

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF091117\
 Data File : BF098432.D
 Acq On : 12 Sep 2017 5:10
 Operator : SJ/JU
 Sample : I5138-18
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-WM-TS004-01

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-BF091117.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	4.869	469	473	488	rVB	406654	679717	9.69%	1.266%
2	5.287	539	544	547	rBV	6341387	6653448	94.82%	12.388%
3	6.328	715	721	736	rBV	5329522	7017254	100.00%	13.066%
4	6.445	736	741	744	rBV	6405907	6358811	90.62%	11.840%
5	6.669	775	779	789	rBV	1711076	1513029	21.56%	2.817%
6	6.822	801	805	809	rBV	4310256	4211434	60.02%	7.841%
7	7.239	870	876	879	rBV	4152515	4196668	59.80%	7.814%
8	7.269	879	881	888	rVB3	56661	71003	1.01%	0.132%
9	7.951	991	997	1000	rBV	2291260	1972267	28.11%	3.672%
10	9.027	1174	1180	1183	rBV	5174679	4327428	61.67%	8.057%
11	9.110	1190	1194	1198	rBV	107291	104941	1.50%	0.195%
12	9.422	1244	1247	1251	rVB	1626976	1447524	20.63%	2.695%
13	9.639	1281	1284	1288	rVB	75159	70313	1.00%	0.131%
14	9.698	1288	1294	1298	rBV	2073026	1935800	27.59%	3.604%
15	10.416	1411	1416	1422	rBV2	77296	97181	1.38%	0.181%
16	10.492	1424	1429	1432	rBV	3100030	2862475	40.79%	5.330%
17	10.763	1472	1475	1480	rVB2	183903	175566	2.50%	0.327%
18	10.839	1485	1488	1491	rVB	106075	89172	1.27%	0.166%
19	10.886	1493	1496	1499	rBV	137462	121117	1.73%	0.226%
20	10.957	1501	1508	1517	rVB3	114086	189401	2.70%	0.353%
21	11.069	1522	1527	1536	rVB2	113277	132748	1.89%	0.247%
22	11.180	1541	1546	1553	rBV	1940592	1636455	23.32%	3.047%
23	11.251	1553	1558	1561	rBV2	108458	125439	1.79%	0.234%
24	11.592	1611	1616	1619	rVB2	66850	92056	1.31%	0.171%
25	12.263	1725	1730	1732	rBV2	148362	158686	2.26%	0.295%
26	12.280	1732	1733	1739	rVB3	109431	87995	1.25%	0.164%
27	12.427	1755	1758	1764	rVB	125006	118001	1.68%	0.220%
28	12.663	1793	1798	1804	rBV3	51689	79100	1.13%	0.147%
29	12.763	1811	1815	1818	rBV	2923131	2525753	35.99%	4.703%
30	13.662	1964	1968	1972	rBV	274400	310422	4.42%	0.578%
31	13.810	1989	1993	1997	rBV	1182028	1136886	16.20%	2.117%
32	14.286	2071	2074	2083	rVB2	85196	118357	1.69%	0.220%
33	14.886	2173	2176	2184	rBV	299165	340291	4.85%	0.634%
34	15.215	2227	2232	2240	rVB	901692	1145980	16.33%	2.134%

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EME-WM-TS004-01

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-BF091117.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.627	2298	2302	2307	rBV	298792	424742	6.05%	0.791%
36	16.615	2466	2470	2477	rVB	126990	226892	3.23%	0.422%
37	17.080	2541	2549	2557	rVB9	134266	404027	5.76%	0.752%
38	17.468	2611	2615	2622	rVB5	59288	121066	1.73%	0.225%
39	17.774	2662	2667	2683	rVB5	106212	427814	6.10%	0.797%

