

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100620\
 Data File : VX018806.D
 Acq On : 06 Oct 2020 18:05
 Operator : JC/SP
 Sample : L4239-05DL2 400X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N5DL2

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	44	49	55	rBV	53024	54259	9.36%	0.786%
2	1.697	96	100	105	rBV	47451	53184	9.17%	0.770%
3	2.343	198	206	220	rVB	97932	157909	27.23%	2.286%
4	2.831	279	286	287	rBV6	5236	8533	1.47%	0.124%
5	2.867	288	292	301	rVB	13400	28147	4.85%	0.407%
6	3.148	328	338	347	rBV2	26488	64413	11.11%	0.933%
7	3.501	387	396	404	rBV	48106	114695	19.78%	1.660%
8	3.672	419	424	430	rVB3	2973	6561	1.13%	0.095%
9	4.373	529	539	551	rVB3	32575	97089	16.74%	1.406%
10	4.556	558	569	580	rBV3	85803	292528	50.45%	4.235%
11	5.141	653	665	677	rBV	69414	213053	36.74%	3.084%
12	5.458	706	717	727	rBV2	37296	122048	21.05%	1.767%
13	6.044	799	813	826	rBV3	159542	466019	80.36%	6.747%
14	6.830	932	942	952	rBV	116947	280112	48.30%	4.055%
15	7.367	1019	1030	1040	rBV	98589	219001	37.77%	3.171%
16	8.372	1189	1195	1204	rBV	60908	103529	17.85%	1.499%
17	8.696	1239	1248	1254	rBV	237238	371351	64.04%	5.376%
18	8.763	1254	1259	1272	rVB	382668	579883	100.00%	8.395%
19	8.994	1292	1297	1305	rBV2	45640	67581	11.65%	0.978%
20	9.427	1362	1368	1388	rBV	366030	504643	87.02%	7.306%
21	9.738	1414	1419	1425	rVB	15880	23054	3.98%	0.334%
22	10.098	1472	1478	1487	rBV	242422	338651	58.40%	4.903%
23	10.238	1494	1501	1506	rBV	37872	55180	9.52%	0.799%
24	10.342	1512	1518	1524	rBV	142083	190623	32.87%	2.760%
25	10.683	1569	1574	1579	rBV	74307	96446	16.63%	1.396%
26	11.000	1620	1626	1633	rVB	15399	20746	3.58%	0.300%
27	11.232	1656	1664	1669	rBV	108479	146115	25.20%	2.115%
28	11.341	1678	1682	1688	rBV	39864	54260	9.36%	0.786%
29	11.421	1688	1695	1702	rVV	173642	312616	53.91%	4.526%
30	11.488	1702	1706	1714	rVB	90371	116009	20.01%	1.679%
31	11.610	1720	1726	1729	rBV	207255	247907	42.75%	3.589%
32	11.652	1729	1733	1740	rVB	99307	127362	21.96%	1.844%
33	11.793	1749	1756	1762	rBV	331676	398299	68.69%	5.766%
34	11.866	1762	1768	1773	rVV2	6100	8608	1.48%	0.125%

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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	12.061	1795	1800	1806	rBV	274730	339935	58.62%	4.921%
36	12.128	1806	1811	1816	rVB	89243	113247	19.53%	1.640%
37	12.280	1830	1836	1844	rBV2	40725	56853	9.80%	0.823%
38	12.360	1844	1849	1858	rVB	306146	381441	65.78%	5.522%
39	12.573	1878	1884	1886	rVV3	6034	8937	1.54%	0.129%
40	12.652	1893	1897	1902	rVV	7614	9669	1.67%	0.140%
41	12.957	1942	1947	1951	rBV2	5877	7739	1.33%	0.112%
42	13.000	1951	1954	1960	rVB	8563	10689	1.84%	0.155%
43	13.329	2002	2008	2014	rVB5	7574	11851	2.04%	0.172%
44	13.817	2084	2088	2093	rBV	20165	26608	4.59%	0.385%

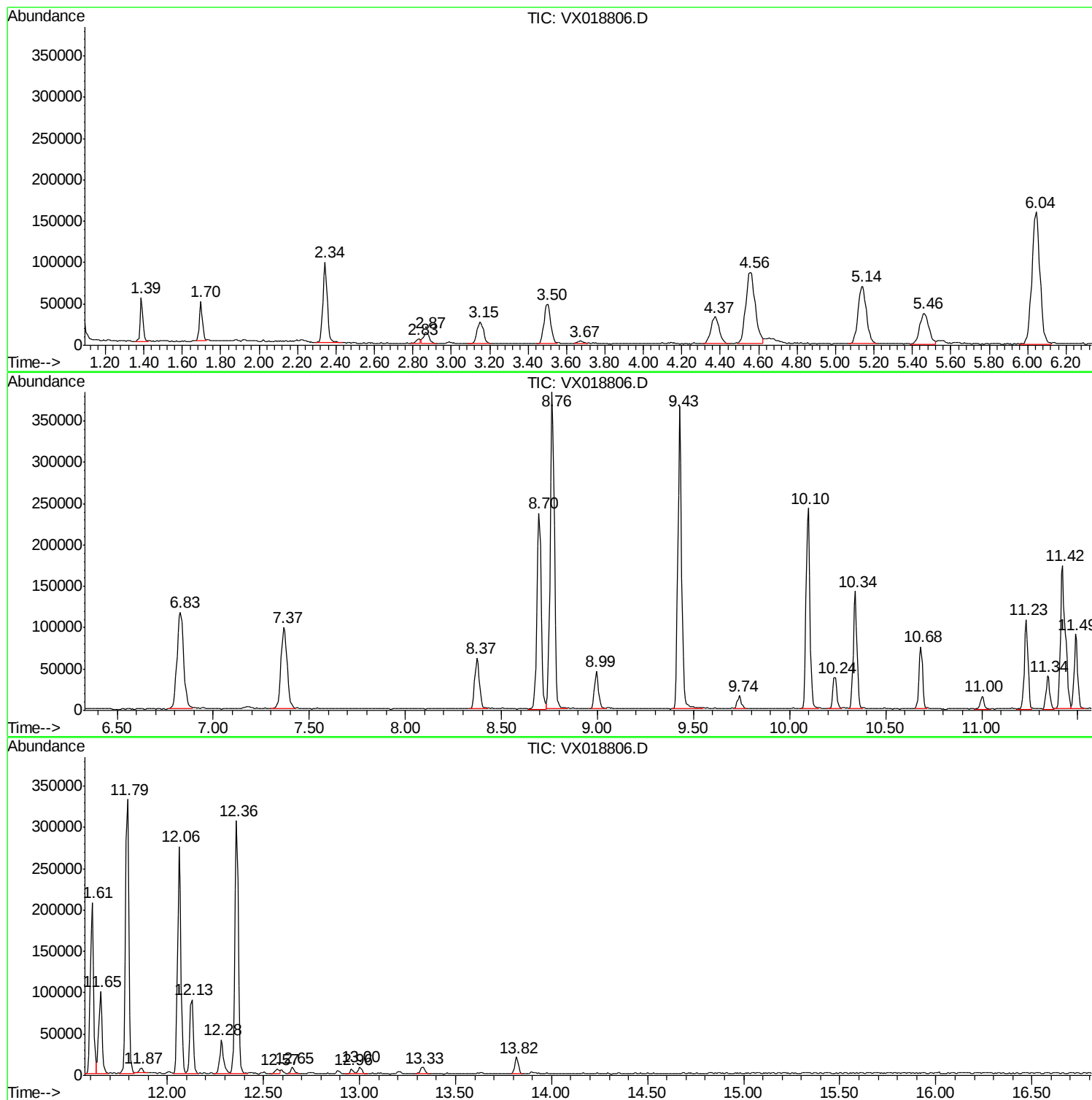
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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



