

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100520\
 Data File : VX018764.D
 Acq On : 05 Oct 2020 14:26
 Operator : JC/SP
 Sample : L4239-02DL 40X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N0DL

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	199	206	220	rBV	121699	208202	1.81%	0.437%
2	2.825	276	285	287	rBV	87352	173151	1.50%	0.364%
3	2.867	287	292	304	rVB	423686	873997	7.58%	1.837%
4	3.148	328	338	353	rBV	809585	1830513	15.87%	3.846%
5	3.495	385	395	408	rBV	1109064	2532068	21.95%	5.321%
6	4.373	528	539	553	rVB	1200634	3451938	29.93%	7.254%
7	4.562	553	570	585	rBV2	366582	1102749	9.56%	2.317%
8	5.135	650	664	679	rBV	76108	238696	2.07%	0.502%
9	5.458	704	717	726	rBV	501325	1557680	13.50%	3.273%
10	5.550	726	732	743	rVB2	119542	339260	2.94%	0.713%
11	6.044	799	813	835	rBV3	183417	552087	4.79%	1.160%
12	6.830	931	942	955	rBV	163715	401080	3.48%	0.843%
13	7.367	1020	1030	1042	rBV	116852	254085	2.20%	0.534%
14	8.373	1188	1195	1204	rBV	77012	127043	1.10%	0.267%
15	8.696	1240	1248	1253	rBV	286786	459385	3.98%	0.965%
16	8.763	1253	1259	1277	rVB	7399988	11534688	100.00%	24.238%
17	9.427	1362	1368	1380	rBV	380905	551189	4.78%	1.158%
18	10.092	1469	1477	1487	rBV	338780	504811	4.38%	1.061%
19	10.232	1494	1500	1506	rBV	379271	497829	4.32%	1.046%
20	10.342	1511	1518	1531	rVB	1383724	1847049	16.01%	3.881%
21	10.683	1567	1574	1586	rBV	735089	942836	8.17%	1.981%
22	11.000	1618	1626	1633	rBV	146723	187508	1.63%	0.394%
23	11.232	1658	1664	1671	rVB	122043	159020	1.38%	0.334%
24	11.342	1677	1682	1689	rBV	483179	593084	5.14%	1.246%
25	11.421	1689	1695	1702	rVV	1948430	3404886	29.52%	7.155%
26	11.488	1702	1706	1720	rVB	921769	1120951	9.72%	2.355%
27	11.610	1720	1726	1730	rBV	3155155	3885335	33.68%	8.164%
28	11.652	1730	1733	1744	rVB	899692	1070855	9.28%	2.250%
29	11.793	1750	1756	1762	rBV	3290767	3852527	33.40%	8.095%
30	12.061	1795	1800	1806	rBV2	410696	562150	4.87%	1.181%
31	12.128	1806	1811	1817	rVB	839551	1003110	8.70%	2.108%
32	12.280	1830	1836	1844	rBV2	341591	495921	4.30%	1.042%
33	12.360	1844	1849	1860	rVB	450024	617515	5.35%	1.298%
34	12.652	1892	1897	1901	rBV	100706	123403	1.07%	0.259%

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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

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Title : TRACE VOA SOM01.0

35	13.000	1951	1954	1961	rVB	100696	121206	1.05%	0.255%
36	13.329	1999	2008	2014	rBV2	90403	131263	1.14%	0.276%
37	13.817	2082	2088	2093	rBV	230335	280879	2.44%	0.590%

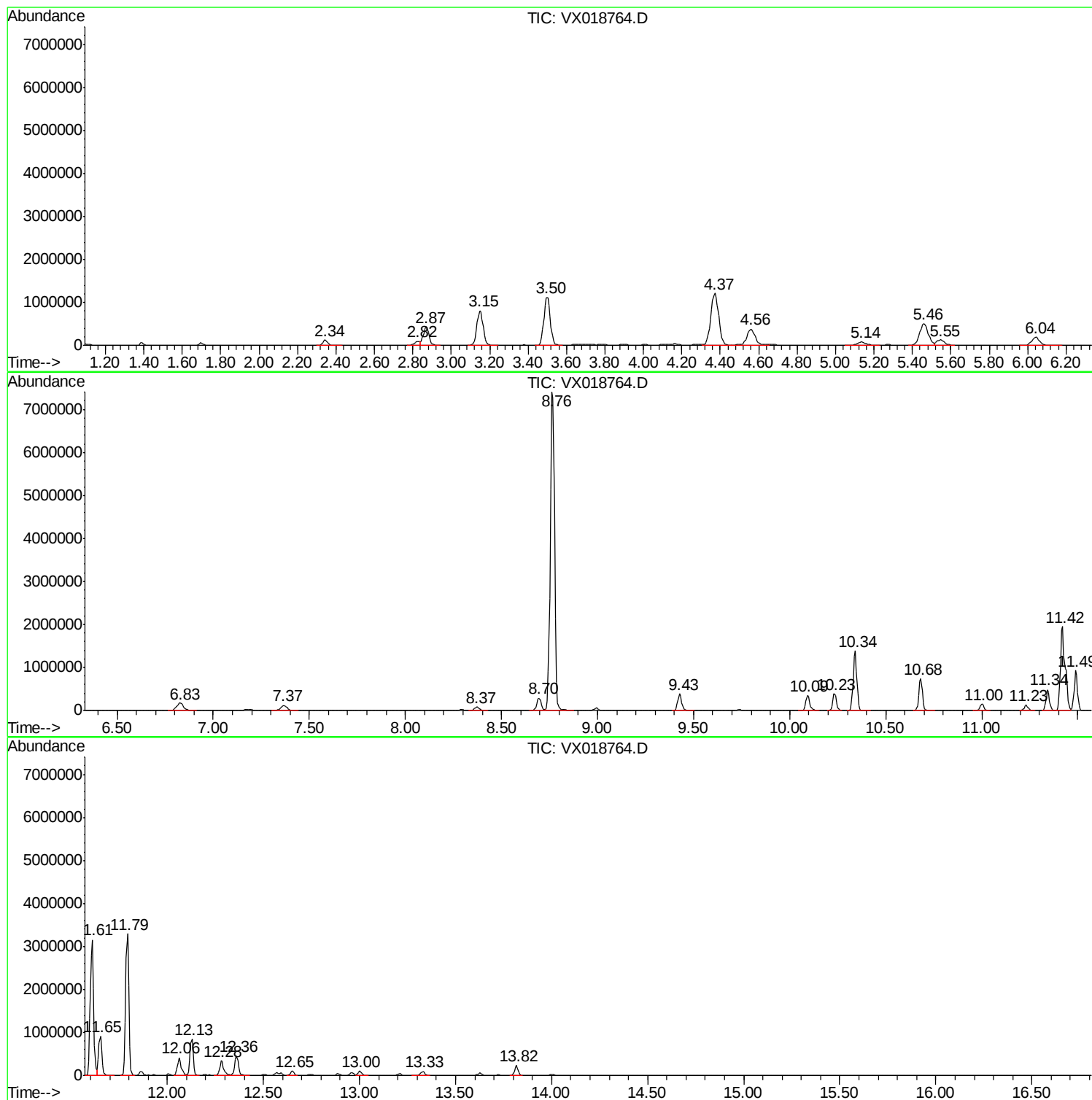
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Quant Title : TRACE VOA SOM01.0

