

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100918\
 Data File : BF109706.D
 Acq On : 9 Oct 2018 20:27
 Operator : JU/SJ
 Sample : J5331-02 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM-STONE-DRUM-3

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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	5.316	546	549	565	rBV2	28015	78783	8.84%	1.255%
2	6.351	722	725	736	rBV2	33971	103265	11.59%	1.644%
3	6.469	742	745	758	rBV	93652	140507	15.76%	2.237%
4	6.698	780	784	806	rBV	266761	416352	46.71%	6.630%
5	6.851	806	810	821	rVB	120708	144030	16.16%	2.294%
6	7.269	877	881	897	rBV2	20574	74868	8.40%	1.192%
7	7.975	997	1001	1009	rBV	434388	489346	54.90%	7.792%
8	8.269	1047	1051	1054	rBV3	12391	14977	1.68%	0.238%
9	8.298	1054	1056	1059	rBV4	12498	12935	1.45%	0.206%
10	8.422	1073	1077	1083	rVB7	17795	36984	4.15%	0.589%
11	8.481	1083	1087	1090	rBV4	22054	28578	3.21%	0.455%
12	8.522	1090	1094	1095	rBV3	17694	21420	2.40%	0.341%
13	8.604	1104	1108	1111	rBV4	21672	33910	3.80%	0.540%
14	8.639	1111	1114	1118	rVV7	15802	26855	3.01%	0.428%
15	8.722	1125	1128	1130	rBV3	24741	27335	3.07%	0.435%
16	8.751	1130	1133	1136	rVV4	22491	34583	3.88%	0.551%
17	8.798	1139	1141	1145	rVV4	29194	40420	4.53%	0.644%
18	8.857	1149	1151	1154	rVB3	29311	26409	2.96%	0.421%
19	8.969	1167	1170	1172	rBV	36157	40503	4.54%	0.645%
20	9.004	1172	1176	1179	rVB5	39715	53592	6.01%	0.853%
21	9.051	1179	1184	1187	rBV	183429	237380	26.63%	3.780%
22	9.363	1234	1237	1241	rBV5	69291	105060	11.79%	1.673%
23	9.457	1251	1253	1256	rBV	66013	63927	7.17%	1.018%
24	9.592	1273	1276	1278	rBV3	98920	115473	12.95%	1.839%
25	9.639	1281	1284	1287	rVB3	208959	212263	23.81%	3.380%
