

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
 Data File : BF082906.D
 Acq On : 13 Nov 2015 19:29
 Operator : UM/IZ
 Sample : G4370-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-TS003-01

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-BF110715.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.430	27	30	33	rBV	91674	129550	3.09%	0.268%
2	4.967	249	252	255	rVB	1709338	1972227	47.02%	4.077%
3	5.413	288	291	293	rBV	3195391	3674345	87.60%	7.596%
4	6.453	379	382	384	rBV	3618069	4194377	100.00%	8.671%
5	6.567	390	392	395	rBV	3011529	3418869	81.51%	7.067%
6	6.796	410	412	415	rBV	1643672	1572291	37.49%	3.250%
7	6.956	424	426	428	rVB	3236798	2716922	64.78%	5.616%
8	7.367	459	462	463	rBV	2256553	2585455	61.64%	5.345%
9	7.836	501	503	511	rVB3	40868	94119	2.24%	0.195%
10	7.996	515	517	519	rBV	44481	44620	1.06%	0.092%
11	8.087	523	525	527	rBV	2592499	2110960	50.33%	4.364%
12	8.853	589	592	595	rBV	49768	81762	1.95%	0.169%
13	9.173	617	620	622	rBV	2788562	3622560	86.37%	7.489%
14	9.299	630	631	638	rVB2	26442	47483	1.13%	0.098%
15	9.562	652	654	656	rBV	1036363	827402	19.73%	1.710%
16	9.665	659	663	665	rBV4	23945	43813	1.04%	0.091%
17	9.790	672	674	676	rVB	105575	85196	2.03%	0.176%
18	9.848	676	679	681	rBV	2018933	2270698	54.14%	4.694%
19	10.053	696	697	701	rVB3	39355	63133	1.51%	0.131%
20	10.259	711	715	716	rBV2	37322	61293	1.46%	0.127%
21	10.293	716	718	719	rVB	99812	121259	2.89%	0.251%
22	10.362	722	724	730	rVB2	27624	46971	1.12%	0.097%
23	10.511	734	737	740	rBV	77076	138251	3.30%	0.286%
24	10.625	745	747	750	rBV2	1899944	2110236	50.31%	4.362%
25	10.762	756	759	763	rBV	225038	270114	6.44%	0.558%
26	10.853	765	767	769	rVV2	53556	76775	1.83%	0.159%
27	10.888	769	770	773	rVV2	48607	71449	1.70%	0.148%
28	10.956	773	776	778	rVV2	37845	68805	1.64%	0.142%
29	11.002	778	780	782	rVV	70999	91111	2.17%	0.188%
30	11.036	782	783	785	rVV	53548	64639	1.54%	0.134%
31	11.082	785	787	790	rVB2	147175	193942	4.62%	0.401%
32	11.208	796	798	799	rBV	96227	85932	2.05%	0.178%
33	11.288	803	805	806	rBV	36929	47757	1.14%	0.099%
