

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
 Data File : BF085022.D
 Acq On : 11 Feb 2016 17:15
 Operator : SJ/IZ
 Sample : H1377-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB04-0-1

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-BF020116.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	1.701	8	10	11	rVB	155869	207206	3.75%	0.387%
2	1.735	11	13	56	rVB	2103592	4263136	77.13%	7.970%
3	4.741	274	276	285	rBV	769270	1522687	27.55%	2.847%
4	5.199	313	316	319	rBV	3967066	4717630	85.35%	8.820%
5	6.147	397	399	402	rBV	484546	508574	9.20%	0.951%
6	6.273	406	410	412	rBV	4424162	5353383	96.85%	10.008%
7	6.376	417	419	422	rBV	3984033	4912843	88.88%	9.185%
8	6.604	437	439	442	rBV	1042968	948697	17.16%	1.774%
9	6.764	450	453	455	rBV	4109255	3734952	67.57%	6.983%
10	7.187	486	490	492	rBV	2889778	3695756	66.86%	6.909%
11	7.656	529	531	536	rBV	51155	66591	1.20%	0.124%
12	7.896	549	552	554	rBV	1596838	1374576	24.87%	2.570%
13	8.650	616	618	627	rVB	349851	347507	6.29%	0.650%
14	8.982	643	647	649	rBV	5140684	5527314	100.00%	10.333%
15	9.382	679	682	684	rBV	793383	850497	15.39%	1.590%
16	9.599	698	701	702	rBV	176659	167123	3.02%	0.312%
17	9.645	702	705	708	rVB	1707329	1510019	27.32%	2.823%
18	9.828	718	721	725	rBV2	35165	79932	1.45%	0.149%
19	10.091	742	744	746	rVB	199121	222900	4.03%	0.417%
20	10.308	760	763	767	rBV	75707	149642	2.71%	0.280%
21	10.433	771	774	777	rVB	3141875	3435569	62.16%	6.423%
22	10.571	784	786	790	rBV	195991	305354	5.52%	0.571%
23	10.753	800	802	807	rBV4	28361	67946	1.23%	0.127%
