

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
 Data File : VX018791.D
 Acq On : 06 Oct 2020 11:43
 Operator : JC/SP
 Sample : L4240-01
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3Q7

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	2.343	201	206	217	rBV	111774	192828	1.04%	0.217%
2	2.825	277	285	287	rBV	554338	1026819	5.53%	1.153%
3	2.873	287	293	303	rVB	2951498	6106803	32.90%	6.860%
4	3.154	328	339	358	rBV	5403310	12110078	65.24%	13.604%
5	3.508	386	397	413	rBV	7745041	17591252	94.77%	19.761%
6	3.727	426	433	441	rVB2	96774	232321	1.25%	0.261%
7	4.117	487	497	511	rVV	333514	1020649	5.50%	1.147%
8	4.282	512	524	529	rVV	257949	774795	4.17%	0.870%
9	4.379	529	540	555	rVV	6425203	18562927	100.00%	20.852%
10	4.544	555	567	581	rVB3	79381	303723	1.64%	0.341%
11	5.141	651	665	674	rBV	70434	232012	1.25%	0.261%
12	5.550	722	732	746	rVB	643139	1945255	10.48%	2.185%
13	6.044	800	813	824	rBV2	163629	479943	2.59%	0.539%
14	6.830	932	942	949	rBV	131731	316715	1.71%	0.356%
15	7.367	1020	1030	1038	rBV	102949	227221	1.22%	0.255%
16	8.696	1233	1248	1253	rBV	259668	433957	2.34%	0.487%
17	8.763	1253	1259	1273	rVB	2828179	4250601	22.90%	4.775%
18	9.427	1361	1368	1377	rBV	330374	457582	2.47%	0.514%
19	10.098	1472	1478	1488	rBV2	269029	506307	2.73%	0.569%
20	10.232	1495	1500	1505	rBV	337919	448879	2.42%	0.504%
21	10.342	1511	1518	1525	rBV	1185001	1597340	8.61%	1.794%
22	10.683	1568	1574	1581	rVB	456796	582497	3.14%	0.654%
23	11.000	1621	1626	1632	rBV	161992	205659	1.11%	0.231%
24	11.342	1676	1682	1689	rBV	613036	758709	4.09%	0.852%
25	11.421	1689	1695	1702	rVV	2114445	3691195	19.88%	4.146%
26	11.488	1702	1706	1716	rVB	1025312	1270861	6.85%	1.428%
27	11.610	1721	1726	1729	rVV	736637	894212	4.82%	1.004%
28	11.652	1729	1733	1740	rVB	877315	1084557	5.84%	1.218%
29	11.793	1750	1756	1761	rBV	3299371	3918161	21.11%	4.401%
30	12.006	1786	1791	1795	rBV	189226	233196	1.26%	0.262%
31	12.079	1795	1803	1807	rVV2	663938	1195323	6.44%	1.343%
32	12.128	1807	1811	1816	rVB	746270	899510	4.85%	1.010%
33	12.280	1831	1836	1844	rBV2	302498	528593	2.85%	0.594%
34	12.360	1844	1849	1858	rVB3	538925	950419	5.12%	1.068%

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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_X\METHOD\SOMXTR100220WMA.M
Title : TRACE VOA SOM01.0

35	12.652	1892	1897	1901	rBV	187568	219388	1.18%	0.246%
36	13.000	1951	1954	1960	rVB	197517	228090	1.23%	0.256%
37	13.329	2004	2008	2013	rVB2	162096	214679	1.16%	0.241%
38	13.628	2051	2057	2067	rBV	2044832	2460531	13.26%	2.764%
39	14.000	2106	2118	2124	rBV	648276	867439	4.67%	0.974%

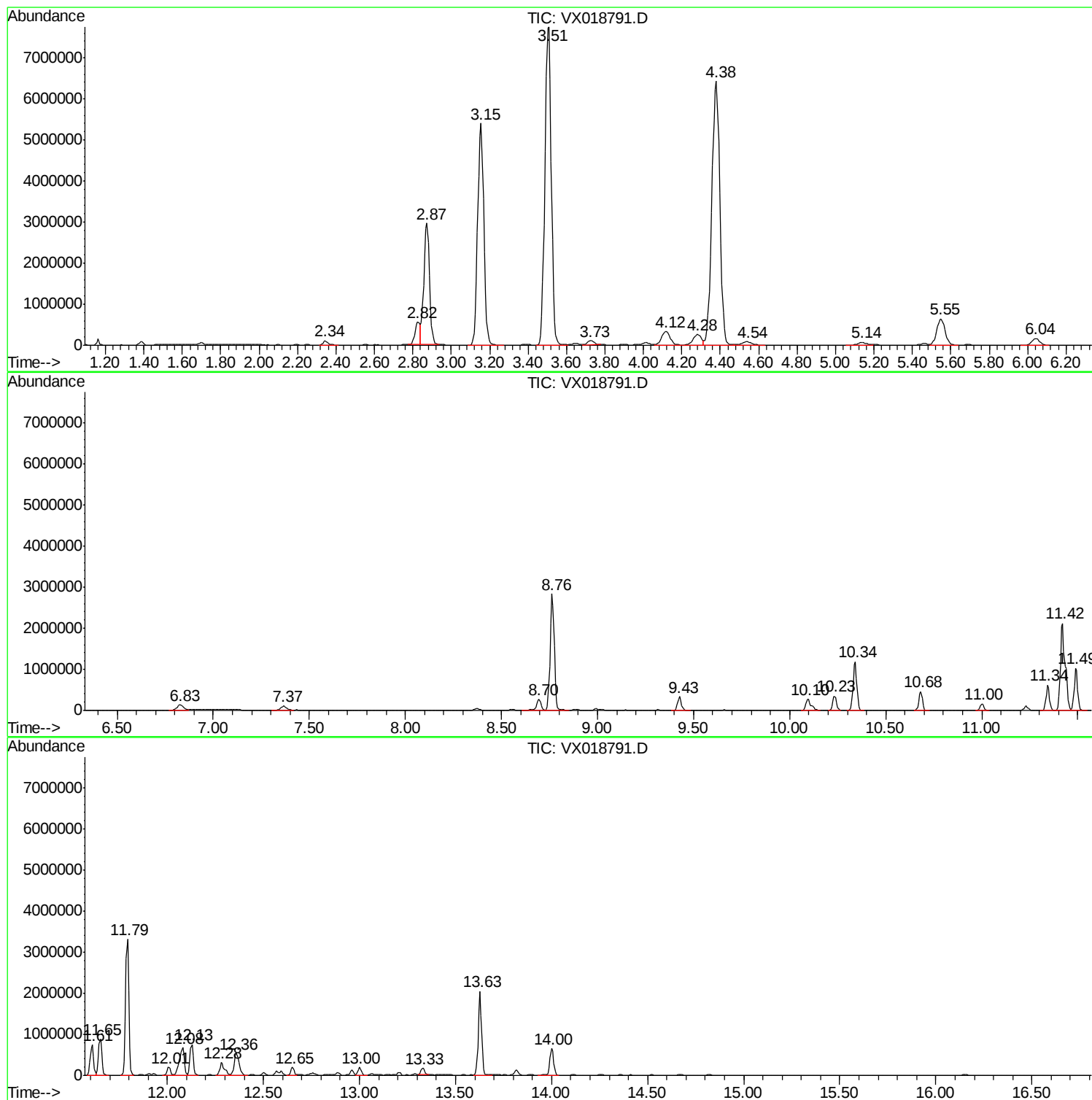
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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



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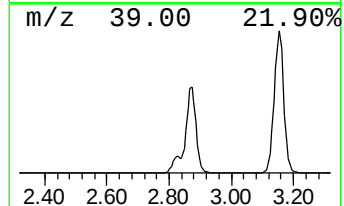
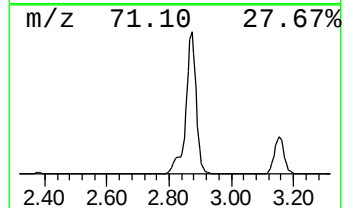
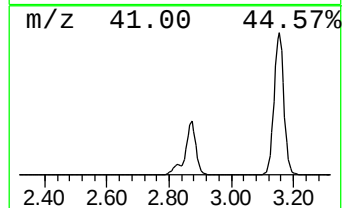
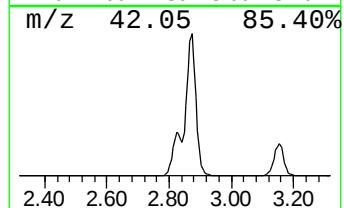
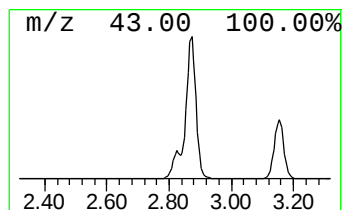
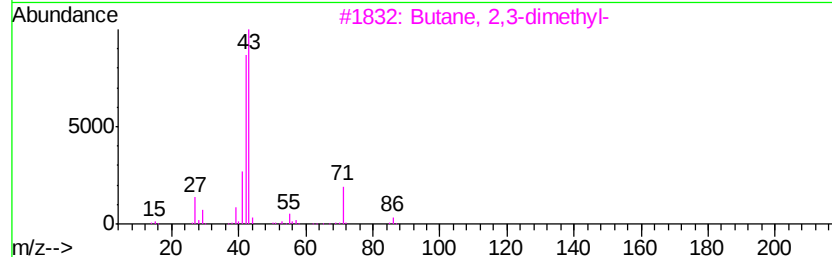
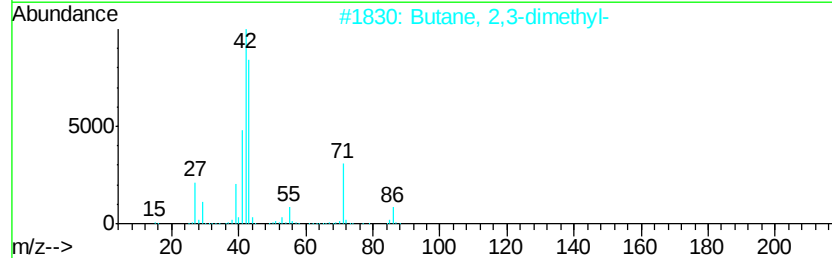
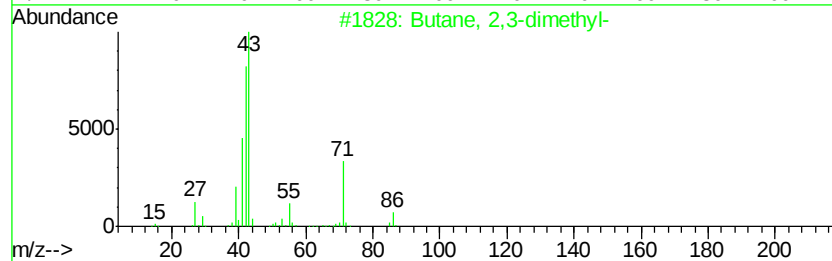
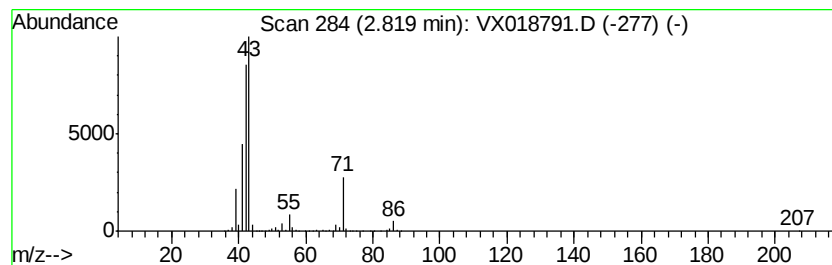
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.82	16.21 ug/L	1026820	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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Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3Q7

