

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085606.D
 Acq On : 20 Mar 2016 8:58
 Operator : UM/SJ
 Sample : H1815-08
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-04-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	239	242	249	rVB	243587	427076	10.46%	1.131%
2	5.453	274	277	280	rBV	2604314	3843412	94.17%	10.177%
3	6.493	364	368	370	rBV	3044298	3923366	96.13%	10.389%
4	6.619	376	379	382	rBV	2954191	3833591	93.93%	10.151%
5	6.847	397	399	402	rBV	806866	849446	20.81%	2.249%
6	7.007	410	413	416	rBV	2863886	2874740	70.43%	7.612%
7	7.419	446	449	451	rBV	2279951	2495531	61.14%	6.608%
8	8.139	509	512	514	rBV	1168348	1168419	28.63%	3.094%
9	9.213	603	606	609	rBV	3720387	4081475	100.00%	10.808%
10	9.602	638	640	642	rBV	422631	506393	12.41%	1.341%
11	9.819	657	659	663	rBV	57201	70162	1.72%	0.186%
12	9.887	663	665	668	rVB	1306808	1228200	30.09%	3.252%
13	10.322	699	703	705	rVB2	151412	146594	3.59%	0.388%
14	10.539	719	722	725	rBV	45476	75137	1.84%	0.199%
15	10.688	731	735	737	rBV2	1925972	2917665	71.49%	7.726%
16	10.790	743	744	751	rVB	123429	188341	4.61%	0.499%
17	11.248	782	784	785	rBV	65752	87285	2.14%	0.231%
18	11.373	793	795	797	rBV	1474418	1269646	31.11%	3.362%
19	11.442	797	801	804	rVB2	24971	49326	1.21%	0.131%
20	11.899	839	841	849	rBV3	141690	233004	5.71%	0.617%
21	12.688	908	910	916	rBV	62043	91824	2.25%	0.243%
22	12.779	916	918	920	rBV	553082	471414	11.55%	1.248%
23	12.836	920	923	924	rBV2	49790	54180	1.33%	0.143%
24	12.962	931	934	936	rBV	3854046	3903654	95.64%	10.337%
