

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
 Data File : BF081757.D
 Acq On : 23 Sep 2015 5:52
 Operator : UM/IZ
 Sample : G3748-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MGP202-30-32

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-BF090115.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	1.247	10	14	17	rVB	1069631	1211886	27.78%	5.447%
2	1.430	29	30	36	rBV	238136	618592	14.18%	2.780%
3	1.521	36	38	79	rVB	1183033	4362215	100.00%	19.607%
4	4.436	288	293	314	rBV	249462	793697	18.19%	3.567%
5	5.316	366	370	395	rBV	658827	1445190	33.13%	6.496%
6	7.579	564	568	572	rBV	617571	1397517	32.04%	6.281%
7	7.659	572	575	583	rVV	904368	1653432	37.90%	7.432%
8	7.796	583	587	591	rVB	156392	246203	5.64%	1.107%
9	8.128	613	616	620	rBV	165167	275524	6.32%	1.238%
10	8.459	641	645	648	rBV	650879	1096661	25.14%	4.929%
11	8.516	648	650	655	rVB	101501	162562	3.73%	0.731%
12	8.688	662	665	672	rBB	176687	315892	7.24%	1.420%
13	9.133	702	704	709	rBV	43385	72039	1.65%	0.324%
14	9.442	726	731	744	rVB	502984	1038796	23.81%	4.669%
15	9.819	762	764	774	rVB	75452	145350	3.33%	0.653%
16	10.505	821	824	836	rVB	48913	136588	3.13%	0.614%
17	10.859	849	855	862	rBV	32582	94240	2.16%	0.424%
18	11.031	867	870	873	rVV	214350	364710	8.36%	1.639%
19	11.236	882	888	895	rBB2	51468	92743	2.13%	0.417%
20	12.185	968	971	979	rVB	20495	46508	1.07%	0.209%
21	12.505	996	999	1006	rVB	42186	76814	1.76%	0.345%
22	13.682	1096	1102	1105	rBV	790034	1550671	35.55%	6.970%
23	14.734	1190	1194	1204	rBB	154553	261245	5.99%	1.174%
24	15.157	1227	1231	1237	rBB2	222558	373687	8.57%	1.680%
25	16.437	1339	1343	1347	rBB	55505	90843	2.08%	0.408%
26	17.066	1394	1398	1409	rVB	496766	976588	22.39%	4.389%
27	17.980	1466	1478	1481	rBV	83464	400452	9.18%	1.800%
28	18.666	1534	1538	1542	rBV	229738	389553	8.93%	1.751%
29	19.626	1619	1622	1630	rBB	74253	136946	3.14%	0.616%
30	20.437	1689	1693	1703	rBV	37500	85336	1.96%	0.384%
31	21.386	1773	1776	1782	rBV	36574	54603	1.25%	0.245%
32	22.049	1831	1834	1837	rBV	1046107	1391699	31.90%	6.255%
33	23.295	1940	1943	1944	rBV	188780	213279	4.89%	0.959%
34	23.329	1944	1946	1953	rVB	282613	328442	7.53%	1.476%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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-BF090115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	24.769	2070	2072	2077	rBV	191211	215994	4.95%	0.971%
36	25.192	2104	2109	2112	rBV2	28381	62114	1.42%	0.279%
37	27.490	2302	2310	2314	rVB2	26725	69972	1.60%	0.315%

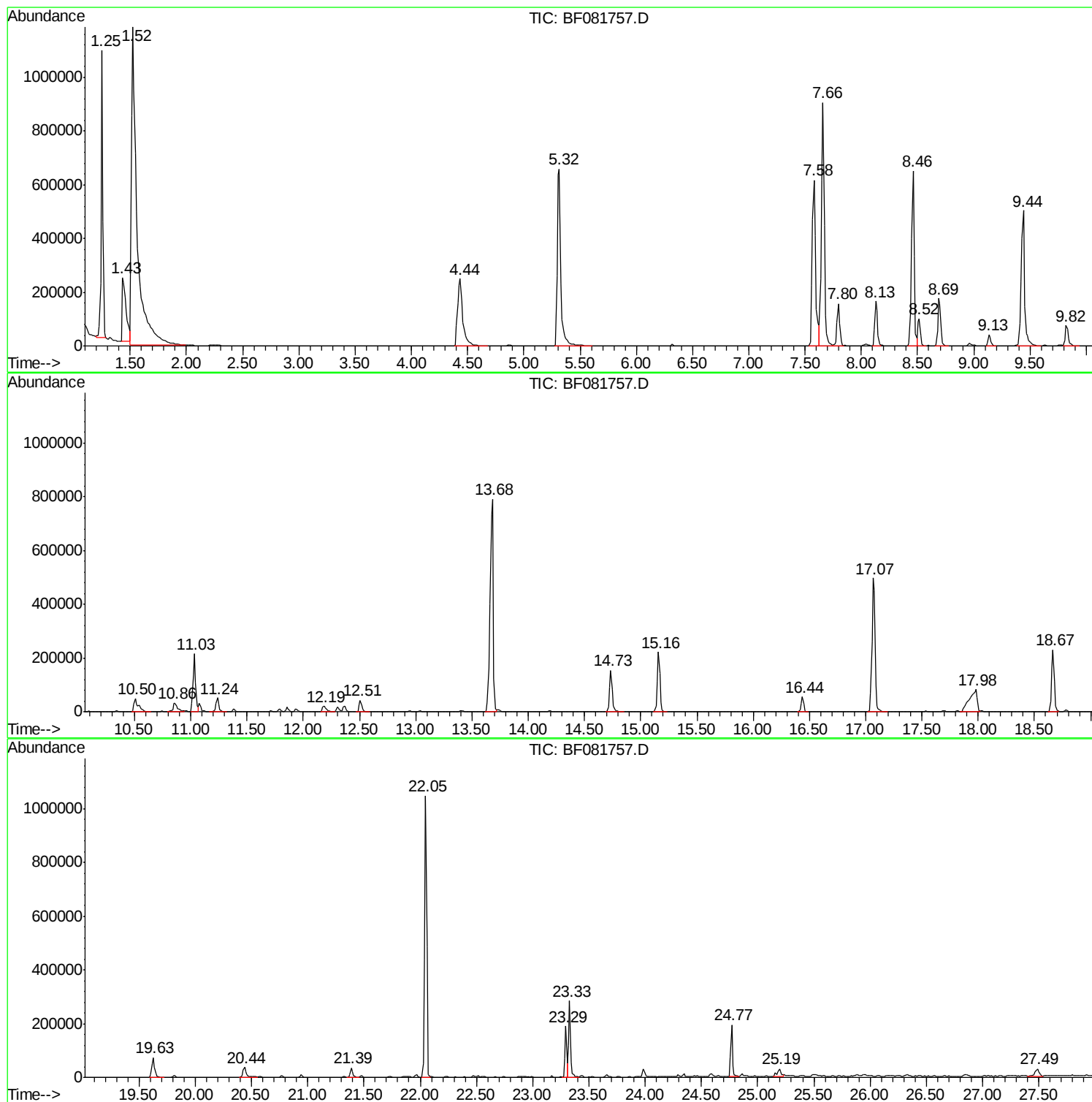
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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MGP202-30-32

