

Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102168.D
 Acq On : 17 Jan 2018 23:28
 Operator : SJ/JU
 Sample : J1163-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-(0-4)

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	478	483	491	rVB	163886	225940	7.24%	0.751%
2	5.334	546	552	555	rBV	3093420	3023392	96.88%	10.050%
3	6.357	720	726	731	rBV	2783507	3120771	100.00%	10.374%
4	6.487	744	748	752	rBV	3018935	3108799	99.62%	10.334%
5	6.722	784	788	791	rBV	1009812	888369	28.47%	2.953%
6	6.875	810	814	817	rBV	2720399	2388154	76.52%	7.939%
7	7.287	878	884	887	rBV	2048488	1893043	60.66%	6.293%
8	7.745	959	962	965	rVB	43881	37240	1.19%	0.124%
9	8.004	1001	1006	1008	rBV	1243299	1155900	37.04%	3.842%
10	8.557	1097	1100	1104	rBV	284612	237803	7.62%	0.790%
11	9.081	1183	1189	1192	rBV	3640862	3022663	96.86%	10.048%
12	9.475	1252	1256	1259	rBV	597482	478634	15.34%	1.591%
13	9.751	1300	1303	1307	rBV	1260327	1125213	36.06%	3.740%
14	10.469	1419	1425	1428	rBV	1250269	976365	31.29%	3.246%
15	10.539	1433	1437	1441	rBV	1734525	1607169	51.50%	5.342%
16	11.233	1549	1555	1559	rBV	1104572	928795	29.76%	3.087%
17	12.645	1792	1795	1797	rBV	53468	42371	1.36%	0.141%
18	12.821	1820	1825	1828	rBV	2089329	1901399	60.93%	6.321%
19	13.715	1973	1977	1984	rBV	302812	323622	10.37%	1.076%
20	13.868	1998	2003	2006	rBV	728080	725393	23.24%	2.411%
21	14.039	2028	2032	2037	rBV4	30846	37016	1.19%	0.123%
22	14.345	2080	2084	2090	rBV2	174092	191851	6.15%	0.638%
23	14.774	2154	2157	2161	rVB2	35829	36497	1.17%	0.121%
24	14.957	2184	2188	2198	rVB	386550	442294	14.17%	1.470%
25	15.286	2238	2244	2247	rBV	583278	720853	23.10%	2.396%
26	15.315	2247	2249	2258	rVB2	55807	69874	2.24%	0.232%
27	15.504	2277	2281	2287	rVB	79811	101010	3.24%	0.336%
28	15.721	2313	2318	2323	rBV	237696	356619	11.43%	1.185%
29	15.768	2323	2326	2333	rVV2	79332	131043	4.20%	0.436%
30	15.839	2335	2338	2347	rVB9	22116	53012	1.70%	0.176%
31	16.192	2395	2398	2404	rVB2	34281	53728	1.72%	0.179%
32	16.468	2440	2445	2452	rVB2	91022	154687	4.96%	0.514%
33	16.751	2488	2493	2497	rVB4	47558	77507	2.48%	0.258%
34	16.809	2499	2503	2513	rVB5	74643	175118	5.61%	0.582%

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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-1-(0-4)

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	16.915	2519	2521	2527	rVB6	20615	36372	1.17%	0.121%
36	17.192	2564	2568	2580	rVB6	58664	149801	4.80%	0.498%
37	18.168	2729	2734	2746	rVB6	32032	84592	2.71%	0.281%

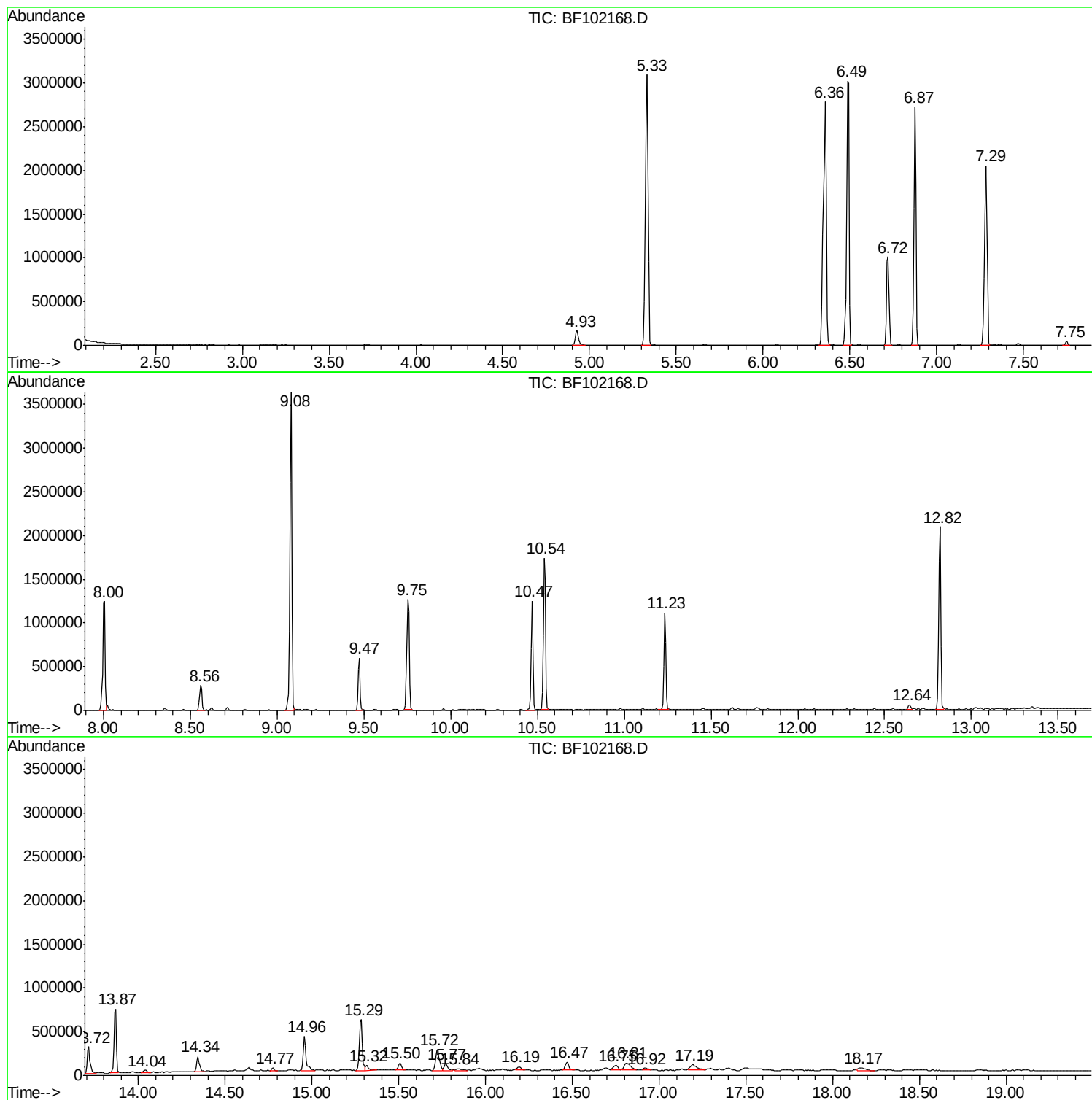
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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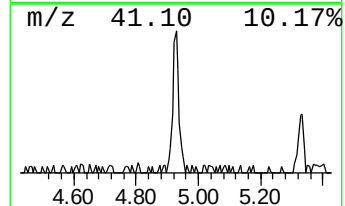
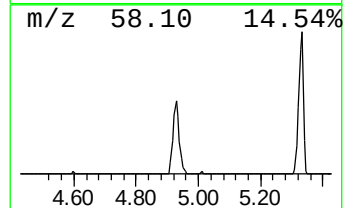
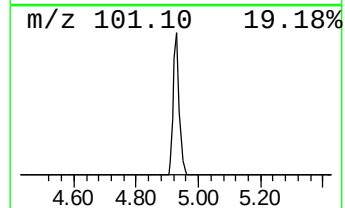
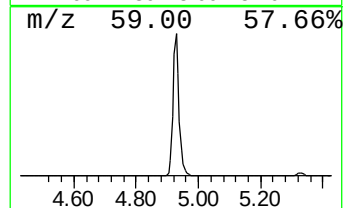
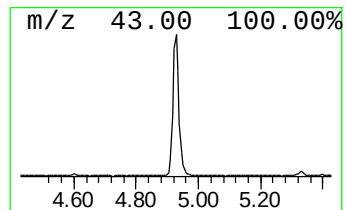
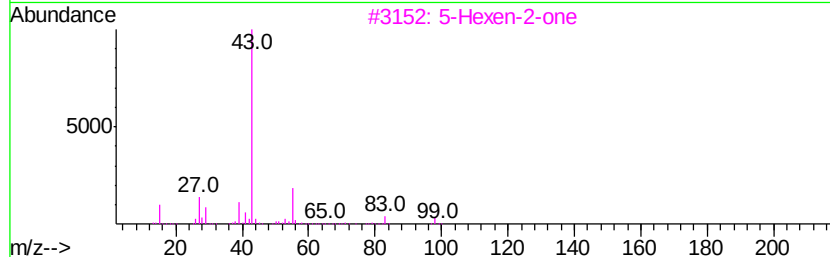
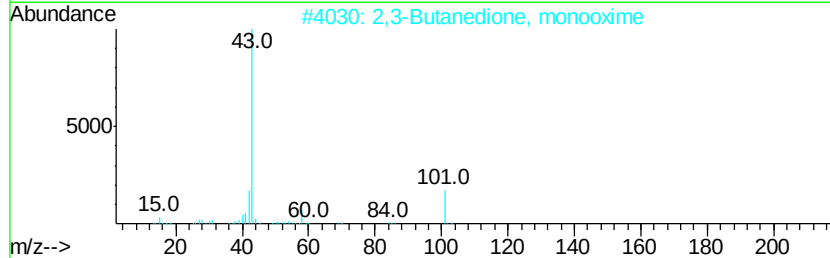
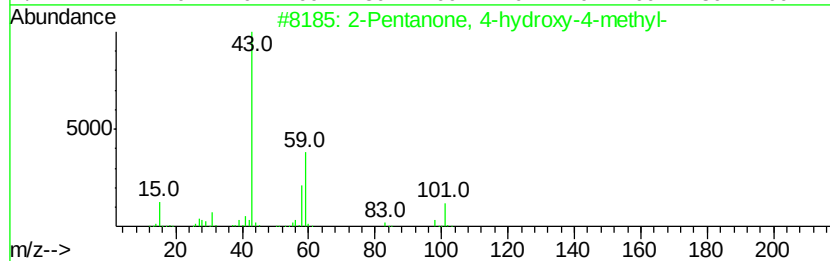
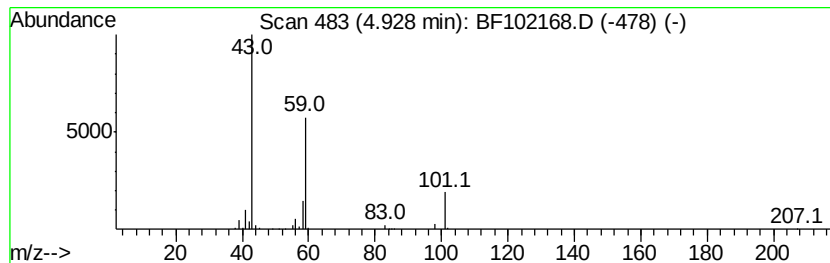
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	5.09 ng	225940	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			.alpha.-D-Galactopyranoside, met...	249	C10H19NO6	017296-07-0	9
5			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9