24	11.016	823	825	826	rBV	99790	104690	1.89%	0.196%
25	11.119	832	834	837	rVB	1337874	1353271	24.48%	2.530%
26	11.679	881	883	885	rBV	87055	86297	1.56%	0.161%
27	12.548	957	959	962	rVB2	89135	88343	1.60%	0.165%
28	12.719	971	974	976	rBV	3249863	4402772	79.65%	8.231%
29	13.622	1050	1053	1056	rBV2	318755	465564	8.42%	0.870%
30	13.748	1062	1064	1067	rBV	1028791	1140582	20.64%	2.132%
31	14.251	1106	1108	1112	rBV	64238	91946	1.66%	0.172%
32	14.605	1137	1139	1141	rBV	59691	60443	1.09%	0.113%
33	14.834	1157	1159	1164	rBV	66757	115493	2.09%	0.216%
34	15.108	1180	1183	1186	rBV	745161	877199	15.87%	1.640%

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SB04-0-1

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.531	1217	1220	1222	rBV	52125	90170	1.63%	0.169%
36	15.634	1227	1229	1232	rVB2	38166	60048	1.09%	0.112%
37	16.445	1296	1300	1303	rBV	28384	64592	1.17%	0.121%
38	16.834	1330	1334	1341	rBV4	60721	177650	3.21%	0.332%
39	17.017	1347	1350	1356	rVB2	23522	64584	1.17%	0.121%
40	17.531	1388	1395	1399	rBV9	15913	62913	1.14%	0.118%
41	17.691	1404	1409	1415	rBV5	43530	148445	2.69%	0.278%
42	18.526	1477	1482	1486	rBV5	18914	64849	1.17%	0.121%

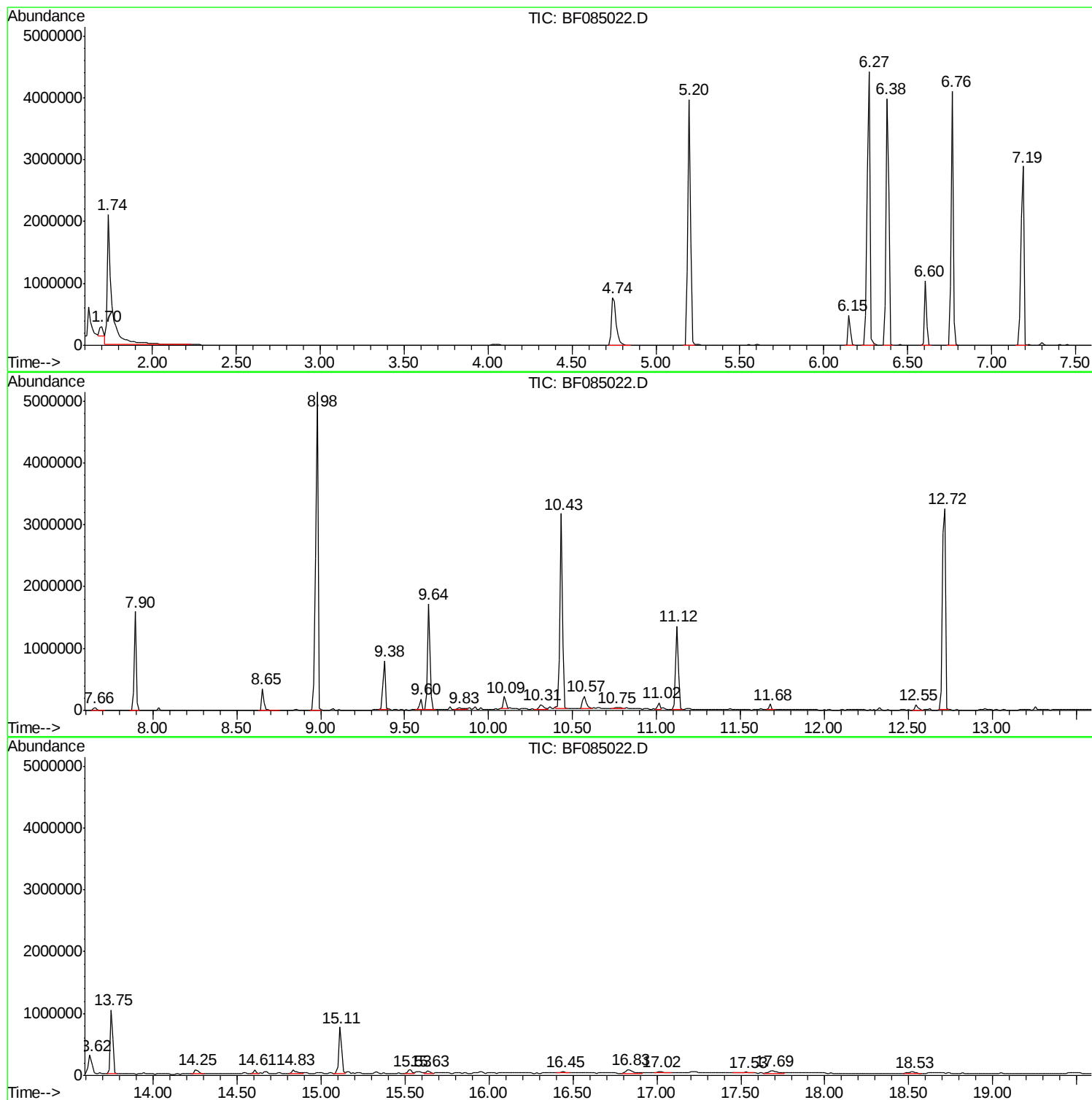
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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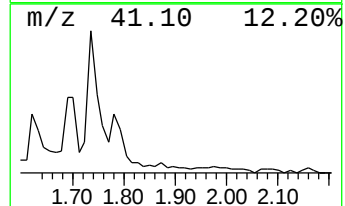
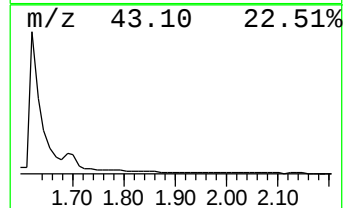
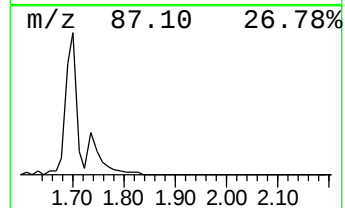
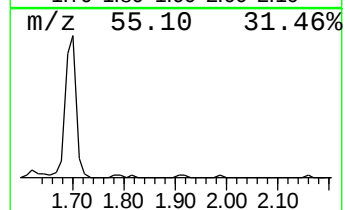
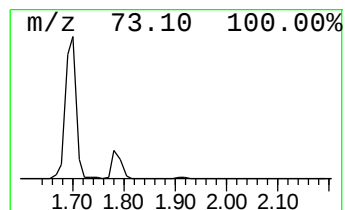
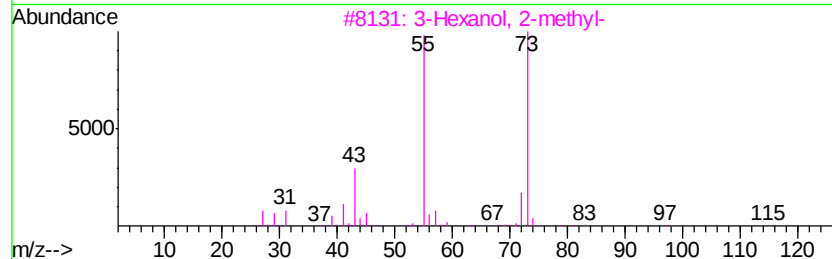
