

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012516\
 Data File : BF084609.D
 Acq On : 25 Jan 2016 16:09
 Operator : SJ/IZ
 Sample : H1201-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-02

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.201	183	185	191	rBV	380761	592240	7.19%	0.984%
2	4.361	197	199	202	rVB	106628	145215	1.76%	0.241%
3	4.898	243	246	249	rBV	1954414	2528871	30.72%	4.202%
4	5.333	282	284	287	rBV	3432184	3370952	40.95%	5.601%
5	5.733	317	319	322	rBV	504773	476551	5.79%	0.792%
6	6.384	373	376	378	rBV	2118593	2415899	29.35%	4.014%
7	6.510	384	387	389	rVB	5748922	6374783	77.44%	10.591%
8	6.579	391	393	395	rVB	222956	234016	2.84%	0.389%
9	6.739	404	407	409	rBV	1080830	1301628	15.81%	2.163%
10	6.887	418	420	423	rBV	3753310	5163588	62.73%	8.579%
11	7.310	453	457	459	rBV	4301133	5178551	62.91%	8.604%
12	8.019	517	519	522	rBV	1883189	1746937	21.22%	2.902%
13	9.105	611	614	616	rBV	7767779	7826027	95.07%	13.002%
14	9.493	646	648	654	rBV	66671	104815	1.27%	0.174%
15	9.768	670	672	675	rBV	1536275	2041096	24.80%	3.391%
16	9.813	675	676	680	rBV3	74482	102130	1.24%	0.170%
