

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
 Data File : BF092222.D
 Acq On : 12 Jan 2017 22:05
 Operator : UM/SJ
 Sample : I1030-13
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-0017

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-BF122116.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.721	245	248	267	rVB	1439963	2127180	44.44%	4.887%
2	5.190	287	289	303	rBV	3360718	3831744	80.05%	8.803%
3	6.264	380	383	385	rBV	3742748	4191226	87.56%	9.629%
4	6.367	390	392	395	rVV	4426779	4027742	84.14%	9.253%
5	6.596	409	412	414	rBV	1138613	1129777	23.60%	2.595%
6	6.744	423	425	428	rBV	2315783	2935798	61.33%	6.744%
7	7.167	459	462	464	rBV	2541460	2665421	55.68%	6.123%
8	7.876	522	524	527	rBV	1087947	1327880	27.74%	3.051%
9	8.962	616	619	621	rBV	4865714	4786901	100.00%	10.997%
10	9.362	652	654	656	rBV	1208640	1022410	21.36%	2.349%
11	9.579	671	673	675	rBV	59490	54222	1.13%	0.125%
12	9.625	675	677	680	rVV	1679261	1757789	36.72%	4.038%
13	10.082	713	717	718	rBV	89870	115157	2.41%	0.265%
14	10.299	733	736	739	rBV	42890	78091	1.63%	0.179%
15	10.413	744	746	749	rBV2	2176304	2776249	58.00%	6.378%
16	10.550	755	758	762	rBV	154017	197910	4.13%	0.455%
17	10.996	795	797	798	rBV2	96508	91023	1.90%	0.209%
18	11.099	804	806	809	rBV	1230364	1758829	36.74%	4.041%
19	11.351	826	828	832	rVB2	28723	48969	1.02%	0.112%
20	11.659	853	855	858	rVV2	494771	493173	10.30%	1.133%
21	11.716	858	860	866	rVB2	38939	80943	1.69%	0.186%
22	12.311	910	912	914	rVB	170189	147921	3.09%	0.340%
23	12.436	922	923	929	rBV	76381	113246	2.37%	0.260%
24	12.539	929	932	934	rVV	139377	150120	3.14%	0.345%
25	12.688	943	945	948	rVB	2636519	3162395	66.06%	7.265%
26	12.939	963	967	968	rBV2	49344	62670	1.31%	0.144%
27	13.602	1022	1025	1034	rBV2	261906	403167	8.42%	0.926%
28	13.728	1034	1036	1039	rBV	832093	1043888	21.81%	2.398%
29	14.242	1076	1081	1086	rBV3	54952	141346	2.95%	0.325%
30	14.517	1101	1105	1108	rBV	29964	56679	1.18%	0.130%
31	14.585	1108	1111	1114	rVB	93384	101496	2.12%	0.233%
32	14.642	1114	1116	1119	rVB	70862	70098	1.46%	0.161%
33	14.734	1121	1124	1127	rBV	55088	110820	2.32%	0.255%
34	14.814	1129	1131	1132	rBV	49054	60976	1.27%	0.140%

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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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.837	1132	1133	1138	rVB3	40804	76813	1.60%	0.176%
36	15.042	1149	1151	1153	rBV	43480	55607	1.16%	0.128%
37	15.088	1153	1155	1158	rBV	704699	877172	18.32%	2.015%
38	15.499	1188	1191	1193	rBV	58124	86353	1.80%	0.198%
39	15.557	1193	1196	1199	rBV3	30489	81249	1.70%	0.187%
40	16.162	1246	1249	1253	rBV4	26142	53801	1.12%	0.124%
41	16.345	1261	1265	1268	rBV	45382	104662	2.19%	0.240%
42	16.471	1272	1276	1279	rVV6	30235	120581	2.52%	0.277%
43	16.551	1279	1283	1284	rVV2	101842	231122	4.83%	0.531%
44	16.585	1284	1286	1292	rVB	245792	504548	10.54%	1.159%
45	16.814	1302	1306	1312	rBV3	67093	165102	3.45%	0.379%
46	17.648	1376	1379	1384	rVB4	16678	48515	1.01%	0.111%

