

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085611.D
 Acq On : 20 Mar 2016 11:16
 Operator : UM/SJ
 Sample : H1815-14
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-02-COMP

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-BF031816.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	5.053	240	242	251	rBV	274756	436148	10.39%	1.165%
2	5.464	275	278	280	rBV	3067227	3482803	82.94%	9.301%
3	6.493	365	368	371	rBV	2746510	3349535	79.76%	8.945%
4	6.630	377	380	382	rBV	2938382	3993569	95.10%	10.665%
5	6.859	397	400	402	rBV	599002	825589	19.66%	2.205%
6	7.007	411	413	416	rBV	2551834	2985575	71.10%	7.973%
7	7.419	446	449	452	rBV	2047956	2394509	57.02%	6.395%
8	8.139	509	512	514	rBV	1306302	1130777	26.93%	3.020%
9	9.213	603	606	609	rBV	3314056	4199288	100.00%	11.214%
10	9.613	638	641	642	rBV	494875	528310	12.58%	1.411%
11	9.831	658	660	661	rBV	40256	47652	1.13%	0.127%
12	9.888	663	665	668	rVB	942193	1124960	26.79%	3.004%
13	10.322	702	703	705	rVB	98857	83656	1.99%	0.223%
14	10.539	719	722	726	rBV	33725	62881	1.50%	0.168%
15	10.688	732	735	737	rBV	2028725	2505117	59.66%	6.690%
16	10.802	743	745	751	rVB	84973	149430	3.56%	0.399%
17	11.248	782	784	785	rBV	59381	59751	1.42%	0.160%
18	11.374	793	795	799	rBV	1189811	1121857	26.72%	2.996%
19	11.911	840	842	844	rBV	162187	233590	5.56%	0.624%
20	12.585	898	901	903	rBV	114054	110572	2.63%	0.295%
21	12.688	908	910	916	rBV	76370	114443	2.73%	0.306%
22	12.779	916	918	920	rBV	196885	205277	4.89%	0.548%
23	12.814	920	921	922	rVB	76833	53855	1.28%	0.144%
24	12.962	931	934	936	rBV	3406808	3273959	77.96%	8.743%
25	13.191	951	954	957	rBV	93154	127378	3.03%	0.340%
26	13.488	978	980	982	rBV	133707	137732	3.28%	0.368%
27	13.534	982	984	985	rBV	166235	151900	3.62%	0.406%
28	13.854	1010	1012	1020	rBV	850412	1146710	27.31%	3.062%
29	14.014	1022	1026	1027	rBV2	606683	957015	22.79%	2.556%
30	14.117	1032	1035	1036	rVV3	31736	64452	1.53%	0.172%