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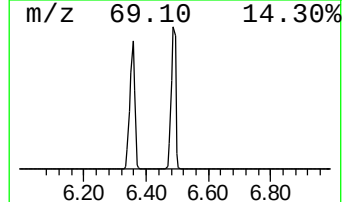
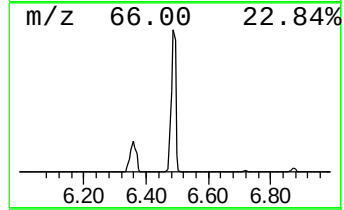
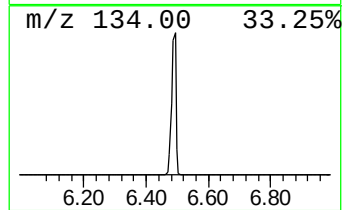
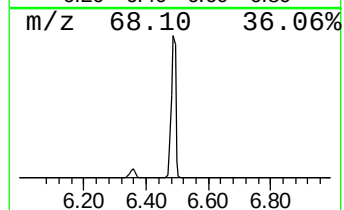
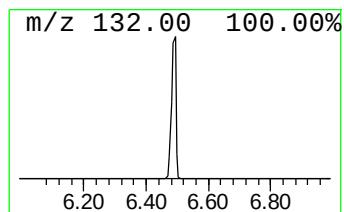
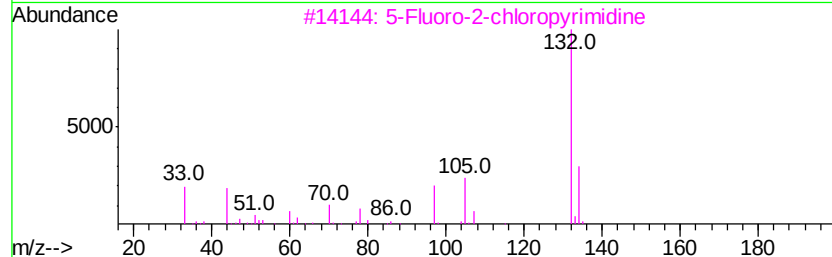
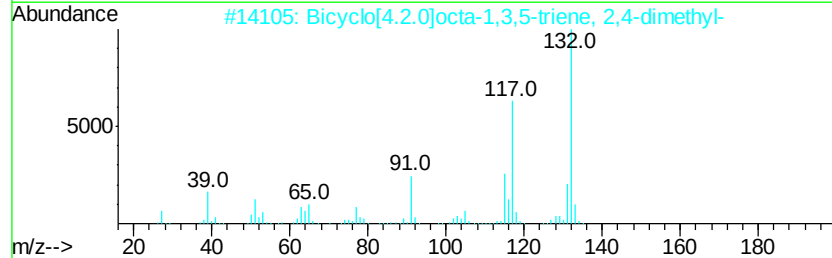
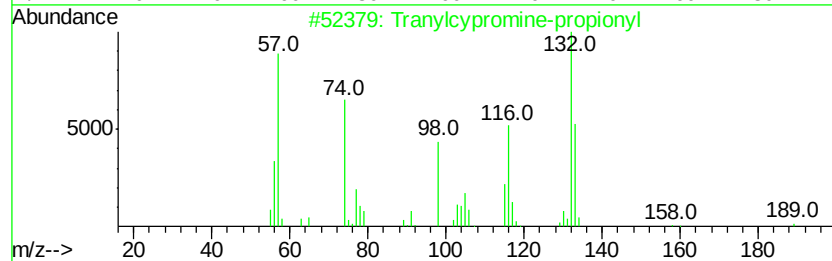
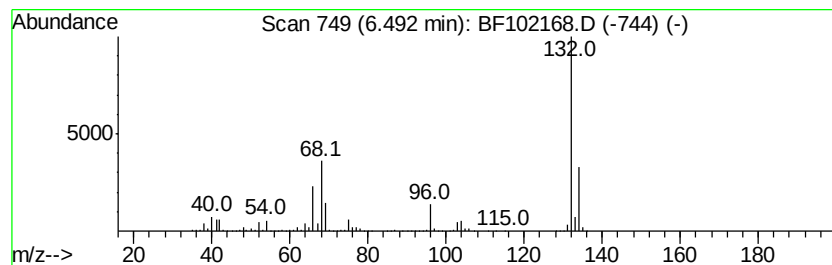
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	69.99 ng	3108800	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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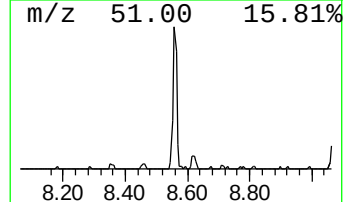
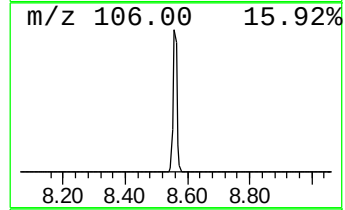
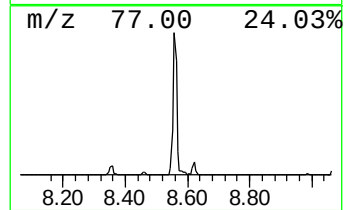
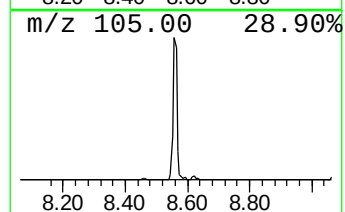
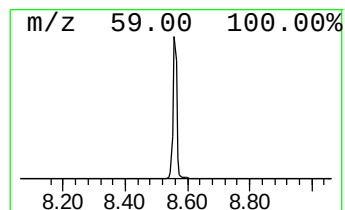
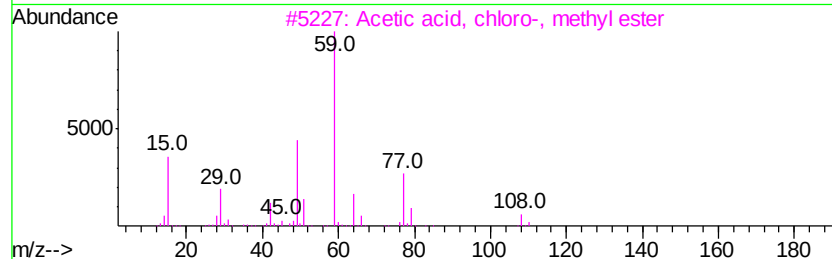
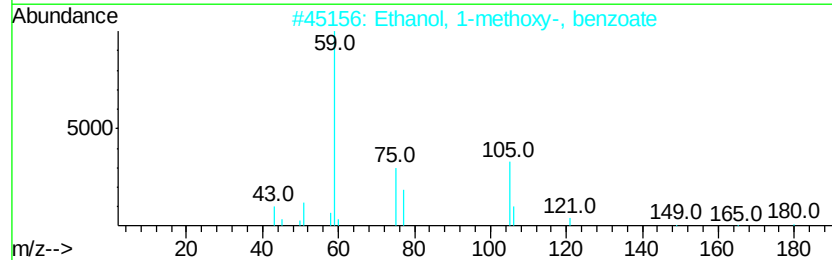
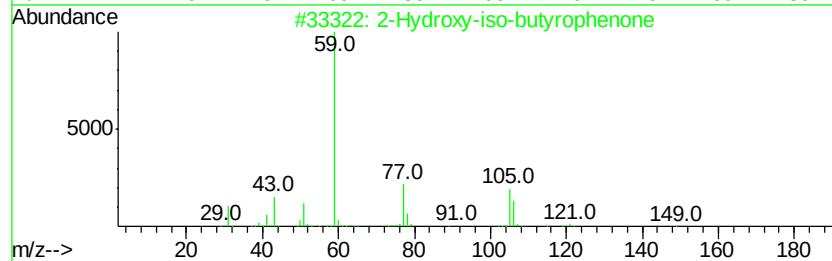
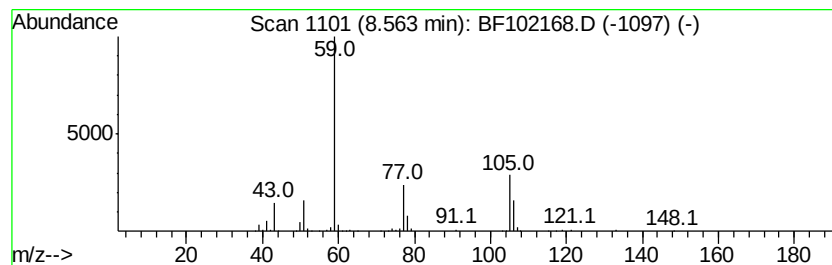
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Hydroxy-iso-butyrophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	4.11 ng	237803	Napthalene-d8	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hydroxy-iso-butyrophenone	164	C10H12O2	007473-98-5	83
2		Ethanol, 1-methoxy-, benzoate	180	C10H12O3	051835-44-0	72
3		Acetic acid, chloro-, methyl ester	108	C3H5ClO2	000096-34-4	9
4		Guanidine	59	CH5N3	000113-00-8	7
5		Formamide, N-methyl-	59	C2H5NO	000123-39-7	7