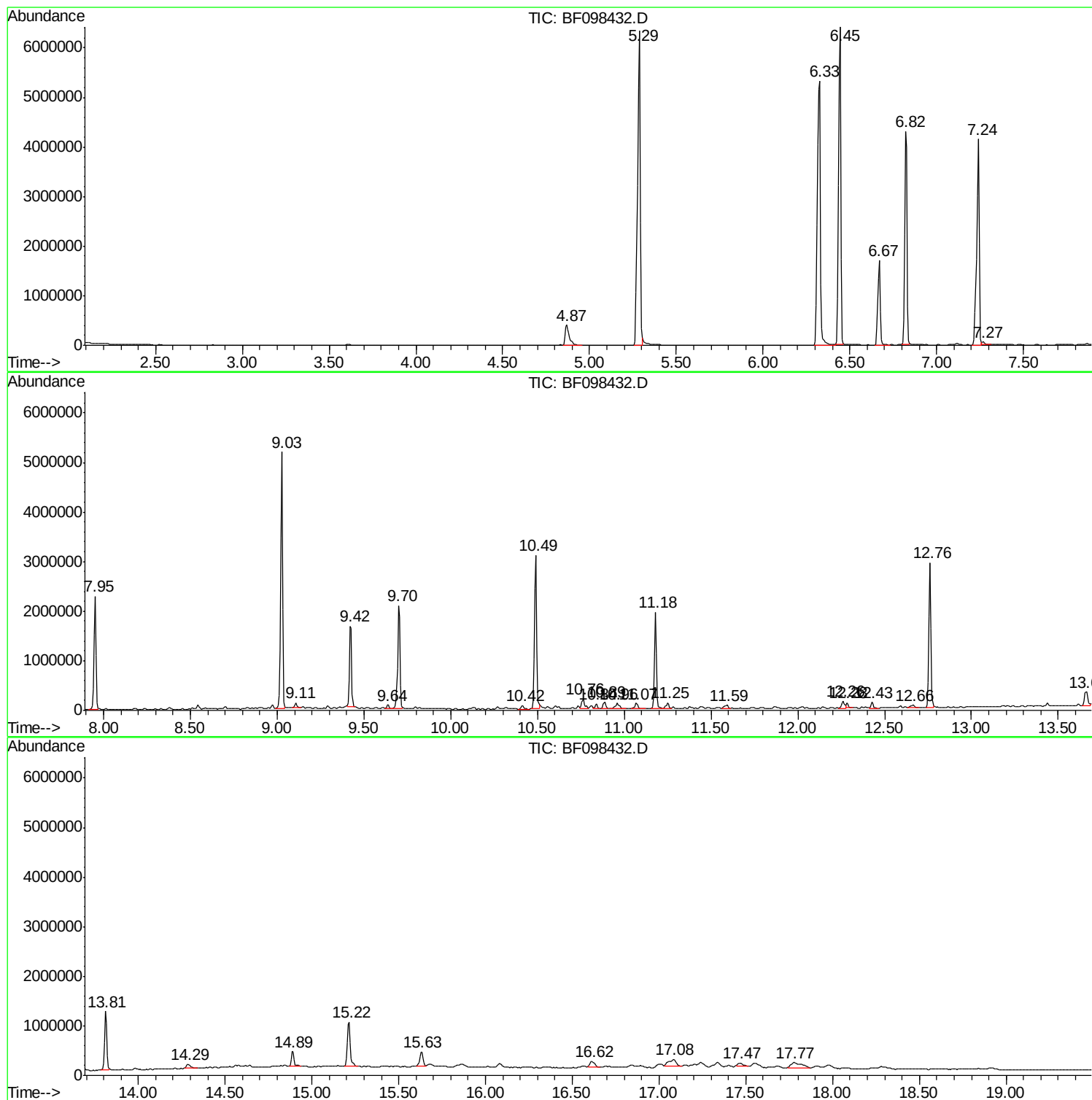
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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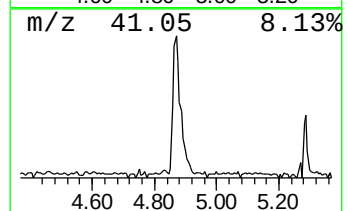
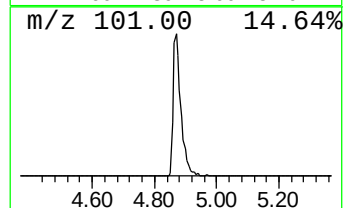
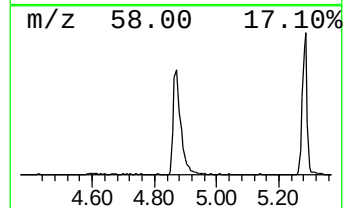
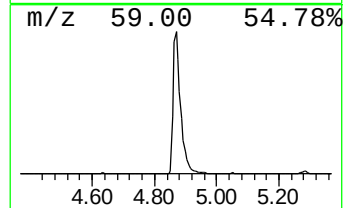
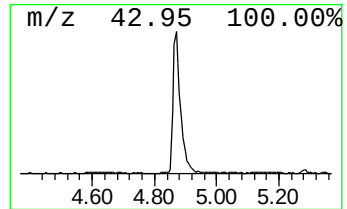
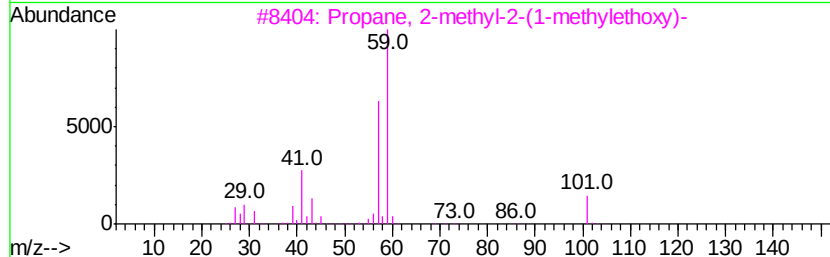
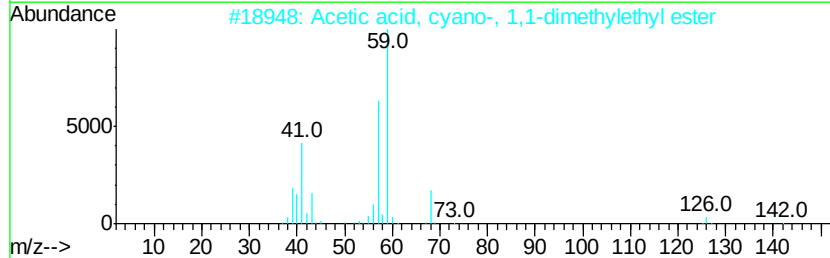
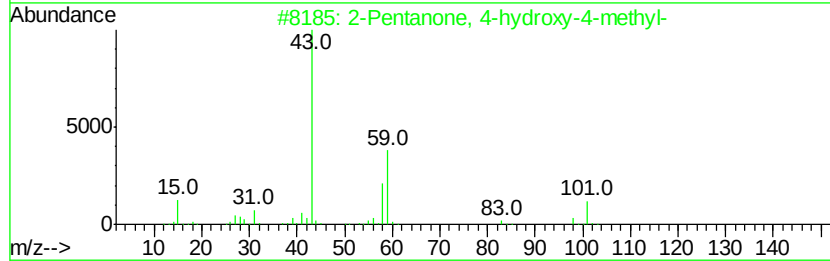
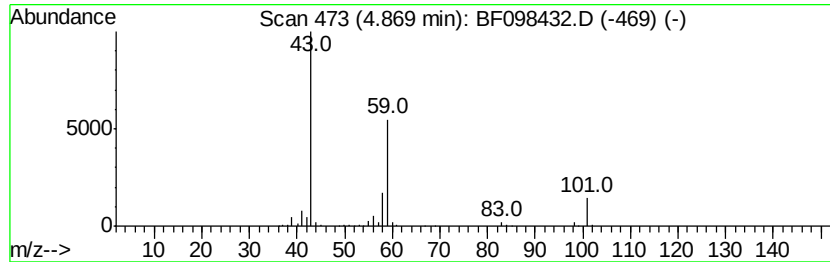
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
4.87	8.98 ng	679717	1,4-Dichlorobenzene-d4	6.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	37
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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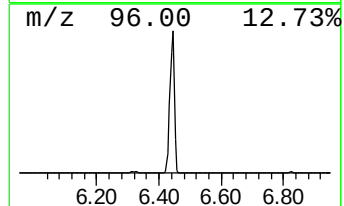
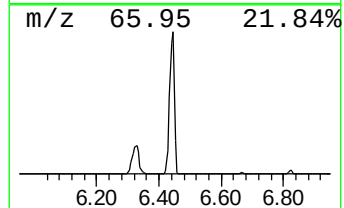
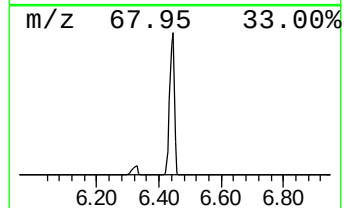
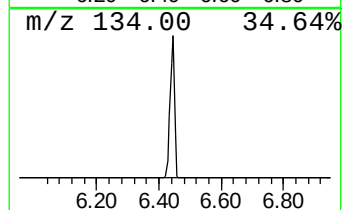
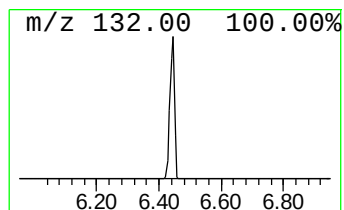
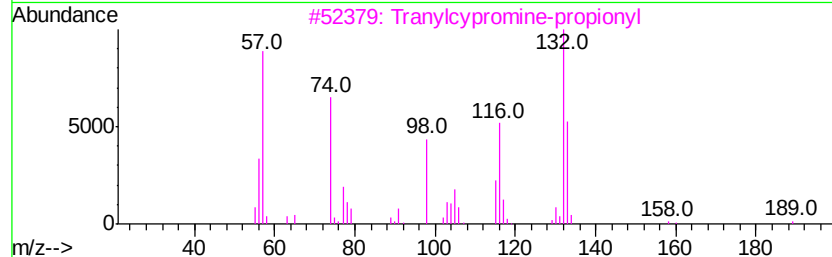
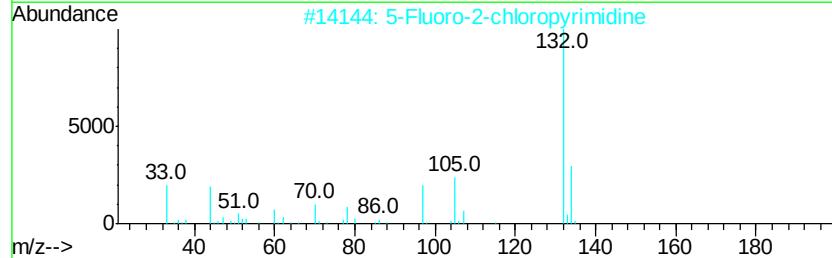
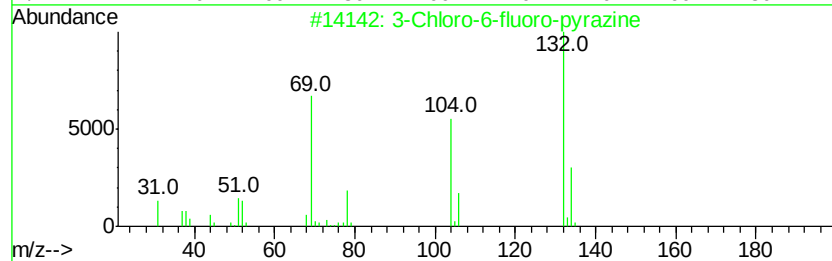
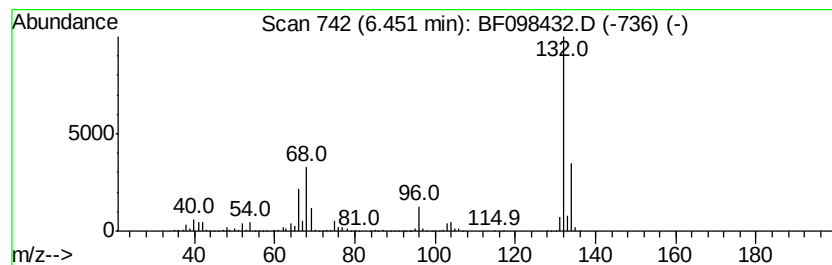
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	84.05 ng	6358810	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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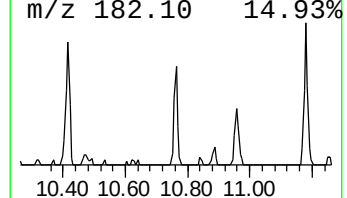
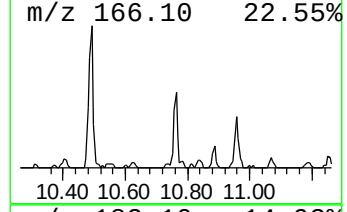
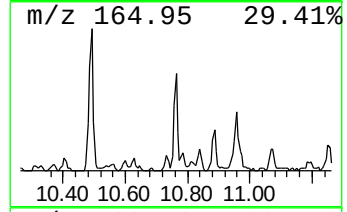
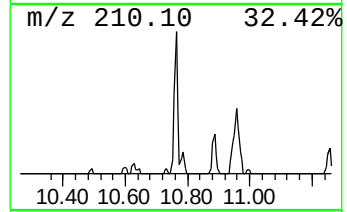
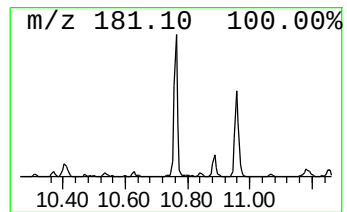
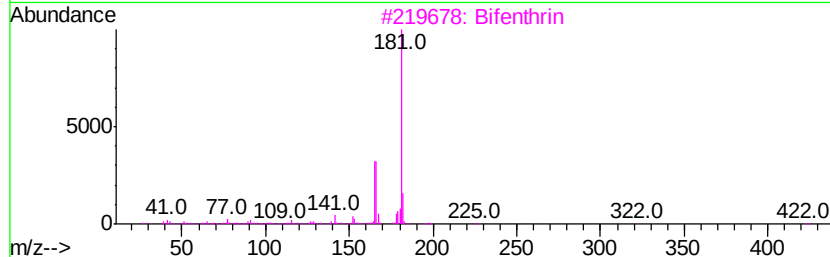
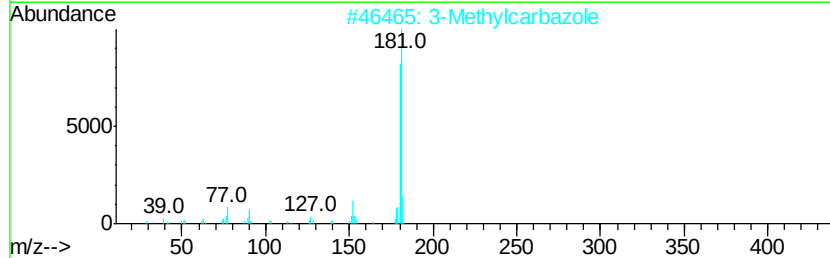
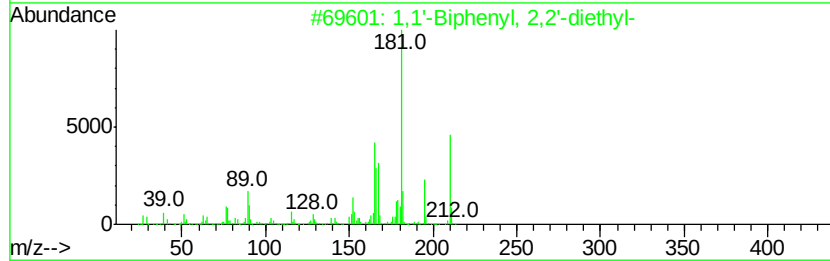
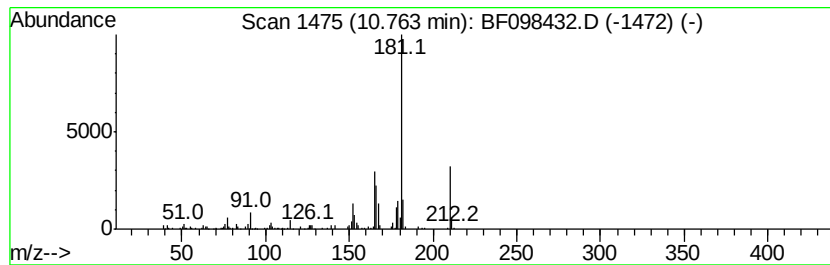
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 2,2'-diethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.76	2.15 ng	175566	Phenanthrene-d10	11.18	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-diethyl-	210	C16H18	013049-35-9	76
2	3-Methylcarbazole	181	C13H11N	004630-20-0	60
3	Bifenthrin	422	C23H22ClF3O2	082657-04-3	59
4	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	55
5	2,2-Diphenylpropylamine	211	C15H17N	034611-07-9	52



