

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078506.D
 Acq On : 14 Apr 2015 18:33
 Operator : TP/IZ
 Sample : G1828-02 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB05B

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-BF040115.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	1.495	26	27	34	rVB2	10068	22830	1.30%	0.360%
2	1.598	34	36	63	rVB	693242	1757057	100.00%	27.732%
3	5.062	337	339	345	rBV	64261	83429	4.75%	1.317%
4	6.410	455	457	464	rBV	86009	96730	5.51%	1.527%
5	6.525	464	467	469	rBV	73056	96210	5.48%	1.518%
6	6.799	489	491	494	rBV	111479	112667	6.41%	1.778%
7	6.982	505	507	510	rBV	54576	66347	3.78%	1.047%
8	7.508	551	553	555	rBV	57850	60236	3.43%	0.951%
9	8.376	627	629	632	rBV	142281	173332	9.86%	2.736%
10	8.810	665	667	670	rBV	13130	18064	1.03%	0.285%
11	9.725	746	747	750	rVB	113416	126597	7.21%	1.998%
12	10.365	801	803	806	rBV	18059	26902	1.53%	0.425%
13	10.434	806	809	811	rVB2	15830	27221	1.55%	0.430%
14	10.536	816	818	821	rVV	215436	223501	12.72%	3.528%
15	10.856	843	846	847	rBV2	12113	18347	1.04%	0.290%
16	10.948	850	854	856	rVV3	9859	19963	1.14%	0.315%
17	11.439	893	897	899	rBV2	14905	24405	1.39%	0.385%
18	11.508	901	903	906	rBV2	75029	86873	4.94%	1.371%
19	11.748	921	924	930	rBV	13999	31020	1.77%	0.490%
20	12.365	976	978	979	rBV	212154	243236	13.84%	3.839%
21	12.388	979	980	983	rVV	61593	55840	3.18%	0.881%
22	12.457	983	986	988	rVV	15729	18943	1.08%	0.299%
23	12.559	993	995	997	rBV2	8875	17744	1.01%	0.280%
24	12.674	1003	1005	1008	rBV3	11019	24507	1.39%	0.387%
25	12.982	1030	1032	1034	rBV	13957	31021	1.77%	0.490%
26	13.017	1034	1035	1039	rVB2	20407	29790	1.70%	0.470%
27	13.119	1042	1044	1045	rBV	24451	34033	1.94%	0.537%
28	13.851	1106	1108	1110	rBV	163393	175970	10.02%	2.777%
29	13.908	1110	1113	1115	rBV2	25231	41880	2.38%	0.661%
30	14.125	1130	1132	1134	rVB	191440	197092	11.22%	3.111%
31	14.240	1140	1142	1144	rBV	83348	110231	6.27%	1.740%
32	14.354	1149	1152	1154	rBV	180137	202400	11.52%	3.194%
33	15.623	1261	1263	1265	rBV2	279170	399506	22.74%	6.305%
34	17.371	1414	1416	1418	rBV	175466	275542	15.68%	4.349%

