

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
 Data File : BF085686.D
 Acq On : 23 Mar 2016 5:04
 Operator : UM/SJ
 Sample : H1886-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-6

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	5.018	236	239	247	rBV	764490	1251456	29.95%	3.138%
2	5.441	273	276	278	rBV	3116599	3814961	91.30%	9.564%
3	5.830	309	310	315	rVB	80994	121395	2.91%	0.304%
4	6.470	363	366	369	rBV	2627063	3722168	89.08%	9.332%
5	6.607	375	378	380	rBV	3149032	4089701	97.88%	10.253%
6	6.836	395	398	400	rBV	651298	811987	19.43%	2.036%
7	6.984	409	411	414	rBV	2400027	2978627	71.29%	7.468%
8	7.396	444	447	449	rBV	1905921	2781520	66.57%	6.974%
9	7.853	484	487	491	rBV	38861	54815	1.31%	0.137%
10	8.116	507	510	512	rBV	1355213	1179125	28.22%	2.956%
11	9.190	601	604	607	rBV	2946008	4178386	100.00%	10.476%
12	9.270	609	611	613	rVB	42296	45754	1.10%	0.115%
13	9.533	629	634	637	rBV2	16898	45998	1.10%	0.115%
14	9.590	637	639	640	rBV	723063	686190	16.42%	1.720%
15	9.807	655	658	660	rVB	125481	145065	3.47%	0.364%
16	9.865	661	663	666	rVB2	1110581	1175770	28.14%	2.948%
17	10.036	674	678	679	rBV	36918	51949	1.24%	0.130%
18	10.128	684	686	687	rBV	41148	46000	1.10%	0.115%
19	10.299	699	701	703	rVB	234090	202246	4.84%	0.507%
20	10.516	717	720	724	rBV	74158	136018	3.26%	0.341%
21	10.665	729	733	735	rBV2	2144387	2867412	68.62%	7.189%
22	10.768	741	742	749	rVB	162258	283878	6.79%	0.712%
23	10.962	757	759	763	rBV2	32231	83901	2.01%	0.210%
24	11.225	779	782	783	rBV	86863	102059	2.44%	0.256%
25	11.248	783	784	788	rVB	39467	42713	1.02%	0.107%
26	11.350	791	793	794	rBV	1373896	1172495	28.06%	2.940%
27	11.373	794	795	796	rVB	124588	86229	2.06%	0.216%
28	11.522	804	808	810	rVB3	18499	46993	1.12%	0.118%
29	11.648	816	819	821	rBV	40980	46371	1.11%	0.116%
30	11.876	838	839	842	rVV	145458	185591	4.44%	0.465%
31	11.956	844	846	850	rVV	35257	69722	1.67%	0.175%
32	12.562	896	899	901	rBV	187918	236401	5.66%	0.593%
33	12.665	906	908	910	rBV2	93170	134332	3.21%	0.337%
