

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
 Data File : BF082069.D
 Acq On : 6 Oct 2015 3:31
 Operator : UM/IZ
 Sample : G3891-16
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-2(0-24)

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-BF092815.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	5.086	255	258	269	rVB	831880	1238828	44.00%	5.161%
2	5.497	291	294	296	rBV	2092167	2485429	88.28%	10.355%
3	6.526	381	384	386	rBV	2108016	2419961	85.96%	10.083%
4	6.652	393	395	398	rBV	1986510	2345939	83.33%	9.774%
5	6.880	413	415	418	rBV	415506	593529	21.08%	2.473%
6	7.040	427	429	432	rBV	1909697	1902348	67.57%	7.926%
7	7.452	462	465	468	rBV	1078065	1380889	49.05%	5.753%
8	7.532	471	472	479	rVB	20483	37421	1.33%	0.156%
9	8.172	526	528	531	rBV	885067	762151	27.07%	3.175%
10	9.258	620	623	625	rBV	2029498	2815380	100.00%	11.730%
11	9.646	655	657	659	rBV	504375	400562	14.23%	1.669%
12	9.932	679	682	684	rBV	934170	826077	29.34%	3.442%
13	10.355	718	719	721	rVB	63968	45942	1.63%	0.191%
14	10.572	735	738	741	rBV	21156	29197	1.04%	0.122%
15	10.721	748	751	754	rBV	1370394	1347981	47.88%	5.616%
16	10.823	758	760	765	rVB	52968	86990	3.09%	0.362%
17	11.281	797	800	801	rBV	43564	45962	1.63%	0.191%
18	11.418	810	812	814	rBV	760041	671506	23.85%	2.798%
19	11.703	835	837	839	rBV	38702	31713	1.13%	0.132%
20	11.943	856	858	865	rBV2	97314	115257	4.09%	0.480%
21	13.006	948	951	954	rBV	2150737	2051169	72.86%	8.546%
22	13.898	1027	1029	1033	rBV	294311	342432	12.16%	1.427%
23	14.058	1041	1043	1049	rVV	529619	598353	21.25%	2.493%
24	14.161	1049	1052	1054	rVV3	20232	37508	1.33%	0.156%
25	14.218	1055	1057	1061	rVV2	18631	40463	1.44%	0.169%
26	14.538	1082	1085	1087	rBV2	58860	111972	3.98%	0.467%
27	14.812	1107	1109	1112	rBV2	47240	69838	2.48%	0.291%
28	14.972	1121	1123	1126	rBV	125873	131612	4.67%	0.548%
29	15.155	1137	1139	1141	rBV	46724	76649	2.72%	0.319%
30	15.190	1141	1142	1145	rVB	79556	101396	3.60%	0.422%
31	15.521	1168	1171	1174	rBV	258092	436342	15.50%	1.818%