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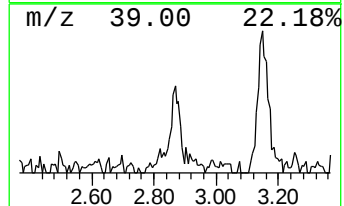
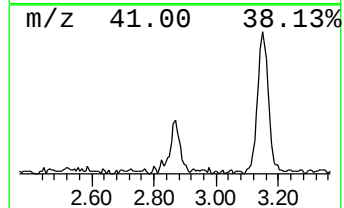
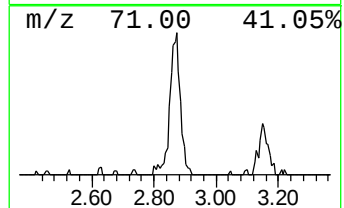
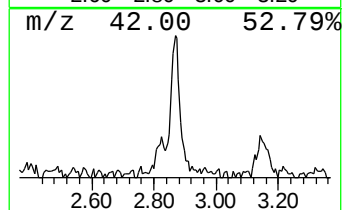
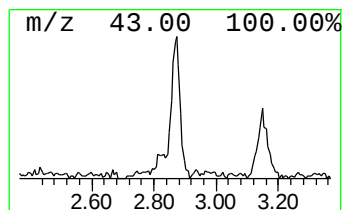
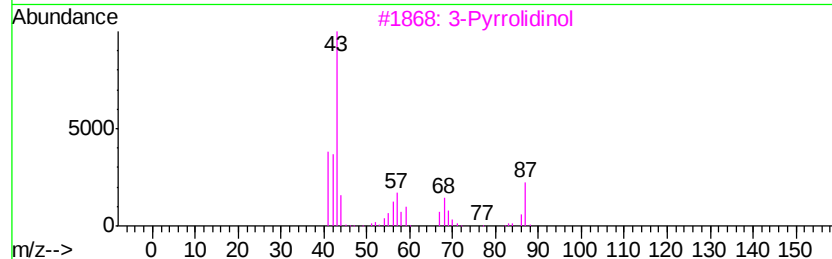
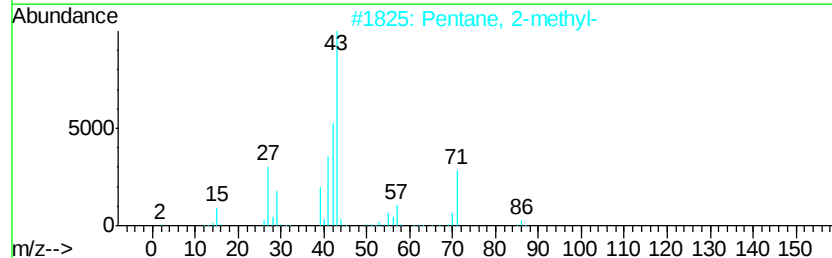
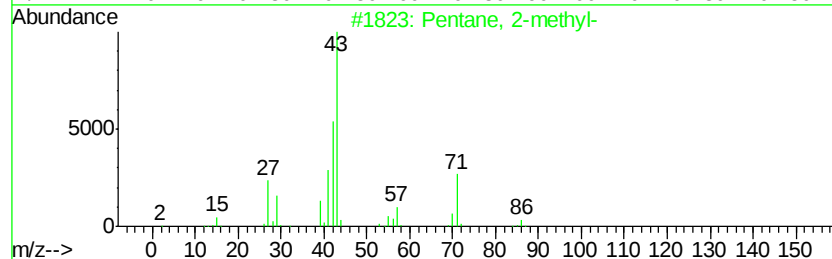
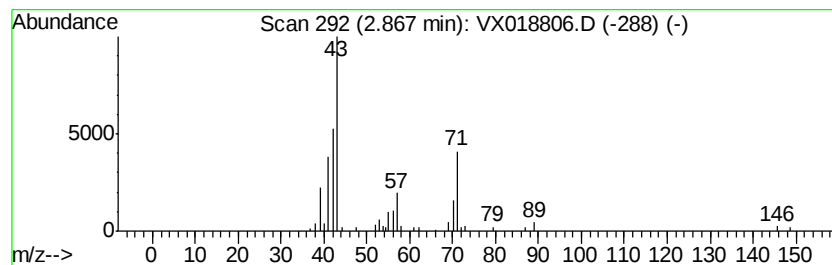
Instrument :
MSVOA_X
ClientSampleId :
BG3N5DL2

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 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.87	0.50 ug/L	28147	1,4-Difluorobenzene	6.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	56
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	56
3	3-Pyrrolidinol	87	C4H9NO	040499-83-0	32
4	Sulfurous acid, pentyl 2-propyl ...	194	C8H18O3S	1000309-11-5	25
5	Propanoic acid, 2-methyl-, 2-eth...	200	C12H24O2	035061-61-1	23



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Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

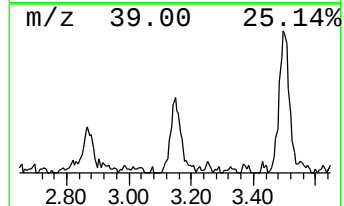
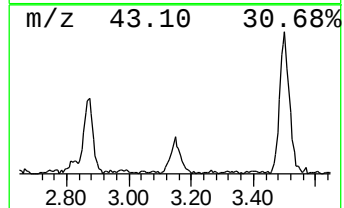
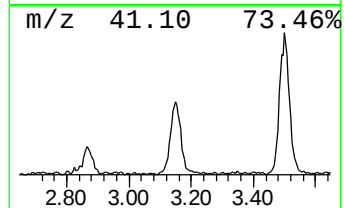
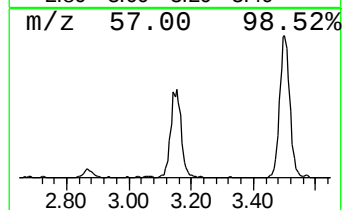
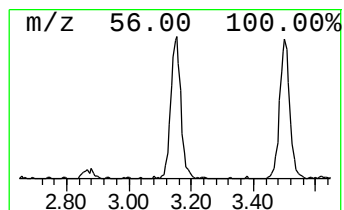
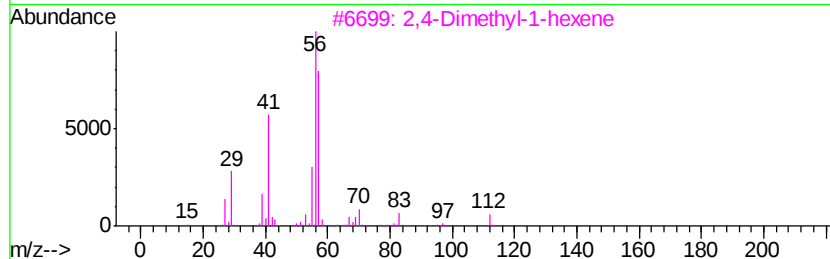
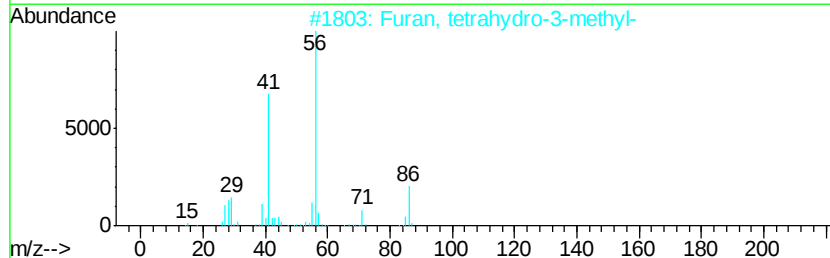
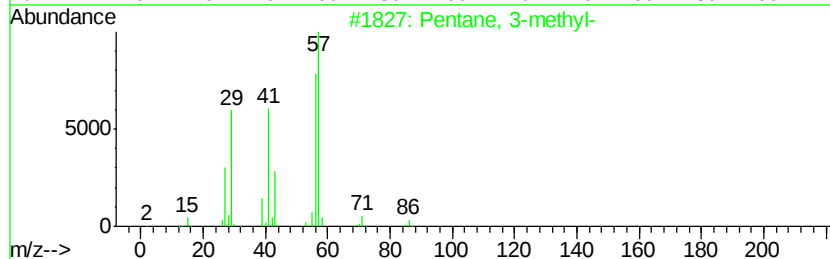
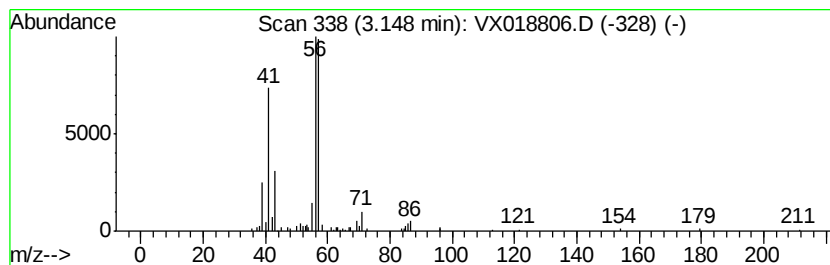
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	1.15 ug/L	64413	1,4-Difluorobenzene	6.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	78
2			Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	64
3			2,4-Dimethyl-1-hexene	112	C8H16	016746-87-5	64
4			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	58
5			Pentane, 3-methyl-	86	C6H14	000096-14-0	56