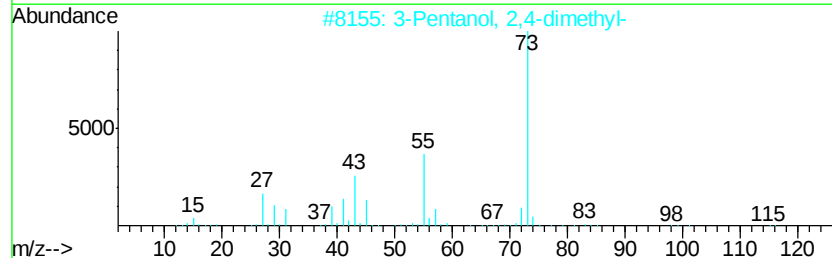
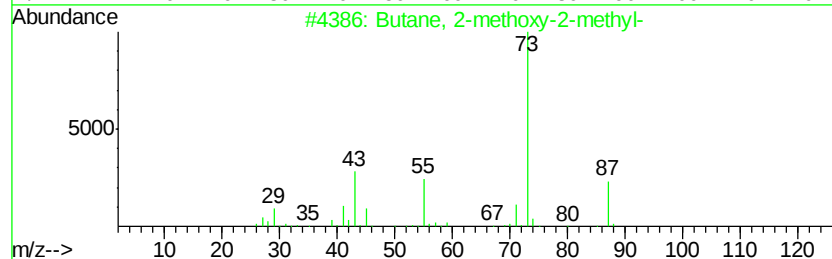
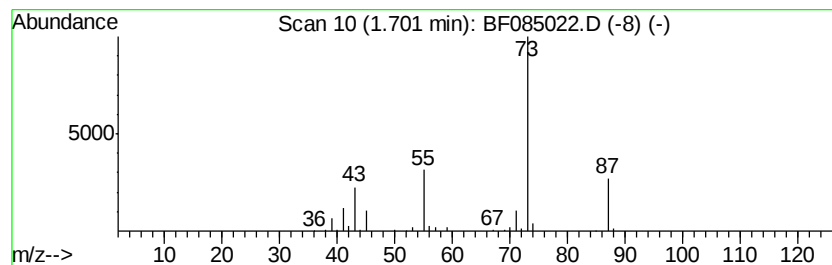
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.70	4.37 ng	207206	1,4-Dichlorobenzene-d4	6.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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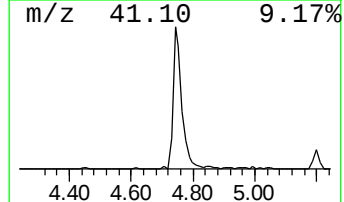
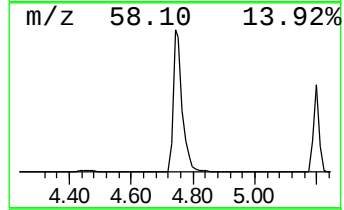
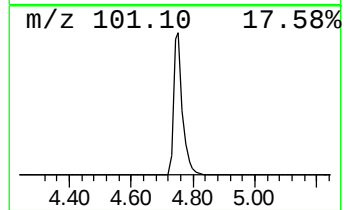
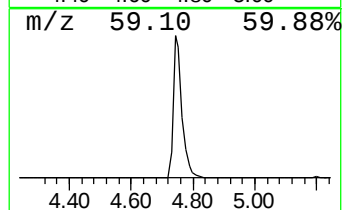
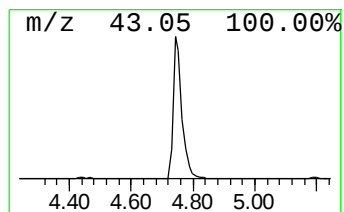
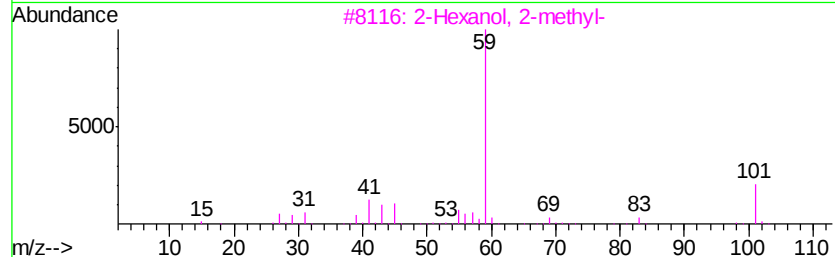
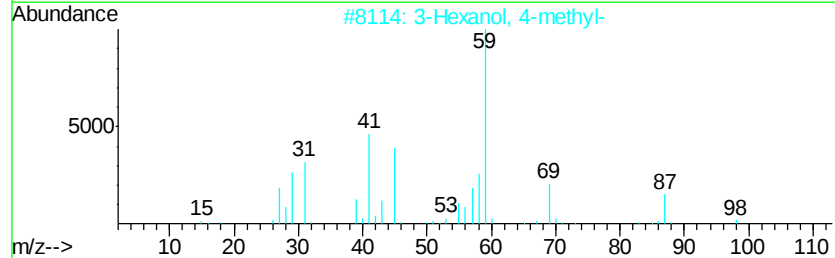
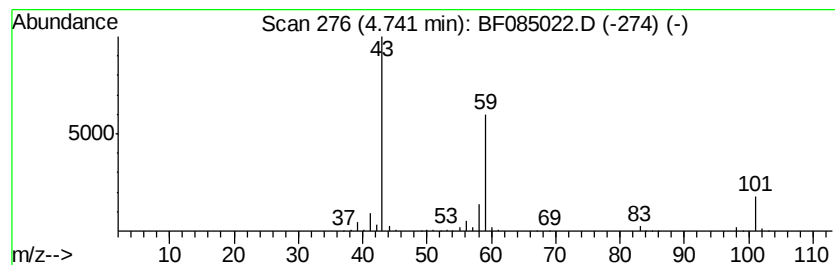
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.74	32.10 ng	1522690	1,4-Dichlorobenzene-d4	6.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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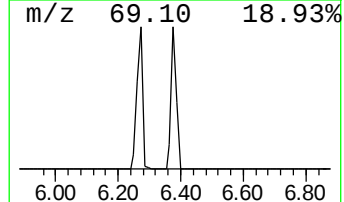
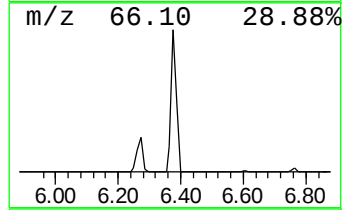
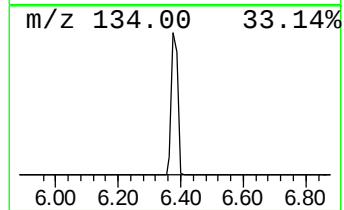
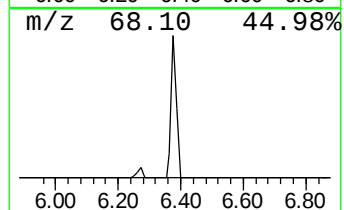
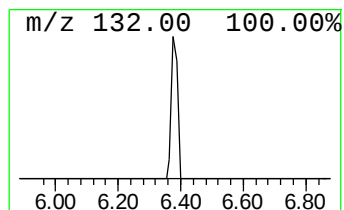
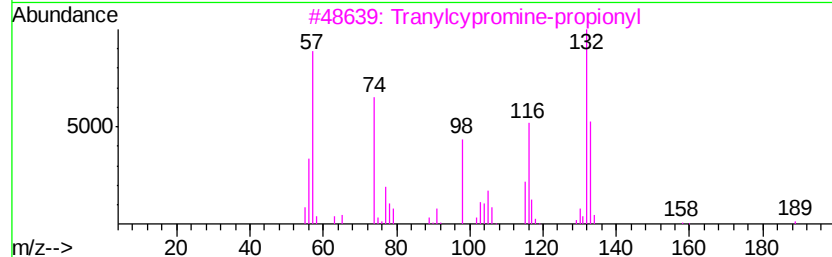
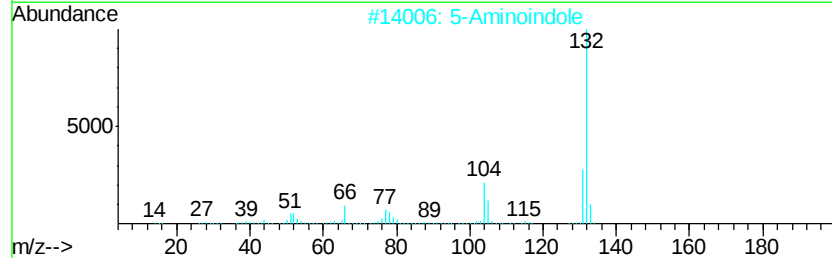
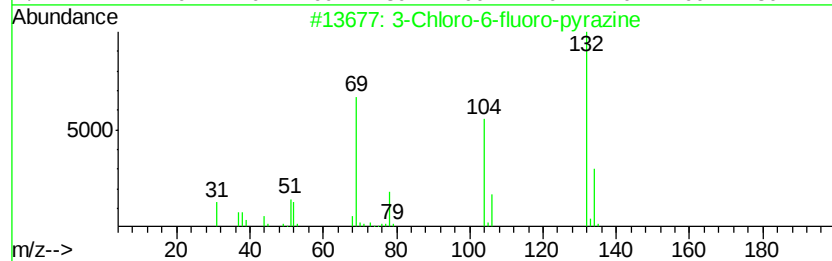
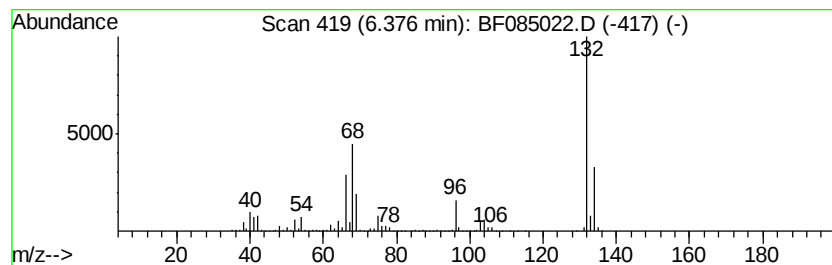
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	103.57 ng	4912840	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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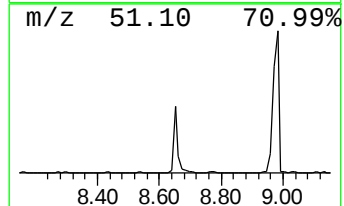
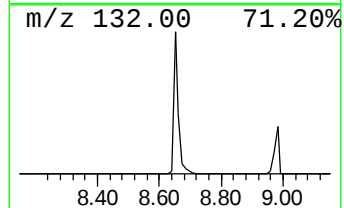
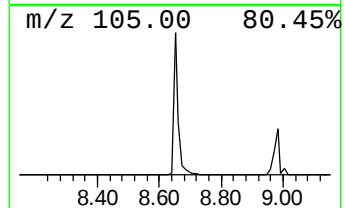