26	9.722	1293	1298	1305	rBV	549891	891344	100.00%	14.194%
27	10.133	1365	1368	1371	rBV	175080	195343	21.92%	3.111%
28	10.386	1409	1411	1414	rVB	163341	149644	16.79%	2.383%
29	10.651	1454	1456	1462	rVB	311789	288182	32.33%	4.589%
30	11.204	1548	1550	1556	rVB	305671	313704	35.19%	4.995%
31	12.780	1816	1818	1827	rBV3	197127	290992	32.65%	4.634%
32	13.827	1993	1996	2005	rVB	432391	512091	57.45%	8.154%
33	14.598	2124	2127	2137	rVB	66177	94923	10.65%	1.512%
34	14.904	2176	2179	2186	rVB	81864	99319	11.14%	1.582%

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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.227	2230	2234	2246	rVB2	311378	581250	65.21%	9.256%
36	15.651	2301	2306	2311	rVB	89526	133162	14.94%	2.120%
37	16.104	2378	2383	2391	rVB2	54504	103165	11.57%	1.643%
38	16.639	2470	2474	2483	rVB5	22679	47044	5.28%	0.749%

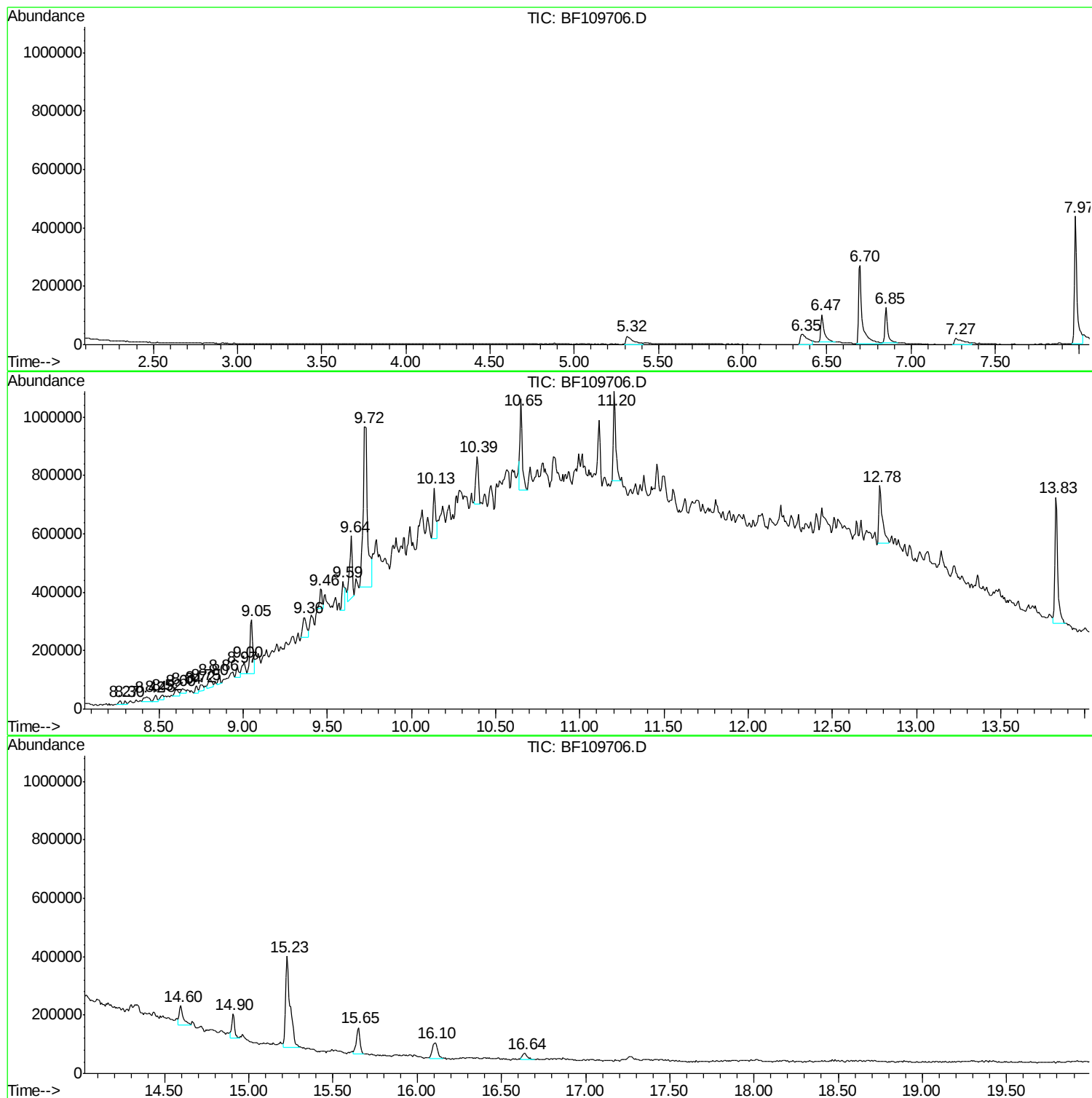
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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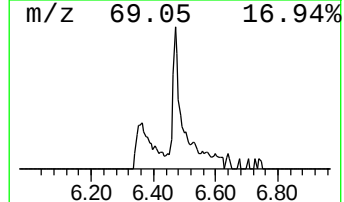
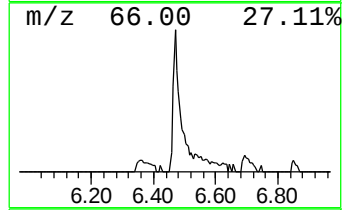
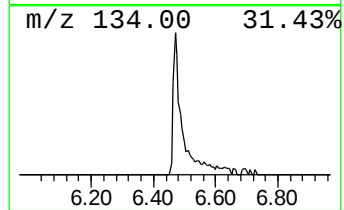
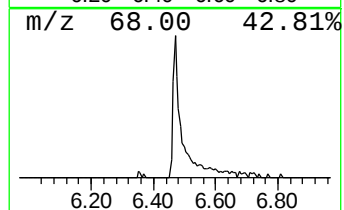
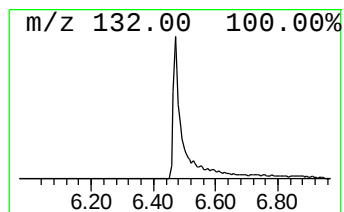
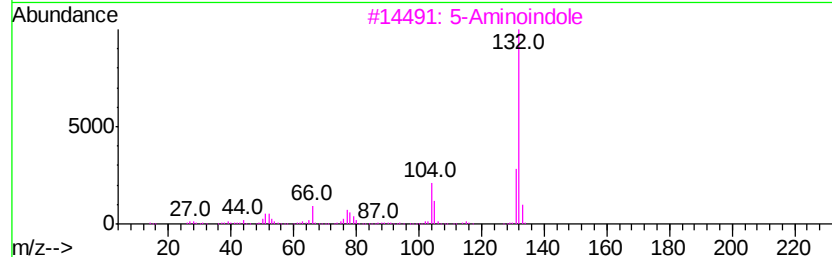
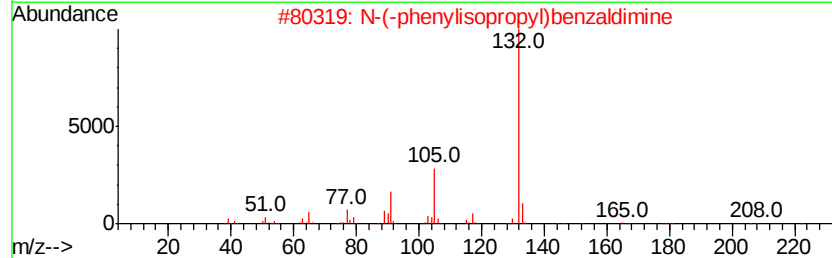
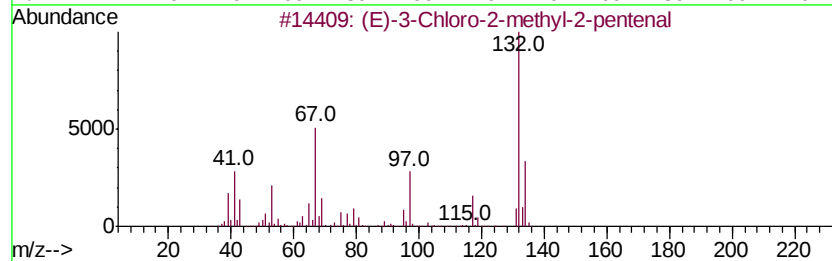
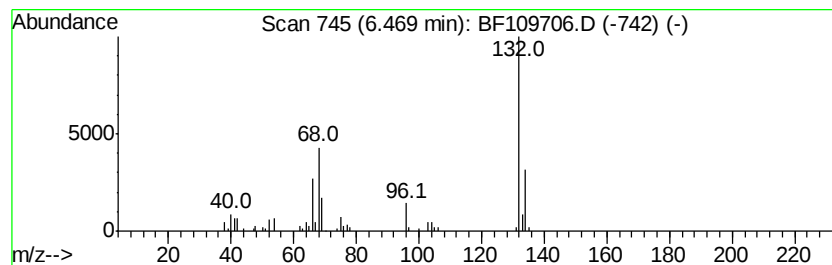
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	6.75 ng	140507	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		Benzeneethanamine, .alpha.-methy...	223	C16H17N	002980-02-1	10
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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ALS Vial : 10 Sample Multiplier: 1

