

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
 Data File : BF085749.D
 Acq On : 25 Mar 2016 8:29
 Operator : UM/SJ
 Sample : H1885-04
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-GRID-15B

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-BF032416.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	4.390	179	184	188	rBV2	35150	62973	2.01%	0.240%
2	5.007	235	238	246	rBV	309267	473509	15.10%	1.803%
3	5.430	272	275	277	rBV	2124035	2594071	82.74%	9.880%
4	5.716	298	300	303	rBV	43784	44898	1.43%	0.171%
5	5.830	307	310	312	rVB	48758	49922	1.59%	0.190%
6	6.459	362	365	368	rBV	1851153	2617631	83.49%	9.970%
7	6.596	374	377	379	rBV	2692155	2746641	87.61%	10.461%
8	6.664	379	383	385	rVB	37200	45895	1.46%	0.175%
9	6.824	395	397	399	rBV	787544	671549	21.42%	2.558%
10	6.984	408	411	413	rBV	2062805	2208449	70.44%	8.411%
11	7.396	444	447	449	rBV	1501795	1828955	58.34%	6.966%
12	7.499	452	456	462	rVB	78041	109703	3.50%	0.418%
13	8.105	507	509	512	rBV	730449	838587	26.75%	3.194%
14	9.190	601	604	606	rBV	3253904	3135166	100.00%	11.941%
15	9.579	636	638	641	rBV	616325	561491	17.91%	2.139%
16	9.796	655	657	659	rVB	126642	95608	3.05%	0.364%
17	9.865	661	663	665	rVB	943849	850634	27.13%	3.240%
18	10.025	675	677	681	rBV2	20978	43642	1.39%	0.166%
19	10.299	697	701	702	rBV	123087	159064	5.07%	0.606%
20	10.516	717	720	723	rBV	48000	83378	2.66%	0.318%
21	10.653	729	732	734	rBV	1465808	1493955	47.65%	5.690%
22	10.768	740	742	749	rVB	156193	185258	5.91%	0.706%
23	10.962	756	759	761	rBV2	21787	31718	1.01%	0.121%
24	11.213	779	781	782	rBV	97296	77276	2.46%	0.294%
25	11.248	782	784	787	rVB	26267	36564	1.17%	0.139%
26	11.339	790	792	795	rBV2	542052	668494	21.32%	2.546%
27	11.568	810	812	817	rBV	159041	232512	7.42%	0.886%
28	11.636	817	818	821	rVB	47926	43542	1.39%	0.166%
29	11.876	837	839	841	rBV	32777	41456	1.32%	0.158%
30	12.128	858	861	866	rVB2	25247	33729	1.08%	0.128%
31	12.391	882	884	887	rBV	43276	65252	2.08%	0.249%
32	12.756	914	916	918	rBV	42605	61514	1.96%	0.234%
33	12.802	918	920	925	rVB	36913	36680	1.17%	0.140%
34	12.928	929	931	933	rBV	2347839	2064773	65.86%	7.864%

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35	13.156	949	951	953	rVB	34889	31571	1.01%	0.120%
36	13.408	968	973	975	rBV	289413	455189	14.52%	1.734%
37	13.488	978	980	985	rVB2	34429	72459	2.31%	0.276%
38	13.831	1007	1010	1014	rBV	56874	105350	3.36%	0.401%
39	13.979	1021	1023	1025	rBV	611096	596960	19.04%	2.274%
40	14.094	1030	1033	1036	rBV	33262	76348	2.44%	0.291%
41	15.397	1144	1147	1152	rVB	512323	623704	19.89%	2.375%

