

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
 Data File : BF087080.D
 Acq On : 16 May 2016 14:39
 Operator : UM/SJ
 Sample : H2968-07 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUPB

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-BF050916.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	1.735	11	13	62	rVB	3362169	7024551	100.00%	23.496%
2	4.719	272	274	287	rBV	61991	117849	1.68%	0.394%
3	5.187	313	315	329	rBV	1472439	1794884	25.55%	6.004%
4	6.273	408	410	417	rBV	1946729	1924039	27.39%	6.436%
5	6.376	417	419	422	rBV	1967627	1800957	25.64%	6.024%
6	6.604	436	439	445	rVB	639496	585648	8.34%	1.959%
7	6.765	450	453	455	rBV	1021428	1330883	18.95%	4.452%
8	7.176	487	489	495	rBV	1211872	1142692	16.27%	3.822%
9	7.896	549	552	554	rBV	763977	731835	10.42%	2.448%
10	8.970	644	646	649	rBV	1969722	1718997	24.47%	5.750%
11	9.370	680	681	684	rBV	116120	133454	1.90%	0.446%
12	9.508	690	693	695	rVB	77271	76530	1.09%	0.256%
13	9.588	698	700	702	rBV	105176	84846	1.21%	0.284%
14	9.645	702	705	707	rBV	726967	695322	9.90%	2.326%
15	10.091	742	744	745	rVB	63903	84604	1.20%	0.283%
16	10.433	771	774	776	rBV	862611	917699	13.06%	3.070%
17	10.559	783	785	789	rVB	103945	126606	1.80%	0.423%
18	10.742	798	801	806	rBV2	23961	73699	1.05%	0.247%
19	11.005	822	824	830	rVB2	55780	98605	1.40%	0.330%
20	11.119	832	834	835	rBV	701688	615645	8.76%	2.059%
21	11.199	838	841	843	rVV	120534	139926	1.99%	0.468%
22	11.622	875	878	879	rBV	96969	97411	1.39%	0.326%
23	11.668	879	882	883	rVV2	84432	165716	2.36%	0.554%
24	11.691	883	884	886	rVV	58466	77201	1.10%	0.258%
25	11.725	886	887	893	rVB	144724	258419	3.68%	0.864%
26	11.919	901	904	905	rBV	63363	73087	1.04%	0.244%
27	12.171	923	926	927	rBV	82985	93565	1.33%	0.313%
28	12.205	927	929	932	rVB3	52041	83655	1.19%	0.280%
29	12.331	938	940	942	rBV	1104710	1103810	15.71%	3.692%
30	12.422	947	948	949	rBV	97084	71378	1.02%	0.239%
31	12.548	956	959	962	rBV2	685545	1040044	14.81%	3.479%
