

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
 Data File : BF080359.D
 Acq On : 6 Jul 2015 22:21
 Operator : TP/IZ
 Sample : G2877-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DCW-WW-155-7215

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-BF070615.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	2.212	8	11	27	rVV	1105560	2017428	100.00%	16.675%
2	2.418	27	29	33	rVB	579154	895894	44.41%	7.405%
3	5.436	290	293	297	rBV	101428	147171	7.29%	1.216%
4	5.836	326	328	330	rBV	309663	335068	16.61%	2.769%
5	6.236	361	363	365	rBV	45170	37008	1.83%	0.306%
6	6.830	413	415	418	rBV	267942	230842	11.44%	1.908%
7	6.990	427	429	432	rBV	785632	687056	34.06%	5.679%
8	7.230	448	450	452	rVB	240147	195804	9.71%	1.618%
9	7.390	462	464	466	rVB	804580	748103	37.08%	6.183%
10	7.790	496	499	501	rBV	403834	540426	26.79%	4.467%
11	8.522	561	563	565	rBV	287411	254685	12.62%	2.105%
12	8.796	584	587	589	rBV	210973	176286	8.74%	1.457%
13	9.470	644	646	650	rBV	19984	20773	1.03%	0.172%
14	9.596	655	657	659	rVB	1709735	1337765	66.31%	11.057%
15	9.985	689	691	693	rVB	82525	61918	3.07%	0.512%
16	10.076	697	699	703	rVV	20178	31330	1.55%	0.259%
17	10.179	706	708	712	rVV3	10093	23713	1.18%	0.196%
18	10.293	716	718	720	rVB	227569	291438	14.45%	2.409%
19	10.693	749	753	755	rBV2	9657	23151	1.15%	0.191%
20	11.082	784	787	789	rBV	726461	711103	35.25%	5.877%
21	11.791	846	849	851	rBV	361928	335060	16.61%	2.769%
22	12.008	865	868	872	rBV	17887	35575	1.76%	0.294%
23	12.865	940	943	948	rBV2	39208	78192	3.88%	0.646%
24	13.459	992	995	997	rBV	1692114	1650856	81.83%	13.645%
25	13.882	1030	1032	1035	rVB	54559	49675	2.46%	0.411%
26	14.499	1083	1086	1095	rVB	57870	116304	5.76%	0.961%
27	14.671	1099	1101	1102	rBV	13466	21553	1.07%	0.178%
28	14.694	1102	1103	1106	rVB	261195	359630	17.83%	2.972%
29	14.922	1120	1123	1130	rVB5	9830	28082	1.39%	0.232%
30	15.791	1196	1199	1201	rBV	43577	57982	2.87%	0.479%
31	16.362	1246	1249	1252	rVB2	15520	21993	1.09%	0.182%
32	16.511	1259	1262	1266	rBV	221272	349566	17.33%	2.889%
