

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
 Data File : BF080696.D
 Acq On : 28 Jul 2015 3:57
 Operator : TP/UM
 Sample : G3084-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUMBO-EP-7

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

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:\HPCHEM1\BNA_F\METHODS\8270-BF072715.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	4.453	205	207	214	rVB	17348	23454	2.24%	0.282%
2	5.093	261	263	266	rBV	596797	605148	57.72%	7.266%
3	5.516	297	300	302	rBV	826987	1048369	100.00%	12.587%
4	5.927	334	336	338	rVB	49980	45063	4.30%	0.541%
5	6.533	387	389	392	rBV	1005148	964764	92.03%	11.583%
6	6.682	400	402	405	rBV	1029382	1025318	97.80%	12.310%
7	6.922	421	423	426	rVB	247349	204399	19.50%	2.454%
8	7.082	434	437	439	rVB	926307	794781	75.81%	9.542%
9	7.482	470	472	474	rVB	812115	634302	60.50%	7.616%
10	8.213	534	536	538	rBV	188852	206828	19.73%	2.483%
11	9.288	627	630	632	rBV	899722	793155	75.66%	9.523%
12	9.676	662	664	666	rBV	204282	160036	15.27%	1.921%
13	9.962	687	689	691	rBV	183080	156368	14.92%	1.877%
14	10.751	756	758	760	rBV	349043	328068	31.29%	3.939%
15	11.448	817	819	821	rBV	149014	133903	12.77%	1.608%
16	11.676	837	839	843	rBV	32370	46943	4.48%	0.564%
17	11.974	864	865	868	rBV	18052	20755	1.98%	0.249%
18	12.236	886	888	890	rVB	54888	43620	4.16%	0.524%
19	12.579	916	918	921	rVV	81837	70762	6.75%	0.850%
20	12.671	923	926	928	rBV	19686	20160	1.92%	0.242%
21	12.899	943	946	948	rBV	13455	18590	1.77%	0.223%
22	13.059	957	960	962	rBV	328659	386737	36.89%	4.643%
23	14.042	1043	1046	1048	rBV2	7379	12824	1.22%	0.154%
24	14.191	1056	1059	1061	rBV	68925	77240	7.37%	0.927%
25	15.220	1146	1149	1151	rVB	67343	80128	7.64%	0.962%
26	15.368	1158	1162	1166	rBV6	6380	16641	1.59%	0.200%
27	15.802	1198	1200	1204	rVB	31790	42973	4.10%	0.516%
28	15.917	1208	1210	1215	rBV6	4980	11549	1.10%	0.139%
29	16.740	1279	1282	1286	rBV6	5830	19939	1.90%	0.239%
30	19.151	1489	1493	1498	rVB3	12390	41868	3.99%	0.503%
31	21.392	1682	1689	1702	rBV3	63834	294237	28.07%	3.533%

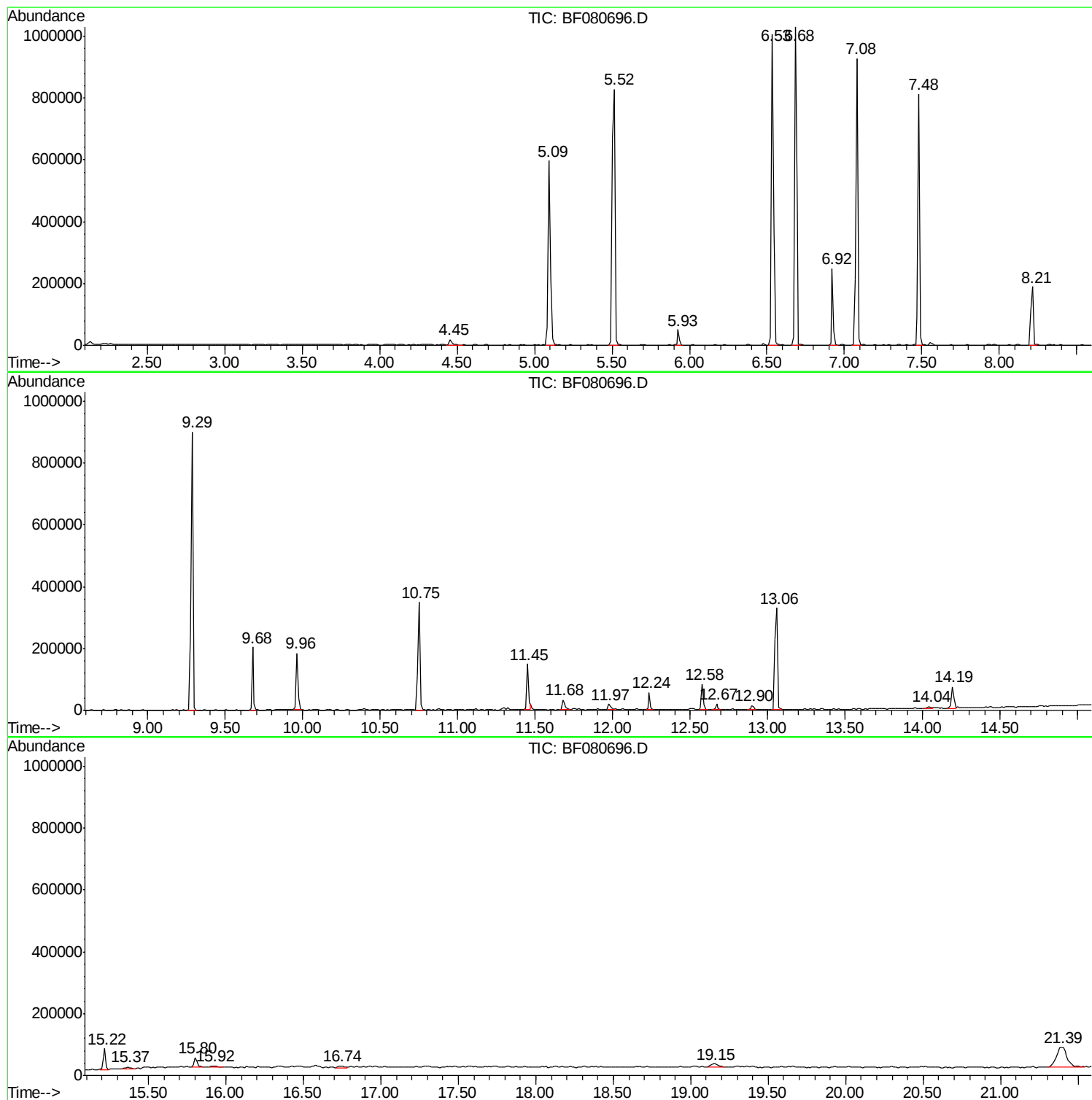
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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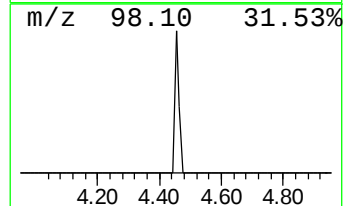
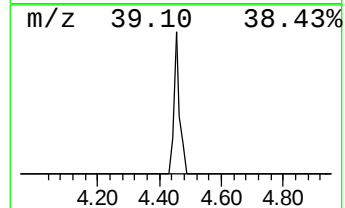
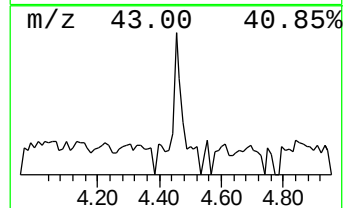
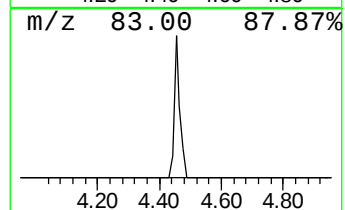
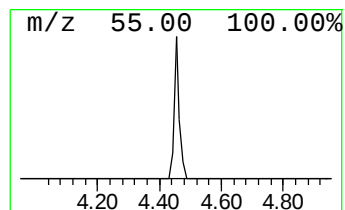
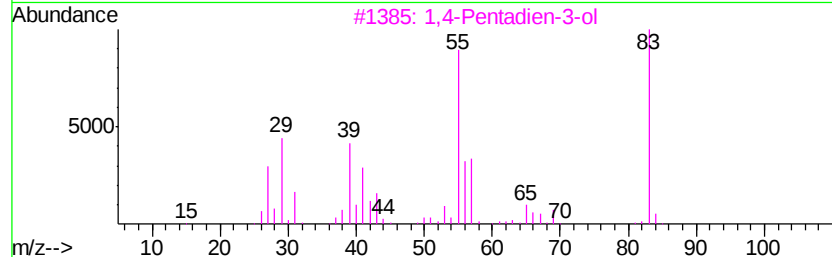
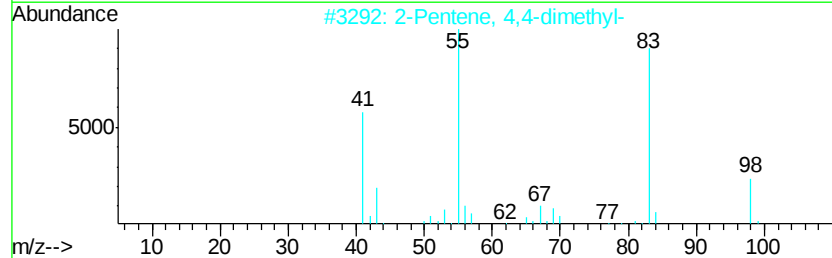
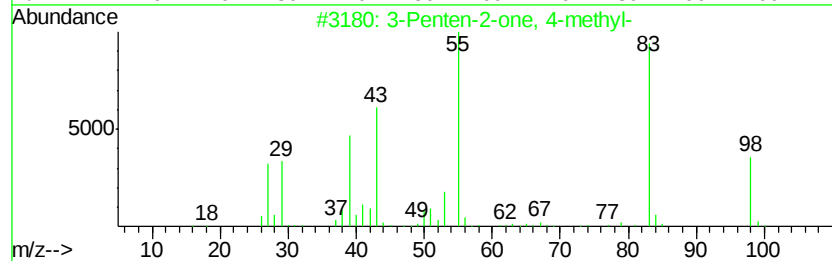
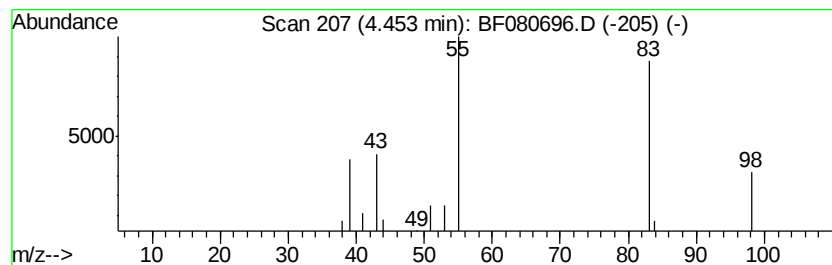
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.45	2.29 ng	23454	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	90
2			2-Pentene, 4,4-dimethyl-	98	C7H14	026232-98-4	83
3			1,4-Pentadien-3-ol	84	C5H8O	000922-65-6	64
4			3,4-Dimethyl-2-pentene(c,t)	98	C7H14	1000118-16-5	64
5			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	59