32	15.738	1187	1190	1196	rBV2	96793	163064	5.79%	0.679%
33	15.955	1207	1209	1212	rBV	37627	65303	2.32%	0.272%
34	16.024	1212	1215	1218	rVV	59381	121344	4.31%	0.506%

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ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-2(0-24)

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-BF092815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.744	1276	1278	1283	rVB3	36752	71082	2.52%	0.296%
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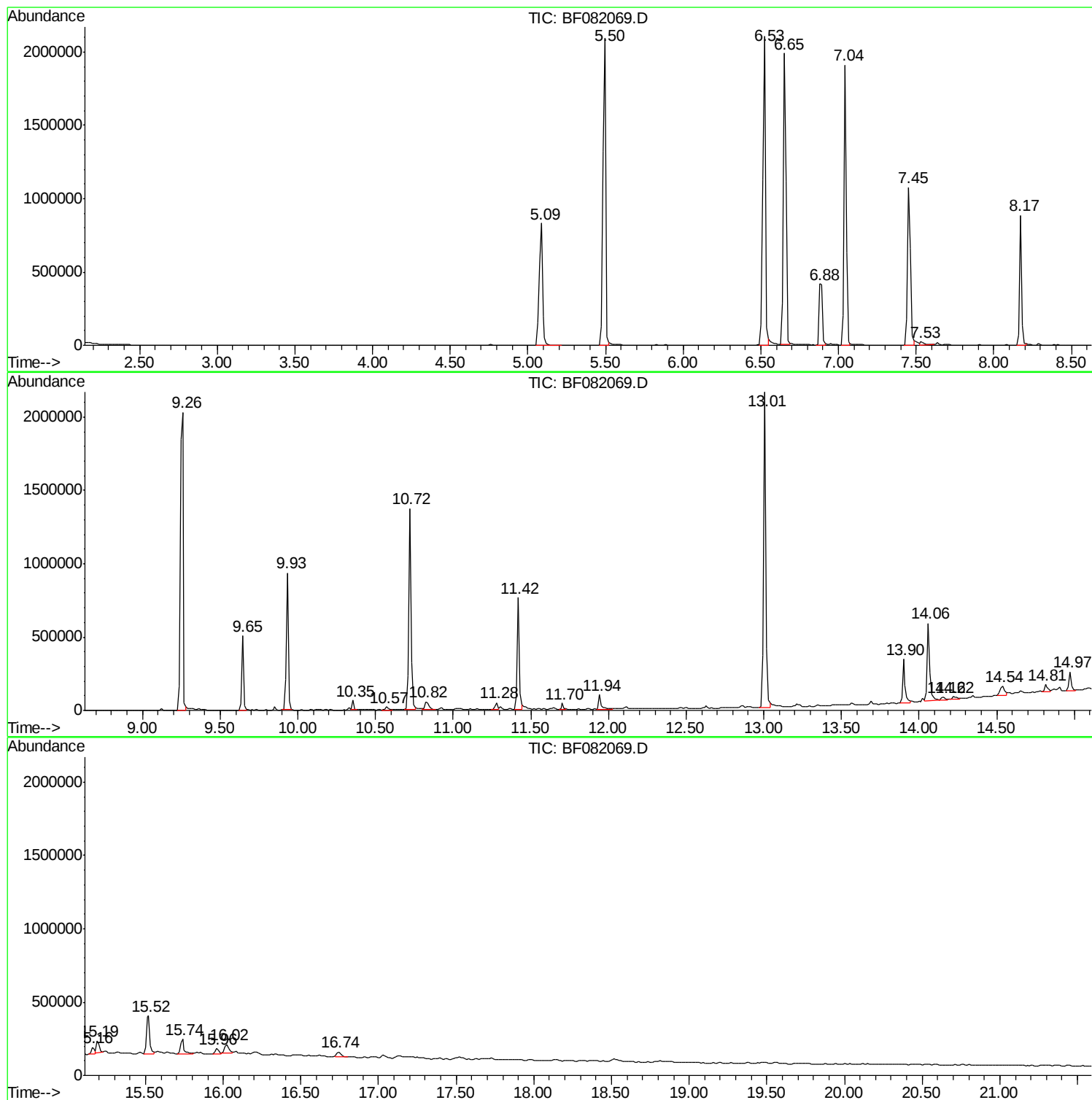
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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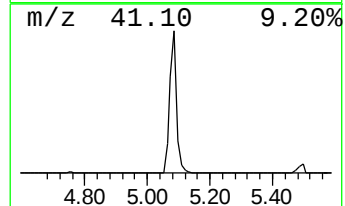
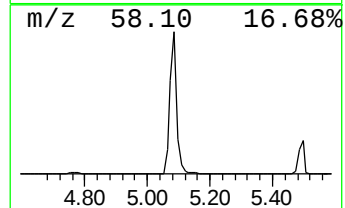
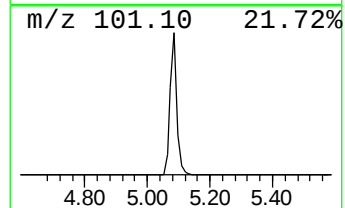
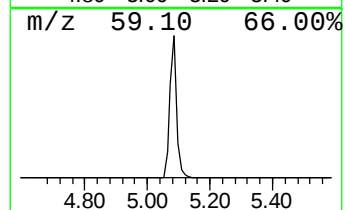
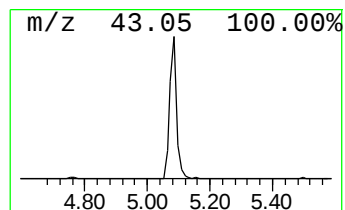
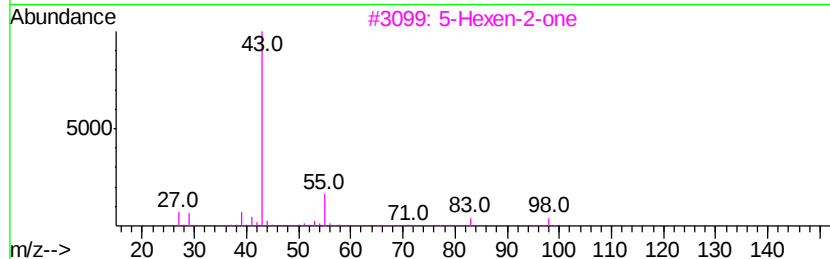
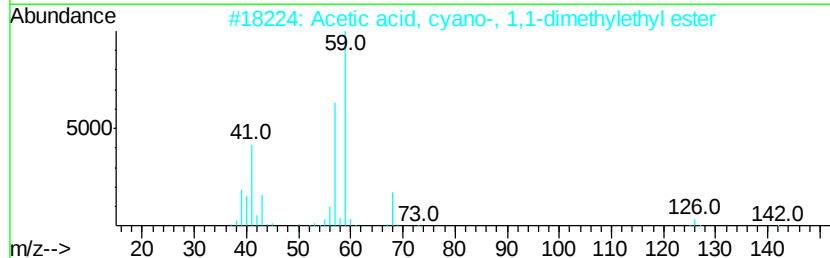
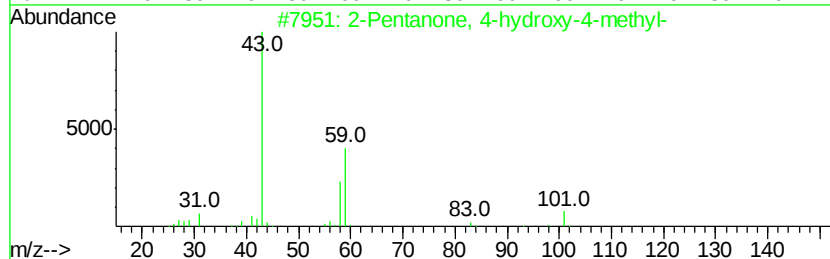
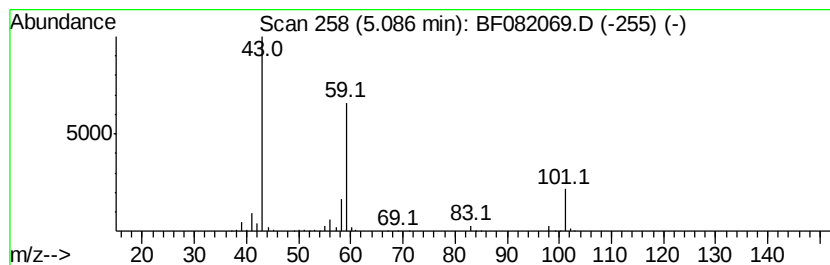
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	41.74 ng	1238830	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Acetone	58	C3H6O	000067-64-1	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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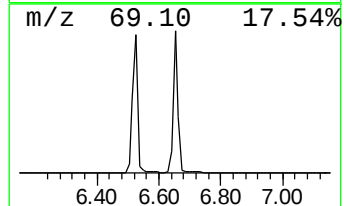
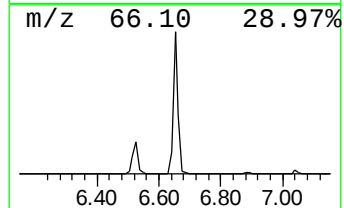
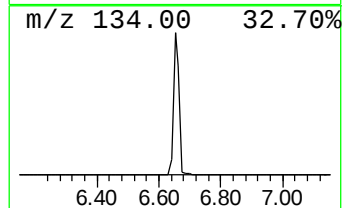
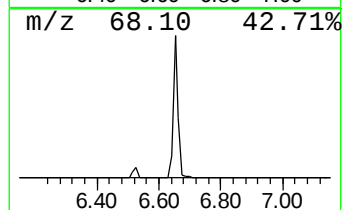
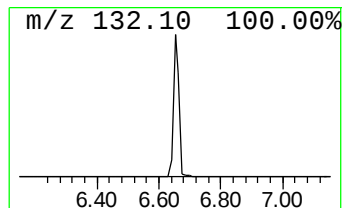
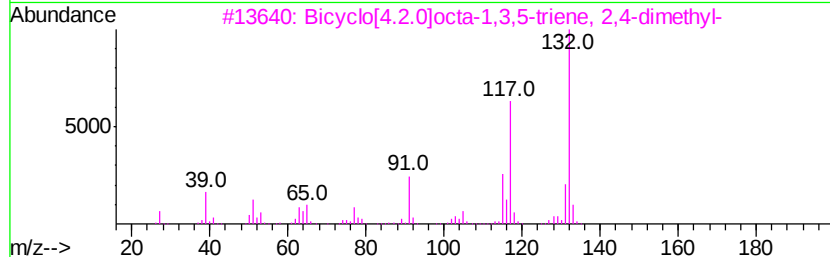
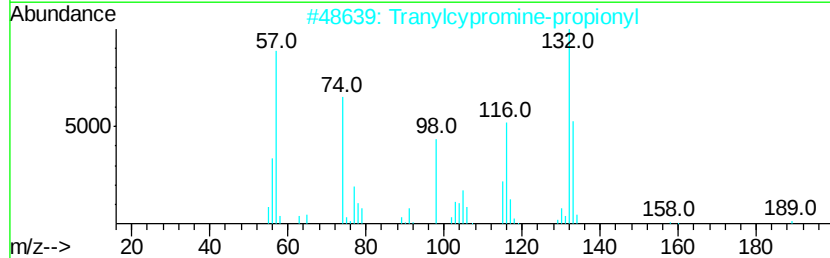
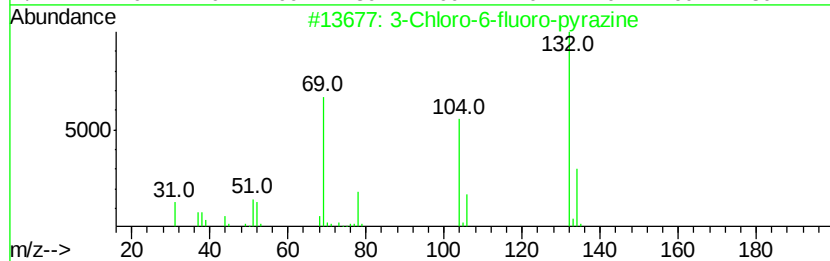
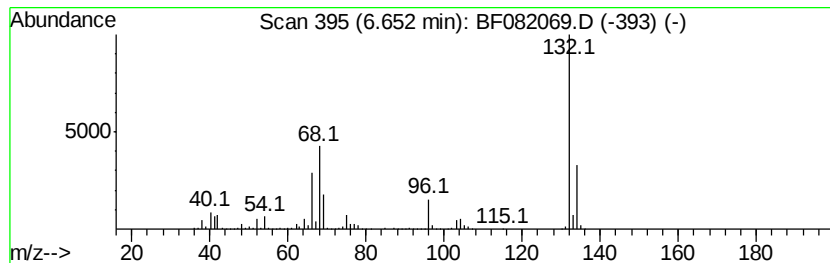
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	79.05 ng	2345940	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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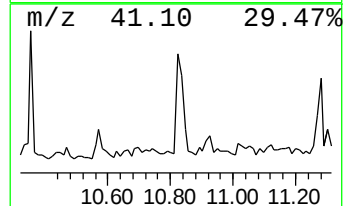
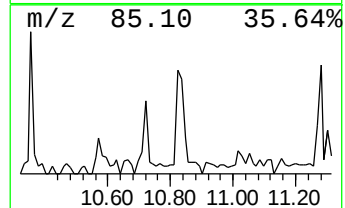
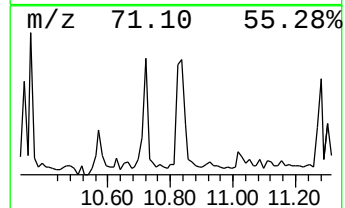
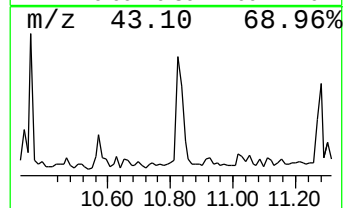