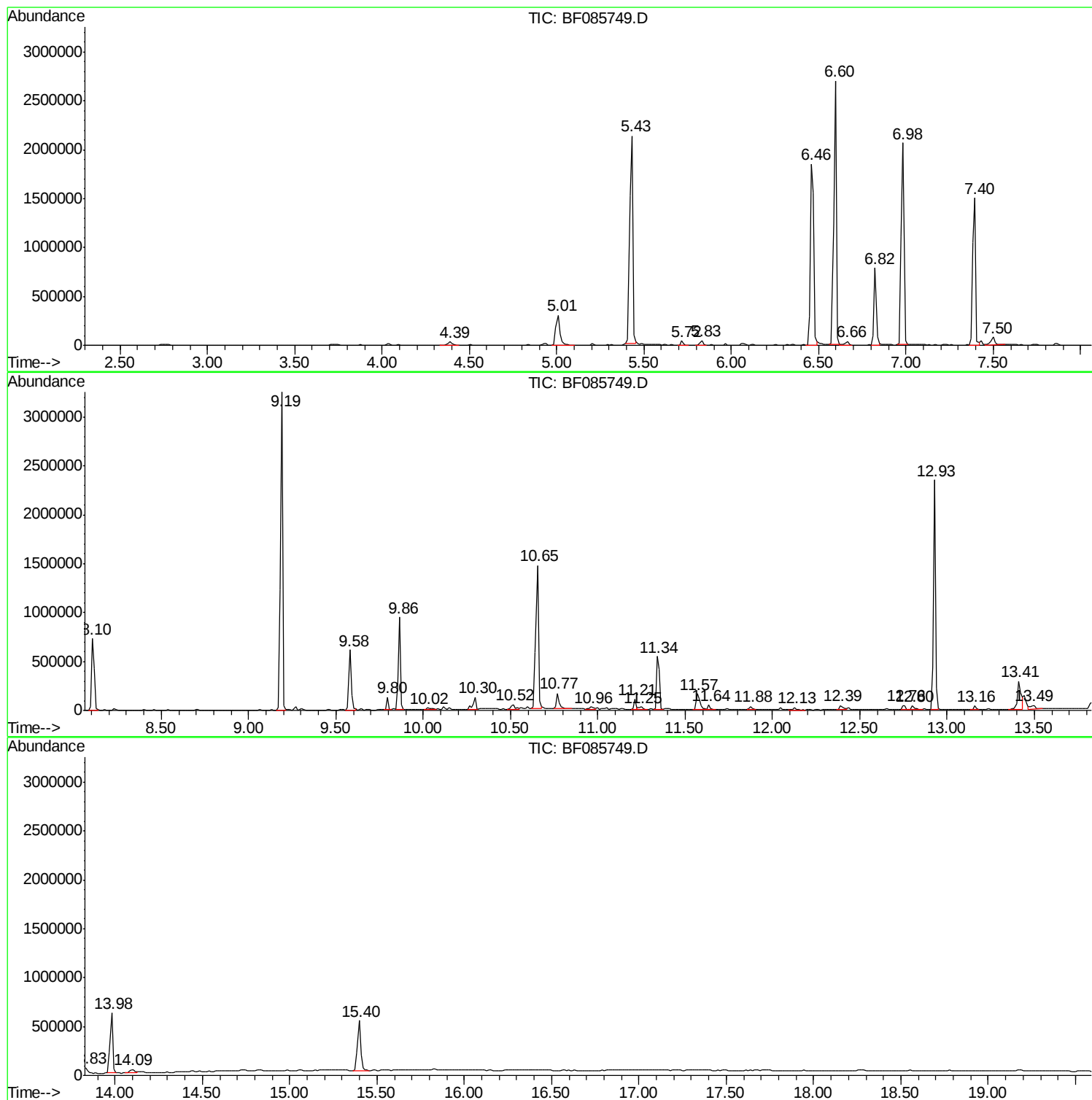
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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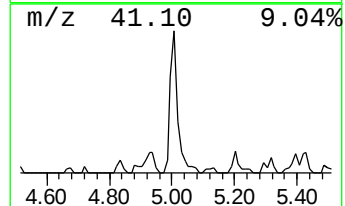
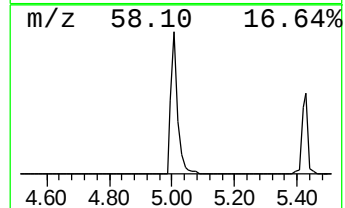
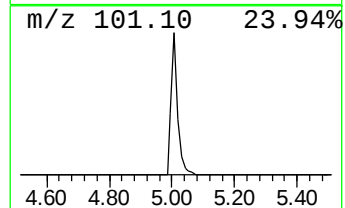
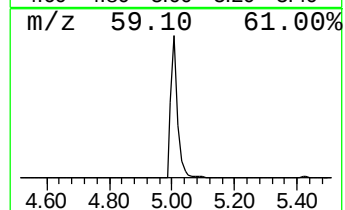
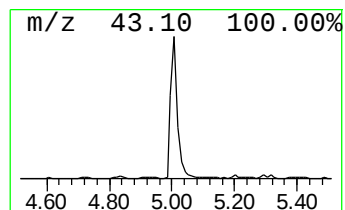
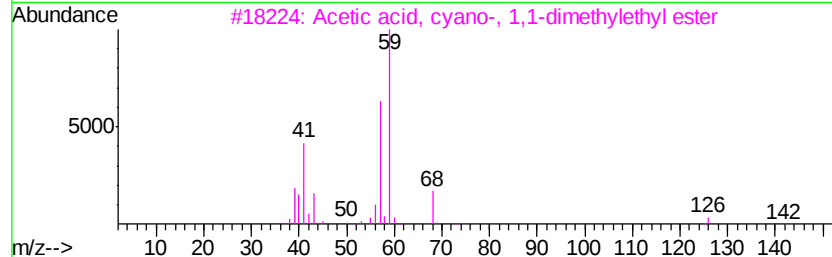
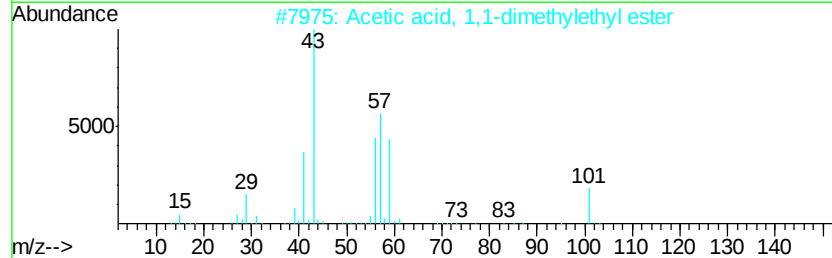
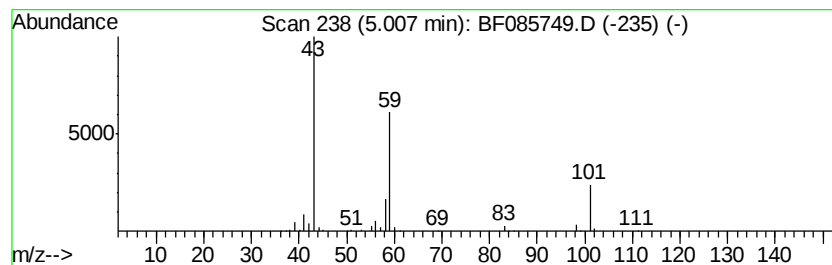
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	14.10 ng	473509	1,4-Dichlorobenzene-d4	6.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	



