

Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102159.D
 Acq On : 17 Jan 2018 19:27
 Operator : SJ/JU
 Sample : J1144-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GR-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-BF010918.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	4.928	473	483	490	rBV	179408	270590	7.80%	0.845%
2	5.334	546	552	555	rVB	3326666	3373265	97.23%	10.538%
3	5.993	660	664	666	rBV	38553	35895	1.03%	0.112%
4	6.363	721	727	731	rBV	2886192	3469256	100.00%	10.837%
5	6.493	744	749	752	rBV	3546446	3456170	99.62%	10.796%
6	6.545	756	758	764	rVB2	34640	37553	1.08%	0.117%
7	6.716	784	787	791	rBV	1000871	889392	25.64%	2.778%
8	6.875	810	814	817	rBV	2804587	2468082	71.14%	7.710%
9	7.287	876	884	887	rBV	2270244	2099531	60.52%	6.559%
10	7.998	1001	1005	1011	rBV	1302468	1213386	34.98%	3.790%
11	8.516	1089	1093	1097	rBV	524040	460416	13.27%	1.438%
12	8.557	1097	1100	1104	rBV	252706	191802	5.53%	0.599%
13	8.845	1145	1149	1157	rBV2	171686	203520	5.87%	0.636%
14	9.081	1183	1189	1192	rBV	3228859	2826669	81.48%	8.830%
15	9.434	1246	1249	1252	rBV	751794	648418	18.69%	2.026%
16	9.475	1252	1256	1259	rVB	411662	359524	10.36%	1.123%
17	9.751	1300	1303	1307	rVB	1313430	1115752	32.16%	3.485%
18	10.245	1384	1387	1390	rBV	163071	131765	3.80%	0.412%
19	10.469	1421	1425	1428	rBV	842387	675211	19.46%	2.109%
20	10.539	1433	1437	1441	rBV	1868743	1749773	50.44%	5.466%
21	10.657	1454	1457	1461	rBV3	47226	59564	1.72%	0.186%
22	10.904	1496	1499	1502	rVB	67112	56177	1.62%	0.175%
23	11.104	1530	1533	1536	rBV2	41146	35944	1.04%	0.112%
24	11.233	1551	1555	1558	rBV	1100106	890781	25.68%	2.783%
25	11.775	1642	1647	1653	rBV	492408	517139	14.91%	1.615%