34	12.791	915	919	920	rVV	143878	246130	5.89%	0.617%

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Sampling : 1 Min Area: 3 % of largest Peak
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35	12.836	920	923	927	rVB3	67089	103433	2.48%	0.259%
36	12.939	929	932	934	rVV	3094102	3412876	81.68%	8.556%
37	13.785	1004	1006	1008	rBV	252420	227057	5.43%	0.569%
38	13.831	1008	1010	1015	rVV3	111177	167226	4.00%	0.419%
39	13.979	1020	1023	1029	rVV	801375	972482	23.27%	2.438%
40	14.288	1048	1050	1053	rBV2	40814	64853	1.55%	0.163%
41	14.345	1053	1055	1059	rVB3	90140	150033	3.59%	0.376%
42	14.414	1059	1061	1062	rBV2	63078	72548	1.74%	0.182%
43	14.459	1062	1065	1067	rBV2	100142	141650	3.39%	0.355%
44	14.997	1110	1112	1116	rBV	108937	202189	4.84%	0.507%
45	15.282	1135	1137	1140	rVB	83963	103158	2.47%	0.259%
46	15.339	1140	1142	1145	rVV	88412	122292	2.93%	0.307%
47	15.397	1145	1147	1157	rVB	509222	777841	18.62%	1.950%
48	16.780	1265	1268	1273	rVB2	45985	122055	2.92%	0.306%
49	17.191	1301	1304	1310	rVB	53751	123690	2.96%	0.310%

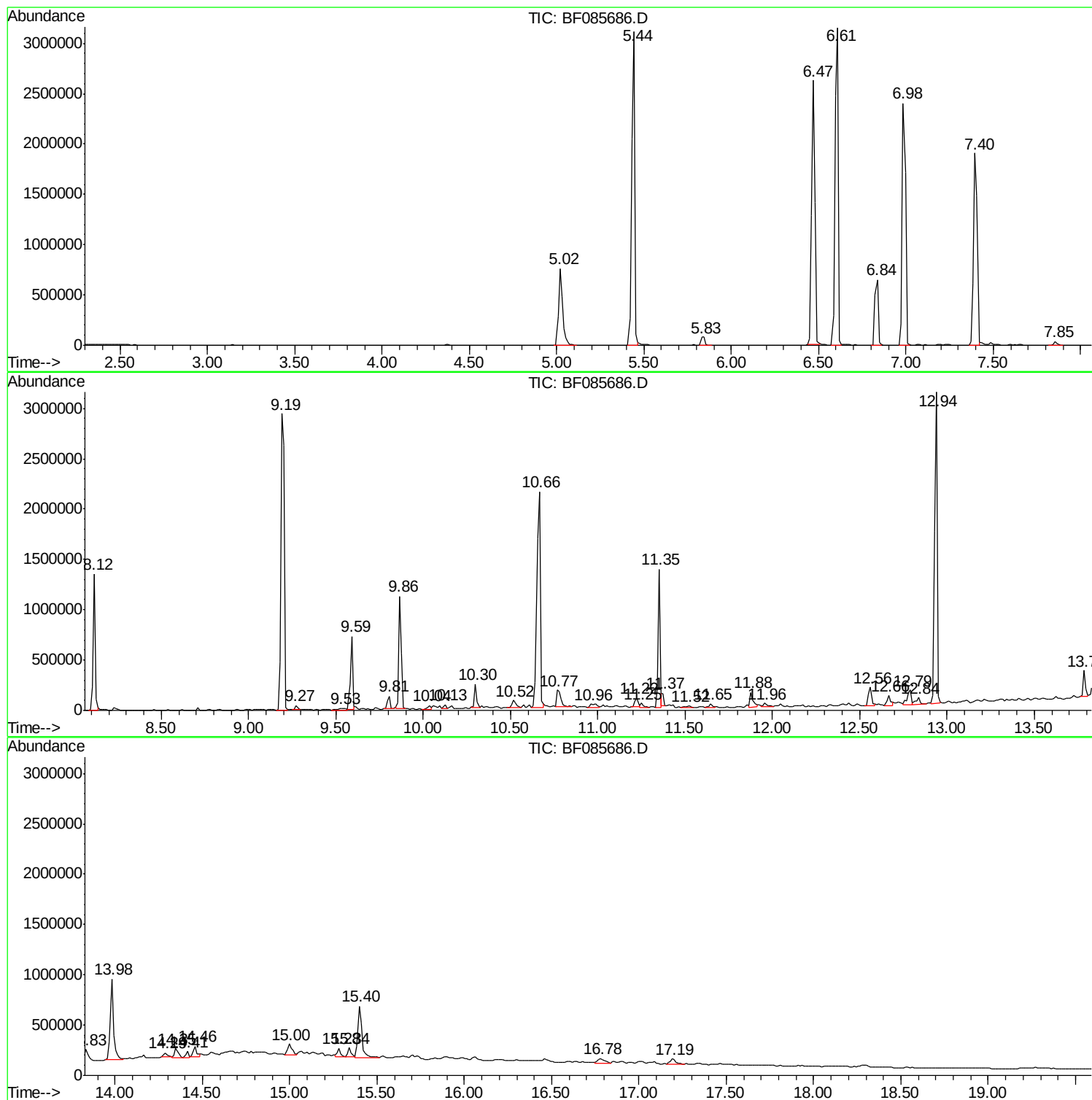
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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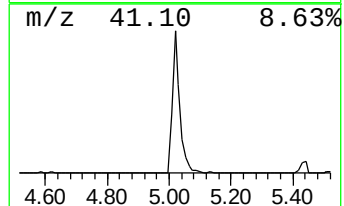
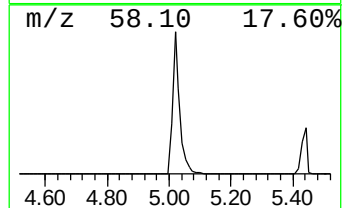
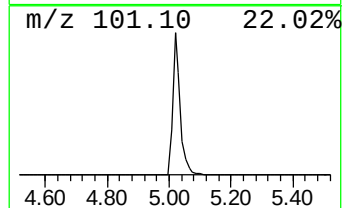
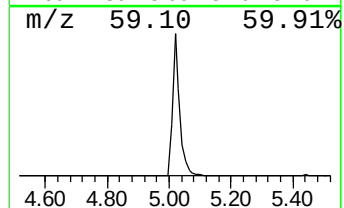
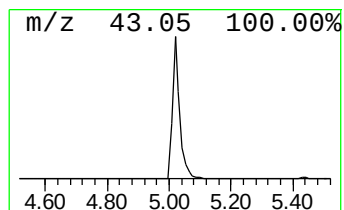
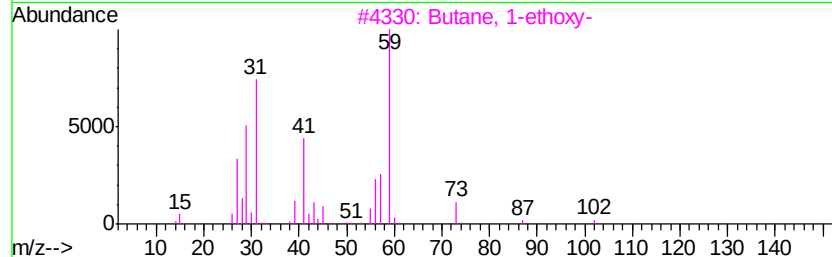
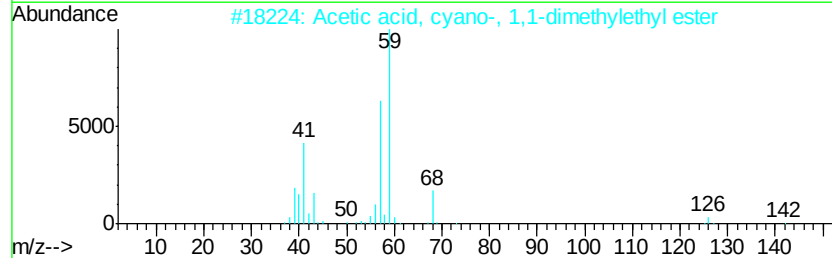
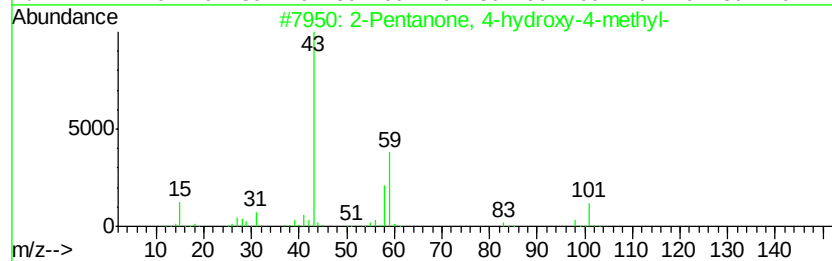
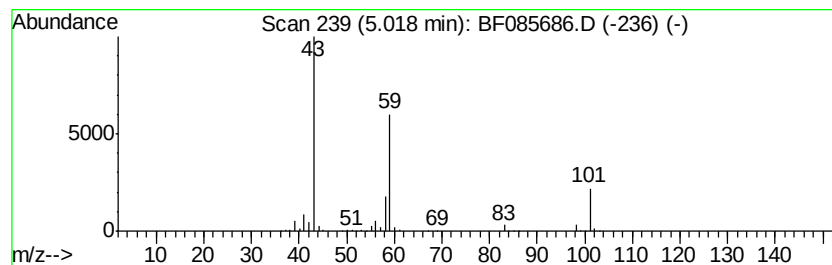
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	30.82 ng	1251460	1,4-Dichlorobenzene-d4	6.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23	
3	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9	
5	Acetone	58	C3H6O	000067-64-1	9	



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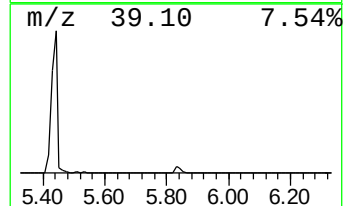
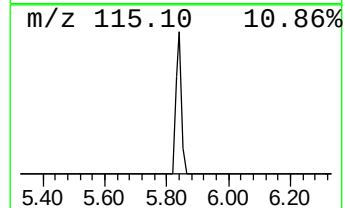
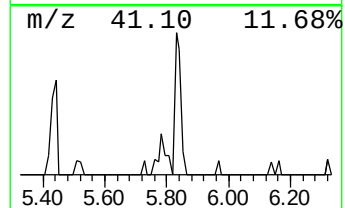