32	12.605	962	964	967	rVB2	41751	70871	1.01%	0.237%
33	12.674	968	970	971	rBV	70656	84272	1.20%	0.282%
34	12.708	971	973	975	rVB	1100076	1245131	17.73%	4.165%

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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-BF050916.M
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35	12.777	976	979	981	rBV2	90988	155146	2.21%	0.519%
36	12.879	985	988	990	rBV	178896	280768	4.00%	0.939%
37	12.948	992	994	996	rBV	182901	177702	2.53%	0.594%
38	12.982	996	997	1001	rVB	116748	134983	1.92%	0.452%
39	13.497	1040	1042	1043	rBV2	91739	117924	1.68%	0.394%
40	13.748	1061	1064	1068	rBV2	829287	1641107	23.36%	5.489%
41	14.754	1149	1152	1158	rBV	320146	752415	10.71%	2.517%
42	15.062	1177	1179	1182	rBV	305857	386922	5.51%	1.294%
43	15.120	1182	1184	1190	rVB2	339477	565603	8.05%	1.892%

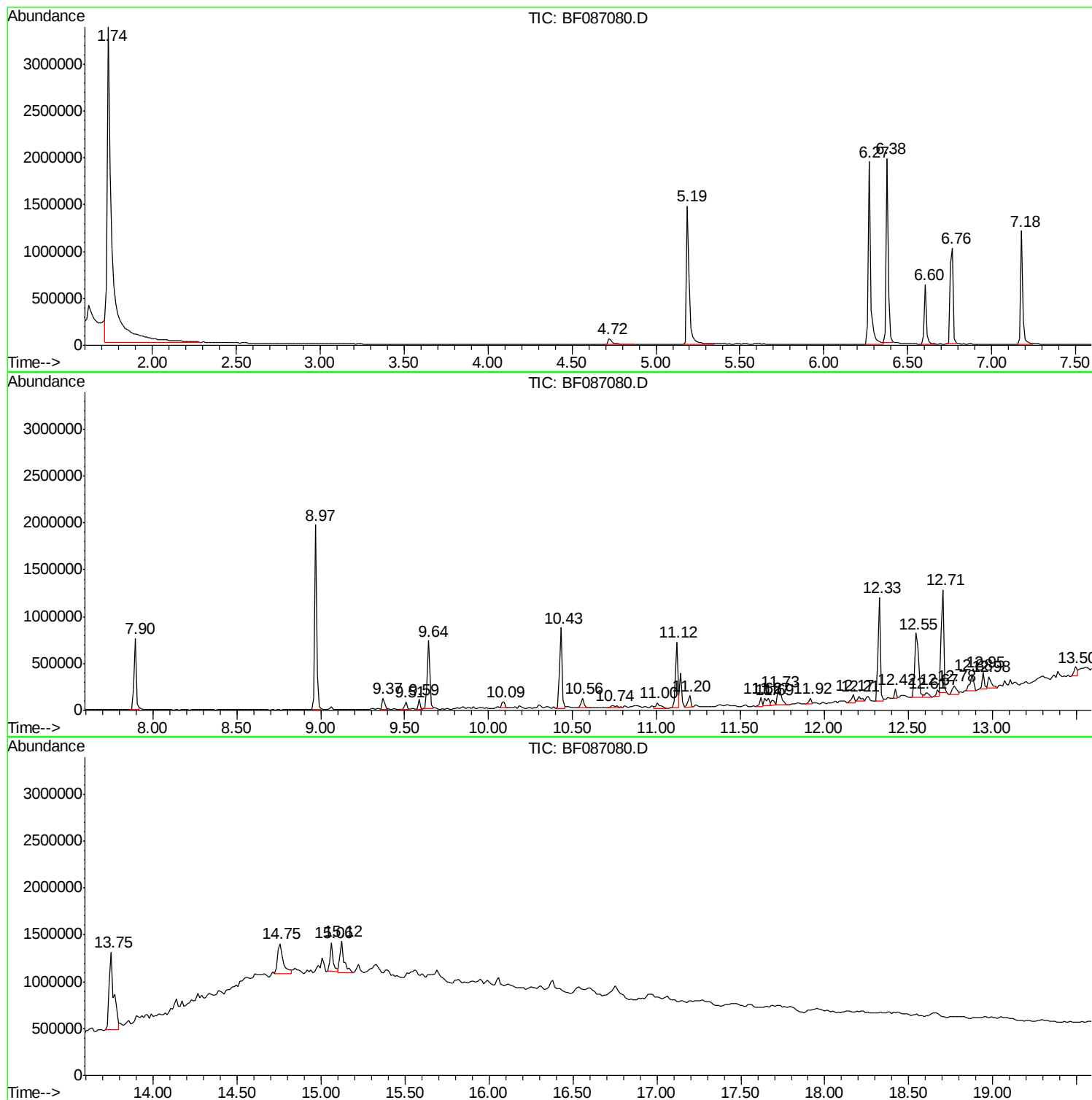
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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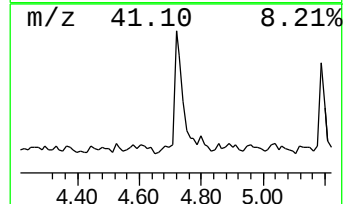
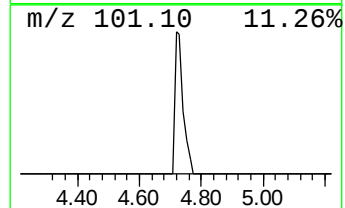
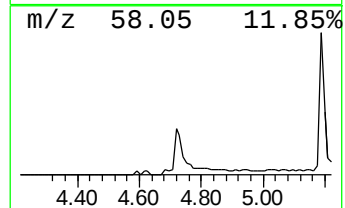
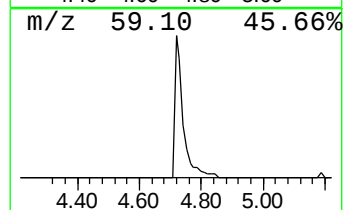
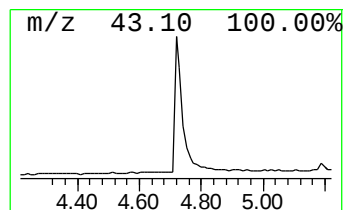
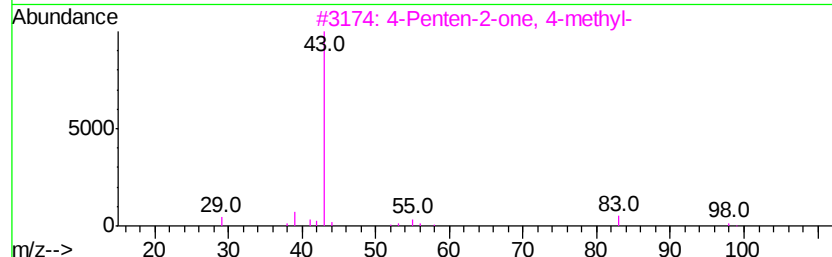
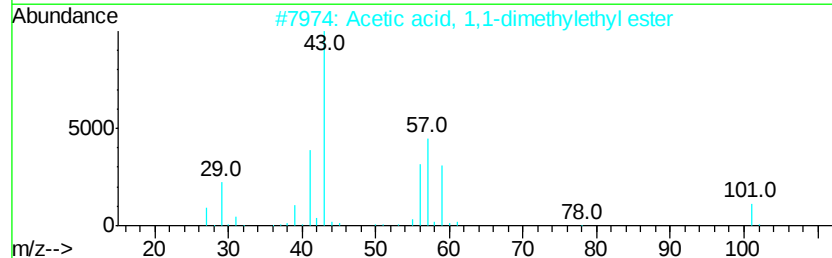
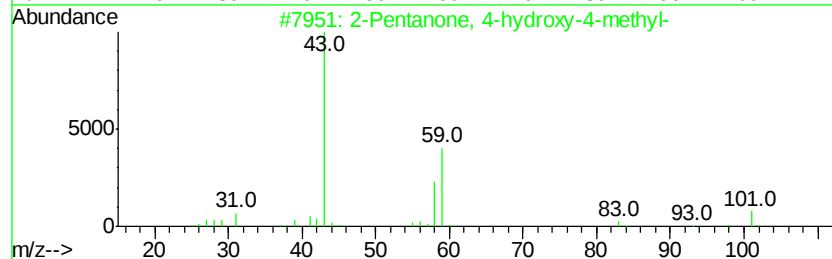
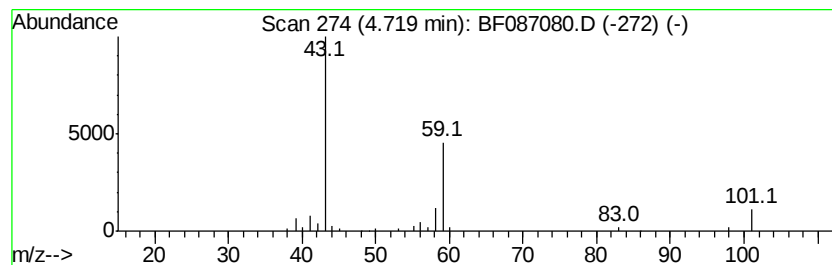
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	4.02 ng	117849	1,4-Dichlorobenzene-d4	6.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	14
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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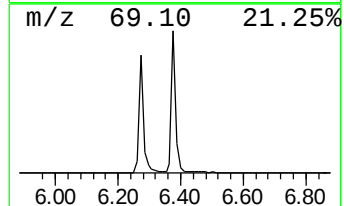
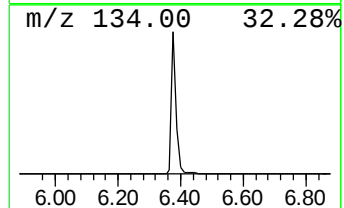
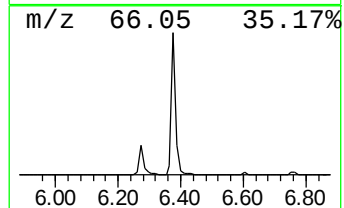
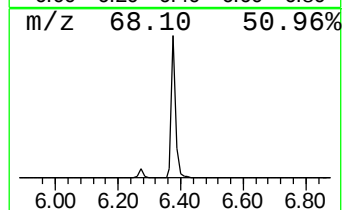
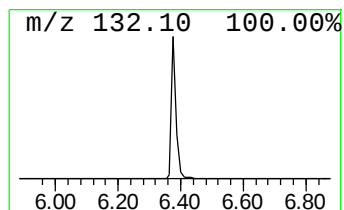
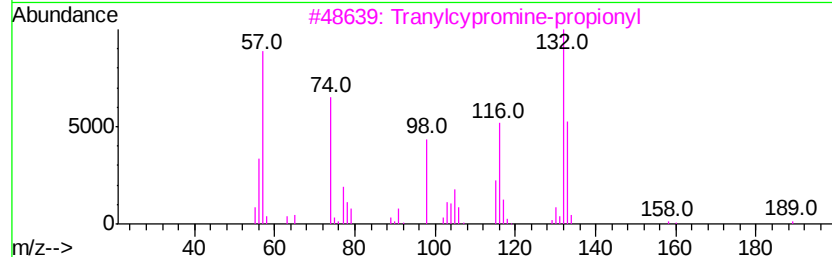
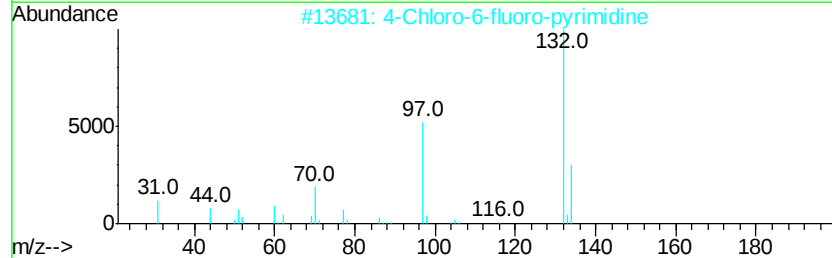
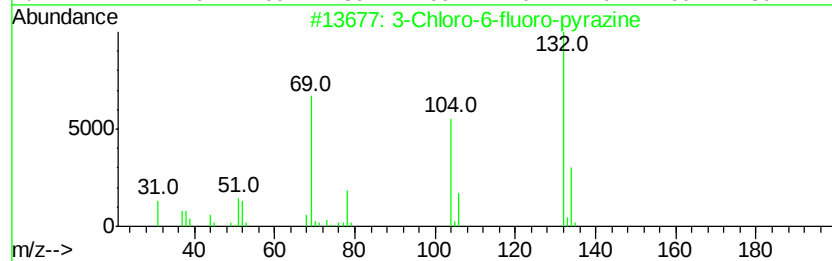