26	12.551	1776	1779	1784	rBV	88647	103289	2.98%	0.323%
27	12.822	1820	1825	1828	rBV	1710786	1681984	48.48%	5.254%
28	13.716	1974	1977	1981	rBV	281356	275510	7.94%	0.861%
29	13.857	1997	2001	2006	rBV2	881102	1287574	37.11%	4.022%
30	14.963	2186	2189	2192	rVB	146965	142776	4.12%	0.446%
31	15.286	2240	2244	2247	rBV	593905	874937	25.22%	2.733%
32	15.945	2353	2356	2367	rVB2	226344	410348	11.83%	1.282%

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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

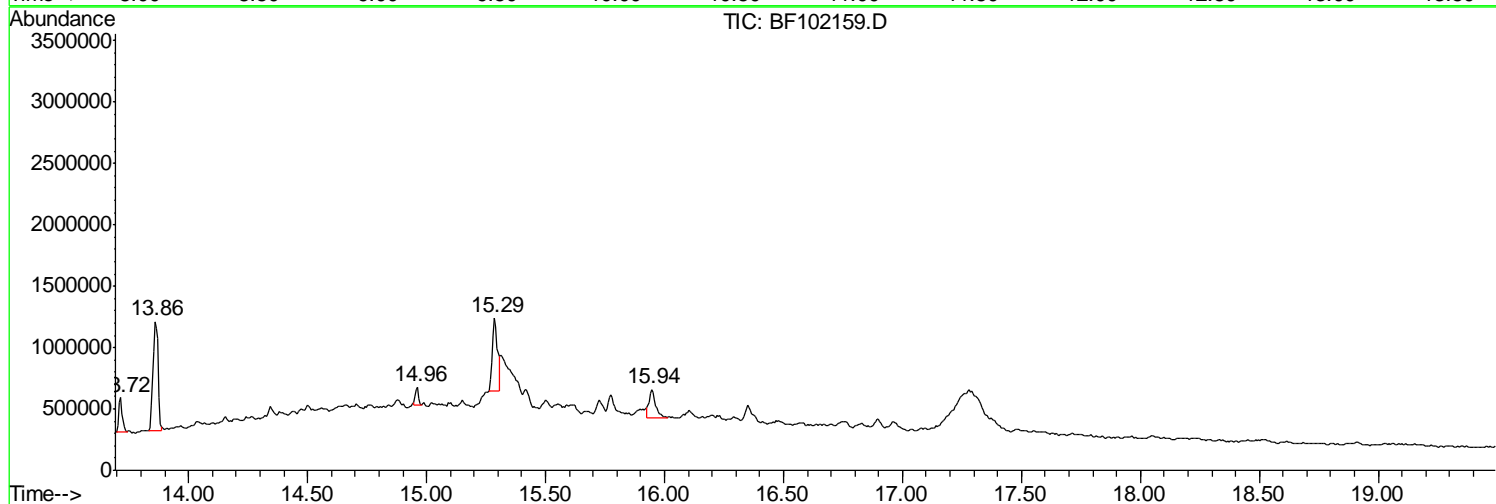
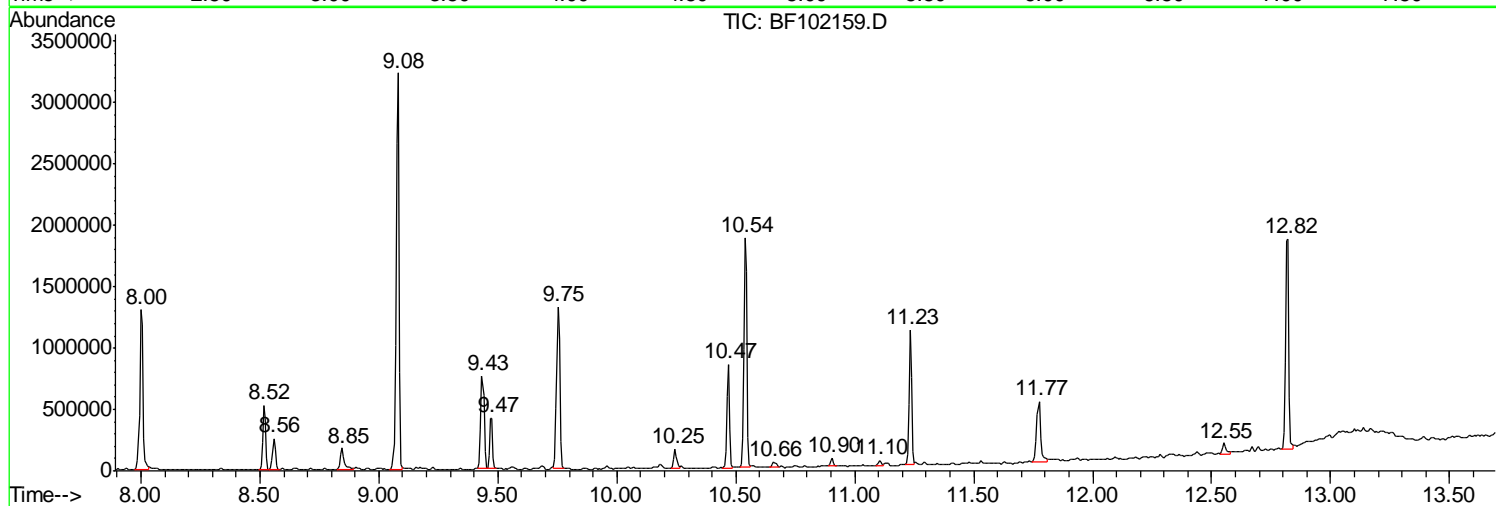
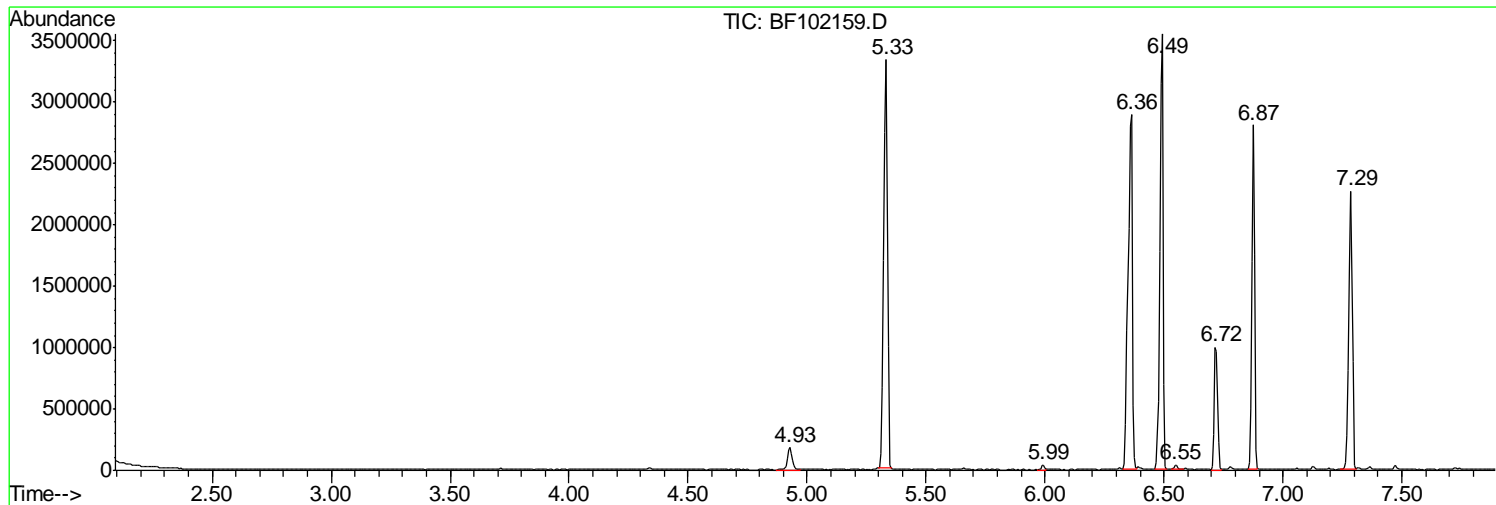
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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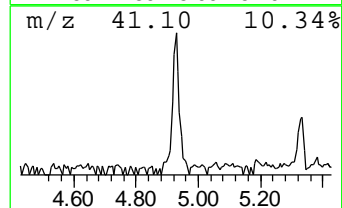
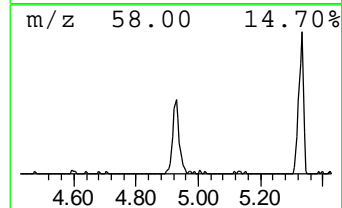
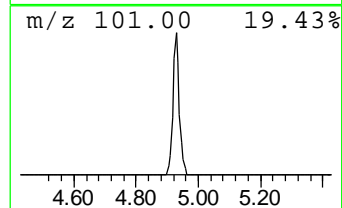
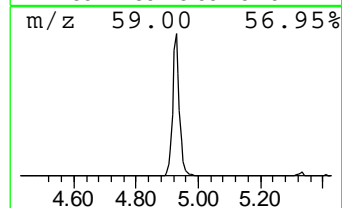
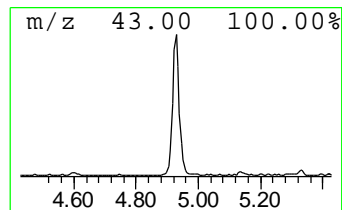
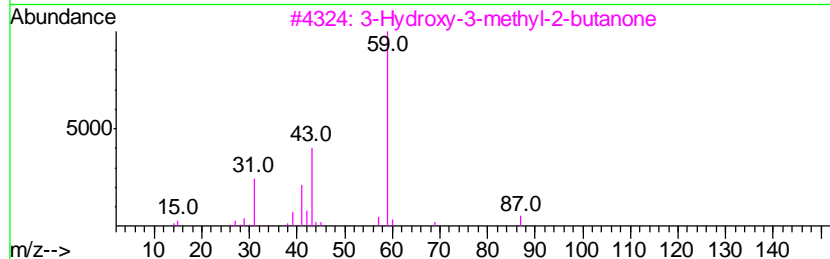
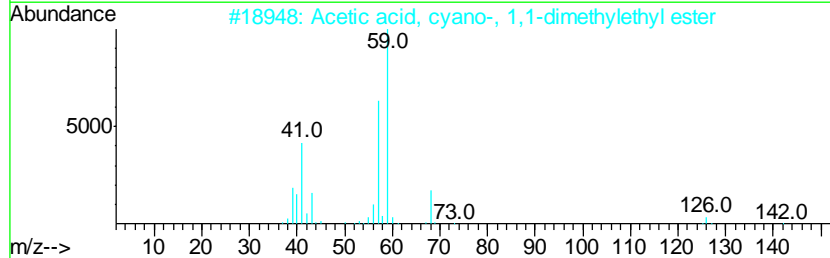
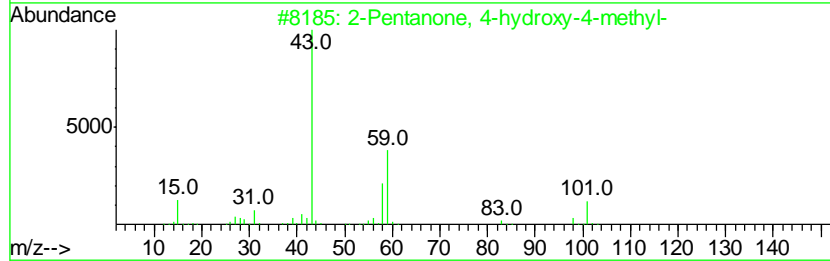
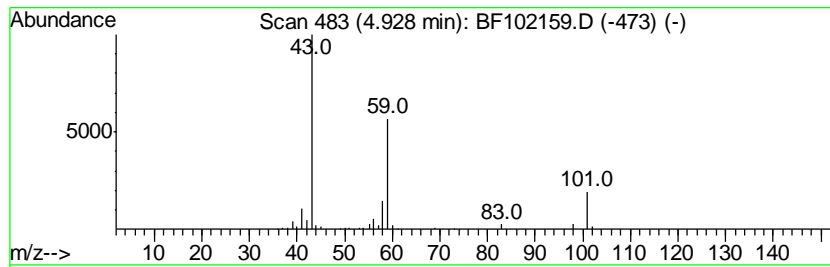
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	6.08 ng	270590	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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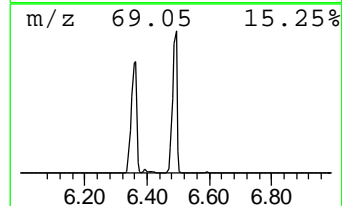
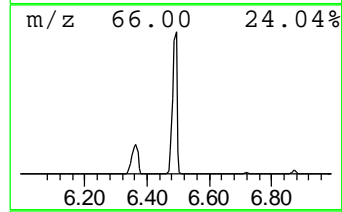
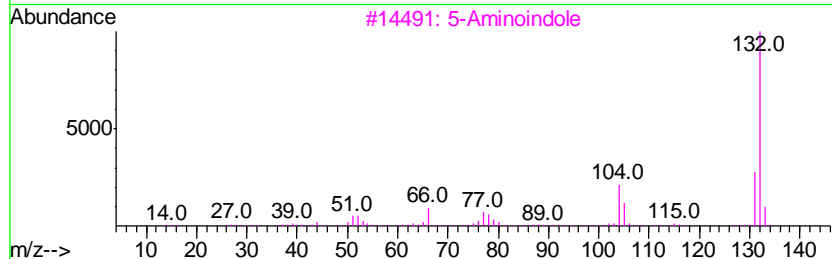
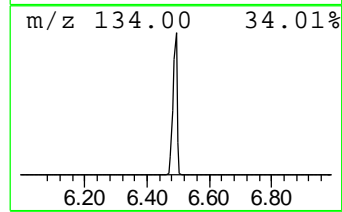
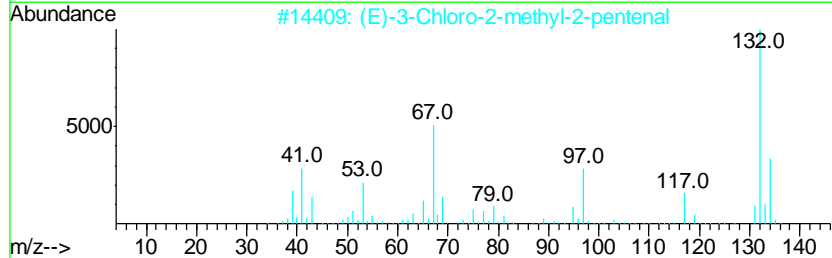
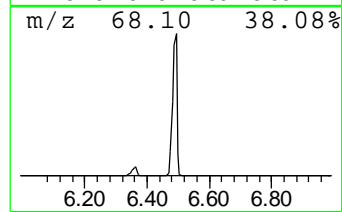
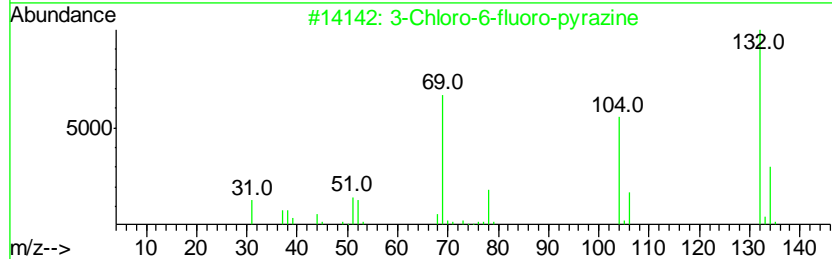
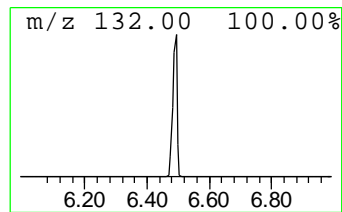
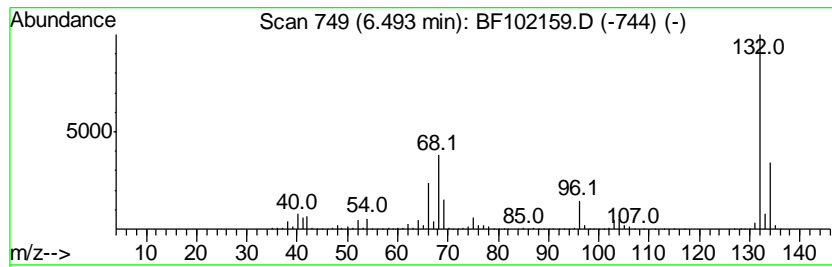
