

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092718\
 Data File : VV007935.D
 Acq On : 28 Sep 2018 04:02
 Operator : SY/MD
 Sample : J5062-05
 Misc : 25.0 mL/MSVOA V/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAC05

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_V\METHOD\SOMVTR092718WMA.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	1.315	63	70	86	rVB	39863	49004	4.44%	1.003%
2	1.405	91	98	115	rBV2	11307	24171	2.19%	0.495%
3	1.582	146	153	165	rVB	34501	40030	3.62%	0.819%
4	2.129	314	323	335	rBV	77868	114416	10.36%	2.341%
5	2.482	423	433	468	rVB	63003	165948	15.02%	3.395%
6	3.154	633	642	655	rVB3	8841	17225	1.56%	0.352%
7	3.859	849	861	878	rVB5	4975	13138	1.19%	0.269%
8	3.994	884	903	936	rBV	20432	93682	8.48%	1.917%
9	4.164	944	956	973	rVB	8358	22936	2.08%	0.469%
10	4.408	1015	1032	1054	rBV2	55751	153146	13.87%	3.133%
11	4.579	1070	1085	1106	rBV2	14824	40490	3.67%	0.828%
12	5.097	1228	1246	1269	rBV2	131664	293794	26.60%	6.011%
13	5.270	1276	1300	1331	rBV3	12018	82792	7.50%	1.694%
14	5.666	1410	1423	1440	rVB	107484	203916	18.46%	4.172%
15	5.849	1465	1480	1495	rBV3	8321	17666	1.60%	0.361%
16	5.968	1499	1517	1519	rBV6	5531	12705	1.15%	0.260%
17	6.122	1555	1565	1591	rVB	76877	162824	14.74%	3.331%
18	6.415	1648	1656	1679	rVB	64131	126529	11.46%	2.589%
19	6.666	1725	1734	1751	rVB2	15175	28711	2.60%	0.587%
20	7.035	1834	1849	1867	rVB	33284	59567	5.39%	1.219%
21	7.363	1937	1951	1966	rBV	152538	259686	23.51%	5.313%
22	7.669	2033	2046	2056	rBV	20242	37476	3.39%	0.767%
23	7.775	2075	2079	2100	rVB6	4729	11395	1.03%	0.233%
24	7.997	2134	2148	2182	rBV4	39319	145912	13.21%	2.985%
25	8.145	2183	2194	2225	rVB	165903	319698	28.94%	6.541%
26	8.293	2231	2240	2253	rVB4	13800	23019	2.08%	0.471%
27	8.495	2292	2303	2316	rBV2	15969	27176	2.46%	0.556%
28	8.900	2416	2429	2453	rVB	157564	252551	22.87%	5.167%
29	9.534	2613	2626	2665	rBV	550108	1104524	100.00%	22.597%
30	9.752	2688	2694	2709	rVB2	7309	11404	1.03%	0.233%
31	9.977	2753	2764	2783	rBV	61793	91635	8.30%	1.875%
32	10.267	2843	2854	2868	rVB2	53714	83246	7.54%	1.703%
33	10.865	3029	3040	3055	rBV	72624	128719	11.65%	2.633%
34	10.929	3055	3060	3075	rVB6	6894	12355	1.12%	0.253%

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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	11.299	3157	3175	3191	rVB	170417	275108	24.91%	5.628%
36	11.579	3252	3262	3281	rBV3	32819	59349	5.37%	1.214%
37	11.675	3281	3292	3310	rVB	164802	256339	23.21%	5.244%
38	12.058	3399	3411	3430	rBV	35688	65662	5.94%	1.343%

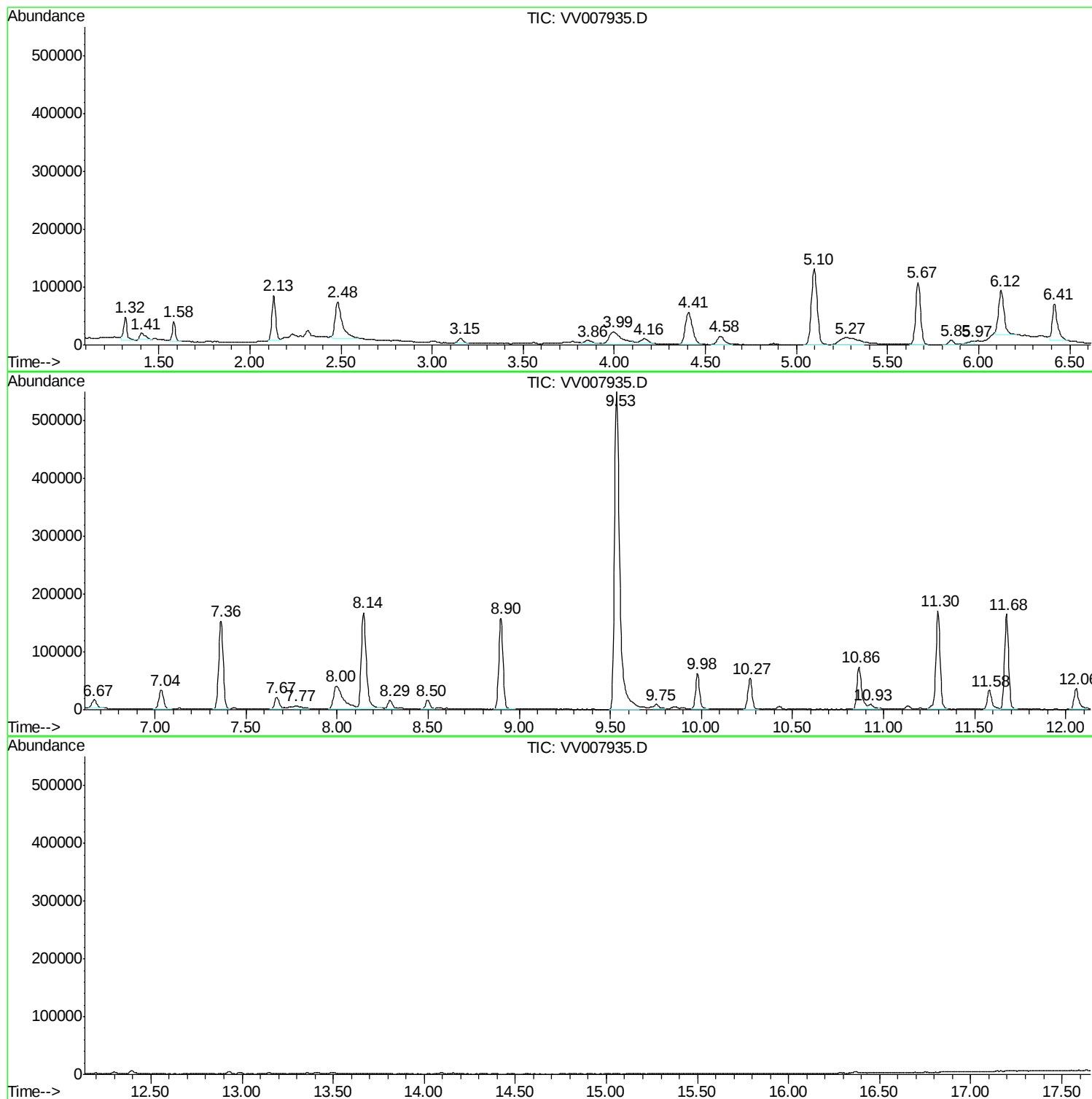
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Quant Title : TRACE VOA SOM01.0