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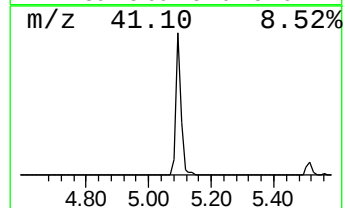
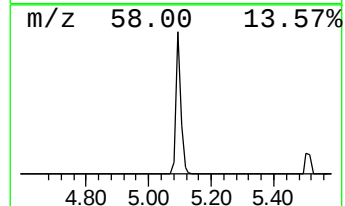
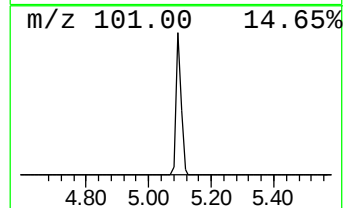
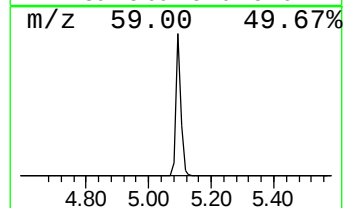
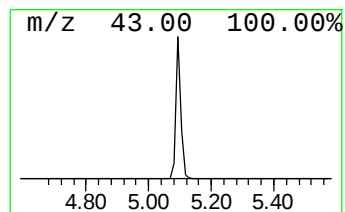
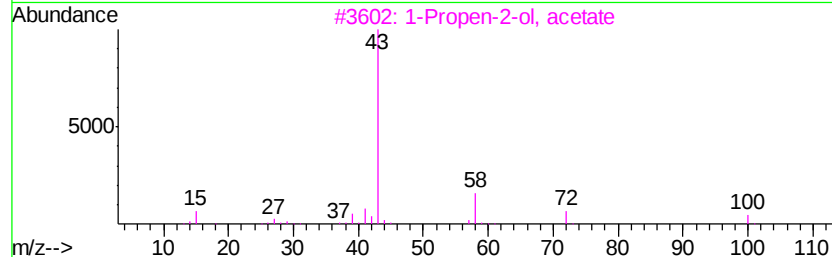
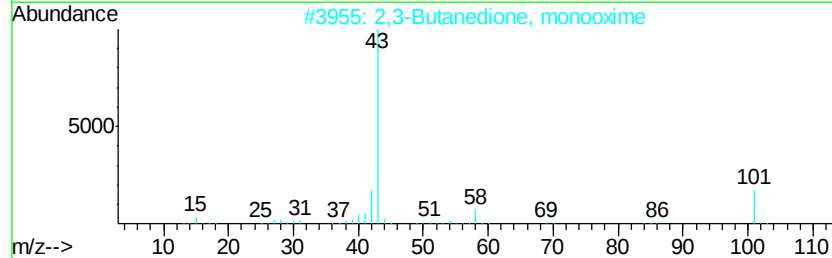
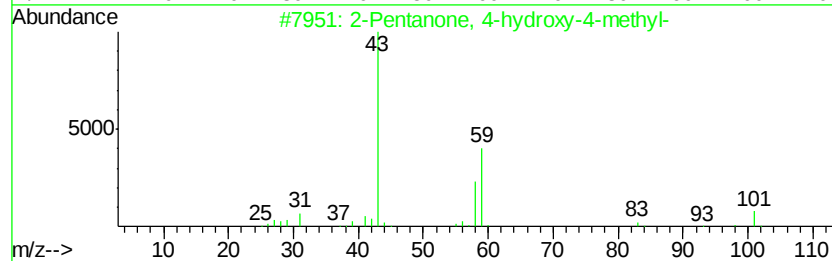
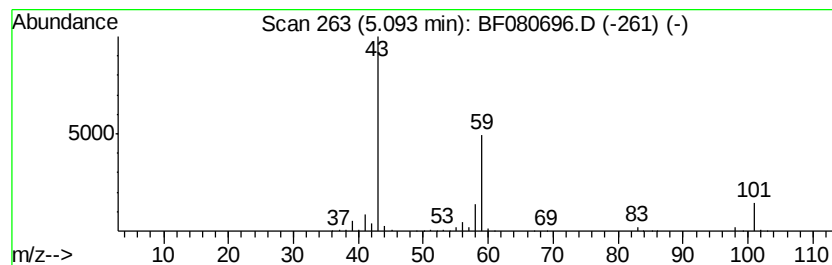
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	59.21 ng	605148	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9