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 Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	77.72 ng	3456170	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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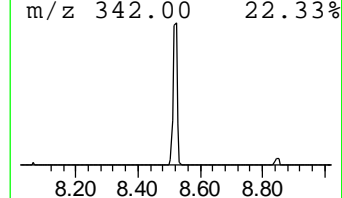
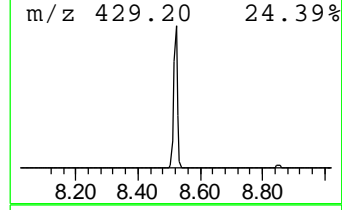
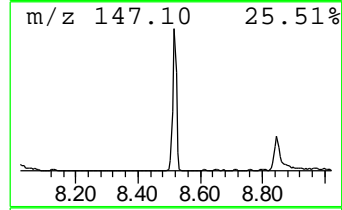
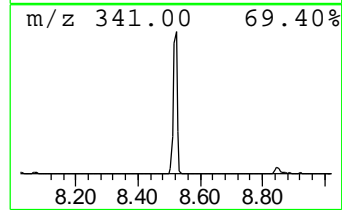
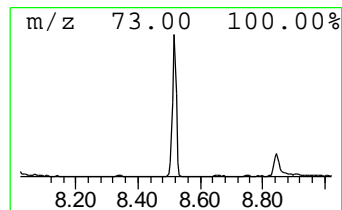
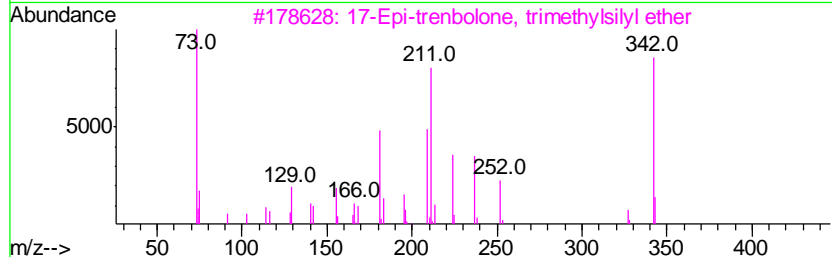
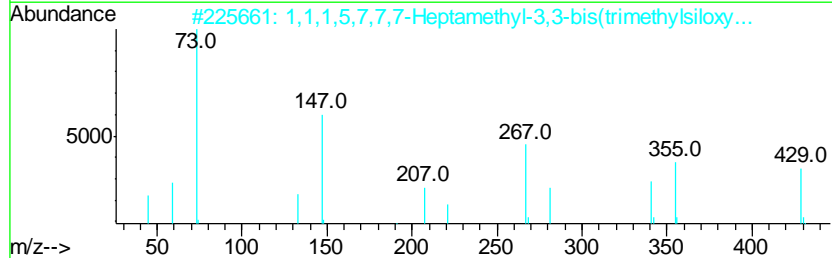
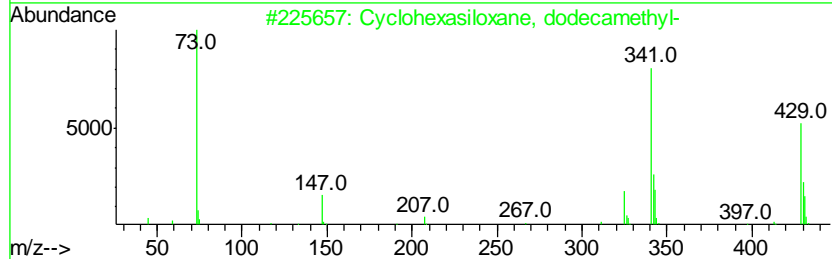
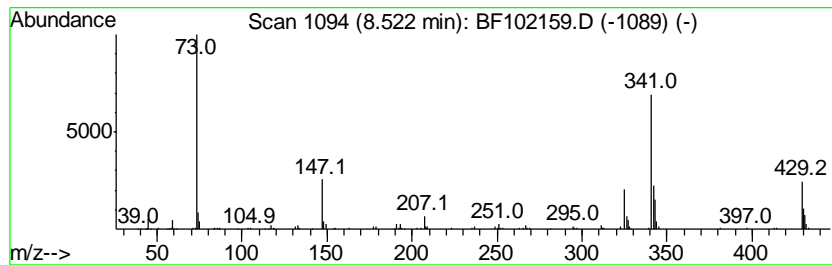
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 Peak Number 4 Cyclohexasiloxane, dodecane... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.52	7.59 ng	460416	Naphthalene-d8	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	90
2		1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	37
3		17-Epi-trenbolone, trimethylsily...	342	C21H30O2Si	1000293-11-0	22
4		2H-1,4-Benzodiazepin-2-one, 7-ch...	342	C18H19ClN2OSi	055299-24-6	22
5		Guaifenesin di-tms derivative	342	C16H30O4Si2	107966-19-8	17



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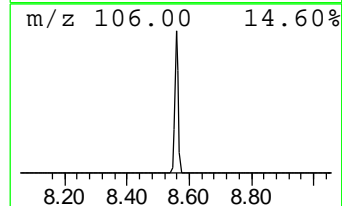
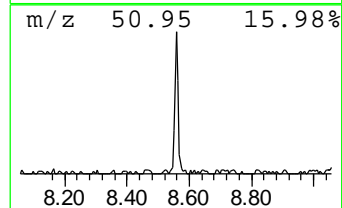
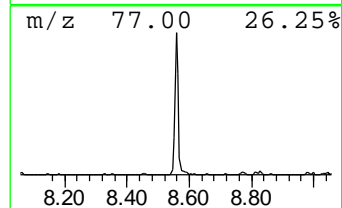
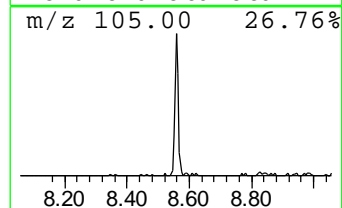
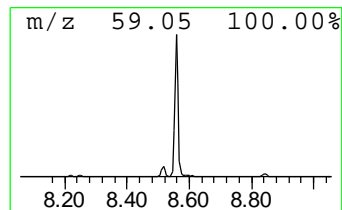
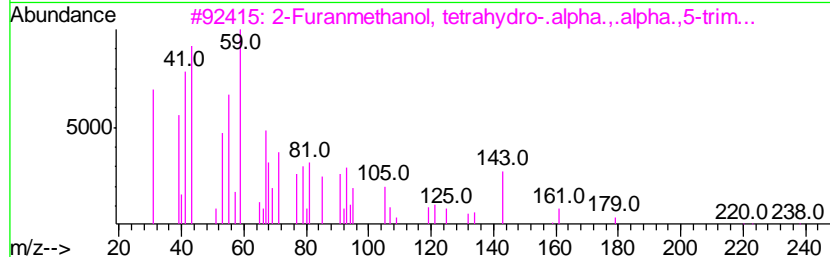
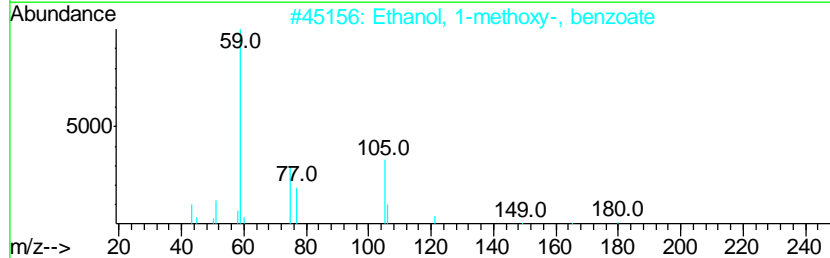
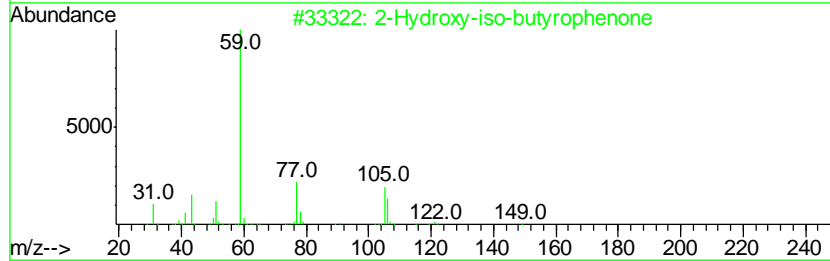
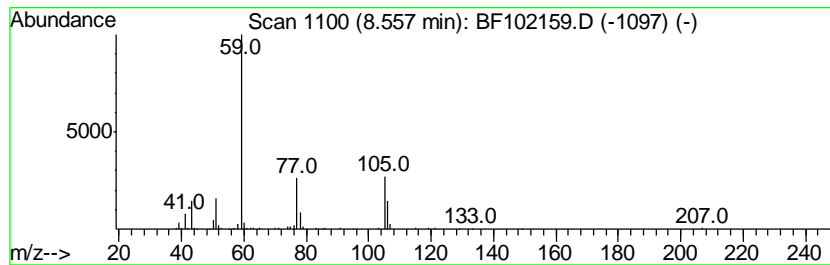
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 Peak Number 5 2-Hydroxy-iso-butyrophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	3.16 ng	191802	Napthalene-d8	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hydroxy-iso-butyrophenone	164	C10H12O2	007473-98-5	90