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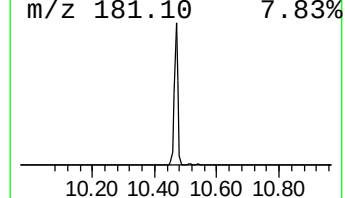
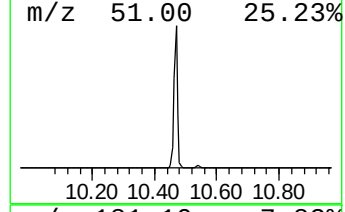
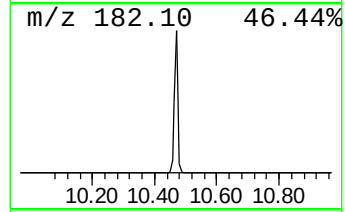
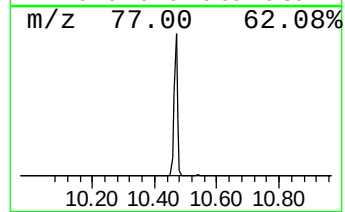
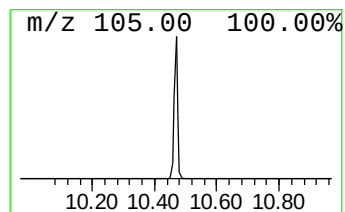
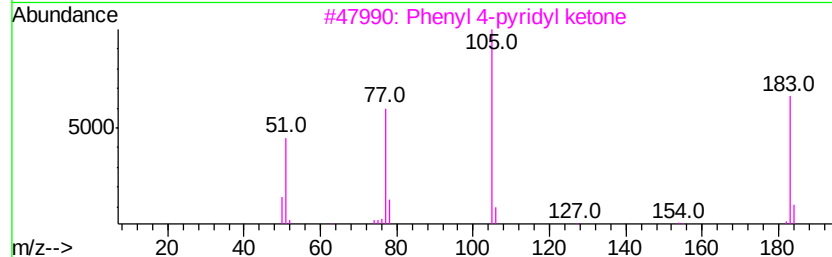
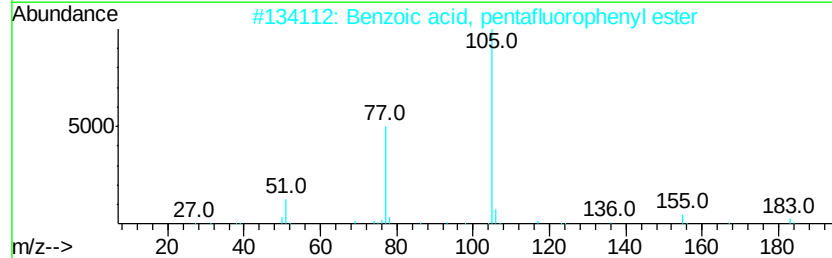
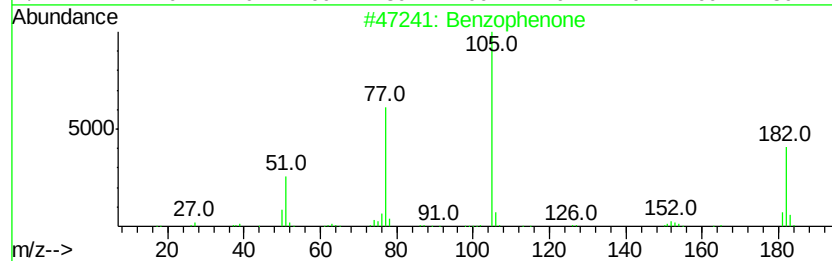
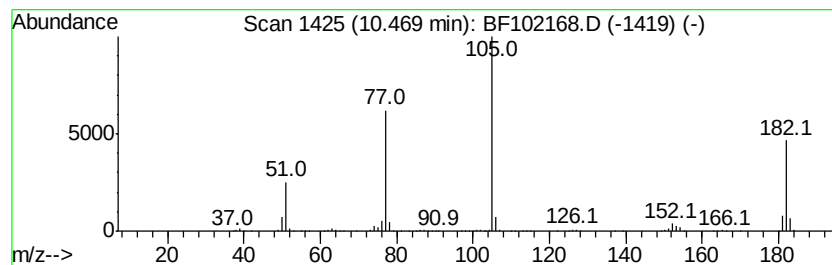
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	17.35 ng	976365	Acenaphthene-d10	9.75
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzophenone	182 C13H10O	000119-61-9 97
2		Benzoic acid, pentafluorophenyl ...	288 C13H5F5O2	1000360-67-9 50
3		Phenvl 4-pyridyl ketone	183 C12H9NO	014548-46-0 50
4		Benzeneacetic acid, .alpha.-oxo-...	164 C9H8O3	015206-55-0 49
5		1-Propanone, 2-bromo-1-phenyl-	212 C9H9BrO	002114-00-3 47



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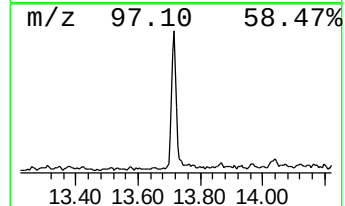
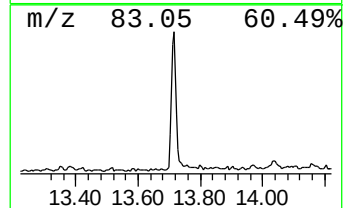
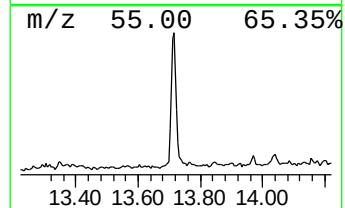
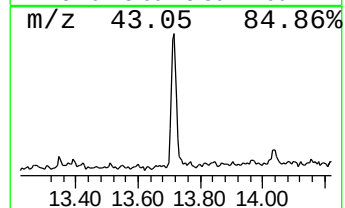
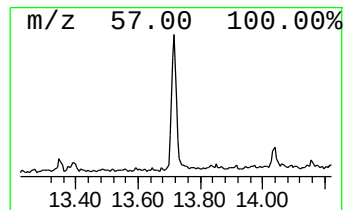
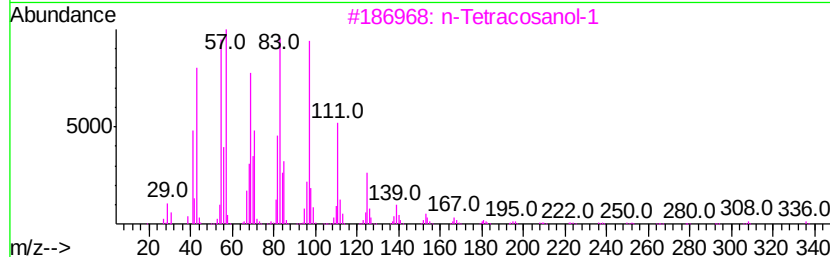
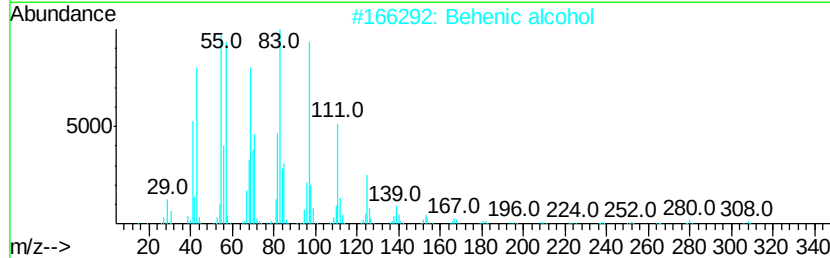
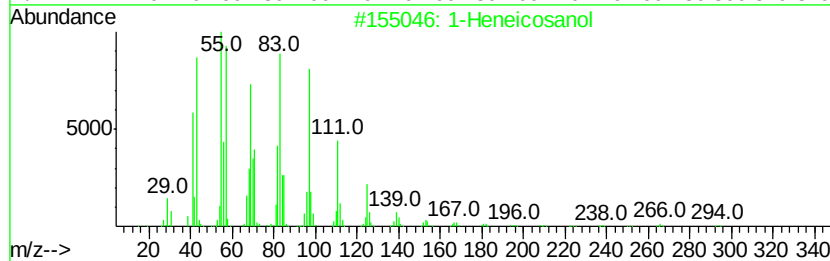
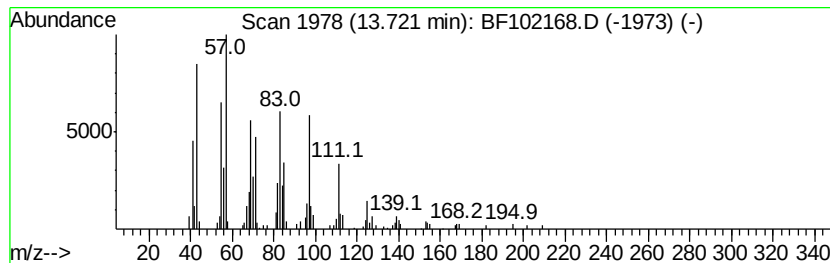
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.72	8.92 ng	323622	Chrysene-d12	13.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	94
2	Behenic alcohol	326	C22H46O	000661-19-8	93
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4	Octacosanol	410	C28H58O	000557-61-9	91
5	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91



