

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
 Data File : BF083185.D
 Acq On : 23 Nov 2015 4:02
 Operator : UM/IZ
 Sample : G4475-03
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 H3

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-BF112115.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	2.227	20	21	35	rVB2	41868	133418	2.09%	0.236%
2	4.799	242	246	262	rVB	2217756	3624655	56.65%	6.420%
3	5.256	284	286	289	rBV	5553454	6110445	95.50%	10.823%
4	6.319	376	379	382	rBV	5557935	6398567	100.00%	11.333%
5	6.422	386	388	391	rBV	5174868	5852505	91.47%	10.366%
6	6.650	406	408	410	rBV	1318969	1234801	19.30%	2.187%
7	6.810	419	422	424	rBV	4347476	4679691	73.14%	8.289%
8	7.222	455	458	460	rBV	3828100	3963067	61.94%	7.019%
9	7.942	518	521	523	rBV	1124244	1483517	23.19%	2.628%
10	8.593	575	578	582	rVB2	49964	74807	1.17%	0.132%
11	8.890	602	604	608	rBV	46227	78184	1.22%	0.138%
12	9.016	612	615	618	rBV	5511754	5402230	84.43%	9.568%
13	9.416	648	650	652	rBV	1736774	1483698	23.19%	2.628%
14	9.633	667	669	671	rBV	110429	98561	1.54%	0.175%
15	9.679	671	673	676	rBV	1222067	1585476	24.78%	2.808%
16	10.136	709	713	714	rBV	169508	193306	3.02%	0.342%
17	10.353	729	732	733	rBV	66492	91367	1.43%	0.162%
18	10.468	740	742	745	rBV2	2973879	3319149	51.87%	5.879%
19	10.605	751	754	758	rBV	237164	278351	4.35%	0.493%
20	11.051	789	793	794	rBV	101042	101528	1.59%	0.180%
21	11.154	800	802	805	rVV	999989	1345229	21.02%	2.383%
22	11.211	805	807	813	rVB3	32794	84311	1.32%	0.149%
23	11.634	841	844	847	rBV	82087	112121	1.75%	0.199%
24	11.714	849	851	855	rBV	758479	751251	11.74%	1.331%
25	12.205	892	894	896	rBV	66956	81492	1.27%	0.144%
26	12.411	908	912	917	rBV5	62651	185477	2.90%	0.329%
27	12.491	917	919	922	rBV	640655	688347	10.76%	1.219%
28	12.548	922	924	925	rVV	83455	96638	1.51%	0.171%
29	12.582	925	927	931	rVB	168044	180595	2.82%	0.320%
30	12.742	938	941	944	rBV	3728102	3747135	58.56%	6.637%
31	13.222	980	983	986	rBV2	40262	74165	1.16%	0.131%
32	13.291	986	989	990	rVV	68072	87846	1.37%	0.156%
33	13.657	1018	1021	1029	rBV	178432	295353	4.62%	0.523%
34	13.782	1029	1032	1035	rBV	1413922	1342755	20.99%	2.378%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.977	1047	1049	1052	rbV	46310	64882	1.01%	0.115%
36	14.274	1072	1075	1079	rbV4	70121	152735	2.39%	0.271%
37	15.154	1149	1152	1154	rbV	654337	981161	15.33%	1.738%