2		Ethanol, 1-methoxy-, benzoate	180	C10H12O3	051835-44-0	50
3		2-Furanmethanol, tetrahydro-.alp...	238	C15H26O2	026184-88-3	28
4		Ethanone, 2-amino-1-phenyl-	135	C8H9NO	000613-89-8	10
5		Guanidine	59	CH5N3	000113-00-8	9



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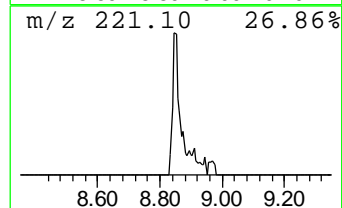
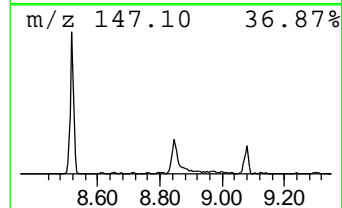
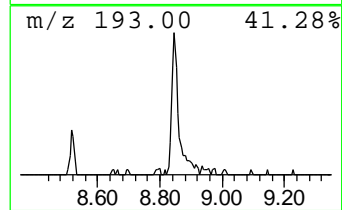
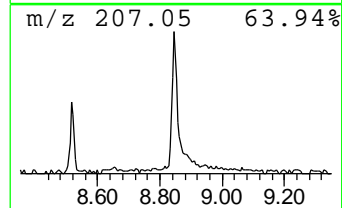
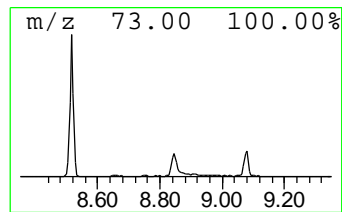
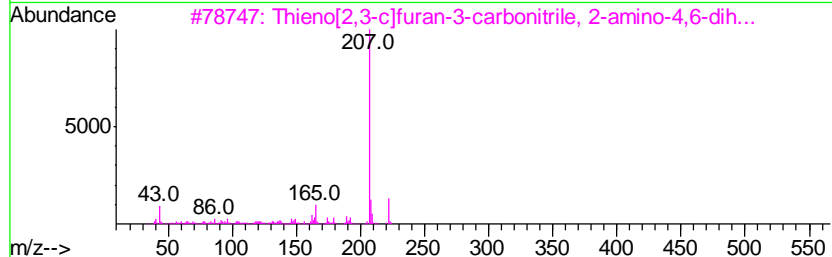
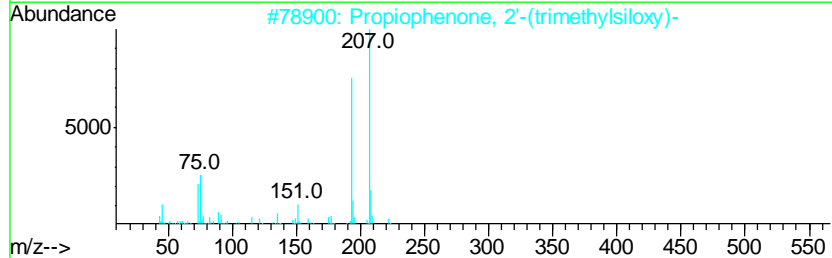
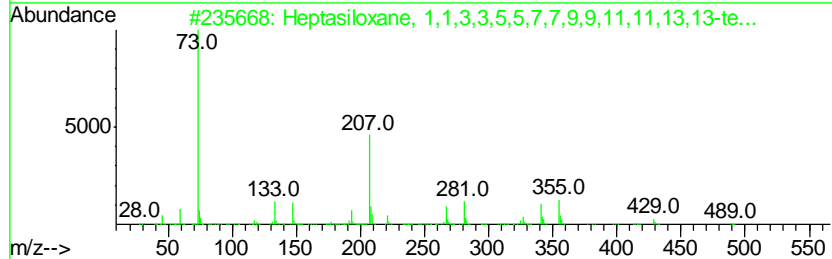
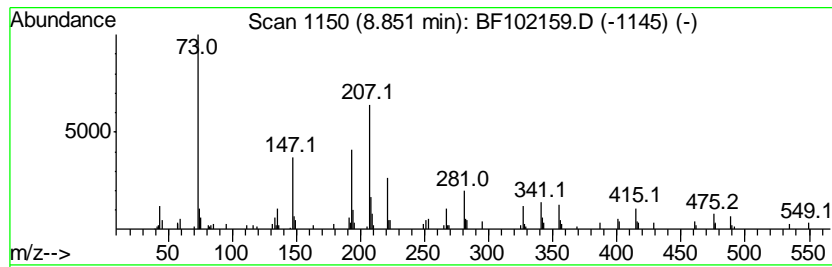
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 Peak Number 6 unknown8.85 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.85	3.35 ng	203520	Naphthalene-d8	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptasiloxane, 1,1,3,3,5,5,7,7,9...	504	C14H44O6Si7	019095-23-9	47
2		Propiophenone, 2'-(trimethylsilo...	222	C12H18O2Si	033342-87-9	43
3		Thieno[2.3-c]furan-3-carbonitril...	222	C11H14N2OS	447412-24-0	30
4		1,4-Bis(trimethylsilyl)benzene	222	C12H22Si2	013183-70-5	27
5		Propanamide, N-(4-methoxyphenyl)...	207	C12H17NO2	056619-94-4	27



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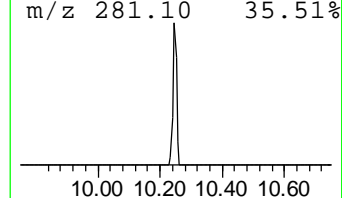
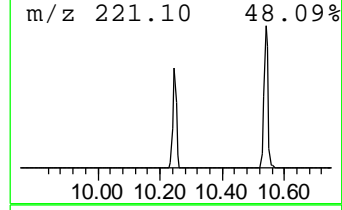
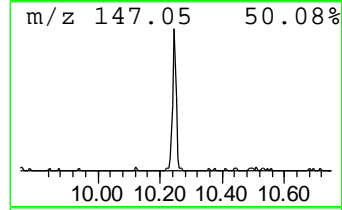
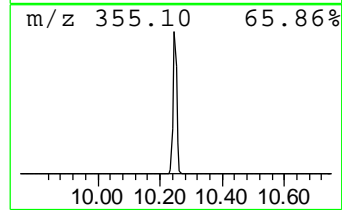
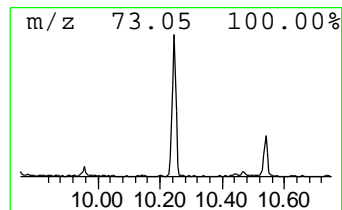
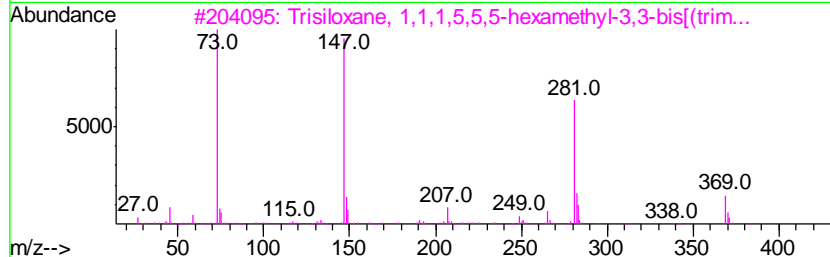
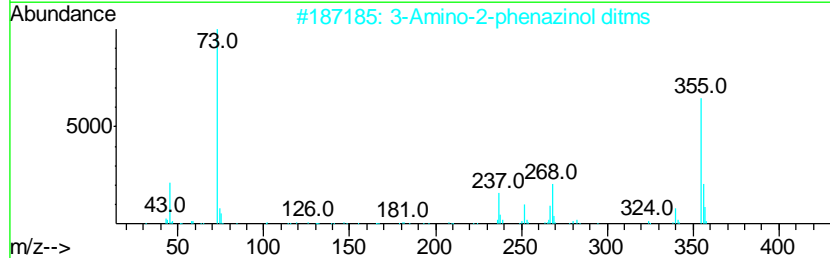
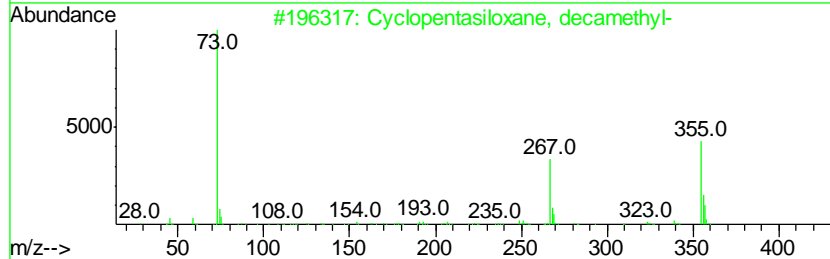
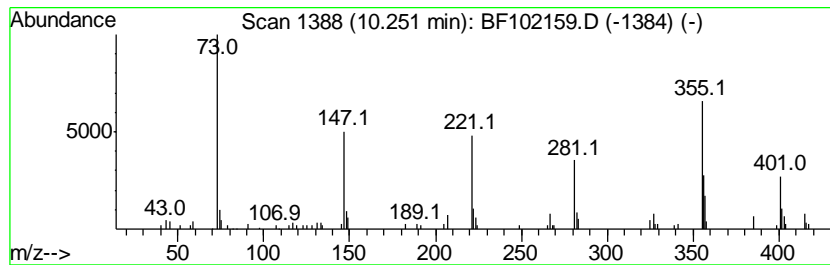