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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	18.023	1471	1473	1481	rVB	322065	898037	51.11%	14.174%
36	18.377	1502	1504	1507	rVB	352905	508395	28.93%	8.024%

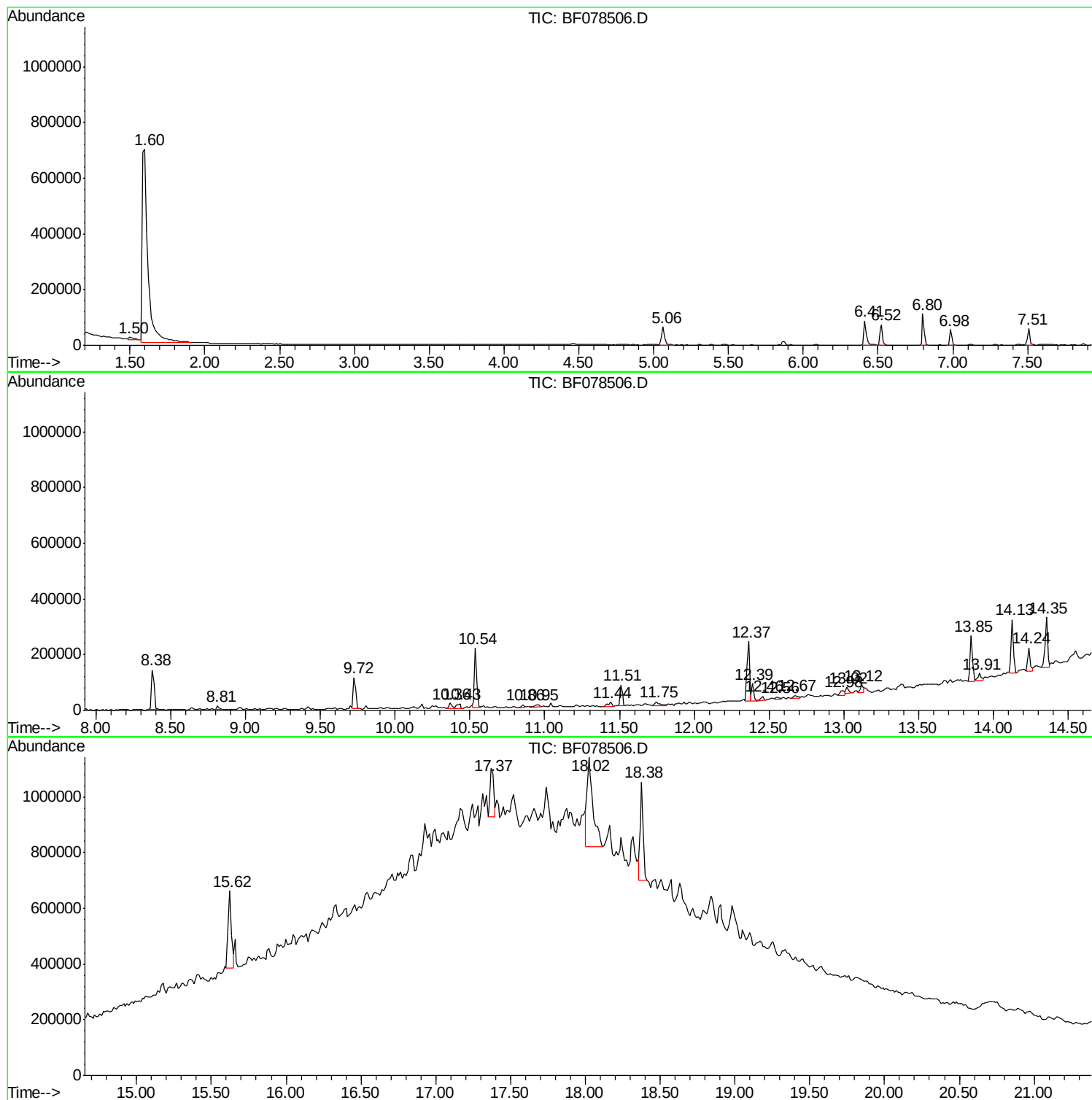
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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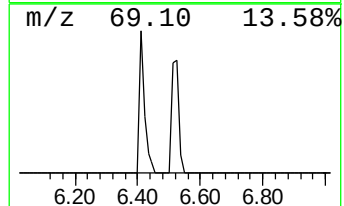
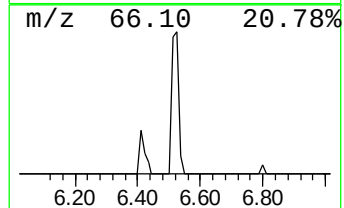
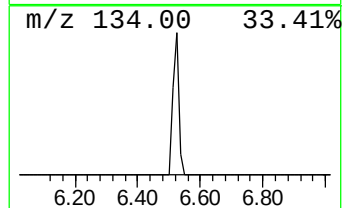
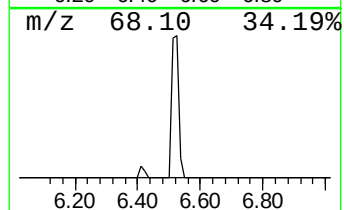
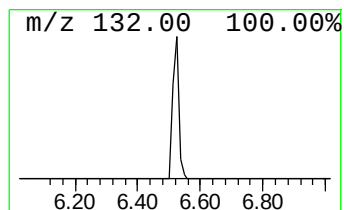
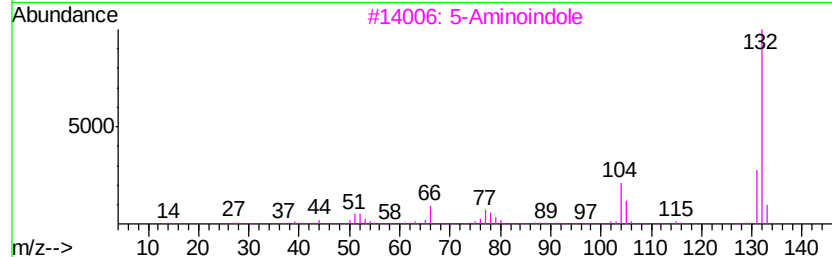
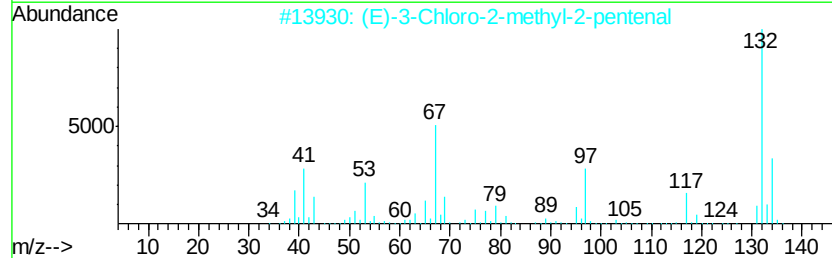
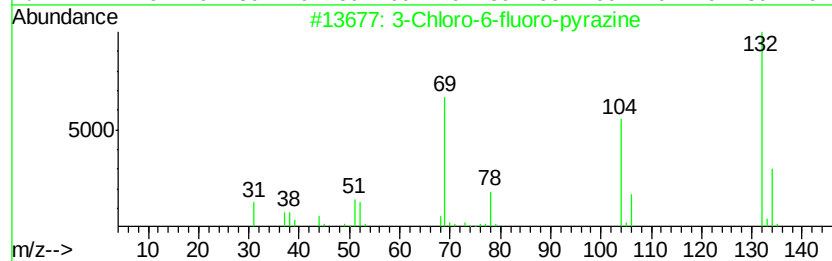
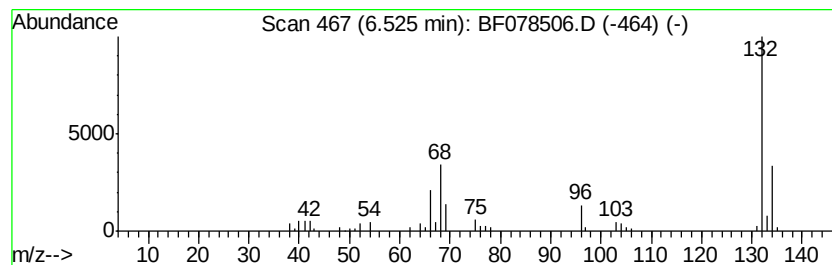
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.52 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	17.08 ng	96210	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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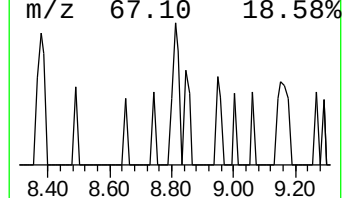
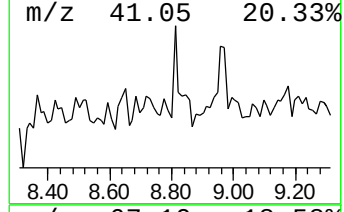
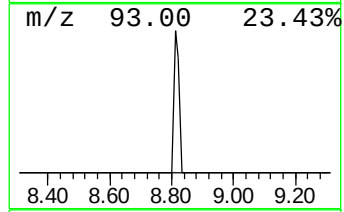
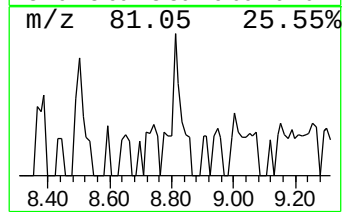
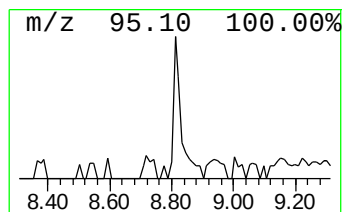
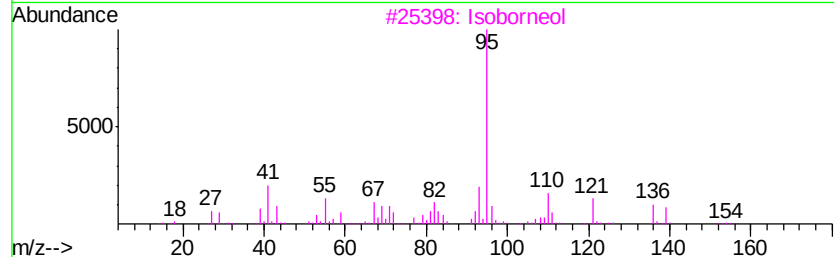
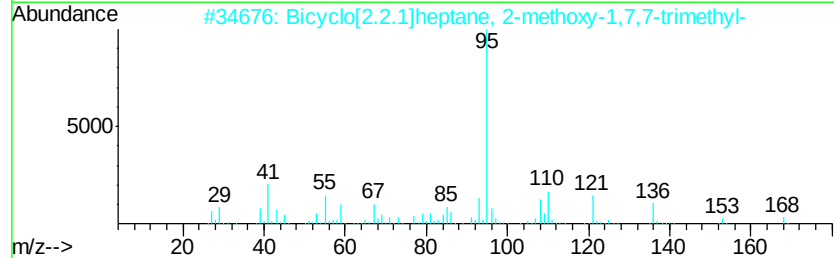
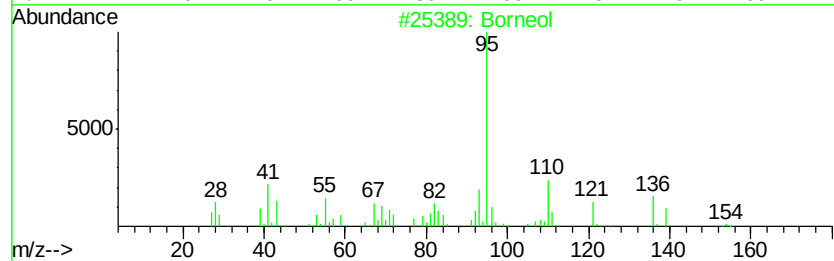
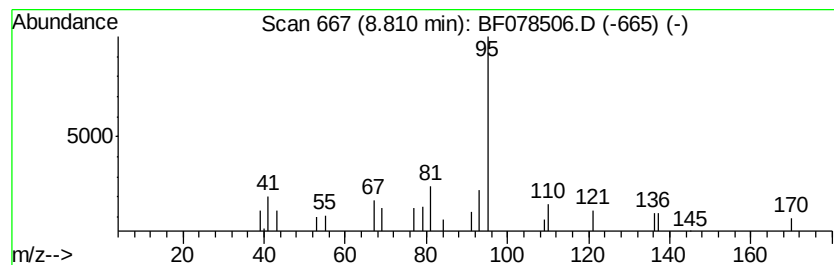
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Borneol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.81	2.08 ng	18064	Naphthalene-d8	8.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Borneol	154	C10H18O	010385-78-1	59
2		Bicyclo[2.2.1]heptane, 2-methoxy...	168	C11H20O	004443-51-0	59
3		Isoborneol	154	C10H18O	000124-76-5	45
4		5-Methyl-5-hexen-3-yn-2-ol	110	C7H10O	068017-33-4	43
5		1,3-Hexadiene, 2,5-dimethyl-	110	C8H14	029253-64-3	38



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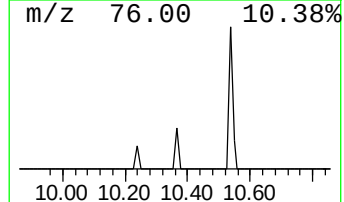
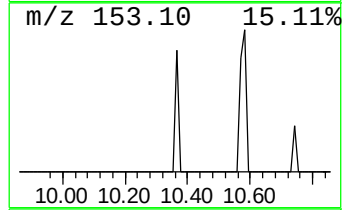
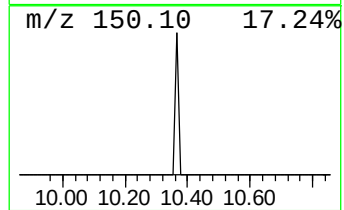
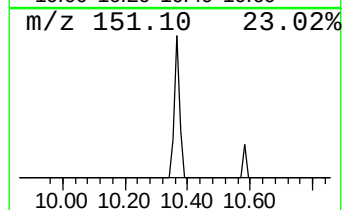
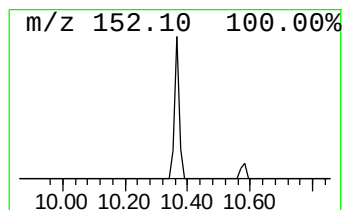