17	10.568	738	742	744	rBV	5068302	5971633	72.55%	9.921%
18	11.253	799	802	804	rBV	2548803	2190971	26.62%	3.640%
19	11.802	847	850	860	rBV2	792183	899935	10.93%	1.495%
20	12.591	917	919	925	rBV	155702	250591	3.04%	0.416%
21	12.842	938	941	944	rVB	5692445	8231605	100.00%	13.676%
22	13.745	1018	1020	1024	rBV	109064	167167	2.03%	0.278%
23	13.882	1030	1032	1035	rBV	1631530	1725980	20.97%	2.868%
24	15.277	1151	1154	1156	rBV	896387	1148452	13.95%	1.908%

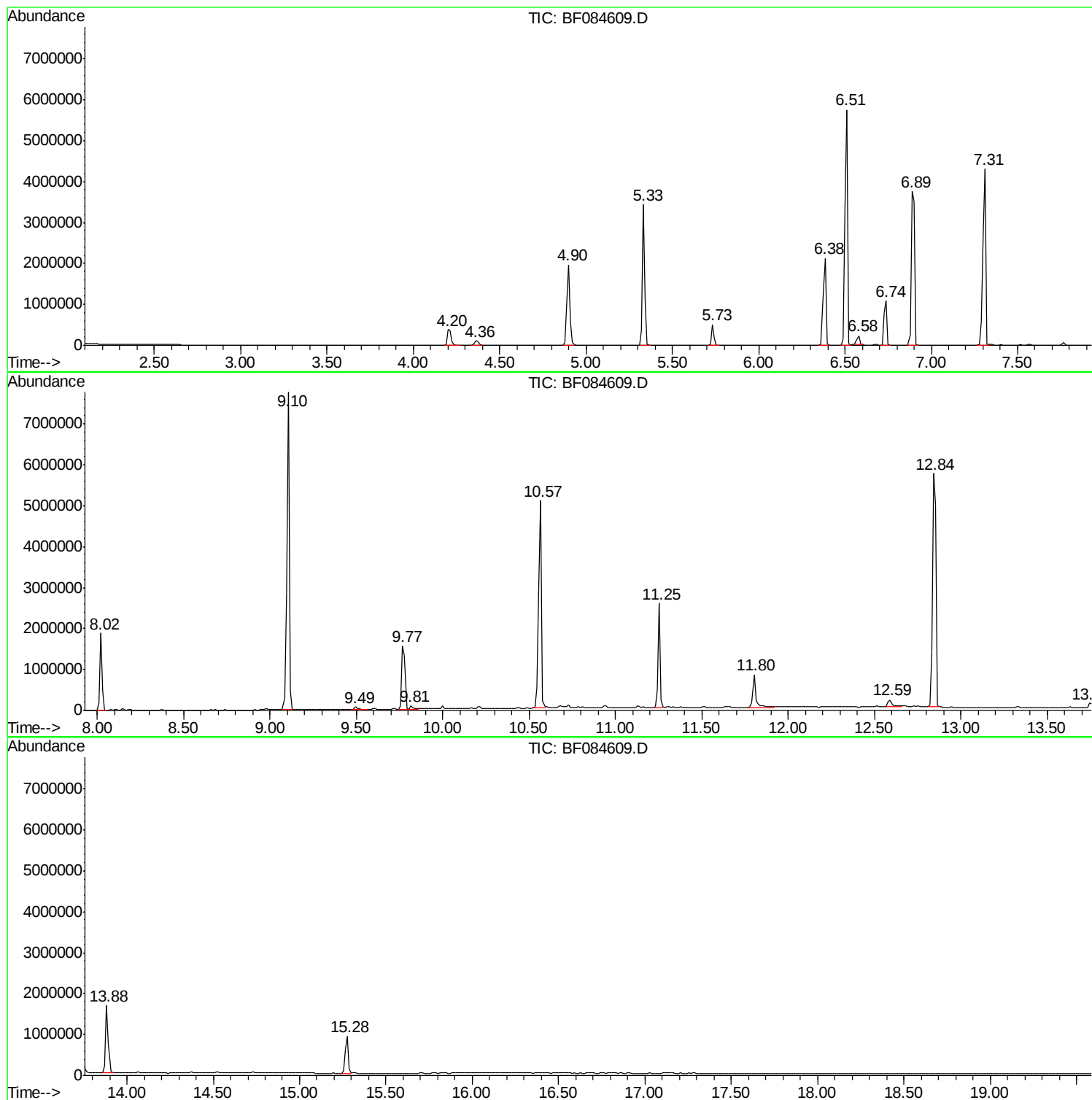
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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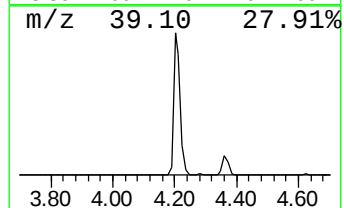
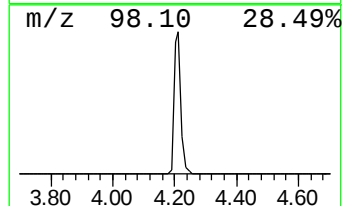
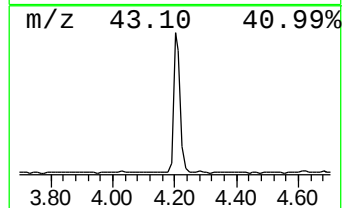
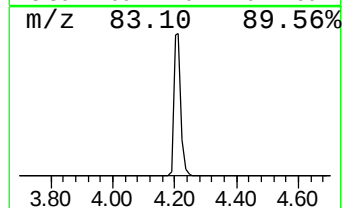
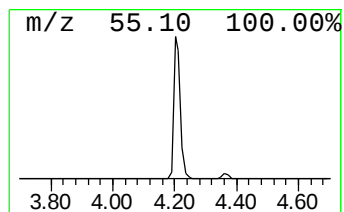
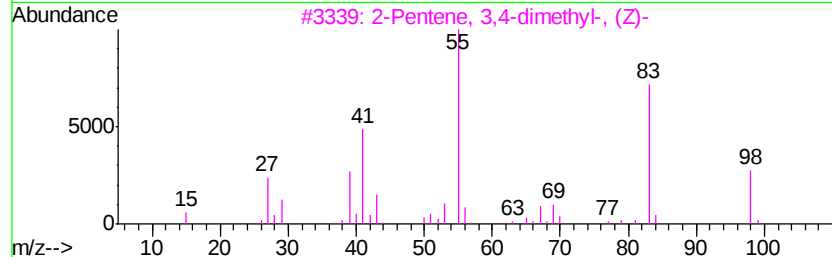
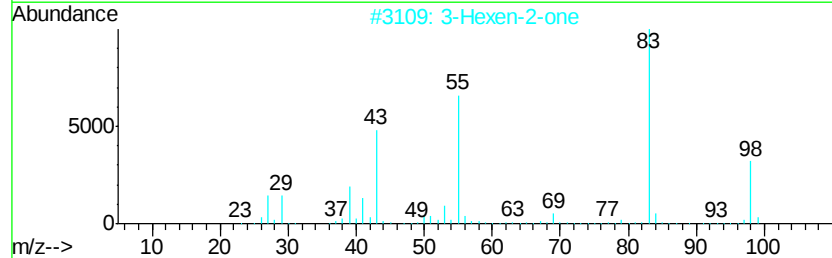
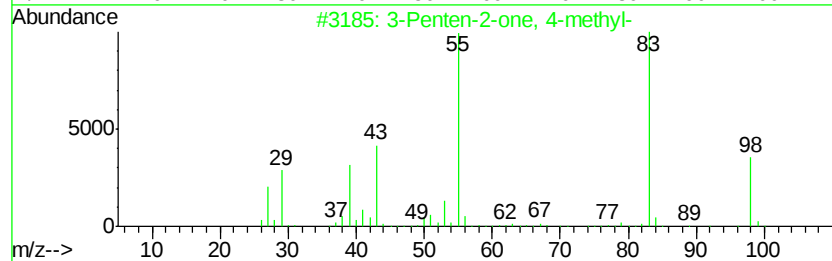
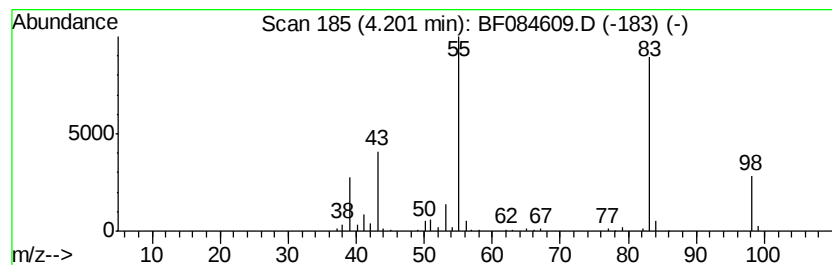
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	9.10 ng	592240	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2			3-Hexen-2-one	98	C6H10O	000763-93-9	90