31	14.174	1038	1040	1045	rVB	317317	329340	7.84%	0.880%
32	14.482	1064	1067	1073	rBV	278755	372536	8.87%	0.995%
33	14.780	1091	1093	1095	rBV	187582	187334	4.46%	0.500%
34	15.020	1112	1114	1118	rBV2	40866	113298	2.70%	0.303%

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Integration Parameters: rteint.p
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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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35	15.100	1119	1121	1123	rVV	122630	135936	3.24%	0.363%
36	15.442	1148	1151	1156	rBV	471154	776077	18.48%	2.073%
37	15.877	1187	1189	1191	rBV	39615	57950	1.38%	0.155%
38	15.934	1192	1194	1204	rVB3	61164	230091	5.48%	0.614%
39	16.357	1228	1231	1238	rVB3	51129	184857	4.40%	0.494%

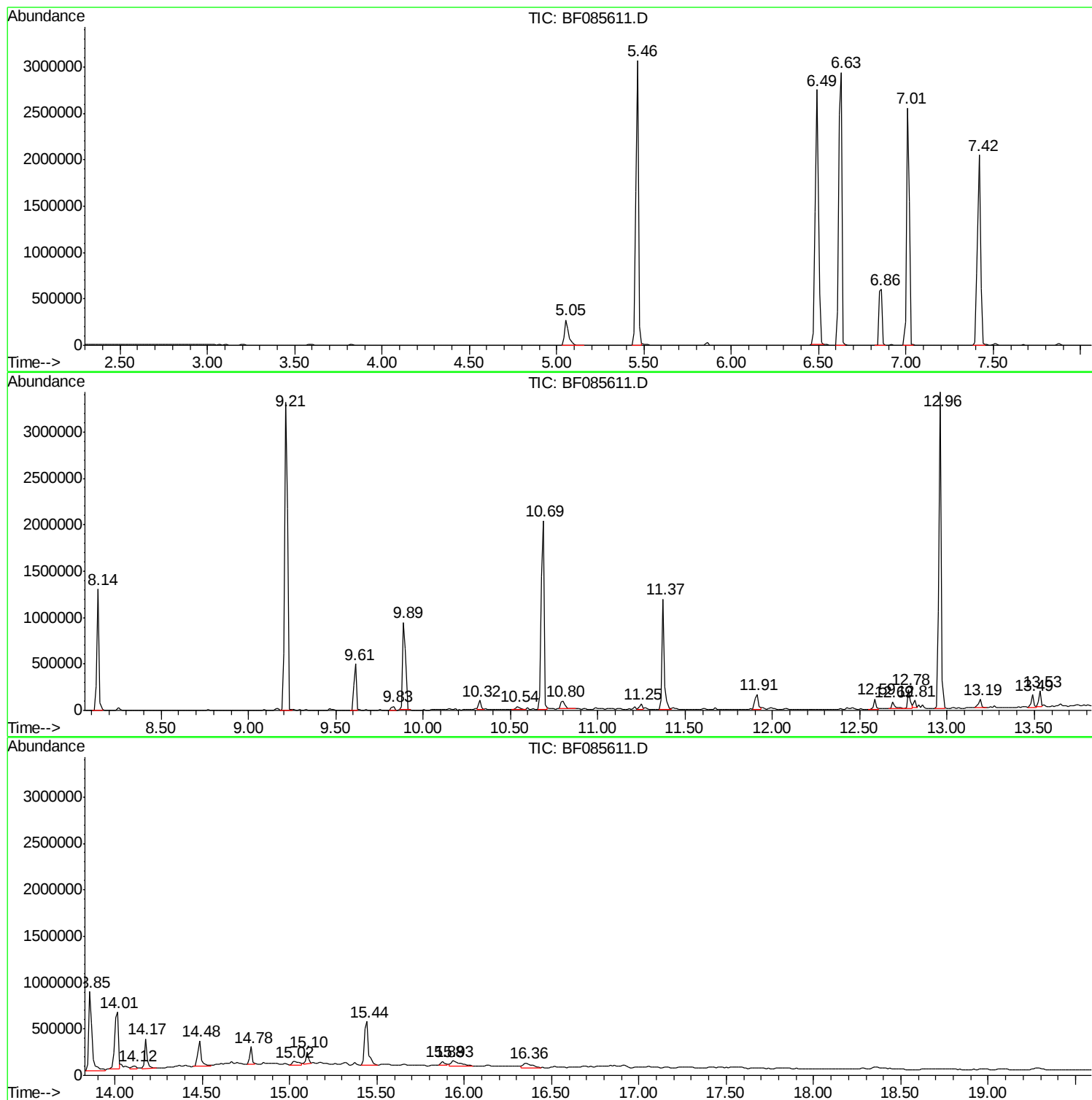
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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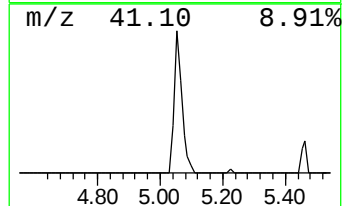
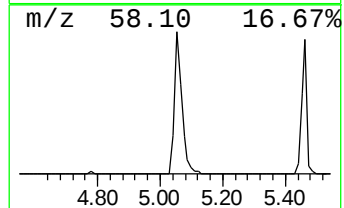
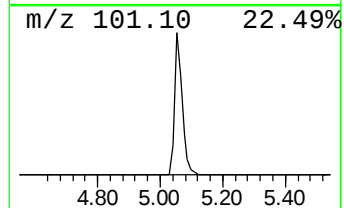
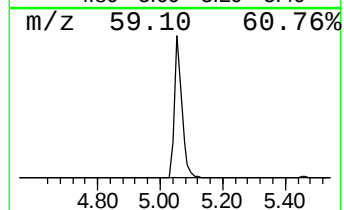
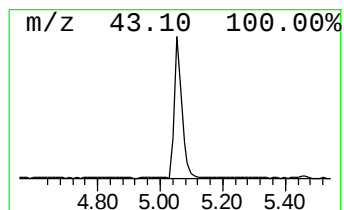
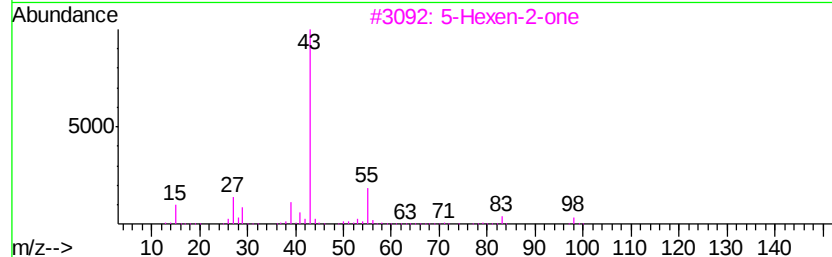
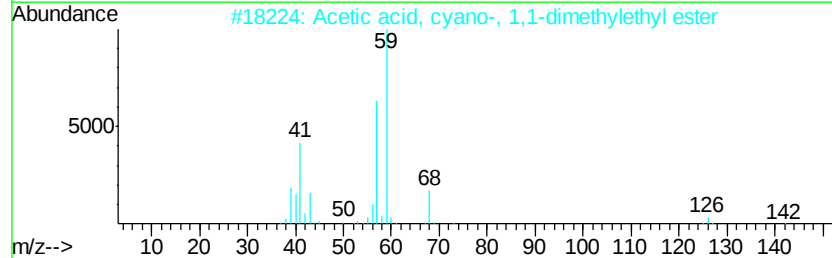
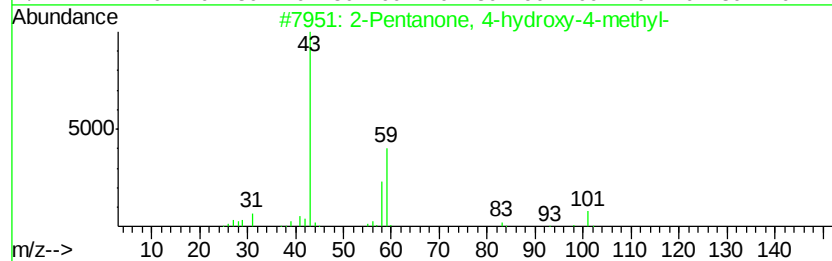
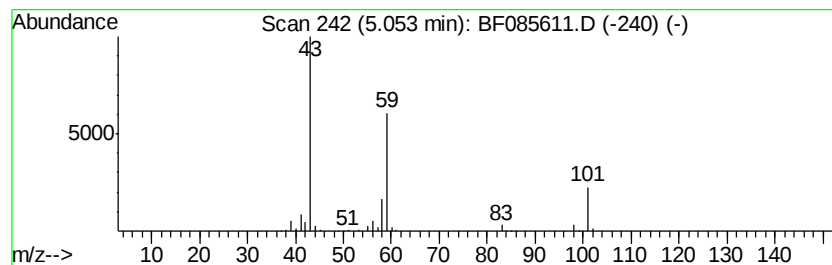
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	10.57 ng	436148	1,4-Dichlorobenzene-d4	6.86

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, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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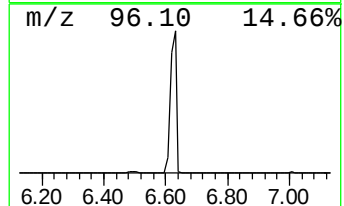
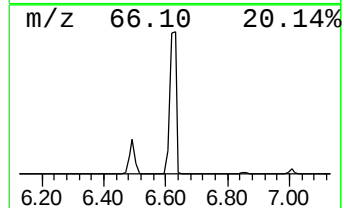
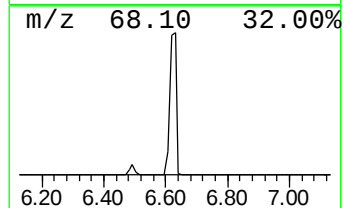
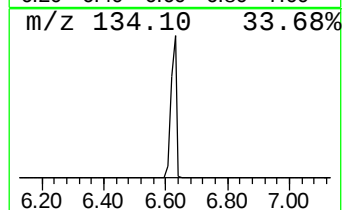
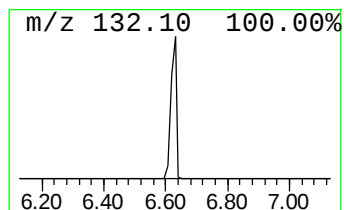
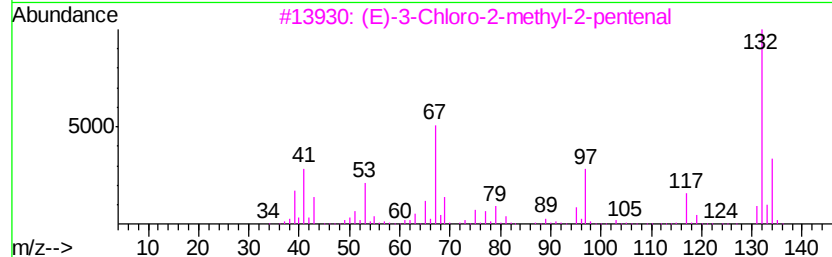
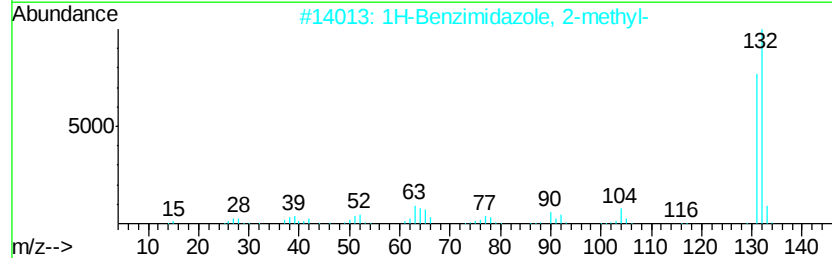
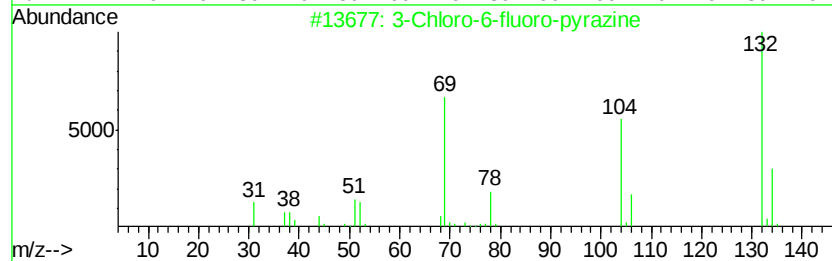
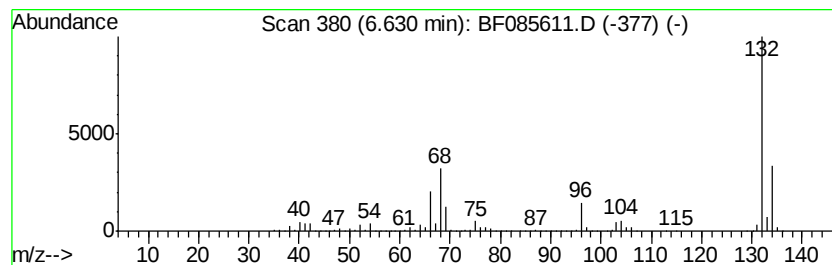