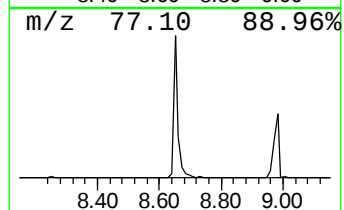
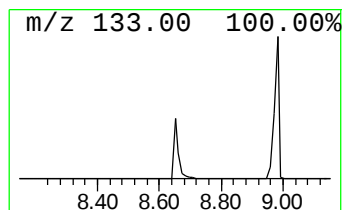
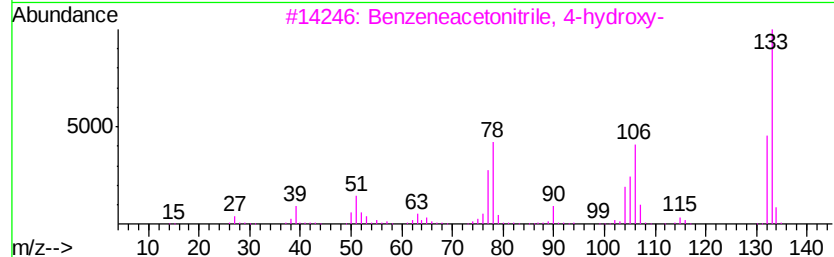
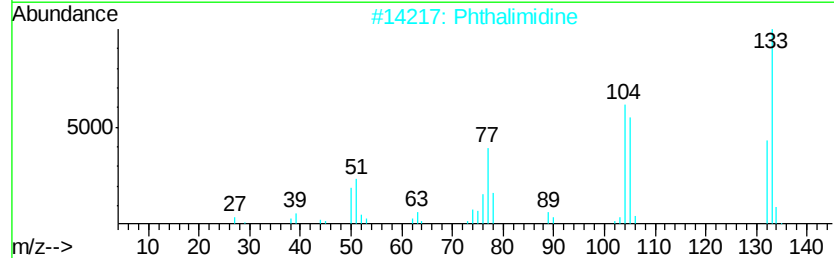
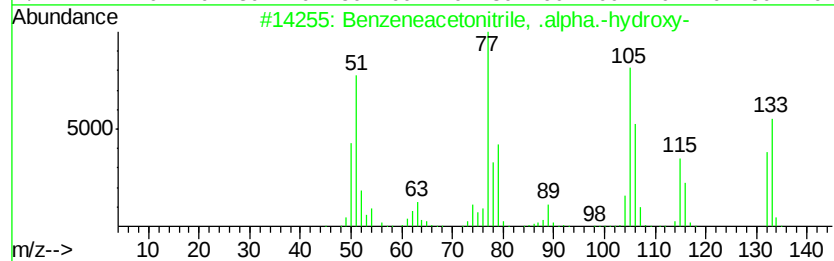
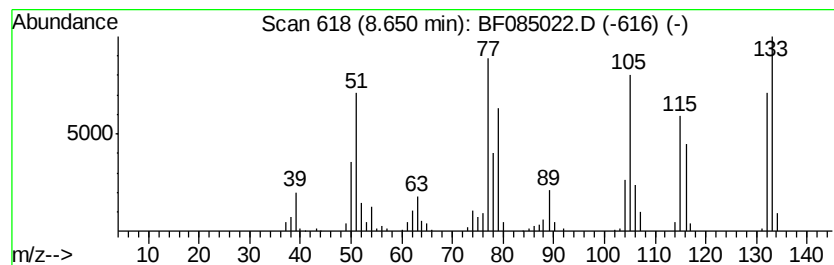
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzeneacetonitrile, .alpha... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.65	5.06 ng	347507	Naphthalene-d8	7.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetonitrile, .alpha.-hyd...	133	C8H7NO	000532-28-5	64
2		Phthalimidine	133	C8H7NO	000480-91-1	49
3		Benzeneacetonitrile, 4-hydroxy-	133	C8H7NO	014191-95-8	27
4		Furo[2,3-c]pyridine, 2-methyl-	133	C8H7NO	069022-76-0	25
5		3-(4-Pyridyl)acrylaldehyde	133	C8H7NO	026505-36-2	22



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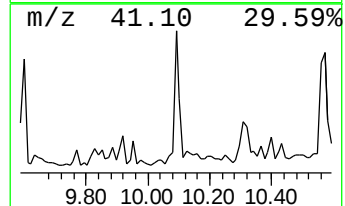
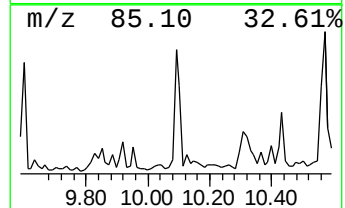
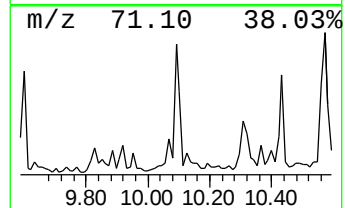
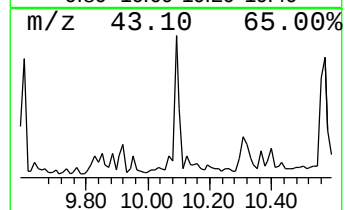
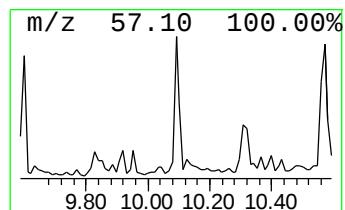
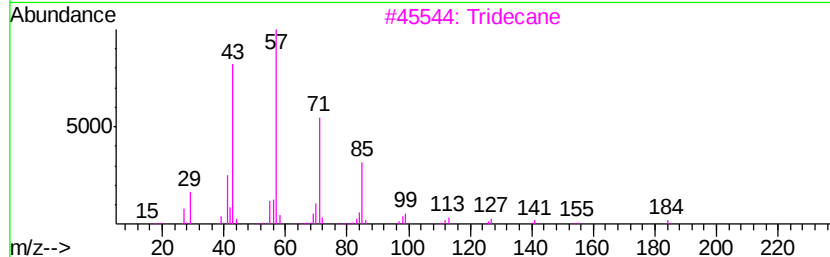
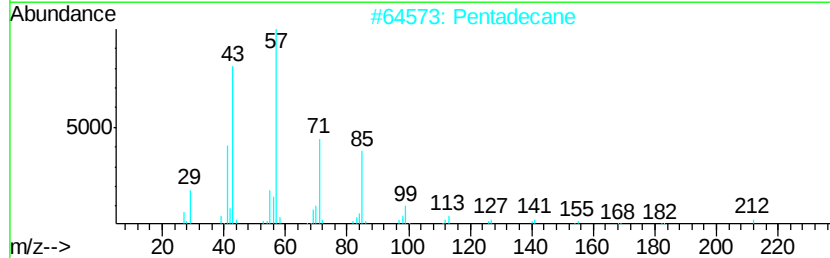
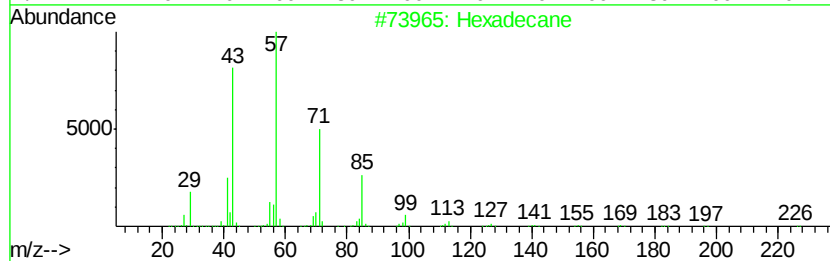
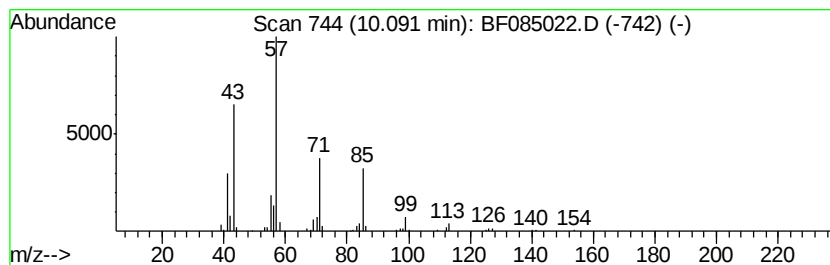
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.09	2.95 ng	222900	Acenaphthene-d10	9.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226	C16H34	000544-76-3 90
2	Pentadecane	212	C15H32	000629-62-9 83
3	Tridecane	184	C13H28	000629-50-5 83
4	Heptadecane	240	C17H36	000629-78-7 78
5	Dotriacontane	451	C32H66	000544-85-4 64



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Sample : H1377-06