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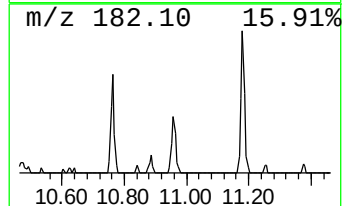
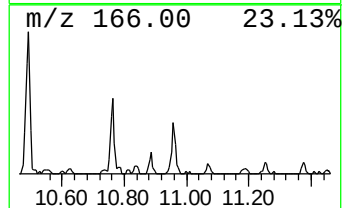
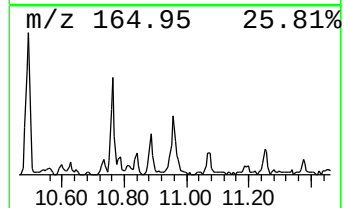
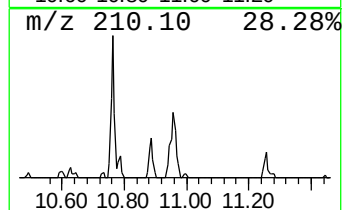
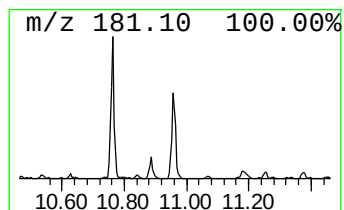
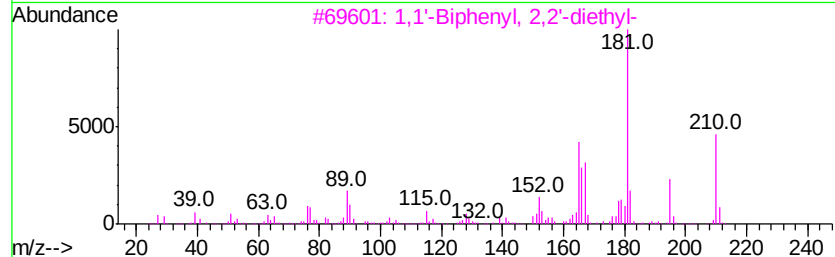
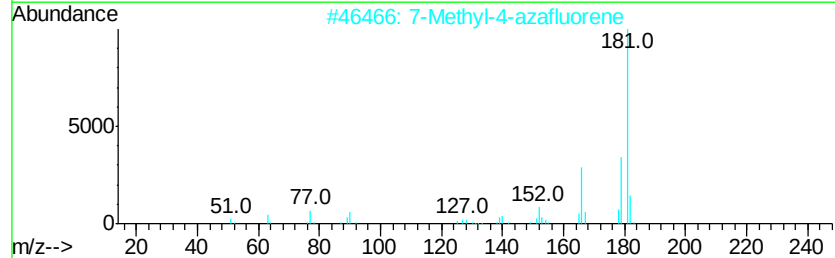
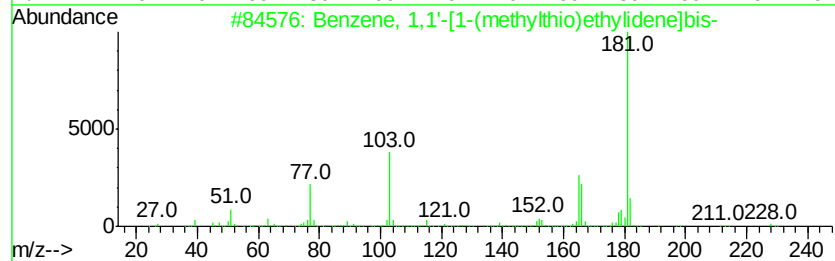
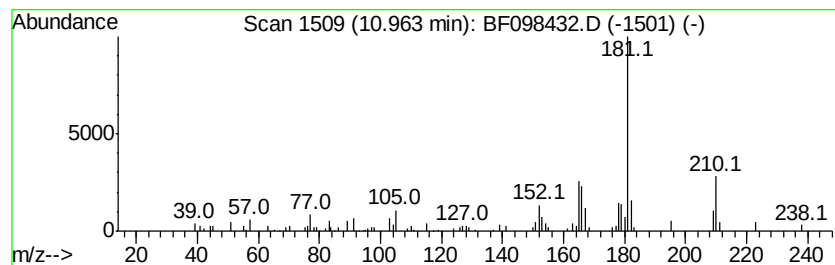
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Peak Number 5 Benzene, 1,1'-[1-(methylthio)eth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.96	2.31 ng	189401	Phenanthrene-d10	11.18		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,1'-[1-(methylthio)eth...	228	C15H16S	054851-63-7	72
2		7-Methyl-4-azafluorene	181	C13H11N	064291-99-2	70
3		1,1'-Biphenyl, 2,2'-diethyl-	210	C16H18	013049-35-9	70
4		Benzo[b]benzo[3,4]cyclobuta[1,2-...	210	C14H10O2	042896-18-4	60
5		Bifenthrin	422	C23H22ClF3O2	082657-04-3	59



