

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
 Data File : BF085746.D
 Acq On : 25 Mar 2016 7:04
 Operator : UM/SJ
 Sample : H1885-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-GRID-14A

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	180	184	188	rBV2	20228	39137	1.23%	0.139%
2	5.007	235	238	244	rBV	301742	466371	14.60%	1.654%
3	5.430	272	275	277	rBV	1988236	2649865	82.98%	9.399%
4	5.830	307	310	313	rVB	41473	42944	1.34%	0.152%
5	6.459	362	365	368	rBV	1845789	2666763	83.51%	9.459%
6	6.596	374	377	379	rBV	2662747	2808988	87.96%	9.963%
7	6.824	395	397	399	rBV	849151	711953	22.29%	2.525%
8	6.984	408	411	413	rBV	2133785	2291719	71.76%	8.128%
9	7.396	444	447	449	rBV	1647049	1913684	59.93%	6.788%
10	7.499	453	456	462	rVB	56835	76966	2.41%	0.273%
11	8.104	507	509	512	rBV	708358	878017	27.49%	3.114%
12	9.190	601	604	606	rBV	3341874	3193389	100.00%	11.327%
13	9.579	636	638	641	rBV	527598	476479	14.92%	1.690%
14	9.796	655	657	659	rVB	109928	83573	2.62%	0.296%
15	9.865	661	663	665	rVB	1066165	946999	29.65%	3.359%
16	10.299	697	701	702	rBV	112424	138043	4.32%	0.490%
17	10.516	717	720	721	rBV	44545	59643	1.87%	0.212%
18	10.653	729	732	734	rBV	1709943	1662773	52.07%	5.898%
19	10.768	740	742	746	rVB	132856	150301	4.71%	0.533%
20	11.213	779	781	782	rBV	82068	69508	2.18%	0.247%
21	11.248	782	784	787	rVB	27974	34828	1.09%	0.124%
22	11.339	790	792	795	rBV	619829	828847	25.96%	2.940%
23	11.579	810	813	816	rBV3	20183	37279	1.17%	0.132%
24	11.636	816	818	820	rVB	36048	34353	1.08%	0.122%
25	11.876	837	839	844	rBV2	116988	123495	3.87%	0.438%
26	11.956	844	846	850	rVB	19929	36831	1.15%	0.131%
27	12.551	896	898	900	rBV	48490	55653	1.74%	0.197%
28	12.653	905	907	909	rBV	32334	47987	1.50%	0.170%
29	12.756	913	916	917	rVV2	20179	35232	1.10%	0.125%
30	12.779	917	918	922	rVB2	59859	80322	2.52%	0.285%
31	12.928	928	931	934	rBV	2357496	2144488	67.15%	7.606%
32	13.156	949	951	954	rBV2	30645	46927	1.47%	0.166%
33	13.408	971	973	975	rBV	176953	261112	8.18%	0.926%