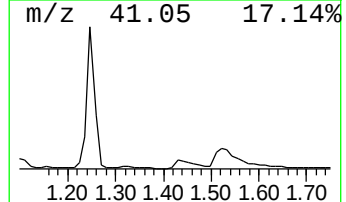
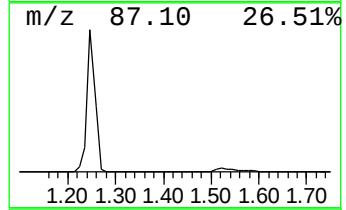
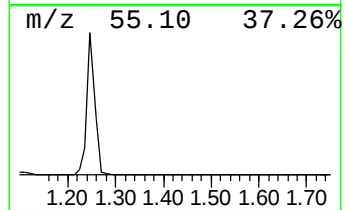
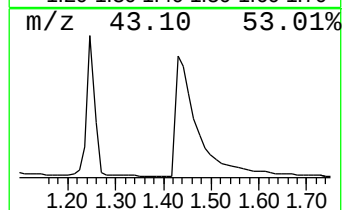
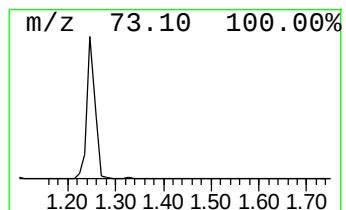
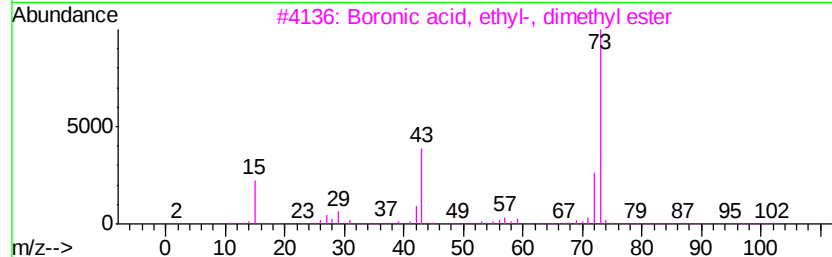
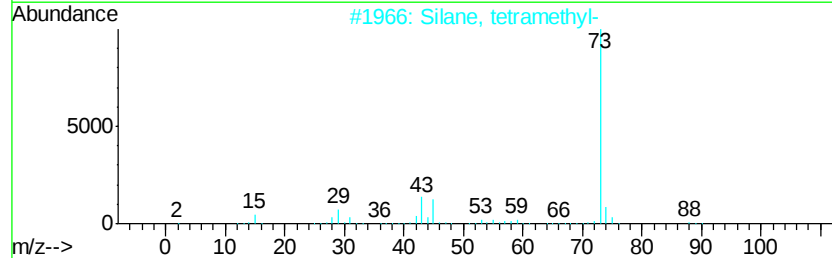
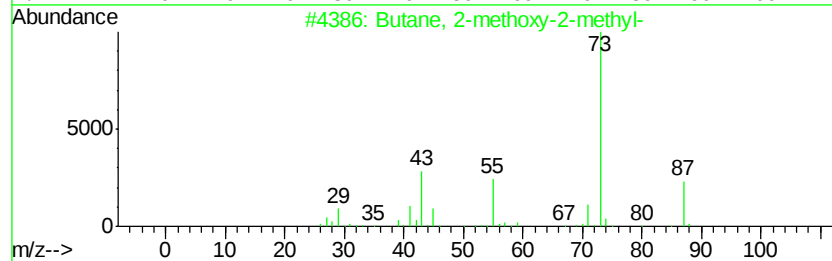
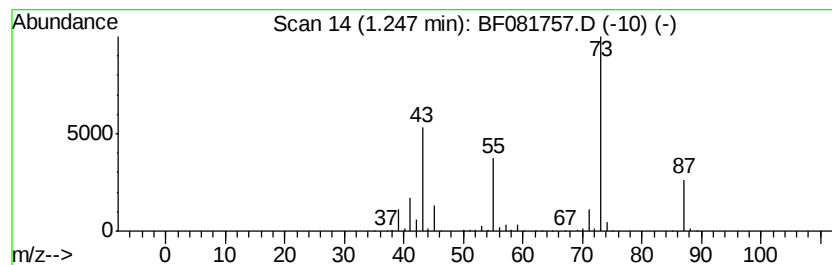
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.25	87.97 ng	1211890	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3		Boronic acid, ethyl-, dimethyl e...	102	C4H11B02	007318-82-3	10
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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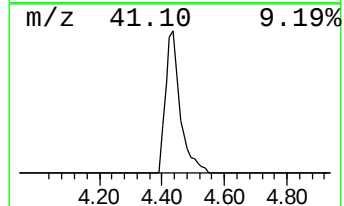
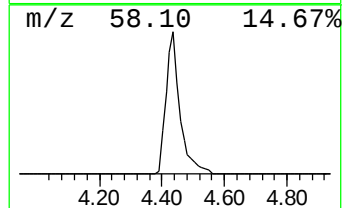
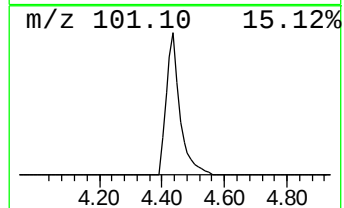
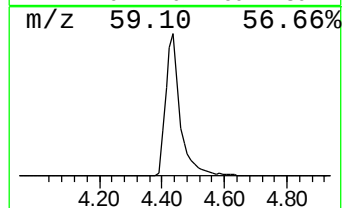
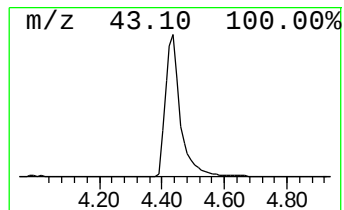
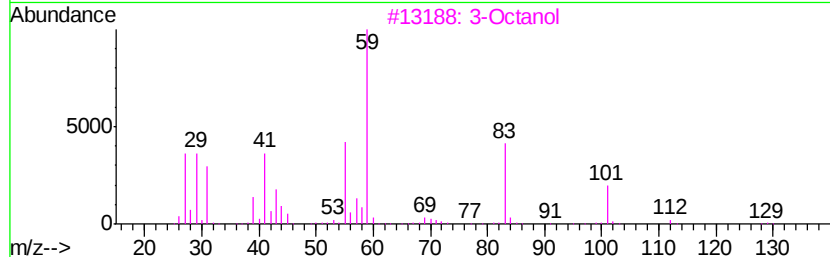
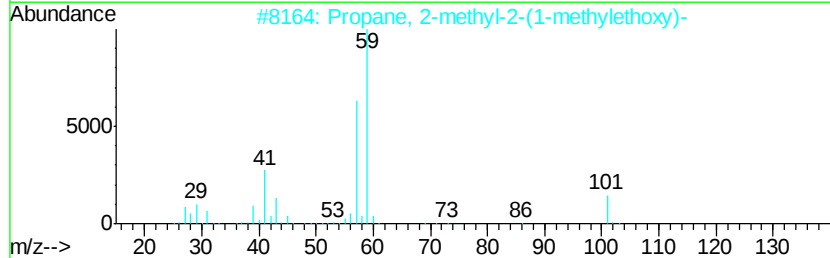
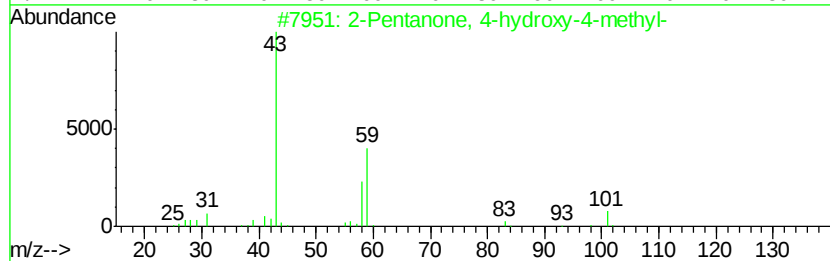
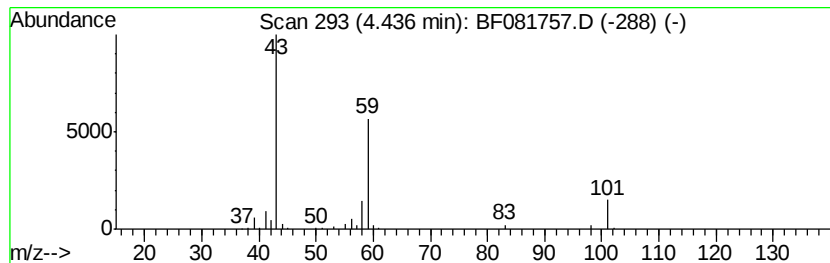
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.44	57.61 ng	793697	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		3-Octanol	130	C8H18O	000589-98-0	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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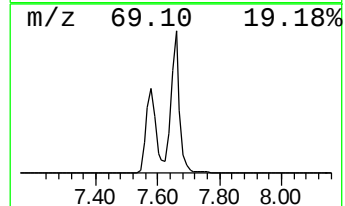
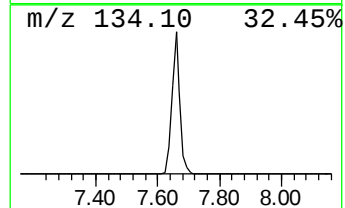
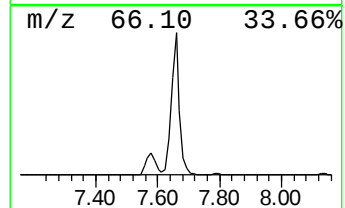
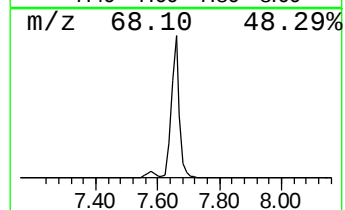
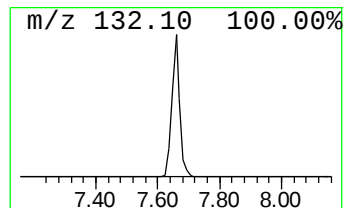
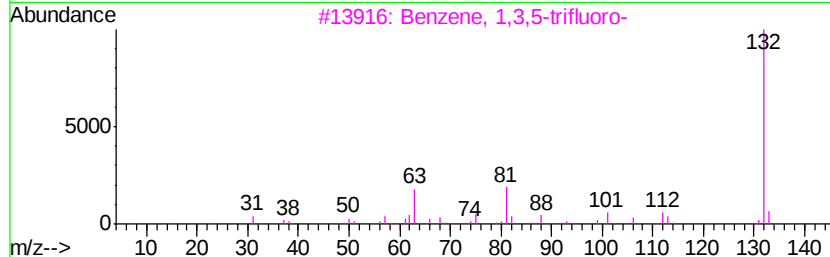
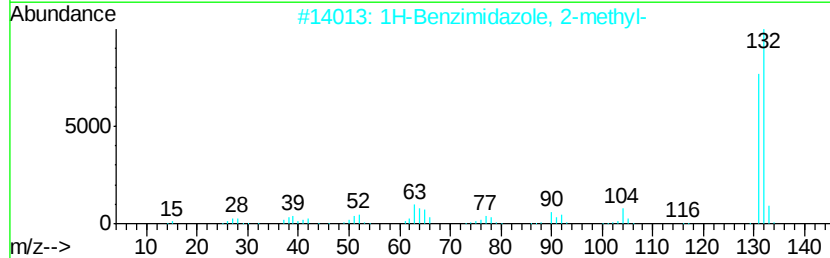
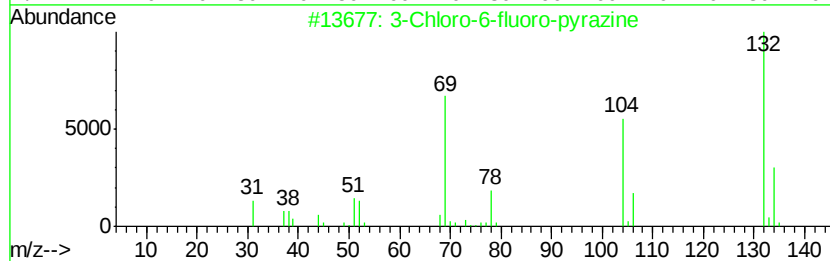
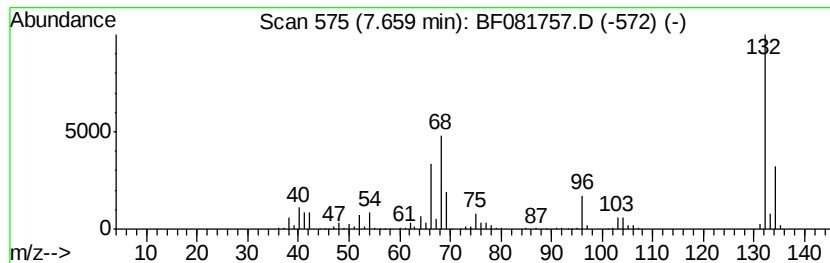
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	120.02 ng	1653430	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	14
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	10



