

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
 Data File : BF081034.D
 Acq On : 18 Aug 2015 7:25
 Operator : UM/IZ
 Sample : G3290-11
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2A(0-2)-5

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-BF081715.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	4.033	177	179	183	rBV	11660	18250	1.12%	0.124%
2	4.753	239	242	252	rBV	580628	1019633	62.81%	6.947%
3	5.210	279	282	284	rBV	1522108	1623269	100.00%	11.059%
4	6.273	372	375	378	rBV	1360929	1416561	87.27%	9.651%
5	6.399	383	386	388	rBV	1288318	1563966	96.35%	10.655%
6	6.627	403	406	408	rBV	454604	373622	23.02%	2.545%
7	6.787	417	420	422	rBV	859968	961234	59.22%	6.549%
8	7.199	453	456	458	rBV	1062794	1092521	67.30%	7.443%
9	7.907	516	518	521	rBV	334638	437232	26.94%	2.979%
10	8.993	610	613	615	rBV	1452132	1214674	74.83%	8.275%
11	9.393	646	648	650	rBV	214436	210839	12.99%	1.436%
12	9.656	669	671	673	rBV	481238	433469	26.70%	2.953%
13	10.445	737	740	742	rBV	772248	765695	47.17%	5.217%
14	10.582	748	752	755	rBV	36024	45028	2.77%	0.307%
15	10.673	758	760	761	rVV2	14531	18871	1.16%	0.129%
16	10.708	761	763	766	rVB3	9377	18139	1.12%	0.124%
17	11.028	787	791	792	rBV	30814	30511	1.88%	0.208%
18	11.131	798	800	802	rBV	461993	410889	25.31%	2.799%
19	11.679	847	848	852	rVV	88230	108356	6.68%	0.738%
20	11.736	852	853	859	rVB4	9161	18386	1.13%	0.125%
21	12.331	904	905	907	rBV	62983	55983	3.45%	0.381%
22	12.559	922	925	927	rBV	72532	70650	4.35%	0.481%
23	12.605	927	929	934	rVB4	15056	25351	1.56%	0.173%
24	12.719	936	939	941	rVB	961069	918417	56.58%	6.257%
25	12.948	956	959	962	rBV3	11375	28154	1.73%	0.192%
26	13.199	978	981	983	rBV	130479	123677	7.62%	0.843%
27	13.576	1011	1014	1016	rBV	9317	17879	1.10%	0.122%
28	13.634	1016	1019	1026	rVV	145029	233016	14.35%	1.588%
29	13.748	1027	1029	1035	rVB	392941	485853	29.93%	3.310%
30	13.874	1036	1040	1042	rBV2	9204	21154	1.30%	0.144%
31	13.954	1045	1047	1050	rVB2	10745	17825	1.10%	0.121%
32	14.262	1071	1074	1078	rBV	47858	92687	5.71%	0.631%
33	14.548	1097	1099	1101	rBV3	12657	17074	1.05%	0.116%