34	13.499	979	981	983	rVB2	66899	77637	2.43%	0.275%

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

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Peak separation: 5

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35	13.819	1007	1009	1015	rBV	328413	562262	17.61%	1.994%
36	13.979	1020	1023	1025	rBV	664637	684559	21.44%	2.428%
37	14.139	1036	1037	1041	rVB	99273	138726	4.34%	0.492%
38	14.448	1062	1064	1066	rBV	151417	164423	5.15%	0.583%
39	14.745	1088	1090	1092	rVB	102303	84442	2.64%	0.300%
40	15.065	1116	1118	1120	rVB	82540	95864	3.00%	0.340%
41	15.397	1145	1147	1152	rVB	425681	622516	19.49%	2.208%
42	16.048	1202	1204	1215	rVB	141820	340119	10.65%	1.206%
43	16.460	1238	1240	1249	rVB	110251	225166	7.05%	0.799%
44	16.997	1285	1287	1291	rVB2	56427	103447	3.24%	0.367%

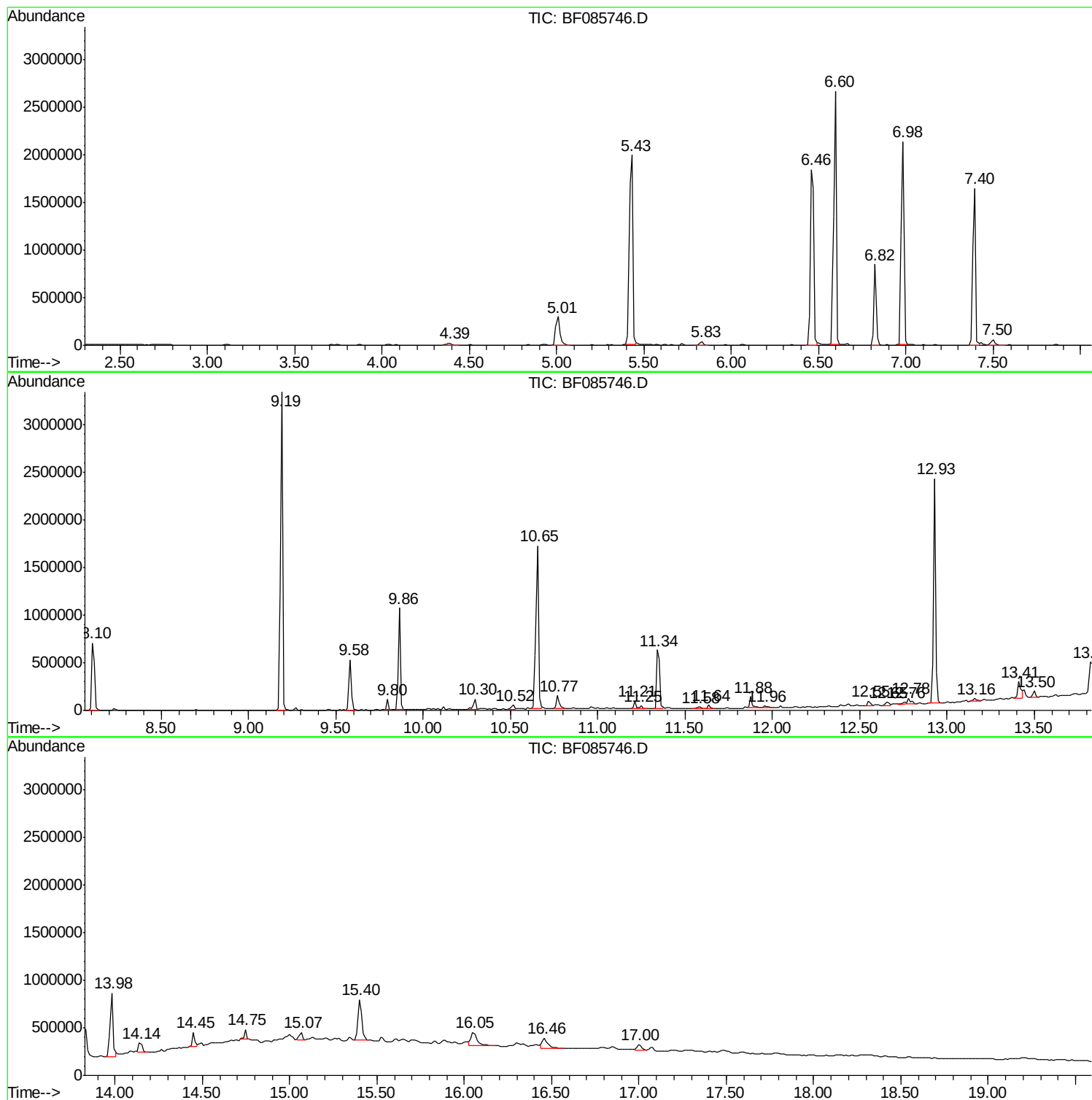
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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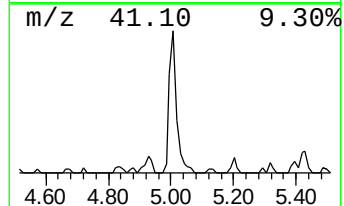
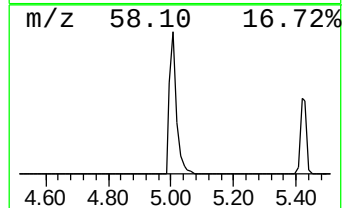
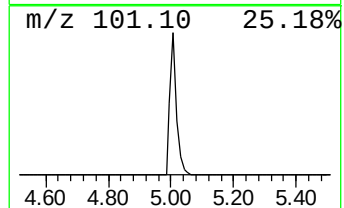
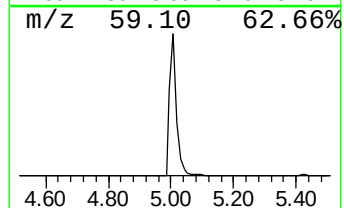
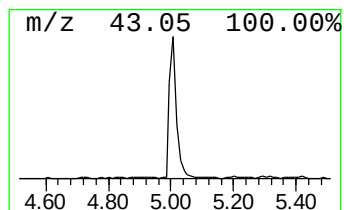
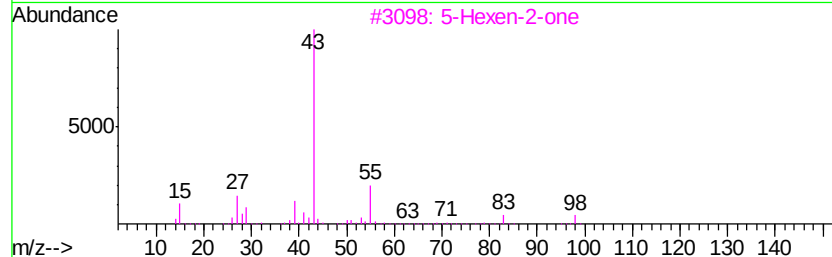
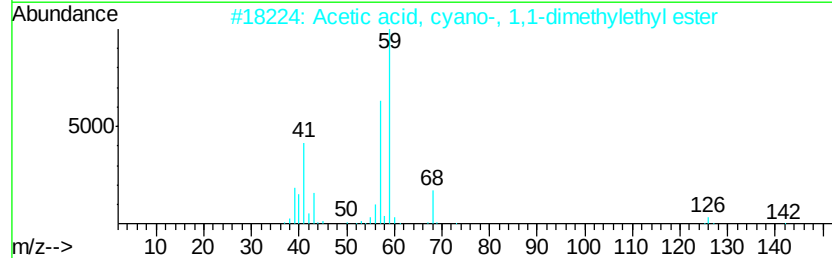
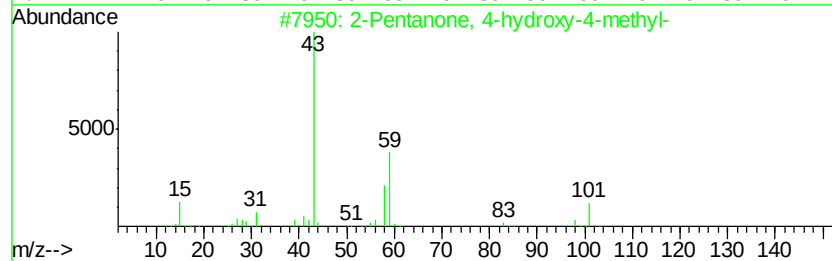
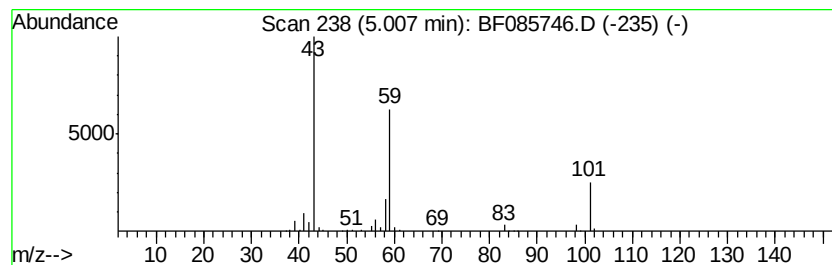
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	13.10 ng	466371	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, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			5-Hexen-2-one	98	C6H10O	000109-49-9	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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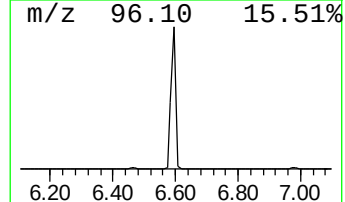
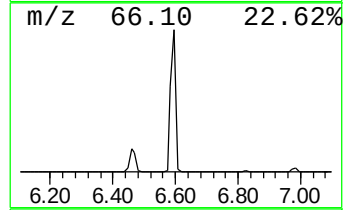
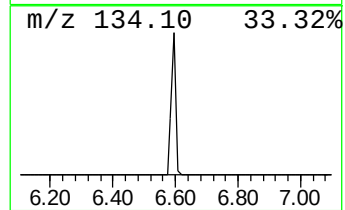
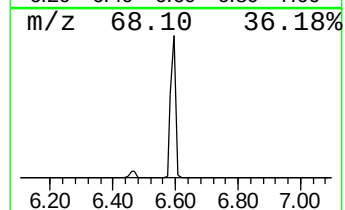
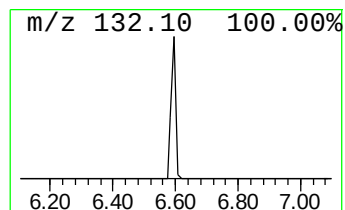
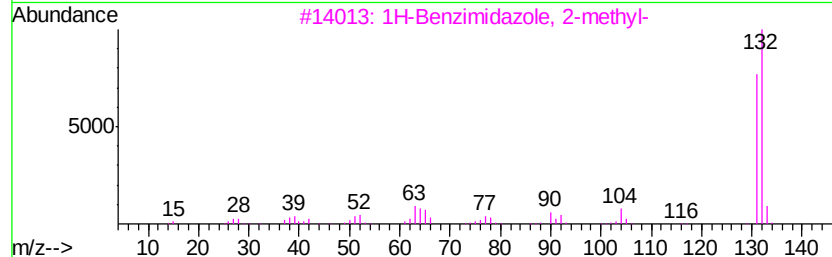
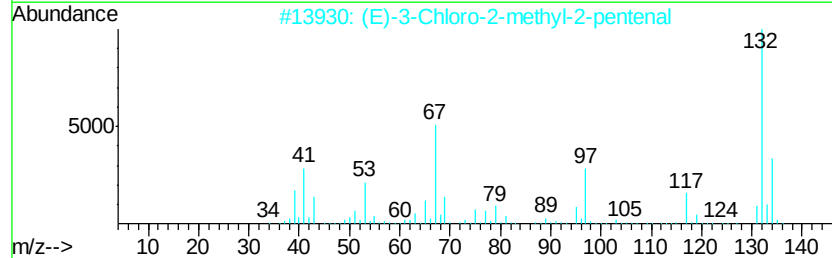
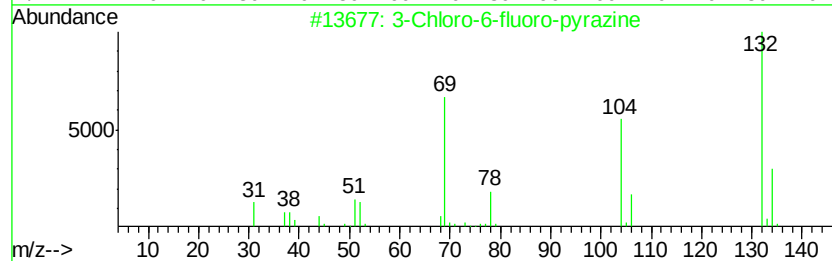
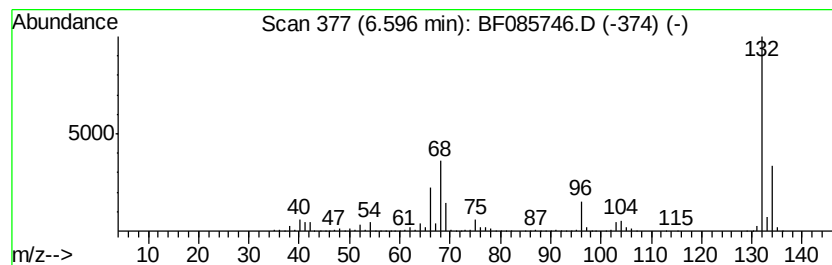