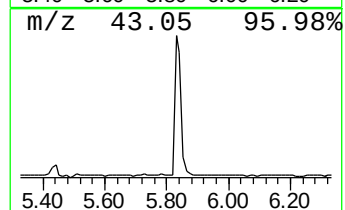
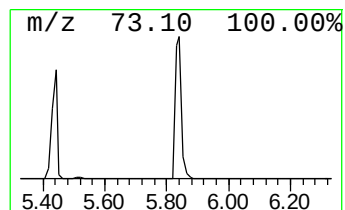
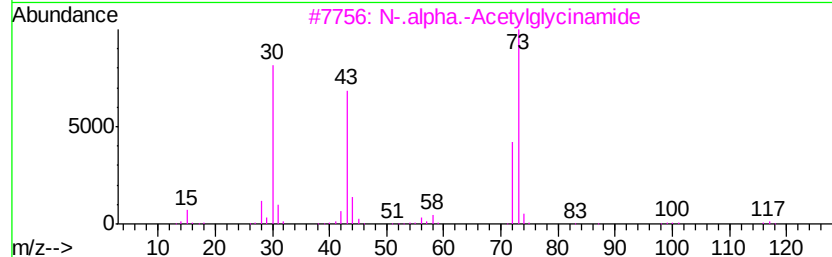
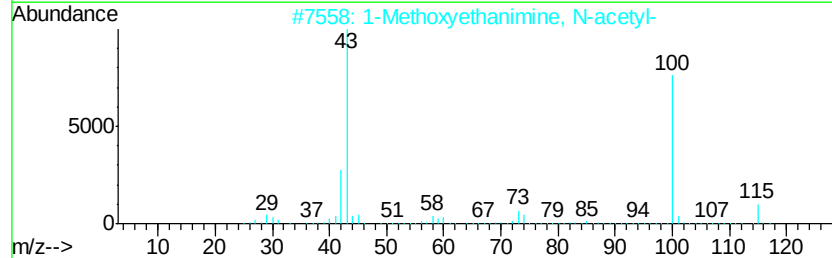
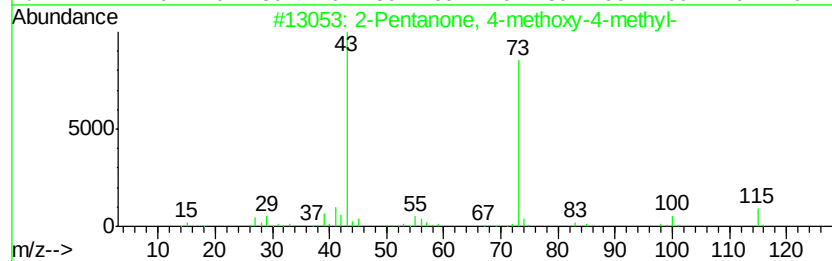
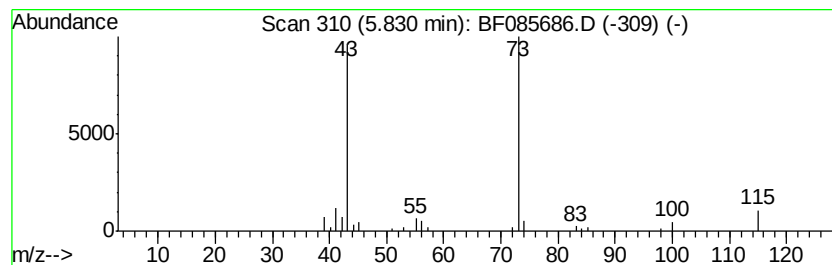
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.83	2.99 ng	121395	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		1-Methoxyethanimine, N-acetyl-	115	C5H9NO2	099028-43-0	38
3		N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	33
4		Butane, 2-methoxy-2,3,3-trimethyl-	130	C8H18O	027705-21-1	25
5		Hexane, 2-methyl-	100	C7H16	000591-76-4	9



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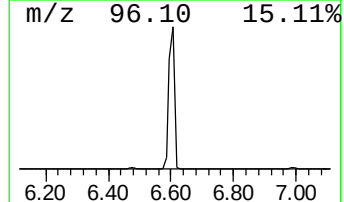
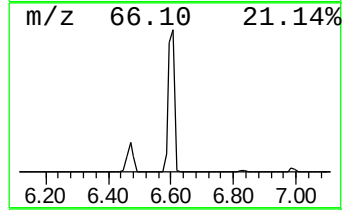
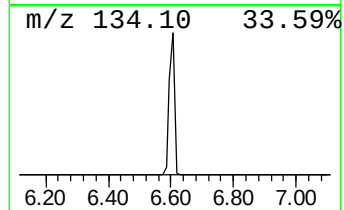
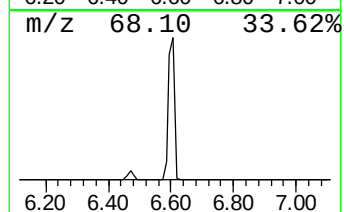
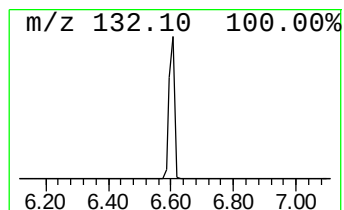
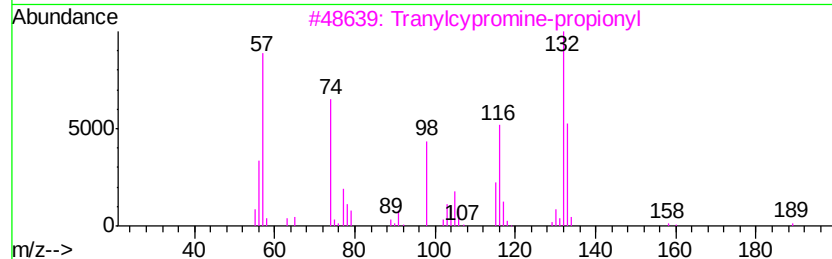
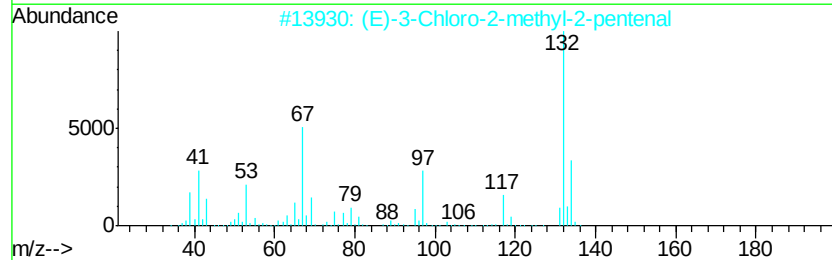
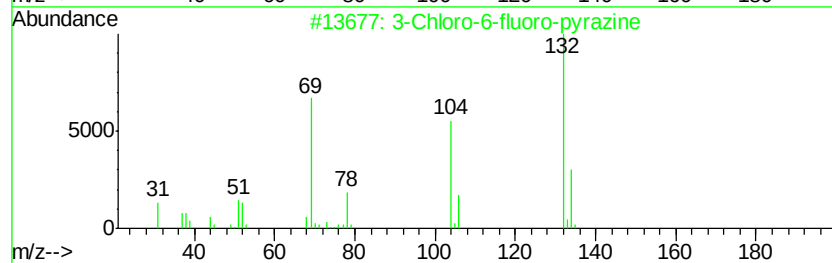
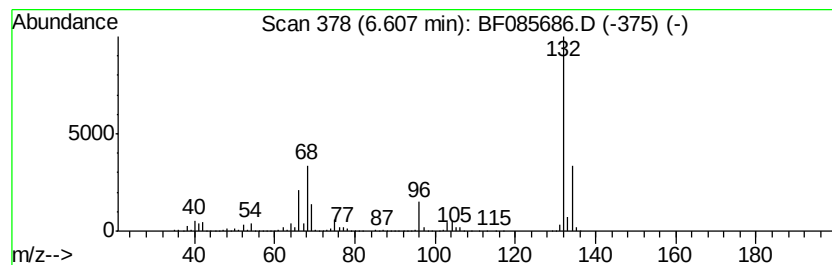
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	100.73 ng	4089700	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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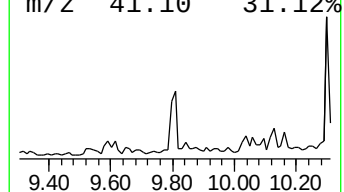
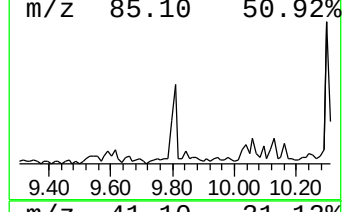
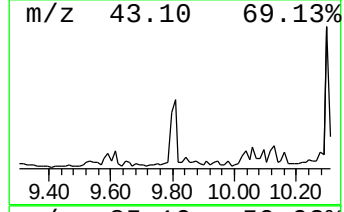
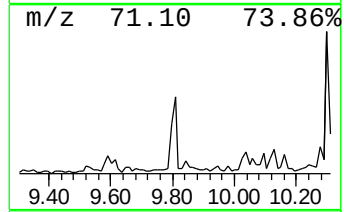
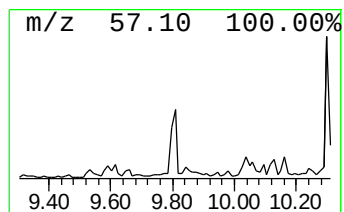
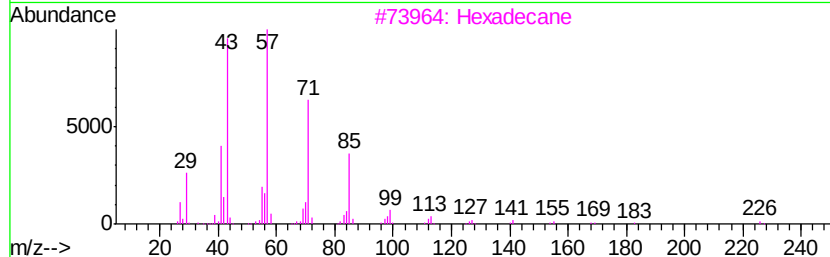
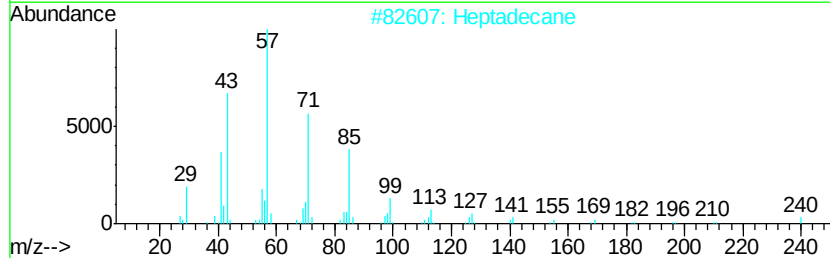
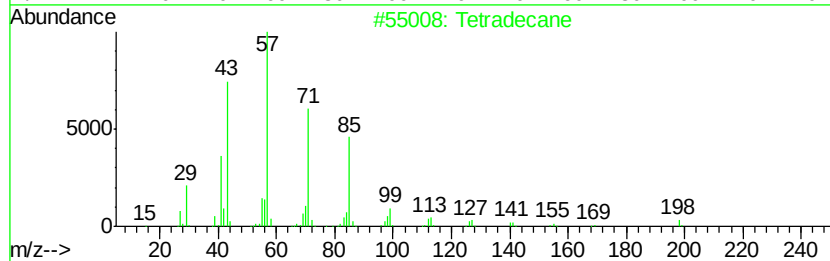
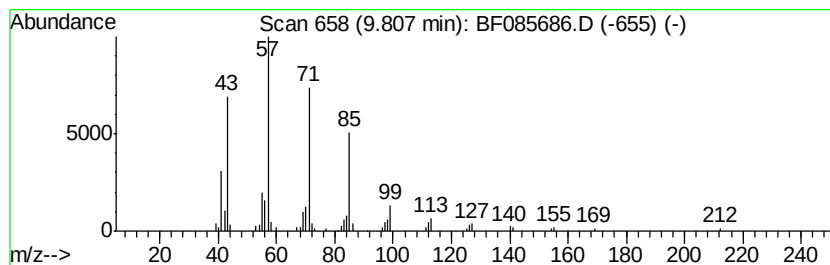
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
9.81	2.47 ng	145065	Acenaphthene-d10		9.86
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	91
2	Heptadecane	240	C17H36	000629-78-7	90
3	Hexadecane	226	C16H34	000544-76-3	86
4	Pentadecane	212	C15H32	000629-62-9	68
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64



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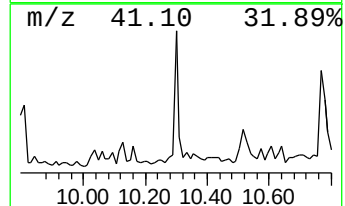