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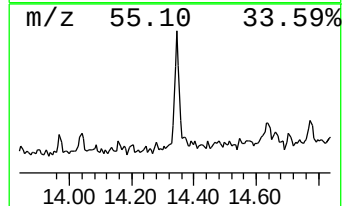
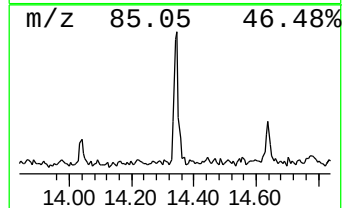
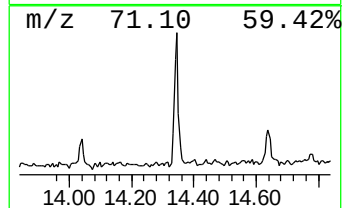
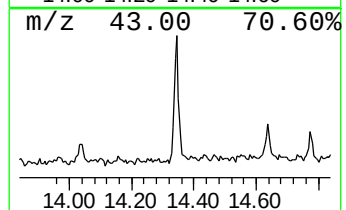
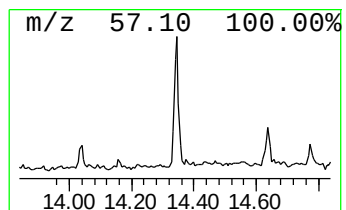
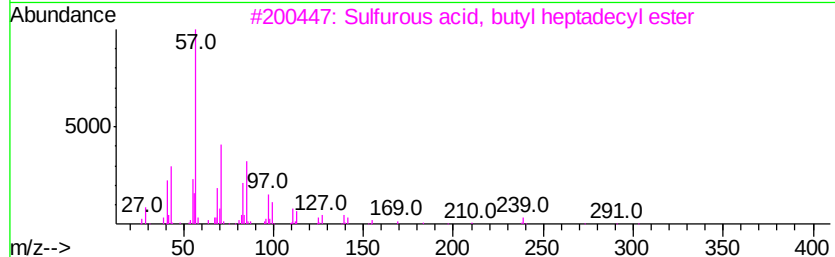
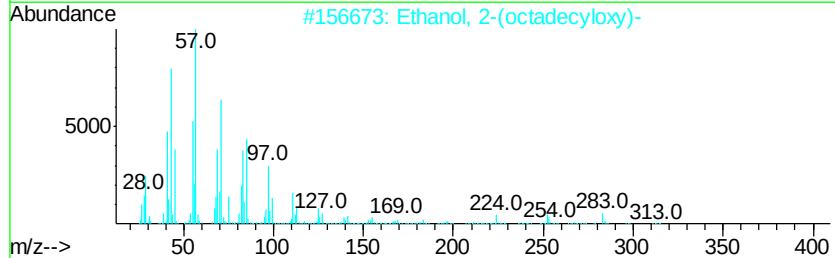
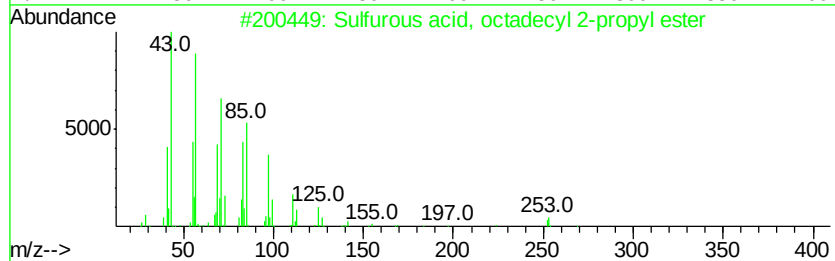
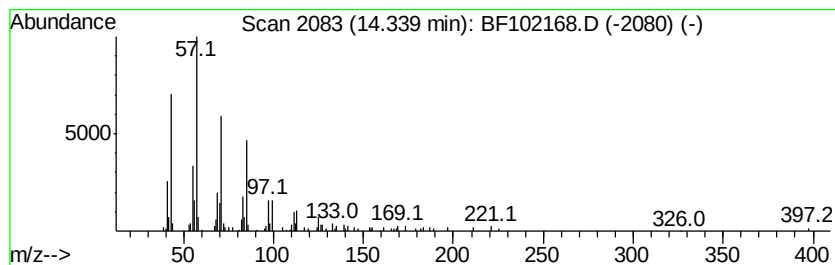
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Sulfurous acid, octadecyl 2... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	5.29 ng	191851	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	91
2		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	90
3		Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	87
4		Hexacosane	366	C26H54	000630-01-3	68
5		Octadecane	254	C18H38	000593-45-3	62



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
Data File : BF102168.D
Acq On : 17 Jan 2018 23:28
Operator : SJ/JU
Sample : J1163-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

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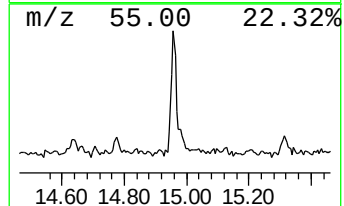
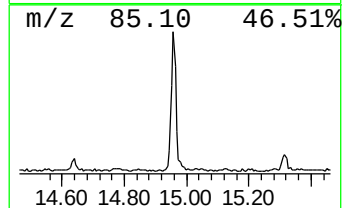
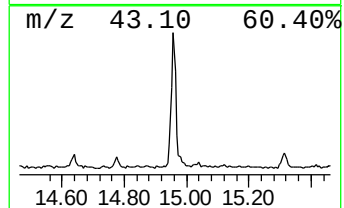
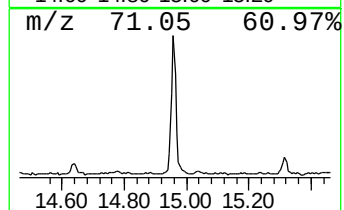
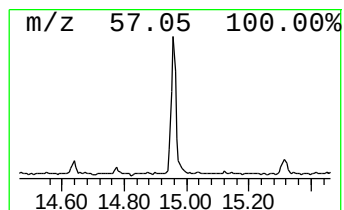
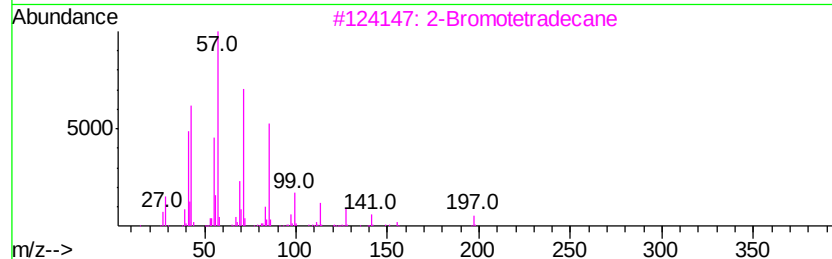
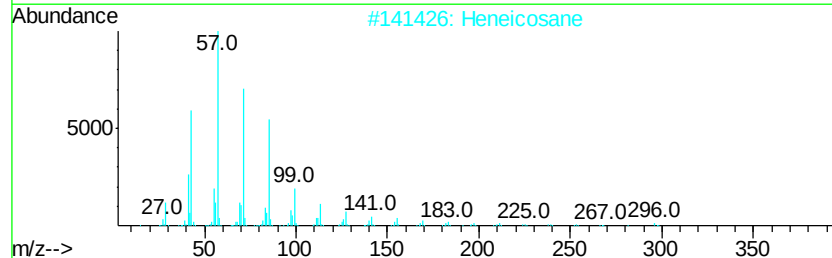
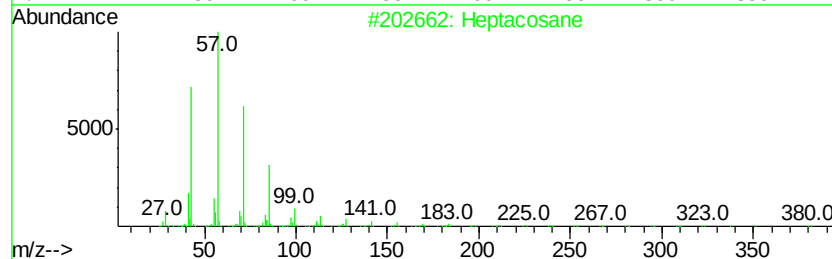
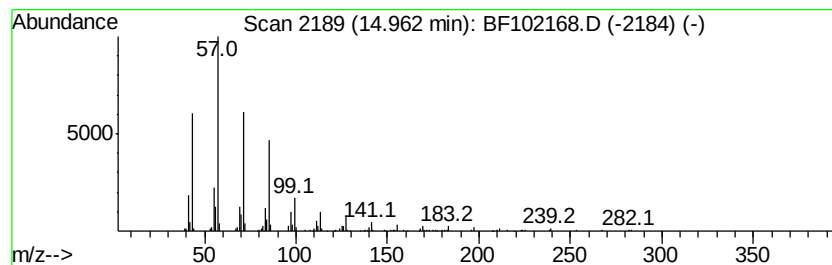
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	12.27 ng	442294	Perylene-d12	15.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Heneicosane		296	C21H44	000629-94-7	90
3	2-Bromotetradecane		276	C14H29Br	074036-95-6	90
4	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90
5	Triacontane		422	C30H62	000638-68-6	90



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
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Sample : J1163-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