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Peak Number 2 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	78.91 ng	2808990	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	(E)-3	Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	1H	Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4	Benzene, 1,3,5	trifluoro-	132	C6H3F3	000372-38-3	9
5	4	Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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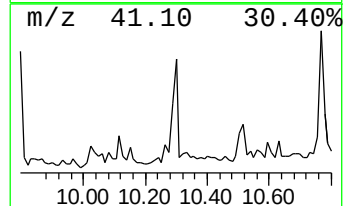
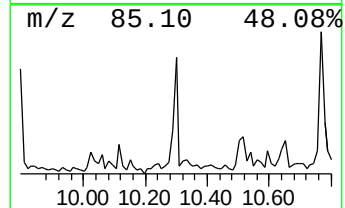
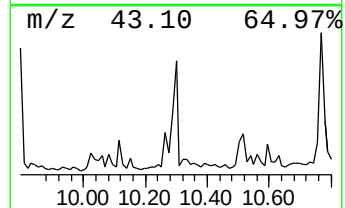
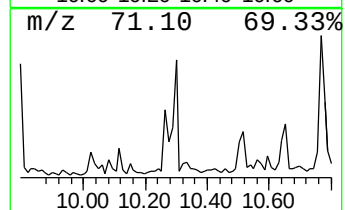
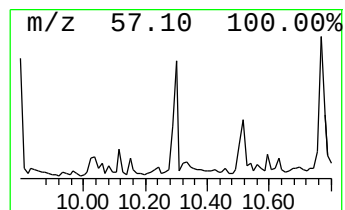
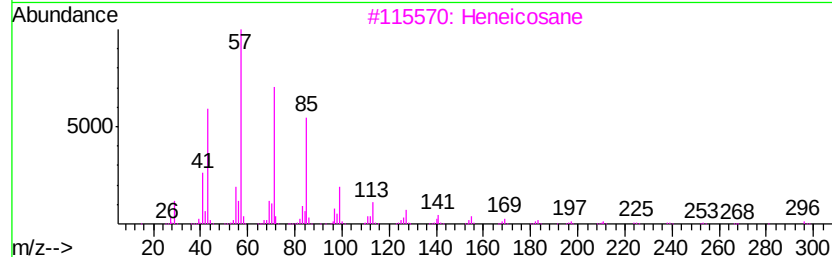
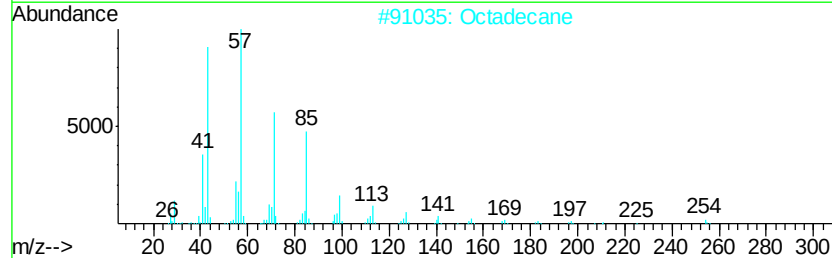
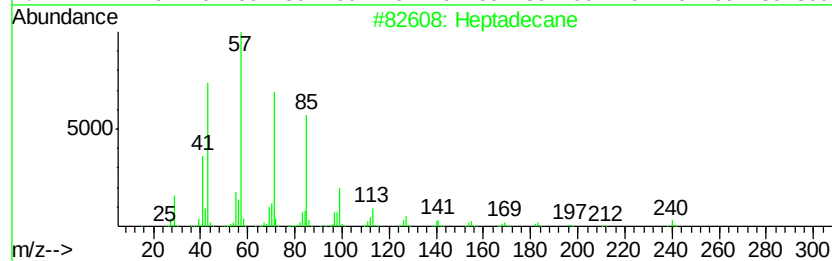
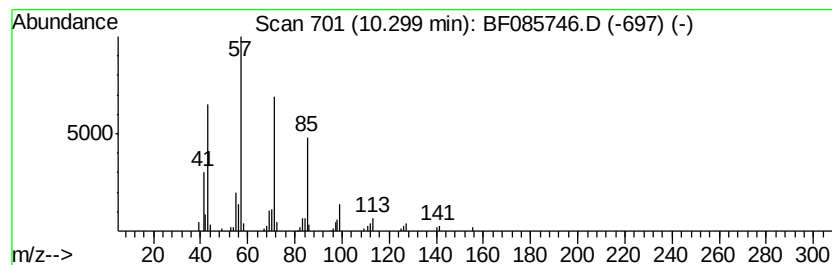
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	2.92 ng	138043	Acenaphthene-d10	9.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	91
2	Octadecane	254 C18H38	000593-45-3	90
3	Heneicosane	296 C21H44	000629-94-7	90
4	Hexadecane	226 C16H34	000544-76-3	90
5	Nonadecane	268 C19H40	000629-92-5	87



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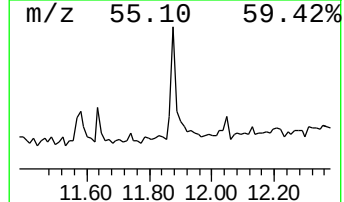
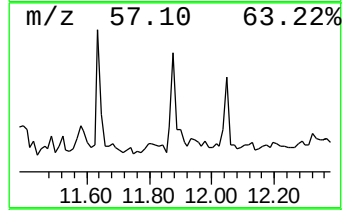
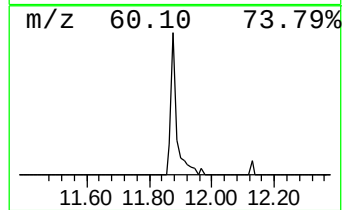
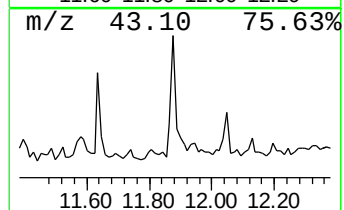
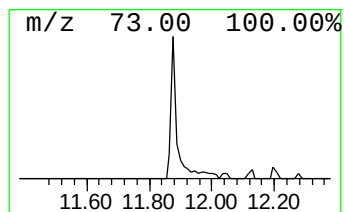
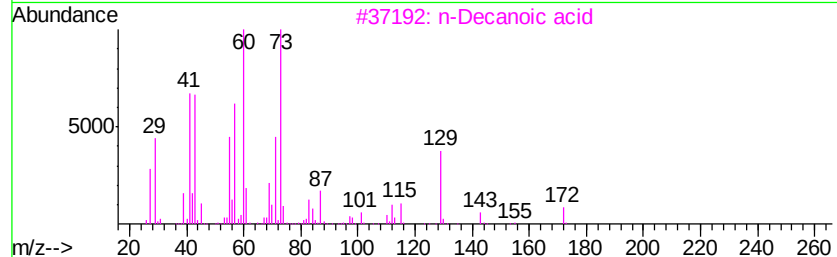
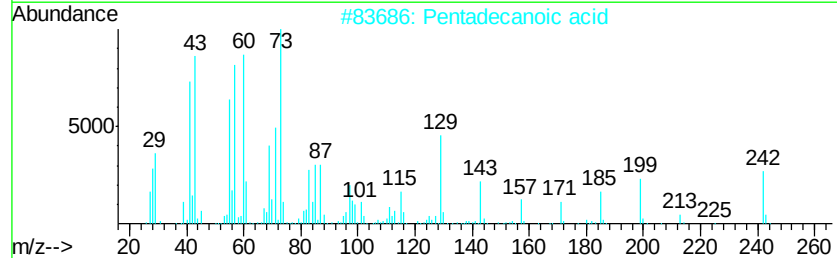
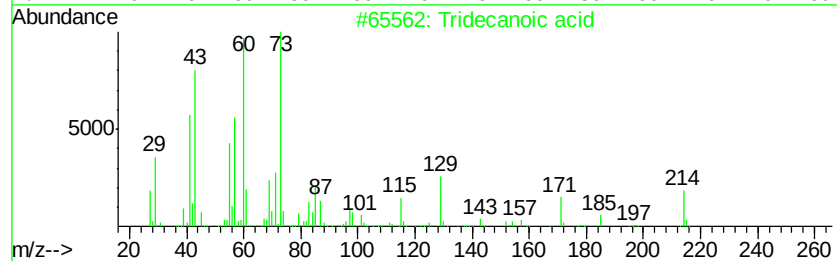
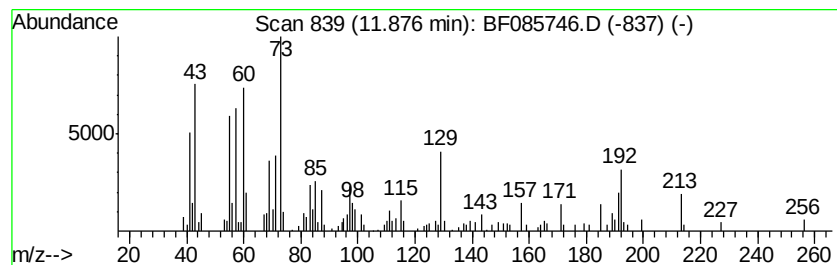
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tridecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.88	2.98 ng	123495	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	76
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
3	n-Decanoic acid	172	C10H20O2	000334-48-5	64
4	Dodecane	170	C12H26	000112-40-3	11
5	Tritetracontane	605	C43H88	007098-21-7	10



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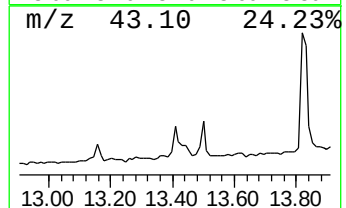