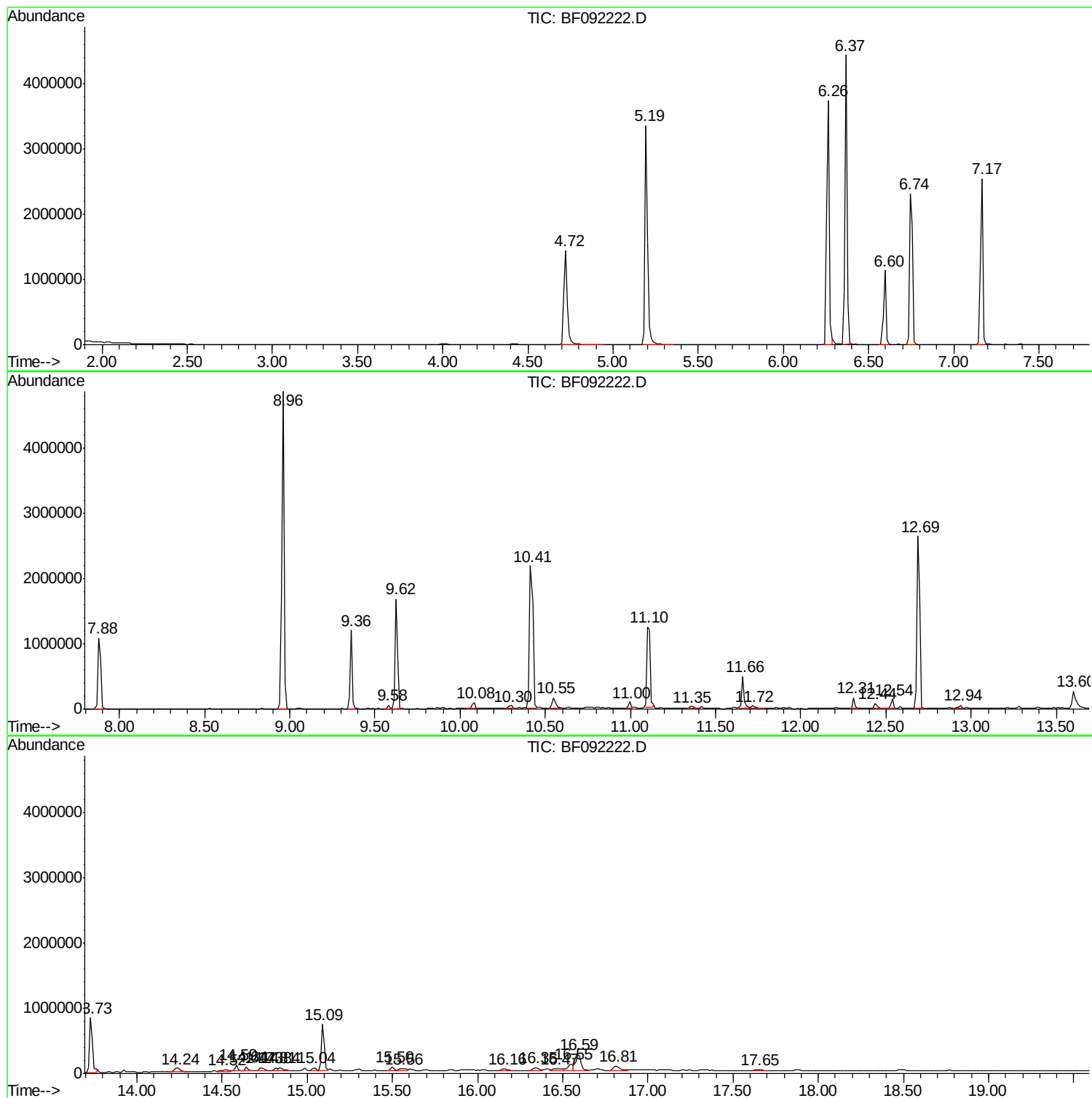
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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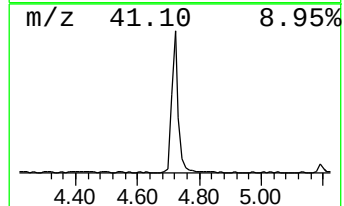
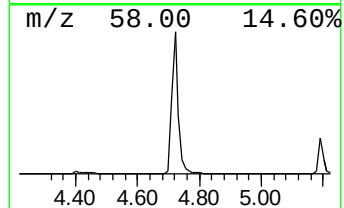
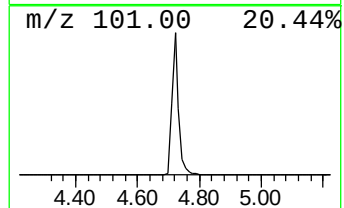
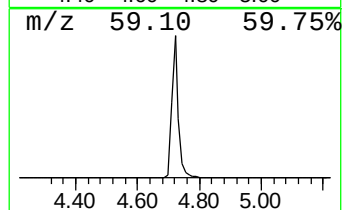
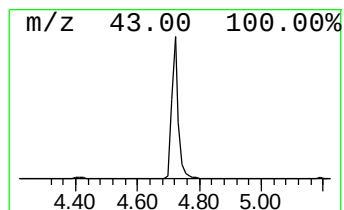
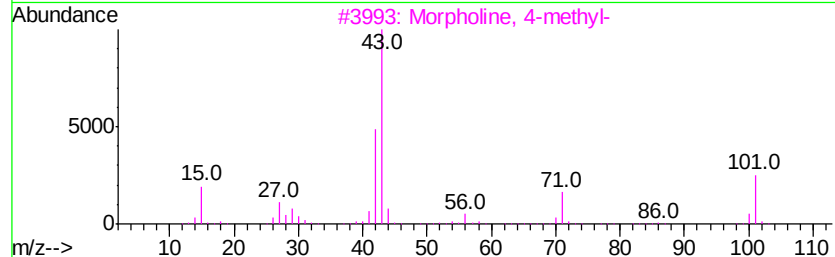
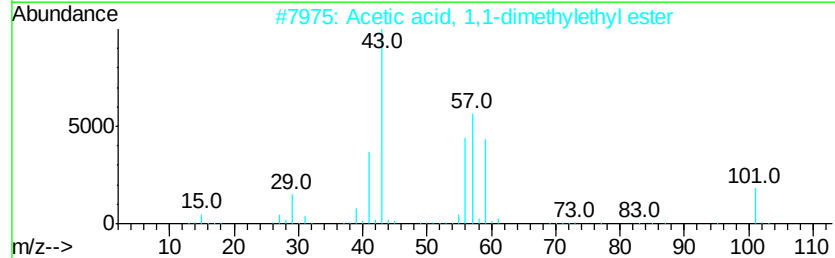
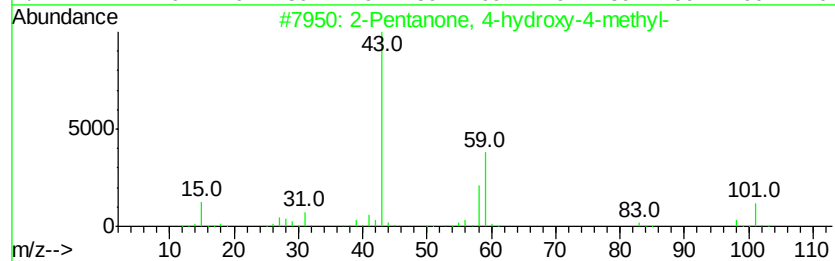
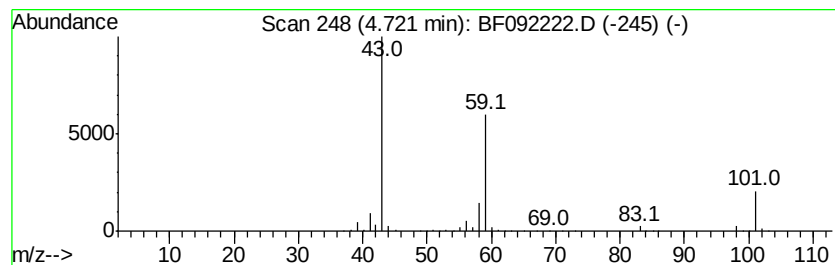
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	37.66 ng	2127180	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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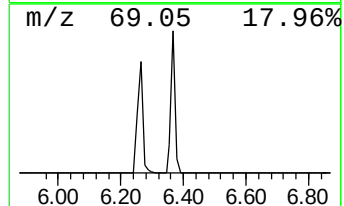
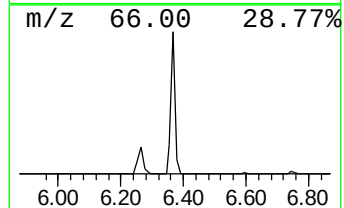
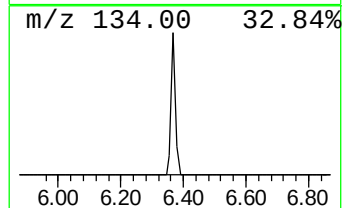
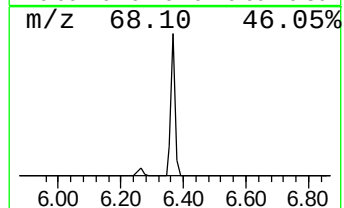
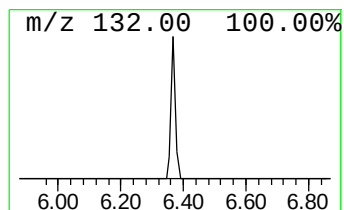
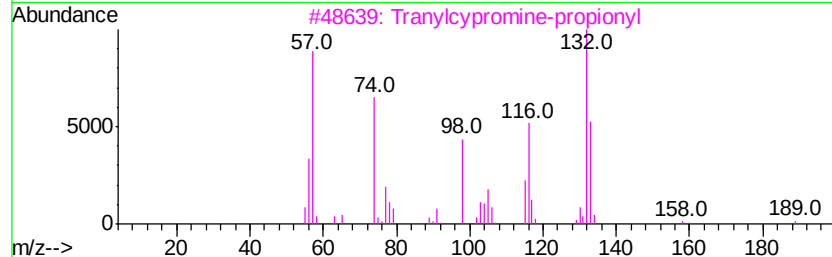
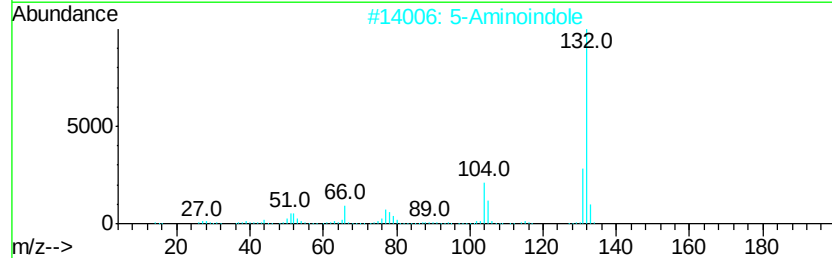
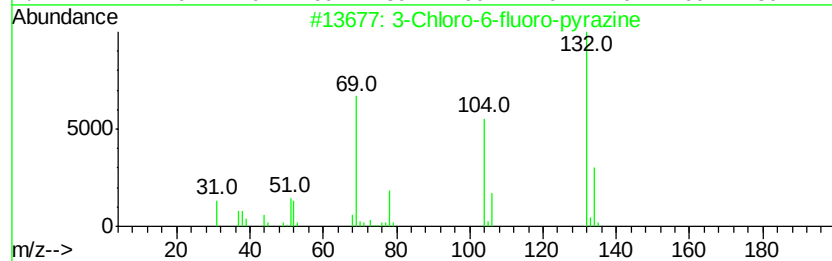
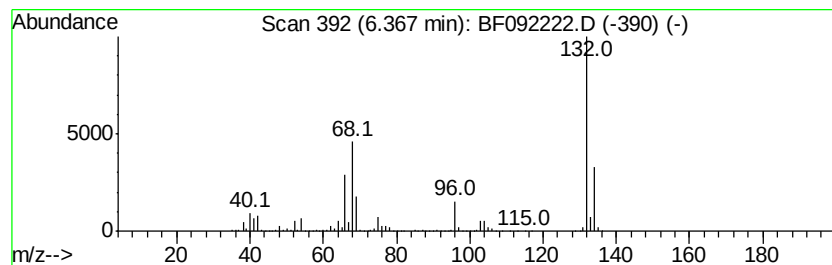
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Peak Number 2 unknown6.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.37	71.30 ng	4027740	1,4-Dichlorobenzene-d4	6.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5		Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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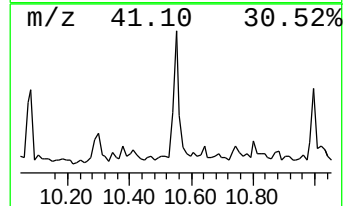
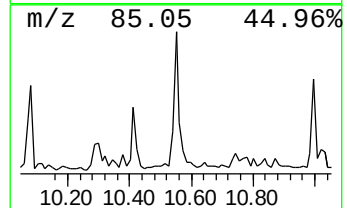
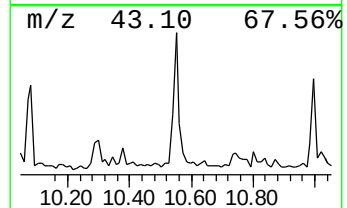
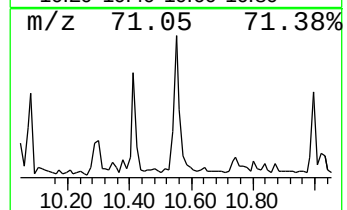
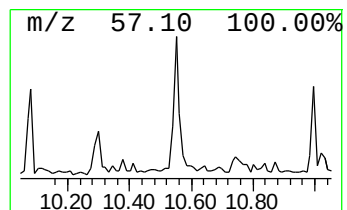
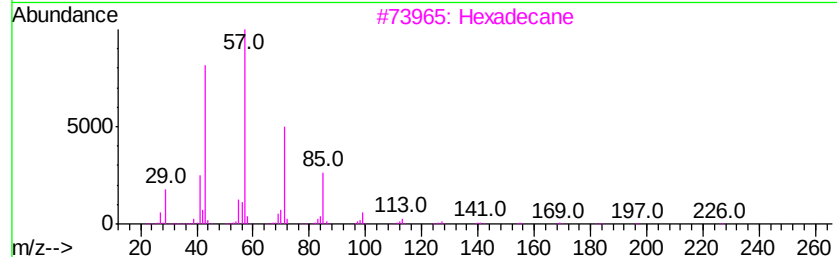
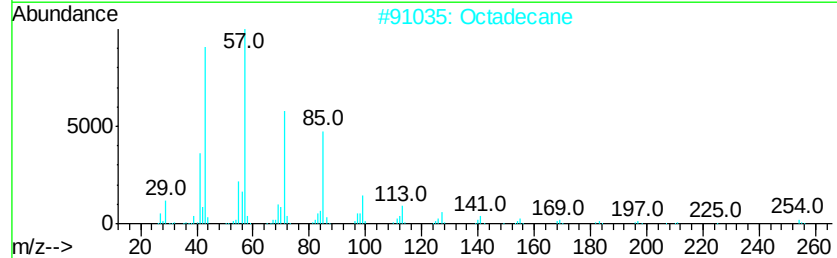
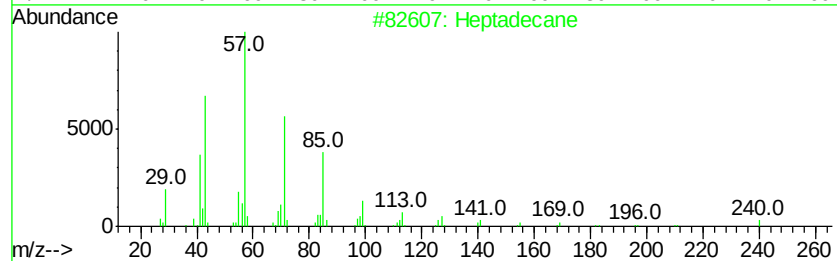
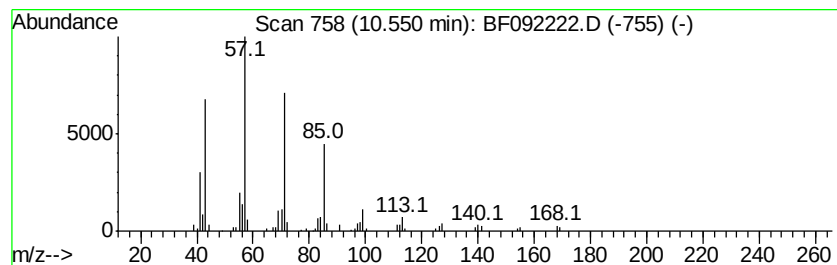
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.55	2.25 ng	197910	Phenanthrene-d10		11.11
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Octadecane	254	C18H38	000593-45-3	91
3	Hexadecane	226	C16H34	000544-76-3	90
4	Pentadecane	212	C15H32	000629-62-9	87
5	Heptacosane	380	C27H56	000593-49-7	86