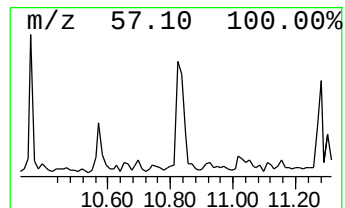
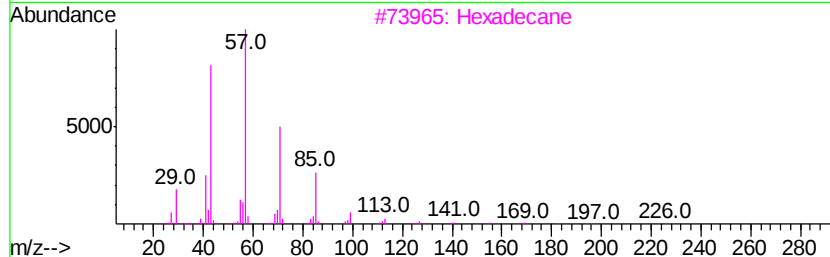
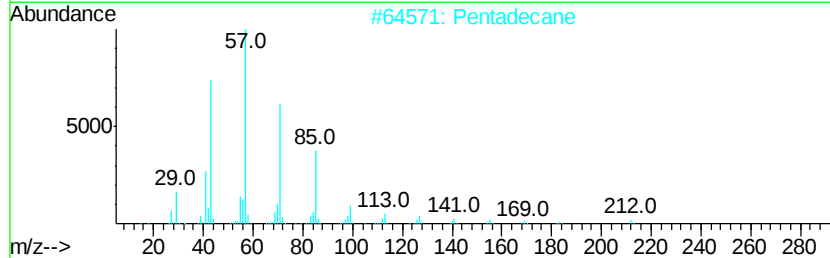
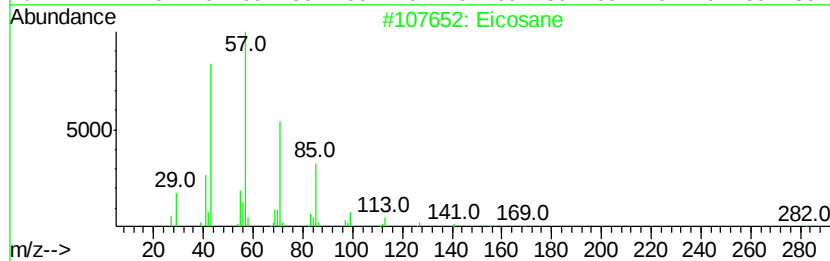
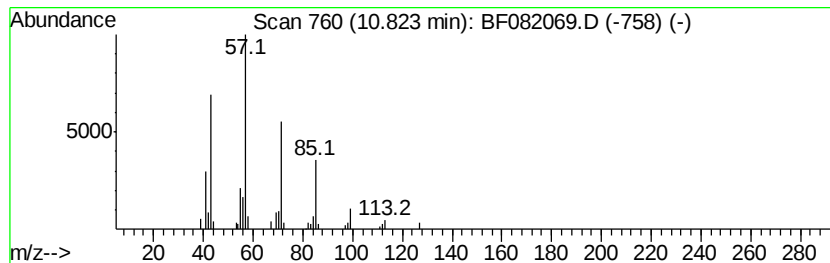
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Eicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.82	2.59 ng	86990	Phenanthrene-d10	11.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Pentadecane	212	C15H32	000629-62-9	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tetradecane	198	C14H30	000629-59-4	86
5	Tridecane	184	C13H28	000629-50-5	83



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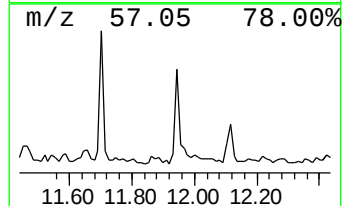
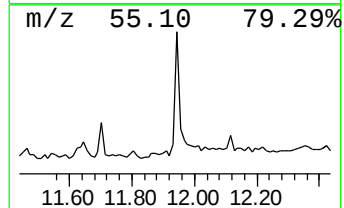
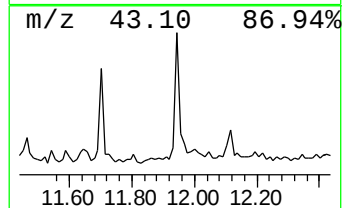
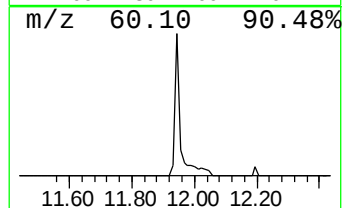
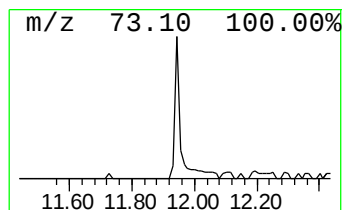
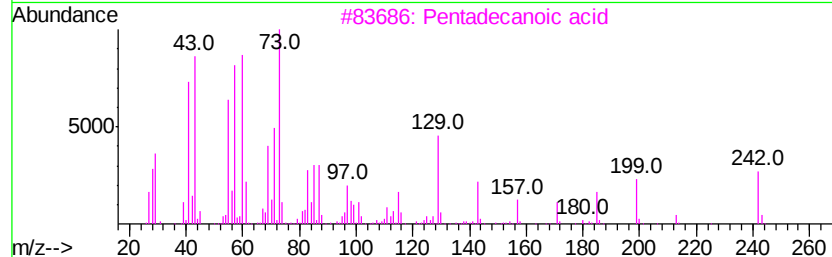
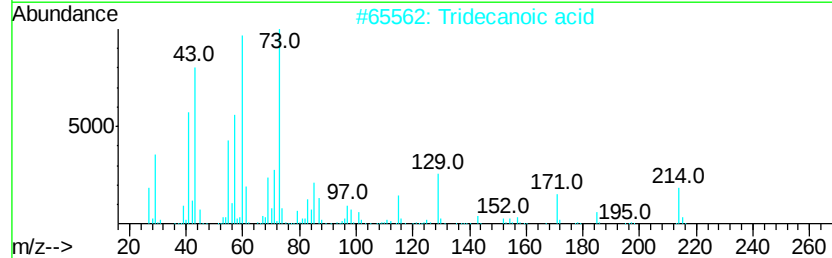
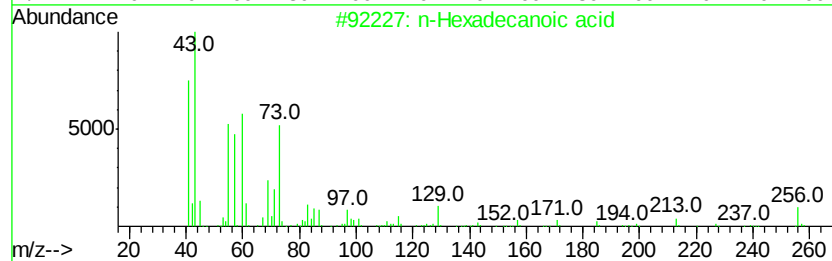
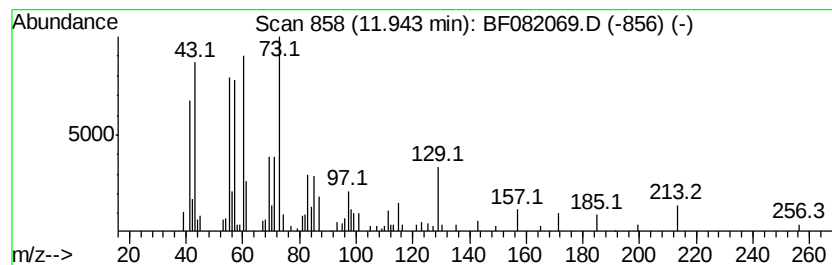
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.94	3.43 ng	115257	Phenanthrene-d10		11.42
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	Tridecanoic acid	214	C13H26O2	000638-53-9	80
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	74
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	Dodecane, 1,2-dibromo-	326	C12H24Br2	055334-42-4	18



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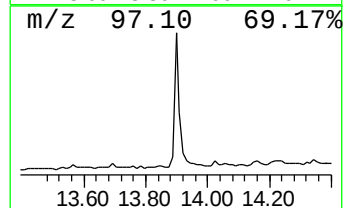
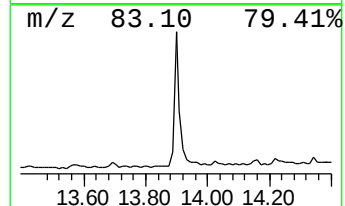
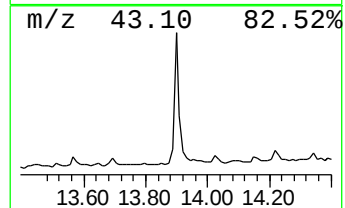
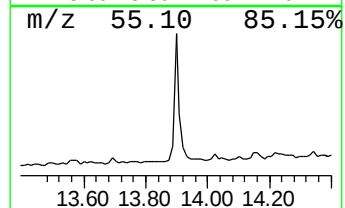
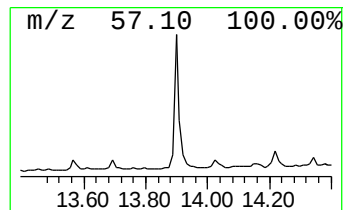
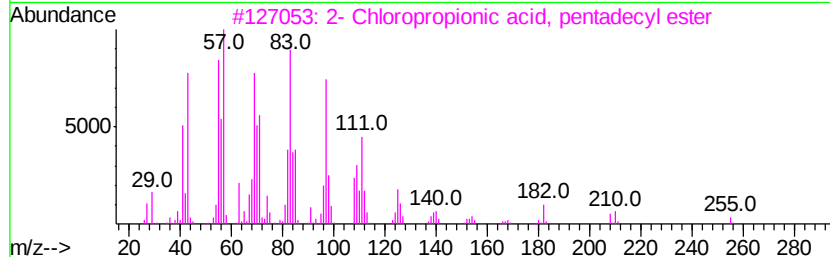
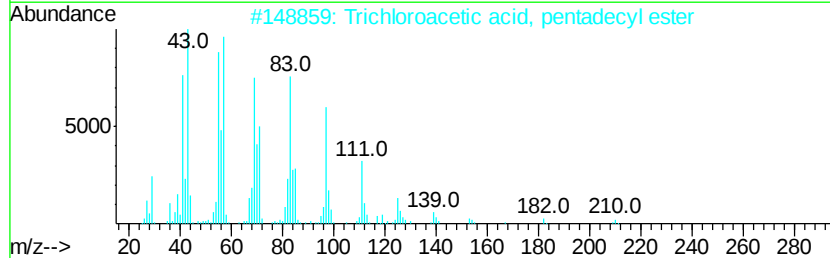
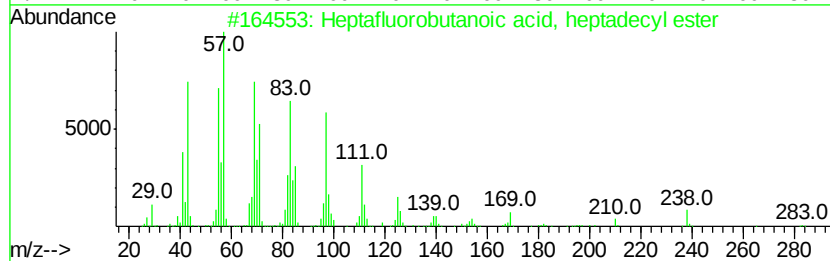
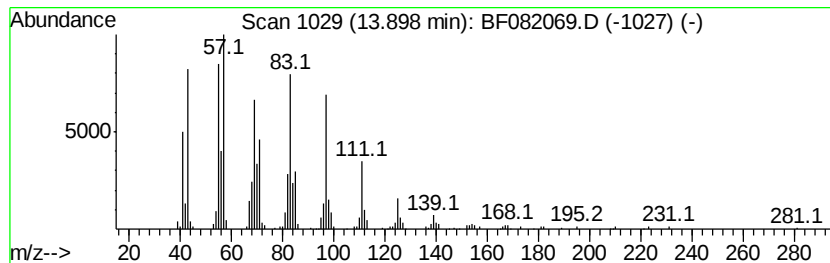
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptafluorobutanoic acid, h... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	11.45 ng	342432	Chrysene-d12	14.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