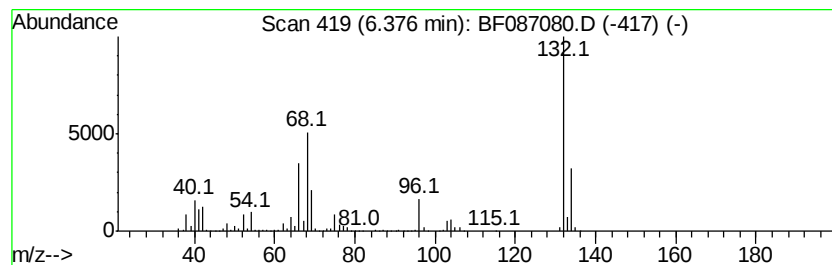
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.38 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	61.50 ng	1800960	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	22
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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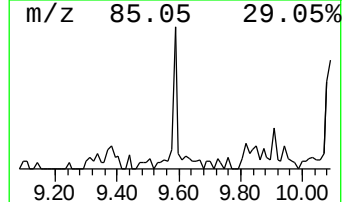
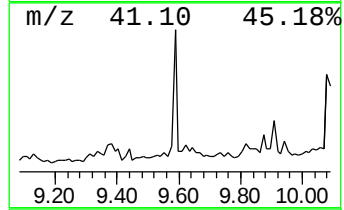
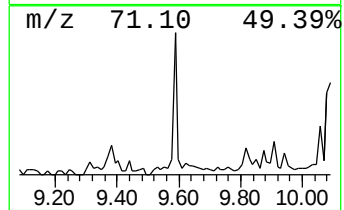
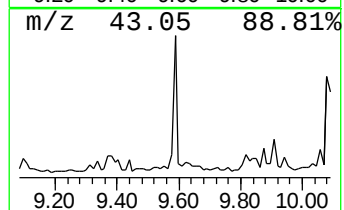
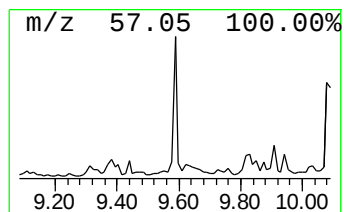
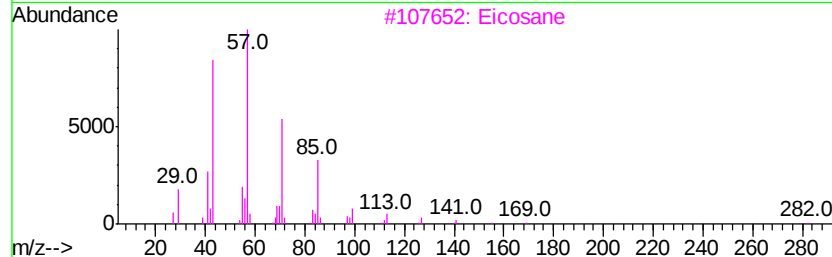
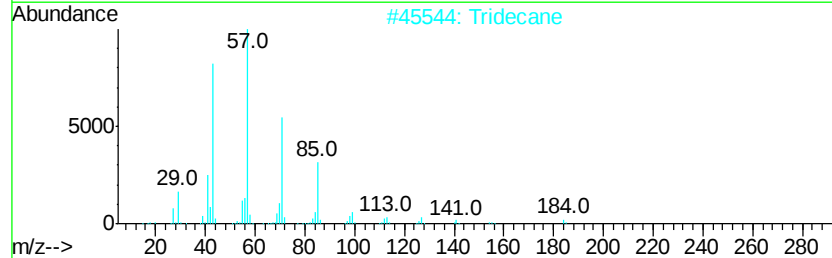
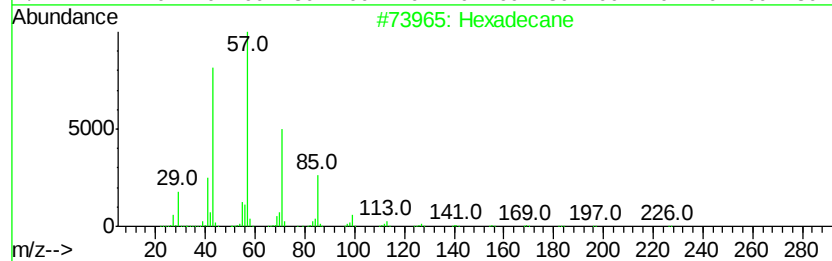
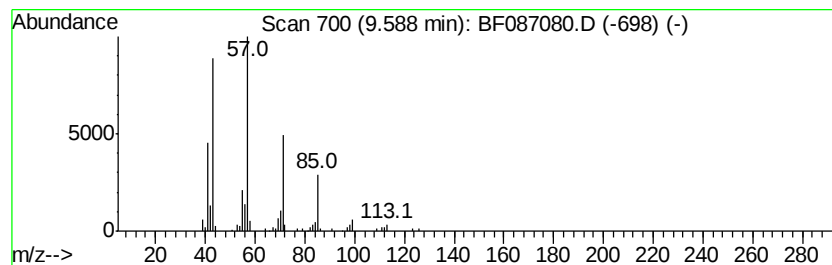
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.59	2.44 ng	84846	Acenaphthene-d10	9.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tridecane		184	C13H28	000629-50-5	90
3	Eicosane		282	C20H42	000112-95-8	83
4	Pentadecane		212	C15H32	000629-62-9	83
5	Hexatriacontane		507	C36H74	000630-06-8	83



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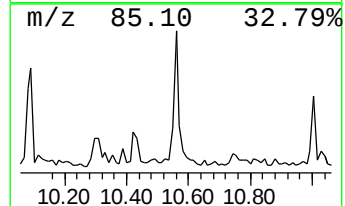
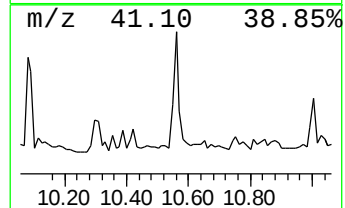
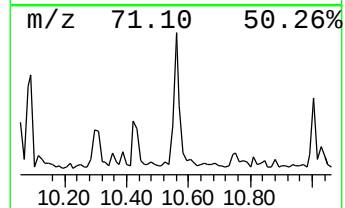
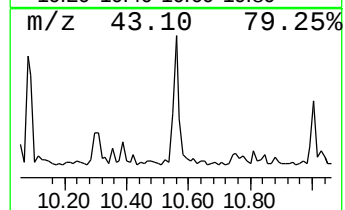
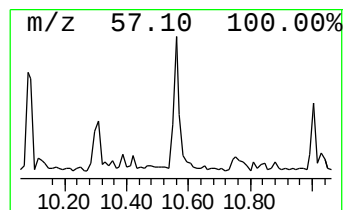
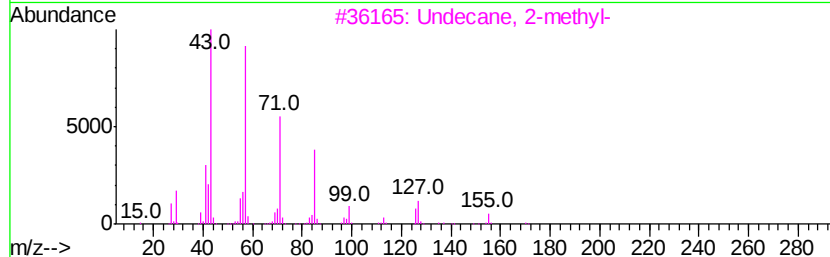
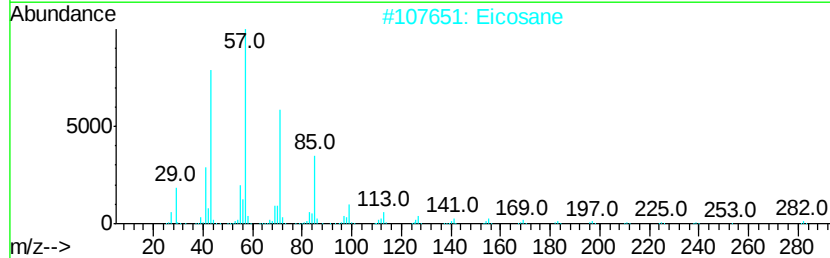
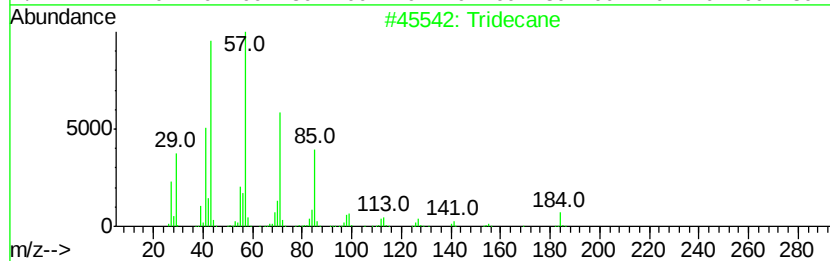
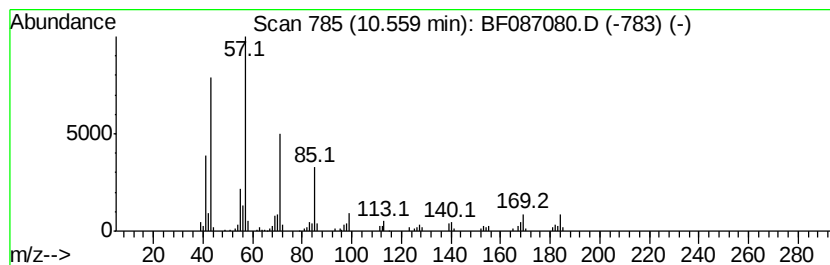
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.56	4.11 ng	126606	Phenanthrene-d10	11.12	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	Eicosane	282	C20H42	000112-95-8	74
3	Undecane, 2-methyl-	170	C12H26	007045-71-8	68
4	Octadecane	254	C18H38	000593-45-3	64
5	Octacosane	394	C28H58	000630-02-4	64



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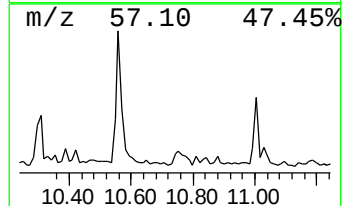
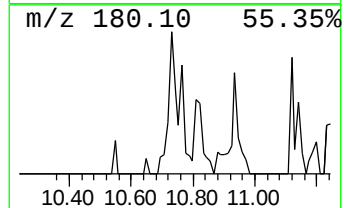
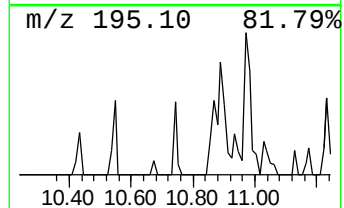
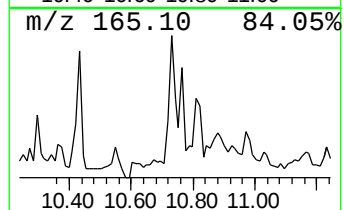