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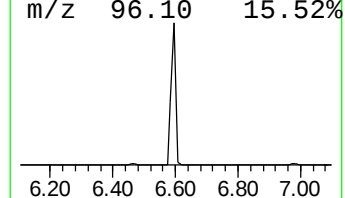
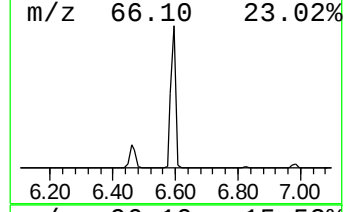
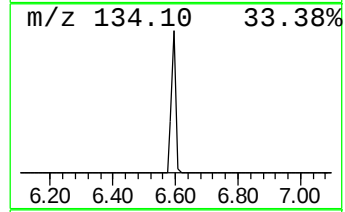
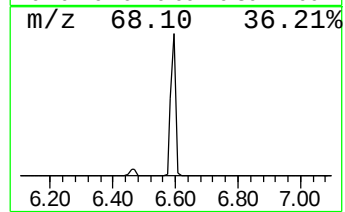
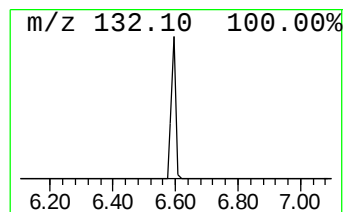
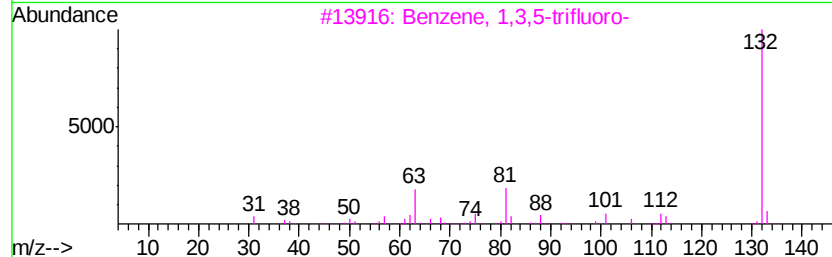
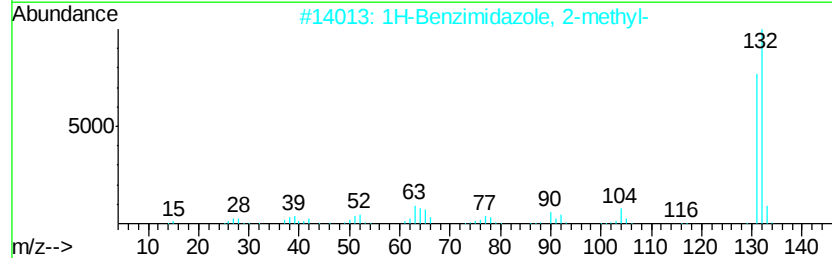
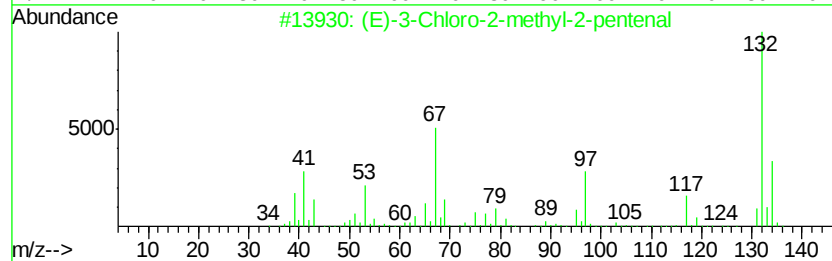
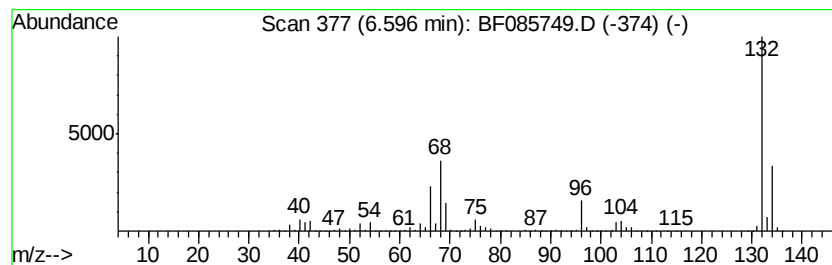
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	81.80 ng	2746640	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Xenon	132	Xe	007440-63-3	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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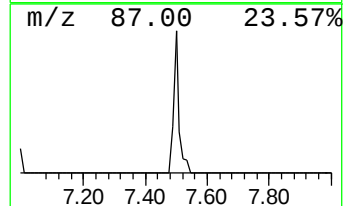
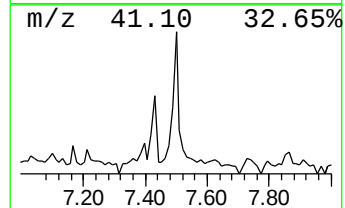
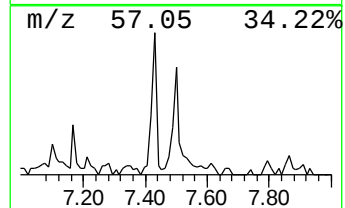
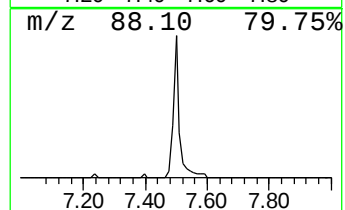
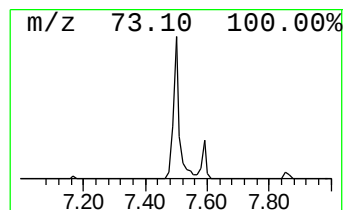
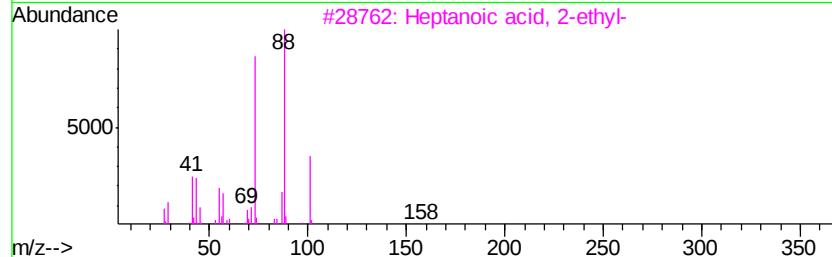
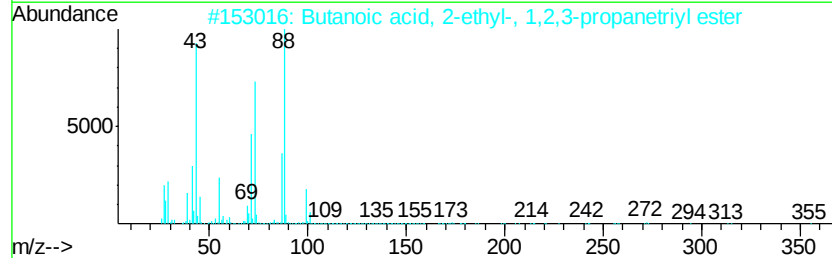
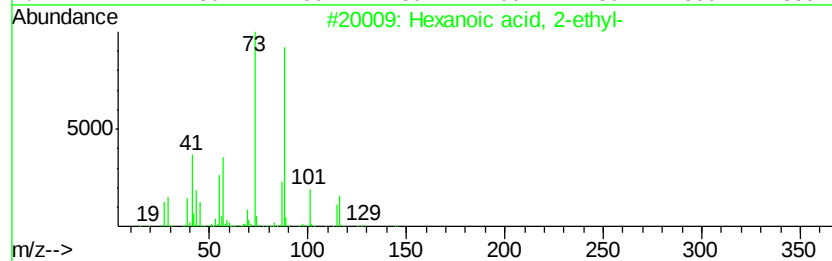
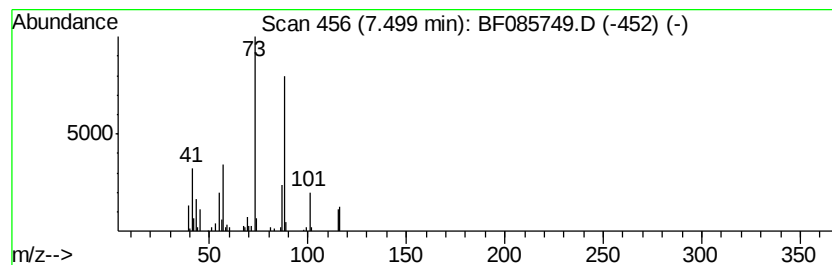
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Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	2.62 ng	109703	Naphthalene-d8	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	50
3		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	45
4		Octanoic acid, 2-methyl-, methyl...	172	C10H20O2	002177-86-8	40
5		1,4-Dioxane	88	C4H8O2	000123-91-1	37