3			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
4			2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86
5			2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	72



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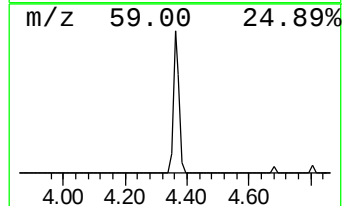
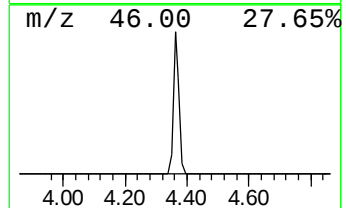
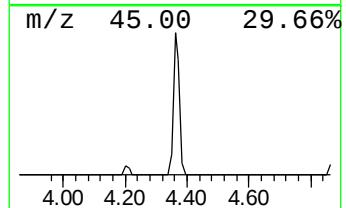
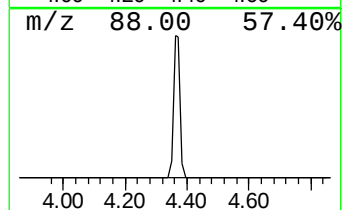
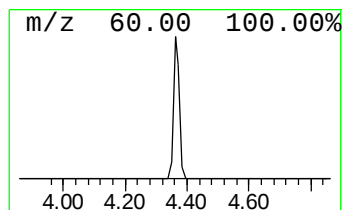
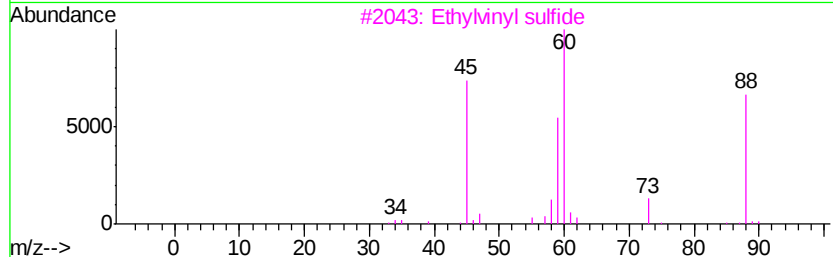
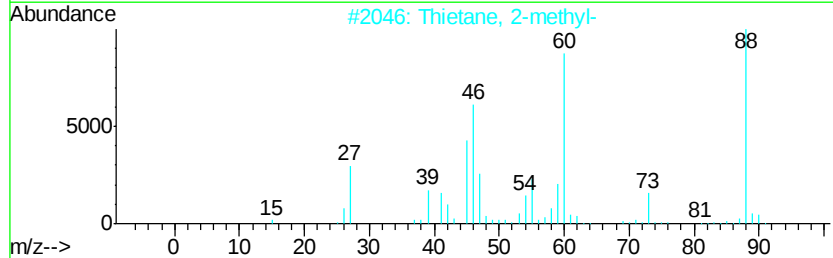
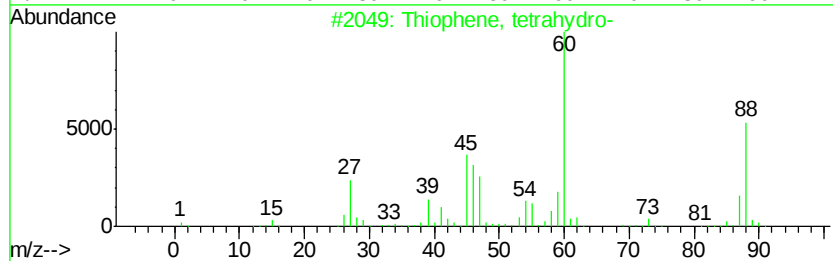
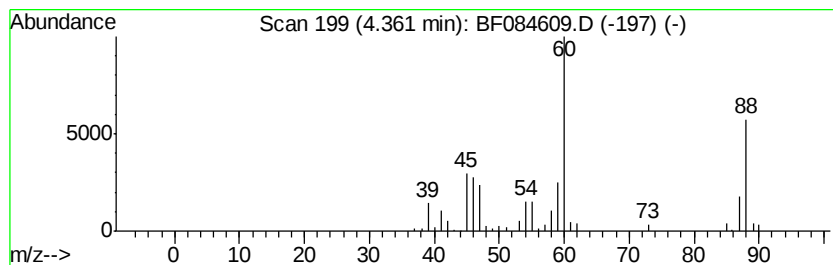
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Thiophene, tetrahydro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.36	2.23 ng	145215	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thiophene, tetrahydro-	88	C4H8S	000110-01-0	95
2		Thietane, 2-methyl-	88	C4H8S	017837-41-1	52
3		Ethylvinyl sulfide	88	C4H8S	000627-50-9	40
4		1,2-Hydrazinedicarboxaldehyde	88	C2H4N2O2	000628-36-4	38