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Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

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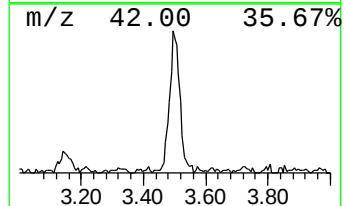
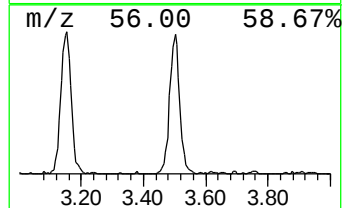
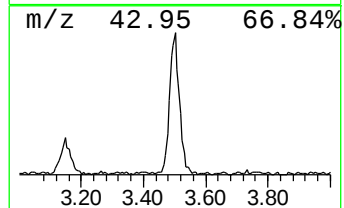
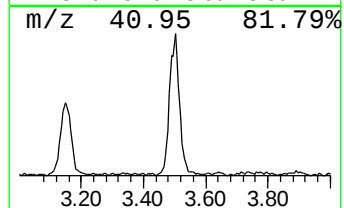
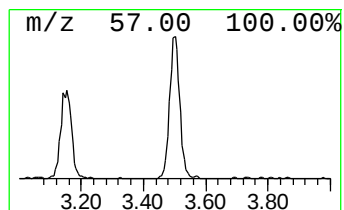
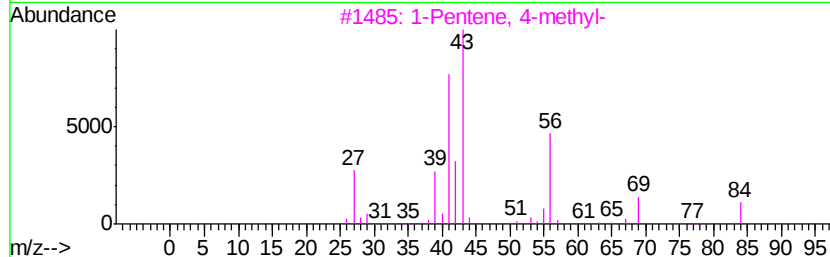
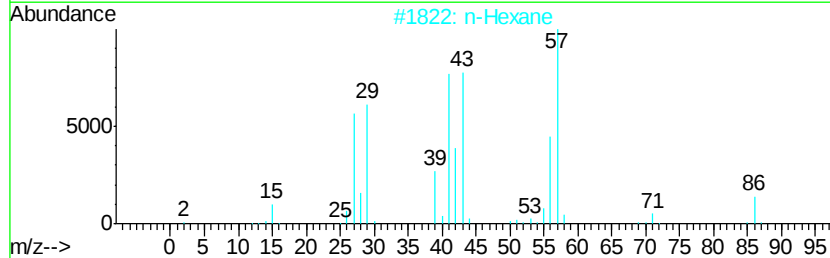
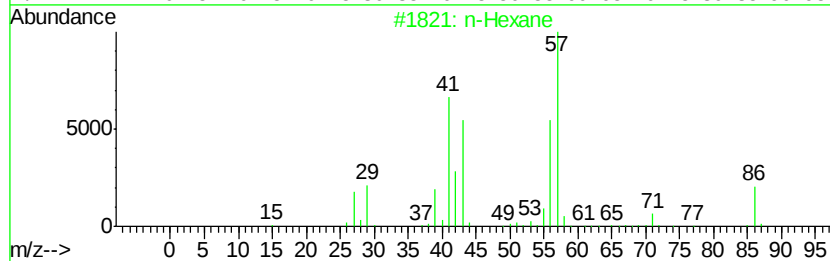
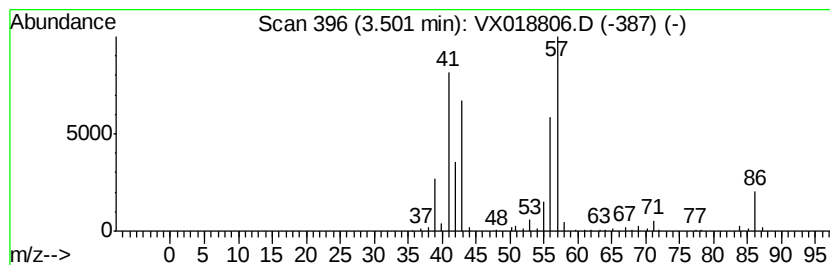
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	2.05 ug/L	114695	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	91
2		n-Hexane	86	C6H14	000110-54-3	59
3		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	47
4		n-Hexane	86	C6H14	000110-54-3	45
5		Butanal, 2-methyl-	86	C5H10O	000096-17-3	43