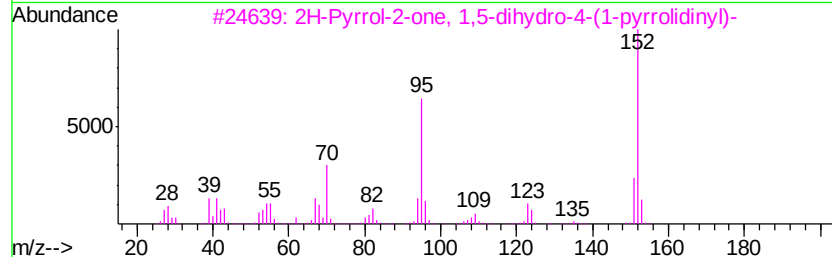
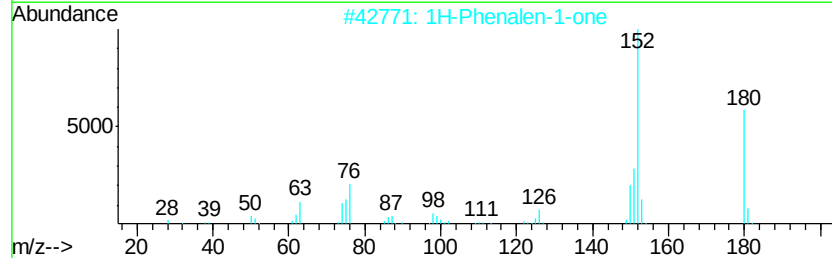
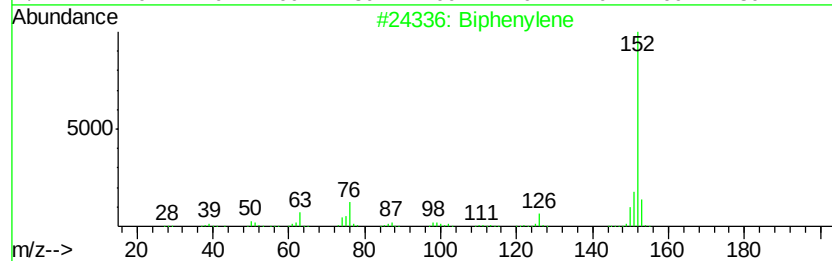
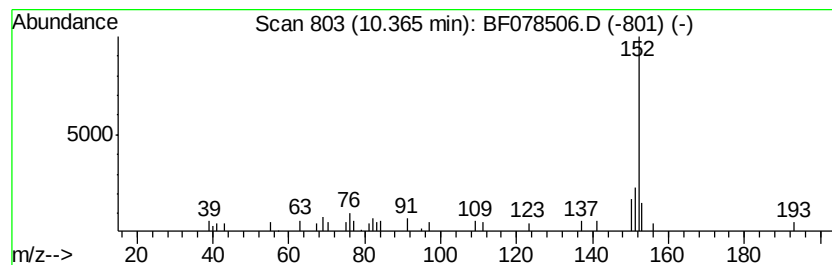
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Biphenylene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.36	2.41 ng	26902	Acenaphthene-d10	10.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenylene	152	C12H8	000259-79-0	72
2		1H-Phenalen-1-one	180	C13H8O	000548-39-0	64
3		2H-Pyrrol-2-one, 1,5-dihydro-4-(...	152	C8H12N2O	105776-93-0	64
4		Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	64
5		Acenaphthylene	152	C12H8	000208-96-8	59



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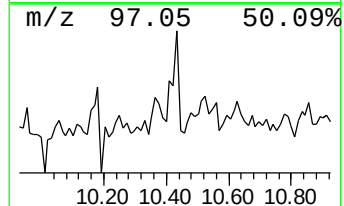
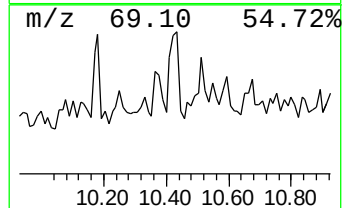
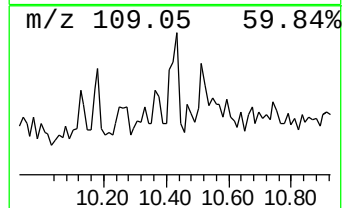
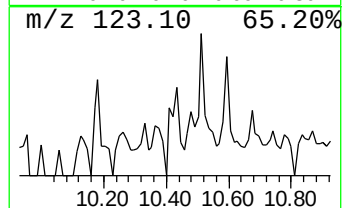
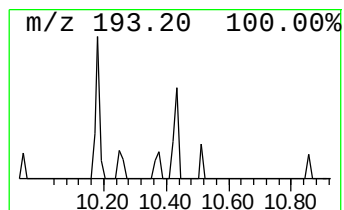
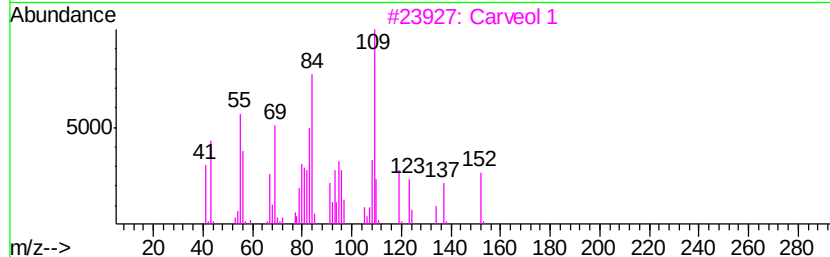
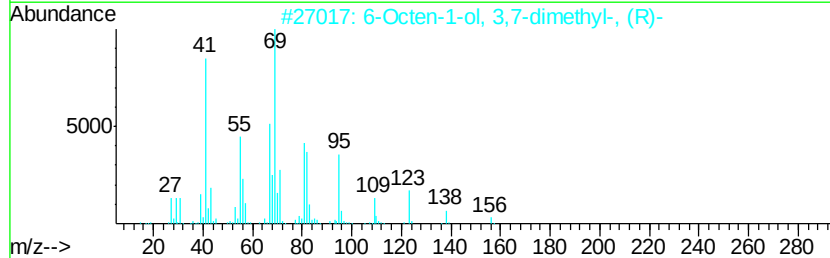
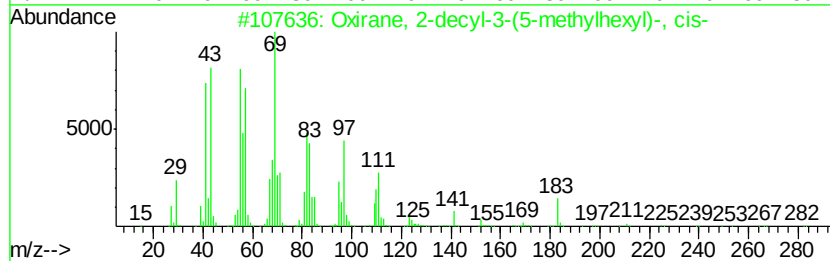
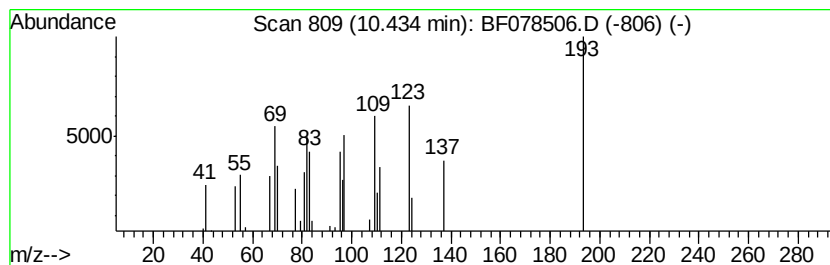
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown10.43 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	2.44 ng	27221	Acenaphthene-d10	10.54
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Oxirane, 2-decyl-3-(5-methylhexy...	282 C19H380	029804-22-6 35
2		6-Octen-1-ol, 3,7-dimethyl-, (R)-	156 C10H200	001117-61-9 17
3		Carveol 1	152 C10H160	1000285-09-9 12
4		3,3,5,5-Tetramethylcyclohexanol	156 C10H200	002650-40-0 12
5		1-Cyclohexene-1-carboxaldehyde, ...	152 C10H160	021391-98-0 10



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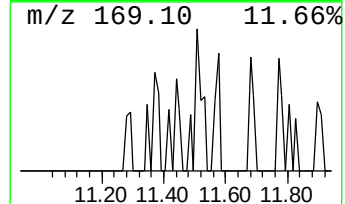
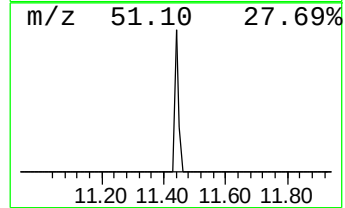
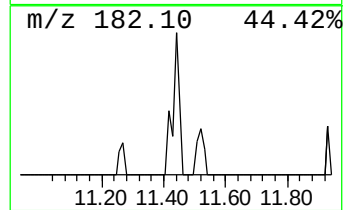
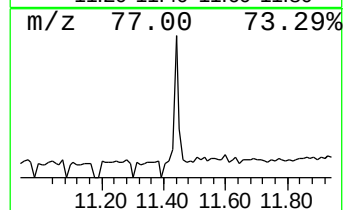
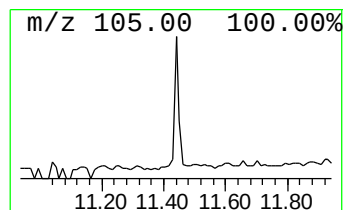
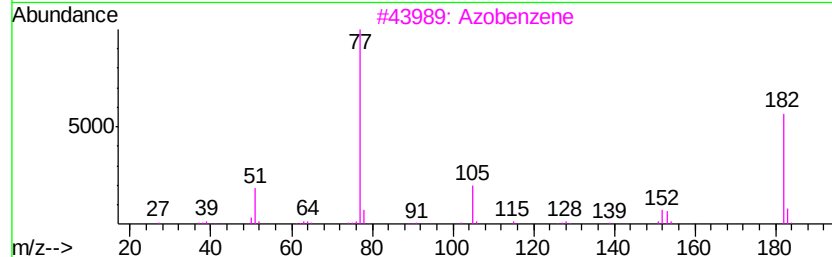
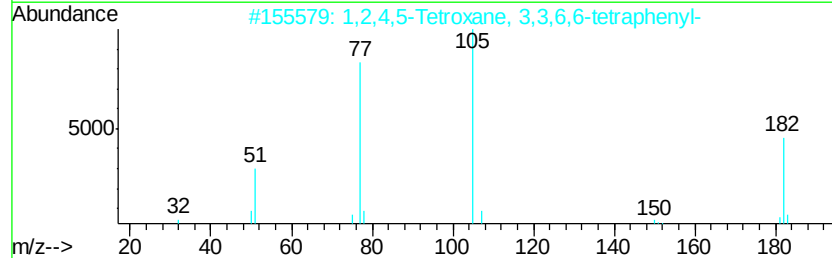
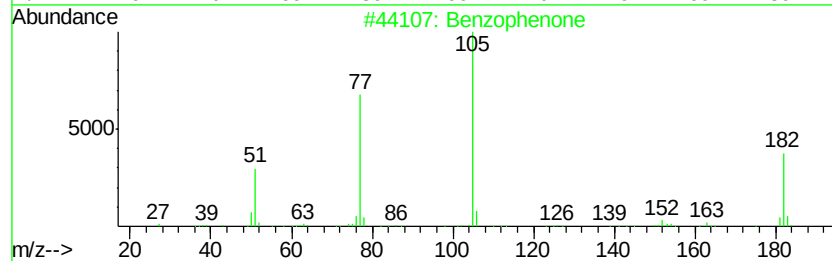
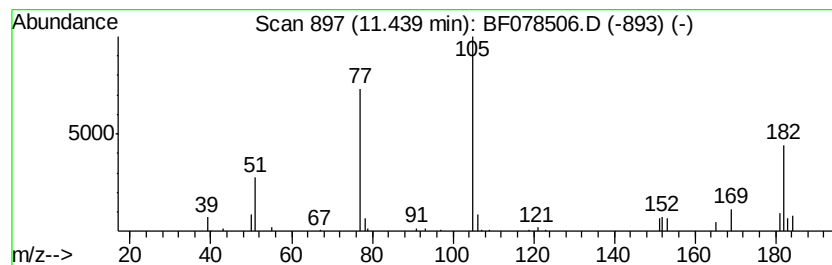
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 Benzophenone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	2.18 ng	24405	Acenaphthene-d10	10.54