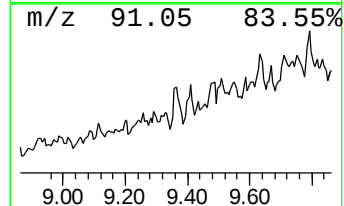
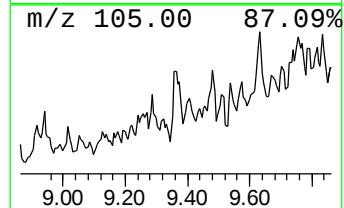
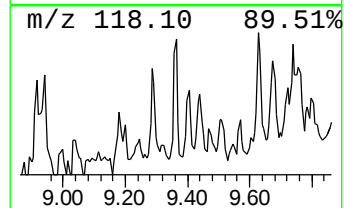
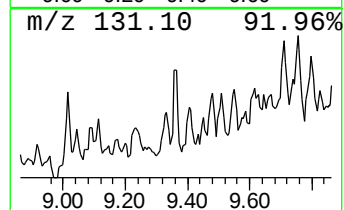
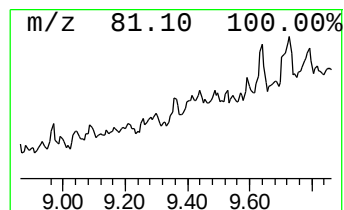
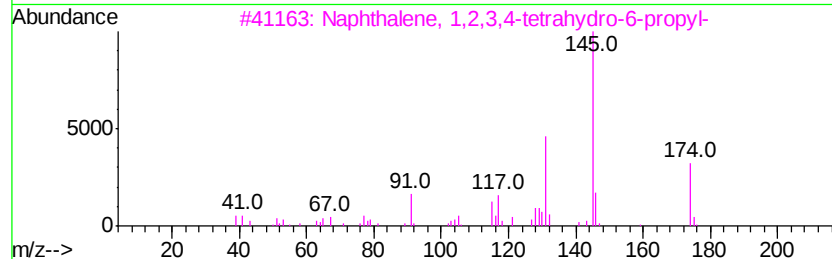
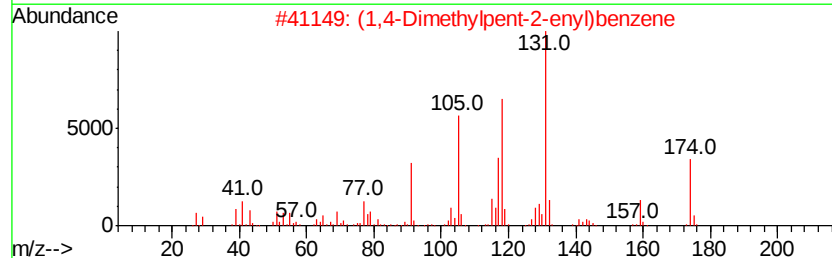
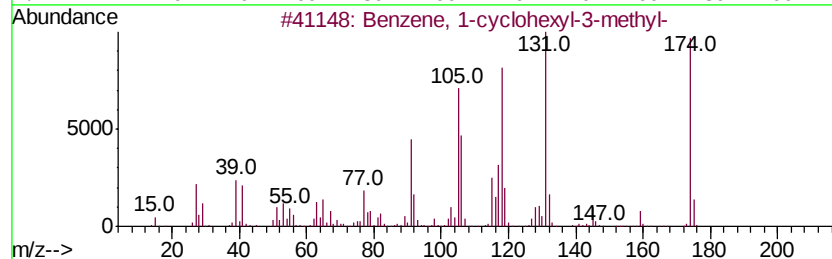
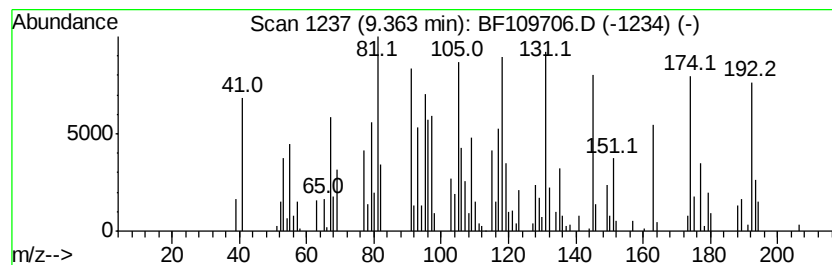
Instrument :
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ClientSampleId :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-cyclohexyl-3-met... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.36	2.36 ng	105060	Acenaphthene-d10	9.72	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	50
2	(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	41
3	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	042775-77-9	38
4	3-Methyl-2-(0-methylbenzyl)butanol	192	C13H20O	029107-38-8	35
5	2H-Indeno[2,1-b]furan-2-one, 3,3...	174	C11H10O2	080343-98-2	30



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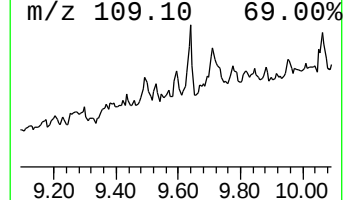
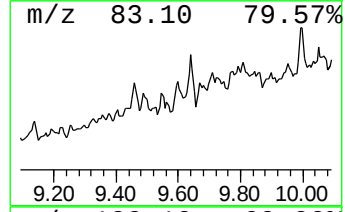
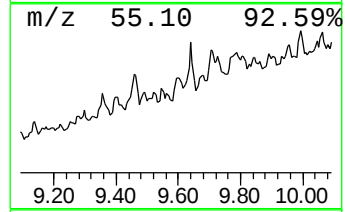
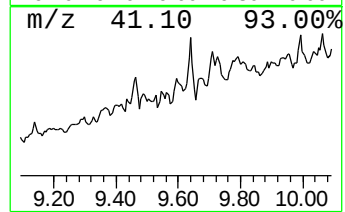
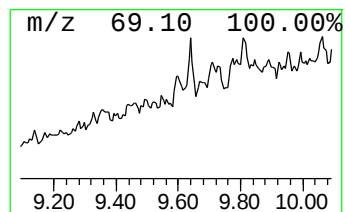
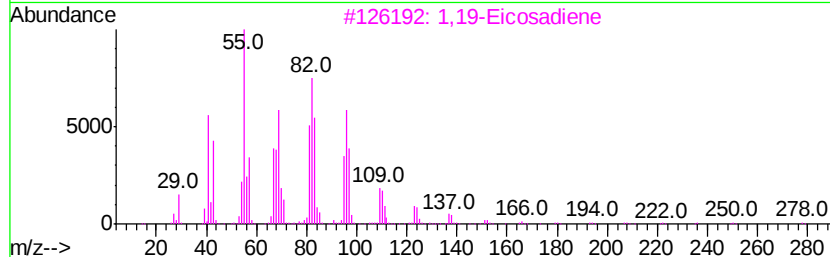
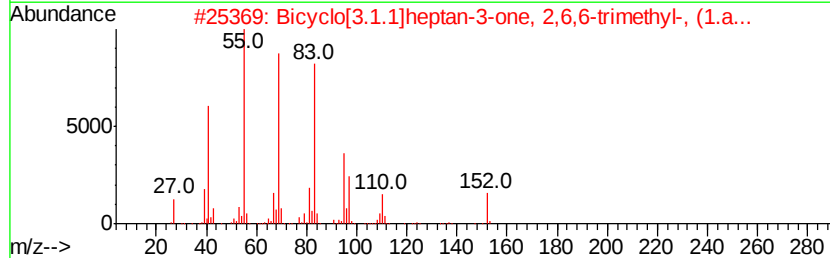
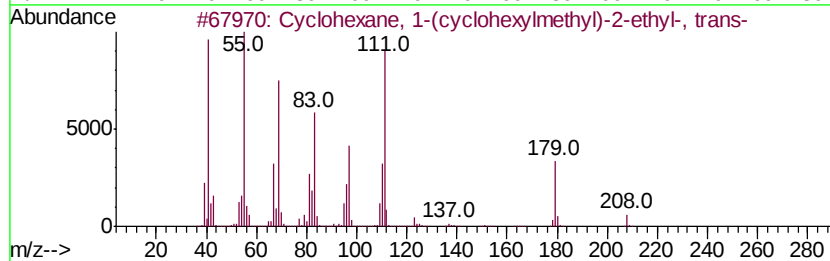
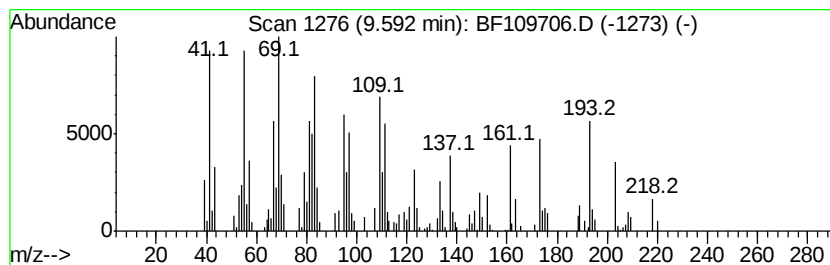
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown9.59 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.59	2.59 ng	115473	Acenaphthene-d10	9.72
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 1-(cyclohexylmethyl)...	208 C15H28	054934-92-8 41