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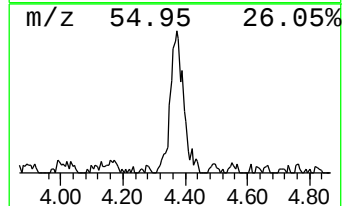
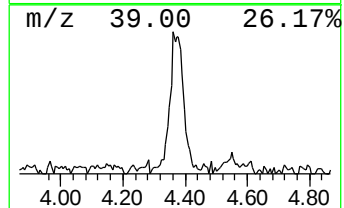
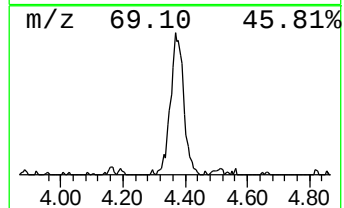
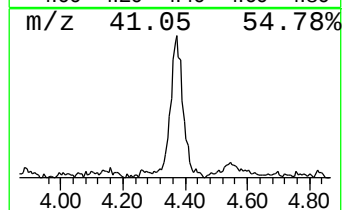
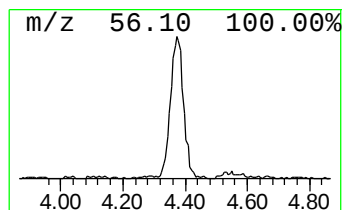
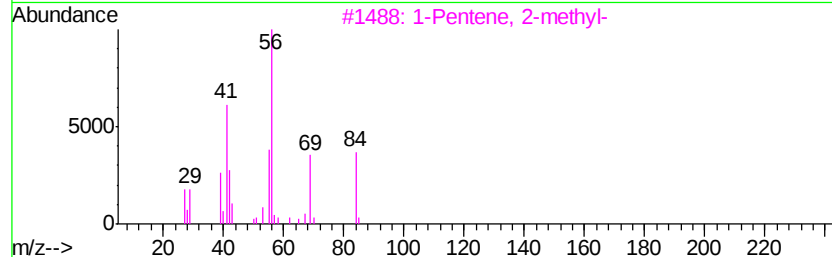
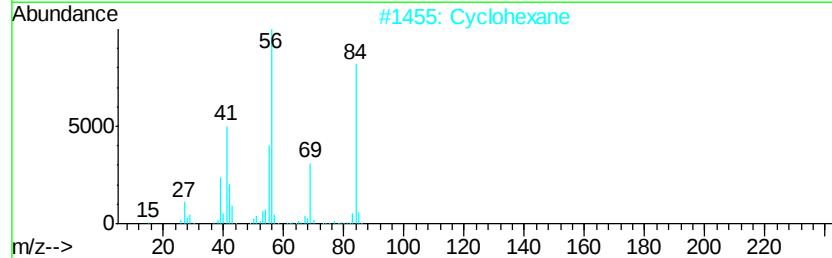
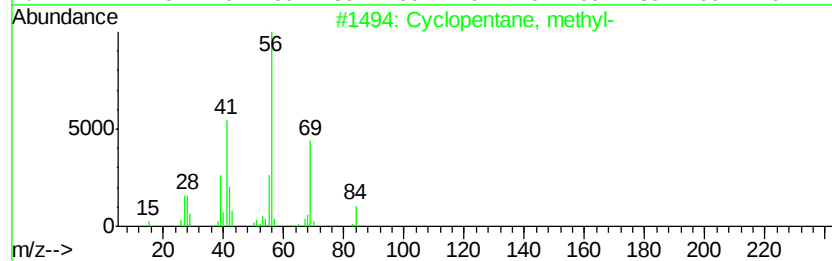
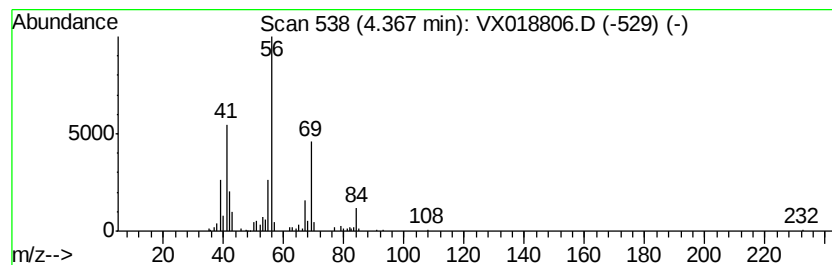
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Cyclic4.37 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	1.73 ug/L	97089	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2		Cyclohexane	84	C6H12	000110-82-7	72
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59



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ClientSampled :
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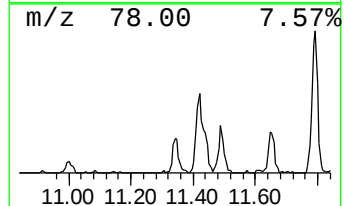
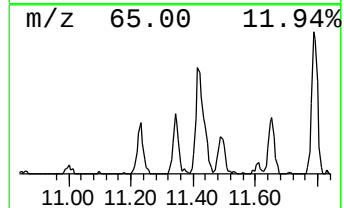
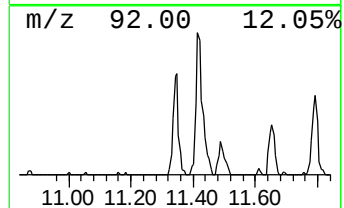
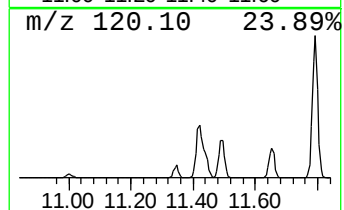
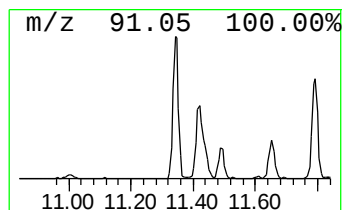
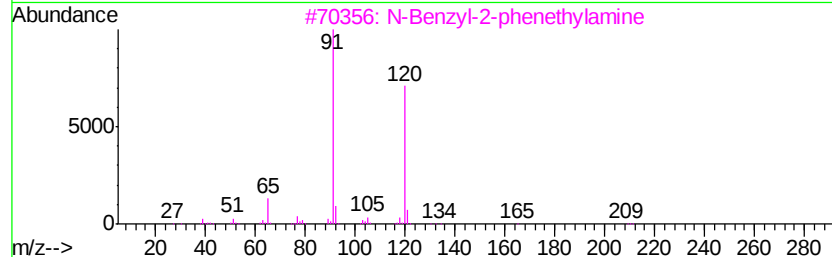
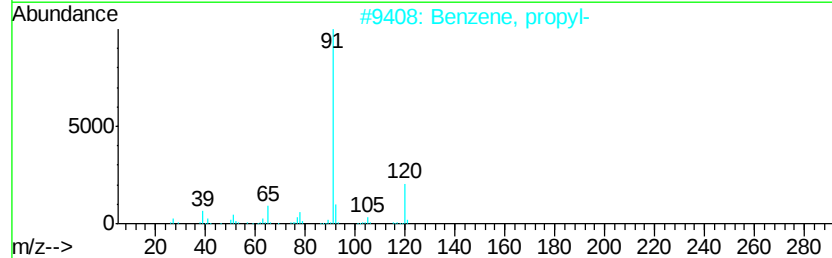
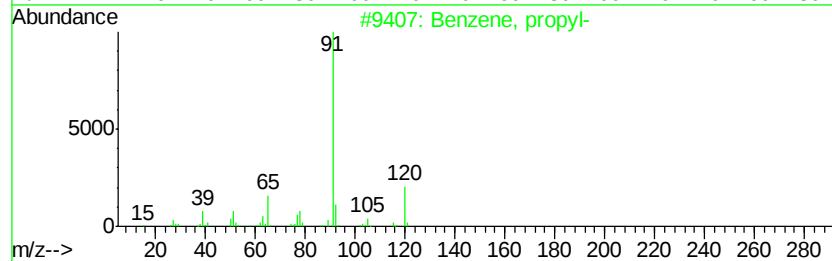
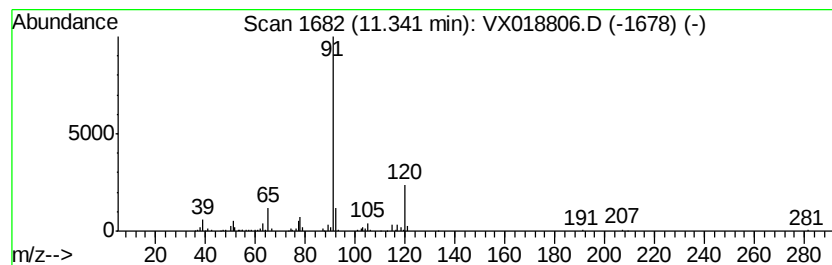
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	0.80 ug/L	54260	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	87
2		Benzene, propyl-	120	C9H12	000103-65-1	81
3		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
4		n-Butylbenzylamine	163	C11H17N	002403-22-7	78
5		Benzene, propyl-	120	C9H12	000103-65-1	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