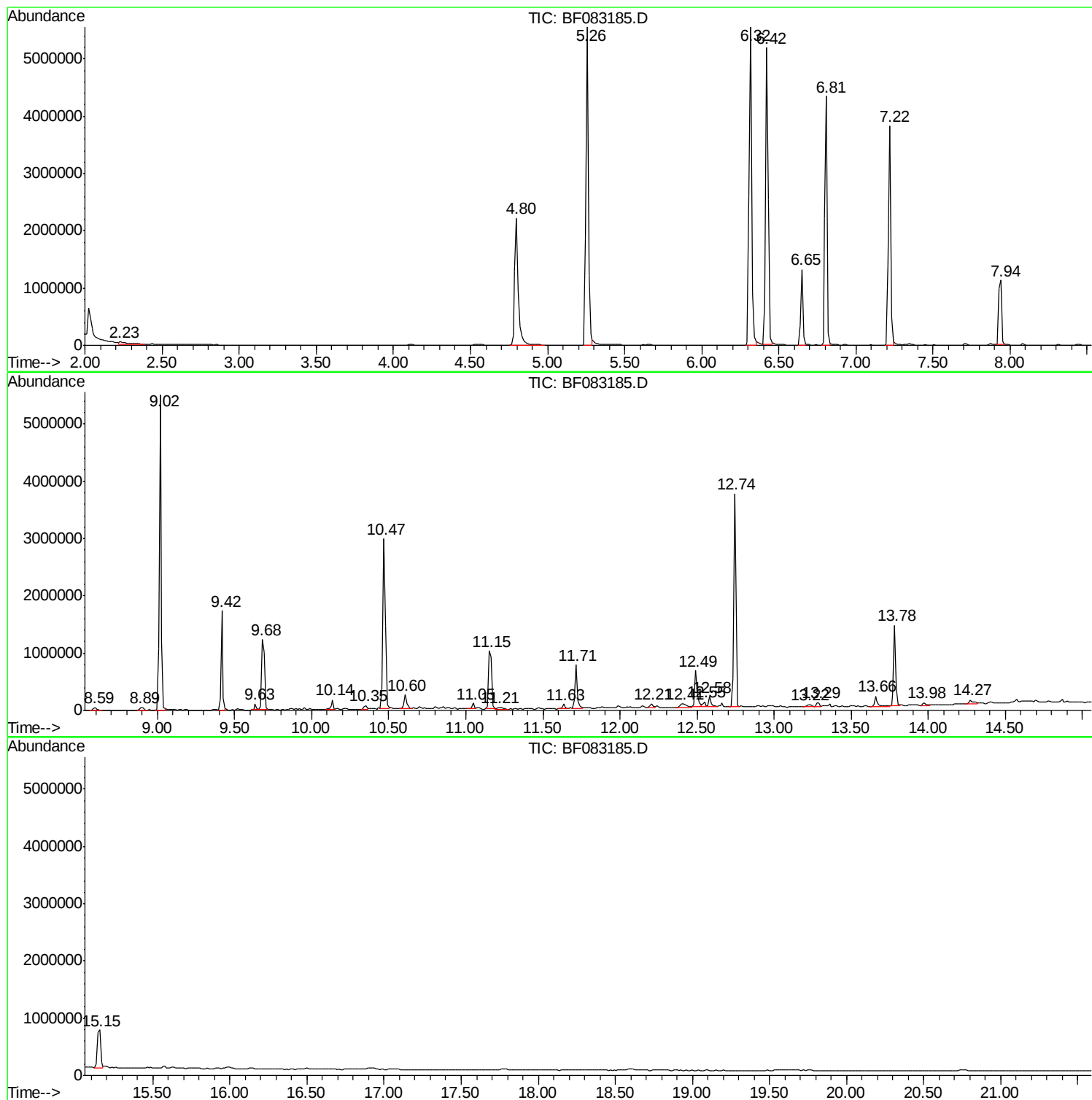
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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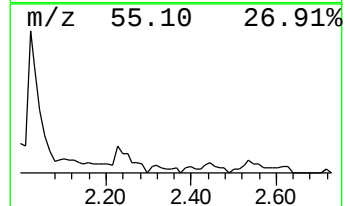
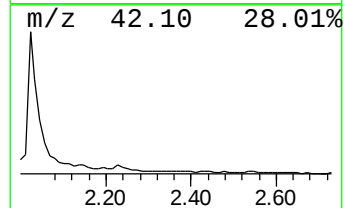
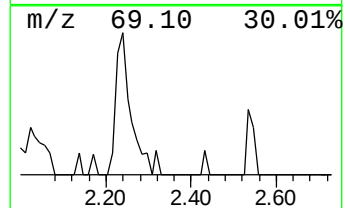
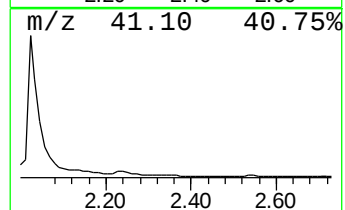
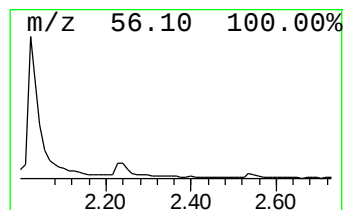
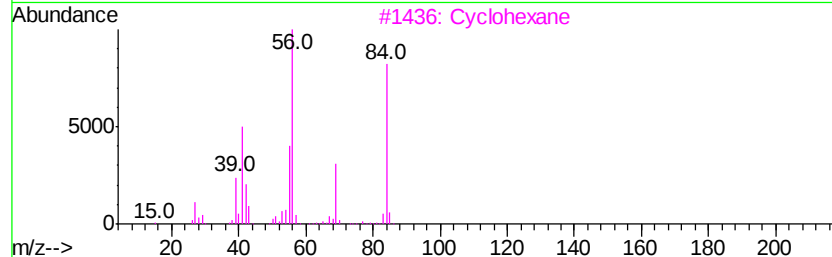
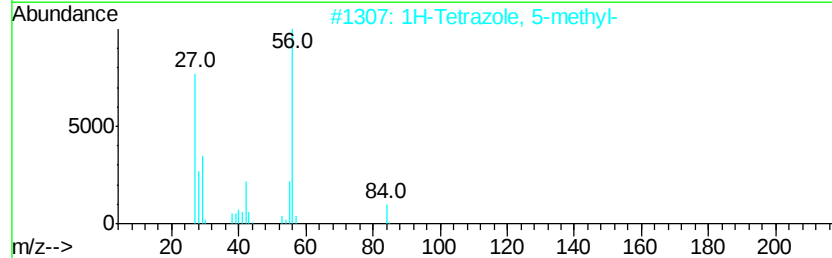
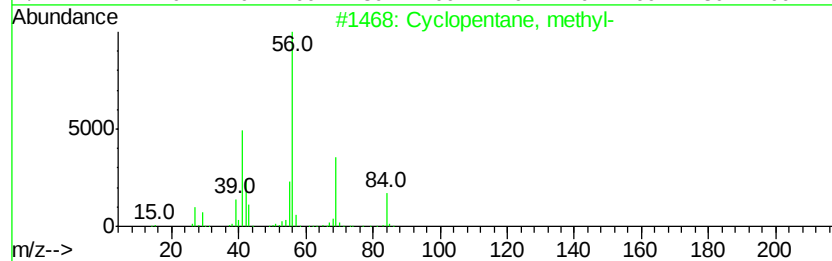
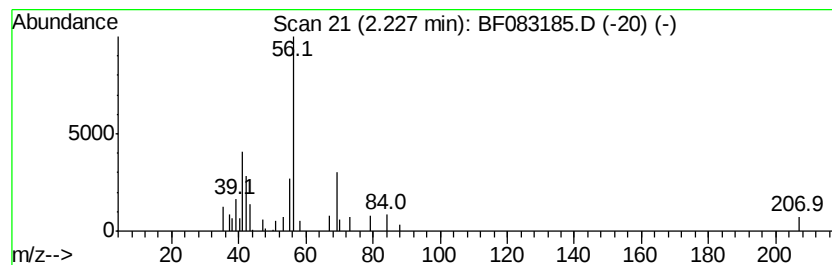
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	2.16 ng	133418	1,4-Dichlorobenzene-d4	6.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	59
2			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	59
3			Cyclohexane	84	C6H12	000110-82-7	50
4			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50
5			2-Methyl-1-nonene	140	C10H20	002980-71-4	40



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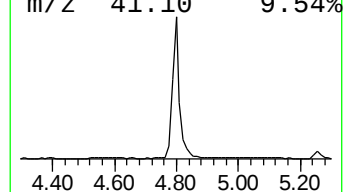
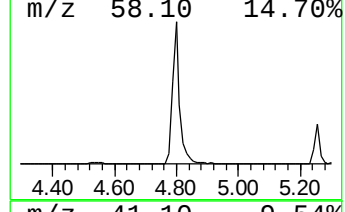
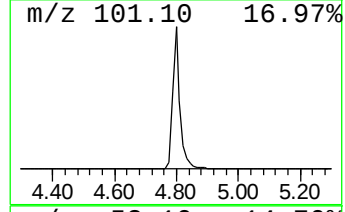
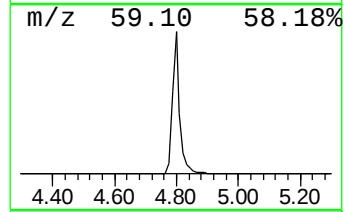
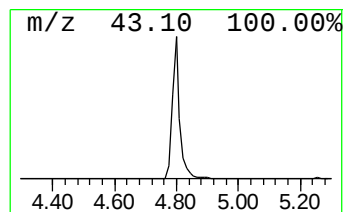
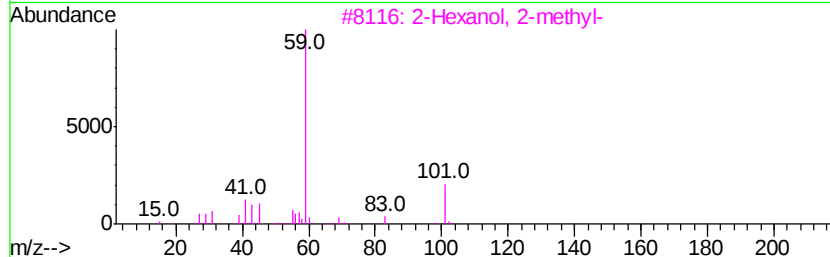
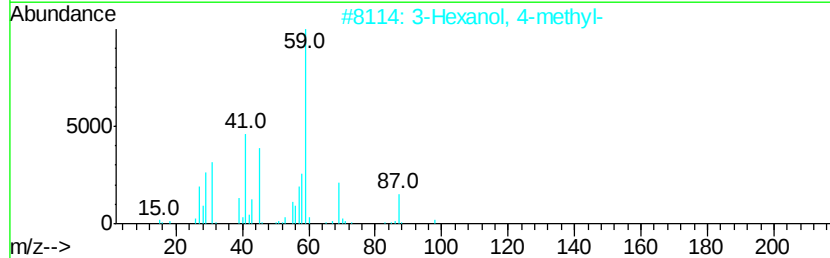
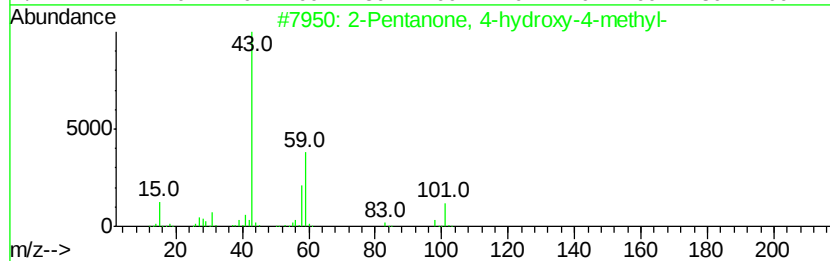
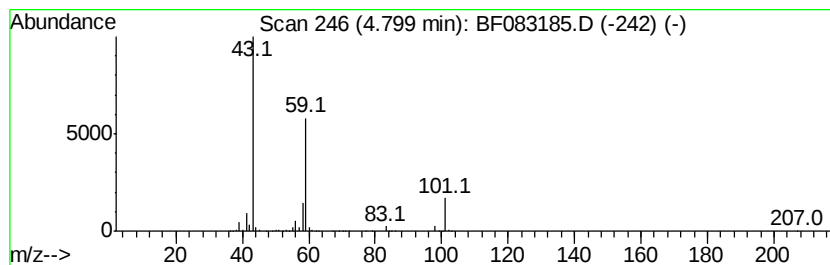
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	58.71 ng	3624660	1,4-Dichlorobenzene-d4	6.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33		
4	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32		
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9		