25	13.191	951	954	956	rBV	96296	90053	2.21%	0.238%
26	13.488	978	980	982	rBV	292557	297642	7.29%	0.788%
27	13.534	982	984	985	rBV	101911	104945	2.57%	0.278%
28	13.865	1010	1013	1021	rBV	241927	517684	12.68%	1.371%
29	14.002	1023	1025	1028	rVB	821975	868764	21.29%	2.300%
30	14.174	1038	1040	1043	rBV	100350	93501	2.29%	0.248%
31	14.482	1064	1067	1068	rBV	55241	64921	1.59%	0.172%
32	14.779	1090	1093	1095	rVB	52756	63335	1.55%	0.168%
33	15.099	1118	1121	1124	rBV	48204	53249	1.30%	0.141%
34	15.431	1147	1150	1157	rVB	583478	771970	18.91%	2.044%

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Misc :
ALS Vial : 64 Sample Multiplier: 1

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ClientSampleId :
SB-04-COMP

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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35	15.980	1193	1198	1201	rBV6	14825	48464	1.19%	0.128%
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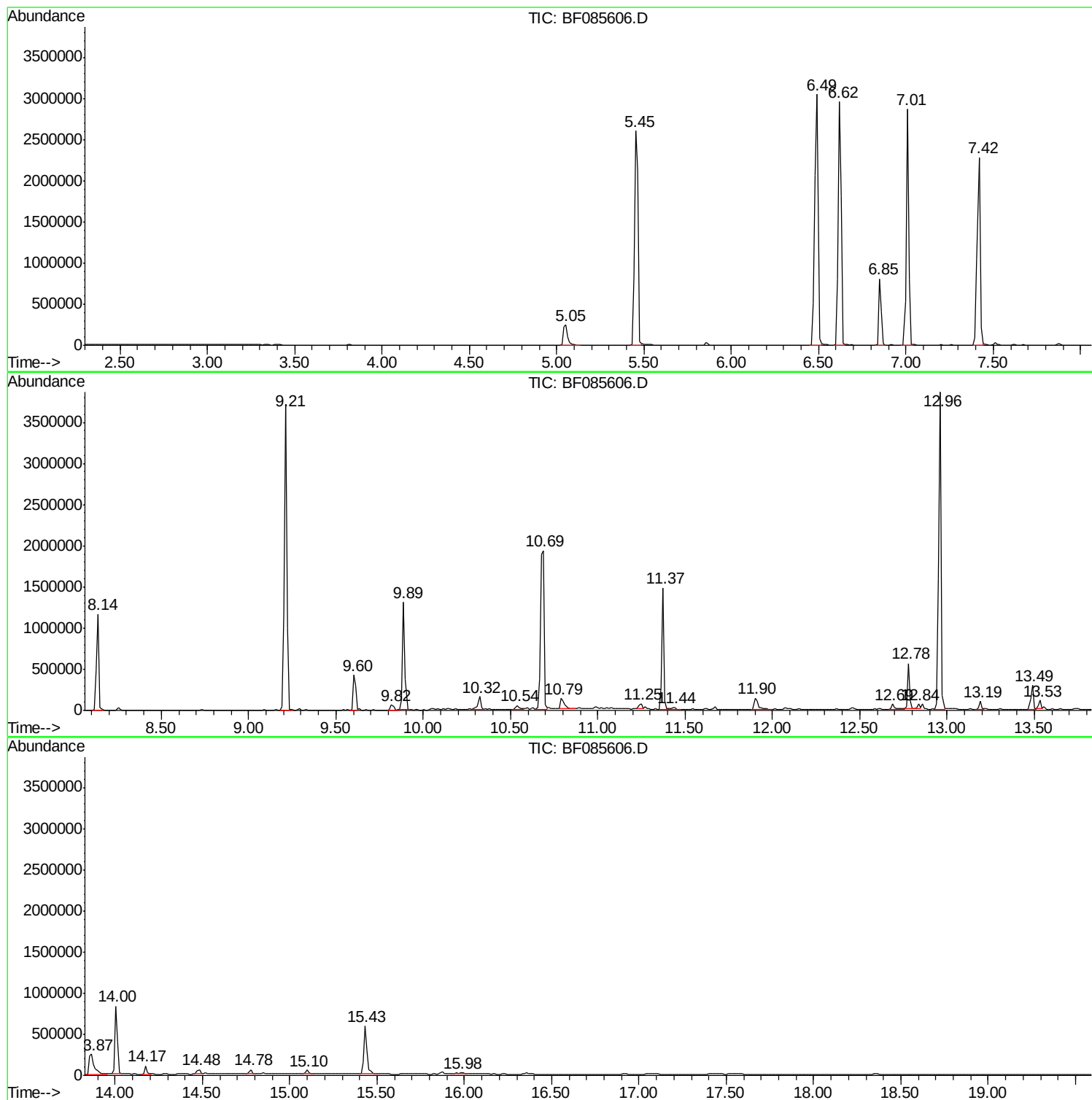
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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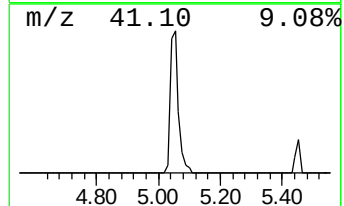
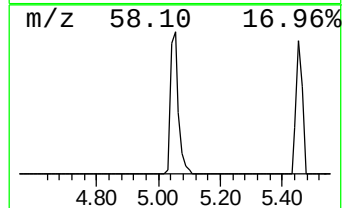
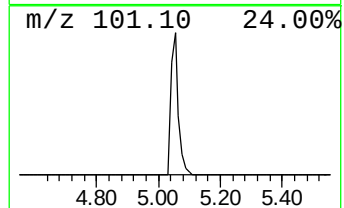
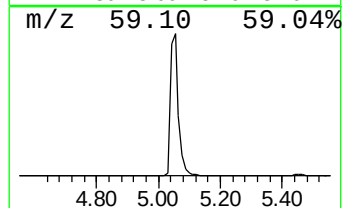
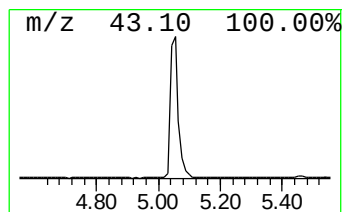
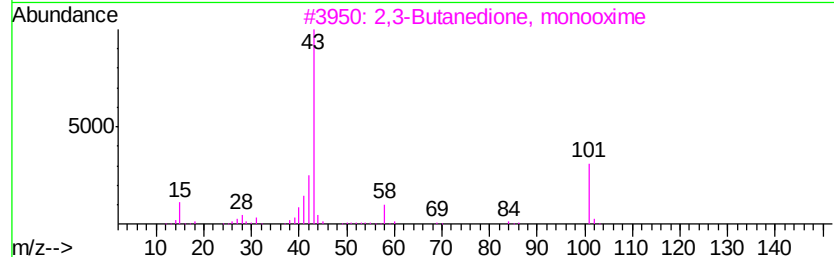
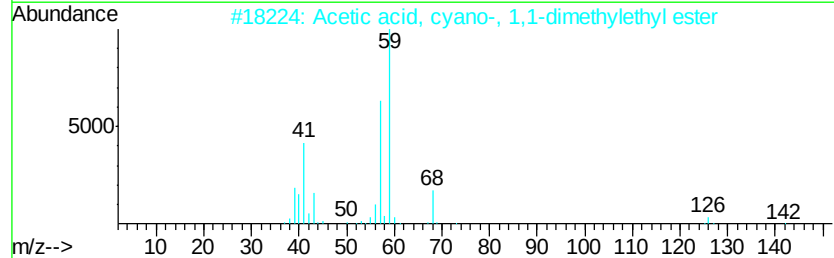
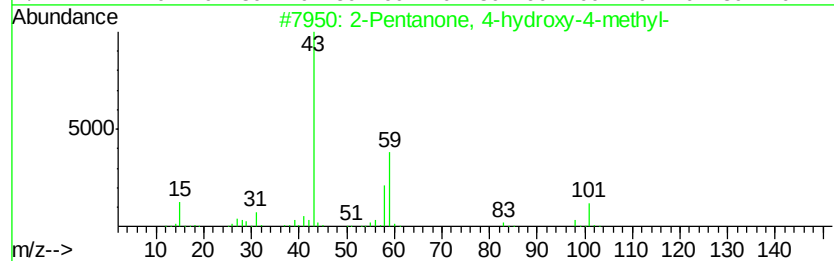
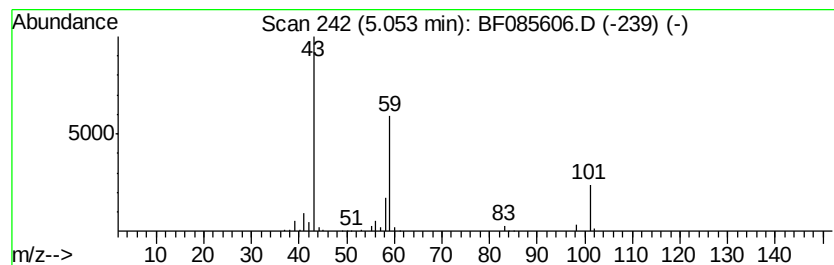