2		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152 C10H16O	015358-88-0 27
3		1,19-Eicosadiene	278 C20H38	014811-95-1 27
4		Cyclopentane, (2-methylbutyl)-	140 C10H20	053366-38-4 18
5		2(1H)-Naphthalenone, octahydro-4...	208 C14H24O	054594-42-2 15



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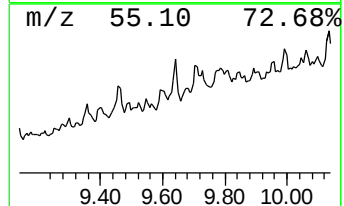
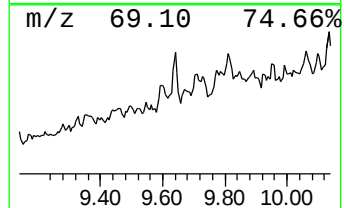
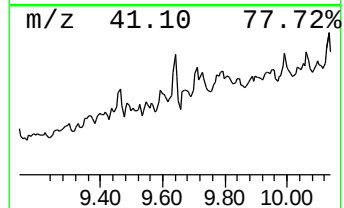
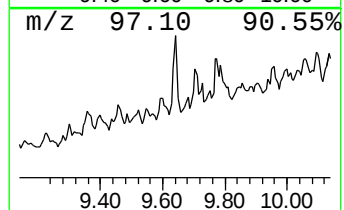
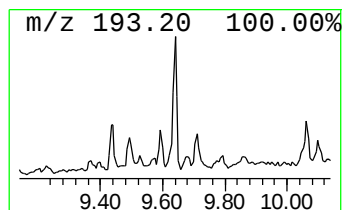
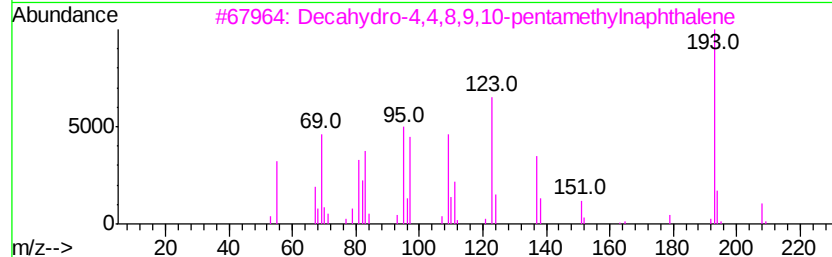
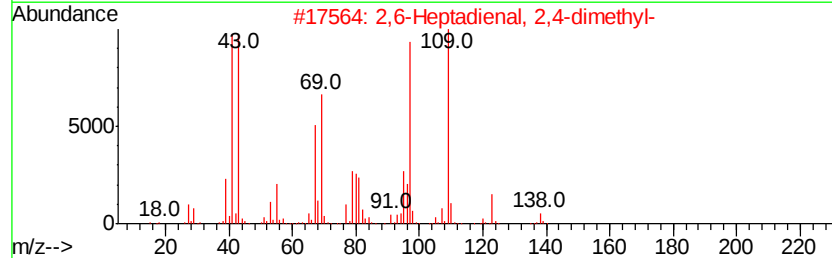
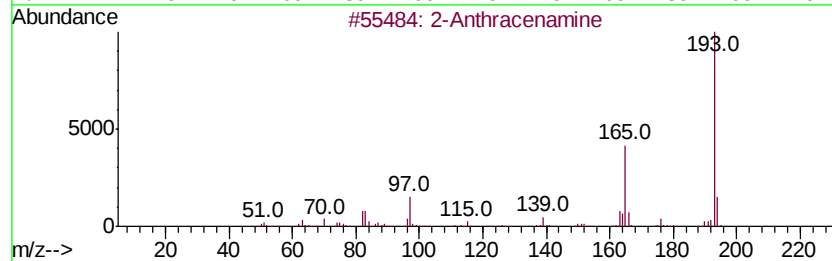
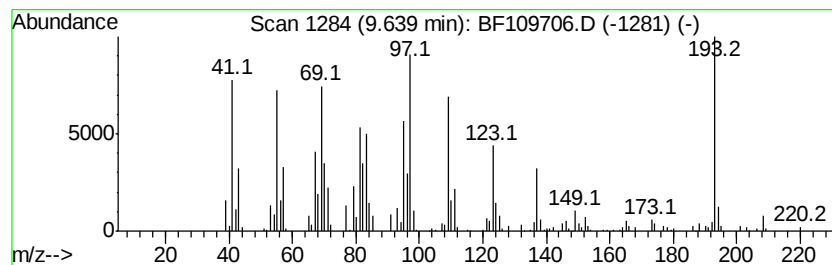
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Peak Number 5 unknown9.64 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.64	4.76 ng	212263	Acenaphthene-d10	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2		Anthracenamine	193	C14H11N	000613-13-8	46
2	2,6		Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	38
3	Decahydro-4,4,8,9,10-pentamethyl...			208	C15H28	080655-44-3	38
4	1-Hexyl-1-nitrocyclohexane			213	C12H23NO2	118252-09-8	22
5	Cyclohexane, 1-(cyclohexylmethyl...			194	C14H26	054823-96-0	20



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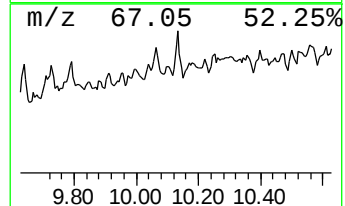
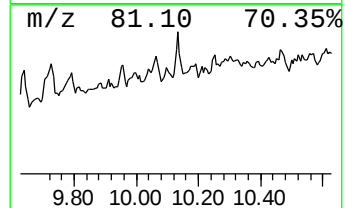
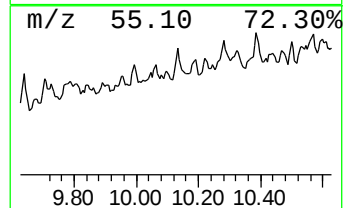
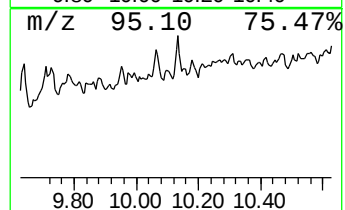
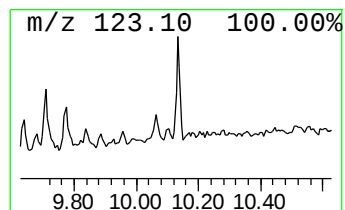
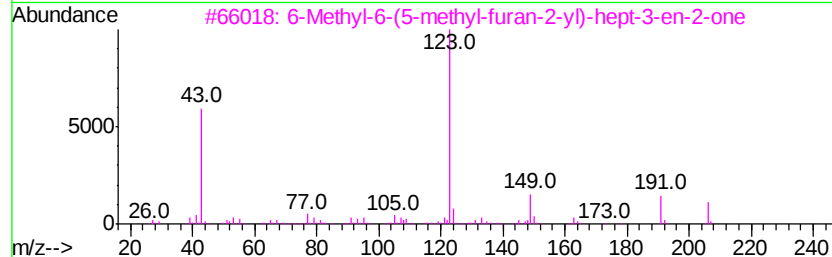
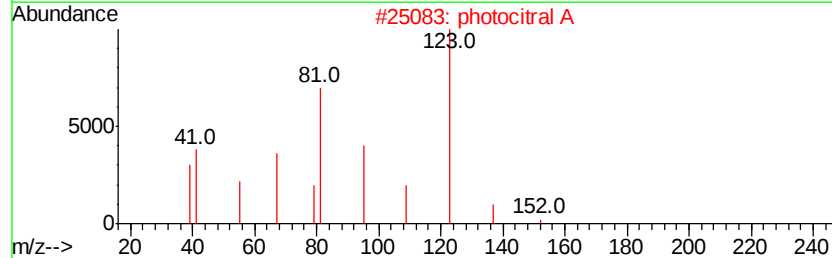
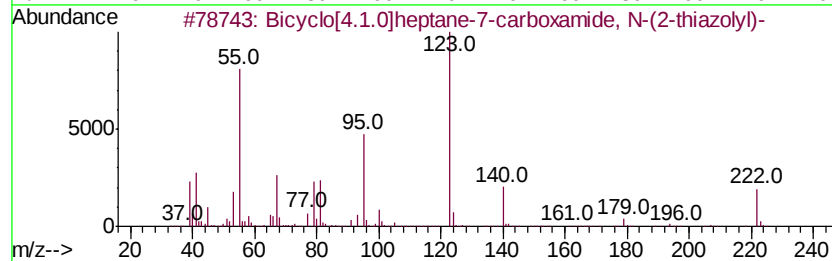
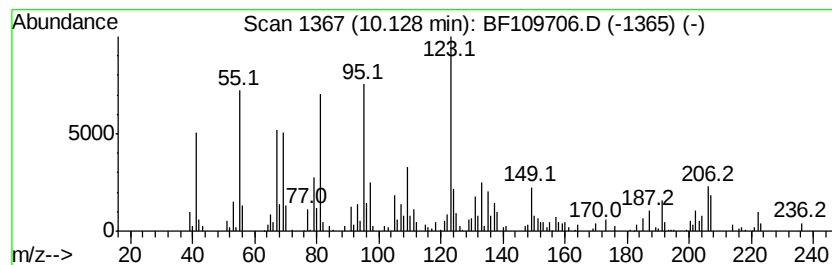