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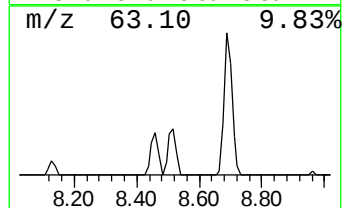
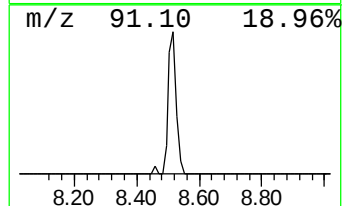
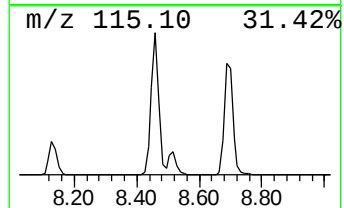
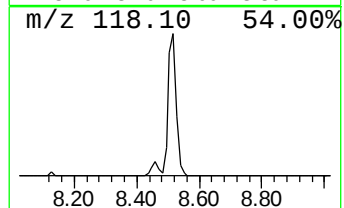
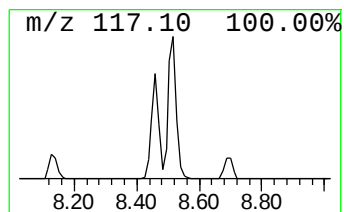
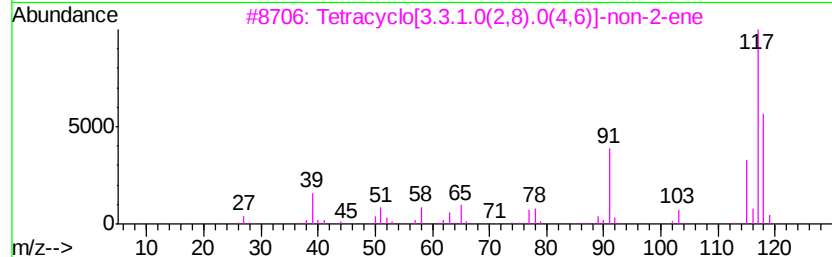
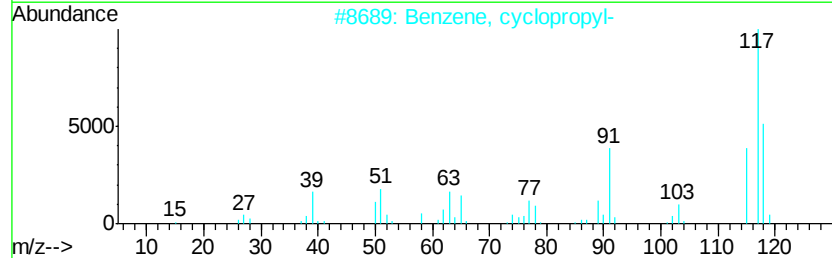
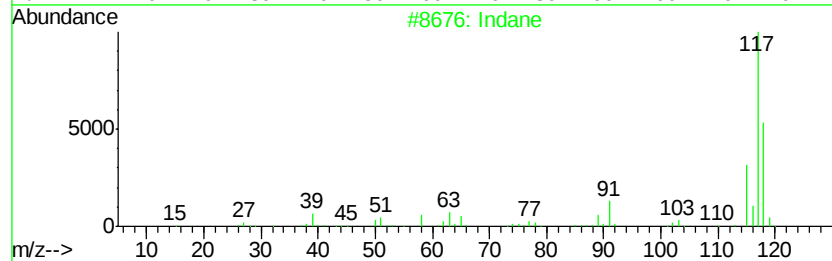
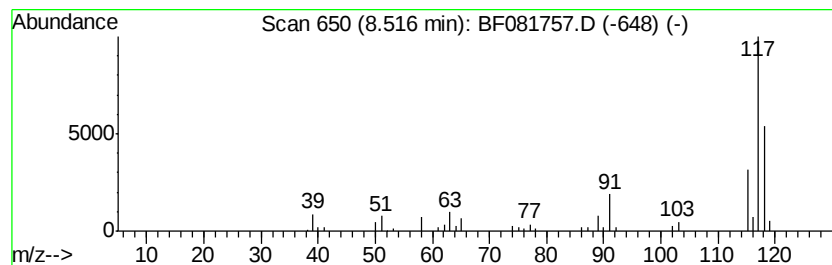
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.52	11.80 ng	162562	1,4-Dichlorobenzene-d4	8.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	87
2	Benzene, cyclopropyl-		118	C9H10	000873-49-4	81
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-	...	118	C9H10	1000191-13-7	80
4	Benzene, 1-propenyl-		118	C9H10	000637-50-3	76
5	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	64



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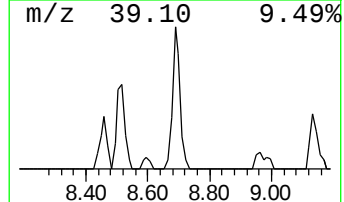
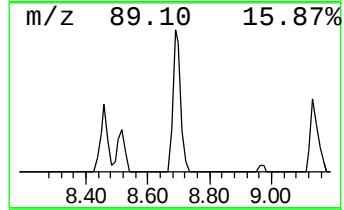
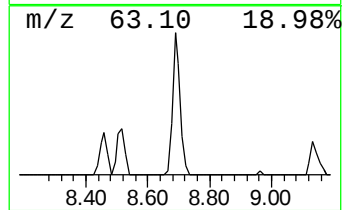
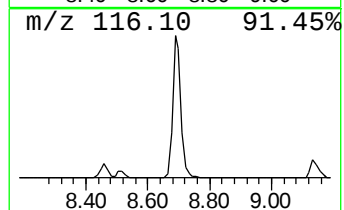
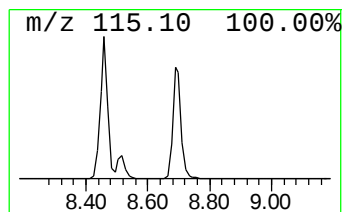
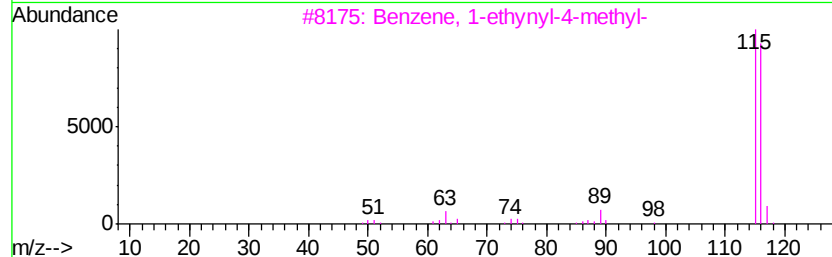
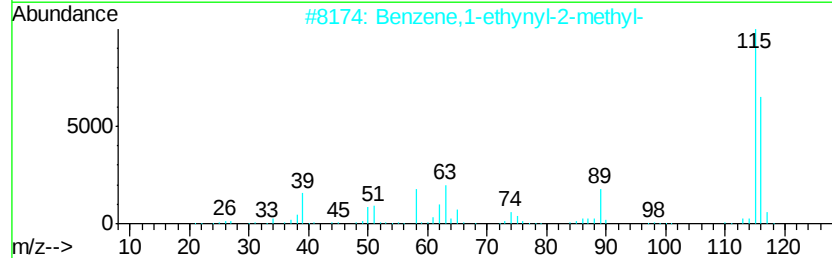
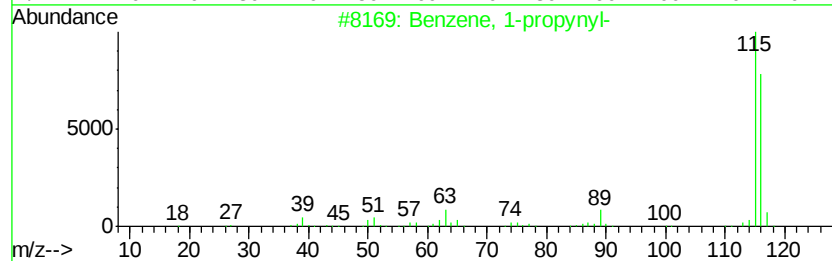
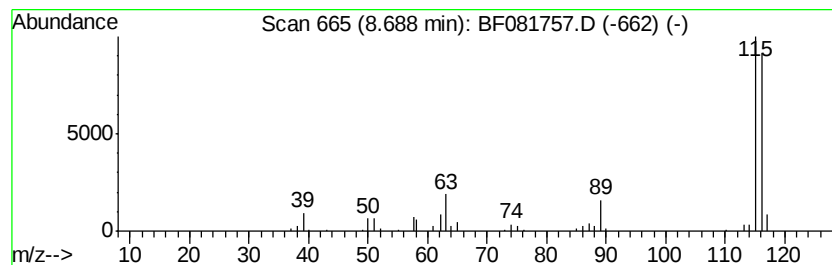
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-propynyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.69	22.93 ng	315892	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
2			Benzene,1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91
3			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
4			Indene	116	C9H8	000095-13-6	90
5			1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	56



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081757.D
Acq On : 23 Sep 2015 5:52
Operator : UM/IZ
Sample : G3748-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MGP202-30-32