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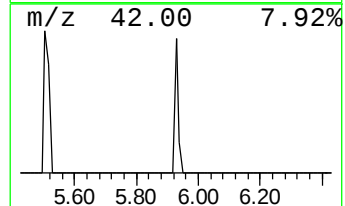
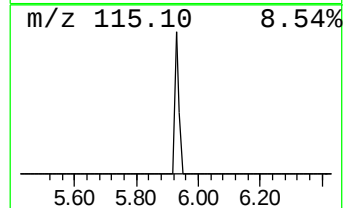
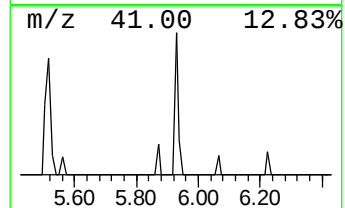
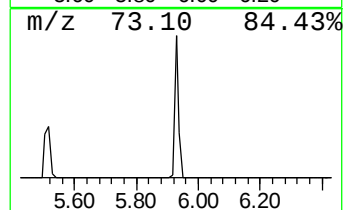
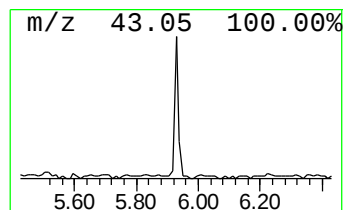
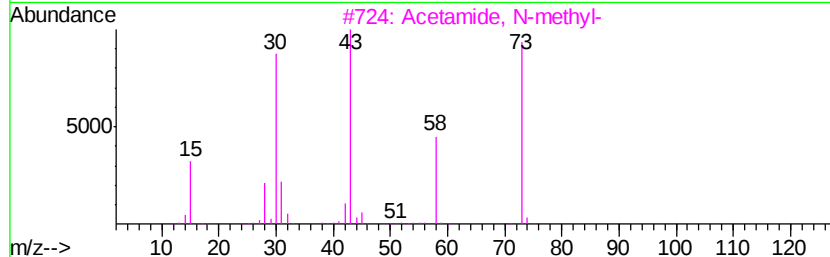
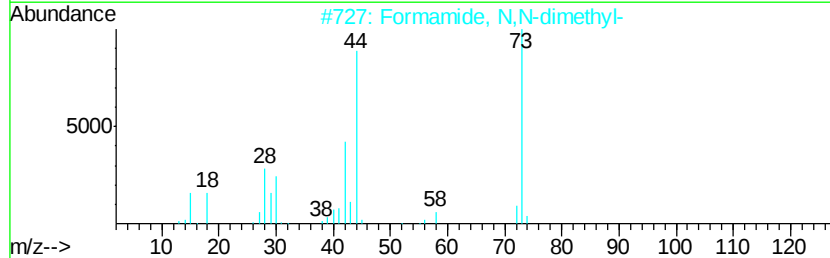
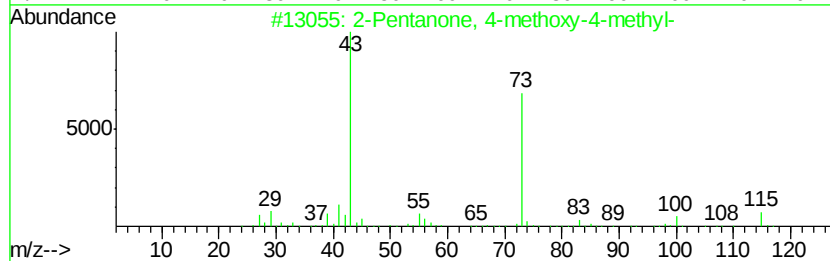
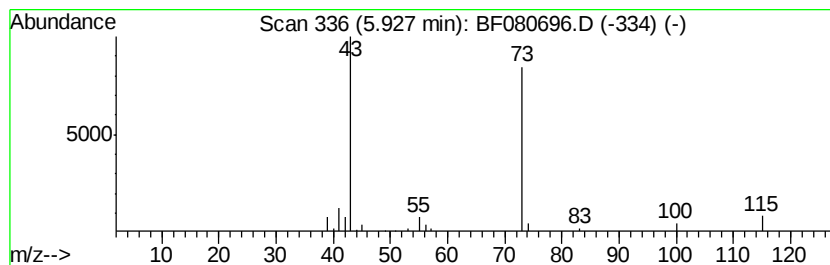
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	4.41 ng	45063	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	9
3		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9
4		Guanidine, methyl-	73	C2H7N3	000471-29-4	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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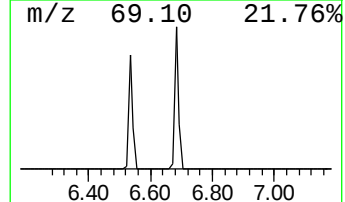
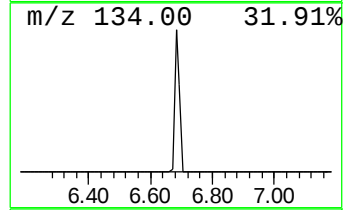
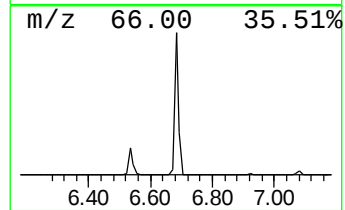
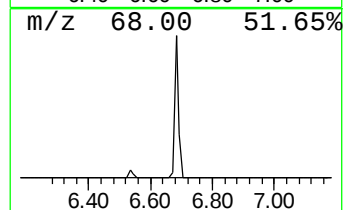
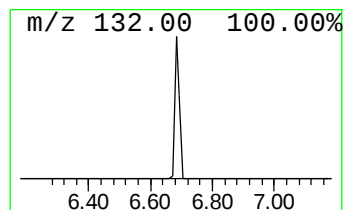
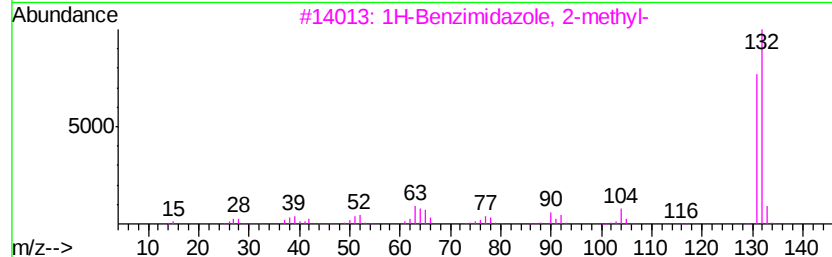
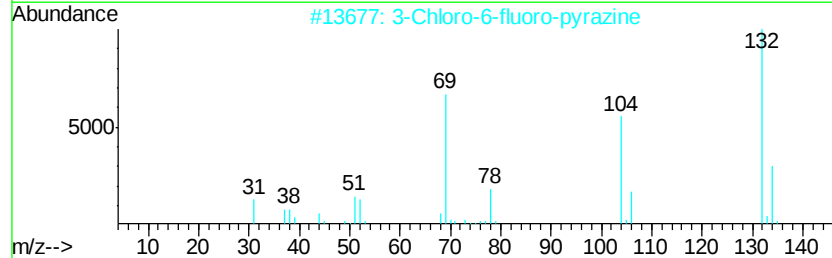
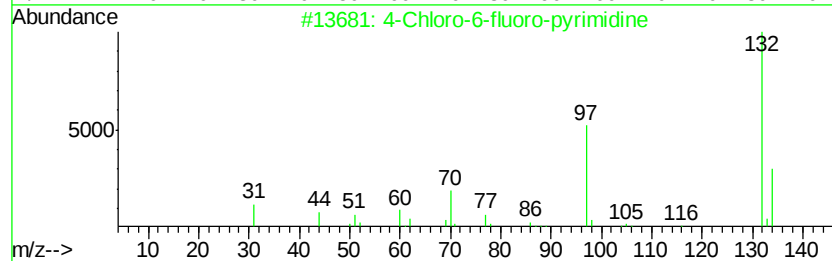
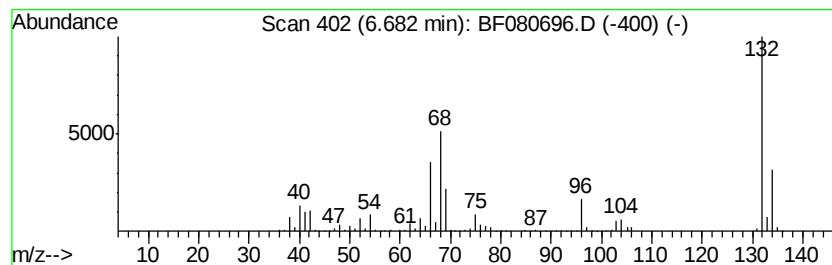
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 unknown6.68 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	100.33 ng	1025320	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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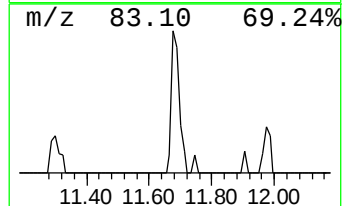
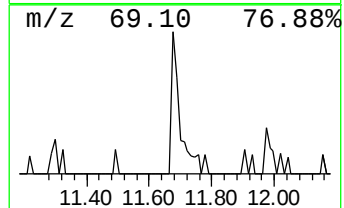
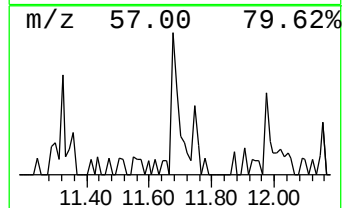
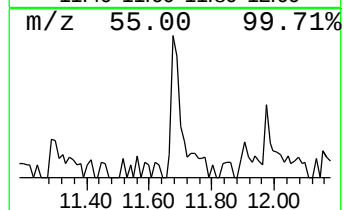
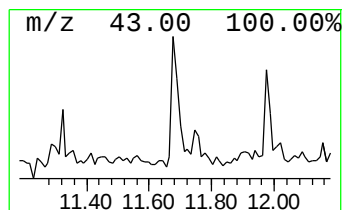
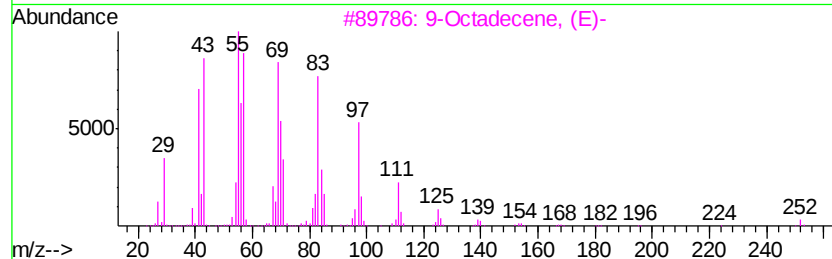
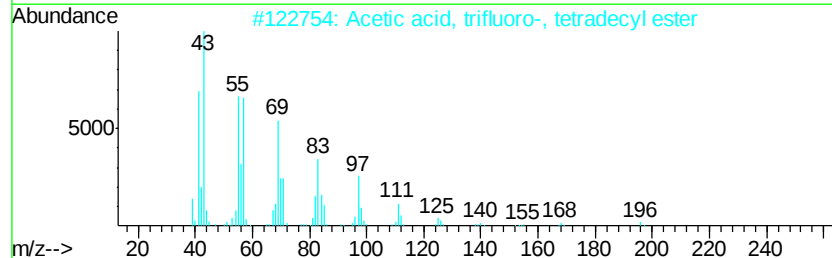
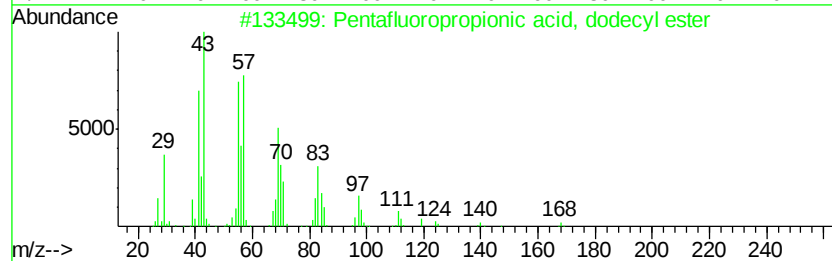
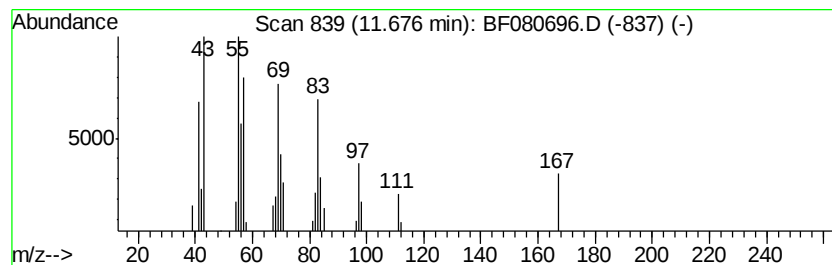
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Pentafluoropropionic acid, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.68	7.01 ng	46943	Phenanthrene-d10	11.45

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, dodec...	332	C15H25F5O2	006222-04-4	90
2			Acetic acid, trifluoro-, tetradec...	310	C16H29F3O2	006222-02-2	87
3			9-Octadecene, (E)-	252	C18H36	007206-25-9	87
4			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	87
5			5-Octadecene, (E)-	252	C18H36	007206-21-5	87