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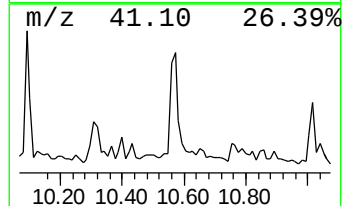
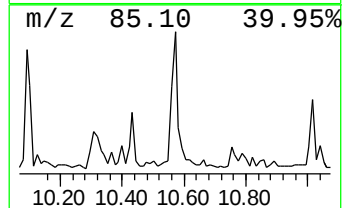
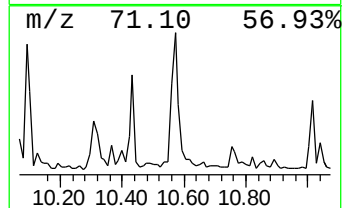
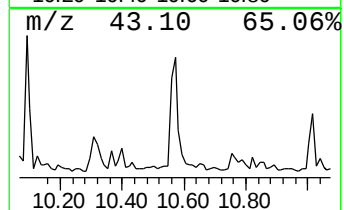
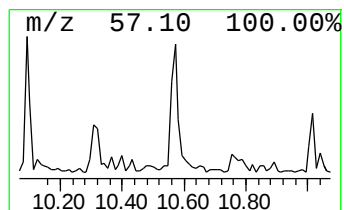
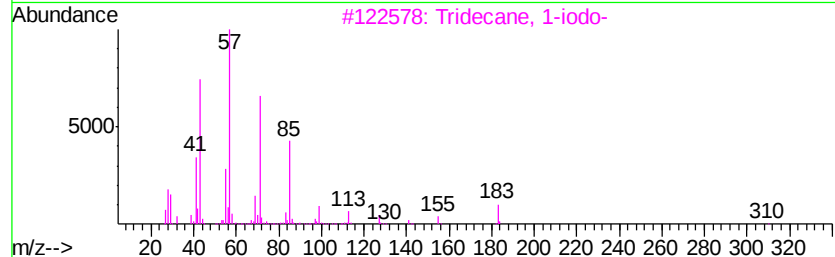
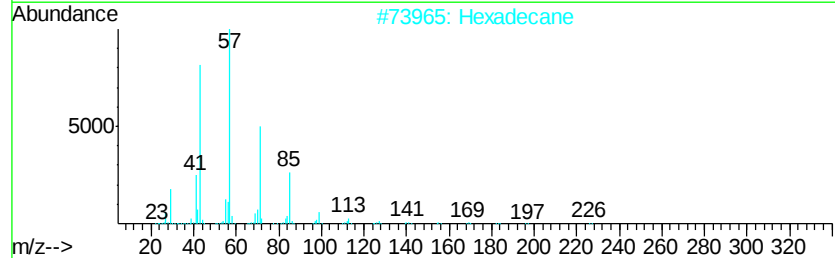
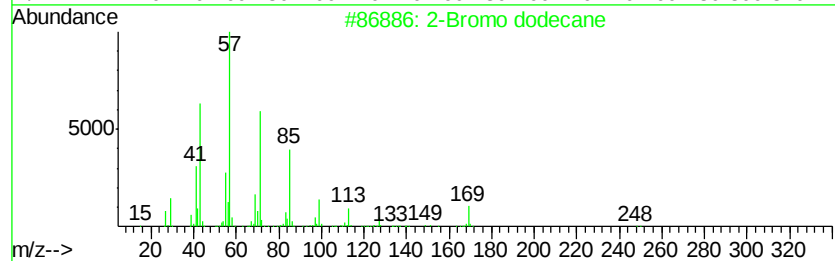
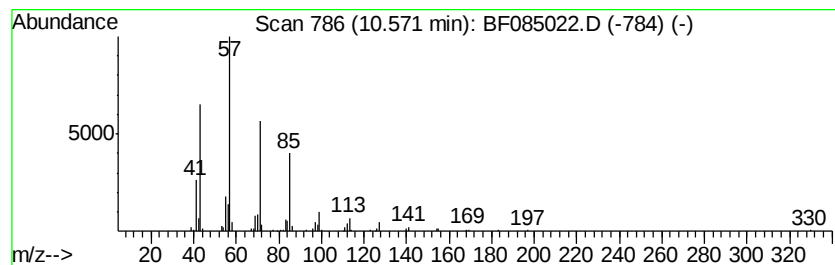
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-Bromo dodecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.57	4.51 ng	305354	Phenanthrene-d10		11.12	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane		248	C12H25Br	013187-99-0	90
2	Hexadecane		226	C16H34	000544-76-3	90
3	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90
4	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	90
5	Eicosane		282	C20H42	000112-95-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
Data File : BF085022.D
Acq On : 11 Feb 2016 17:15
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ClientSampleId :
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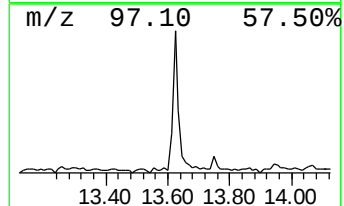
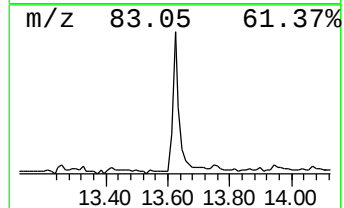
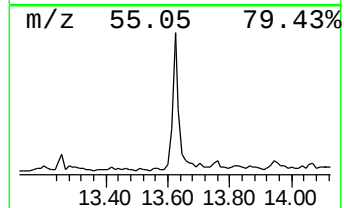
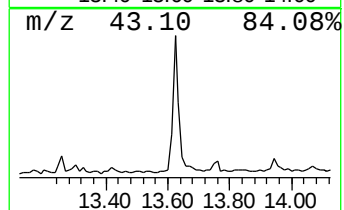
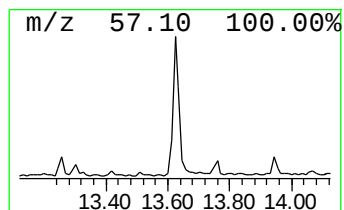
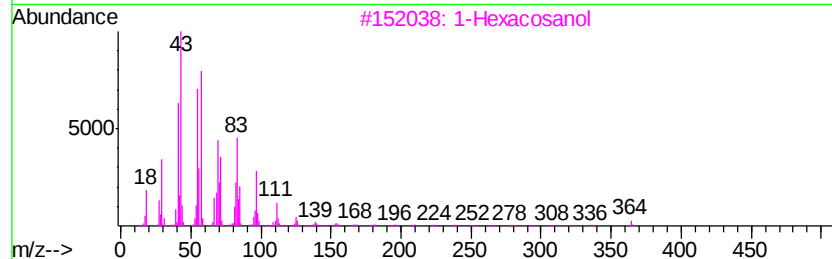
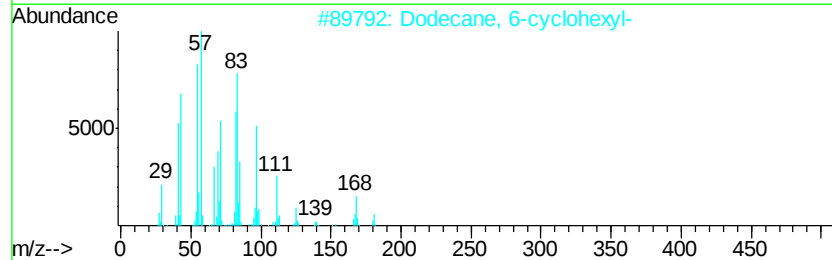
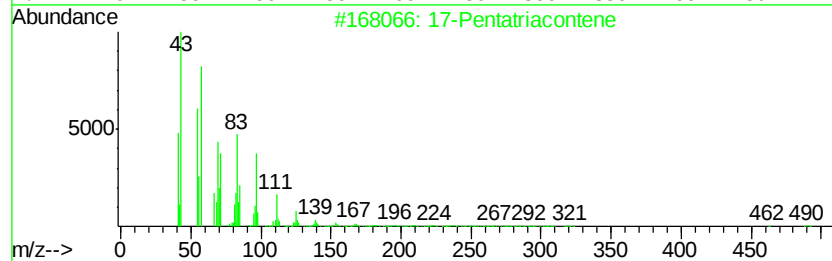
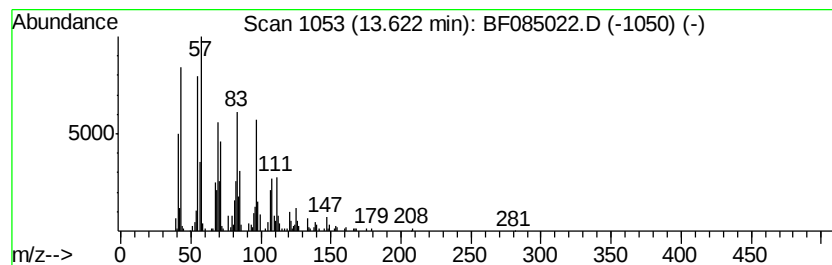