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



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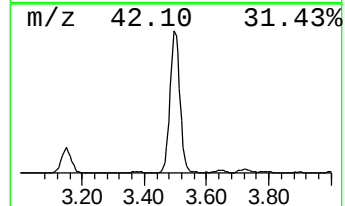
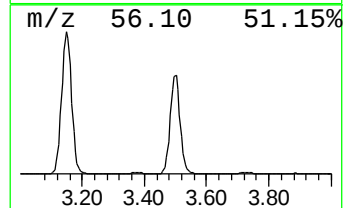
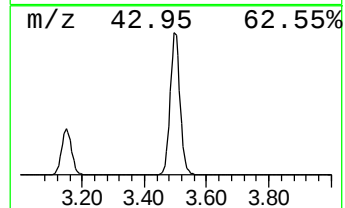
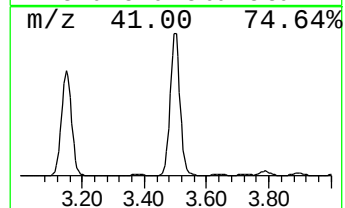
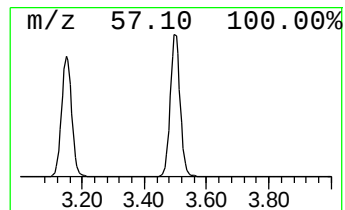
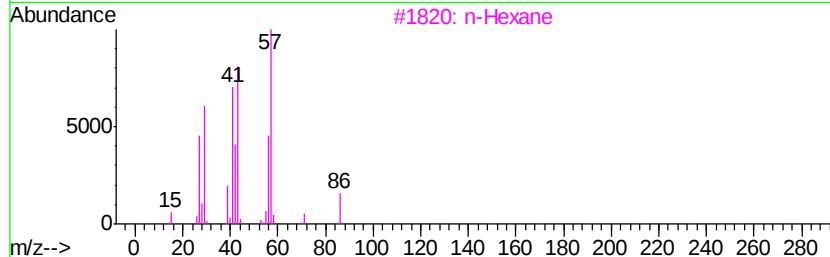
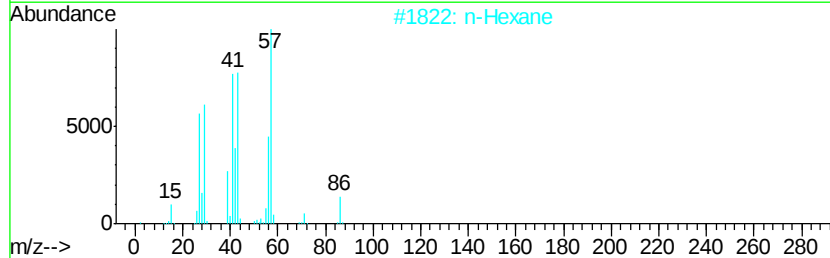
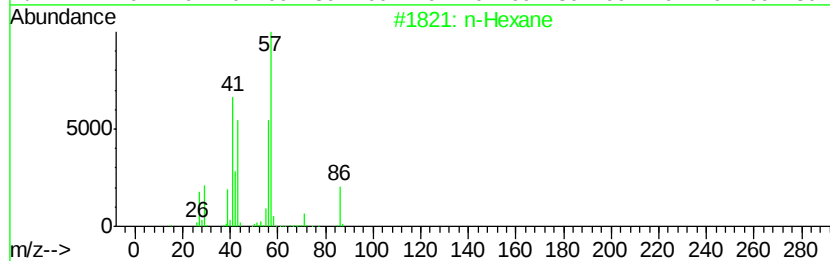
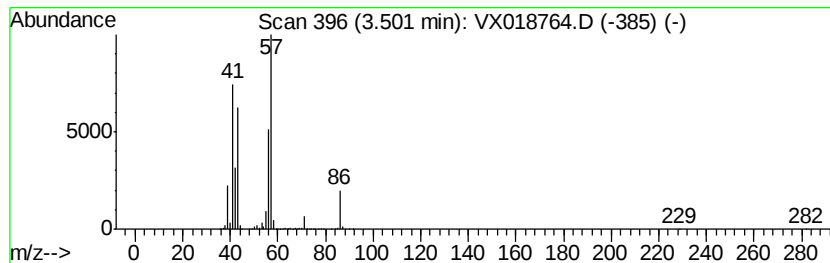
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	31.57 ug/L	2532070	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		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	47
5		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43



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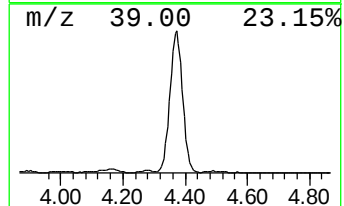
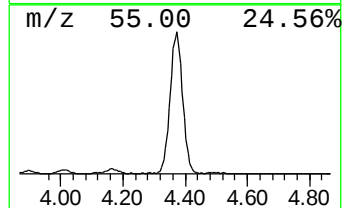
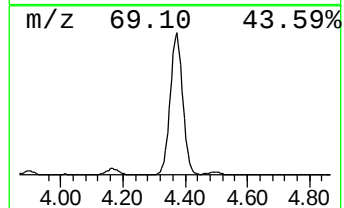
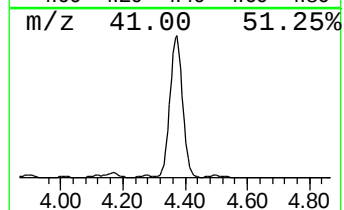
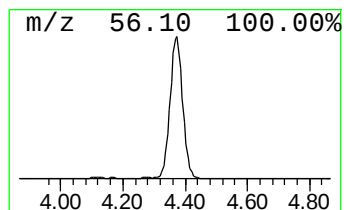
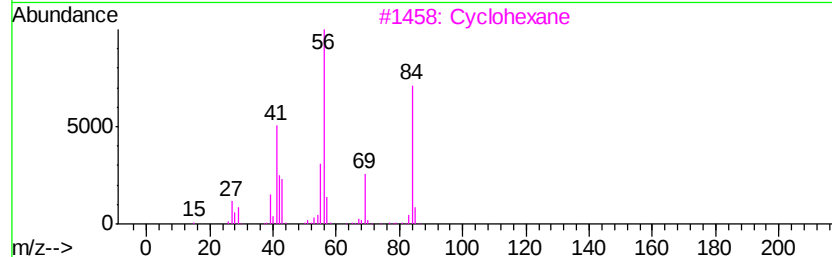
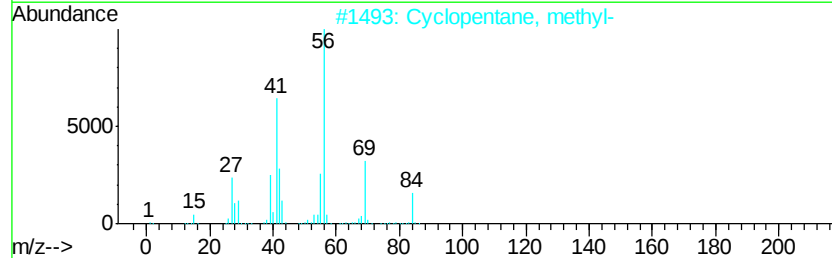
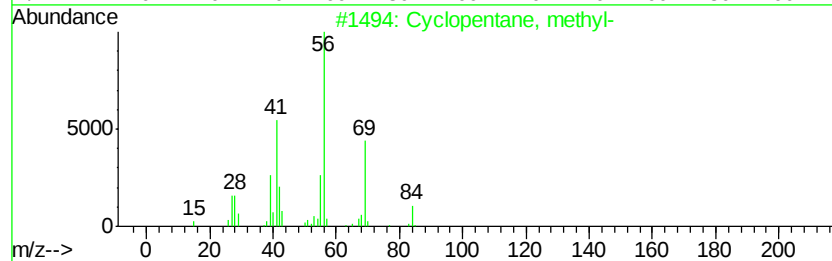
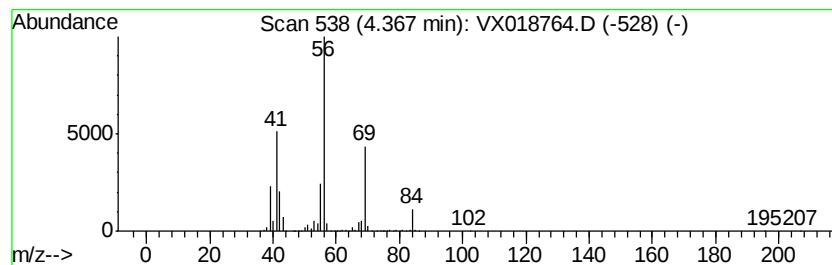
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 (DEL) Alkane: Cyclic4.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	43.03 ug/L	3451940	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		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64



