

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100520\
 Data File : VX018772.D
 Acq On : 05 Oct 2020 17:54
 Operator : JC/SP
 Sample : L4239-03DL 100X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N2DL

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	1.386	45	49	57	rBV	72800	76423	1.53%	0.321%
2	1.697	95	100	108	rBV	58356	68858	1.38%	0.289%
3	2.343	200	206	214	rBV	127962	208011	4.18%	0.873%
4	2.867	288	292	302	rVB2	37588	76142	1.53%	0.319%
5	3.148	329	338	349	rBV	83747	195895	3.93%	0.822%
6	3.501	386	396	407	rBV	118610	274181	5.51%	1.150%
7	3.672	415	424	432	rBV	40423	103018	2.07%	0.432%
8	4.373	527	539	551	rBV	192149	552836	11.10%	2.319%
9	4.562	556	570	583	rBV3	326542	979862	19.68%	4.110%
10	5.141	653	665	677	rBV2	80094	239789	4.82%	1.006%
11	5.464	704	718	728	rBV2	291829	922258	18.52%	3.869%
12	5.550	728	732	742	rVB3	24700	61053	1.23%	0.256%
13	6.043	798	813	827	rBV2	193481	563800	11.32%	2.365%
14	6.830	931	942	953	rBV	148433	358274	7.20%	1.503%
15	7.366	1020	1030	1043	rBV	120850	263302	5.29%	1.104%
16	8.372	1188	1195	1208	rBV	77624	131379	2.64%	0.551%
17	8.695	1239	1248	1253	rBV	296047	464528	9.33%	1.949%
18	8.762	1253	1259	1276	rVB	3251944	4979199	100.00%	20.886%
19	8.994	1291	1297	1303	rBV	53272	79272	1.59%	0.333%
20	9.427	1362	1368	1378	rBV	393485	568794	11.42%	2.386%
21	10.098	1472	1478	1491	rVB	321791	452034	9.08%	1.896%
22	10.232	1494	1500	1506	rBV	285558	388486	7.80%	1.630%
23	10.341	1512	1518	1527	rVB	1119395	1478339	29.69%	6.201%
24	10.683	1568	1574	1582	rBV	584569	736683	14.80%	3.090%
25	11.000	1621	1626	1636	rVB	84599	110023	2.21%	0.462%
26	11.232	1659	1664	1673	rVB	131239	166684	3.35%	0.699%
27	11.341	1675	1682	1689	rBV	204913	257980	5.18%	1.082%
28	11.421	1689	1695	1702	rVV	1048464	1766390	35.48%	7.409%
29	11.488	1702	1706	1713	rVB	454355	572553	11.50%	2.402%
30	11.610	1721	1726	1730	rBV	1553264	1861580	37.39%	7.809%
31	11.652	1730	1733	1740	rVB	523199	625799	12.57%	2.625%
32	11.792	1750	1756	1764	rBV	1740054	2081627	41.81%	8.732%
33	12.061	1795	1800	1806	rBV	350751	442210	8.88%	1.855%
34	12.128	1806	1811	1816	rVB	520302	619453	12.44%	2.598%

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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	12.280	1831	1836	1844	rBV	221194	297683	5.98%	1.249%
36	12.359	1844	1849	1857	rVB	399982	535323	10.75%	2.245%
37	13.000	1951	1954	1962	rVB	40189	51827	1.04%	0.217%
38	13.329	2002	2008	2014	rVB2	38868	56991	1.14%	0.239%
39	13.816	2081	2088	2094	rBV	138858	171572	3.45%	0.720%

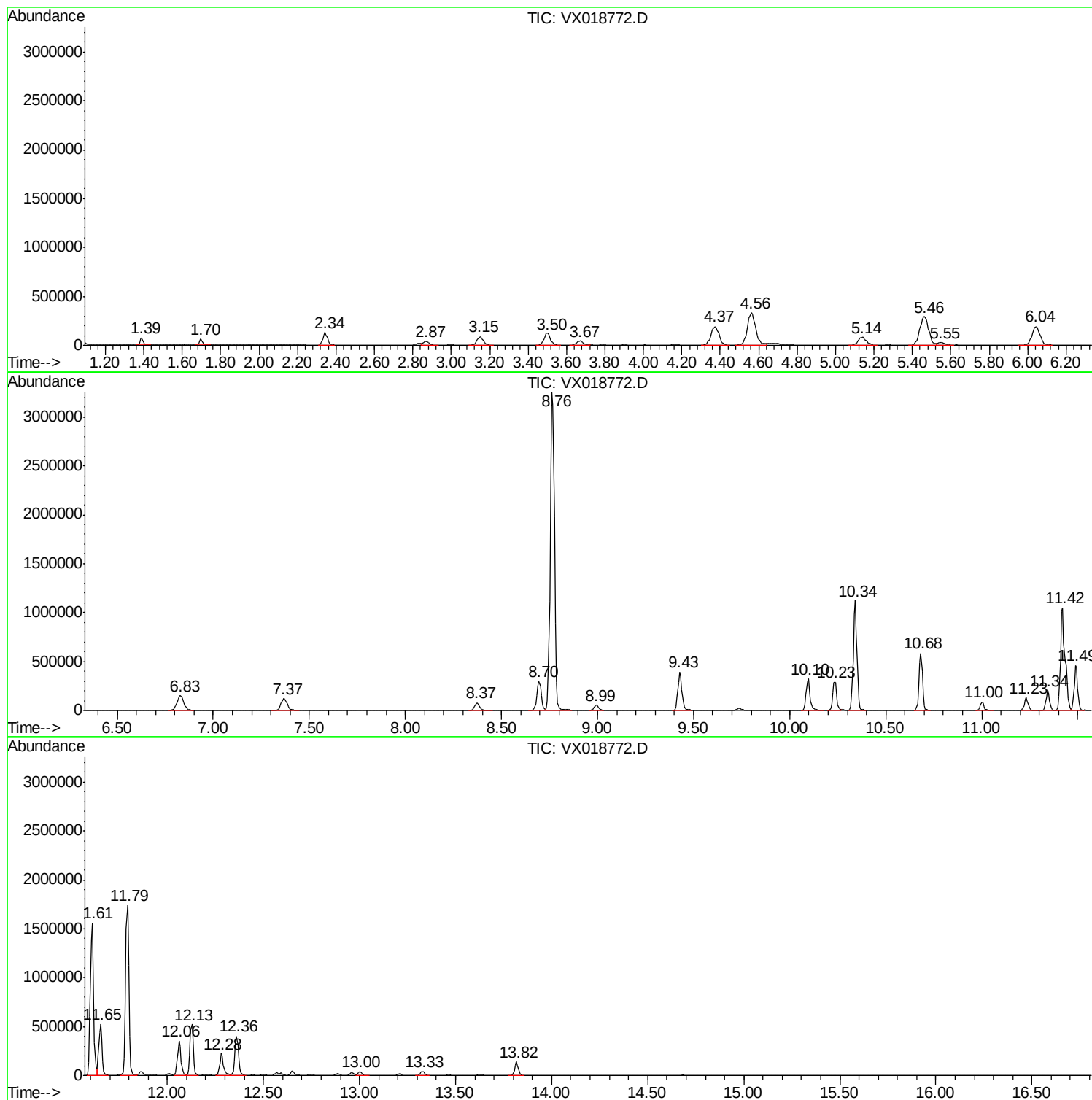
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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