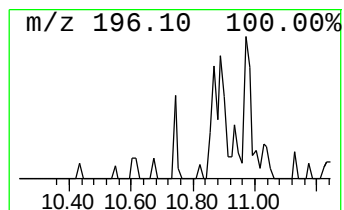
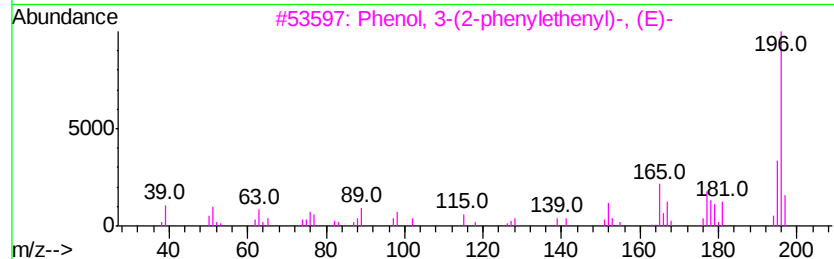
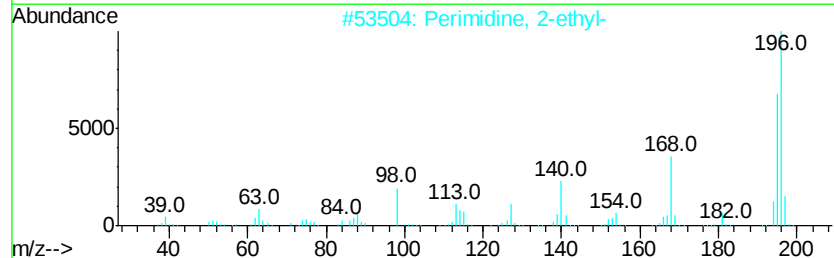
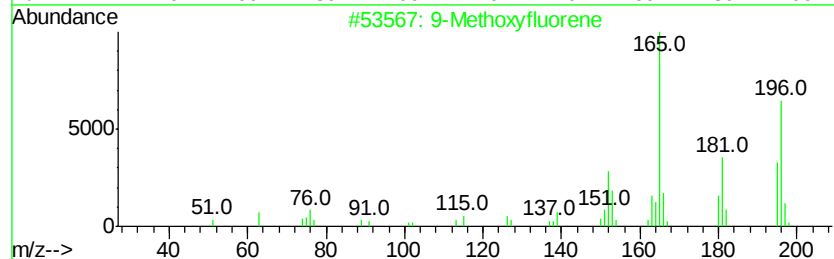
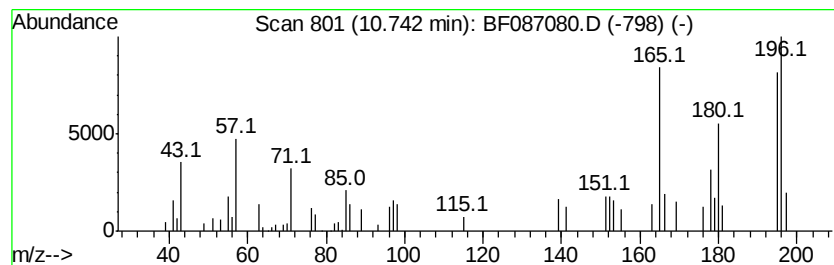
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 9-Methoxyfluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.74	2.39 ng	73699	Phenanthrene-d10	11.12		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Methoxyfluorene	196	C14H12O	019126-15-9	58
2		Perimidine, 2-ethyl-	196	C13H12N2	028478-13-9	43
3		Phenol, 3-(2-phenylethenyl)-, (E)-	196	C14H12O	017861-18-6	38
4		Pyridine, 2-[2-(4-aminophenyl)et...	196	C13H12N2	001694-46-8	38
5		8-Dimethylaminonaphthalene-1-car...	196	C13H12N2	128644-69-9	38



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Operator : UM/SJ
Sample : H2968-07 2X
Misc :
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ClientSampleId :
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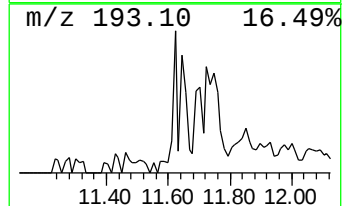
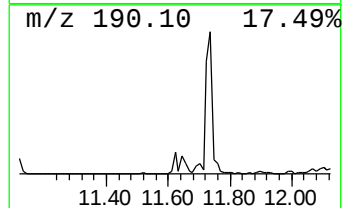
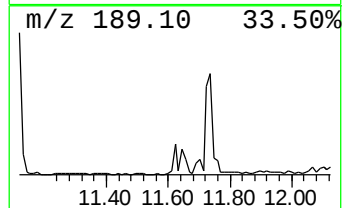
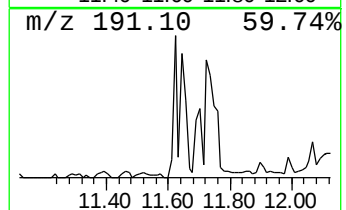
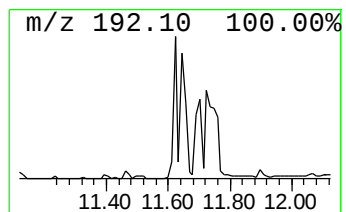
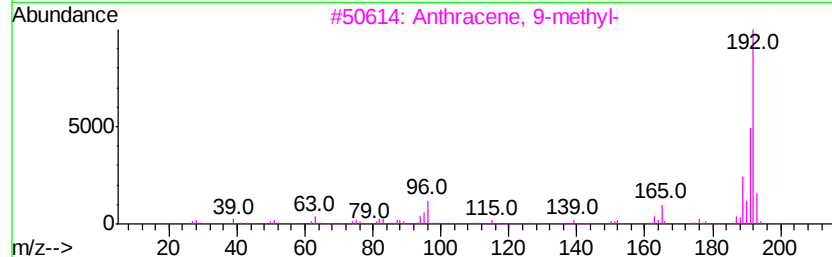
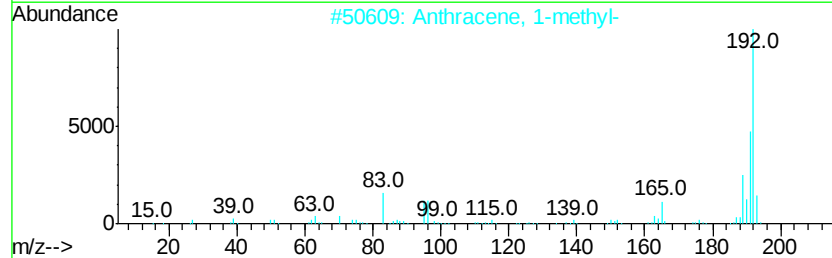
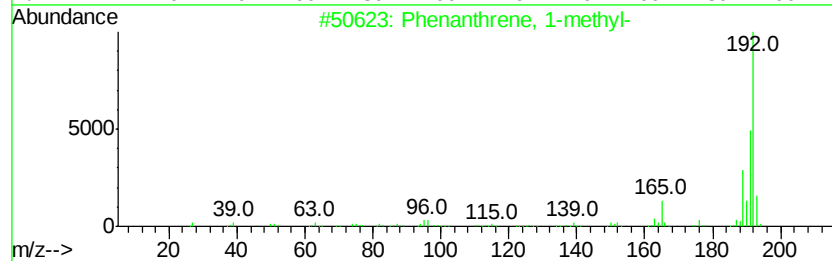
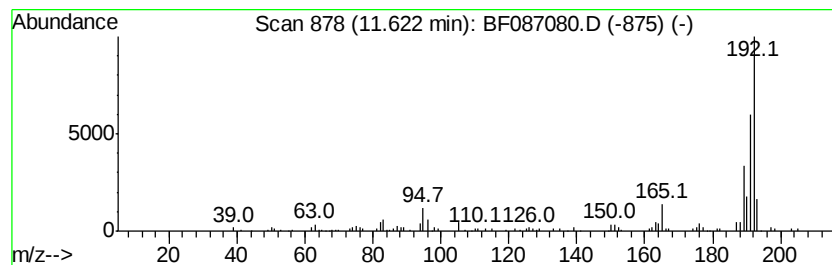
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	3.16 ng	97411	Phenanthrene-d10	11.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
Data File : BF087080.D
Acq On : 16 May 2016 14:39
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Sample : H2968-07 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

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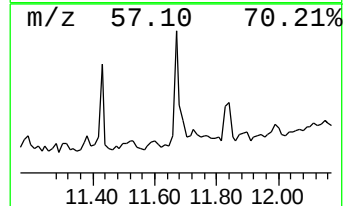
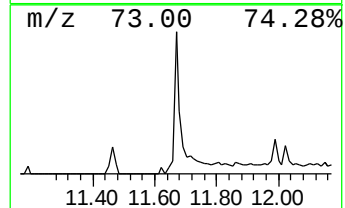
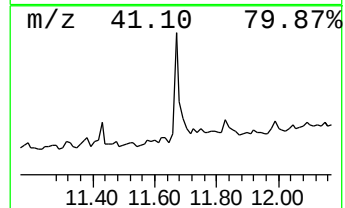
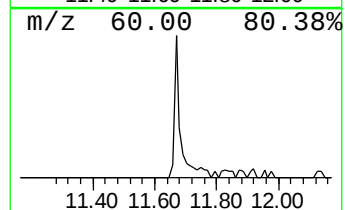
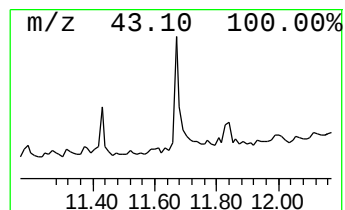
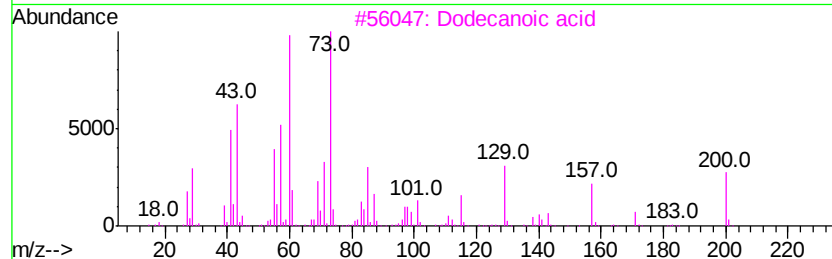