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	10.06 ng	427076	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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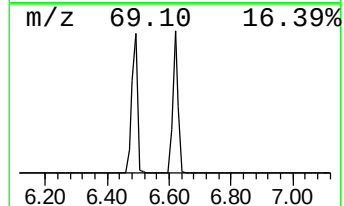
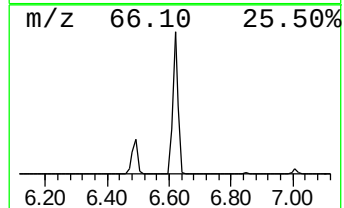
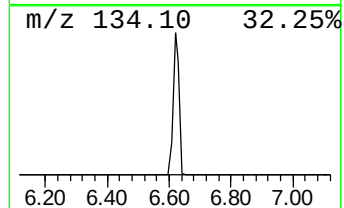
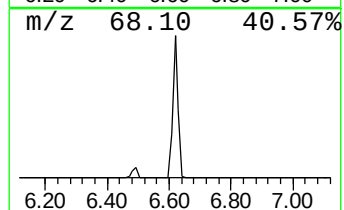
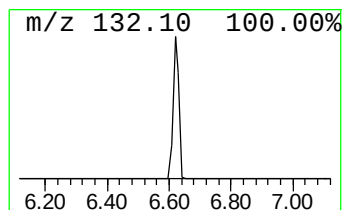
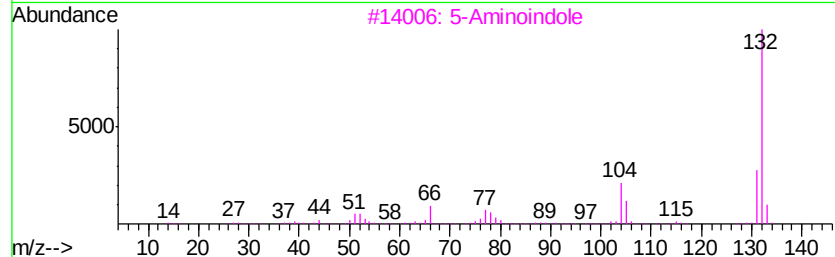
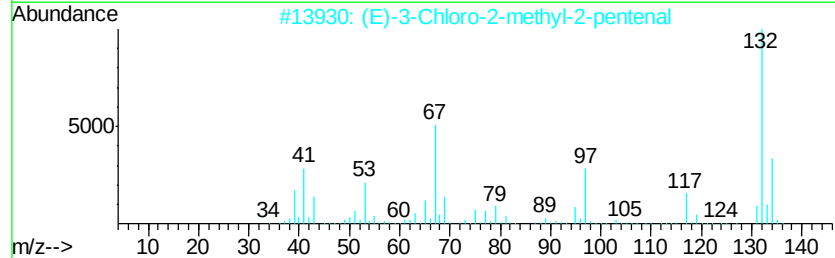
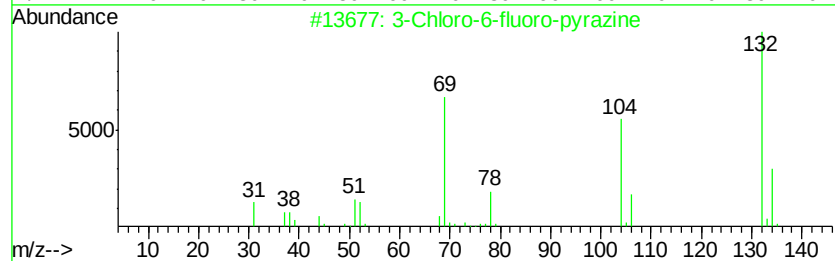
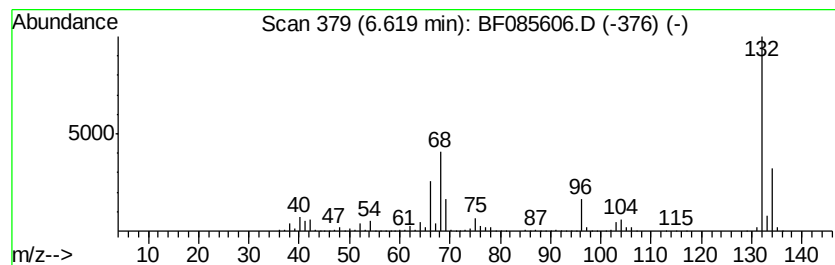
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	90.26 ng	3833590	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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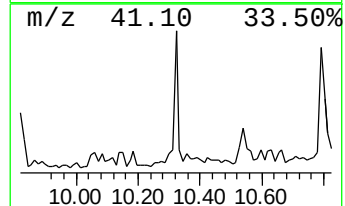
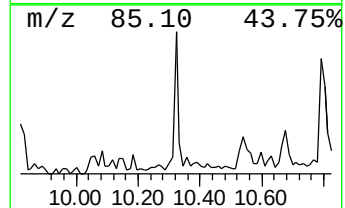
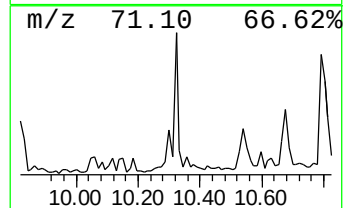
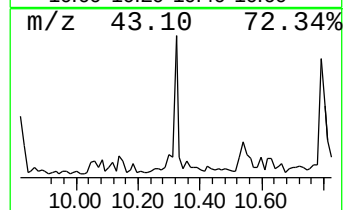
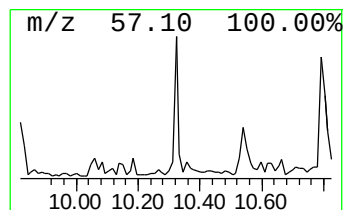
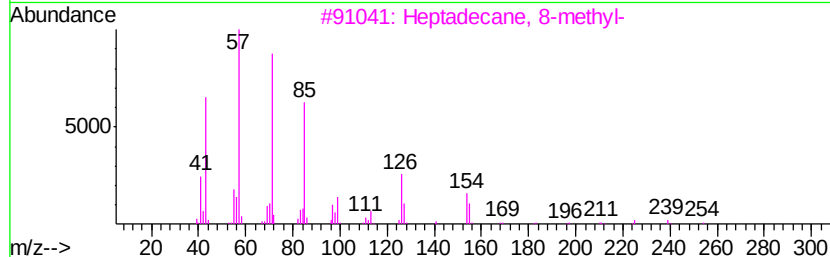
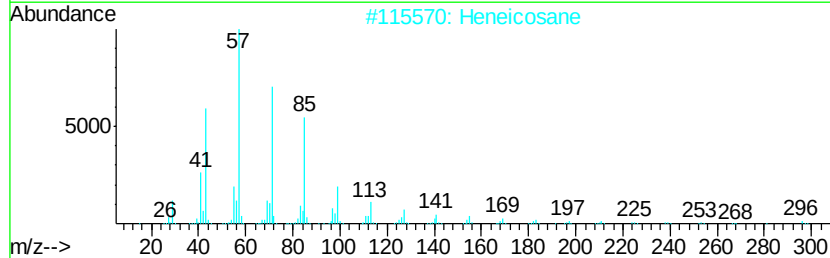
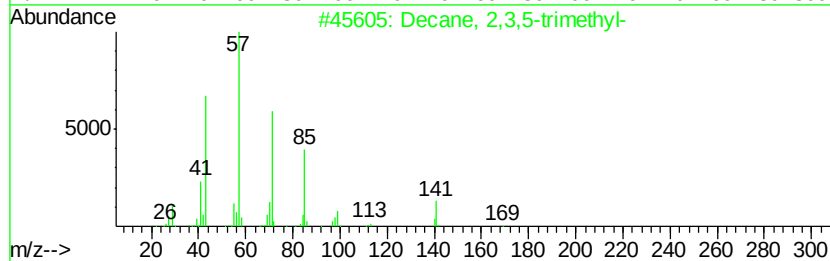
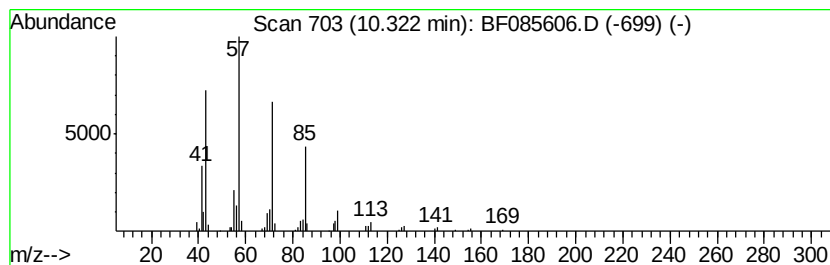
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Decane, 2,3,5-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.32	2.39 ng	146594	Acenaphthene-d10		9.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
2			Heneicosane	296	C21H44	000629-94-7	90
3			Heptadecane, 8-methyl-	254	C18H38	013287-23-5	90
4			Heptadecane	240	C17H36	000629-78-7	90
5			Pentadecane	212	C15H32	000629-62-9	86



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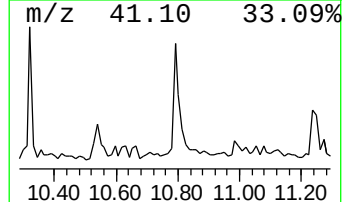