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 7 unknown10.25 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	2.36 ng	131765	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	27
2		3-Amino-2-phenazinol ditms	355	C18H25N3OSi2	1000332-15-5	27
3		Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	27
4		Mercaptoacetic acid, bis(trimeth...	236	C8H20O2SSi2	006398-62-5	25
5		N-(Trifluoroacetyl)-N,O,O',O''-t...	553	C22H42F3NO4Si4	1000072-26-7	22



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102159.D
 Acq On : 17 Jan 2018 19:27
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Instrument :
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 ClientSampled :
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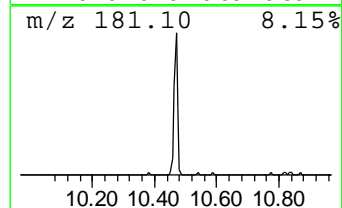
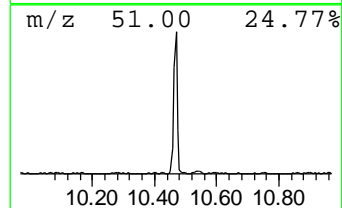
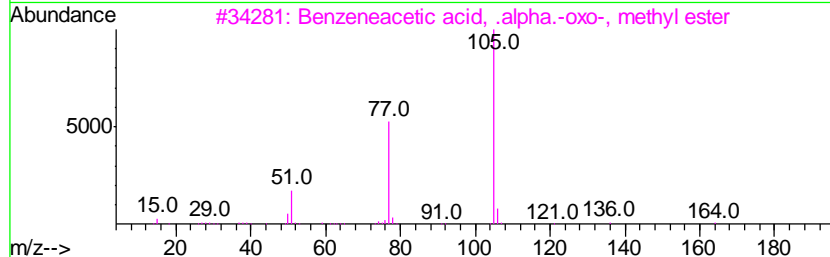
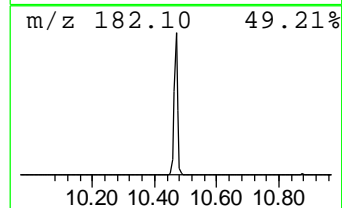
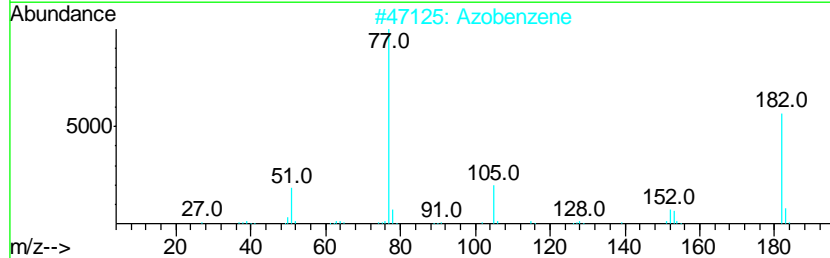
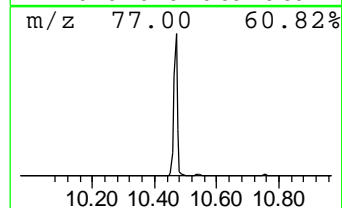
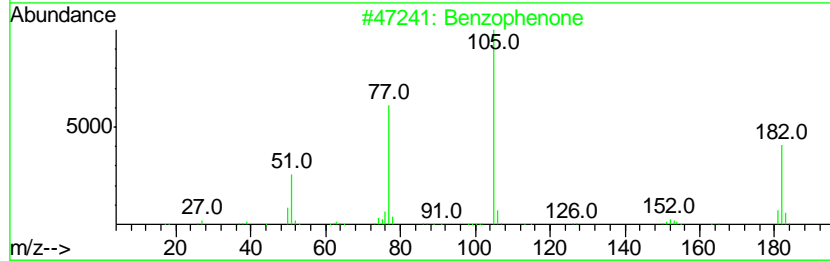
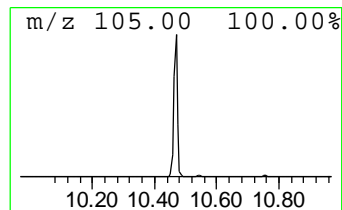
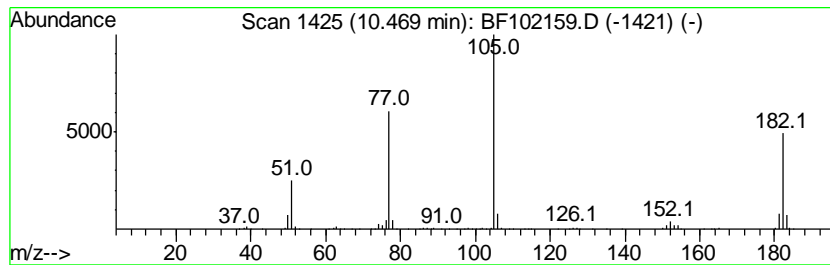
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	12.10 ng	675211	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	97
2		Azobenzene	182	C12H10N2	000103-33-3	72
3		Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	49
4		1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	47
5		Vinyl benzoate	148	C9H8O2	000769-78-8	47



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102159.D
 Acq On : 17 Jan 2018 19:27
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Instrument :
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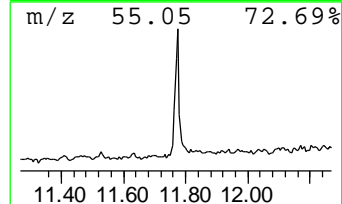