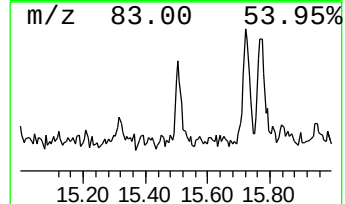
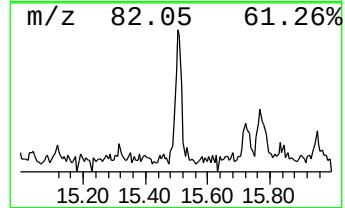
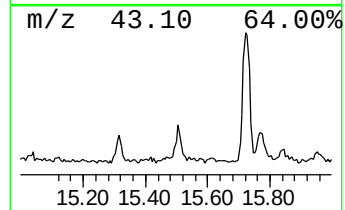
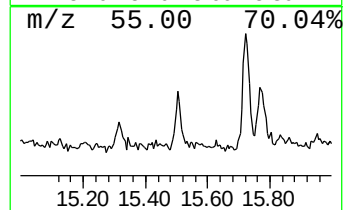
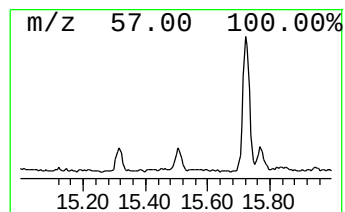
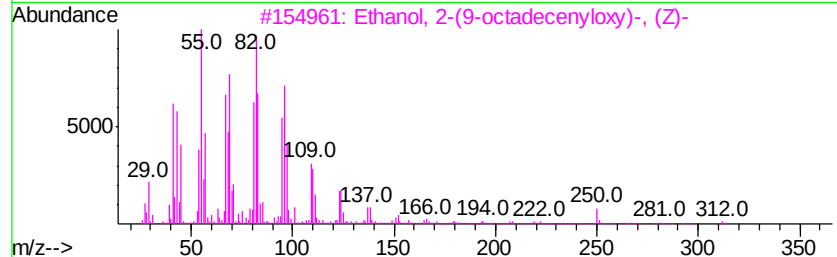
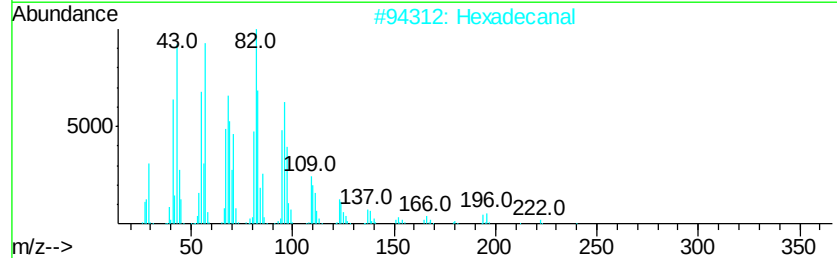
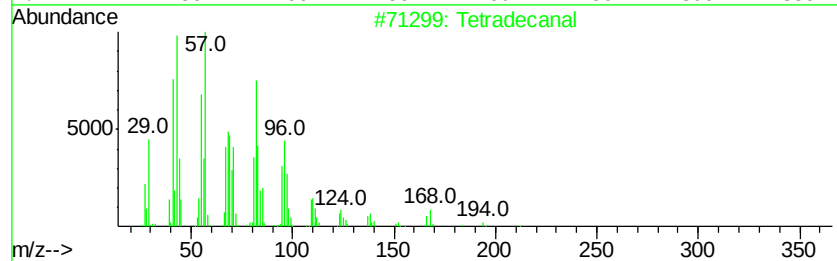
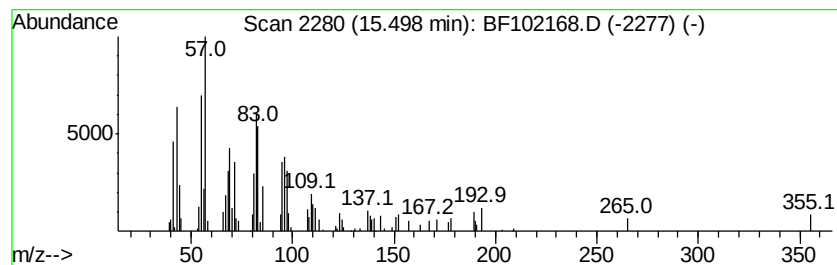
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetradecanal Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.50	2.80 ng	101010	Perylene-d12	15.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanal	212	C14H28O	000124-25-4	86
2	Hexadecanal	240	C16H32O	000629-80-1	72
3	Ethanol, 2-(9-octadecenyloxy)-, ...	312	C20H40O2	005353-25-3	68
4	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	64
5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	58



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
Data File : BF102168.D
Acq On : 17 Jan 2018 23:28
Operator : SJ/JU
Sample : J1163-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

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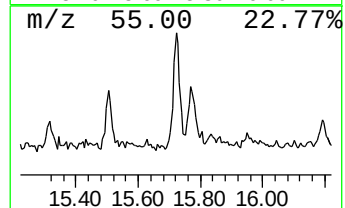
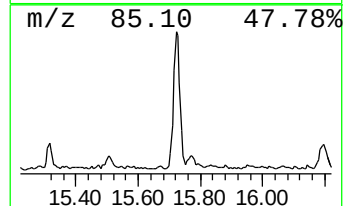
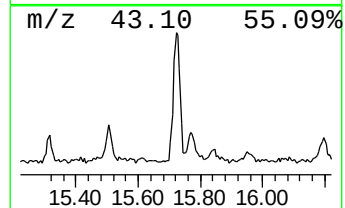
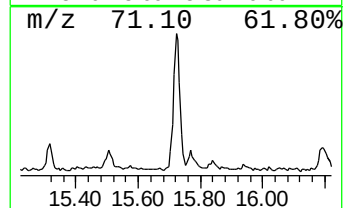
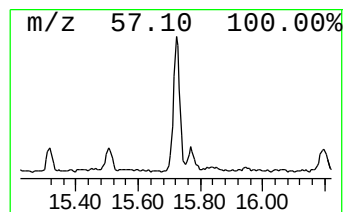
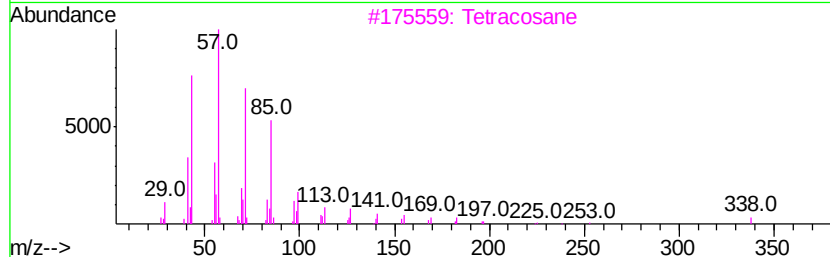
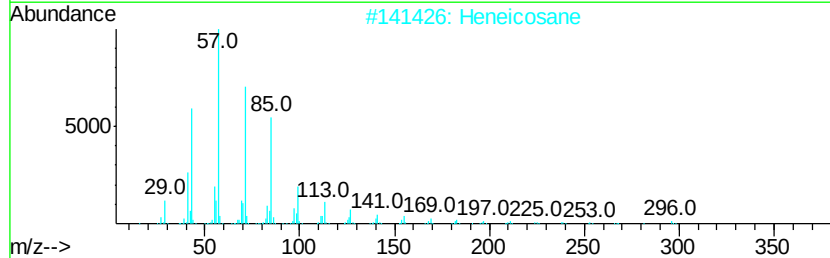
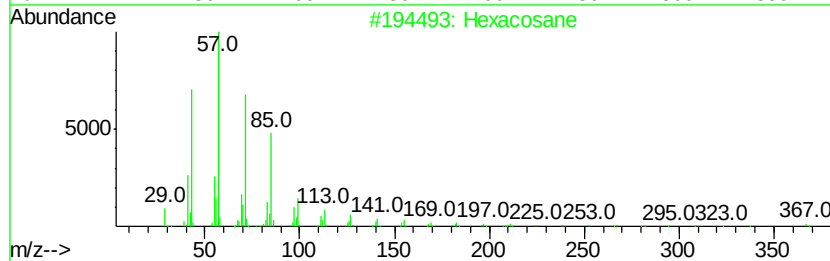
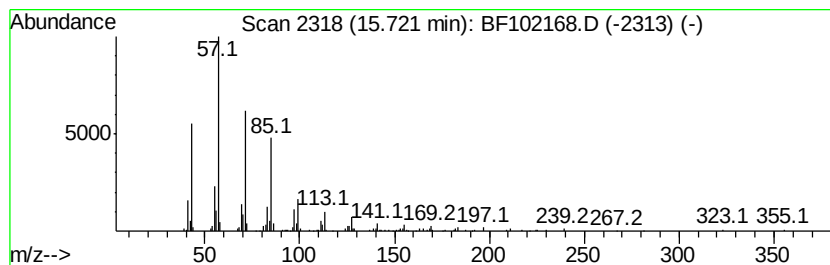
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	9.89 ng	356619	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetracosane	338	C24H50	000646-31-1	90
4		Heptadecane	240	C17H36	000629-78-7	87
5		Docosane	310	C22H46	000629-97-0	87