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	96.74 ng	3993570	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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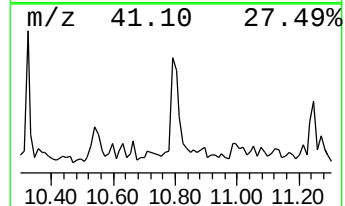
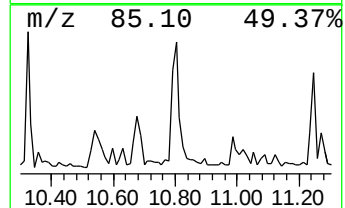
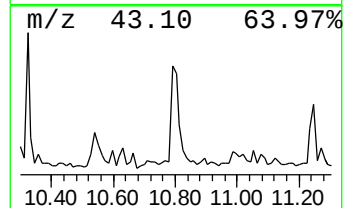
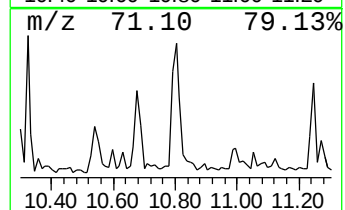
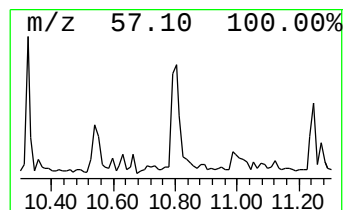
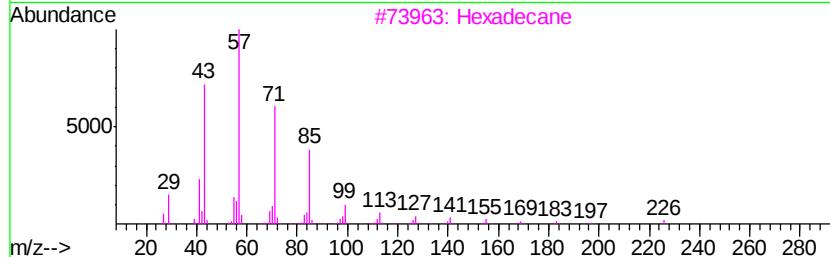
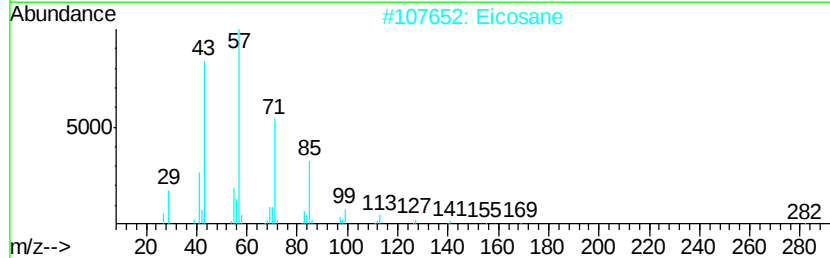
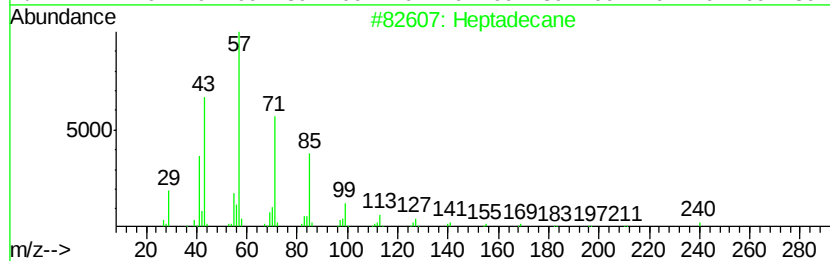
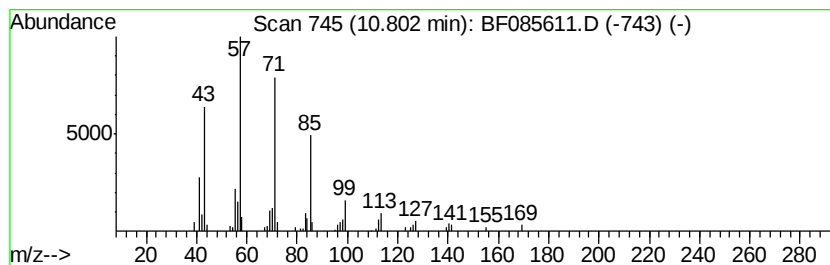
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.80	2.66 ng	149430	Phenanthrene-d10		11.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Eicosane	282	C20H42	000112-95-8	91
3	Hexadecane	226	C16H34	000544-76-3	90
4	Pentadecane	212	C15H32	000629-62-9	90
5	Tritetracontane	605	C43H88	007098-21-7	90



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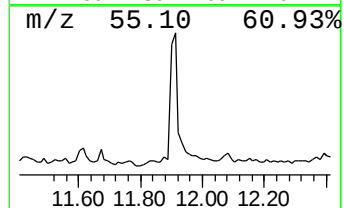
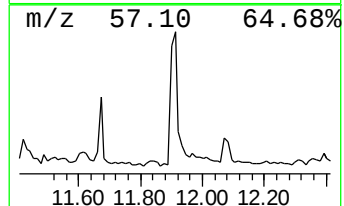
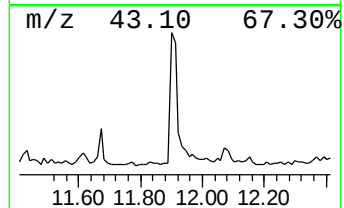
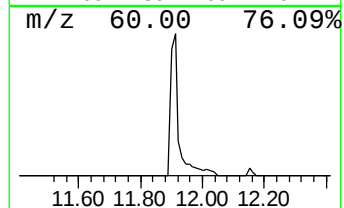
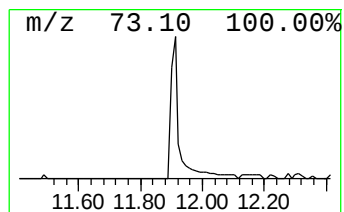
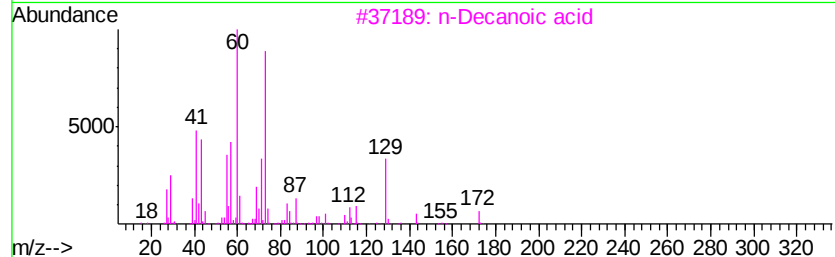
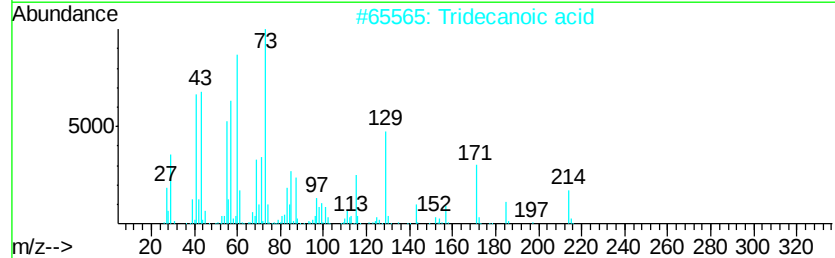
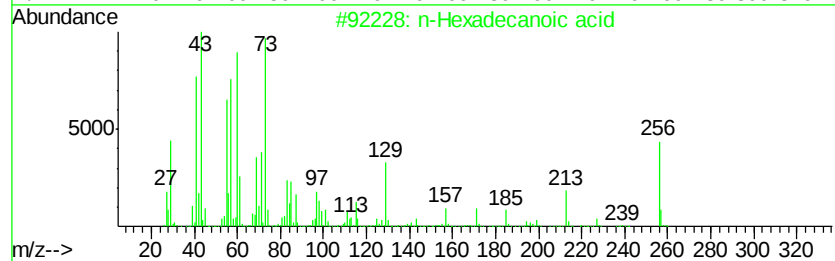
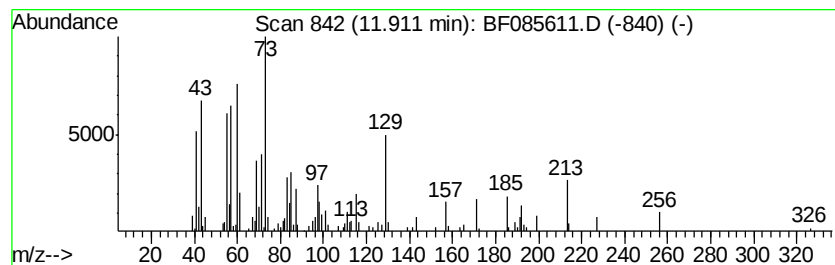
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.91	4.16 ng	233590	Phenanthrene-d10		11.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tridecanoic acid	214	C13H26O2	000638-53-9	89
3	n-Decanoic acid	172	C10H20O2	000334-48-5	76
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	52
5	Nonanoic acid	158	C9H18O2	000112-05-0	18



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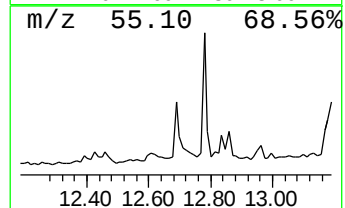
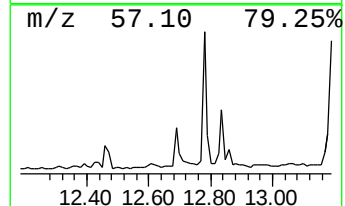
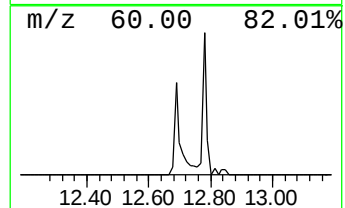
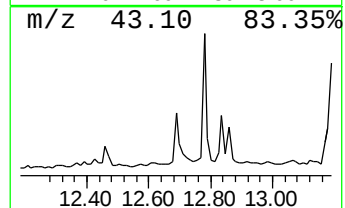
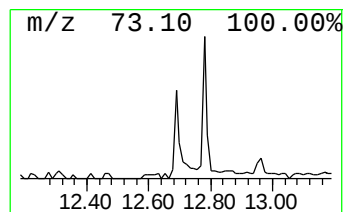
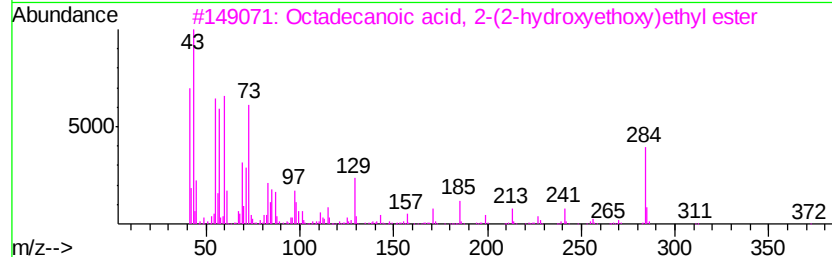