34	11.322	806	808	811	rVV	2221418	2033715	48.49%	4.204%

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35	11.391	811	814	816	rVB4	56734	120486	2.87%	0.249%
36	11.791	847	849	851	rBV	59054	104173	2.48%	0.215%
37	11.871	853	856	858	rBV	606744	645751	15.40%	1.335%
38	12.156	879	881	883	rBV	30266	43199	1.03%	0.089%
39	12.214	884	886	887	rVB2	46733	52132	1.24%	0.108%
40	12.316	893	895	897	rBV	49915	81030	1.93%	0.168%
41	12.374	898	900	902	rVV	96160	102921	2.45%	0.213%
42	12.568	912	917	922	rBV2	1637011	3563980	84.97%	7.367%
43	12.659	922	925	929	rVV	341609	610100	14.55%	1.261%
44	12.717	929	930	932	rVV	106142	97590	2.33%	0.202%
45	12.762	932	934	935	rVV	119740	126823	3.02%	0.262%
46	12.911	945	947	949	rBV	2612796	2244405	53.51%	4.640%
47	13.071	959	961	964	rBV2	101869	172318	4.11%	0.356%
48	13.402	988	990	992	rVB2	86098	122504	2.92%	0.253%
49	13.574	1003	1005	1006	rVB	99794	74990	1.79%	0.155%
50	13.757	1019	1021	1023	rVB2	74001	86047	2.05%	0.178%
51	13.825	1024	1027	1029	rBV2	191955	315212	7.52%	0.652%
52	13.962	1036	1039	1041	rBV	1512052	1639292	39.08%	3.389%
53	14.454	1080	1082	1090	rBV2	133157	360325	8.59%	0.745%
54	15.082	1134	1137	1140	rBV2	115030	298906	7.13%	0.618%
55	15.380	1160	1163	1166	rBV	1161355	1446953	34.50%	2.991%
56	15.848	1202	1204	1206	rVB	106959	134175	3.20%	0.277%
57	16.054	1220	1222	1228	rVB5	121612	372800	8.89%	0.771%
58	17.083	1310	1312	1316	rVB2	113040	243575	5.81%	0.504%
59	17.311	1329	1332	1333	rBV2	146478	276819	6.60%	0.572%

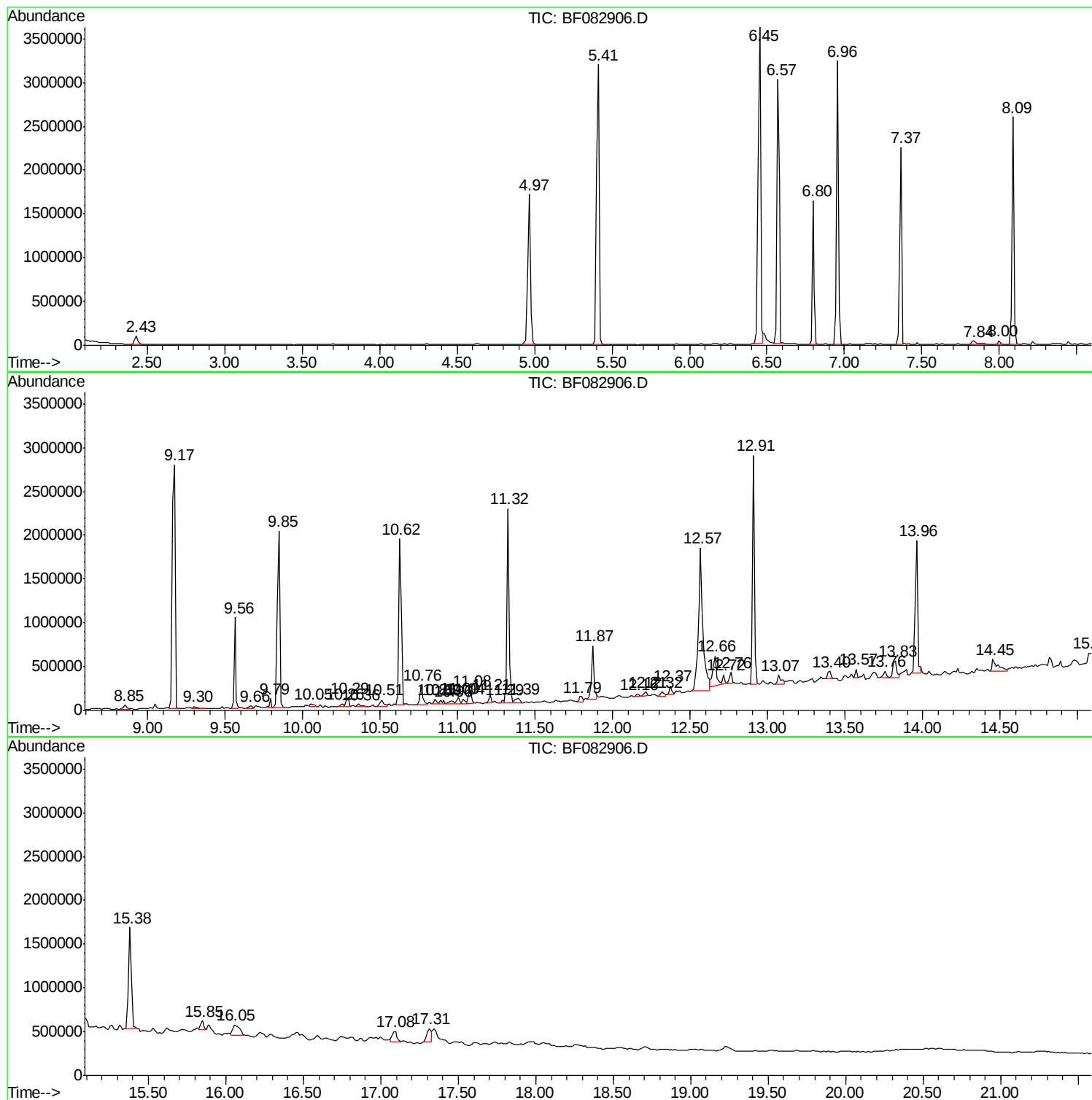
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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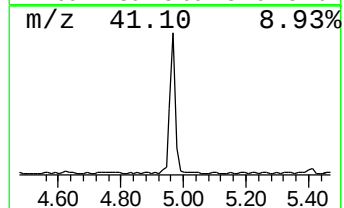
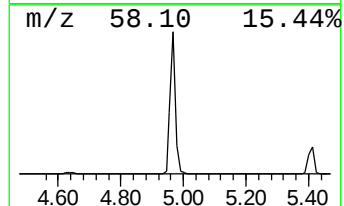
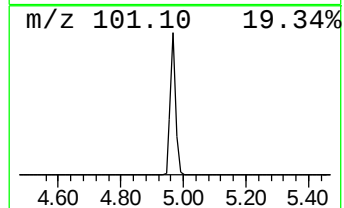
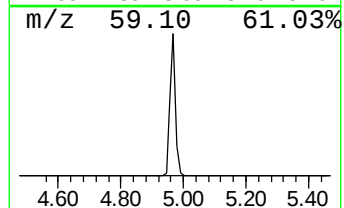
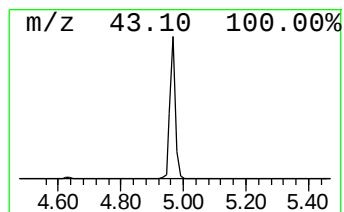
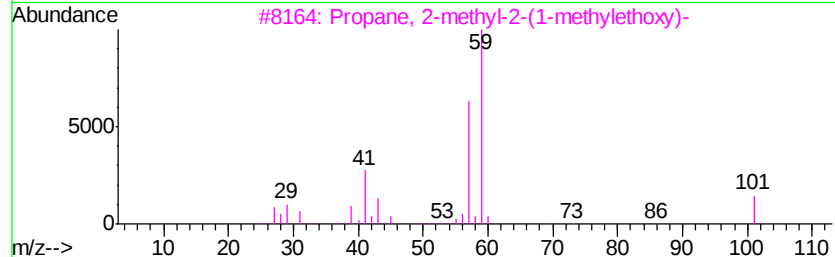
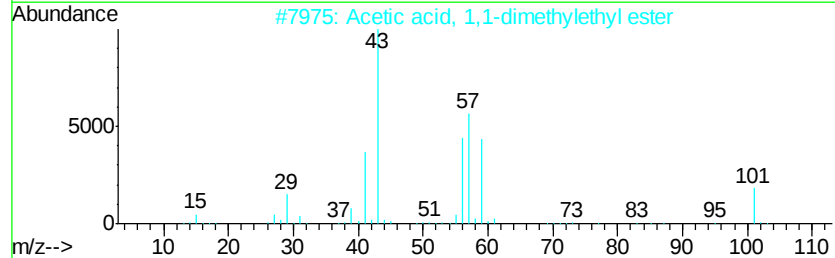
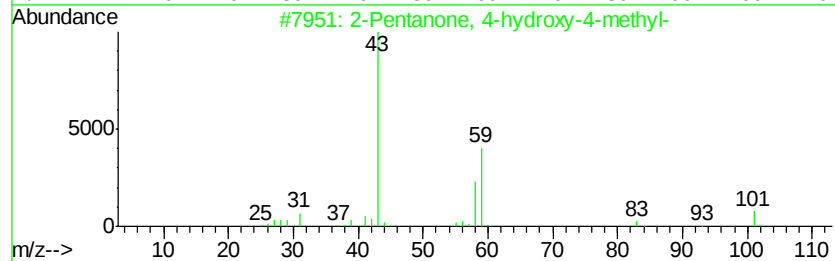
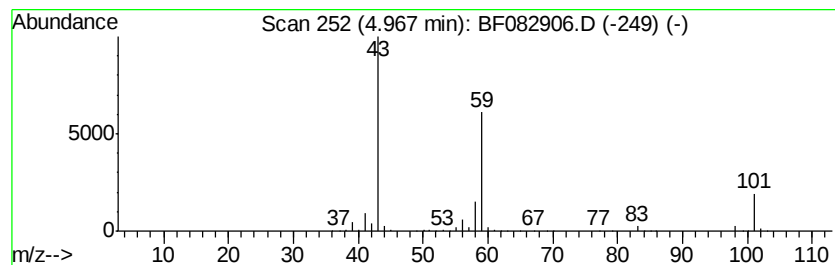