34	14.605	1103	1104	1106	rVB	30353	28206	1.74%	0.192%

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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

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35	14.674	1108	1110	1112	rBV2	15458	19791	1.22%	0.135%
36	14.742	1113	1116	1121	rVB	27431	66009	4.07%	0.450%
37	14.834	1122	1124	1126	rBV	30843	45973	2.83%	0.313%
38	14.994	1136	1138	1141	rVB	26023	36511	2.25%	0.249%
39	15.051	1141	1143	1145	rBV	22591	34942	2.15%	0.238%
40	15.108	1145	1148	1150	rBV	196568	251166	15.47%	1.711%
41	15.337	1166	1168	1172	rVB3	22655	33273	2.05%	0.227%
42	15.531	1183	1185	1187	rBV	34354	52254	3.22%	0.356%
43	15.588	1187	1190	1192	rVV	25422	50846	3.13%	0.346%
44	16.068	1230	1232	1237	rVB5	17833	38436	2.37%	0.262%
45	16.194	1241	1243	1248	rVB2	22411	39108	2.41%	0.266%
46	16.845	1296	1300	1307	rBV8	17164	57301	3.53%	0.390%
47	17.691	1370	1374	1379	rBV8	10749	35411	2.18%	0.241%

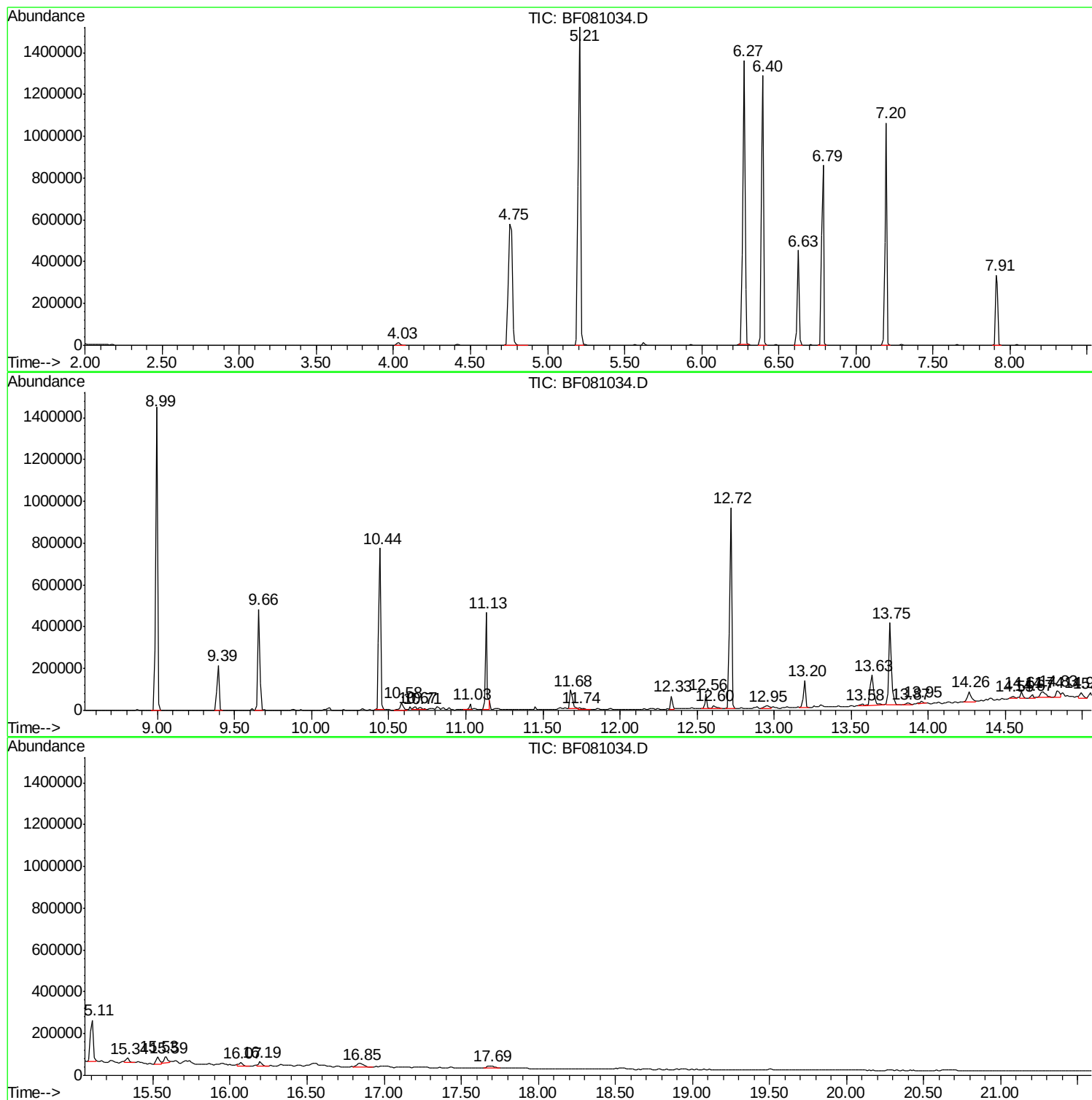
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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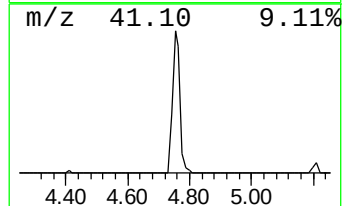
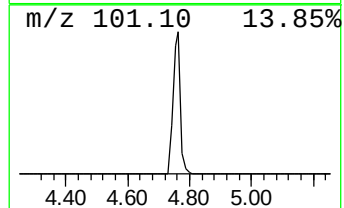
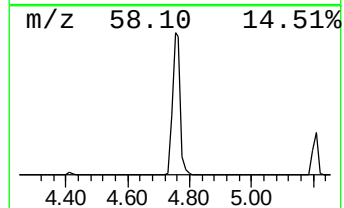
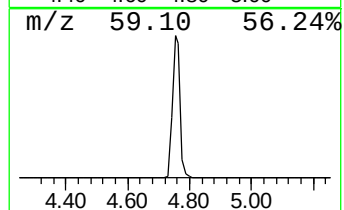
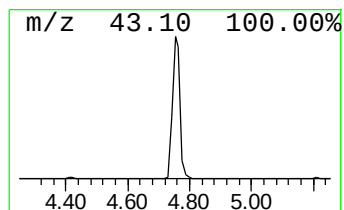
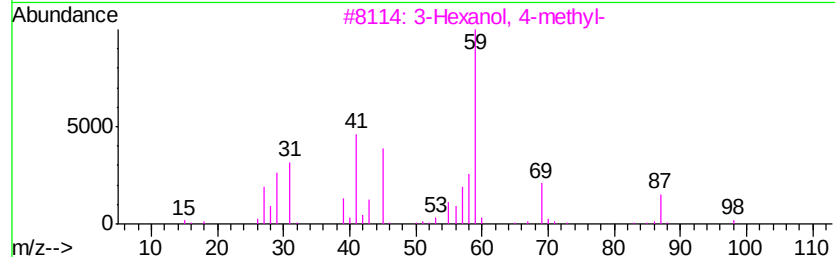
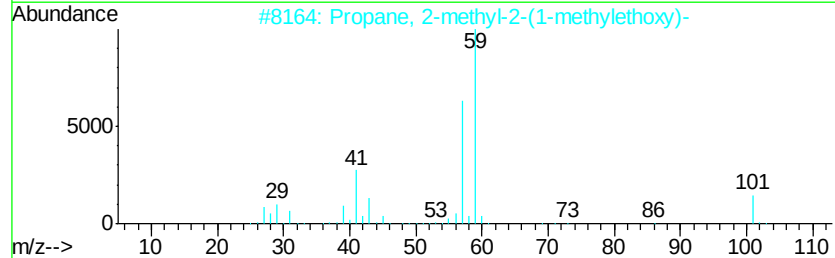
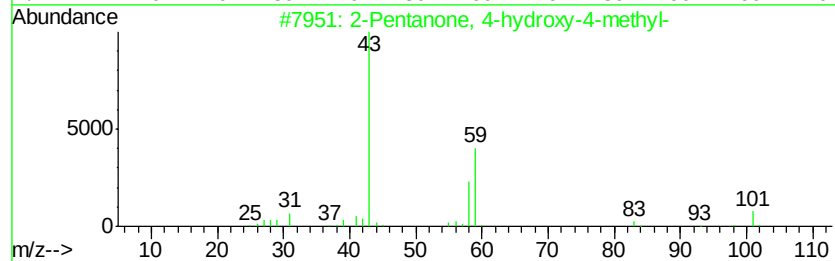
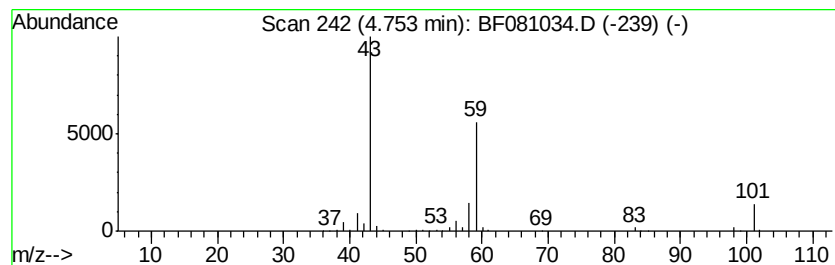
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.75	54.58 ng	1019630	1,4-Dichlorobenzene-d4	6.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42	
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
4	3-Hexanol	102	C6H14O	000623-37-0	33	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32	



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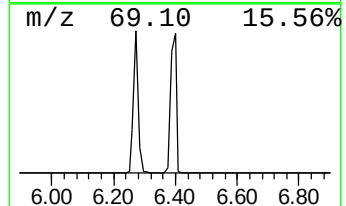
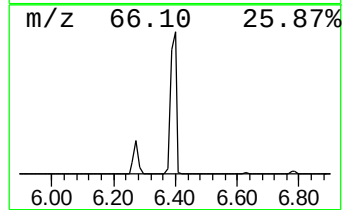
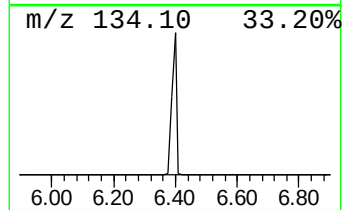
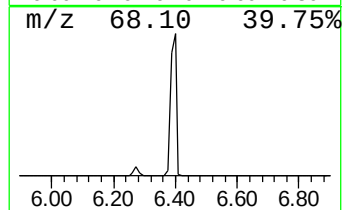
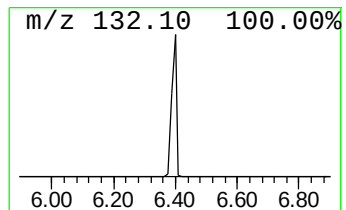