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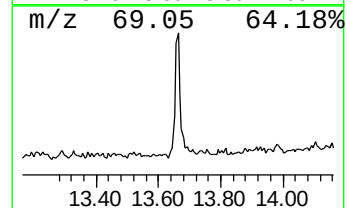
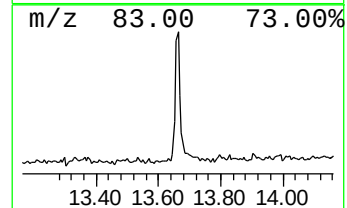
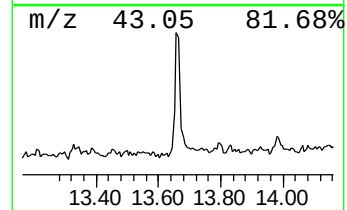
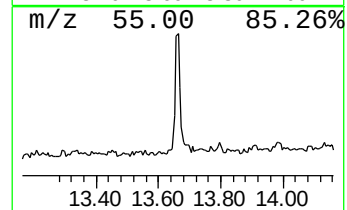
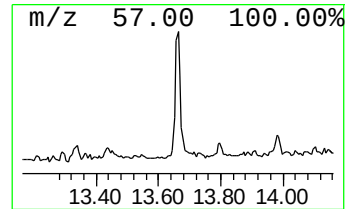
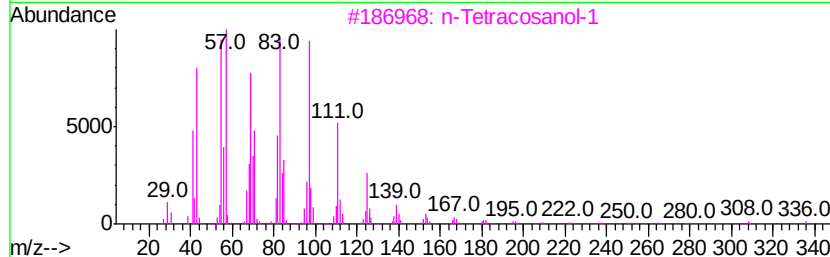
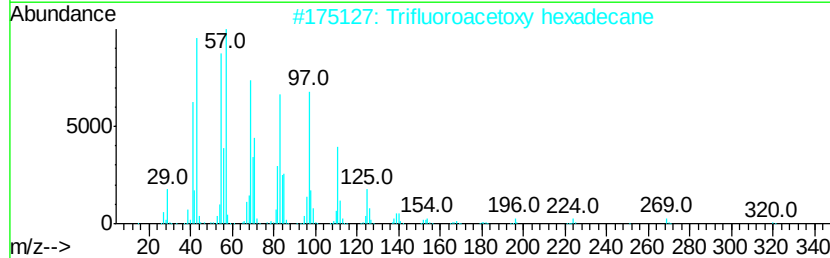
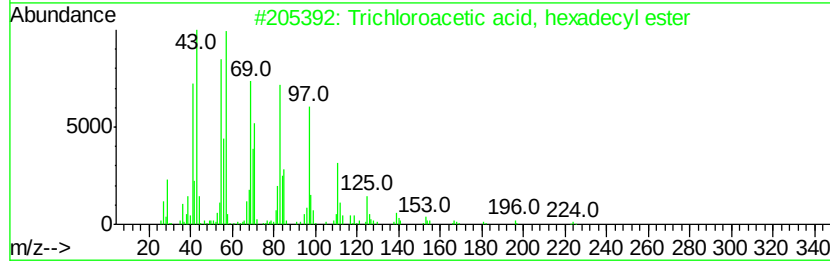
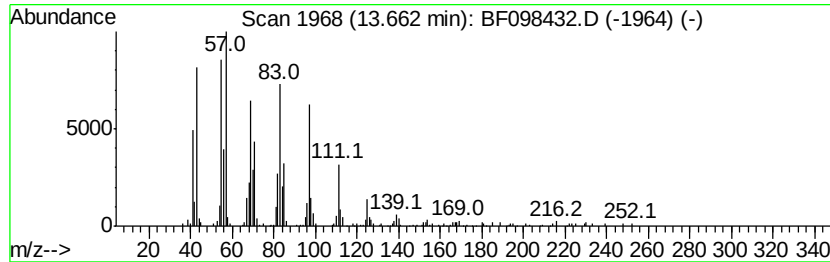
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Trichloroacetic acid, hexad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.66	5.46 ng	310422	Chrysene-d12	13.81	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2	Trifluoroacetoxyl hexadecane	338	C18H33F3O2	006222-03-3	91
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
4	Pentafluoropropionic acid, octadecyl ...	416	C21H37F5O2	959261-25-7	91
5	Pentafluoropropionic acid, pentadecyl ...	374	C18H31F5O2	959092-08-1	91