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	25.09 ng	1972230	1,4-Dichlorobenzene-d4	6.80

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	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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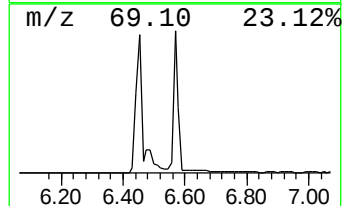
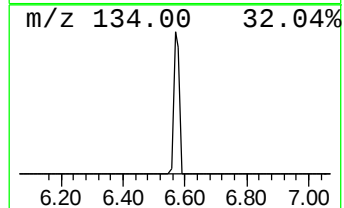
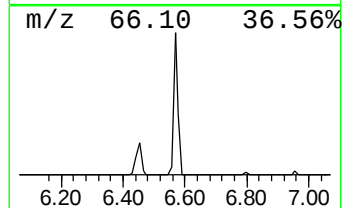
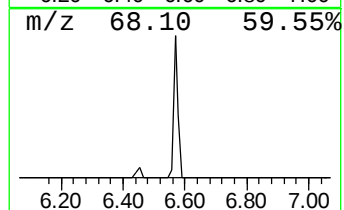
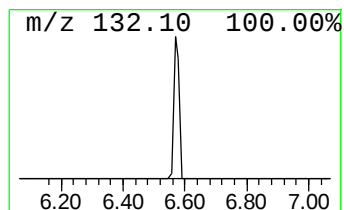
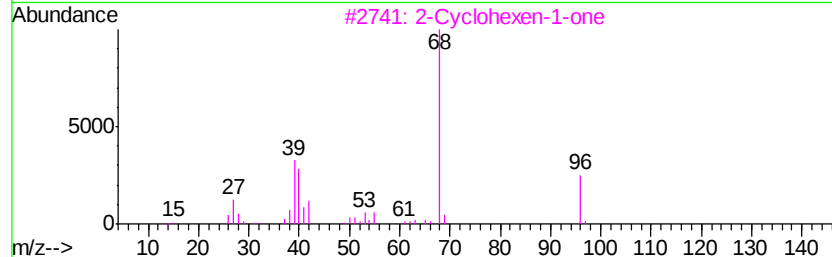
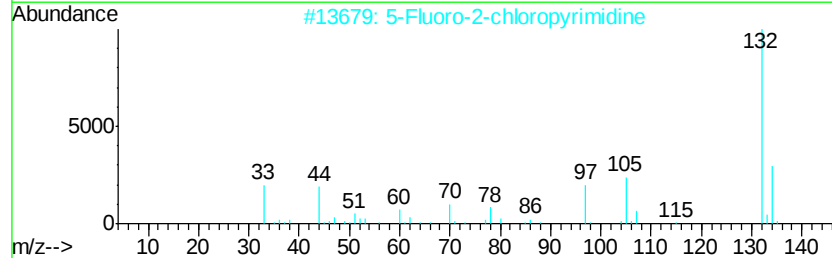
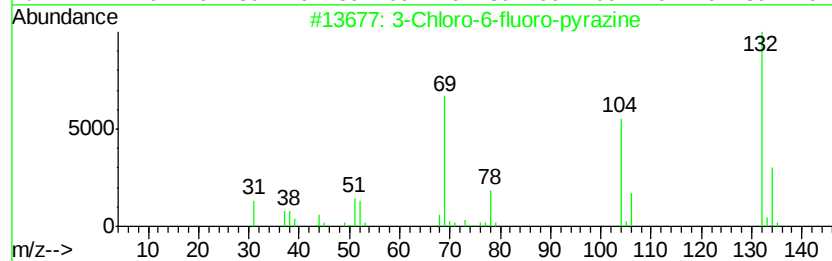
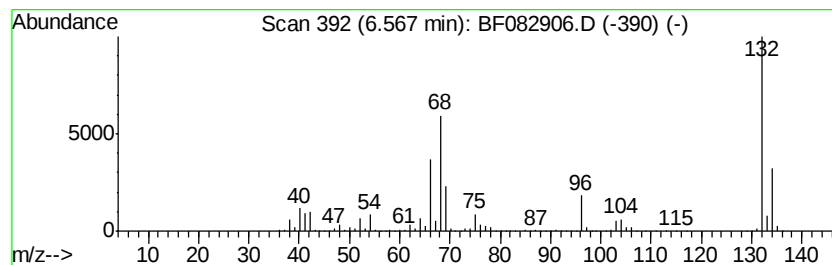
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	43.49 ng	3418870	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	11
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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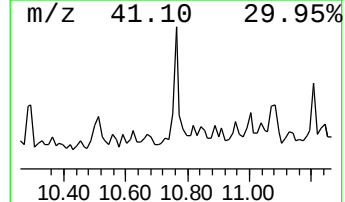
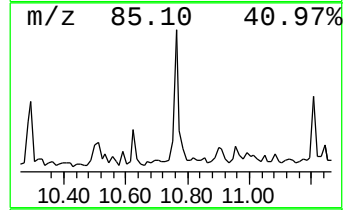
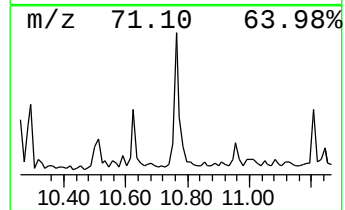
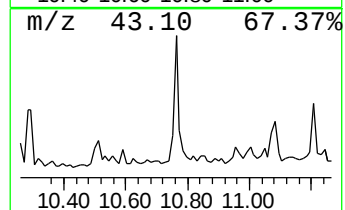
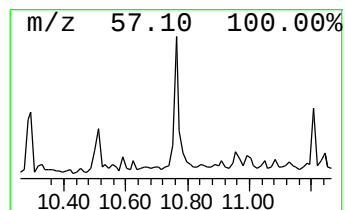
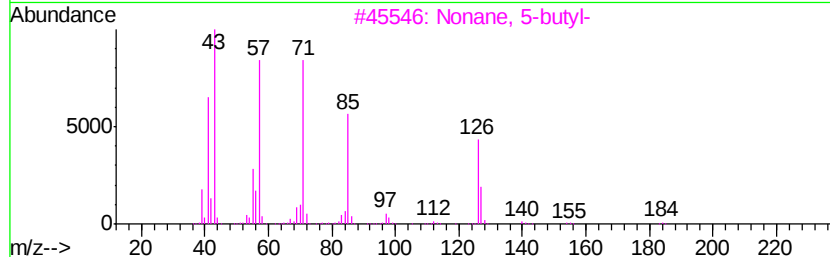
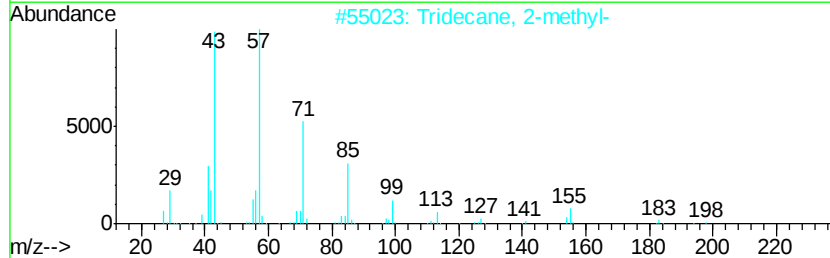
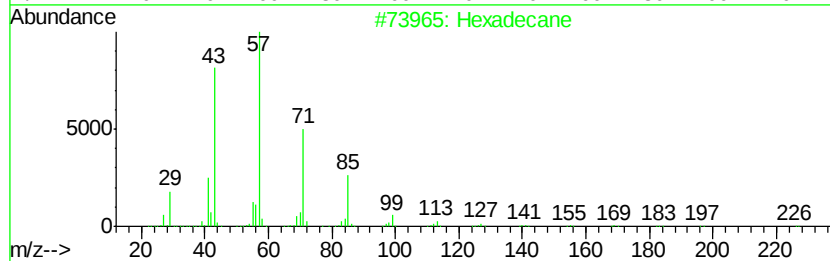
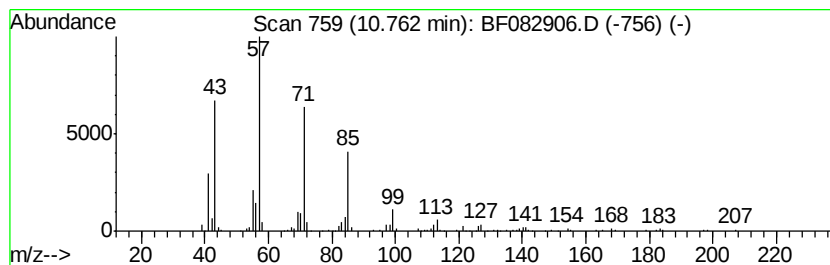
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.76	2.66 ng	270114	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	83
2	Tridecane, 2-methyl-	198	C14H30	001560-96-9	81
3	Nonane, 5-butyl-	184	C13H28	017312-63-9	70
4	Dodecane, 2-methyl-	184	C13H28	001560-97-0	68
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64



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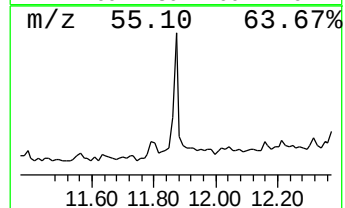
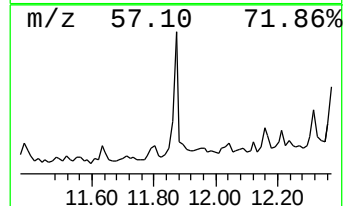
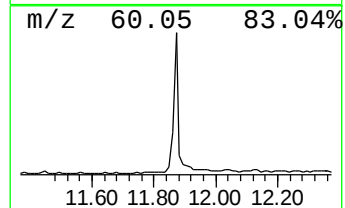
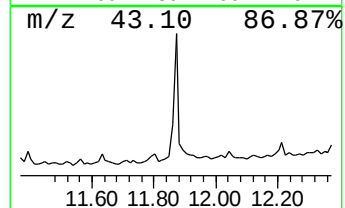
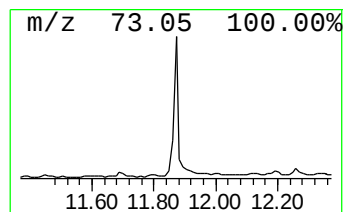
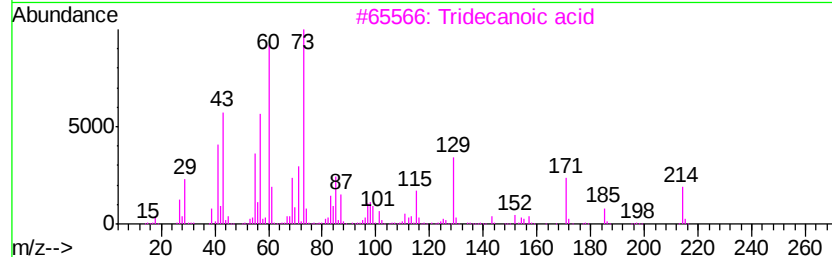
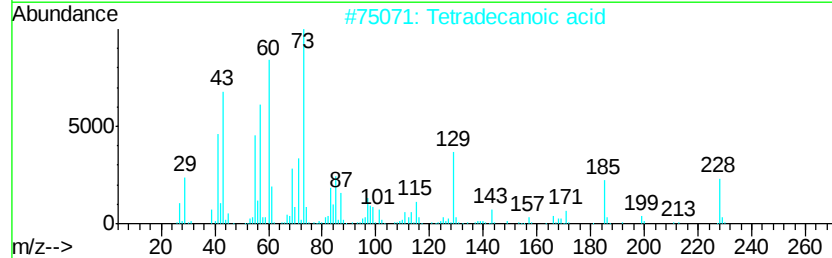
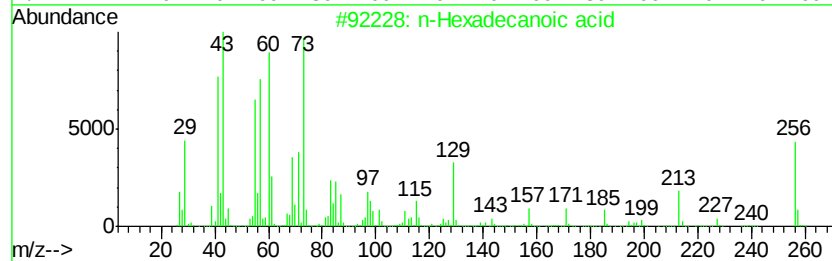
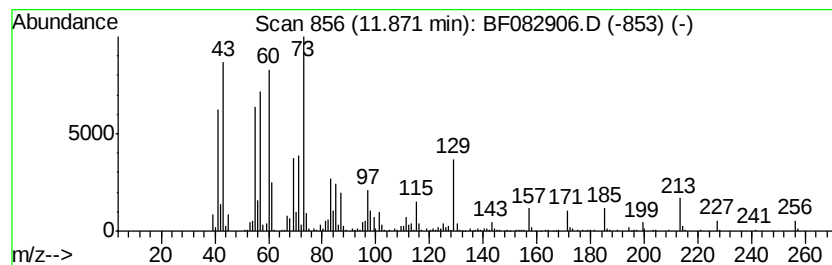