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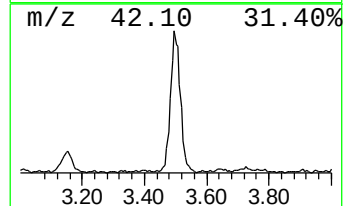
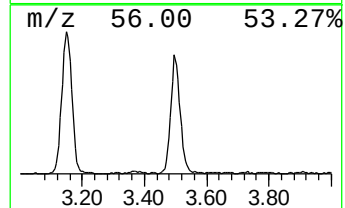
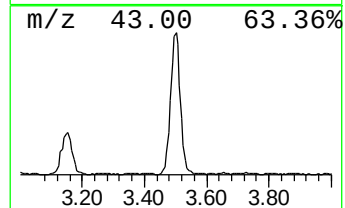
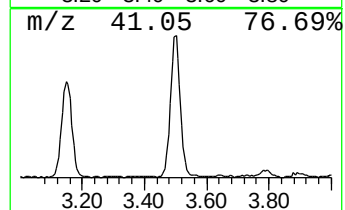
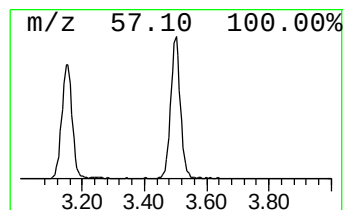
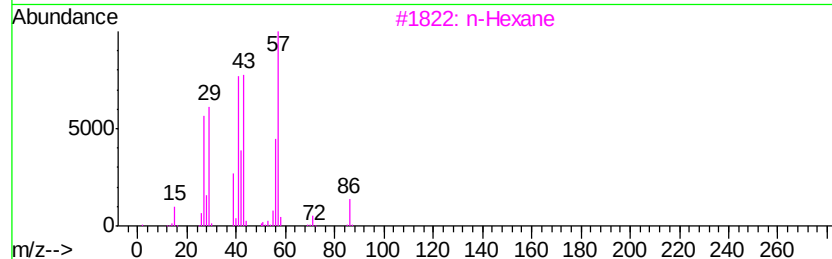
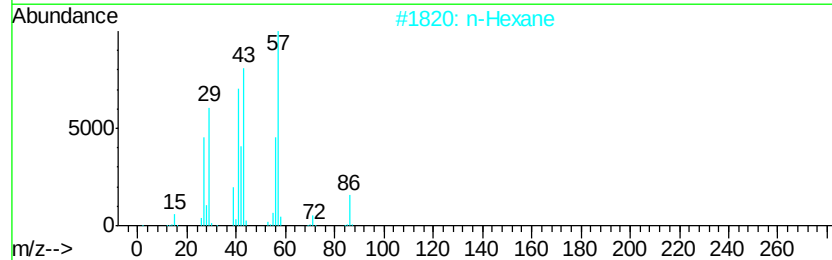
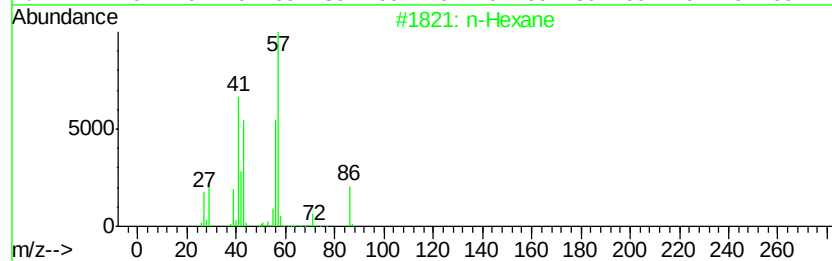
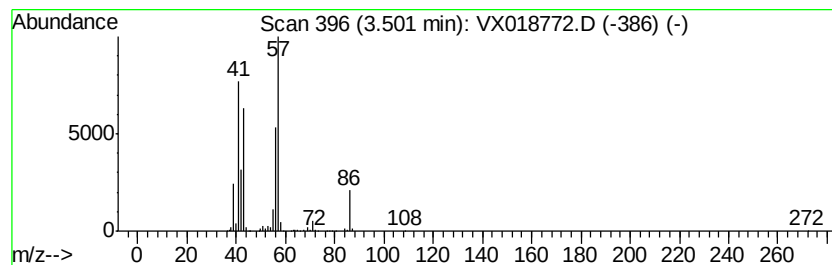
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	3.83 ug/L	274181	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	64
4		Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	53
5		2(3H)-Furanone, dihydro-3-methyl-	100	C5H8O2	001679-47-6	47



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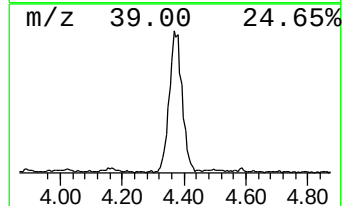
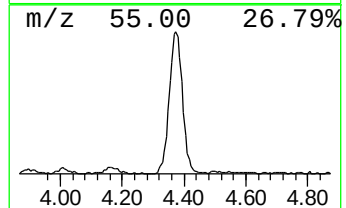
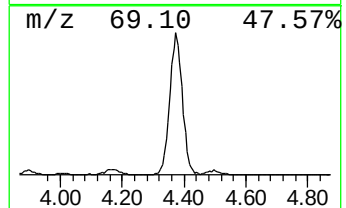
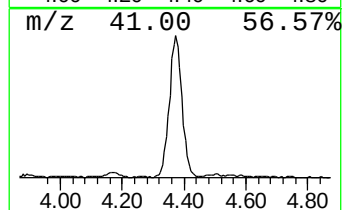
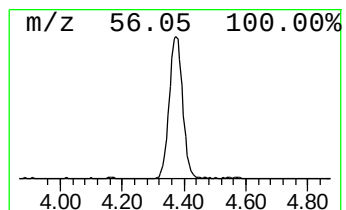
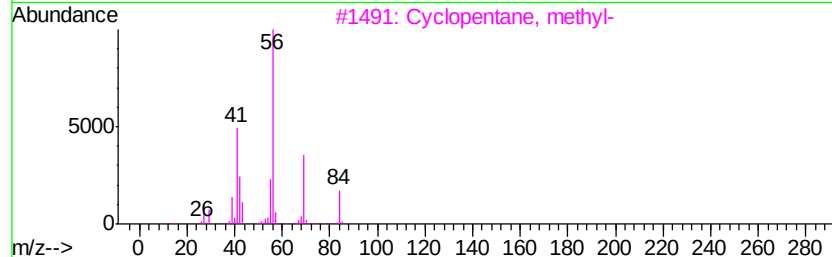
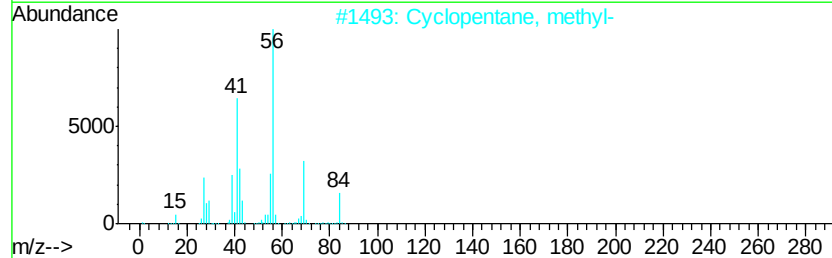
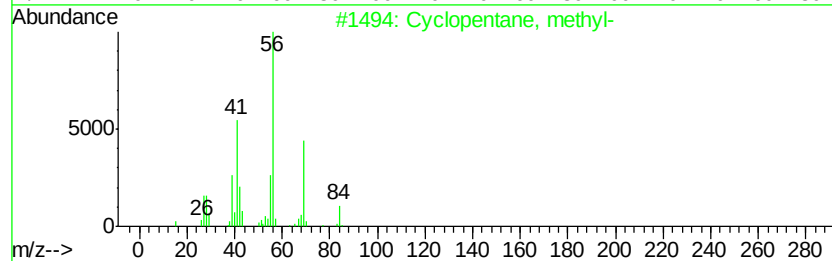
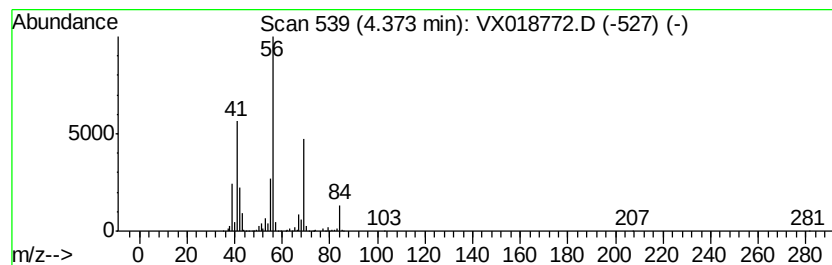
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 (DEL) Alkane: Cyclic4.37 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	7.72 ug/L	552836	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		Cyclopentane, methyl-	84	C6H12	000096-37-7	86
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80