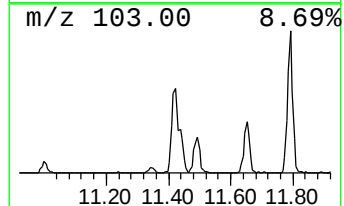
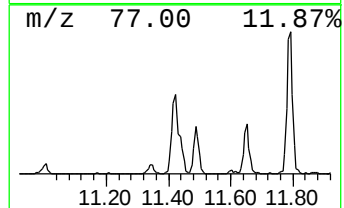
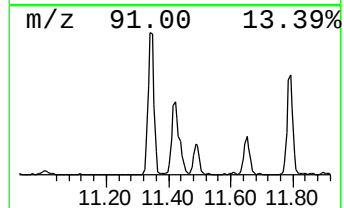
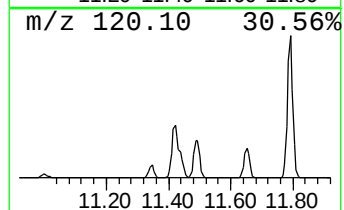
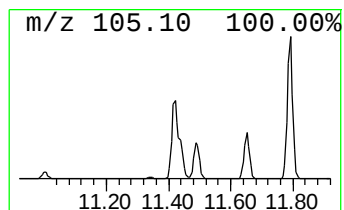
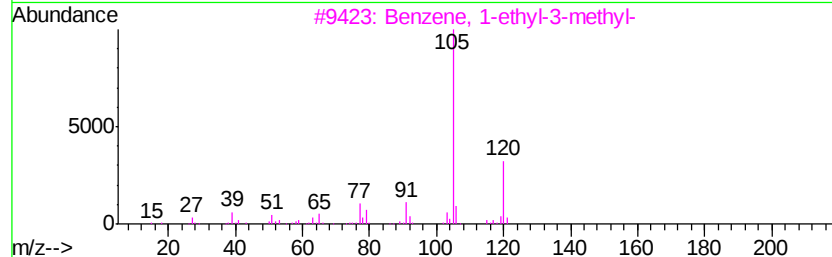
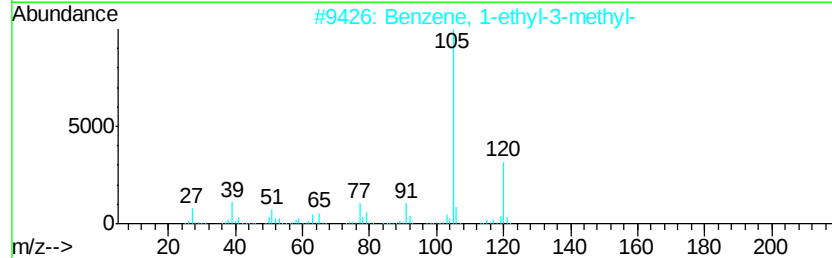
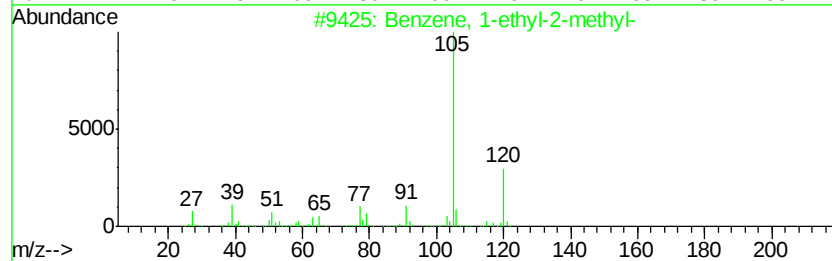
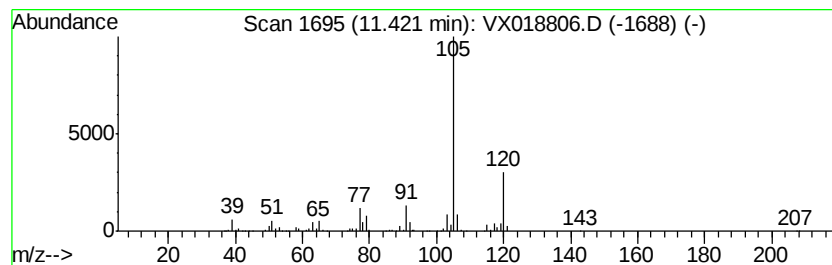
Instrument :
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ClientSampled :
BG3N5DL2

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 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	4.60 ug/L	312616	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4	94
3	Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4	94
4	Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N5DL2

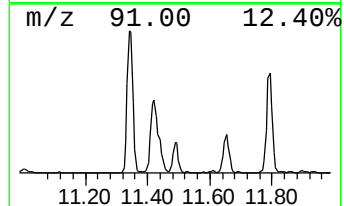
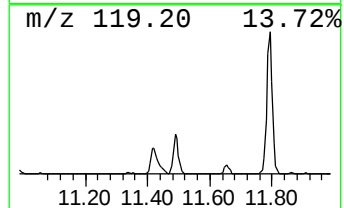
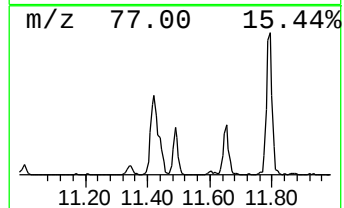
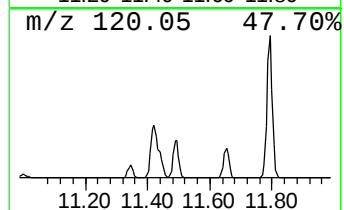
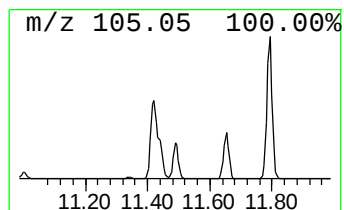
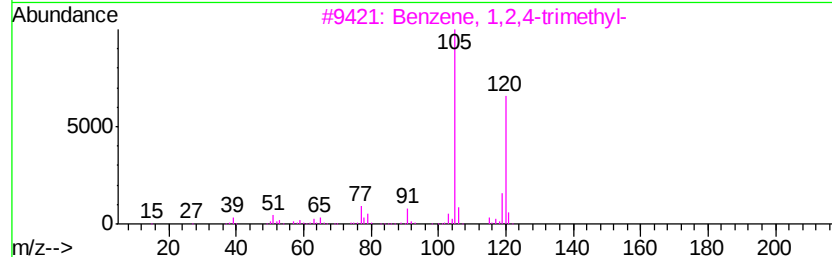
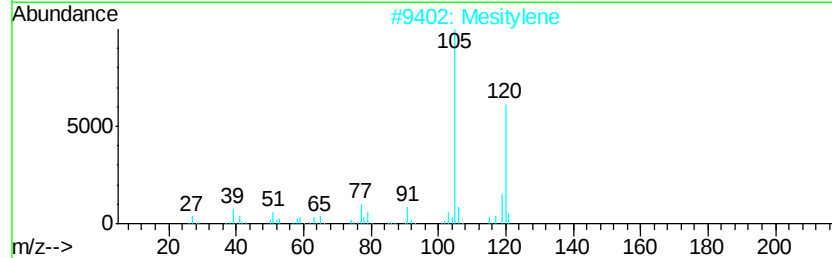
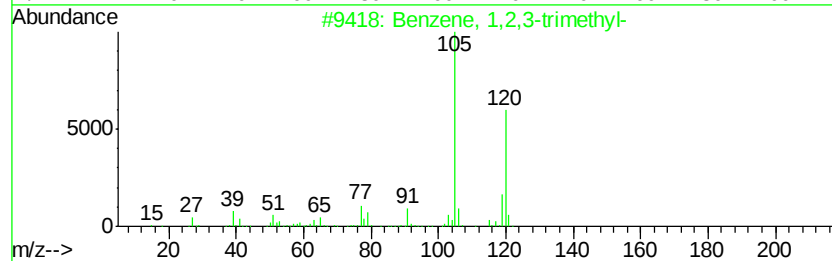
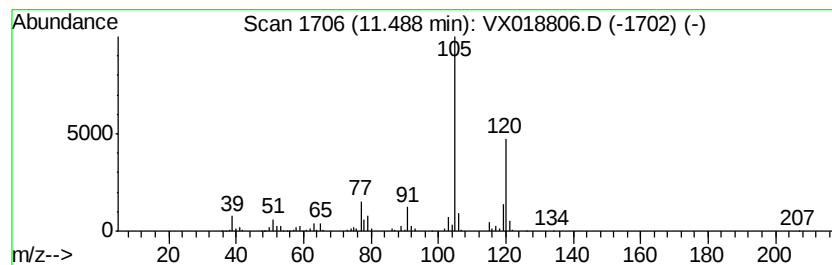
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	1.71 ug/L	116009	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	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Mesitylene	120	C9H12	000108-67-8	95
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