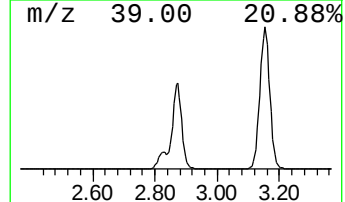
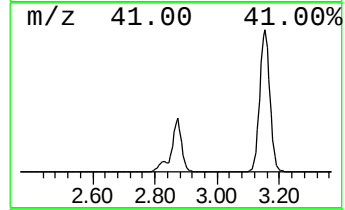
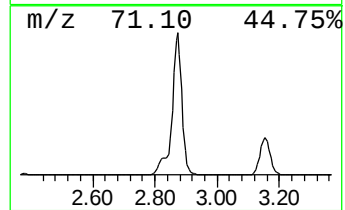
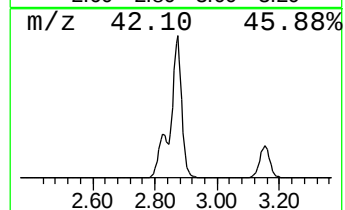
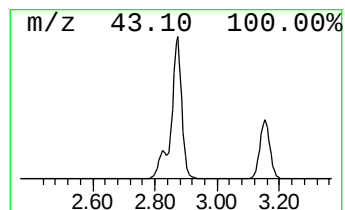
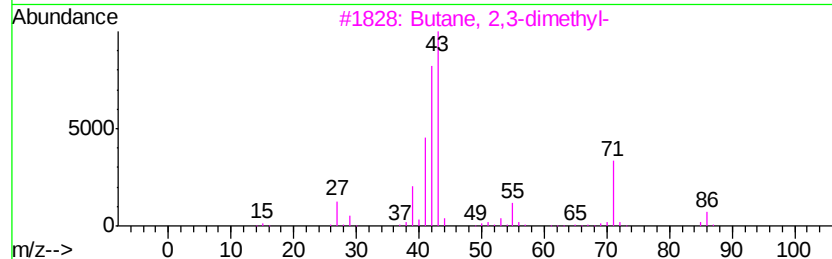
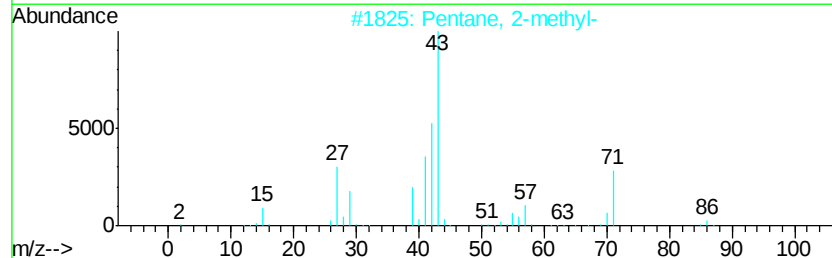
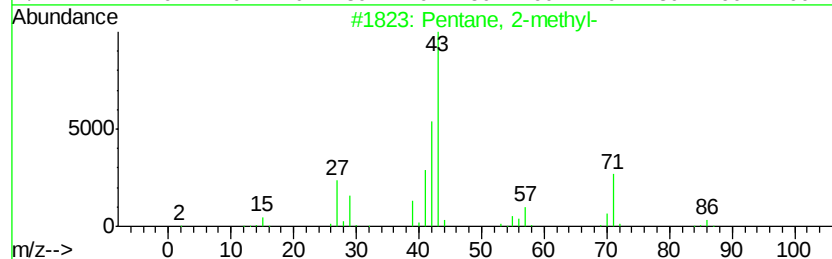
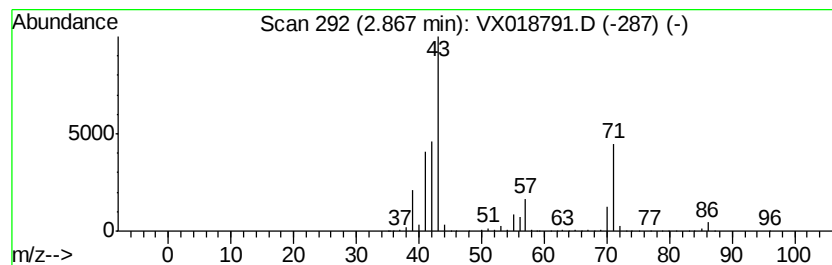
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Quant Title : TRACE VOA SOM01.0

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	96.41 ug/L	6106800	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		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
4		Pentane	72	C5H12	000109-66-0	46
5		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45



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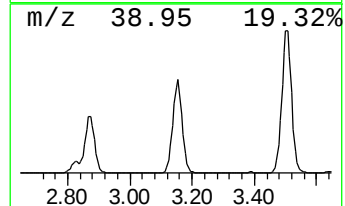
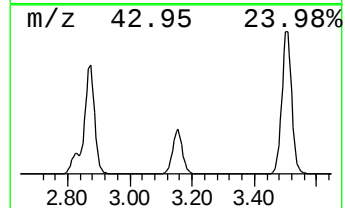
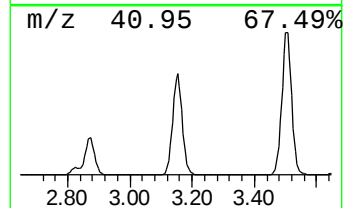
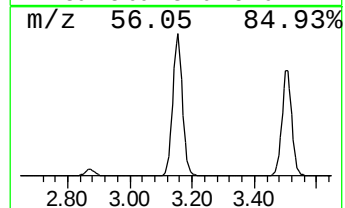
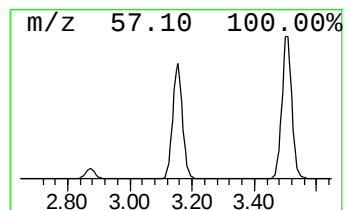
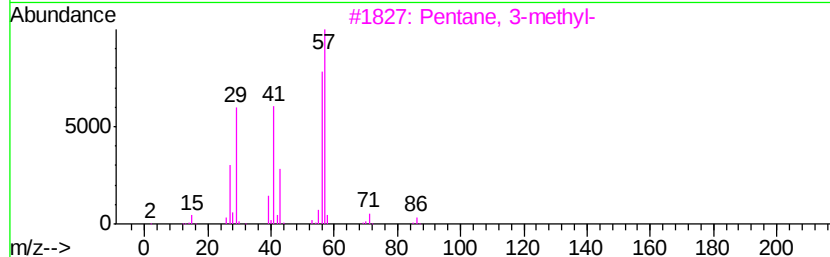
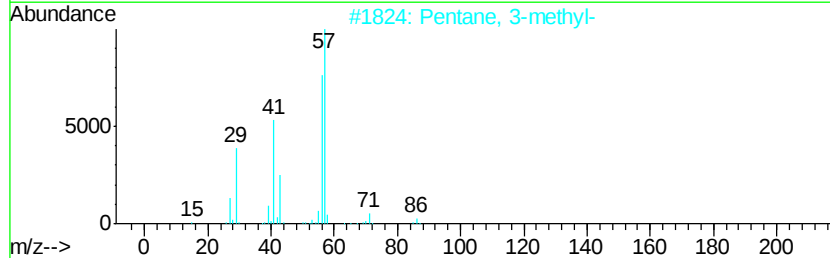
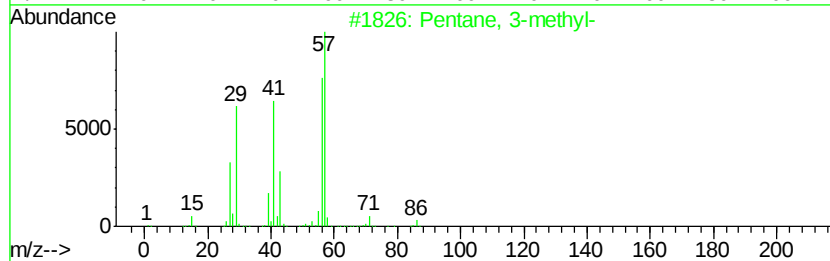
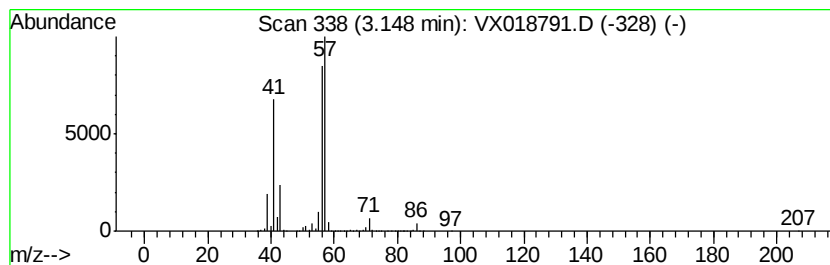
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	191.18 ug/L	12110100	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	91
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	90
4		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64