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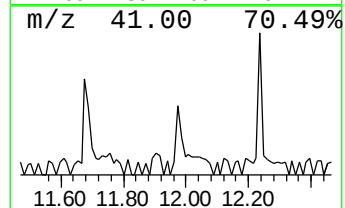
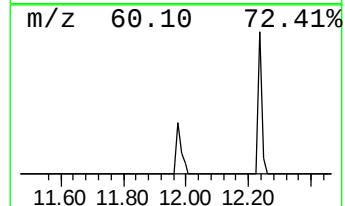
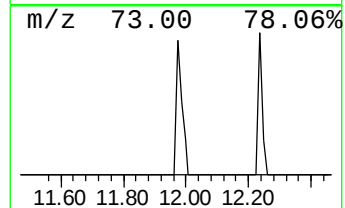
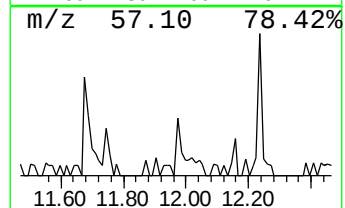
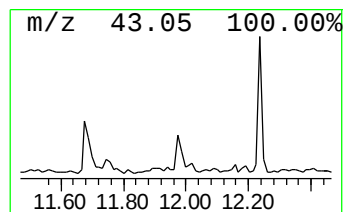
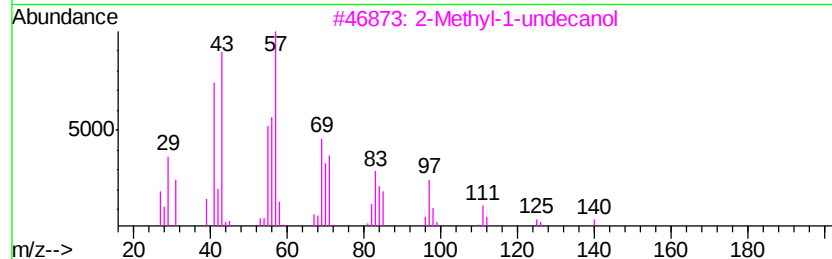
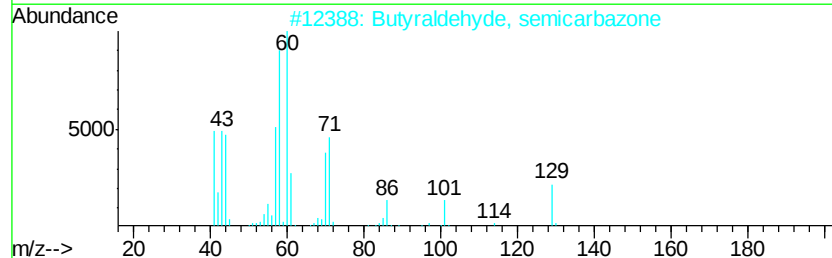
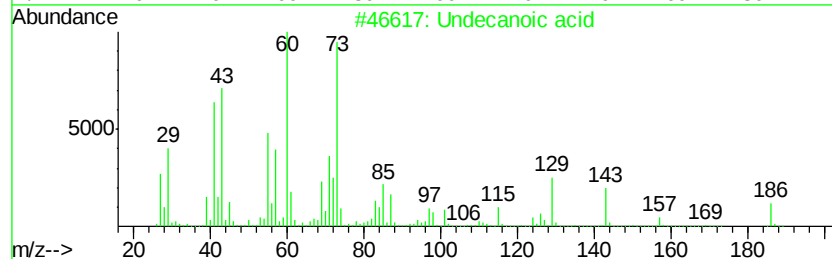
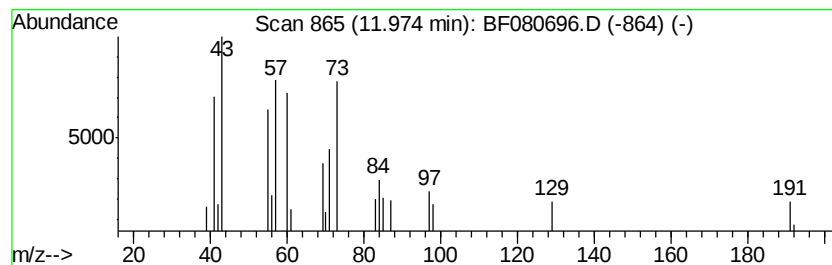
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown11.97 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	3.10 ng	20755	Phenanthrene-d10	11.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecanoic acid	186	C11H22O2	000112-37-8	47
2		Butyraldehyde, semicarbazone	129	C5H11N3O	013183-21-6	35
3		2-Methyl-1-undecanol	186	C12H26O	010522-26-6	35
4		Amvl Nitrite	117	C5H11NO2	000110-46-3	27
5		Dodecane, 1-chloro-	204	C12H25Cl	000112-52-7	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

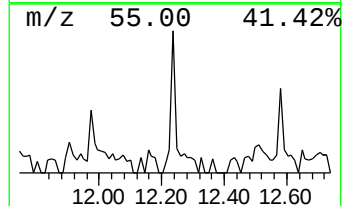
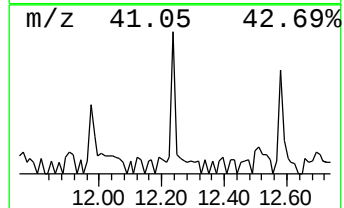
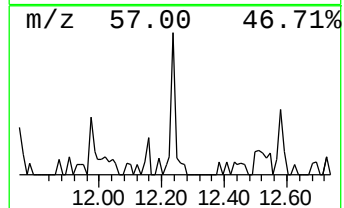
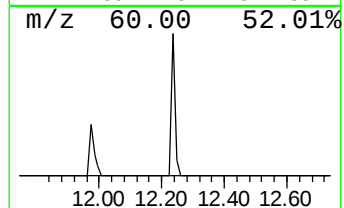
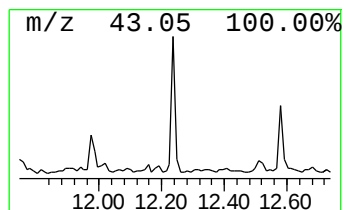
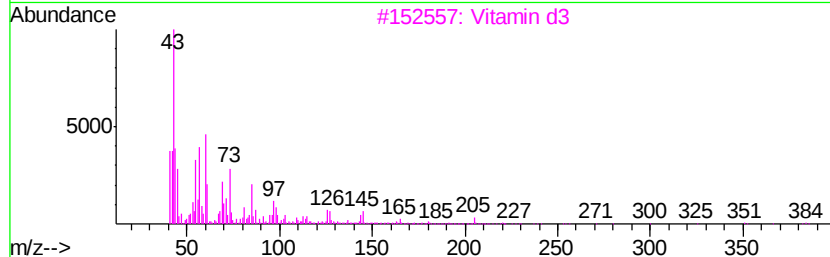
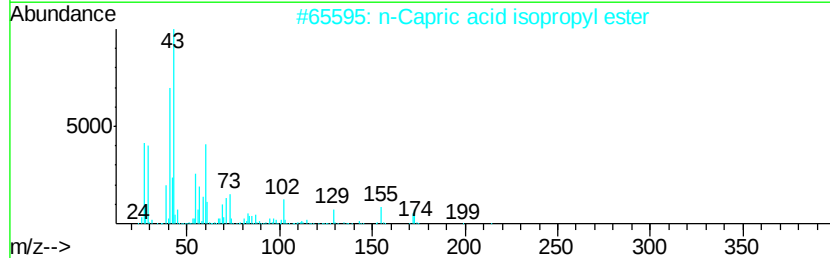
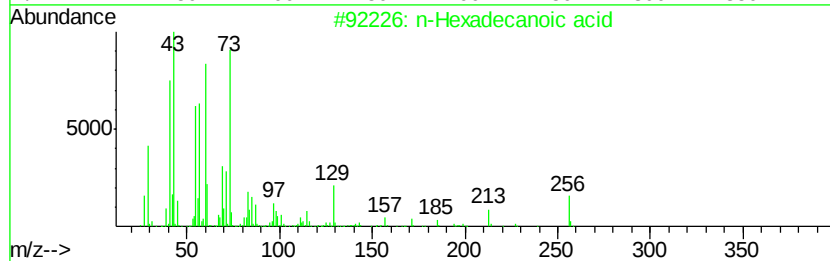
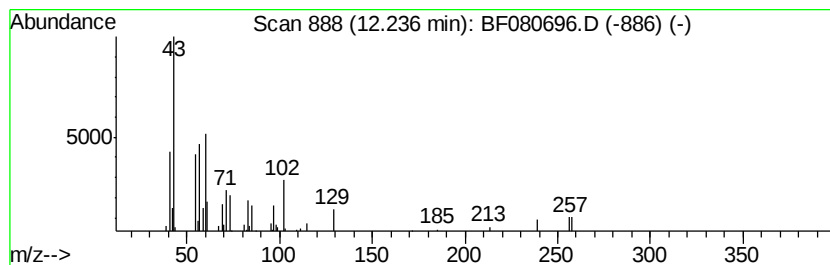
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.24	6.52 ng	43620	Phenanthrene-d10	11.45	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	78
2	n-Capric acid isopropyl ester	214	C13H26O2	002311-59-3	40
3	Vitamin d3	384	C27H44O	001406-16-2	38
4	Isopropyl Myristate	270	C17H34O2	000110-27-0	38
5	Nonanoic acid, 1-methylethyl ester	200	C12H24O2	028267-32-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUMBO-EP-7