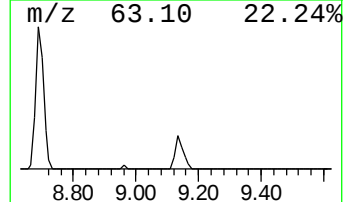
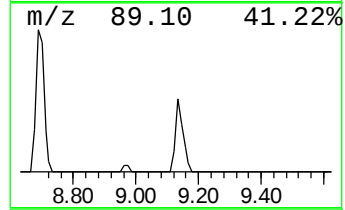
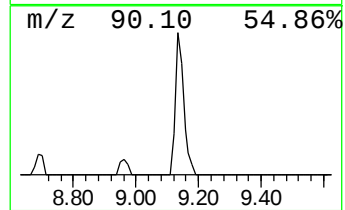
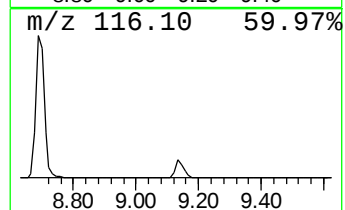
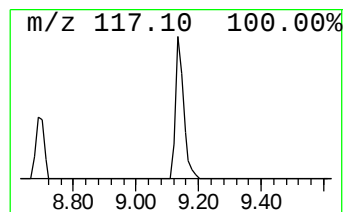
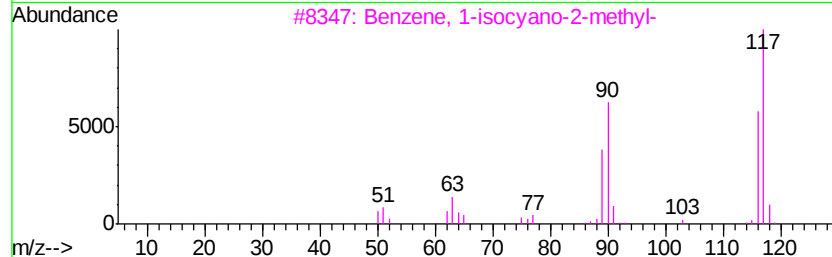
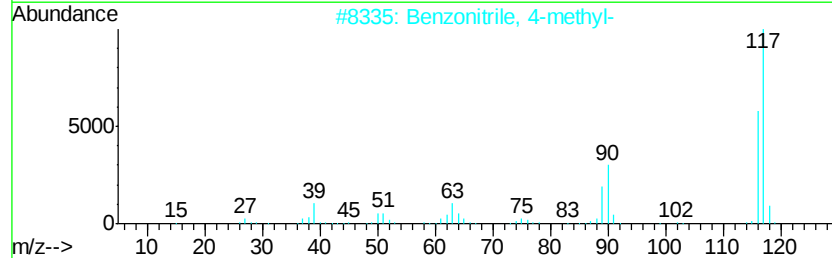
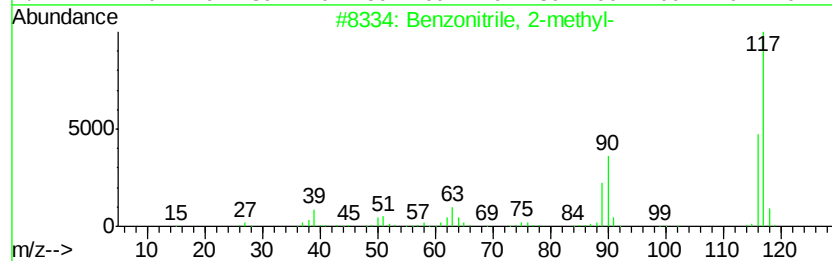
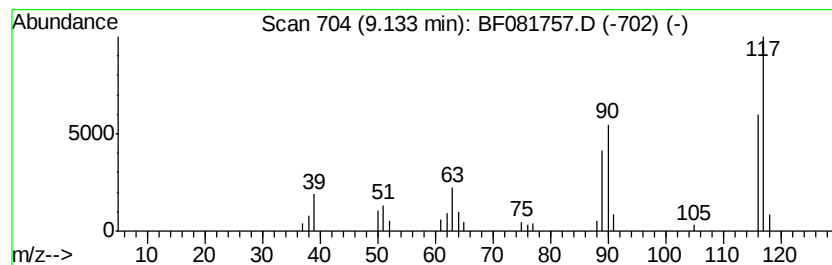
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzonitrile, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.13	5.23 ng	72039	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	96
2			Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	95
3			Benzene, 1-isocvano-2-methyl-	117	C8H7N	010468-64-1	93
4			Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	91
5			5H-1-Pyridine	117	C8H7N	000270-91-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092215\
Data File : BF081757.D
Acq On : 23 Sep 2015 5:52
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Misc :
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ClientSampleId :
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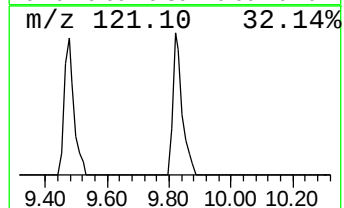
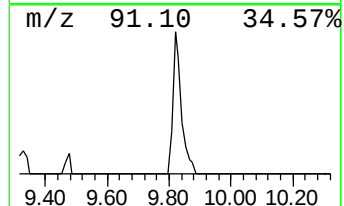
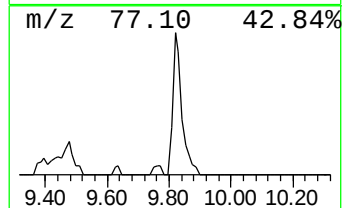
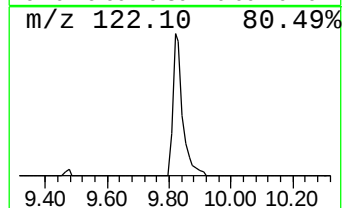
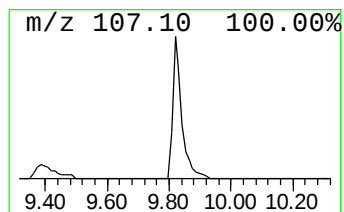
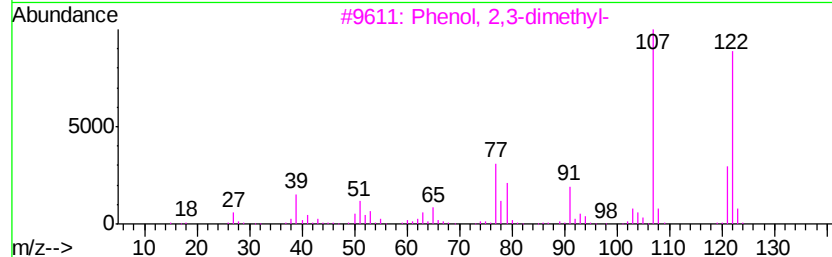
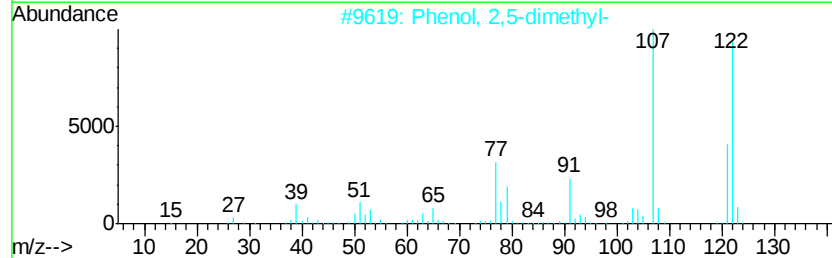
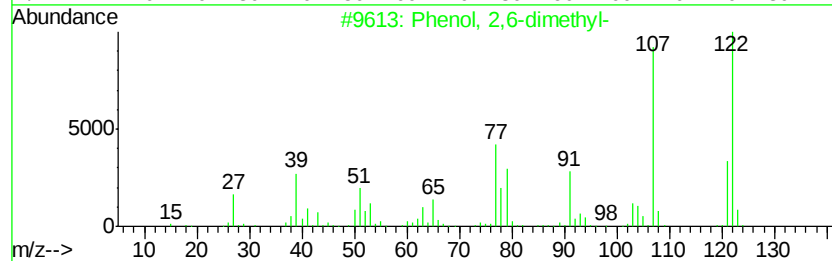
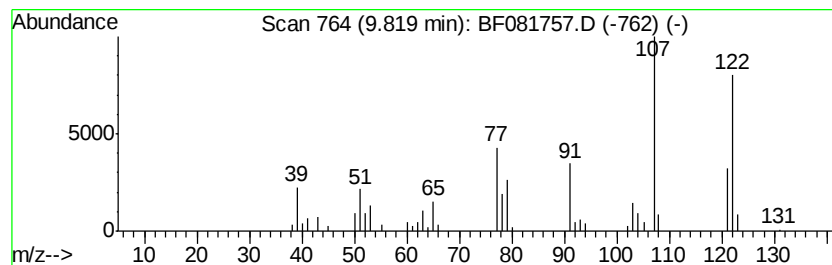
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenol, 2,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	7.97 ng	145350	Napthalene-d8	11.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	97
2		Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	97
3		Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	95
4		Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	93
5		Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081757.D
Acq On : 23 Sep 2015 5:52
Operator : UM/IZ
Sample : G3748-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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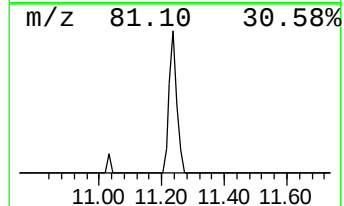
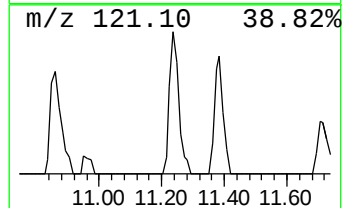
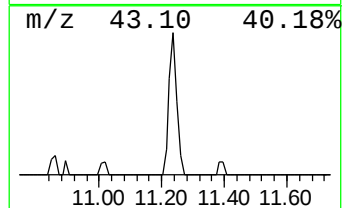
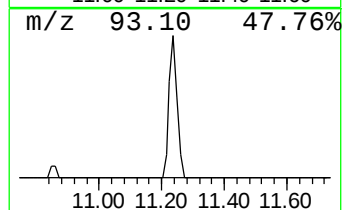
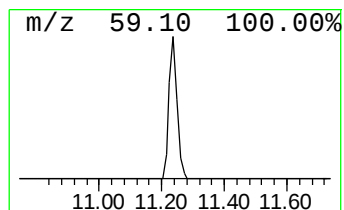
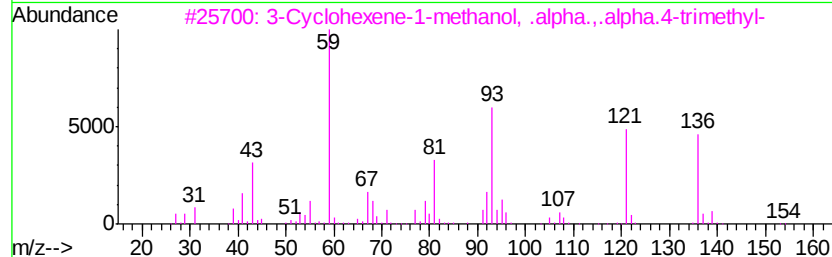
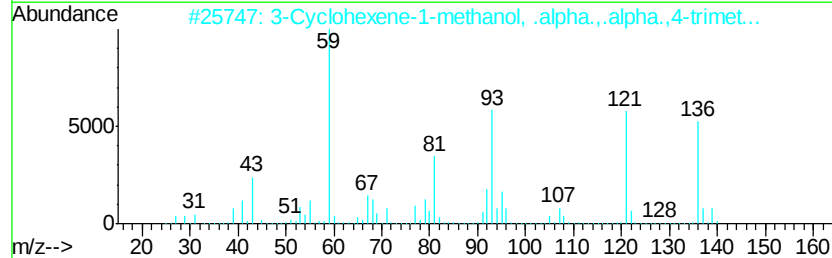
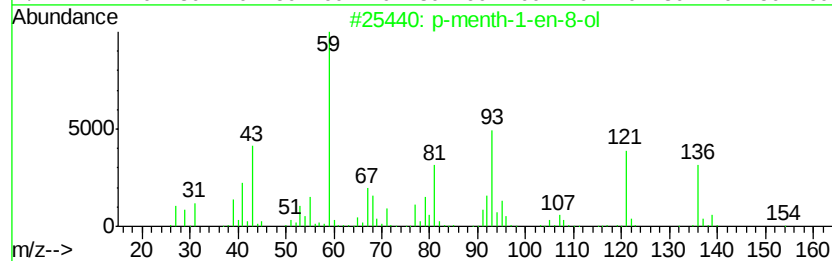
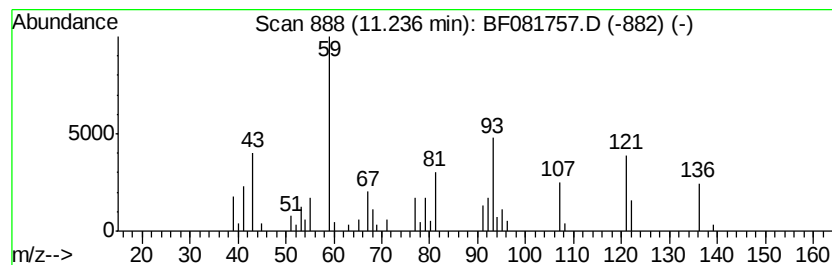
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 p-menth-1-en-8-ol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.24	5.09 ng	92743	Napthalene-d8	11.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-menth-1-en-8-ol	154	C10H18O	1000157-89-9	78
2		3-Cyclohexene-1-methanol, .alpha...	154	C10H18O	010482-56-1	64
3		3-Cyclohexene-1-methanol, .alpha...	154	C10H18O	000098-55-5	64
4		7-Octen-2-ol, 2-methyl-6-methylene-	154	C10H18O	000543-39-5	45
5		p-menth-1-en-8-ol	154	C10H18O	1000151-92-4	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092215\
Data File : BF081757.D
Acq On : 23 Sep 2015 5:52
Operator : UM/IZ
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MGP202-30-32

