

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
 Data File : BF098750.D
 Acq On : 23 Sep 2017 20:29
 Operator : SJ/JU
 Sample : I5340-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LT-TS-012

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-BF092317.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	4.087	328	340	343	rBV	6798110	14166772	100.00%	26.970%
2	5.163	518	523	535	rVB	581558	796034	5.62%	1.515%
3	5.545	582	588	591	rBV	3649852	4132970	29.17%	7.868%
4	6.545	751	758	761	rBV	3504737	4095639	28.91%	7.797%
5	6.692	778	783	786	rBV	4059967	4126785	29.13%	7.856%
6	6.922	818	822	825	rVB	1177320	943227	6.66%	1.796%
7	7.081	844	849	852	rVB	3310570	3146521	22.21%	5.990%
8	7.486	911	918	921	rBV	2311038	2601072	18.36%	4.952%
9	8.204	1035	1040	1043	rBV	1376914	1210672	8.55%	2.305%
10	9.280	1217	1223	1226	rBV	4315045	4221786	29.80%	8.037%
11	9.669	1283	1289	1292	rBV	507052	475811	3.36%	0.906%
12	9.957	1335	1338	1341	rVB	1367188	1113250	7.86%	2.119%
13	10.745	1464	1472	1475	rBV	2101685	1964408	13.87%	3.740%
14	11.445	1587	1591	1593	rBV	1047843	930472	6.57%	1.771%
15	11.957	1674	1678	1686	rVB2	346347	372135	2.63%	0.708%
16	12.651	1793	1796	1800	rBV	248669	255278	1.80%	0.486%
17	12.745	1808	1812	1821	rBV3	183046	193123	1.36%	0.368%
18	12.886	1830	1836	1840	rBV	224798	245965	1.74%	0.468%
19	13.027	1856	1860	1863	rVB	2914193	2604917	18.39%	4.959%
20	13.786	1985	1989	1992	rBV	176681	152192	1.07%	0.290%
21	13.910	2007	2010	2018	rVB2	453735	416796	2.94%	0.793%