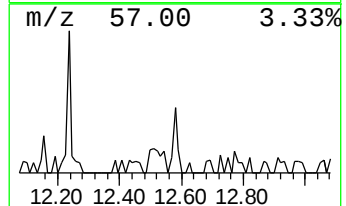
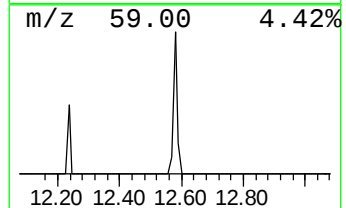
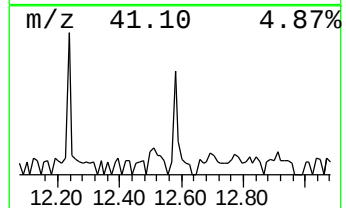
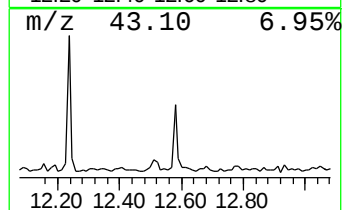
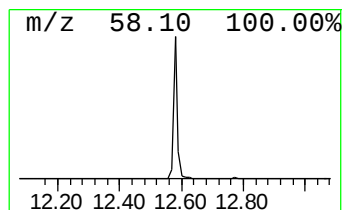
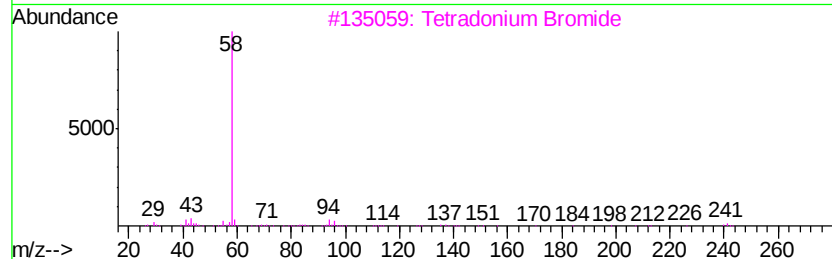
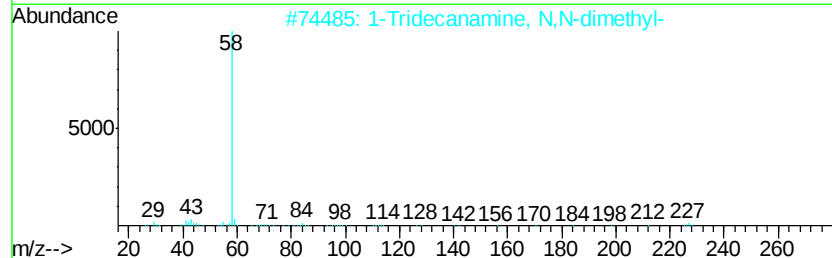
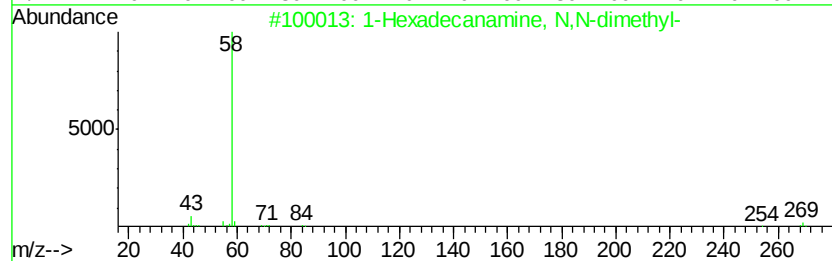
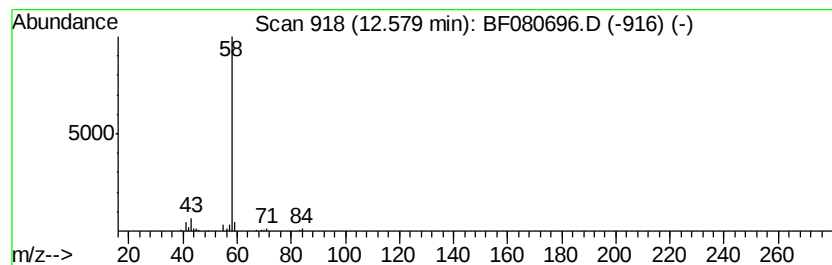
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Hexadecanamine, N,N-dimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	10.57 ng	70762	Phenanthrene-d10	11.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexadecanamine, N,N-dimethyl-	269	C18H39N	000112-69-6	83
2		1-Tridecanamine, N,N-dimethyl-	227	C15H33N	017373-29-4	83
3		Tetradonium Bromide	335	C17H38BrN	001119-97-7	78
4		1-Pentadecanamine, N,N-dimethyl-	255	C17H37N	017678-60-3	74
5		1-Undecanamine, N,N-dimethyl-	199	C13H29N	017373-28-3	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DUMBO-EP-7

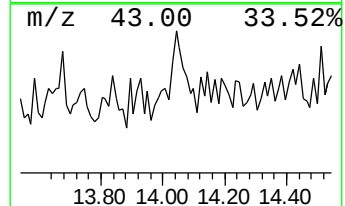
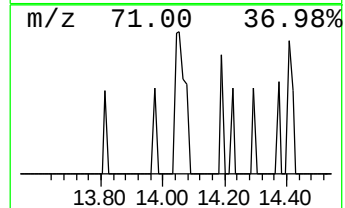
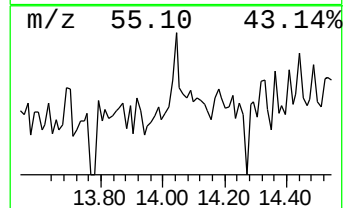
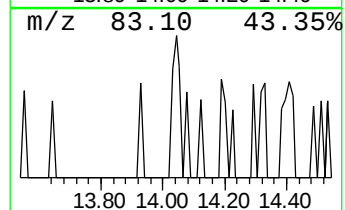
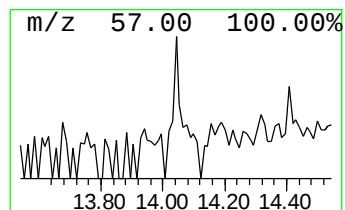
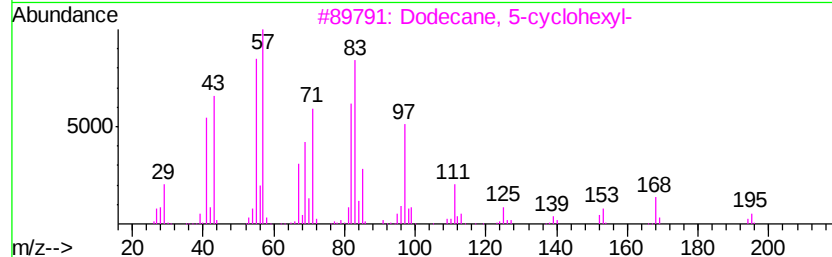
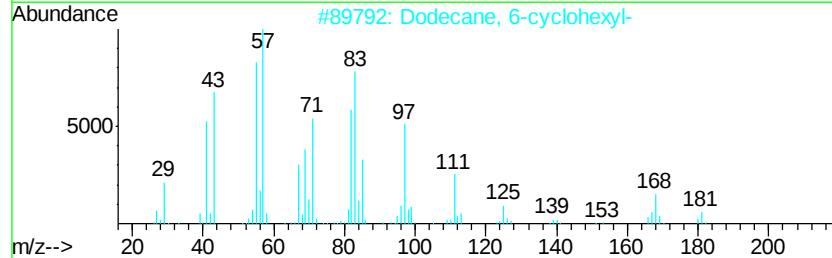
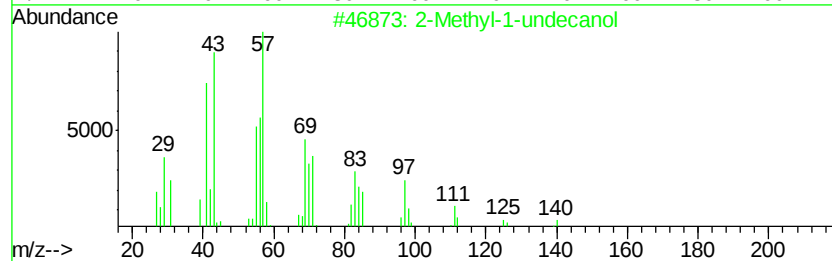
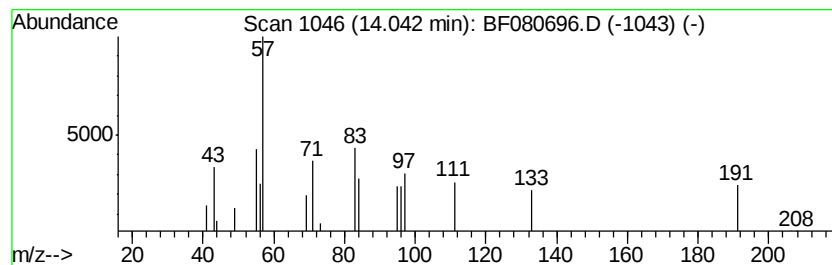
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown14.04 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	3.32 ng	12824	Chrysene-d12	14.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-1-undecanol	186	C12H26O	010522-26-6	32
2			Dodecane, 6-cyclohexyl-	252	C18H36	013151-86-5	17
3			Dodecane, 5-cyclohexyl-	252	C18H36	013151-85-4	17
4			2-Decen-1-ol, (E)-	156	C10H20O	018409-18-2	9
5			1-Octadecene	252	C18H36	000112-88-9	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
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Acq On : 28 Jul 2015 3:57
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Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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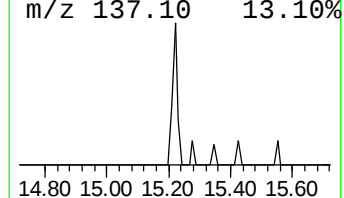
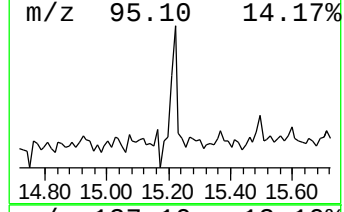
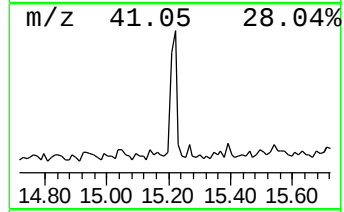
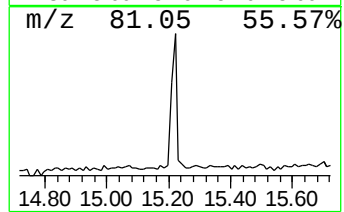
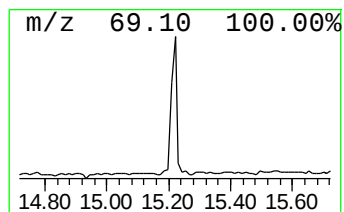
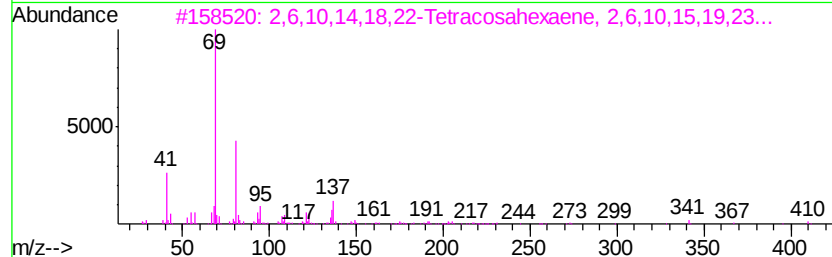
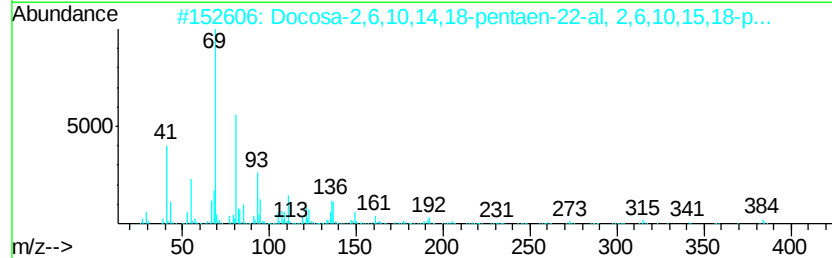
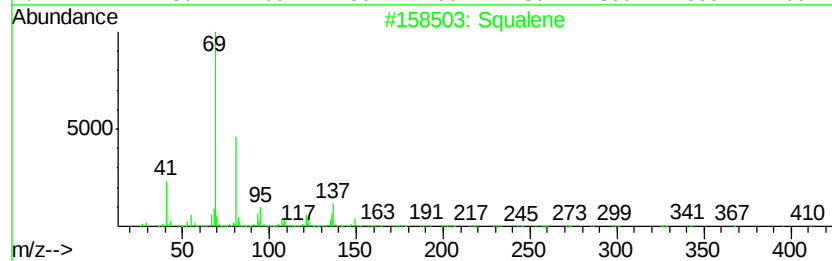
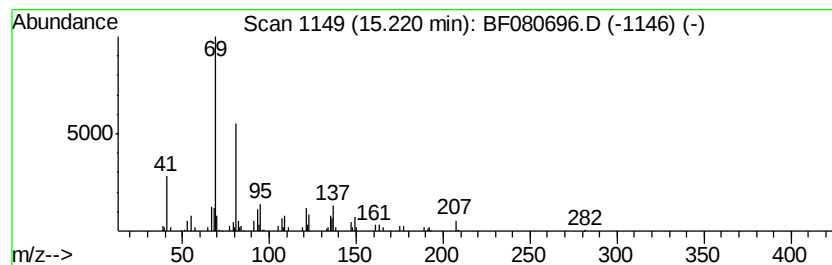
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	37.29 ng	80128	Perylene-d12	15.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	90
2		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	86
3		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	86
4		1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	81
5		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DUMBO-EP-7