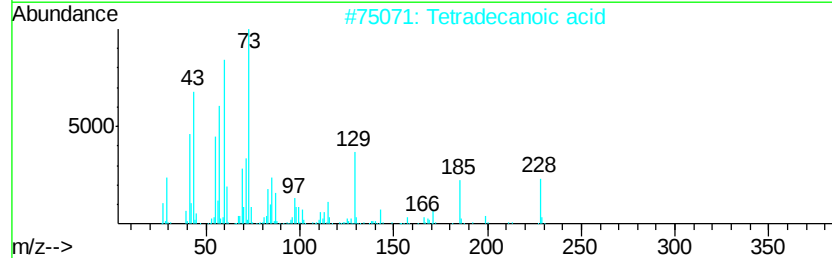
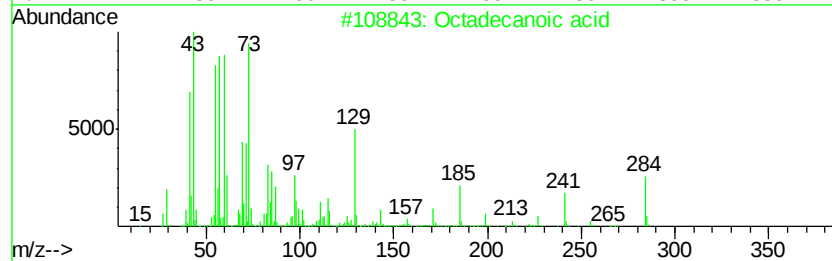
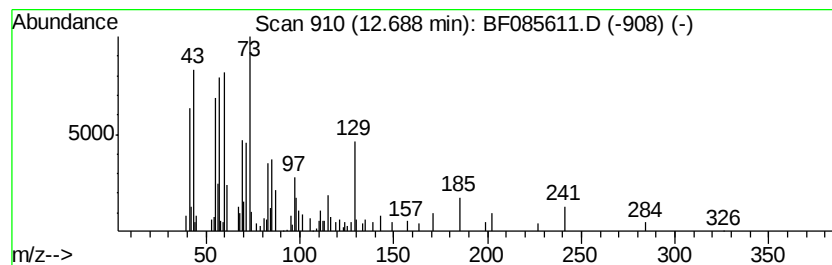
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	2.04 ng	114443	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
3		Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
4		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	18
5		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085611.D
Acq On : 20 Mar 2016 11:16
Operator : UM/SJ
Sample : H1815-14
Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-02-COMP

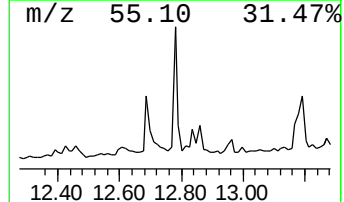
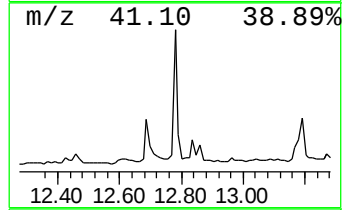
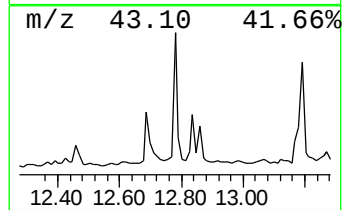
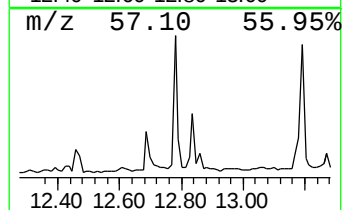
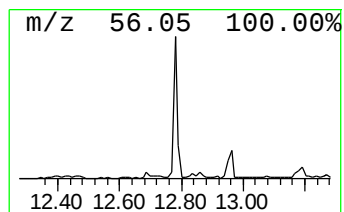
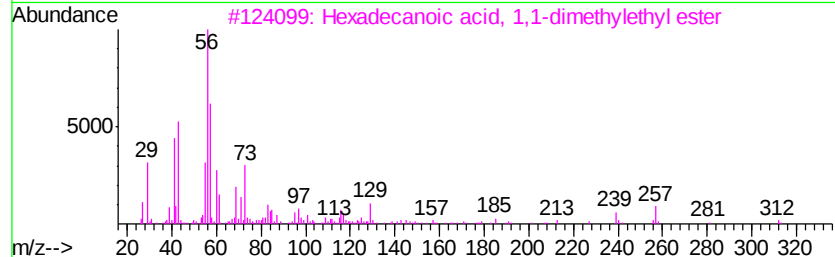
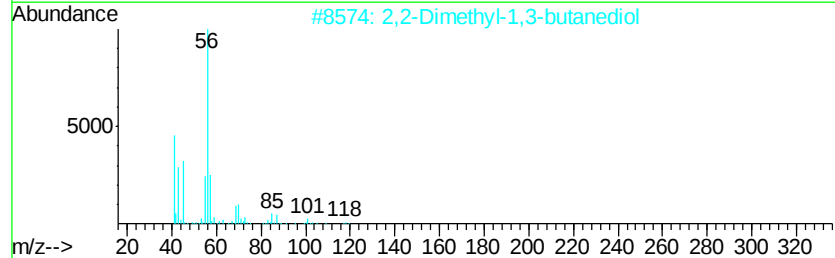
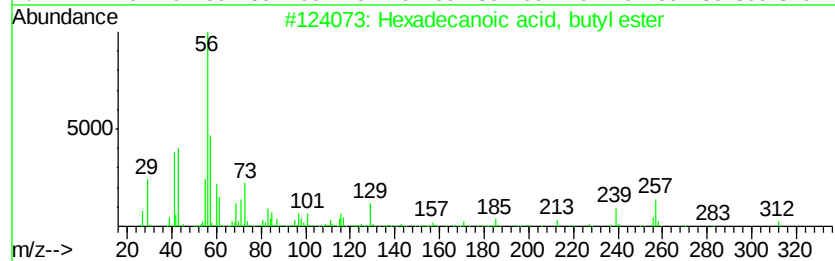
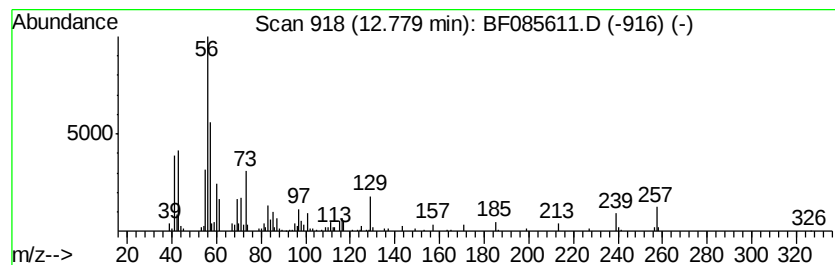
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.78	4.29 ng	205277	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		2,2-Dimethyl-1,3-butanediol	118	C6H14O2	000076-35-7	43
3		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	43
4		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	38
5		Cyclohexanol,2-amino-,trans-	115	C6H13NO	006982-39-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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Acq On : 20 Mar 2016 11:16
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Sample : H1815-14
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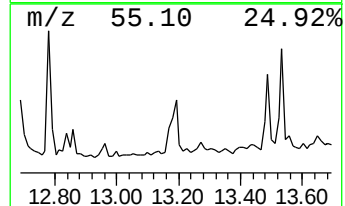
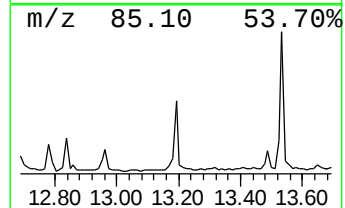
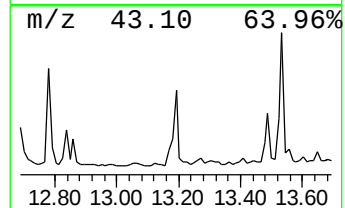
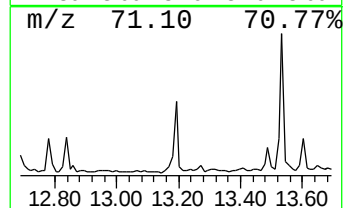
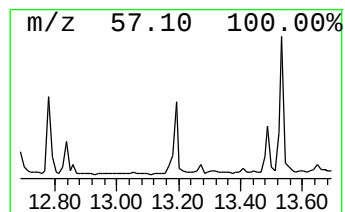
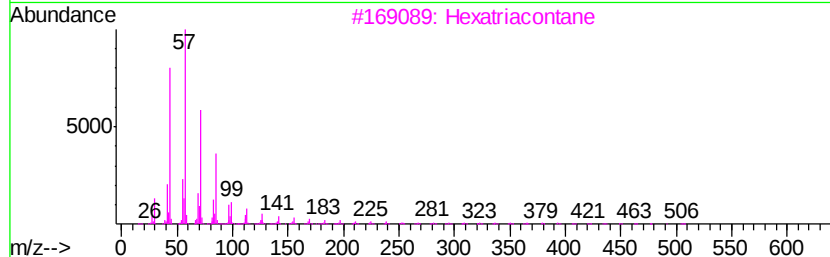
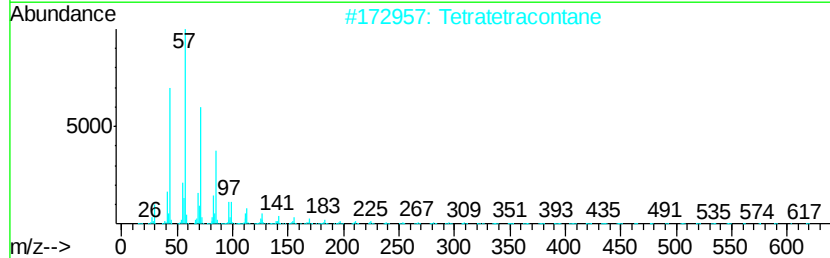
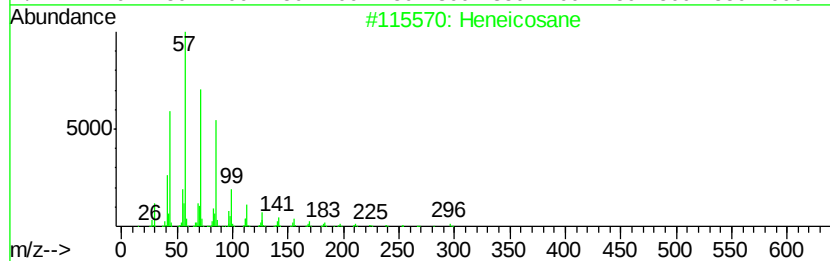
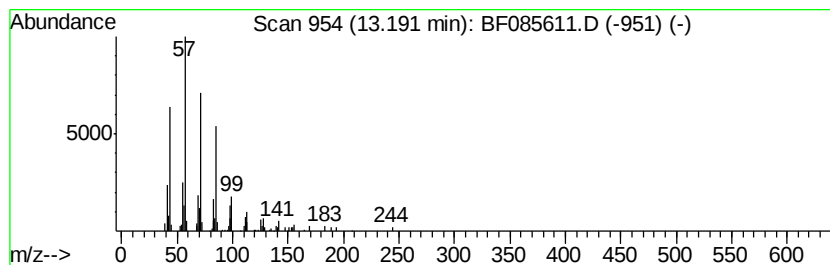