33	16.980	1300	1303	1309	rVB	20960	39258	1.95%	0.324%
34	17.723	1364	1368	1371	rBV	22389	54092	2.68%	0.447%

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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

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35	18.603	1442	1445	1452	rVB2	20838	54803	2.72%	0.453%
36	19.700	1536	1541	1547	rVB	15028	50467	2.50%	0.417%
37	21.026	1655	1657	1665	rVB2	8620	28828	1.43%	0.238%

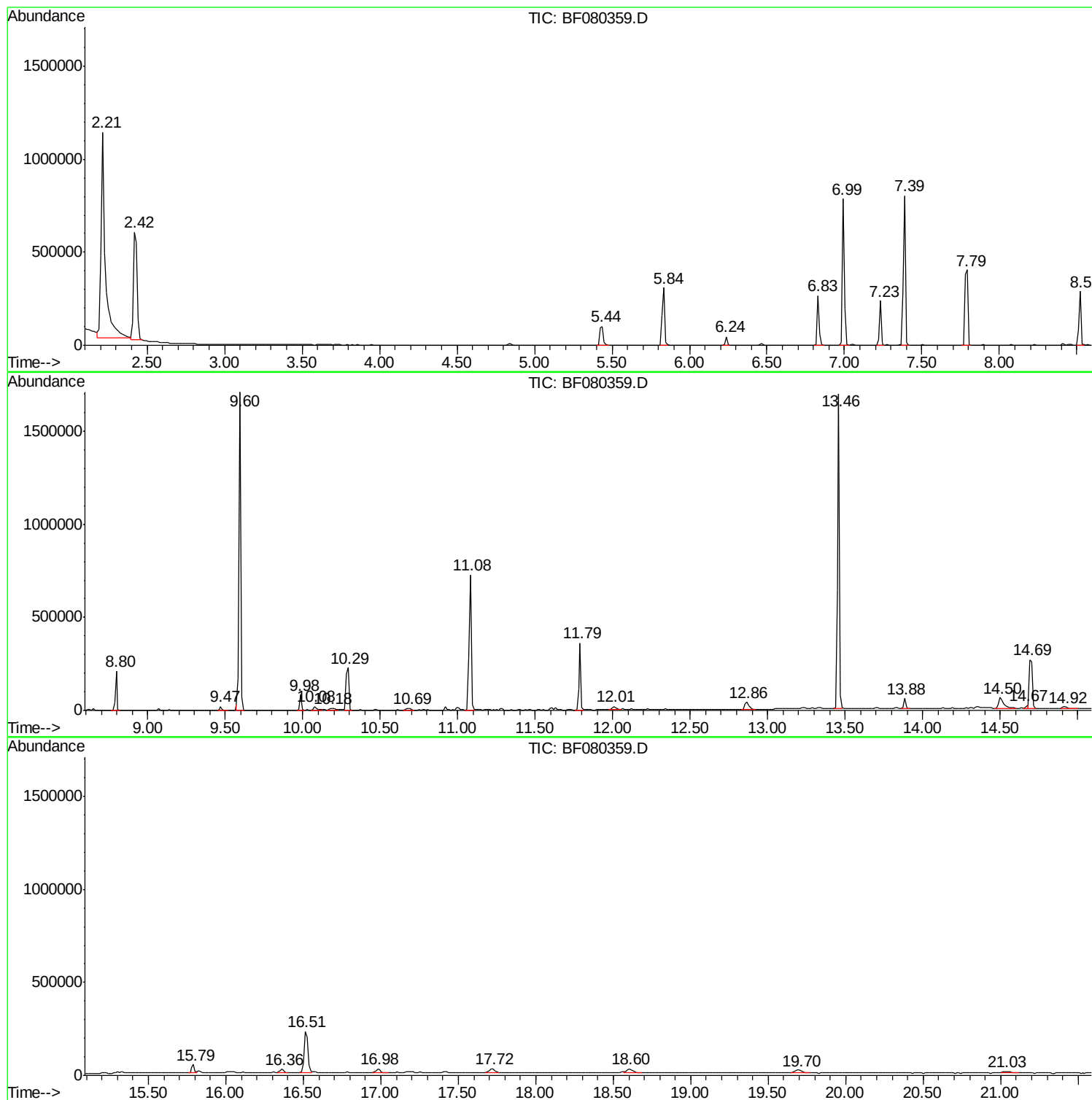
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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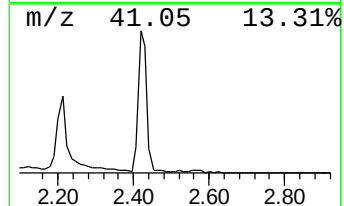
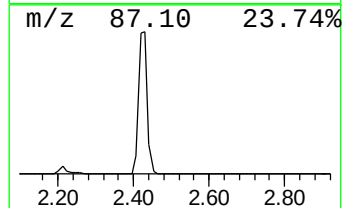
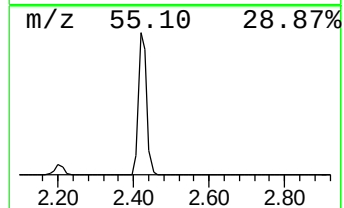
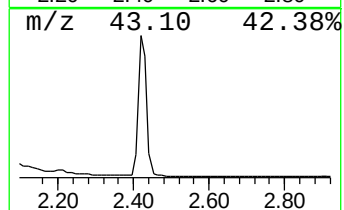
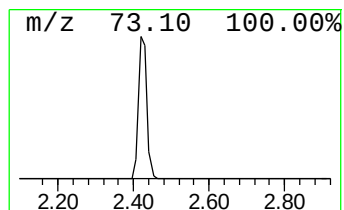
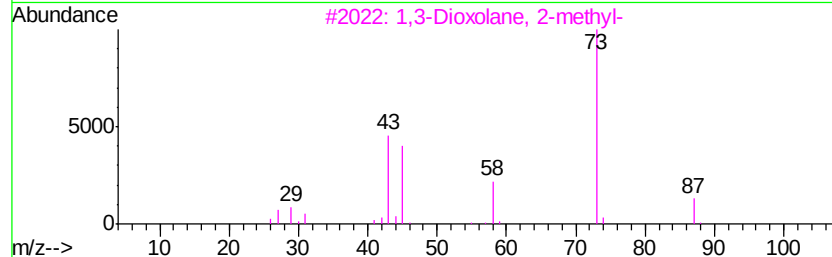
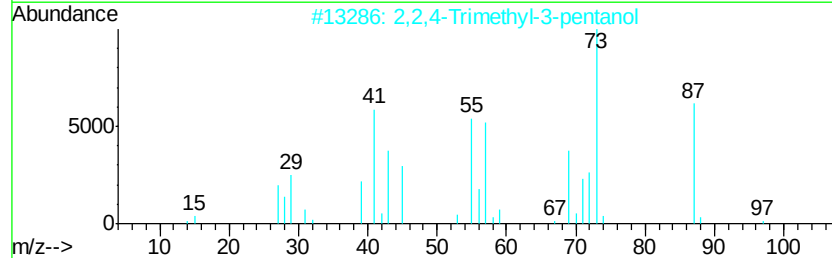
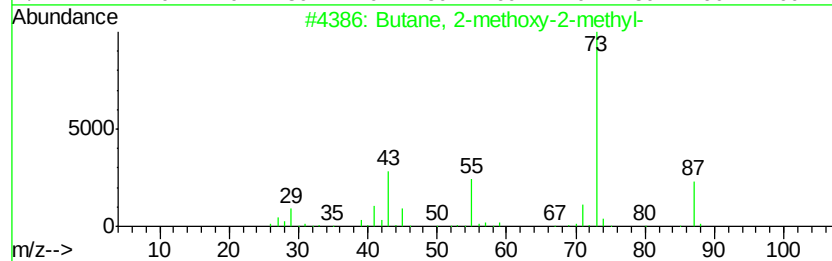
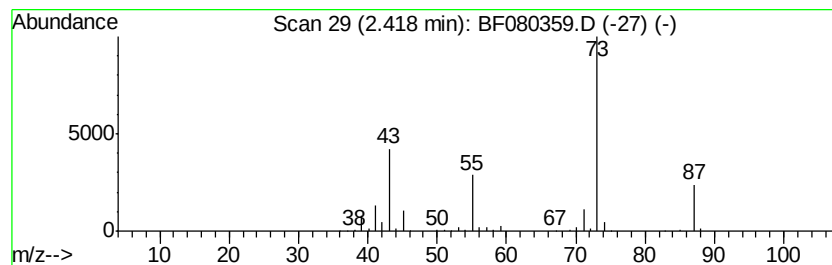
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.42	91.51 ng	895894	1,4-Dichlorobenzene-d4	7.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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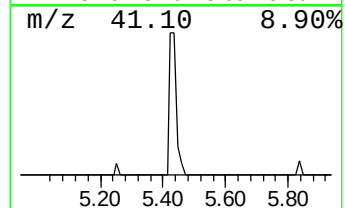
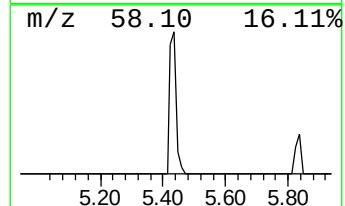
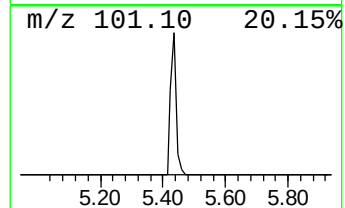
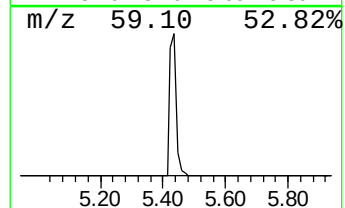
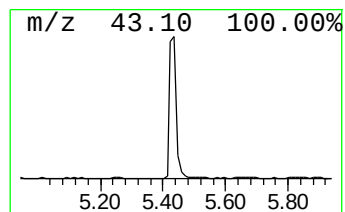
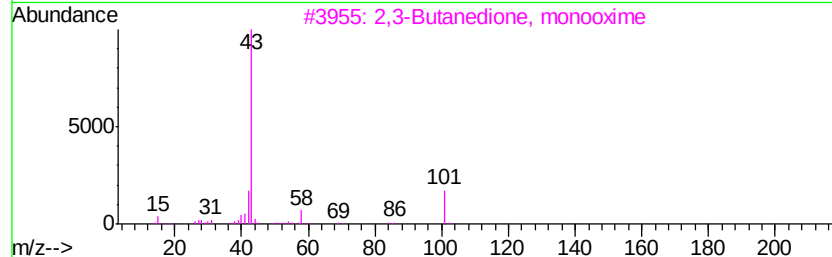
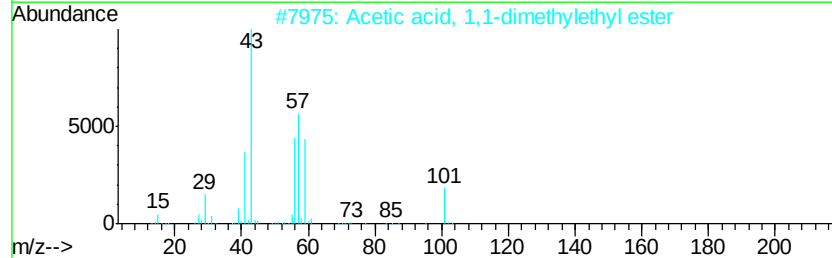
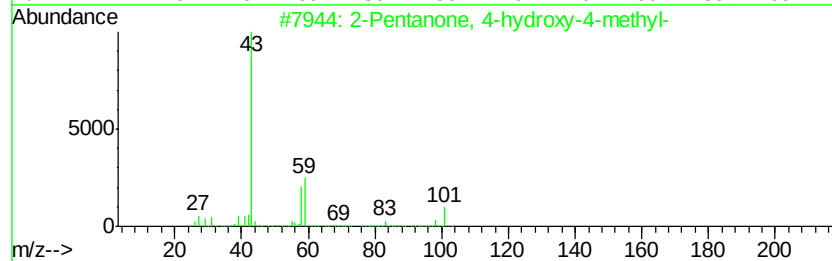
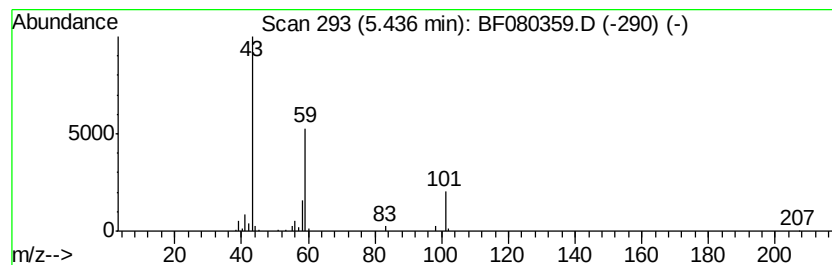
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.44	15.03 ng	147171	1,4-Dichlorobenzene-d4	7.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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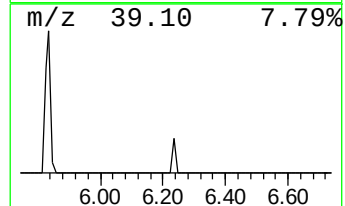
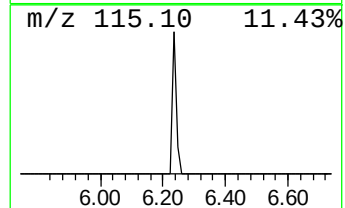
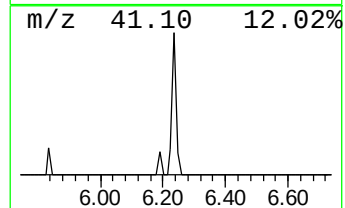
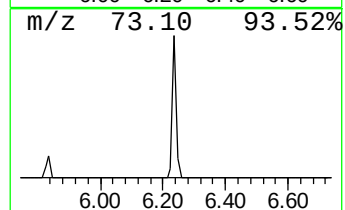
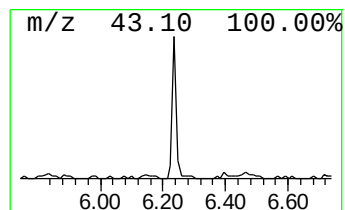
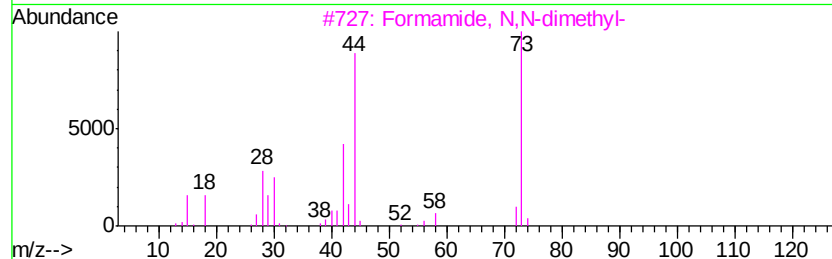
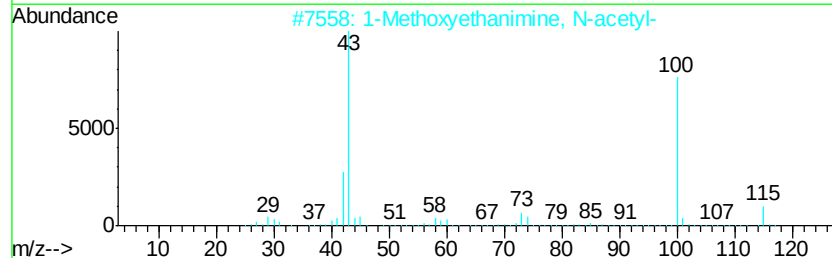
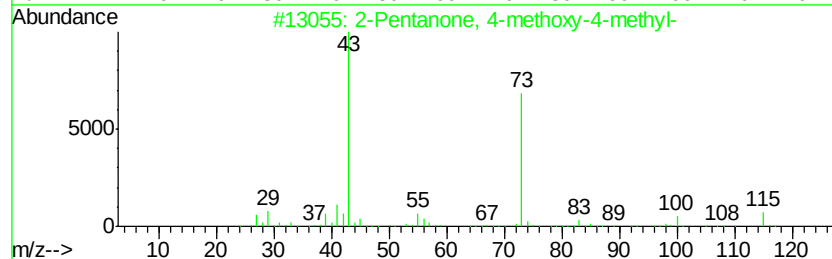