3		2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	91
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
5		1-Docosene	308	C22H44	001599-67-3	91



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Misc :
ALS Vial : 26 Sample Multiplier: 1

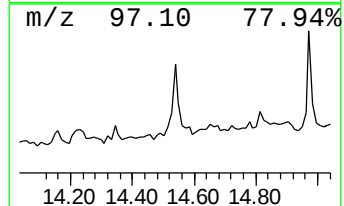
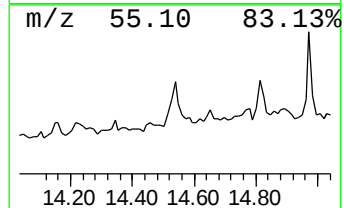
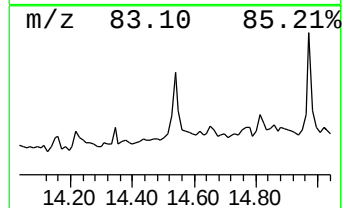
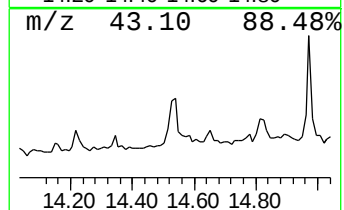
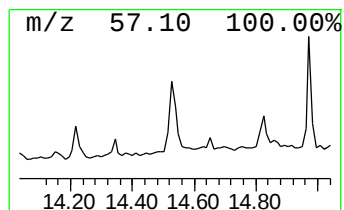
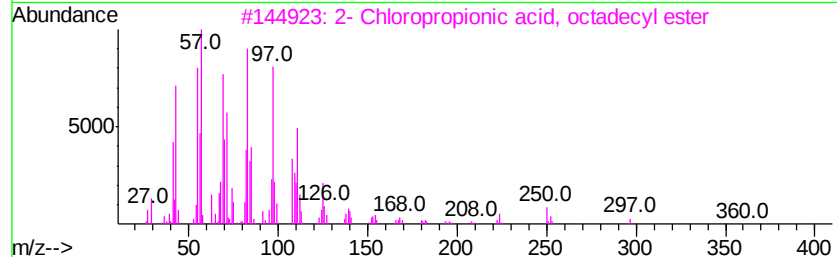
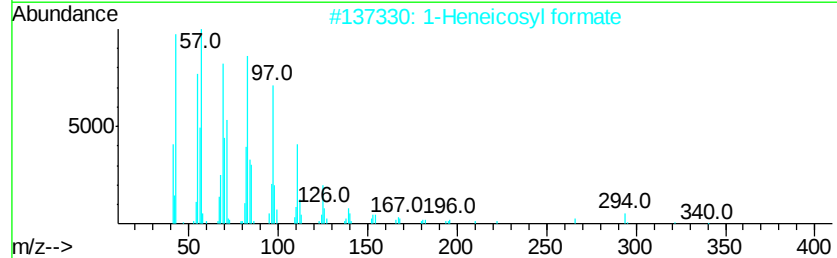
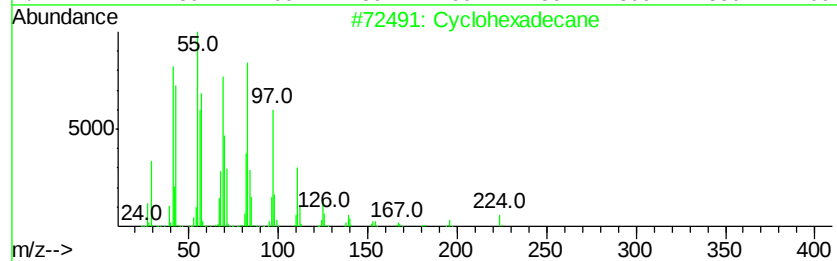
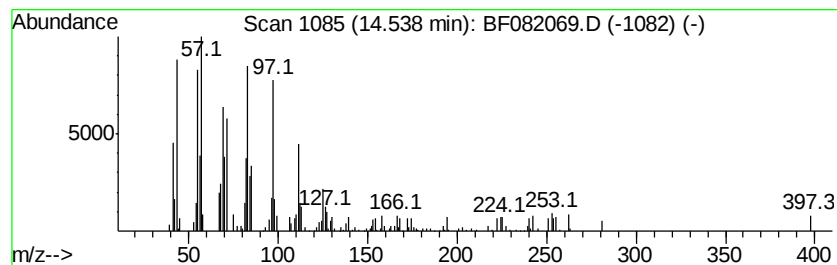
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ClientSampleId :
SS-2(0-24)

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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 Cyclohexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	3.74 ng	111972	Chrysene-d12	14.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 98
2		1-Heneicosyl formate	340 C22H44O2	077899-03-7 94
3		2- Chloropropionic acid, octadec...	360 C21H41ClO2	088104-31-8 94
4		1-Octadecanol	270 C18H38O	000112-92-5 93
5		Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1 91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
Data File : BF082069.D
Acq On : 6 Oct 2015 3:31
Operator : UM/IZ
Sample : G3891-16
Misc :
ALS Vial : 26 Sample Multiplier: 1

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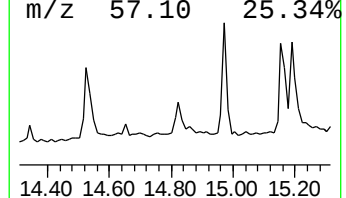
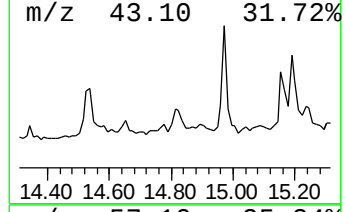
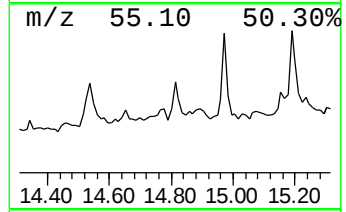
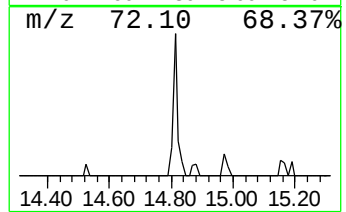
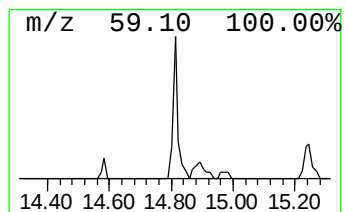
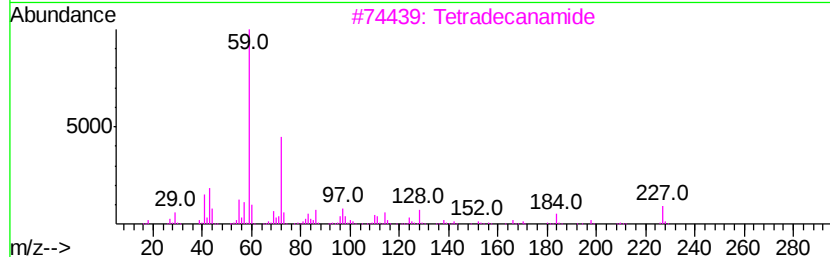
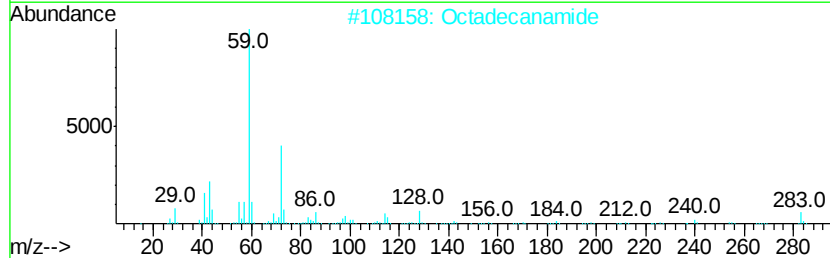
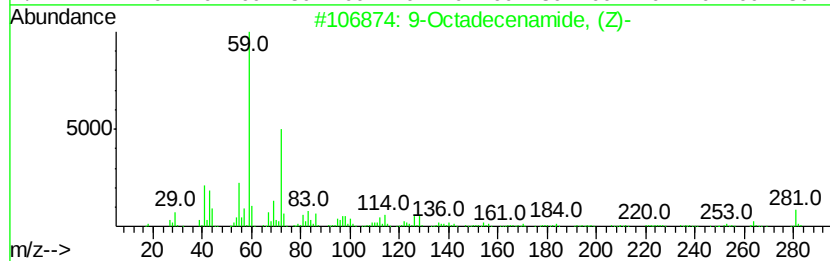
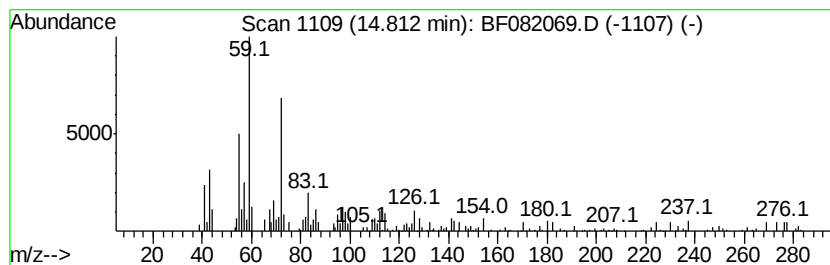
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	3.20 ng	69838	Perylene-d12	15.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
2		Octadecanamide	283	C18H37NO	000124-26-5	59