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.87	6.35 ng	645751	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3	Tridecanoic acid	214	C13H26O2	000638-53-9	86
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5	Hexadecane	226	C16H34	000544-76-3	11



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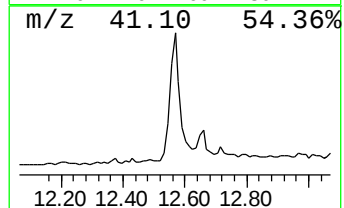
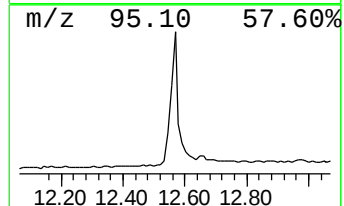
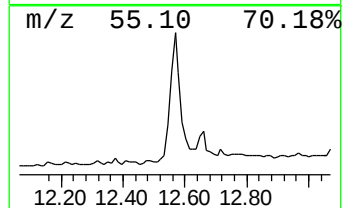
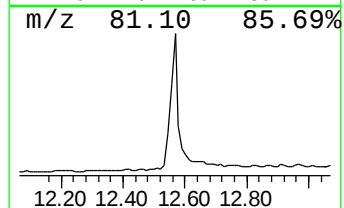
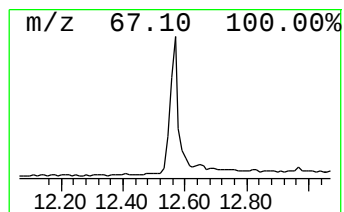
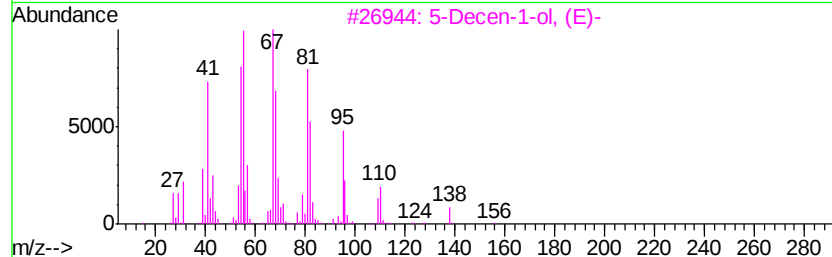
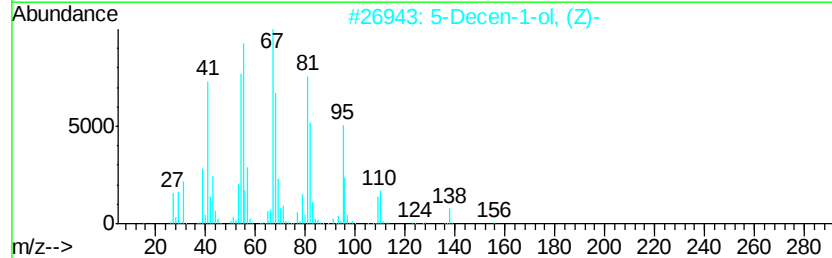
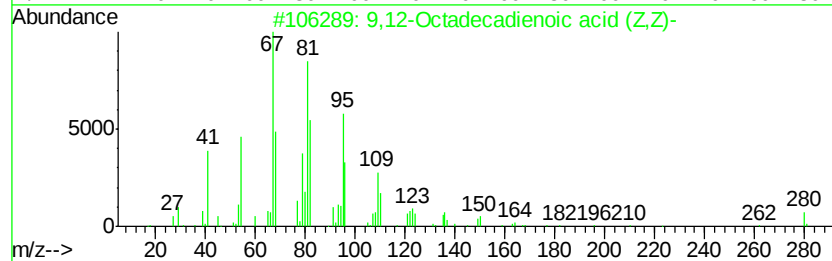
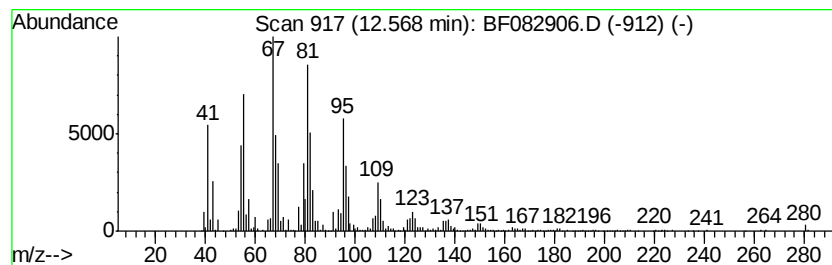
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9,12-Octadecadienoic acid (... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	35.05 ng	3563980	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
2		5-Decen-1-ol, (Z)-	156	C10H20O	051652-47-2	91
3		5-Decen-1-ol, (E)-	156	C10H20O	056578-18-8	91
4		7-Hexadecyne	222	C16H30	074685-28-2	90
5		1-Tetradecyne	194	C14H26	000765-10-6	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082906.D
Acq On : 13 Nov 2015 19:29
Operator : UM/IZ
Sample : G4370-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

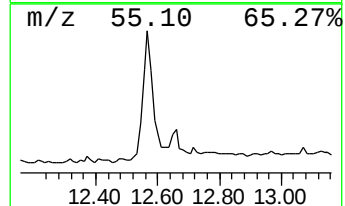
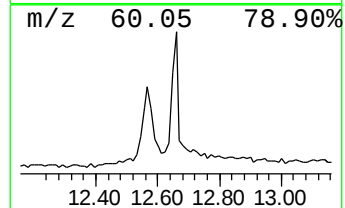
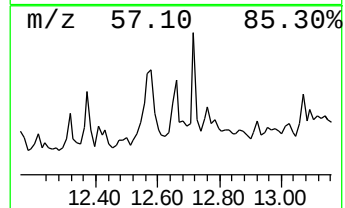
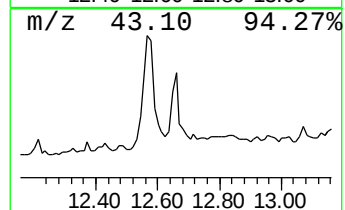
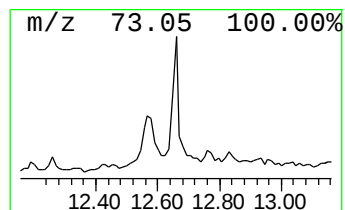
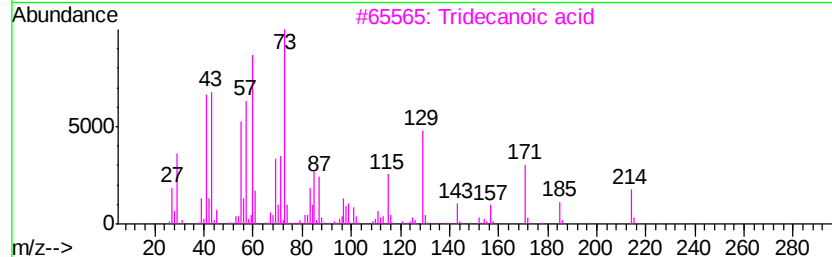
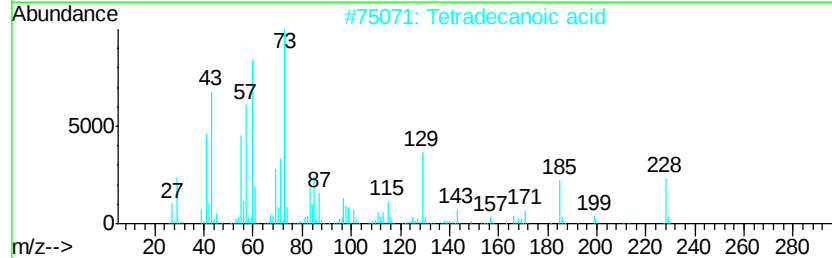
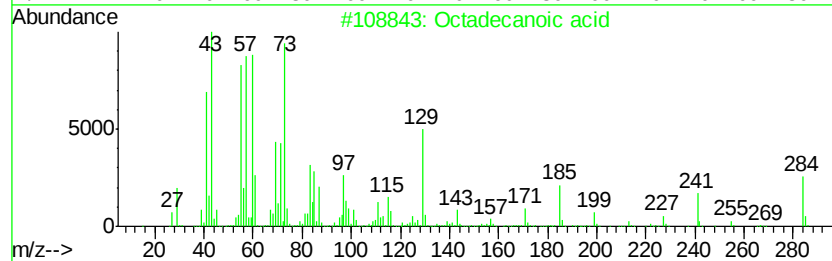
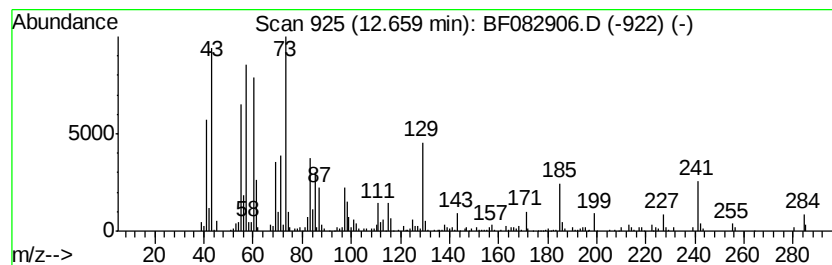
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.66	7.44 ng	610100	Chrysene-d12	13.96	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	90
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3	Tridecanoic acid	214	C13H26O2	000638-53-9	74
4	n-Decanoic acid	172	C10H20O2	000334-48-5	70
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	62



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082906.D
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Instrument :