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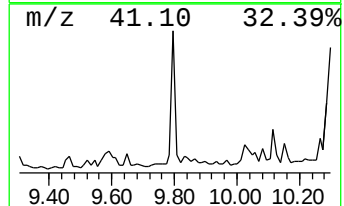
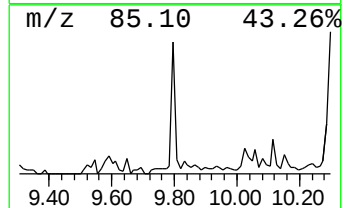
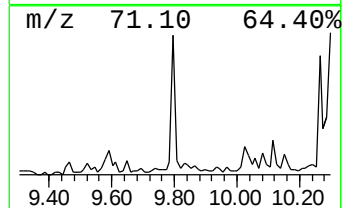
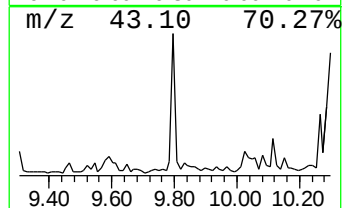
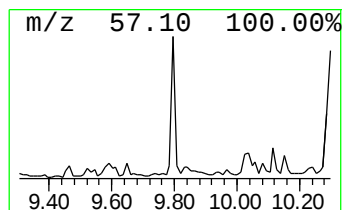
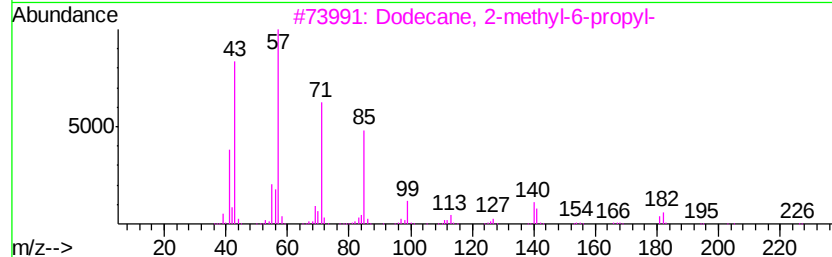
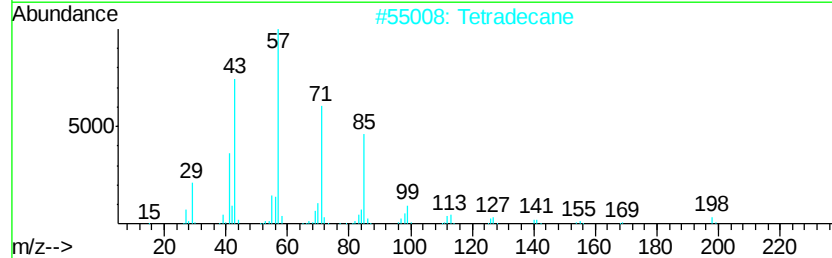
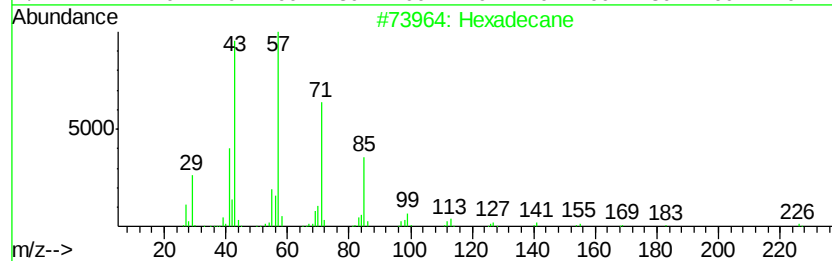
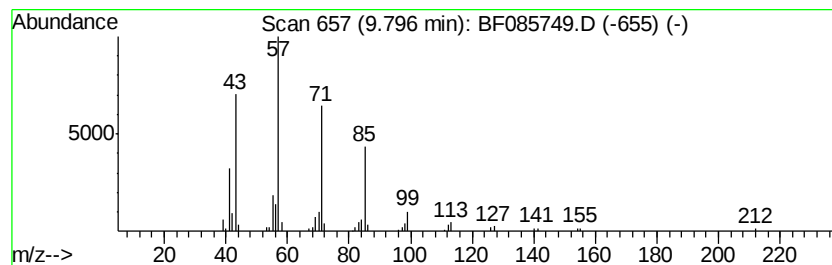
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.80	2.25 ng	95608	Acenaphthene-d10	9.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	90
2		Tetradecane	198	C14H30	000629-59-4	87
3		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4		10-Methylnonadecane	282	C20H42	056862-62-5	83
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83



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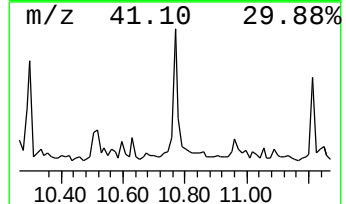
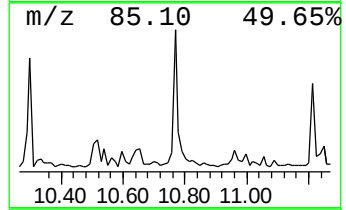
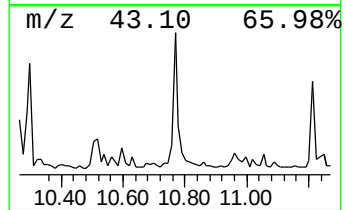
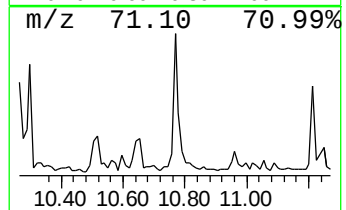
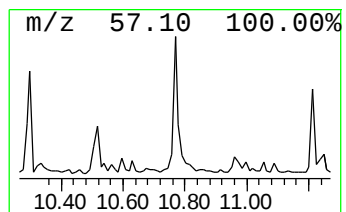
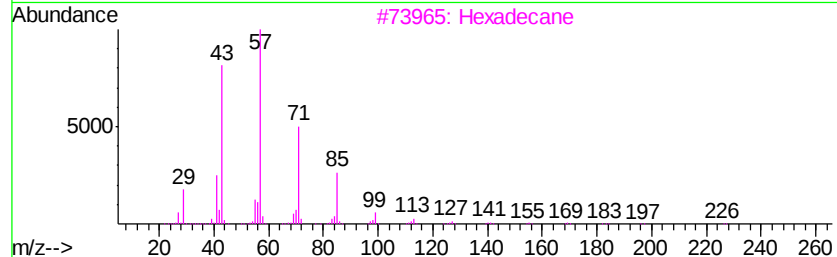
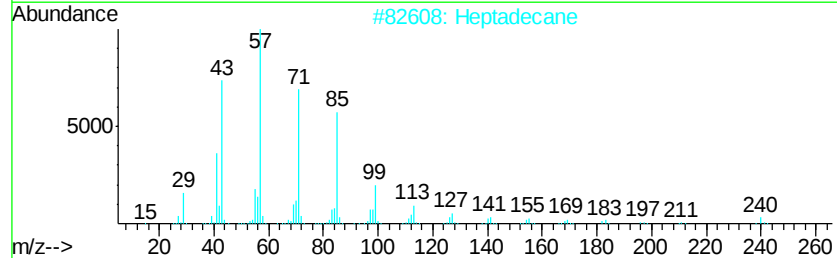
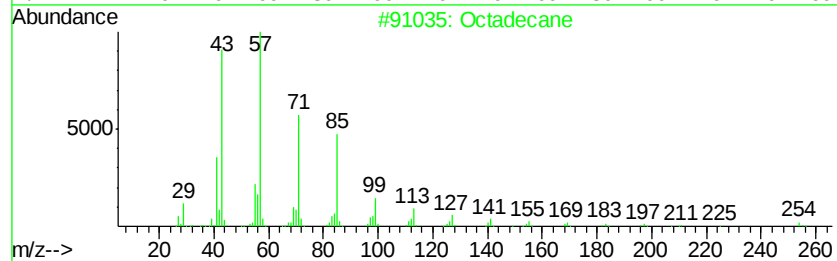
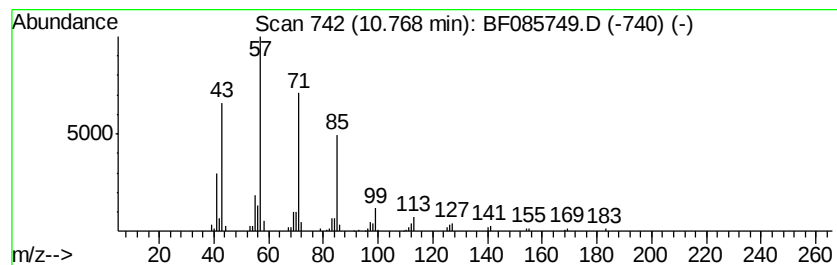
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Peak Number 7 Octadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	5.54 ng	185258	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Hexadecane	226	C16H34	000544-76-3	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90