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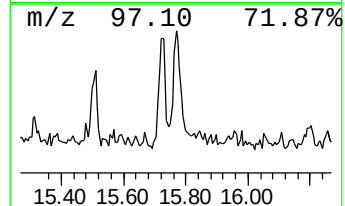
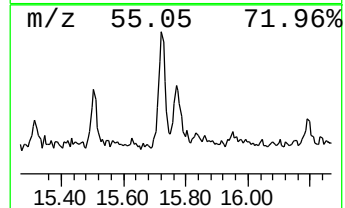
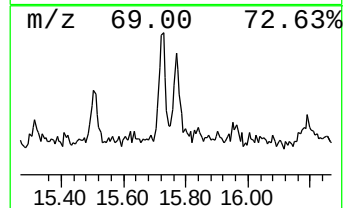
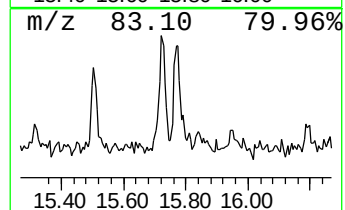
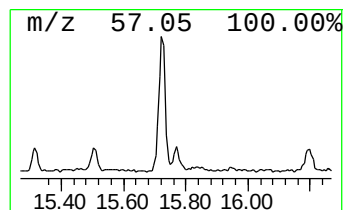
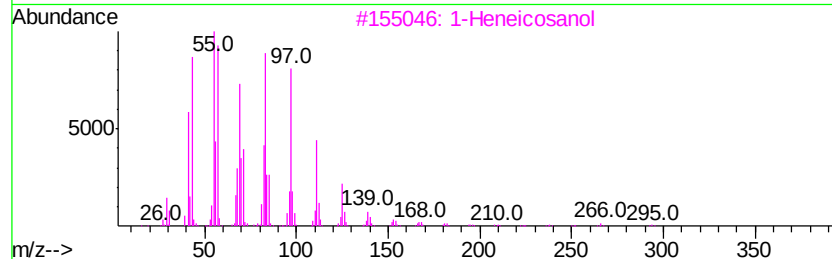
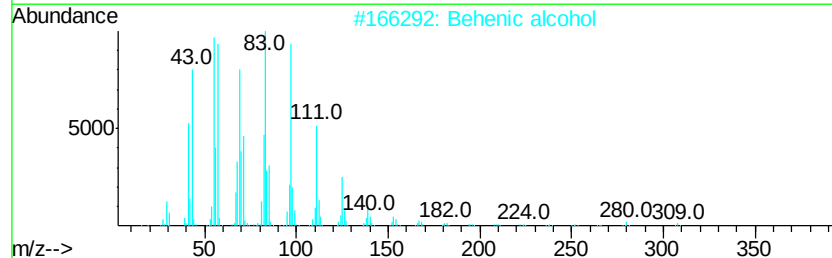
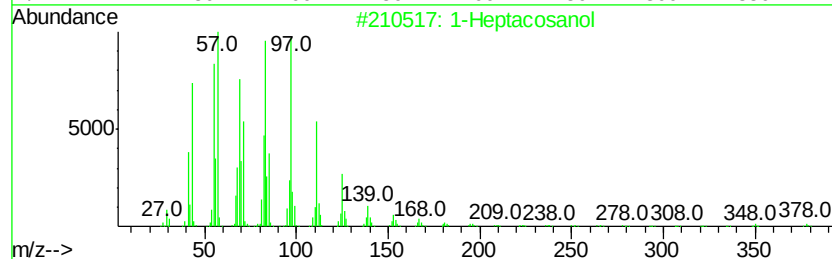
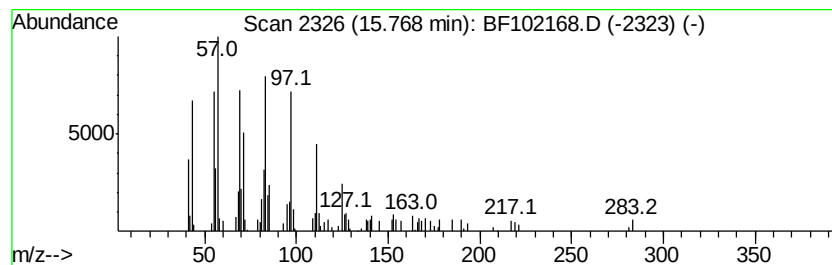
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Heptacosanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.77	3.64 ng	131043	Perylene-d12	15.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	74
2	Behenic alcohol	326	C22H46O	000661-19-8	70
3	1-Heneicosanol	312	C21H44O	015594-90-8	62
4	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	62
5	Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	62



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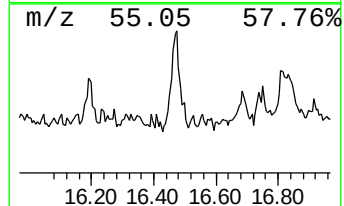
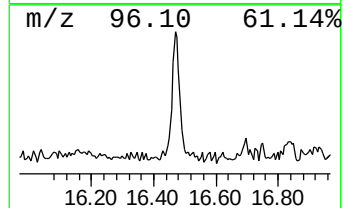
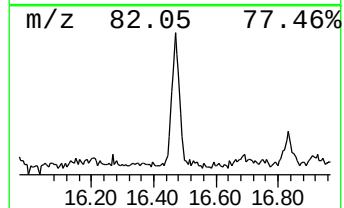
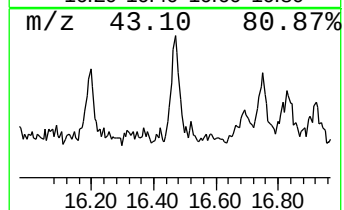
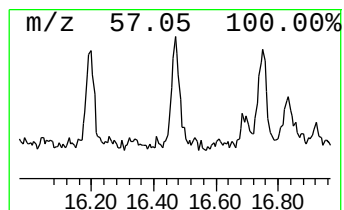
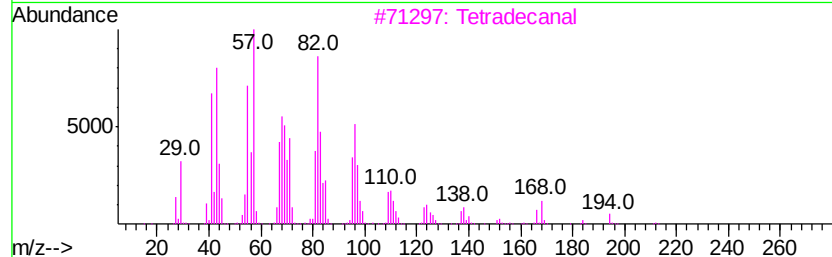
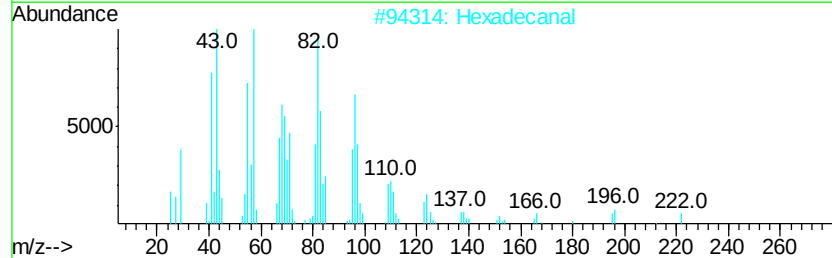
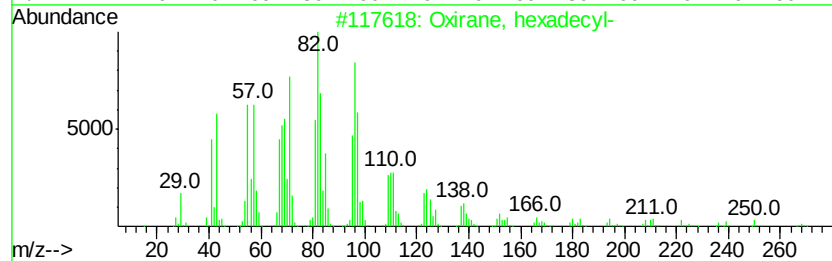
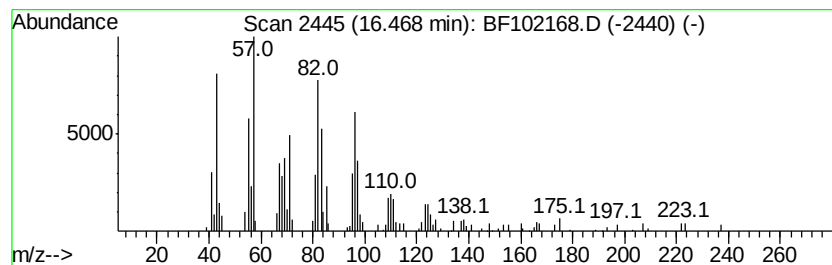
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Oxirane, hexadecyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	4.29 ng	154687	Perylene-d12	15.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2			Hexadecanal	240	C16H32O	000629-80-1	91
3			Tetradecanal	212	C14H28O	000124-25-4	90
4			1,19-Eicosadiene	278	C20H38	014811-95-1	76
5			18-Nonadecen-1-ol	282	C19H38O	1000142-89-2	72