3		Tetradecanamide	227	C14H29NO	000638-58-4	59
4		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	59
5		Dodecanamide	199	C12H25NO	001120-16-7	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
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Acq On : 6 Oct 2015 3:31
Operator : UM/IZ
Sample : G3891-16
Misc :
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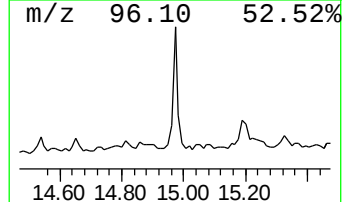
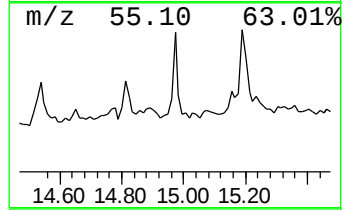
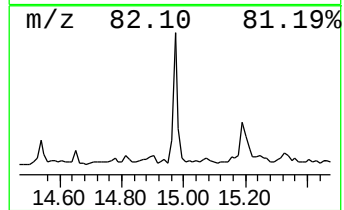
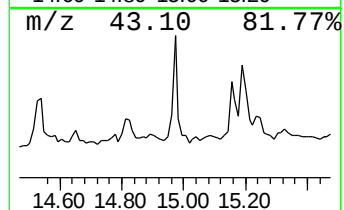
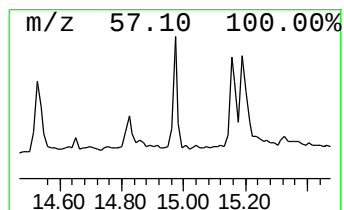
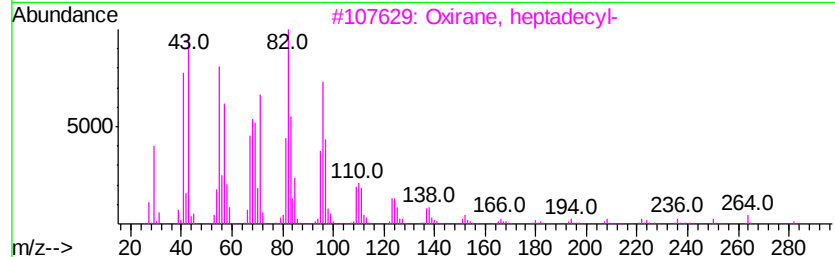
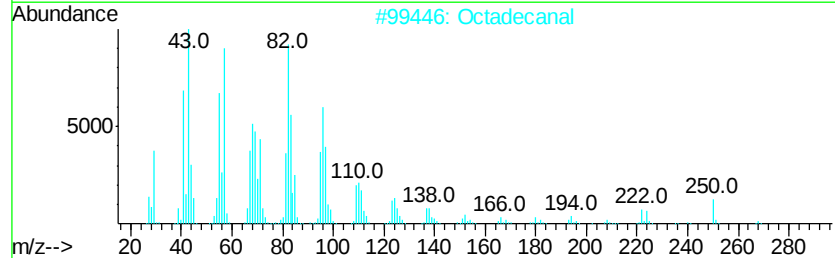
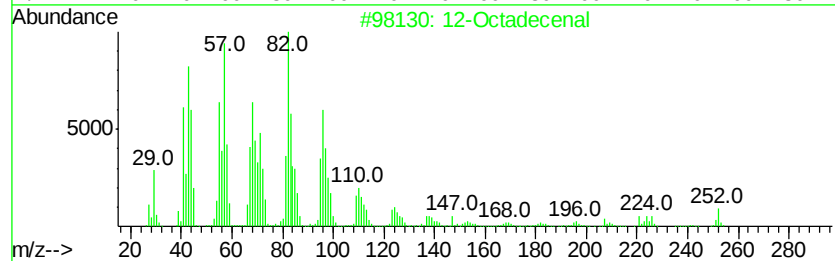
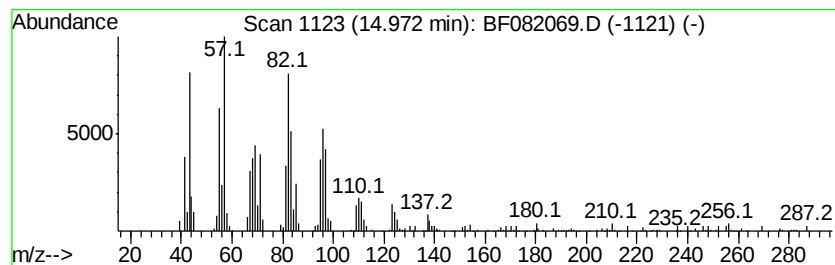
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 12-Octadecenal Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.97	6.03 ng	131612	Perylene-d12		15.52
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	12-Octadecenal	266	C18H34O	056554-91-7	90
2	Octadecanal	268	C18H36O	000638-66-4	87
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	74
4	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	68
5	11-Tetradecyn-1-ol acetate	252	C16H28O2	033925-72-3	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
Data File : BF082069.D
Acq On : 6 Oct 2015 3:31
Operator : UM/IZ
Sample : G3891-16
Misc :
ALS Vial : 26 Sample Multiplier: 1

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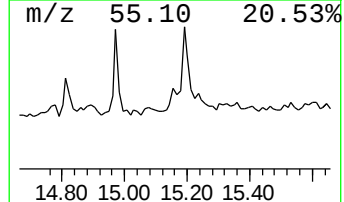
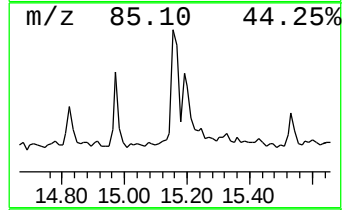
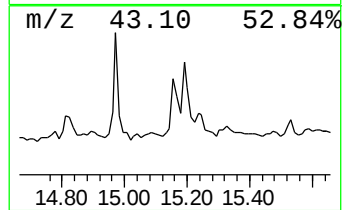
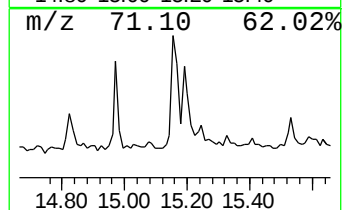
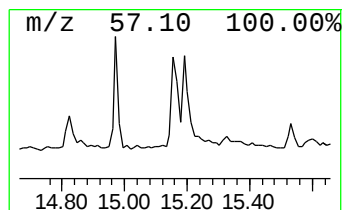
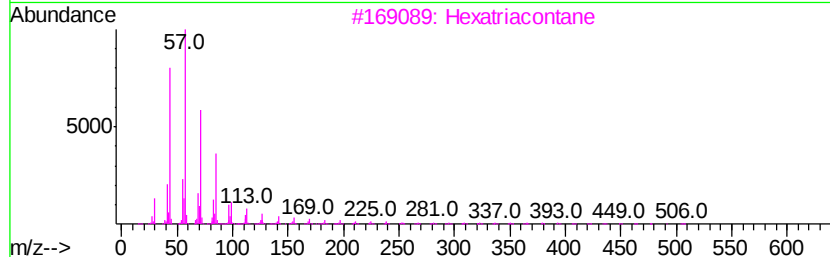
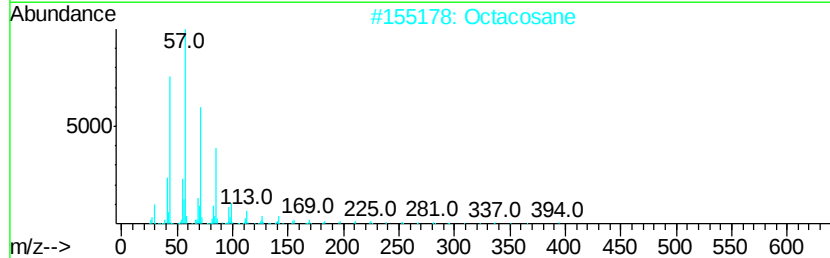
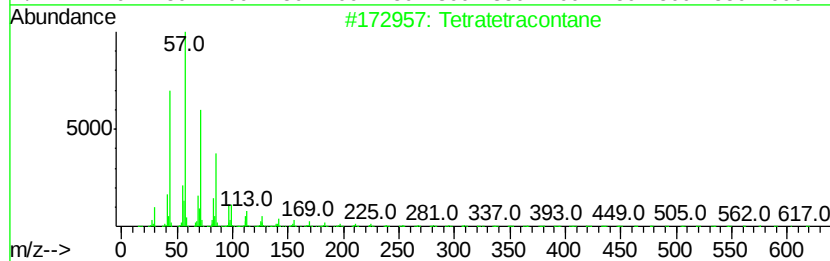
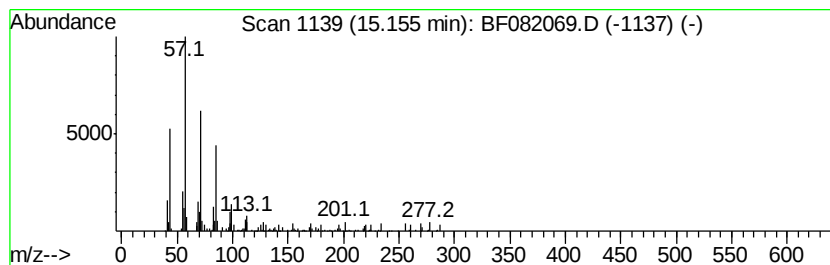
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetratetracontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	3.51 ng	76649	Perylene-d12	15.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	83