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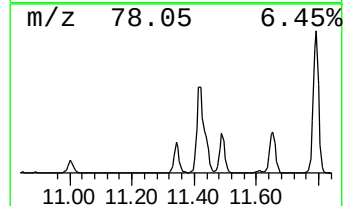
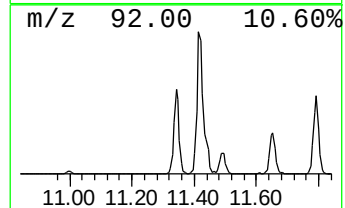
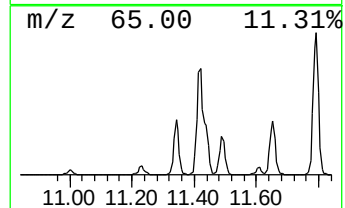
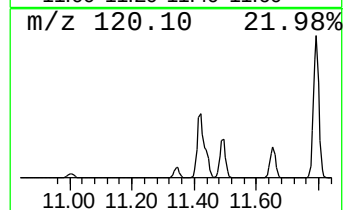
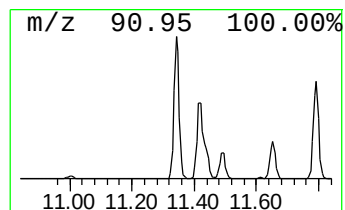
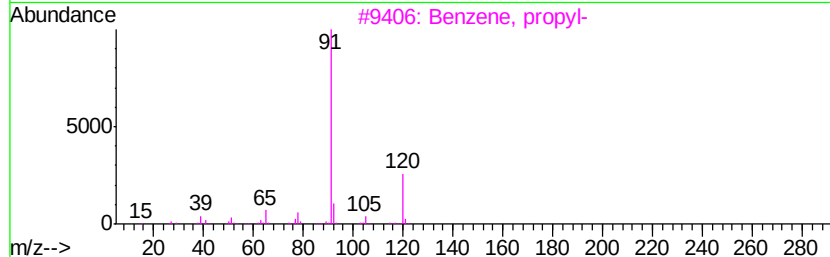
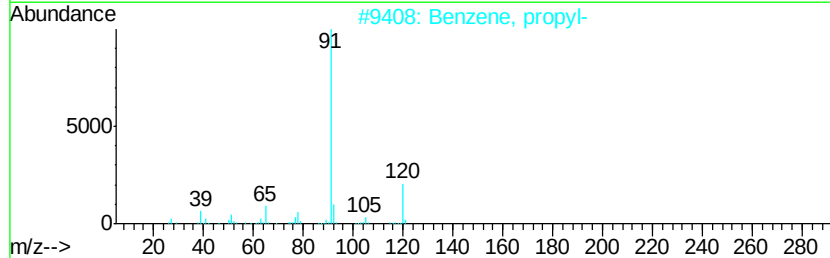
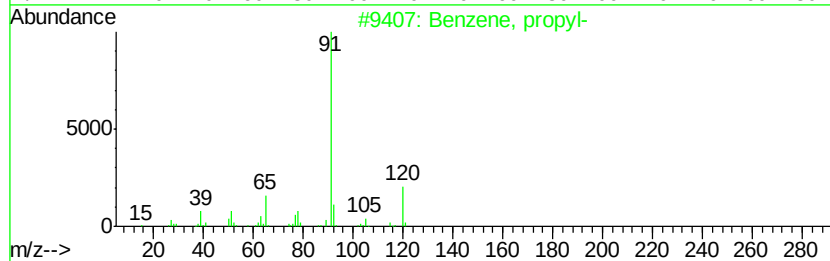
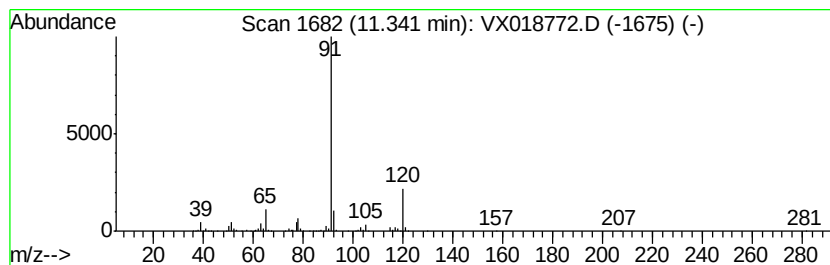
Instrument :
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ClientSampleId :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	2.92 ug/L	257980	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, propyl-	120 C9H12	000103-65-1	93
2	Benzene, propyl-	120 C9H12	000103-65-1	87
3	Benzene, propyl-	120 C9H12	000103-65-1	83
4	1,2-Ethanediamine, N-(phenylmeth...	150 C9H14N2	004152-09-4	72
5	N-Benzyl-2-phenethylamine	211 C15H17N	003647-71-0	72



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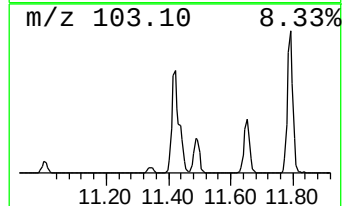
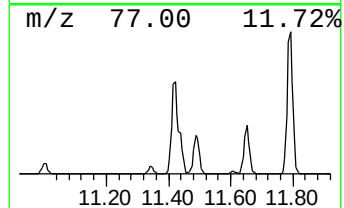
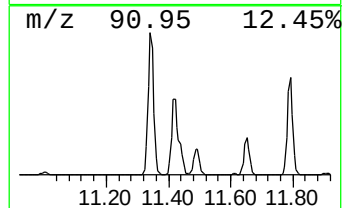
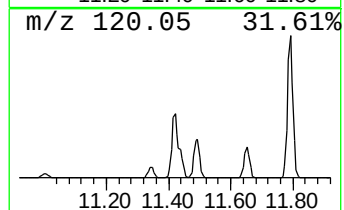
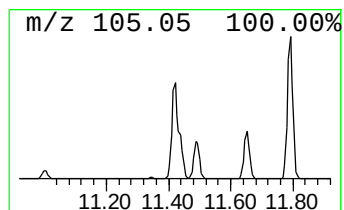
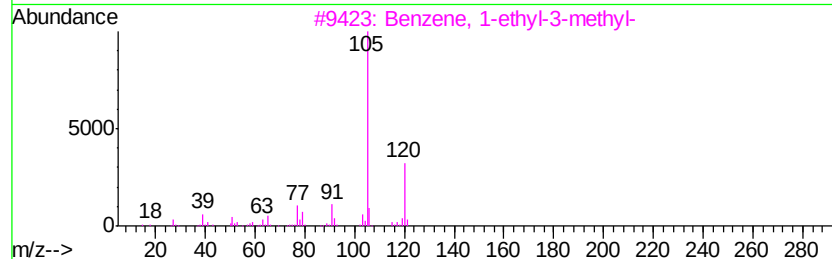
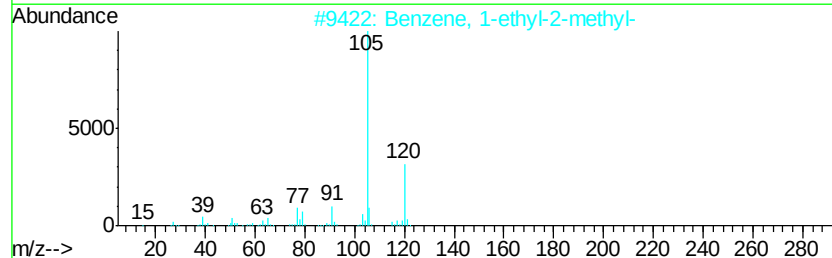
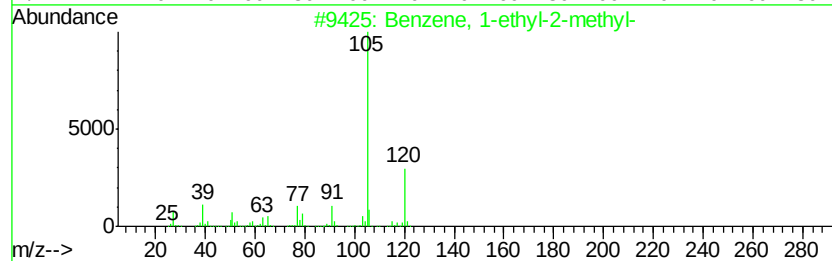
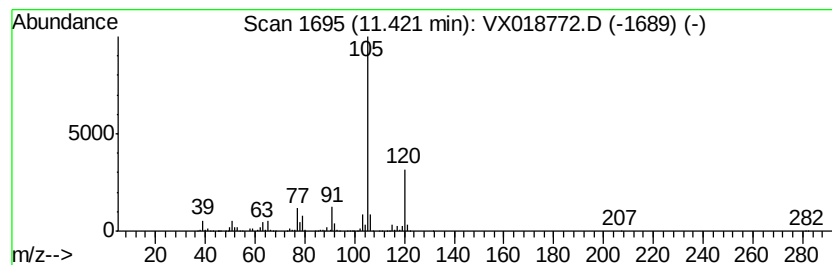
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	19.97 ug/L	1766390	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		Mesitylene	120	C9H12	000108-67-8	91