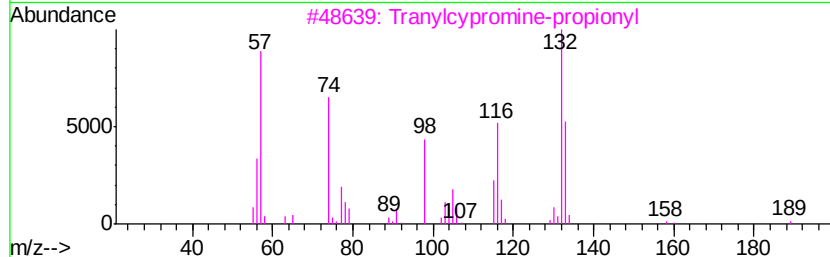
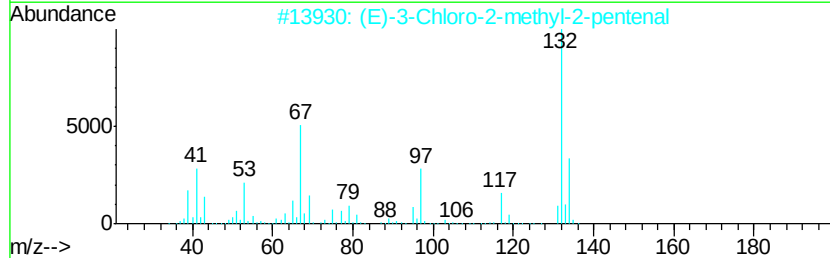
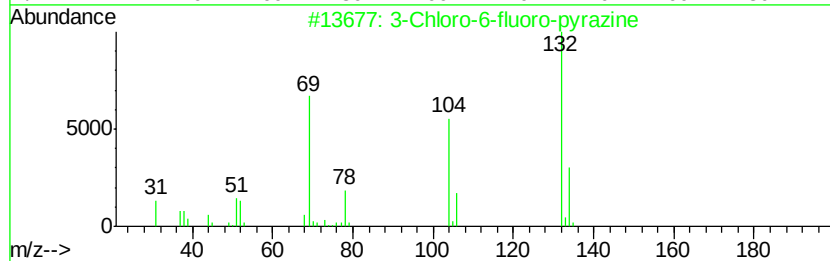
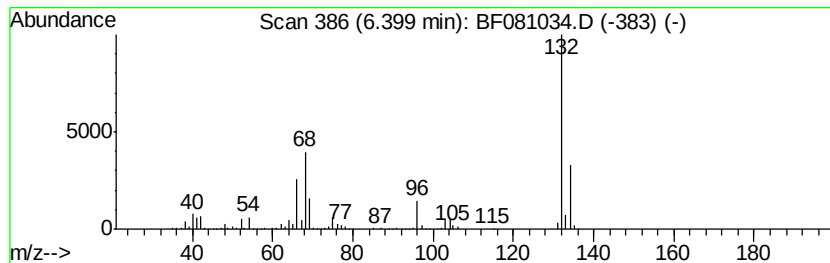
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	83.72 ng	1563970	1,4-Dichlorobenzene-d4	6.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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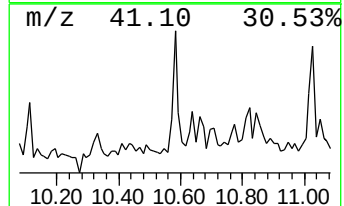
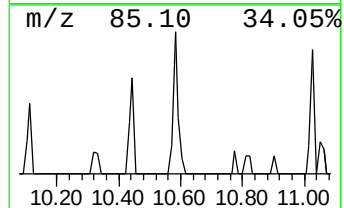
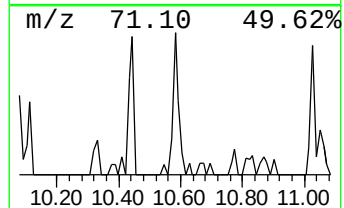
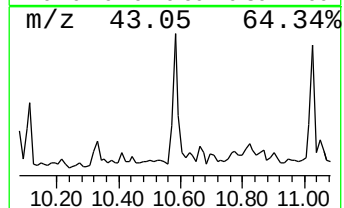
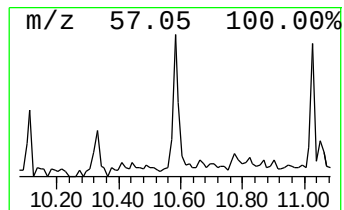
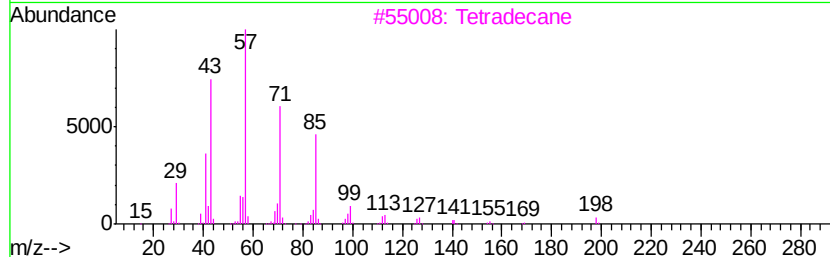
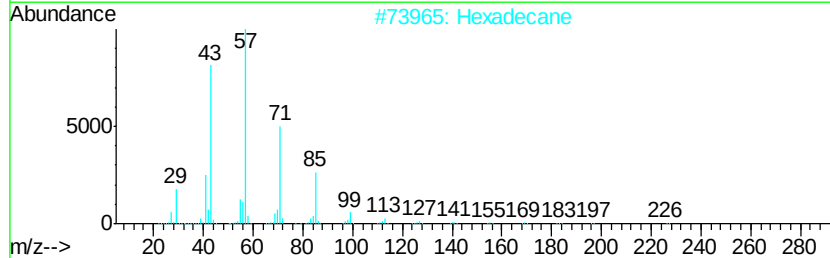
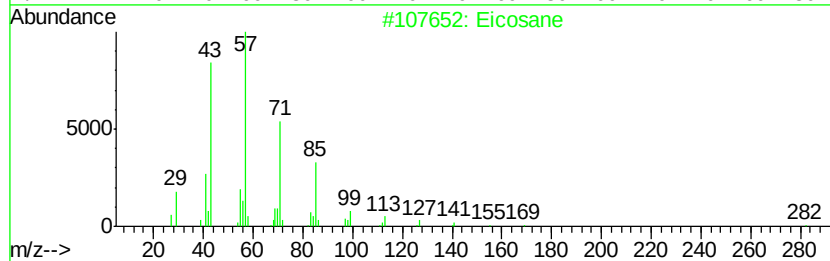
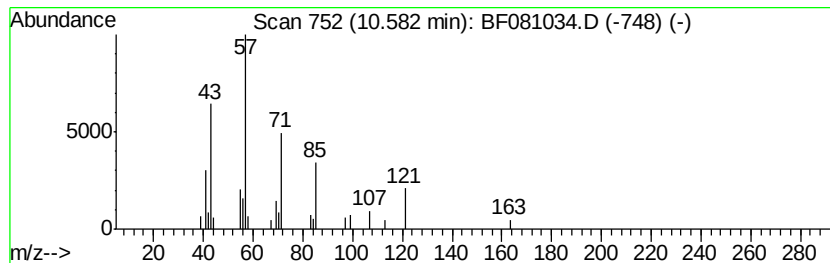
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Eicosane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	2.19 ng	45028	Phenanthrene-d10	11.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	72
2		Hexadecane	226	C16H34	000544-76-3	64
3		Tetradecane	198	C14H30	000629-59-4	64
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	64
5		Pentadecane	212	C15H32	000629-62-9	64



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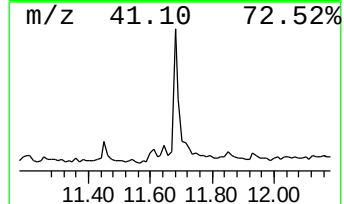
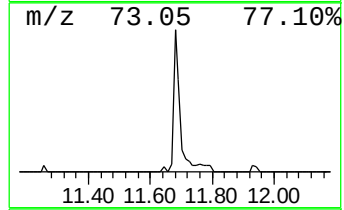
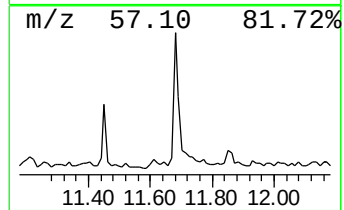
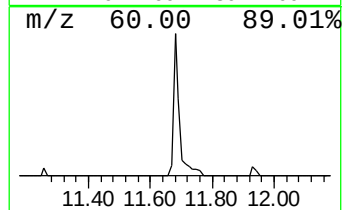
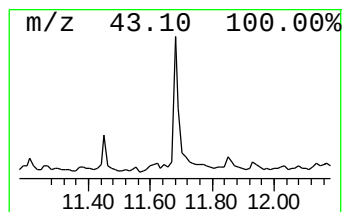
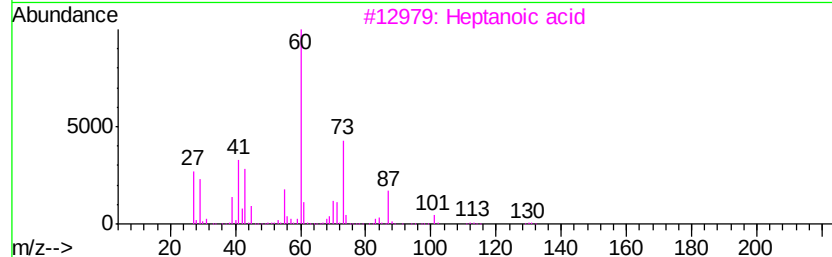
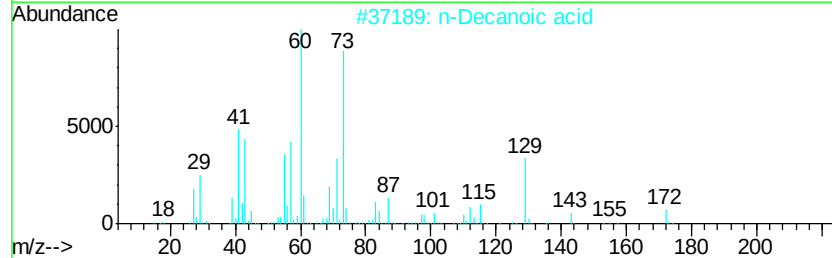
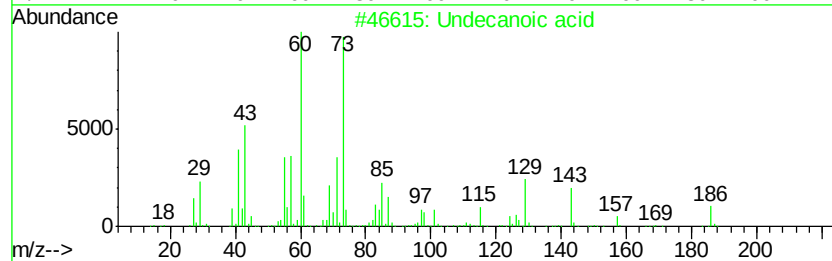
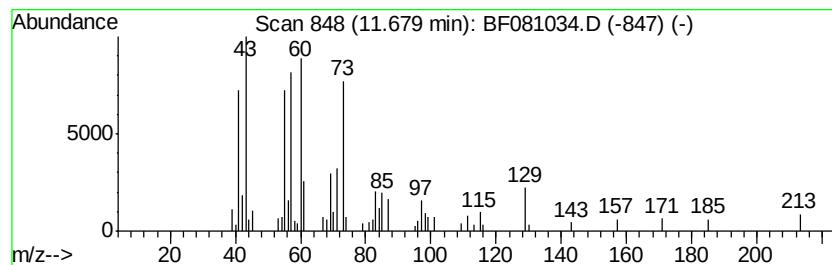
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Undecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.68	5.27 ng	108356	Phenanthrene-d10	11.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecanoic acid		186	C11H22O2	000112-37-8	80
2	n-Decanoic acid		172	C10H20O2	000334-48-5	50
3	Heptanoic acid		130	C7H14O2	000111-14-8	43
4	.beta.-D-Mannofuranoside, 1-O-(1...		332	C17H32O6	1000155-29-9	22
5	5-Acetoxypentadecane		270	C17H34O2	1000245-62-3	14



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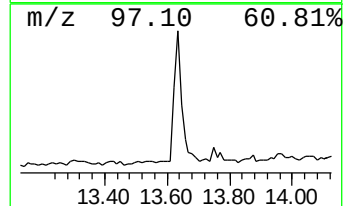