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



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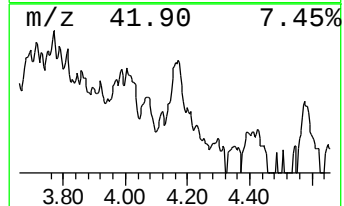
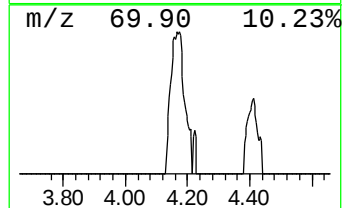
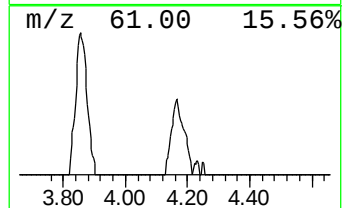
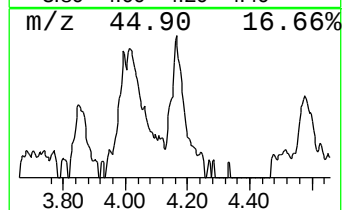
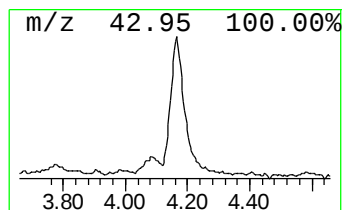
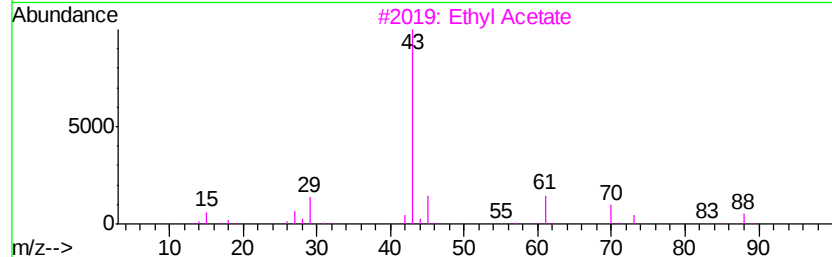
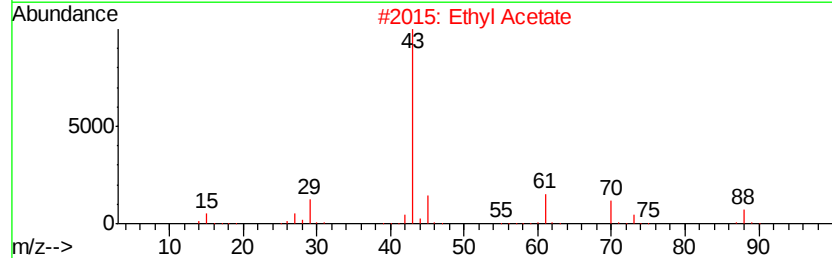
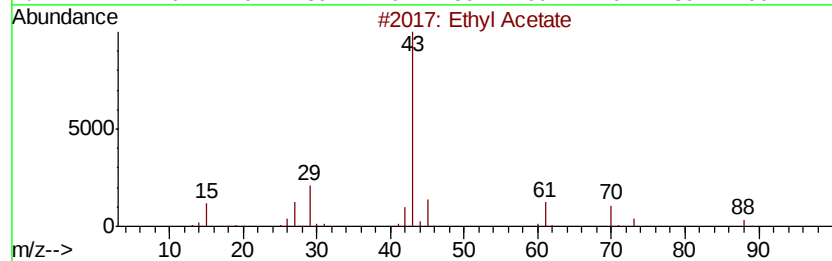
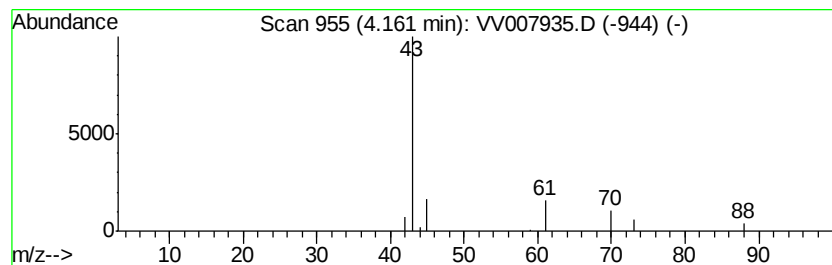
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Ethyl Acetate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.16	0.56 ug/L	22936	1,4-Difluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl Acetate	88	C4H8O2	000141-78-6	83
2			Ethyl Acetate	88	C4H8O2	000141-78-6	83
3			Ethyl Acetate	88	C4H8O2	000141-78-6	83
4			Ethyl Acetate	88	C4H8O2	000141-78-6	83
5			CH3C(O)CH2CH2OH	88	C4H8O2	000590-90-9	9



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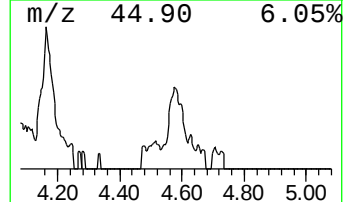
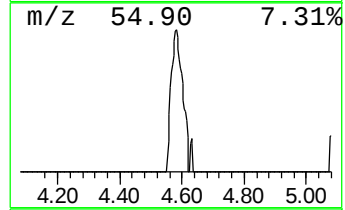
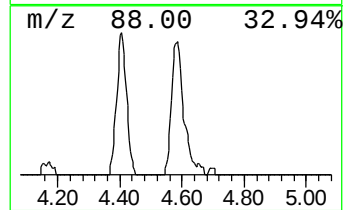
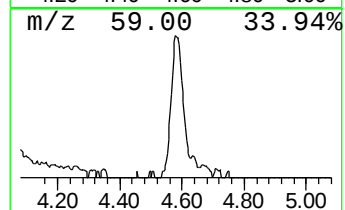
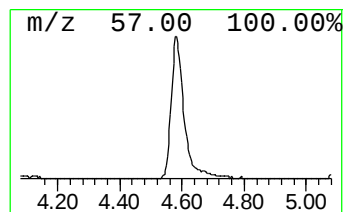
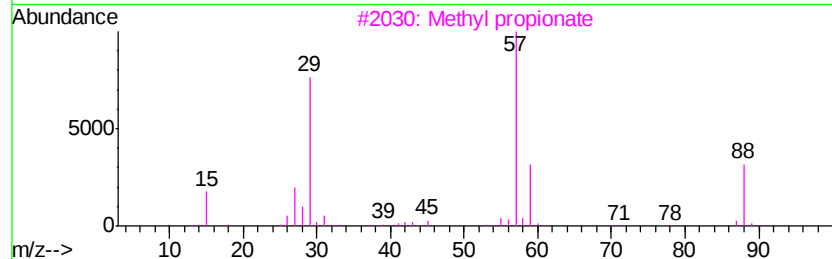
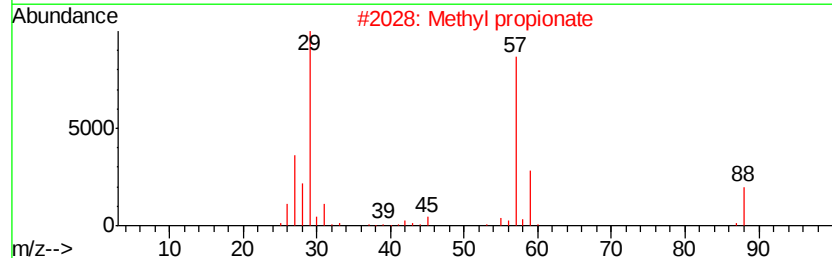
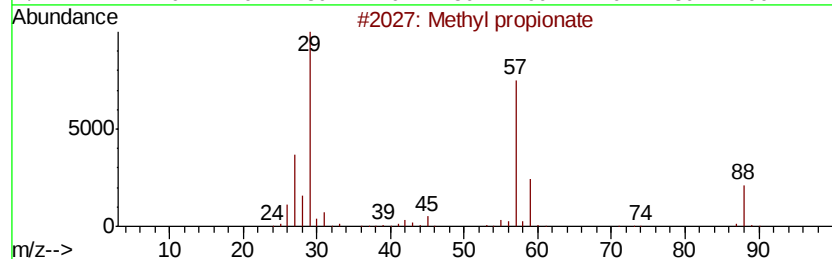
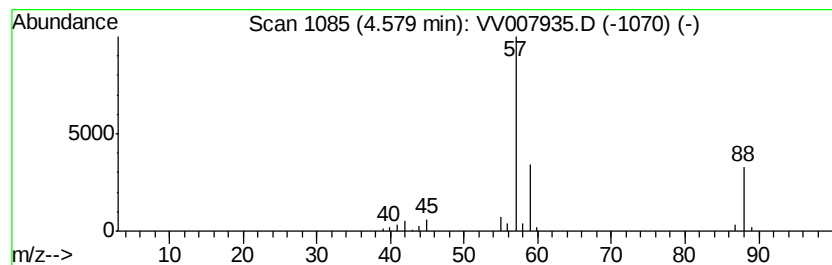
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Methyl propionate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.58	0.99 ug/L	40490	1,4-Difluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl propionate	88	C4H8O2	000554-12-1	90
2			Methyl propionate	88	C4H8O2	000554-12-1	90
3			Methyl propionate	88	C4H8O2	000554-12-1	80
4			Methyl propionate	88	C4H8O2	000554-12-1	72
5			Propylamine	59	C3H9N	000107-10-8	9