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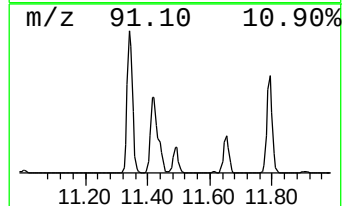
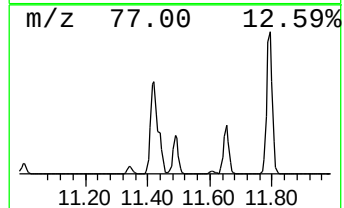
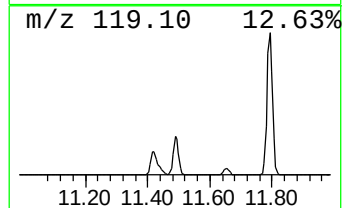
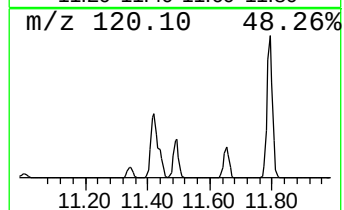
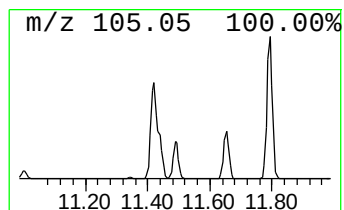
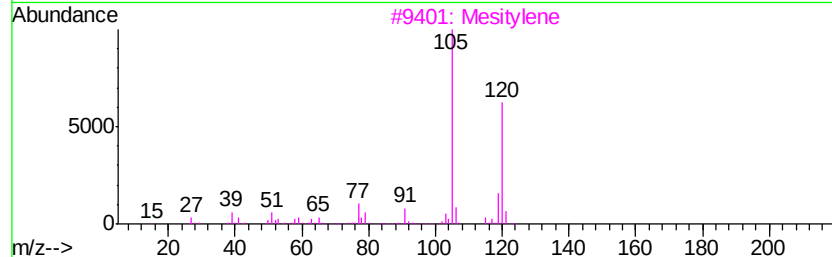
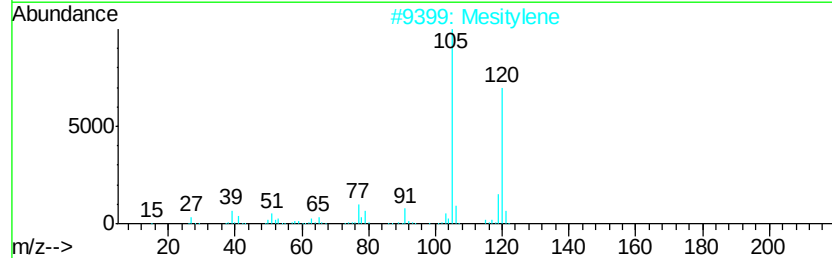
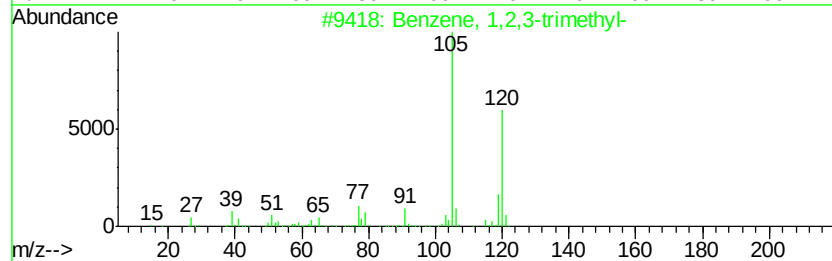
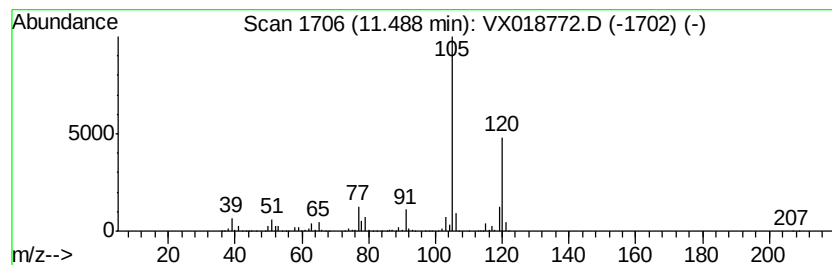
Instrument :
MSVOA_X
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.49	6.47 ug/L	572553	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	Mesitylene	120	C9H12	000108-67-8	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\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