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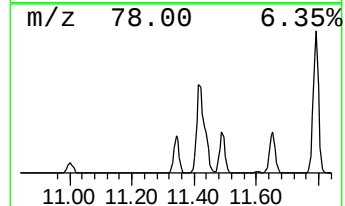
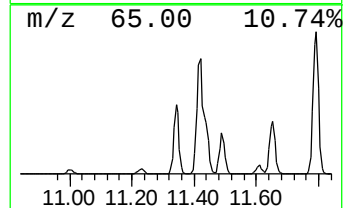
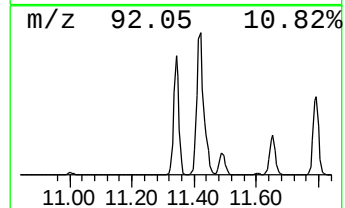
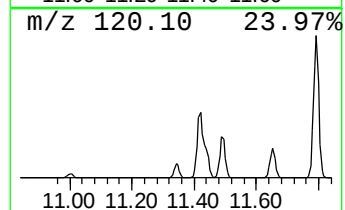
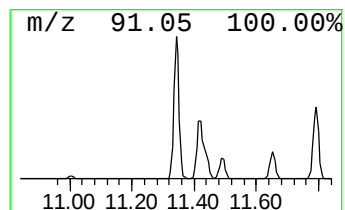
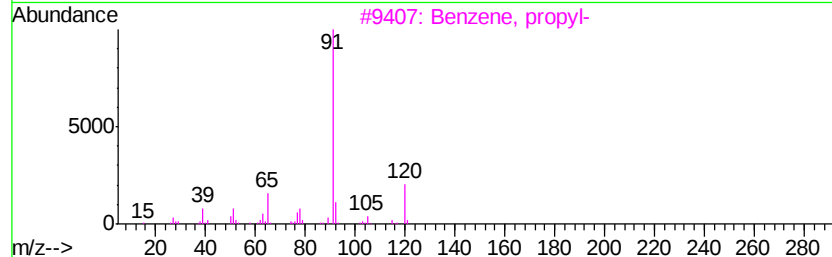
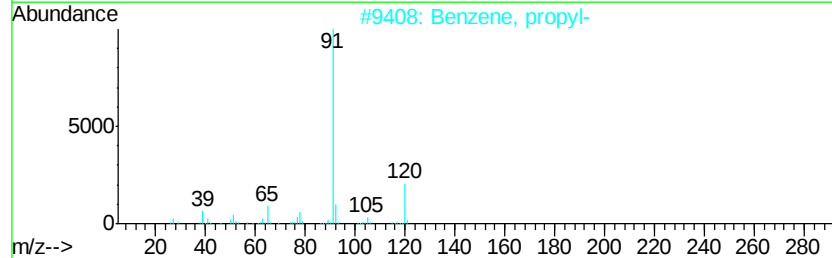
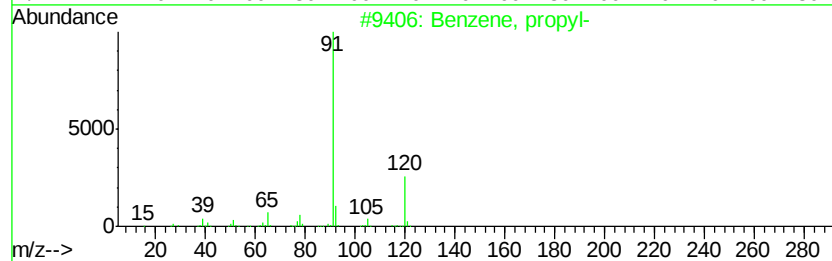
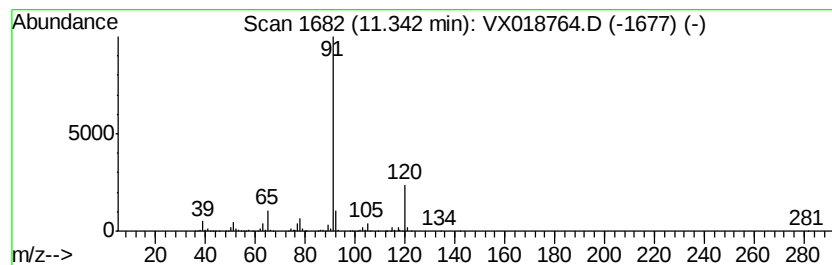
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 Benzene, propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	5.28 ug/L	593084	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	86
4		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	78
5		Benzeneacetaldehyde	120	C8H8O	000122-78-1	53



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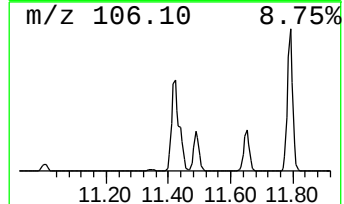
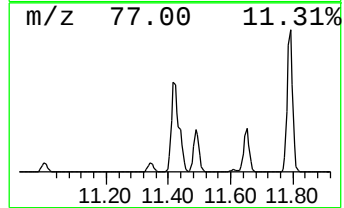
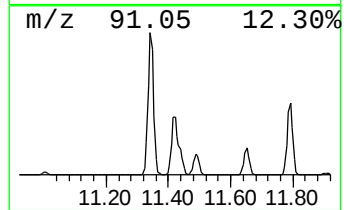
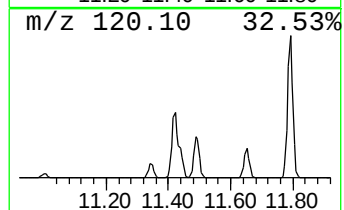
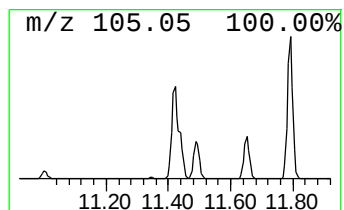
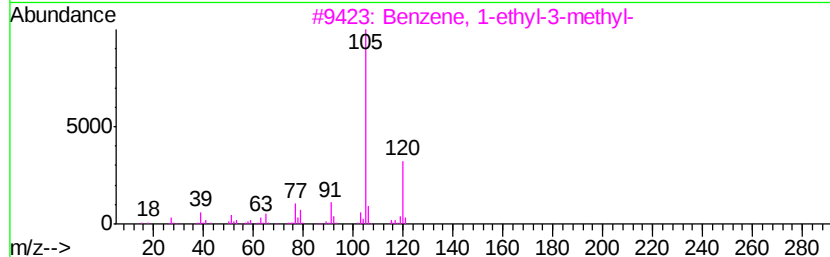
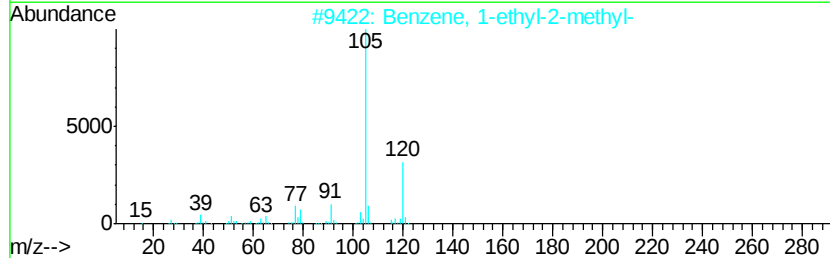
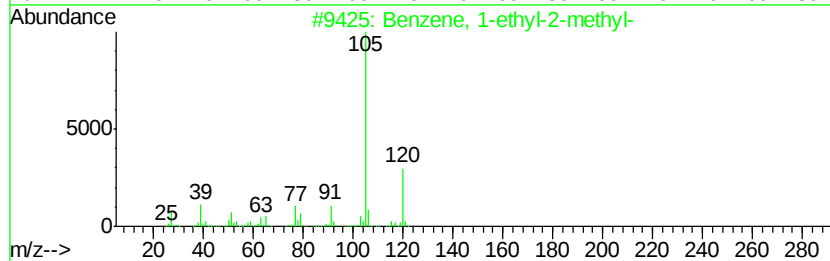
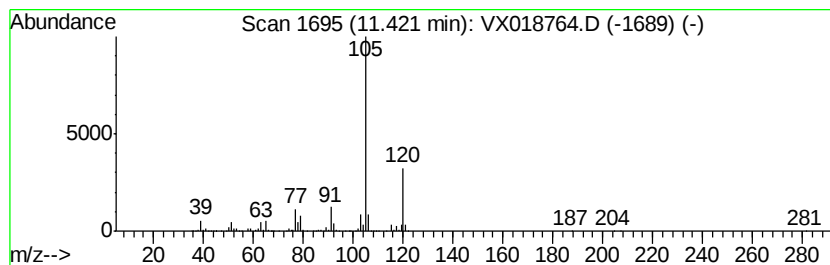
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	30.28 ug/L	3404890	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