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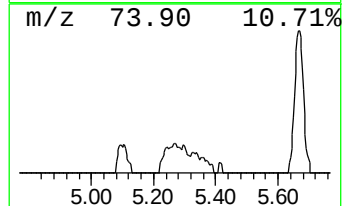
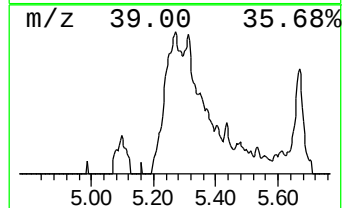
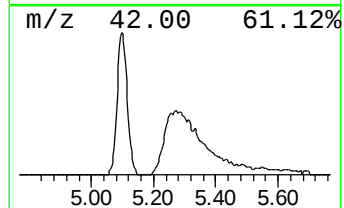
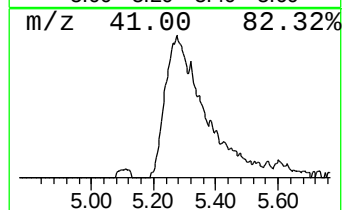
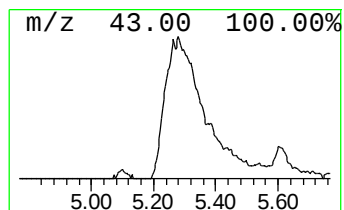
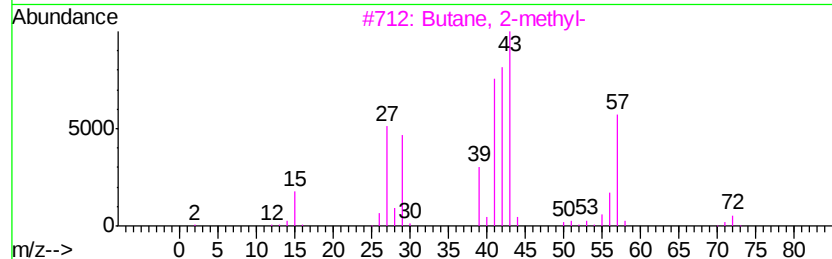
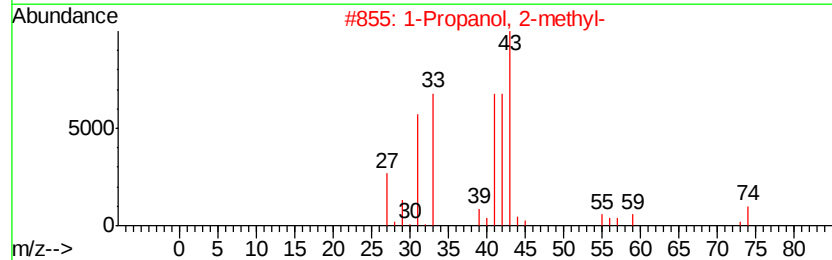
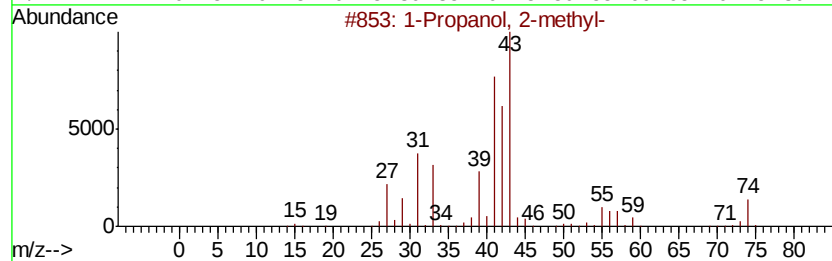
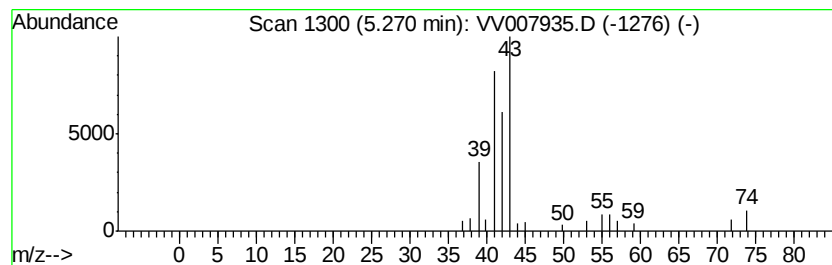
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 1-Propanol, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.27	2.03 ug/L	82792	1,4-Difluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	72
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	72
3		Butane, 2-methyl-	72	C5H12	000078-78-4	59
4		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	50
5		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	50



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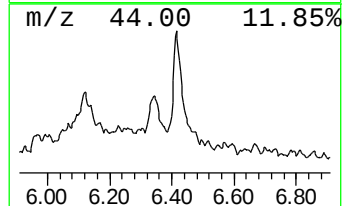
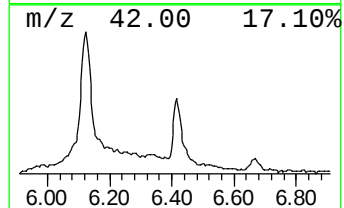
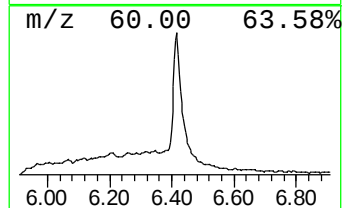
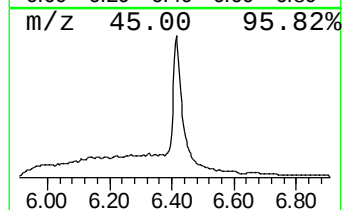
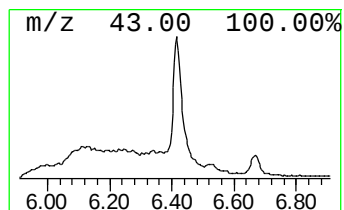
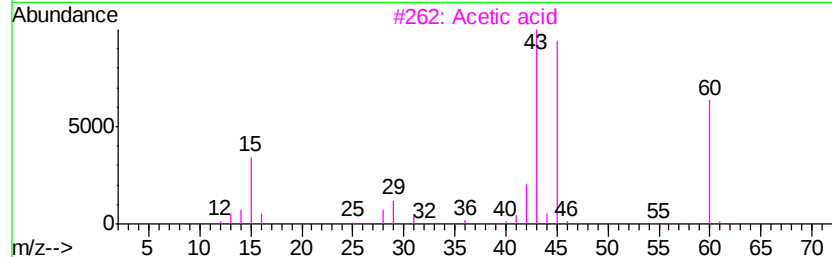
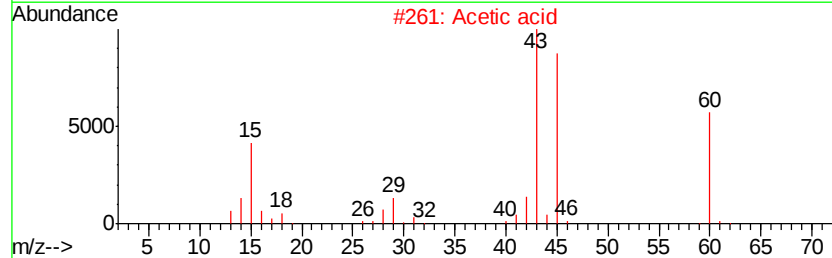
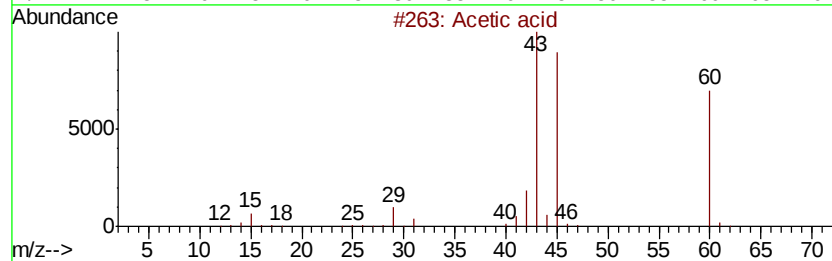
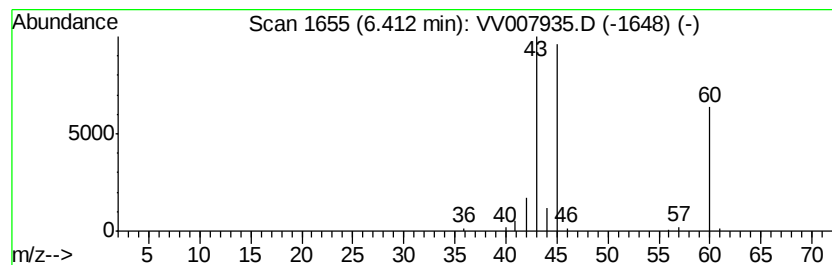
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Acetic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	3.10 ug/L	126529	1,4-Difluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid			60	C2H4O2	000064-19-7	90
2	Acetic acid			60	C2H4O2	000064-19-7	90
3	Acetic acid			60	C2H4O2	000064-19-7	90
4	Acetic acid			60	C2H4O2	000064-19-7	78
5	Acetic acid			60	C2H4O2	000064-19-7	78