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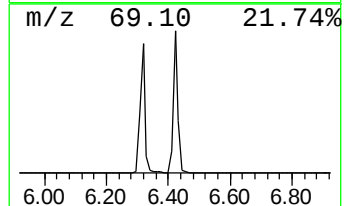
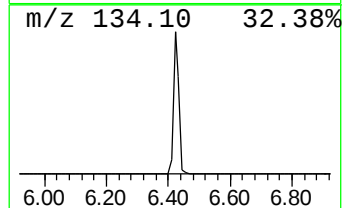
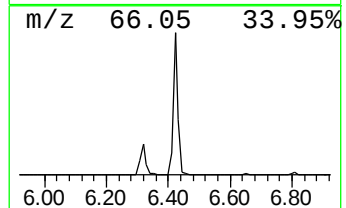
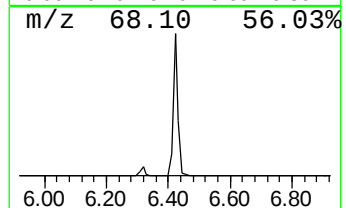
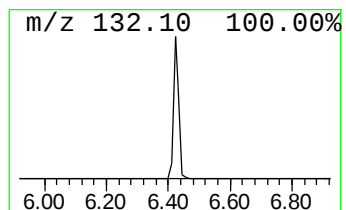
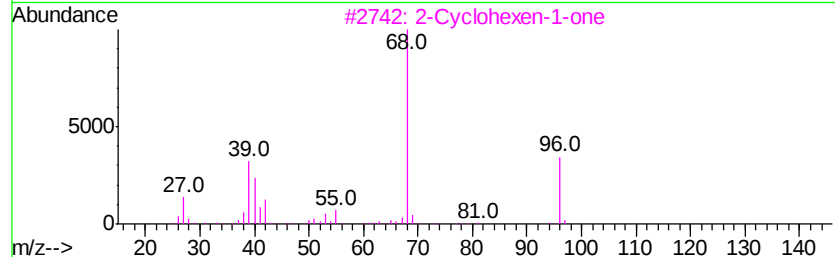
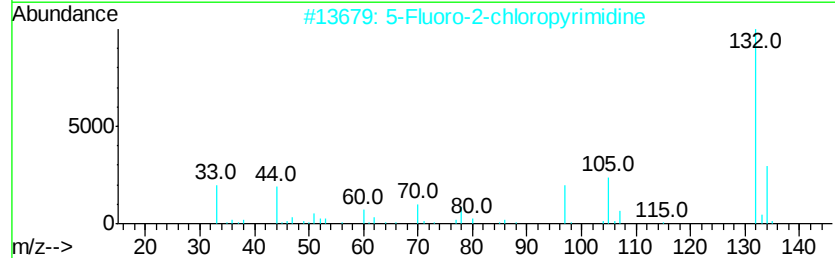
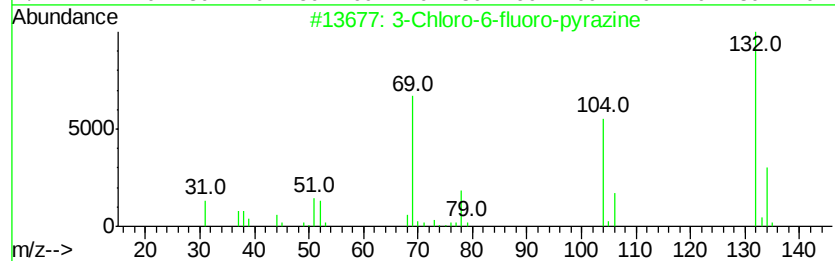
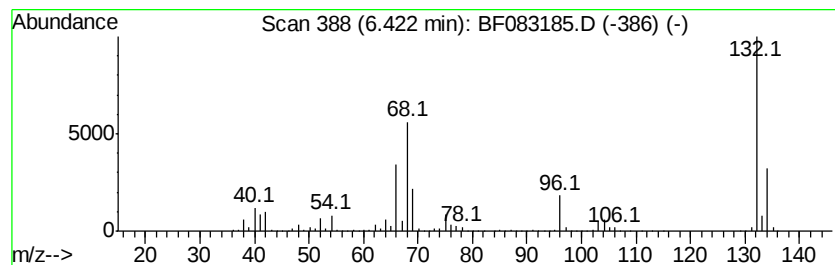
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	94.79 ng	5852510	1,4-Dichlorobenzene-d4	6.65

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	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		5-Aminoindole	132	C8H8N2	005192-03-0	10