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heneicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	2.66 ng	127378	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296	C21H44	000629-94-7 91
2	Tetratetracontane	619	C44H90	007098-22-8 91
3	Hexatriacontane	507	C36H74	000630-06-8 91
4	Octacosane	394	C28H58	000630-02-4 91
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1 91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085611.D
Acq On : 20 Mar 2016 11:16
Operator : UM/SJ
Sample : H1815-14
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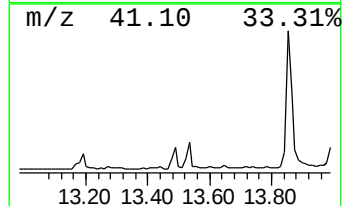
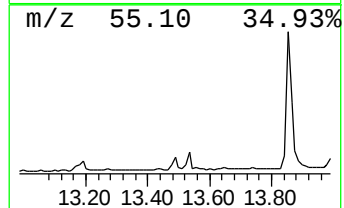
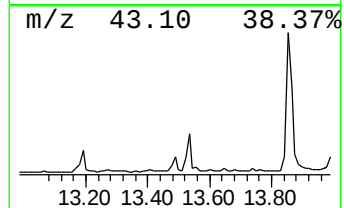
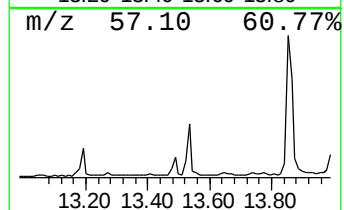
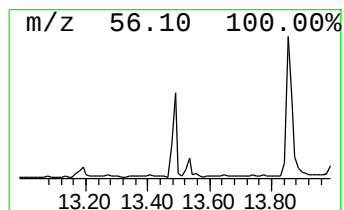
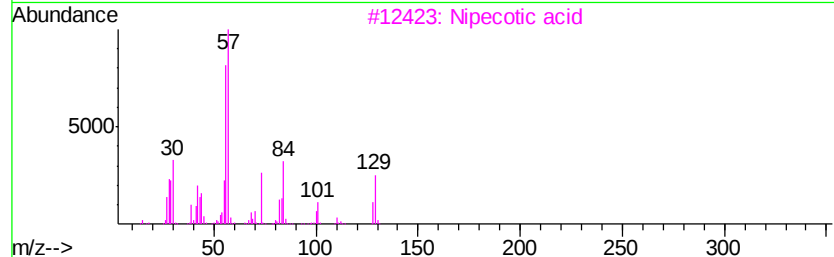
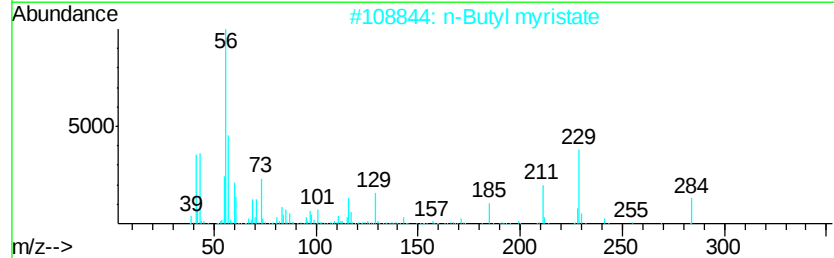
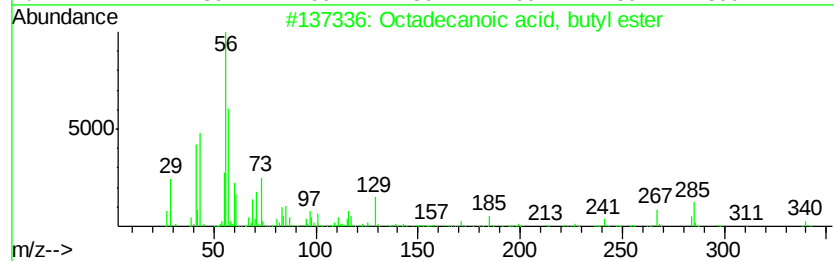
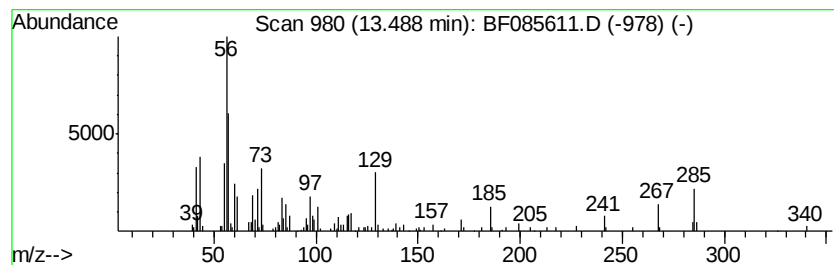
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid, butyl ester Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.49	2.88 ng	137732	Chrysene-d12		14.01
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	98
2	n-Butyl myristate	284	C18H36O2	000110-36-1	64
3	Nipecotic acid	129	C6H11NO2	000498-95-3	47
4	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	30
5	Cyclohexanol, 4-amino-, trans-	115	C6H13NO	027489-62-9	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085611.D
Acq On : 20 Mar 2016 11:16
Operator : UM/SJ
Sample : H1815-14
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ALS Vial : 69 Sample Multiplier: 1

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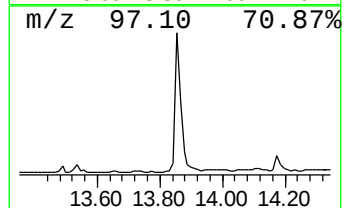
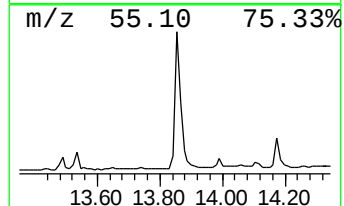
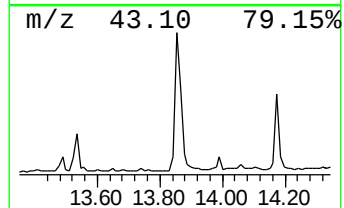
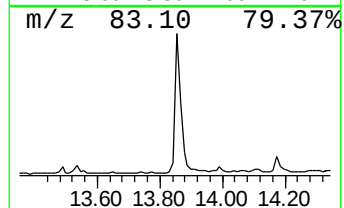
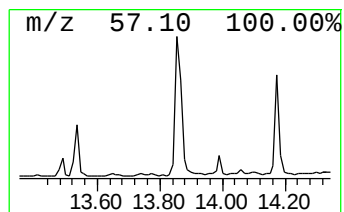
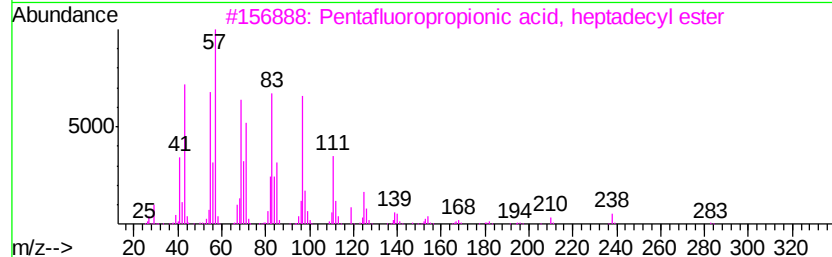
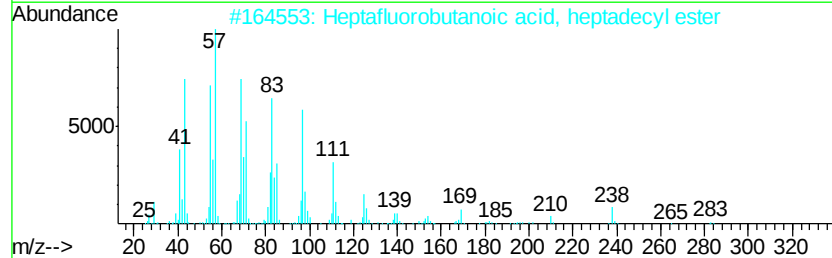