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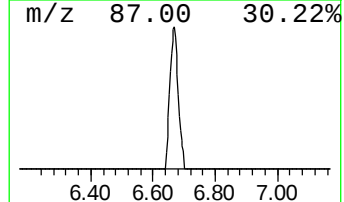
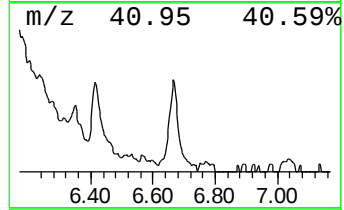
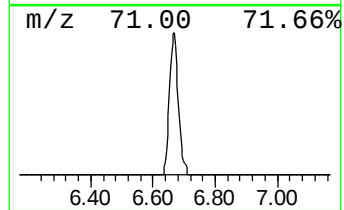
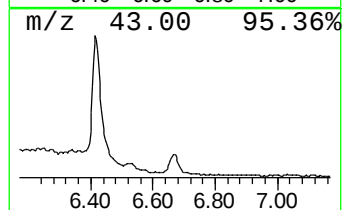
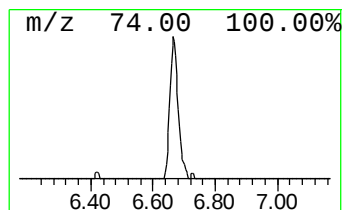
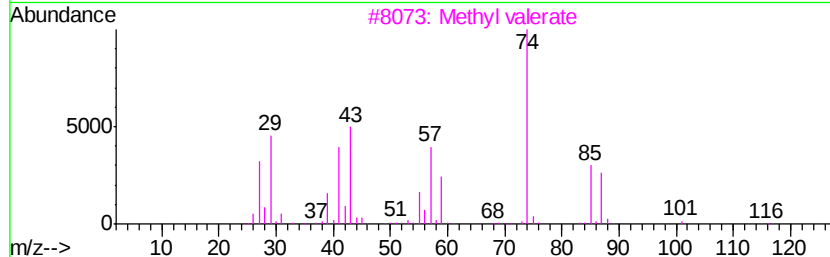
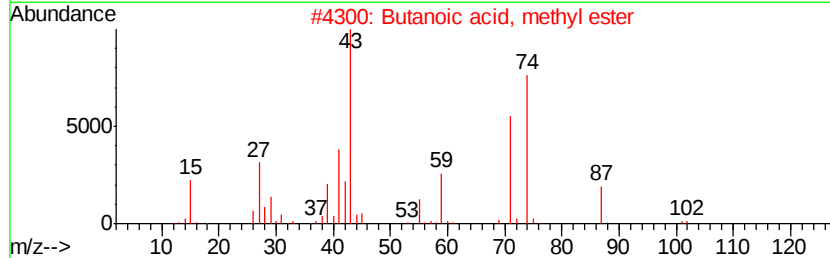
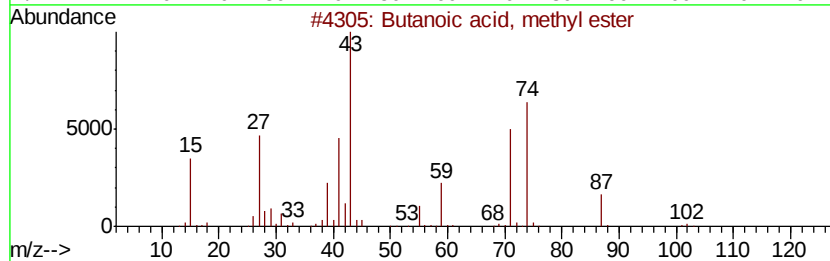
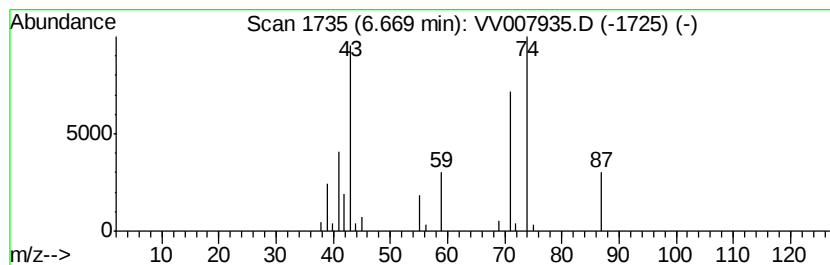
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Peak Number 6 Butanoic acid, methyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	0.70 ug/L	28711	1,4-Difluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	78
2		Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	74
3		Methyl valerate	116	C6H12O2	000624-24-8	28
4		Methyl valerate	116	C6H12O2	000624-24-8	28
5		3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