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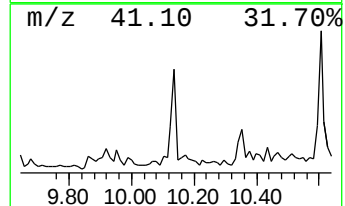
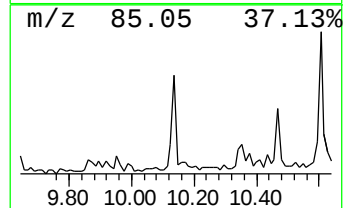
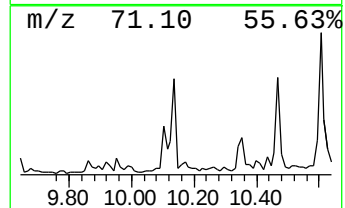
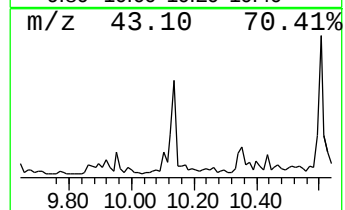
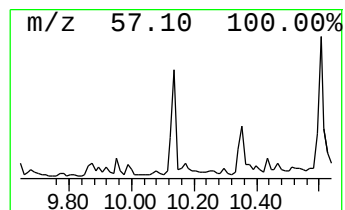
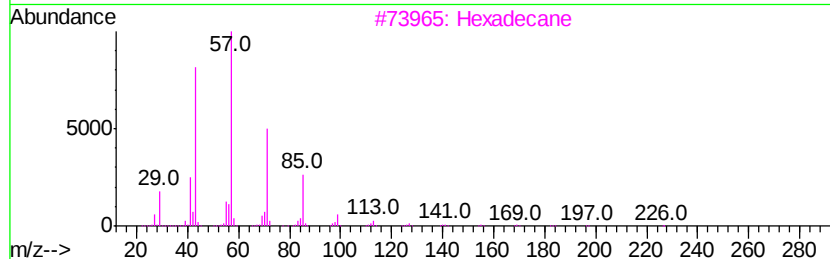
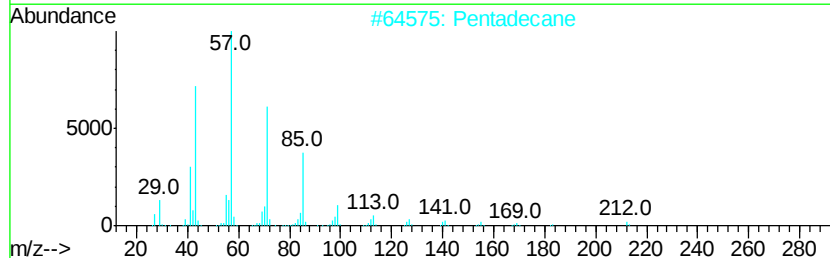
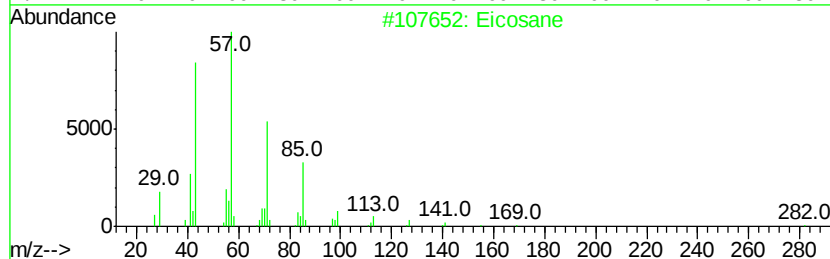
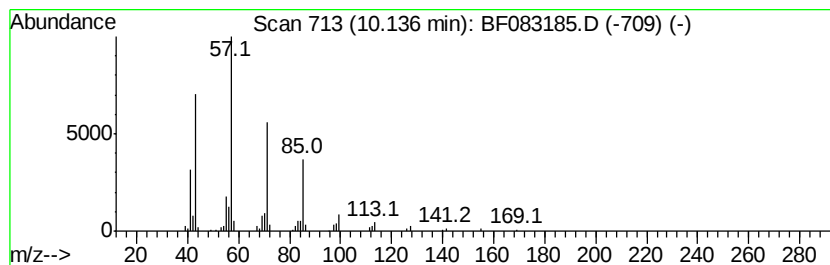
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.44 ng	193306	Acenaphthene-d10	9.68

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	Tridecane		184	C13H28	000629-50-5	90
5	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	78



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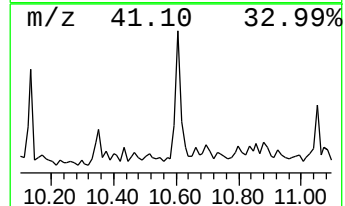
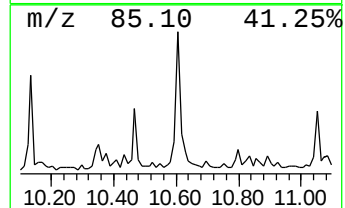
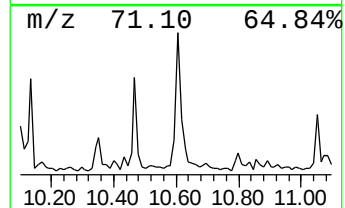
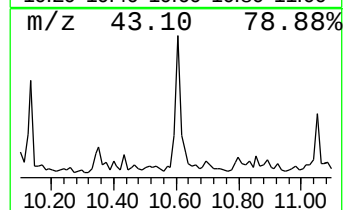
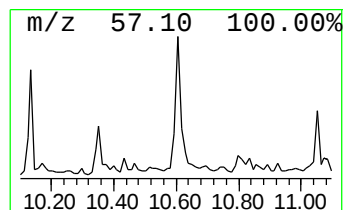
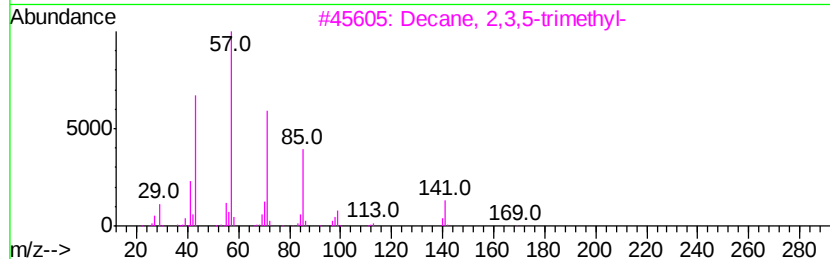
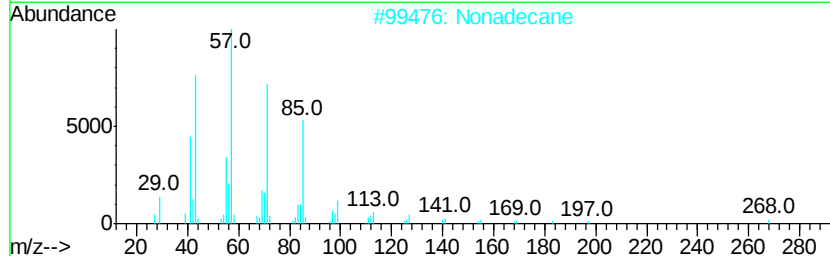
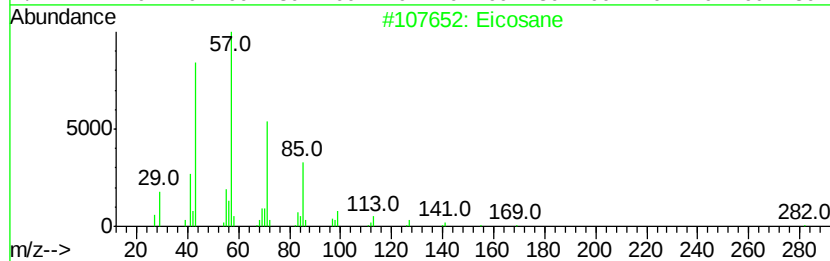
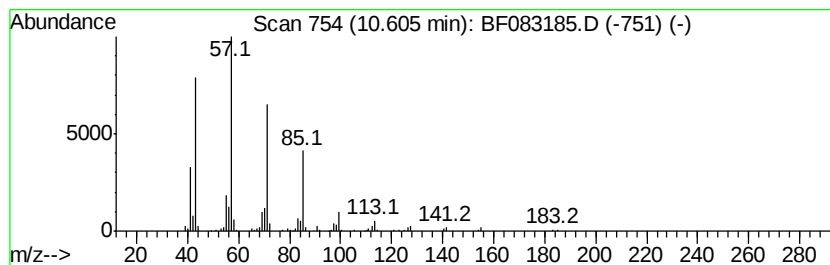
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.60	4.14 ng	278351	Phenanthrene-d10	11.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Nonadecane		268	C19H40	000629-92-5	87
3	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	83
4	Tridecane		184	C13H28	000629-50-5	80
5	Hexadecane		226	C16H34	000544-76-3	80