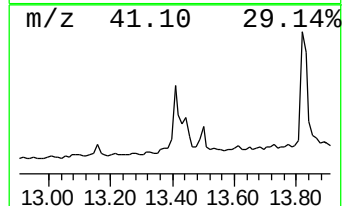
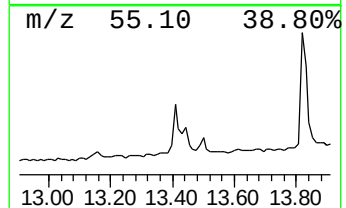
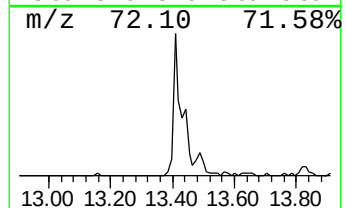
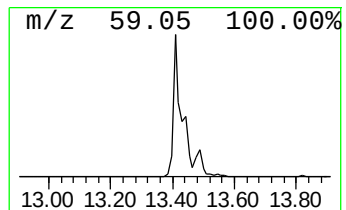
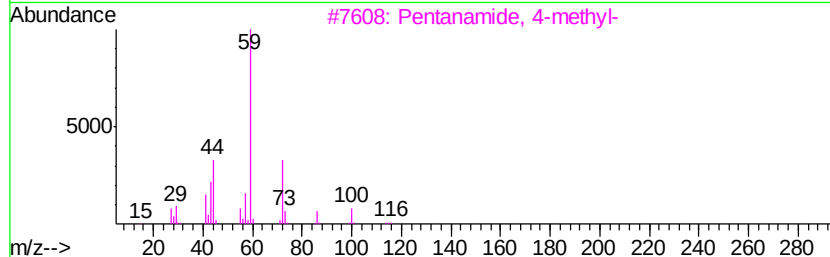
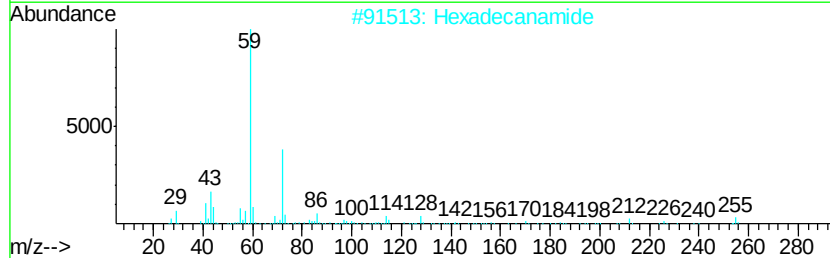
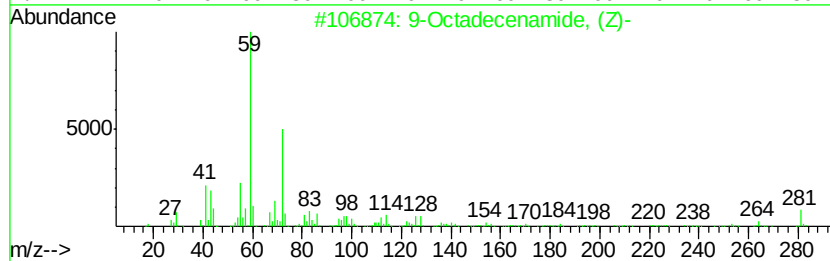
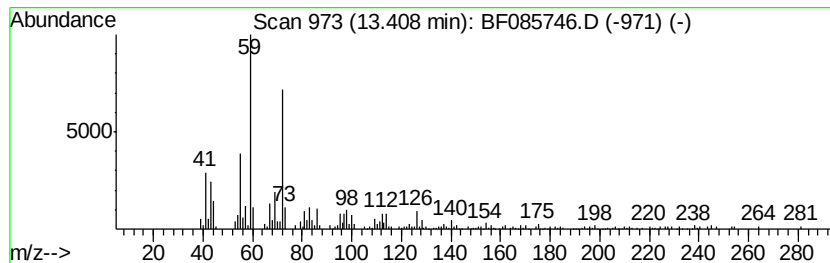
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TIC Integration Parameters: LSCINT.P

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.41	7.63 ng	261112	Chrysene-d12	13.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
2	Hexadecanamide	255	C16H33NO	000629-54-9	72
3	Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	64
4	Dodecanamide	199	C12H25NO	001120-16-7	59
5	Tetradecanamide	227	C14H29NO	000638-58-4	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
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Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
Sample : H1885-01
Misc :
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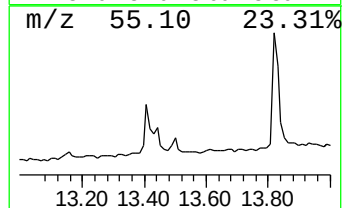
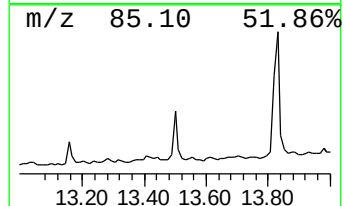
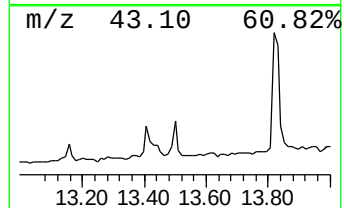
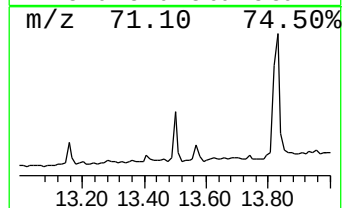
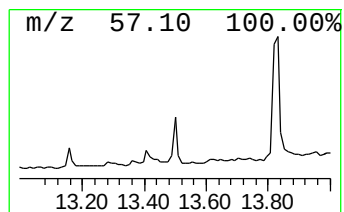
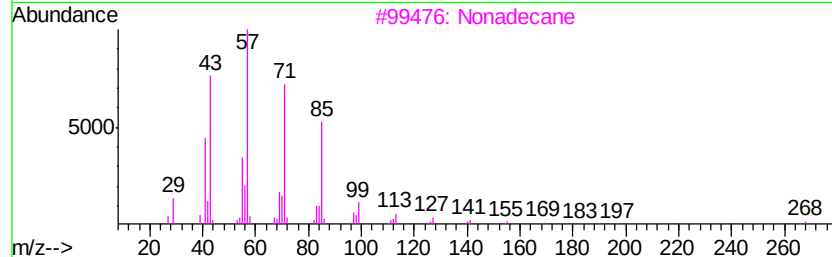
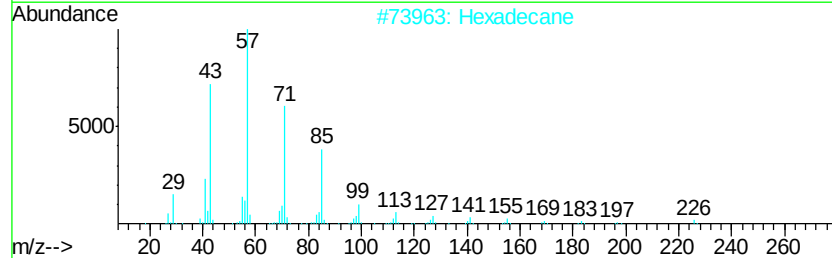
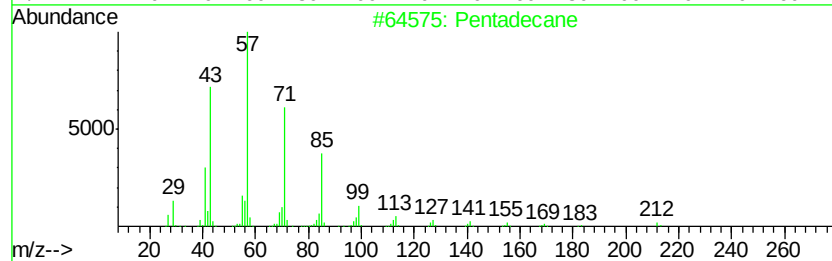
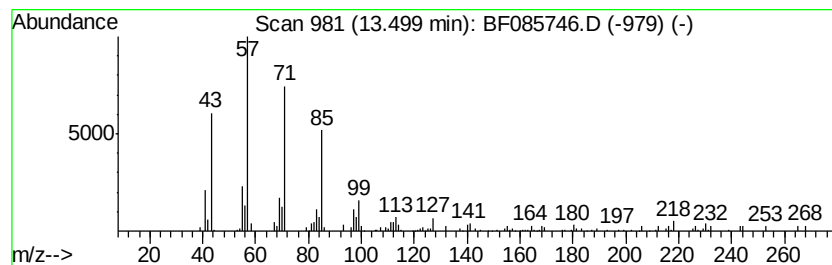
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.50	2.27 ng	77637	Chrysene-d12	13.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane	212 C15H32	000629-62-9	95
2	Hexadecane	226 C16H34	000544-76-3	94
3	Nonadecane	268 C19H40	000629-92-5	93
4	Octadecane, 2-methyl-	268 C19H40	001560-88-9	87
5	Pentadecane, 8-heptyl-	310 C22H46	071005-15-7	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085746.D
Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
Sample : H1885-01
Misc :
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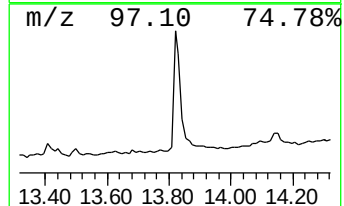
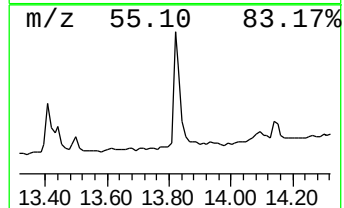
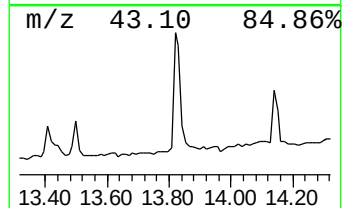