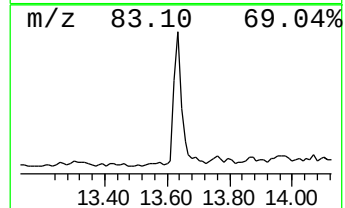
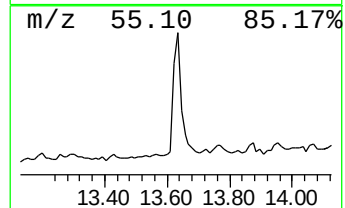
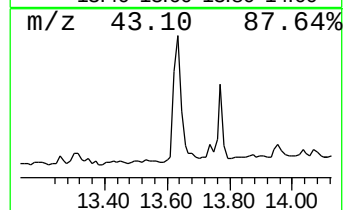
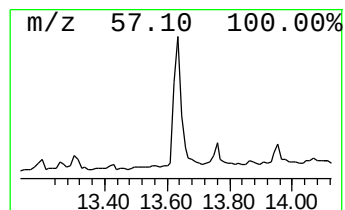
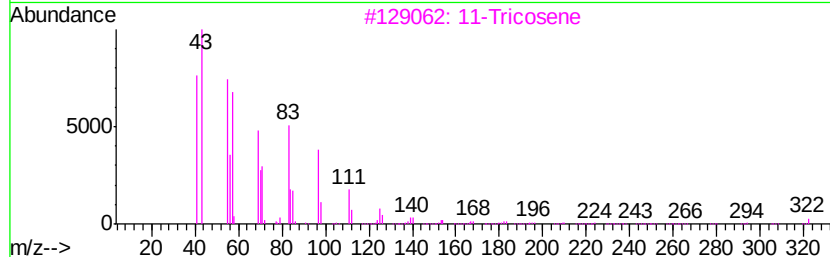
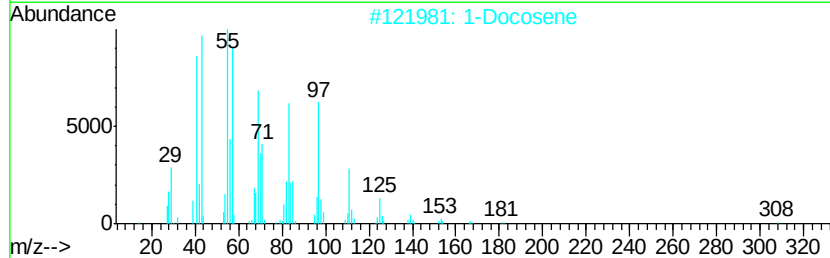
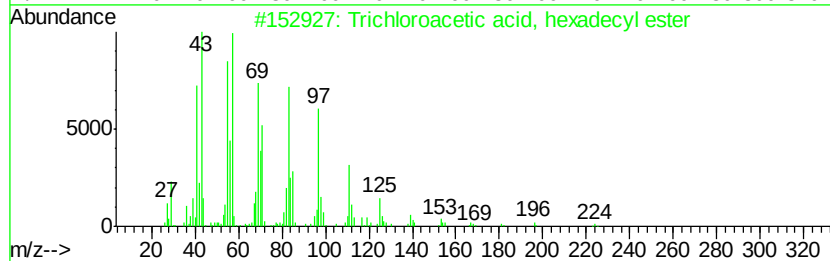
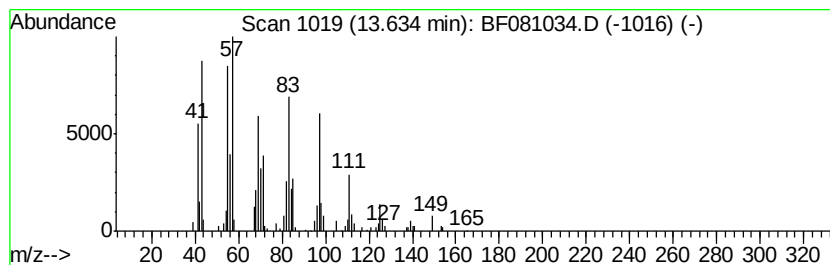
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Trichloroacetic acid, hexadecyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	9.59 ng	233016	Chrysene-d12	13.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	95
2			1-Docosene	308	C22H44	001599-67-3	91
3			11-Tricosene	322	C23H46	052078-56-5	91
4			1-Nonadecene	266	C19H38	018435-45-5	91
5			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91



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Data File : BF081034.D
Acq On : 18 Aug 2015 7:25
Operator : UM/IZ
Sample : G3290-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

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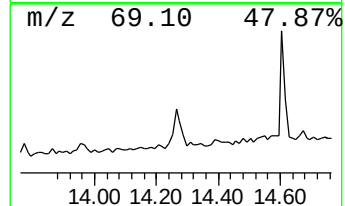
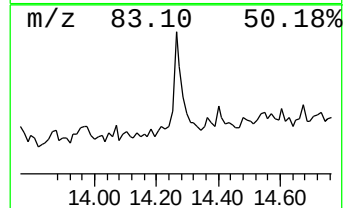
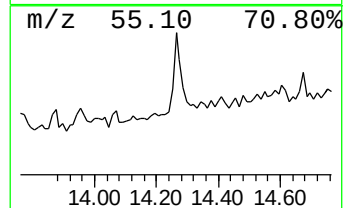
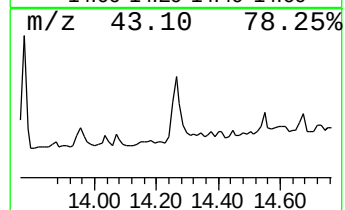
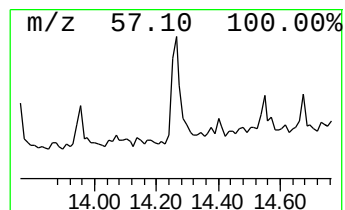
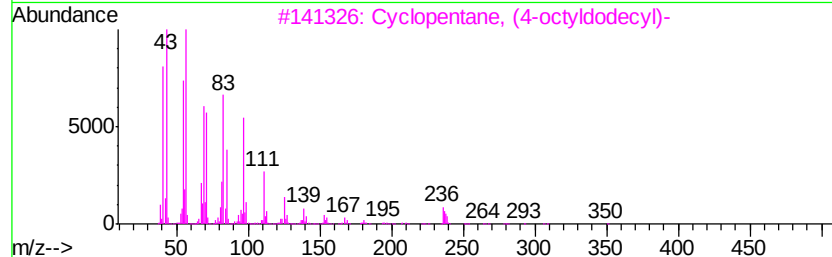
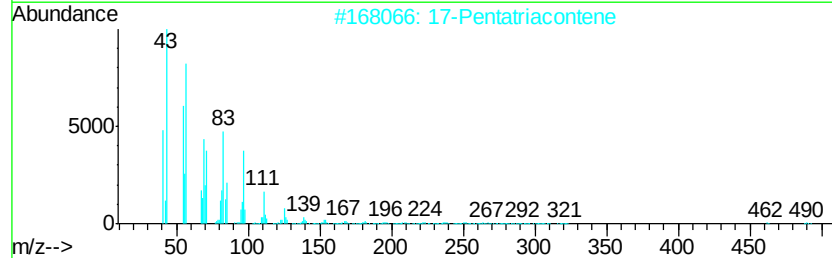
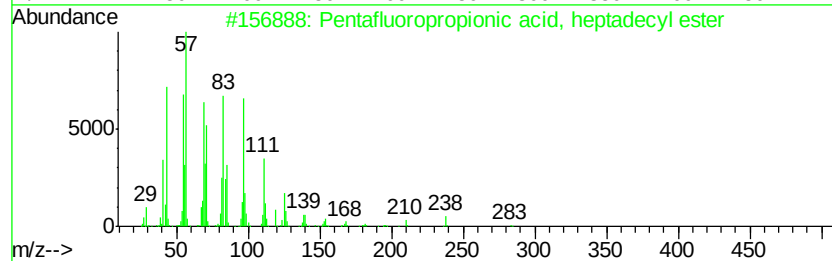
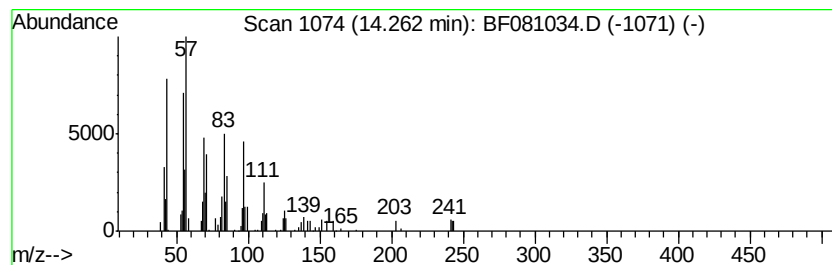
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentafluoropropionic acid, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	3.82 ng	92687	Chrysene-d12	13.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	90
2		17-Pentatriacontene	491	C35H70	006971-40-0	90
3		Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	83
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	83
5		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
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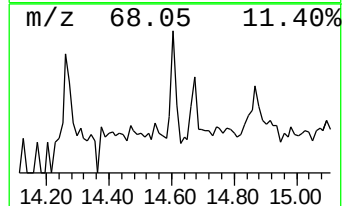
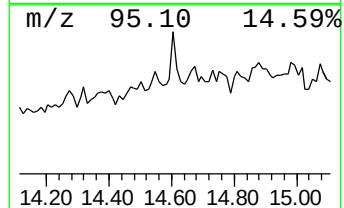
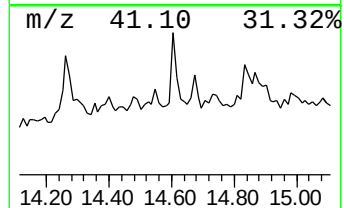
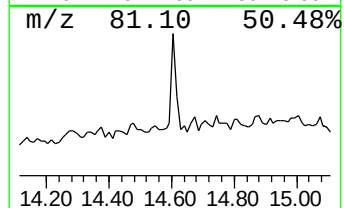