5		Sulfide, allyl methyl	88	C4H8S	010152-76-8	22



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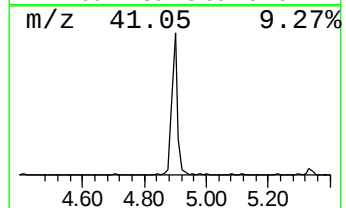
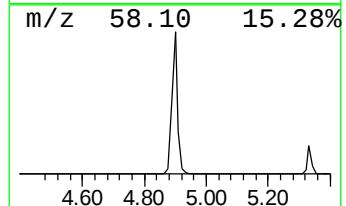
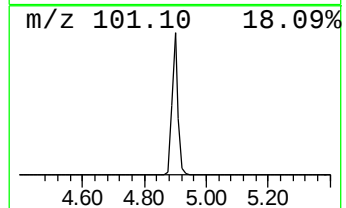
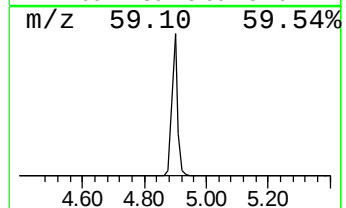
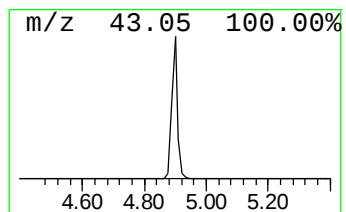
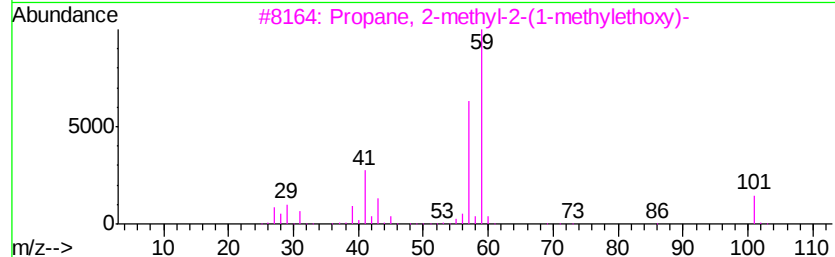
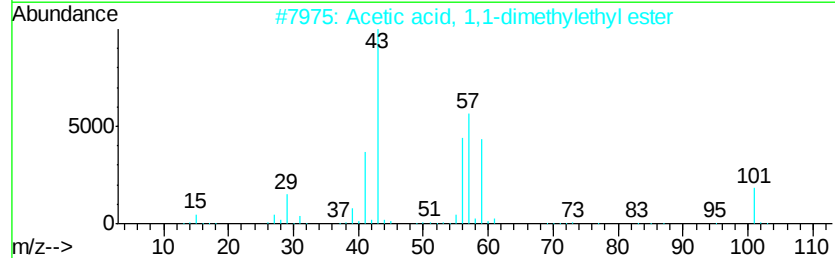
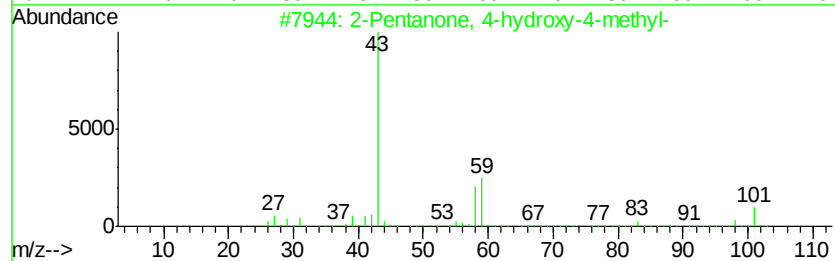
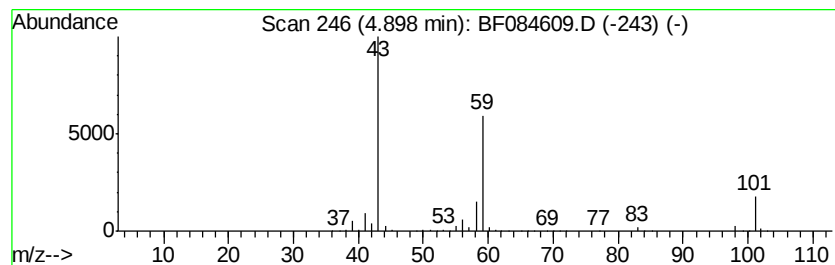
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	38.86 ng	2528870	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



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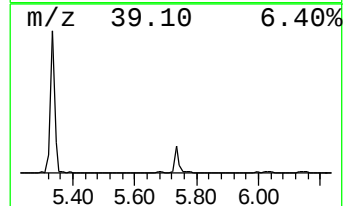
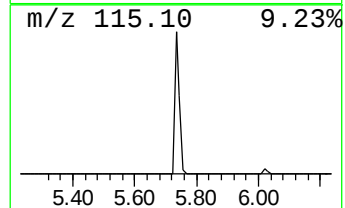
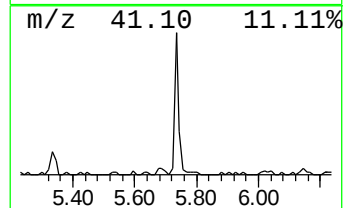
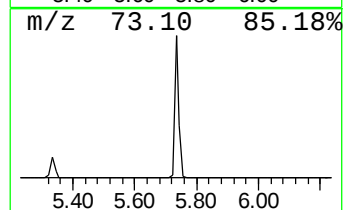
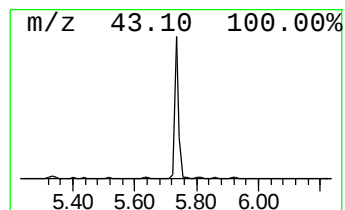
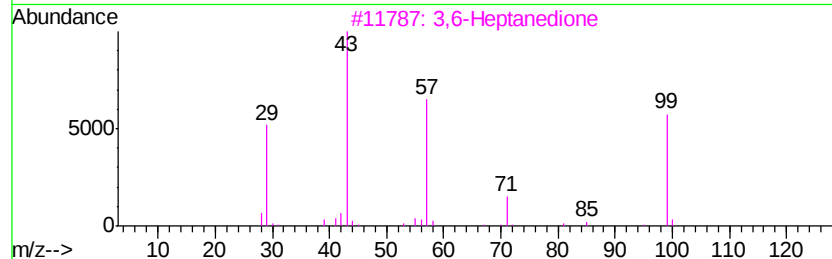
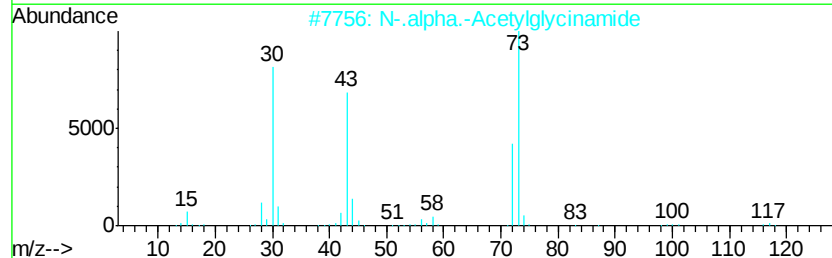
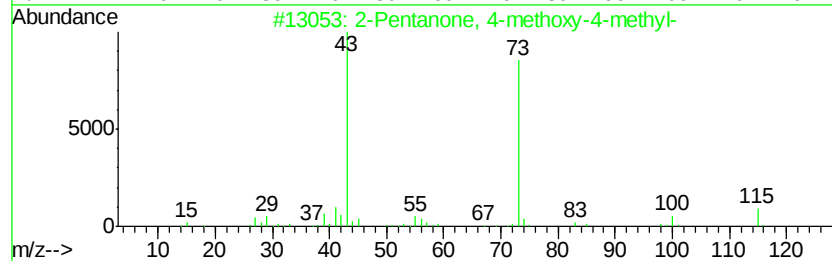
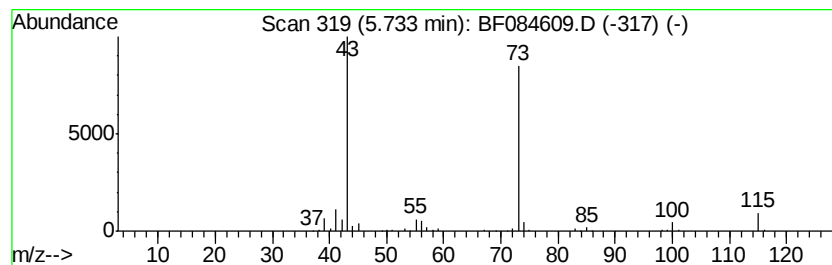