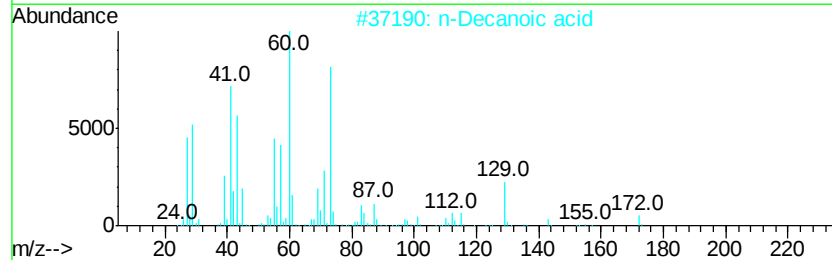
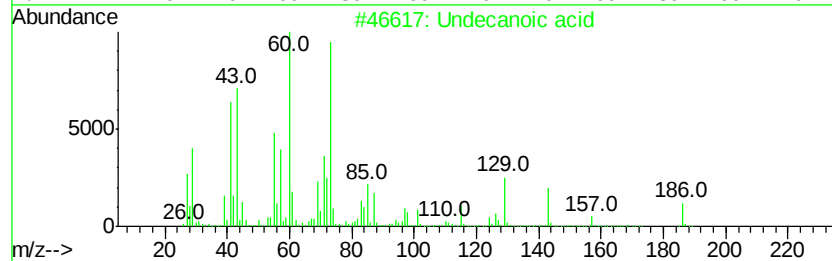
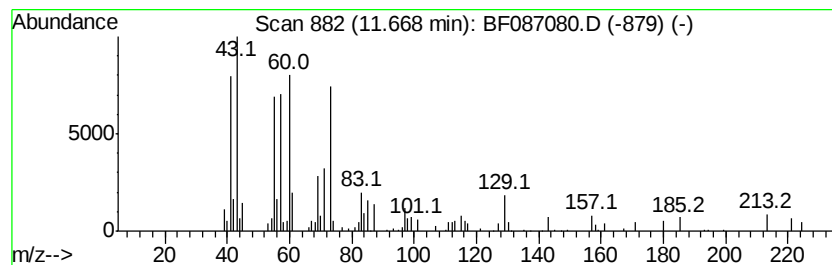
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Undecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	5.38 ng	165716	Phenanthrene-d10	11.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecanoic acid	186	C11H22O2	000112-37-8	72
2		n-Decanoic acid	172	C10H20O2	000334-48-5	59
3		Dodecanoic acid	200	C12H24O2	000143-07-7	59
4		Tridecanoic acid	214	C13H26O2	000638-53-9	59
5		Butyraldehyde, semicarbazone	129	C5H11N3O	013183-21-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
Data File : BF087080.D
Acq On : 16 May 2016 14:39
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

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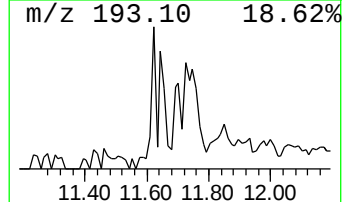
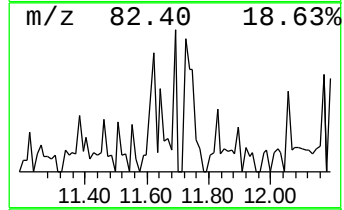
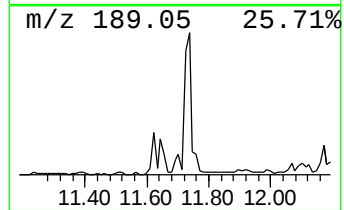
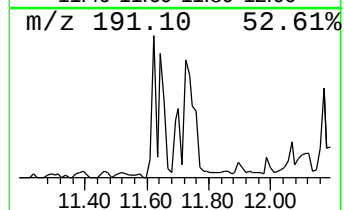
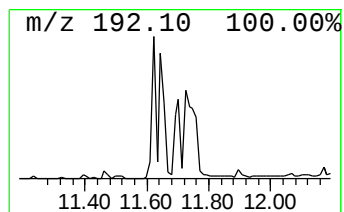
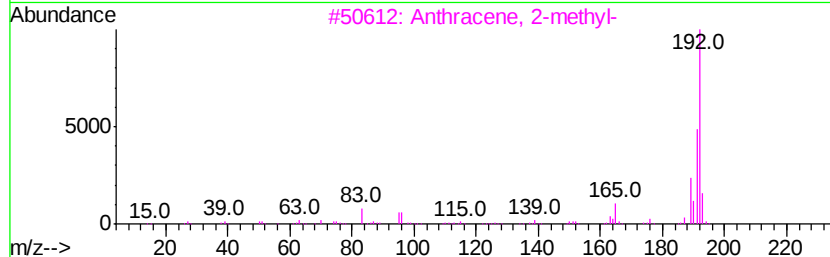
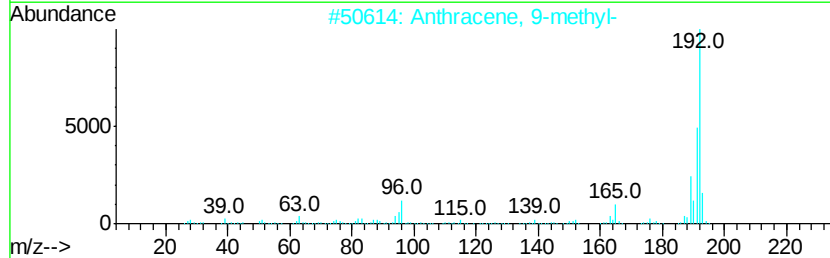
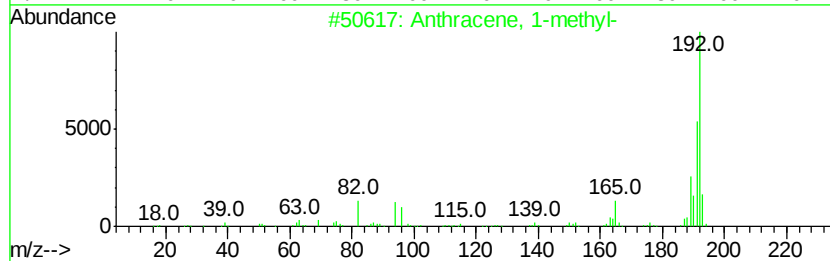
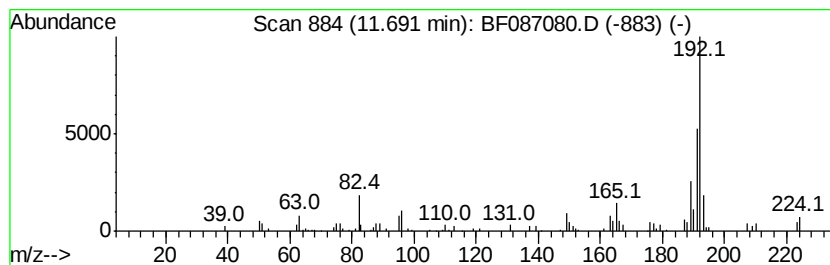
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Anthracene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	2.51 ng	77201	Phenanthrene-d10	11.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	94
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
Data File : BF087080.D
Acq On : 16 May 2016 14:39
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

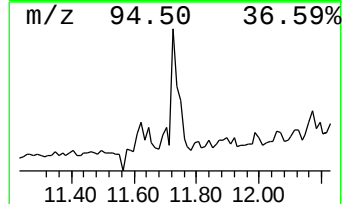
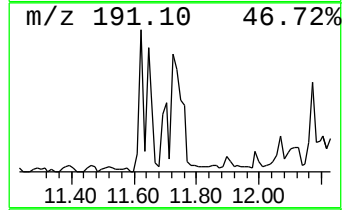
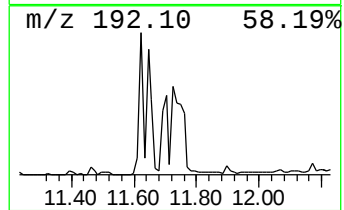
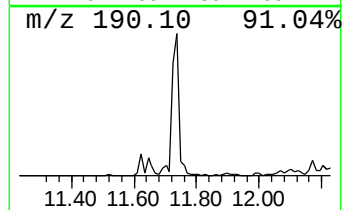
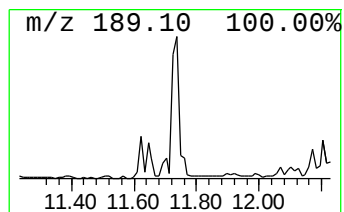
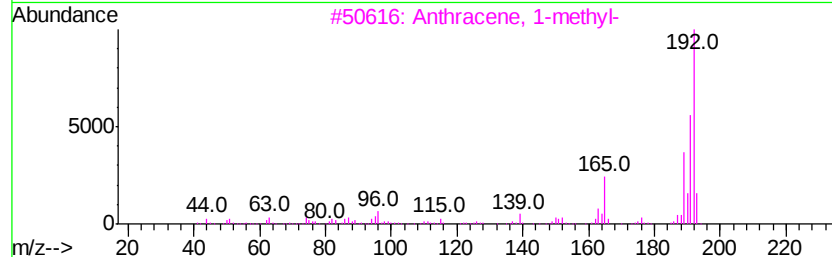
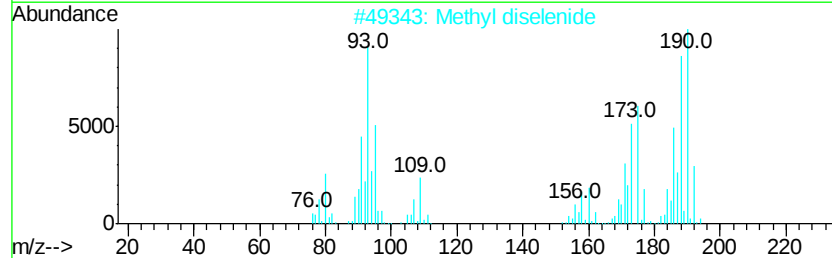
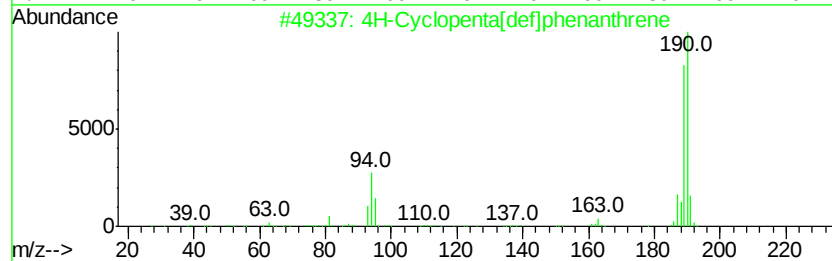
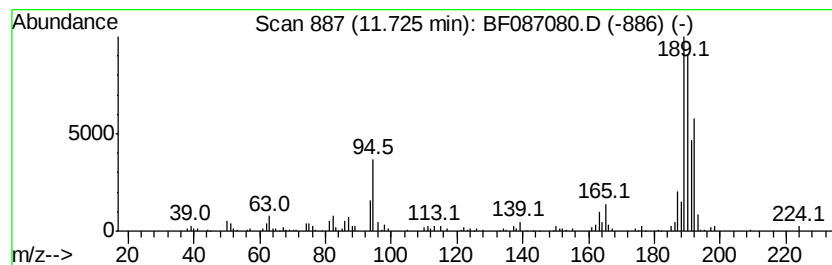