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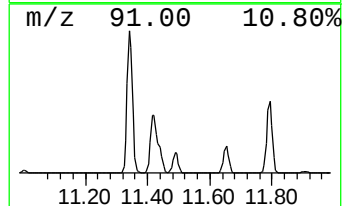
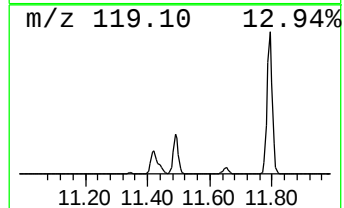
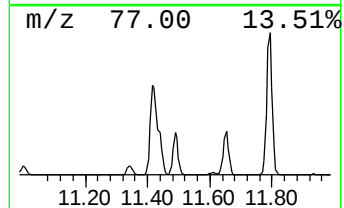
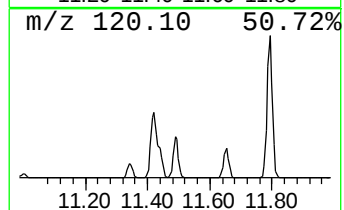
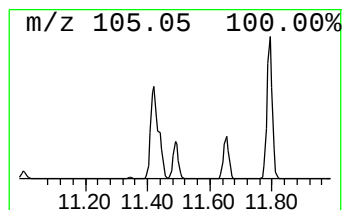
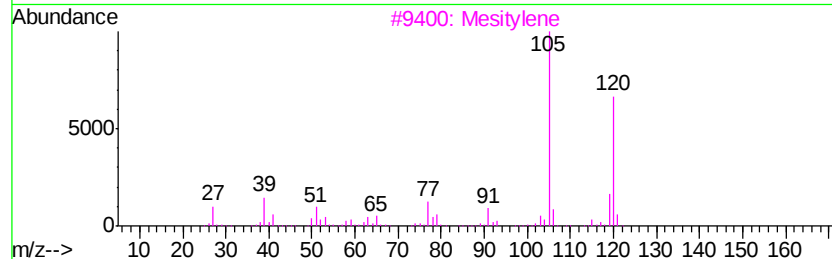
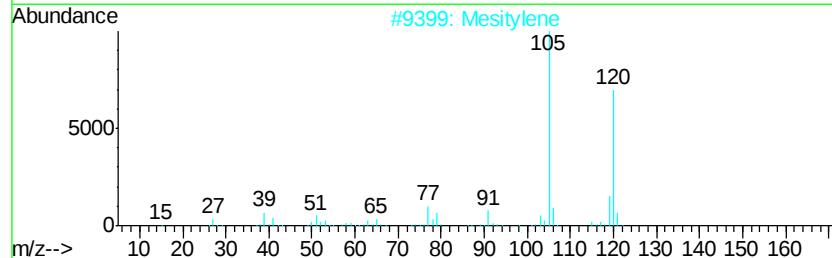
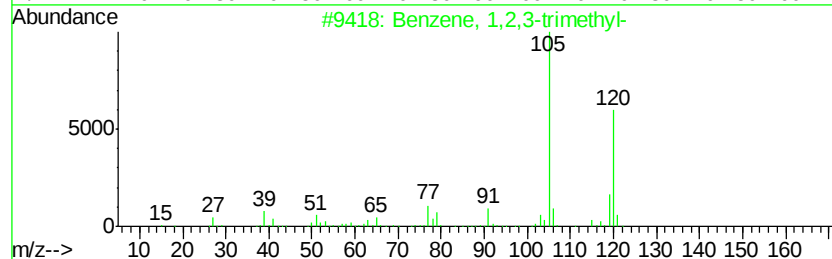
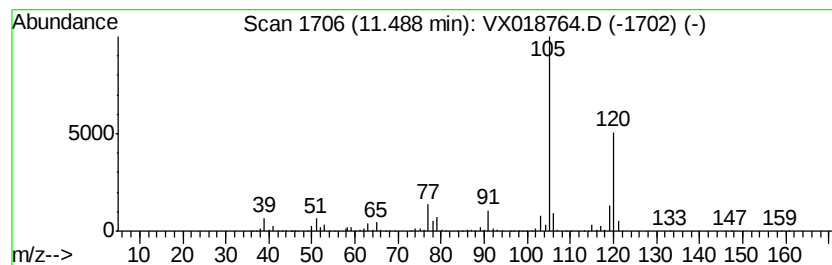
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	9.97 ug/L	1120950	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		Mesitylene	120	C9H12	000108-67-8	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Mesitylene	120	C9H12	000108-67-8	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

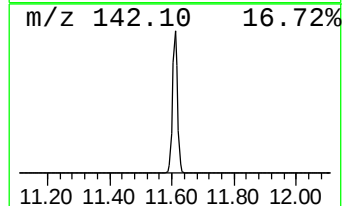
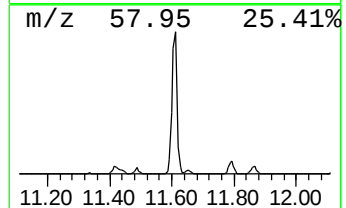
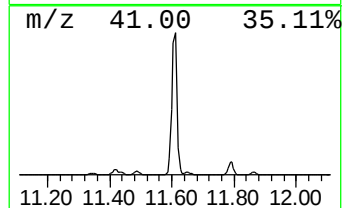
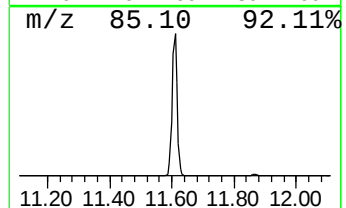
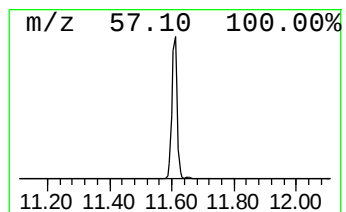
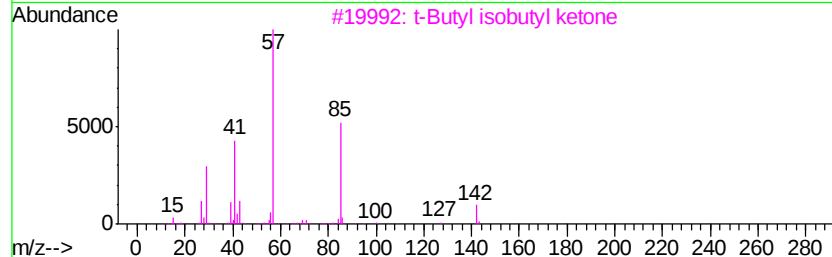
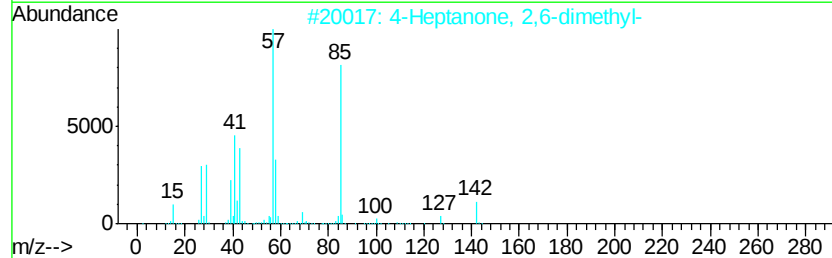
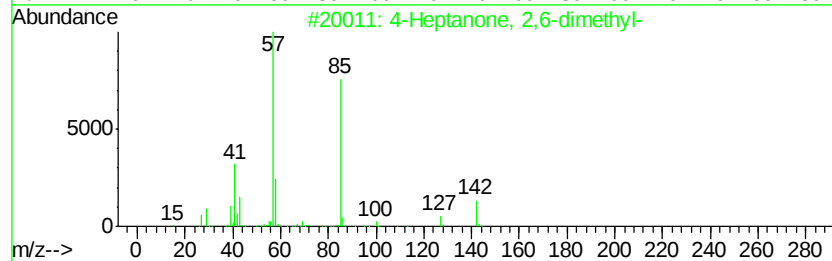
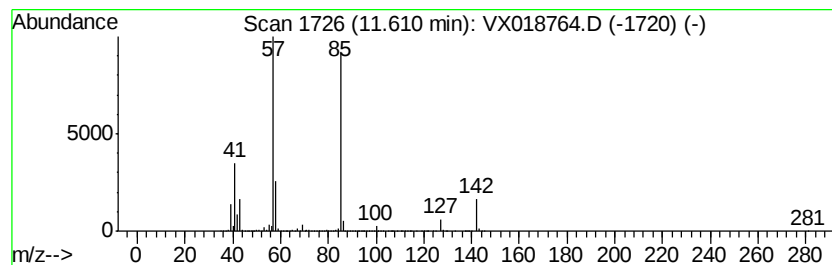
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Quant Title : TRACE VOA SOM01.0