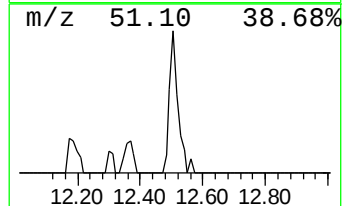
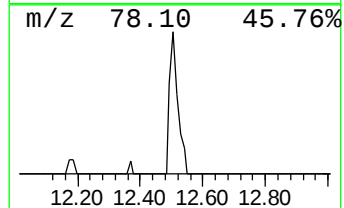
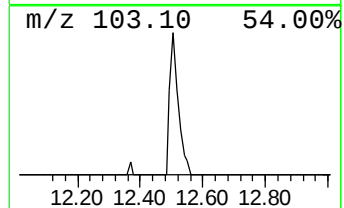
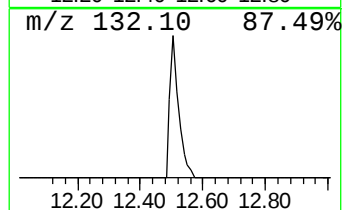
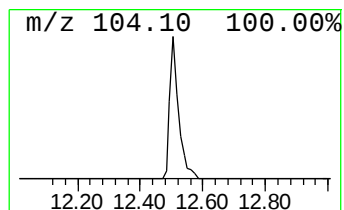
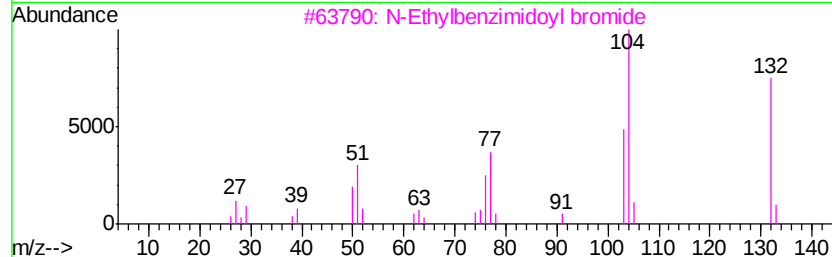
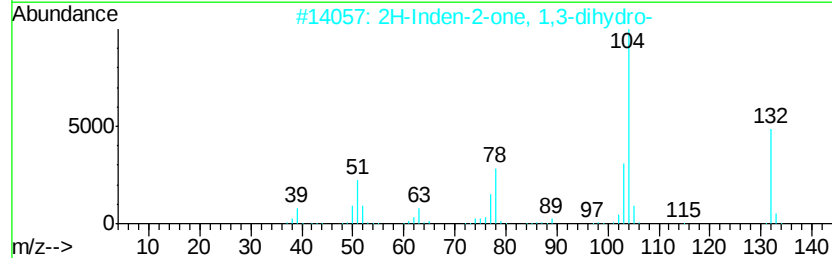
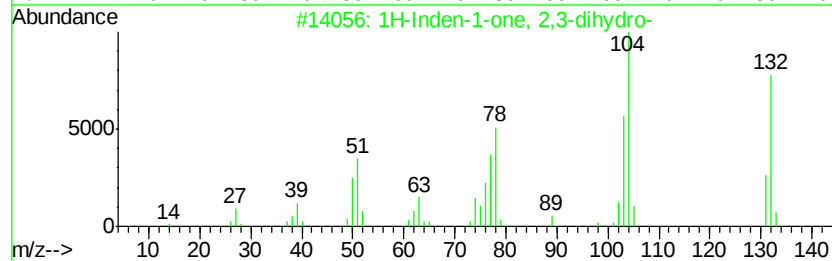
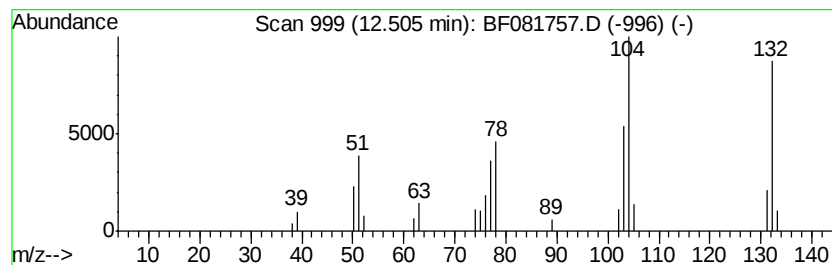
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	4.21 ng	76814	Naphthalene-d8	11.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	98
2			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	70
3			N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	68
4			Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	64
5			Styrene	104	C8H8	000100-42-5	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092215\
Data File : BF081757.D
Acq On : 23 Sep 2015 5:52
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Misc :
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Instrument :
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ClientSampleId :
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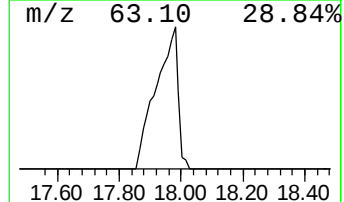
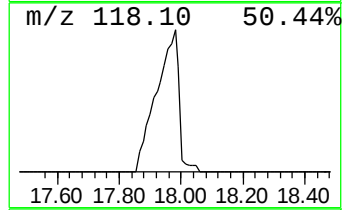
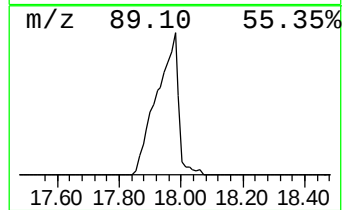
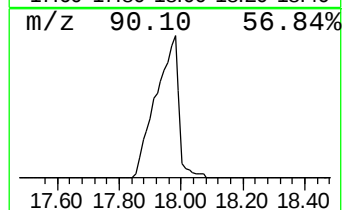
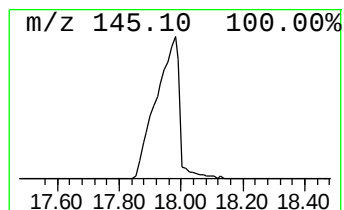
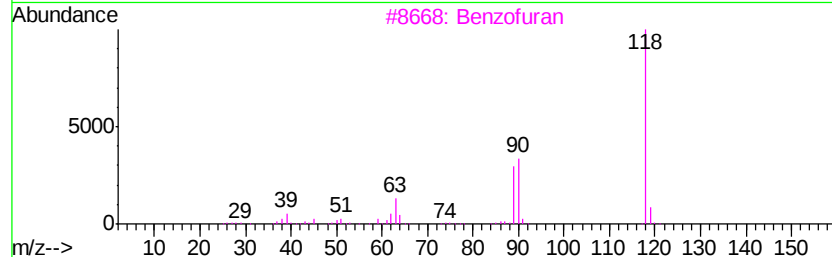
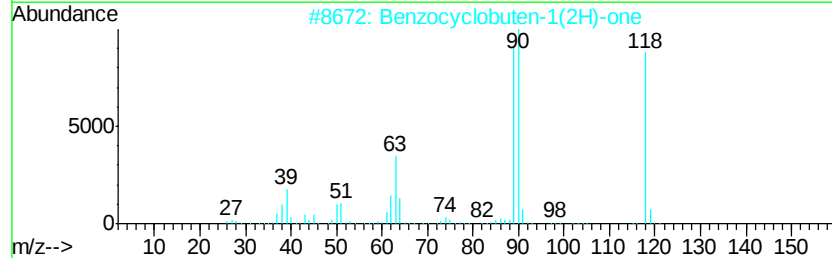
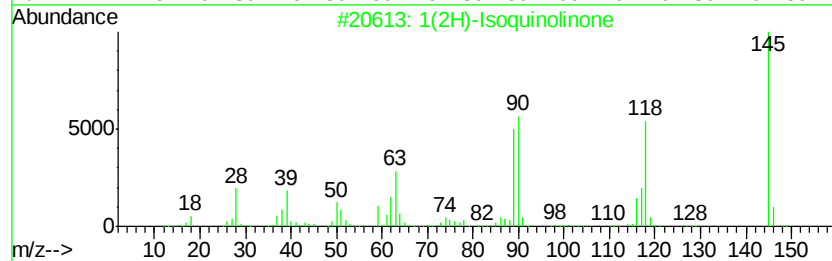
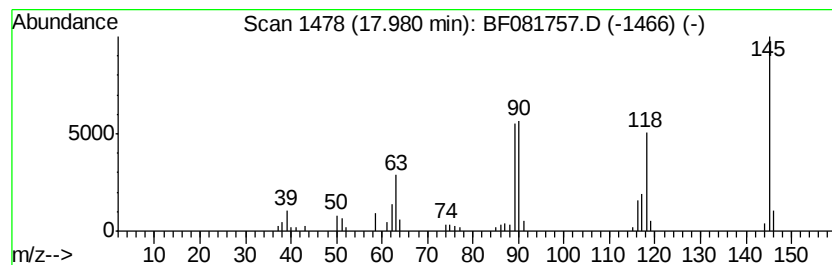
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1(2H)-Isoquinolinone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.98	20.56 ng	400452	Phenanthrene-d10	18.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	96
2			Benzocyclobuten-1(2H)-one	118	C8H6O	003469-06-5	87
3			Benzofuran	118	C8H6O	000271-89-6	78
4			3-Quinolinol	145	C9H7NO	000580-18-7	62
5			1-Phthalazinamine	145	C8H7N3	019064-69-8	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092215\
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Acq On : 23 Sep 2015 5:52
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ClientSampleId :
MGP202-30-32