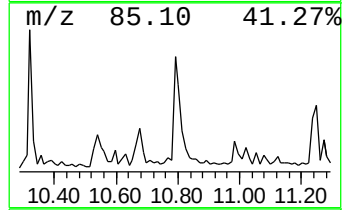
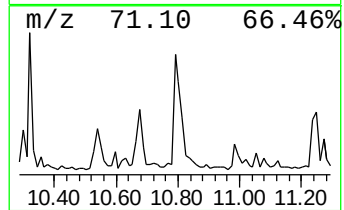
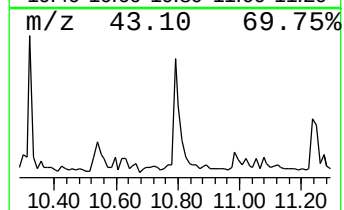
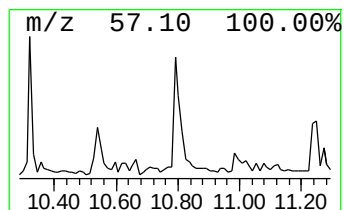
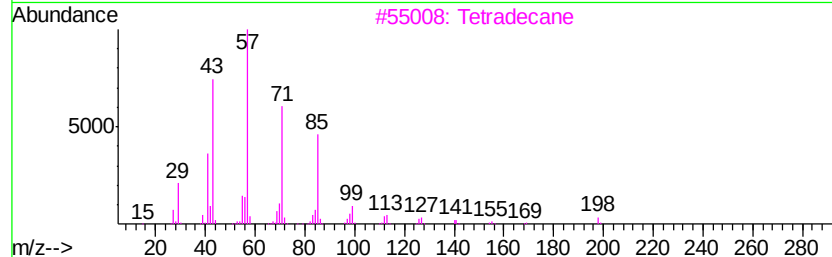
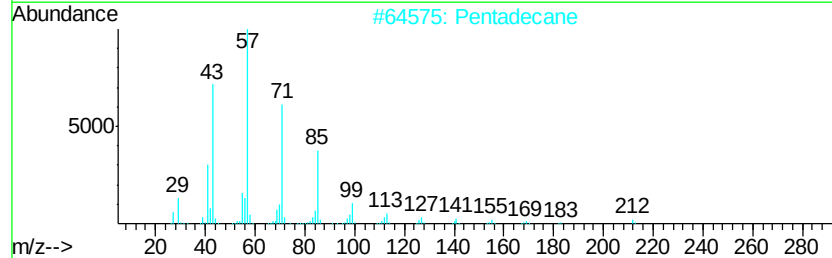
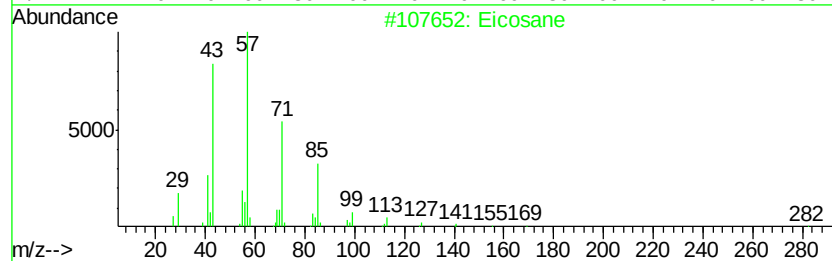
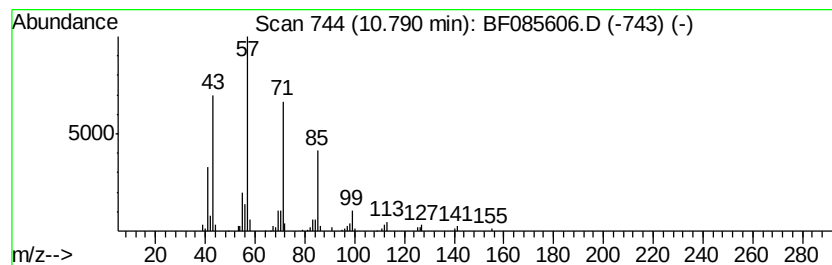
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	2.97 ng	188341	Phenanthrene-d10	11.37

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	Tetradecane		198	C14H30	000629-59-4	90
4	Hexadecane		226	C16H34	000544-76-3	86
5	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	83



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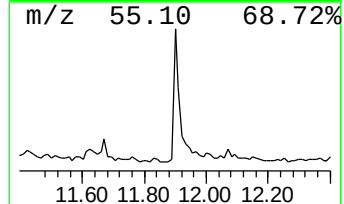
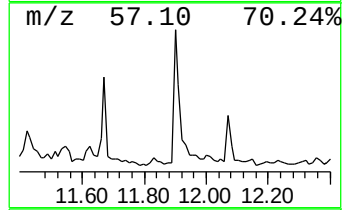
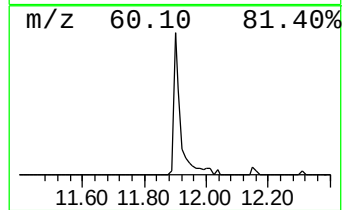
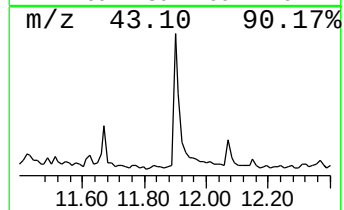
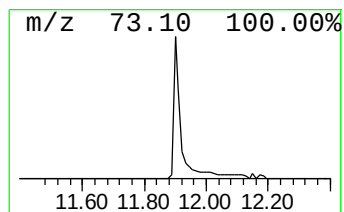
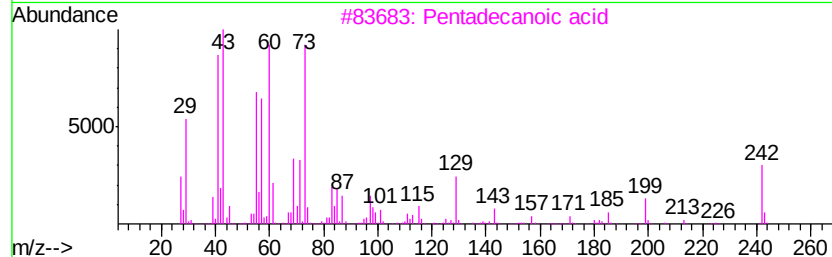
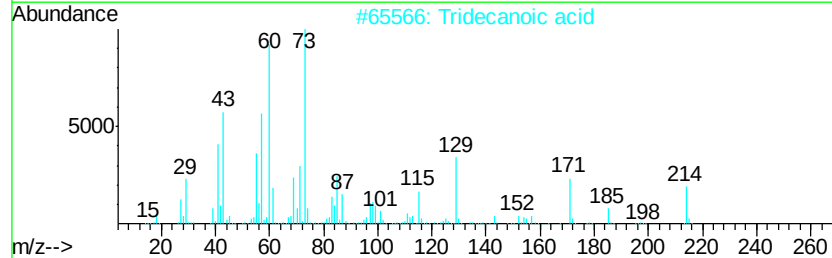
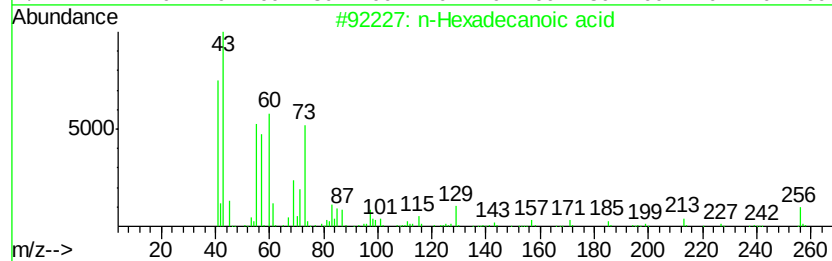
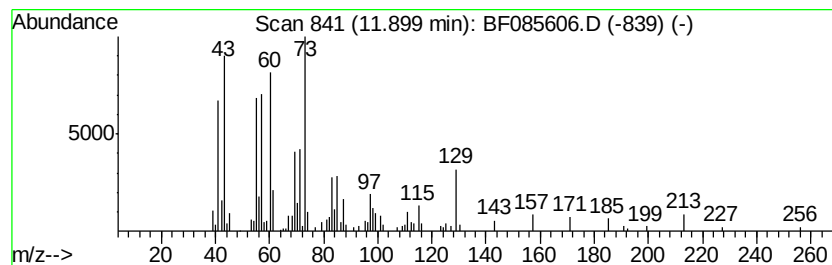
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.90	3.67 ng	233004	Phenanthrene-d10		11.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Tridecanoic acid	214	C13H26O2	000638-53-9	80
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	72
4	n-Decanoic acid	172	C10H20O2	000334-48-5	43
5	Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	25



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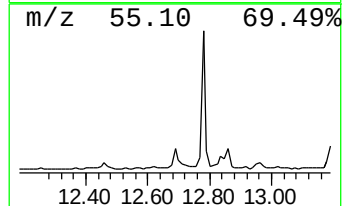