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	34.56 ug/L	3885340	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	94
2		4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	83
3		t-Butyl isobutyl ketone	142	C9H18O	014705-50-1	59
4		5-Nonanone	142	C9H18O	000502-56-7	53
5		4-Octanone, 2-methyl-	142	C9H18O	007492-38-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

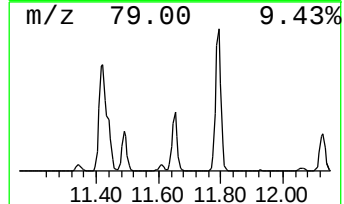
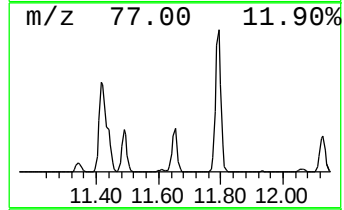
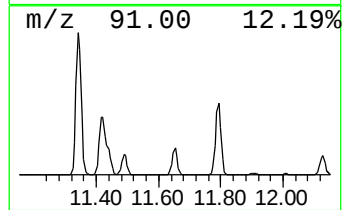
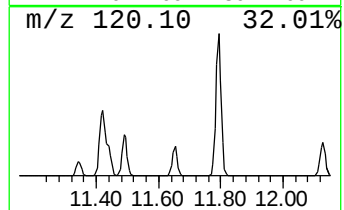
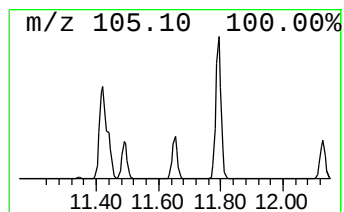
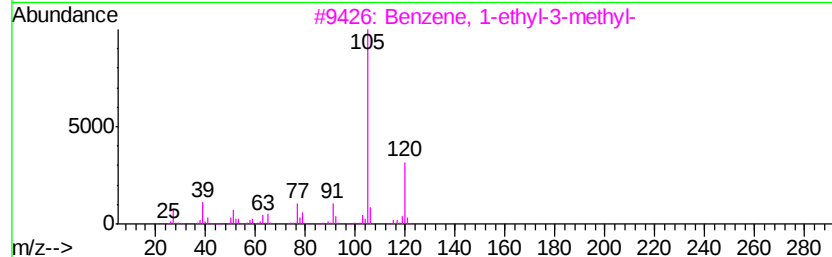
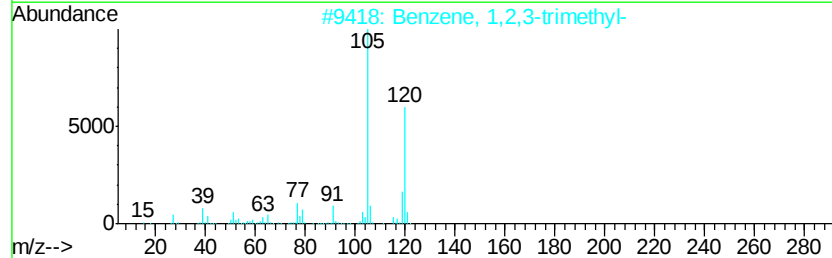
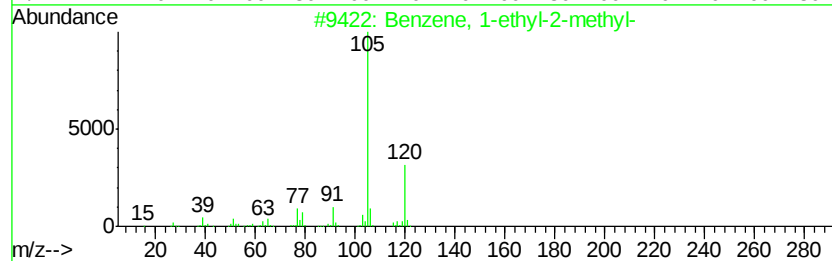
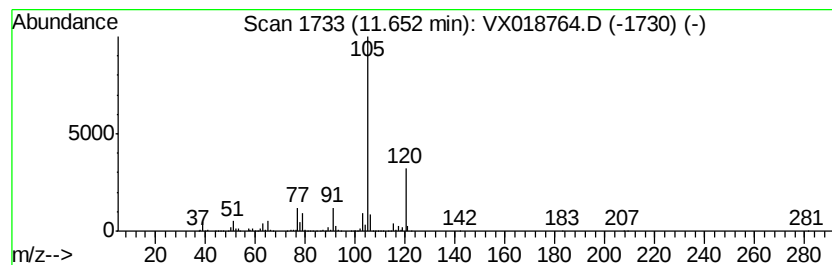
Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	9.52 ug/L	1070860	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, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91
4		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 91
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

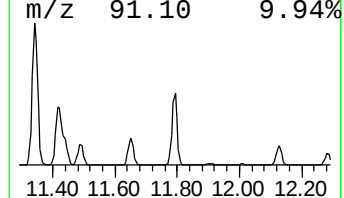
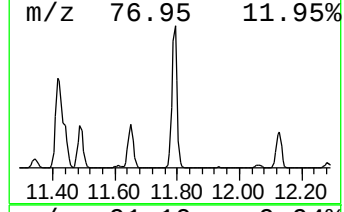
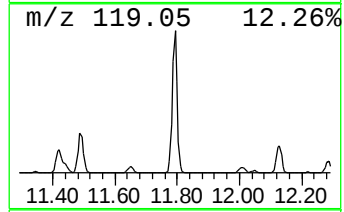
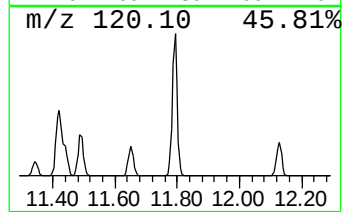
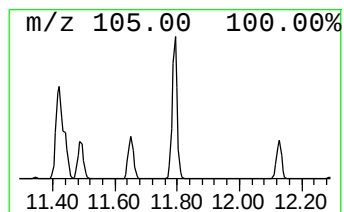
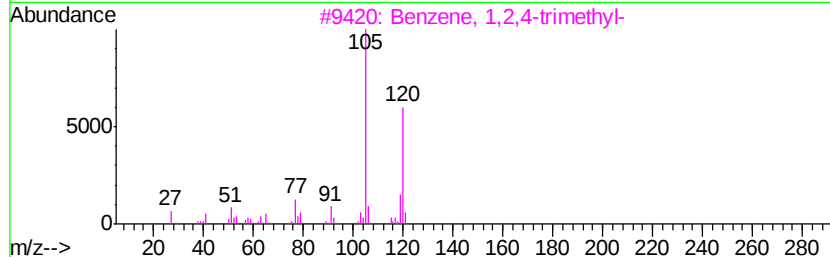
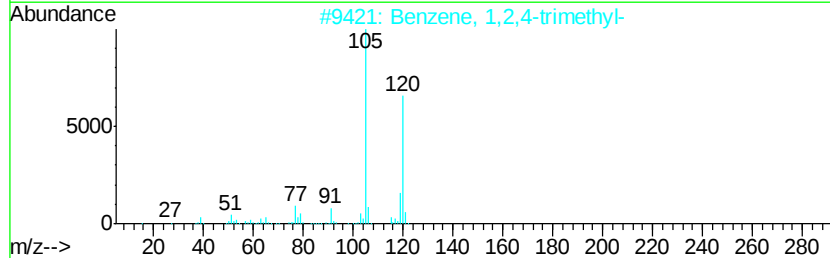
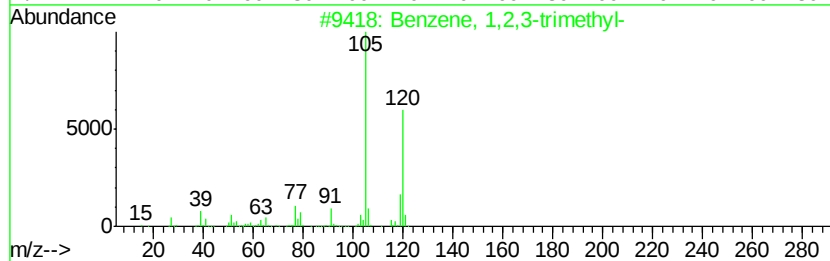
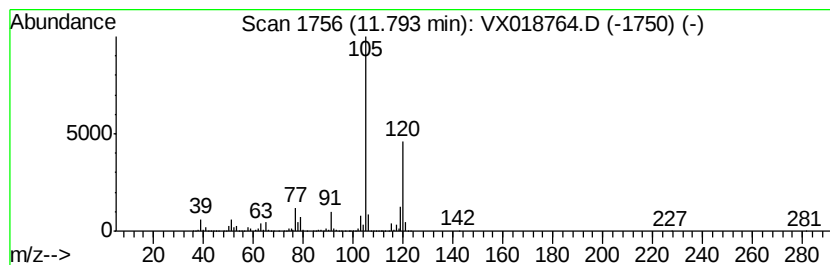
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, 1,2,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	34.27 ug/L	3852530	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		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Mesitylene	120	C9H12	000108-67-8	94
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