2			Octacosane	394	C28H58	000630-02-4	83
3			Hexatriacontane	507	C36H74	000630-06-8	83
4			Tetratriacontane	479	C34H70	014167-59-0	83
5			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
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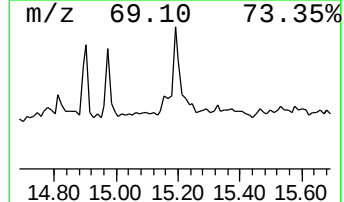
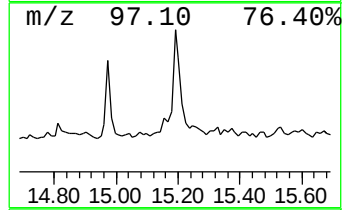
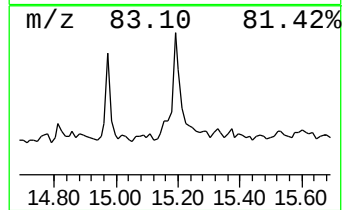
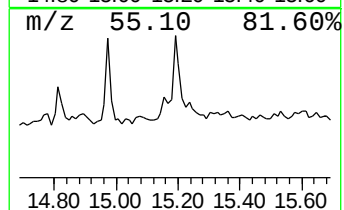
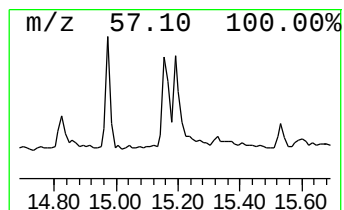
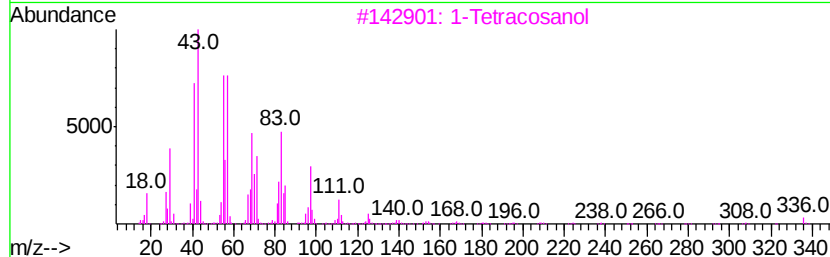
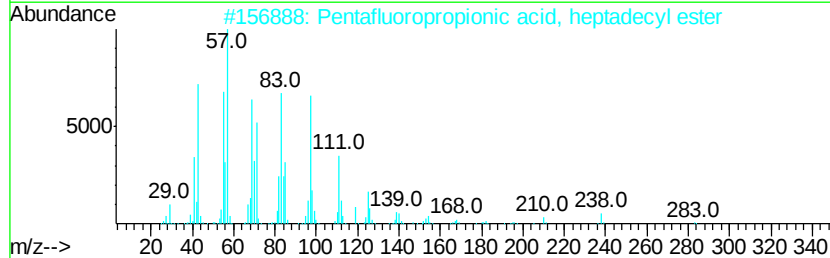
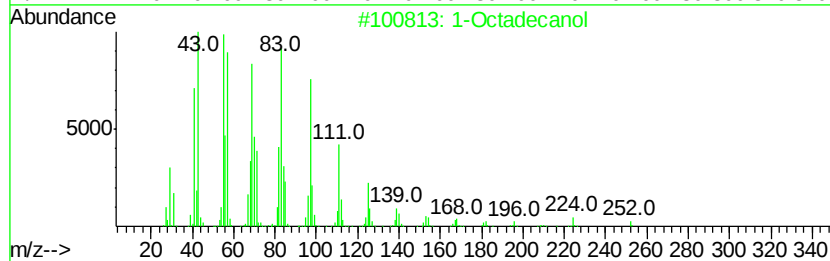
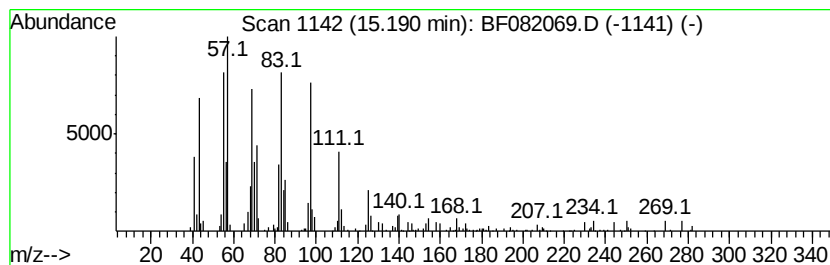
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Octadecanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	4.65 ng	101396	Perylene-d12	15.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Octadecanol	270 C18H38O	000112-92-5	90
2	Pentafluoropropionic acid, hepta...	402 C20H35F5O2	1000283-04-2	76
3	1-Tetracosanol	354 C24H50O	000506-51-4	72
4	Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1	62
5	1-Heptadecanol	256 C17H36O	001454-85-9	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
Data File : BF082069.D
Acq On : 6 Oct 2015 3:31
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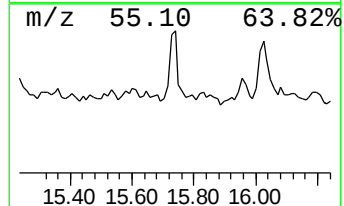
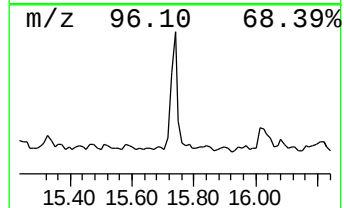
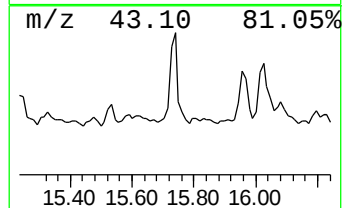
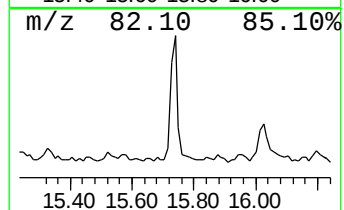
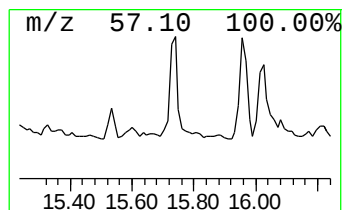
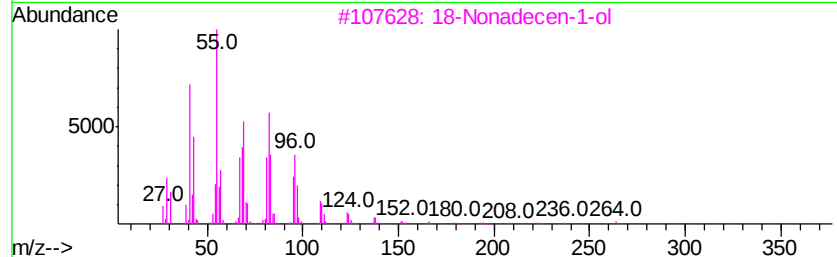
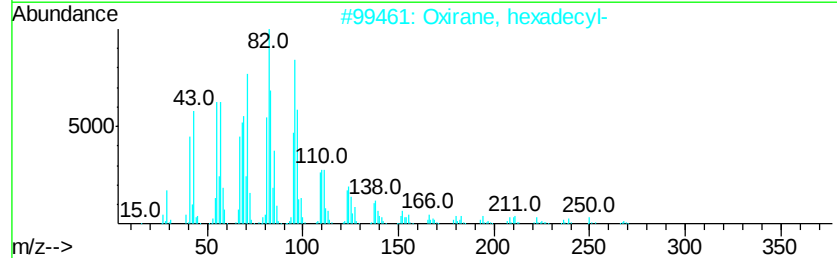
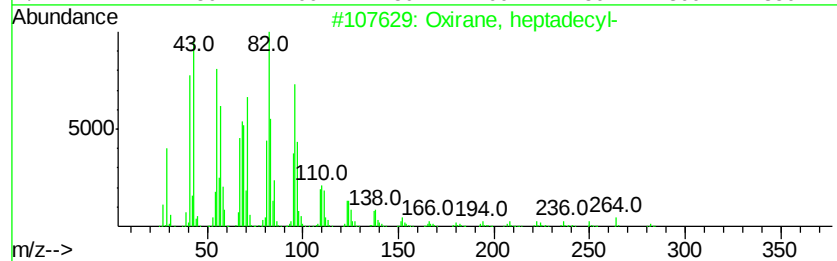
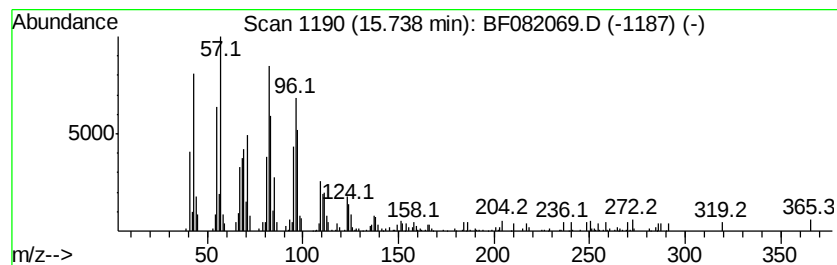
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Oxirane, heptadecyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	7.47 ng	163064	Perylene-d12	15.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
3		18-Nonadecen-1-ol	282	C19H38O	1000142-89-2	72
4		Octadecanal	268	C18H36O	000638-66-4	64
5		E-7-Dodecen-1-ol acetate	226	C14H26O2	1000131-35-3	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