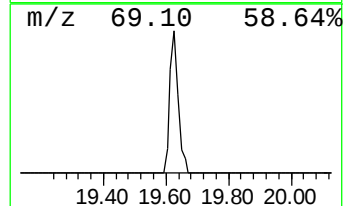
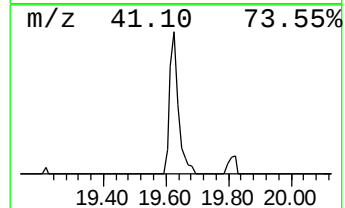
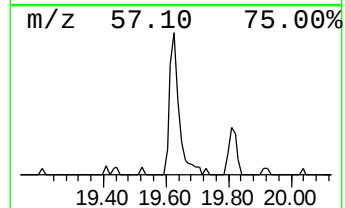
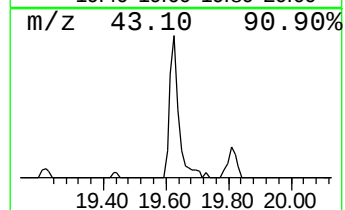
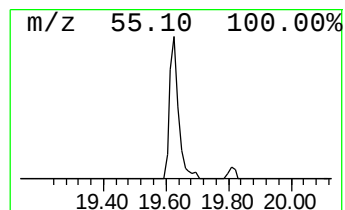
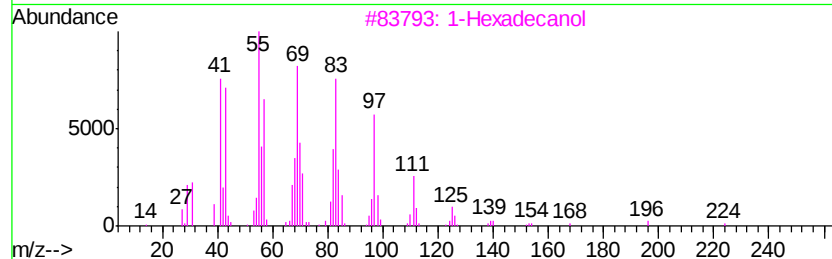
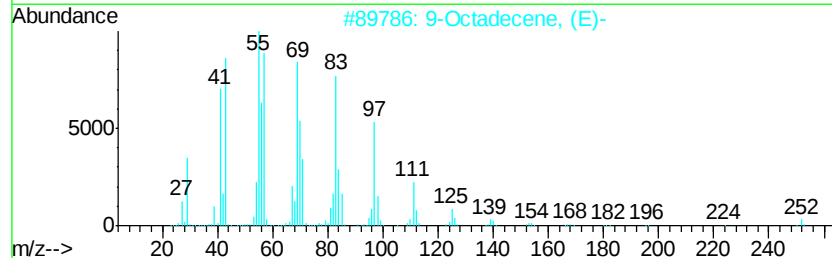
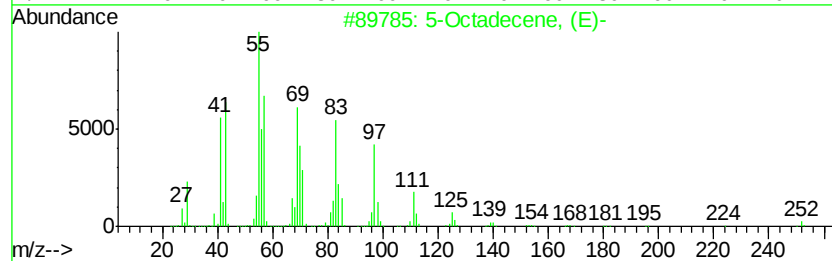
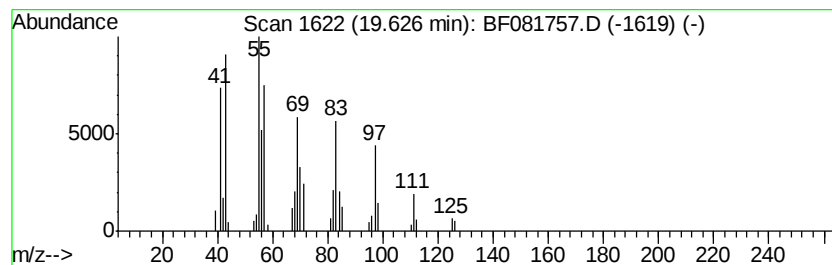
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 5-Octadecene, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.63	7.03 ng	136946	Phenanthrene-d10	18.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	91
2			9-Octadecene, (E)-	252	C18H36	007206-25-9	90
3			1-Hexadecanol	242	C16H34O	036653-82-4	87
4			1-Hexadecene	224	C16H32	000629-73-2	83
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092215\
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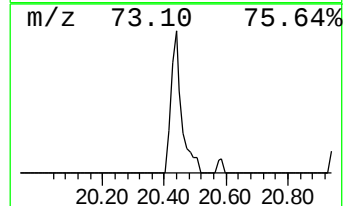
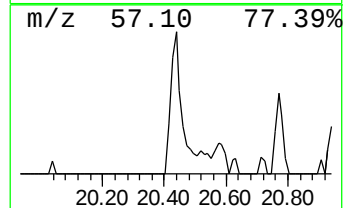
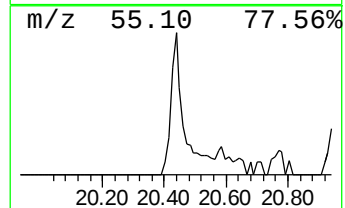
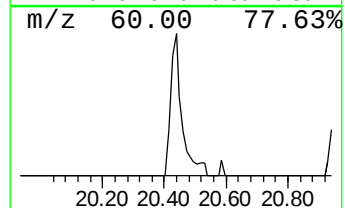
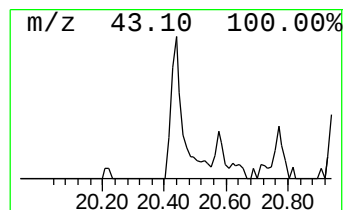
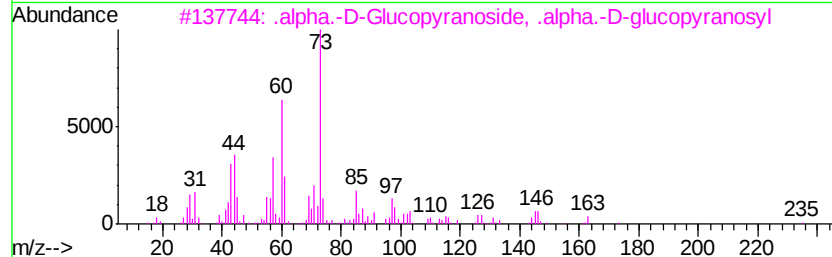
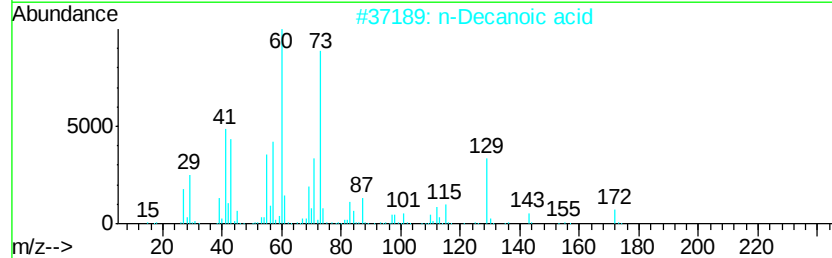
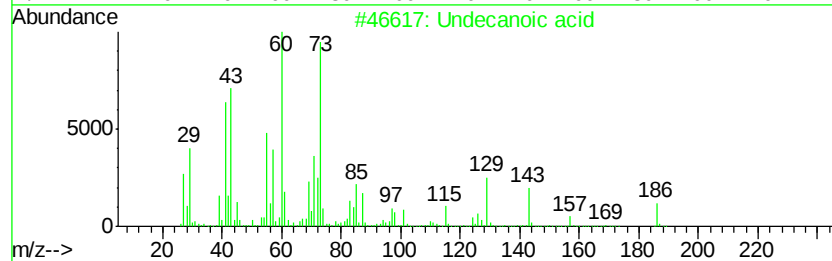
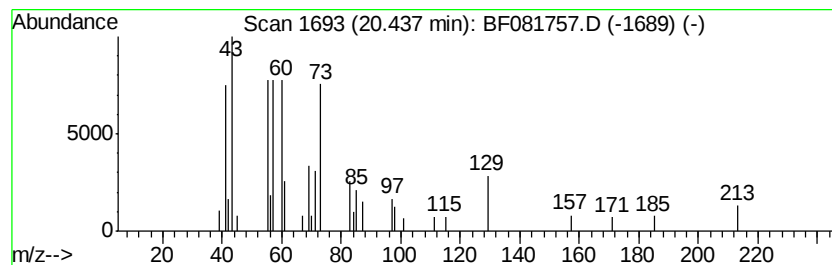
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Undecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.44	4.38 ng	85336	Phenanthrene-d10	18.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecanoic acid		186	C11H22O2	000112-37-8	64
2	n-Decanoic acid		172	C10H20O2	000334-48-5	53
3	.alpha.-D-Glucopyranoside, .alph...		342	C12H22O11	000099-20-7	53
4	Nonanoic acid		158	C9H18O2	000112-05-0	50
5	11-Bromoundecanoic acid		264	C11H21BrO2	002834-05-1	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081757.D
Acq On : 23 Sep 2015 5:52
Operator : UM/IZ
Sample : G3748-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MGP202-30-32