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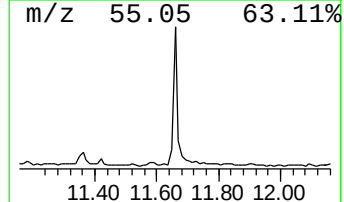
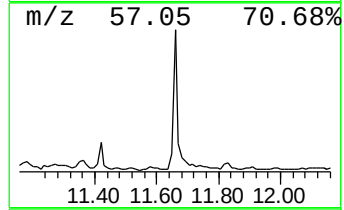
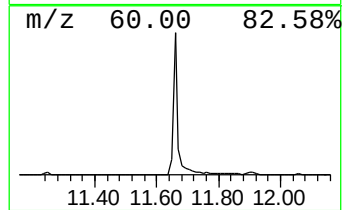
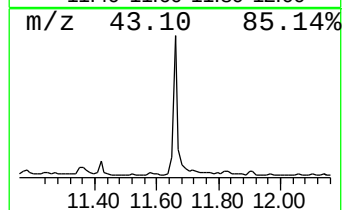
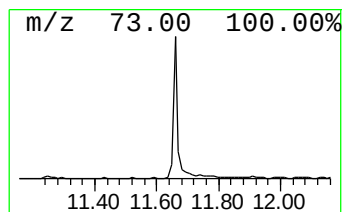
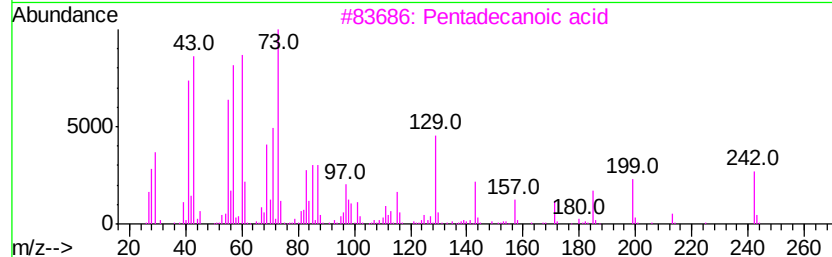
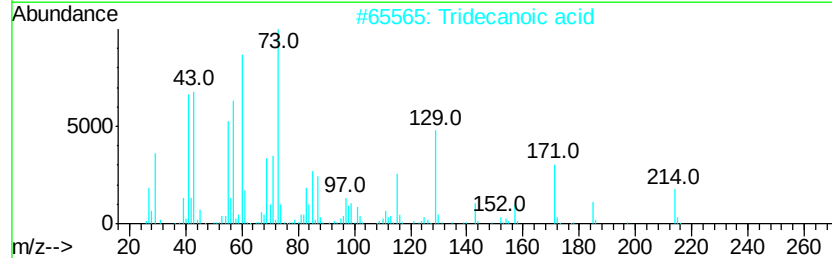
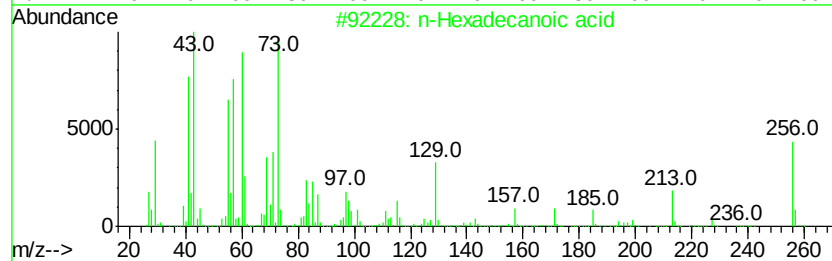
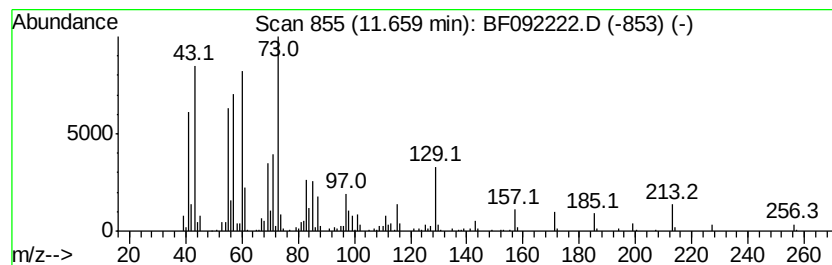
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.66	5.61 ng	493173	Phenanthrene-d10		11.11
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	49
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	11
5	Nonanoic acid	158	C9H18O2	000112-05-0	11