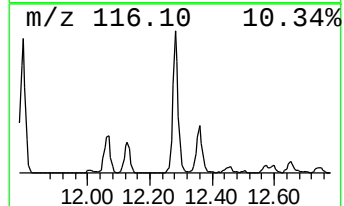
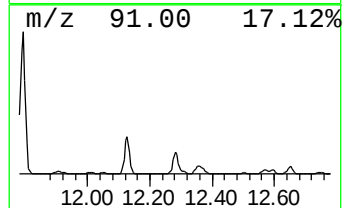
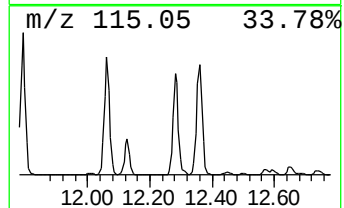
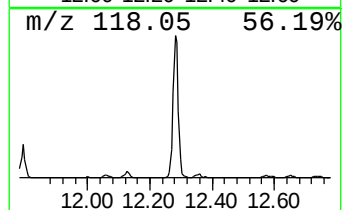
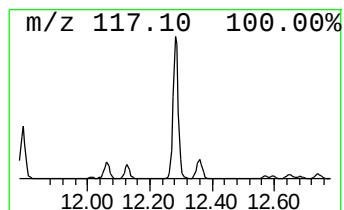
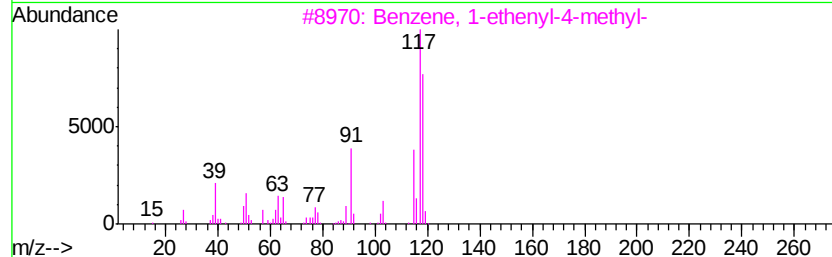
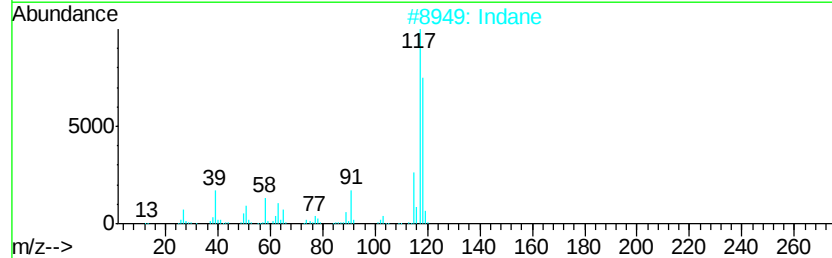
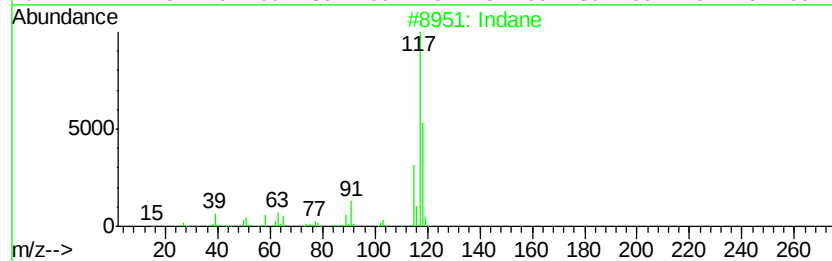
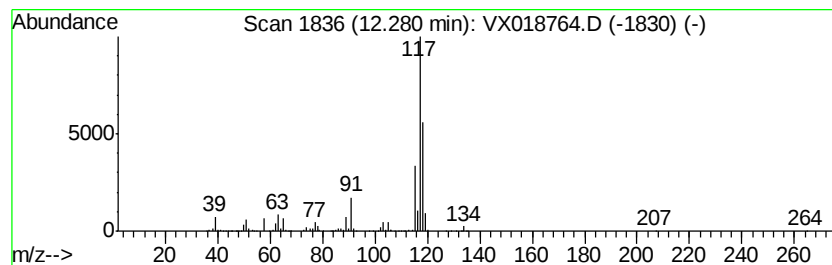
Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

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 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	4.41 ug/L	495921	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	Benzene, 1-ethenyl-4-methyl-		118 C9H10	000622-97-9 83
4	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 81
5	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

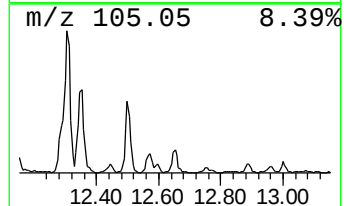
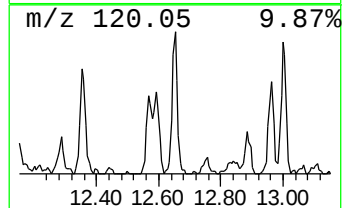
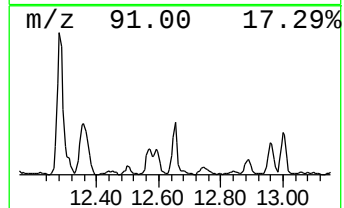
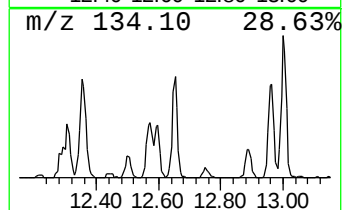
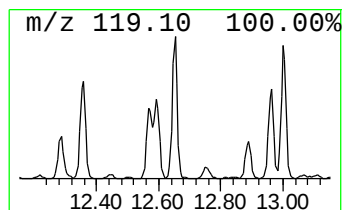
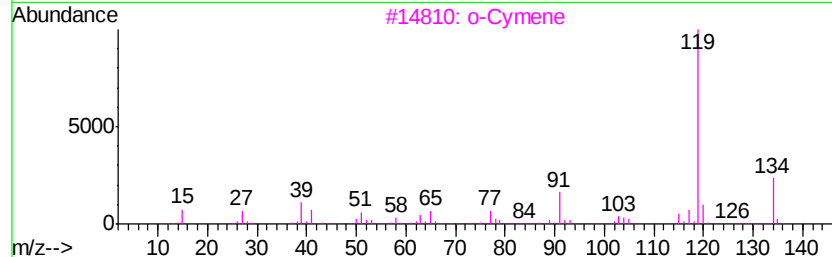
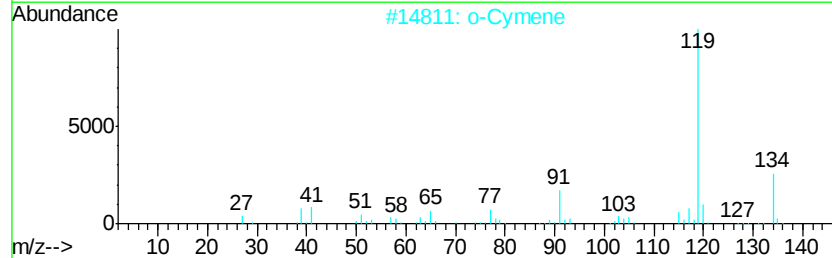
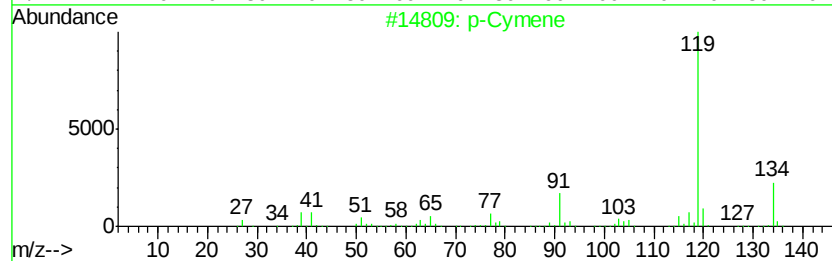
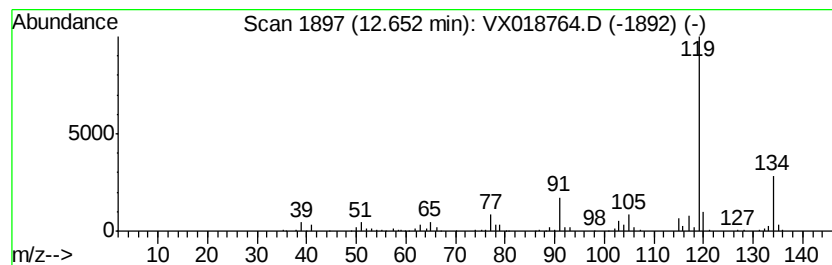
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ClientSampleId :
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 p-Cymene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	1.10 ug/L	123403	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	p-Cymene	134	C10H14	000099-87-6 97
2	o-Cymene	134	C10H14	000527-84-4 97
3	o-Cymene	134	C10H14	000527-84-4 97
4	o-Cymene	134	C10H14	000527-84-4 95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5 95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