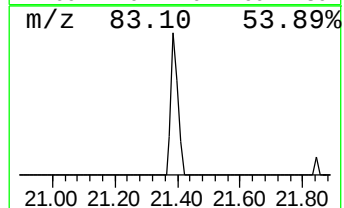
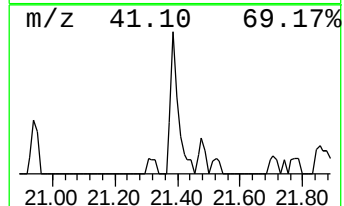
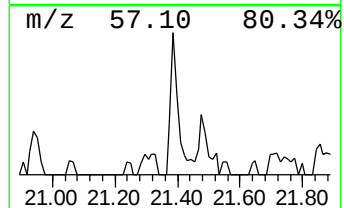
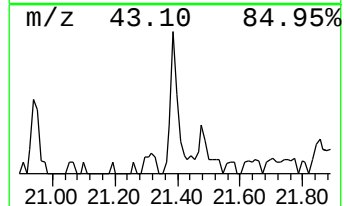
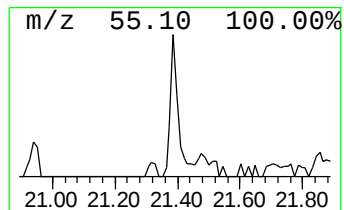
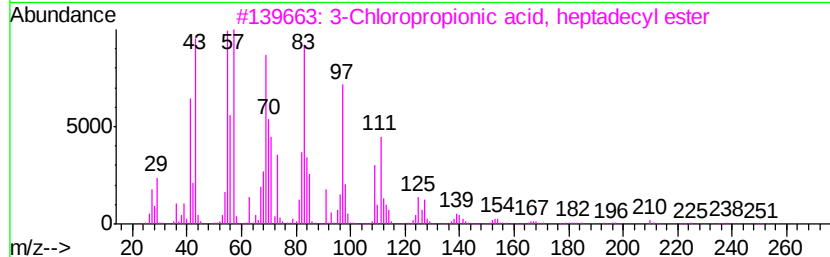
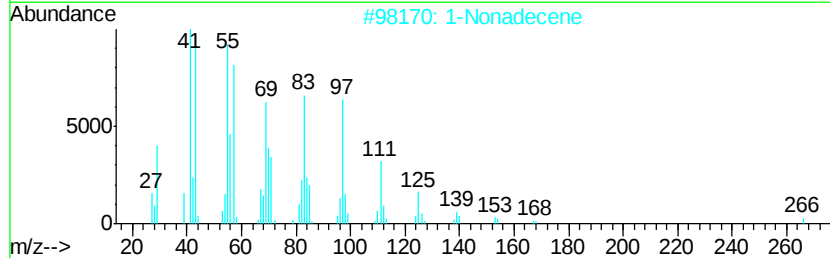
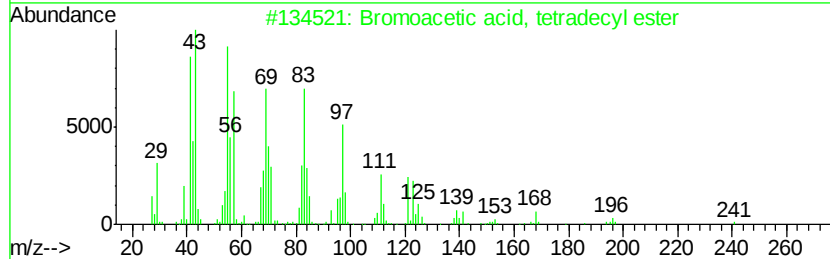
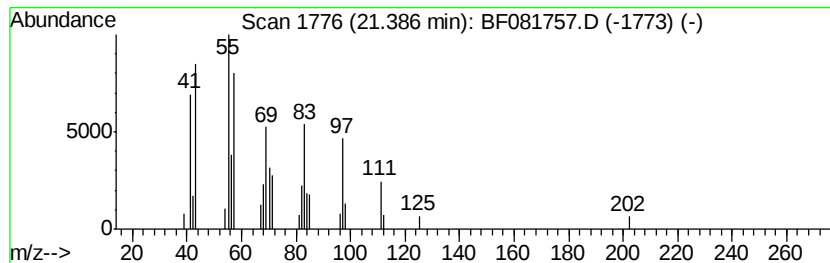
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Bromoacetic acid, tetradecy... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.39	3.32 ng	54603	Chrysene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, tetradecyl ester	334	C16H31BrO2	018992-01-3	91
2		1-Nonadecene	266	C19H38	018435-45-5	91
3		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
4		1-Hexadecene	224	C16H32	000629-73-2	91
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	90



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Acq On : 23 Sep 2015 5:52
Operator : UM/IZ
Sample : G3748-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MGP202-30-32