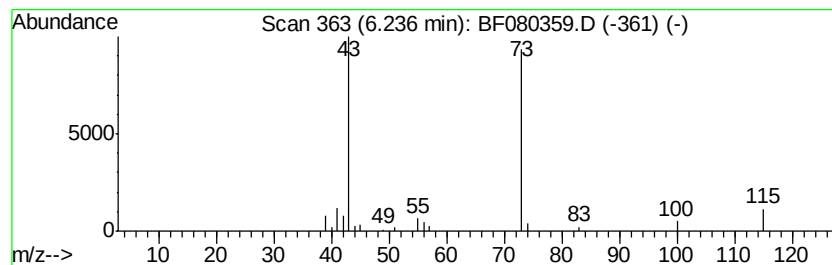
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-methoxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.24	3.78 ng	37008	1,4-Dichlorobenzene-d4	7.23	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
2	1-Methoxyethanimine, N-acetyl-	115	C5H9NO2	099028-43-0	9
3	Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	9
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
5	5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9



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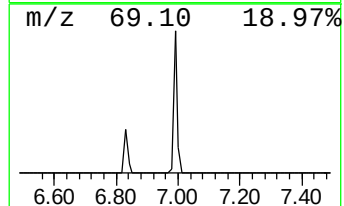
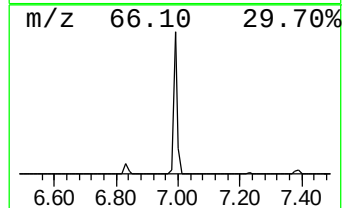
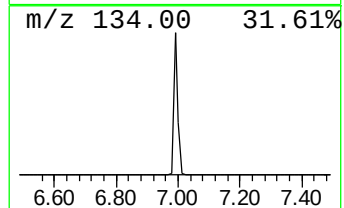
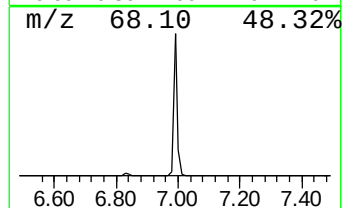
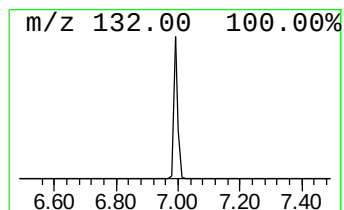
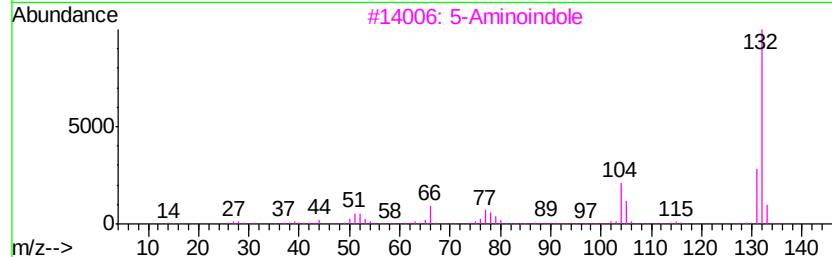
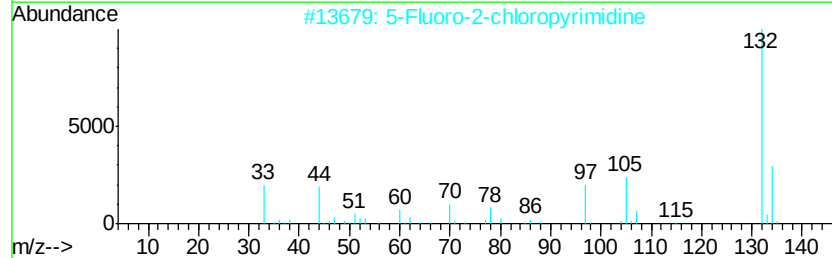
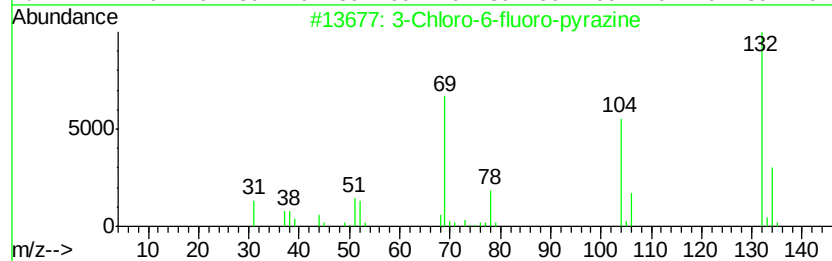
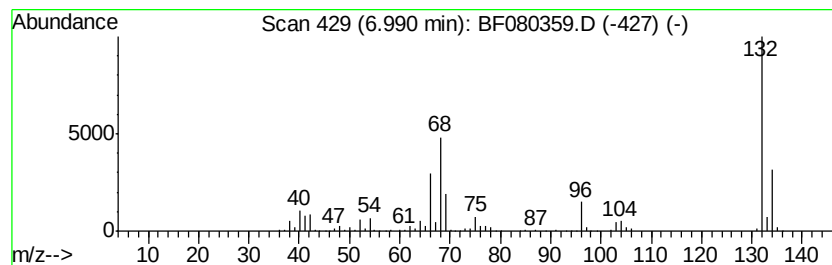
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.99 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.99	70.18 ng	687056	1,4-Dichlorobenzene-d4	7.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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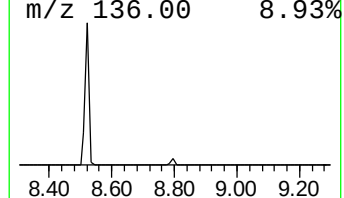
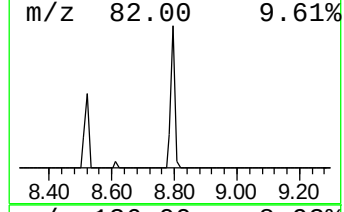
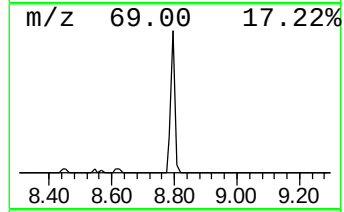
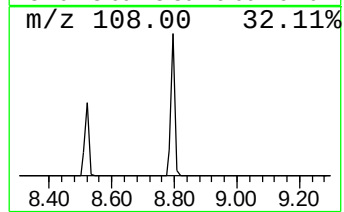
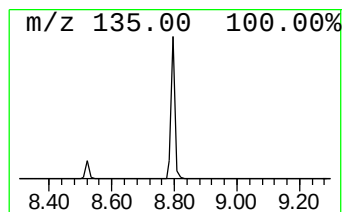
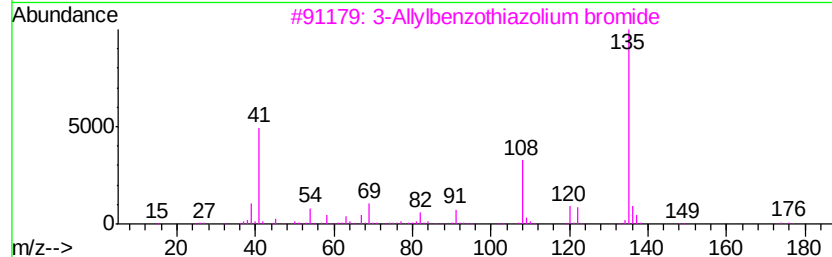
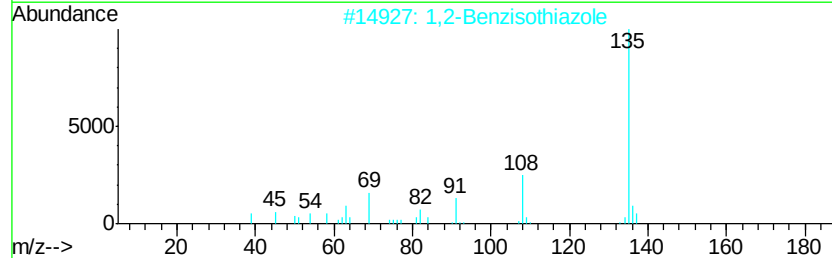
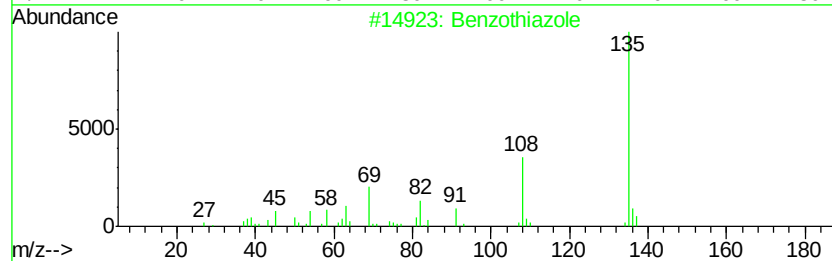
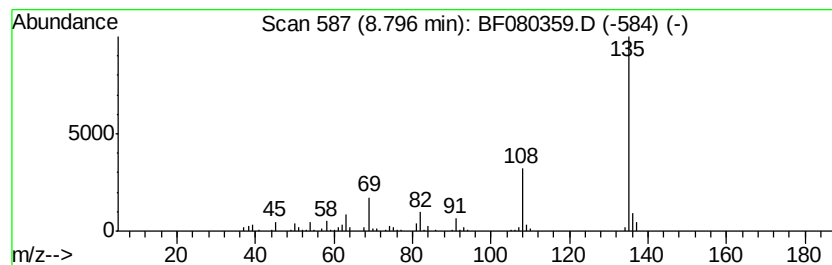
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzothiazole Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	13.84 ng	176286	Naphthalene-d8	8.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzothiazole	135	C7H5NS	000095-16-9	97
2			1,2-Benzisothiazole	135	C7H5NS	000272-16-2	91
3			3-Allylbenzothiazolium bromide	255	C10H10BrNS	016407-55-9	83
4			1H-Pyrazolo[3,4-d]pyrimidin-4-amine	135	C5H5N5	002380-63-4	59
5			4-Aminopyrazolo(3,4-d)pyrimidine	135	C5H5N5	020289-44-5	59