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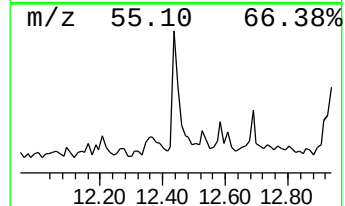
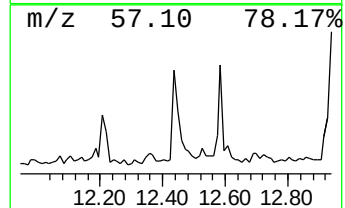
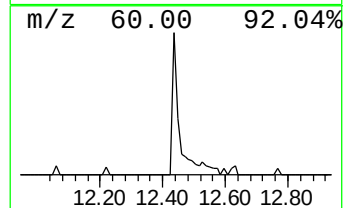
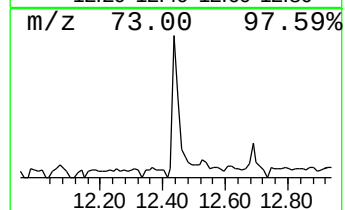
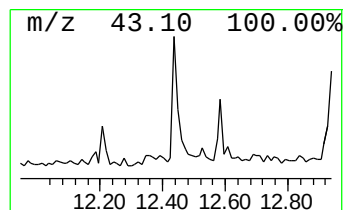
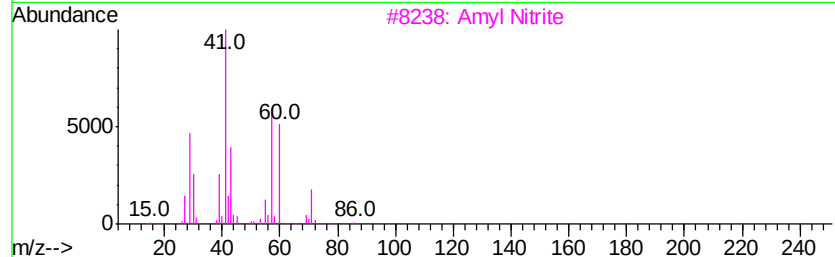
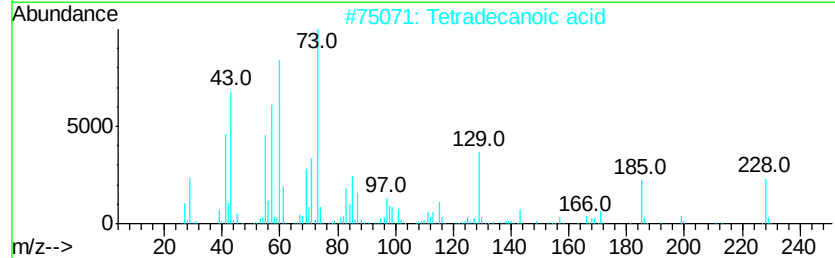
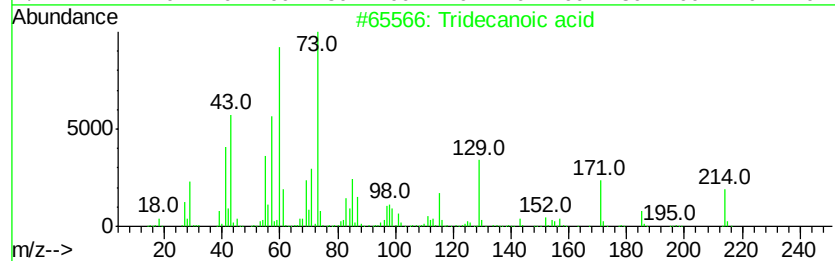
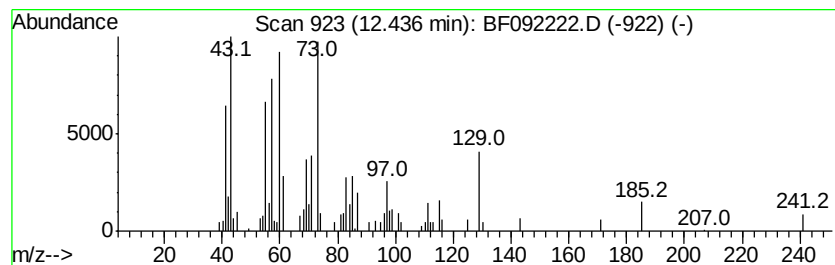
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.44	2.17 ng	113246	Chrysene-d12	13.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	86
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3	Amyl Nitrite	117	C5H11NO2	000110-46-3	27
4	Nonahexacontanoic acid	999	C69H138O2	040710-32-5	18
5	5-Acetoxypentadecane	270	C17H34O2	1000245-62-3	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
Data File : BF092222.D
Acq On : 12 Jan 2017 22:05
Operator : UM/SJ
Sample : I1030-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

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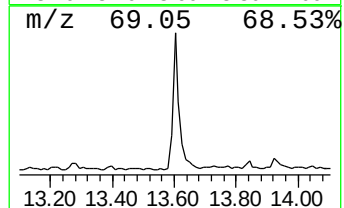
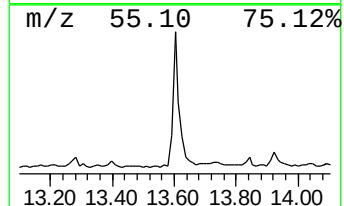
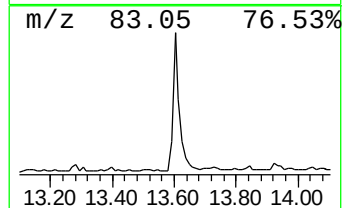
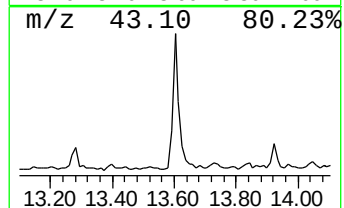
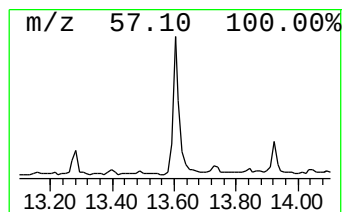
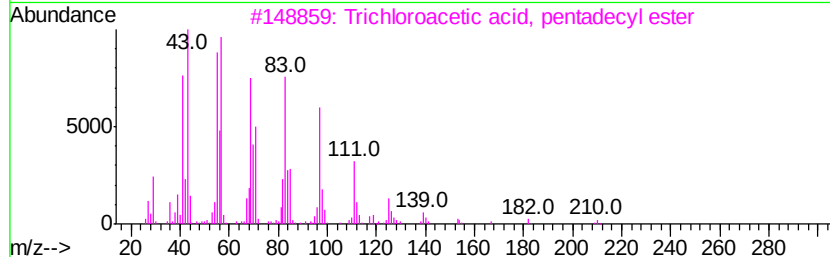
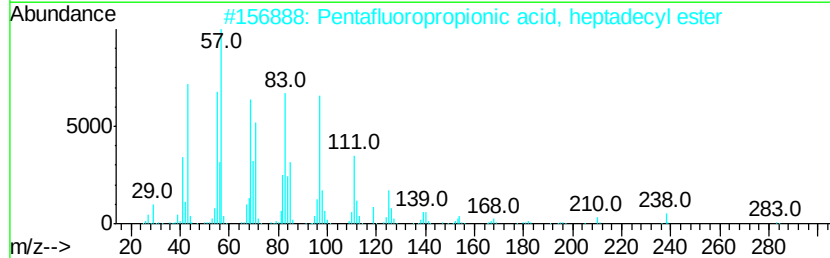
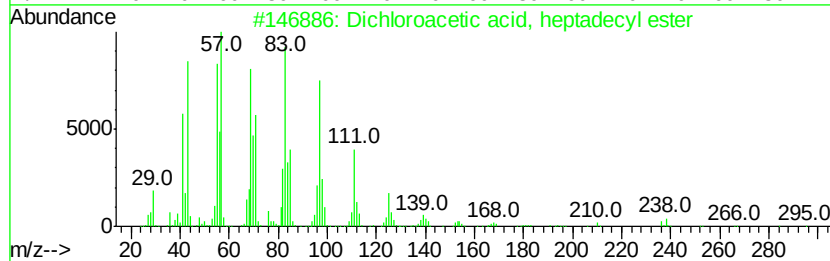
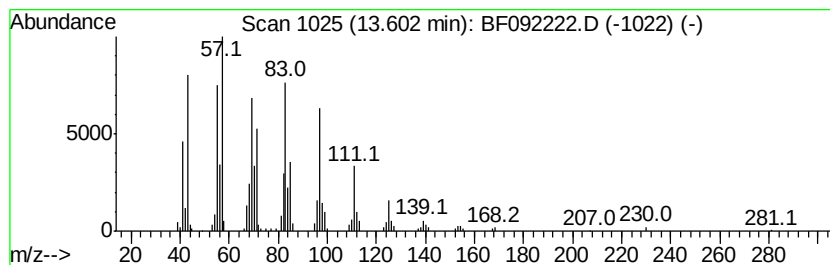
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dichloroacetic acid, heptad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	7.72 ng	403167	Chrysene-d12	13.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	91
3	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
4	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	87
5	Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	87