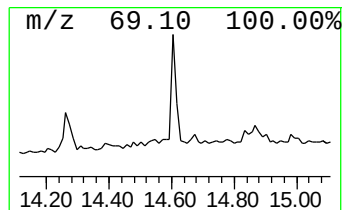
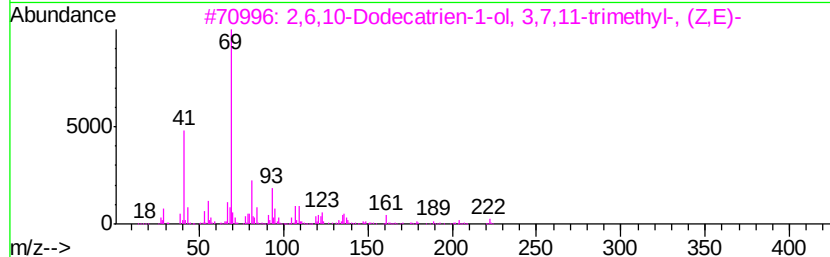
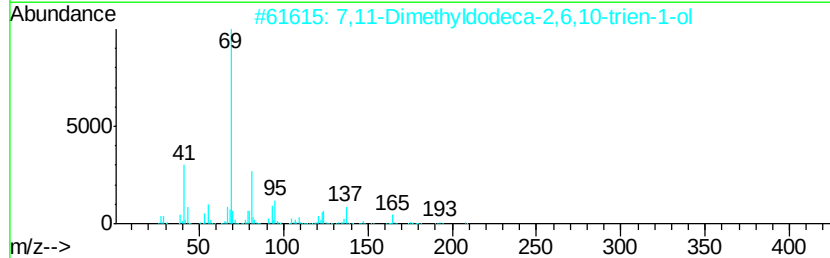
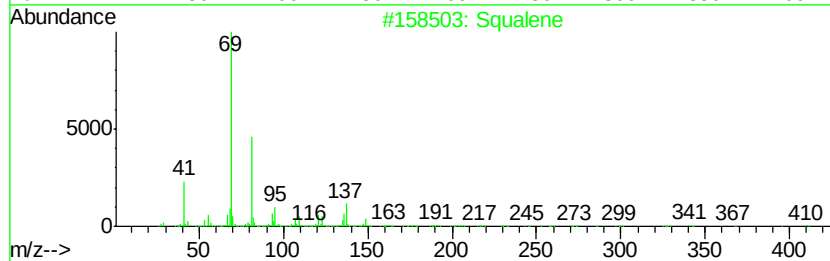
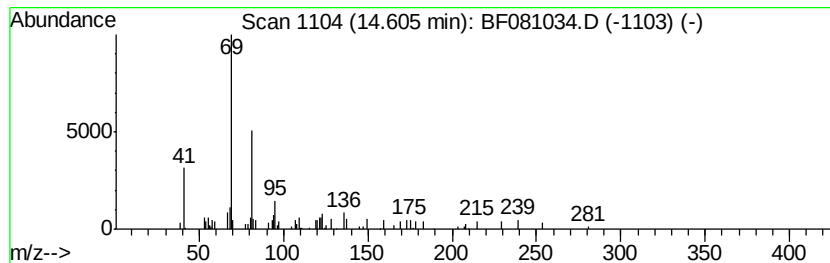
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Squalene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	2.25 ng	28206	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	58
2		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	53
3		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	003790-71-4	53
4		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	42
5		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	42



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Operator : UM/IZ
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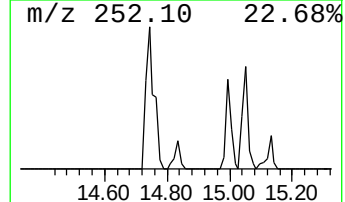
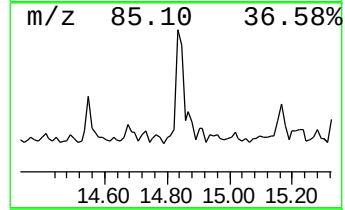
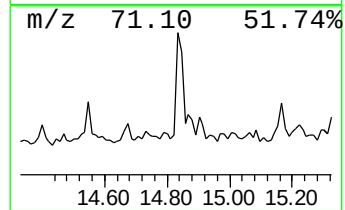
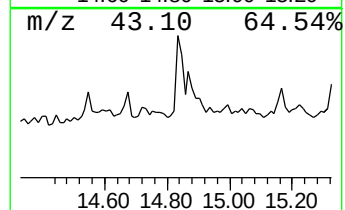
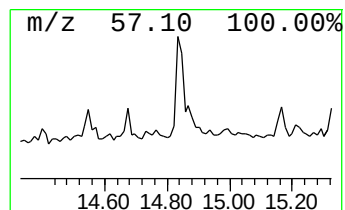
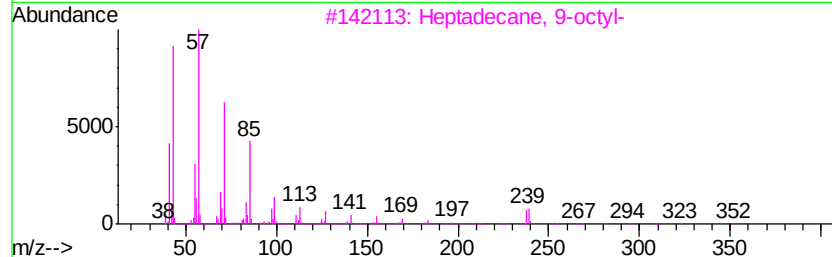
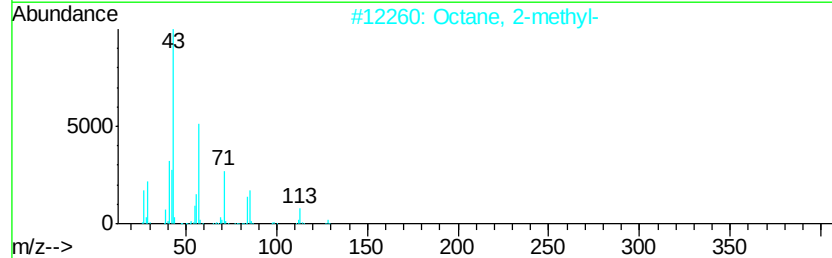
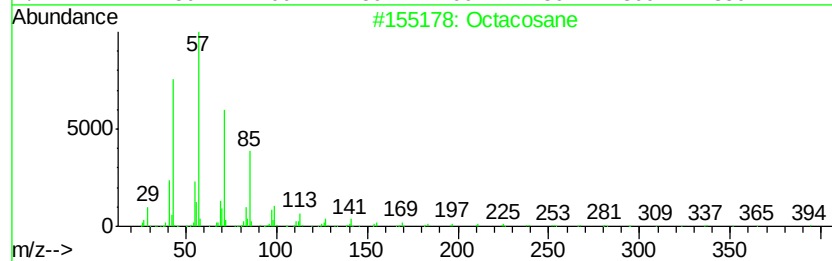
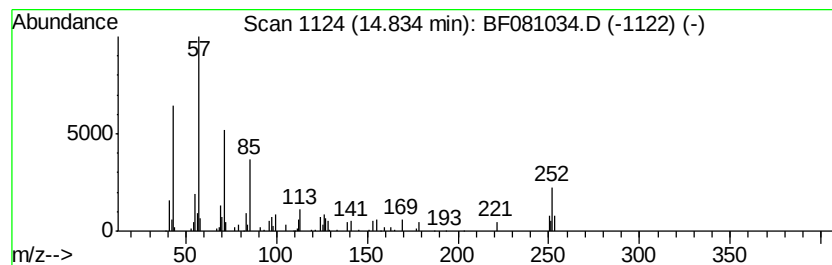
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	3.66 ng	45973	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	53
2		Octane, 2-methyl-	128	C9H20	003221-61-2	52
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	49
4		Eicosane	282	C20H42	000112-95-8	49
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
Data File : BF081034.D
Acq On : 18 Aug 2015 7:25
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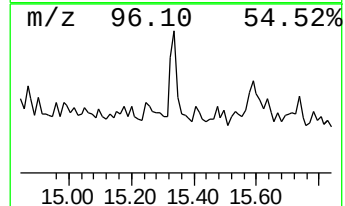
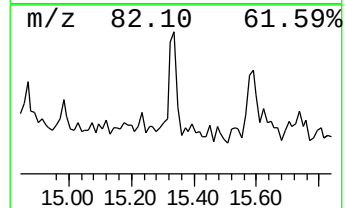
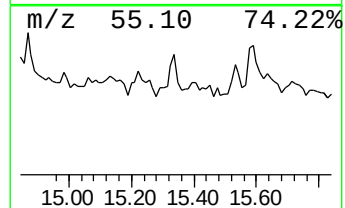
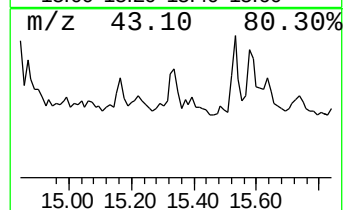
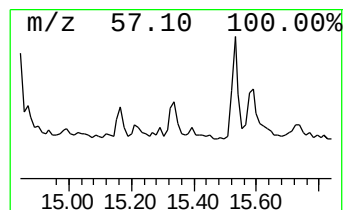
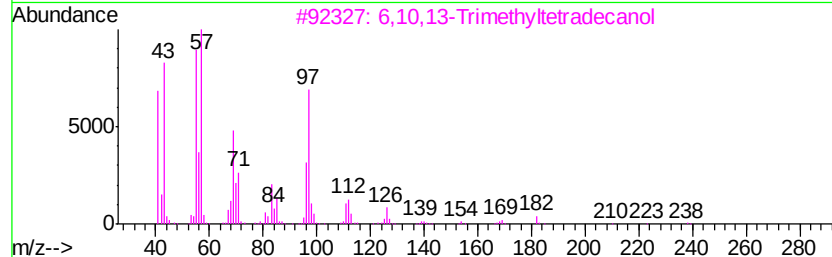
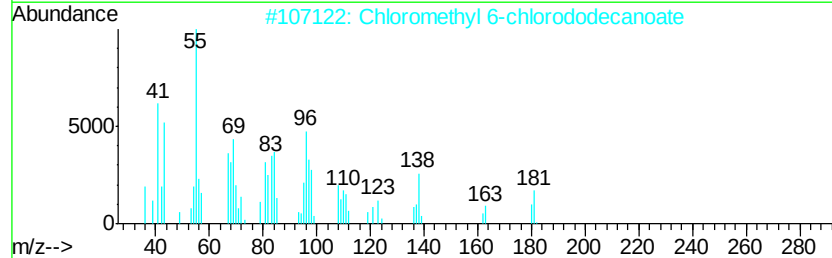
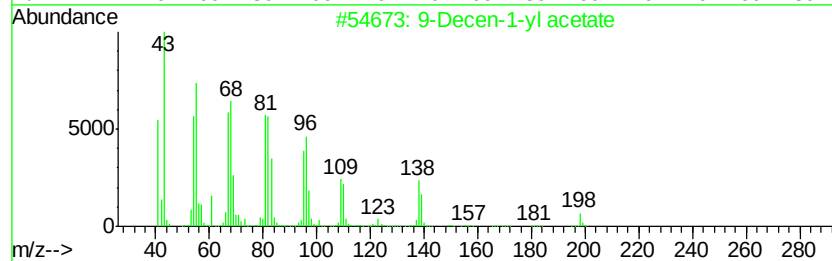
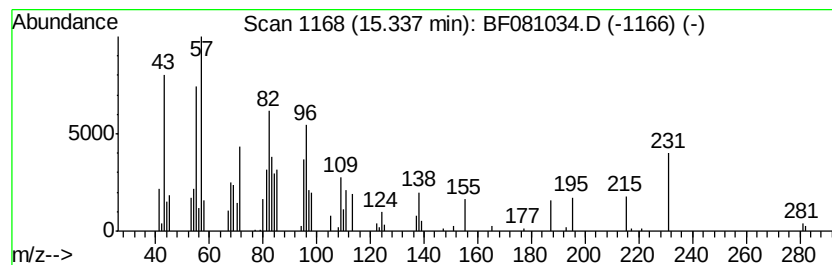