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Sample : G2877-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

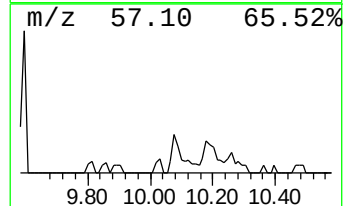
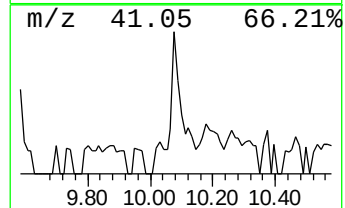
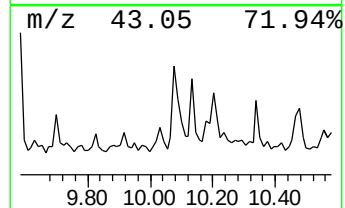
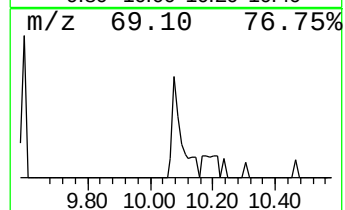
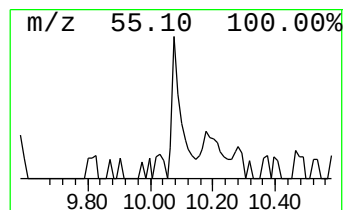
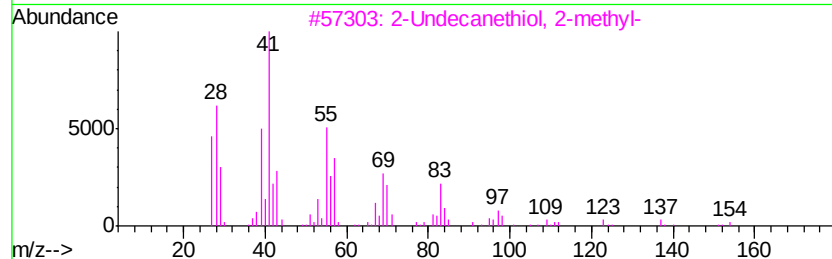
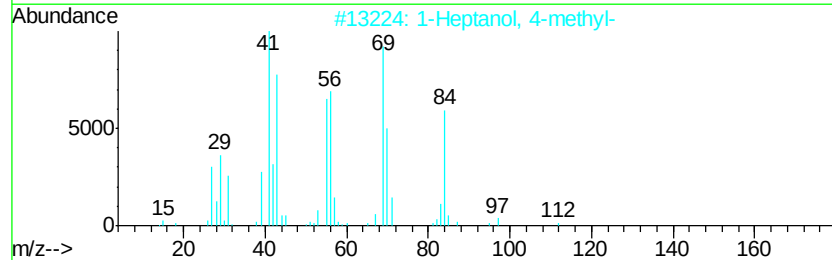
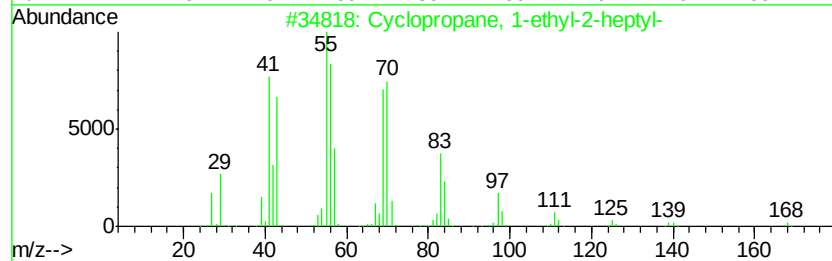
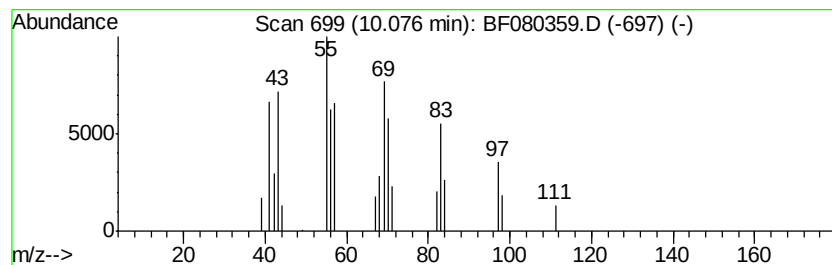
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopropane, 1-ethyl-2-hep... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.08	2.15 ng	31330	Acenaphthene-d10		10.29
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-ethyl-2-heptyl-	168	C12H24	074663-86-8	53
2	1-Heptanol, 4-methyl-	130	C8H18O	000817-91-4	50
3	2-Undecanethiol, 2-methyl-	202	C12H26S	010059-13-9	43
4	1-Hexene, 3-methyl-	98	C7H14	003404-61-3	38
5	Cyclopentane, butyl-	126	C9H18	002040-95-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
Data File : BF080359.D
Acq On : 6 Jul 2015 22:21
Operator : TP/IZ
Sample : G2877-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DCW-WW-155-7215

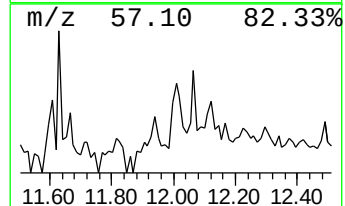
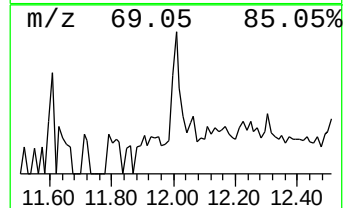
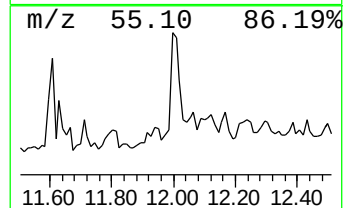
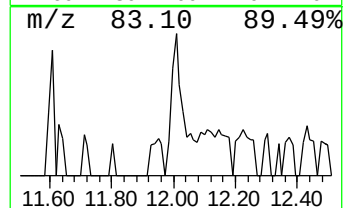
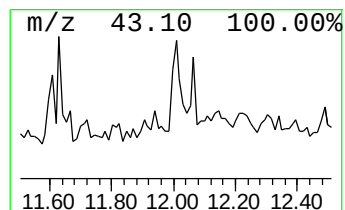
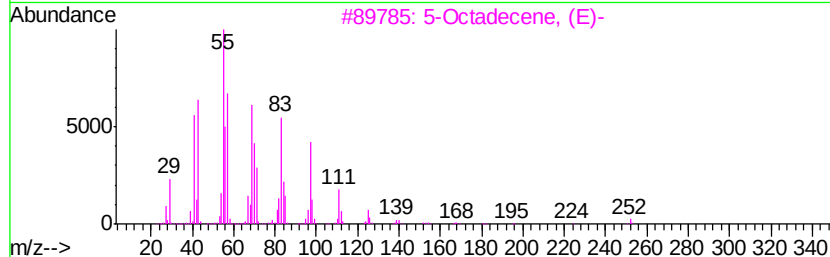
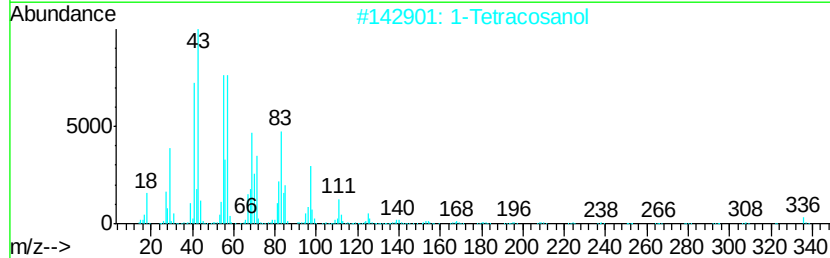
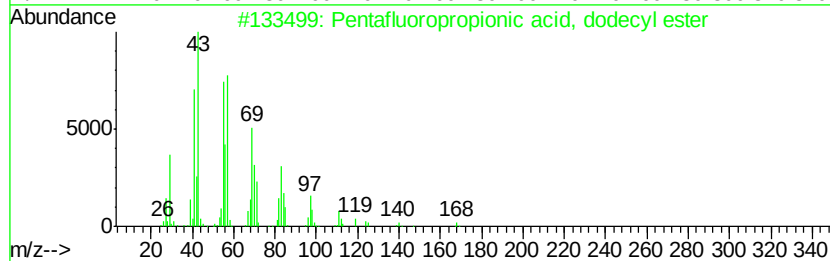
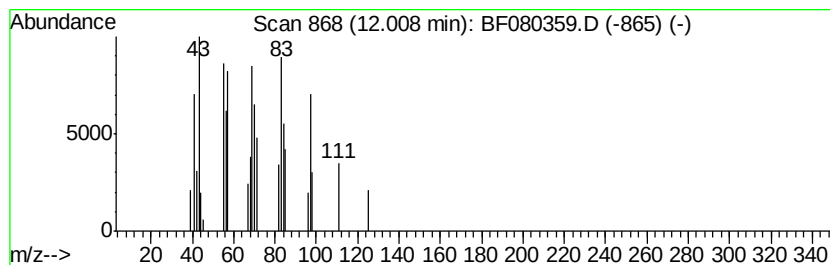
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentafluoropropionic acid, ... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	2.12 ng	35575	Phenanthrene-d10	11.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, dodec...	332	C15H25F5O2	006222-04-4	87
2		1-Tetracosanol	354	C24H50O	000506-51-4	78
3		5-Octadecene, (E)-	252	C18H36	007206-21-5	72
4		Heptafluorobutyric acid,n-tridec...	396	C17H27F7O2	1000216-78-9	64
5		Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