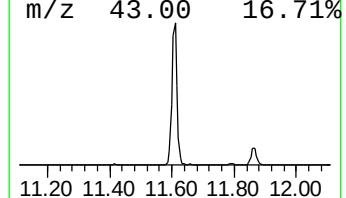
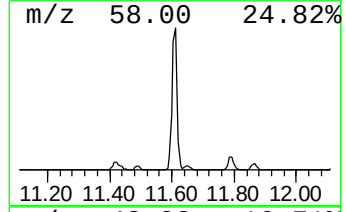
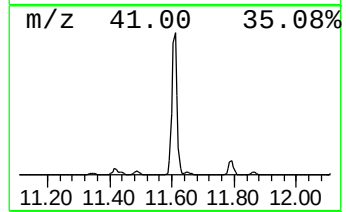
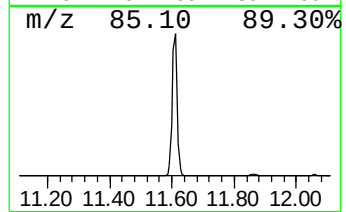
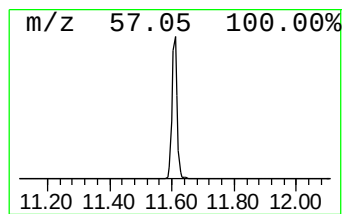
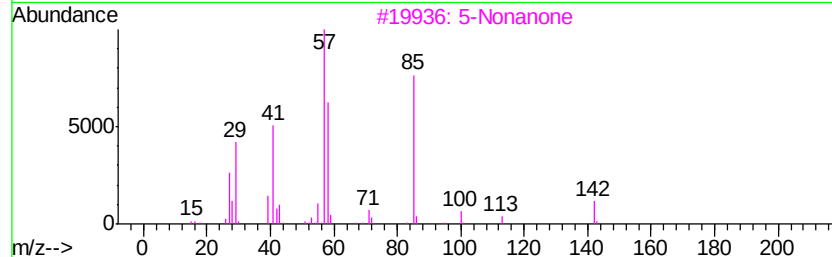
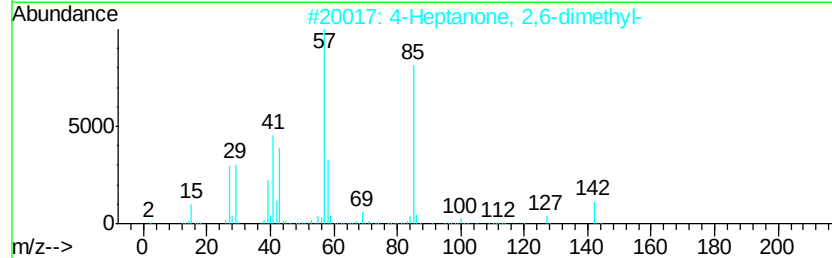
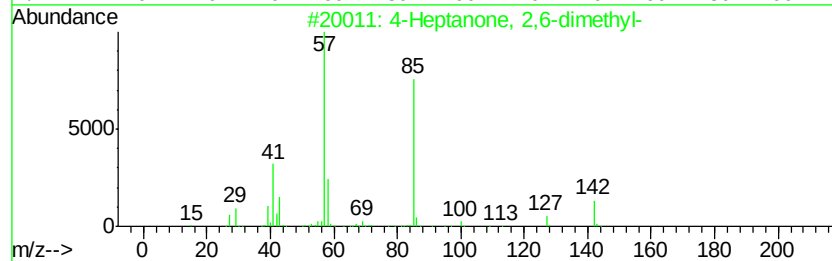
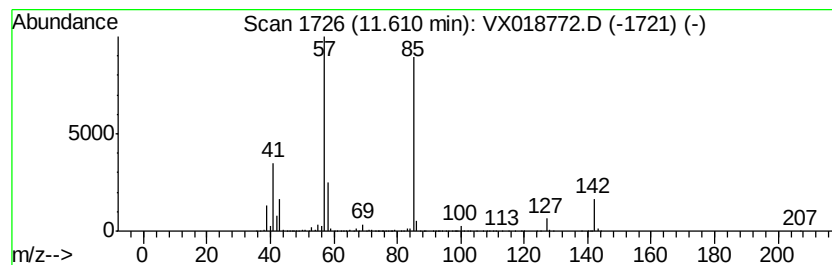
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ClientSampleId :
BG3N2DL

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 7 4-Heptanone, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.61	21.05 ug/L	1861580	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	t-Butyl isobutyl ketone	142	C9H18O	014705-50-1	59
5	5-Nonanone	142	C9H18O	000502-56-7	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N2DL

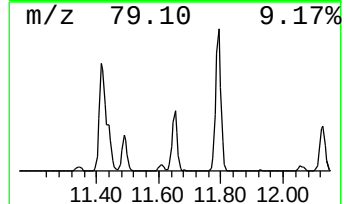
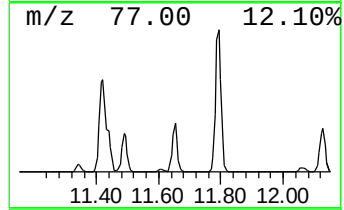
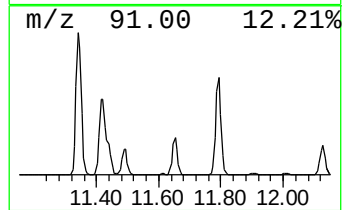
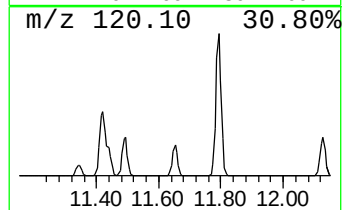
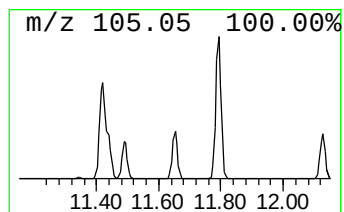
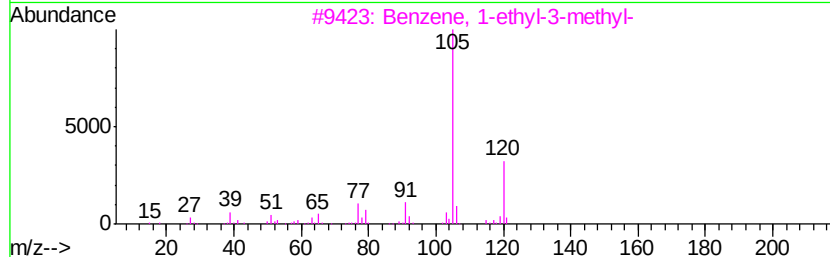
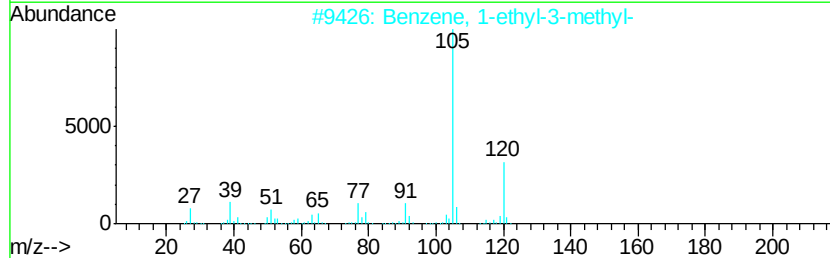
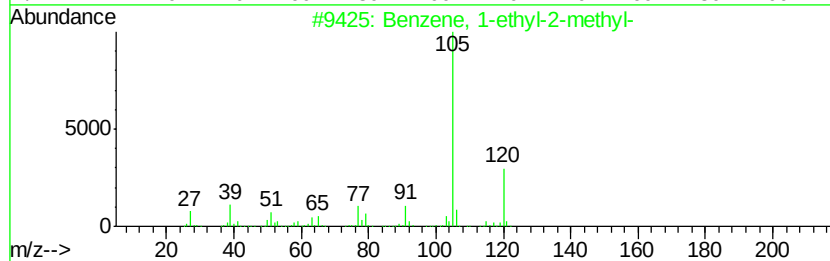
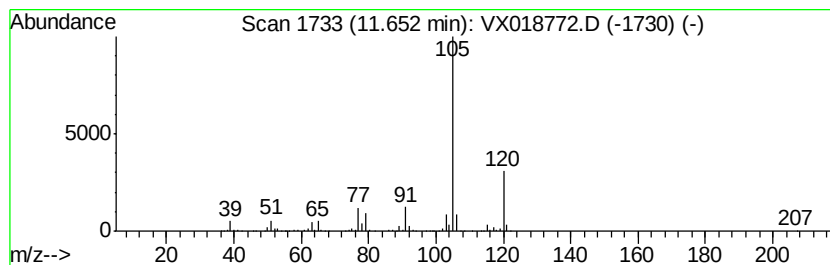
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	7.08 ug/L	625799	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-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N2DL