Data Path : U:\HPCHEM1\BNA_F\DATA\BF011718\
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Acq On : 17 Jan 2018 23:28
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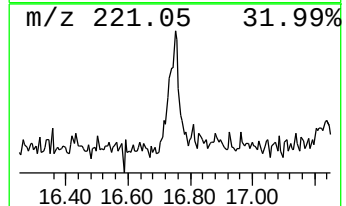
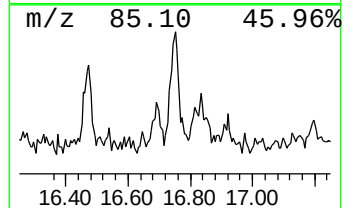
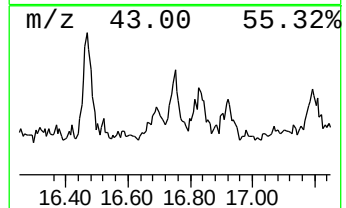
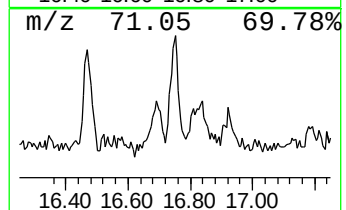
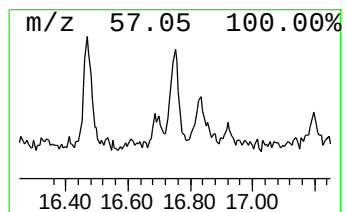
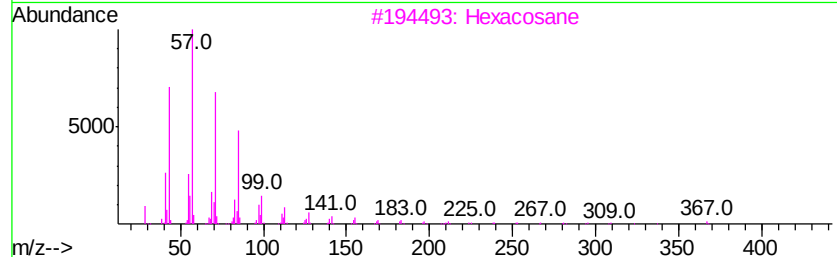
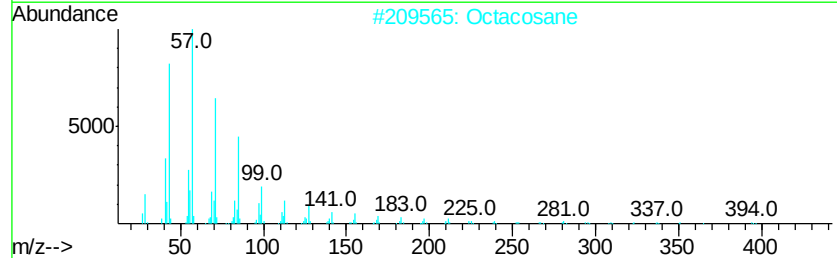
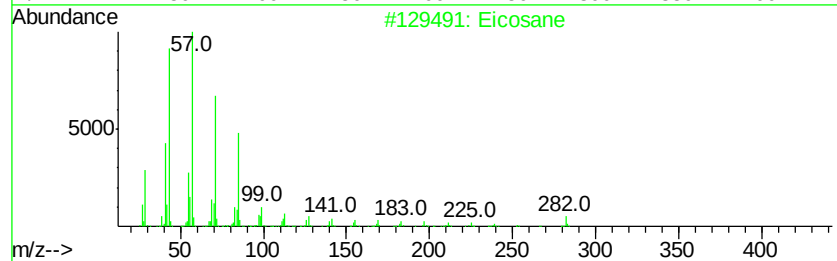
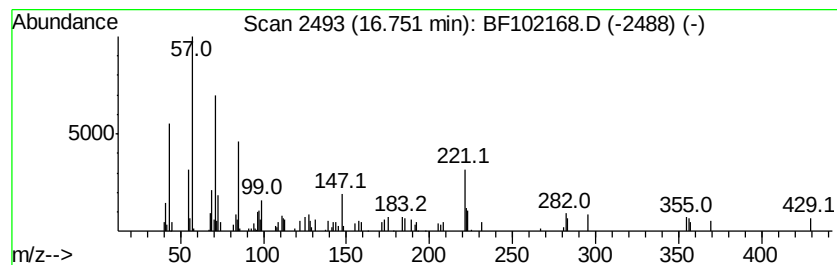
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.75	2.15 ng	77507	Perylene-d12	15.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	70
2	Octacosane	394 C28H58	000630-02-4	43
3	Hexacosane	366 C26H54	000630-01-3	43
4	Pentacosane	352 C25H52	000629-99-2	43
5	Decane, 2,9-dimethyl-	170 C12H26	001002-17-1	38



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
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Acq On : 17 Jan 2018 23:28
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Misc :
ALS Vial : 23 Sample Multiplier: 1

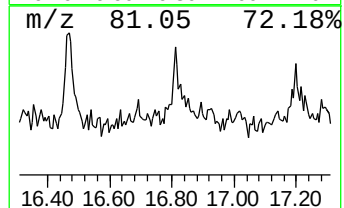
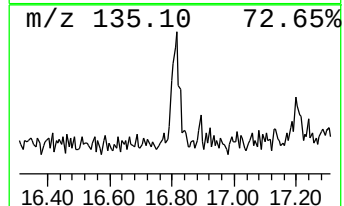
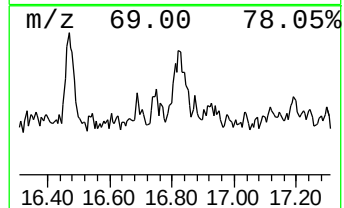
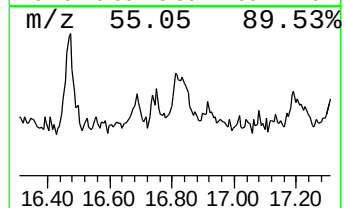
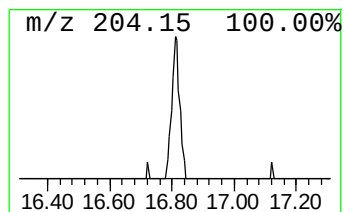
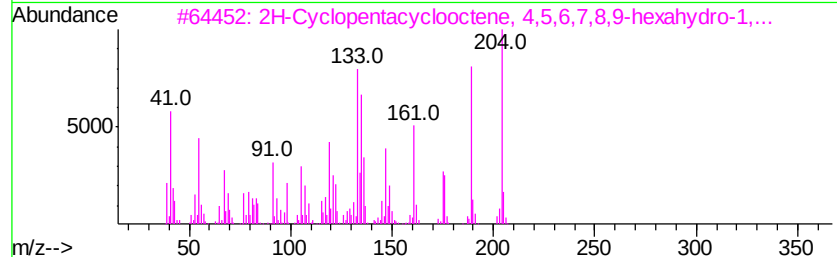
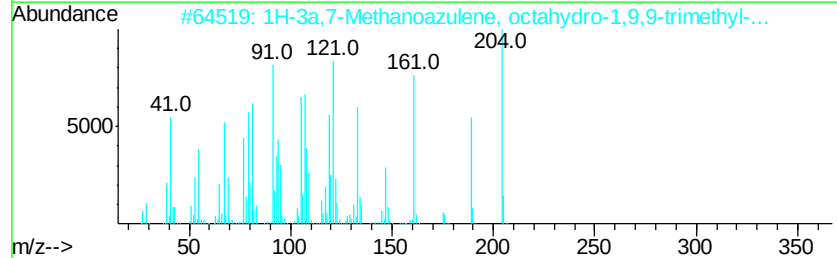
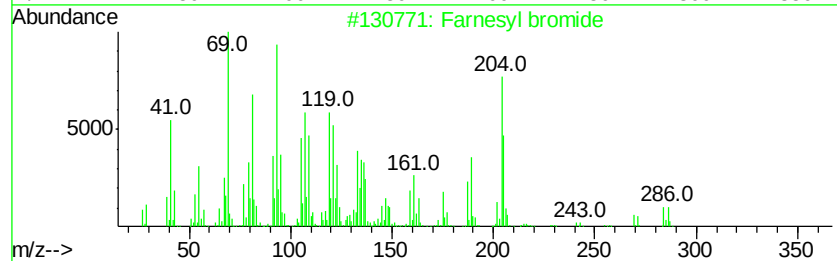
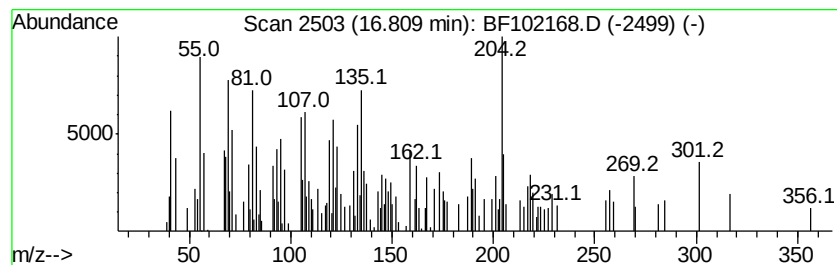
Instrument :
BNA_F
ClientSampleId :
TP-1-(0-4)

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 14 Farnesyl bromide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.81	4.86 ng	175118	Perylene-d12	15.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Farnesyl bromide	284 C15H25Br	006874-67-5	55
2	1H-3a,7-Methanoazulene, octahydr...	204 C15H24	000508-55-4	50
3	2H-Cyclopentacyclooctene, 4,5,6,...	204 C15H24	1000221-85-8	49
4	1,1,4a-Trimethyl-5,6-dimethylene...	204 C15H24	1000193-60-8	49
5	Aromandendrene	204 C15H24	000489-39-4	46



Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
Data File : BF102168.D
Acq On : 17 Jan 2018 23:28
Operator : SJ/JU
Sample : J1163-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