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Operator : UM/SJ
Sample : H1885-04
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-15B

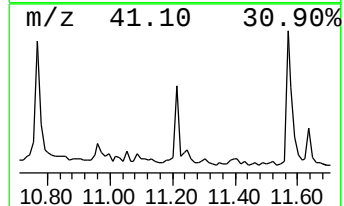
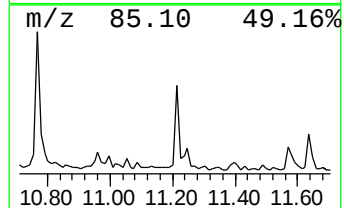
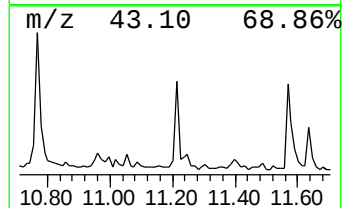
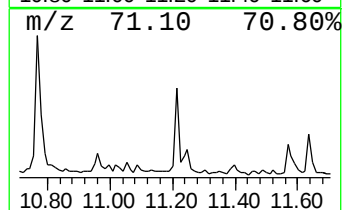
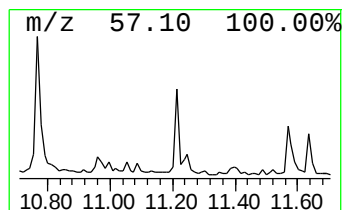
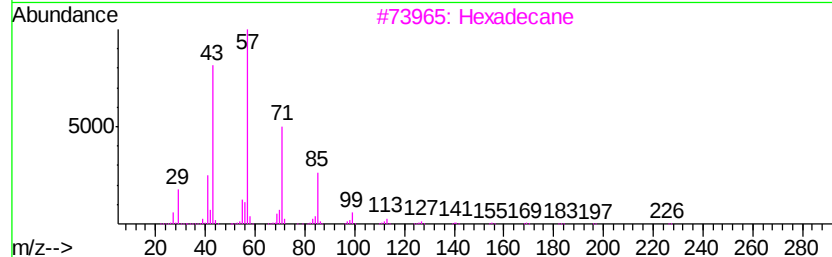
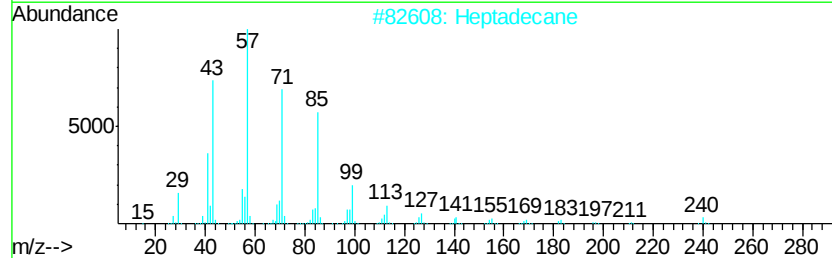
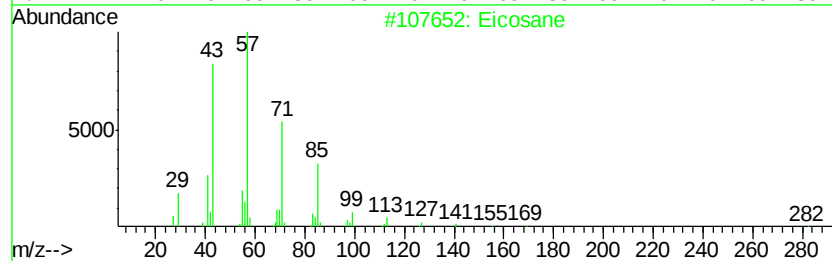
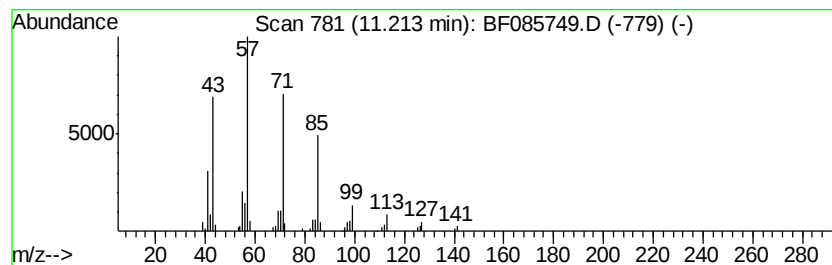
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.21	2.31 ng	77276	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Hexadecane	226	C16H34	000544-76-3	90
4		Octadecane	254	C18H38	000593-45-3	90
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
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Sample : H1885-04
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-15B

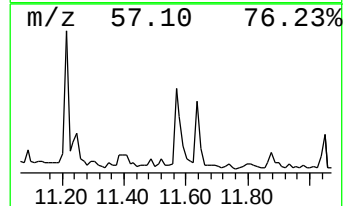
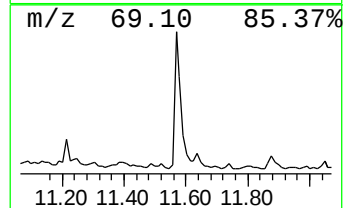
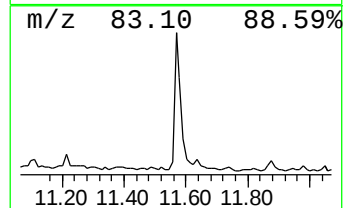
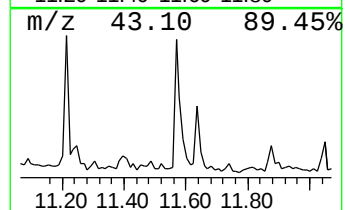
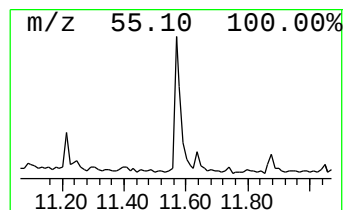
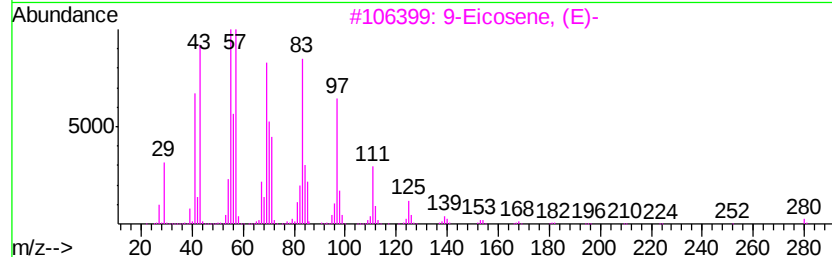
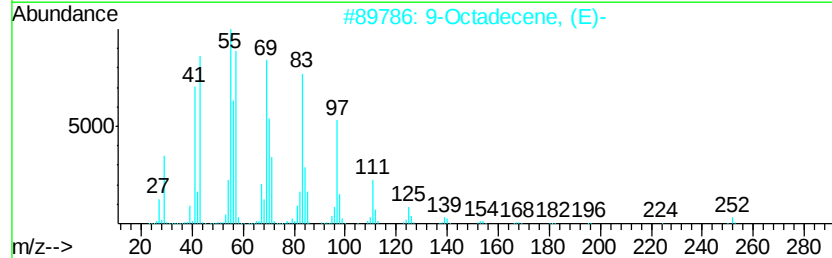
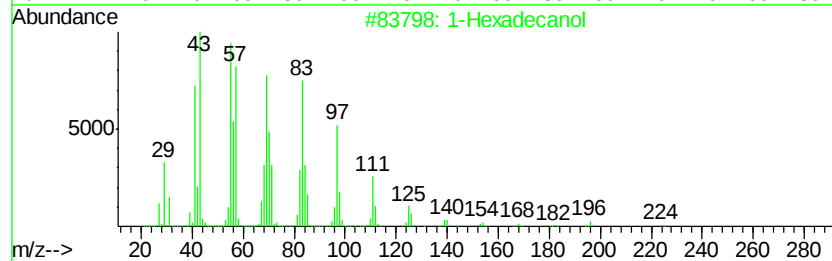
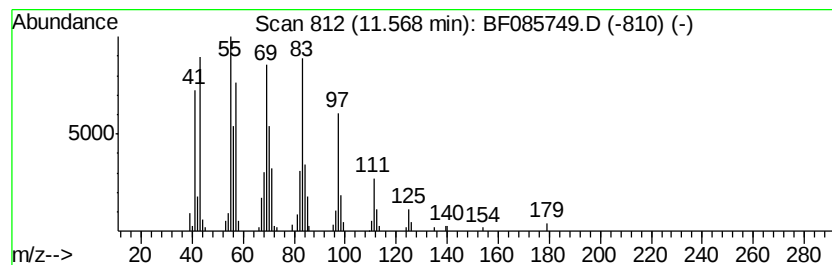
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Hexadecanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	6.96 ng	232512	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexadecanol	242	C16H34O	036653-82-4	95
2		9-Octadecene, (E)-	252	C18H36	007206-25-9	91
3		9-Eicosene, (E)-	280	C20H40	074685-29-3	91
4		5-Octadecene, (E)-	252	C18H36	007206-21-5	91
5		1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
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Sample : H1885-04
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TRACK-D-GRID-15B