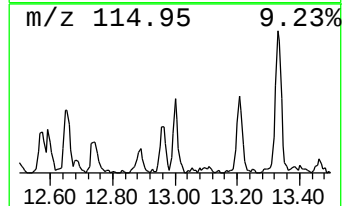
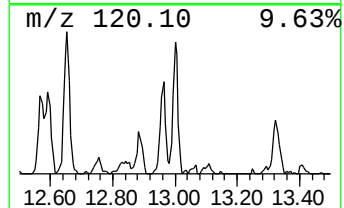
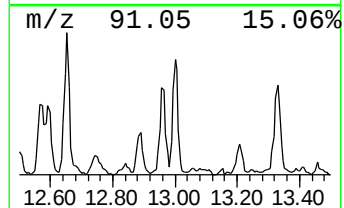
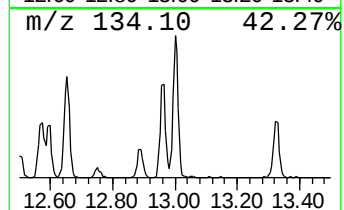
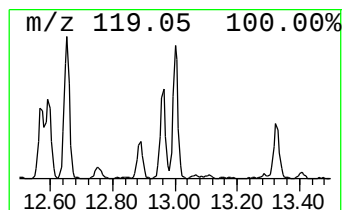
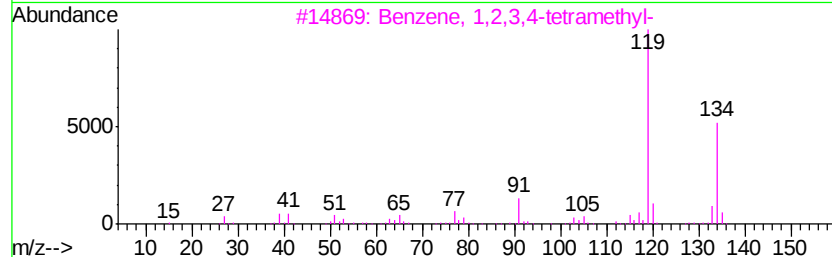
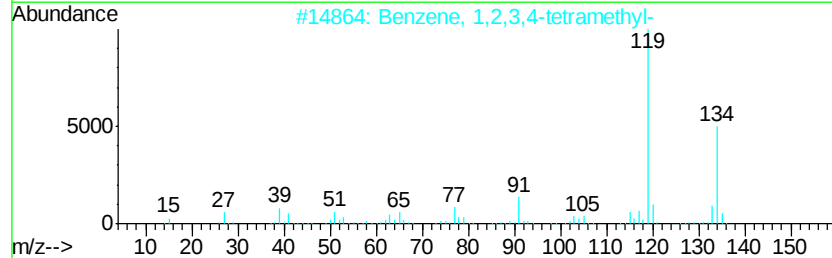
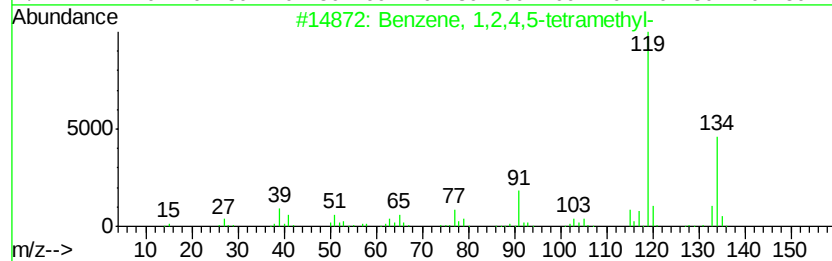
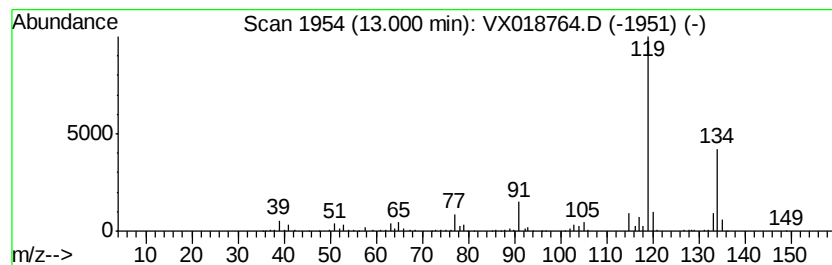
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	1.08 ug/L	121206	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		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

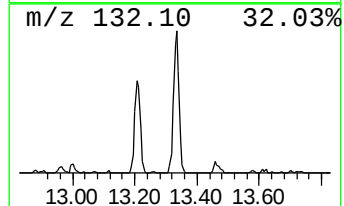
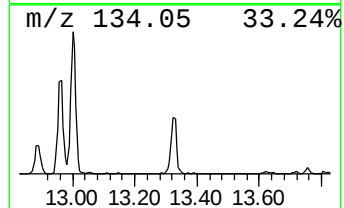
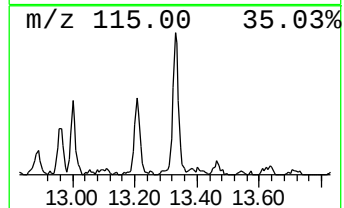
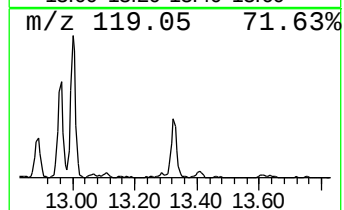
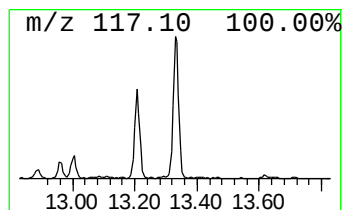
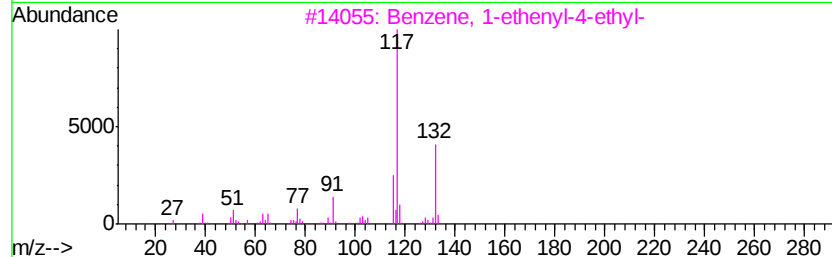
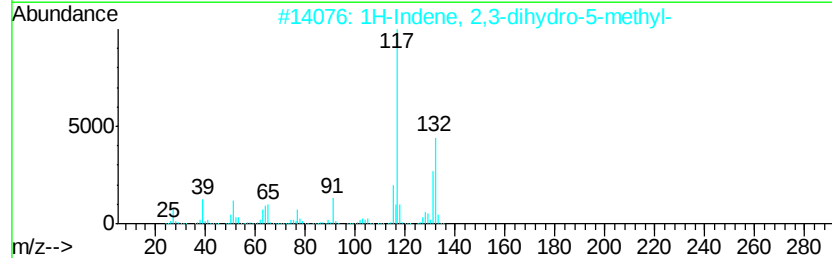
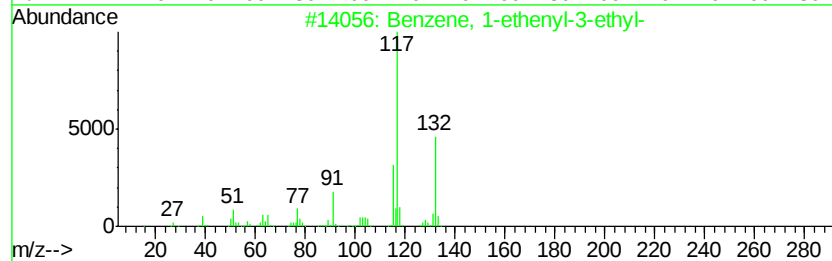
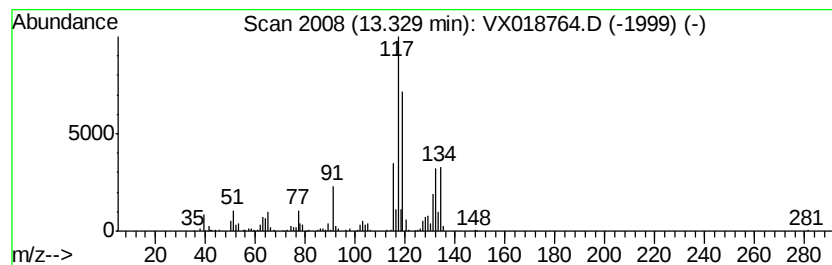
Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