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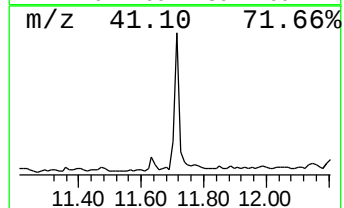
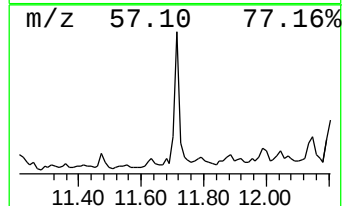
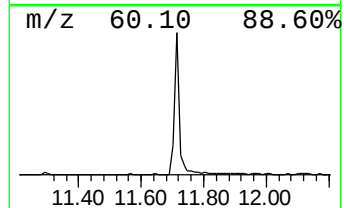
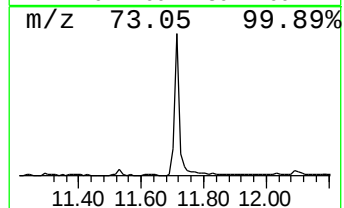
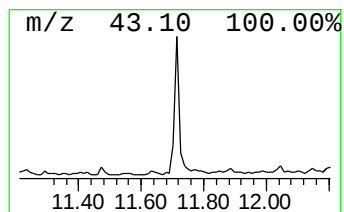
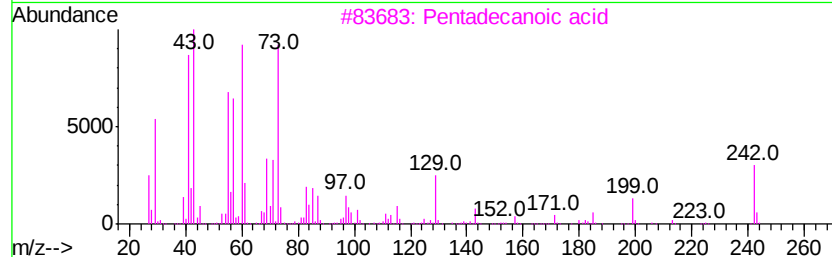
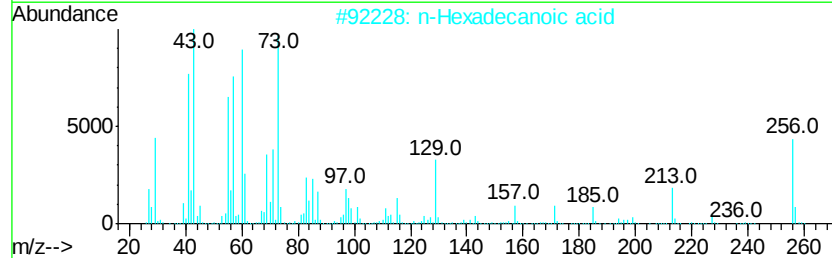
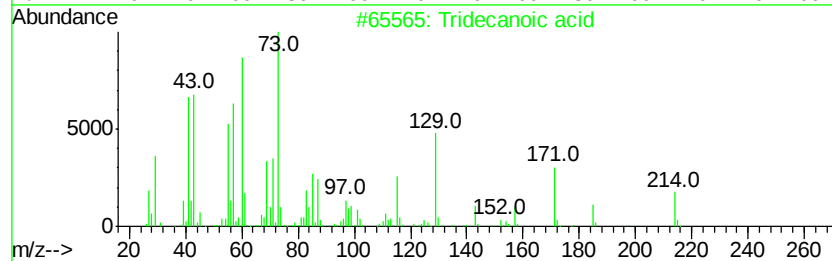
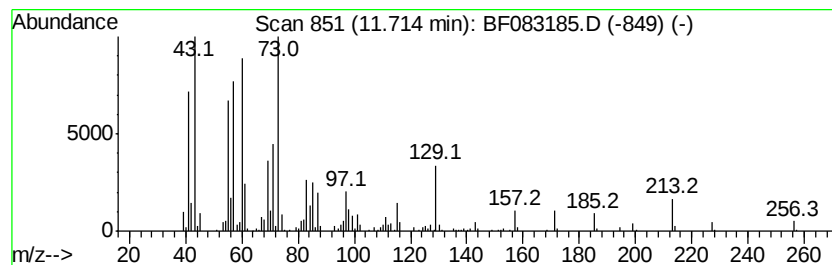
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	11.17 ng	751251	Phenanthrene-d10	11.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Undecane	156	C11H24	001120-21-4	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083185.D
Acq On : 23 Nov 2015 4:02
Operator : UM/IZ
Sample : G4475-03
Misc :
ALS Vial : 71 Sample Multiplier: 1

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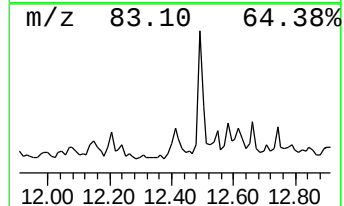
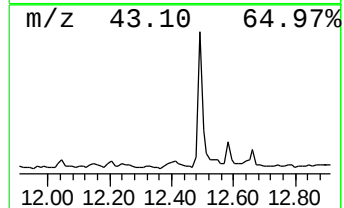
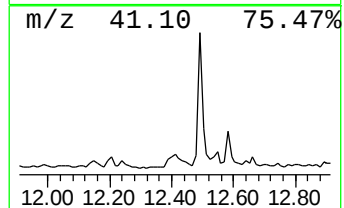
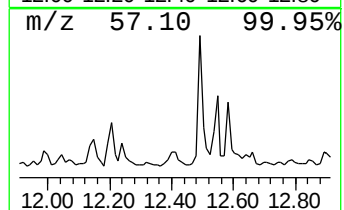
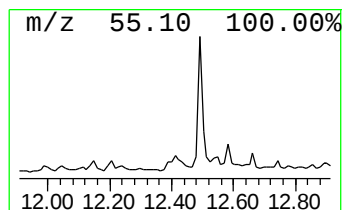
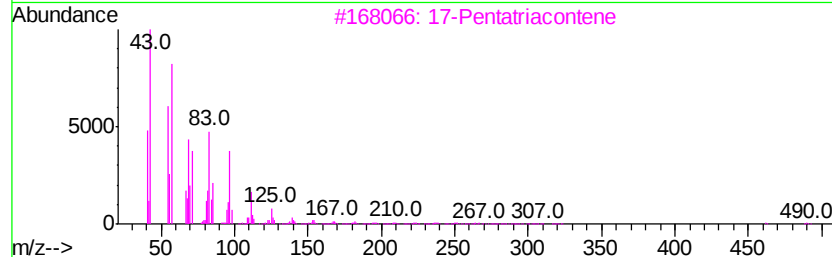
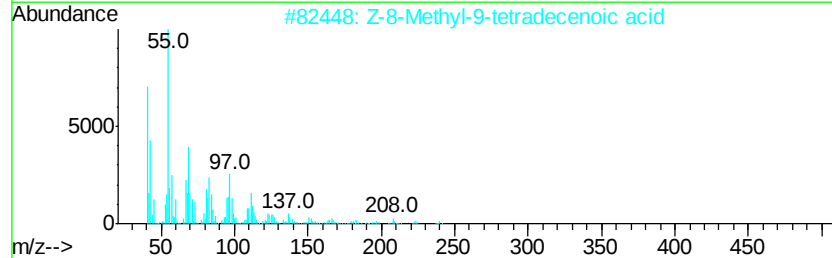
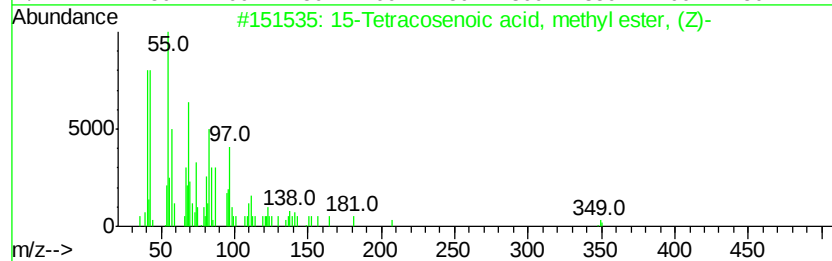
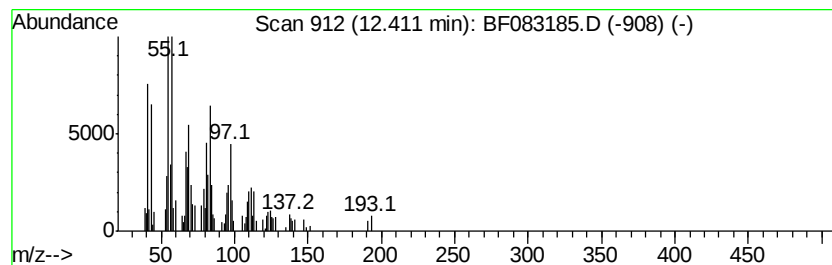
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 15-Tetracosenoic acid, methyl ester, (Z)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	2.76 ng	185477	Phenanthrene-d10	11.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	15-Tetracosenoic acid, methyl ester, (Z)-	380	C25H48O2	002733-88-2	64
2		Z-8-Methyl-9-tetradecenoic acid	240	C15H28O2	1000130-84-5	52
3		17-Pentatriacontene	491	C35H70	006971-40-0	46
4		9-Borabicyclo[3.3.1]nonane, 9-hydroxy-	138	C8H15BO	063366-65-4	46
5		1-Pentadecanol	228	C15H32O	000629-76-5	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
Data File : BF083185.D
Acq On : 23 Nov 2015 4:02
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Sample : G4475-03
Misc :
ALS Vial : 71 Sample Multiplier: 1