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Acq On : 12 Sep 2017 5:10
Operator : SJ/JU
Sample : I5138-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

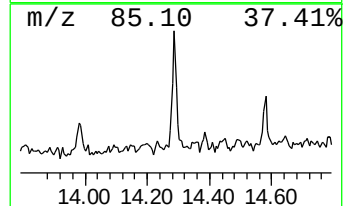
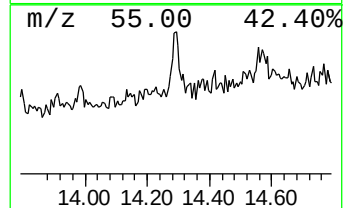
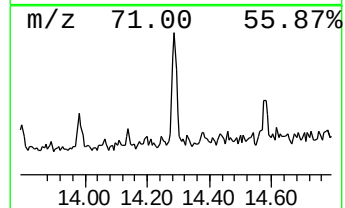
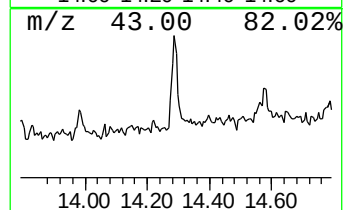
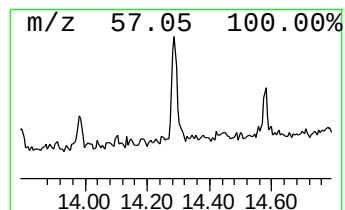
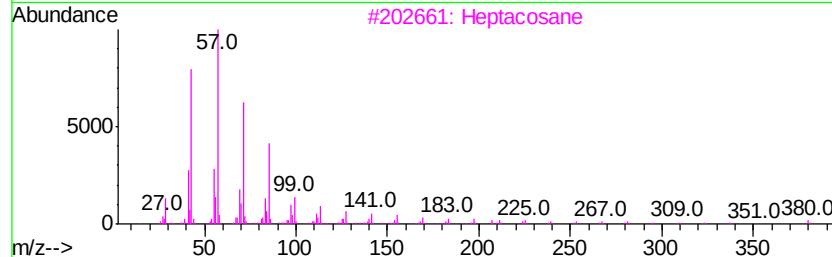
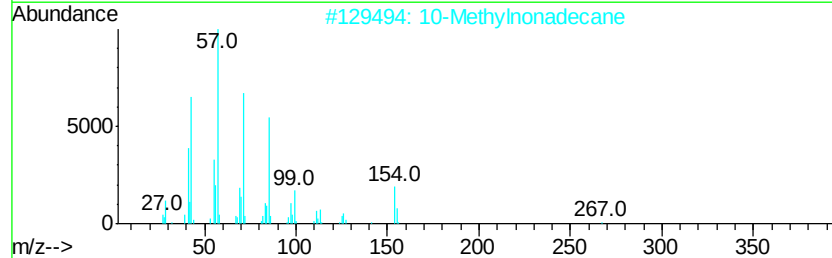
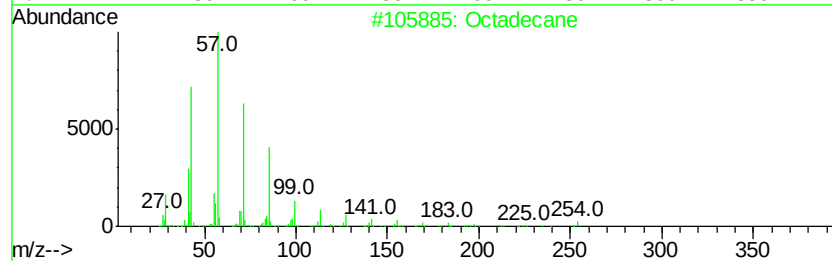
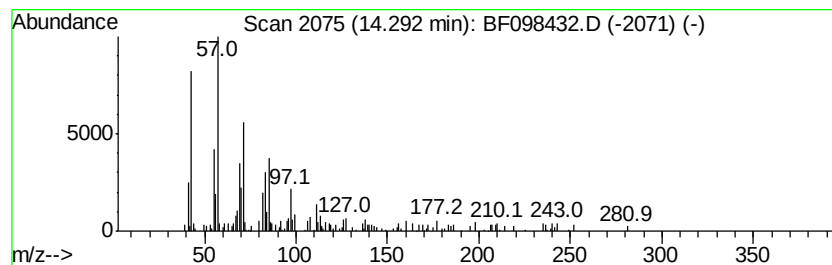
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Octadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	2.08 ng	118357	Chrysene-d12	13.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octadecane	254 C18H38	000593-45-3	93
2	10-Methylnonadecane	282 C20H42	056862-62-5	87
3	Heptacosane	380 C27H56	000593-49-7	87
4	Sulfurous acid, 2-propyl tridecy...	306 C16H34O3S	1000309-12-4	86
5	2-methyloctacosane	408 C29H60	1000376-72-8	83