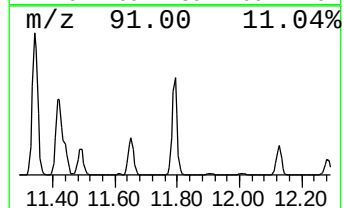
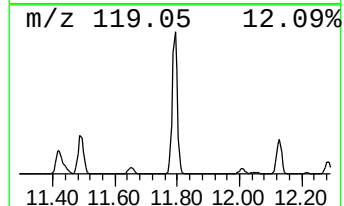
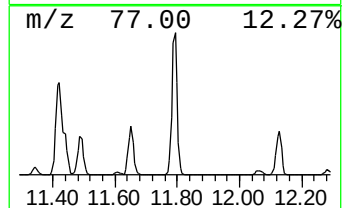
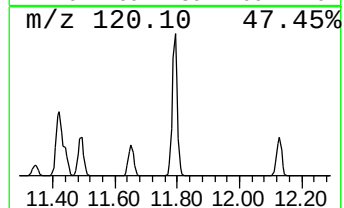
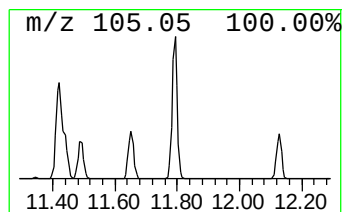
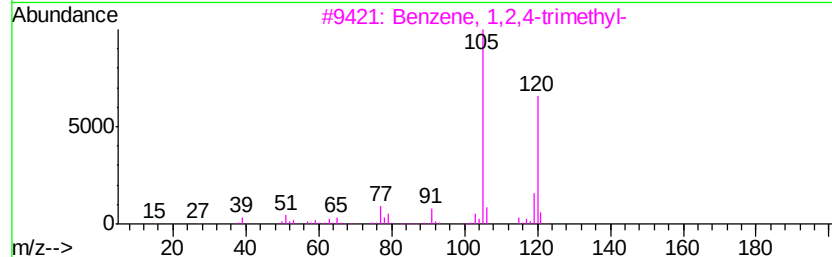
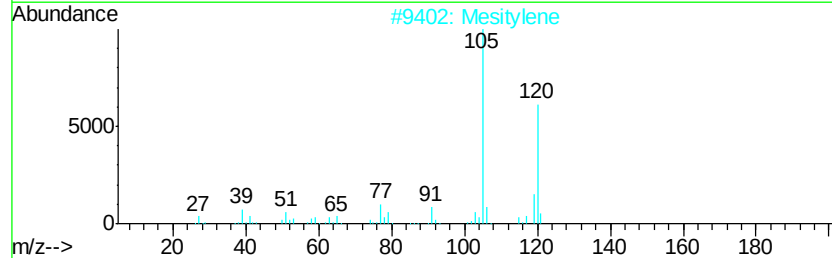
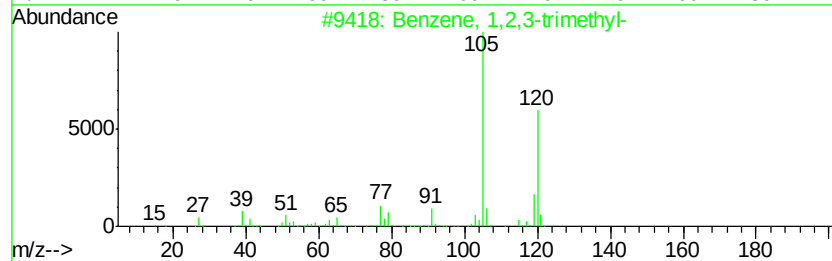
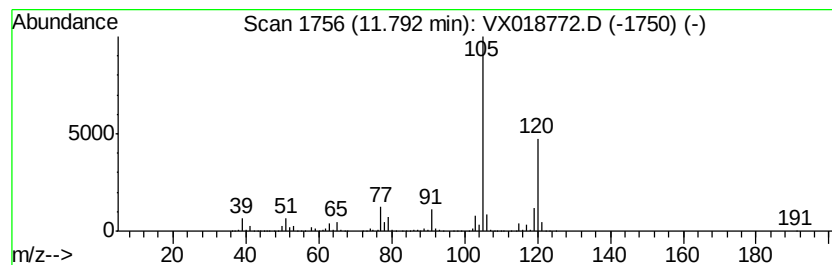
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Mesitylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	23.54 ug/L	2081630	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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

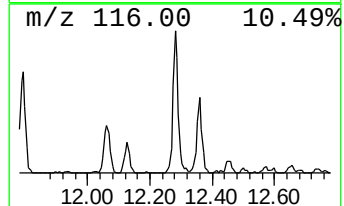
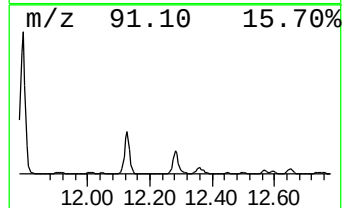
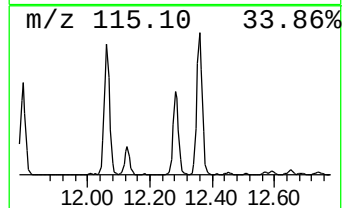
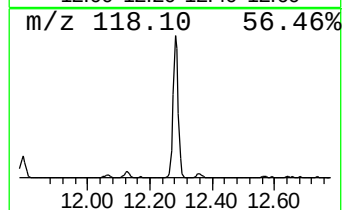
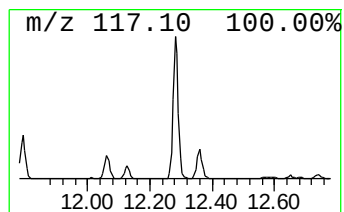
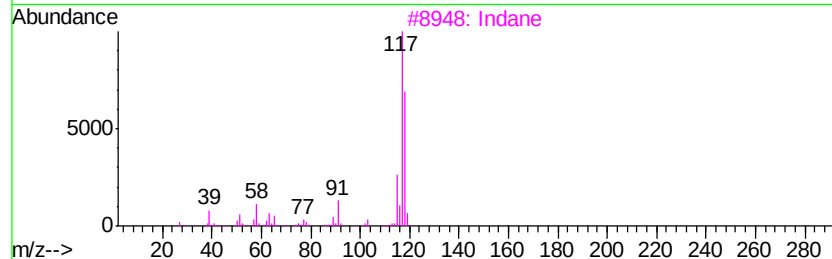
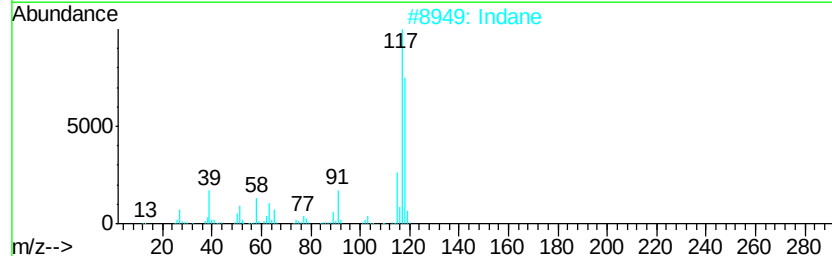
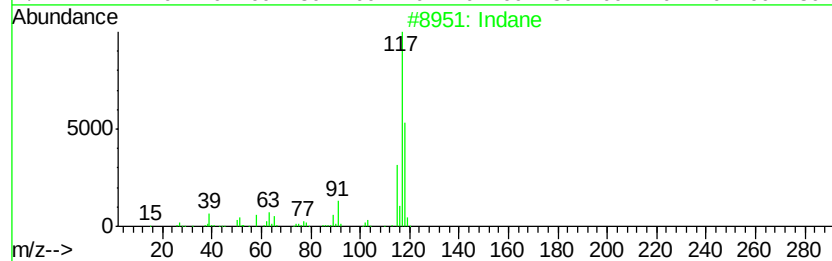
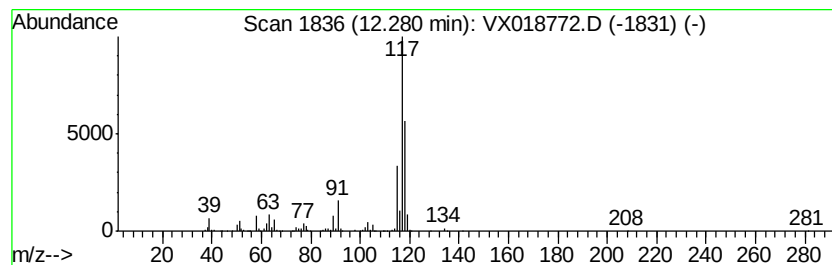
Instrument :
MSVOA_X
ClientSampleId :
BG3N2DL