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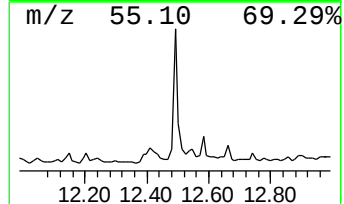
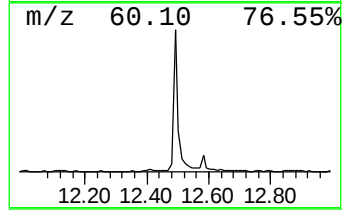
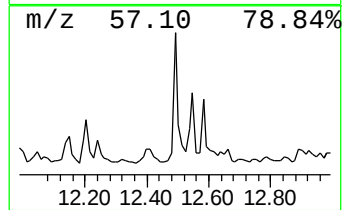
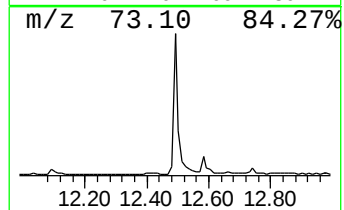
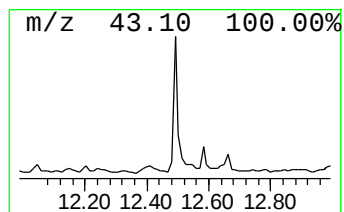
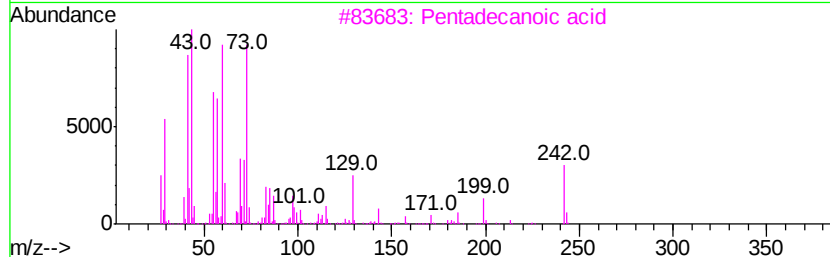
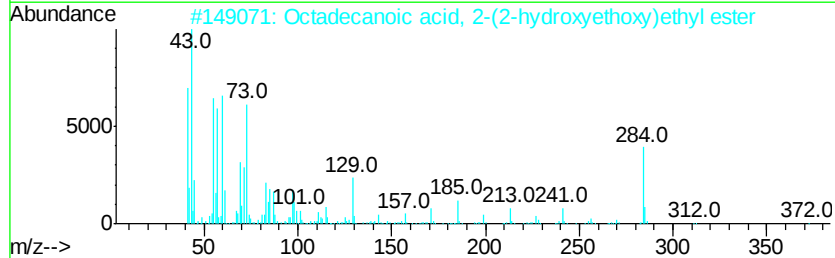
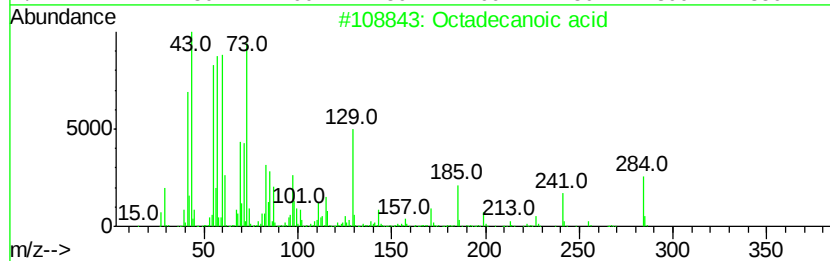
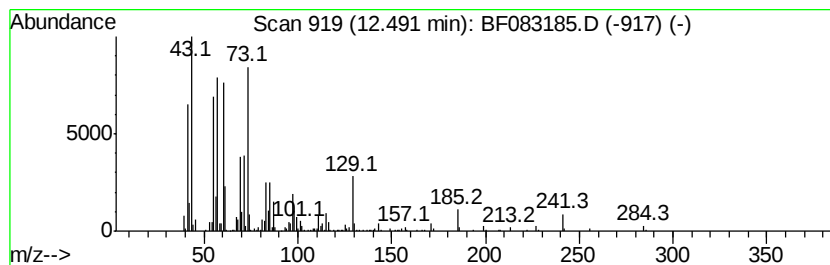
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	10.25 ng	688347	Chrysene-d12	13.78

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Undecane	156	C11H24	001120-21-4	18
5		Hexadecane	226	C16H34	000544-76-3	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083185.D
Acq On : 23 Nov 2015 4:02
Operator : UM/IZ
Sample : G4475-03
Misc :
ALS Vial : 71 Sample Multiplier: 1