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzophenone	182 C13H10O	000119-61-9 83
2		1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396 C26H20O4	016204-36-7 72
3		Azobenzene	182 C12H10N2	000103-33-3 59
4		Azobenzene, (E)-	182 C12H10N2	017082-12-1 59
5		Benzilamide	227 C14H13NO2	004746-87-6 50



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ALS Vial : 4 Sample Multiplier: 1

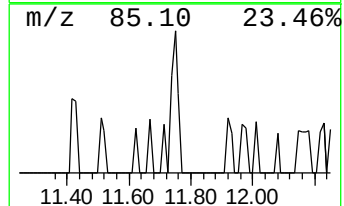
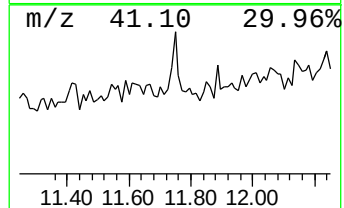
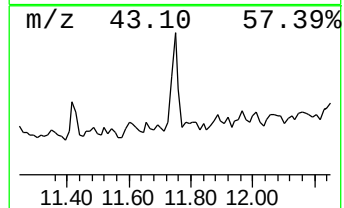
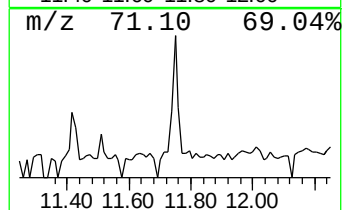
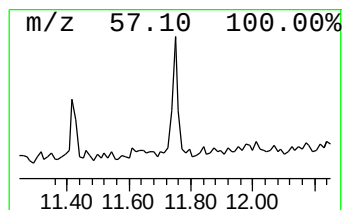
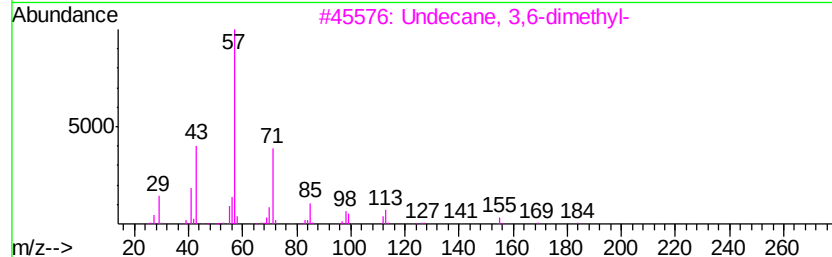
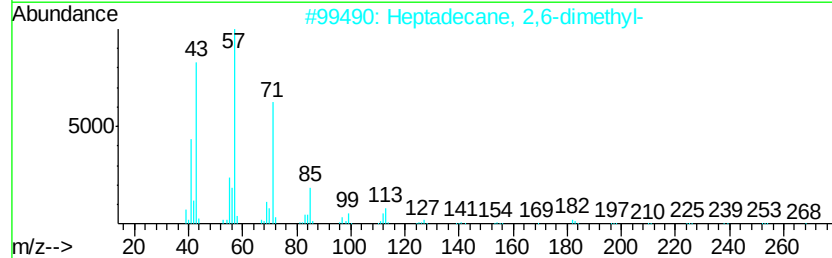
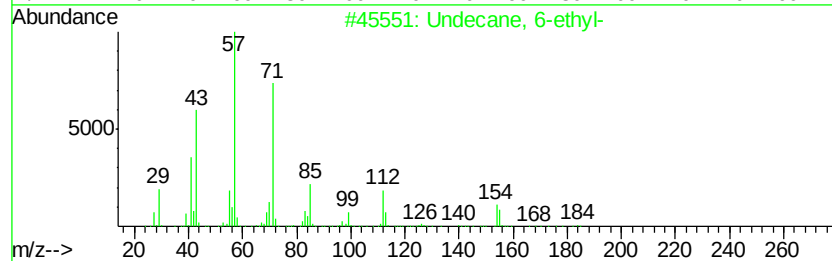
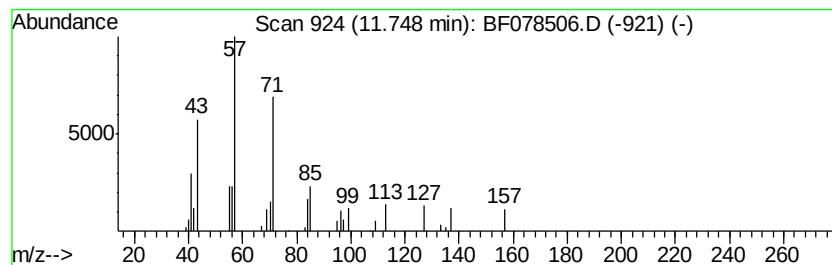
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ClientSampleId :
HS-SB05B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Undecane, 6-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.75	2.55 ng	31020	Phenanthrene-d10		12.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 6-ethyl-	184	C13H28	017312-60-6	53
2	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	53
3	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	53
4	Hexadecane, 7-methyl-	240	C17H36	026730-20-1	53
5	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
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Acq On : 14 Apr 2015 18:33
Operator : TP/IZ
Sample : G1828-02 5X
Misc :
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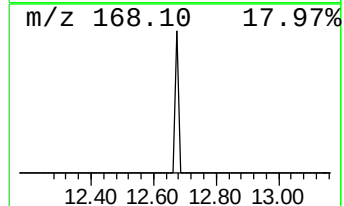
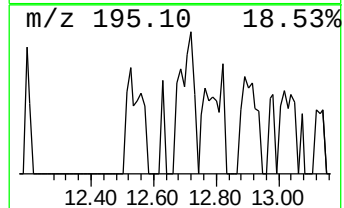
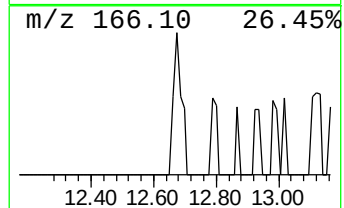
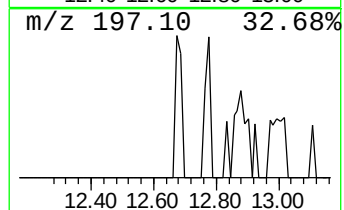
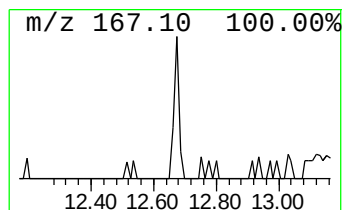
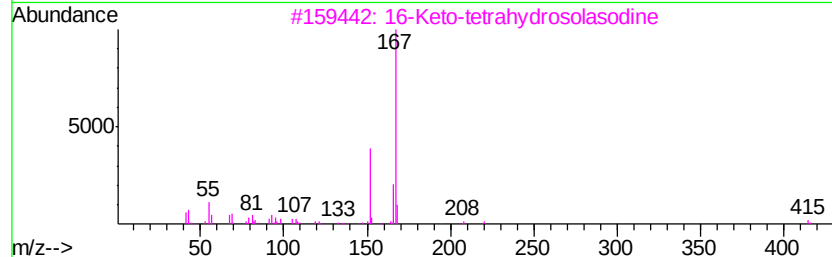
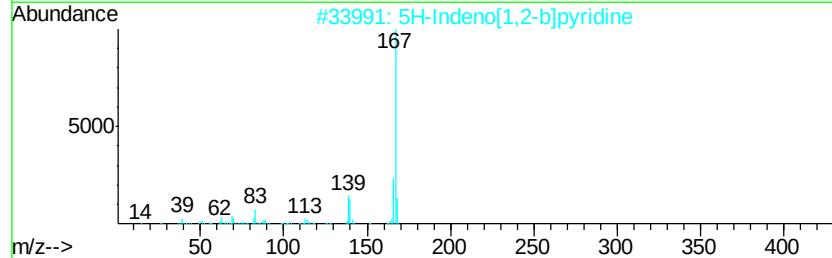
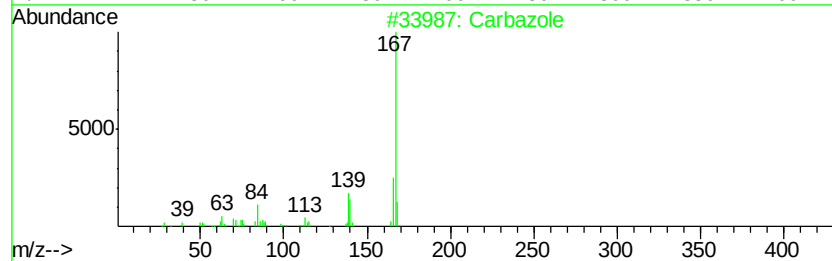
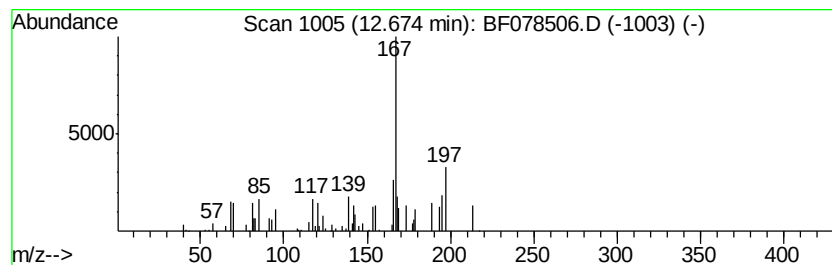
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown12.67 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.67	2.02 ng	24507	Phenanthrene-d10		12.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbazole	167	C12H9N	000086-74-8	43