22	14.080	2034	2039	2042	rBV	899023	867794	6.13%	1.652%
23	14.545	2114	2118	2124	rBV	430649	422892	2.99%	0.805%
24	15.004	2193	2196	2200	rBV	231744	233279	1.65%	0.444%
25	15.127	2213	2217	2221	rBV	113778	182348	1.29%	0.347%
26	15.204	2226	2230	2231	rBV2	168735	197569	1.39%	0.376%
27	15.221	2231	2233	2239	rVB2	384797	485032	3.42%	0.923%
28	15.568	2288	2292	2301	rBV	554314	786658	5.55%	1.498%
29	16.039	2368	2372	2377	rBV	160389	246099	1.74%	0.469%
30	16.092	2377	2381	2388	rVB2	119780	189371	1.34%	0.361%
31	16.868	2507	2513	2520	rBV3	97636	201412	1.42%	0.383%
32	17.262	2575	2580	2586	rBV4	76944	155301	1.10%	0.296%
33	17.680	2644	2651	2660	rBV3	169294	394866	2.79%	0.752%

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ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

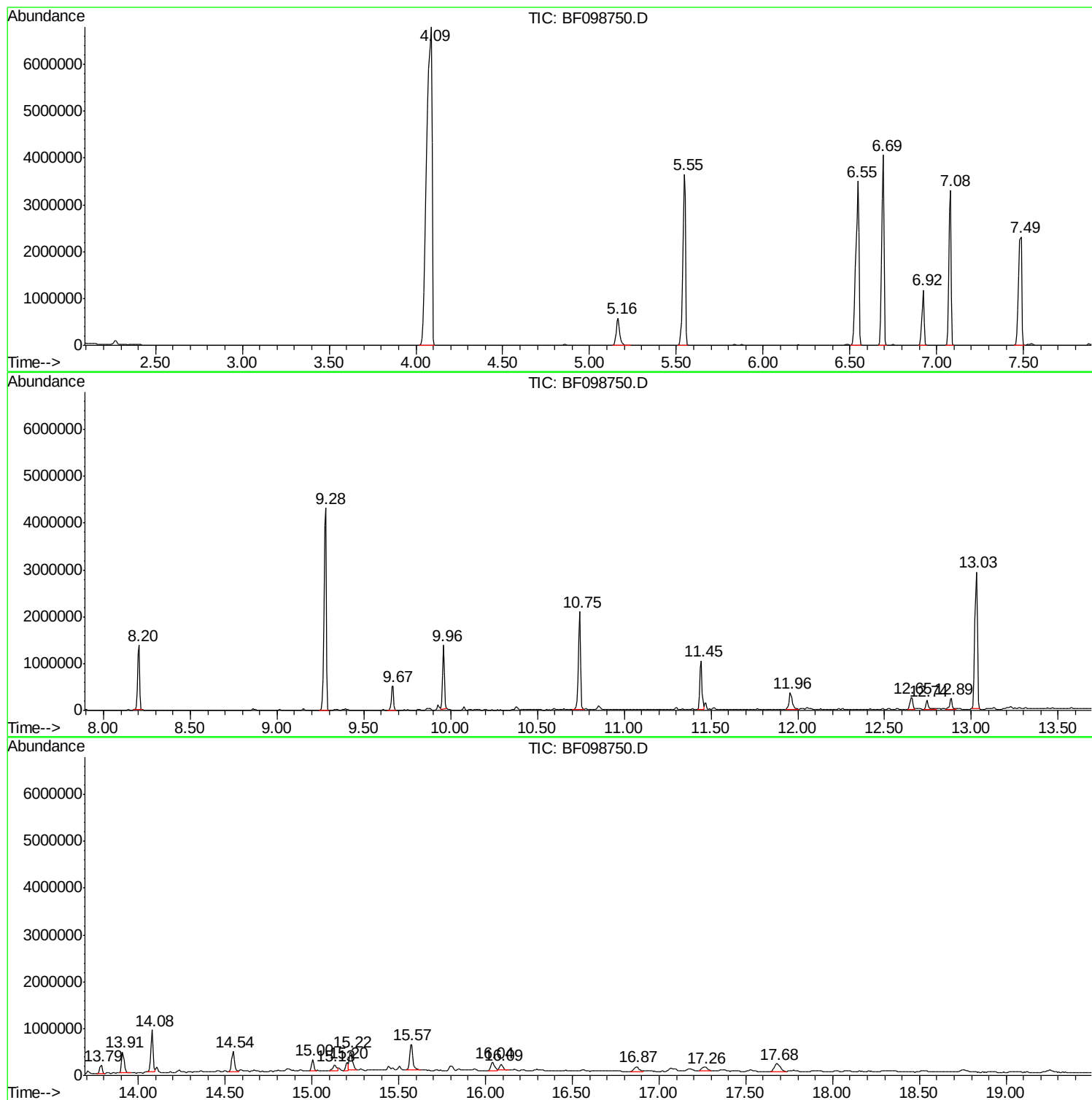
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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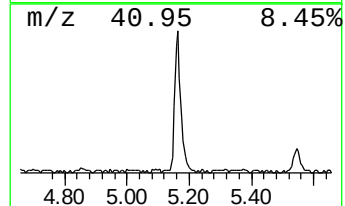
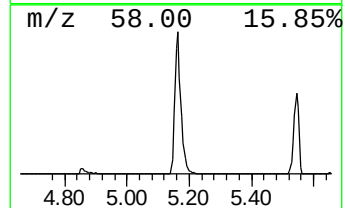
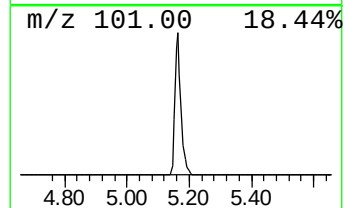
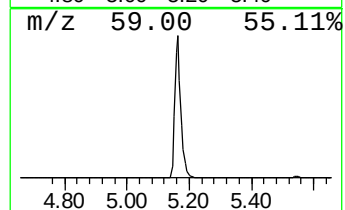
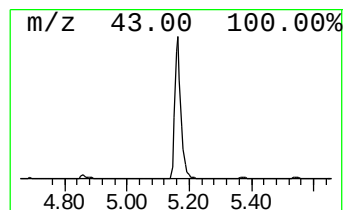
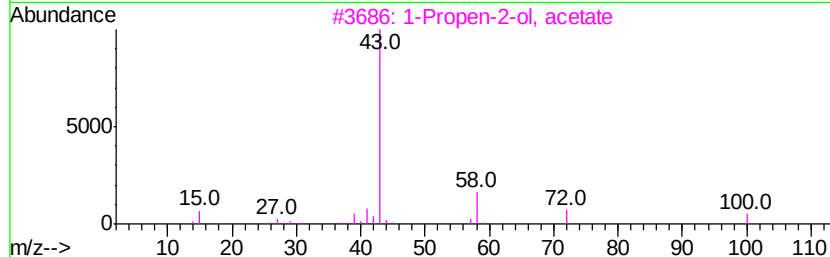
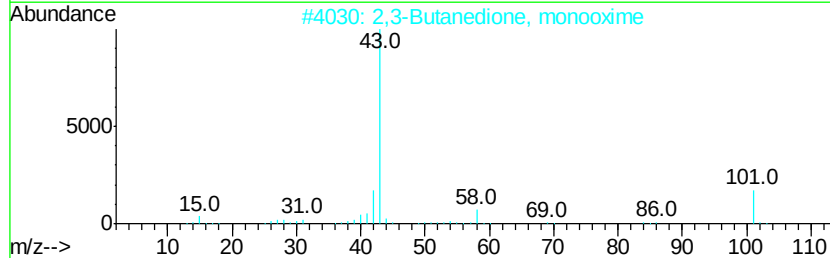
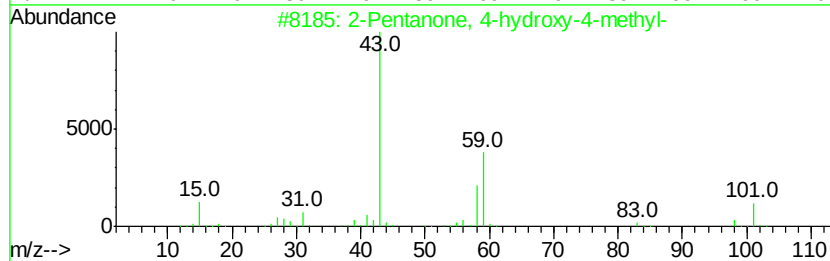
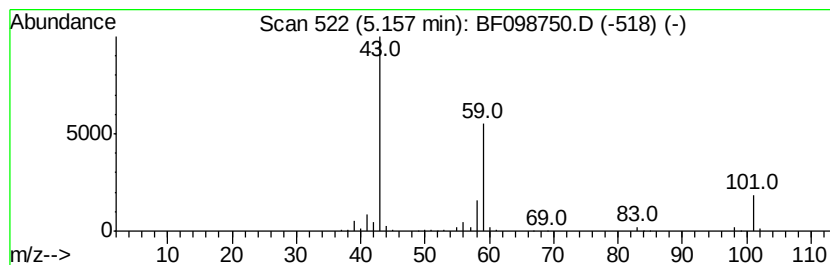
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	16.88 ng	796034	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		N,N'-Bis(2-methyl-2-nitrosopenta...	258	C12H22N2O4	094514-30-4	9