Data File : BF080359.D
Acq On : 6 Jul 2015 22:21
Operator : TP/IZ
Sample : G2877-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

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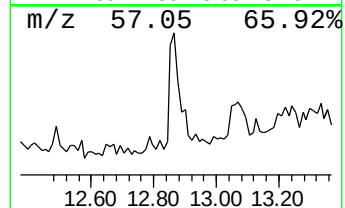
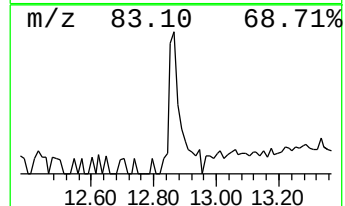
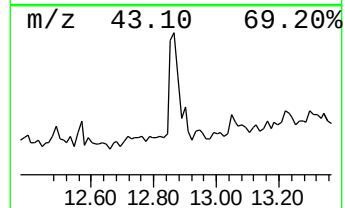
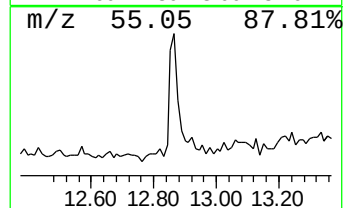
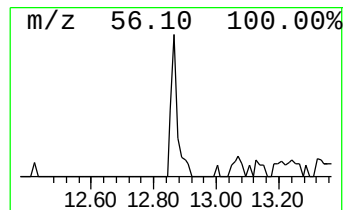
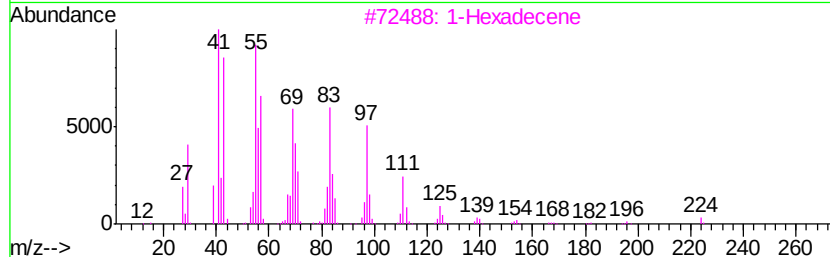
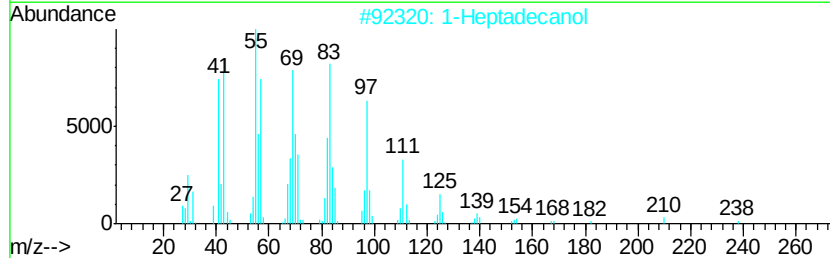
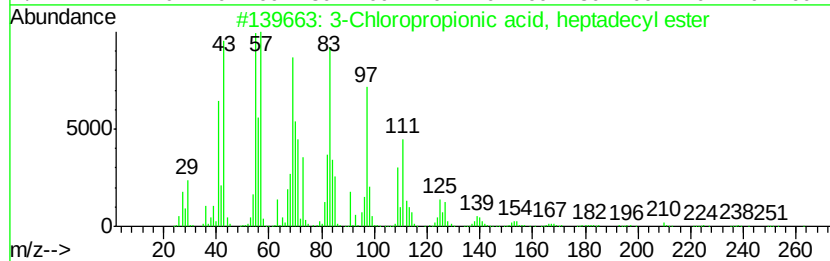
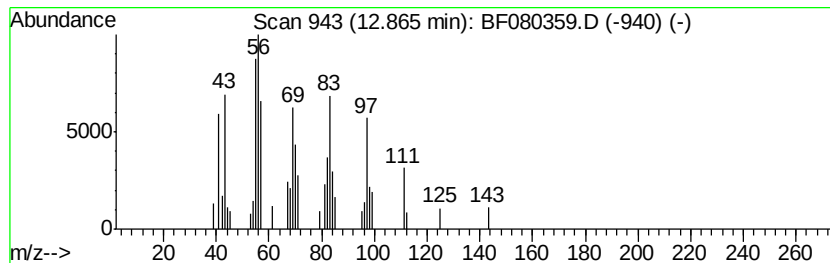
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 3-Chloropropionic acid, hep... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	4.67 ng	78192	Phenanthrene-d10	11.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	64
2			1-Heptadecanol	256	C17H36O	001454-85-9	60
3			1-Hexadecene	224	C16H32	000629-73-2	58
4			1-Tridecene	182	C13H26	002437-56-1	58
5			9-Eicosene, (E)-	280	C20H40	074685-29-3	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
Data File : BF080359.D
Acq On : 6 Jul 2015 22:21
Operator : TP/IZ
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Misc :
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Instrument :
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ClientSampleId :
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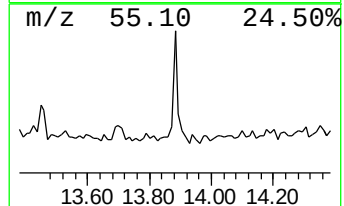
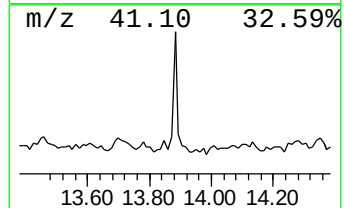
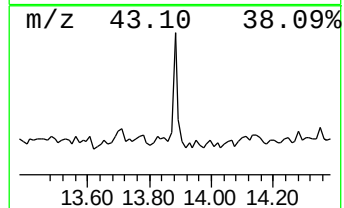
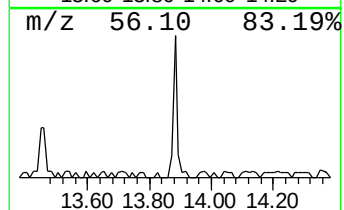
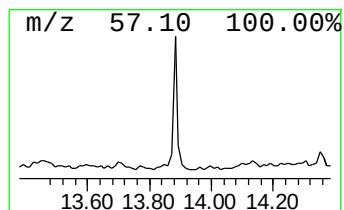
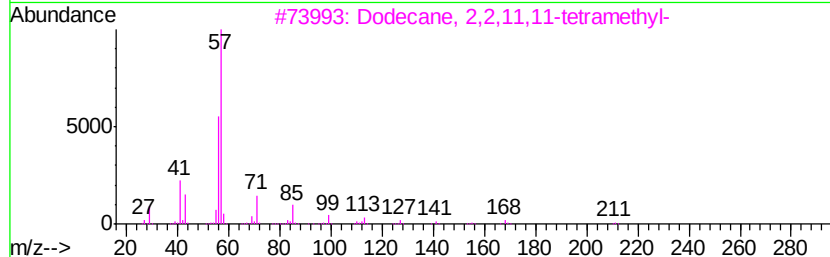
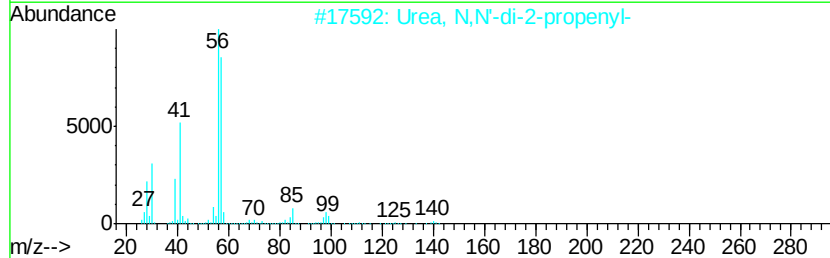
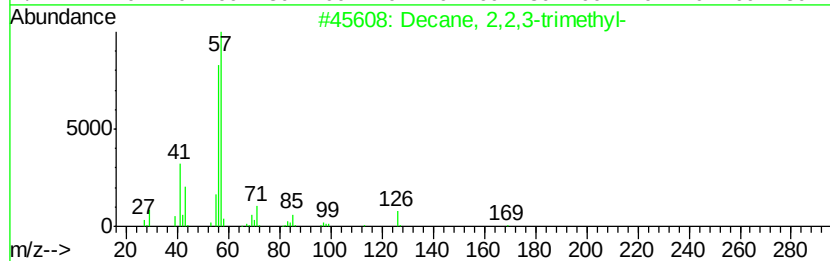
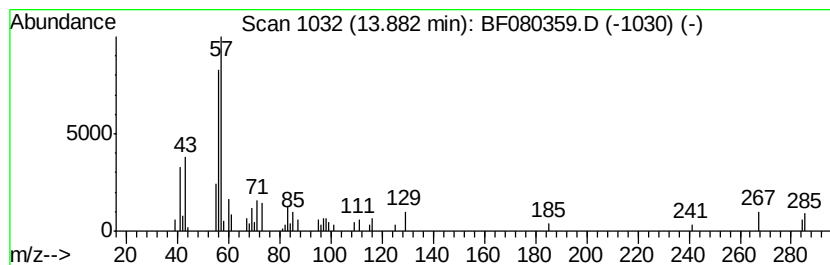
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Decane, 2,2,3-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	2.76 ng	49675	Chrysene-d12	14.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	50
2			Urea, N,N'-di-2-propenyl-	140	C7H12N2O	001801-72-5	50
3			Dodecane, 2,2,11,11-tetramethyl-	226	C16H34	127204-12-0	50