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 2-Pentanone, 4-methoxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.73	7.32 ng	476551	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		N-.alpha.-Acetylglucynamide	116	C4H8N2O2	002620-63-5	38
3		3,6-Heptanedione	128	C7H12O2	001703-51-1	23
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	12



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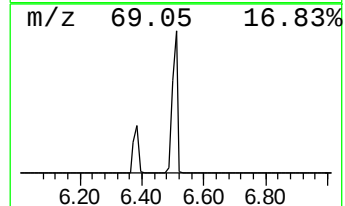
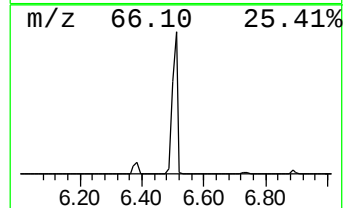
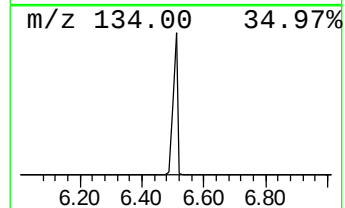
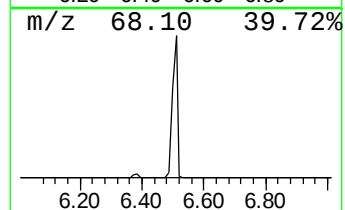
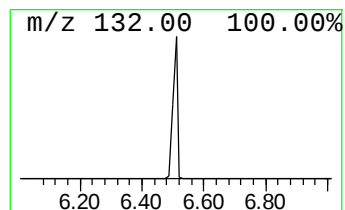
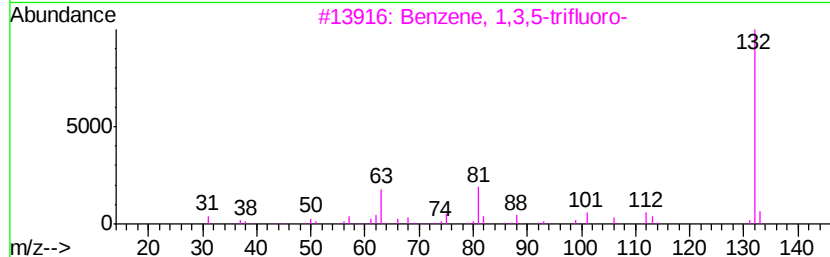
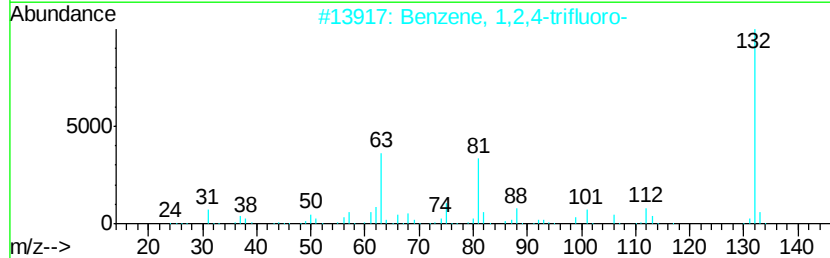
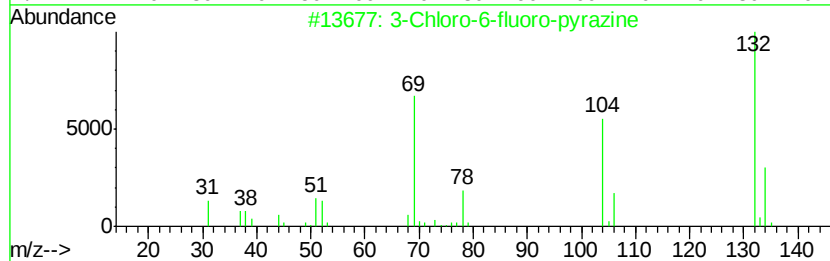
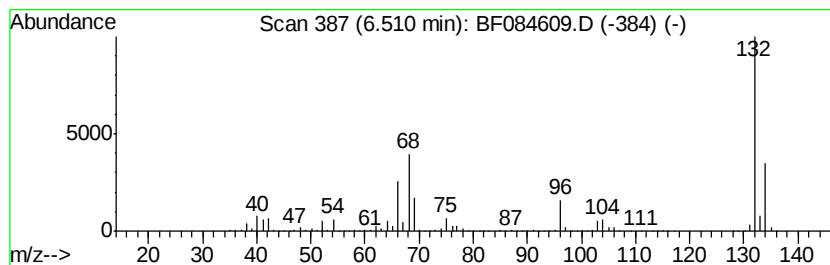
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	97.95 ng	6374780	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9	
3	Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9	
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9	
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9	



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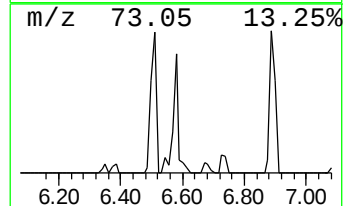