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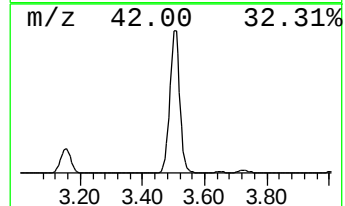
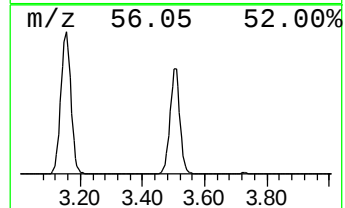
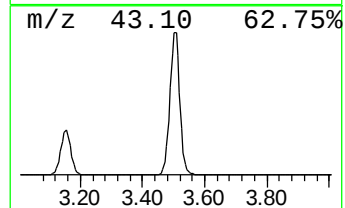
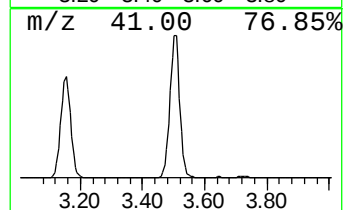
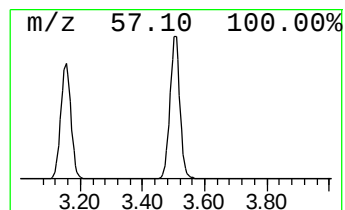
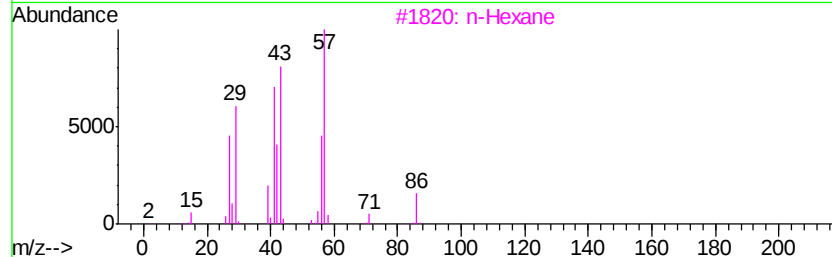
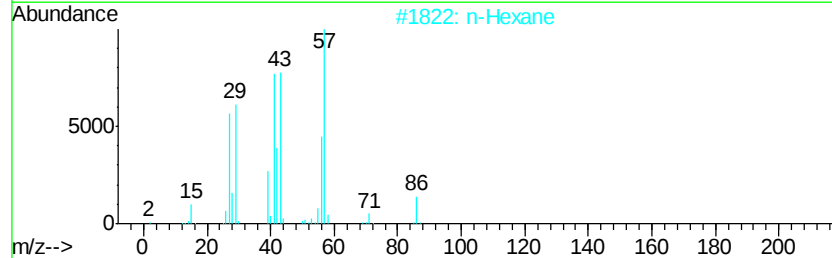
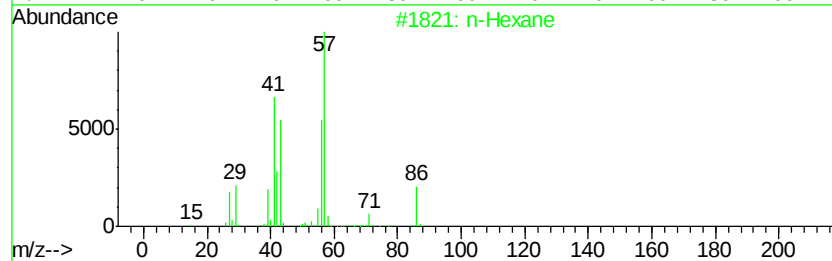
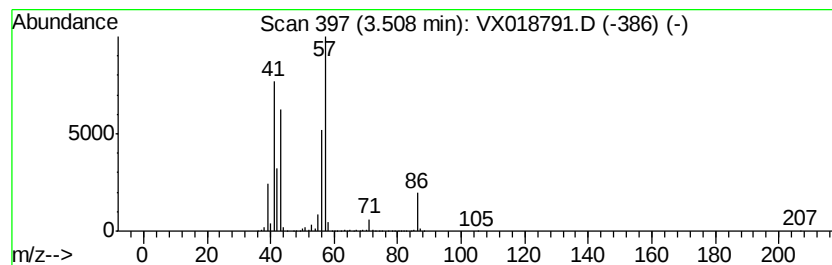
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.51	277.71 ug/L	17591300	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		n-Hexane	86	C6H14	000110-54-3	59
4		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43
5		Butanal, 2-methyl-	86	C5H10O	000096-17-3	43



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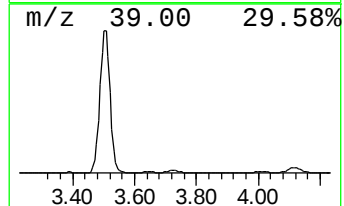
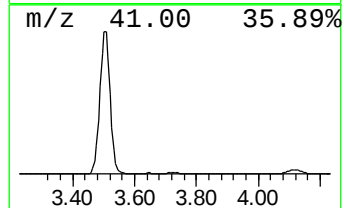
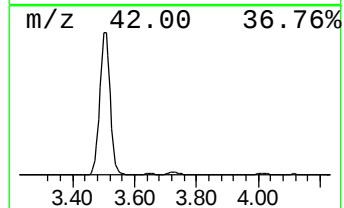
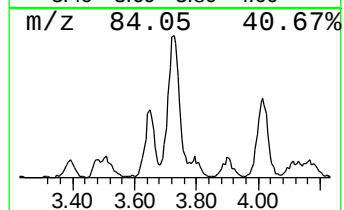
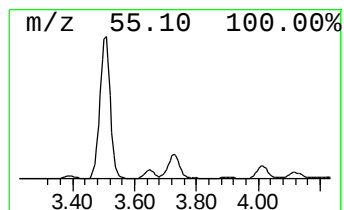
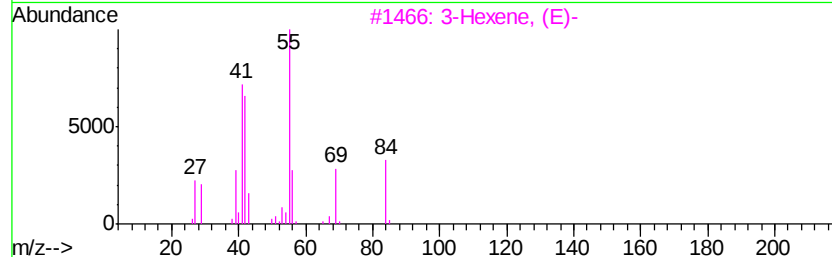
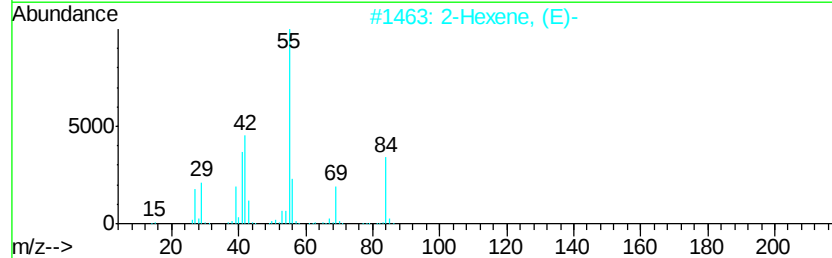
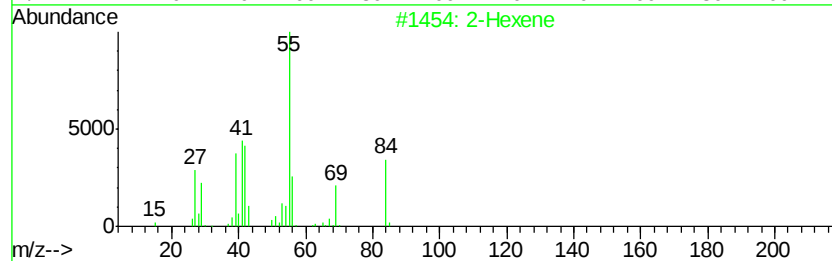
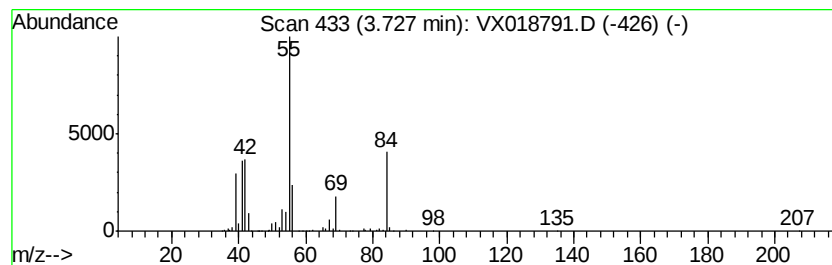
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.73	3.67 ug/L	232321	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hexene	84	C6H12	000592-43-8	94
2		2-Hexene, (E)-	84	C6H12	004050-45-7	90
3		3-Hexene, (E)-	84	C6H12	013269-52-8	86
4		2-Hexene, (Z)-	84	C6H12	007688-21-3	83
5		3-Hexene	84	C6H12	000592-47-2	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3Q7

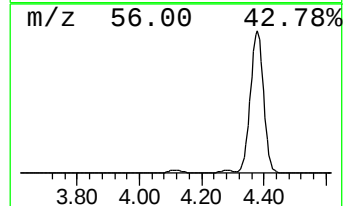
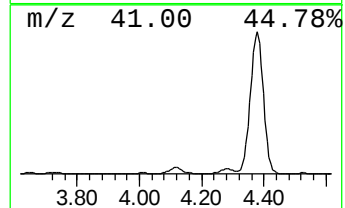
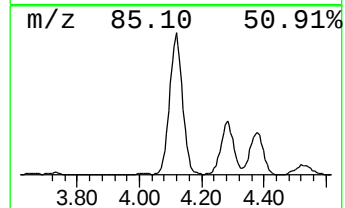
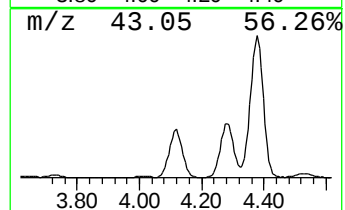
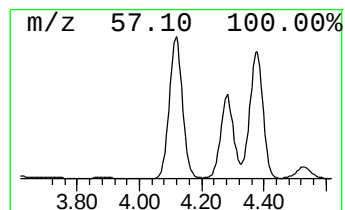
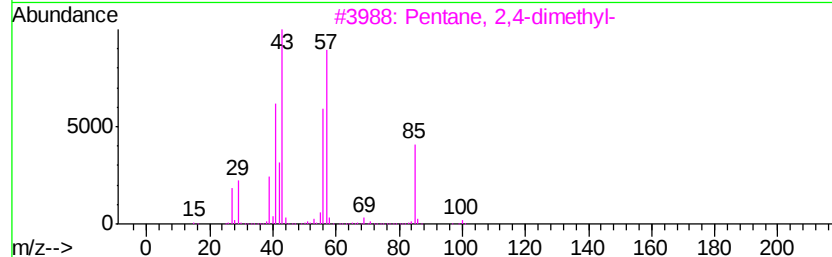
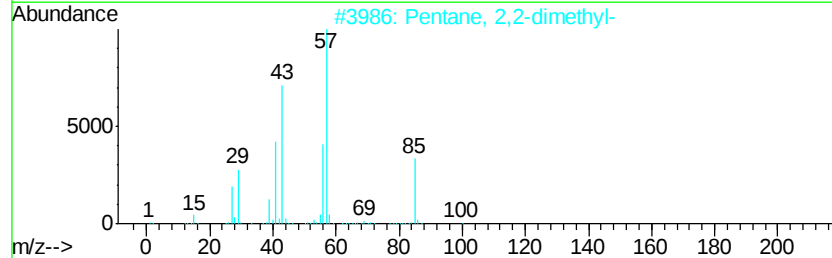
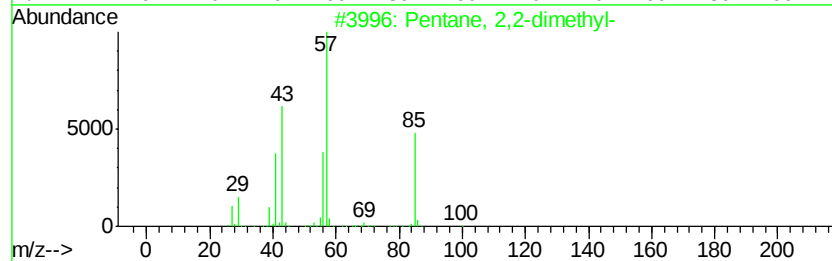
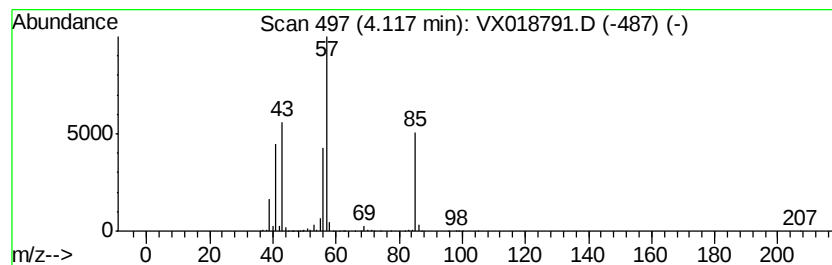
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.12	16.11 ug/L	1020650	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83
2		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	74
3		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
4		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72
5		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3Q7