4			1-Hexene, 4-methyl-	98	C7H14	003769-23-1	47
5			Hexanoic acid, butyl ester	172	C10H20O2	000626-82-4	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
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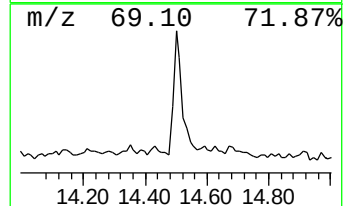
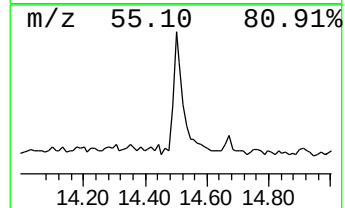
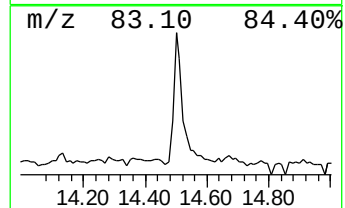
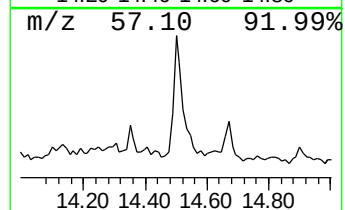
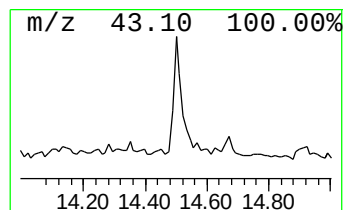
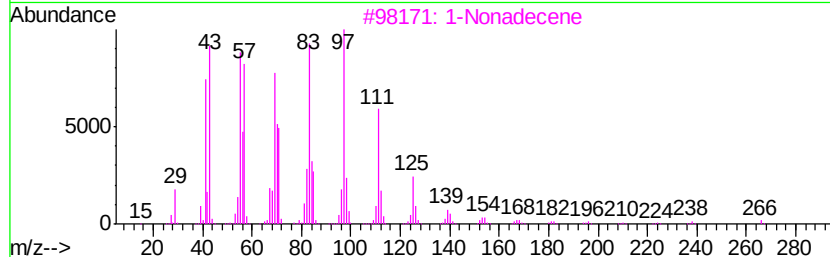
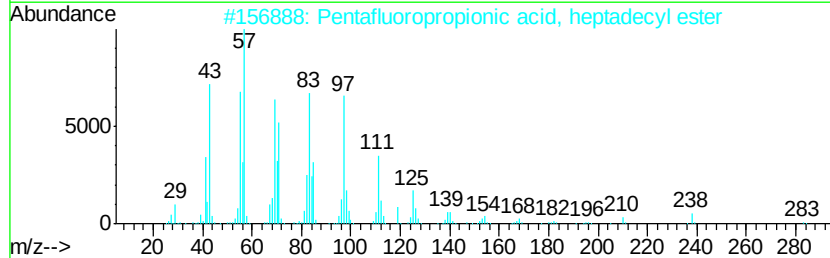
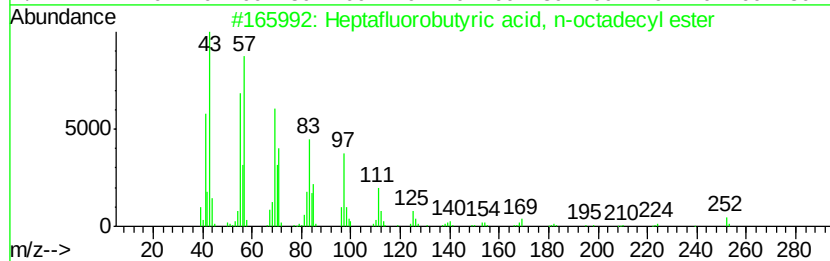
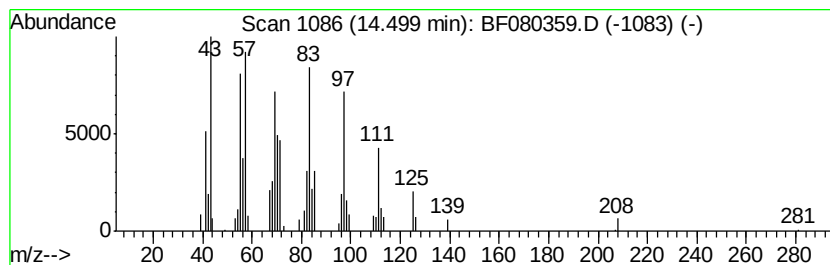
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptafluorobutyric acid, n-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	6.47 ng	116304	Chrysene-d12	14.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91
2			Pentafluoropropionic acid, heptadecyl ester	402	C20H35F5O2	1000283-04-2	90
3			1-Nonadecene	266	C19H38	018435-45-5	90
4			1-Octadecanol	270	C18H38O	000112-92-5	87
5			2-Chloropropionic acid, octadecyl ester	360	C21H41ClO2	088104-31-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
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ClientSampleId :
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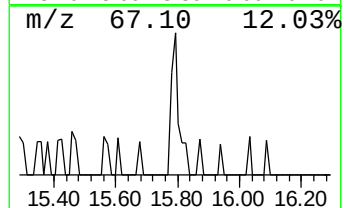
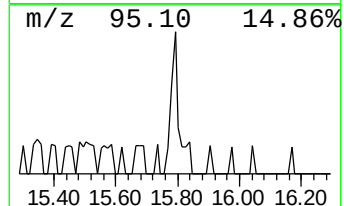
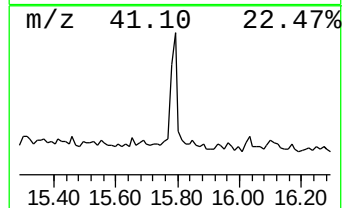
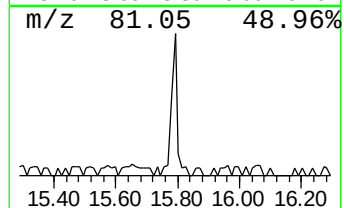
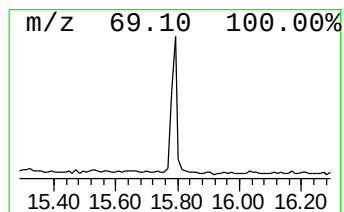
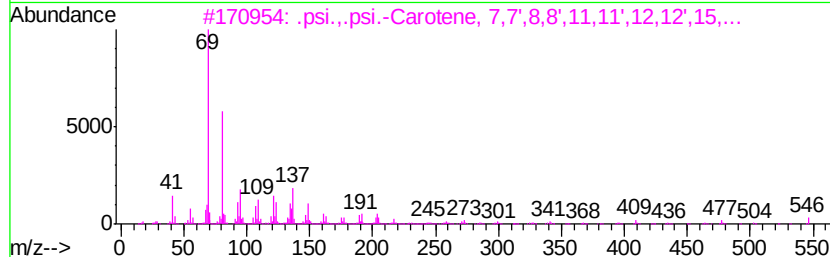
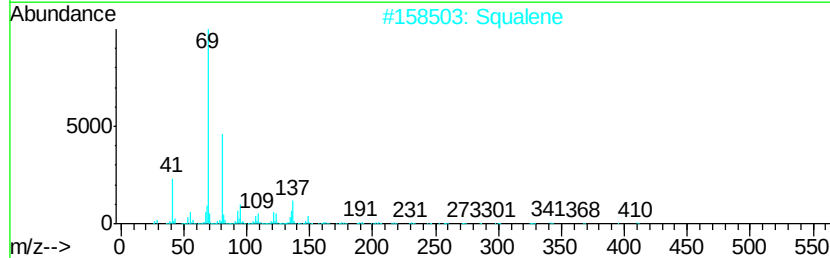
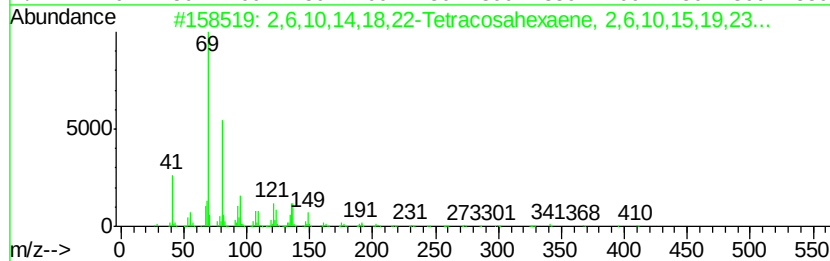
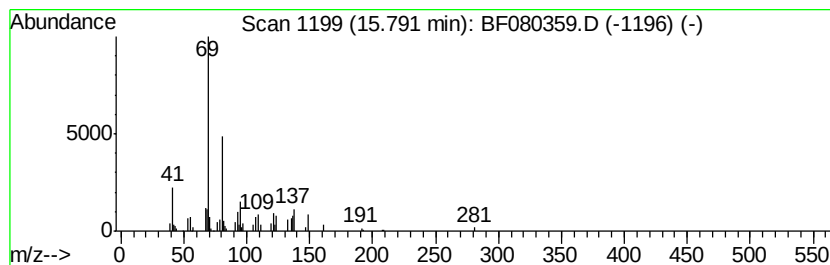
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2,6,10,14,18,22-Tetracosahexaene... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	3.32 ng	57982	Perylene-d12	16.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	91		
2	Squalene	410	C30H50	007683-64-9	90		
3	.psi.,.psi.-Carotene, 7,7',8,8',...	547	C40H66	000502-62-5	83		
4	4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	78		