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 10 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	3.37 ug/L	297683	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 94
2	Indane		118 C9H10	000496-11-7 91
3	Indane		118 C9H10	000496-11-7 87
4	Benzene, cyclopropyl-		118 C9H10	000873-49-4 81
5	Benzene, cyclopropyl-		118 C9H10	000873-49-4 81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N2DL

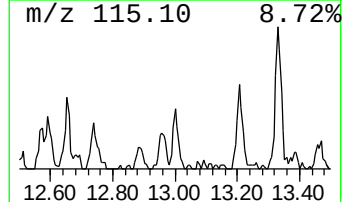
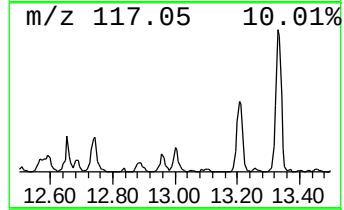
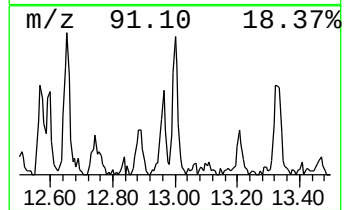
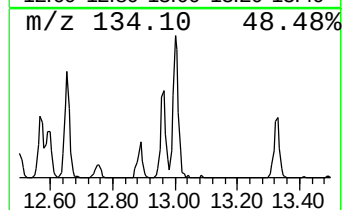
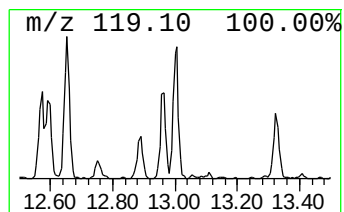
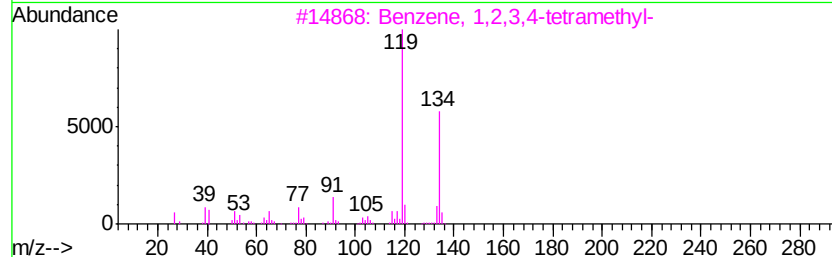
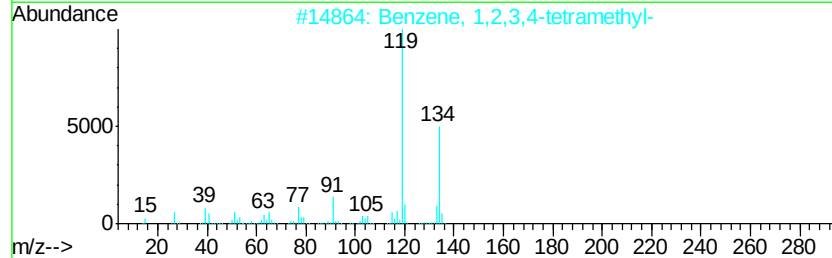
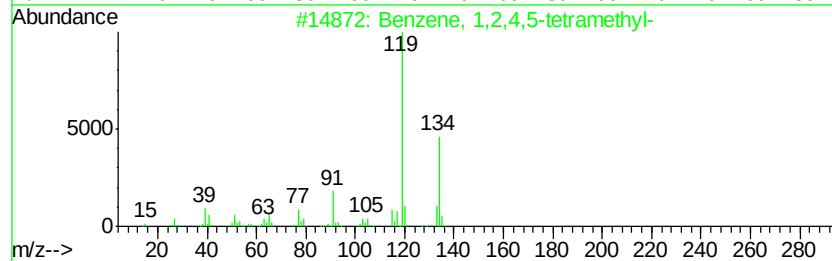
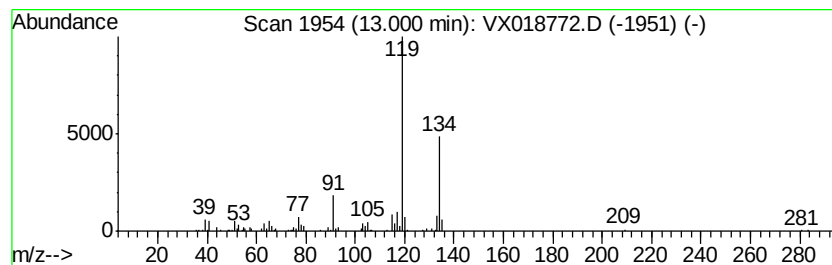
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	0.59 ug/L	51827	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		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N2DL

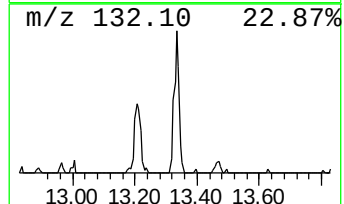
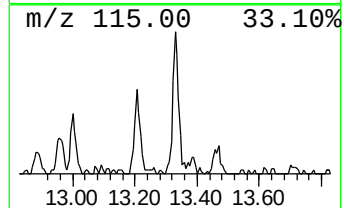
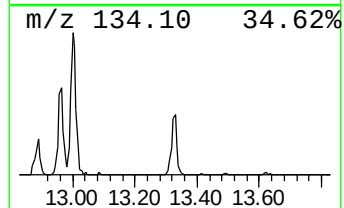
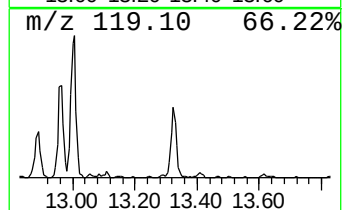
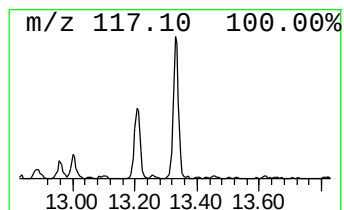
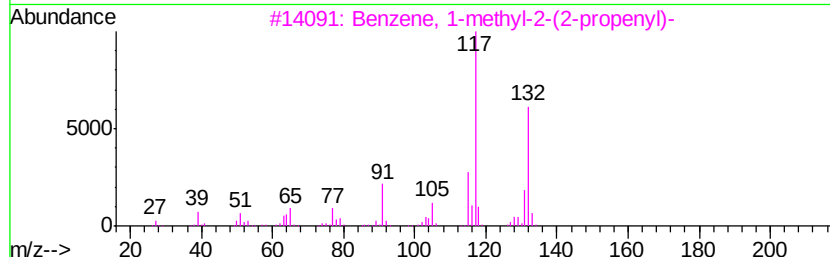
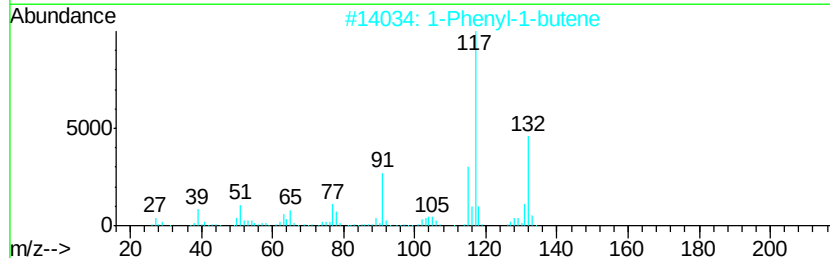
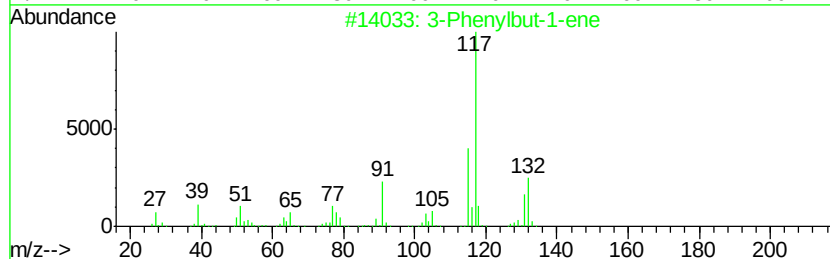
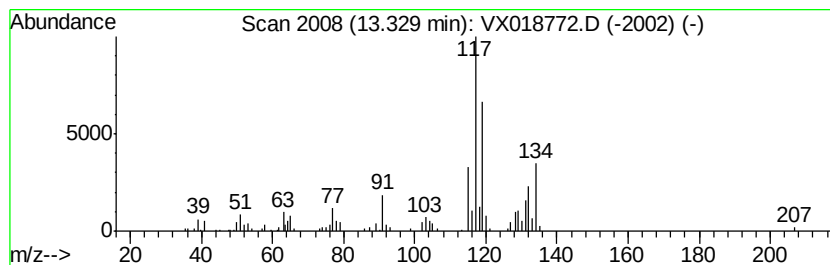
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 3-Phenylbut-1-ene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	0.64 ug/L	56991	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	64
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018772.D
Acq On : 05 Oct 2020 17:54
Operator : JC/SP
Sample : L4239-03DL 100X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N2DL