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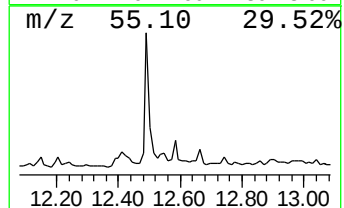
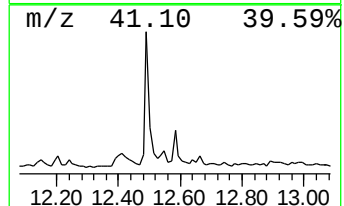
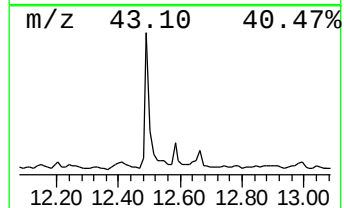
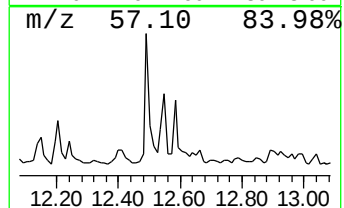
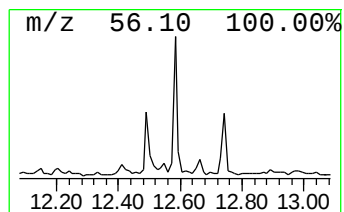
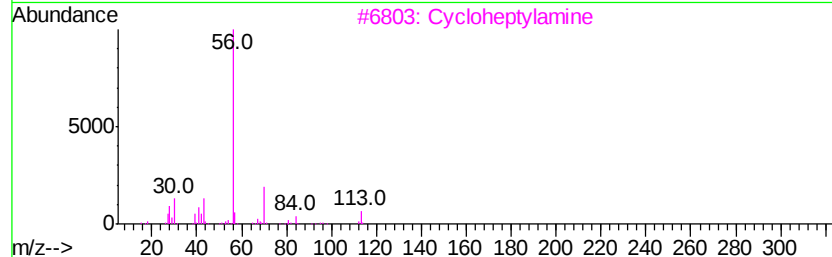
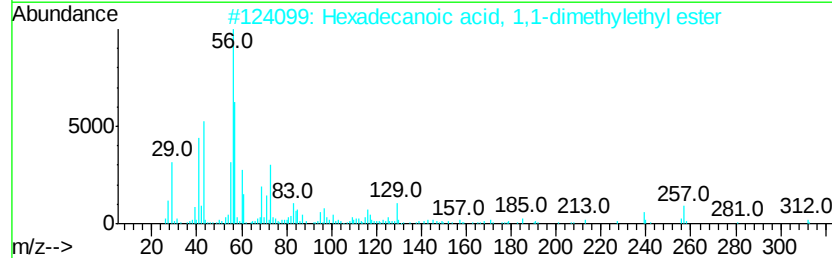
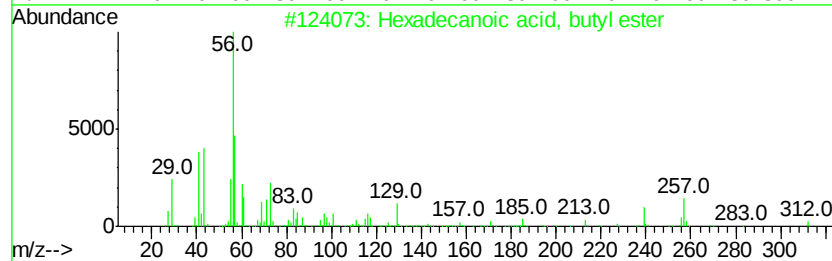
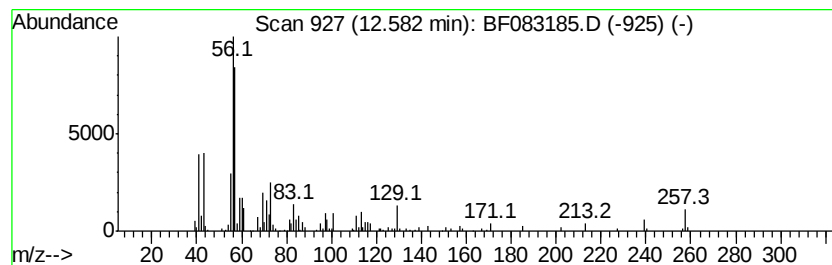
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Hexadecanoic acid, butyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	2.69 ng	180595	Chrysene-d12	13.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	68
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	49
3		Cycloheptylamine	113	C7H15N	005452-35-7	43
4		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	38
5		1-Hexene, 4-methyl-	98	C7H14	003769-23-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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Misc :
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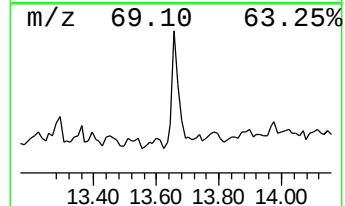
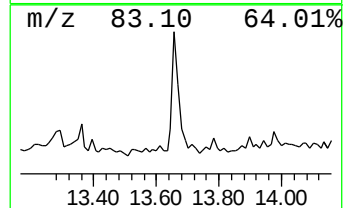
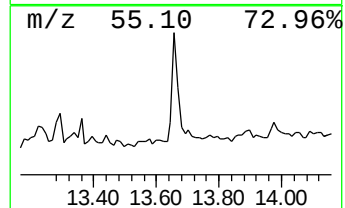
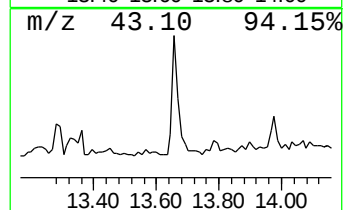
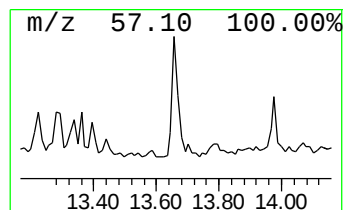
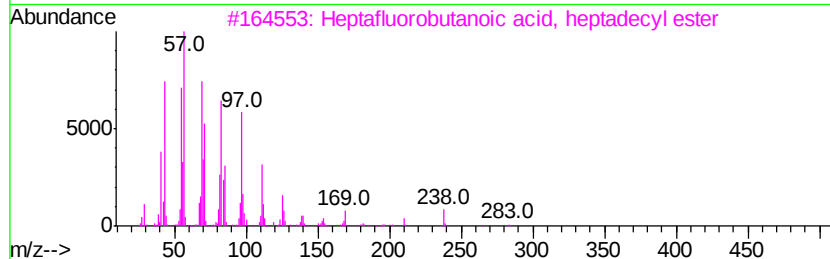
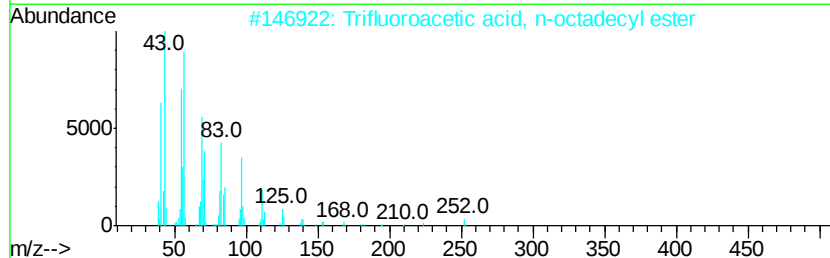
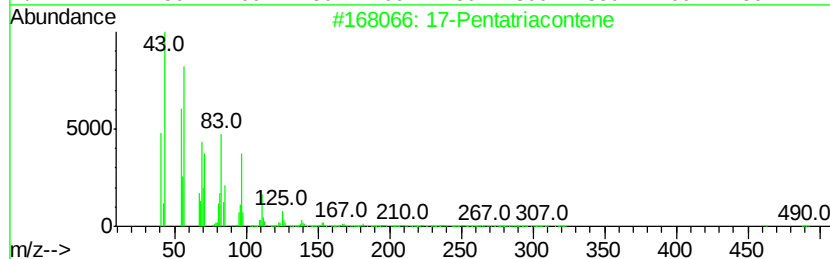
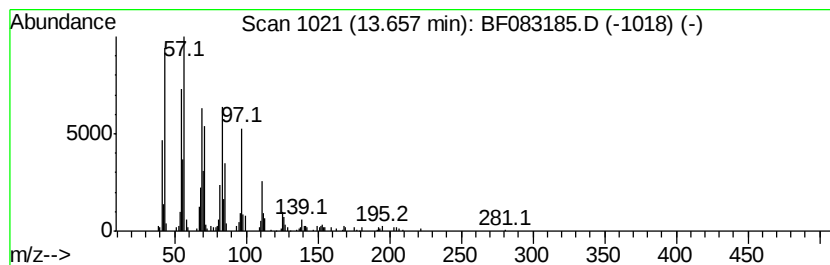
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 17-Pentatriacontene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	4.40 ng	295353	Chrysene-d12	13.78
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		17-Pentatriacontene	491 C35H70	006971-40-0 91
2		Trifluoroacetic acid, n-octadecy...	366 C20H37F3O2	079392-43-1 91
3		Heptafluorobutanoic acid, heptad...	452 C21H35F7O2	1000282-97-3 90
4		1-Hexadecanesulfonyl chloride	324 C16H33ClO2S	038775-38-1 90
5		Trichloroacetic acid, pentadecyl...	372 C17H31Cl3O2	074339-53-0 86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083185.D
Acq On : 23 Nov 2015 4:02
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Misc :
ALS Vial : 71 Sample Multiplier: 1