5	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	78		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
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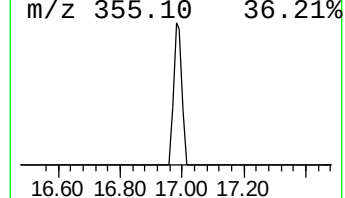
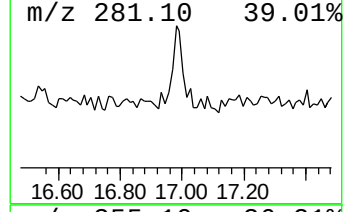
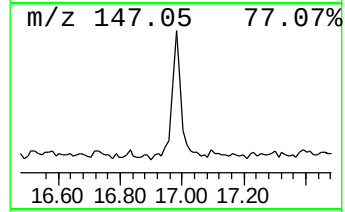
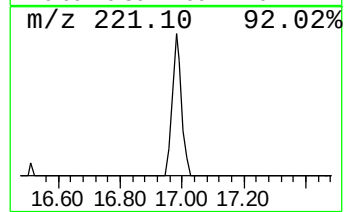
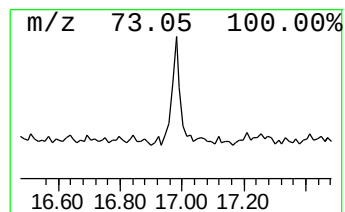
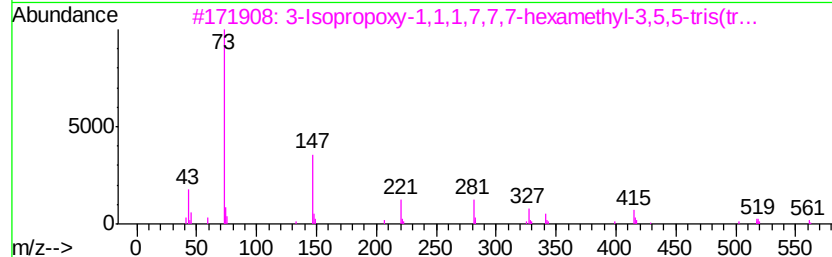
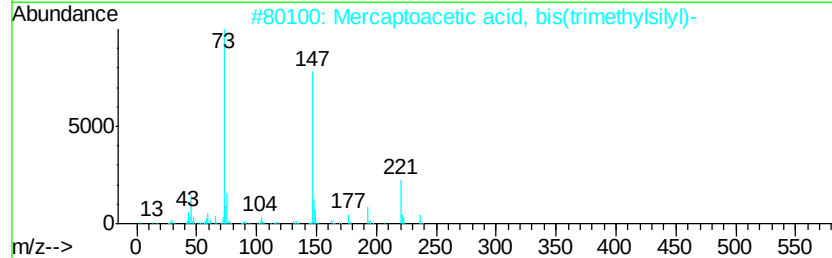
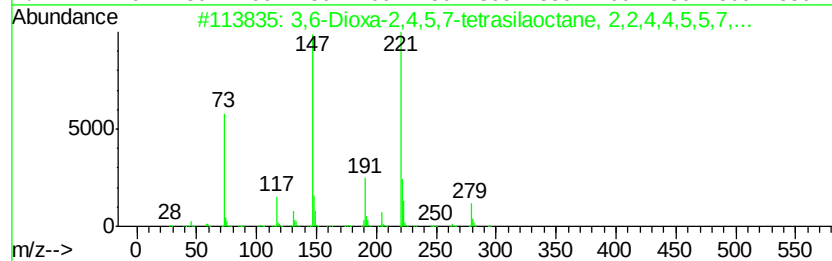
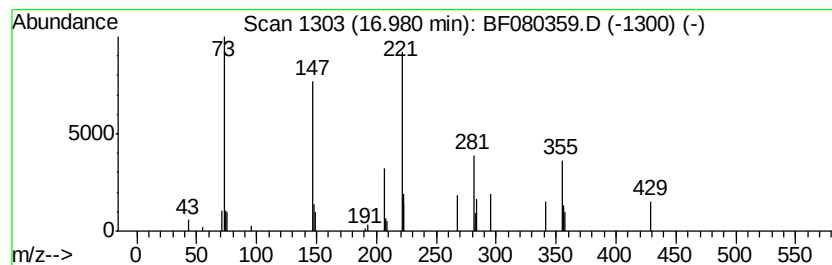
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.98 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.98	2.25 ng	39258	Perylene-d12	16.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	47
2		Mercaptoacetic acid, bis(trimeth...	236	C8H20O2SSi2	006398-62-5	38
3		3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	32
4		Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	27
5		N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
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Acq On : 6 Jul 2015 22:21
Operator : TP/IZ
Sample : G2877-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DCW-WW-155-7215

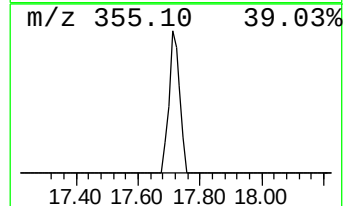
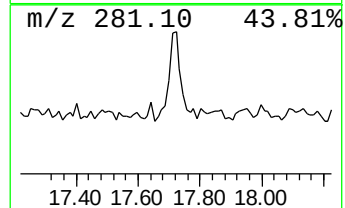
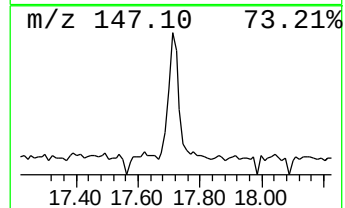
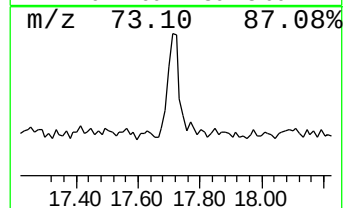
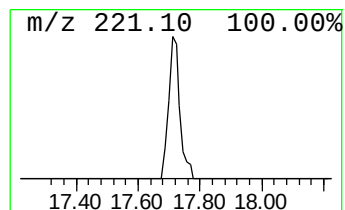
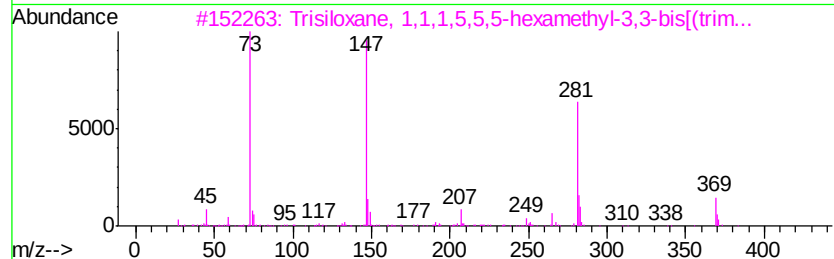
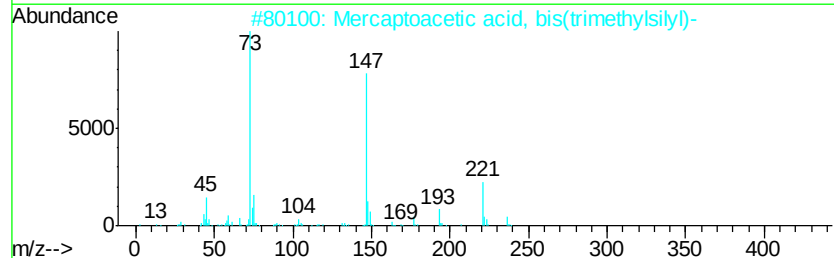
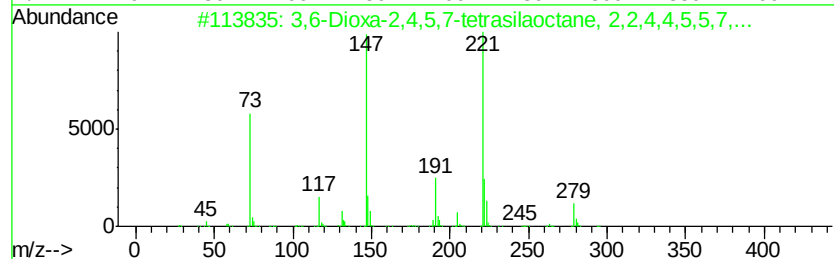
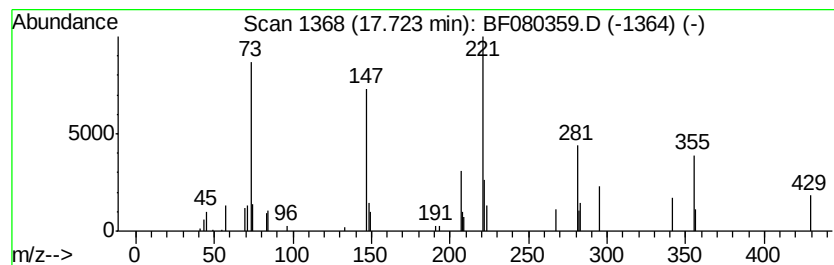
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown17.72 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.72	3.09 ng	54092	Perylene-d12	16.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
2		Mercaptoacetic acid, bis(trimeth...	236	C8H20O2SSi2	006398-62-5	32
3		Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	27
4		Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	27
5		N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	16



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
Data File : BF080359.D
Acq On : 6 Jul 2015 22:21