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ClientSampleId :
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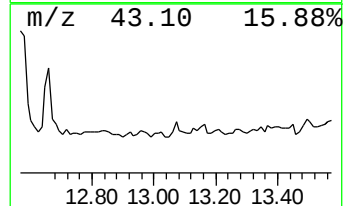
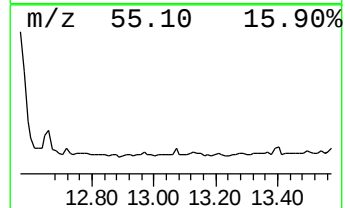
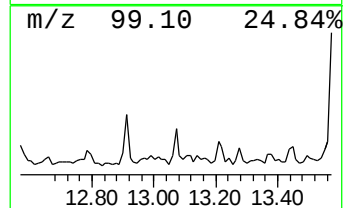
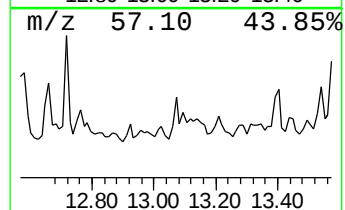
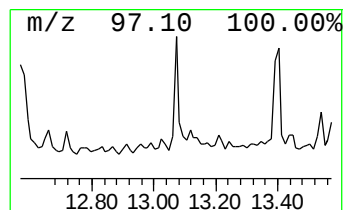
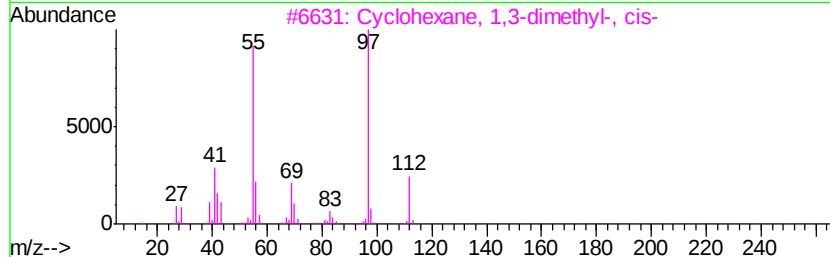
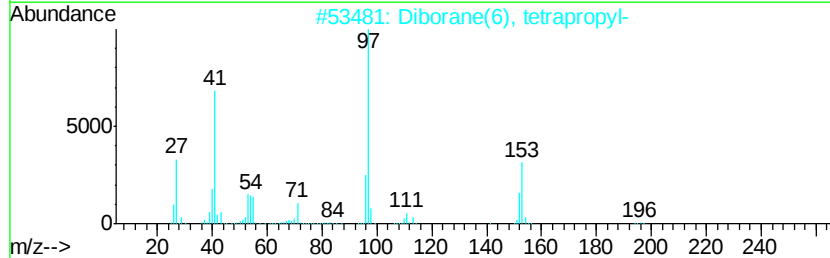
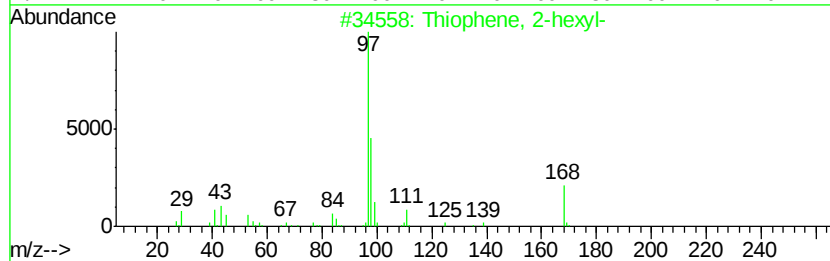
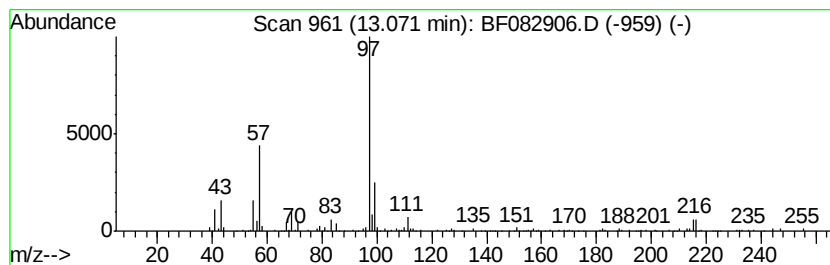
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Thiophene, 2-hexyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.07	2.10 ng	172318	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thiophene, 2-hexyl-	168	C10H16S	018794-77-9	59
2		Diborane(6), tetrapropyl-	196	C12H30B2	022784-01-6	50
3		Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	47
4		2H-Pyran-4-carboxylic acid, 3,4-...	170	C9H14O3	038858-64-9	47
5		Isoxazole, 3,5-dimethyl-	97	C5H7NO	000300-87-8	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111315\
Data File : BF082906.D
Acq On : 13 Nov 2015 19:29
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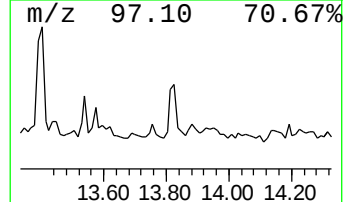
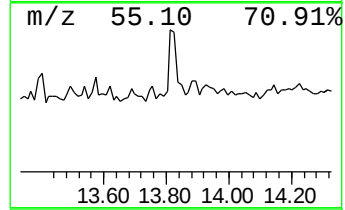
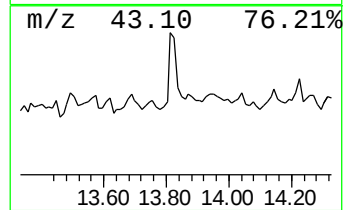
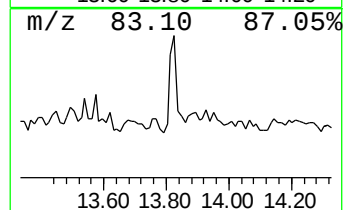
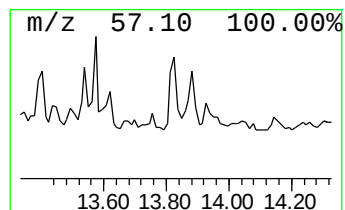
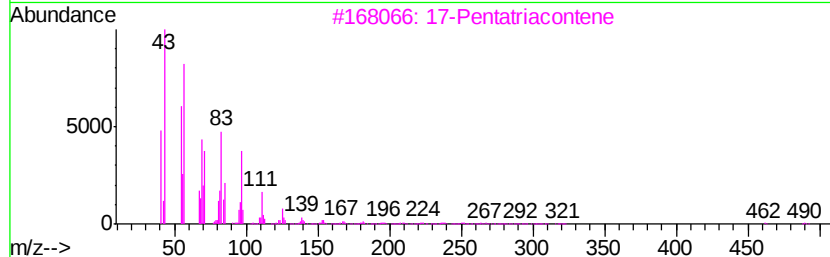
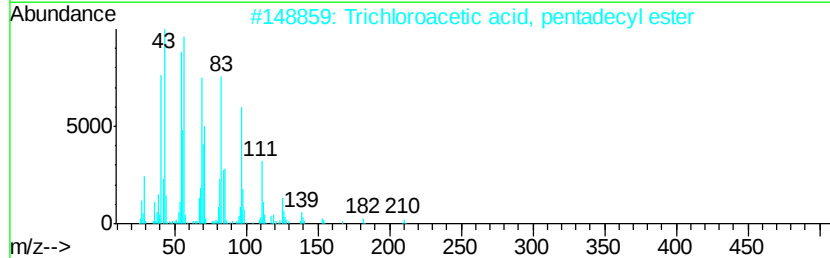
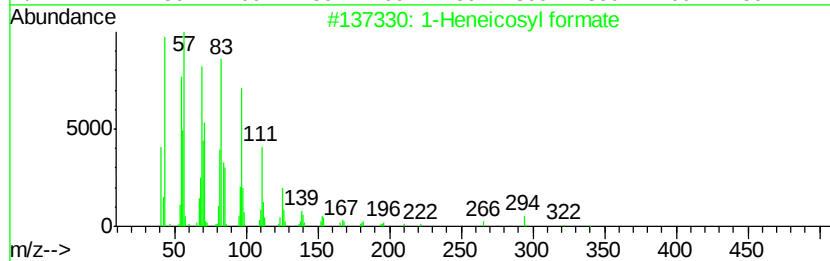
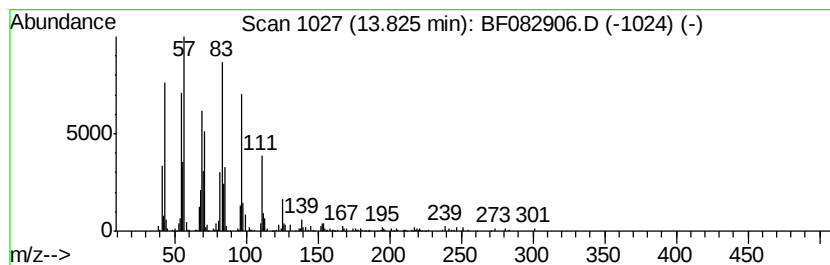
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Heneicosyl formate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	3.85 ng	315212	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	86
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	86
5		1-Octadecanol	270	C18H38O	000112-92-5	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082906.D
Acq On : 13 Nov 2015 19:29
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Misc :
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ClientSampleId :