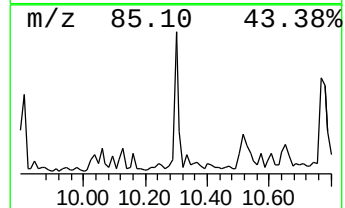
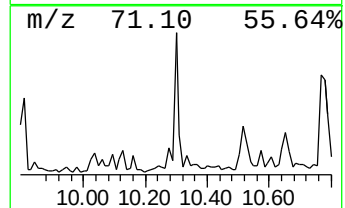
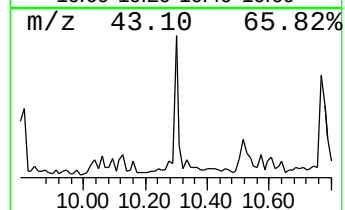
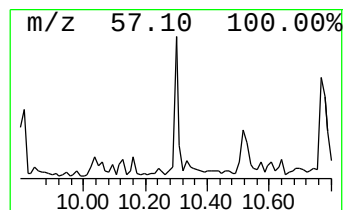
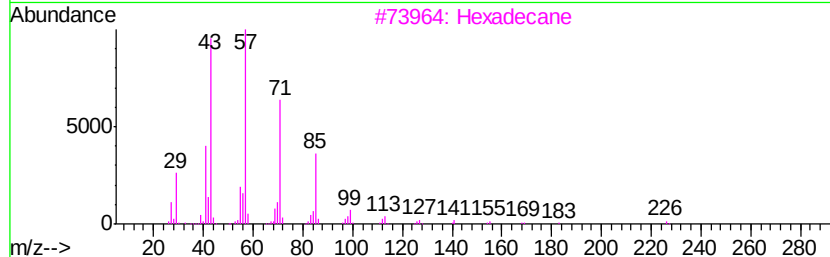
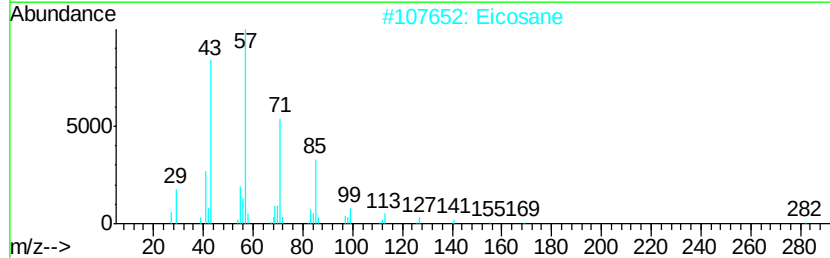
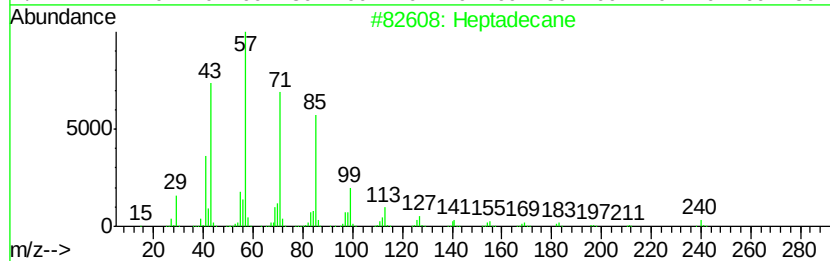
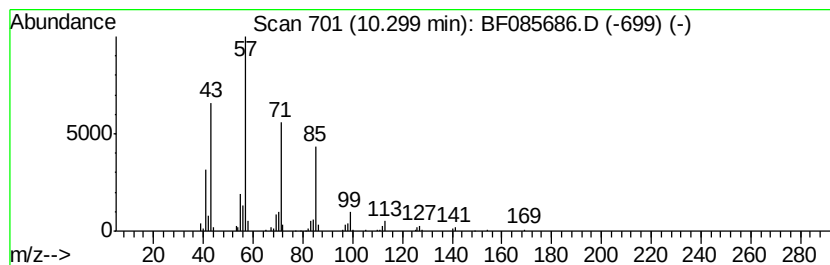
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TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	3.44 ng	202246	Acenaphthene-d10	9.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240	C17H36	000629-78-7 90
2	Eicosane	282	C20H42	000112-95-8 83
3	Hexadecane	226	C16H34	000544-76-3 83
4	Tridecane	184	C13H28	000629-50-5 80
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3 78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085686.D
Acq On : 23 Mar 2016 5:04
Operator : UM/SJ
Sample : H1886-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

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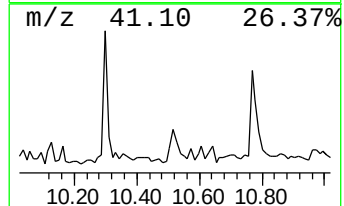
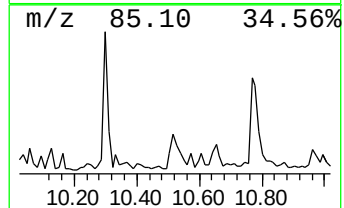
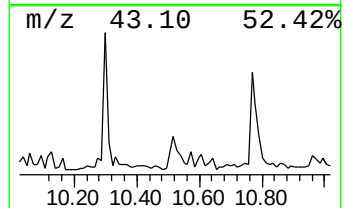
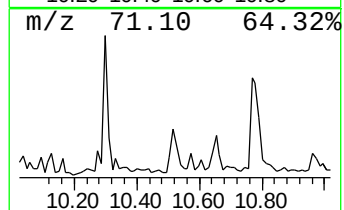
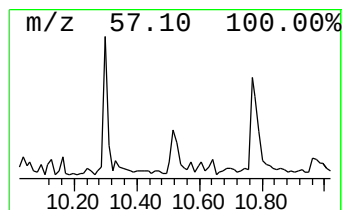
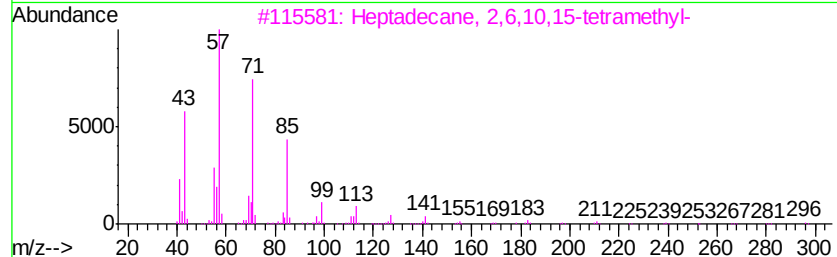
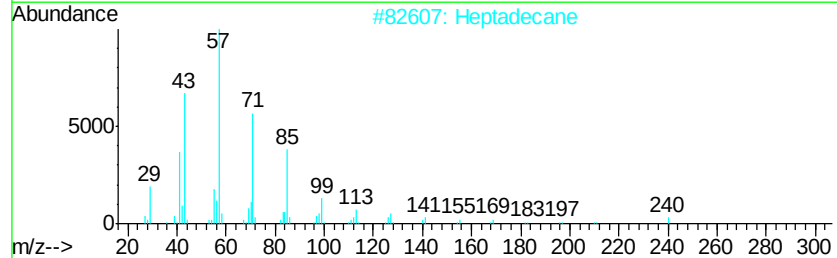
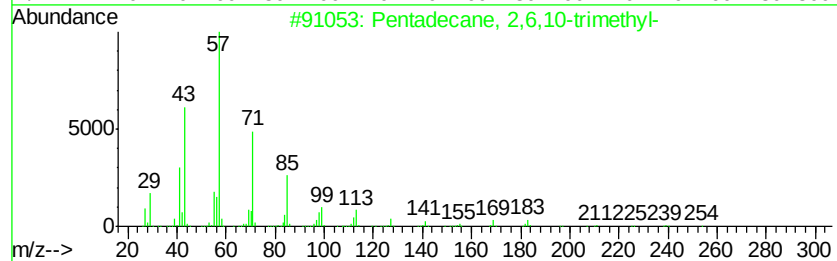
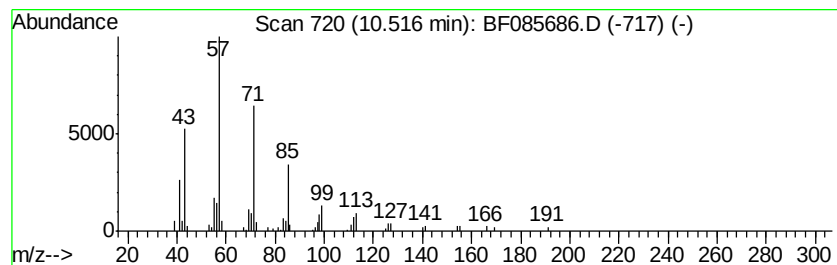
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane, 2,6,10-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.52	2.31 ng	136018	Acenaphthene-d10	9.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
4		Heptacosane	380	C27H56	000593-49-7	87
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
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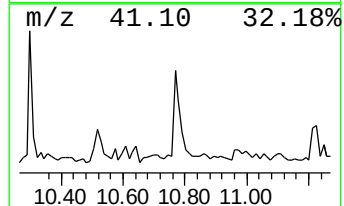
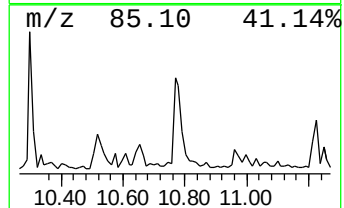
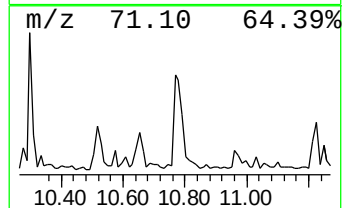
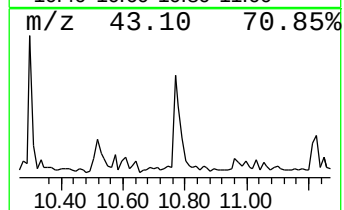
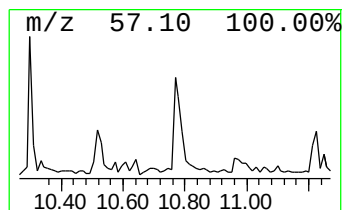
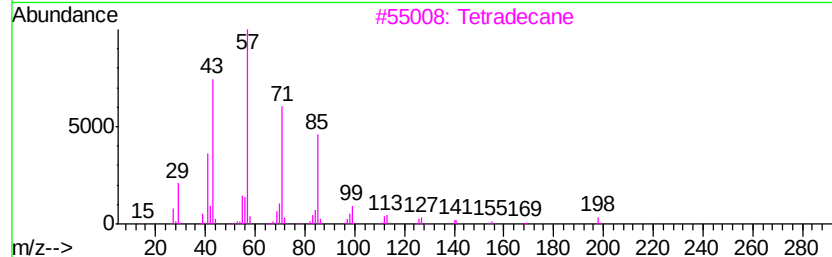
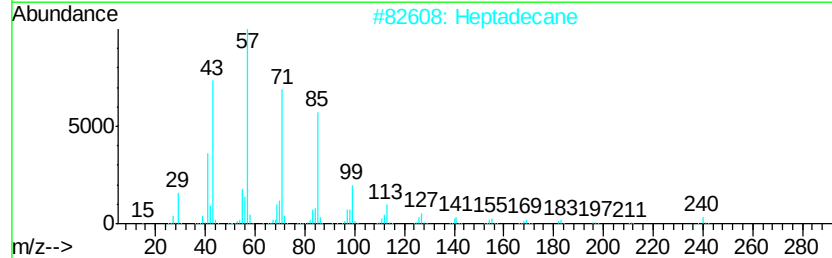
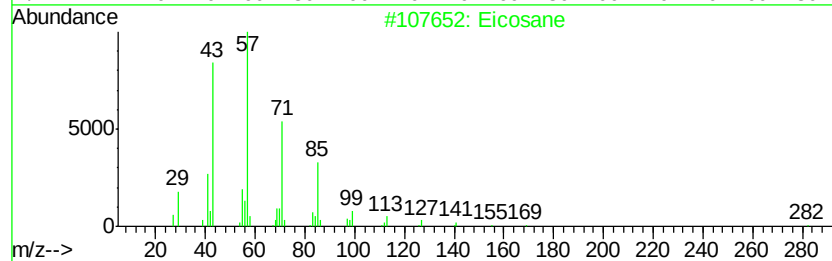