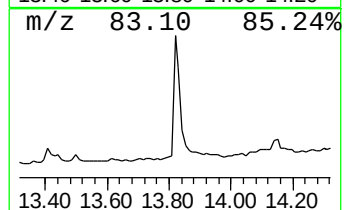
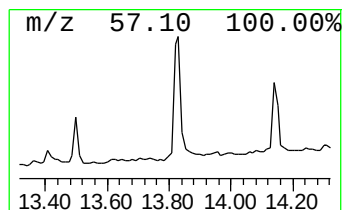
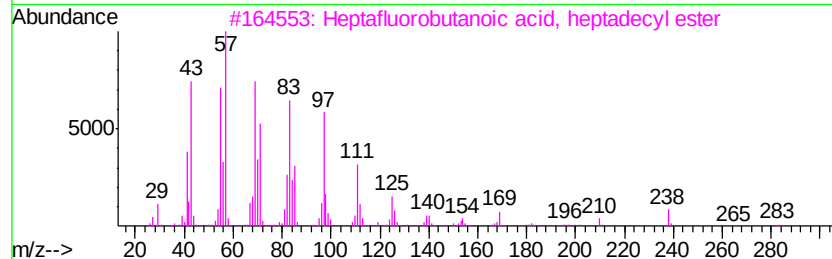
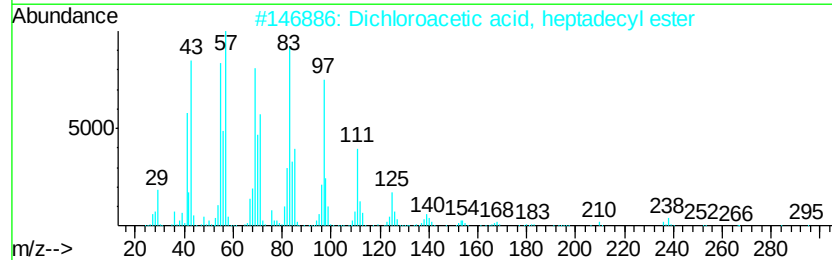
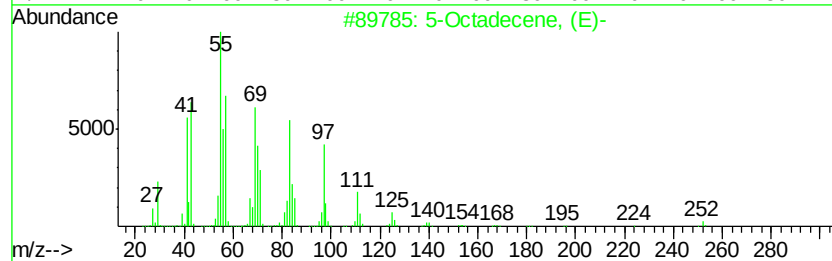
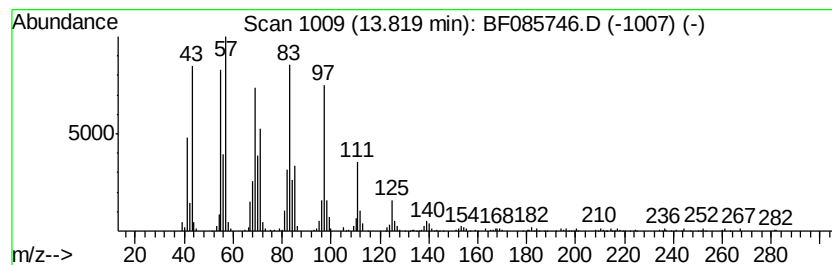
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 5-Octadecene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.82	16.43 ng	562262	Chrysene-d12	13.98		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Octadecene, (E)-	252	C18H36	007206-21-5	96
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085746.D
Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
Sample : H1885-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

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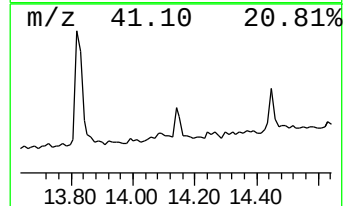
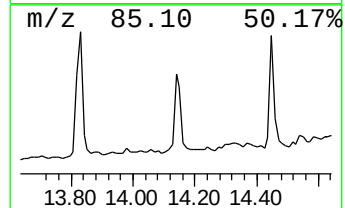
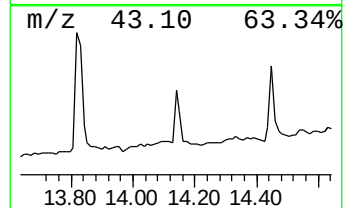
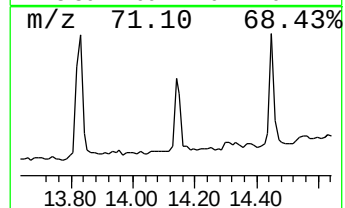
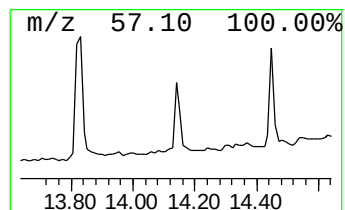
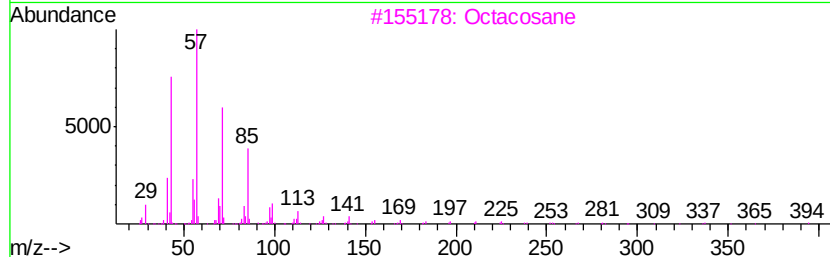
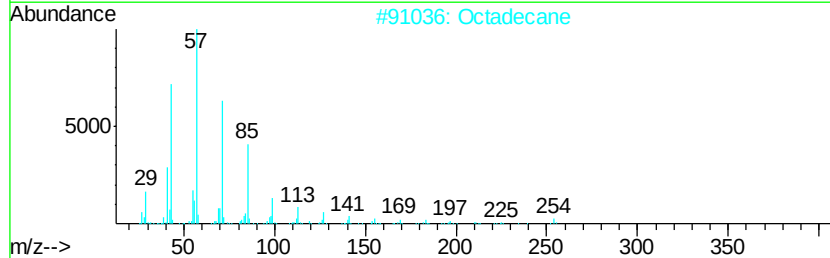
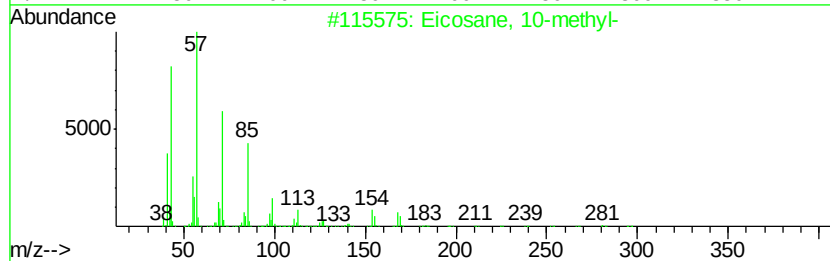
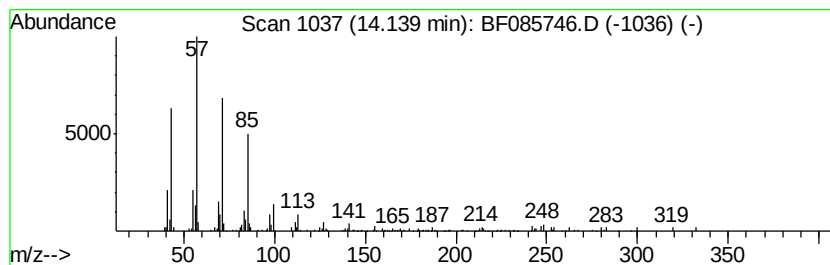
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane, 10-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	4.05 ng	138726	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
2		Octadecane	254	C18H38	000593-45-3	86
3		Octacosane	394	C28H58	000630-02-4	83
4		Heneicosane	296	C21H44	000629-94-7	83
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085746.D
Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
Sample : H1885-01
Misc :
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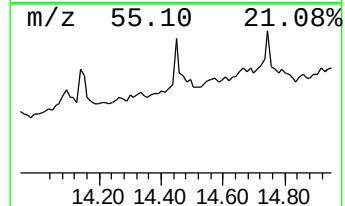
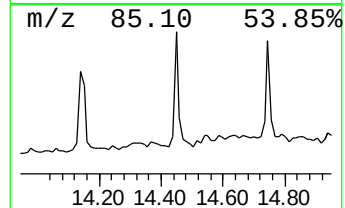
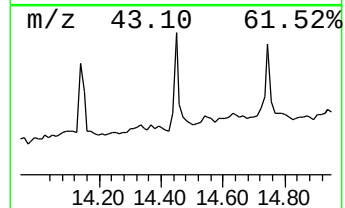
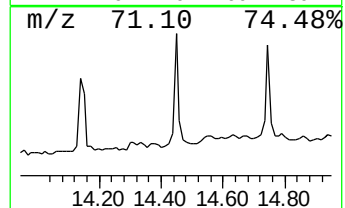
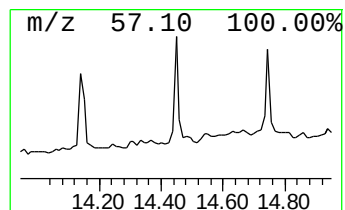
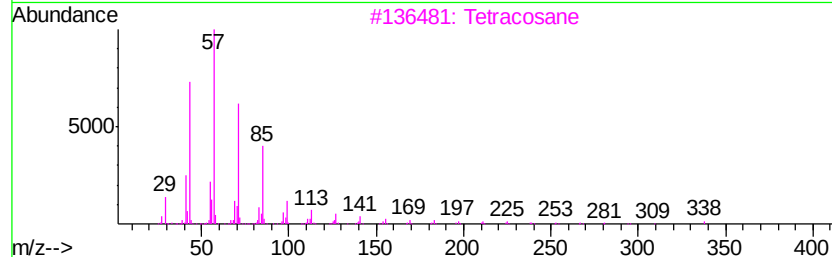
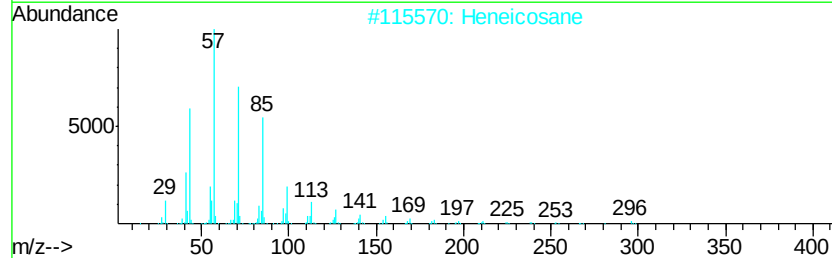
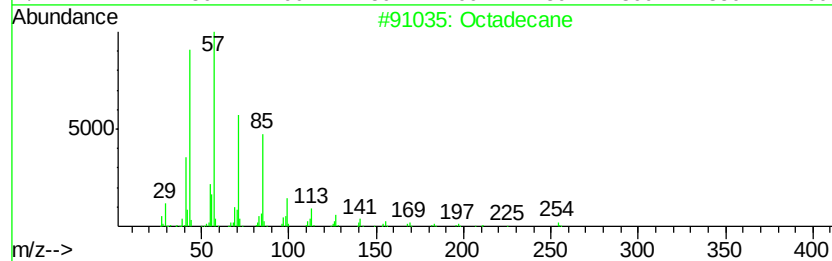
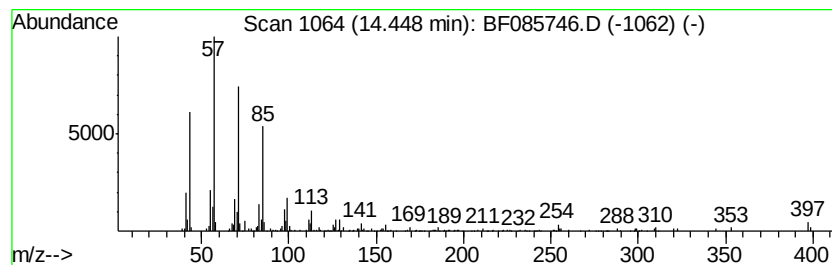