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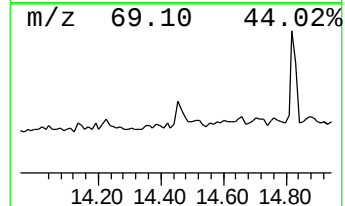
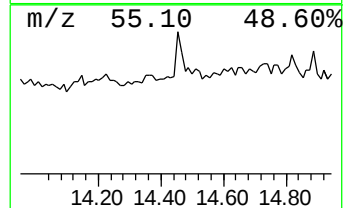
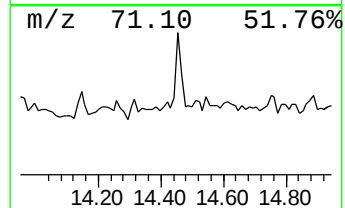
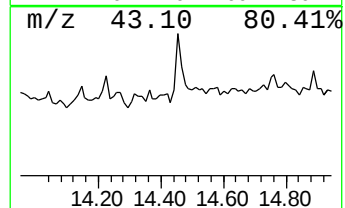
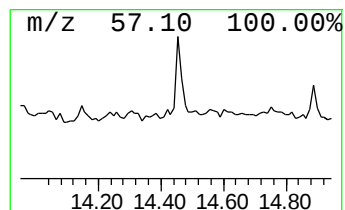
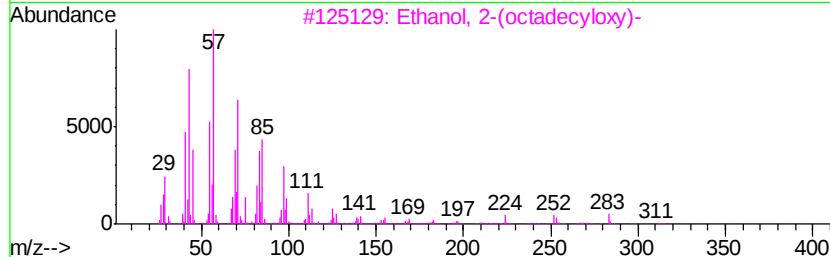
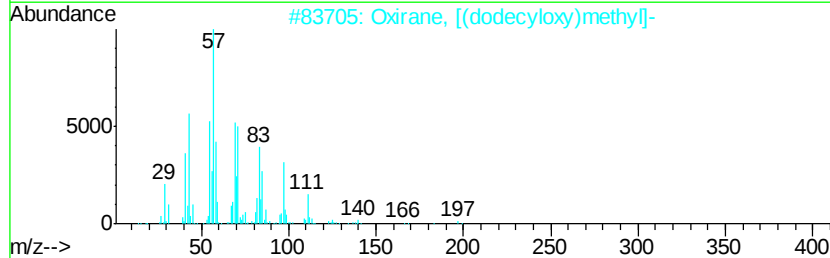
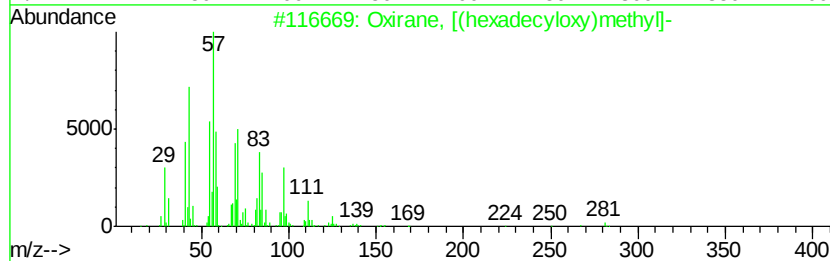
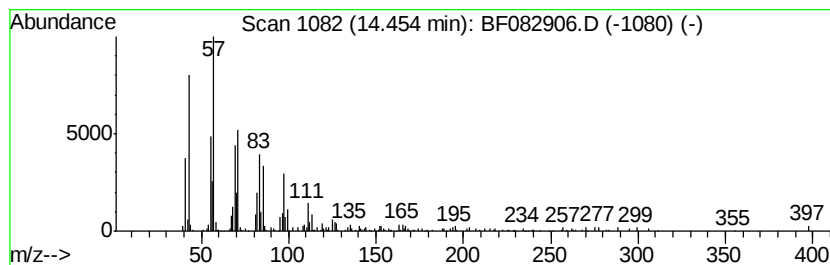
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Oxirane, [(hexadecyloxy)met... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	4.40 ng	360325	Chrysene-d12	13.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, [(hexadecyloxy)methyl]-	298	C19H38O2	015965-99-8	91
2			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	87
3			Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	83
4			17-Pentatriacontene	491	C35H70	006971-40-0	76
5			1-Docosanol	326	C22H46O	000661-19-8	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111315\
Data File : BF082906.D
Acq On : 13 Nov 2015 19:29
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Misc :
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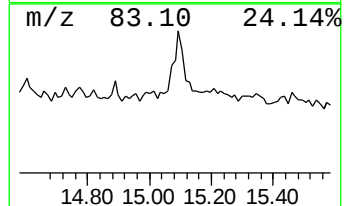
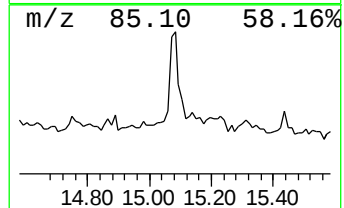
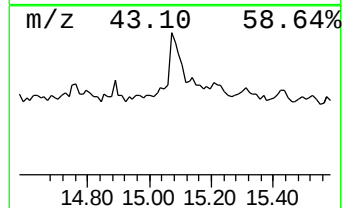
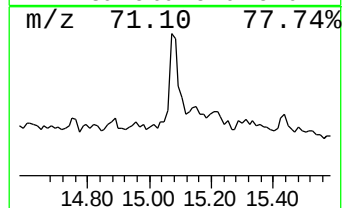
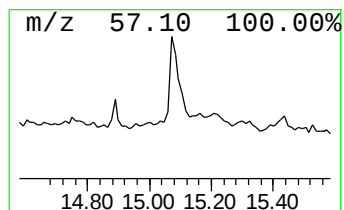
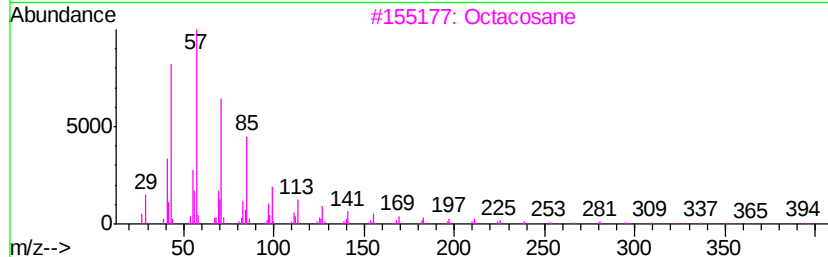
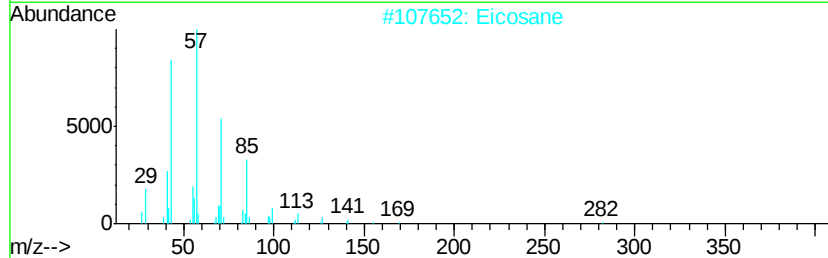
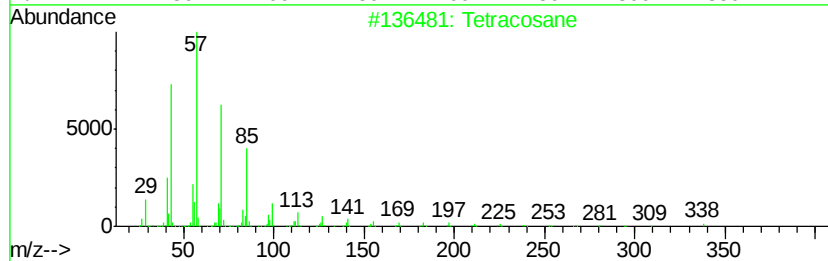
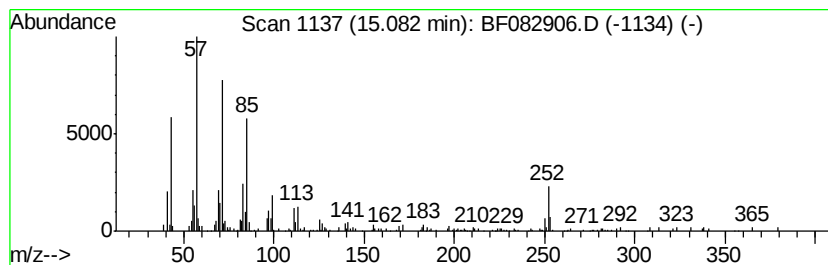
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	4.13 ng	298906	Perylene-d12	15.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	93
2		Eicosane	282	C20H42	000112-95-8	90
3		Octacosane	394	C28H58	000630-02-4	87
4		10-Methylnonadecane	282	C20H42	056862-62-5	83
5		Heptadecane	240	C17H36	000629-78-7	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
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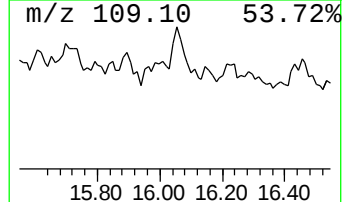
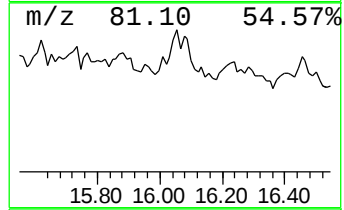
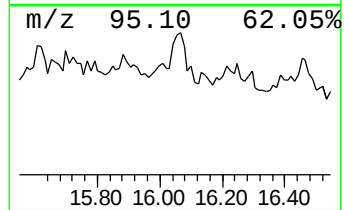