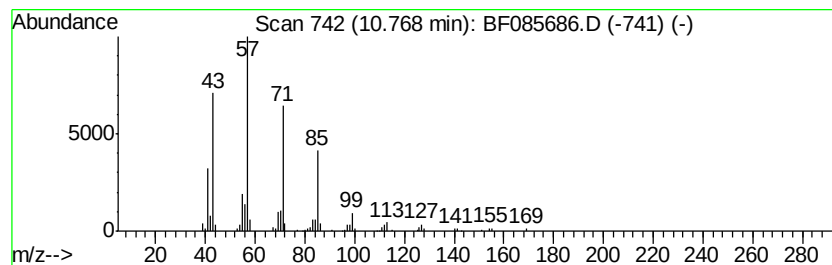
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.77	4.84 ng	283878	Phenanthrene-d10		11.35	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Heptadecane		240	C17H36	000629-78-7	90
3	Tetradecane		198	C14H30	000629-59-4	86
4	Hexadecane		226	C16H34	000544-76-3	78
5	Nonane, 2-methyl-5-propyl-		184	C13H28	031081-17-1	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
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Sample : H1886-03
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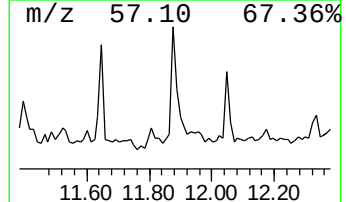
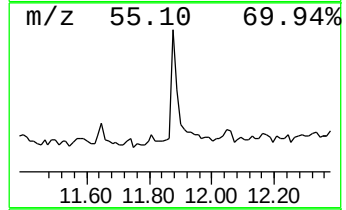
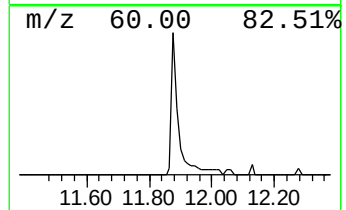
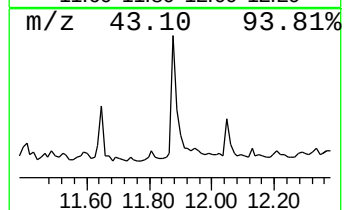
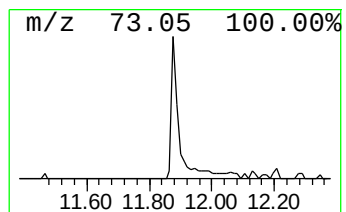
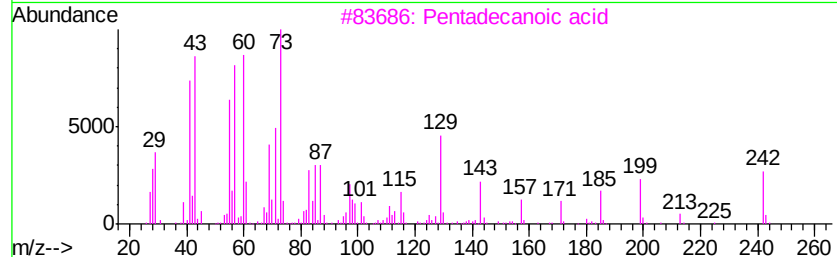
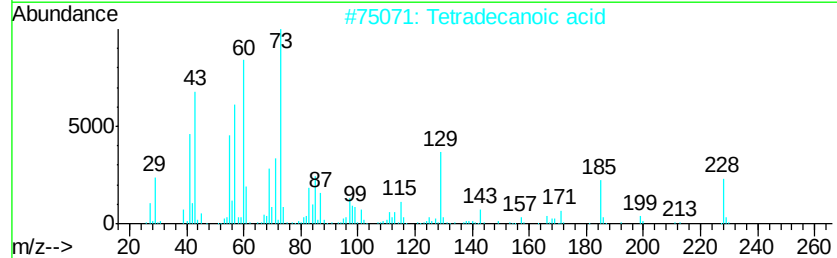
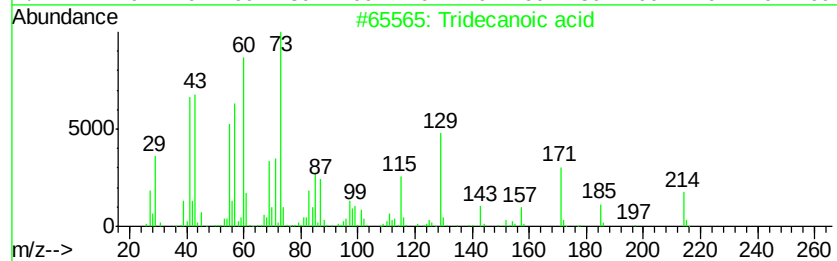
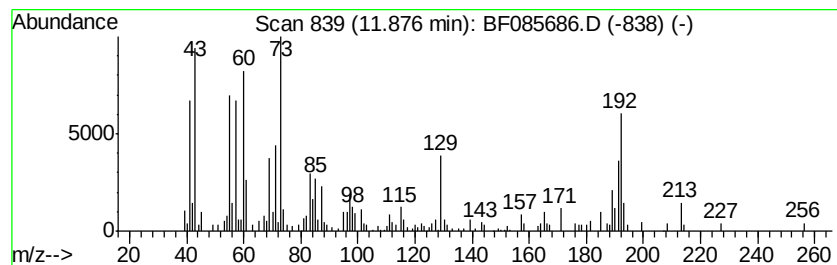
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tridecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.88	3.17 ng	185591	Phenanthrene-d10		11.35
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	86
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	68
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	62
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58
5	n-Decanoic acid	172	C10H20O2	000334-48-5	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
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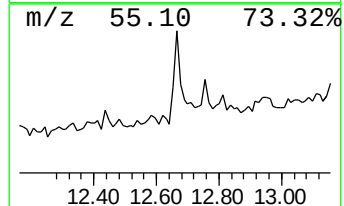
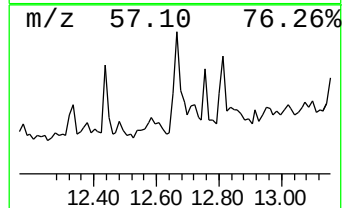
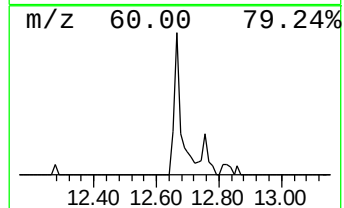
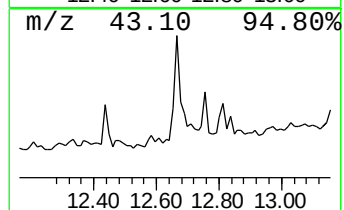
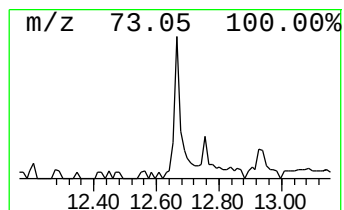
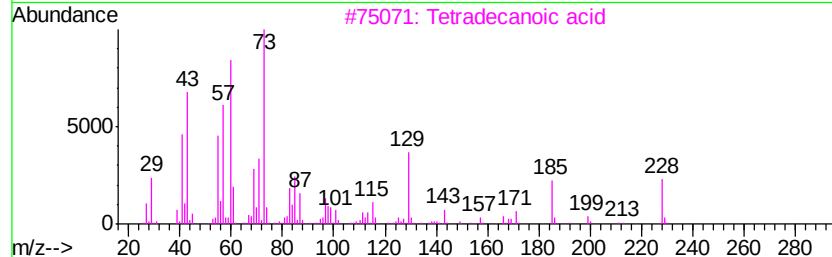
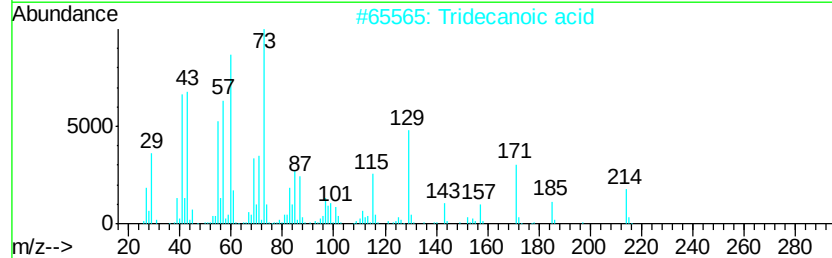
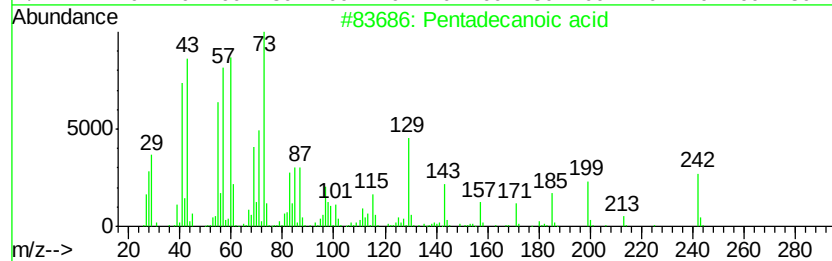
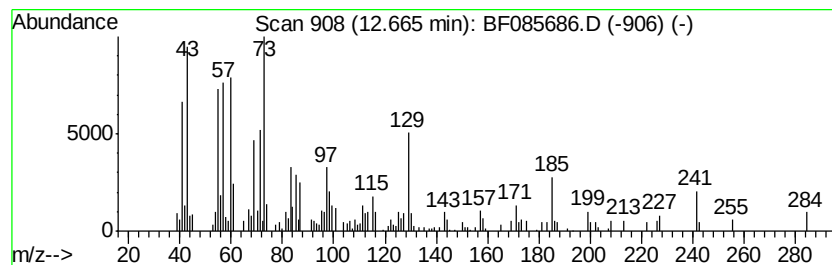
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.66	2.76 ng	134332	Chrysene-d12		13.98
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	89
2	Tridecanoic acid	214	C13H26O2	000638-53-9	64