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 17-Pentatriacontene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.62	8.16 ng	465564	Chrysene-d12	13.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-Pentatriacontene	491	C35H70	006971-40-0	89
2		Dodecane, 6-cyclohexyl-	252	C18H36	013151-86-5	68
3		1-Hexacosanol	382	C26H54O	000506-52-5	60
4		1-Tetracosanol	354	C24H50O	000506-51-4	53
5		1-Nonadecene	266	C19H38	018435-45-5	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
Data File : BF085022.D
Acq On : 11 Feb 2016 17:15
Operator : SJ/IZ
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Misc :
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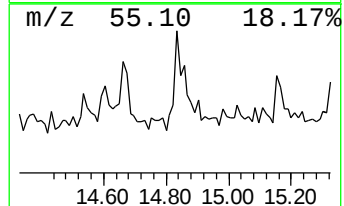
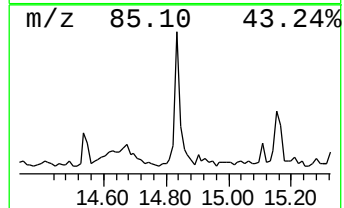
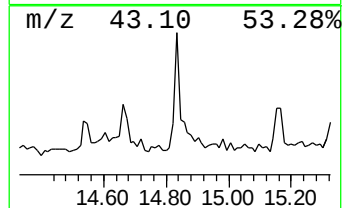
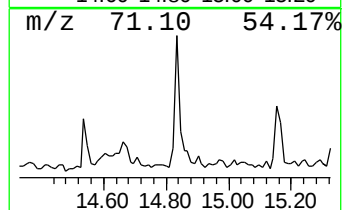
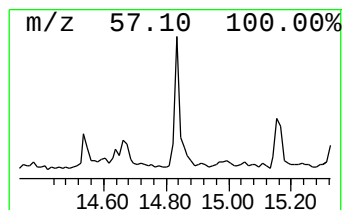
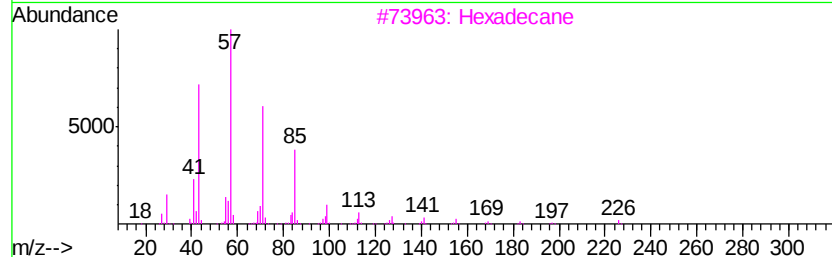
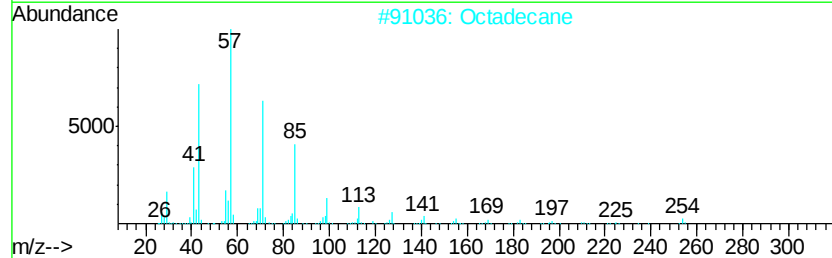
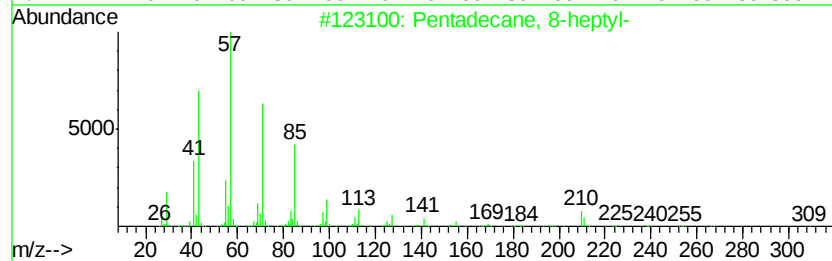
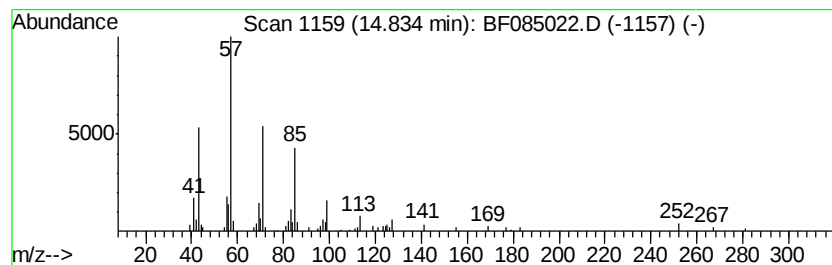
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane, 8-heptyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	2.63 ng	115493	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
2		Octadecane	254	C18H38	000593-45-3	86
3		Hexadecane	226	C16H34	000544-76-3	86
4		Eicosane	282	C20H42	000112-95-8	86
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
Data File : BF085022.D
Acq On : 11 Feb 2016 17:15
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Misc :
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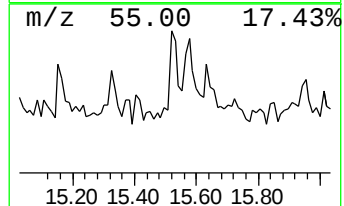
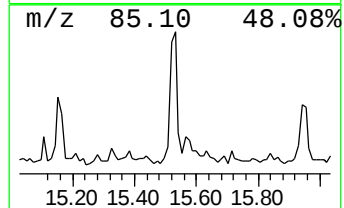