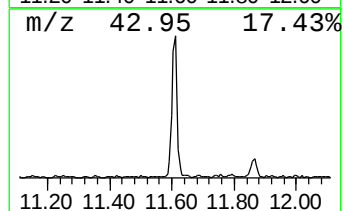
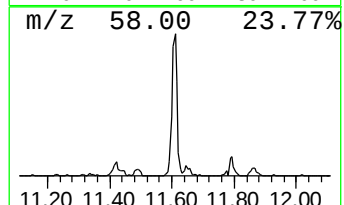
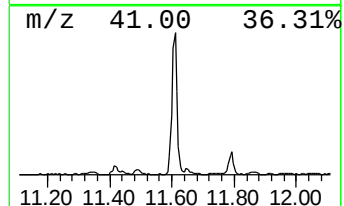
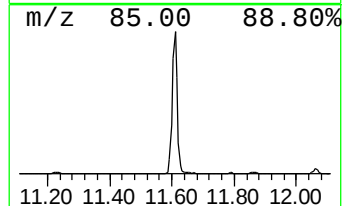
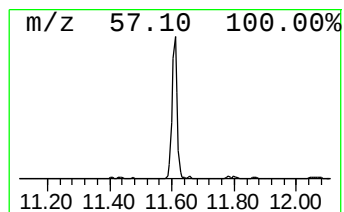
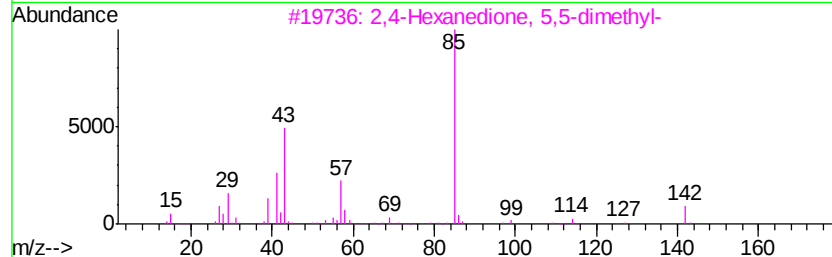
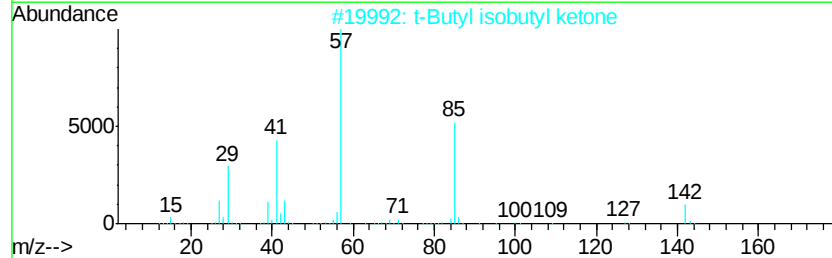
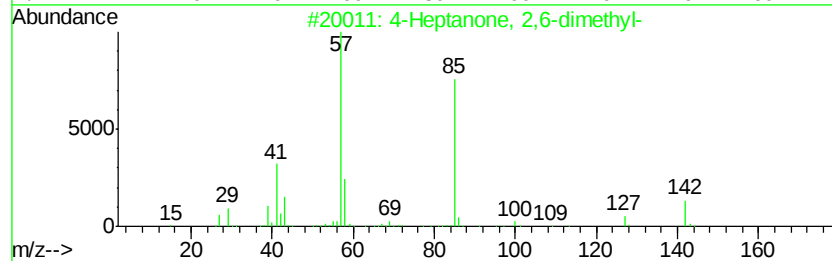
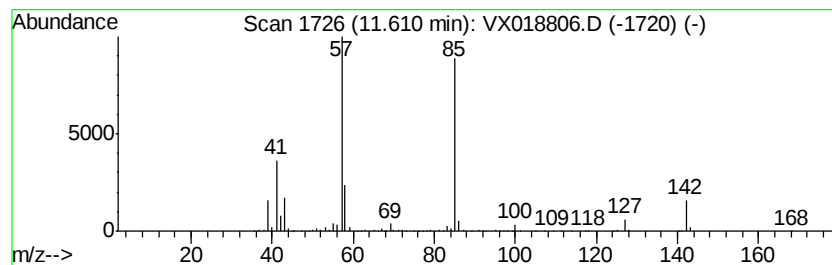
Instrument :
MSVOA_X
ClientSampleId :
BG3N5DL2

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 9 4-Heptanone, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.61	3.65 ug/L	247907	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	t-Butyl isobutyl ketone	142	C9H18O	014705-50-1	64
3	2,4-Hexanedione, 5,5-dimethyl-	142	C8H14O2	007307-04-2	59
4	5-Nonanone	142	C9H18O	000502-56-7	59
5	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

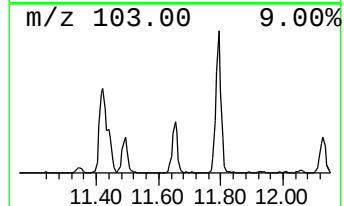
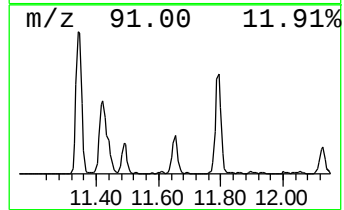
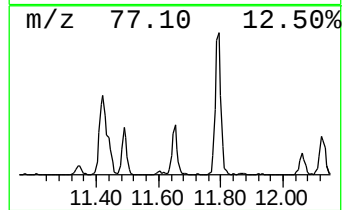
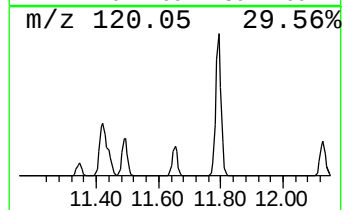
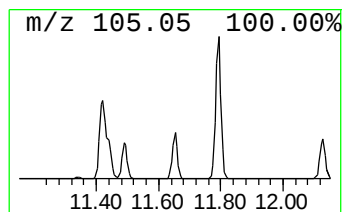
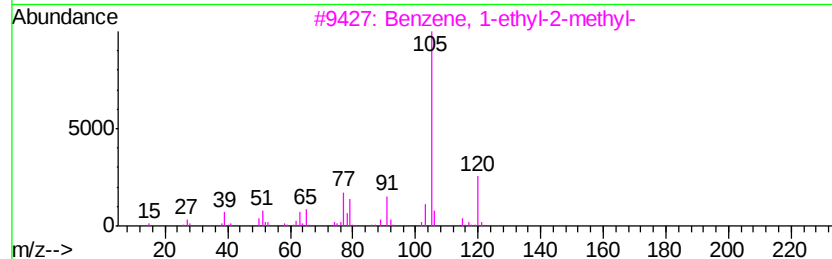
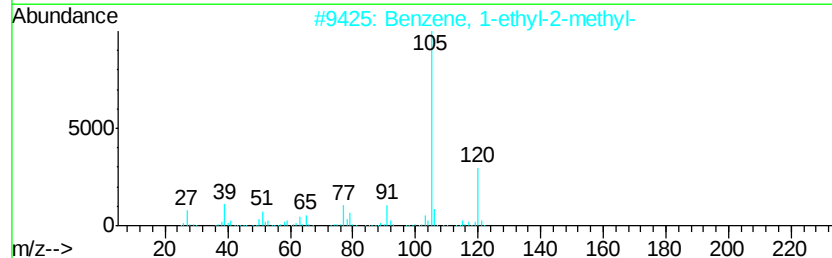
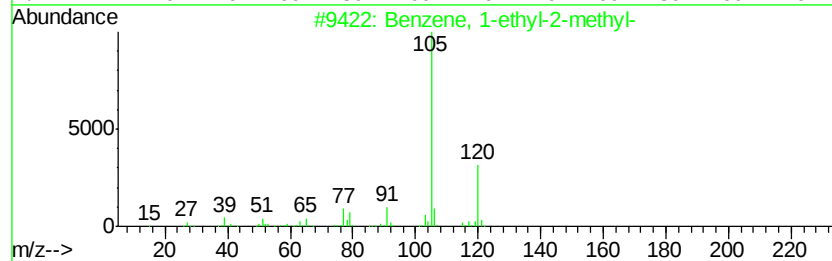
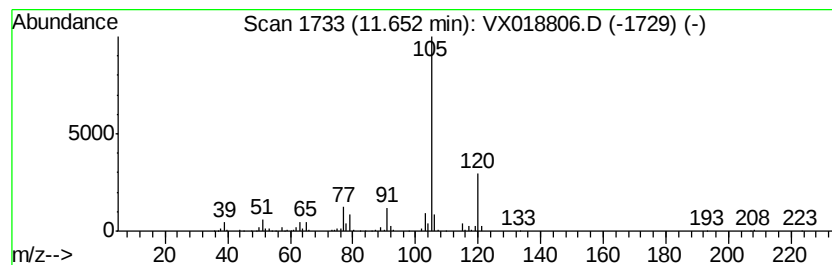
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Benzene, 1-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	1.87 ug/L	127362	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-2-methyl-	120	C9H12	000611-14-3	93
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