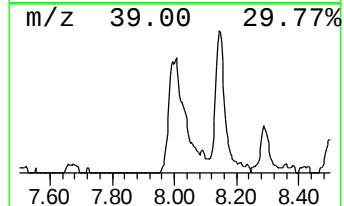
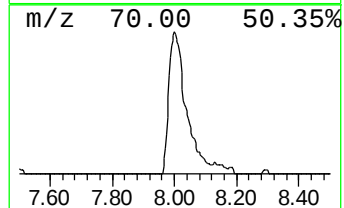
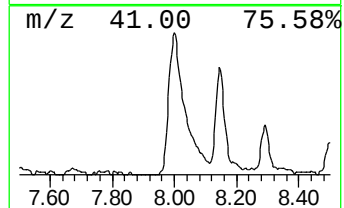
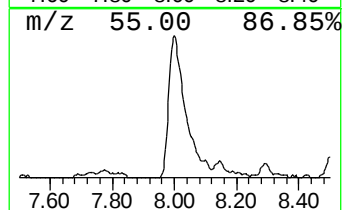
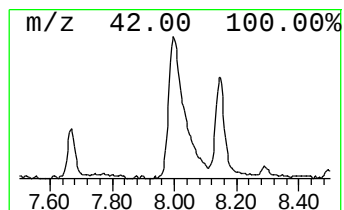
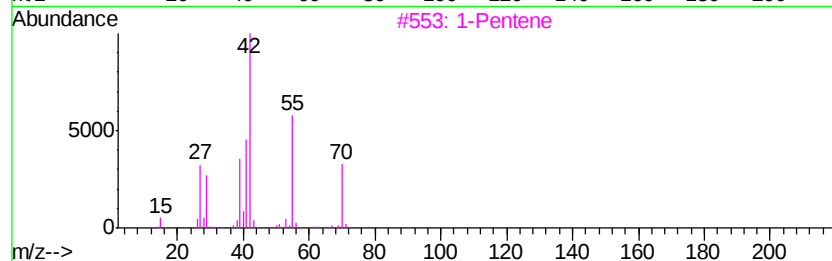
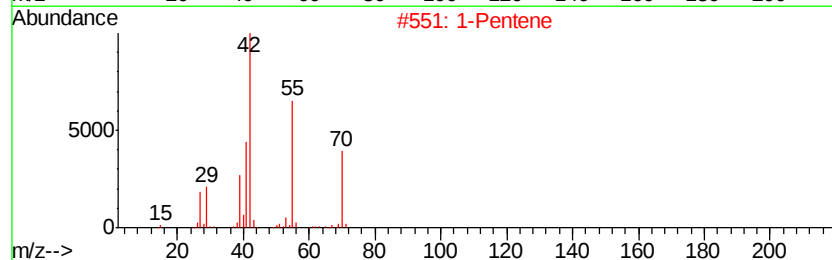
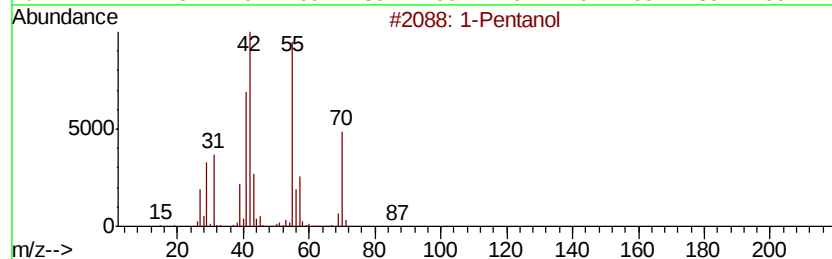
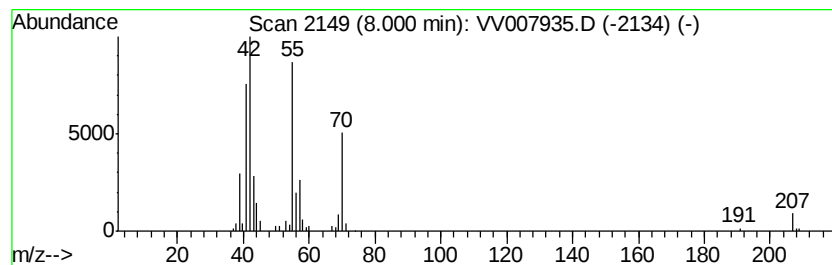
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 1-Pentanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.00	2.89 ug/L	145912	Chlorobenzene-d5	8.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentanol	88	C5H12O	000071-41-0	72
2		1-Pentene	70	C5H10	000109-67-1	72
3		1-Pentene	70	C5H10	000109-67-1	72
4		1-Pentanol	88	C5H12O	000071-41-0	72
5		1-Pentanol	88	C5H12O	000071-41-0	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

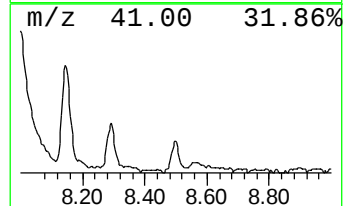
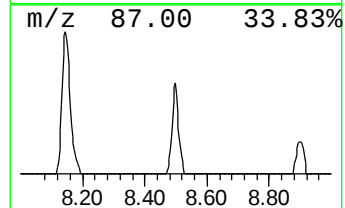
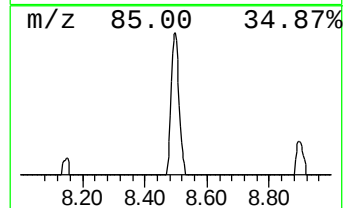
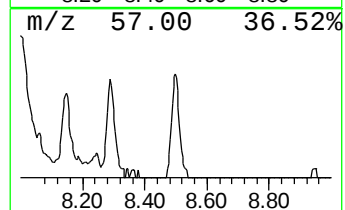
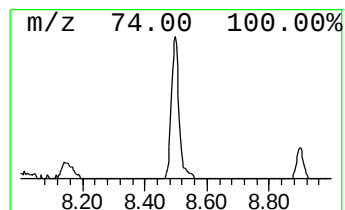
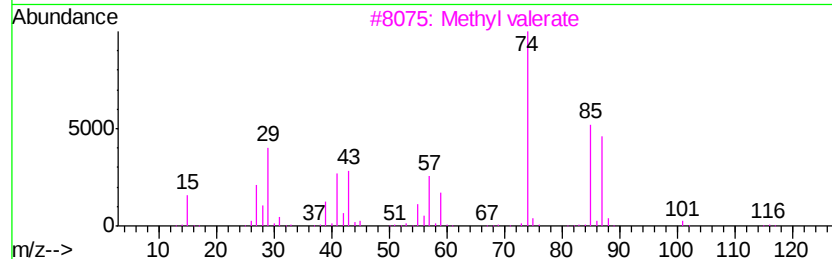
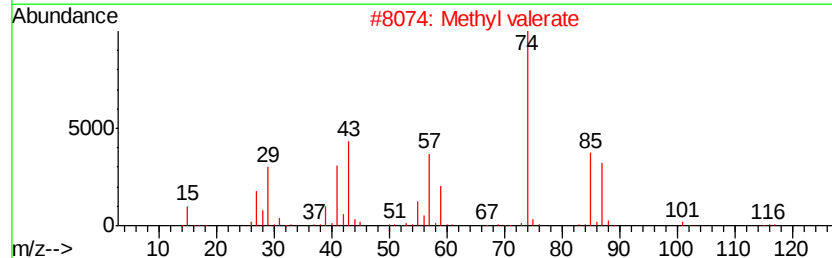
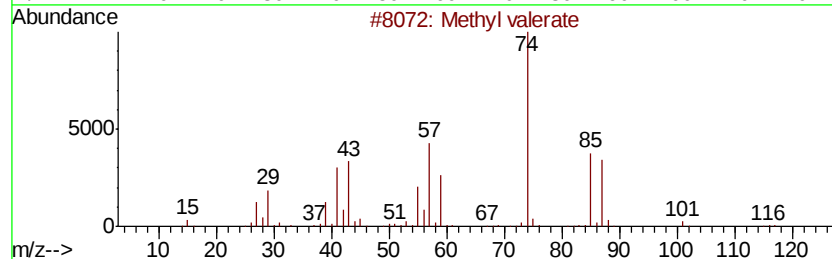
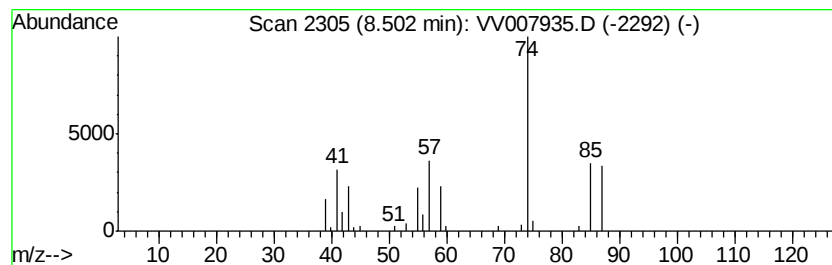
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Methyl valerate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.50	0.54 ug/L	27176	Chlorobenzene-d5	8.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl valerate	116	C6H12O2	000624-24-8	83
2		Methyl valerate	116	C6H12O2	000624-24-8	78
3		Methyl valerate	116	C6H12O2	000624-24-8	72
4		Methyl valerate	116	C6H12O2	000624-24-8	64
5		Propanoic acid	74	C3H6O2	000079-09-4	23