2	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	38
3	16-Keto-tetrahydrosolasodine	415	C27H45N02	1000255-34-4	38
4	1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	37
5	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078506.D
Acq On : 14 Apr 2015 18:33
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Sample : G1828-02 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

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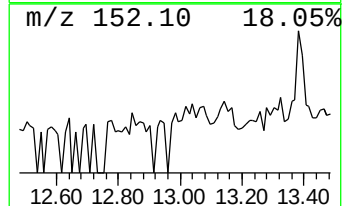
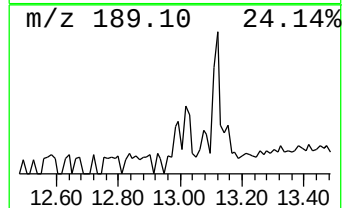
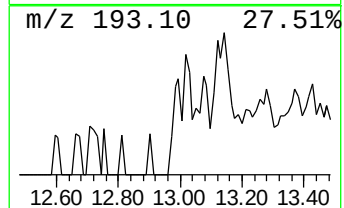
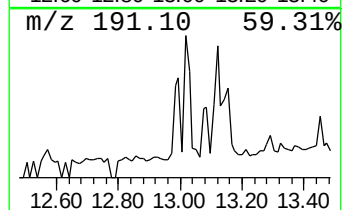
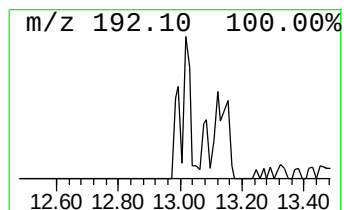
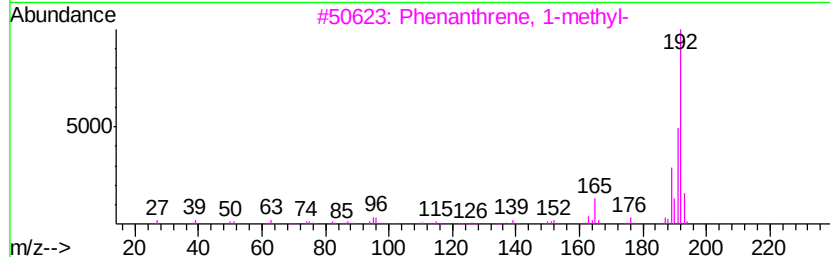
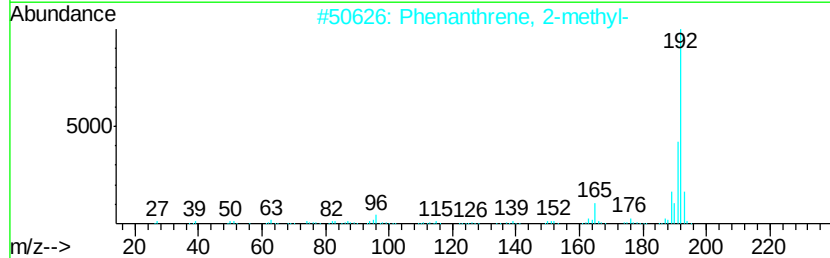
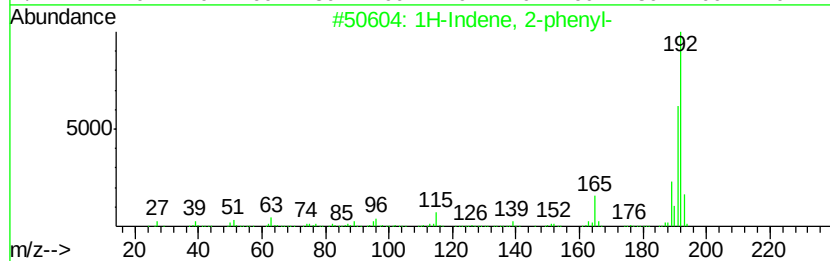
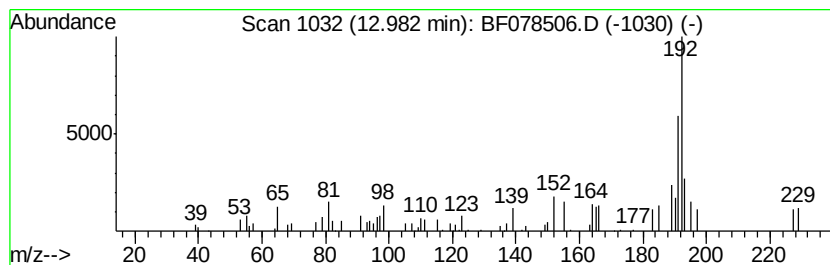
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown12.98 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	2.55 ng	31021	Phenanthrene-d10	12.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	49
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	47
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	47
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	47
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078506.D
Acq On : 14 Apr 2015 18:33
Operator : TP/IZ
Sample : G1828-02 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

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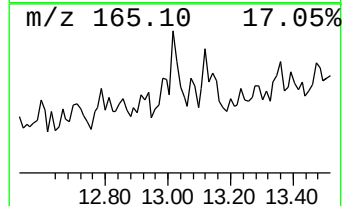
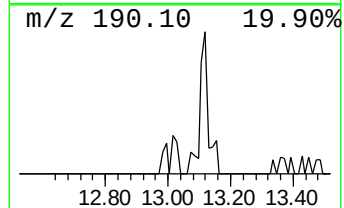
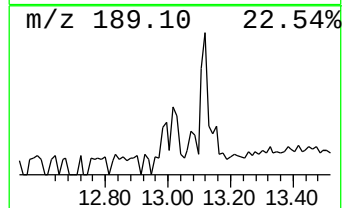
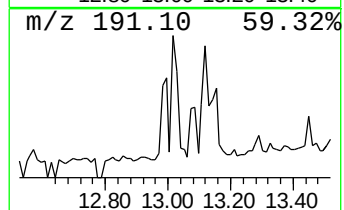
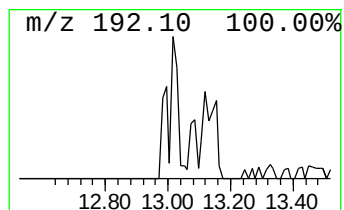
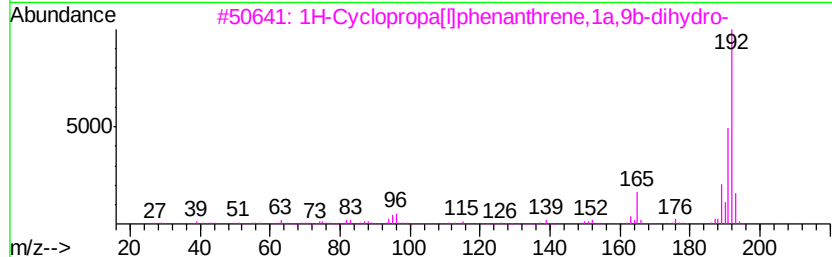
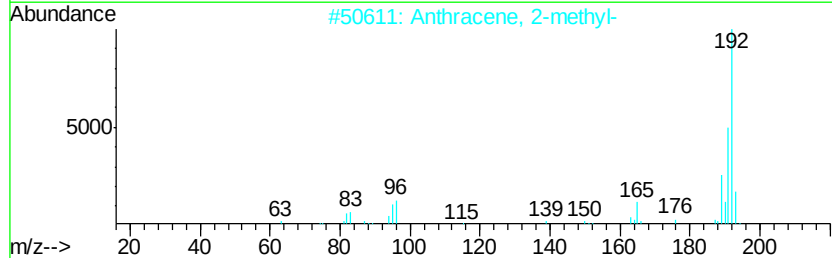
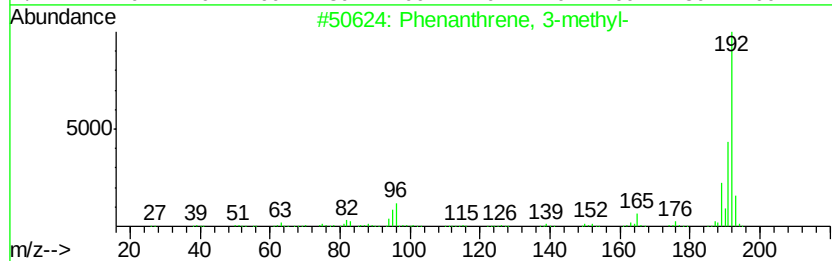
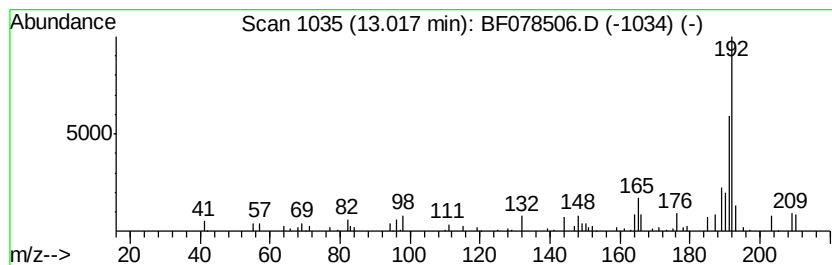
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Phenanthrene, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	2.45 ng	29790	Phenanthrene-d10	12.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	59