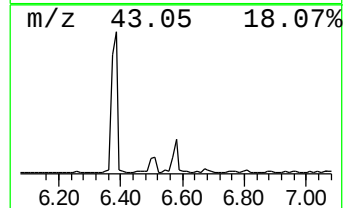
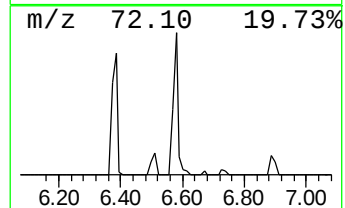
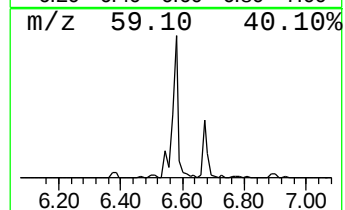
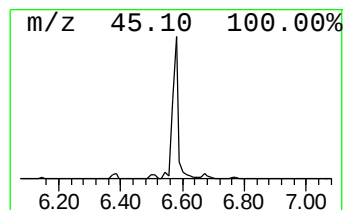
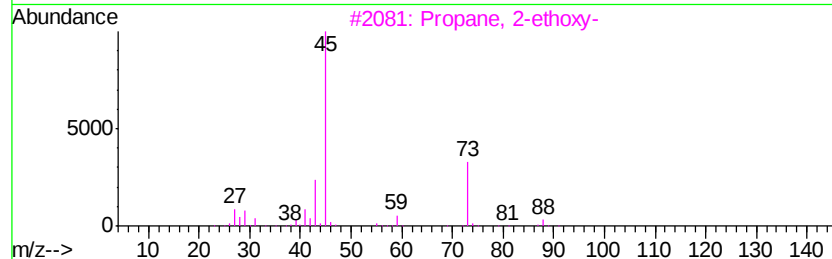
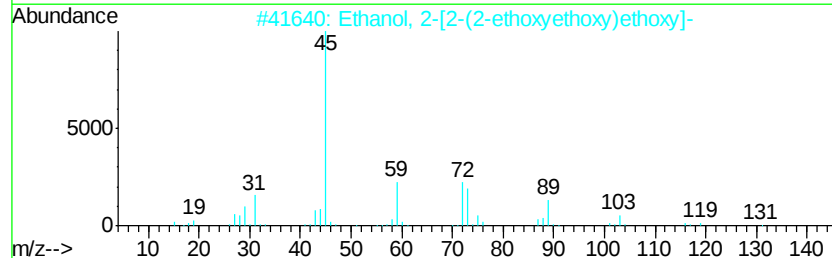
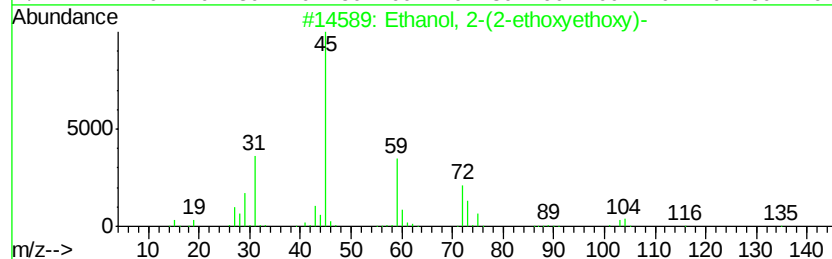
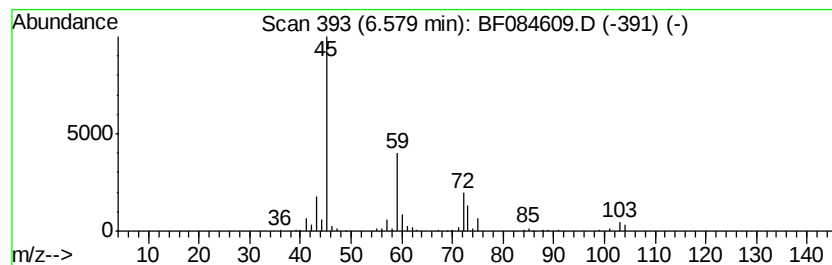
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	3.60 ng	234016	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-ethoxvethoxy)-	134	C6H14O3	000111-90-0	83
2		Ethanol, 2-[2-(2-ethoxvethoxy)et...	178	C8H18O4	000112-50-5	59
3		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	43
4		Ethanol, 2-(2-methoxvethoxy)-	120	C5H12O3	000111-77-3	42
5		2-Methoxyisopropyl cyanoacetate	157	C7H11NO3	032804-79-8	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012516\
Data File : BF084609.D
Acq On : 25 Jan 2016 16:09
Operator : SJ/IZ
Sample : H1201-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-02

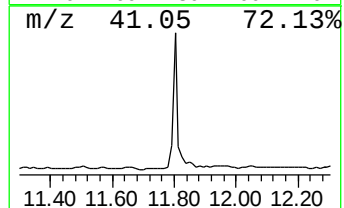
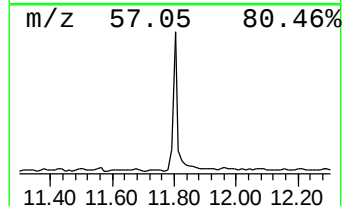
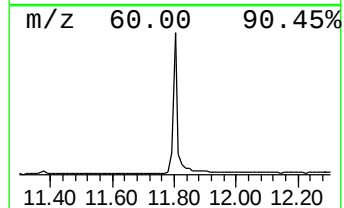