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown10.13 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.13	4.38 ng	195343	Acenaphthene-d10	9.72		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bicvclo[4.1.0]heptane-7-carboxam...	222	C11H14N2OS	329912-72-3	43
2		photocitral A	152	C10H16O	055253-28-6	43
3		6-Methyl-6-(5-methyl-furan-2-yl)...	206	C13H18O2	077822-40-3	38
4		Pentanamide, N-(4-methoxyphenyl)-	207	C12H17NO2	1000306-89-3	35
5		cis-1,3,3-trimethylbicyclo[3.1.0]...	152	C10H16O	1000365-94-2	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
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Operator : JU/SJ
Sample : J5331-02 10X
Misc :
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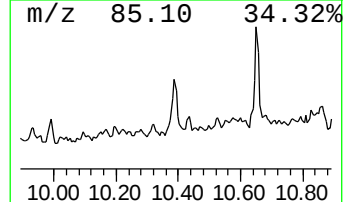
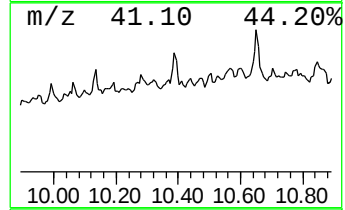
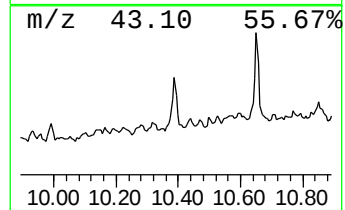
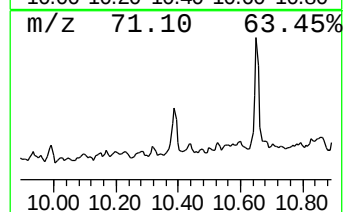
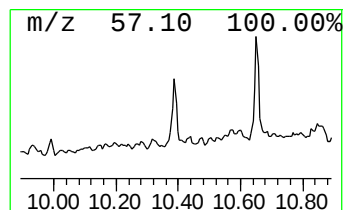
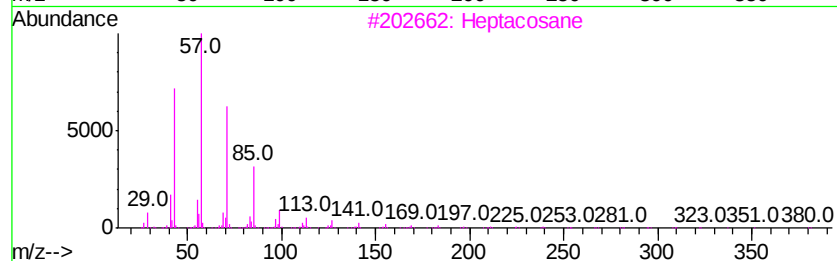
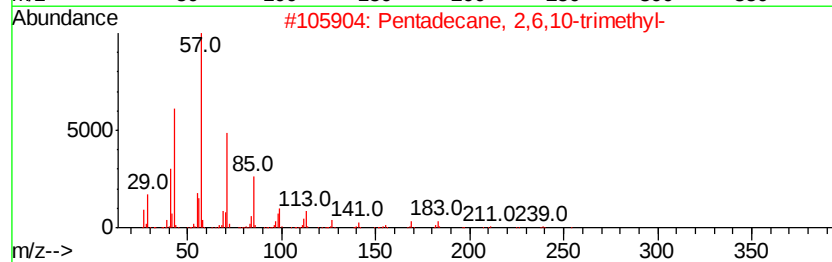
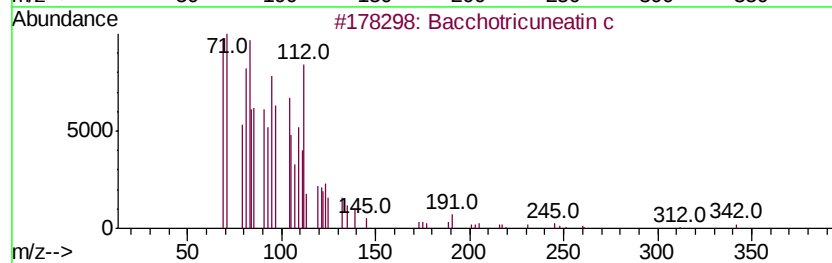
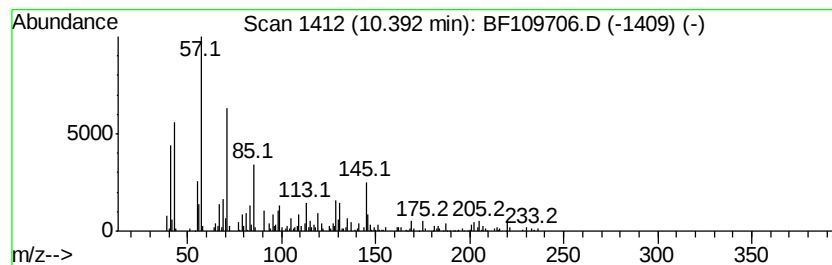
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Bacchotricuneatin c Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.39	3.36 ng	149644	Acenaphthene-d10		9.72
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bacchotricuneatin c	342	C20H22O5	066563-30-2	95
2	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
3	Heptacosane	380	C27H56	000593-49-7	72
4	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	62
5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
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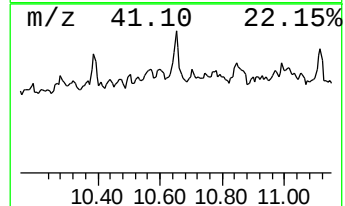