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091117\
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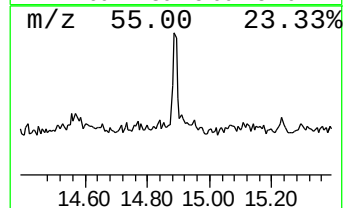
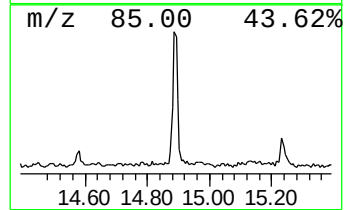
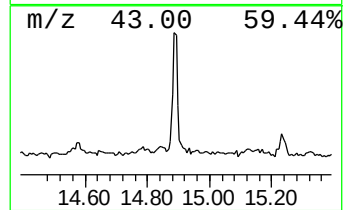
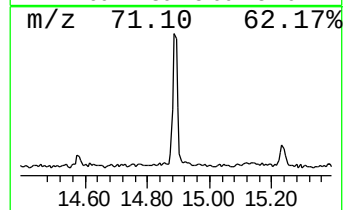
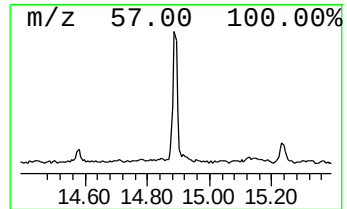
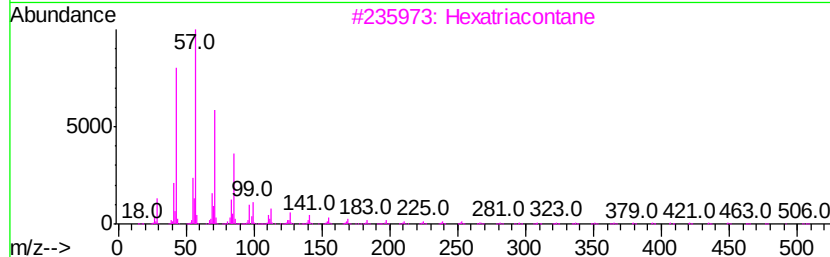
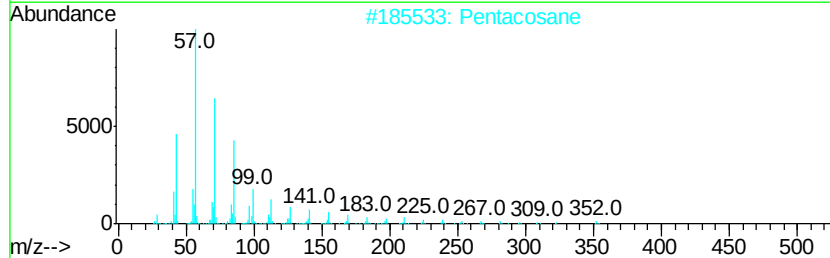
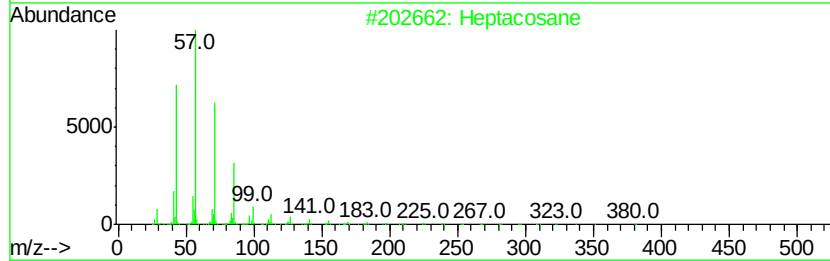
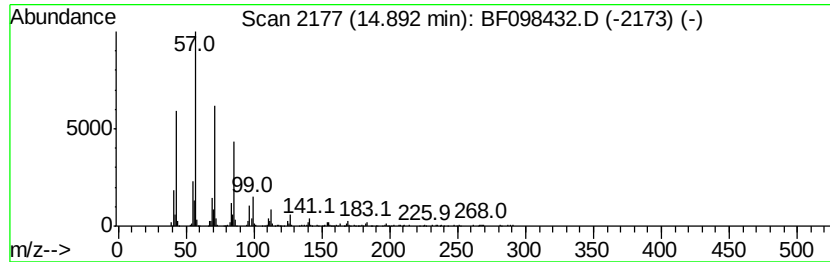
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.89	5.94 ng	340291	Perylene-d12		15.22
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Pentacosane	352	C25H52	000629-99-2	90
3	Hexatriacontane	507	C36H74	000630-06-8	90
4	Tetratetracontane	619	C44H90	007098-22-8	90
5	Octadecane	254	C18H38	000593-45-3	86



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091117\
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 Acq On : 12 Sep 2017 5:10
 Operator : SJ/JU
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 ClientSampleId :
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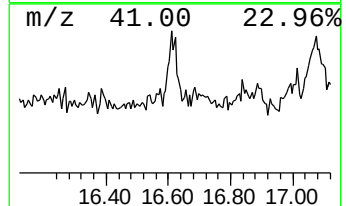
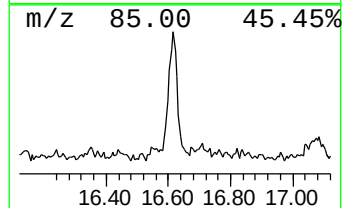
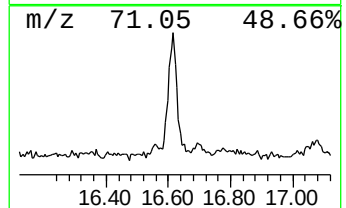
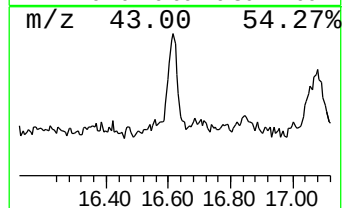
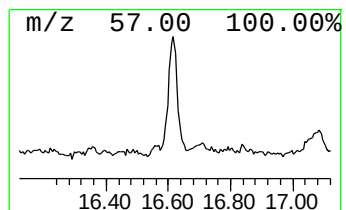
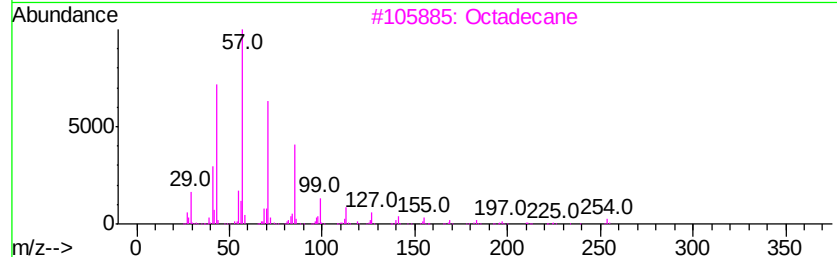
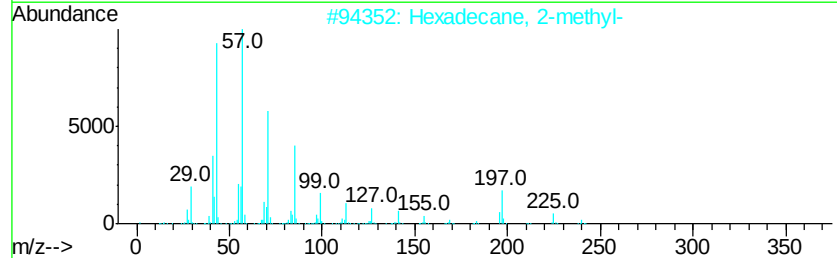
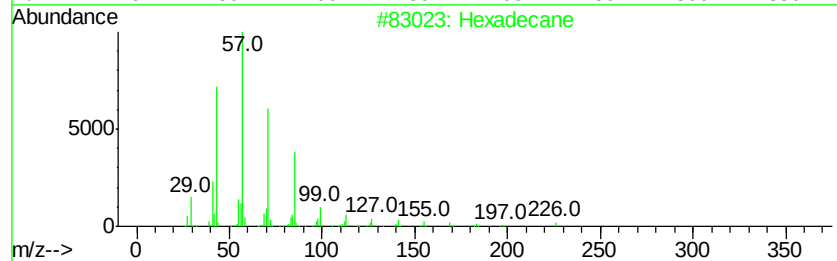
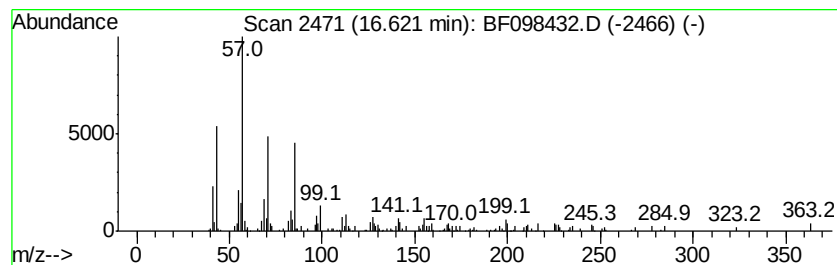
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 10 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.62	3.96 ng	226892	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Hexadecane, 2-methyl-	240	C17H36	001560-92-5	86
3		Octadecane	254	C18H38	000593-45-3	83
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	83
5		Tetratetracontane	619	C44H90	007098-22-8	83



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
Data File : BF098432.D
Acq On : 12 Sep 2017 5:10
Operator : SJ/JU
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Misc :
ALS Vial : 22 Sample Multiplier: 1