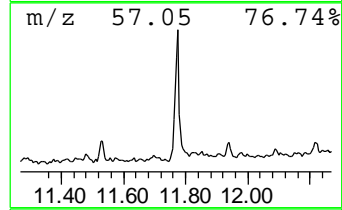
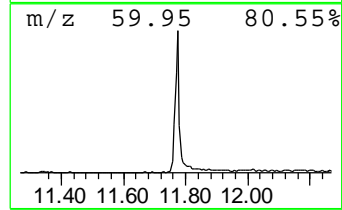
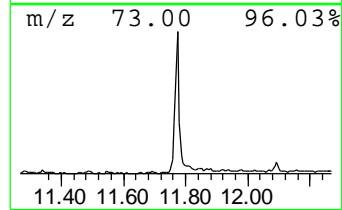
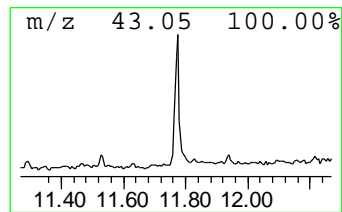
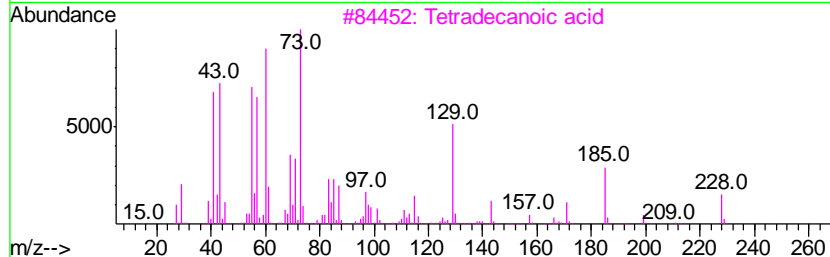
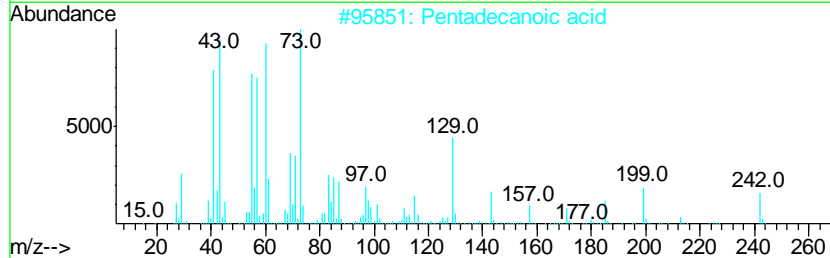
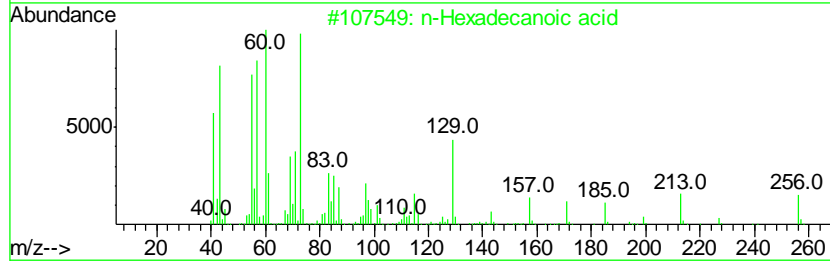
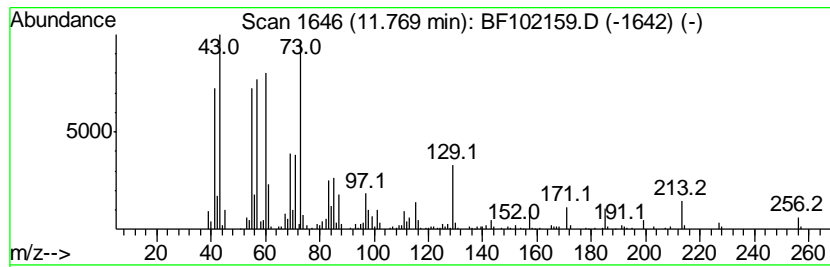
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	11.61 ng	517139	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	94
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4		Tridecanoic acid	214	C13H26O2	000638-53-9	90
5		Undecanoic acid	186	C11H22O2	000112-37-8	68



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
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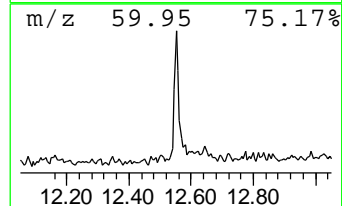
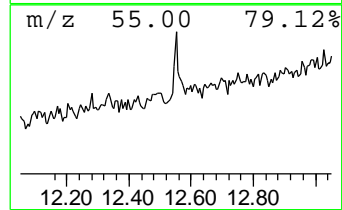
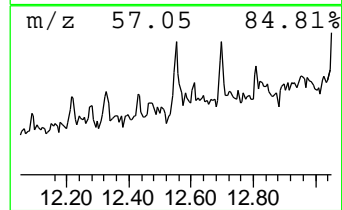
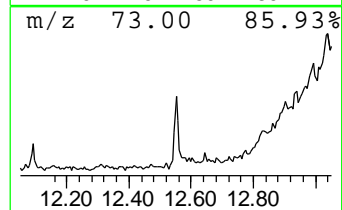
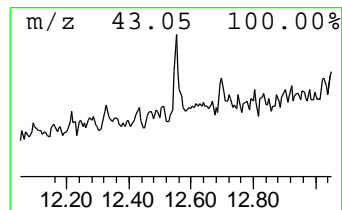
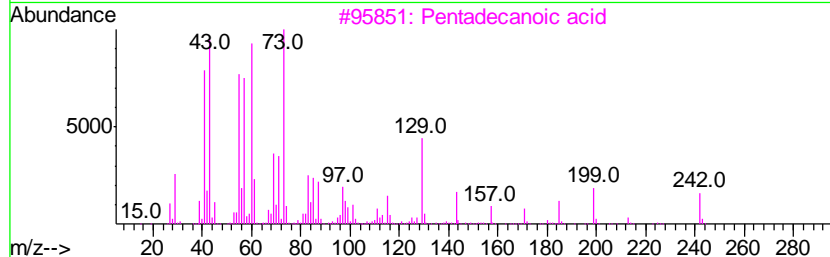
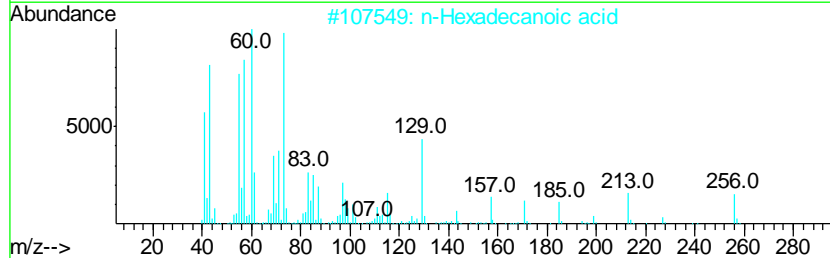
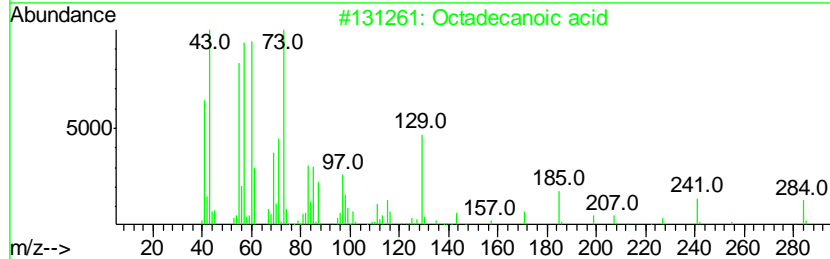
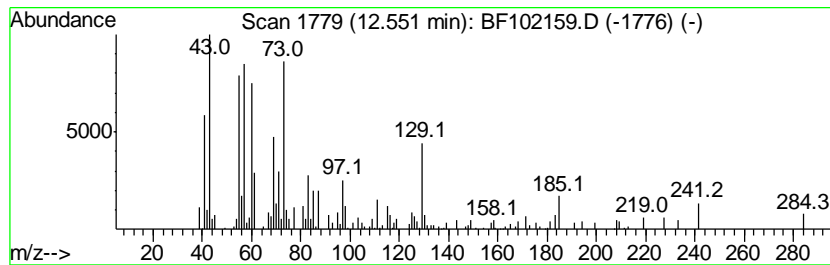
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	2.32 ng	103289	Phenanthrene-d10	11.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octadecanoic acid	284 C18H36O2	000057-11-4 91
2		n-Hexadecanoic acid	256 C16H32O2	000057-10-3 83
3		Pentadecanoic acid	242 C15H30O2	001002-84-2 72
4		Undecanoic acid	186 C11H22O2	000112-37-8 45
5		Trehalose	342 C12H22O11	000099-20-7 43



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102159.D
 Acq On : 17 Jan 2018 19:27
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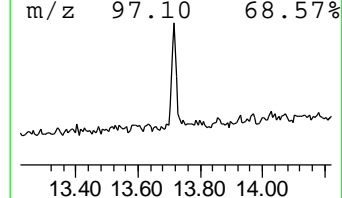
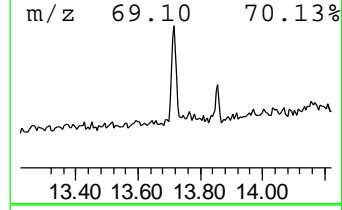