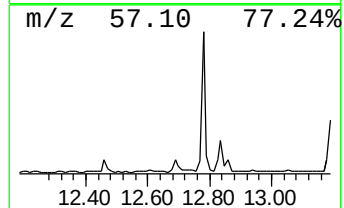
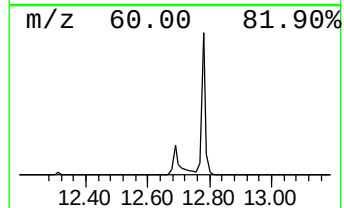
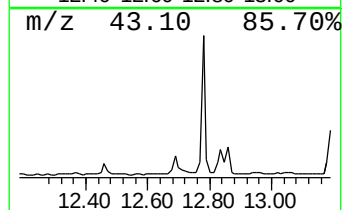
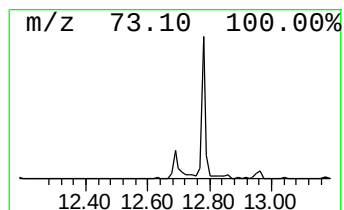
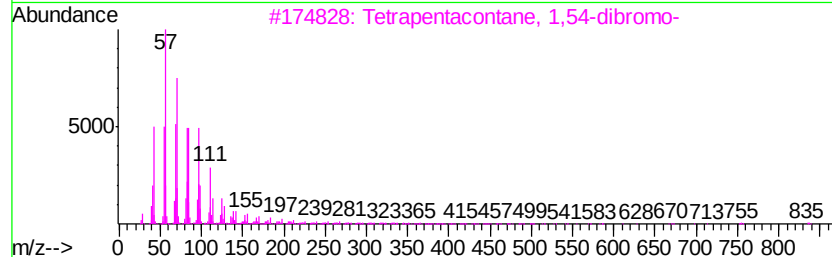
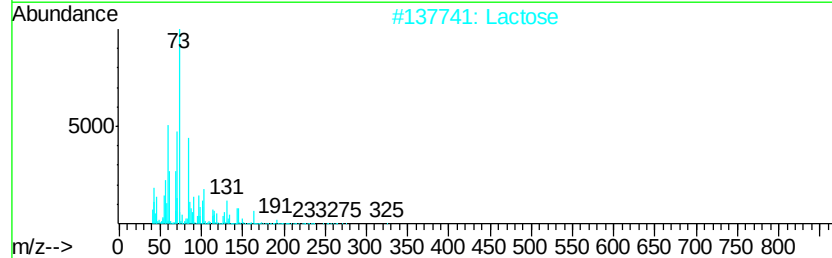
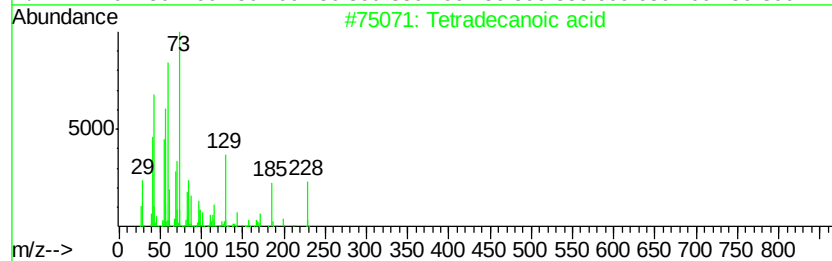
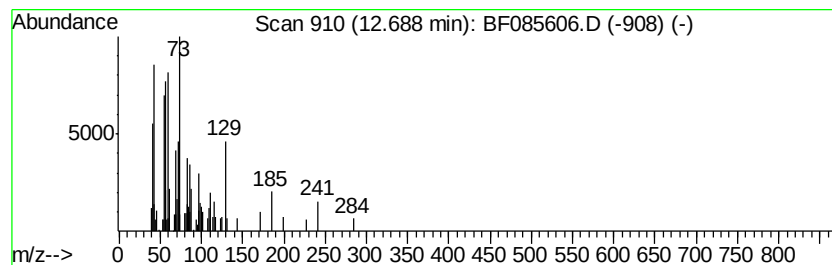
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	2.11 ng	91824	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
2		Lactose	342	C12H22O11	000063-42-3	38
3		Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	18
4		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	18
5		Acetic acid, octadecyl ester	312	C20H40O2	000822-23-1	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085606.D
Acq On : 20 Mar 2016 8:58
Operator : UM/SJ
Sample : H1815-08
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-04-COMP

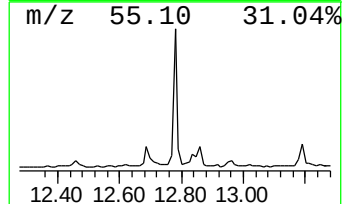
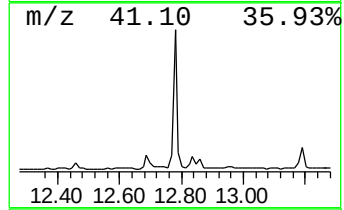
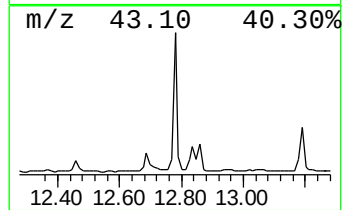
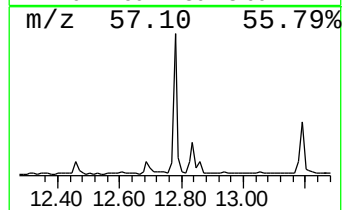
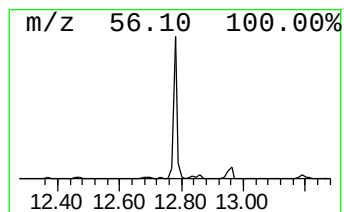
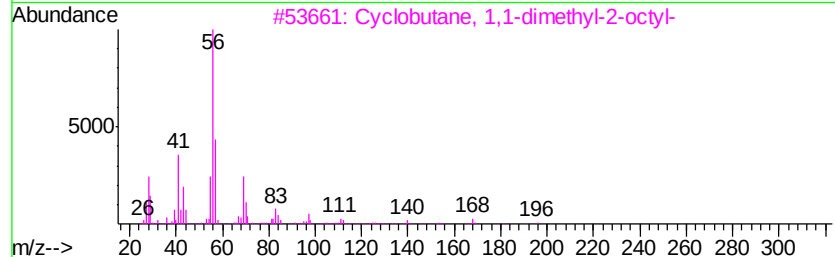
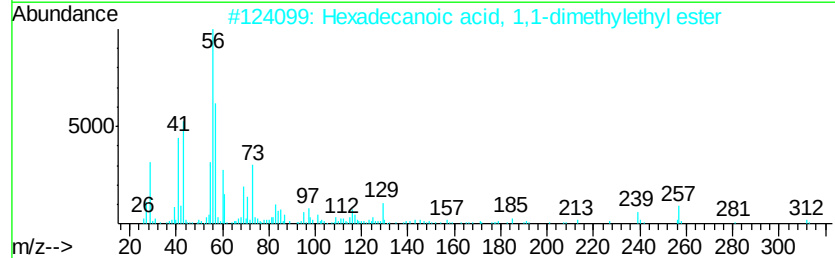
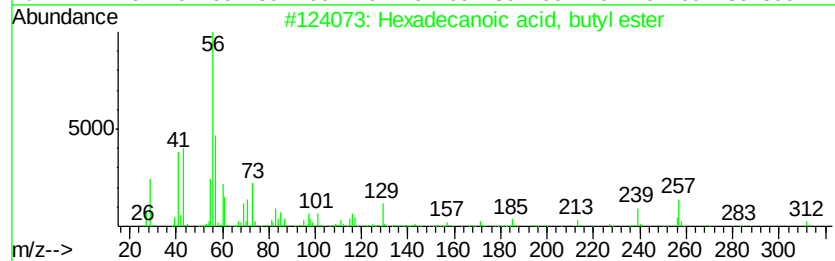
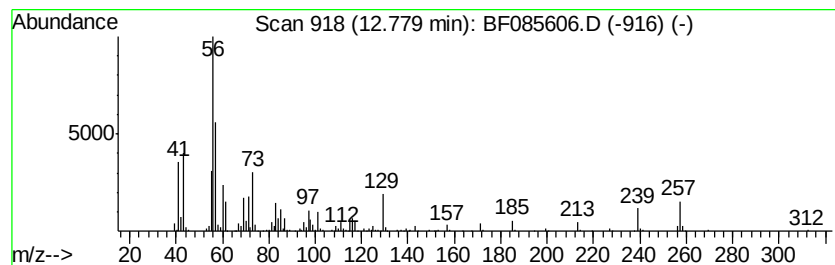
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, butyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.78	10.85 ng	471414	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	83
3		Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	35
4		Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	35