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4	Undecanoic acid	186	C11H22O2	000112-37-8	62
5	n-Decanoic acid	172	C10H20O2	000334-48-5	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
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ClientSampleId :
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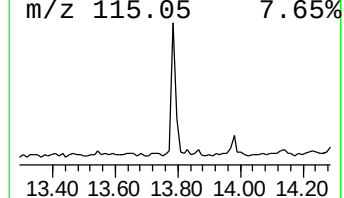
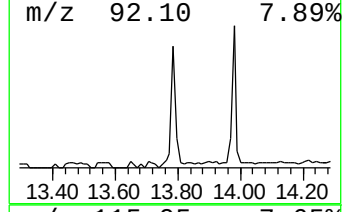
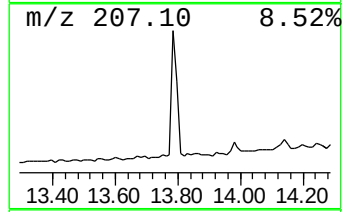
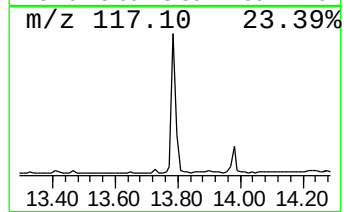
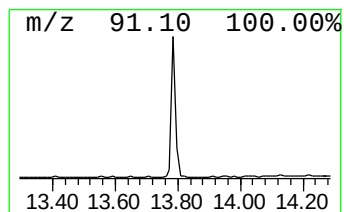
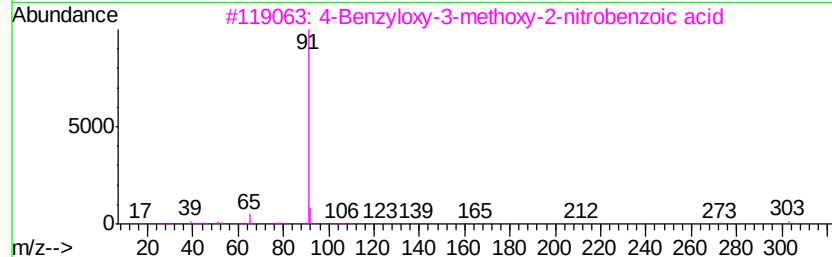
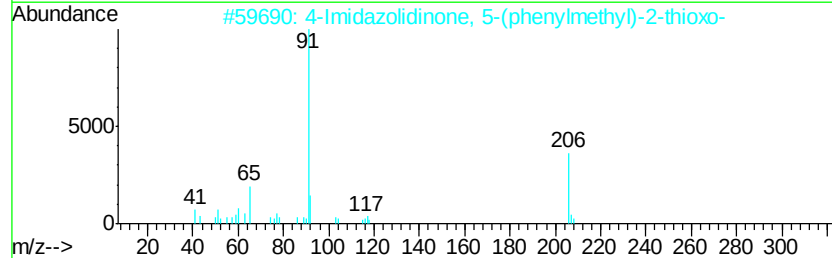
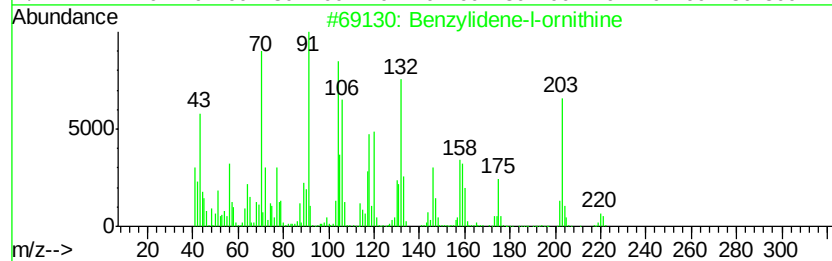
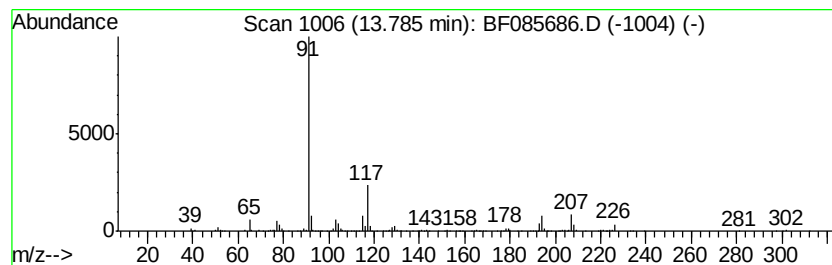
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.79 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	4.67 ng	227057	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzyldene-l-ornithine	220	C12H16N2O2	025693-00-9	38
2		4-Imidazolidinone, 5-(phenylmeth...	206	C10H10N2OS	006330-09-2	37
3		4-Benzyloxy-3-methoxy-2-nitroben...	303	C15H13N06	003584-32-5	14
4		Diazene, bis(4-methylphenyl)-, 1...	226	C14H14N2O	000955-98-6	14
5		(+)-N-Benzyl-.alpha.-phenethylamine	211	C15H17N	038235-77-7	12



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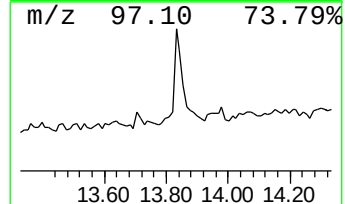
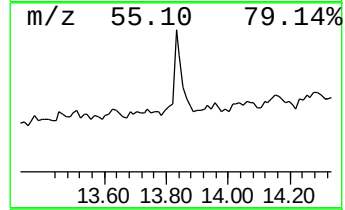
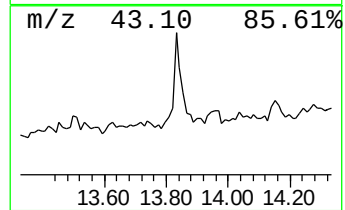
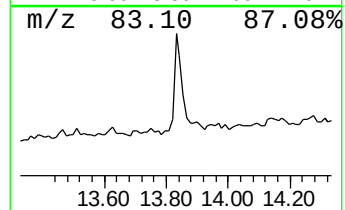
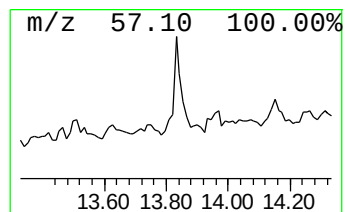
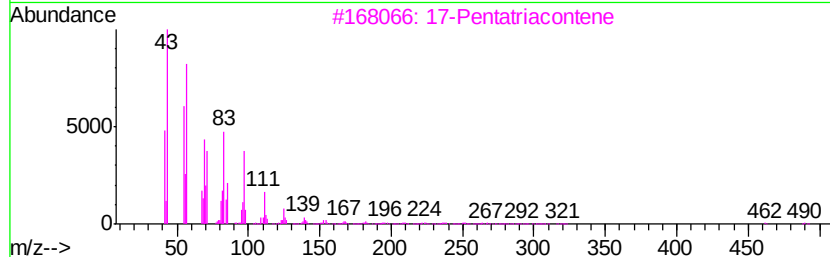
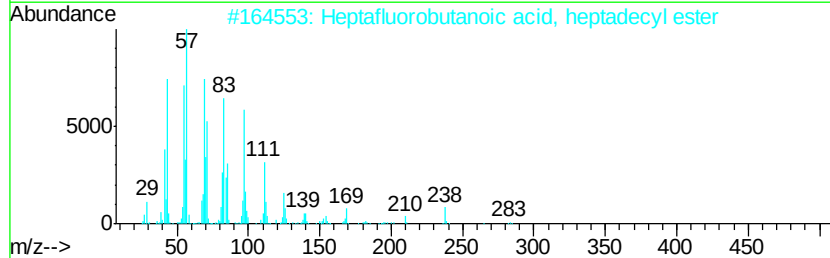
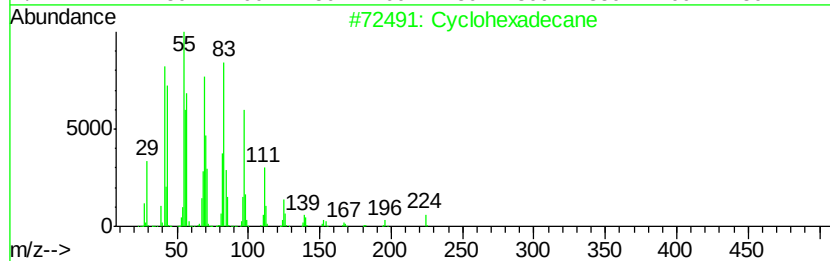
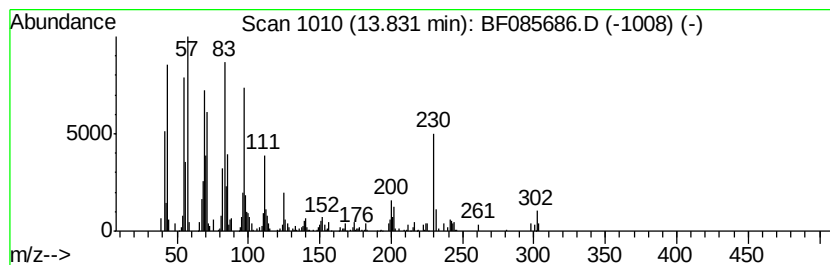
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cyclohexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	3.44 ng	167226	Chrysene-d12	13.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 97
2		Heptafluorobutanoic acid, heptad...	452 C21H35F7O2	1000282-97-3 91
3		17-Pentatriacontene	491 C35H70	006971-40-0 87
4		Pentafluoropropionic acid, hepta...	402 C20H35F5O2	1000283-04-2 87