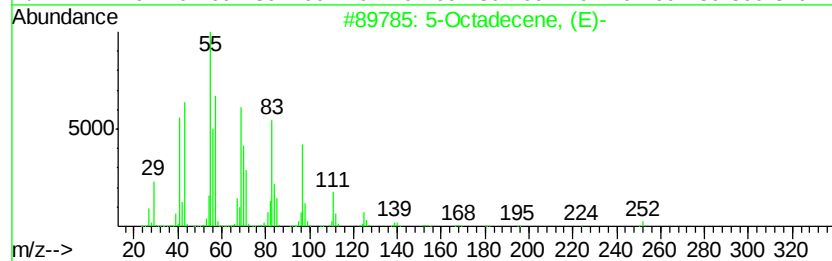
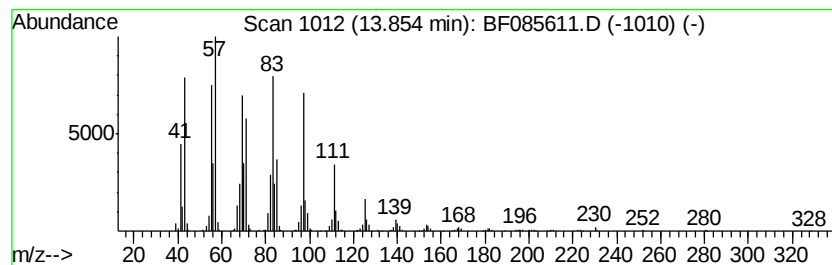
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 5-Octadecene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	23.96 ng	1146710	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Octadecene, (E)-	252	C18H36	007206-21-5	94
2		Heptafluorobutanoic acid, heptadecyl ester	452	C21H35F7O2	1000282-97-3	94
3		Pentafluoropropionic acid, heptadecyl ester	402	C20H35F5O2	1000283-04-2	91
4		Trifluoroacetic acid, n-octadecyl ester	366	C20H37F3O2	079392-43-1	91
5		Cyclohexadecane	224	C16H32	000295-65-8	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085611.D
Acq On : 20 Mar 2016 11:16
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Misc :
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ClientSampleId :
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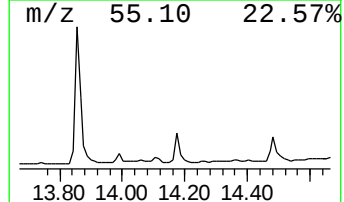
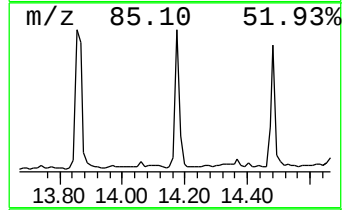
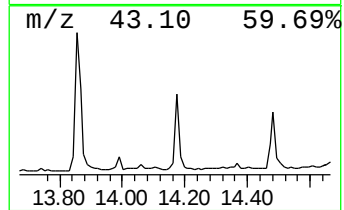
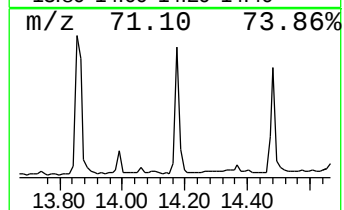
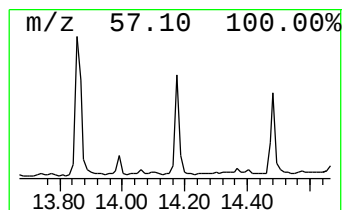
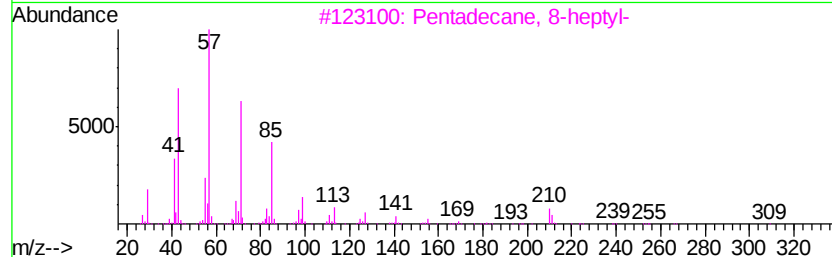
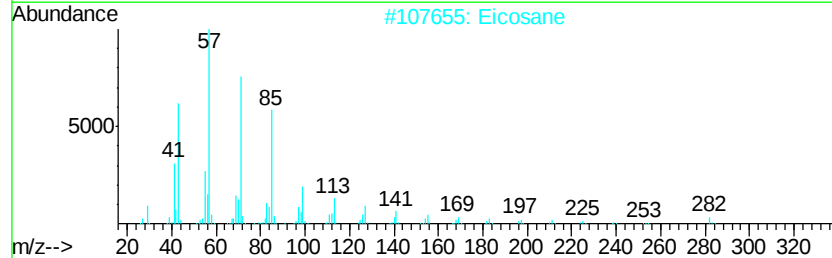
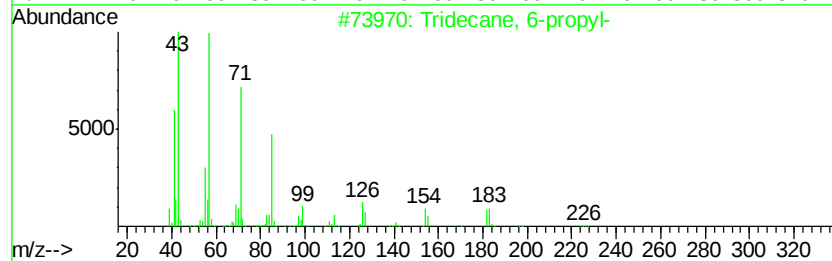
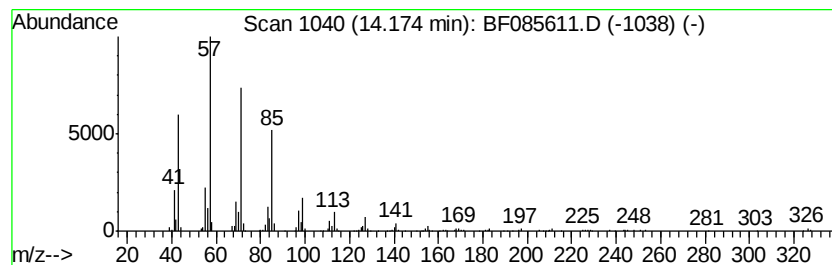
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecane, 6-propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	6.88 ng	329340	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 6-propyl-	226	C16H34	055045-10-8	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4		Tricosane	324	C23H48	000638-67-5	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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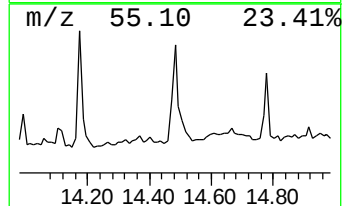
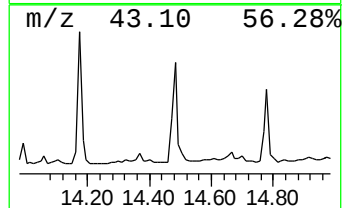
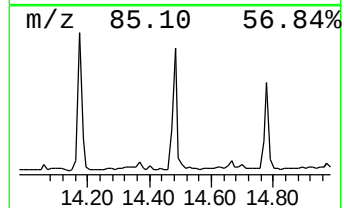
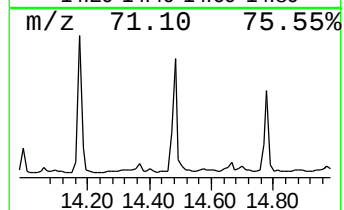
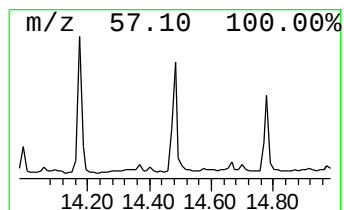
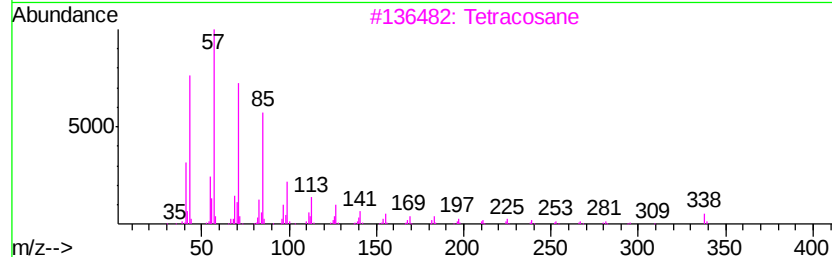
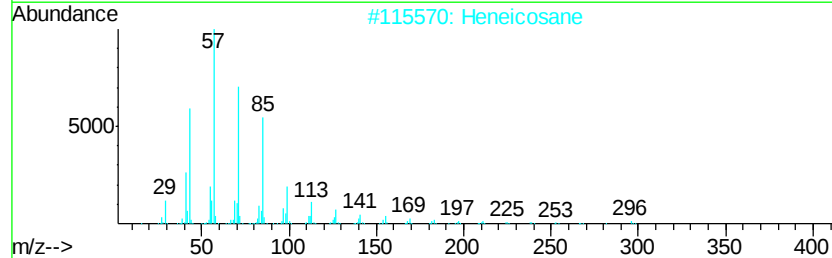