5		2-Methyl-n-1-tridecene	196	C14H28	018094-01-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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Operator : UM/SJ
Sample : H1815-08
Misc :
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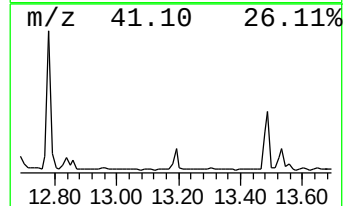
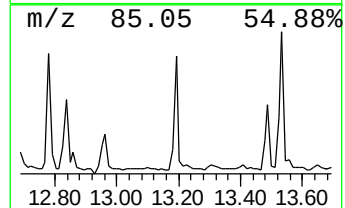
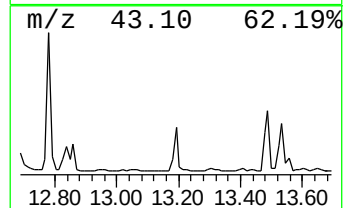
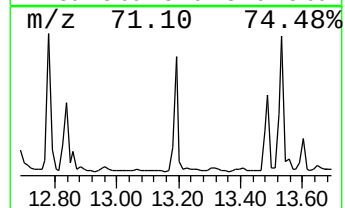
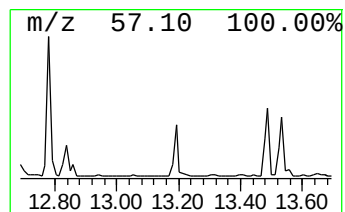
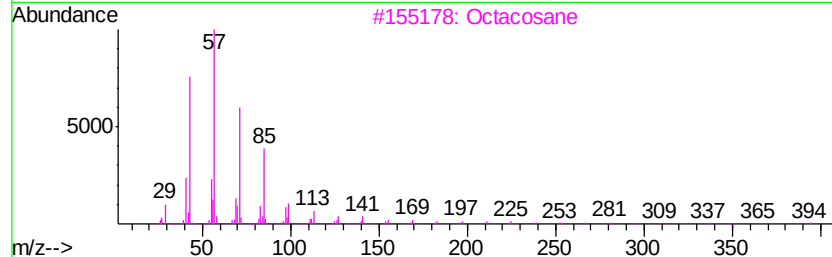
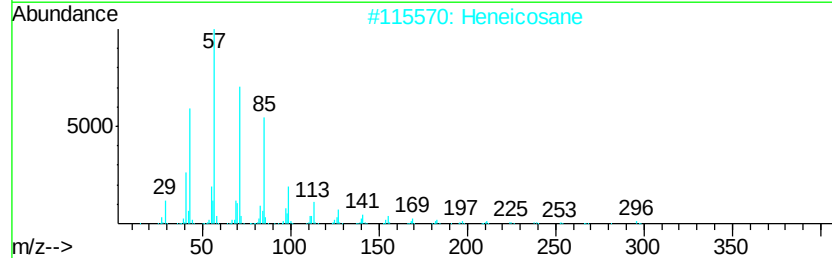
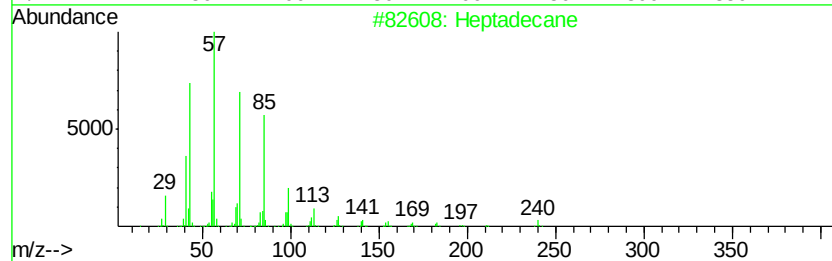
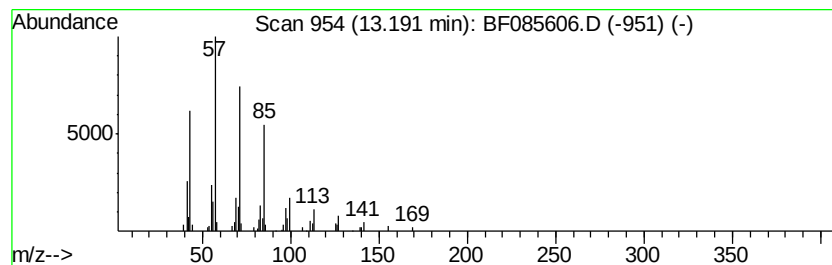
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	2.07 ng	90053	Chrysene-d12	14.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	91
2	Heneicosane	296 C21H44	000629-94-7	91
3	Octacosane	394 C28H58	000630-02-4	91
4	Heptacosane	380 C27H56	000593-49-7	91
5	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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Acq On : 20 Mar 2016 8:58
Operator : UM/SJ
Sample : H1815-08
Misc :
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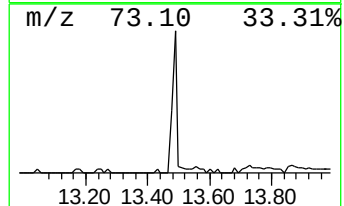
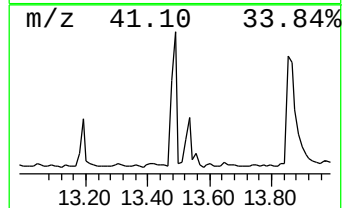
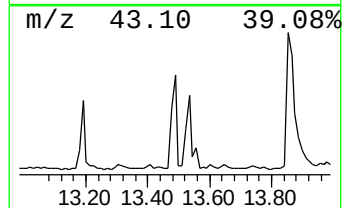
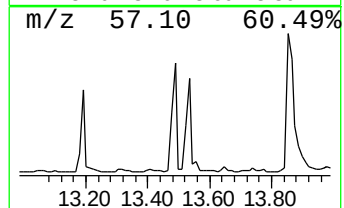
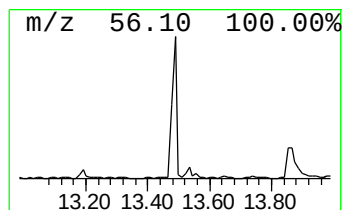
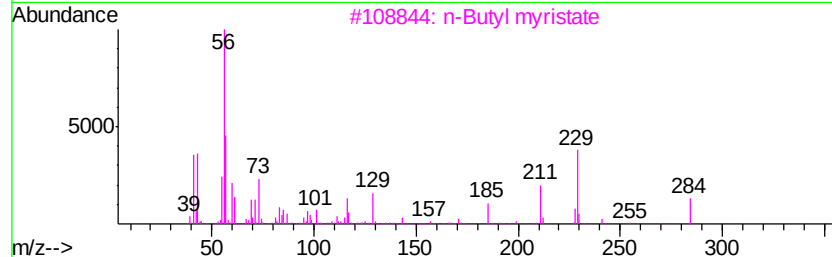
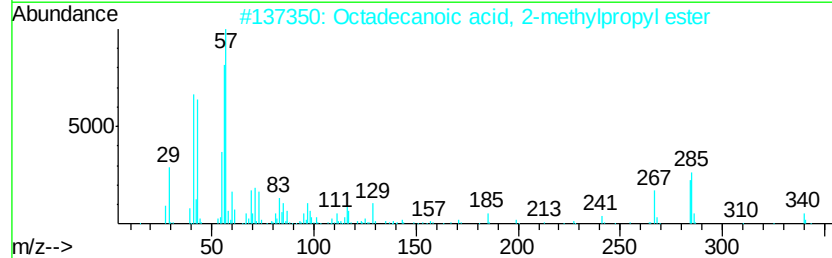
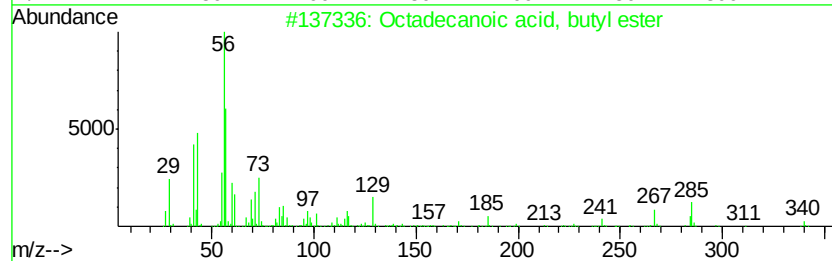
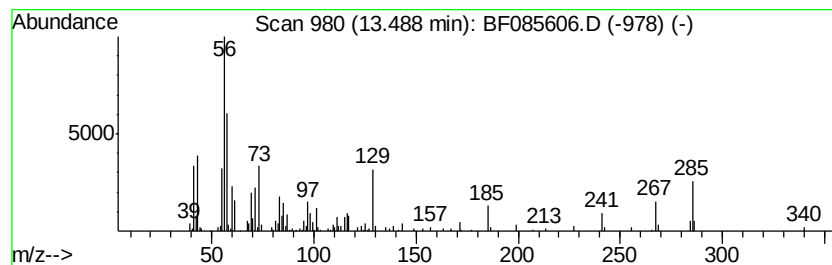