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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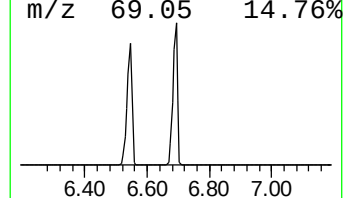
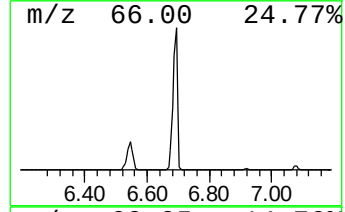
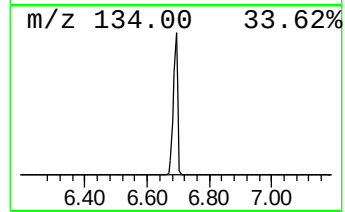
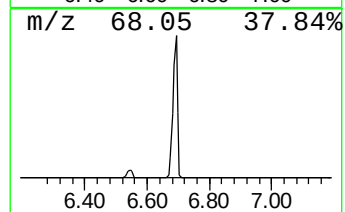
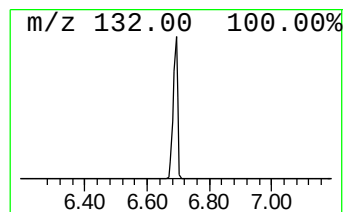
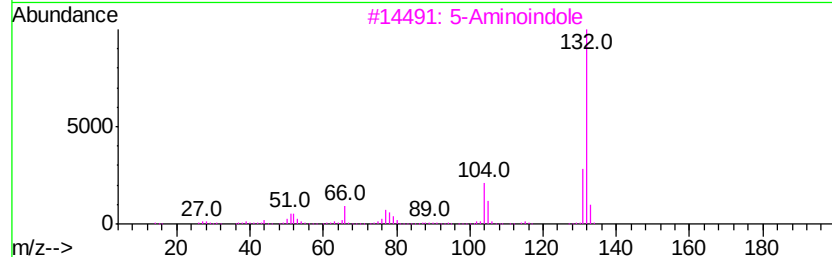
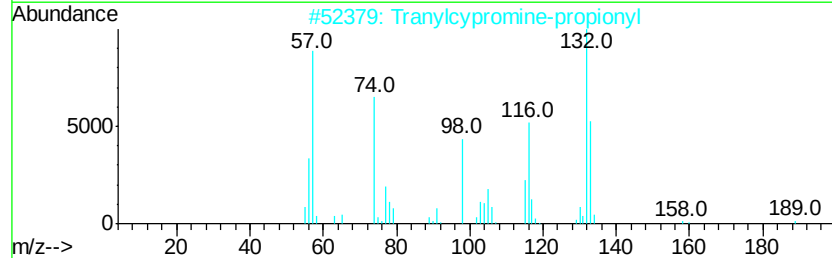
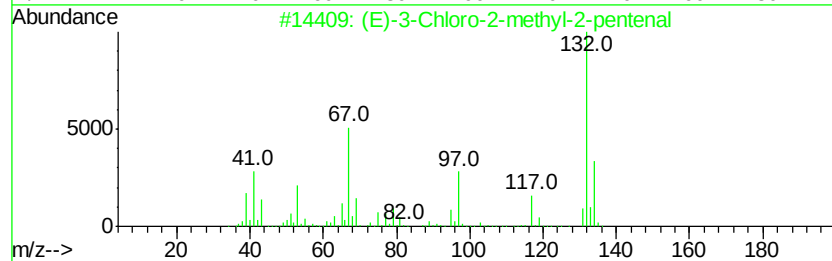
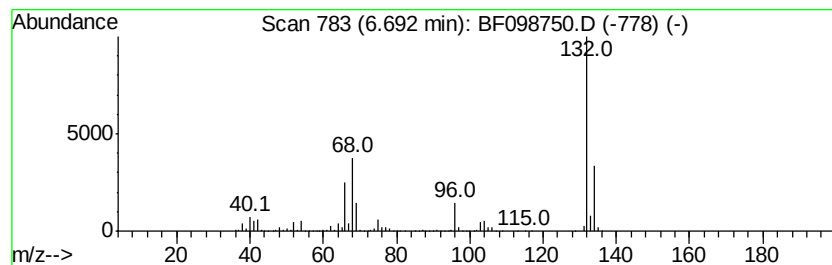
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.69 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	87.50 ng	4126790	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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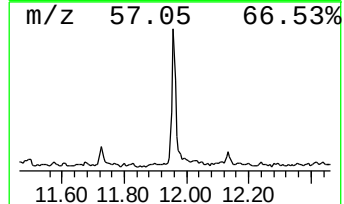
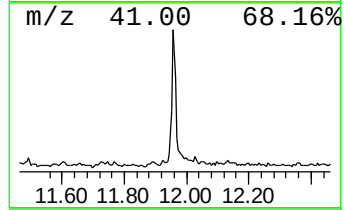