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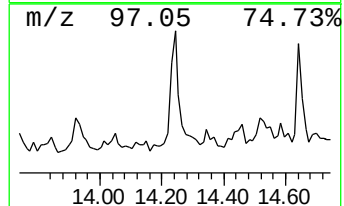
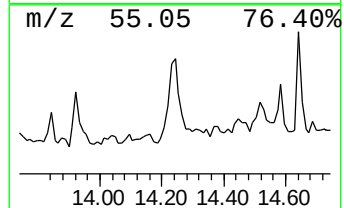
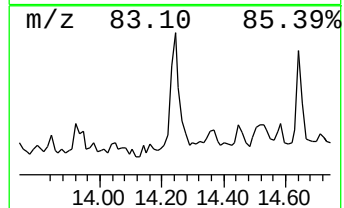
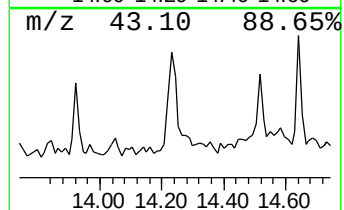
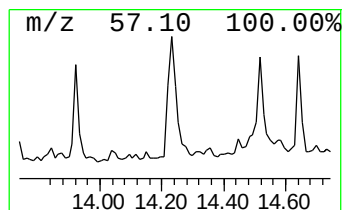
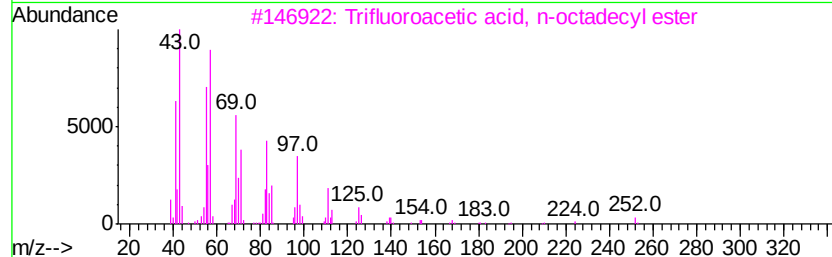
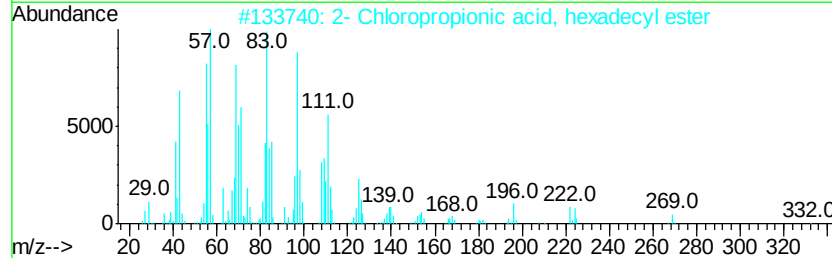
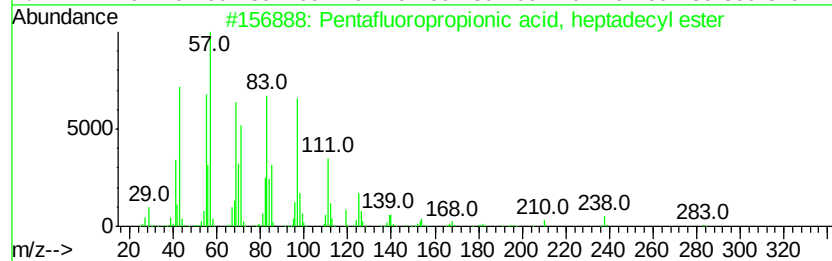
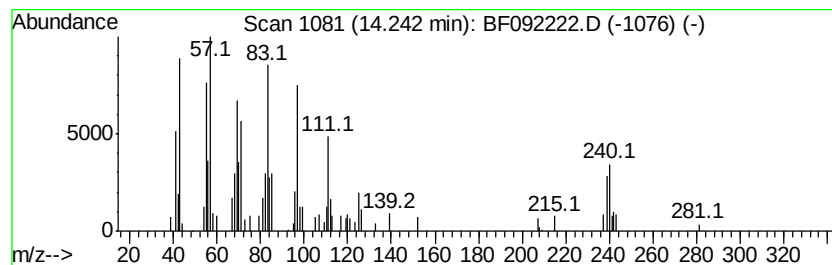
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentafluoropropionic acid, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	2.71 ng	141346	Chrysene-d12	13.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87
2			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	83
3			Trifluoroacetic acid, n-octadecv...	366	C20H37F3O2	079392-43-1	83
4			11-Tricosene	322	C23H46	052078-56-5	81
5			1-Eicosene	280	C20H40	003452-07-1	72



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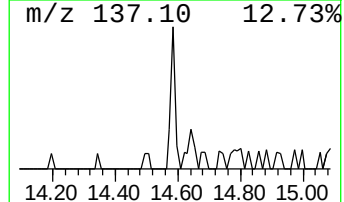
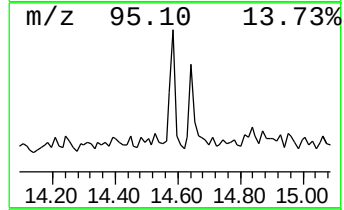
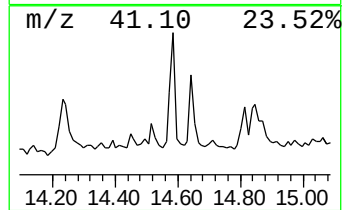
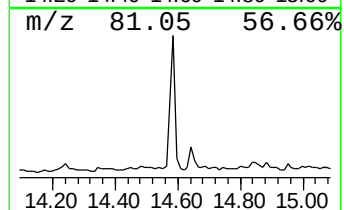
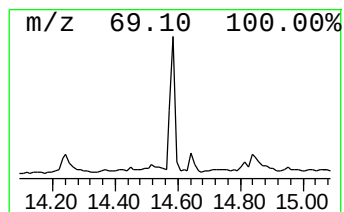
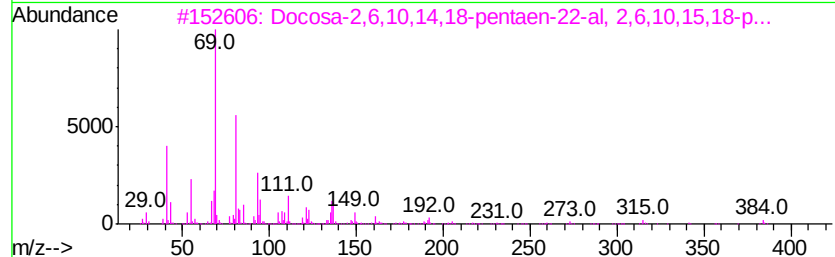
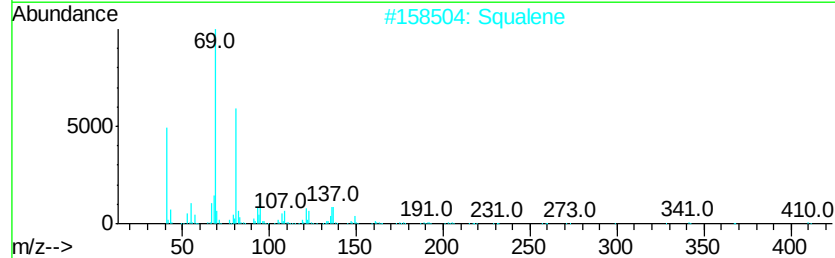
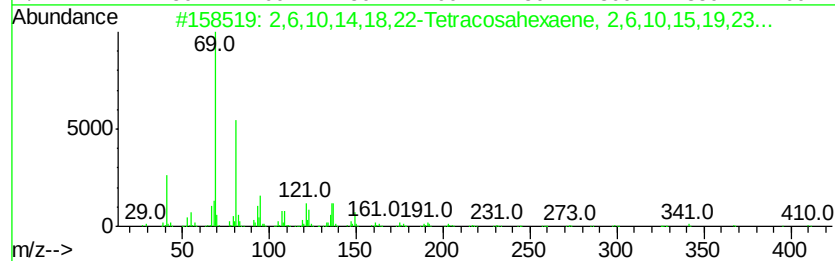
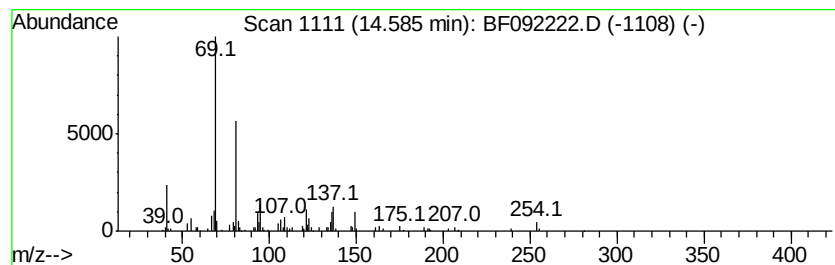
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2,6,10,14,18,22-Tetracosahexaene... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	2.31 ng	101496	Perylene-d12	15.09

Hit#	of	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	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	78	
4	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	74	
5	1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	64	