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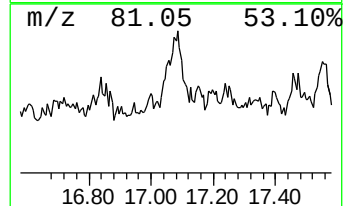
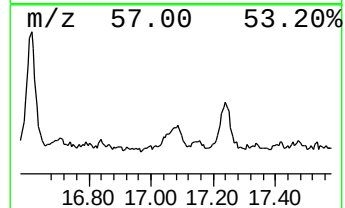
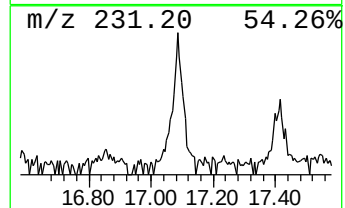
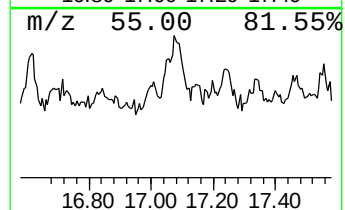
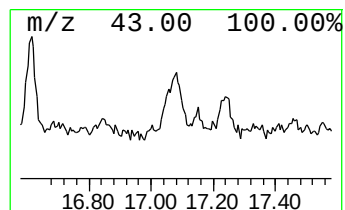
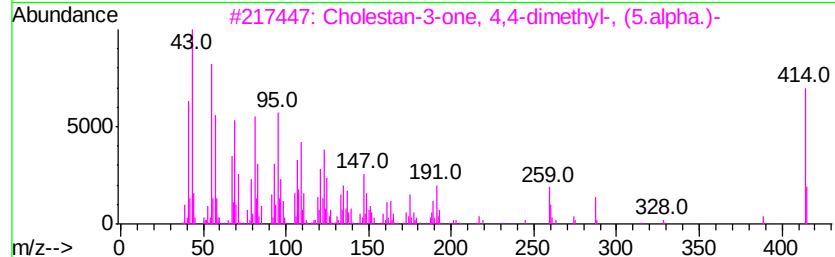
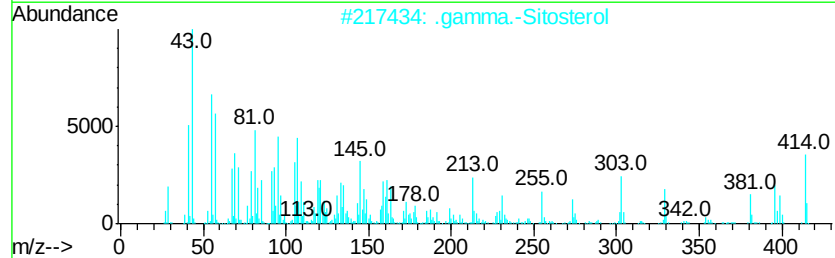
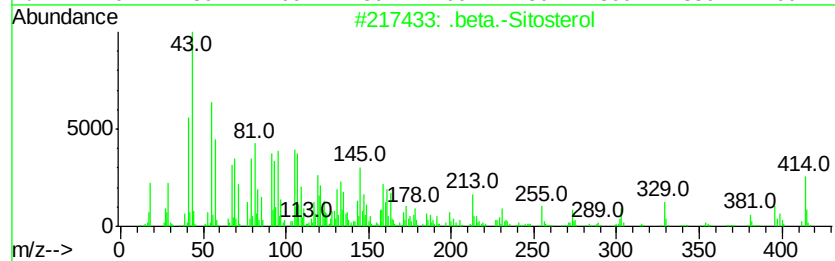
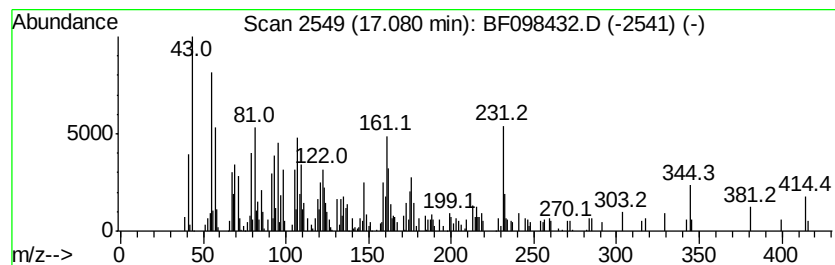
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown17.08 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.08	7.05 ng	404027	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-	Sitosterol	414	C29H50O	000083-46-5	38
2	.	gamma.-	Sitosterol	414	C29H50O	000083-47-6	38
3	Cholestan-3-one, 4,4-dimethyl-, ...	414	C29H50O	002097-85-0	25		
4	1H-Cyclopropylazulen-4-ol, deca...	222	C15H26O	000552-02-3	25		
5	17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	25		



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
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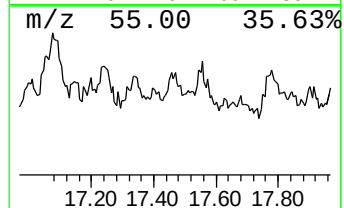
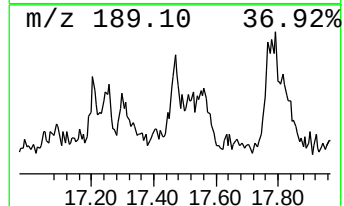
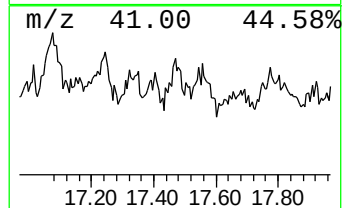
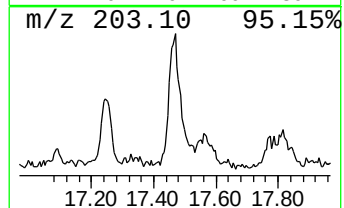
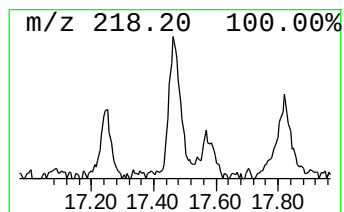
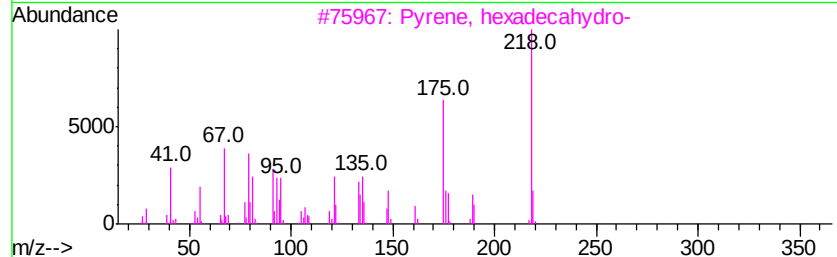
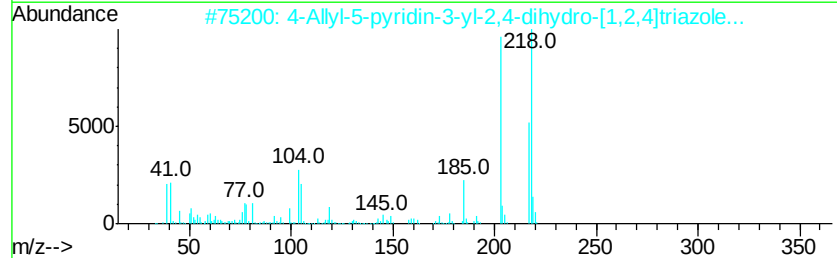
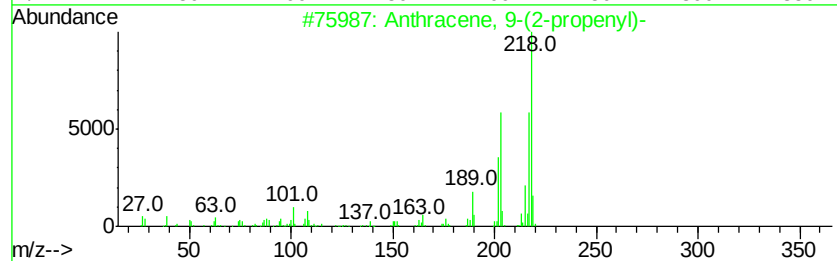
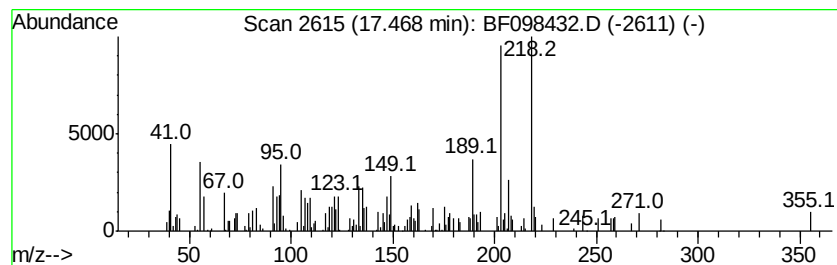
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Anthracene, 9-(2-propenyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.47	2.11 ng	121066	Perylene-d12	15.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	53
2		4-Allyl-5-pyridin-3-yl-2,4-dihyd...	218	C10H10N4S	1000300-40-5	49
3		Pvrene, hexadecahydro-	218	C16H26	002435-85-0	43
4		Thieno[2,3-b]pyridine-2-carbonit...	218	C10H10N4S	147992-88-9	30
5		7b-Phenyl-2a,7b-dihydro-3H-cyclo...	218	C17H14	138089-70-0	30