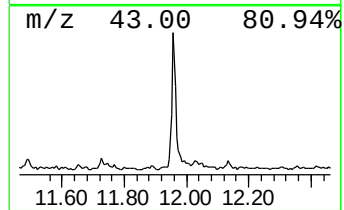
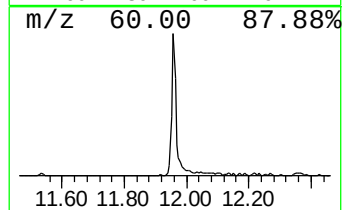
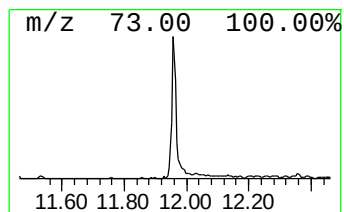
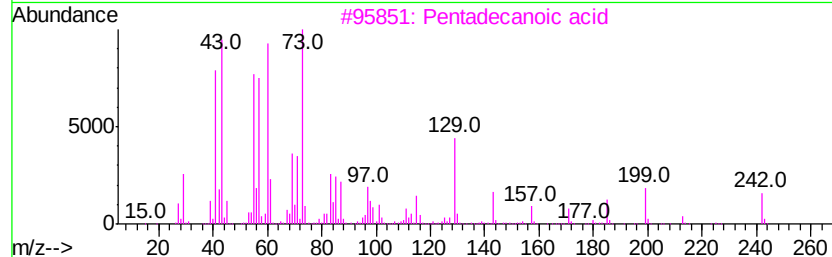
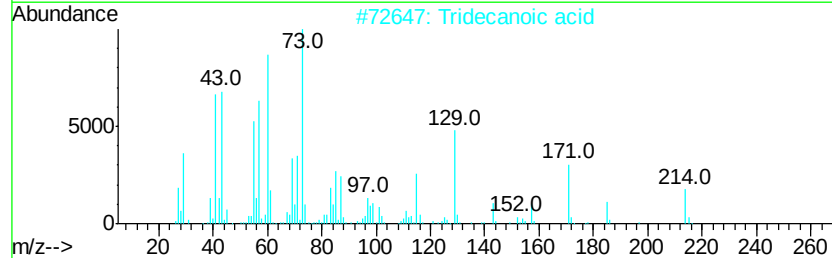
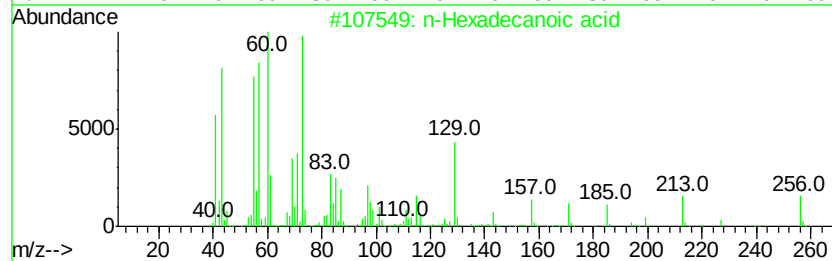
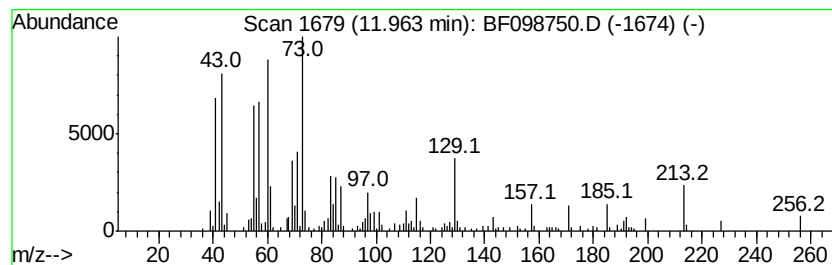
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	8.00 ng	372135	Phenanthrene-d10	11.45

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



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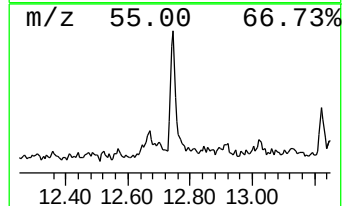
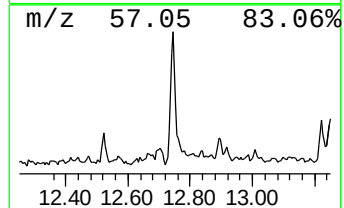
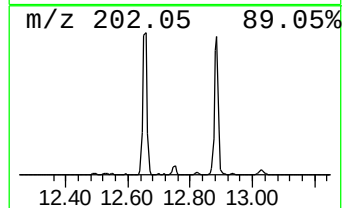
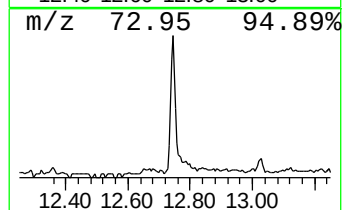
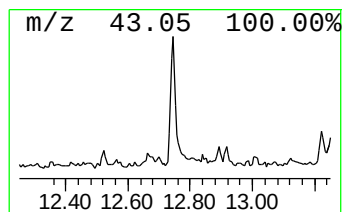
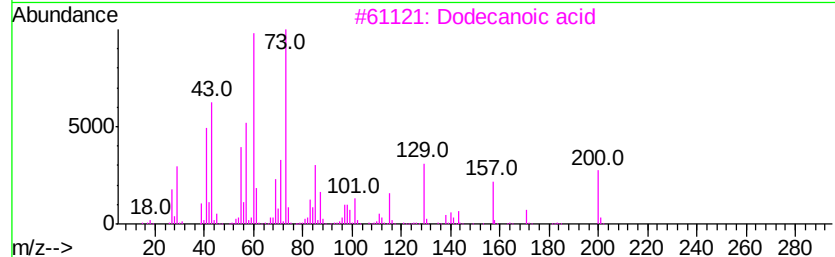