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.73	8.40 ng	258419	Phenanthrene-d10	11.12	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2	Methyl diselenide	190	C2H6Se2	007101-31-7	38
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	35
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
Data File : BF087080.D
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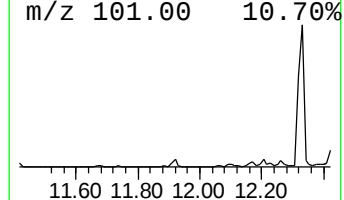
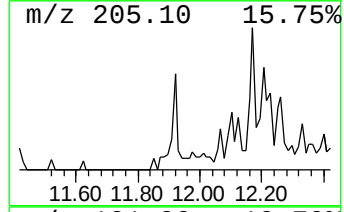
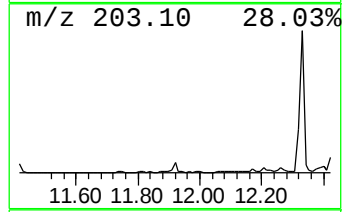
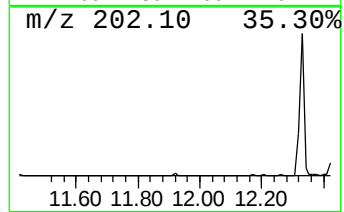
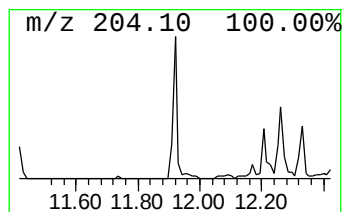
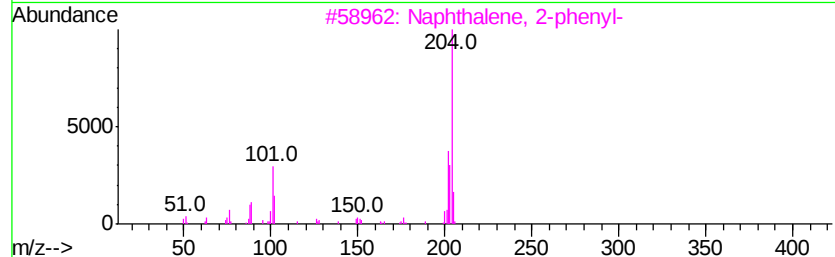
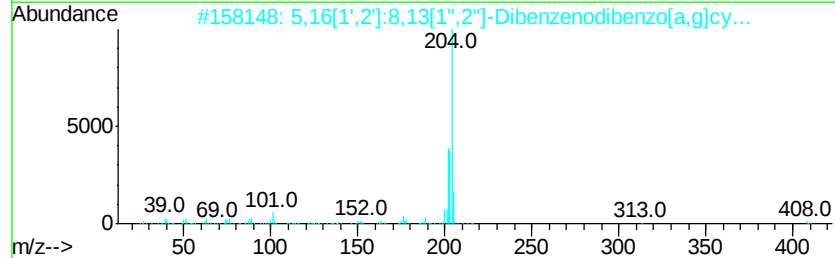
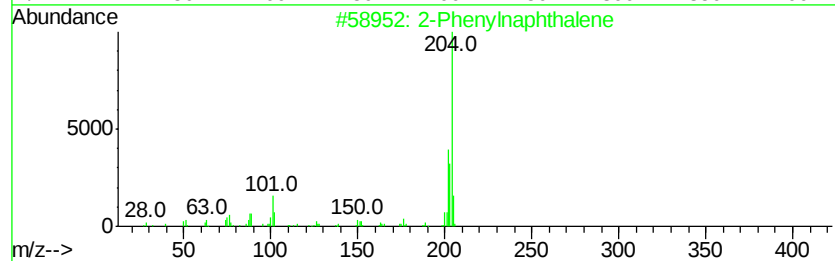
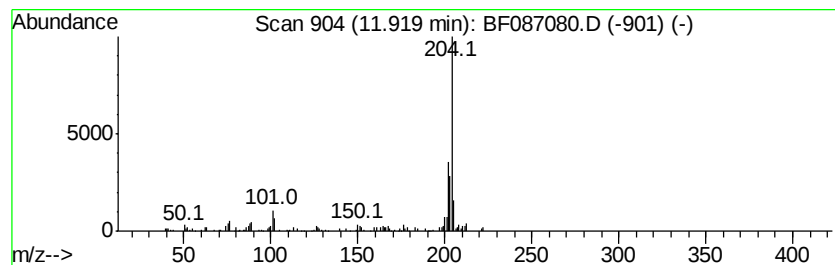
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2-Phenylnaphthalene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	2.37 ng	73087	Phenanthrene-d10	11.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenylnaphthalene	204	C16H12	035465-71-5	93
2		5,16[1',2']: ^{8,13} [1'',2'']-Dibenz...	408	C32H24	005672-97-9	80
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76
4		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	64
5		[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051616\
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Acq On : 16 May 2016 14:39
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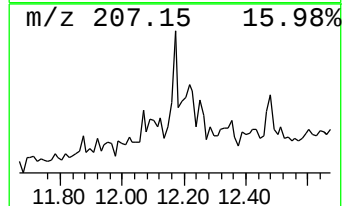
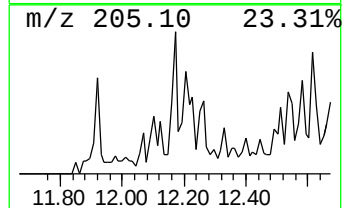
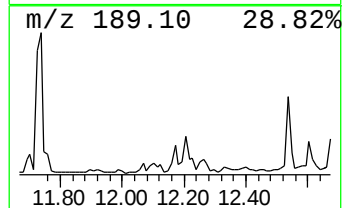
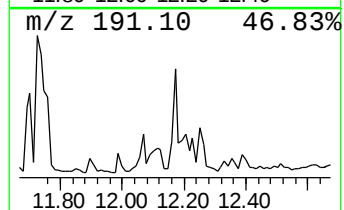
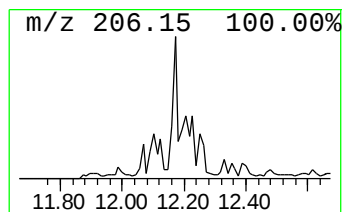
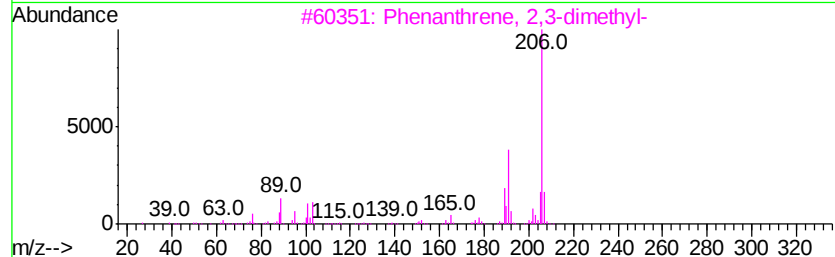
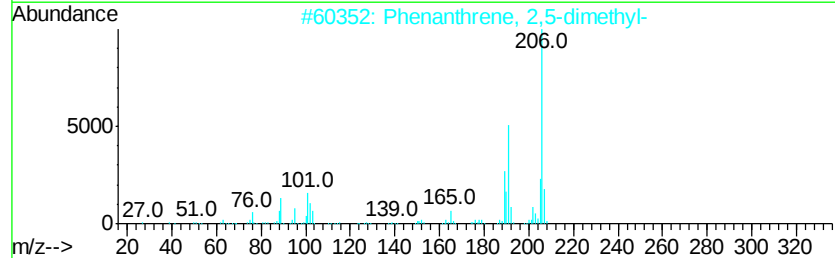
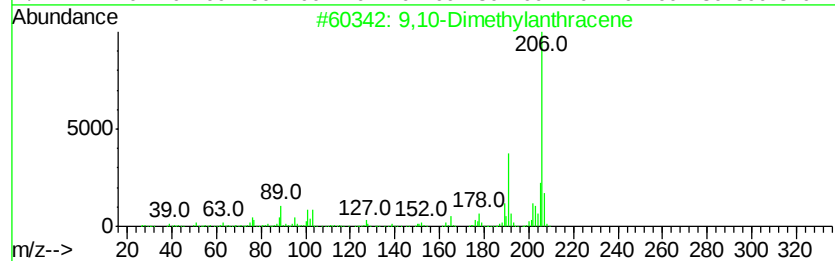
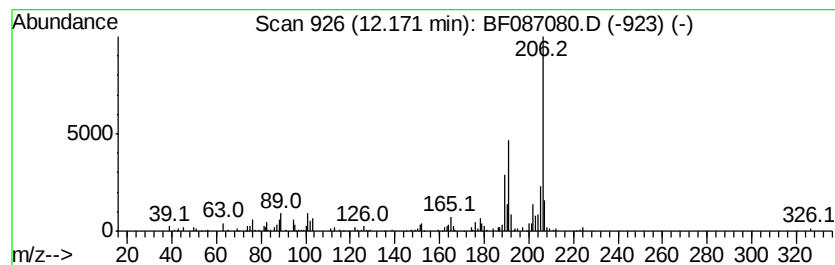
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 9,10-Dimethylantracene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.17	3.04 ng	93565	Phenanthrene-d10	11.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Dimethylantracene	206	C16H14	000781-43-1	96