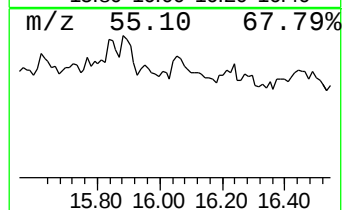
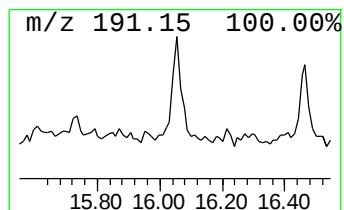
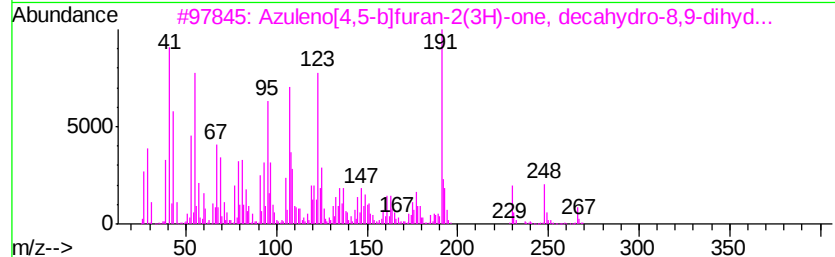
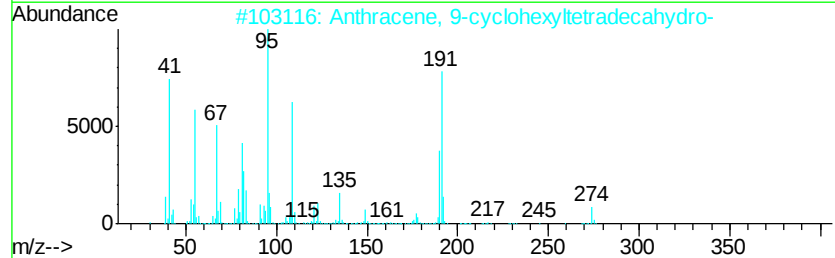
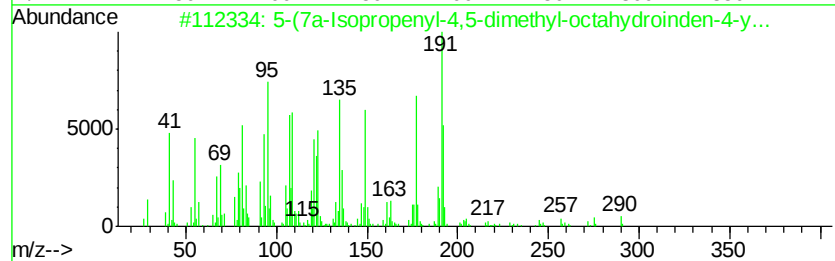
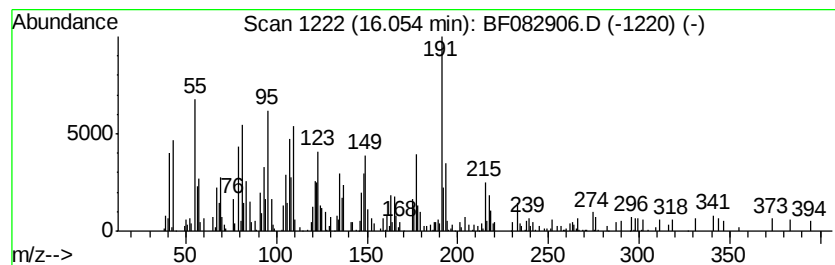
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 5-(7a-Isopropenyl-4,5-dimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.05	5.15 ng	372800	Perylene-d12	15.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(7a-Isopropenyl-4,5-dimethyl-o...	290	C20H34O	1000193-54-0	52
2	Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	47
3	Azuleno[4,5-b]furan-2(3H)-one, d...	266	C15H22O4	005090-67-5	38
4	2-Anthracenylloxirane	220	C16H12O	052643-86-4	30
5	(7a-Isopropenyl-4,5-dimethylocta...	222	C15H26O	1000187-02-9	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
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Acq On : 13 Nov 2015 19:29
Operator : UM/IZ
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P001-TS003-01

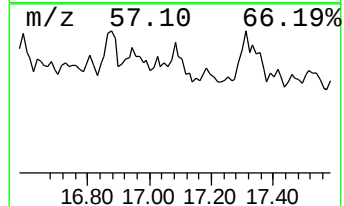
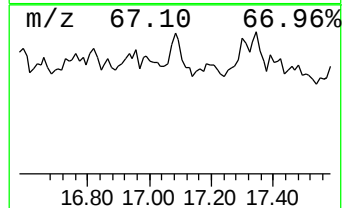
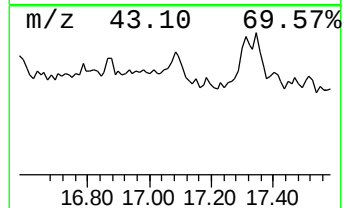
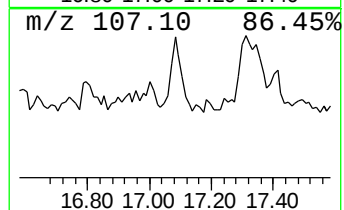
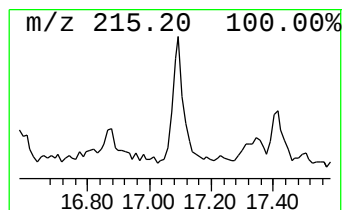
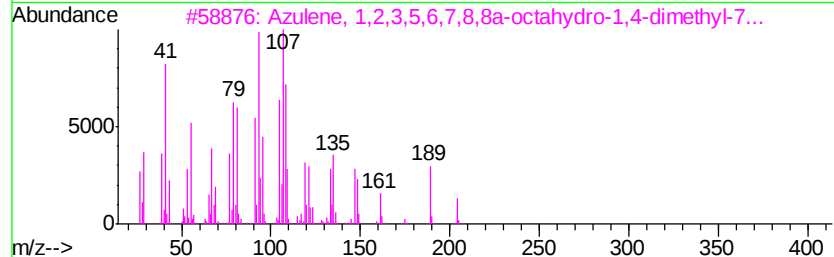
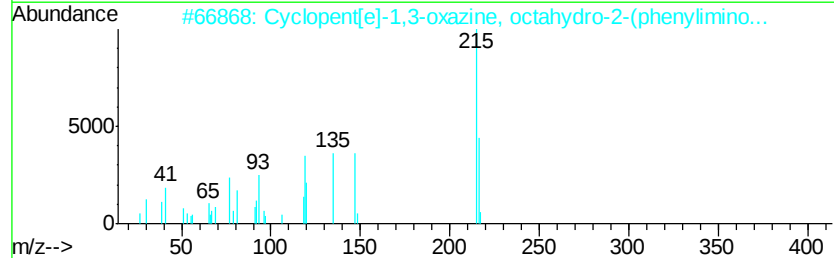
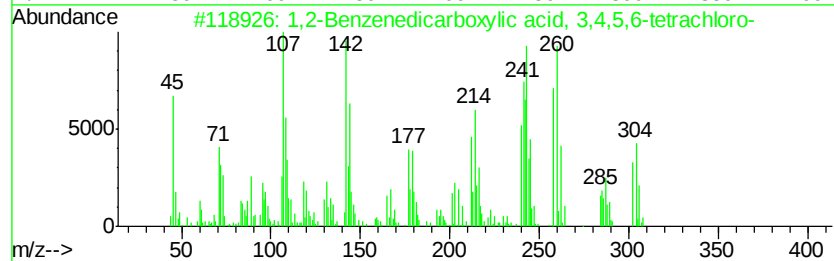
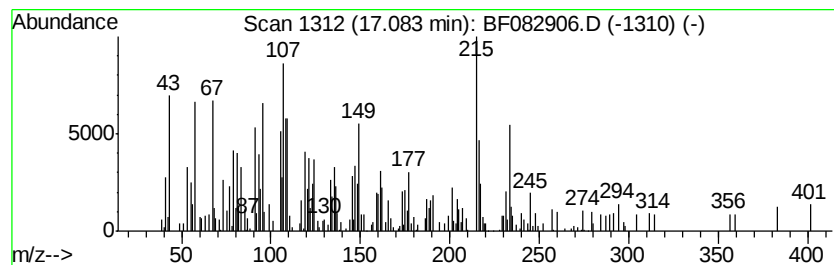
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,2-Benzenedicarboxylic aci... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.08	3.37 ng	243575	Perylene-d12	15.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Benzenedicarboxylic acid, 3,...	302	C8H2Cl4O4	000632-58-6	90
2		Cyclopent[e]-1,3-oxazine, octahy...	216	C13H16N2O	1000138-92-2	35
3		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	20
4		Cyclopentanecarboxylic acid, 3-e...	152	C9H12O2	108451-43-0	15
5		2-Cyclohexen-3-ol-1-one, 2-benzoyl-	216	C13H12O3	061834-43-3	15



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082906.D
Acq On : 13 Nov 2015 19:29
Operator : UM/IZ