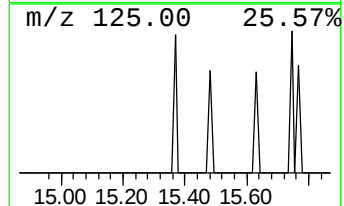
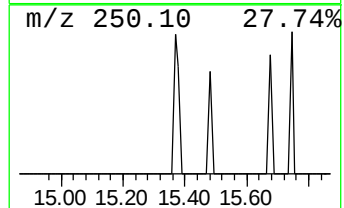
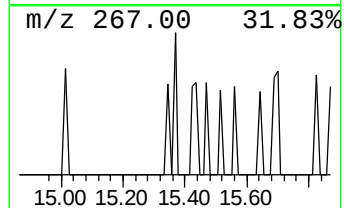
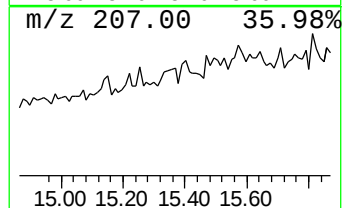
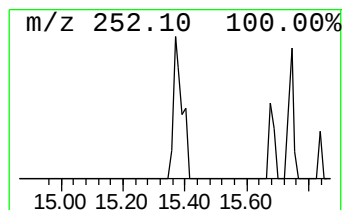
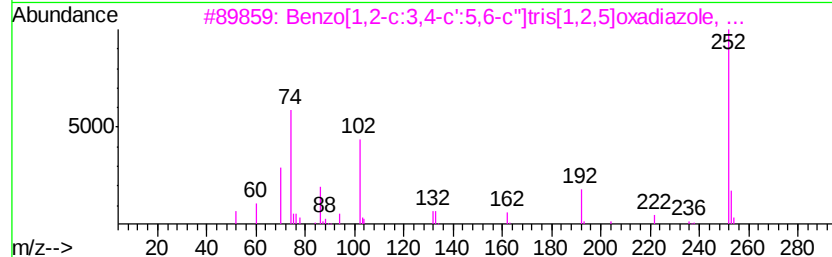
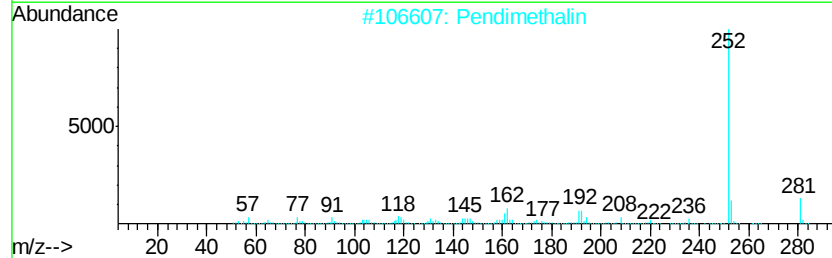
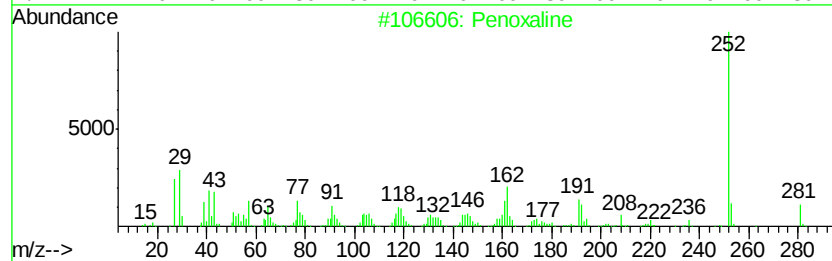
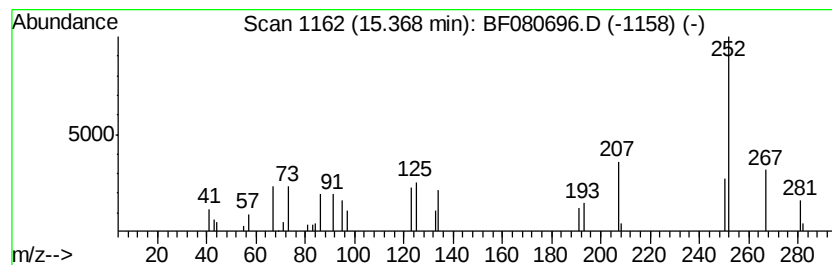
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.37 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	7.74 ng	16641	Perylene-d12	15.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Penoxaline	281	C13H19N3O4	040487-42-1	12
2		Pendimethalin	281	C13H19N3O4	040318-45-4	12
3		Benzo[1,2-c:3,4-c':5,6-c'']tris[...]	252	C6N6O6	003470-17-5	9
4		Benzo[el]pyrene	252	C20H12	000192-97-2	9
5		Benzonitrile, 3-(3-hydroxy-4-met...	252	C15H12N2O2	304683-23-6	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072715\
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Acq On : 28 Jul 2015 3:57
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Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
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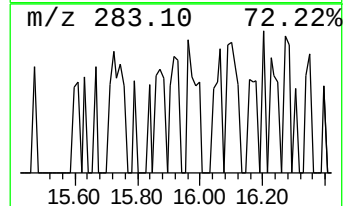
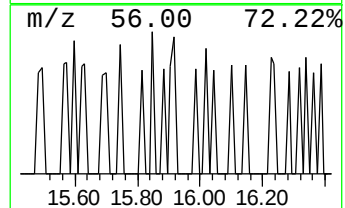
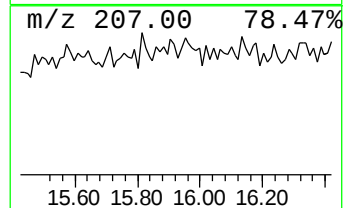
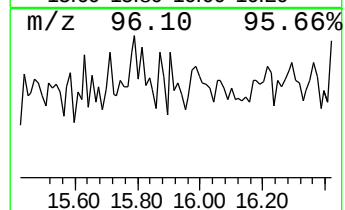
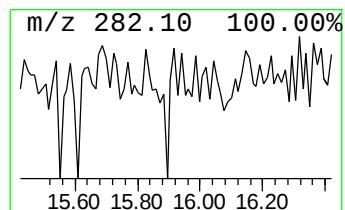
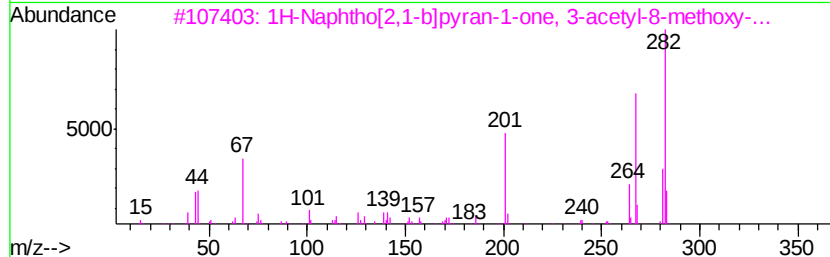
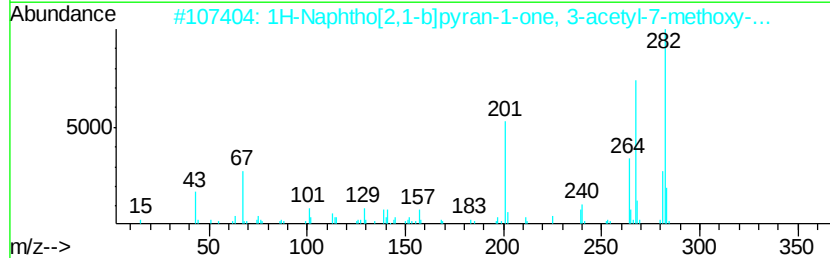
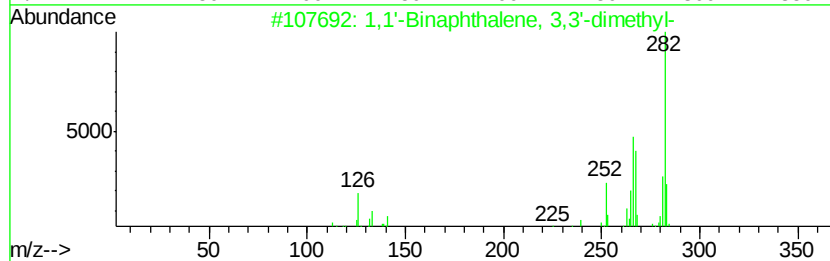
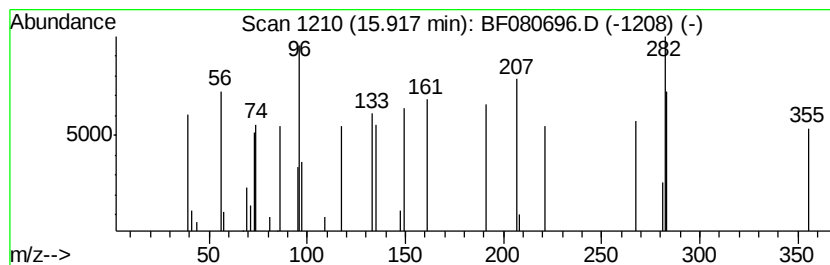
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.92 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	5.38 ng	11549	Perylene-d12	15.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Binaphthalene, 3,3'-dimethyl-	282	C22H18	034042-82-5	12
2		1H-Naphtho[2,1-b]pyran-1-one, 3-...	282	C17H14O4	055044-51-4	10
3		1H-Naphtho[2,1-b]pyran-1-one, 3-...	282	C17H14O4	055044-50-3	10
4		Isoindolo[2,1-f]phenanthridin-10...	283	C20H13NO	063558-92-9	9
5		2,3,8-Trimethyl-1H-pyrrolo[3,2-H...	282	C17H18N2O2	1000296-79-8	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