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	58
3			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	58
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	58
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078506.D
Acq On : 14 Apr 2015 18:33
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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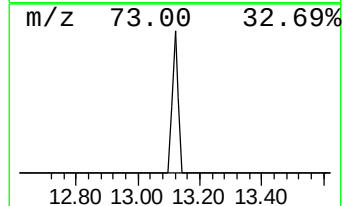
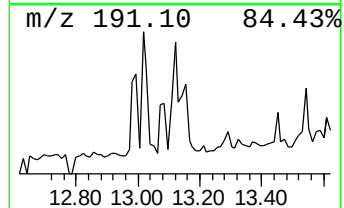
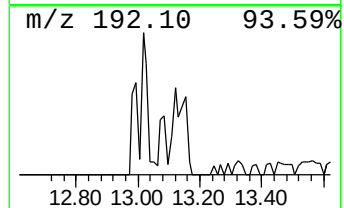
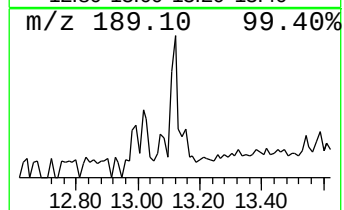
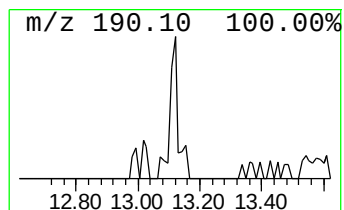
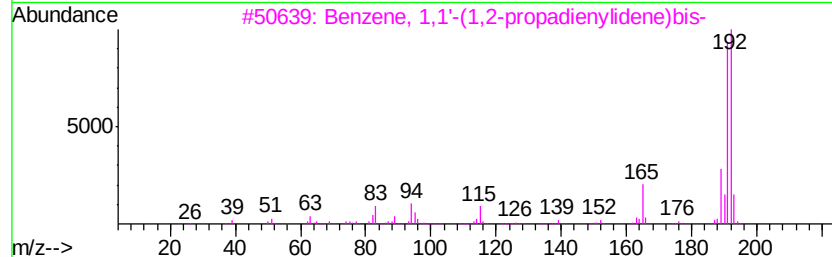
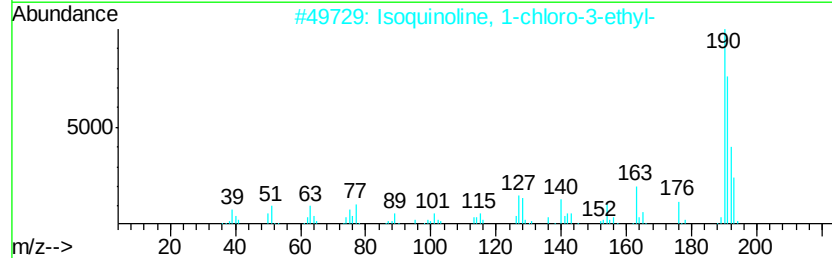
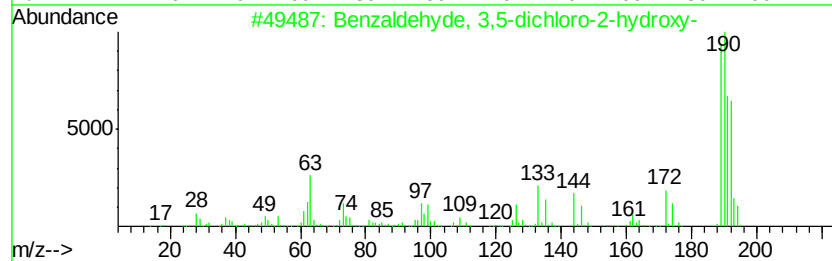
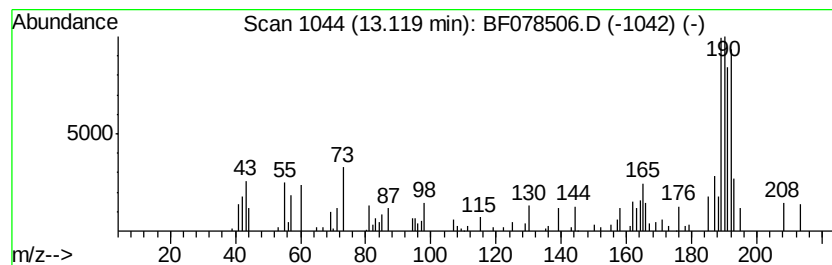
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzaldehyde, 3,5-dichloro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	2.80 ng	34033	Phenanthrene-d10	12.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2		Isoquinoline, 1-chloro-3-ethyl-	191	C11H10ClN	055150-52-2	43
3		Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	38
4		Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	38
5		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078506.D
Acq On : 14 Apr 2015 18:33
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Misc :
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ClientSampleId :
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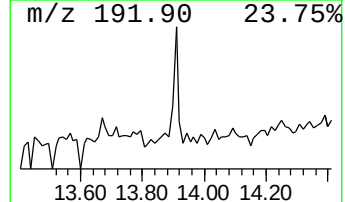
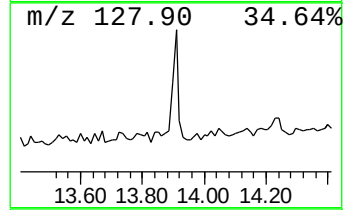
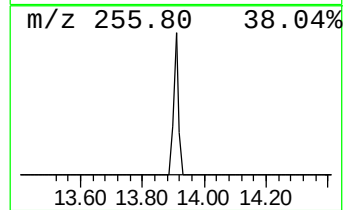
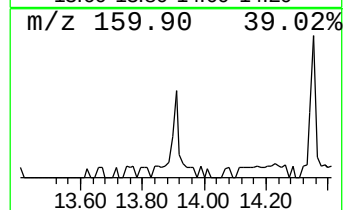
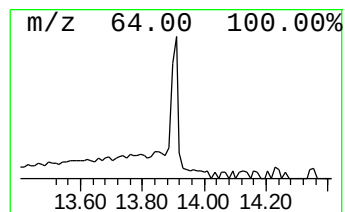
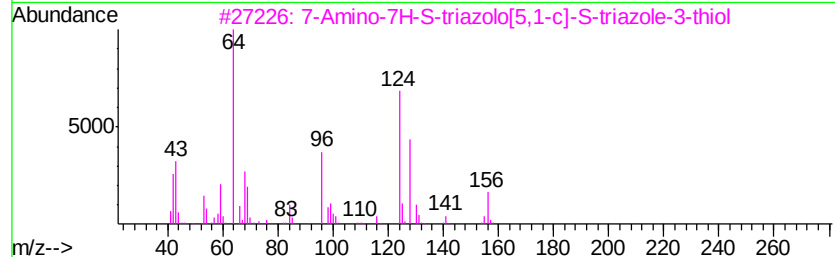
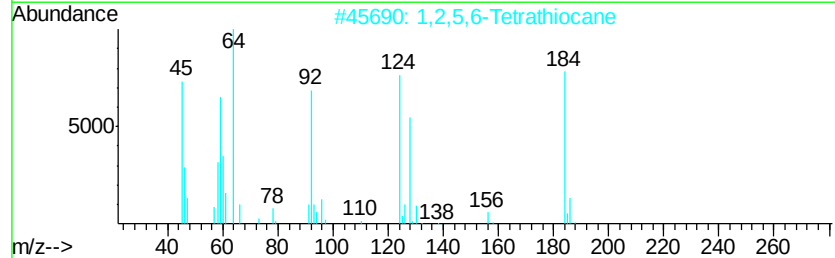
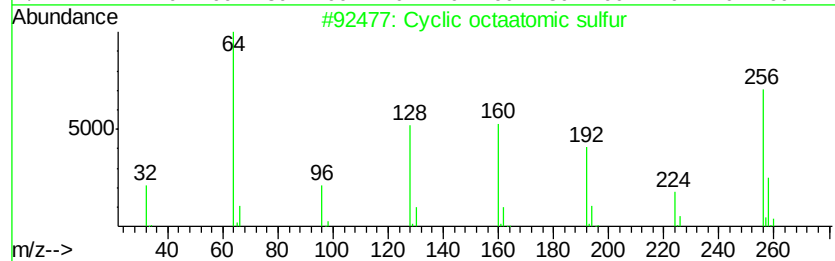
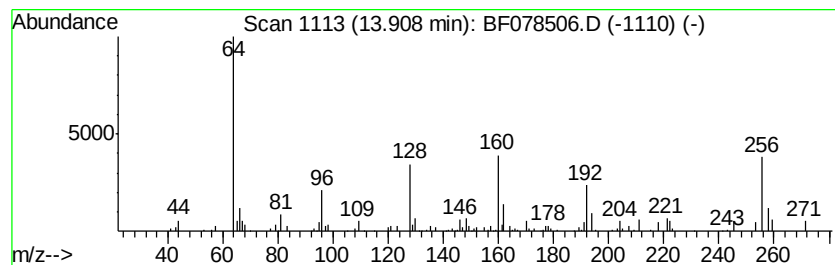
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Cyclic octaatomic sulfur Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	3.44 ng	41880	Phenanthrene-d10	12.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclic octaatomic sulfur	256	S8	010544-50-0	91
2		1,2,5,6-Tetrathiocane	184	C4H8S4	001940-01-8	45