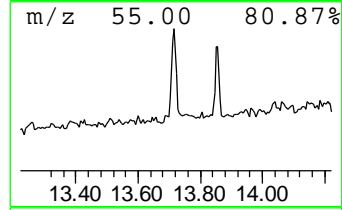
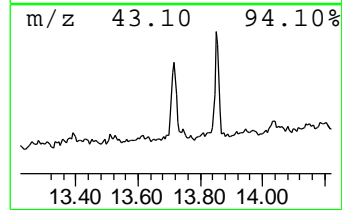
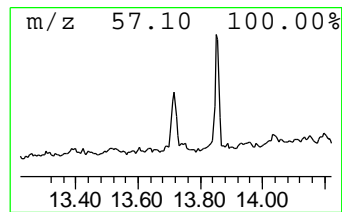
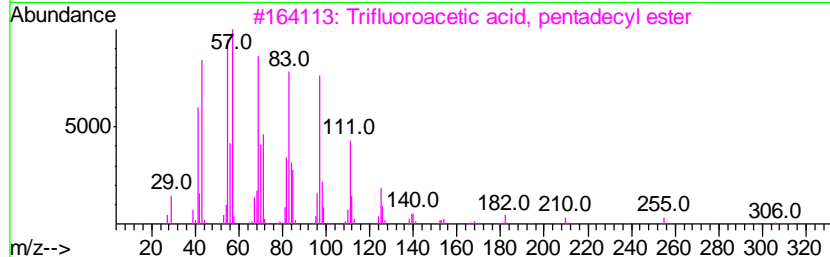
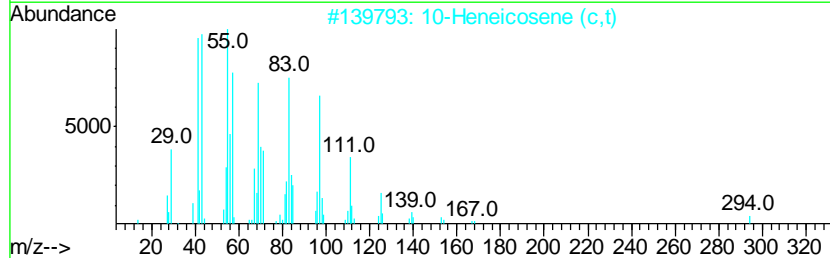
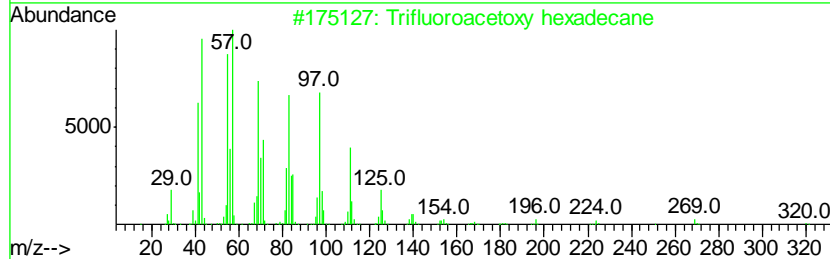
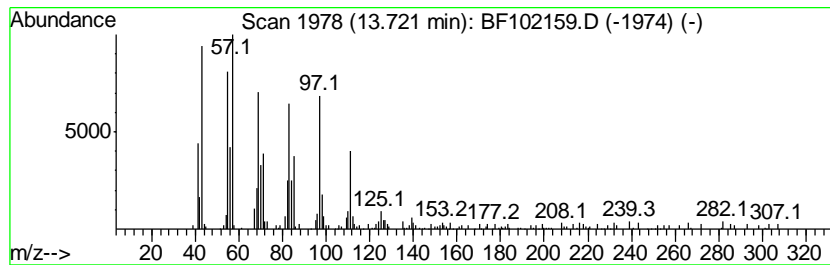
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 11 Trifluoroacetoxy hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	4.28 ng	275510	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2		10-Heneicosene (c,t)	294	C21H42	095008-11-0	92
3		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	91
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5		1-Heneicosyl formate	340	C22H44O2	077899-03-7	90



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
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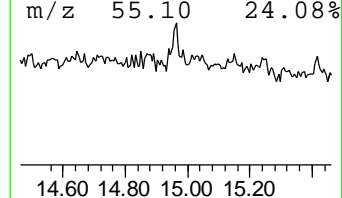
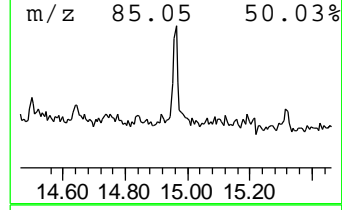
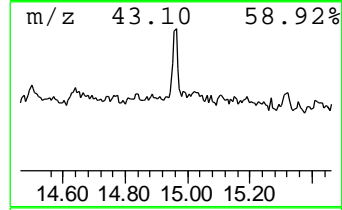
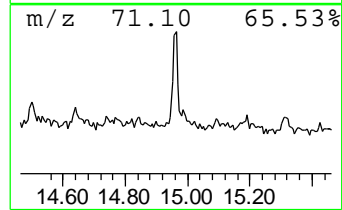
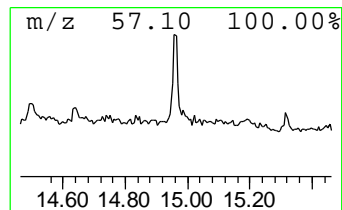
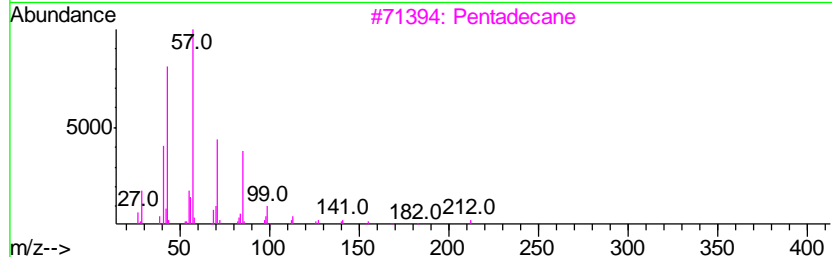
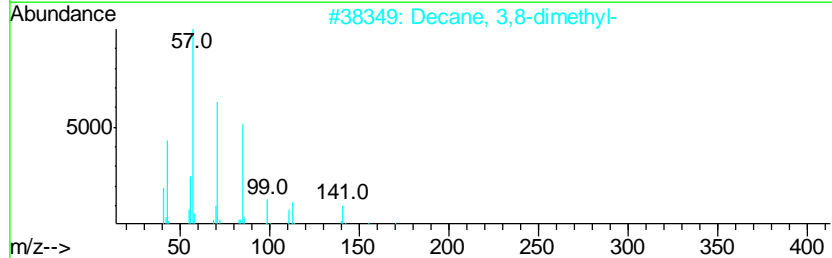
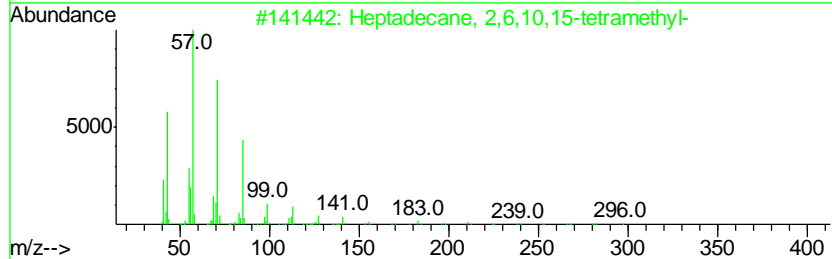
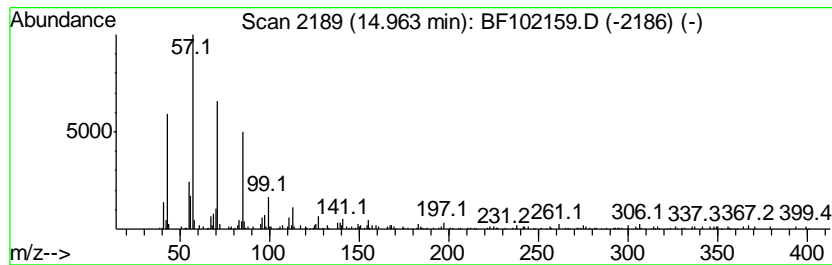
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 12 Heptadecane, 2,6,10,15-tetr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	3.26 ng	142776	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	93
3		Pentadecane	212	C15H32	000629-62-9	90
4		Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
5		Eicosane	282	C20H42	000112-95-8	90



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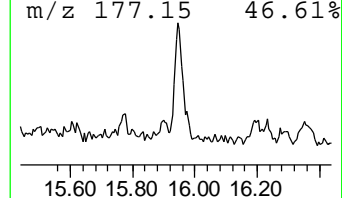
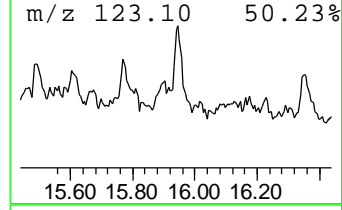