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 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.33	1.17 ug/L	131263	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	58
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	58
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018764.D
Acq On : 05 Oct 2020 14:26
Operator : JC/SP
Sample : L4239-02DL 40X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N0DL

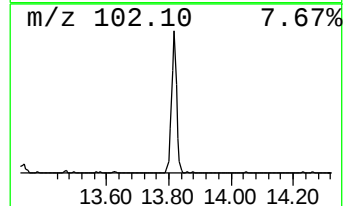
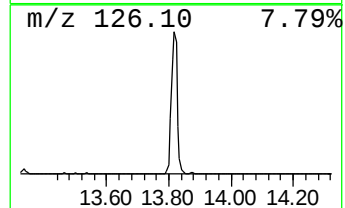
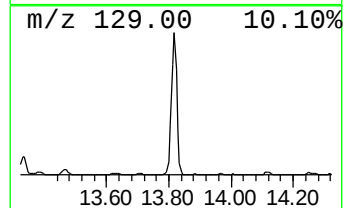
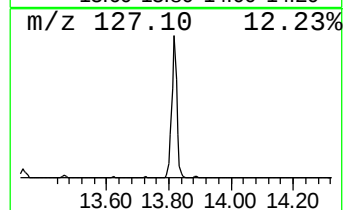
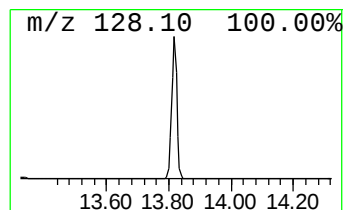
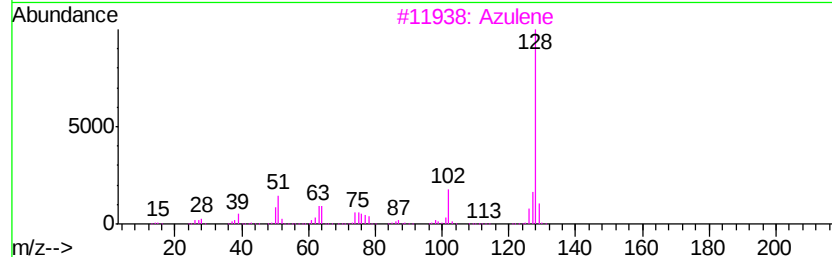
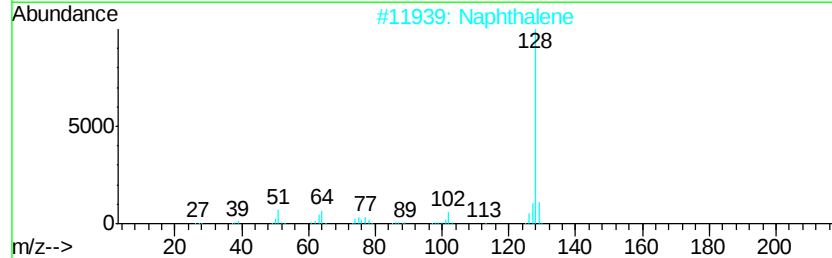
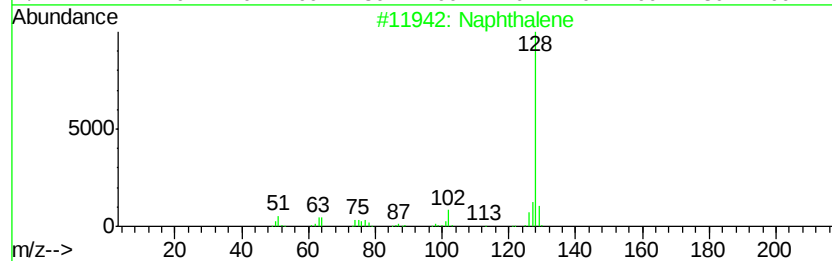
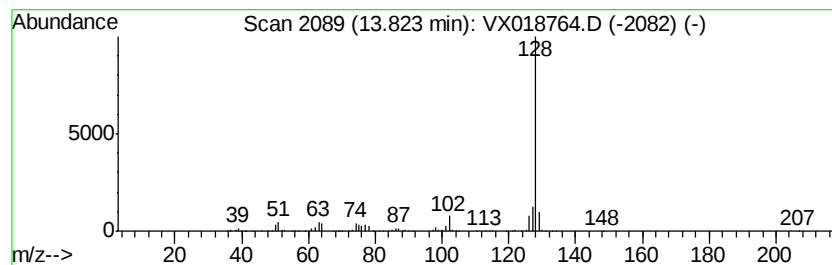
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 Azulene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	2.50 ug/L	280879	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene	128	C10H8	000091-20-3	95
2		Naphthalene	128	C10H8	000091-20-3	91
3		Azulene	128	C10H8	000275-51-4	91
4		Azulene	128	C10H8	000275-51-4	91
5		Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\WX100520\
 Data File : VX018764.D
 Acq On : 05 Oct 2020 14:26
 Operator : JC/SP
 Sample : L4239-02DL 40X
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N0DL

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	31.6	ug/L	2532070	1	6.83	401080	5.0
(DEL) Alkane: Cyc...	4.37	43.0	ug/L	3451940	1	6.83	401080	5.0
Benzene, propyl-	11.34	5.3	ug/L	593084	3	12.06	562150	5.0
Benzene, 1-ethyl-...	11.42	30.3	ug/L	3404890	3	12.06	562150	5.0
Benzene, 1,2,3-tr...	11.49	10.0	ug/L	1120950	3	12.06	562150	5.0
4-Heptanone, 2,6-...	11.61	34.6	ug/L	3885340	3	12.06	562150	5.0
Benzene, 1-ethyl-...	11.65	9.5	ug/L	1070860	3	12.06	562150	5.0
Benzene, 1,2,4-tr...	11.79	34.3	ug/L	3852530	3	12.06	562150	5.0
Indane	12.28	4.4	ug/L	495921	3	12.06	562150	5.0
p-Cymene	12.65	1.1	ug/L	123403	3	12.06	562150	5.0
Benzene, 1,2,4,5-...	13.00	1.1	ug/L	121206	3	12.06	562150	5.0
Benzene, 1-etheny...	13.33	1.2	ug/L	131263	3	12.06	562150	5.0
Azulene	13.82	2.5	ug/L	280879	3	12.06	562150	5.0