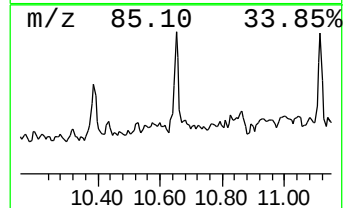
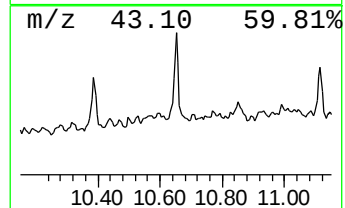
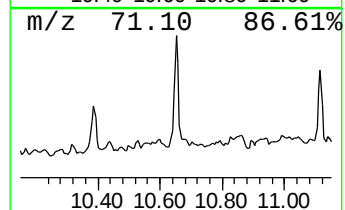
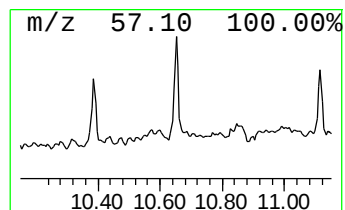
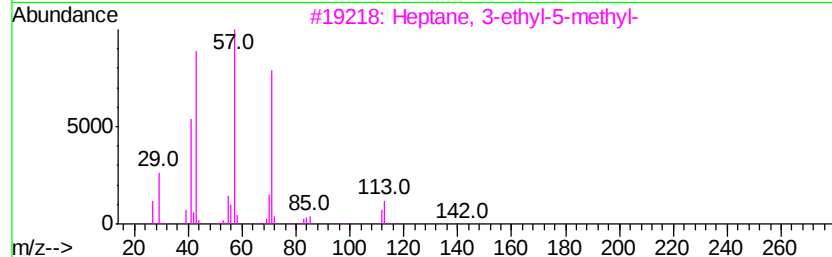
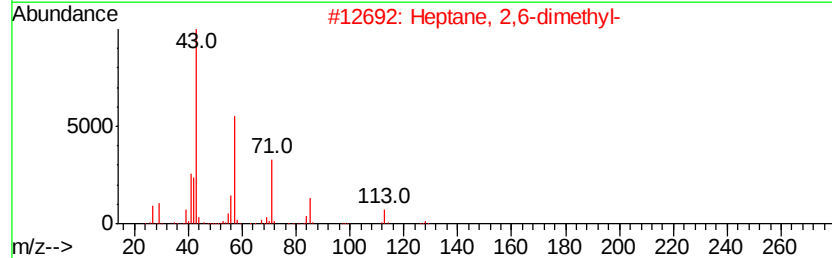
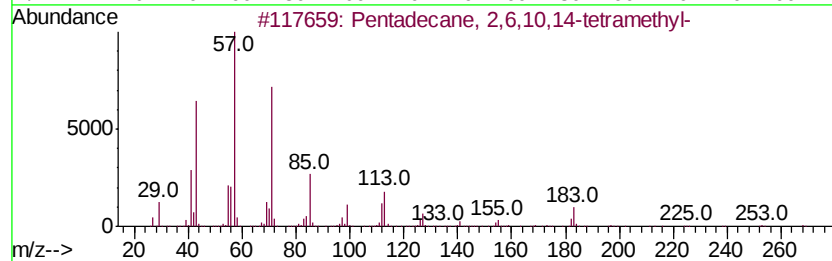
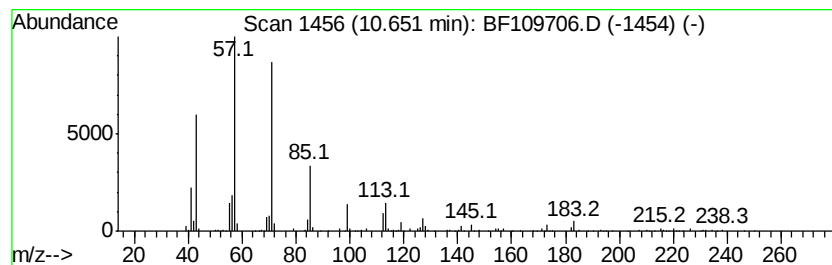
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane, 2,6,10,14-tetr... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.65	18.37 ng	288182	Phenanthrene-d10	11.20		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	78
2		Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	64
3		Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	64
4		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
5		Hexadecane	226	C16H34	000544-76-3	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109706.D
Acq On : 9 Oct 2018 20:27
Operator : JU/SJ
Sample : J5331-02 10X
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Instrument :
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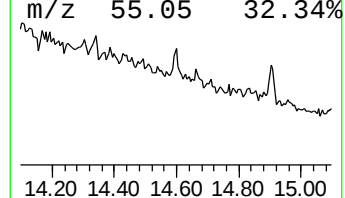
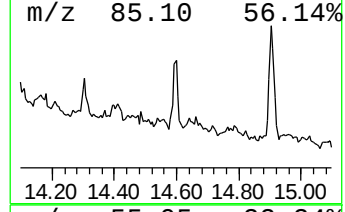
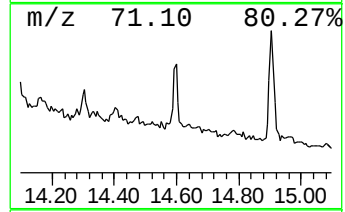
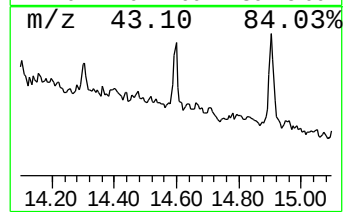
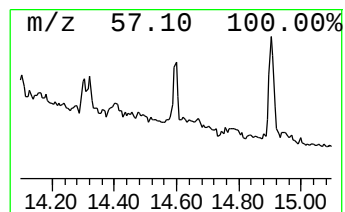
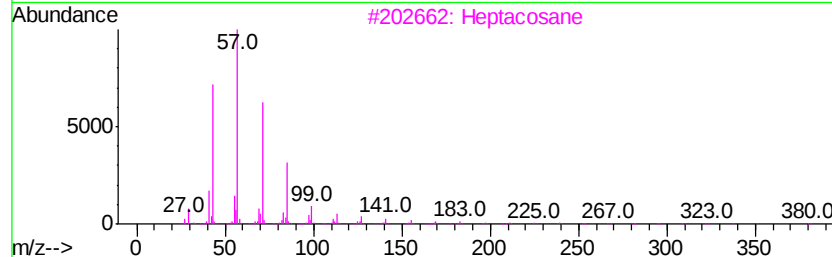
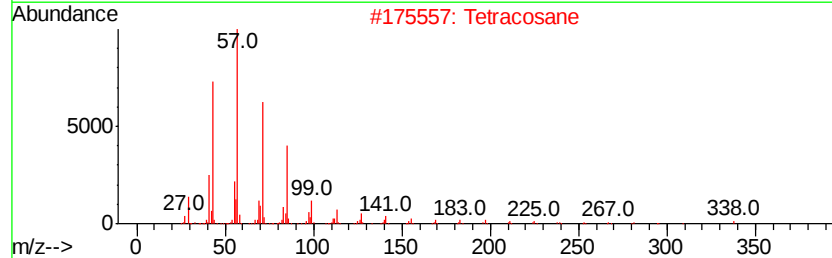
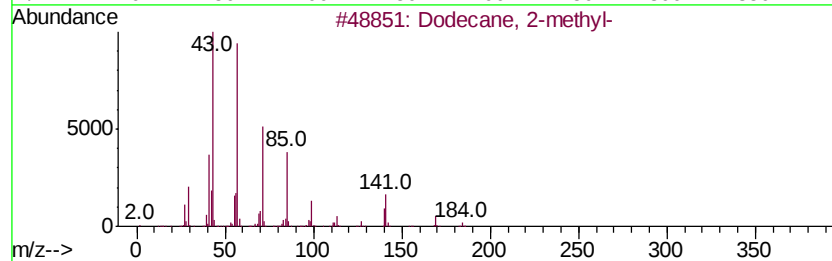
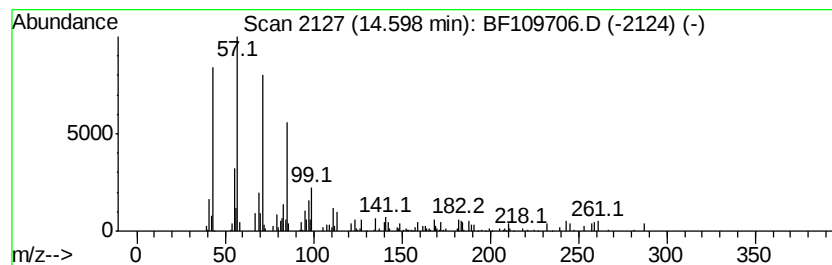
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dodecane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	3.27 ng	94923	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
2		Tetracosane	338	C24H50	000646-31-1	64