3		7-Amino-7H-S-triazolo[5,1-c]-S-t...	156	C3H4N6S	013728-28-4	34
4		Cyclobutane, 1,1'-(1,1,2,2-tetra...	354	C10H6F12	035207-94-4	33
5		Tetrazolo[1,5-b]pyridazine, 6-ch...	155	C4H2ClN5	021413-15-0	7



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
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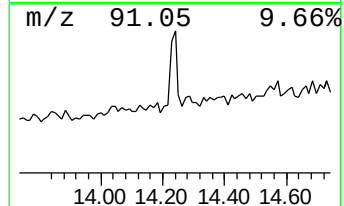
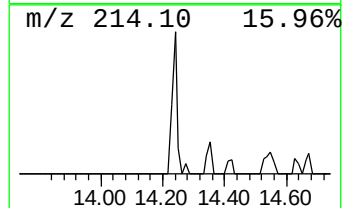
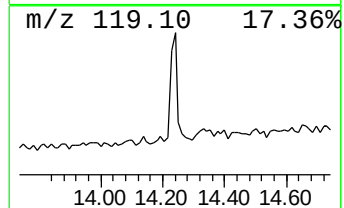
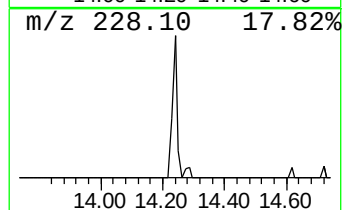
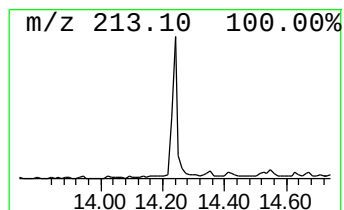
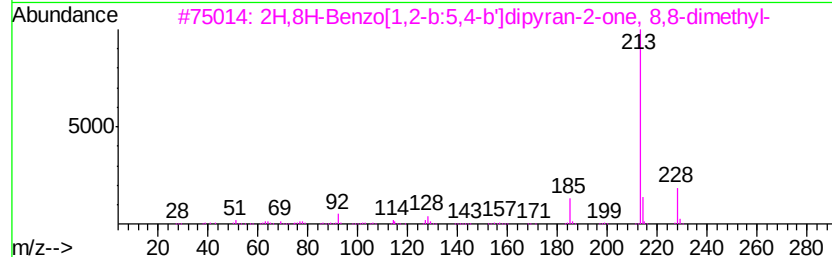
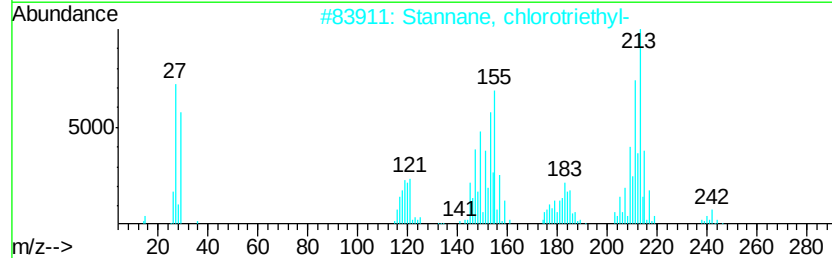
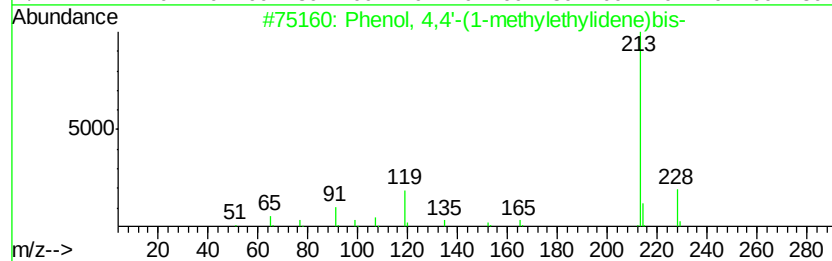
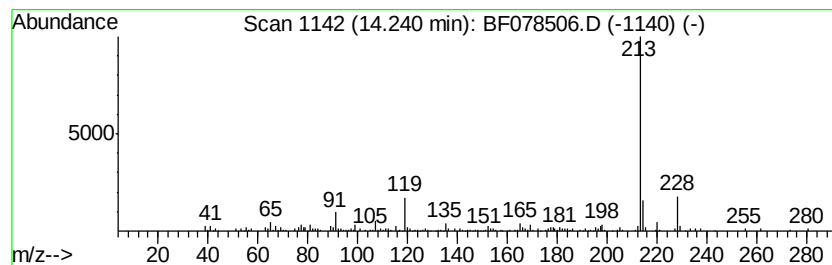
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Phenol, 4,4'-(1-methylethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	5.52 ng	110231	Chrysene-d12	15.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2			Stannane, chlorotriethyl-	242	C6H15ClSn	000994-31-0	70
3			2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	64
4			Phenol, 2-bromo-4-(1,1-dimethyle...	228	C10H13BrO	002198-66-5	64
5			2,2-Bis-(4-hydroxyphenyl)-butane	242	C16H18O2	000077-40-7	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078506.D
Acq On : 14 Apr 2015 18:33
Operator : TP/IZ
Sample : G1828-02 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HS-SB05B

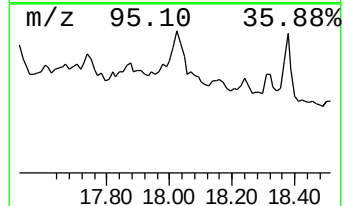
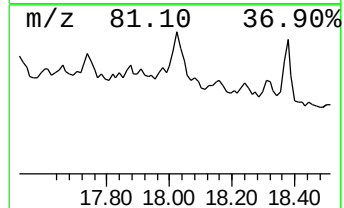
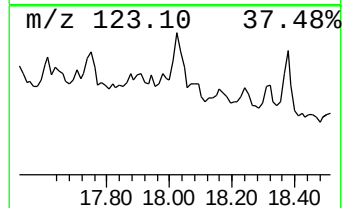
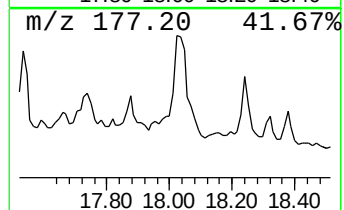
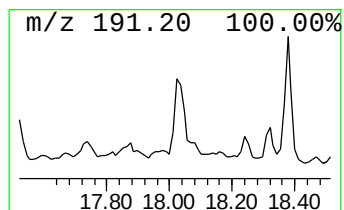
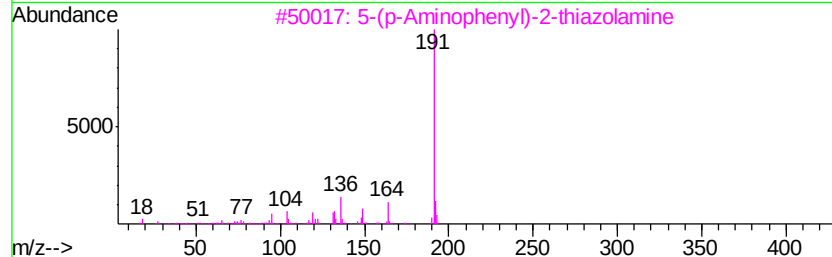
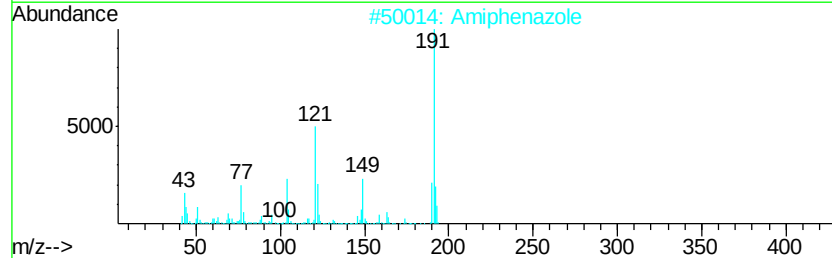
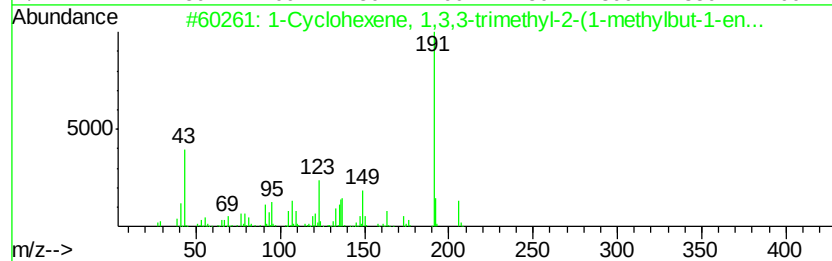
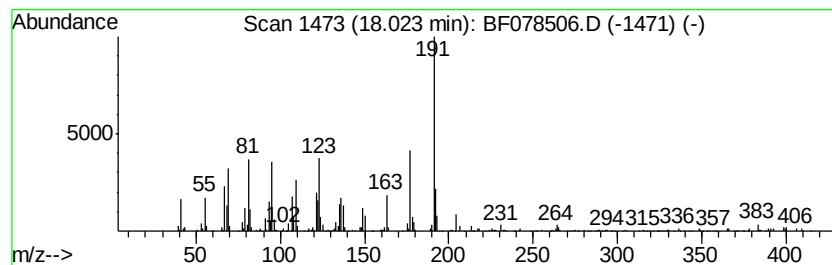
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown18.02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.02	65.18 ng	898037	Perylene-d12	17.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	38
2		Amiphenazole	191	C9H9N3S	000490-55-1	38
3		5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	35
4		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	35