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.34 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.34	2.65 ng	33273	Perylene-d12		15.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Decen-1-yl	acetate	198	C12H22O2	050816-18-7	38	
2	Chloromethyl	6-chlorododecanoate	282	C13H24Cl2O2	080419-02-9	27	
3	6,10,13-Trimethyl	tetradecanol	256	C17H36O	1000131-71-0	25	
4	Z-13-Octadecen-1-yl	acetate	310	C20H38O2	060037-58-3	22	
5	trans-3-tert-butyl	cycloheptanol	170	C11H22O	019790-14-8	22	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
Data File : BF081034.D
Acq On : 18 Aug 2015 7:25
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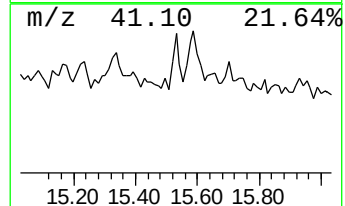
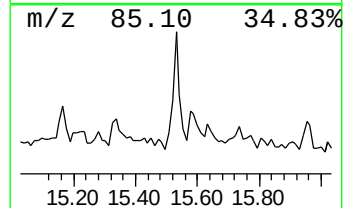
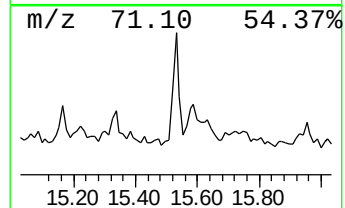
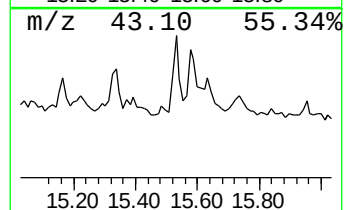
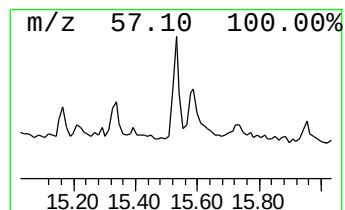
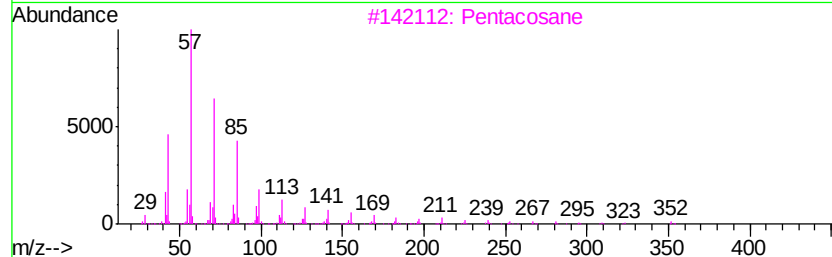
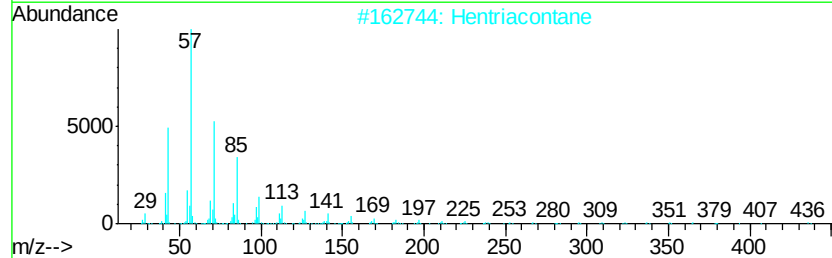
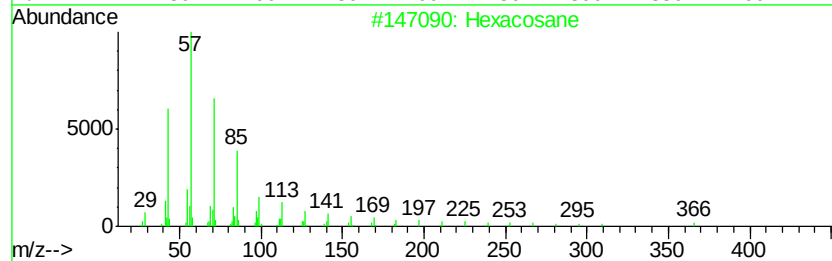
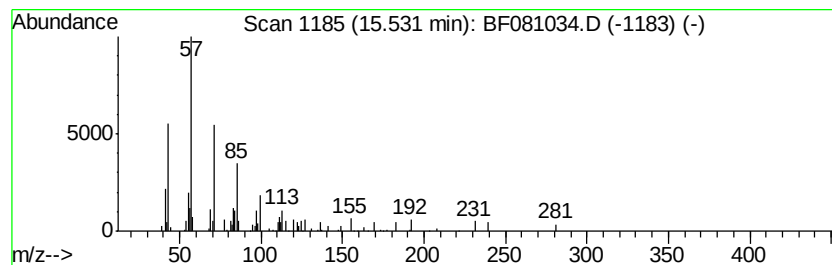
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	4.16 ng	52254	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	80
2		Hentriacontane	437	C31H64	000630-04-6	80
3		Pentacosane	352	C25H52	000629-99-2	80
4		Tetratriacontane	479	C34H70	014167-59-0	80
5		Eicosane	282	C20H42	000112-95-8	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
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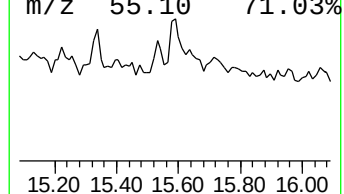
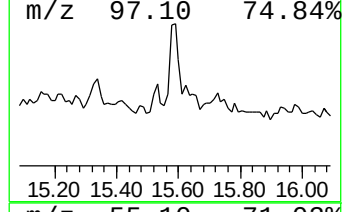
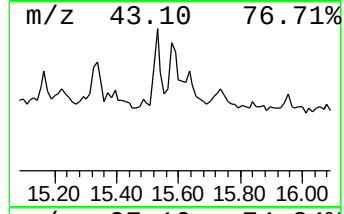
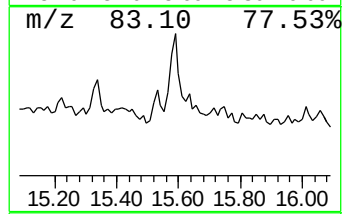
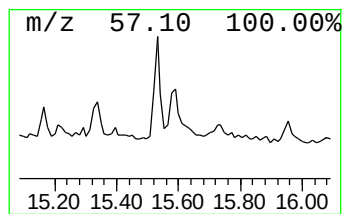
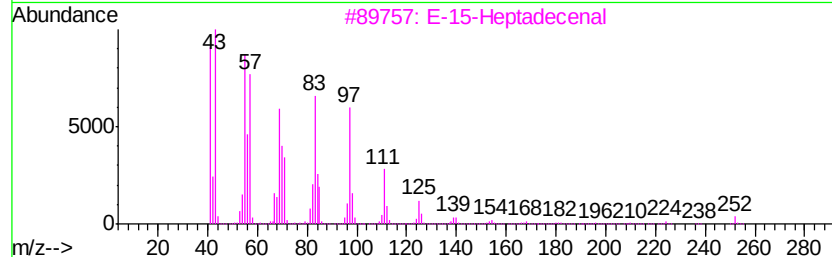
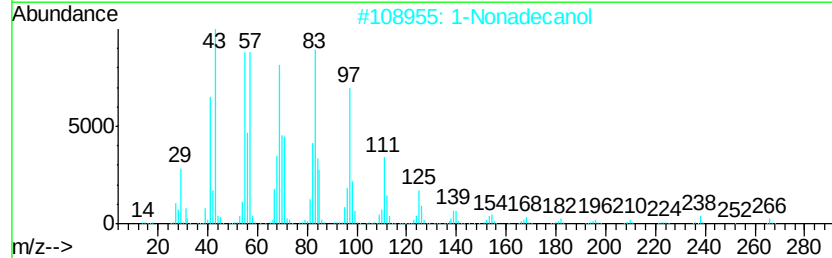
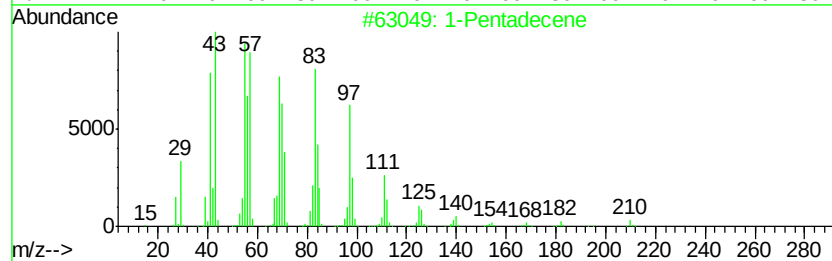
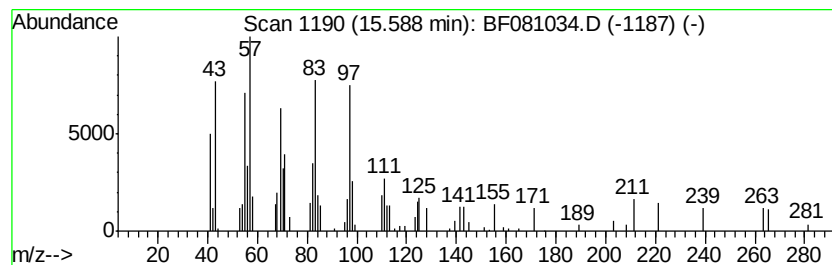
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Pentadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.59	4.05 ng	50846	Perylene-d12		15.11
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentadecene	210	C15H30	013360-61-7	72
2	1-Nonadecanol	284	C19H40O	001454-84-8	72
3	E-15-Heptadecenal	252	C17H32O	1000130-97-9	64
4	1-Tetracosanol	354	C24H50O	000506-51-4	52