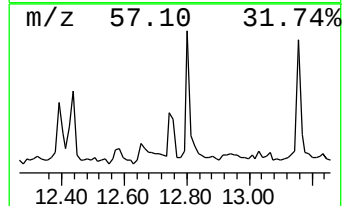
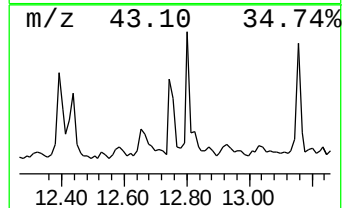
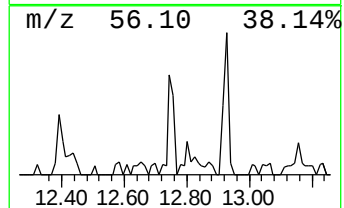
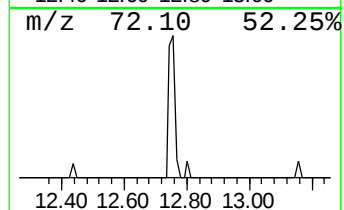
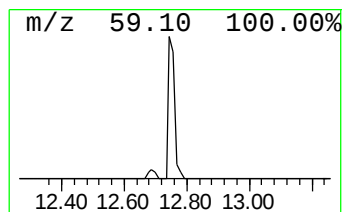
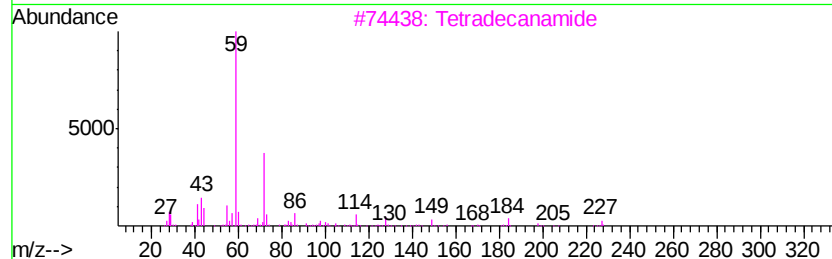
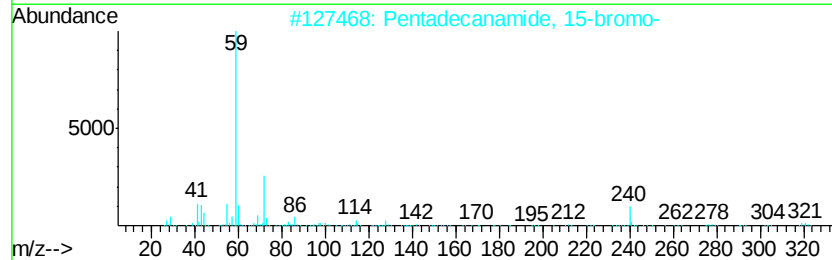
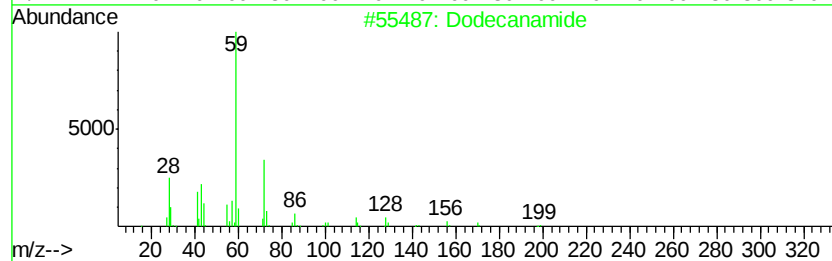
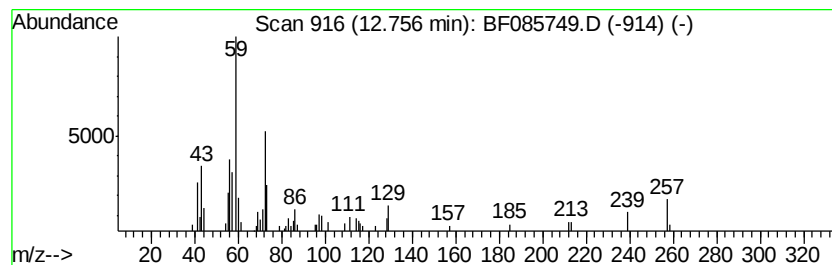
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown12.76 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	2.06 ng	61514	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanamide	199	C12H25NO	001120-16-7	47
2		Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	43
3		Tetradecanamide	227	C14H29NO	000638-58-4	43
4		Decanamide-	171	C10H21NO	002319-29-1	43
5		2-Methyl-2-decanol	172	C11H24O	003396-02-9	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085749.D
Acq On : 25 Mar 2016 8:29
Operator : UM/SJ
Sample : H1885-04
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TRACK-D-GRID-15B

