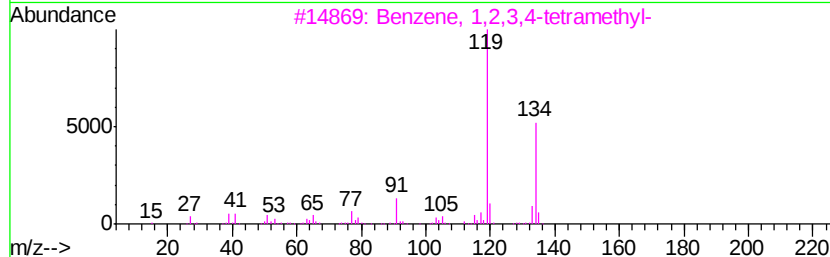
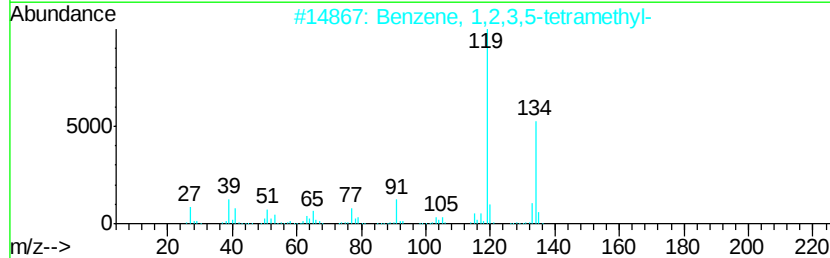
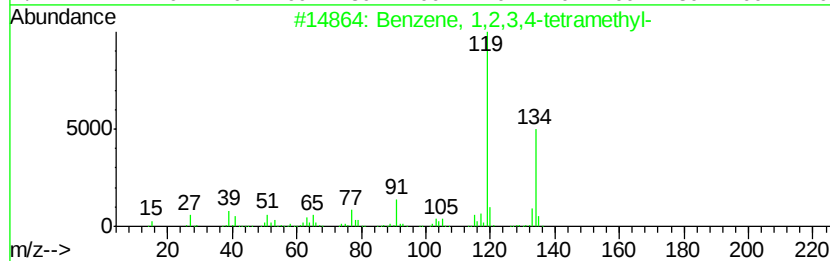
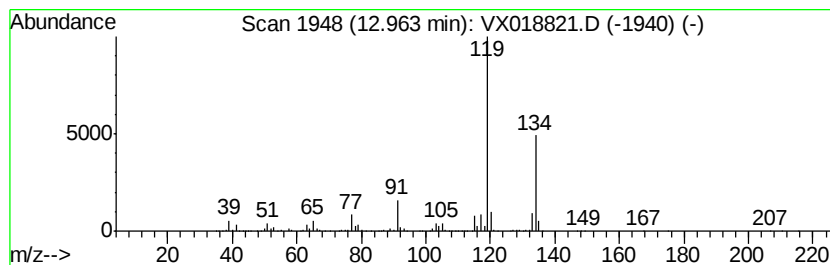
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 15 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	3.30 ug/L	925512	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
OK-01-100620

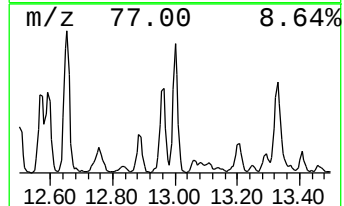
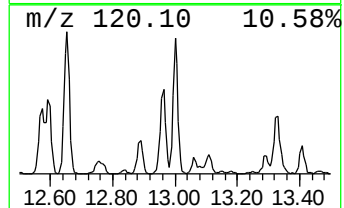
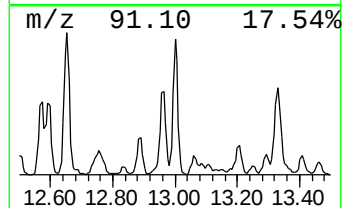
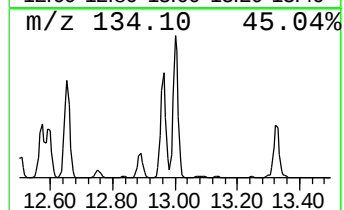
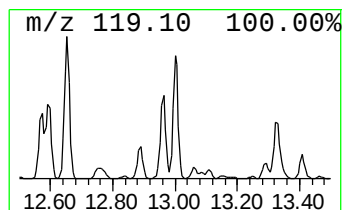
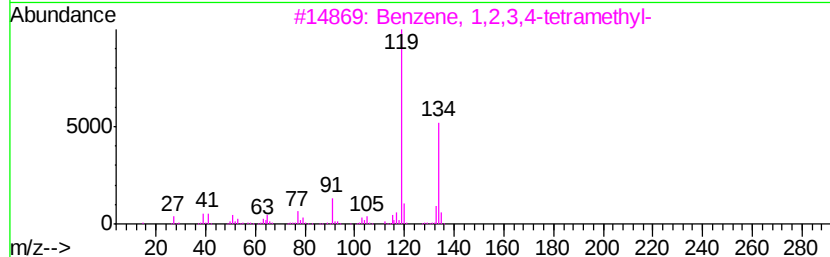
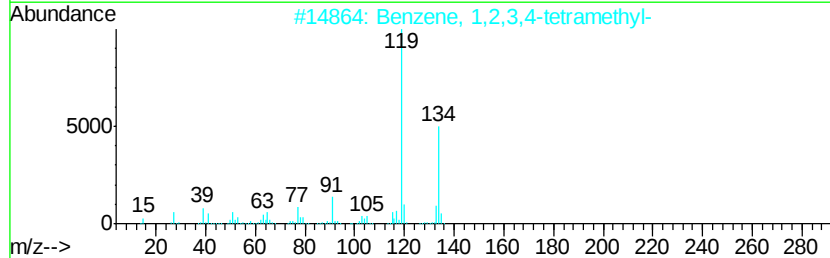
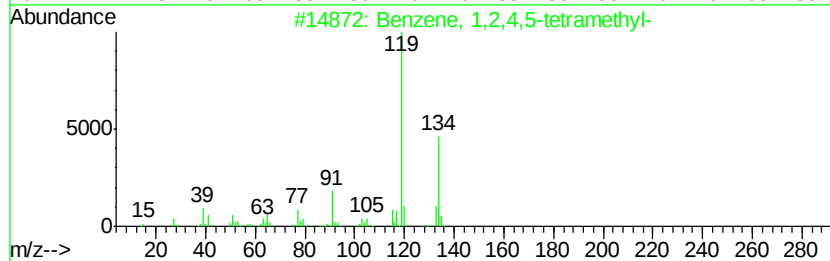
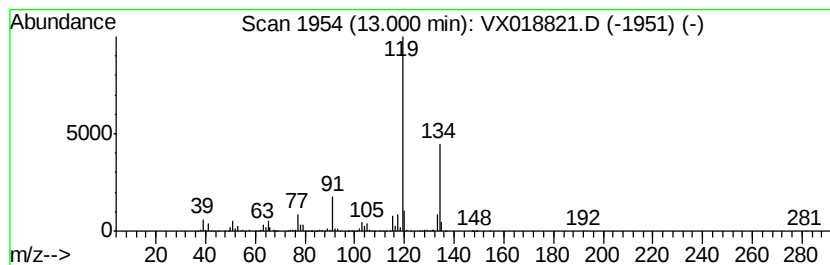
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 16 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	4.52 ug/L	1269360	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100720\
Data File : VX018821.D
Acq On : 07 Oct 2020 19:22
Operator : JC/SP
Sample : L4240-11DL 4X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