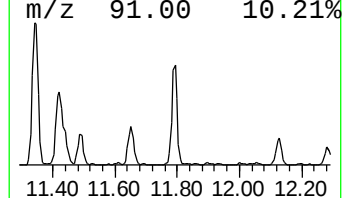
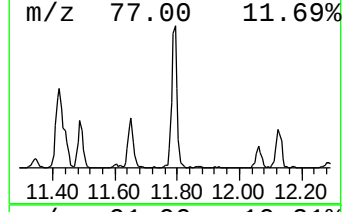
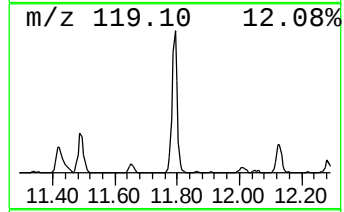
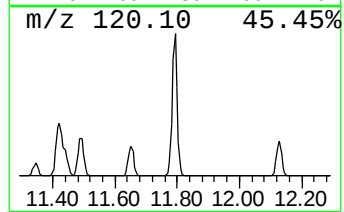
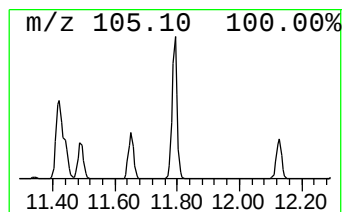
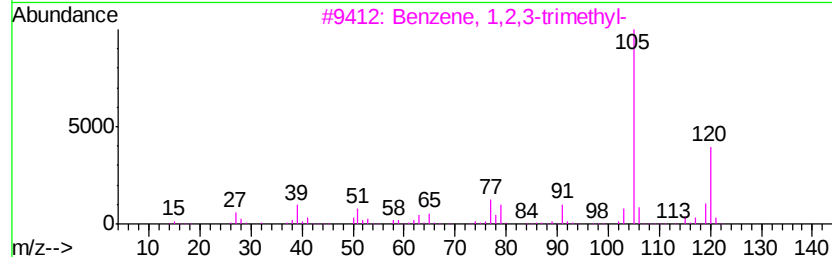
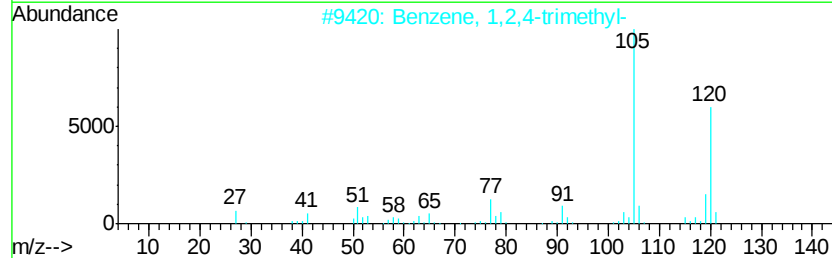
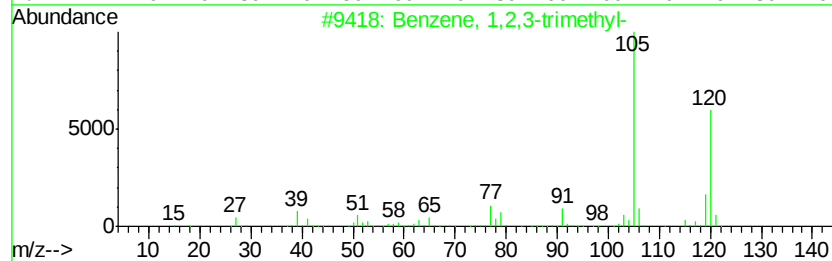
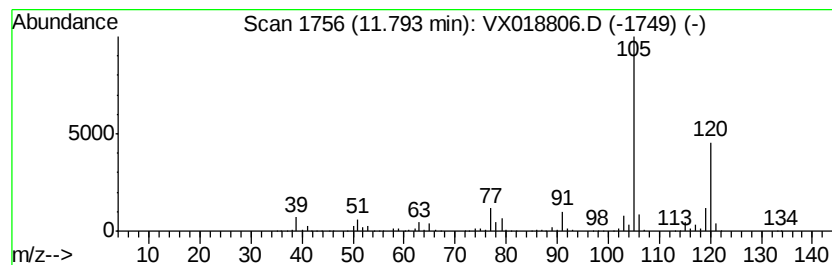
Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

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 11 Benzene, 1,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.79	5.86 ug/L	398299	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	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
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\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

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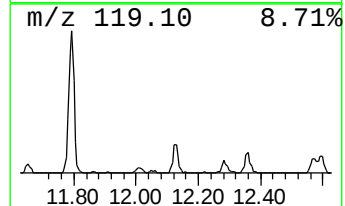
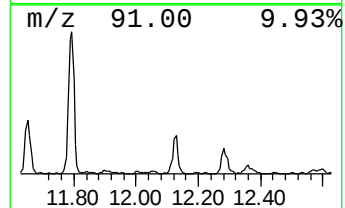
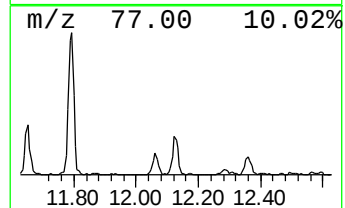
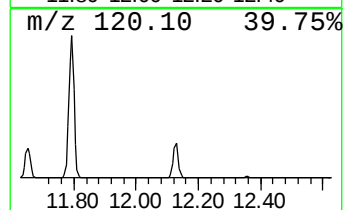
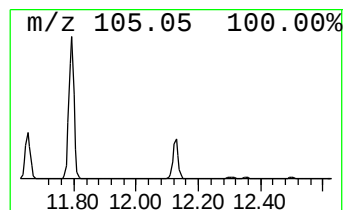
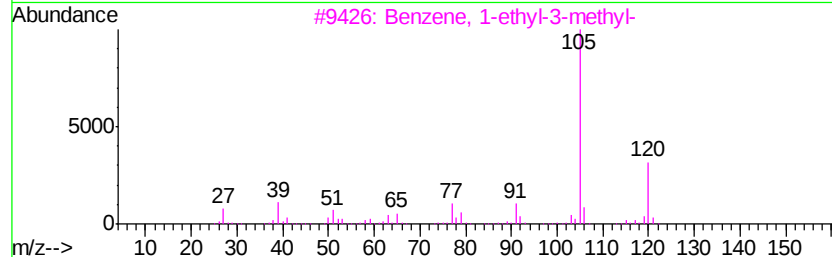
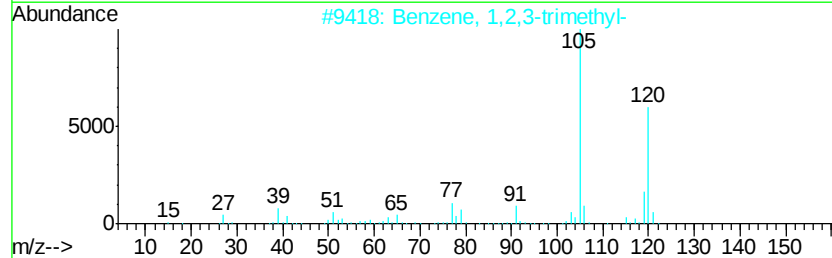
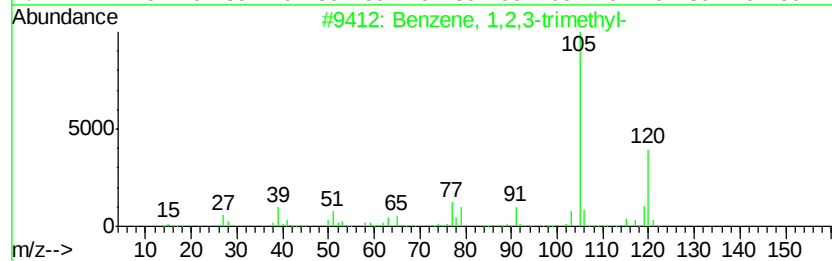
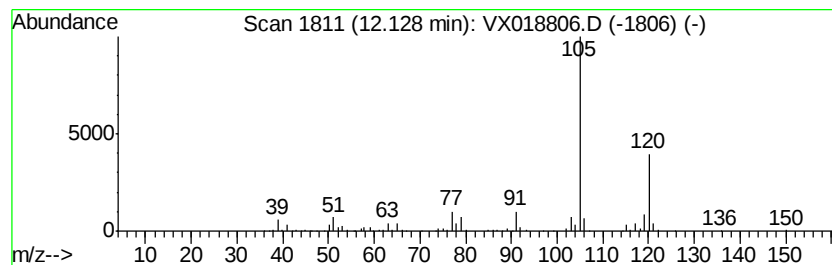
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Mesitylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	1.67 ug/L	113247	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	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100620\
Data File : VX018806.D
Acq On : 06 Oct 2020 18:05
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BG3N5DL2