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecanoic acid, butyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	6.85 ng	297642	Chrysene-d12	14.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	93
2		Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	59
3		n-Butyl myristate	284	C18H36O2	000110-36-1	58
4		Nipecotic acid	129	C6H11NO2	000498-95-3	47
5		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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Operator : UM/SJ
Sample : H1815-08
Misc :
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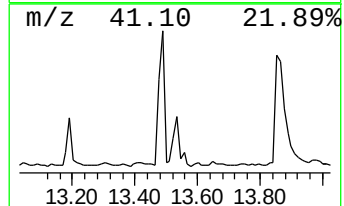
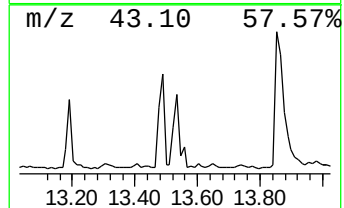
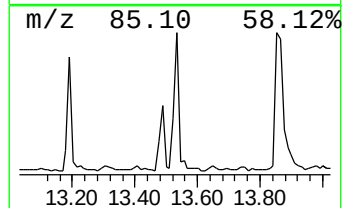
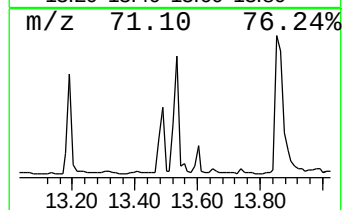
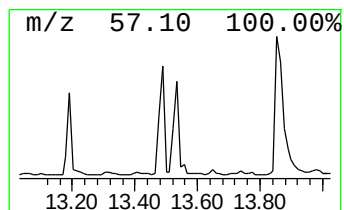
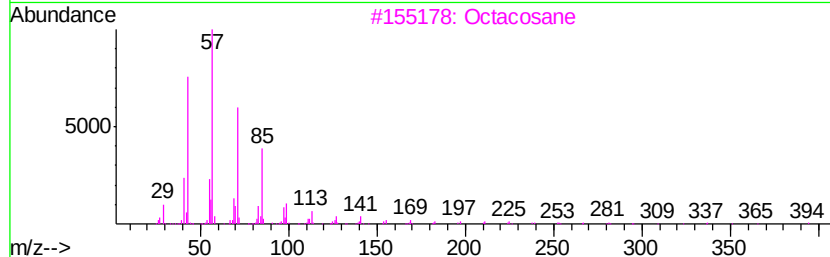
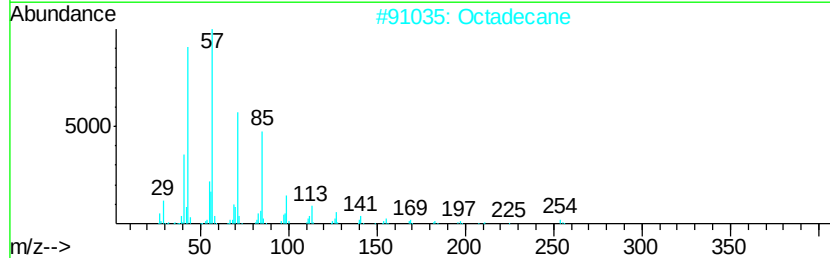
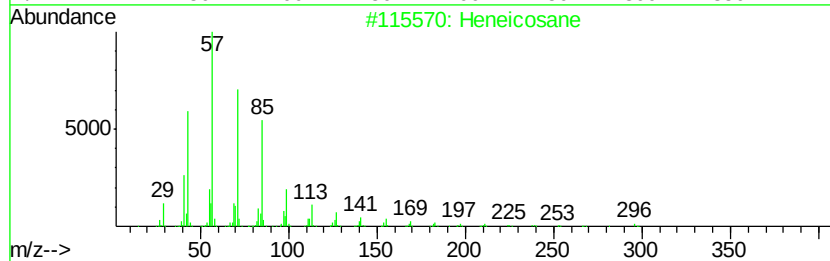
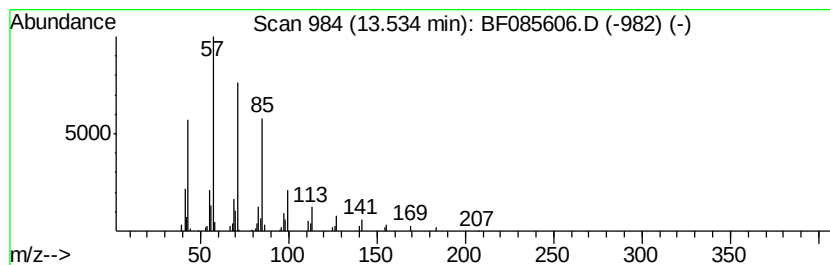
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	2.42 ng	104945	Chrysene-d12	14.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	91
2	Octadecane	254 C18H38	000593-45-3	90
3	Octacosane	394 C28H58	000630-02-4	87
4	Tetracosane	338 C24H50	000646-31-1	86
5	Hexadecane	226 C16H34	000544-76-3	86



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Acq On : 20 Mar 2016 8:58
Operator : UM/SJ
Sample : H1815-08
Misc :
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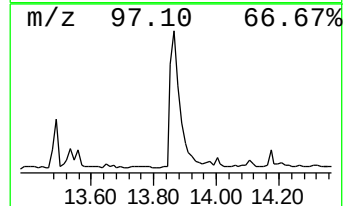
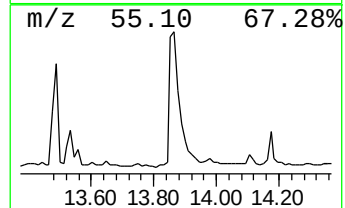
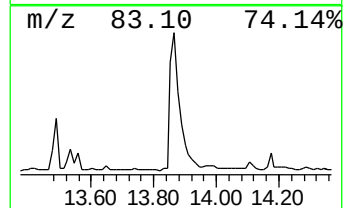