2		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	94
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
5		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
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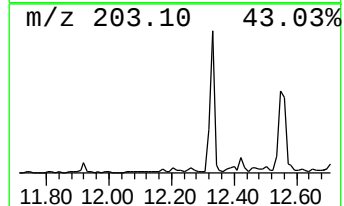
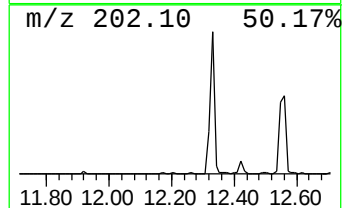
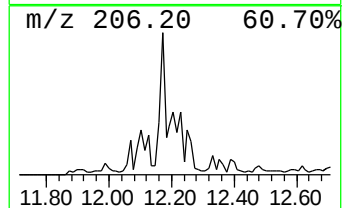
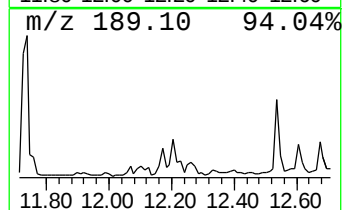
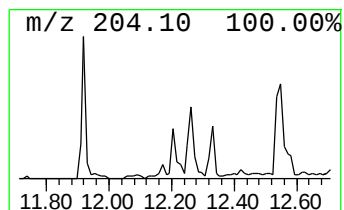
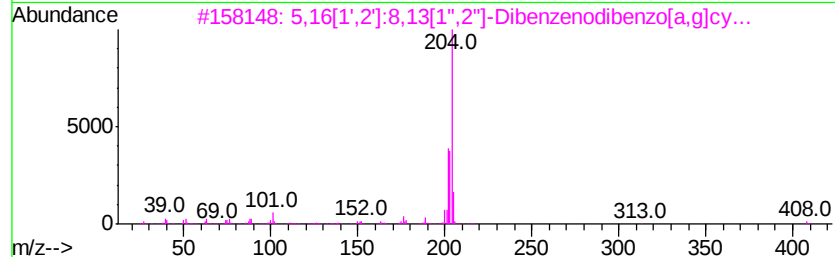
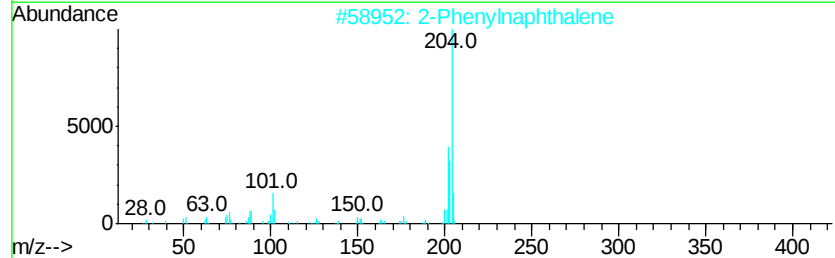
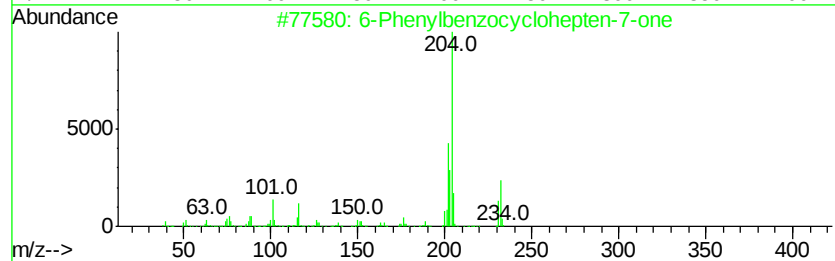
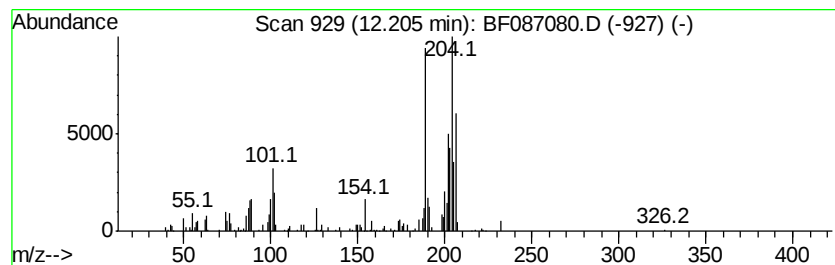
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown12.21 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.21	2.72 ng	83655	Phenanthrene-d10	11.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	43
2		2-Phenylnaphthalene	204	C16H12	035465-71-5	38
3		5,16[1',2':8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	38
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	38
5		[1,1'-Biphenyl]-4-ol, 4'-chloro-	204	C12H9ClO	028034-99-3	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051616\
Data File : BF087080.D
Acq On : 16 May 2016 14:39
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

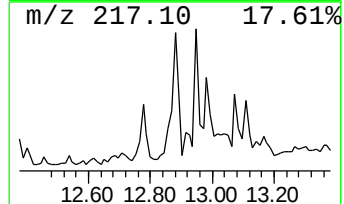
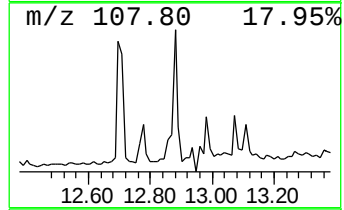
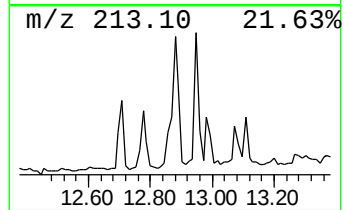
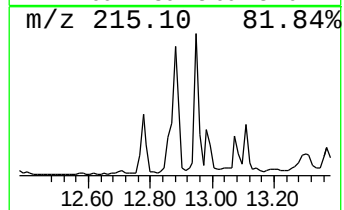
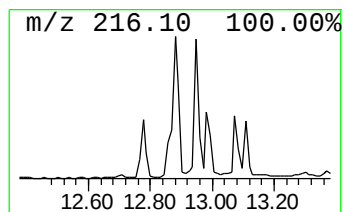
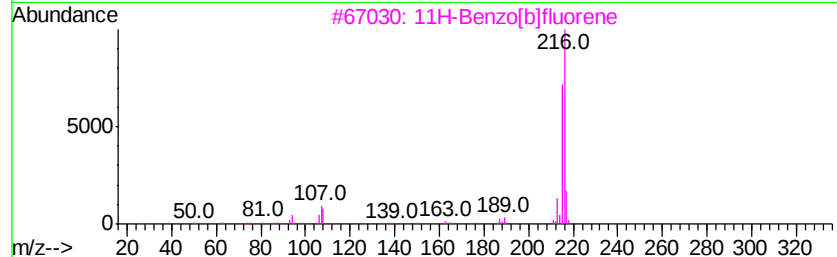
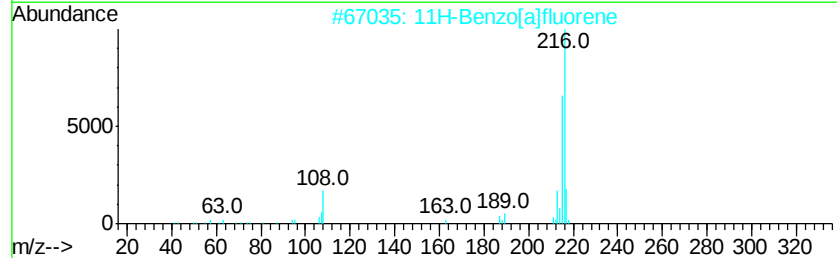
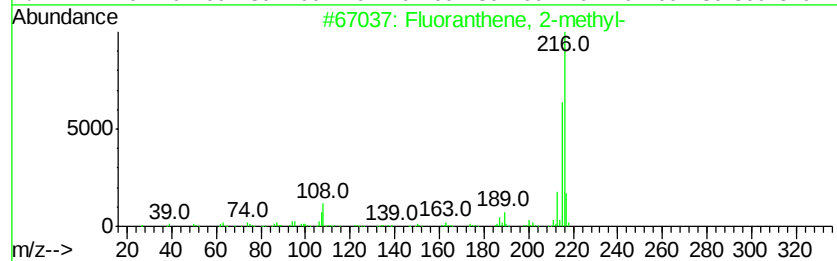
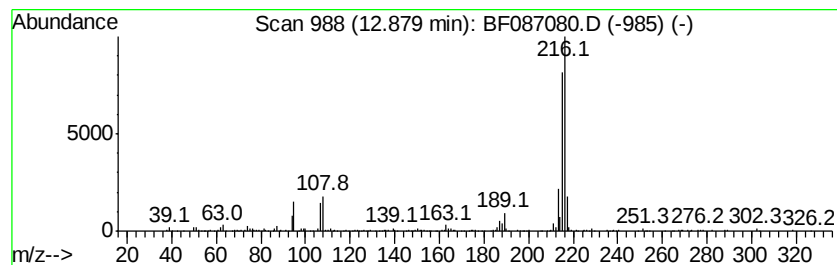
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Fluoranthene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.88	3.42 ng	280768	Chrysene-d12		13.75	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	96
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051616\
Data File : BF087080.D
Acq On : 16 May 2016 14:39