3		Heptacosane	380	C27H56	000593-49-7	59
4		Octacosane	394	C28H58	000630-02-4	49
5		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109706.D
Acq On : 9 Oct 2018 20:27
Operator : JU/SJ
Sample : J5331-02 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

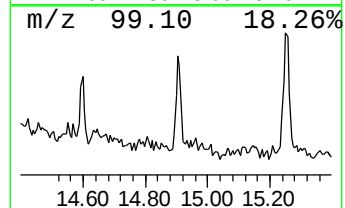
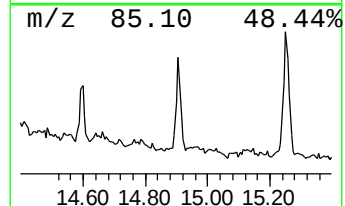
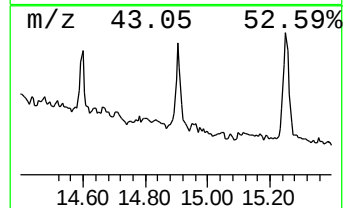
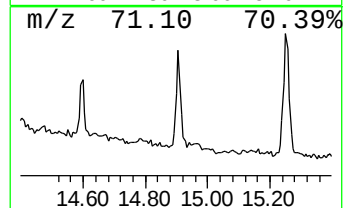
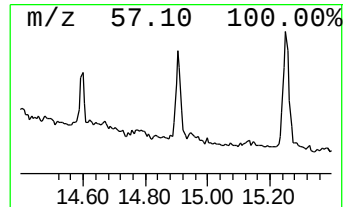
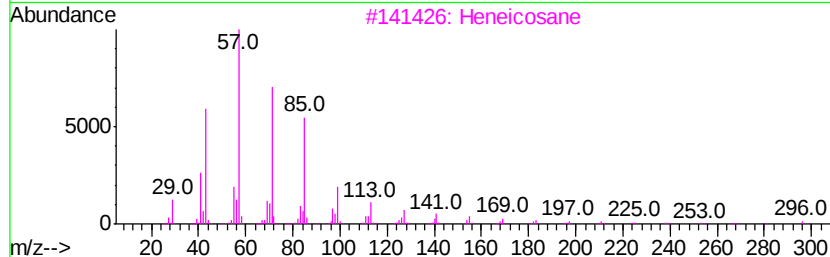
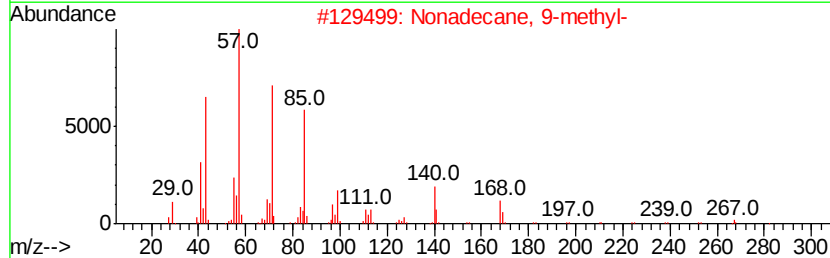
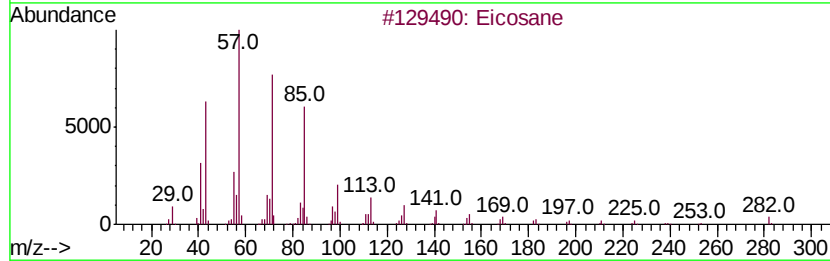
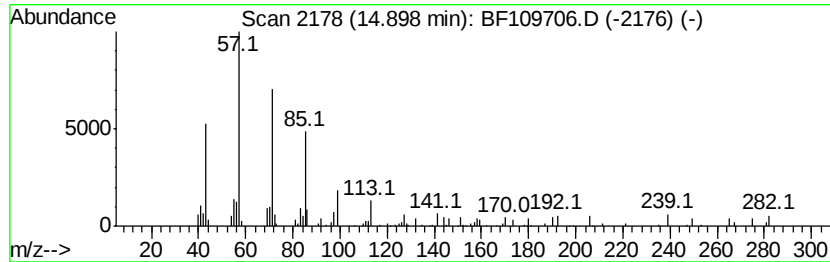
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	3.42 ng	99319	Perylene-d12	15.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 94
2	Nonadecane, 9-methyl-		282 C20H42	013287-24-6 91
3	Heneicosane		296 C21H44	000629-94-7 91
4	Heptadecane		240 C17H36	000629-78-7 91
5	Octacosane		394 C28H58	000630-02-4 90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109706.D
Acq On : 9 Oct 2018 20:27
Operator : JU/SJ
Sample : J5331-02 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

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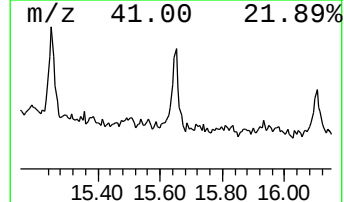
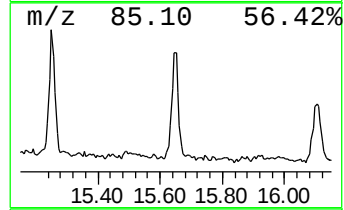
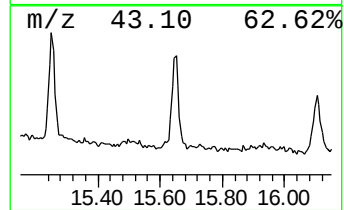
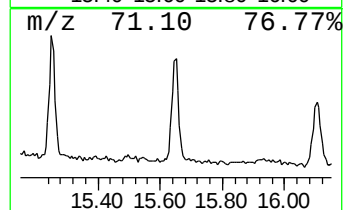
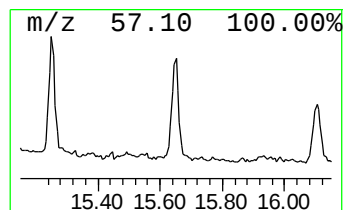
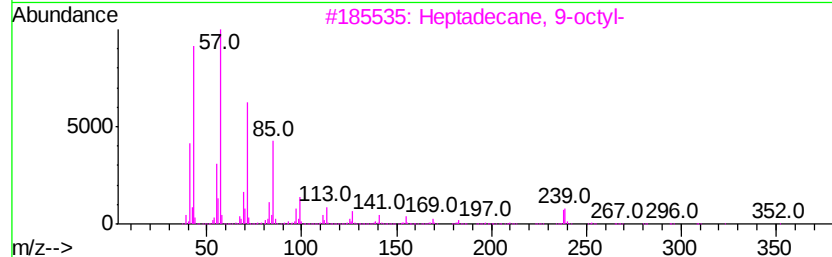
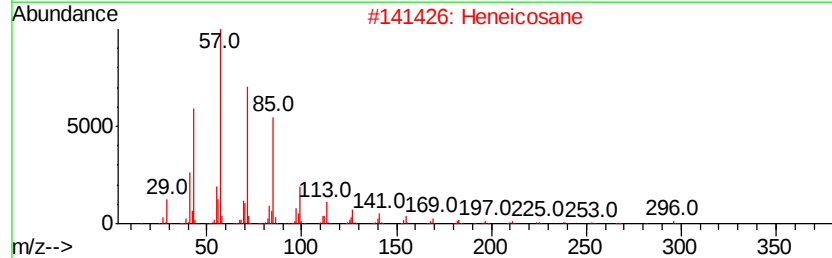
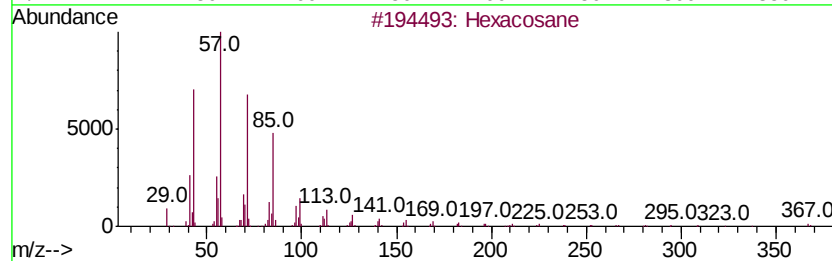
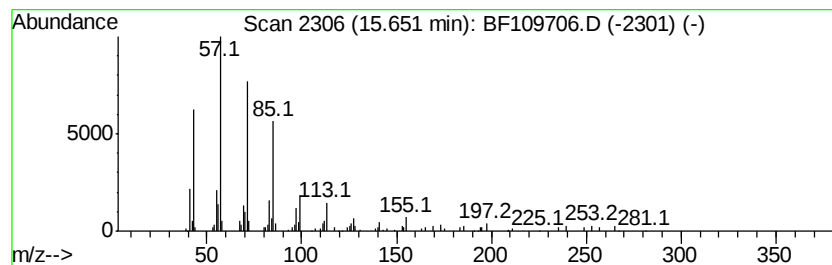
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	4.58 ng	133162	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