5		Dichloroacetic acid, heptadecyl ...	366 C19H36Cl2O2	1000282-98-2 87



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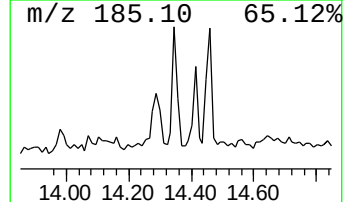
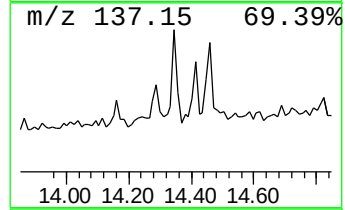
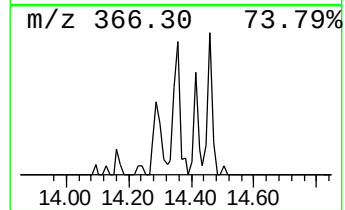
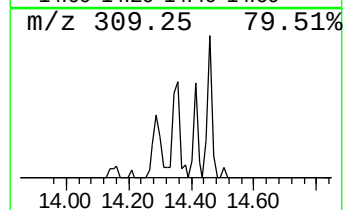
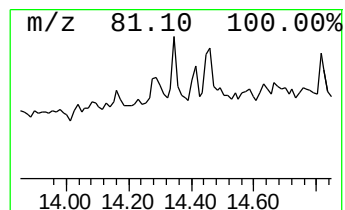
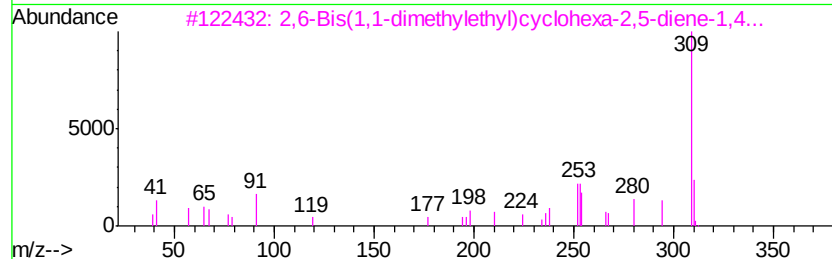
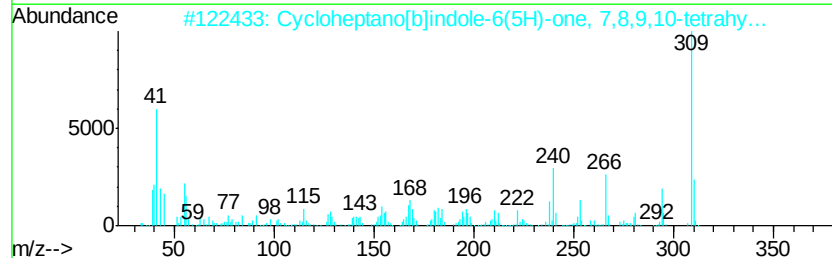
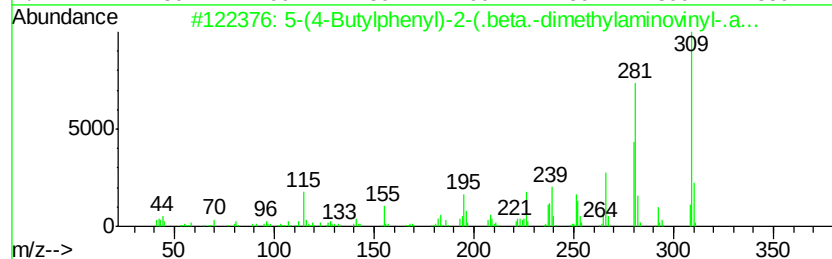
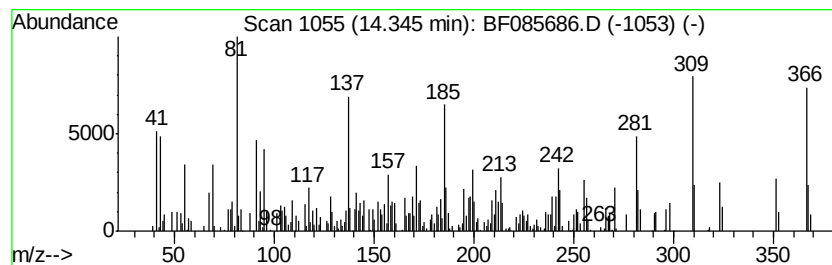
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 5-(4-Butylphenyl)-2-(.beta.... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.35	3.09 ng	150033	Chrysene-d12	13.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(4-Butylphenyl)-2-(.beta.-dime...	309	C19H23N3O	131548-35-1	51
2	Cycloheptano[b]indole-6(5H)-one,...	309	C21H27NO	327990-59-0	18
3	2,6-Bis(1,1-dimethylethyl)cycloh...	309	C21H27NO	017118-97-7	15
4	2-Oxo-5-phenyl-4-(4-methoxypheny...	366	C24H18N2O2	028840-53-1	14
5	1-(4-Anisidino)-2-phenyl-2-(1,2,...	309	C17H15N3OS	153333-55-2	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085686.D
Acq On : 23 Mar 2016 5:04
Operator : UM/SJ
Sample : H1886-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

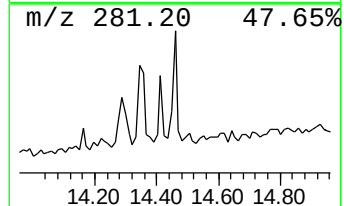
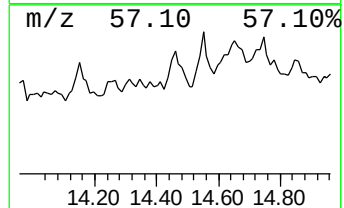
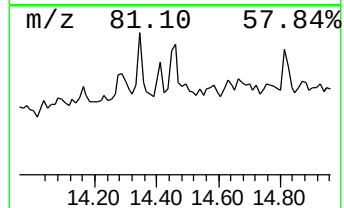
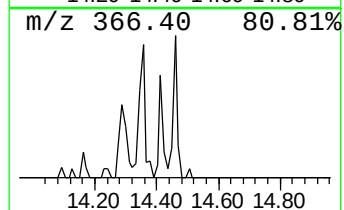
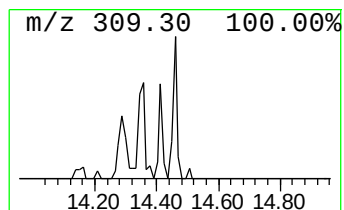
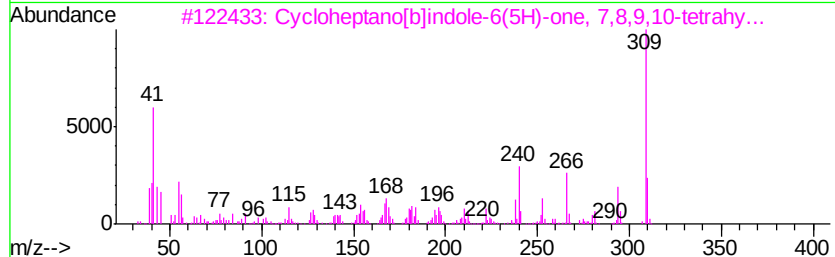
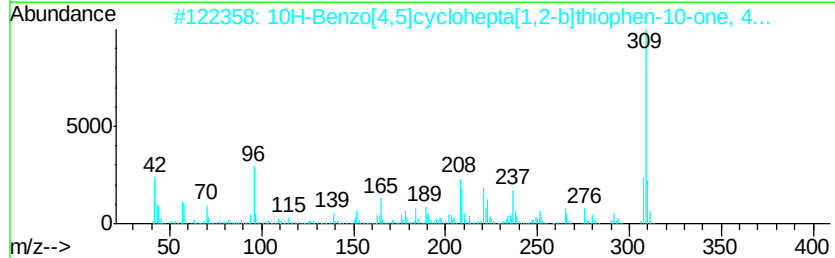
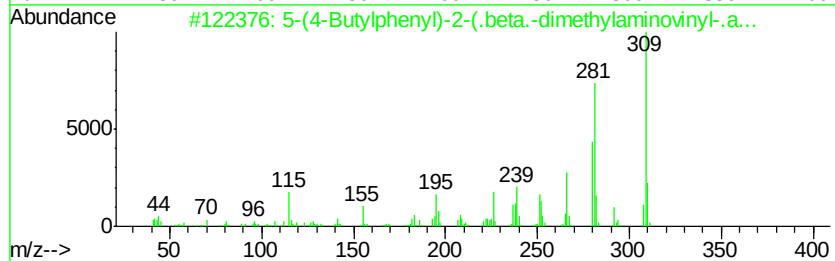
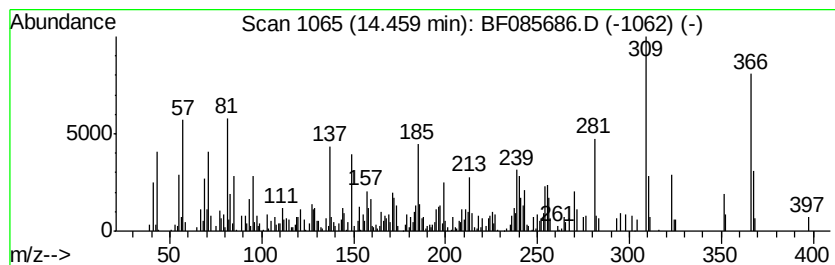
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown14.46 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.46	2.91 ng	141650	Chrysene-d12	13.98		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(4-Butylphenyl)-2-(.beta.-dime...	309	C19H23N3O	131548-35-1	42	
2	10H-Benzo[4,5]cyclohepta[1,2-b]t...	309	C19H19NOS	034580-13-7	25	
3	Cycloheptano[b]indole-6(5H)-one,...	309	C21H27NO	327990-59-0	22	