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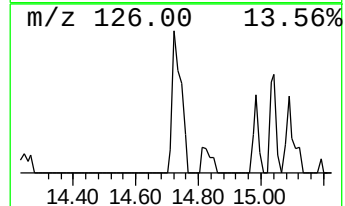
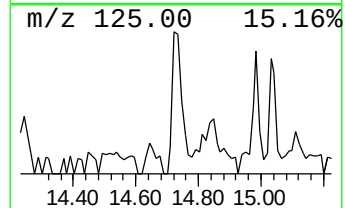
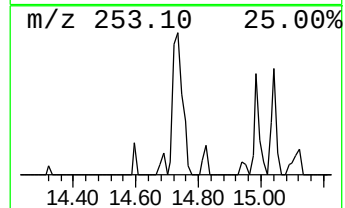
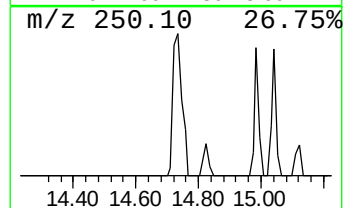
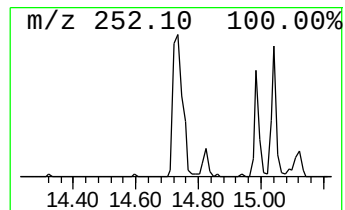
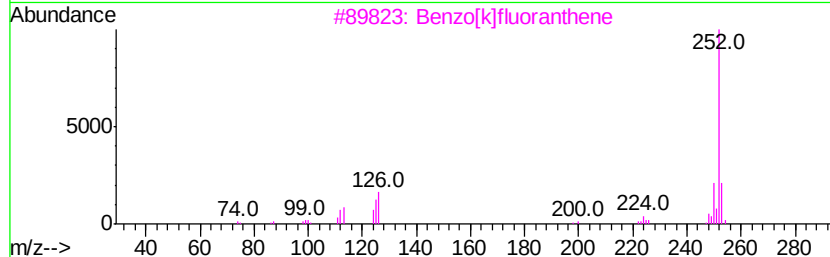
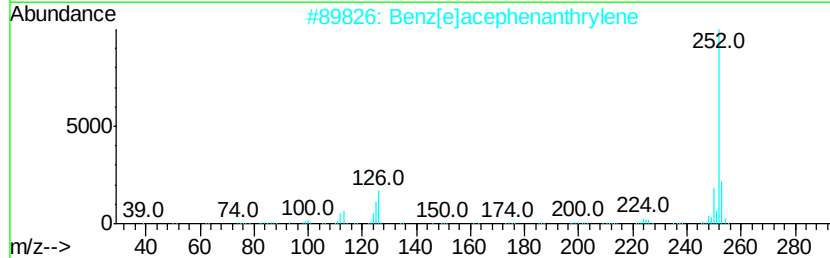
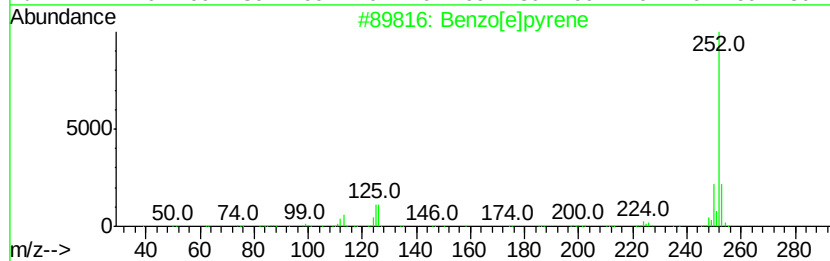
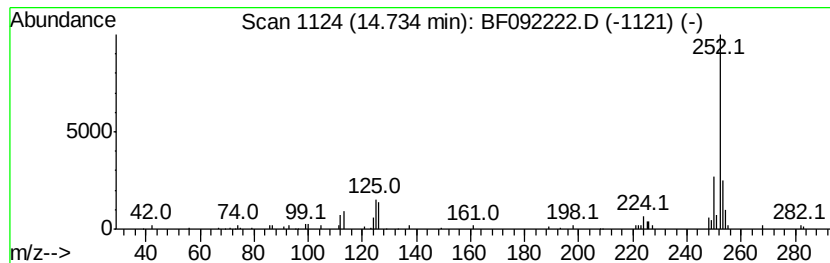
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.73	2.53 ng	110820	Perylene-d12	15.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	93
2	Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
4	Benzo[a]pyrene	252	C20H12	000050-32-8	81
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
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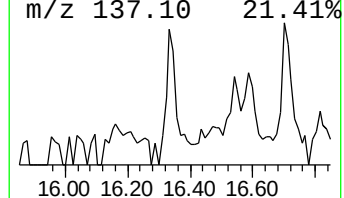
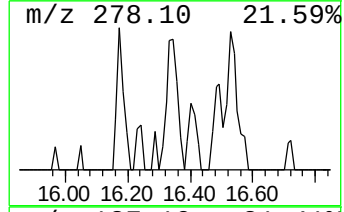
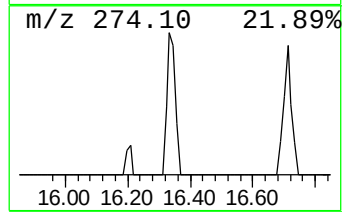
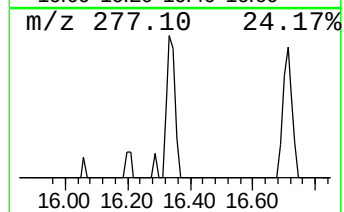
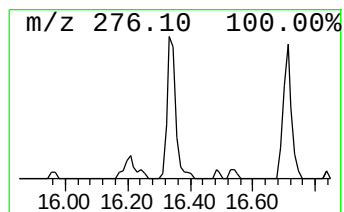
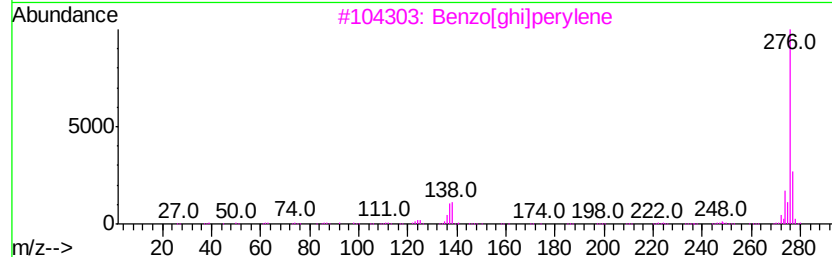
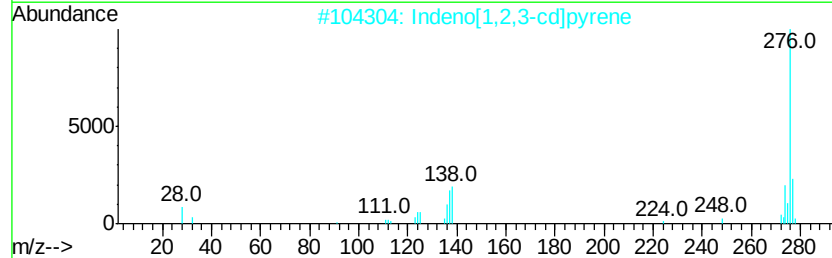
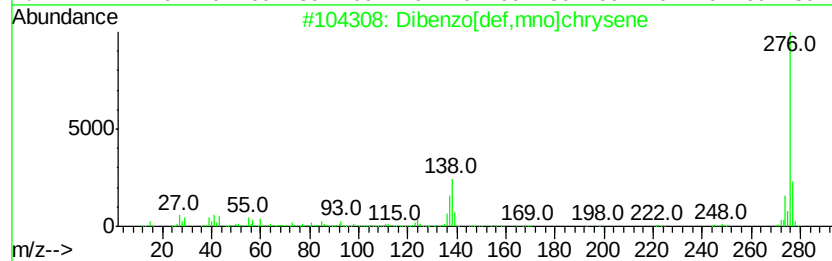
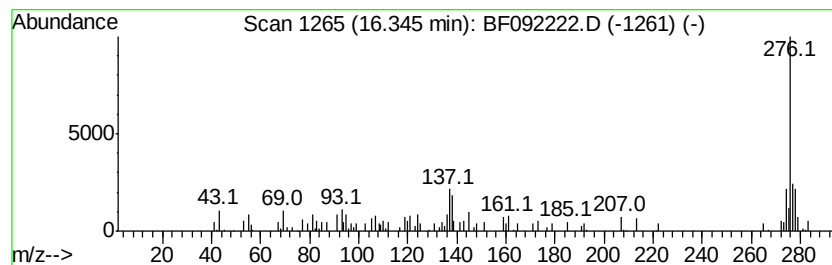
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Dibenzo[def,mno]chrysene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.35	2.39 ng	104662	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	92
2		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	89
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	76
4		Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	64
5		2-Amino-3-cyano-4,5-diphenylthio...	276	C17H12N2S	042160-34-9	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
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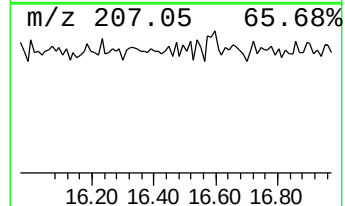
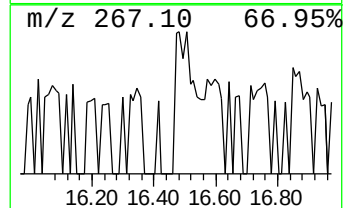
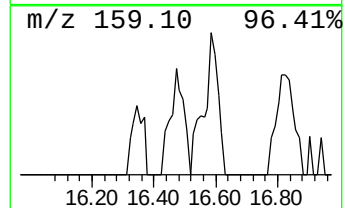
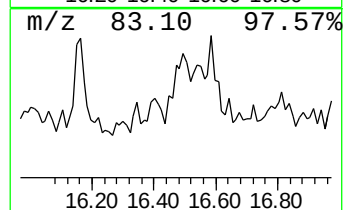
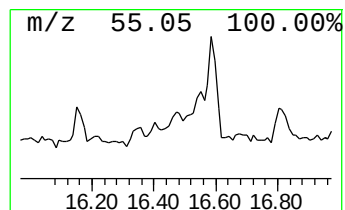
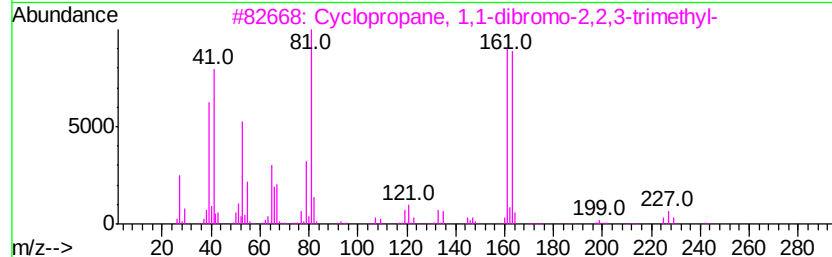
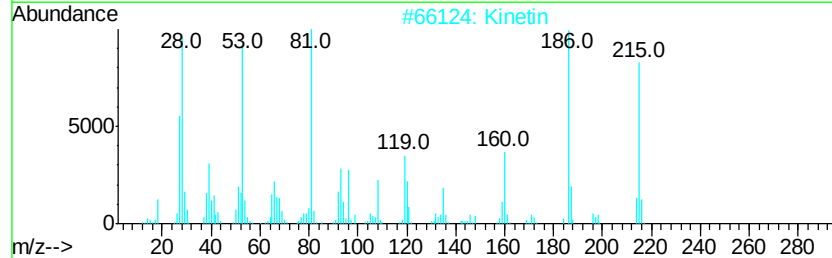
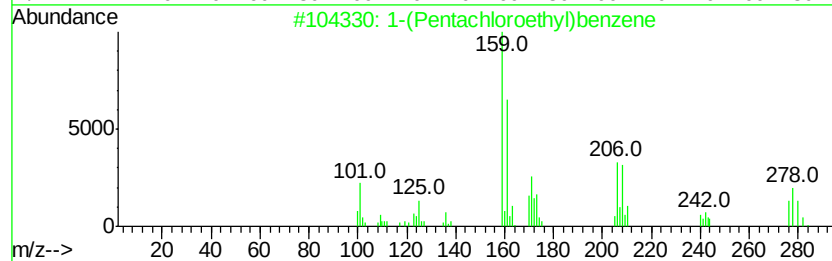
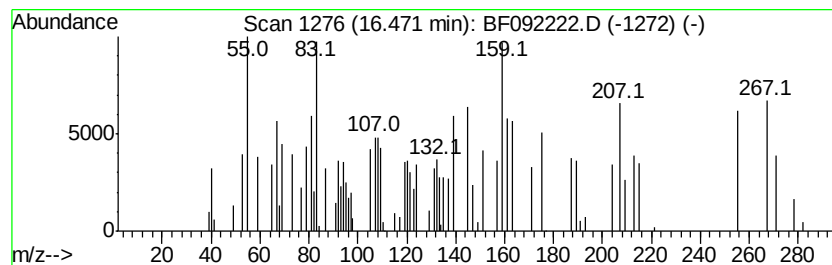
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown16.47 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	2.75 ng	120581	Perylene-d12	15.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(Pentachloroethyl)benzene	276	C8H5Cl5	000706-93-4	14
2			Kinetin	215	C10H9N5O	000525-79-1	14
3			Cyclopropane, 1,1-dibromo-2,2,3-...	240	C6H10Br2	021960-71-4	12
4			Thiazolidin-4-one, 3-furfuryl-2-...	213	C8H7N02S2	004703-95-1	12
5			Pyrrole-2,5-dione, 1-(3-chloro-2...	267	C12H10ClN04	1000273-93-8	12