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

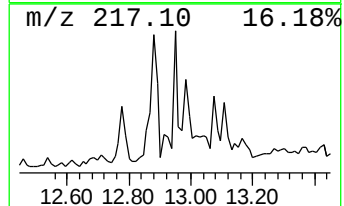
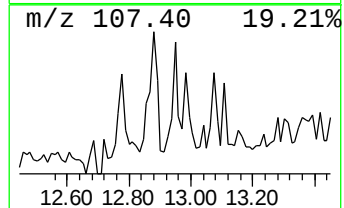
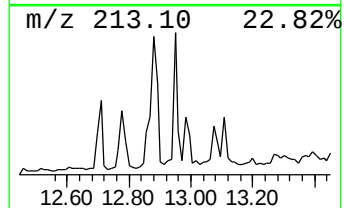
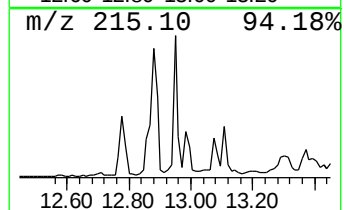
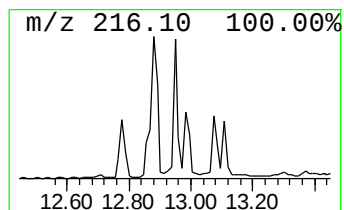
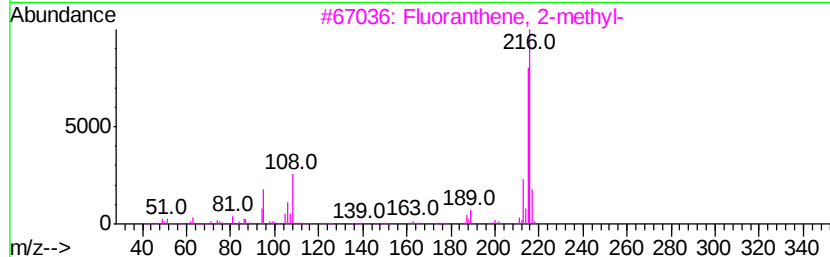
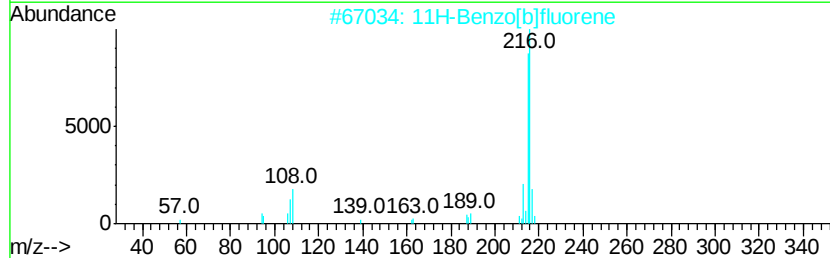
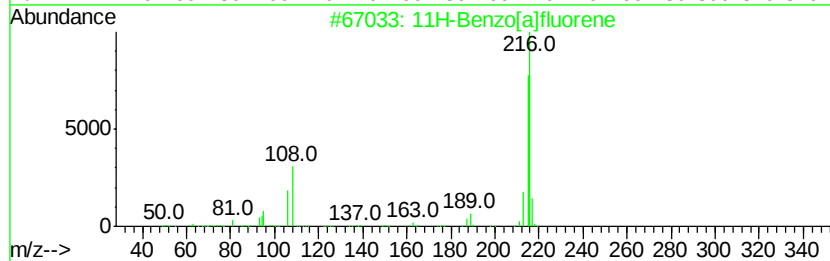
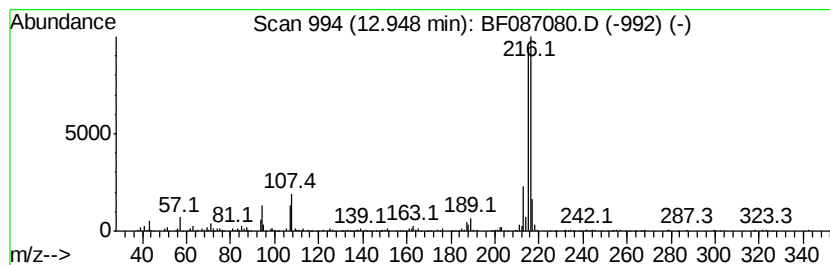
Instrument :
BNA_F
ClientSampleId :
DUPB

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 11H-Benzo[a]fluorene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	2.17 ng	177702	Chrysene-d12	13.75
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 81
2		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 81
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 81
4		7H-Benzo[c]fluorene	216 C17H12	000205-12-9 64
5		Pyrene, 1-methyl-	216 C17H12	002381-21-7 64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051616\
 Data File : BF087080.D
 Acq On : 16 May 2016 14:39
 Operator : UM/SJ
 Sample : H2968-07 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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
2-Pentanone, 4-hy...	4.72	4.0 ng		117849	1	6.60	585648	20.0
unknown6.38	6.38	61.5 ng		1800960	1	6.60	585648	20.0
Hexadecane	9.59	2.4 ng		84846	3	9.64	695322	20.0
Tridecane	10.56	4.1 ng		126606	4	11.12	615645	20.0
9-Methoxyfluorene	10.74	2.4 ng		73699	4	11.12	615645	20.0
Phenanthrene, 1-m...	11.62	3.2 ng		97411	4	11.12	615645	20.0
Undecanoic acid	11.67	5.4 ng		165716	4	11.12	615645	20.0
Anthracene, 1-met...	11.69	2.5 ng		77201	4	11.12	615645	20.0
4H-Cyclopenta[def...	11.73	8.4 ng		258419	4	11.12	615645	20.0
2-Phenylanthracene	11.92	2.4 ng		73087	4	11.12	615645	20.0
9,10-Dimethylantr...	12.17	3.0 ng		93565	4	11.12	615645	20.0
unknown12.21	12.21	2.7 ng		83655	4	11.12	615645	20.0
Fluoranthene, 2-m...	12.88	3.4 ng		280768	5	13.75	1641110	20.0
11H-Benzo[a]fluorene	12.95	2.2 ng		177702	5	13.75	1641110	20.0