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
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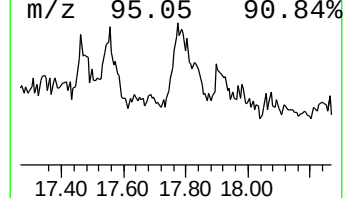
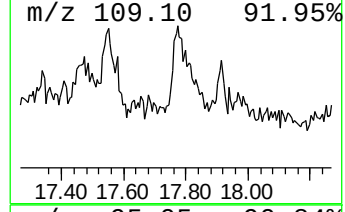
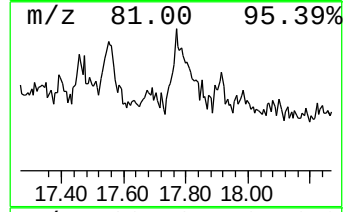
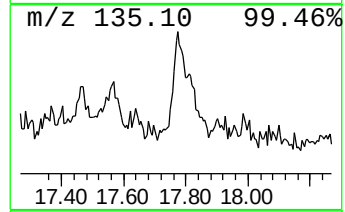
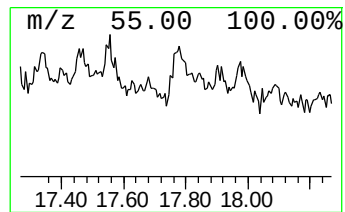
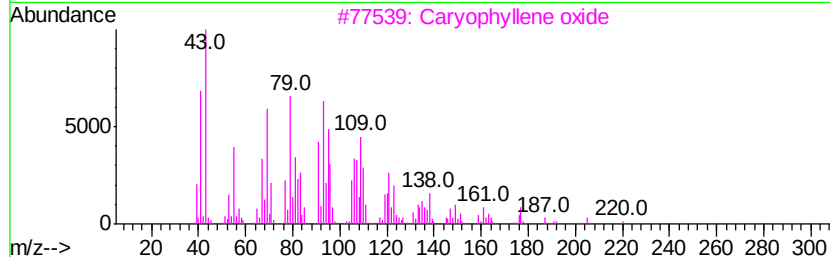
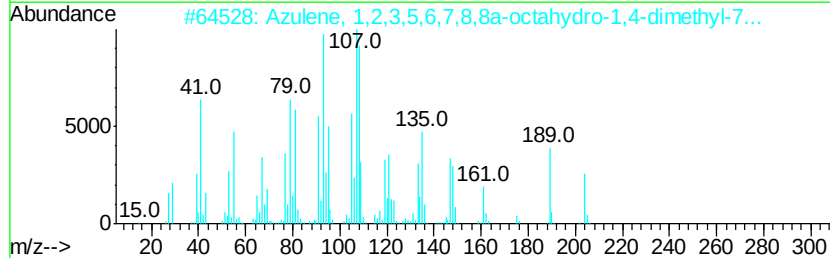
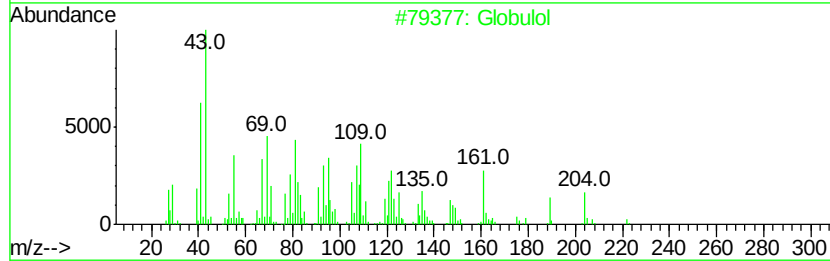
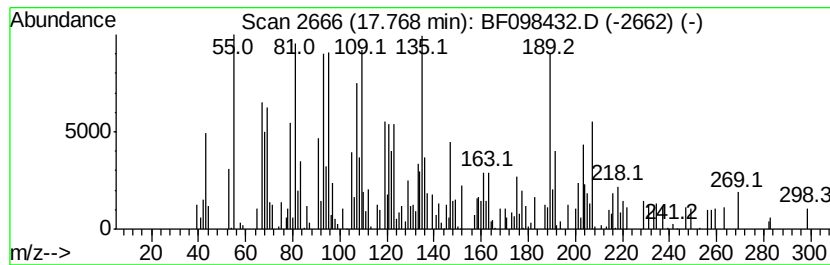
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown17.77 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.77	7.47 ng	427814	Perylene-d12	15.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Globulol	222 C15H26O	051371-47-2 41
2		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204 C15H24	003691-11-0 30
3		Carvophyllene oxide	220 C15H24O	001139-30-6 30
4		.alpha.-Cadinol	222 C15H26O	000481-34-5 25
5		Cyclohexanone, 5-methyl-2-(1-met...	152 C10H16O	029606-79-9 25



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF091117\
Data File : BF098432.D
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Operator : SJ/JU
Sample : I5138-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EME-WM-TS004-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.87	9.0 ng		679717	1	6.67	1513030	20.0
unknown6.45	6.45	84.0 ng		6358810	1	6.67	1513030	20.0
1,1'-Biphenyl, 2,...	10.76	2.1 ng		175566	4	11.18	1636460	20.0
Benzene, 1,1'-[1-...	10.96	2.3 ng		189401	4	11.18	1636460	20.0
Trichloroacetic a...	13.66	5.5 ng		310422	5	13.81	1136890	20.0
Octadecane	14.29	2.1 ng		118357	5	13.81	1136890	20.0
Heptacosane	14.89	5.9 ng		340291	6	15.22	1145980	20.0
Hexadecane	16.62	4.0 ng		226892	6	15.22	1145980	20.0
unknown17.08	17.08	7.0 ng		404027	6	15.22	1145980	20.0
Anthracene, 9-(2-...	17.47	2.1 ng		121066	6	15.22	1145980	20.0
unknown17.77	17.77	7.5 ng		427814	6	15.22	1145980	20.0