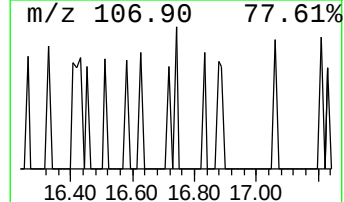
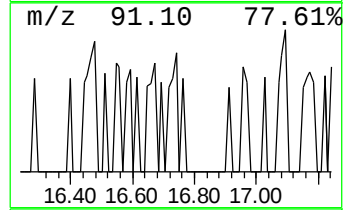
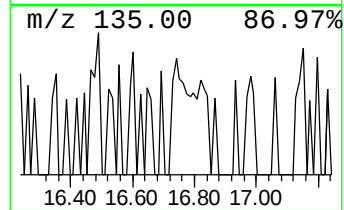
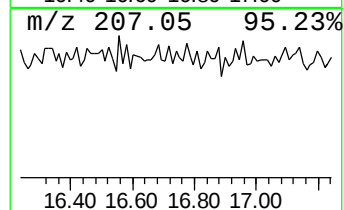
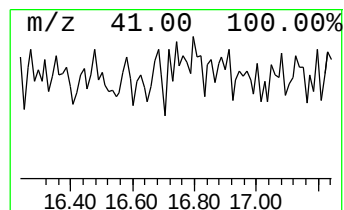
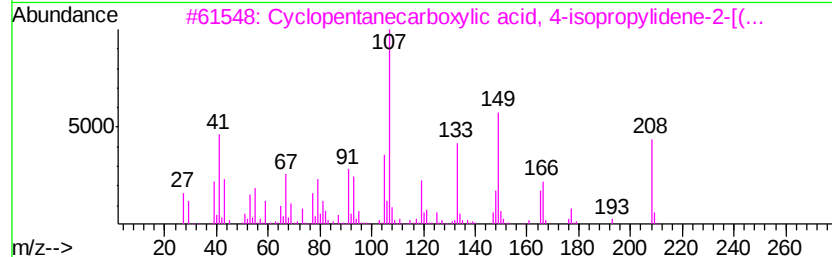
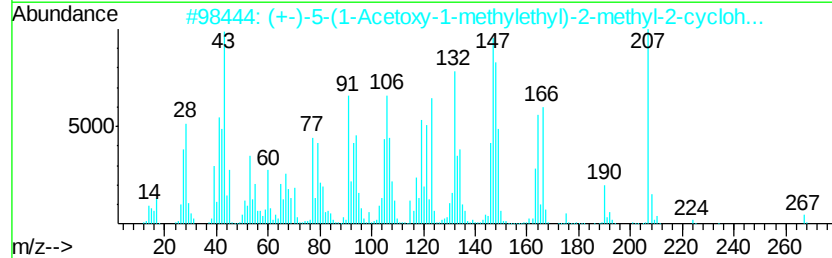
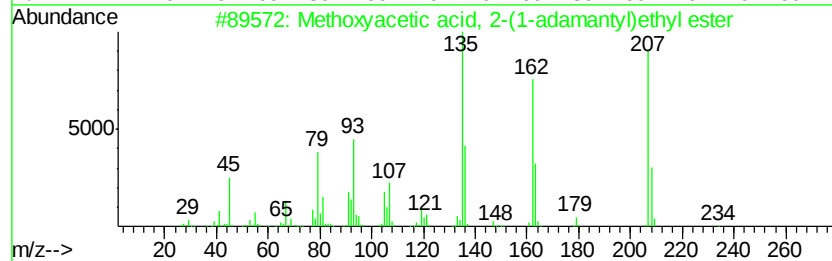
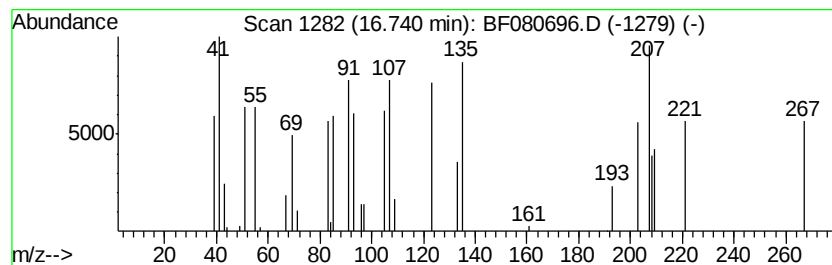
Instrument :
BNA_F
ClientSampleId :
DUMBO-EP-7

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.74 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.74	9.28 ng	19939	Perylene-d12	15.80		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methoxvacetic acid, 2-(1-adamant...	252	C15H24O3	1000282-98-7	22
2		(+)-5-(1-Acetoxy-1-methylethyl)...	267	C13H21N3O3	108904-53-6	12
3		Cyclopentanecarboxylic acid, 4-i...	208	C13H20O2	1000159-44-5	10
4		2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-17-5	10
5		Cyclobutaneacetamide, N-(4-methy...	203	C13H17NO	302567-77-7	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUMBO-EP-7