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.45	4.80 ng	164423	Chrysene-d12		13.98
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	93
2	Heneicosane	296	C21H44	000629-94-7	90
3	Tetracosane	338	C24H50	000646-31-1	87
4	Docosane	310	C22H46	000629-97-0	87
5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
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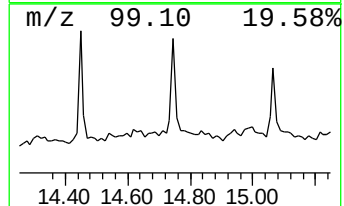
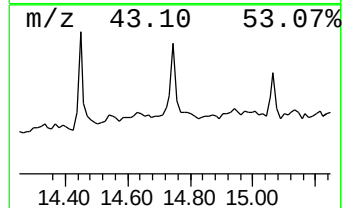
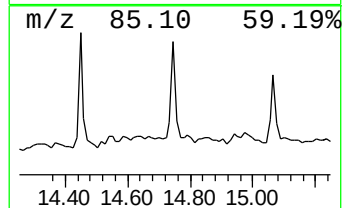
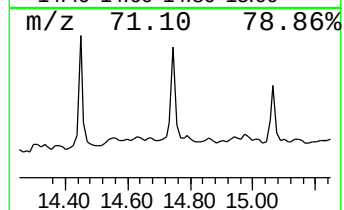
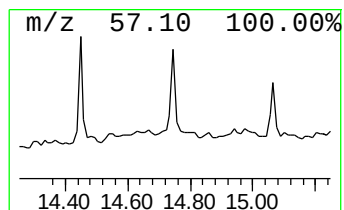
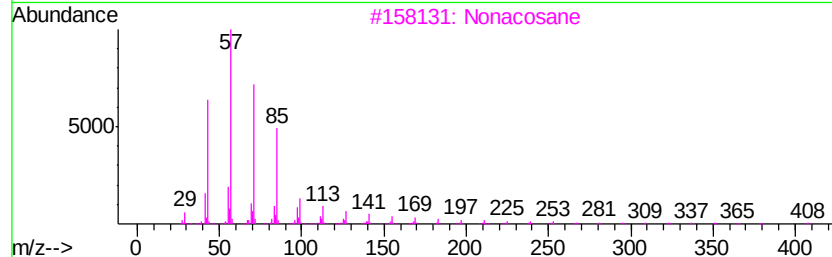
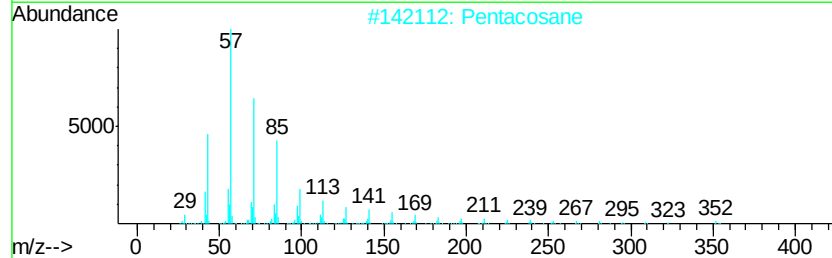
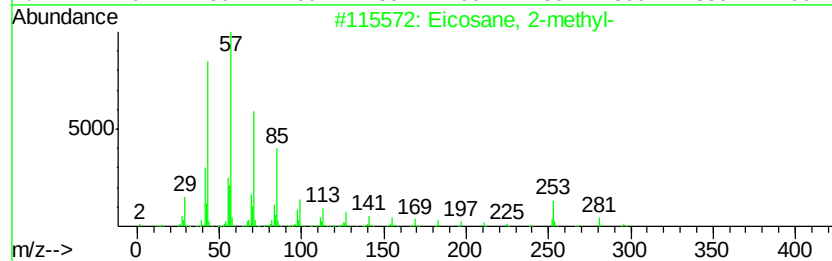
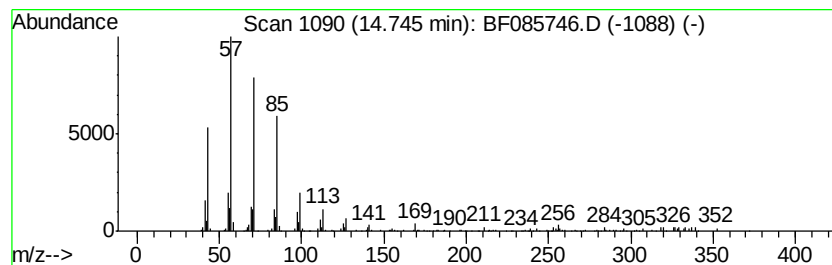
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.75	2.71 ng	84442	Perylene-d12		15.40
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 2-methyl-	296	C21H44	001560-84-5	94
2	Pentacosane	352	C25H52	000629-99-2	91
3	Nonacosane	408	C29H60	000630-03-5	91
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085746.D
Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
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Misc :
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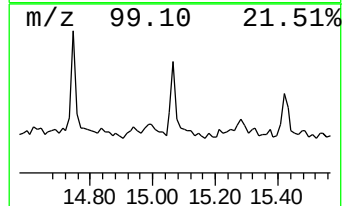
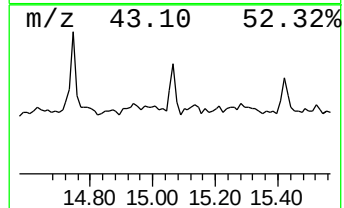
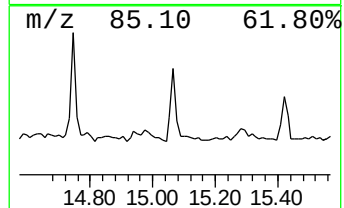
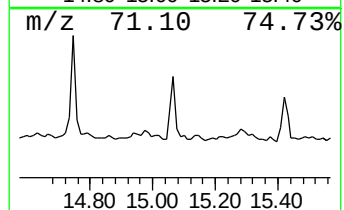
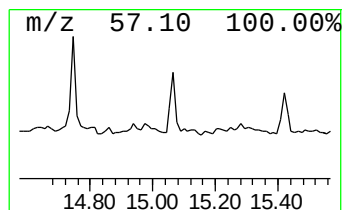
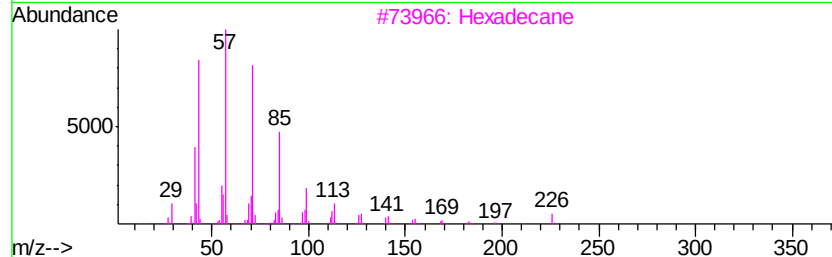
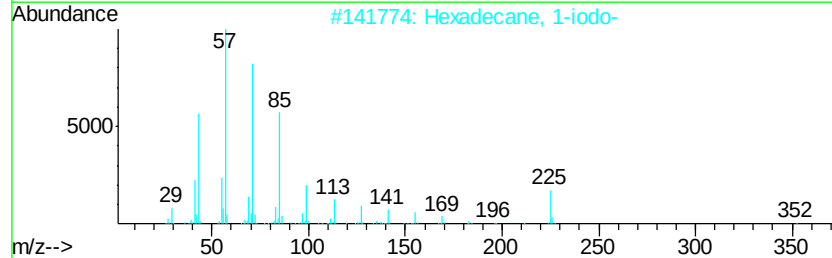
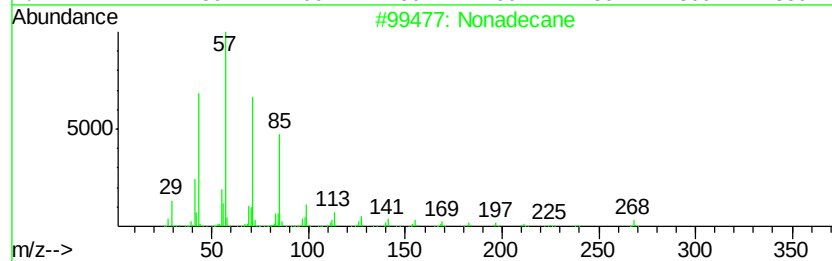
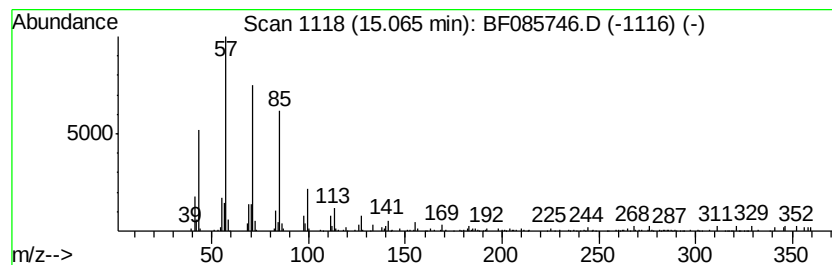
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	3.08 ng	95864	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	93
2		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
3		Hexadecane	226	C16H34	000544-76-3	86
4		triacontane	422	C30H62	000638-68-6	86
5		Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
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Acq On : 25 Mar 2016 7:04