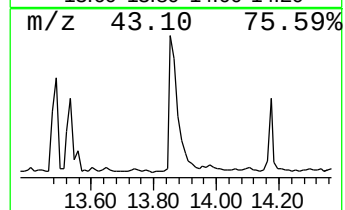
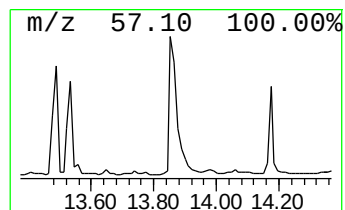
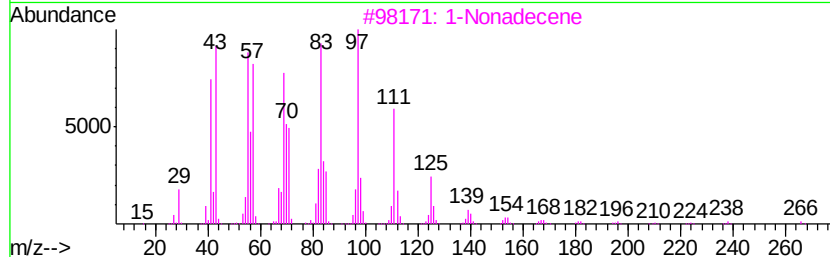
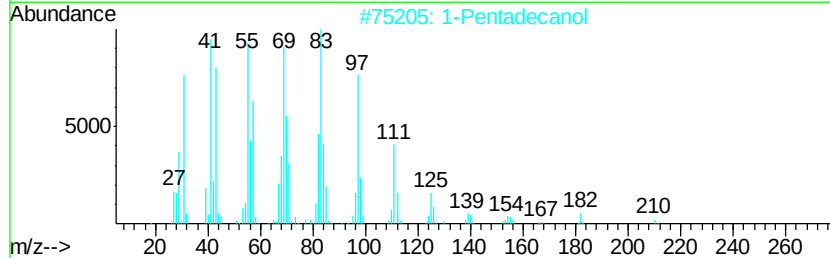
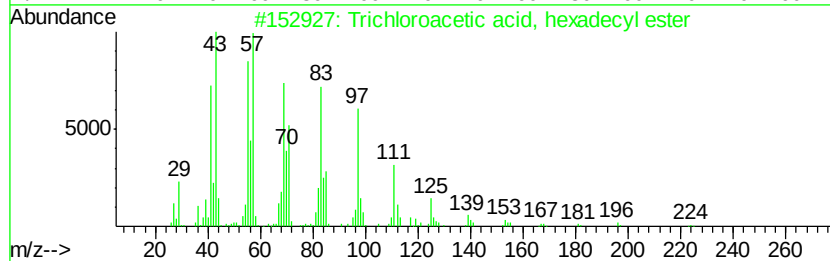
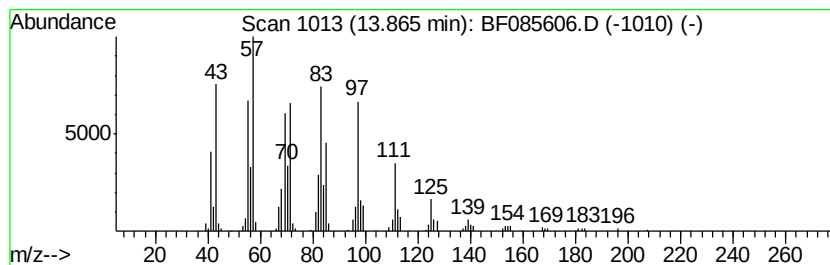
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Trichloroacetic acid, hexad... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	11.92 ng	517684	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
2		1-Pentadecanol	228	C15H32O	000629-76-5	76
3		1-Nonadecene	266	C19H38	018435-45-5	76
4		5-Eicosene, (E)-	280	C20H40	074685-30-6	74
5		1-Docosene	308	C22H44	001599-67-3	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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Acq On : 20 Mar 2016 8:58
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Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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	10.1 ng		427076	1	6.85	849446	20.0
unknown6.62	6.62	90.3 ng		3833590	1	6.85	849446	20.0
Decane, 2,3,5-tri...	10.32	2.4 ng		146594	3	9.89	1228200	20.0
Eicosane	10.79	3.0 ng		188341	4	11.37	1269650	20.0
n-Hexadecanoic acid	11.90	3.7 ng		233004	4	11.37	1269650	20.0
Tetradecanoic acid	12.69	2.1 ng		91824	5	14.00	868764	20.0
Hexadecanoic acid...	12.78	10.9 ng		471414	5	14.00	868764	20.0
Heptadecane	13.19	2.1 ng		90053	5	14.00	868764	20.0
Octadecanoic acid...	13.49	6.8 ng		297642	5	14.00	868764	20.0
Heneicosane	13.53	2.4 ng		104945	5	14.00	868764	20.0
Trichloroacetic a...	13.87	11.9 ng		517684	5	14.00	868764	20.0