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

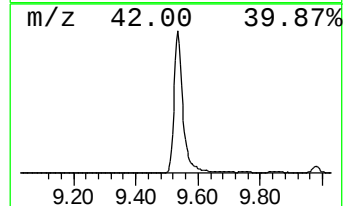
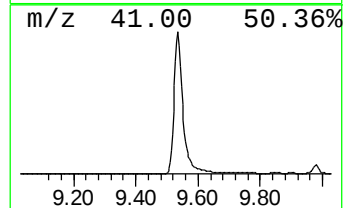
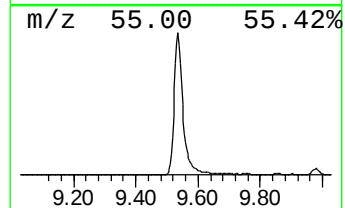
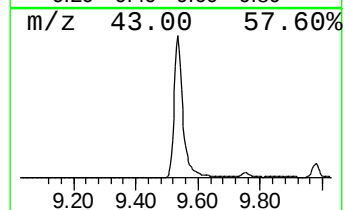
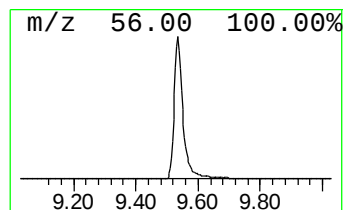
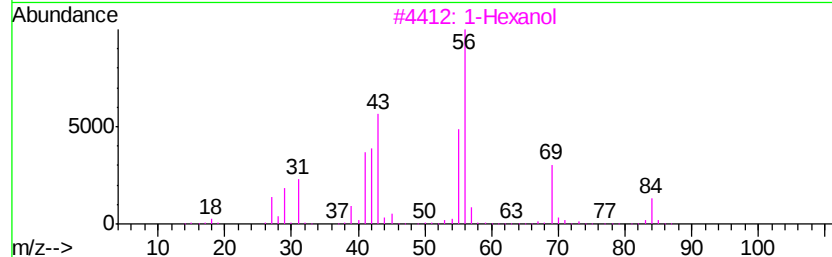
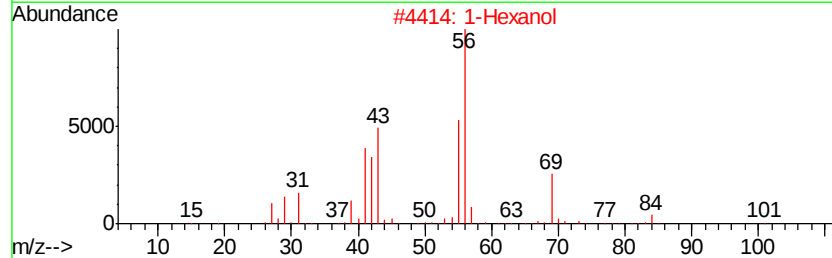
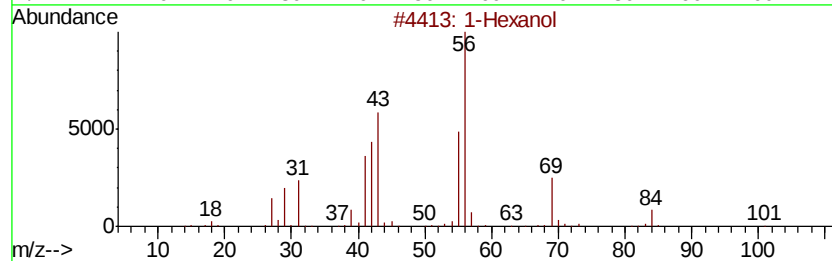
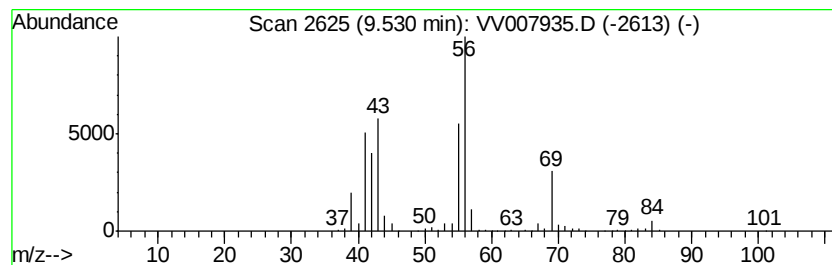
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 1-Hexanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	21.87 ug/L	1104520	Chlorobenzene-d5	8.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol	102	C6H14O	000111-27-3	90
2		1-Hexanol	102	C6H14O	000111-27-3	83
3		1-Hexanol	102	C6H14O	000111-27-3	78
4		1-Hexanol	102	C6H14O	000111-27-3	78
5		1-Hexanol	102	C6H14O	000111-27-3	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

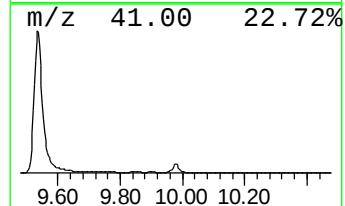
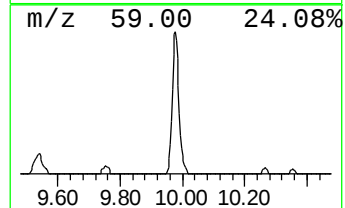
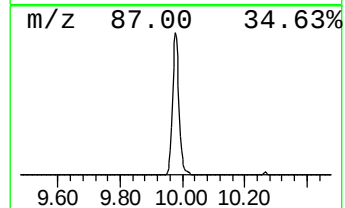
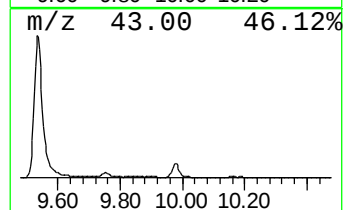
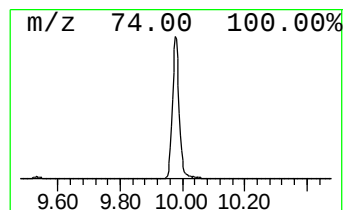
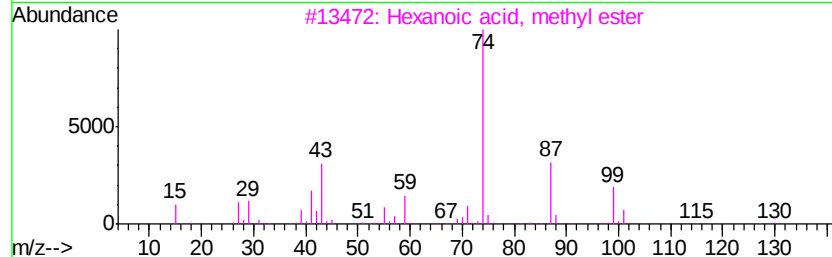
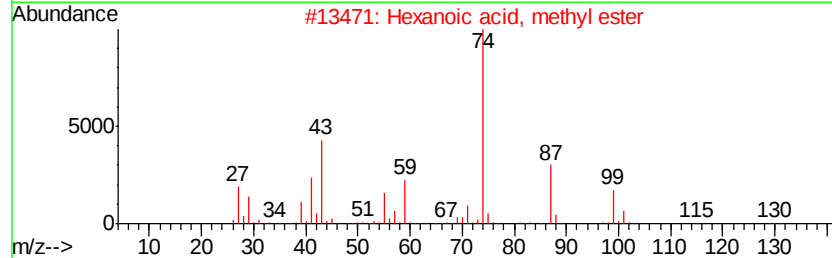
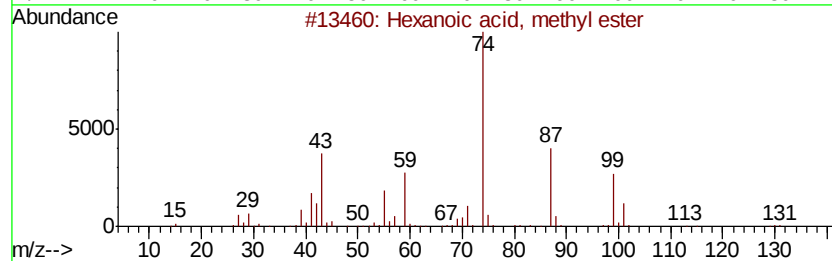
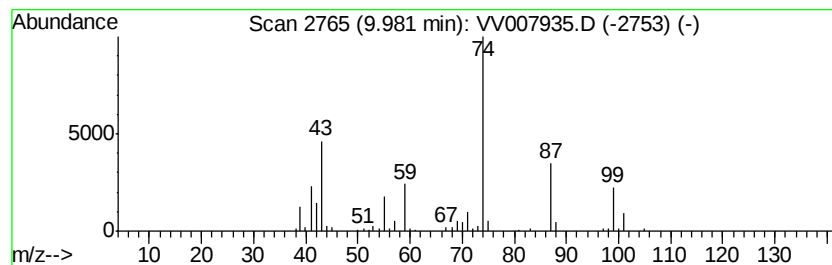
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Hexanoic acid, methyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.98	1.81 ug/L	91635	Chlorobenzene-d5	8.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, methyl ester	130	C7H14O2	000106-70-7	90
2			Hexanoic acid, methyl ester	130	C7H14O2	000106-70-7	78
3			Hexanoic acid, methyl ester	130	C7H14O2	000106-70-7	72
4			Pentanoic acid, 3-methyl-, methyl...	130	C7H14O2	002177-78-8	59
5			Methyl valerate	116	C6H12O2	000624-24-8	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