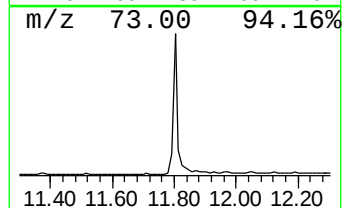
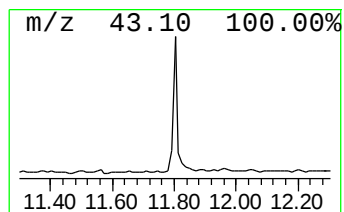
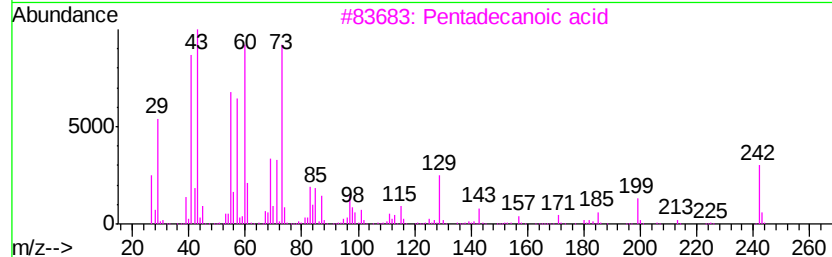
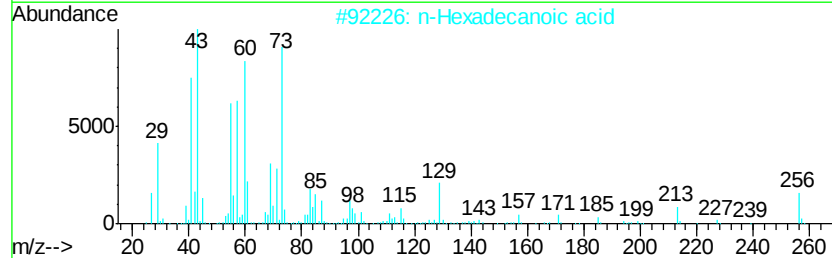
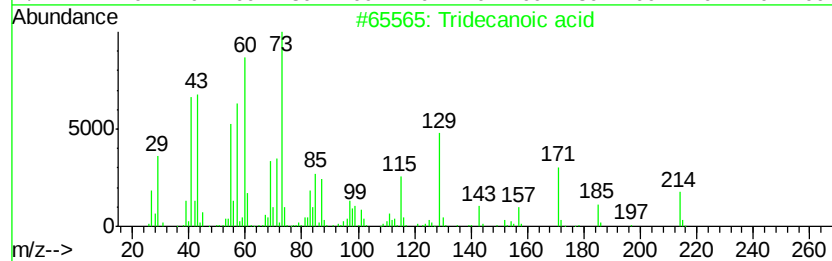
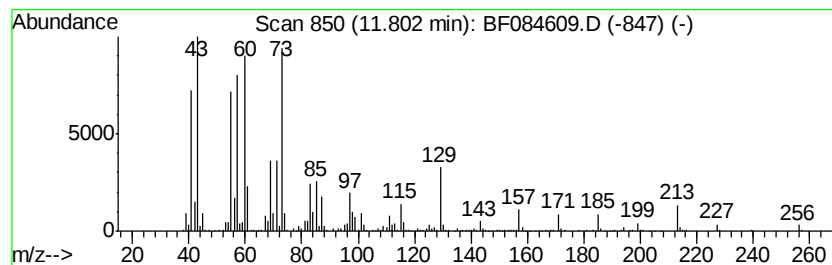
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	8.21 ng	899935	Phenanthrene-d10	11.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		Dodecanoic acid	200	C12H24O2	000143-07-7	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012516\
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Acq On : 25 Jan 2016 16:09
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Sample : H1201-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

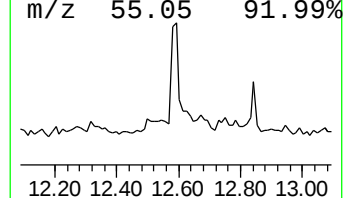
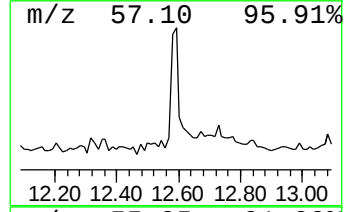
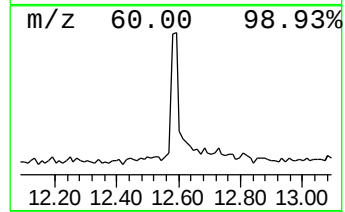
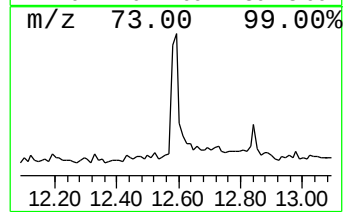
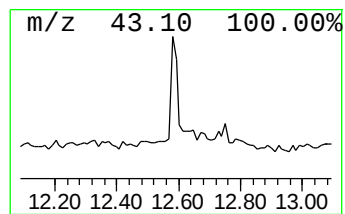
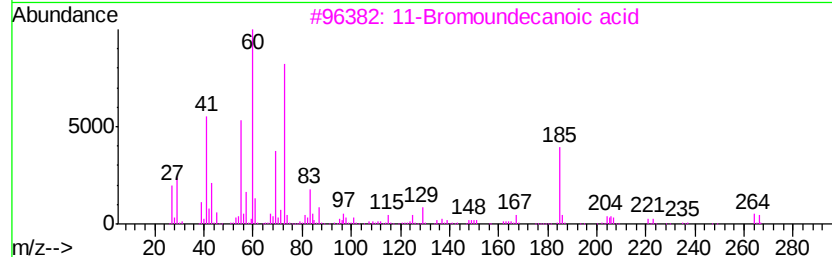
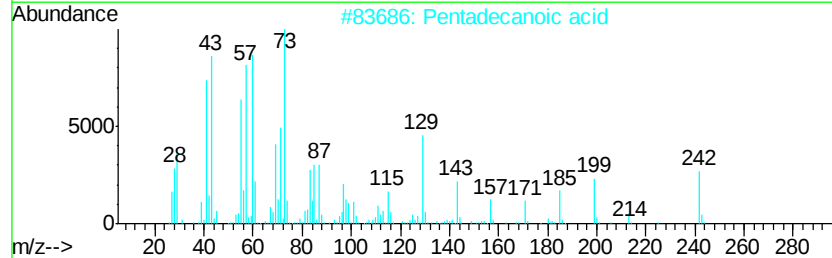
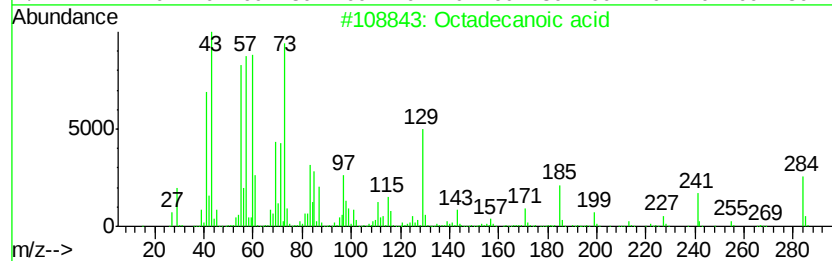
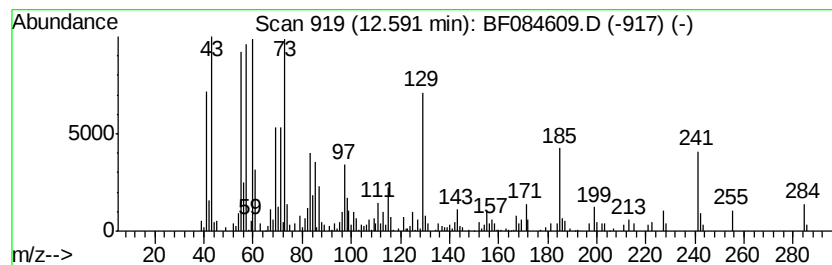
Instrument :
BNA_F
ClientSampleId :
MW-02

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.59	2.90 ng	250591	Chrysene-d12		13.88	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	90
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	45
3		11-Bromoundecanoic acid	264	C11H21BrO2	002834-05-1	38
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	38
5		Undecanoic acid	186	C11H22O2	000112-37-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012516\
Data File : BF084609.D
Acq On : 25 Jan 2016 16:09
Operator : SJ/IZ
Sample : H1201-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-02

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.20	9.1 ng		592240	1	6.74	1301630	20.0
Thiophene, tetrah...	4.36	2.2 ng		145215	1	6.74	1301630	20.0
2-Pentanone, 4-hy...	4.90	38.9 ng		2528870	1	6.74	1301630	20.0
2-Pentanone, 4-me...	5.73	7.3 ng		476551	1	6.74	1301630	20.0
unknown6.51	6.51	98.0 ng		6374780	1	6.74	1301630	20.0
Ethanol, 2-(2-eth...	6.58	3.6 ng		234016	1	6.74	1301630	20.0
Tridecanoic acid	11.80	8.2 ng		899935	4	11.25	2190970	20.0
Octadecanoic acid	12.59	2.9 ng		250591	5	13.88	1725980	20.0