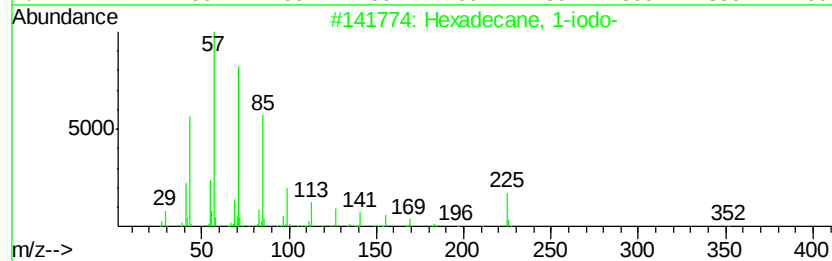
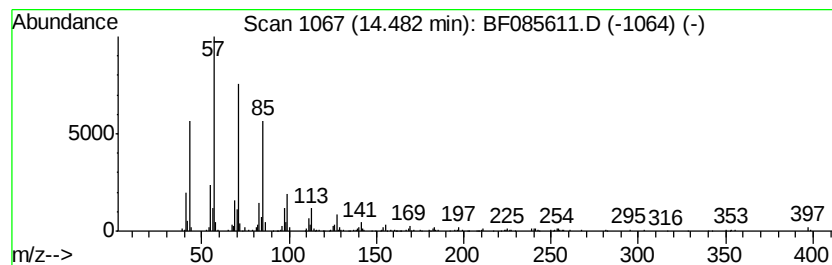
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecane, 1-iodo- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.48	7.79 ng	372536	Chrysene-d12	14.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5	Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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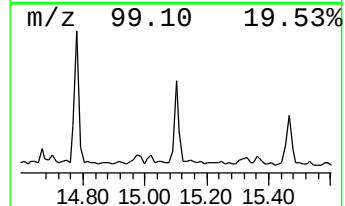
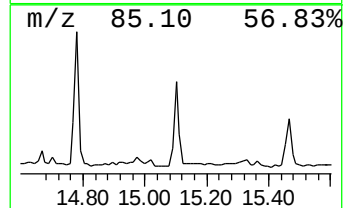
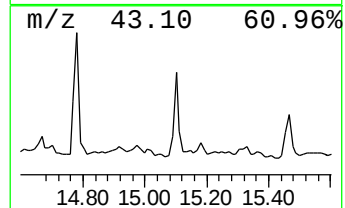
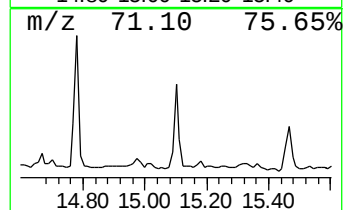
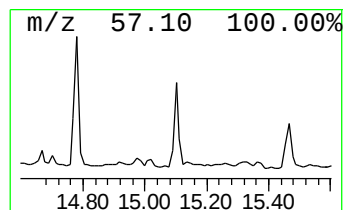
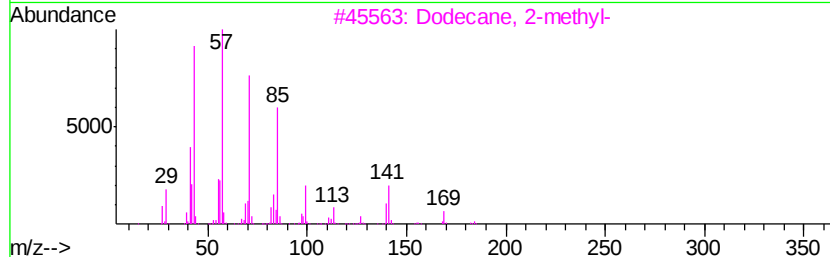
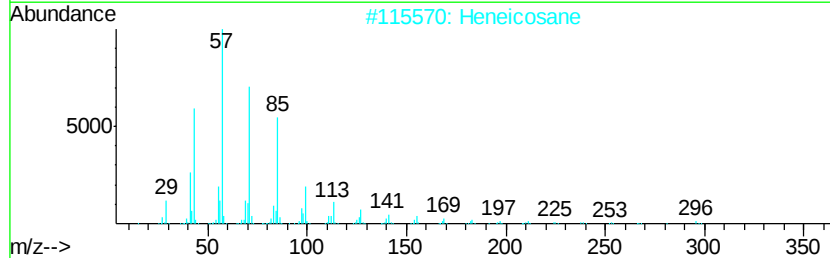
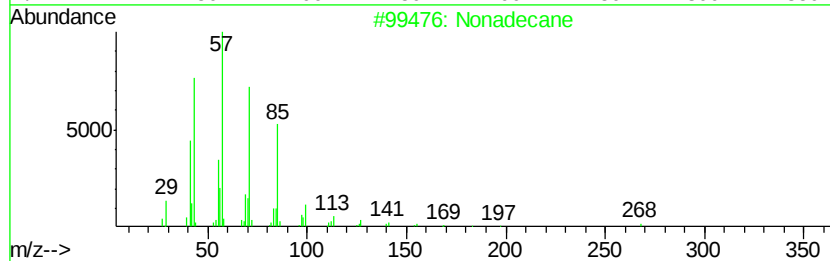
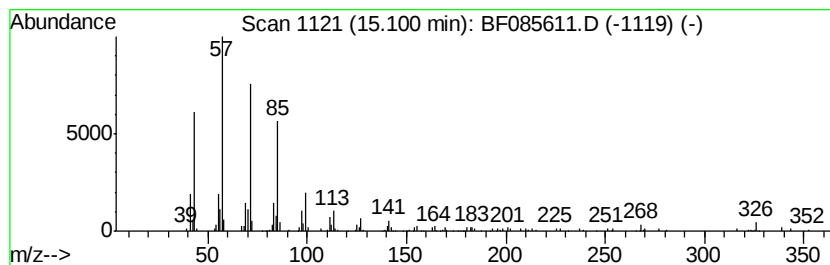
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	3.50 ng	135936	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	93
2		Heneicosane	296	C21H44	000629-94-7	90
3		Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085611.D
Acq On : 20 Mar 2016 11:16
Operator : UM/SJ
Sample : H1815-14
Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-02-COMP

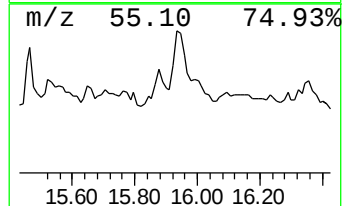
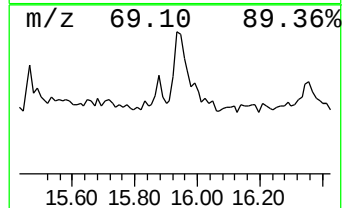
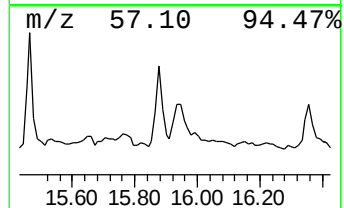
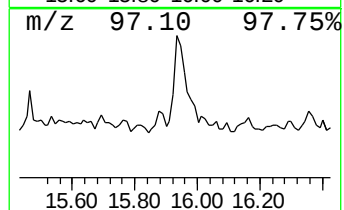
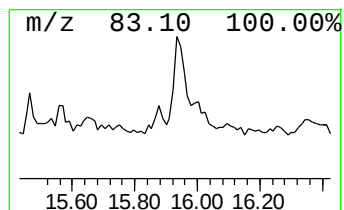
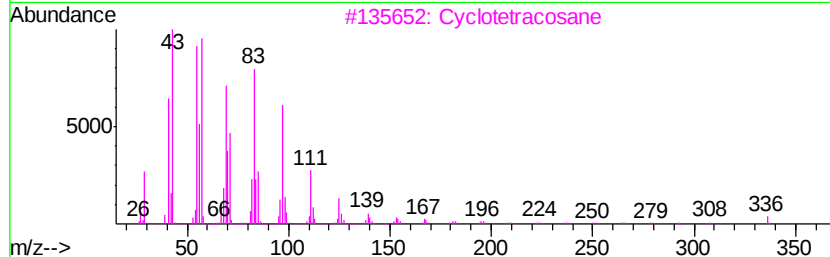
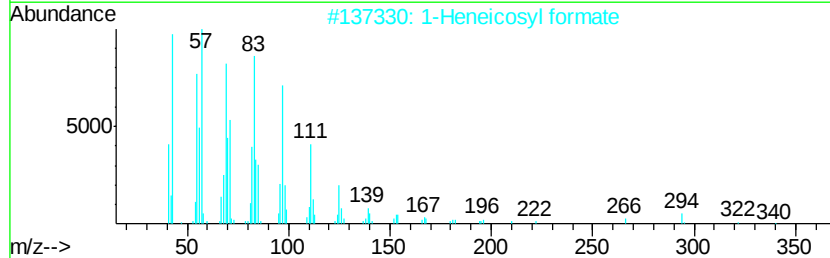
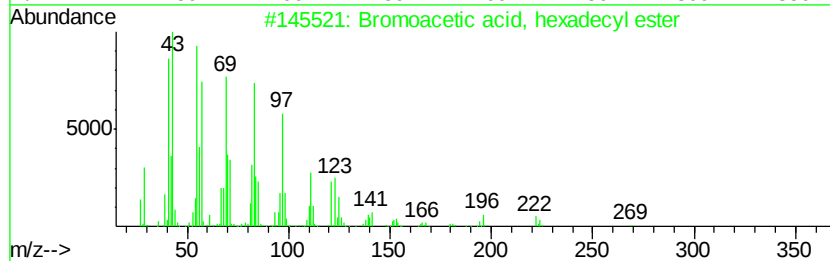
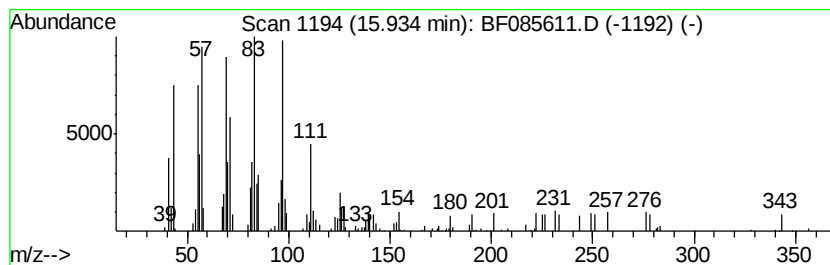
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Bromoacetic acid, hexadecyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	5.93 ng	230091	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	64
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	64
3		Cyclotetracosane	336	C24H48	000297-03-0	64
4		1-Docosene	308	C22H44	001599-67-3	62
5		1-Eicosanol	298	C20H42O	000629-96-9	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085611.D