Sample : G4370-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P001-TS003-01

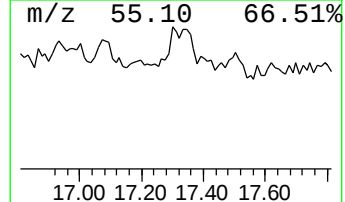
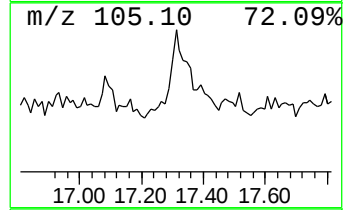
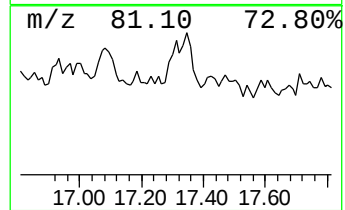
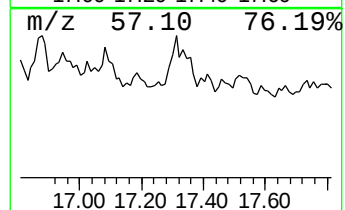
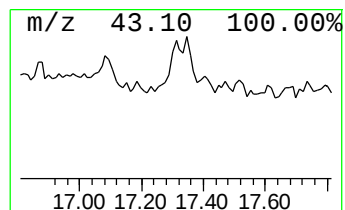
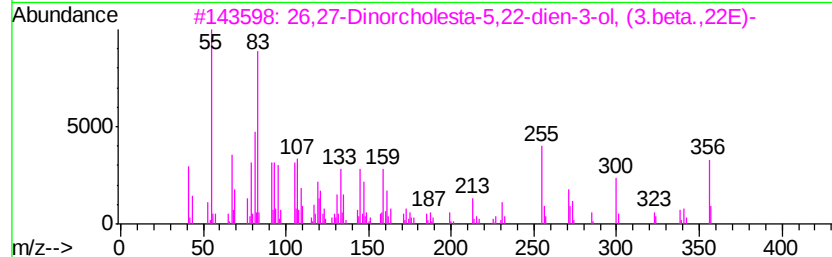
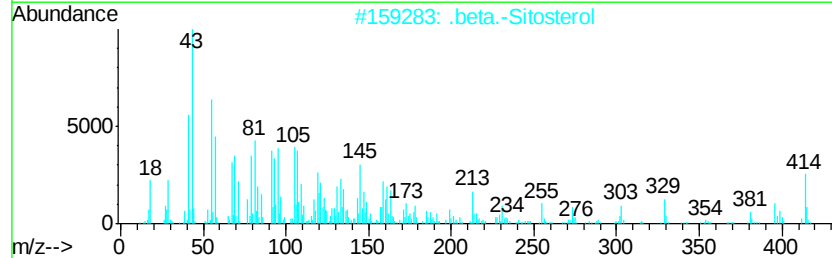
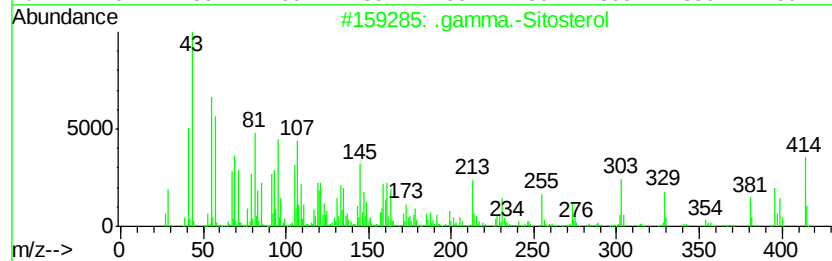
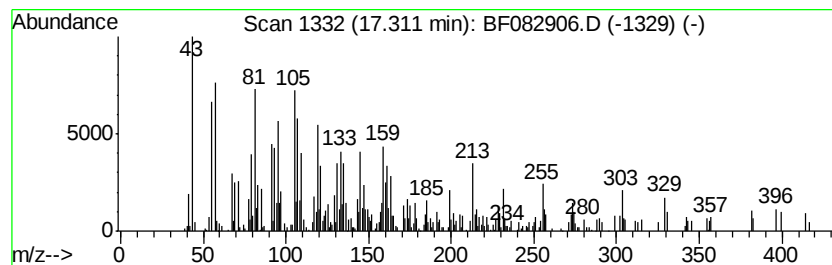
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .gamma.-Sitosterol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.31	3.83 ng	276819	Perylene-d12	15.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Sitosterol	414	C29H50O	000083-47-6	97
2			.beta.-Sitosterol	414	C29H50O	000083-46-5	95
3			26,27-Dinorcholesta-5,22-dien-3-...	356	C25H40O	057597-10-1	37
4			17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	30
5			Cholestan-3-one, 4,4-dimethyl-, ...	414	C29H50O	002097-85-0	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111315\
 Data File : BF082906.D
 Acq On : 13 Nov 2015 19:29
 Operator : UM/IZ
 Sample : G4370-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-TS003-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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...	4.97	25.1 ng		1972230	1	6.80	1572290	20.0
unknown6.57	6.57	43.5 ng		3418870	1	6.80	1572290	20.0
Hexadecane	10.76	2.7 ng		270114	4	11.32	2033720	20.0
n-Hexadecanoic acid	11.87	6.3 ng		645751	4	11.32	2033720	20.0
9,12-Octadecadien...	12.57	35.0 ng		3563980	4	11.32	2033720	20.0
Octadecanoic acid	12.66	7.4 ng		610100	5	13.96	1639290	20.0
Thiophene, 2-hexyl-	13.07	2.1 ng		172318	5	13.96	1639290	20.0
1-Heneicosyl formate	13.83	3.9 ng		315212	5	13.96	1639290	20.0
Oxirane, [(hexade...	14.45	4.4 ng		360325	5	13.96	1639290	20.0
Tetracosane	15.08	4.1 ng		298906	6	15.38	1446950	20.0
5-(7a-Isopropenyl...	16.05	5.2 ng		372800	6	15.38	1446950	20.0
1,2-Benzenedicarb...	17.08	3.4 ng		243575	6	15.38	1446950	20.0
.gamma.-Sitosterol	17.31	3.8 ng		276819	6	15.38	1446950	20.0