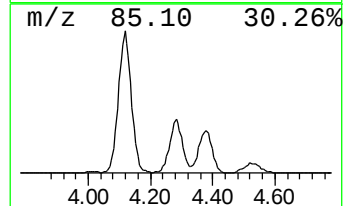
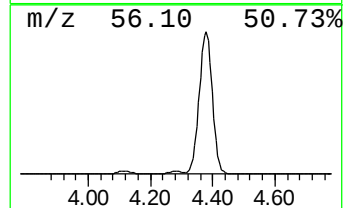
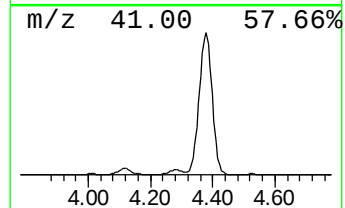
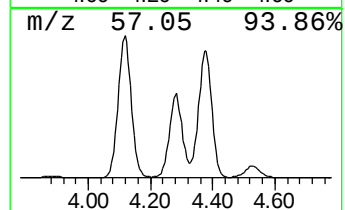
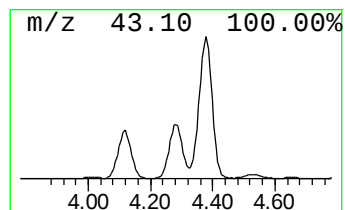
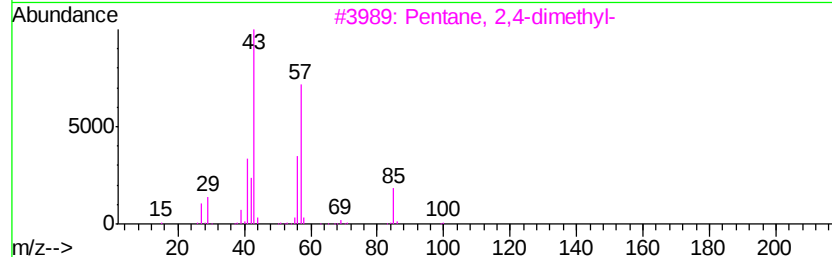
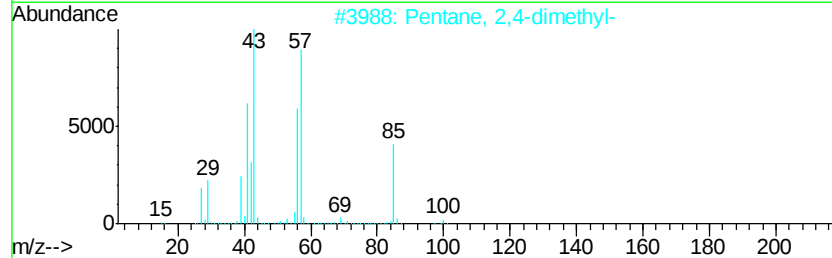
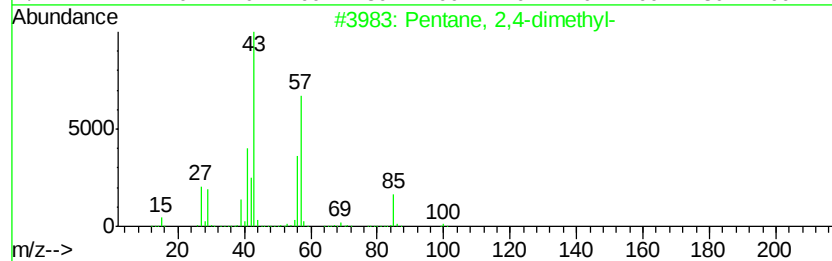
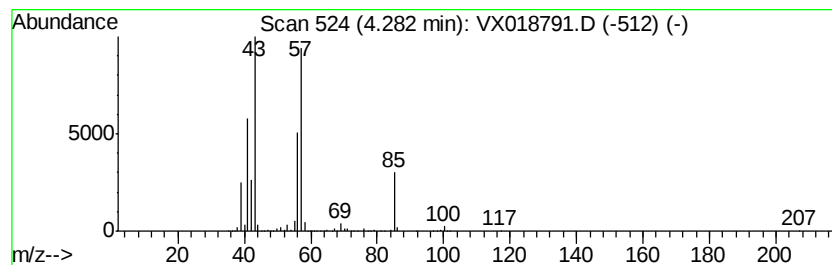
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	12.23 ug/L	774795	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	90
2		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	87
3		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	74
4		Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	64
5		Hexane, 2-methyl-	100	C7H16	000591-76-4	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

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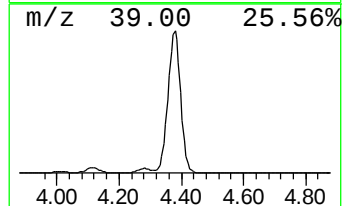
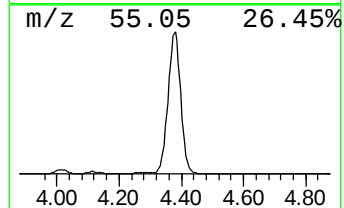
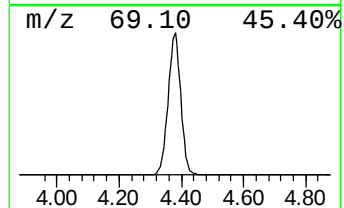
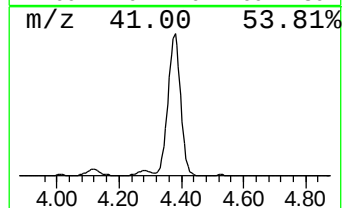
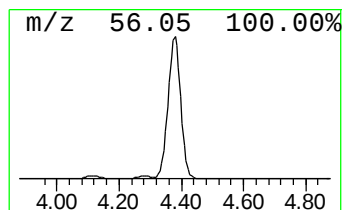
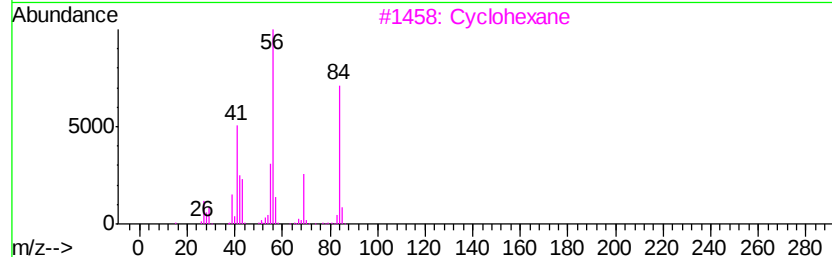
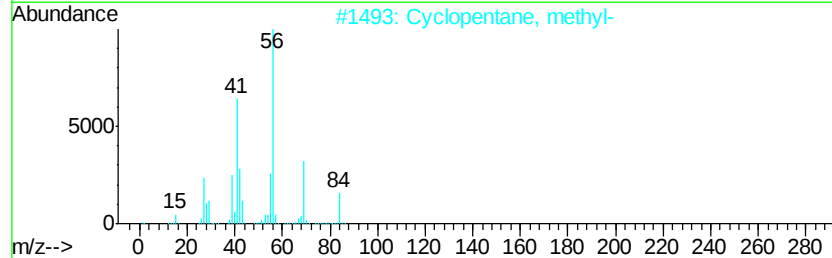
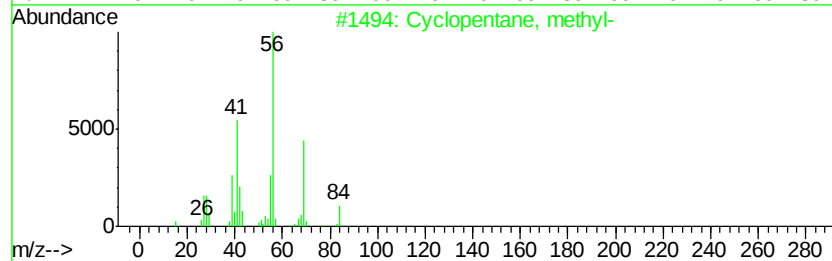
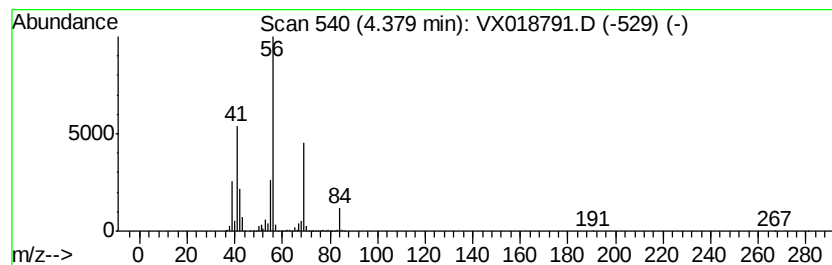
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 (DEL) Alkane: Cyclic4.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	293.05 ug/L	18562900	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	91
3		Cyclohexane	84	C6H12	000110-82-7	78
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