Acq On : 20 Mar 2016 11:16
Operator : UM/SJ
Sample : H1815-14
Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-02-COMP

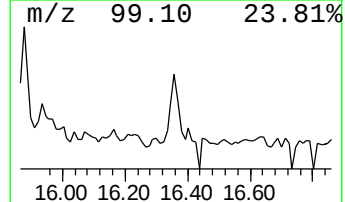
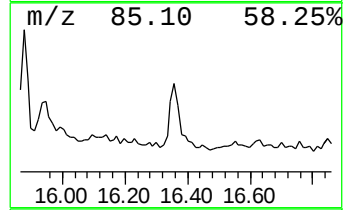
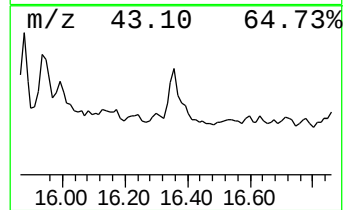
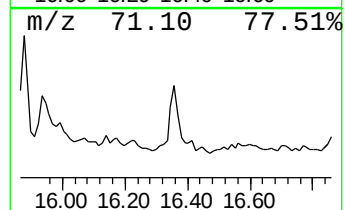
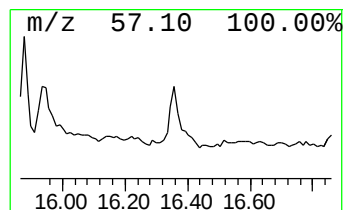
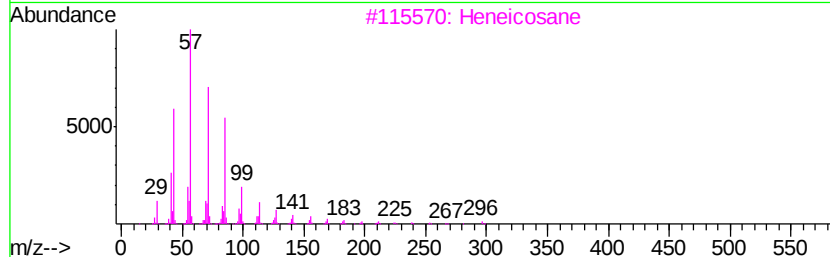
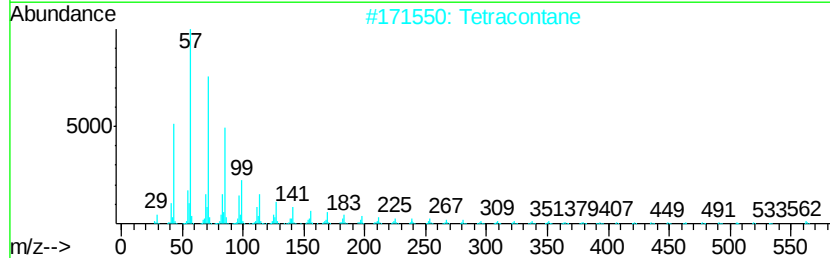
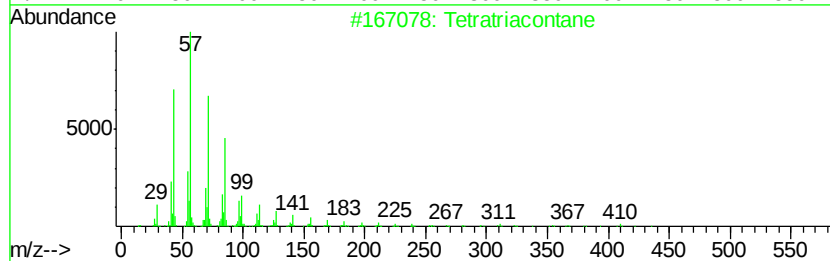
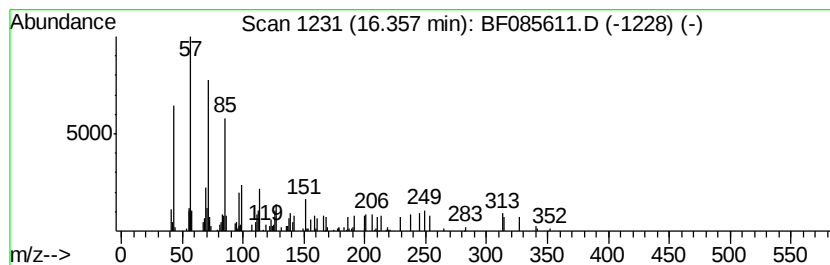
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Tetratriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	4.76 ng	184857	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontane	479	C34H70	014167-59-0	64
2		Tetracontane	563	C40H82	004181-95-7	64
3		Heneicosane	296	C21H44	000629-94-7	58
4		Hentriacontane	437	C31H64	000630-04-6	53
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085611.D
 Acq On : 20 Mar 2016 11:16
 Operator : UM/SJ
 Sample : H1815-14
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.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...	5.05	10.6	ng	436148	1	6.86	825589	20.0
unknown6.63	6.63	96.7	ng	3993570	1	6.86	825589	20.0
Heptadecane	10.80	2.7	ng	149430	4	11.37	1121860	20.0
n-Hexadecanoic acid	11.91	4.2	ng	233590	4	11.37	1121860	20.0
Octadecanoic acid	12.69	2.0	ng	114443	4	11.37	1121860	20.0
Hexadecanoic acid...	12.78	4.3	ng	205277	5	14.01	957015	20.0
Heneicosane	13.19	2.7	ng	127378	5	14.01	957015	20.0
Octadecanoic acid...	13.49	2.9	ng	137732	5	14.01	957015	20.0
5-Octadecene, (E)-	13.85	24.0	ng	1146710	5	14.01	957015	20.0
Tridecane, 6-propyl-	14.17	6.9	ng	329340	5	14.01	957015	20.0
Hexadecane, 1-iodo-	14.48	7.8	ng	372536	5	14.01	957015	20.0
Nonadecane	15.10	3.5	ng	135936	6	15.44	776077	20.0
Bromoacetic acid,...	15.93	5.9	ng	230091	6	15.44	776077	20.0
Tetratriacontane	16.36	4.8	ng	184857	6	15.44	776077	20.0