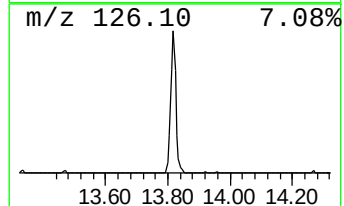
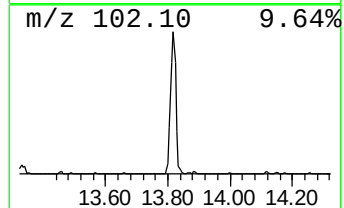
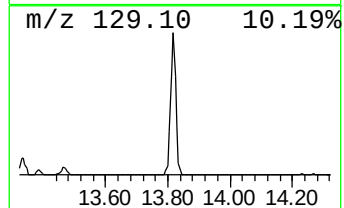
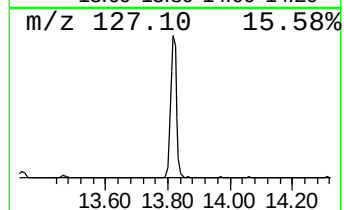
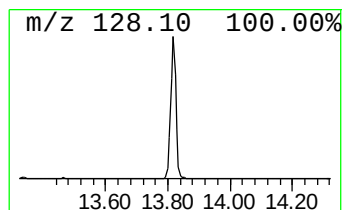
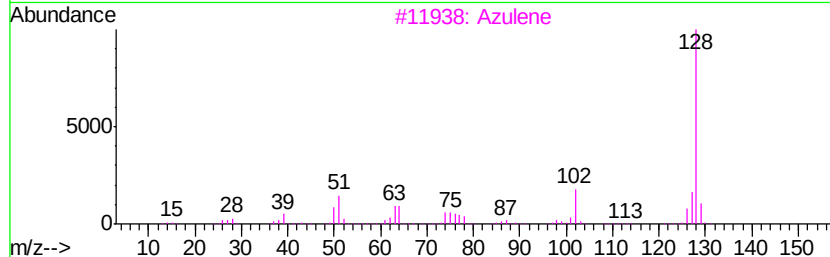
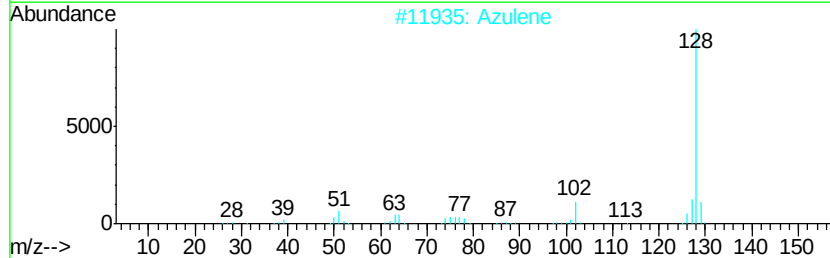
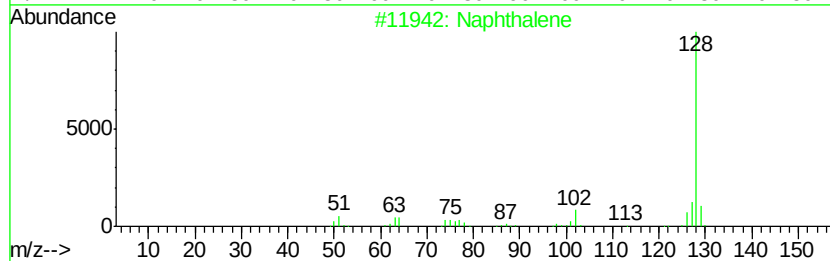
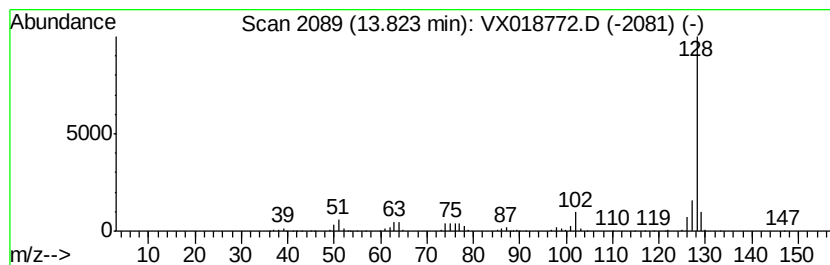
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 Azulene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	1.94 ug/L	171572	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene	128	C10H8	000091-20-3	95
2		Azulene	128	C10H8	000275-51-4	93
3		Azulene	128	C10H8	000275-51-4	91
4		1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
5		Naphthalene	128	C10H8	000091-20-3	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100520\
 Data File : VX018772.D
 Acq On : 05 Oct 2020 17:54
 Operator : JC/SP
 Sample : L4239-03DL 100X
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N2DL

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...	3.50	3.8	ug/L	274181	1	6.83	358274	5.0
(DEL) Alkane: Cyc...	4.37	7.7	ug/L	552836	1	6.83	358274	5.0
Benzene, propyl-	11.34	2.9	ug/L	257980	3	12.06	442210	5.0
Benzene, 1-ethyl-...	11.42	20.0	ug/L	1766390	3	12.06	442210	5.0
Benzene, 1,2,3-tr...	11.49	6.5	ug/L	572553	3	12.06	442210	5.0
4-Heptanone, 2,6-...	11.61	21.1	ug/L	1861580	3	12.06	442210	5.0
Benzene, 1-ethyl-...	11.65	7.1	ug/L	625799	3	12.06	442210	5.0
Mesitylene	11.79	23.5	ug/L	2081630	3	12.06	442210	5.0
Indane	12.28	3.4	ug/L	297683	3	12.06	442210	5.0
Benzene, 1,2,4,5-...	13.00	0.6	ug/L	51827	3	12.06	442210	5.0
3-Phenylbut-1-ene	13.33	0.6	ug/L	56991	3	12.06	442210	5.0
Azulene	13.82	1.9	ug/L	171572	3	12.06	442210	5.0