Operator : TP/IZ
Sample : G2877-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DCW-WW-155-7215

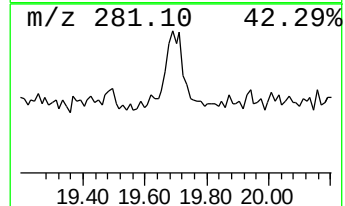
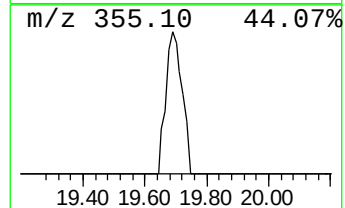
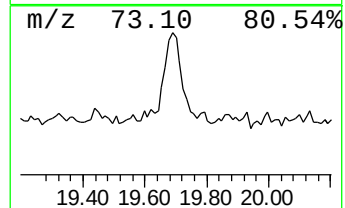
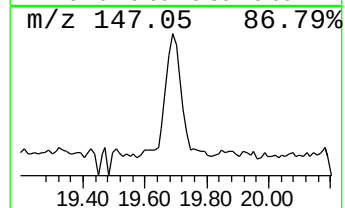
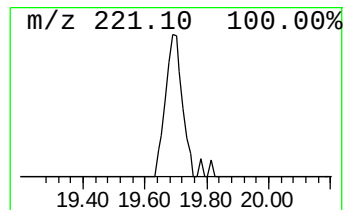
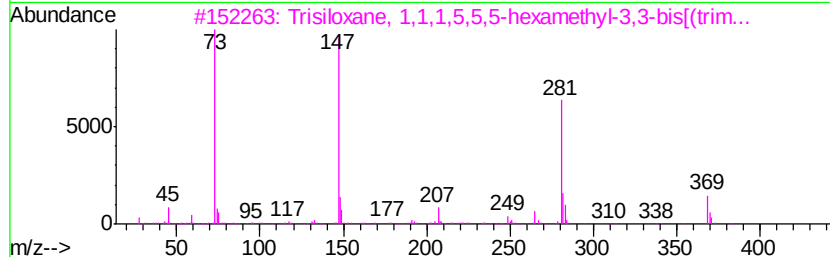
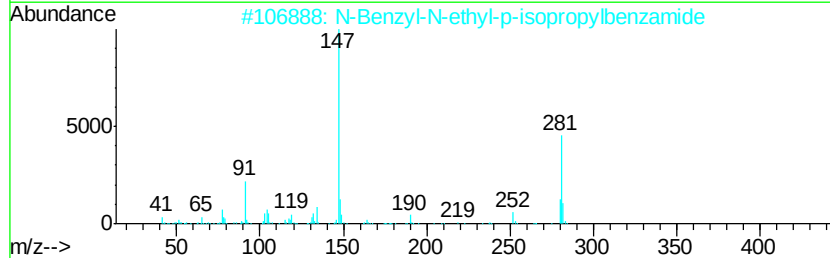
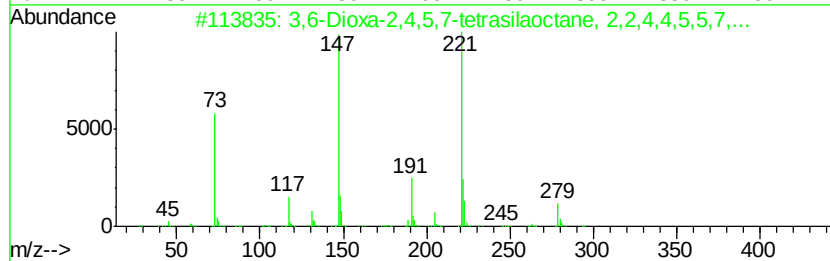
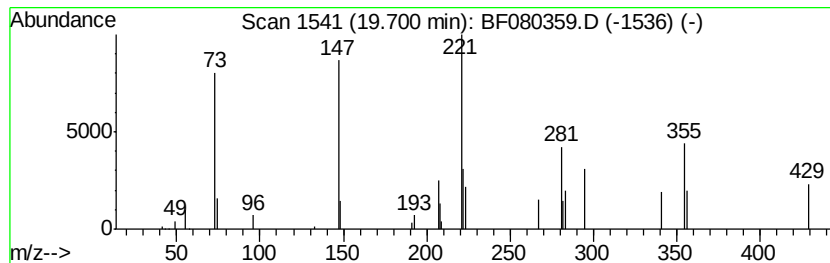
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown19.70 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.70	2.89 ng	50467	Perylene-d12	16.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	38
2		N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	22
3		Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	10
4		Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	10
5		Piperidine, 1-(5-trifluoromethyl...	295	C15H16F3N3	1000268-74-7	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
Data File : BF080359.D
Acq On : 6 Jul 2015 22:21
Operator : TP/IZ
Sample : G2877-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DCW-WW-155-7215

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.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...	2.42	91.5	ng	895894	1	7.23	195804	20.0
2-Pentanone, 4-hy...	5.44	15.0	ng	147171	1	7.23	195804	20.0
2-Pentanone, 4-me...	6.24	3.8	ng	37008	1	7.23	195804	20.0
unknown6.99	6.99	70.2	ng	687056	1	7.23	195804	20.0
Benzothiazole	8.80	13.8	ng	176286	2	8.52	254685	20.0
Cyclopropane, 1-e...	10.08	2.1	ng	31330	3	10.29	291438	20.0
Pentafluoropropio...	12.01	2.1	ng	35575	4	11.79	335060	20.0
3-Chloropropionic...	12.87	4.7	ng	78192	4	11.79	335060	20.0
Decane, 2,2,3-tri...	13.88	2.8	ng	49675	5	14.71	359630	20.0
Heptafluorobutyri...	14.50	6.5	ng	116304	5	14.71	359630	20.0
2,6,10,14,18,22-T...	15.79	3.3	ng	57982	6	16.51	349566	20.0
unknown16.98	16.98	2.3	ng	39258	6	16.51	349566	20.0
unknown17.72	17.72	3.1	ng	54092	6	16.51	349566	20.0
unknown19.70	19.70	2.9	ng	50467	6	16.51	349566	20.0