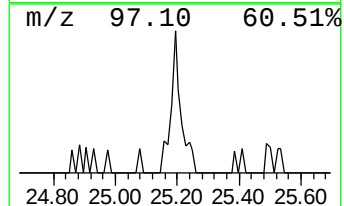
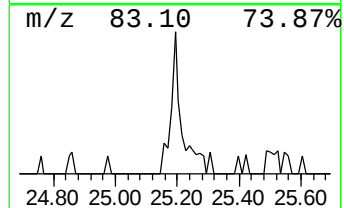
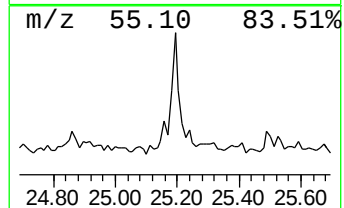
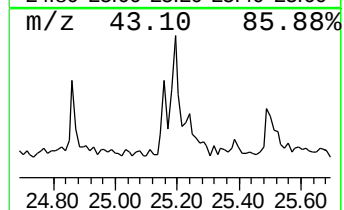
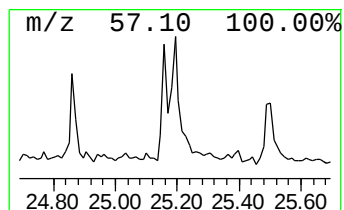
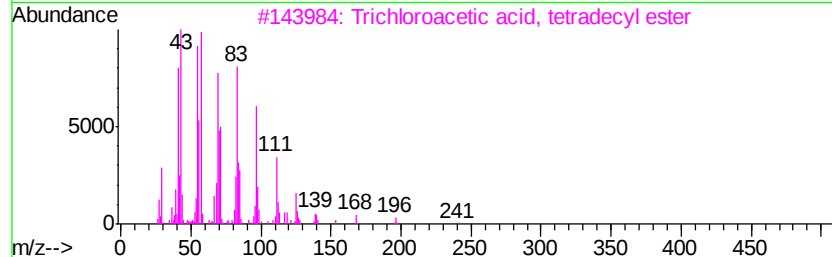
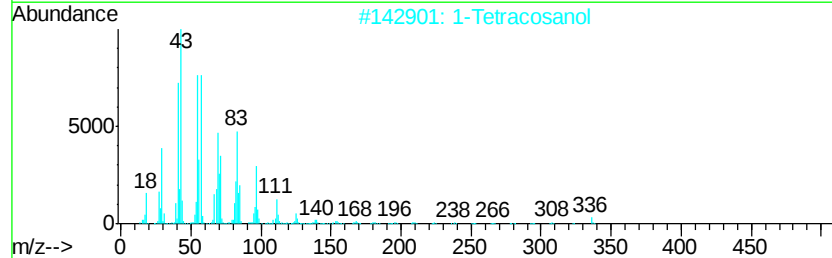
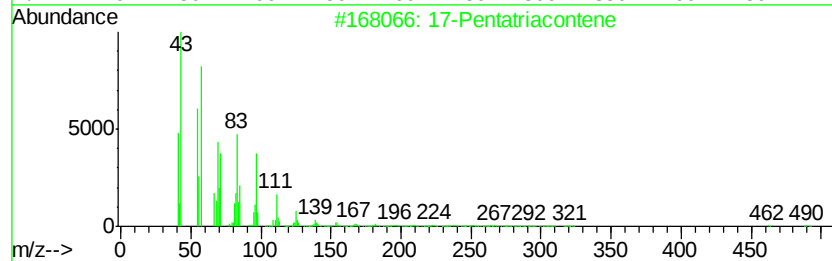
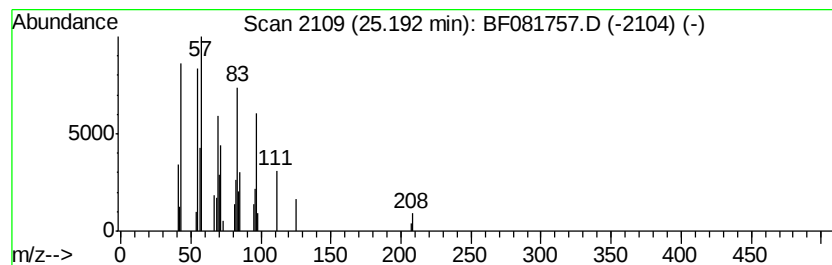
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 17-Pentatriacontene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.19	5.75 ng	62114	Perylene-d12	24.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			17-Pentatriacontene	491	C35H70	006971-40-0	87
2			1-Tetracosanol	354	C24H50O	000506-51-4	74
3			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	72
4			1-Hexacosanol	382	C26H54O	000506-52-5	72
5			1-Hentetracontanol	593	C41H84O	040710-42-7	58



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Acq On : 23 Sep 2015 5:52
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Sample : G3748-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MGP202-30-32

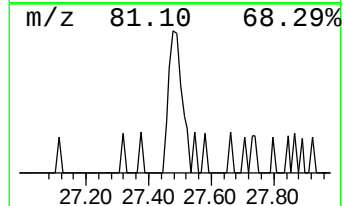
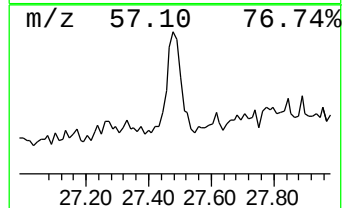
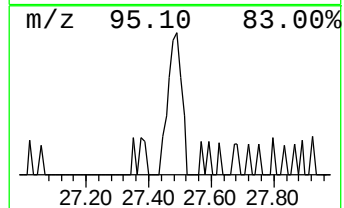
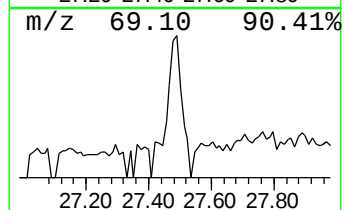
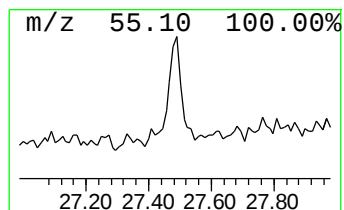
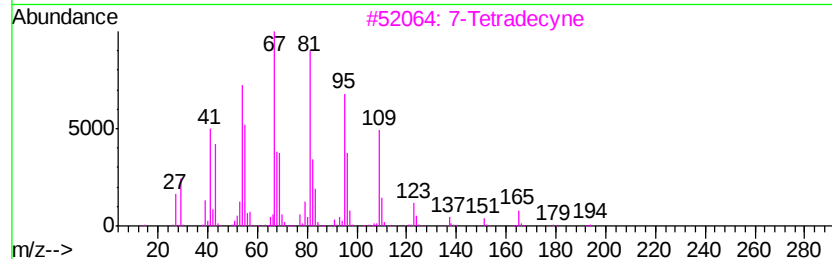
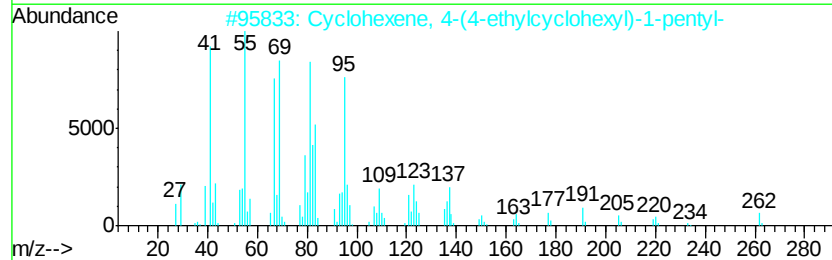
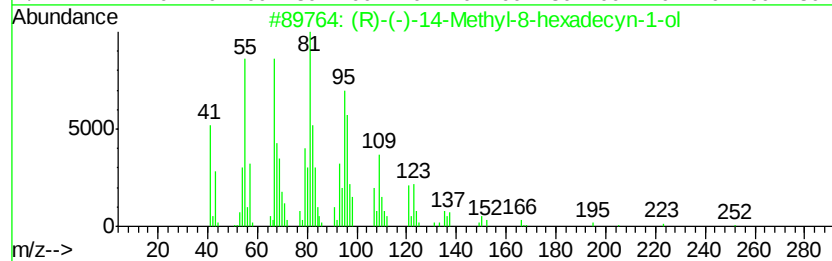
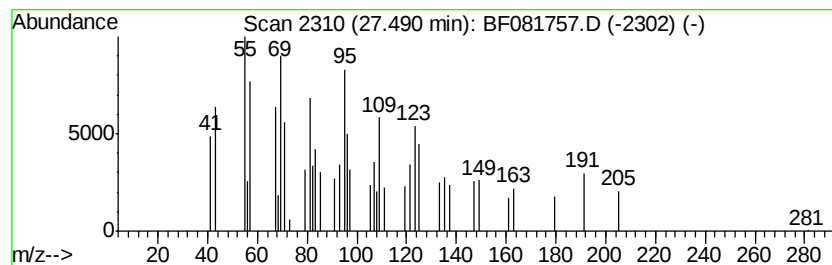
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown27.49 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.49	6.48 ng	69972	Perylene-d12	24.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(R)-(-)-14-Methyl-8-hexadecyn-1-ol	252	C17H32O	064566-18-3	38		
2	Cyclohexene, 4-(4-ethylcyclohexyl)-1-pentyl-	262	C19H34	301643-32-3	38		
3	7-Tetradecyne	194	C14H26	035216-11-6	27		
4	7-Heptadecyne, 17-chloro-	270	C17H31Cl	056554-75-7	27		
5	1-Formyl-2,2-dimethyl-3-trans-(3...	220	C15H24O	1000144-09-7	27		



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 Operator : UM/IZ
 Sample : G3748-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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
Butane, 2-methoxy...	1.25	88.0	ng	1211890	1	8.13	275524	20.0
2-Pentanone, 4-hy...	4.44	57.6	ng	793697	1	8.13	275524	20.0
unknown7.66	7.66	120.0	ng	1653430	1	8.13	275524	20.0
Indane	8.52	11.8	ng	162562	1	8.13	275524	20.0
Benzene, 1-propynyl-	8.69	22.9	ng	315892	1	8.13	275524	20.0
Benzonitrile, 2-m...	9.13	5.2	ng	72039	1	8.13	275524	20.0
Phenol, 2,6-dimet...	9.82	8.0	ng	145350	2	11.03	364710	20.0
p-menth-1-en-8-ol	11.24	5.1	ng	92743	2	11.03	364710	20.0
1H-Inden-1-one, 2...	12.51	4.2	ng	76814	2	11.03	364710	20.0
1(2H)-Isoquinolinone	17.98	20.6	ng	400452	4	18.67	389553	20.0
5-Octadecene, (E)-	19.63	7.0	ng	136946	4	18.67	389553	20.0
Undecanoic acid	20.44	4.4	ng	85336	4	18.67	389553	20.0
Bromoacetic acid,...	21.39	3.3	ng	54603	5	23.33	328442	20.0
17-Pentatriacontene	25.19	5.8	ng	62114	6	24.77	215994	20.0
unknown27.49	27.49	6.5	ng	69972	6	24.77	215994	20.0