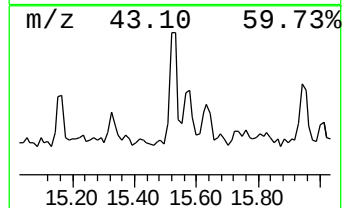
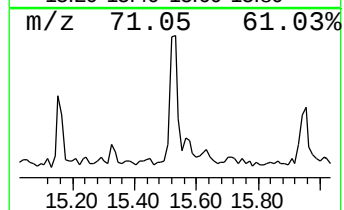
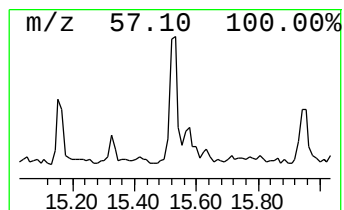
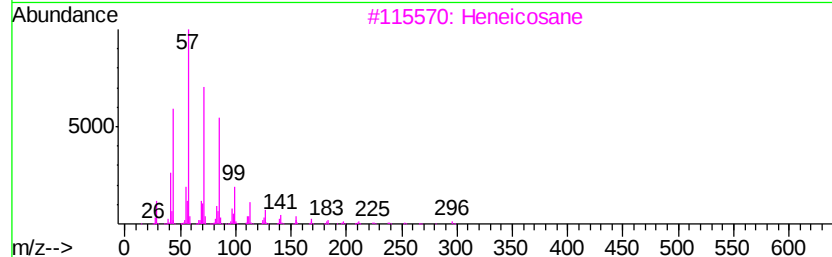
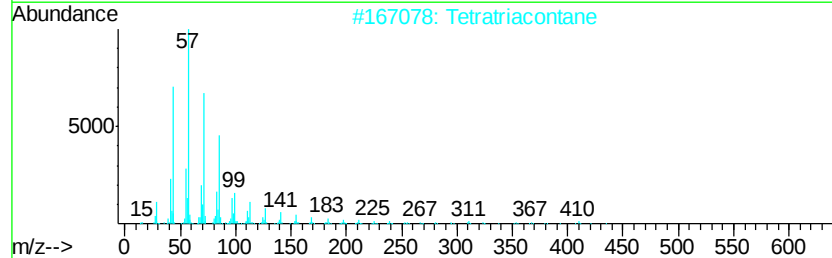
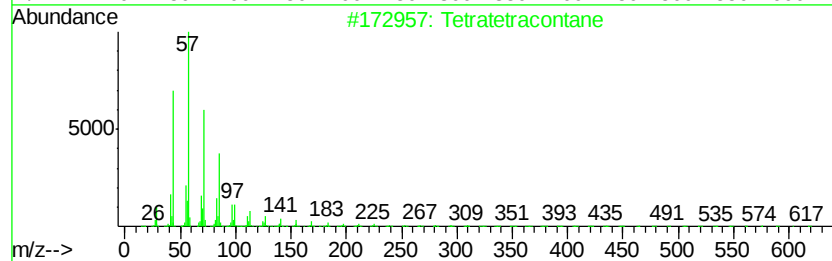
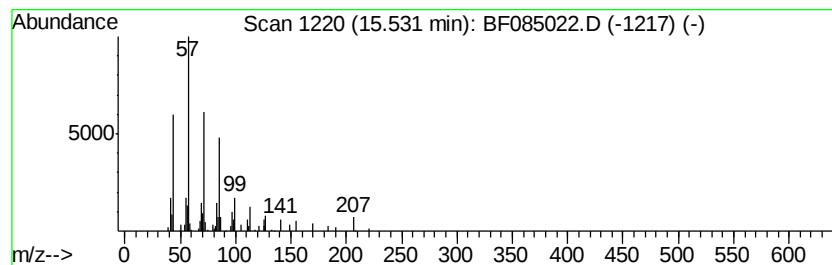
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetratetracontane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	2.06 ng	90170	Perylene-d12	15.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	90
2			Tetratriacontane	479	C34H70	014167-59-0	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Pentacosane	352	C25H52	000629-99-2	87
5			Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
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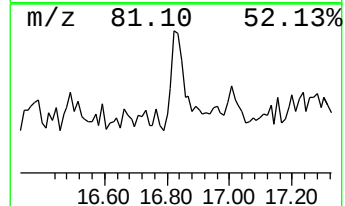
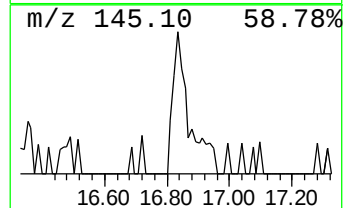
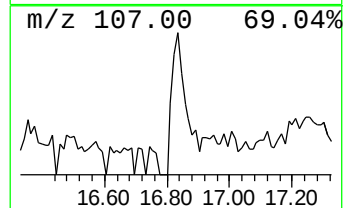
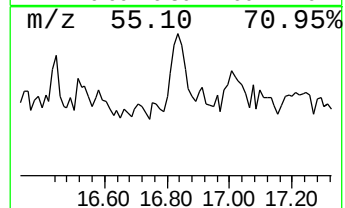
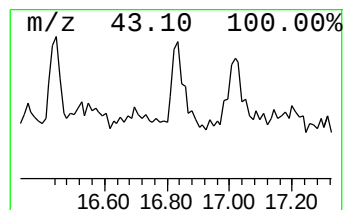
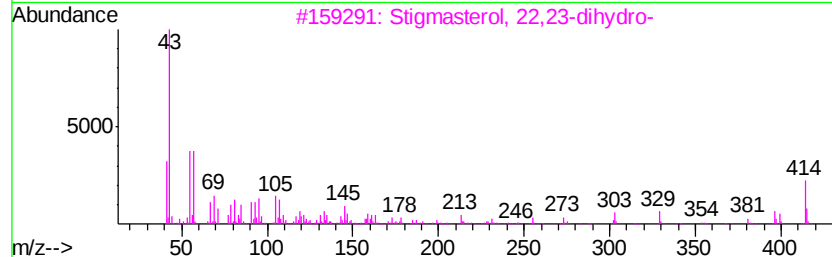
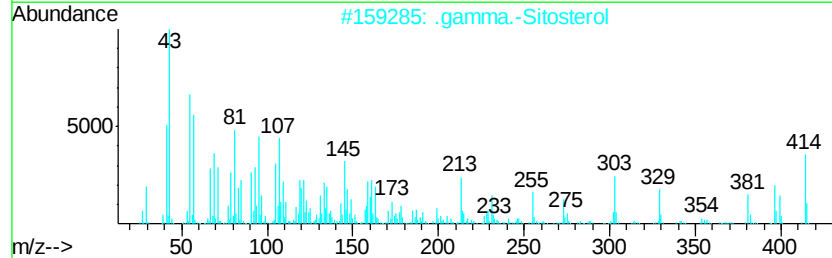
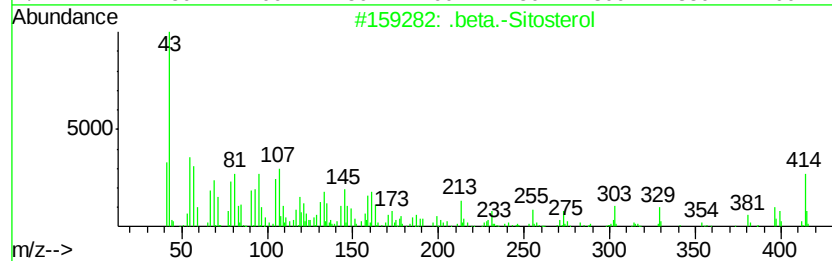
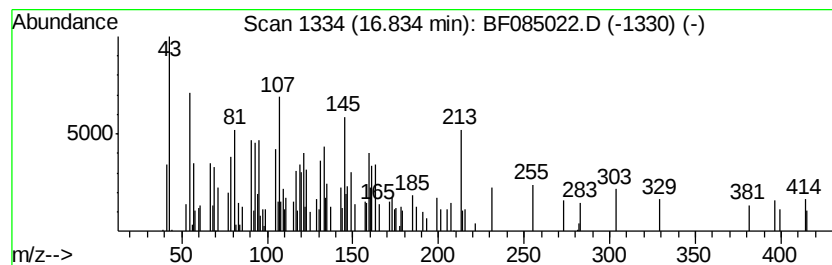
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 .beta.-Sitosterol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.83	4.05 ng	177650	Perylene-d12	15.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.beta.-Sitosterol	414 C29H50O	000083-46-5 95
2		.gamma.-Sitosterol	414 C29H50O	000083-47-6 86
3		Stigmasterol, 22,23-dihydro-	414 C29H50O	1000214-20-7 46
4		Cholest-5-ene, 3-ethoxy-, (3.bet...	414 C29H50O	000986-19-6 17