5	1-Nonadecene	266	C19H38	018435-45-5	43



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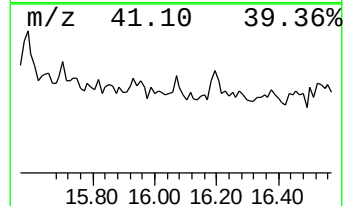
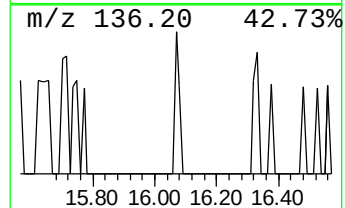
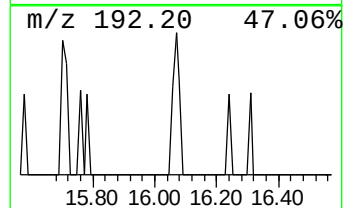
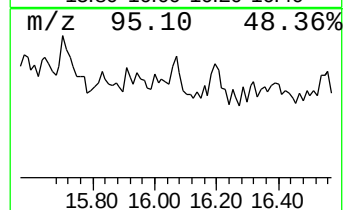
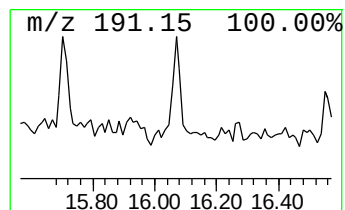
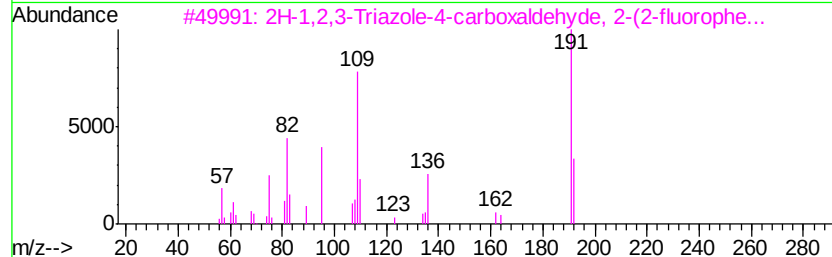
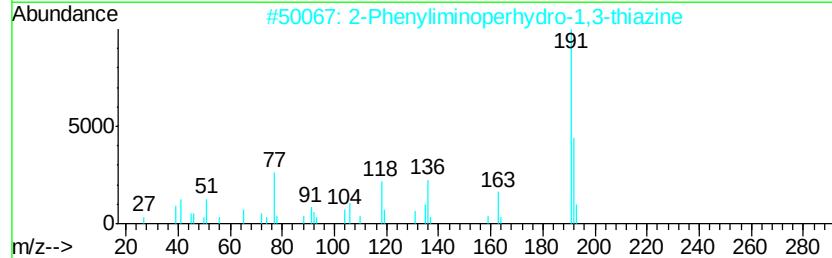
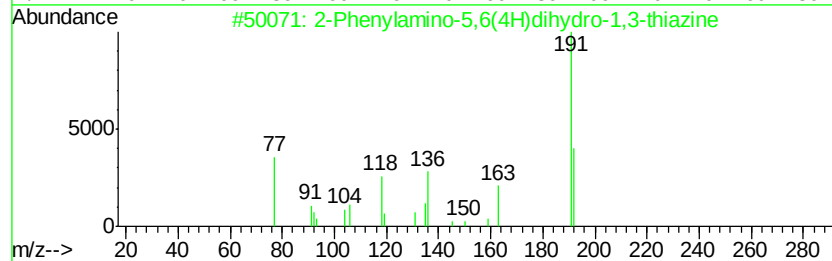
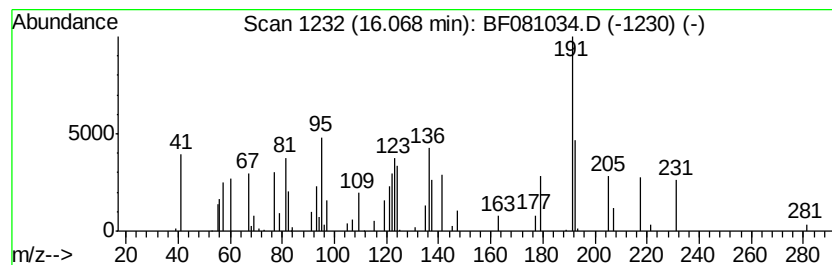
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown16.07 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	3.06 ng	38436	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenylamino-5,6(4H)dihydro-1,3...	192	C10H12N2S	1000147-35-4	27
2		2-Phenyliminoperhydro-1,3-thiazine	192	C10H12N2S	1000138-91-3	27
3		2H-1,2,3-Triazole-4-carboxaldehy...	191	C9H6FN3O	051306-43-5	25
4		7a-Isopropenyl-4,5-dimethyloctah...	236	C15H24O2	1000192-14-7	23
5		Pyrido[3,4-d]pyridazin-1(2H)-one...	191	C8H9N5O	1000263-69-7	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
Data File : BF081034.D
Acq On : 18 Aug 2015 7:25
Operator : UM/IZ
Sample : G3290-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
2A(0-2)-5

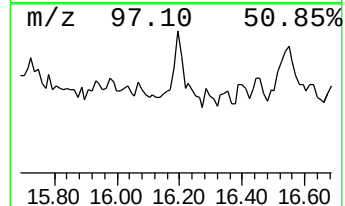
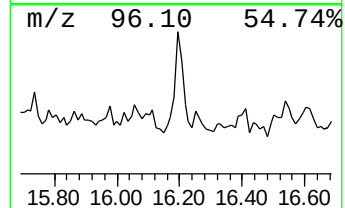
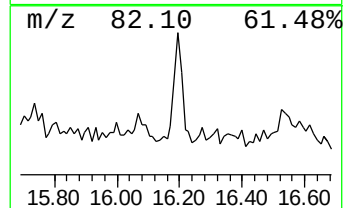
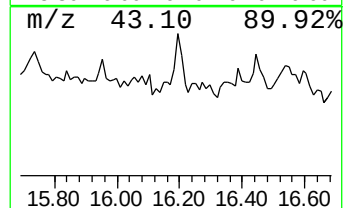
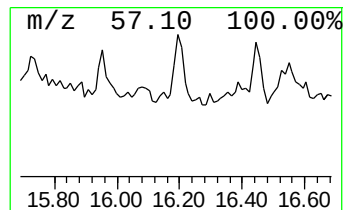
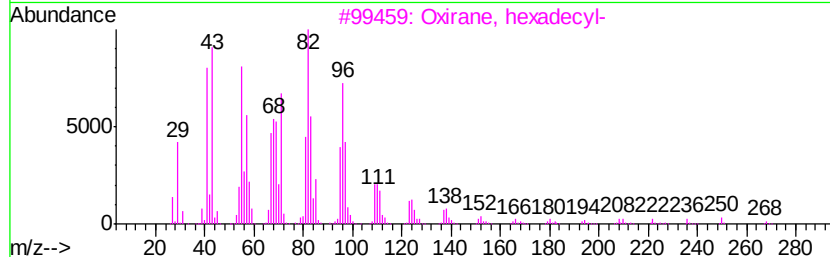
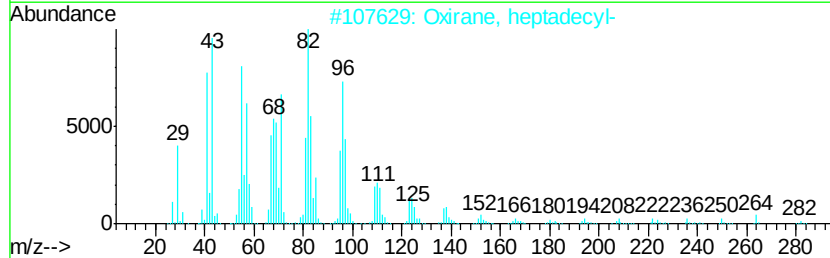
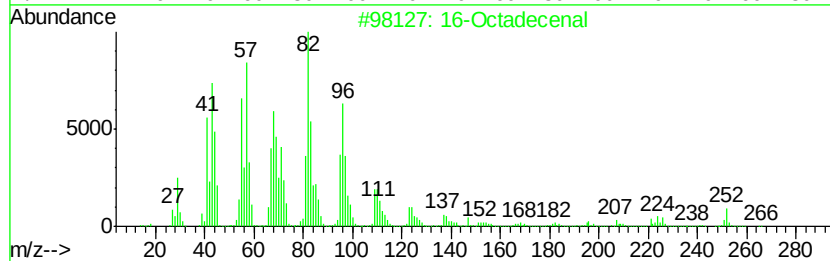
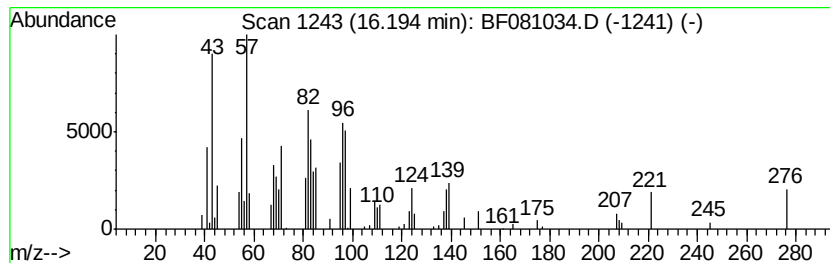
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.19 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.19	3.11 ng	39108	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	16-Octadecenal	266	C18H34O	056554-87-1	38
2		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	30
3		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	27
4		2,5a-Methano-5aH-pyrido[1,2-b][1...	167	C10H17NO	054980-46-0	27
5		Ecgonine	185	C9H15NO3	000481-37-8	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