5		4-(3H-Imidazo[4,5-b]pyridin-2-yl...	342	C15H18N8S	1000276-08-2	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078506.D
Acq On : 14 Apr 2015 18:33
Operator : TP/IZ
Sample : G1828-02 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HS-SB05B

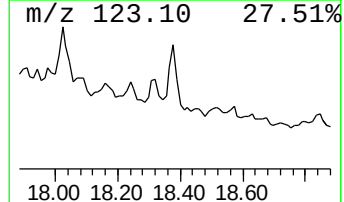
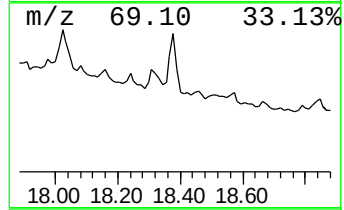
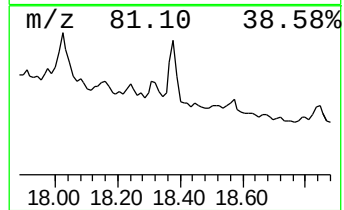
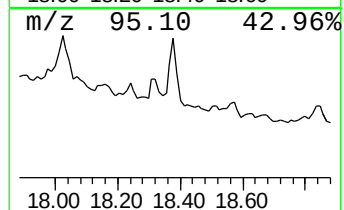
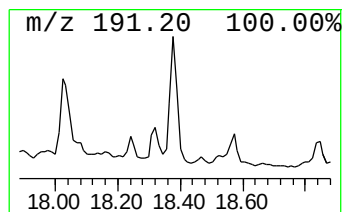
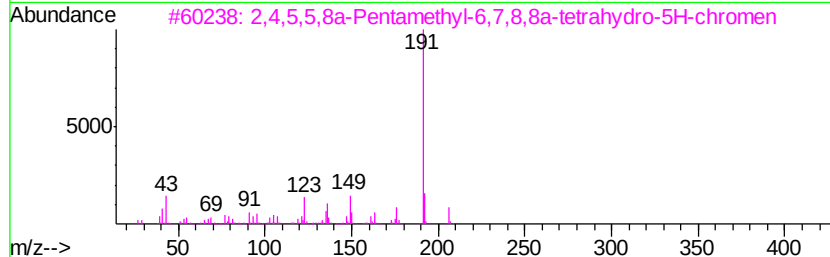
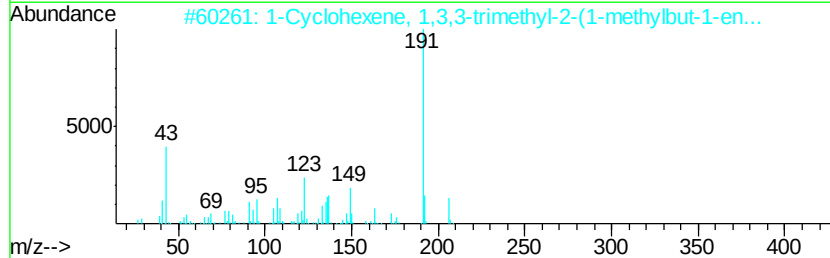
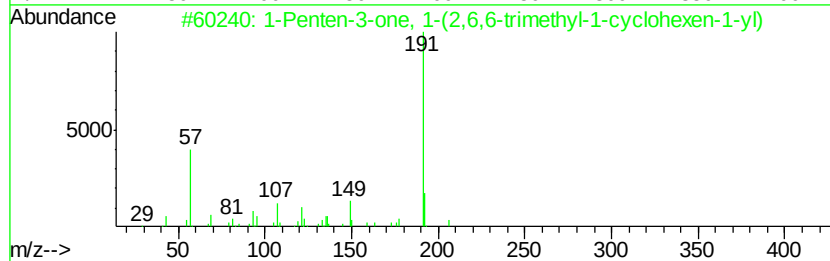
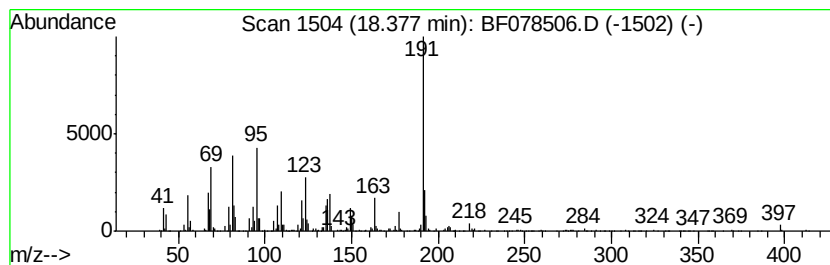
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown18.38 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.38	36.90 ng	508395	Perylene-d12	17.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	43
2		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	42
3		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	38
4		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	38
5		28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078506.D
 Acq On : 14 Apr 2015 18:33
 Operator : TP/IZ
 Sample : G1828-02 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB05B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.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
unknown6.52	6.52	17.1	ng	96210	1	6.80	112667	20.0
Borneol	8.81	2.1	ng	18064	2	8.38	173332	20.0
Biphenylene	10.36	2.4	ng	26902	3	10.54	223501	20.0
unknown10.43	10.43	2.4	ng	27221	3	10.54	223501	20.0
Benzophenone	11.44	2.2	ng	24405	3	10.54	223501	20.0
Undecane, 6-ethyl-	11.75	2.5	ng	31020	4	12.37	243236	20.0
unknown12.67	12.67	2.0	ng	24507	4	12.37	243236	20.0
unknown12.98	12.98	2.5	ng	31021	4	12.37	243236	20.0
Phenanthrene, 3-m...	13.02	2.5	ng	29790	4	12.37	243236	20.0
Benzaldehyde, 3,5...	13.12	2.8	ng	34033	4	12.37	243236	20.0
Cyclic octaatomic...	13.91	3.4	ng	41880	4	12.37	243236	20.0
Phenol, 4,4'-(1-m...	14.24	5.5	ng	110231	5	15.62	399506	20.0
unknown18.02	18.02	65.2	ng	898037	6	17.38	275542	20.0
unknown18.38	18.38	36.9	ng	508395	6	17.38	275542	20.0