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ClientSampleId :
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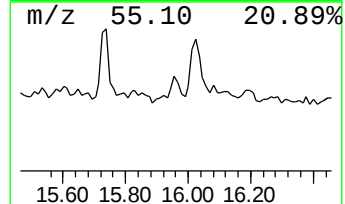
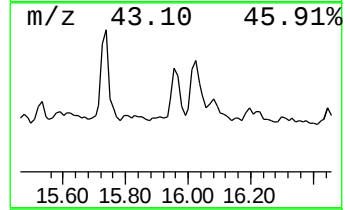
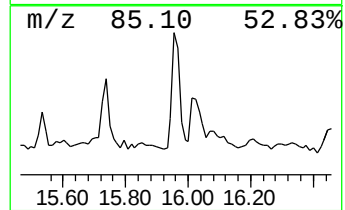
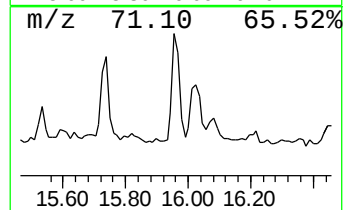
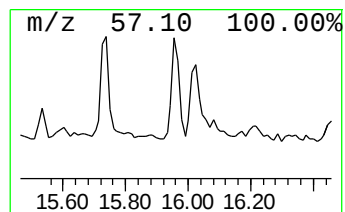
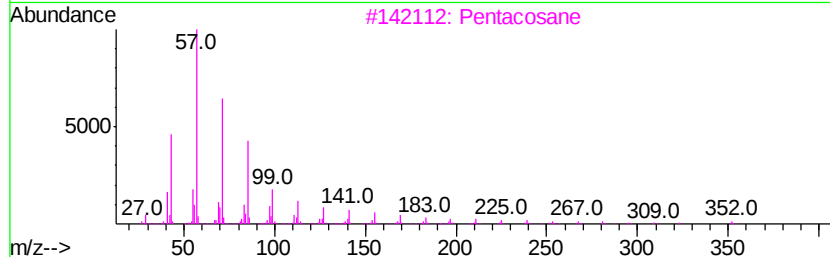
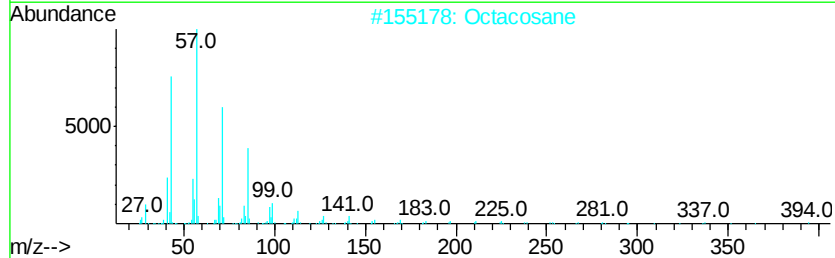
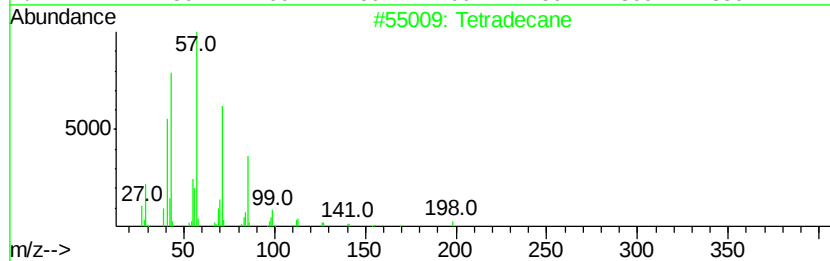
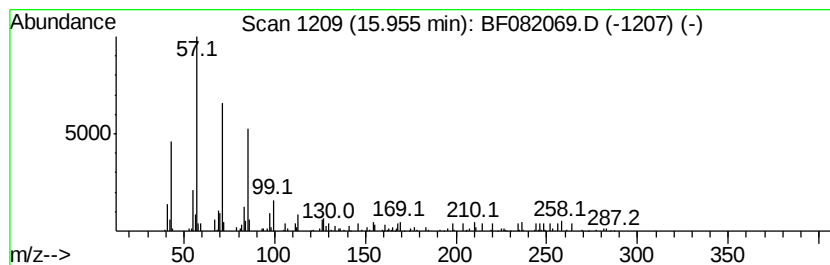
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	2.99 ng	65303	Perylene-d12	15.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	52
2		Octacosane	394	C28H58	000630-02-4	52
3		Pentacosane	352	C25H52	000629-99-2	52
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	52
5		Nonacosane	408	C29H60	000630-03-5	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
Data File : BF082069.D
Acq On : 6 Oct 2015 3:31
Operator : UM/IZ
Sample : G3891-16
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-2(0-24)

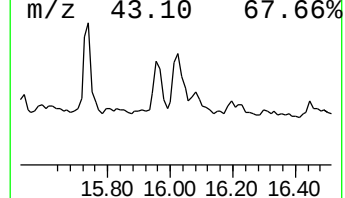
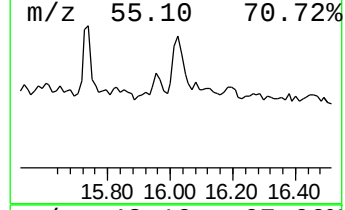
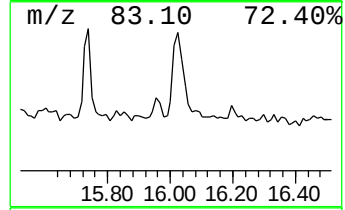
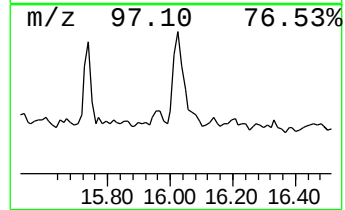
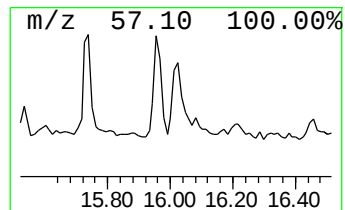
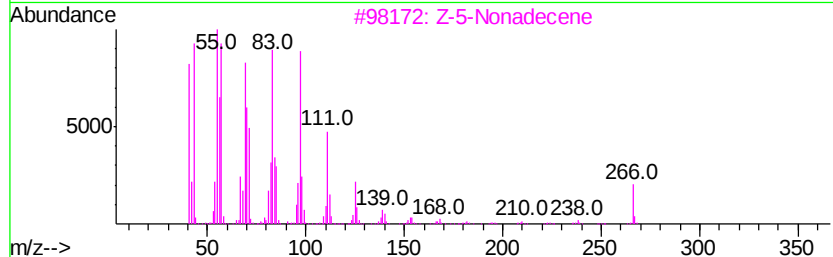
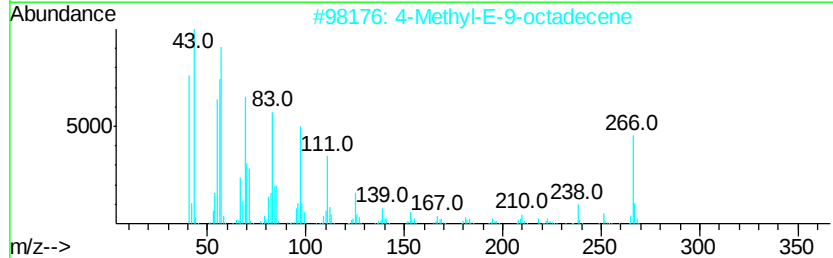
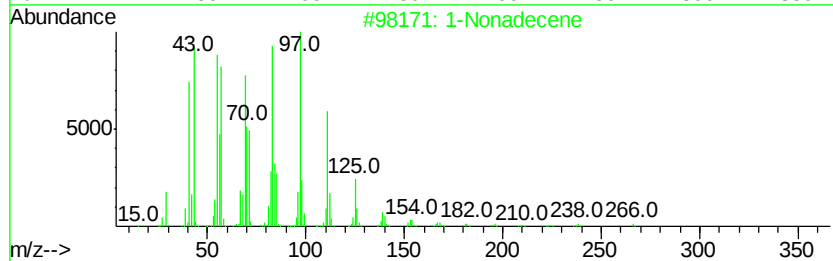
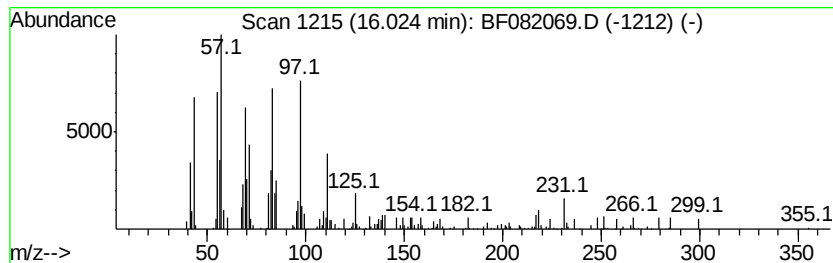
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Nonadecene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.02	5.56 ng	121344	Perylene-d12	15.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	87
2			4-Methyl-E-9-octadecene	266	C19H38	1000130-87-1	87
3			Z-5-Nonadecene	266	C19H38	1000131-11-8	86
4			Cyclooctadecane, ethyl-	280	C20H40	1000151-22-5	74
5			Cyclotetradecane, 1,7,11-trimeth...	280	C20H40	001786-12-5	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
Data File : BF082069.D
Acq On : 6 Oct 2015 3:31
Operator : UM/IZ
Sample : G3891-16
Misc :
ALS Vial : 26 Sample Multiplier: 1