Data File : BF081034.D
Acq On : 18 Aug 2015 7:25
Operator : UM/IZ
Sample : G3290-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
2A(0-2)-5

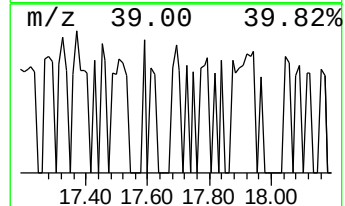
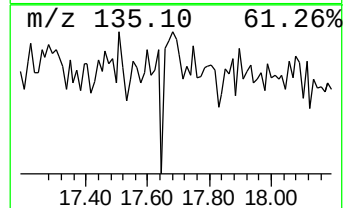
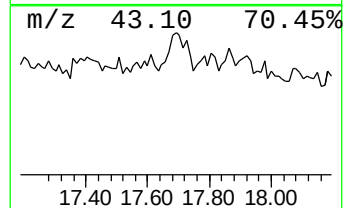
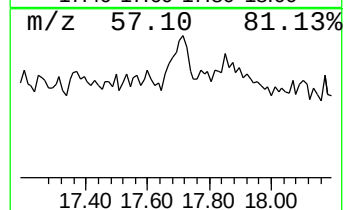
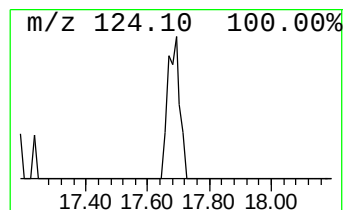
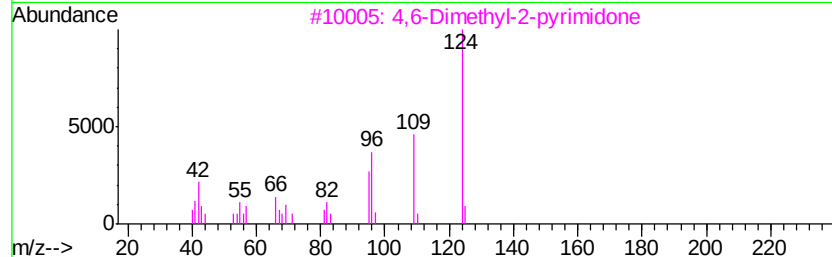
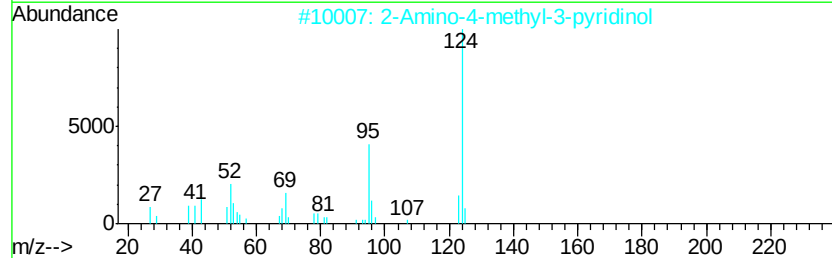
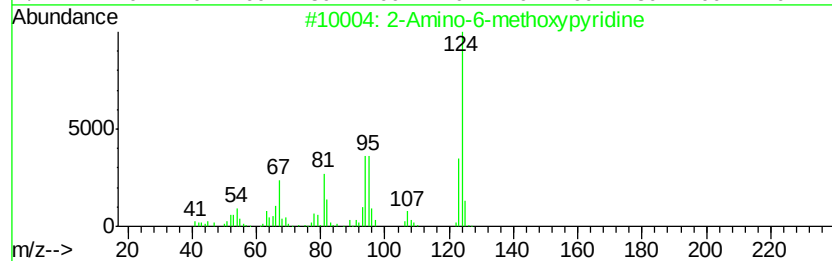
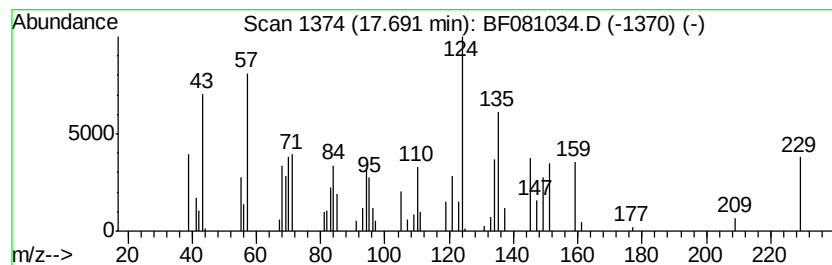
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown17.69 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.69	2.82 ng	35411	Perylene-d12	15.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Amino-6-methoxypyridine	124	C6H8N2O	017920-35-3	25
2			2-Amino-4-methyl-3-pyridinol	124	C6H8N2O	020348-18-9	22
3			4,6-Dimethyl-2-pyrimidone	124	C6H8N2O	000108-79-2	22
4			Methyl 2-[(1H-2-oxo-4-pyrimidinyl...]	183	C7H9N3O3	128786-77-6	16
5			N-Methyl-2-(propionylmethylene)p...	153	C9H15NO	082071-14-5	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
Data File : BF081034.D
Acq On : 18 Aug 2015 7:25
Operator : UM/IZ
Sample : G3290-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
2A(0-2)-5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.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...	4.75	54.6	ng	1019630	1	6.63	373622	20.0
unknown6.40	6.40	83.7	ng	1563970	1	6.63	373622	20.0
Eicosane	10.58	2.2	ng	45028	4	11.13	410889	20.0
Undecanoic acid	11.68	5.3	ng	108356	4	11.13	410889	20.0
Trichloroacetic a...	13.63	9.6	ng	233016	5	13.75	485853	20.0
Pentafluoropropio...	14.26	3.8	ng	92687	5	13.75	485853	20.0
Squalene	14.61	2.3	ng	28206	6	15.11	251166	20.0
Octacosane	14.83	3.7	ng	45973	6	15.11	251166	20.0
unknown15.34	15.34	2.6	ng	33273	6	15.11	251166	20.0
Hexacosane	15.53	4.2	ng	52254	6	15.11	251166	20.0
1-Pentadecene	15.59	4.0	ng	50846	6	15.11	251166	20.0
unknown16.07	16.07	3.1	ng	38436	6	15.11	251166	20.0
unknown16.19	16.19	3.1	ng	39108	6	15.11	251166	20.0
unknown17.69	17.69	2.8	ng	35411	6	15.11	251166	20.0