5		Ethinyl Estradiol	296 C20H24O2	000057-63-6 16



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
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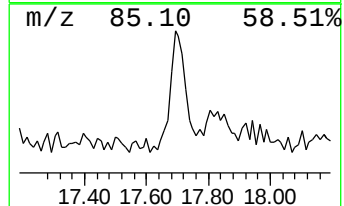
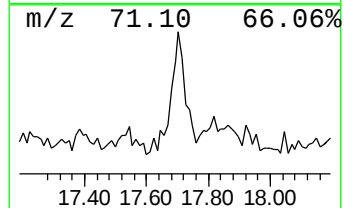
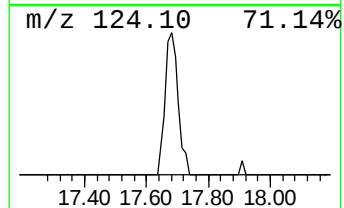
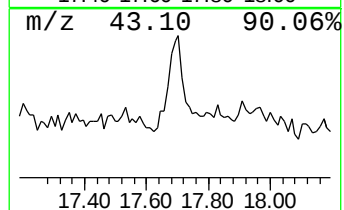
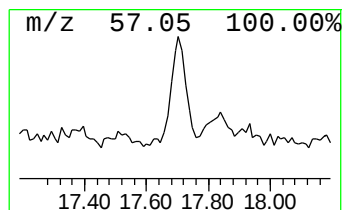
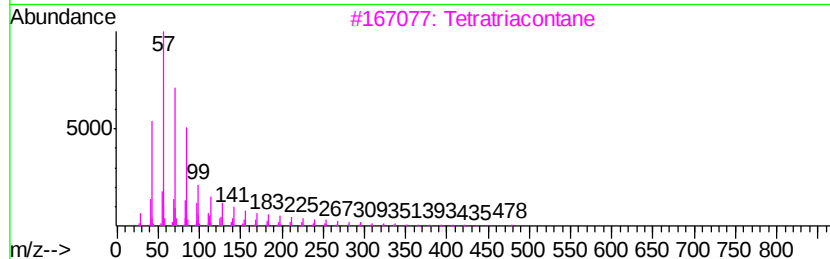
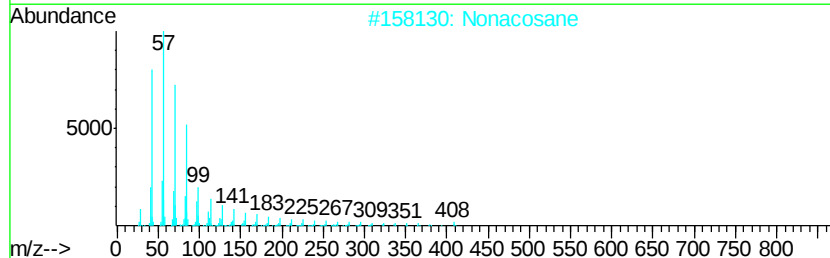
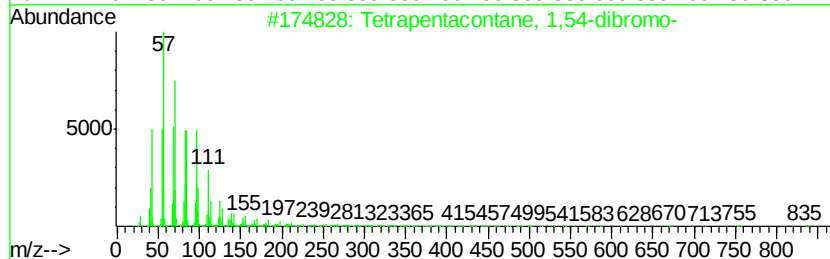
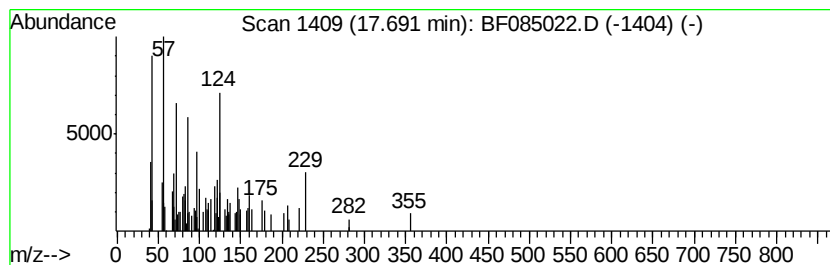
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown17.69 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.69	3.38 ng	148445	Perylene-d12	15.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	35
2		Nonacosane	408	C29H60	000630-03-5	27
3		Tetratriacontane	479	C34H70	014167-59-0	27
4		Tetrapentacontane	759	C54H110	005856-66-6	27
5		Eicosane	282	C20H42	000112-95-8	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Pentanone, 4-hy...	4.74	32.1	ng	1522690	1	6.60	948697	20.0
unknown6.38	6.38	103.6	ng	4912840	1	6.60	948697	20.0
Benzeneacetoneitri...	8.65	5.1	ng	347507	2	7.90	1374580	20.0
Hexadecane	10.09	3.0	ng	222900	3	9.64	1510020	20.0
2-Bromo dodecane	10.57	4.5	ng	305354	4	11.12	1353270	20.0
17-Pentatriacontene	13.62	8.2	ng	465564	5	13.75	1140580	20.0
Pentadecane, 8-he...	14.83	2.6	ng	115493	6	15.11	877199	20.0
Tetratetracontane	15.53	2.1	ng	90170	6	15.11	877199	20.0
.beta.-Sitosterol	16.83	4.0	ng	177650	6	15.11	877199	20.0
unknown17.69	17.69	3.4	ng	148445	6	15.11	877199	20.0