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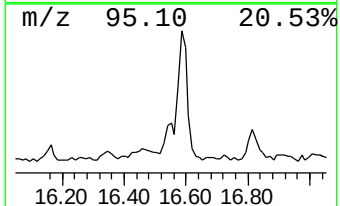
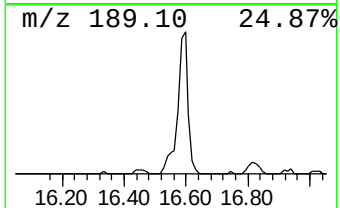
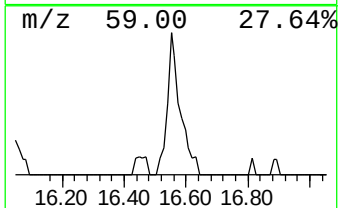
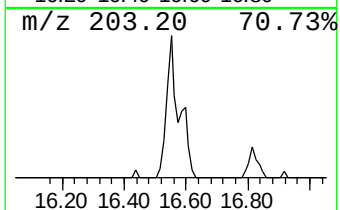
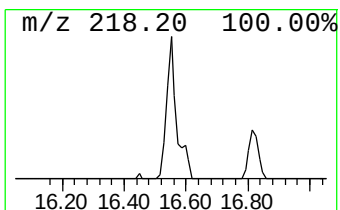
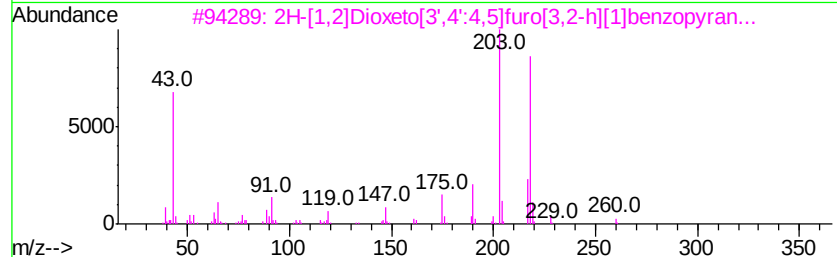
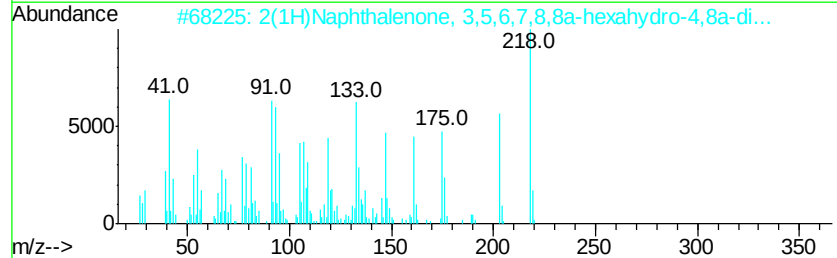
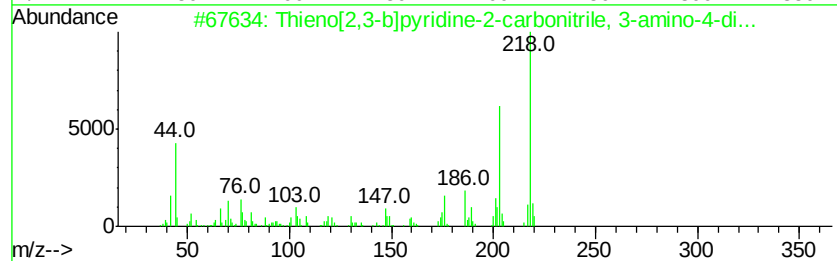
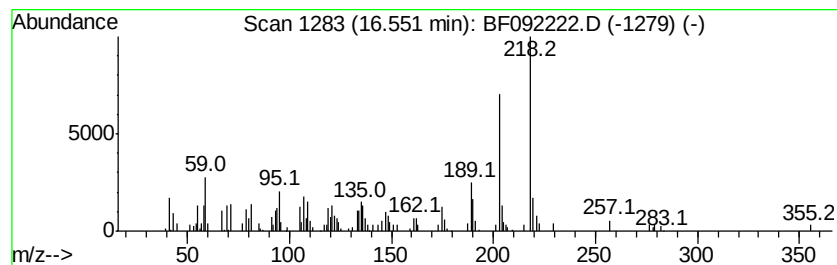
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Thieno[2,3-b]pyridine-2-car... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	5.27 ng	231122	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thieno[2,3-b]pyridine-2-carbonit...	218	C10H10N4S	147992-88-9	53
2		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	53
3		2H-[1,2]Dioxeto[3',4':4,5]furo[3...	260	C14H12O5	129833-00-7	53
4		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	52
5		Pyrene, hexadecahydro-	218	C16H26	002435-85-0	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011217\
Data File : BF092222.D
Acq On : 12 Jan 2017 22:05
Operator : UM/SJ
Sample : I1030-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