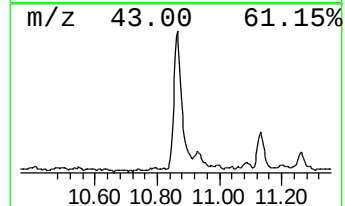
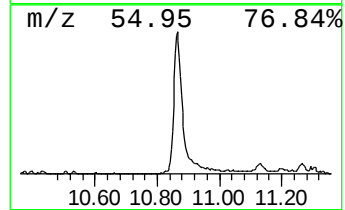
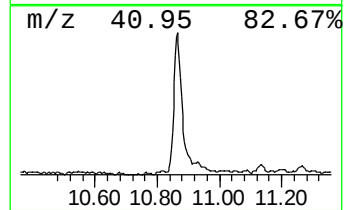
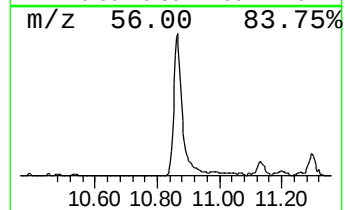
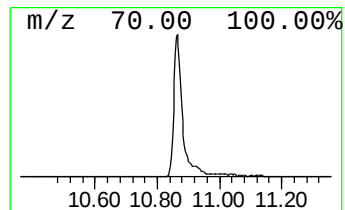
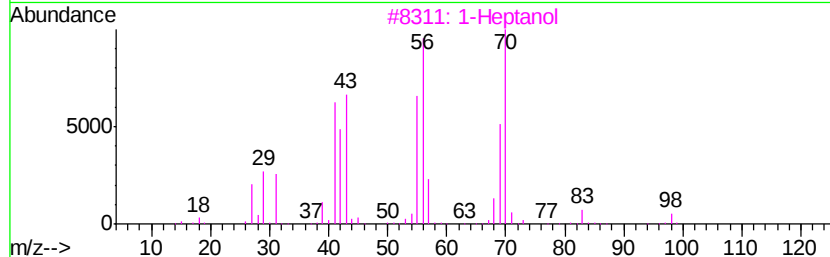
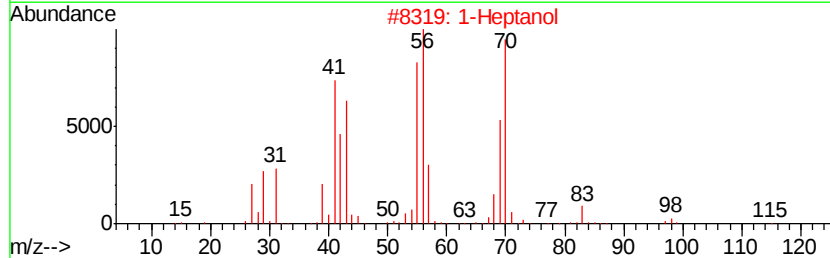
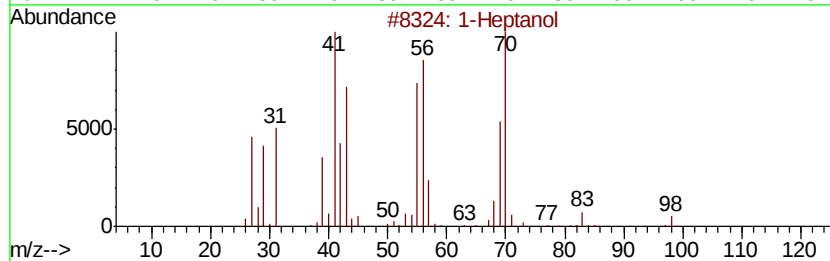
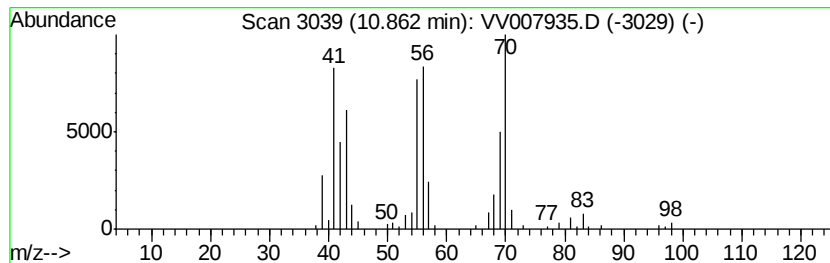
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 1-Heptanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.86	2.34 ug/L	128719	1,4-Dichlorobenzene-d4	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptanol		116	C7H16O	000111-70-6	86
2	1-Heptanol		116	C7H16O	000111-70-6	86
3	1-Heptanol		116	C7H16O	000111-70-6	72
4	Formic acid, heptyl ester		144	C8H16O2	000112-23-2	72
5	Formic acid, heptyl ester		144	C8H16O2	000112-23-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

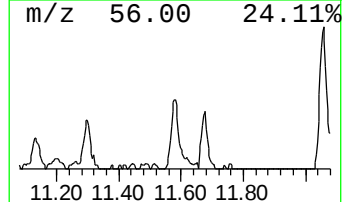
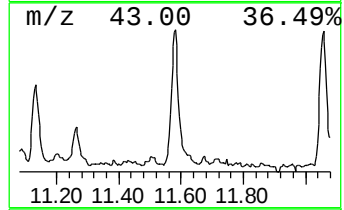
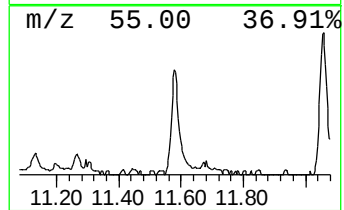
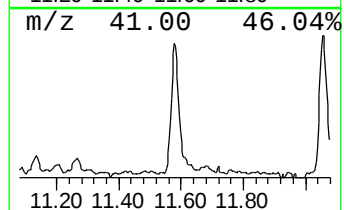
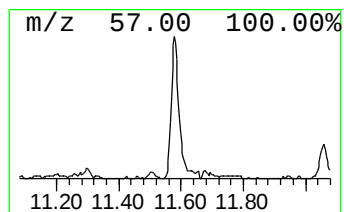
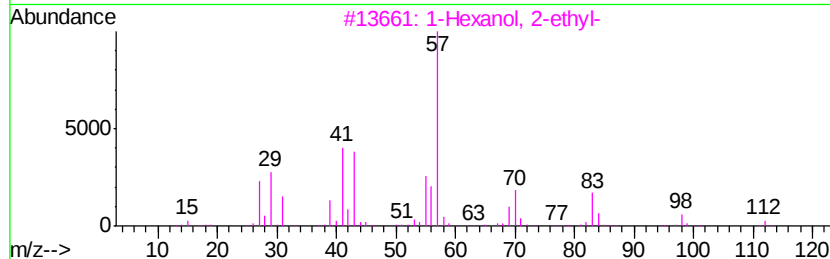
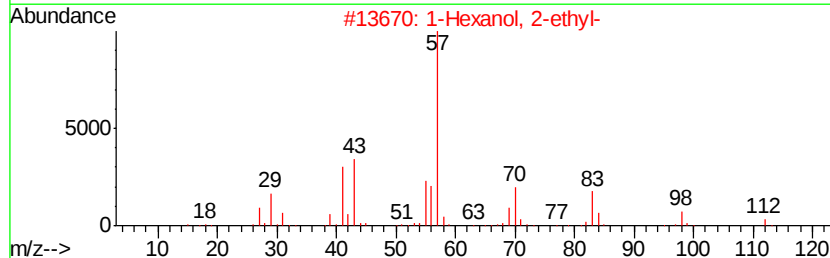
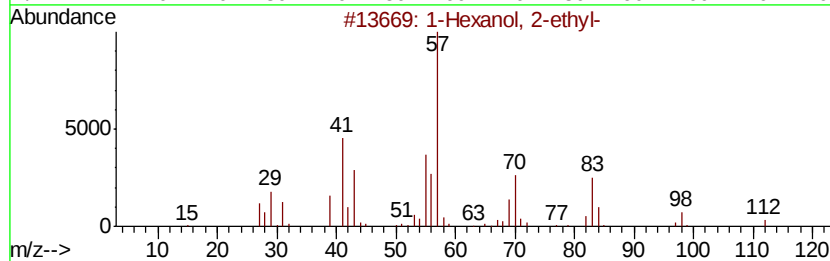
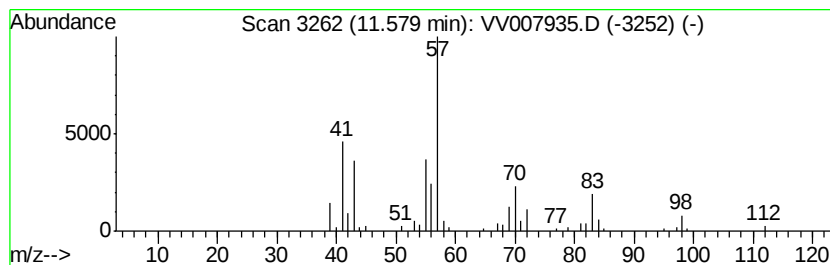
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 1-Hexanol, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.58	1.08 ug/L	59349	1,4-Dichlorobenzene-d4	11.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
4			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	53
5			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092718\
Data File : VV007935.D
Acq On : 28 Sep 2018 04:02
Operator : SY/MD
Sample : J5062-05
Misc : 25.0 mL/MSVOA V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAC05