OK-01-100620

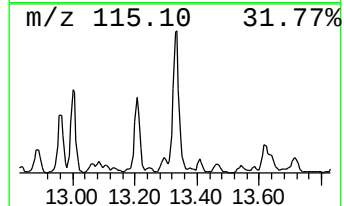
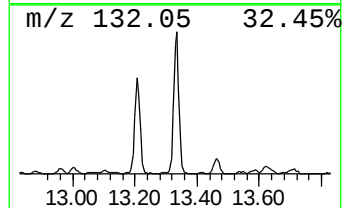
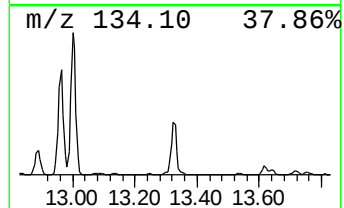
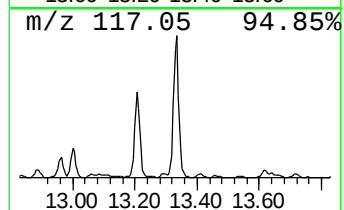
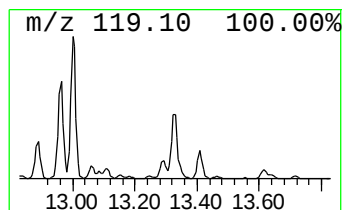
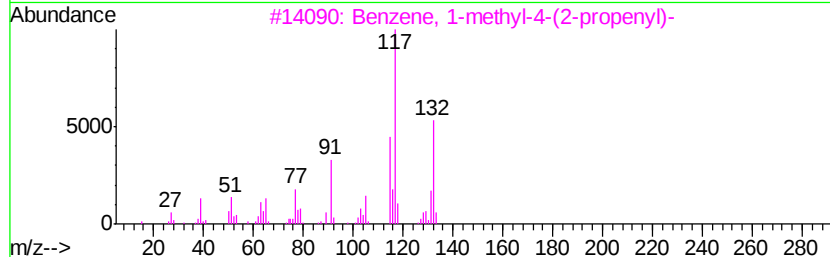
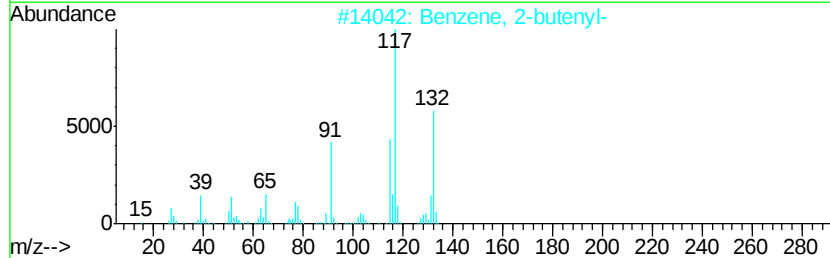
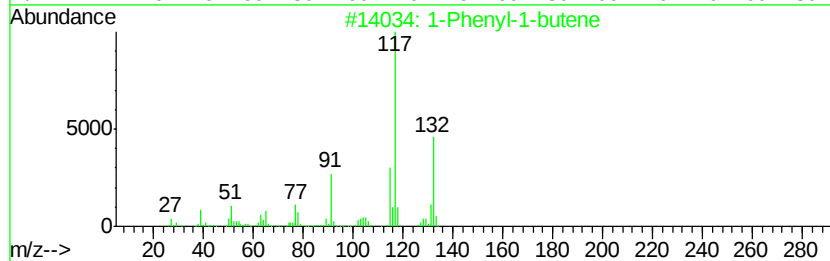
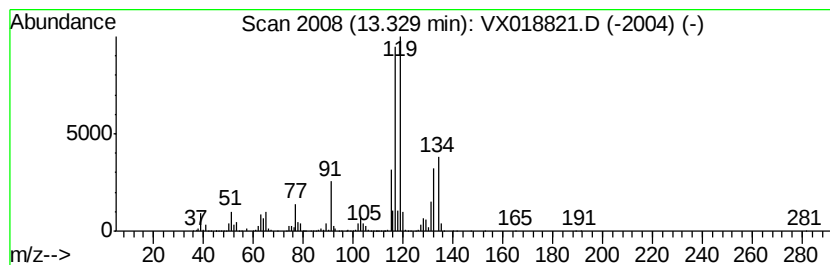
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 17 1-Phenyl-1-butene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	4.49 ug/L	1260580	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	55
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	55



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100720\
 Data File : VX018821.D
 Acq On : 07 Oct 2020 19:22
 Operator : JC/SP
 Sample : L4240-11DL 4X
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OK-01-100620

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.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...	2.82	3.9	ug/L	2409990	1	6.83	3103460	5.0