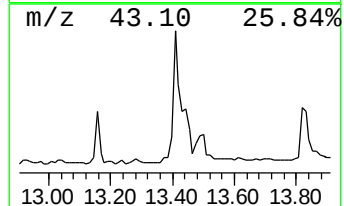
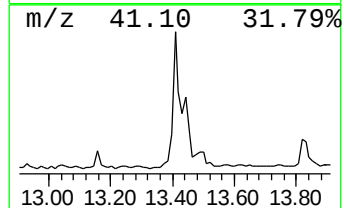
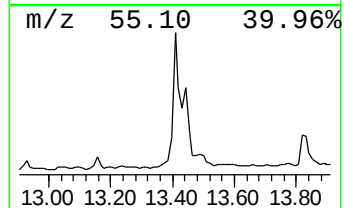
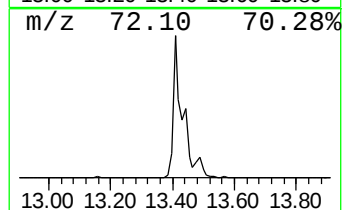
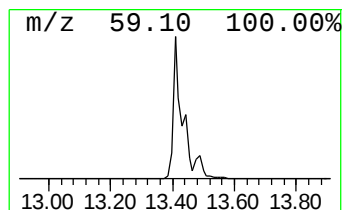
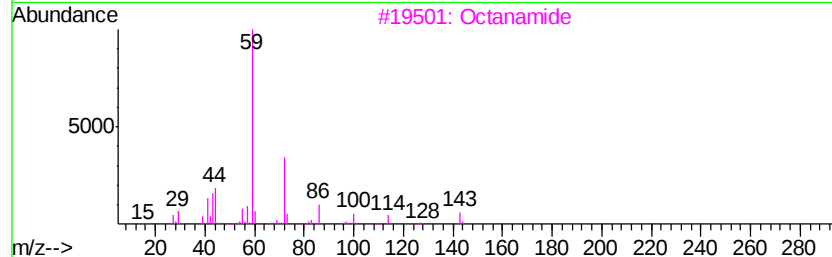
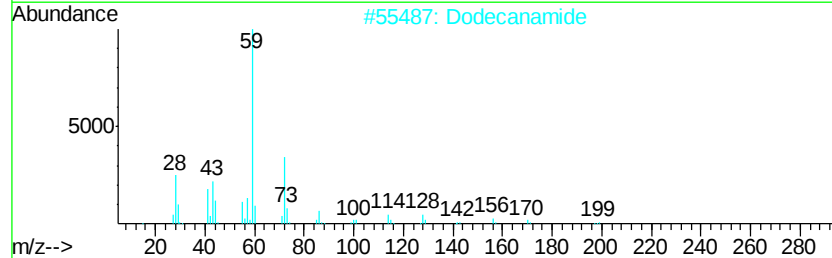
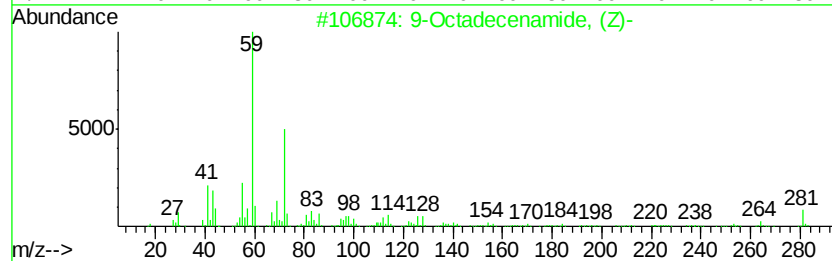
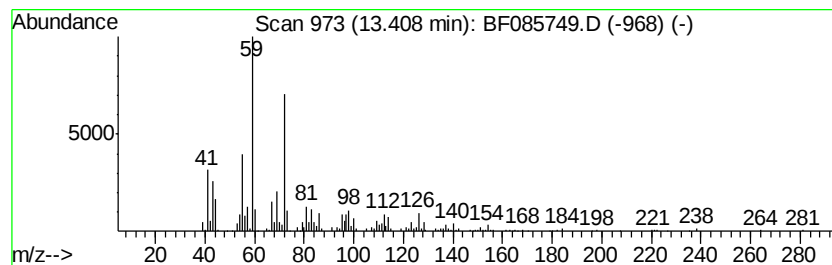
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	15.25 ng	455189	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		Dodecanamide	199	C12H25NO	001120-16-7	53
3		Octanamide	143	C8H17NO	000629-01-6	53
4		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	45
5		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	45



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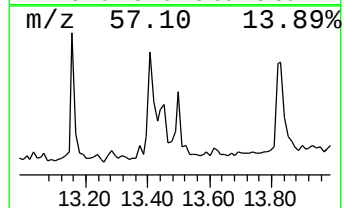
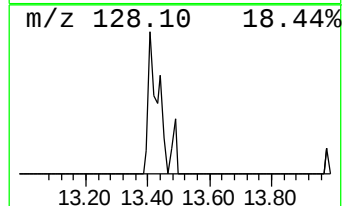
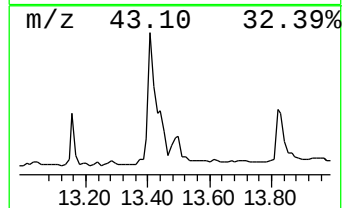
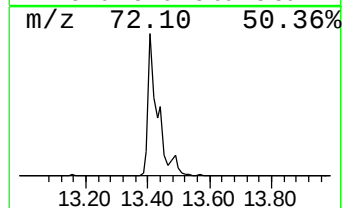
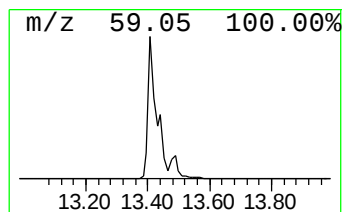
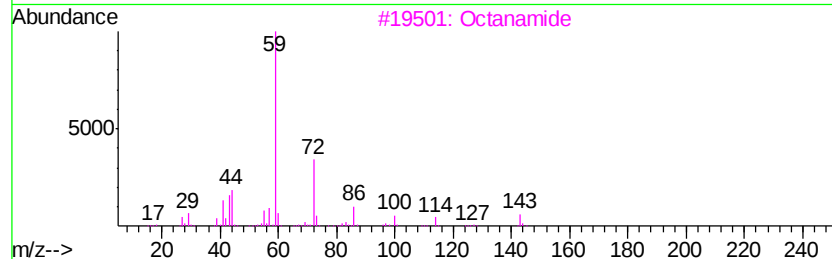
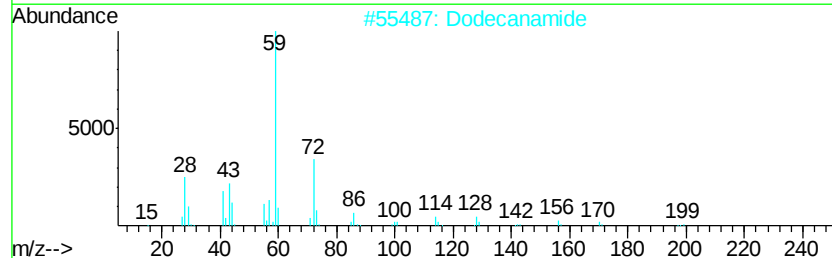
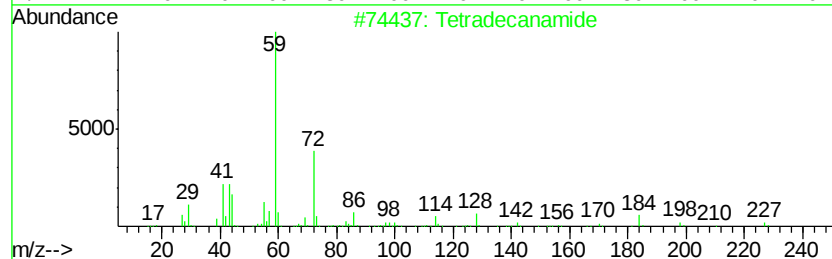
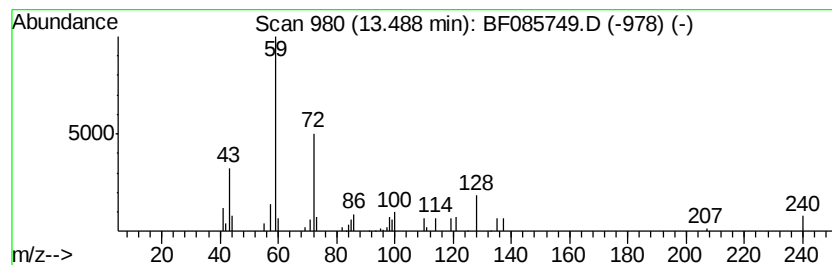
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Tetradecanamide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	2.43 ng	72459	Chrysene-d12	13.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecanamide	227 C14H29NO	000638-58-4 59
2		Dodecanamide	199 C12H25NO	001120-16-7 50
3		Octanamide	143 C8H17NO	000629-01-6 50
4		Hexadecanamide	255 C16H33NO	000629-54-9 45
5		Pentanamide, 5-hydroxy-	117 C5H11NO2	029686-12-2 45