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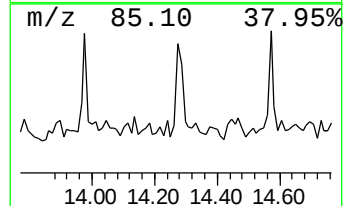
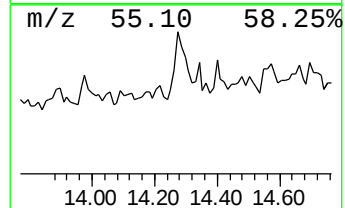
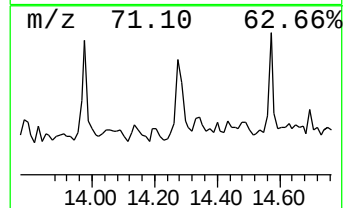
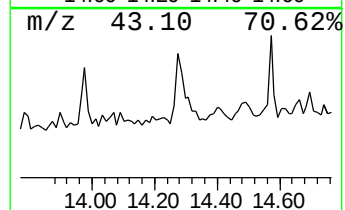
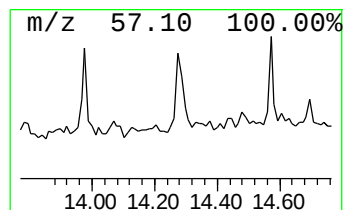
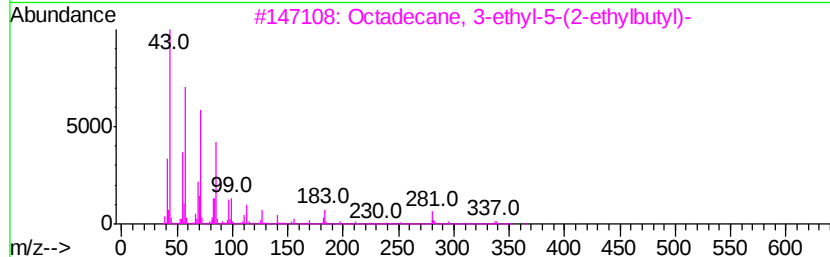
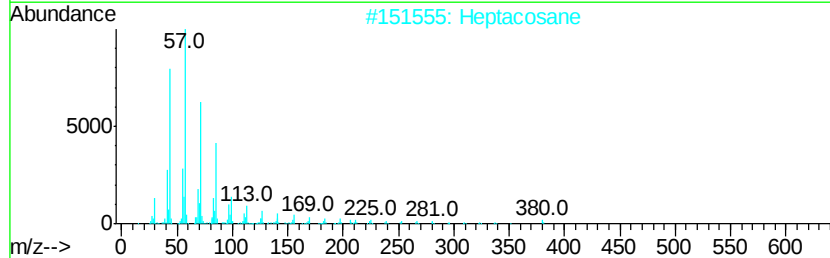
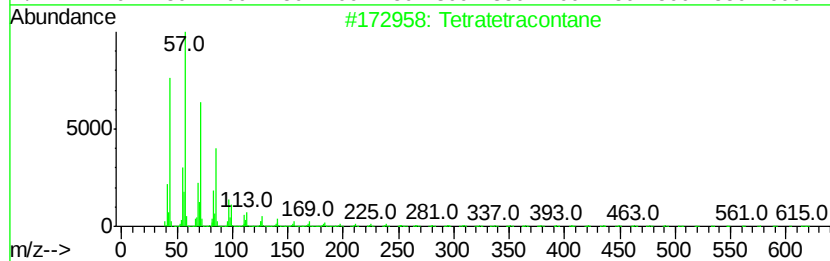
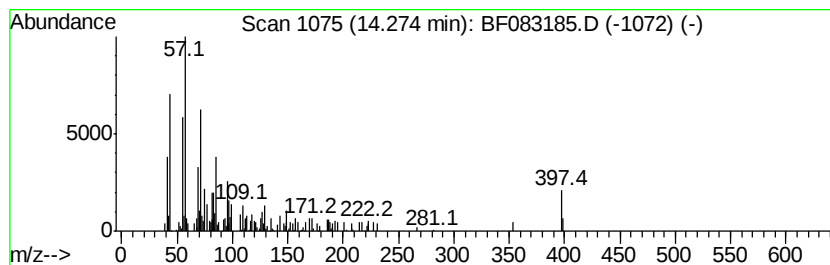
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown14.27 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	2.27 ng	152735	Chrysene-d12	13.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	41
2		Heptacosane	380	C27H56	000593-49-7	38
3		Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	38
4		Methoxyacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	38
5		Hexadecane	226	C16H34	000544-76-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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Acq On : 23 Nov 2015 4:02
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Sample : G4475-03
Misc :
ALS Vial : 71 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Pentanone, 4-hy...	4.80	58.7	ng	3624660	1	6.65	1234800	20.0
unknown6.42	6.42	94.8	ng	5852510	1	6.65	1234800	20.0
Eicosane	10.14	2.4	ng	193306	3	9.68	1585480	20.0
Nonadecane	10.60	4.1	ng	278351	4	11.16	1345230	20.0
Tridecanoic acid	11.71	11.2	ng	751251	4	11.16	1345230	20.0
15-Tetracosenoic ...	12.41	2.8	ng	185477	4	11.16	1345230	20.0
Octadecanoic acid	12.49	10.3	ng	688347	5	13.78	1342760	20.0
Hexadecanoic acid...	12.58	2.7	ng	180595	5	13.78	1342760	20.0
17-Pentatriacontene	13.66	4.4	ng	295353	5	13.78	1342760	20.0
unknown14.27	14.27	2.3	ng	152735	5	13.78	1342760	20.0