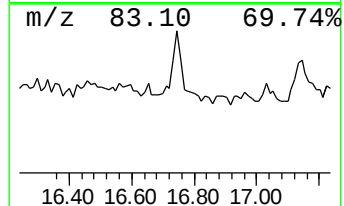
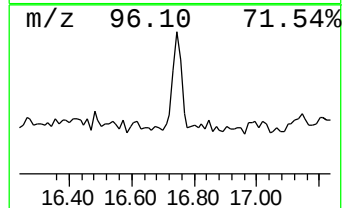
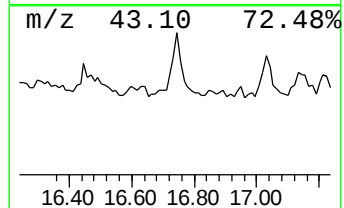
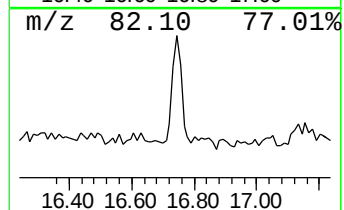
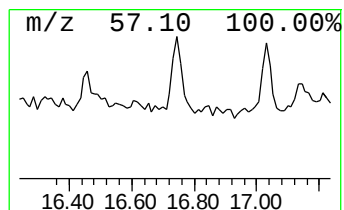
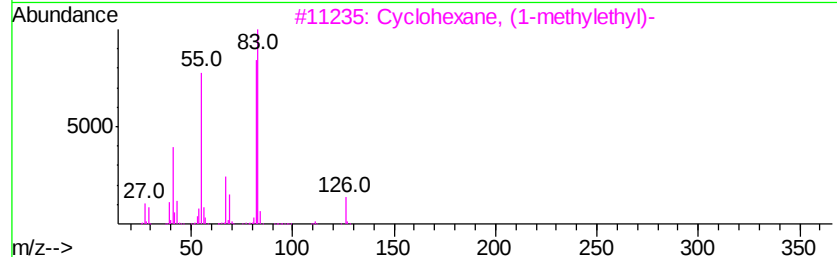
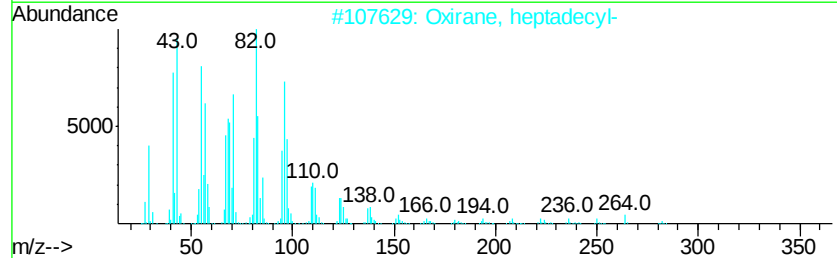
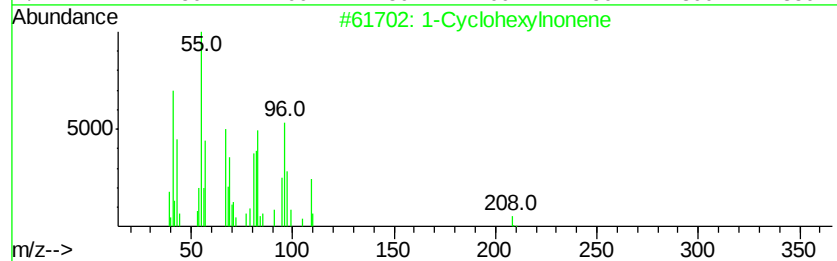
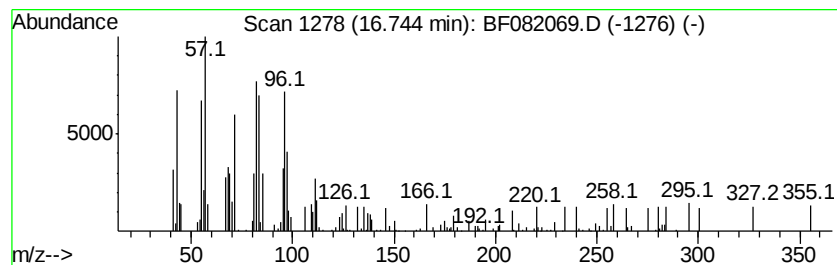
Instrument :
BNA_F
ClientSampleId :
SS-2(0-24)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Cyclohexylnonene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.74	3.26 ng	71082	Perylene-d12	15.52	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Cyclohexyl-nonene	208	C15H28	114614-84-5	50
2	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	49
3	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	38
4	Octadecanal	268	C18H36O	000638-66-4	38
5	Cyclobuta[1,2:3,4]dicyclooctene,...	220	C16H28	061277-45-0	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
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 Operator : UM/IZ
 Sample : G3891-16
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-2(0-24)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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...	5.09	41.7	ng	1238830	1	6.89	593529	20.0
unknown6.65	6.65	79.0	ng	2345940	1	6.89	593529	20.0
Eicosane	10.82	2.6	ng	86990	4	11.42	671506	20.0
n-Hexadecanoic acid	11.94	3.4	ng	115257	4	11.42	671506	20.0
Heptafluorobutano...	13.90	11.4	ng	342432	5	14.06	598353	20.0
Cyclohexadecane	14.54	3.7	ng	111972	5	14.06	598353	20.0
9-Octadecenamide, ...	14.81	3.2	ng	69838	6	15.52	436342	20.0
12-Octadecenal	14.97	6.0	ng	131612	6	15.52	436342	20.0
Tetratetracontane	15.16	3.5	ng	76649	6	15.52	436342	20.0
1-Octadecanol	15.19	4.7	ng	101396	6	15.52	436342	20.0
Oxirane, heptadecyl-	15.74	7.5	ng	163064	6	15.52	436342	20.0
Tetradecane	15.96	3.0	ng	65303	6	15.52	436342	20.0
1-Nonadecene	16.02	5.6	ng	121344	6	15.52	436342	20.0
1-Cyclohexylnonene	16.74	3.3	ng	71082	6	15.52	436342	20.0