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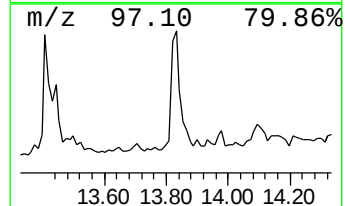
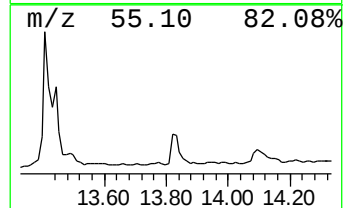
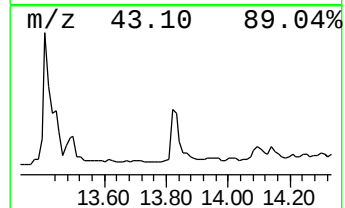
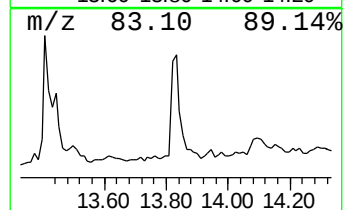
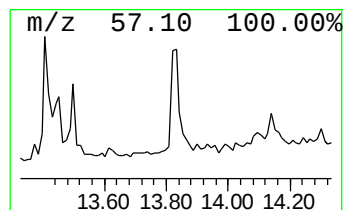
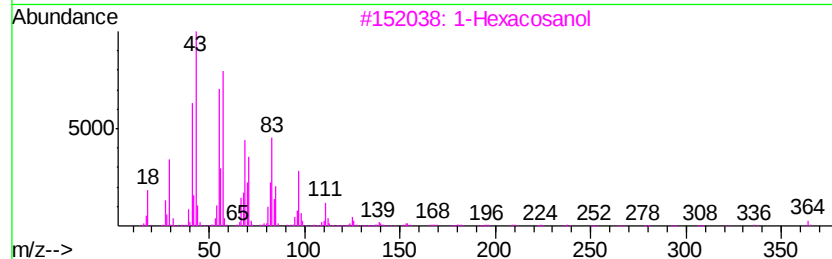
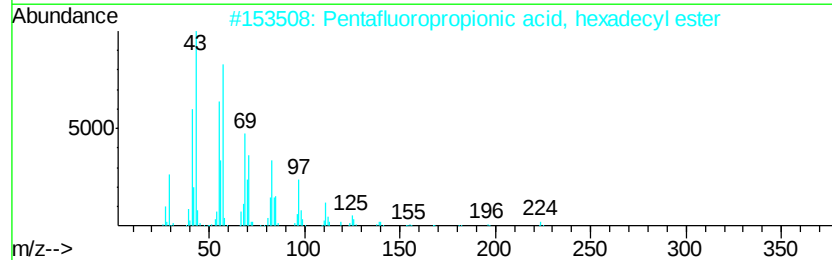
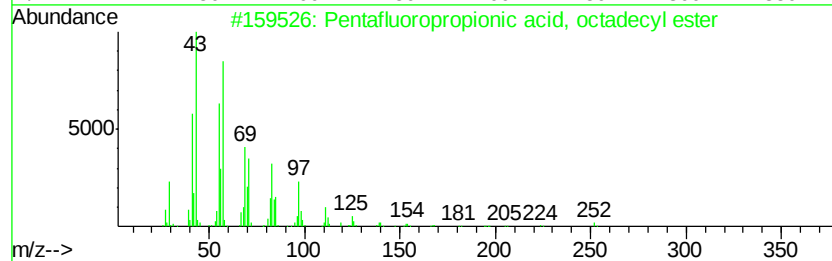
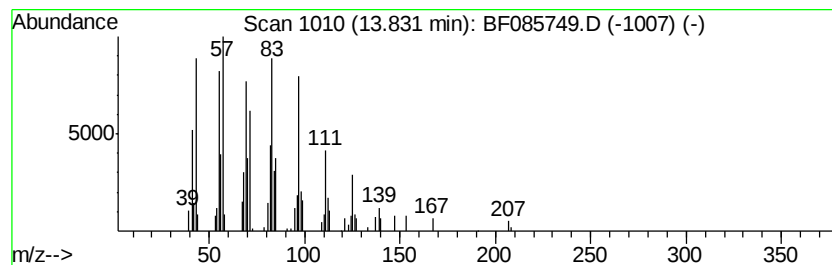
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentafluoropropionic acid, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	3.53 ng	105350	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, octadecyl...	416	C21H37F5O2	1000280-07-7	91
2		Pentafluoropropionic acid, hexadecyl...	388	C19H33F5O2	006222-07-7	91
3		1-Hexacosanol	382	C26H54O	000506-52-5	91
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



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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
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unknown6.60	6.60	81.8	ng	2746640	1	6.82	671549	20.0
Hexanoic acid, 2-...	7.50	2.6	ng	109703	2	8.10	838587	20.0
Hexadecane	9.80	2.3	ng	95608	3	9.86	850634	20.0
Octadecane	10.77	5.5	ng	185258	4	11.34	668494	20.0
Eicosane	11.21	2.3	ng	77276	4	11.34	668494	20.0
1-Hexadecanol	11.57	7.0	ng	232512	4	11.34	668494	20.0
unknown12.76	12.76	2.1	ng	61514	5	13.98	596960	20.0
9-Octadecenamide, ...	13.41	15.3	ng	455189	5	13.98	596960	20.0
Tetradecanamide	13.49	2.4	ng	72459	5	13.98	596960	20.0
Pentafluoropropio...	13.83	3.5	ng	105350	5	13.98	596960	20.0