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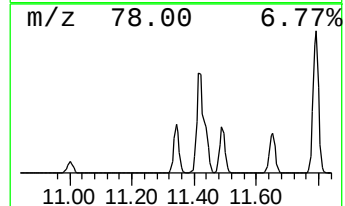
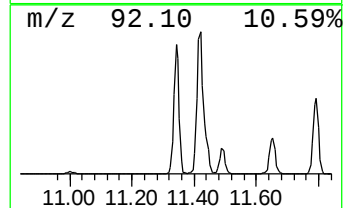
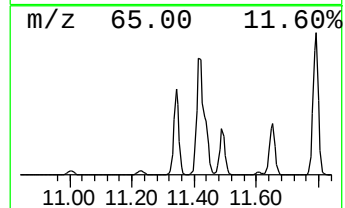
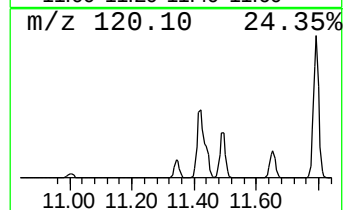
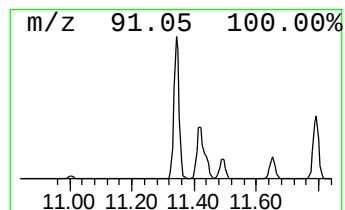
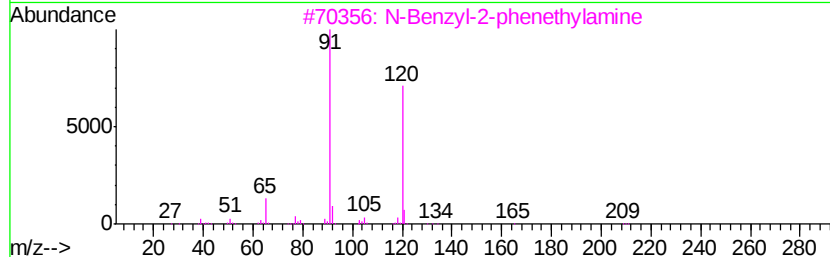
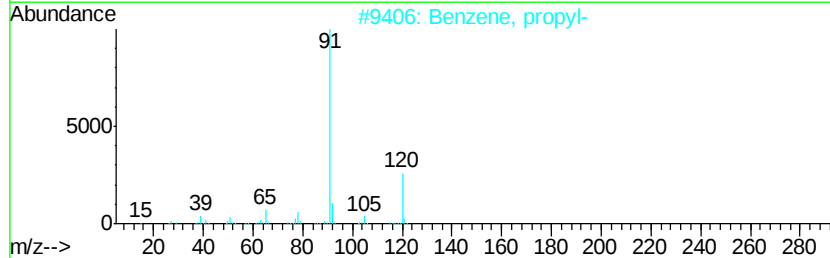
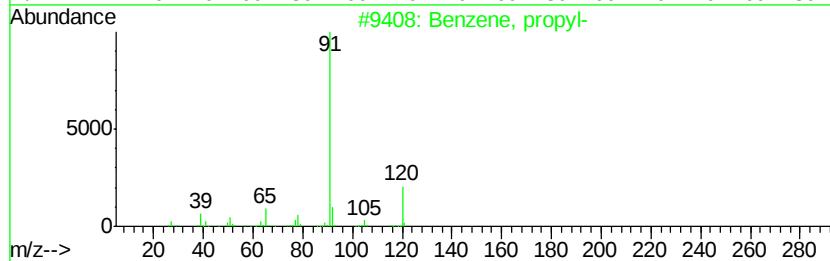
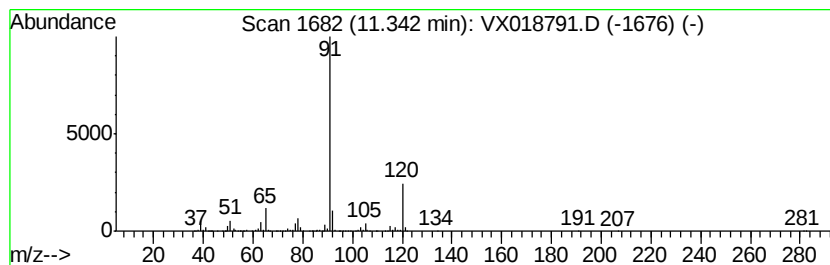
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	3.17 ug/L	758709	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	87
3		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	83
4		Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	78
5		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

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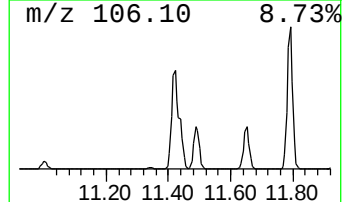
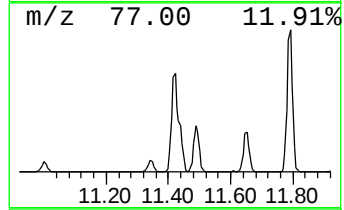
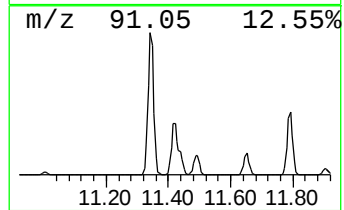
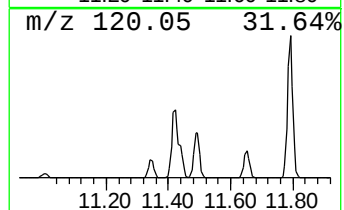
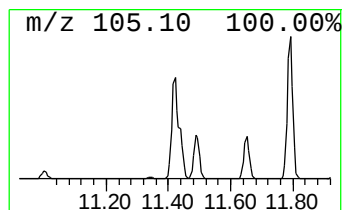
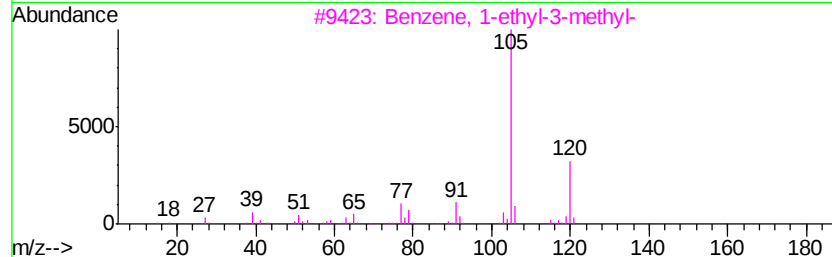
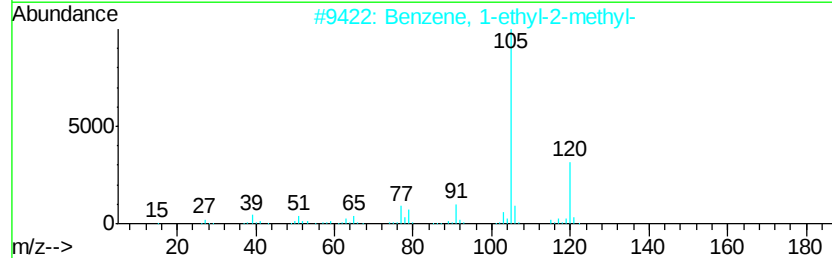
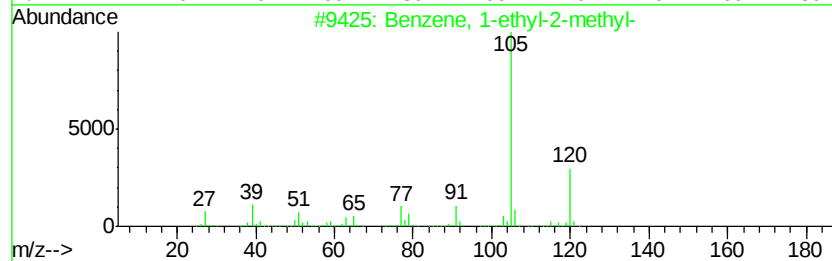
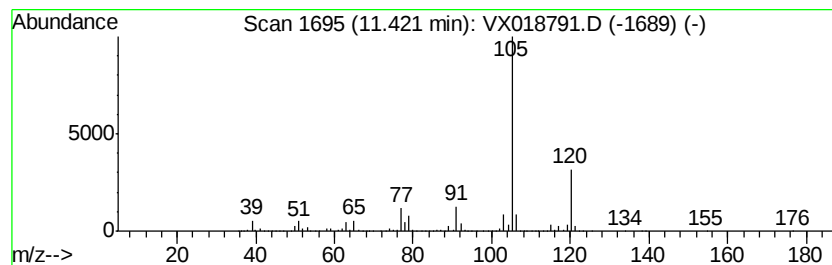
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	15.44 ug/L	3691200	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

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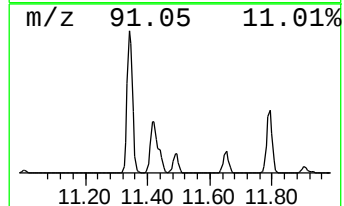
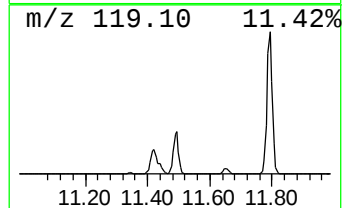
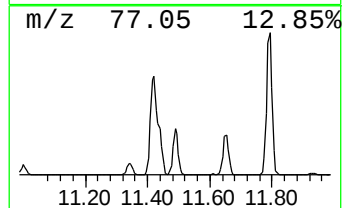
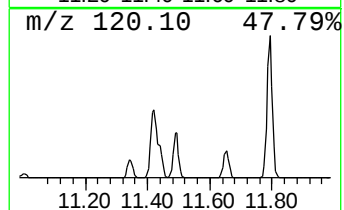
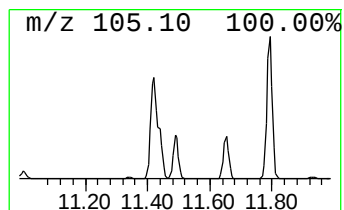
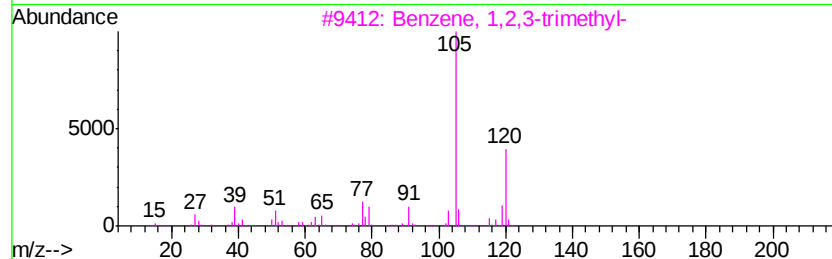
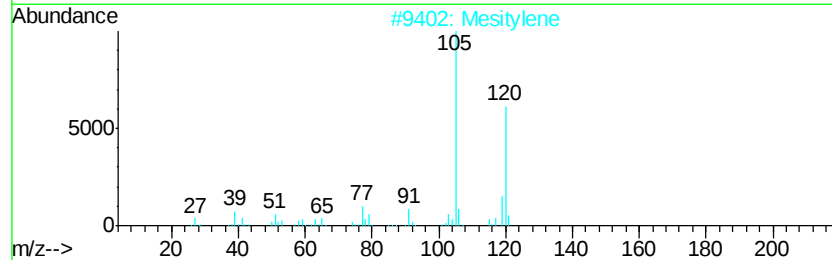
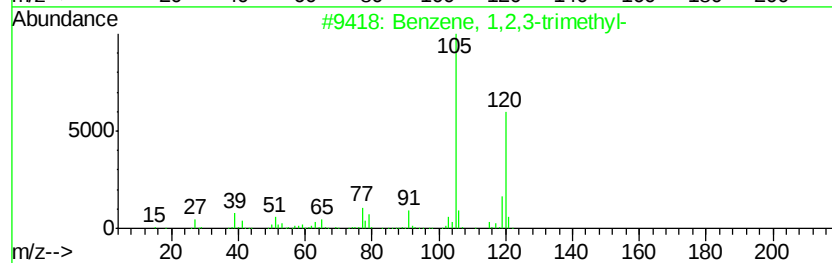
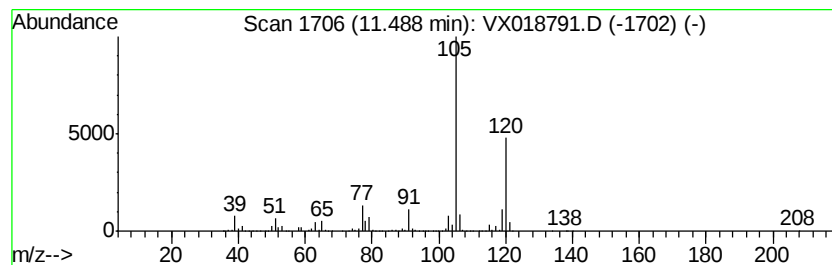
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	5.32 ug/L	1270860	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		Mesitylene	120	C9H12	000108-67-8	97
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4		Mesitylene	120	C9H12	000108-67-8	94
5		Mesitylene	120	C9H12	000108-67-8	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