4		Hexadecane	226	C16H34	000544-76-3	80
5		Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
Data File : BF109706.D
Acq On : 9 Oct 2018 20:27
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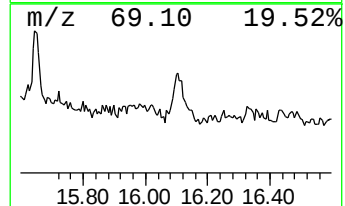
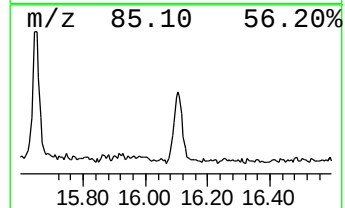
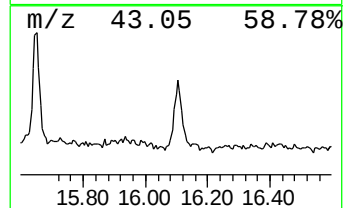
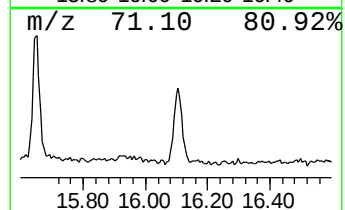
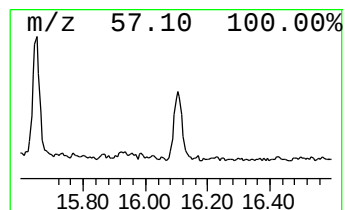
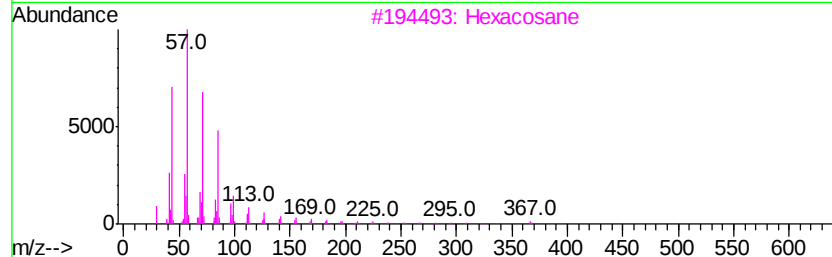
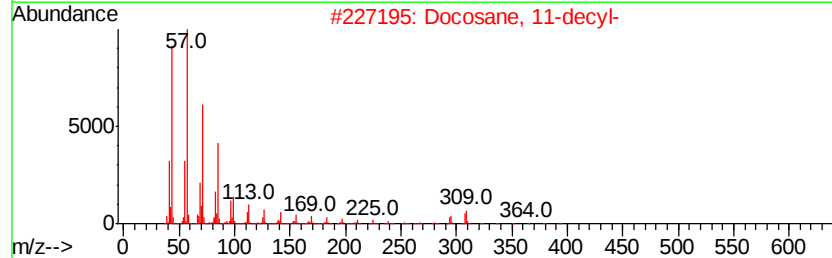
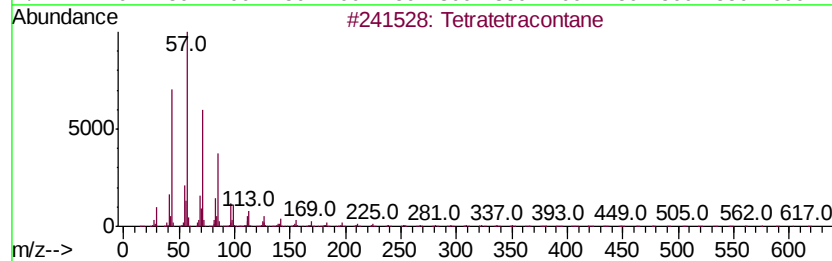
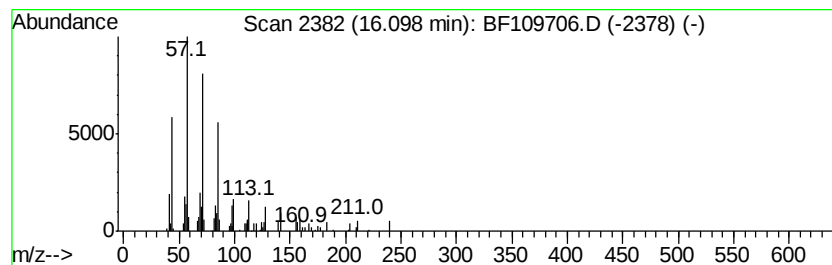
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetratetracontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	3.55 ng	103165	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Docosane, 11-decyl-	451	C32H66	055401-55-3	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100918\
Data File : BF109706.D
Acq On : 9 Oct 2018 20:27
Operator : JU/SJ
Sample : J5331-02 10X
Misc :
ALS Vial : 10 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\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1-cycloh...	9.36	2.4	ng	105060	3	9.72	891344	20.0
unknown9.59	9.59	2.6	ng	115473	3	9.72	891344	20.0
unknown9.64	9.64	4.8	ng	212263	3	9.72	891344	20.0
unknown10.13	10.13	4.4	ng	195343	3	9.72	891344	20.0
Bacchotricuneatin c	10.39	3.4	ng	149644	3	9.72	891344	20.0
Pentadecane, 2,6,...	10.65	18.4	ng	288182	4	11.20	313704	20.0
Dodecane, 2-methyl-	14.60	3.3	ng	94923	6	15.23	581250	20.0
Eicosane	14.90	3.4	ng	99319	6	15.23	581250	20.0
Hexacosane	15.65	4.6	ng	133162	6	15.23	581250	20.0
Tetratetracontane	16.10	3.5	ng	103165	6	15.23	581250	20.0