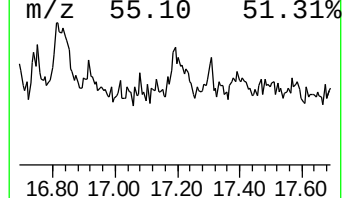
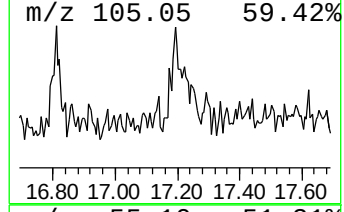
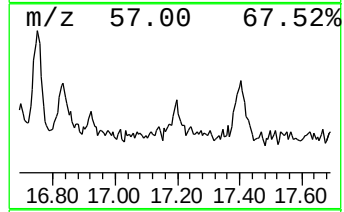
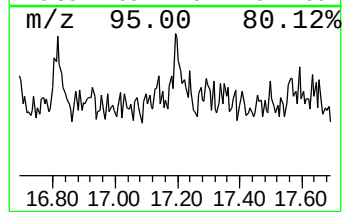
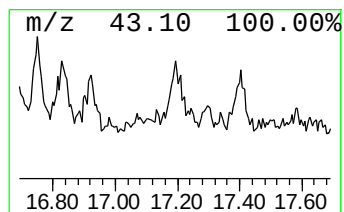
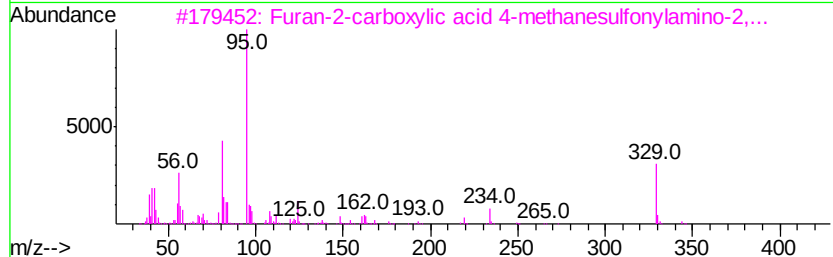
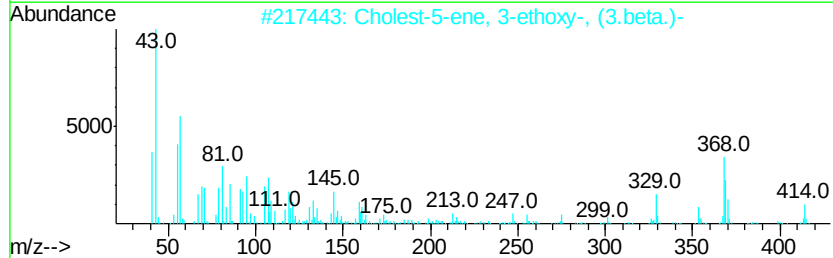
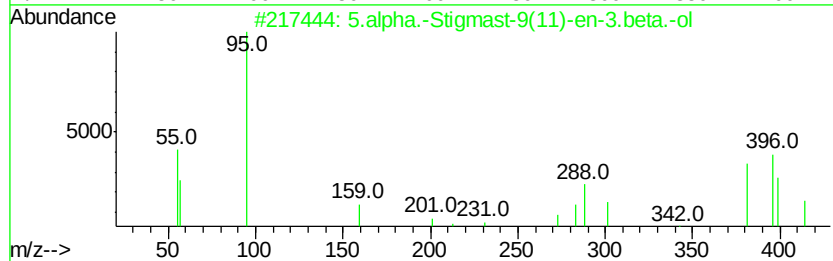
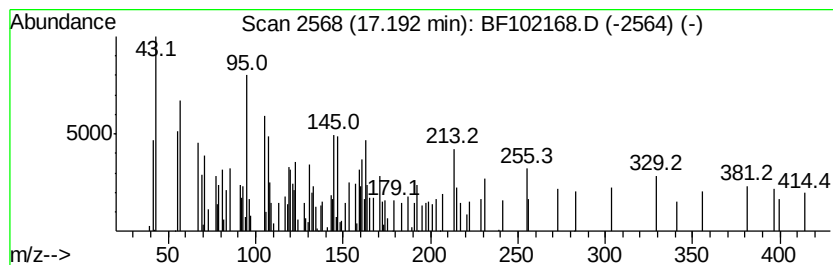
Instrument :
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ClientSampleId :
TP-1-(0-4)

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 15 unknown17.19 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.19	4.16 ng	149801	Perylene-d12	15.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5.alpha.-Stigmast-9(11)-en-3.beta...	414	C29H50O	077794-81-1	25
2	Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	15
3	Furan-2-carboxylic acid 4-methan...	344	C15H24N2O5S	1000300-98-9	10
4	2,20-Cyclo-8,9-secoaspidospermid...	396	C24H32N2O3	056145-37-0	9
5	Stigmastan-3,5-diene	396	C29H48	1000214-16-4	9



Data Path : U:\HPCHEM1\BNA_F\DATA\BF011718\
Data File : BF102168.D
Acq On : 17 Jan 2018 23:28
Operator : SJ/JU
Sample : J1163-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-1-(0-4)

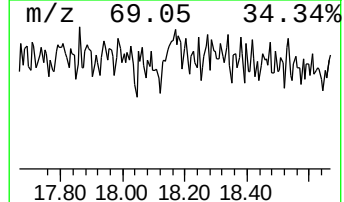
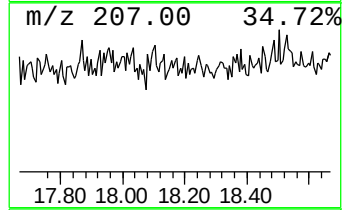
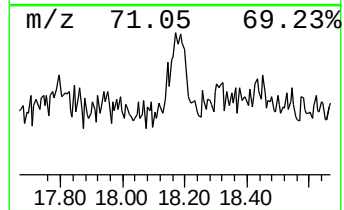
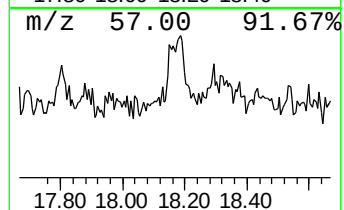
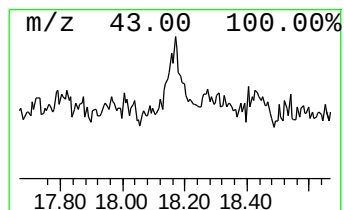
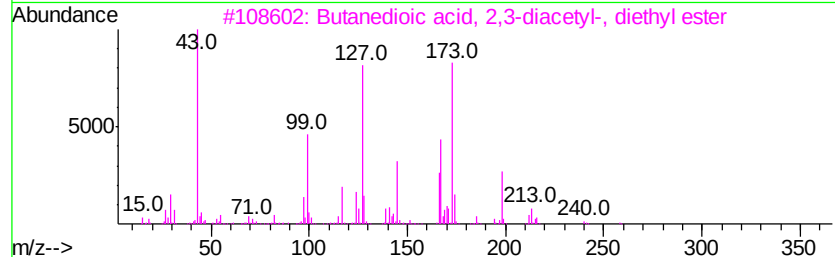
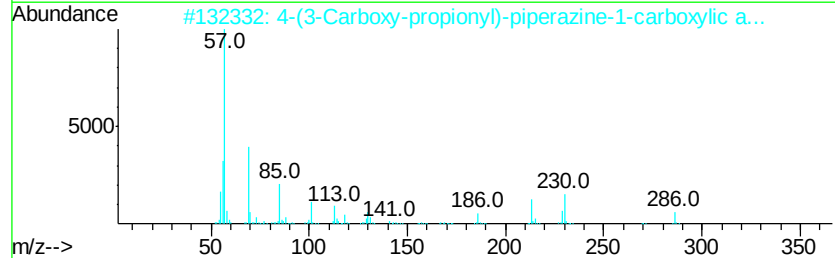
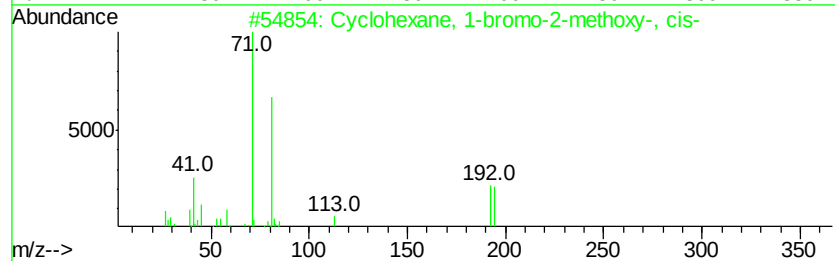
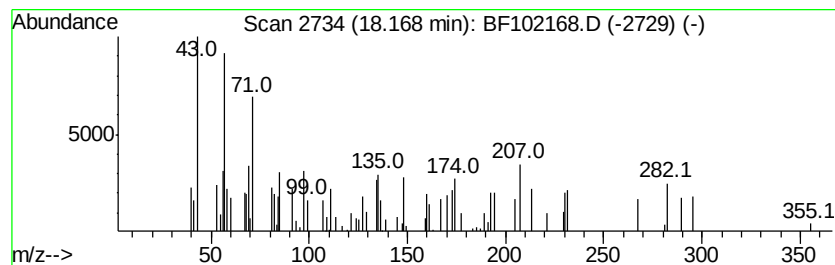
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown18.17 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	2.35 ng	84592	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-bromo-2-methoxy-,...	192	C7H13BrO	051332-48-0	10
2		4-(3-Carboxy-propionyl)-piperazi...	286	C13H22N2O5	288851-44-5	9
3		Butanedioic acid, 2,3-diacetyl-,...	258	C12H18O6	002049-86-7	9
4		Cyclohexane, 1-bromo-2-methoxy-,...	192	C7H13BrO	005927-93-5	9
5		Germane, dichlorodihexyl-	314	C12H26Cl2Ge	035551-43-0	7



Data Path : U:\HPCHEM1\BNA_F\DATA\BF011718\
 Data File : BF102168.D
 Acq On : 17 Jan 2018 23:28
 Operator : SJ/JU
 Sample : J1163-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-(0-4)

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	5.1 ng		225940	1	6.72	888369	20.0
unknown6.49	6.49	70.0 ng		3108800	1	6.72	888369	20.0
2-Hydroxy-iso-but...	8.56	4.1 ng		237803	2	8.00	1155900	20.0
Benzophenone	10.47	17.4 ng		976365	3	9.75	1125210	20.0
1-Heneicosanol	13.72	8.9 ng		323622	5	13.87	725393	20.0
Sulfurous acid, o...	14.34	5.3 ng		191851	5	13.87	725393	20.0
Heptacosane	14.96	12.3 ng		442294	6	15.29	720853	20.0
Tetradecanal	15.50	2.8 ng		101010	6	15.29	720853	20.0
Hexacosane	15.72	9.9 ng		356619	6	15.29	720853	20.0
1-Heptacosanol	15.77	3.6 ng		131043	6	15.29	720853	20.0
Oxirane, hexadecyl-	16.47	4.3 ng		154687	6	15.29	720853	20.0
Eicosane	16.75	2.1 ng		77507	6	15.29	720853	20.0
Farnesyl bromide	16.81	4.9 ng		175118	6	15.29	720853	20.0
unknown17.19	17.19	4.2 ng		149801	6	15.29	720853	20.0
unknown18.17	18.17	2.4 ng		84592	6	15.29	720853	20.0