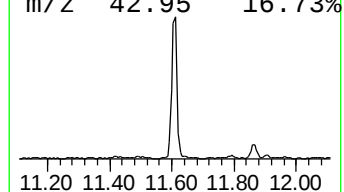
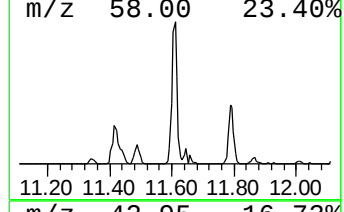
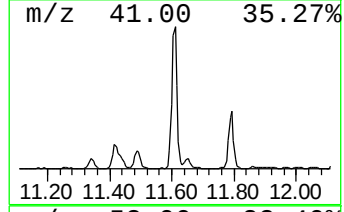
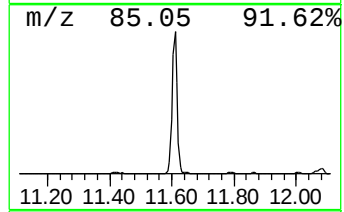
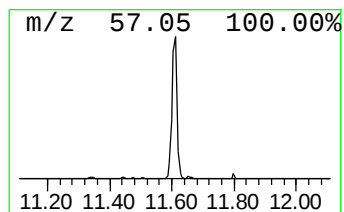
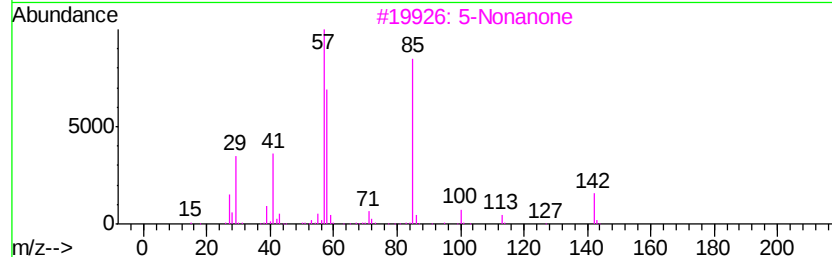
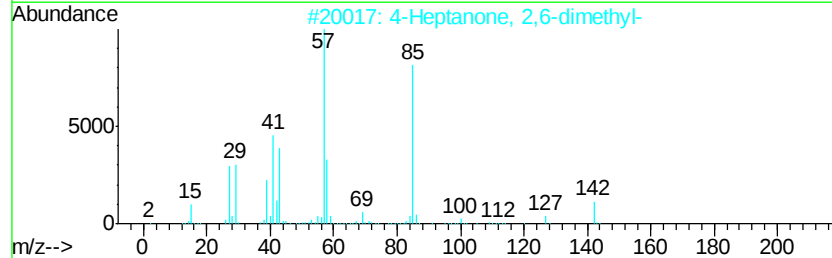
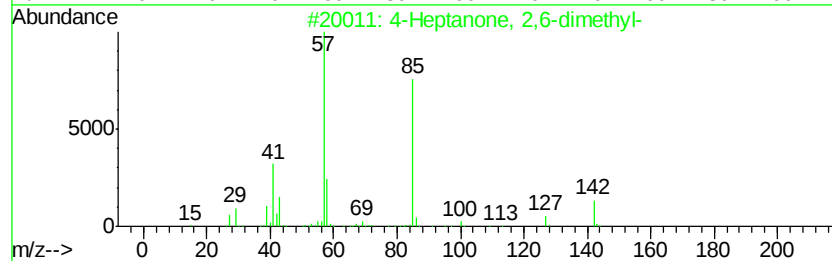
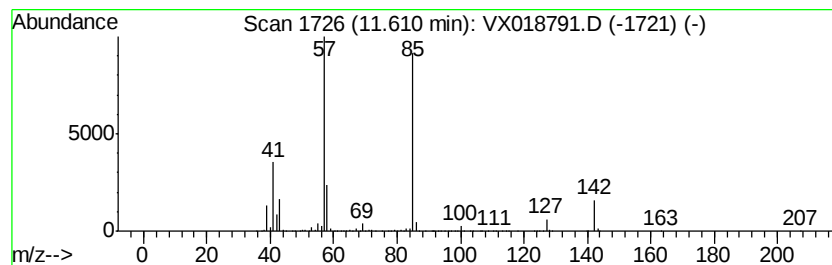
Instrument :
MSVOA_X
ClientSampleId :
BG3Q7

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 12 4-Heptanone, 2,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.61	3.74 ug/L	894212	1,4-Dichlorobenzene-d4		12.06		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	94
2			4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	64
3			5-Nonanone	142	C9H18O	000502-56-7	59
4			5-Nonanone	142	C9H18O	000502-56-7	59
5			3-Buten-2-ol, 2,3-dimethyl-	100	C6H12O	010473-13-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3Q7

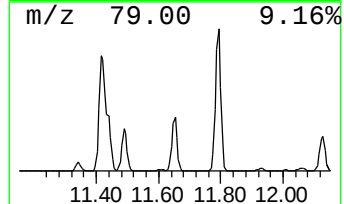
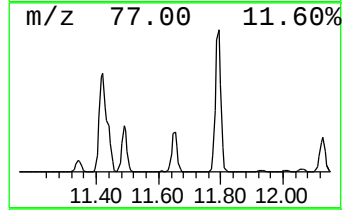
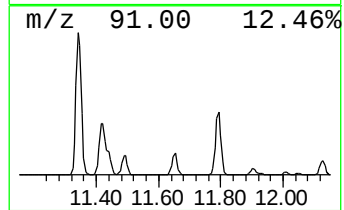
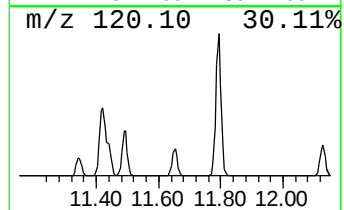
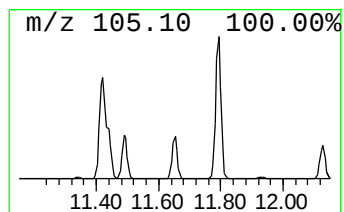
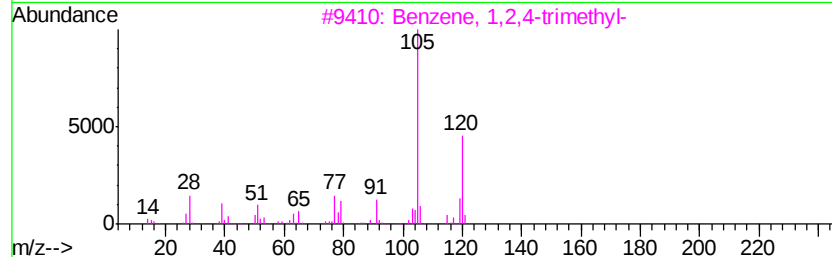
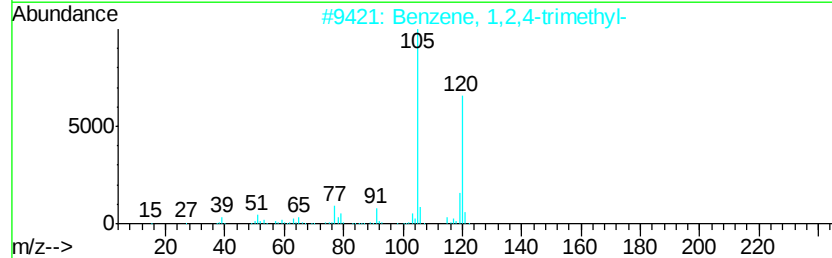
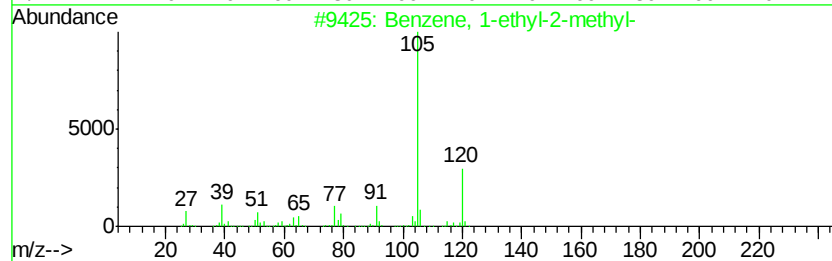
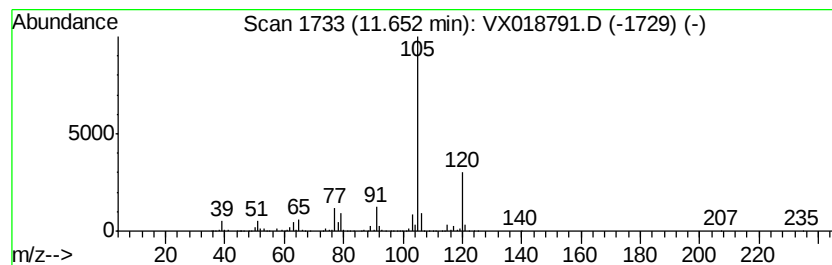
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	4.54 ug/L	1084560	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	94
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
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-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