4	2,6-Bis(1,1-dimethylethyl)cycloh...	309	C21H27NO	017118-97-7	20	
5	Indole-3-carboxylic acid, 5-hydr...	309	C19H19NO3	005165-44-6	20	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
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Acq On : 23 Mar 2016 5:04
Operator : UM/SJ
Sample : H1886-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-6

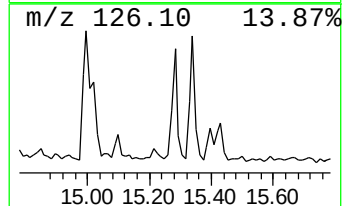
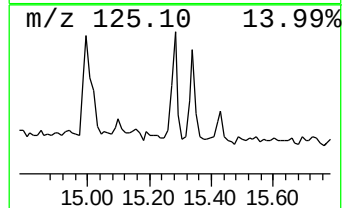
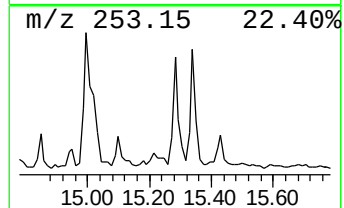
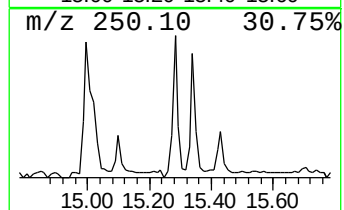
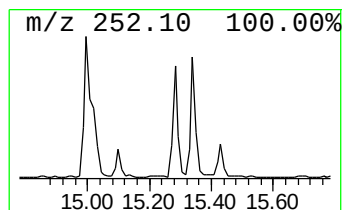
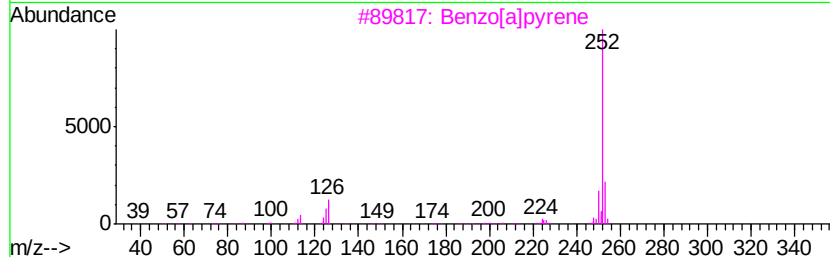
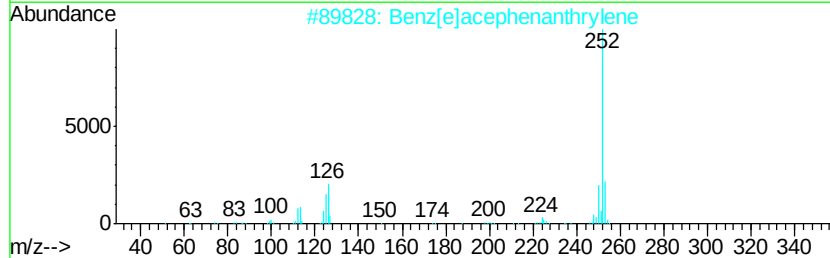
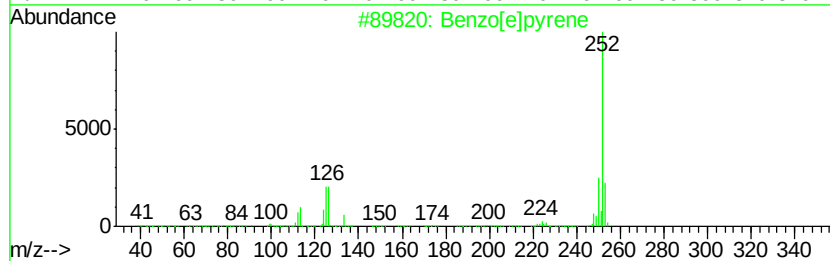
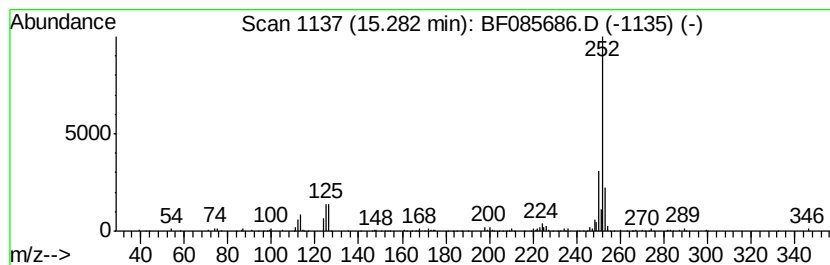
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzo[e]pyrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	2.65 ng	103158	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
3		Benzo[a]pyrene	252	C20H12	000050-32-8	96
4		Perylene	252	C20H12	000198-55-0	95
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93



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 Sample : H1886-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-6

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	30.8	ng	1251460	1	6.84	811987	20.0
2-Pentanone, 4-me...	5.83	3.0	ng	121395	1	6.84	811987	20.0
unknown6.61	6.61	100.7	ng	4089700	1	6.84	811987	20.0
Tetradecane	9.81	2.5	ng	145065	3	9.86	1175770	20.0
Heptadecane	10.30	3.4	ng	202246	3	9.86	1175770	20.0
Pentadecane, 2,6,...	10.52	2.3	ng	136018	3	9.86	1175770	20.0
Eicosane	10.77	4.8	ng	283878	4	11.35	1172500	20.0
Tridecanoic acid	11.88	3.2	ng	185591	4	11.35	1172500	20.0
Pentadecanoic acid	12.66	2.8	ng	134332	5	13.98	972482	20.0
unknown13.79	13.79	4.7	ng	227057	5	13.98	972482	20.0
Cyclohexadecane	13.83	3.4	ng	167226	5	13.98	972482	20.0
5-(4-Butylphenyl)...	14.35	3.1	ng	150033	5	13.98	972482	20.0
unknown14.46	14.46	2.9	ng	141650	5	13.98	972482	20.0
Benzo[e]pyrene	15.28	2.6	ng	103158	6	15.40	777841	20.0