(DEL) Alkane: Str...	2.87	36.3	ug/L	22516400	1	6.83	3103460	5.0
(DEL) Alkane: Str...	3.15	66.5	ug/L	41273600	1	6.83	3103460	5.0
(DEL) Alkane: Str...	3.52	146.7	ug/L	91048600	1	6.83	3103460	5.0
(DEL) Alkane: Str...	4.12	6.0	ug/L	3716530	1	6.83	3103460	5.0
(DEL) Alkane: Str...	4.28	4.0	ug/L	2477310	1	6.83	3103460	5.0
(DEL) Alkane: Cyc...	4.38	53.9	ug/L	33431600	1	6.83	3103460	5.0
Benzene, propyl-	11.34	6.9	ug/L	1937140	3	12.06	1403930	5.0
Benzene, 1-ethyl-...	11.42	30.5	ug/L	8565000	3	12.06	1403930	5.0
Mesitylene	11.49	11.4	ug/L	3210710	3	12.06	1403930	5.0
Benzene, 1,2,4-tr...	11.65	6.3	ug/L	1761920	3	12.06	1403930	5.0
Benzene, 1,2,3-tr...	11.79	30.2	ug/L	8468950	3	12.06	1403930	5.0
Benzene, 1-methyl...	12.30	3.8	ug/L	1066140	3	12.06	1403930	5.0
o-Cymene	12.65	4.6	ug/L	1294550	3	12.06	1403930	5.0
Benzene, 1,2,3,4-...	12.96	3.3	ug/L	925512	3	12.06	1403930	5.0
Benzene, 1,2,4,5-...	13.00	4.5	ug/L	1269360	3	12.06	1403930	5.0
1-Phenyl-1-butene	13.33	4.5	ug/L	1260580	3	12.06	1403930	5.0