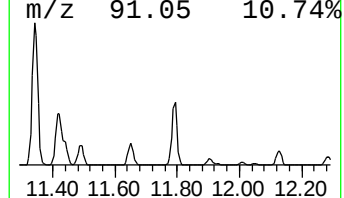
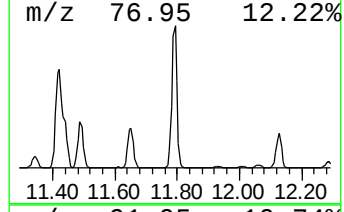
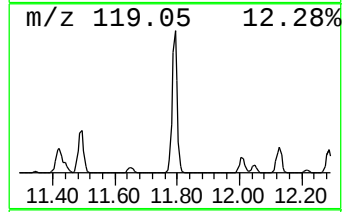
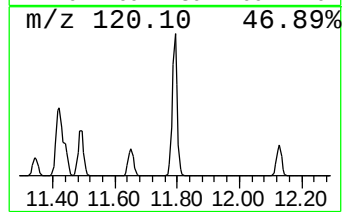
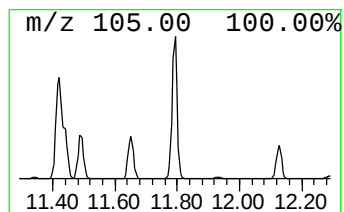
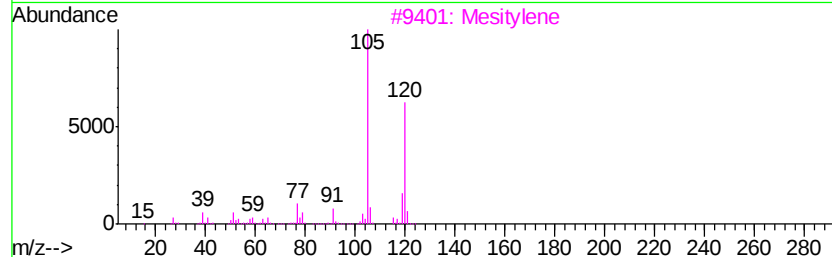
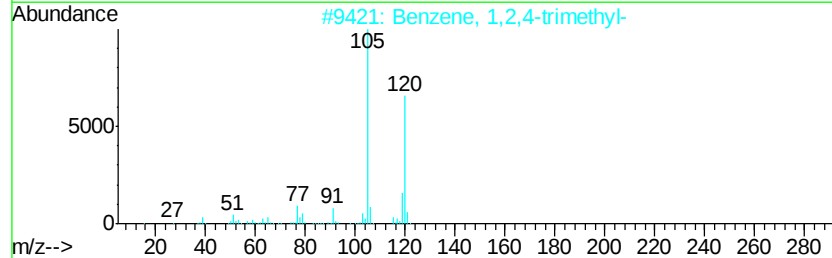
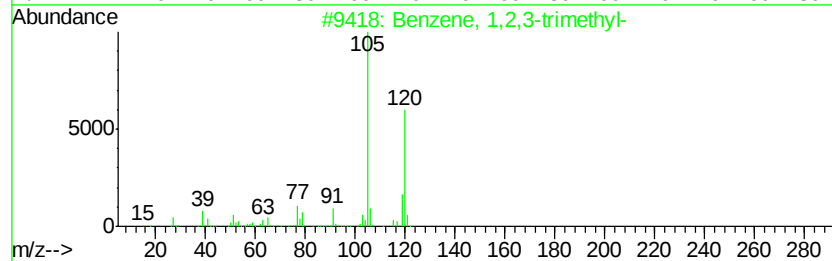
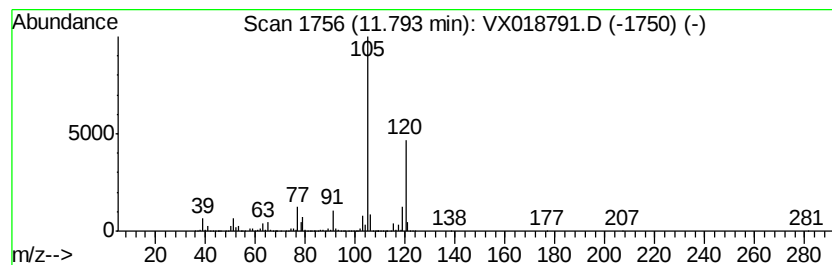
Instrument :
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ClientSampled :
BG3Q7

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Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 Mesitylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.79	16.39 ug/L	3918160	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

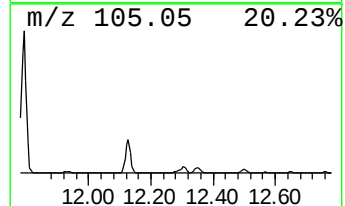
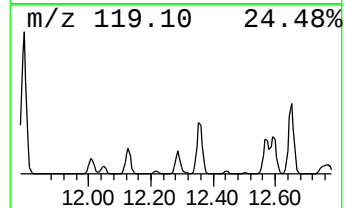
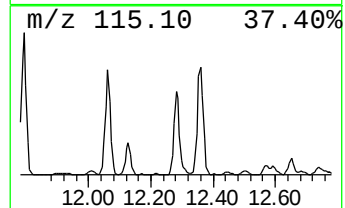
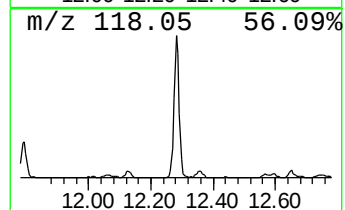
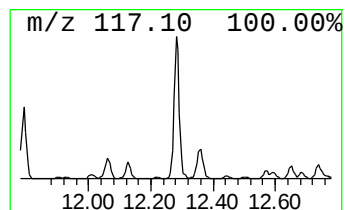
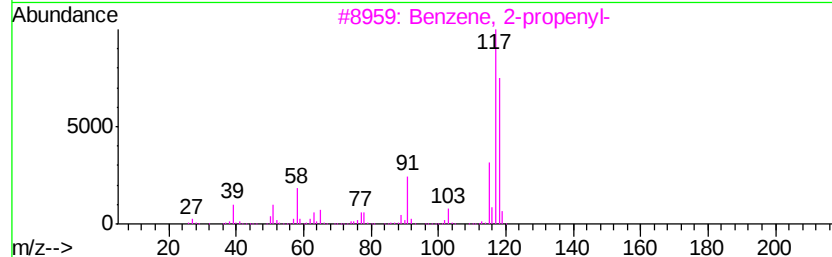
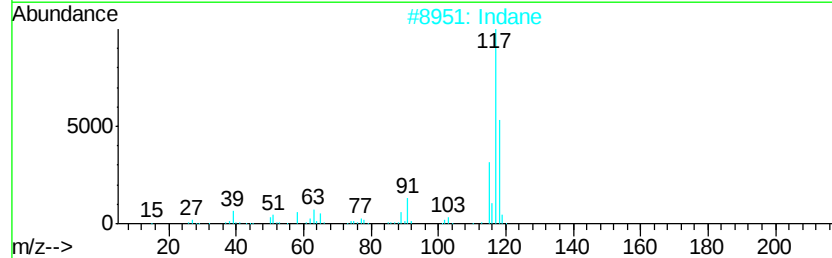
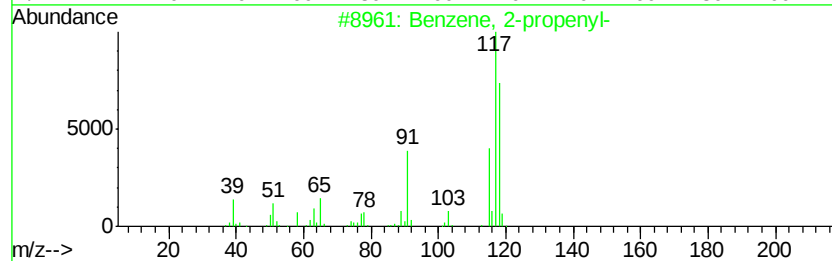
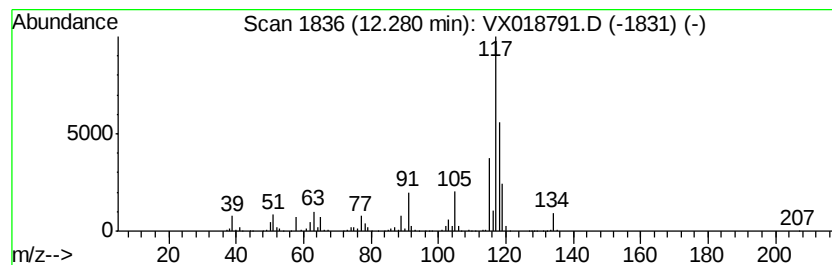
Instrument :
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ClientSampleId :
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 15 Benzene, 2-propenyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	2.21 ug/L	528593	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-propenyl-	118 C9H10	000300-57-2 70
2		Indane	118 C9H10	000496-11-7 70
3		Benzene, 2-propenyl-	118 C9H10	000300-57-2 64
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
5		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3Q7

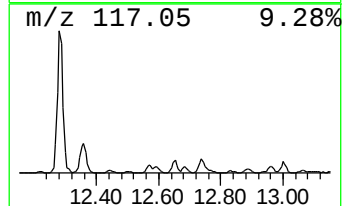
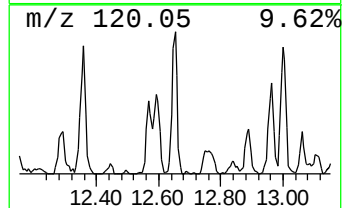
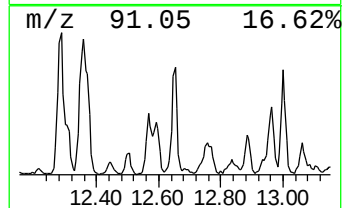
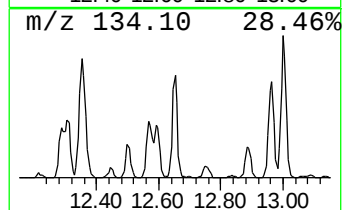
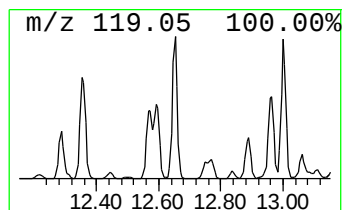
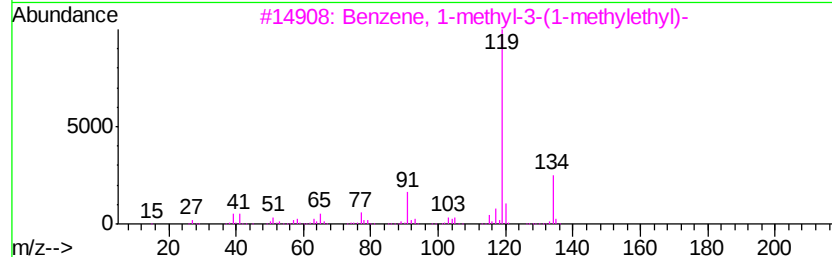
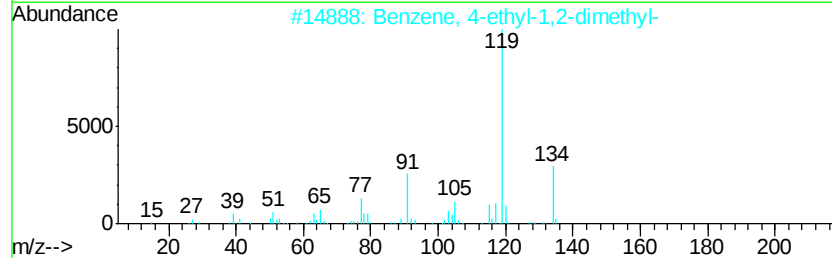
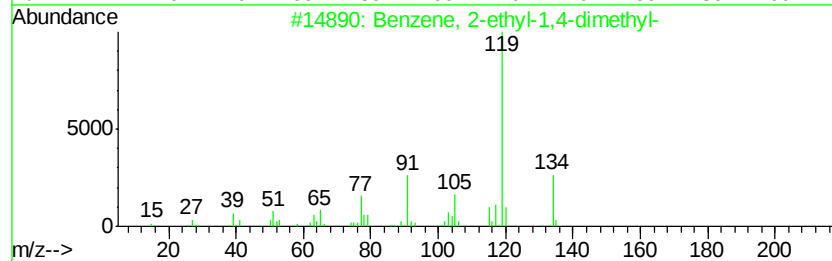
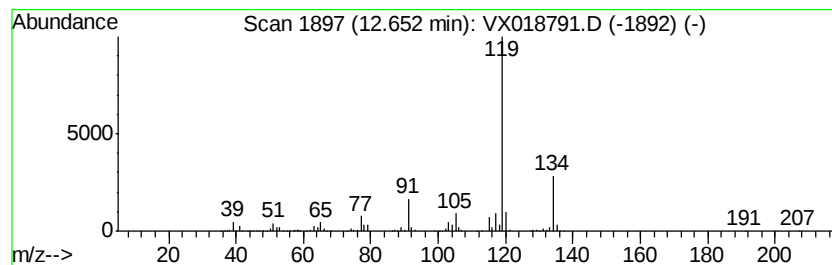
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	0.92 ug/L	219388	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	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 X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3Q7