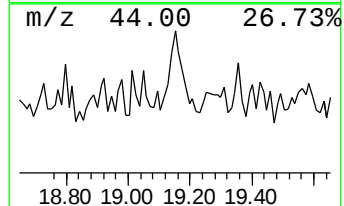
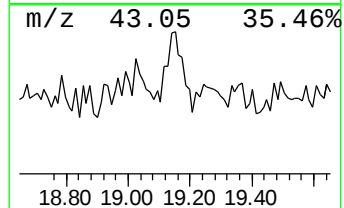
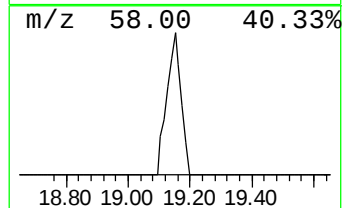
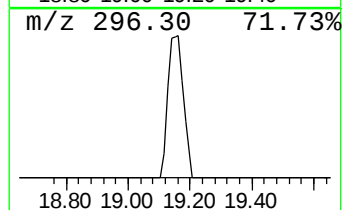
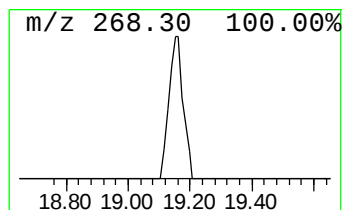
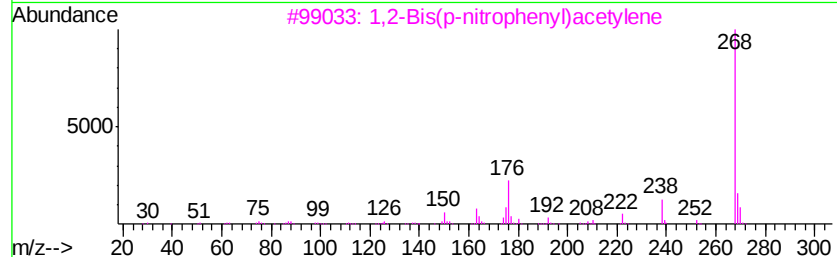
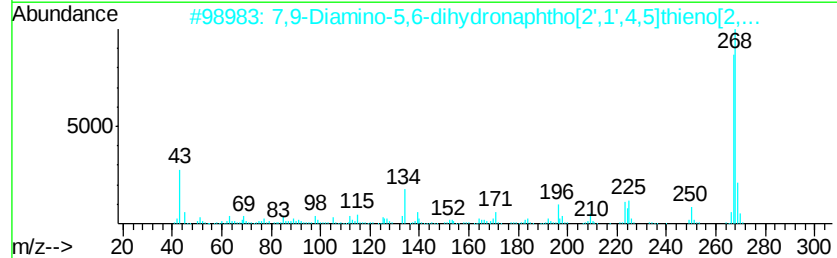
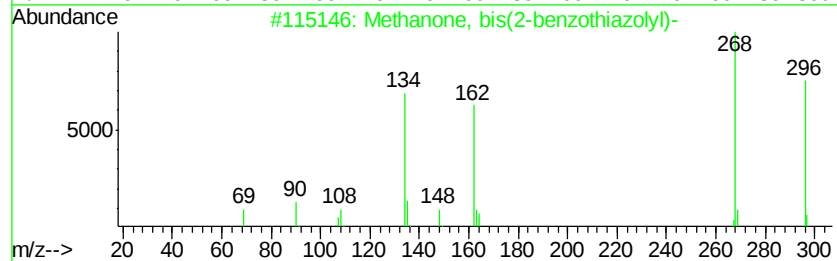
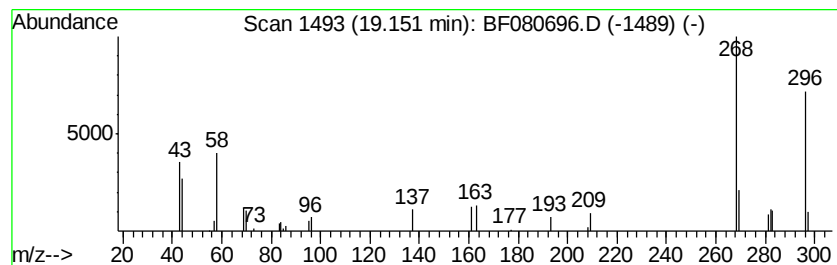
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Methanone, bis(2-benzothiaz... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.15	19.49 ng	41868	Perylene-d12	15.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methanone, bis(2-benzothiazolyl)-	296	C15H8N2OS2	004464-60-2	59
2		7,9-Diamino-5,6-dihydronaphtho[2...	268	C14H12N4S	037071-29-7	25
3		1,2-Bis(p-nitrophenyl)acetylene	268	C14H8N2O4	002735-14-0	17
4		4-Imidazolidinone, 5,5-diphenyl-...	268	C15H12N2OS	021083-47-6	16
5		Indol-2(3H)-one, 3-(2,4-dimethox...	296	C17H16N2O3	294877-12-6	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072715\
Data File : BF080696.D
Acq On : 28 Jul 2015 3:57
Operator : TP/UM
Sample : G3084-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUMBO-EP-7

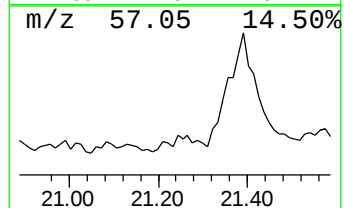
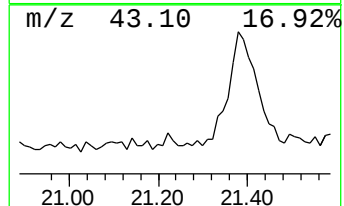
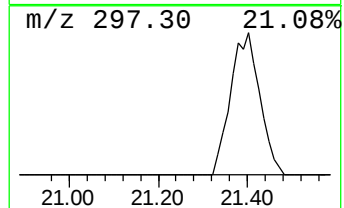
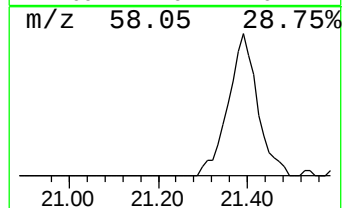
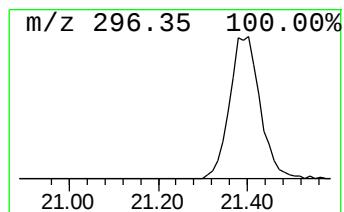
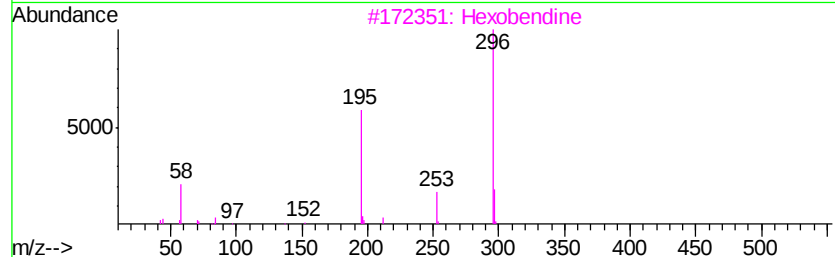
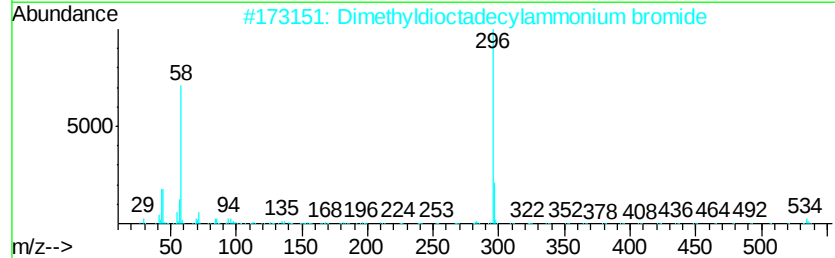
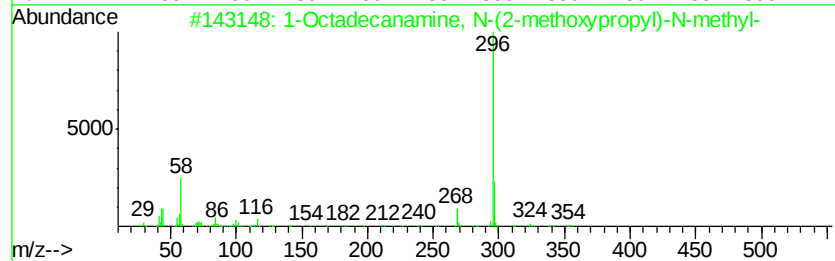
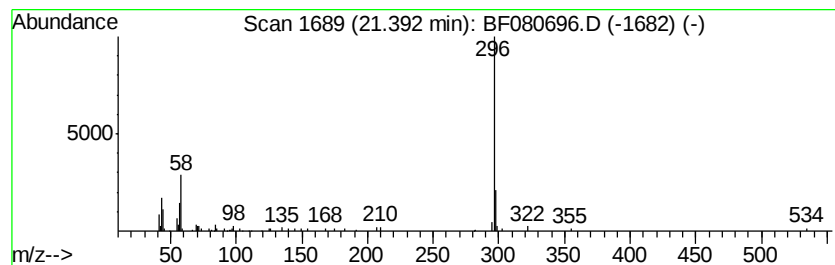
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Octadecanamine, N-(2-meth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.39	136.94 ng	294237	Perylene-d12	15.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecanamine, N-(2-methoxyvpr...	355	C23H49NO	1000196-18-5	78
2		Dimethyldioctadecylammonium bromide	630	C38H80BrN	003700-67-2	64
3		Hexobendine	592	C30H44N2O10	000054-03-5	64
4		1H-Imidazole, 2,4,5-triphenyl-	296	C21H16N2	000484-47-9	59
5		Benzeneethanamine, .beta.-hydrox...	403	C27H49NO	1000197-20-9	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072715\
 Data File : BF080696.D
 Acq On : 28 Jul 2015 3:57
 Operator : TP/UM
 Sample : G3084-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUMBO-EP-7

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072715.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...	4.45	2.3	ng	23454	1	6.92	204399	20.0
2-Pentanone, 4-hy...	5.09	59.2	ng	605148	1	6.92	204399	20.0
2-Pentanone, 4-me...	5.93	4.4	ng	45063	1	6.92	204399	20.0
unknown6.68	6.68	100.3	ng	1025320	1	6.92	204399	20.0
Pentafluoropropio...	11.68	7.0	ng	46943	4	11.45	133903	20.0
unknown11.97	11.97	3.1	ng	20755	4	11.45	133903	20.0
n-Hexadecanoic acid	12.24	6.5	ng	43620	4	11.45	133903	20.0
1-Hexadecanamine,...	12.58	10.6	ng	70762	4	11.45	133903	20.0
unknown14.04	14.04	3.3	ng	12824	5	14.19	77240	20.0
Squalene	15.22	37.3	ng	80128	6	15.80	42973	20.0
unknown15.37	15.37	7.7	ng	16641	6	15.80	42973	20.0
unknown15.92	15.92	5.4	ng	11549	6	15.80	42973	20.0
unknown16.74	16.74	9.3	ng	19939	6	15.80	42973	20.0
Methanone, bis(2-...	19.15	19.5	ng	41868	6	15.80	42973	20.0
1-Octadecanamine,...	21.39	136.9	ng	294237	6	15.80	42973	20.0