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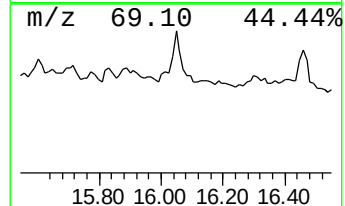
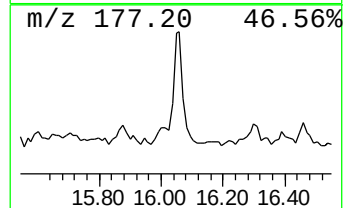
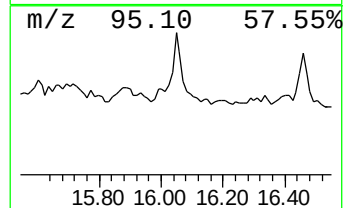
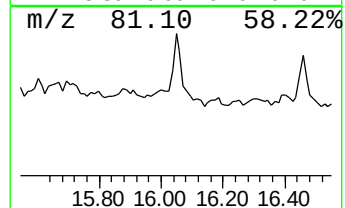
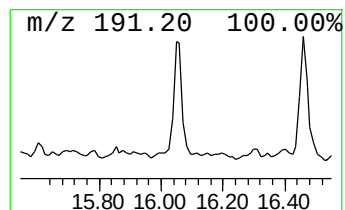
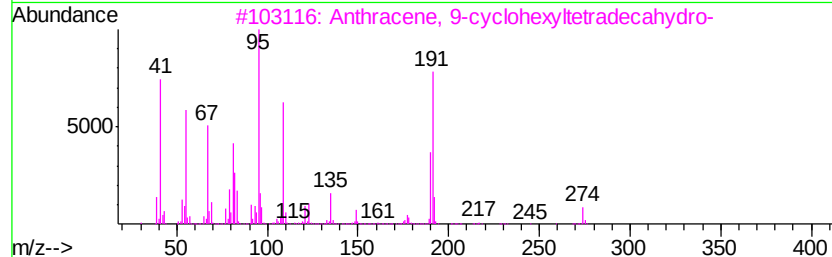
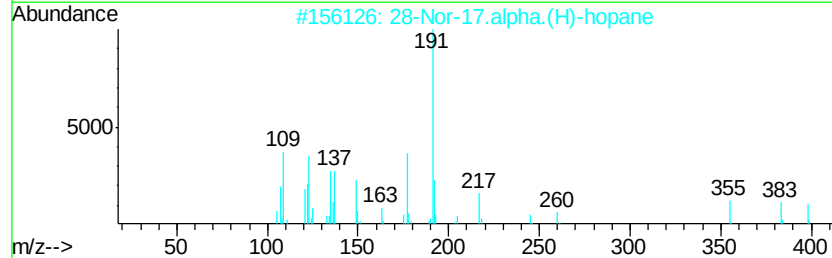
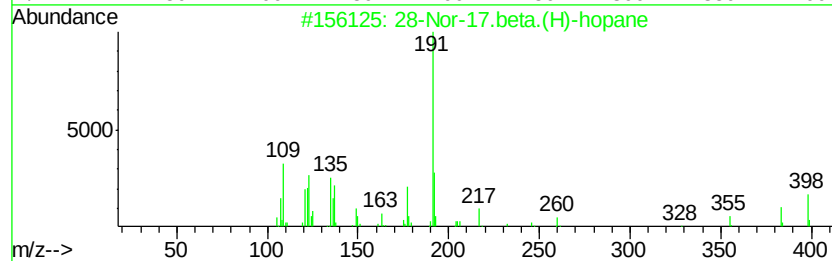
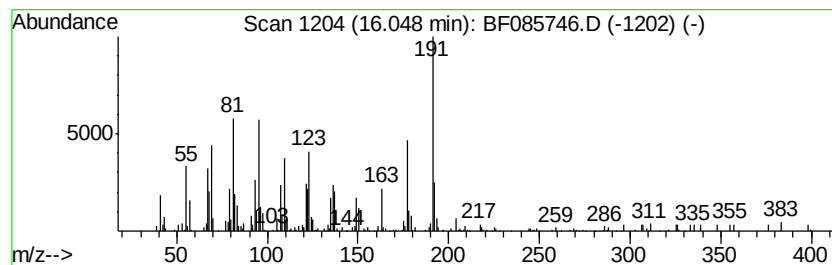
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 28-Nor-17.beta.(H)-hopane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	10.93 ng	340119	Perylene-d12	15.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	28-Nor-17.beta.(H)-hopane	398 C29H50	036728-72-0	68
2	28-Nor-17.alpha.(H)-hopane	398 C29H50	053584-60-4	64
3	Anthracene, 9-cyclohexyltetradec...	274 C20H34	055255-70-4	38
4	Anthracene, 9-dodecyltetradeca...	360 C26H48	055401-75-7	32
5	Acetamide, N-benzyl-2-[2,3-dihyd...	339 C19H21N3OS	1000272-50-7	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085746.D
Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
Sample : H1885-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-14A

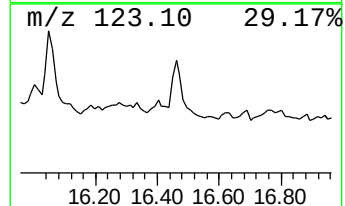
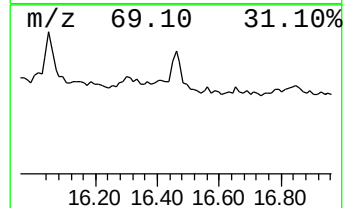
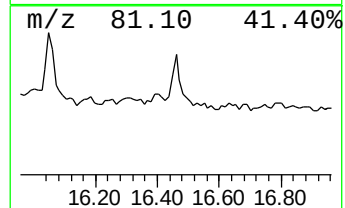
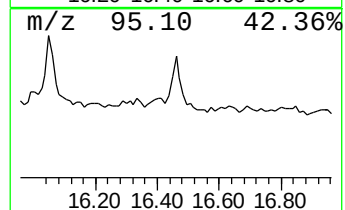
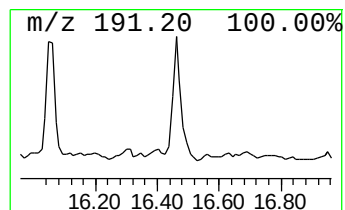
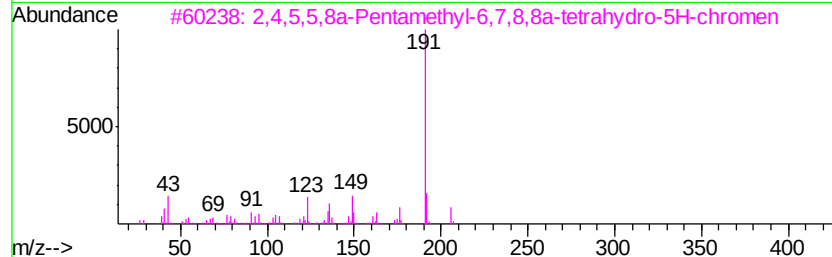
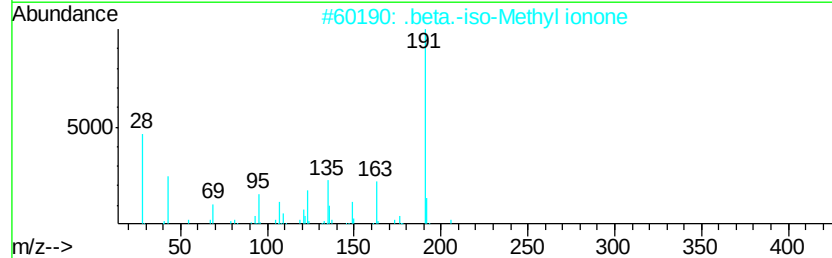
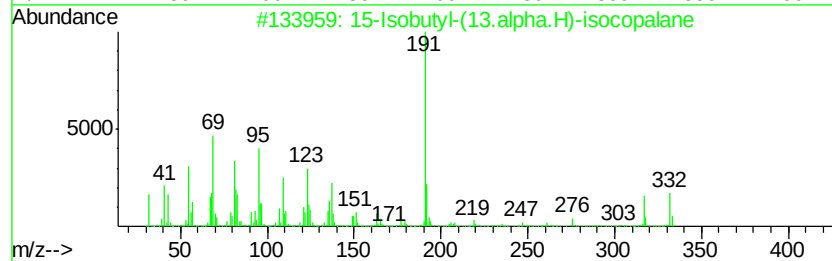
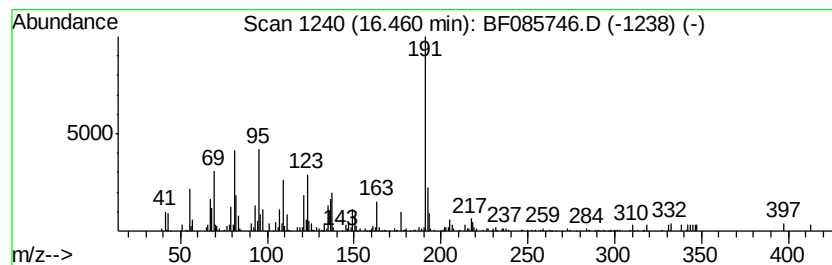
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 15-Isobutyl-(13.alpha.H)-is... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	7.23 ng	225166	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	15-Isobutyl-(13.alpha.H)-isocopa...	332	C24H44	087953-47-7	72
2		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	58
3		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	43
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35
5		4H-1,2,4-Triazole, 3-ethyl-5-(3,...	191	C9H13N5	328532-29-2	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085746.D