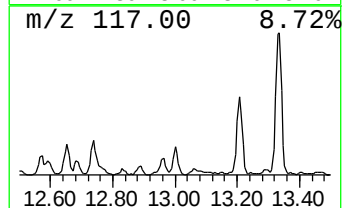
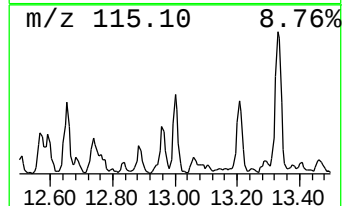
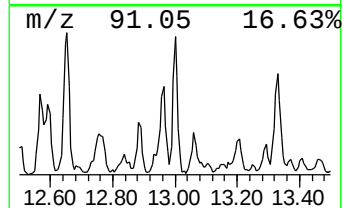
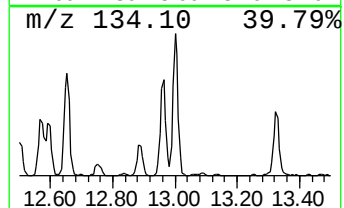
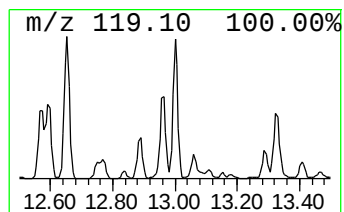
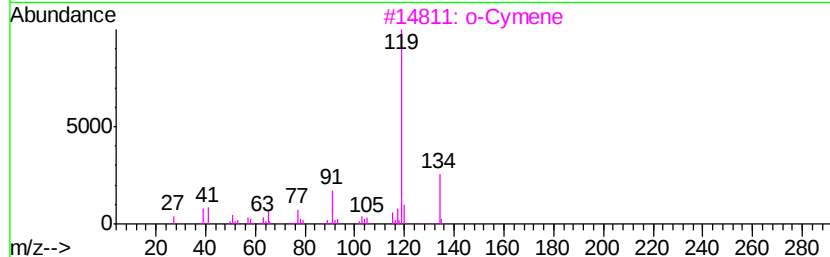
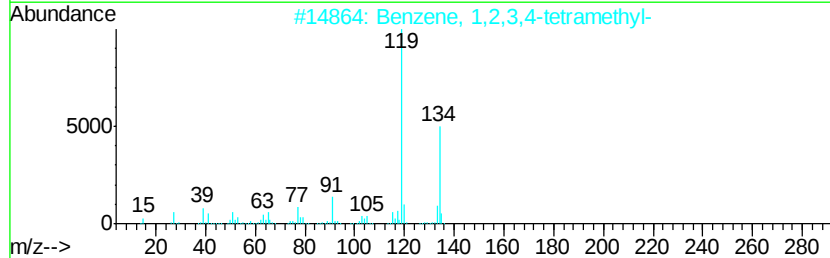
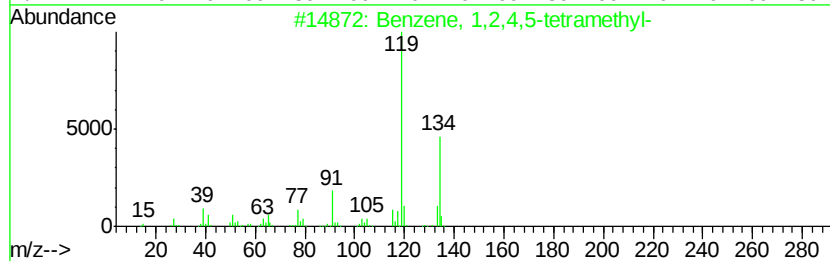
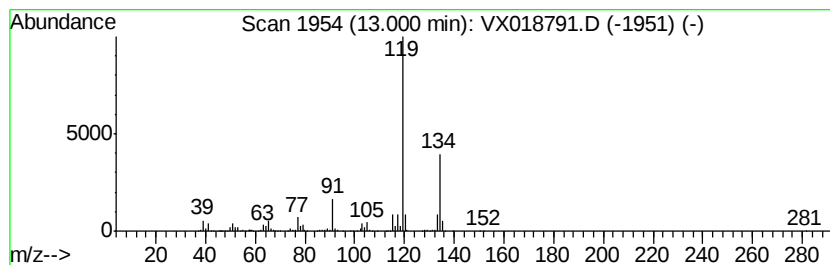
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 17 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	0.95 ug/L	228090	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	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018791.D
Acq On : 06 Oct 2020 11:43
Operator : JC/SP
Sample : L4240-01
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3Q7

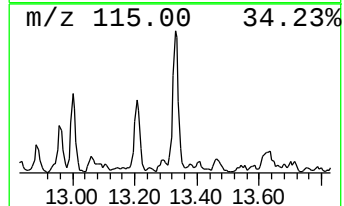
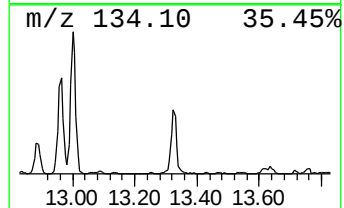
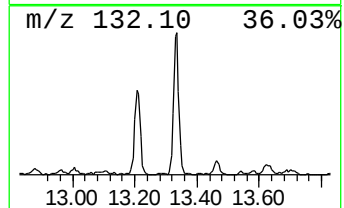
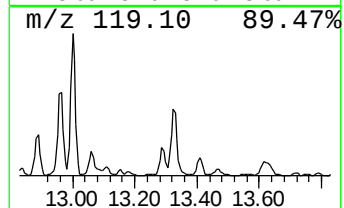
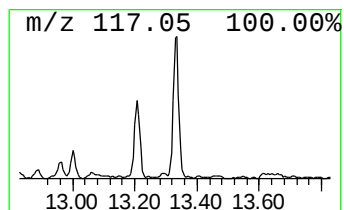
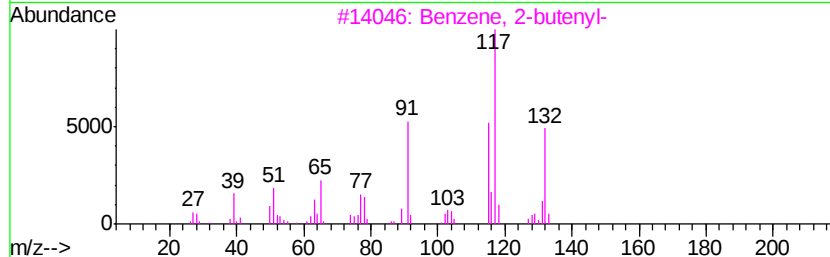
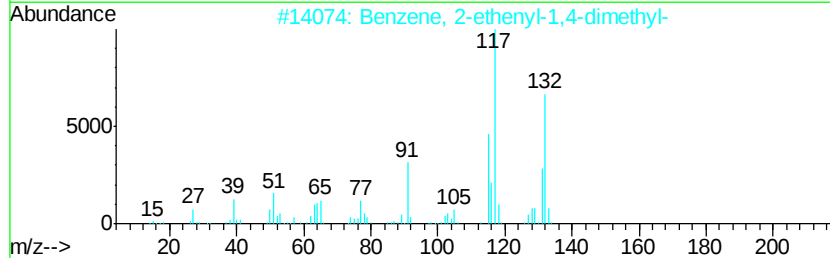
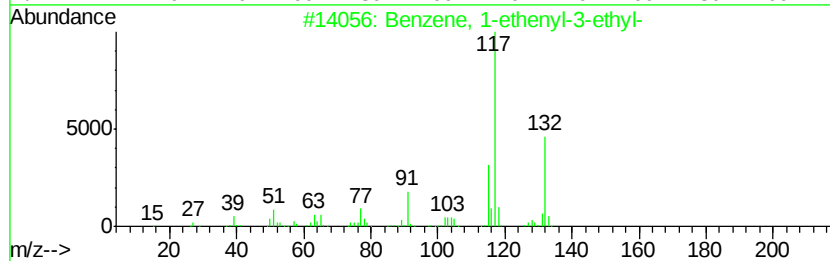
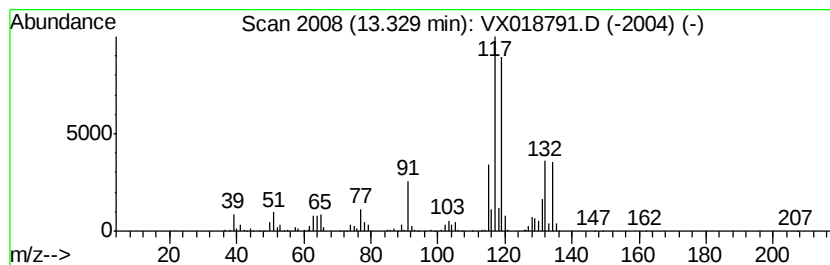
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 18 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	0.90 ug/L	214679	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	84
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	78
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	60



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100620\
 Data File : VX018791.D
 Acq On : 06 Oct 2020 11:43
 Operator : JC/SP
 Sample : L4240-01
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3Q7

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	16.2	ug/L	1026820	1	6.83	316715	5.0
(DEL) Alkane: Str...	2.87	96.4	ug/L	6106800	1	6.83	316715	5.0
(DEL) Alkane: Str...	3.15	191.2	ug/L	12110100	1	6.83	316715	5.0
(DEL) Alkane: Str...	3.51	277.7	ug/L	17591300	1	6.83	316715	5.0
(DEL) Alkane: Str...	3.73	3.7	ug/L	232321	1	6.83	316715	5.0
(DEL) Alkane: Str...	4.12	16.1	ug/L	1020650	1	6.83	316715	5.0
(DEL) Alkane: Str...	4.28	12.2	ug/L	774795	1	6.83	316715	5.0
(DEL) Alkane: Cyc...	4.38	293.1	ug/L	18562900	1	6.83	316715	5.0
Benzene, propyl-	11.34	3.2	ug/L	758709	3	12.06	1195320	5.0
Benzene, 1-ethyl-...	11.42	15.4	ug/L	3691200	3	12.06	1195320	5.0
Benzene, 1,2,3-tr...	11.49	5.3	ug/L	1270860	3	12.06	1195320	5.0
4-Heptanone, 2,6-...	11.61	3.7	ug/L	894212	3	12.06	1195320	5.0
Benzene, 1,2,4-tr...	11.65	4.5	ug/L	1084560	3	12.06	1195320	5.0
Mesitylene	11.79	16.4	ug/L	3918160	3	12.06	1195320	5.0
Benzene, 2-propenyl-	12.28	2.2	ug/L	528593	3	12.06	1195320	5.0
Benzene, 2-ethyl-...	12.65	0.9	ug/L	219388	3	12.06	1195320	5.0
Benzene, 1,2,4,5-...	13.00	0.9	ug/L	228090	3	12.06	1195320	5.0
Benzene, 1-etheny...	13.33	0.9	ug/L	214679	3	12.06	1195320	5.0