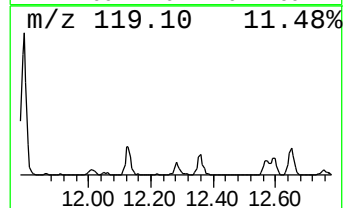
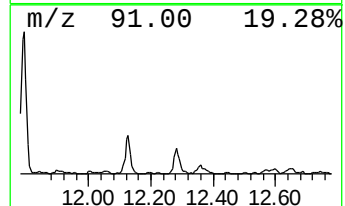
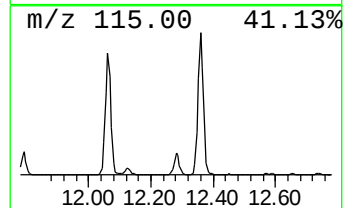
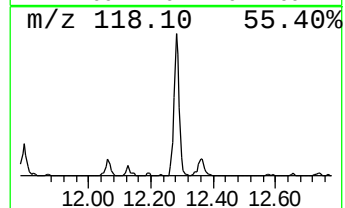
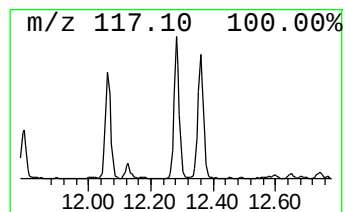
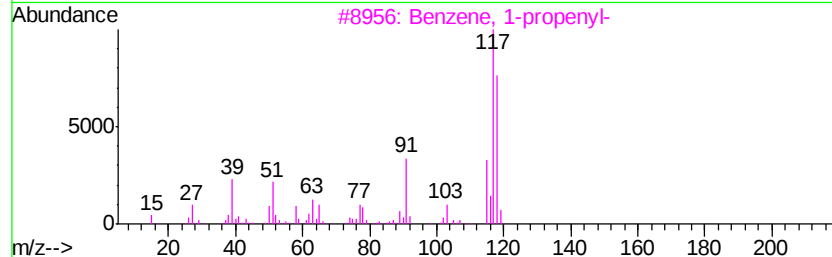
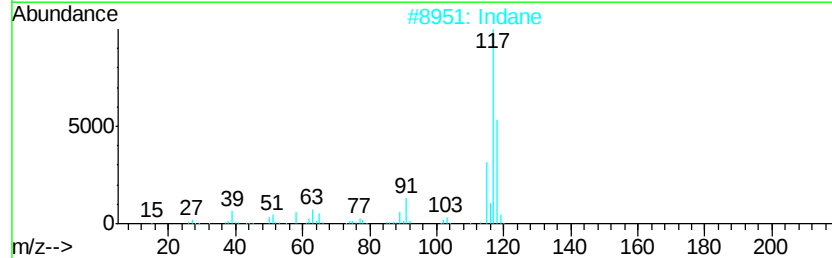
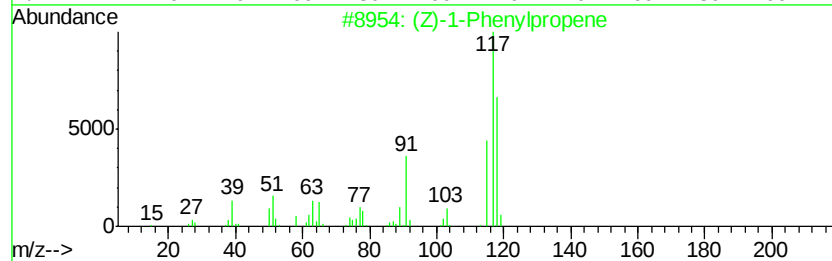
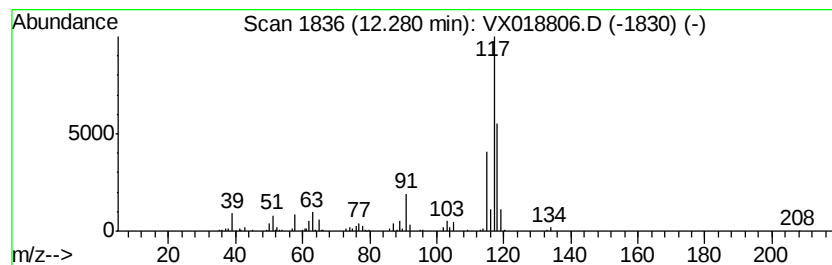
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 (Z)-1-Phenylpropene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	0.84 ug/L	56853	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	70
2		Indane	118	C9H10	000496-11-7	68
3		Benzene, 1-propenyl-	118	C9H10	000637-50-3	62
4		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	58
5		7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	58



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100620\
 Data File : VX018806.D
 Acq On : 06 Oct 2020 18:05
 Operator : JC/SP
 Sample : L4239-05DL2 400X
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N5DL2

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...	2.87	0.5	ug/L	28147	1	6.83	280112	5.0
(DEL) Alkane: Str...	3.15	1.1	ug/L	64413	1	6.83	280112	5.0
(DEL) Alkane: Str...	3.50	2.0	ug/L	114695	1	6.83	280112	5.0
(DEL) Alkane: Cyc...	4.37	1.7	ug/L	97089	1	6.83	280112	5.0
Benzene, propyl-	11.34	0.8	ug/L	54260	3	12.06	339935	5.0
Benzene, 1-ethyl-...	11.42	4.6	ug/L	312616	3	12.06	339935	5.0
Benzene, 1,2,3-tr...	11.49	1.7	ug/L	116009	3	12.06	339935	5.0
4-Heptanone, 2,6-...	11.61	3.6	ug/L	247907	3	12.06	339935	5.0
Benzene, 1-ethyl-...	11.65	1.9	ug/L	127362	3	12.06	339935	5.0
Benzene, 1,2,4-tr...	11.79	5.9	ug/L	398299	3	12.06	339935	5.0
Mesitylene	12.13	1.7	ug/L	113247	3	12.06	339935	5.0
(Z)-1-Phenylpropene	12.28	0.8	ug/L	56853	3	12.06	339935	5.0