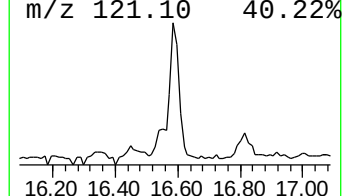
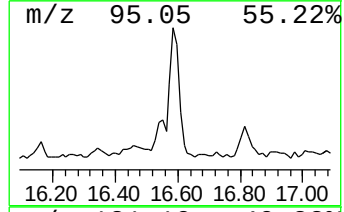
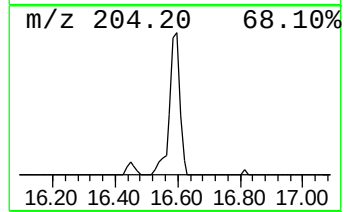
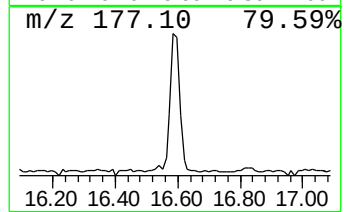
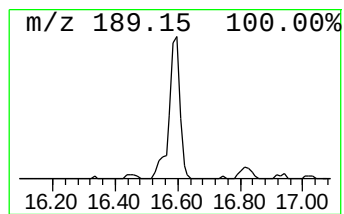
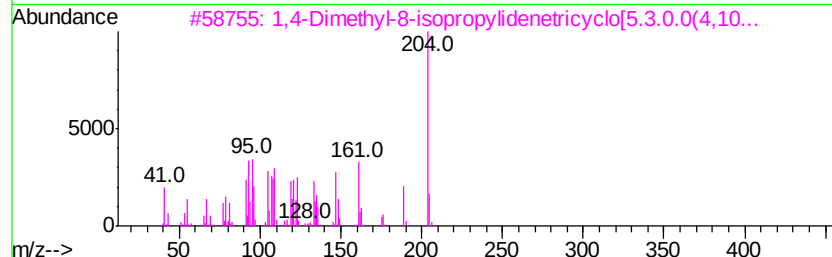
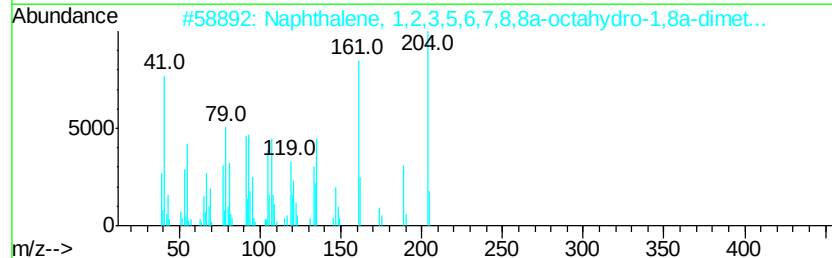
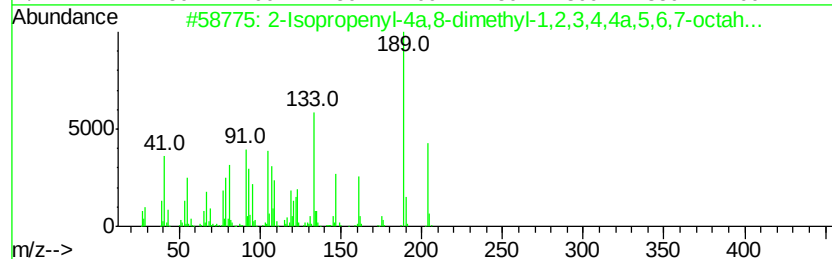
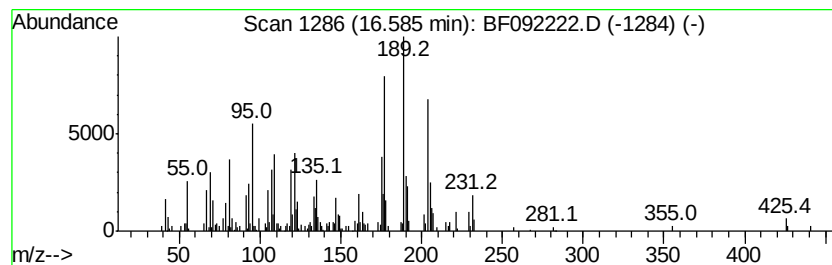
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.59 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	11.50 ng	504548	Perylene-d12	15.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Isopropenyl-4a,8-dimethyl-1,2,...	204 C15H24	1000192-43-5	38
2	Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204 C15H24	004630-07-3	25
3	1,4-Dimethyl-8-isopropylidenetri...	204 C15H24	1000140-07-7	25
4	1,3a-Ethano-3aH-indene, 1,2,3,6,...	204 C15H24	004545-68-0	25
5	3H-1,2-Dithiole-3-thione, 5-tert...	204 C8H12S3	013120-76-8	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011217\
Data File : BF092222.D
Acq On : 12 Jan 2017 22:05
Operator : UM/SJ
Sample : I1030-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-0017

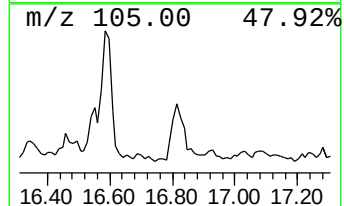
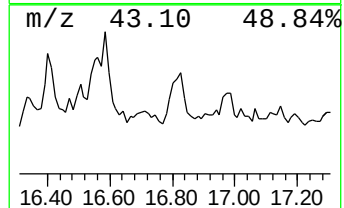
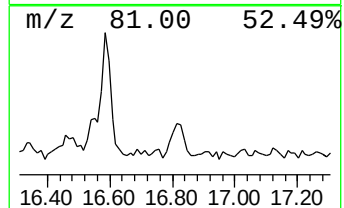
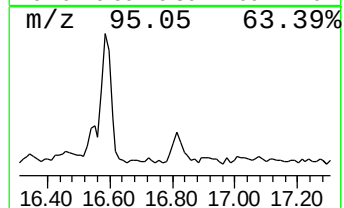
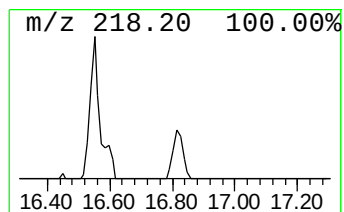
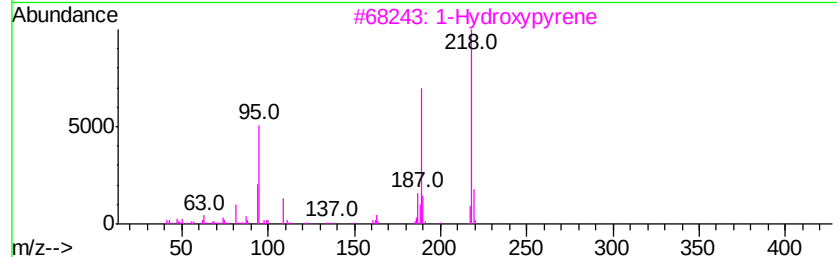
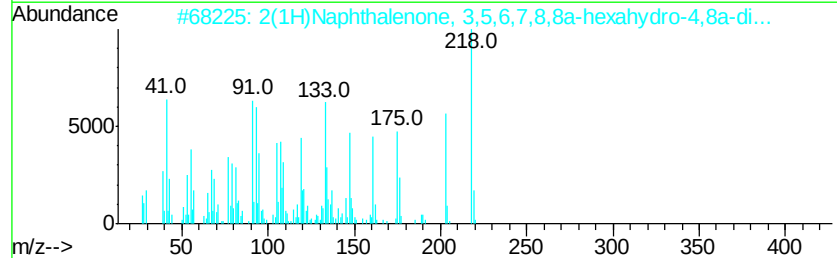
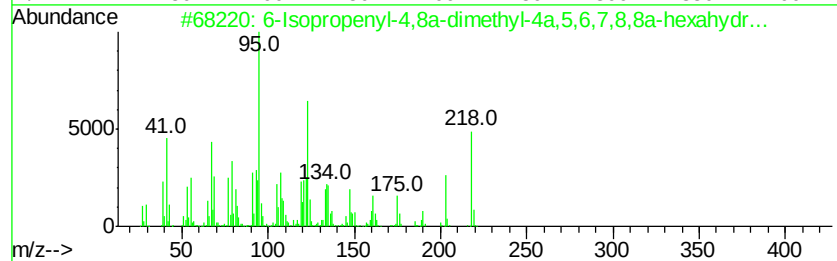
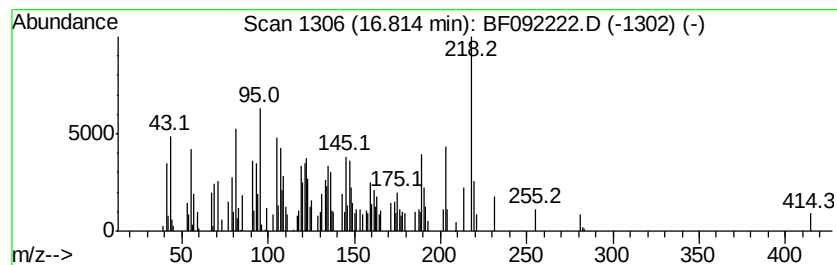
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 6-Isopropenyl-4,8a-dimethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.81	3.76 ng	165102	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	50
2		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	46
3		1-Hydroxypyrene	218	C16H10O	005315-79-7	38
4		Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	38
5		9-Phenyl-5H-benzocycloheptene	218	C17H14	138089-71-1	35



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 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-0017

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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	37.7	ng	2127180	1	6.60	1129780	20.0
unknown6.37	6.37	71.3	ng	4027740	1	6.60	1129780	20.0
Heptadecane	10.55	2.3	ng	197910	4	11.11	1758830	20.0
n-Hexadecanoic acid	11.66	5.6	ng	493173	4	11.11	1758830	20.0
Tridecanoic acid	12.44	2.2	ng	113246	5	13.73	1043890	20.0
Dichloroacetic ac...	13.60	7.7	ng	403167	5	13.73	1043890	20.0
Pentafluoropropio...	14.24	2.7	ng	141346	5	13.73	1043890	20.0
2,6,10,14,18,22-T...	14.59	2.3	ng	101496	6	15.09	877172	20.0
Benzo[e]pyrene	14.73	2.5	ng	110820	6	15.09	877172	20.0
Dibenzo[def,mno]c...	16.35	2.4	ng	104662	6	15.09	877172	20.0
unknown16.47	16.47	2.8	ng	120581	6	15.09	877172	20.0
Thieno[2,3-b]pyri...	16.55	5.3	ng	231122	6	15.09	877172	20.0
unknown16.59	16.59	11.5	ng	504548	6	15.09	877172	20.0
6-Isopropenyl-4,8...	16.81	3.8	ng	165102	6	15.09	877172	20.0