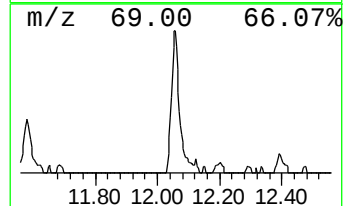
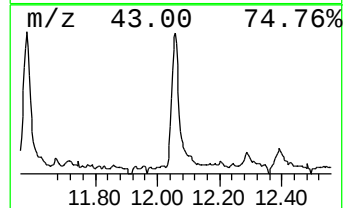
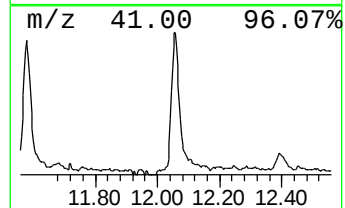
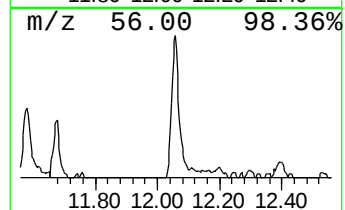
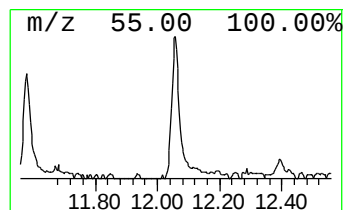
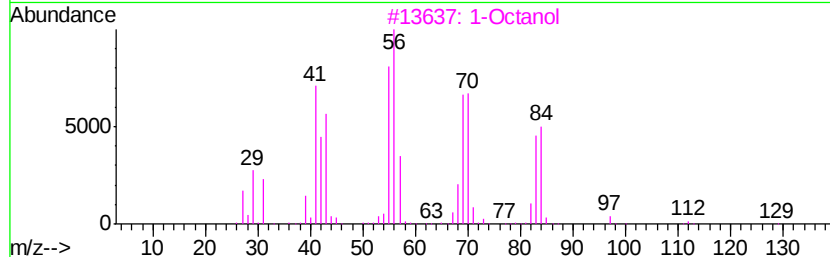
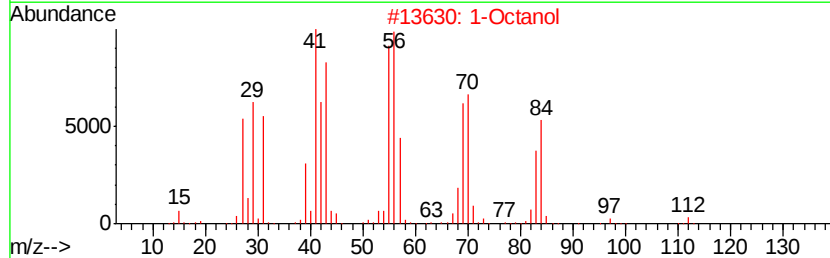
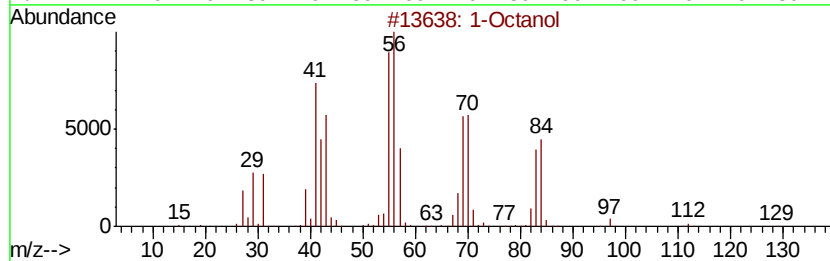
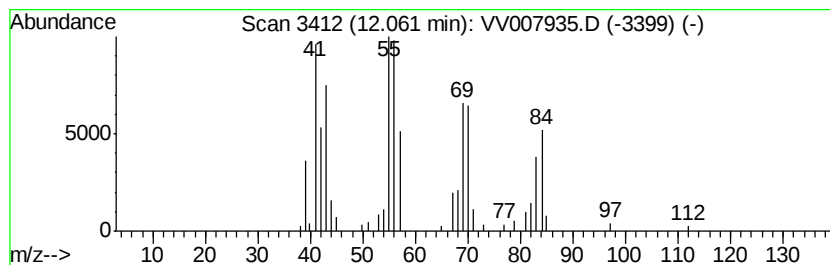
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 1-Octanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	1.19 ug/L	65662	1,4-Dichlorobenzene-d4	11.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octanol	130	C8H18O	000111-87-5	91
2			1-Octanol	130	C8H18O	000111-87-5	90
3			1-Octanol	130	C8H18O	000111-87-5	90
4			Formic acid, octyl ester	158	C9H18O2	000112-32-3	90
5			1-Octanol	130	C8H18O	000111-87-5	86



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092718\
 Data File : VV007935.D
 Acq On : 28 Sep 2018 04:02
 Operator : SY/MD
 Sample : J5062-05
 Misc : 25.0 mL/MSVOA_V/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAC05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.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
Ethyl Acetate	4.16	0.6	ug/L	22936	1	5.67	203916	5.0
Methyl propionate	4.58	1.0	ug/L	40490	1	5.67	203916	5.0
1-Propanol, 2-met...	5.27	2.0	ug/L	82792	1	5.67	203916	5.0
Acetic acid	6.41	3.1	ug/L	126529	1	5.67	203916	5.0
Butanoic acid, me...	6.67	0.7	ug/L	28711	1	5.67	203916	5.0
1-Pentanol	8.00	2.9	ug/L	145912	2	8.90	252551	5.0
Methyl valerate	8.50	0.5	ug/L	27176	2	8.90	252551	5.0
1-Hexanol	9.53	21.9	ug/L	1104520	2	8.90	252551	5.0
Hexanoic acid, me...	9.98	1.8	ug/L	91635	2	8.90	252551	5.0
1-Heptanol	10.86	2.3	ug/L	128719	3	11.30	275108	5.0
1-Hexanol, 2-ethyl-	11.58	1.1	ug/L	59349	3	11.30	275108	5.0
1-Octanol	12.06	1.2	ug/L	65662	3	11.30	275108	5.0