Acq On : 25 Mar 2016 7:04
Operator : UM/SJ
Sample : H1885-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-14A

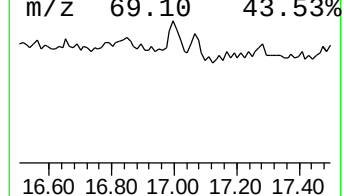
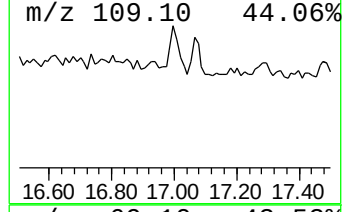
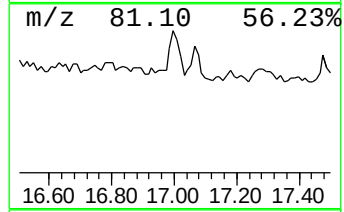
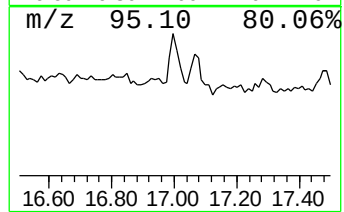
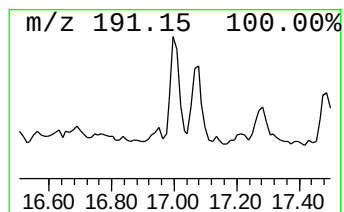
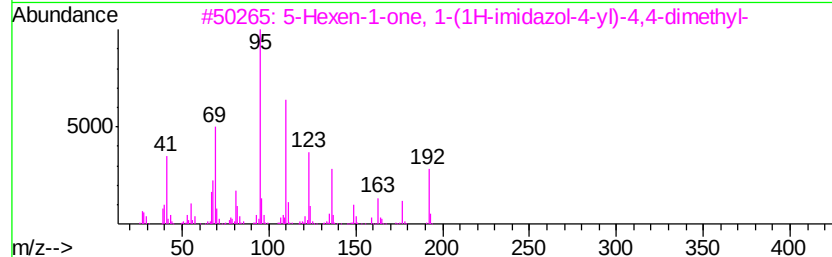
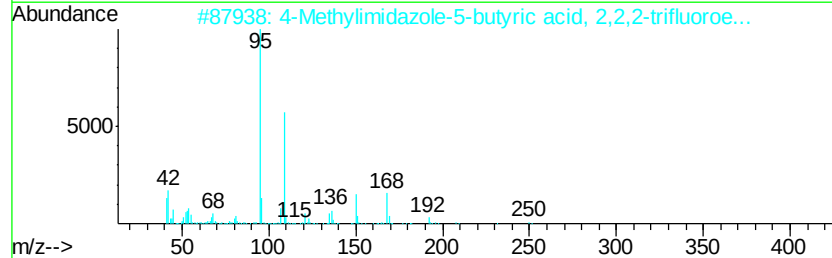
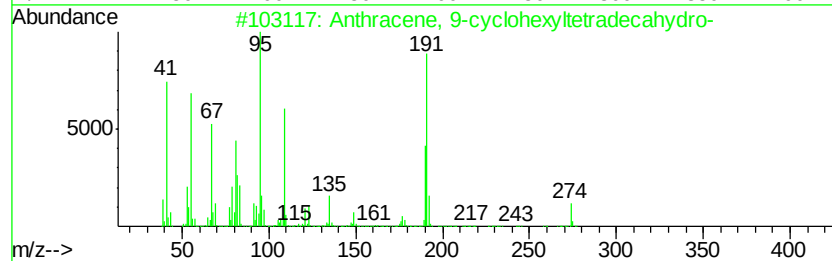
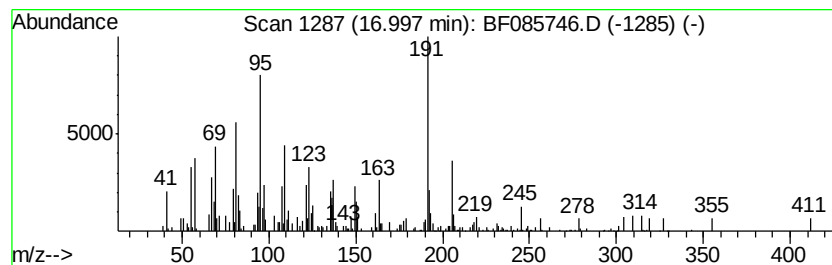
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown17.00 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	3.32 ng	103447	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	43
2		4-Methylimidazole-5-butyric acid...	250	C10H13F3N2O2	1000129-15-9	30
3		5-Hexen-1-one, 1-(1H-imidazol-4-...	192	C11H16N2O	069393-41-5	25
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	22
5		2-(2-((2-Chloro-6-fluorobenzyl)s...	334	C17H16ClFN2S	1000298-18-2	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
 Data File : BF085746.D
 Acq On : 25 Mar 2016 7:04
 Operator : UM/SJ
 Sample : H1885-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-GRID-14A

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	13.1	ng	466371	1	6.82	711953	20.0
unknown6.60	6.60	78.9	ng	2808990	1	6.82	711953	20.0
Heptadecane	10.30	2.9	ng	138043	3	9.86	946999	20.0
Tridecanoic acid	11.88	3.0	ng	123495	4	11.34	828847	20.0
9-Octadecenamide,...	13.41	7.6	ng	261112	5	13.98	684559	20.0
Pentadecane	13.50	2.3	ng	77637	5	13.98	684559	20.0
5-Octadecene, (E)-	13.82	16.4	ng	562262	5	13.98	684559	20.0
Eicosane, 10-methyl-	14.14	4.0	ng	138726	5	13.98	684559	20.0
Octadecane	14.45	4.8	ng	164423	5	13.98	684559	20.0
Eicosane, 2-methyl-	14.75	2.7	ng	84442	6	15.40	622516	20.0
Nonadecane	15.07	3.1	ng	95864	6	15.40	622516	20.0
28-Nor-17.beta.(H...	16.05	10.9	ng	340119	6	15.40	622516	20.0
15-Isobutyl-(13.a...	16.46	7.2	ng	225166	6	15.40	622516	20.0
unknown17.00	17.00	3.3	ng	103447	6	15.40	622516	20.0