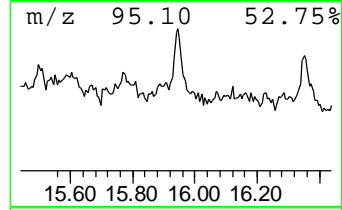
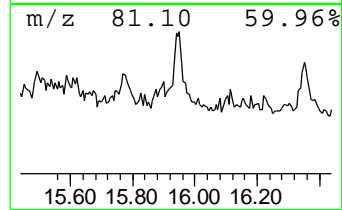
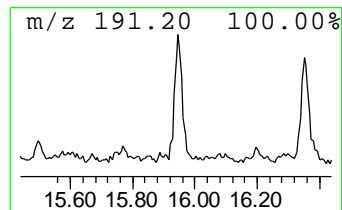
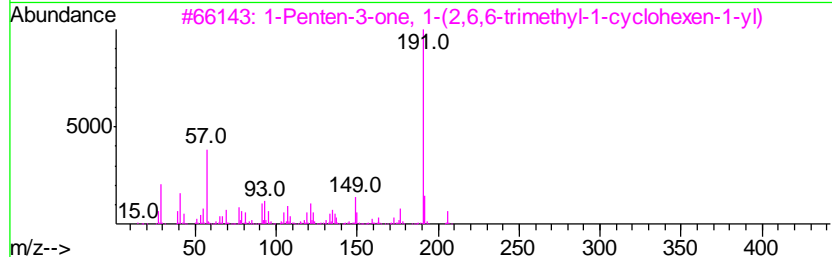
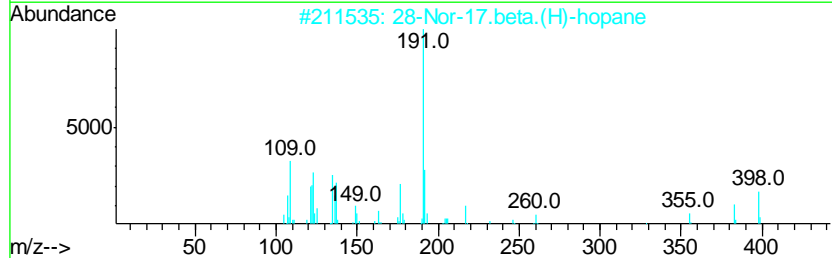
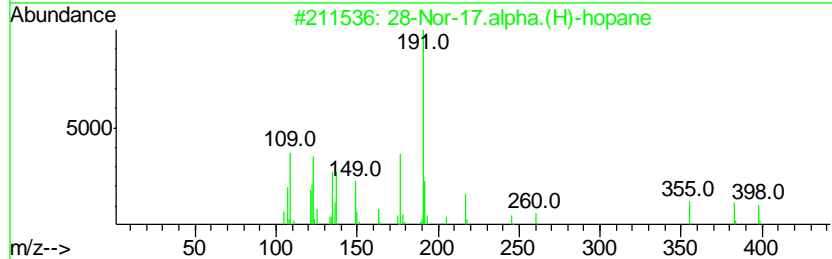
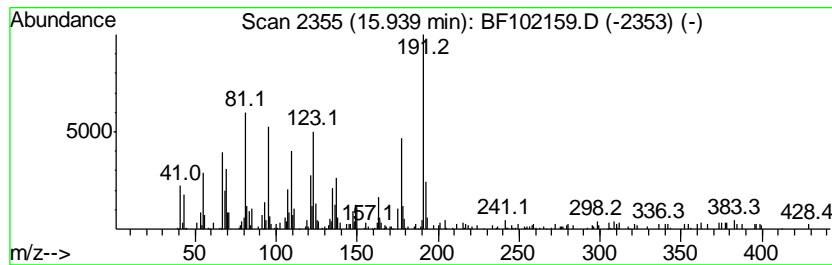
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 13 28-Nor-17.alpha.(H)-hopane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	9.38 ng	410348	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	83
2		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	45
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
4		9-Undecylanthracene	332	C25H32	132034-75-4	27
5		Terephthalic acid, 4-chloro-3-me...	332	C18H17ClO4	1000323-69-7	22



Data Path : U:\HPCHEM1\BNA_F\DATA\BF011718\
 Data File : BF102159.D
 Acq On : 17 Jan 2018 19:27
 Operator : SJ/JU
 Sample : J1144-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GR-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.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...	4.93	6.1	ng	270590	1	6.72	889392	20.0
unknown6.49	6.49	77.7	ng	3456170	1	6.72	889392	20.0
Cyclohexasiloxane...	8.52	7.6	ng	460416	2	8.00	1213390	20.0
2-Hydroxy-iso-but...	8.56	3.2	ng	191802	2	8.00	1213390	20.0
unknown8.85	8.85	3.4	ng	203520	2	8.00	1213390	20.0
unknown10.25	10.25	2.4	ng	131765	3	9.75	1115750	20.0
Benzophenone	10.47	12.1	ng	675211	3	9.75	1115750	20.0
n-Hexadecanoic acid	11.77	11.6	ng	517139	4	11.23	890781	20.0
Octadecanoic acid	12.55	2.3	ng	103289	4	11.23	890781	20.0
Trifluoroacetoxy ...	13.72	4.3	ng	275510	5	13.87	1287570	20.0
Heptadecane, 2,6,...	14.96	3.3	ng	142776	6	15.29	874937	20.0
28-Nor-17.alpha.(...	15.94	9.4	ng	410348	6	15.29	874937	20.0