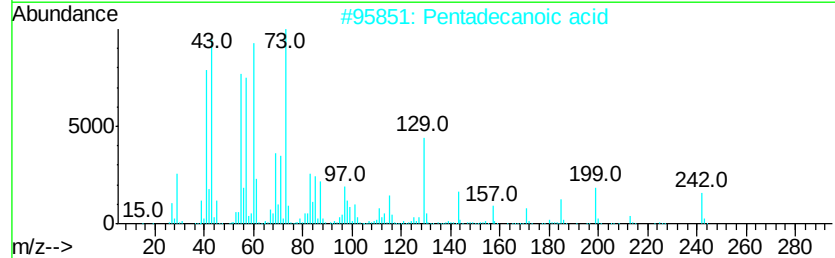
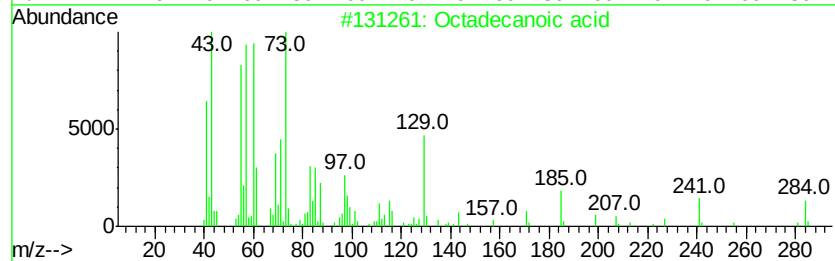
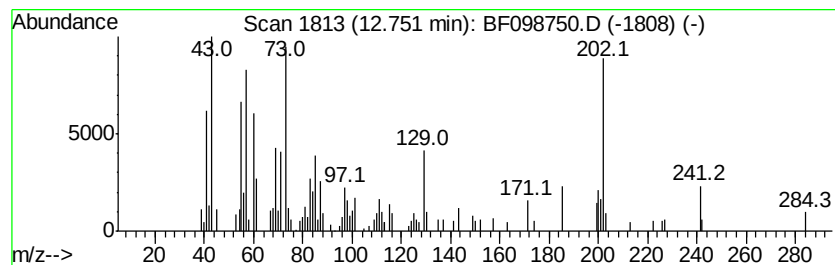
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	4.15 ng	193123	Phenanthrene-d10	11.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Dodecanoic acid	200	C12H24O2	000143-07-7	87
4	Undecanoic acid	186	C11H22O2	000112-37-8	83
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53



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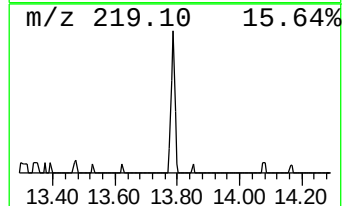
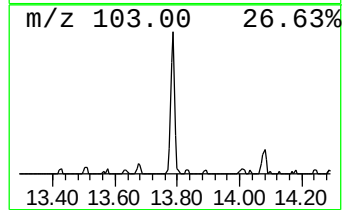
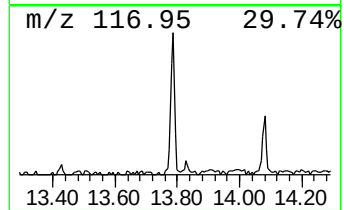
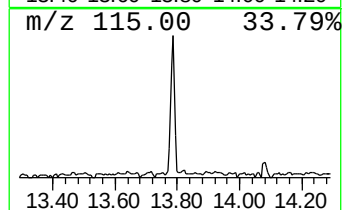
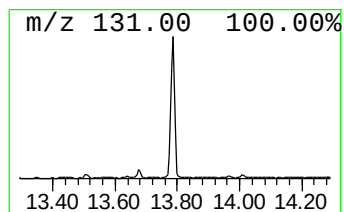
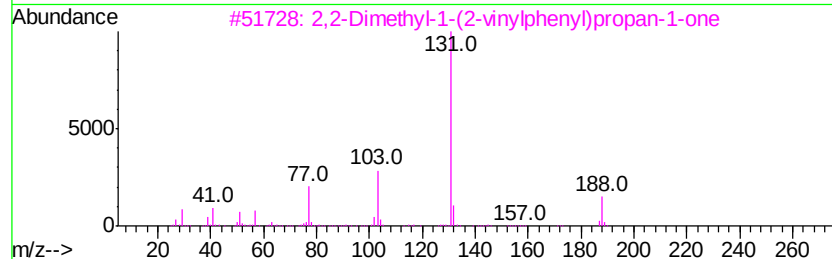
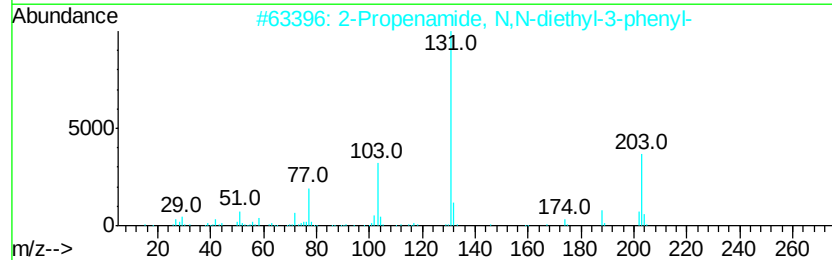
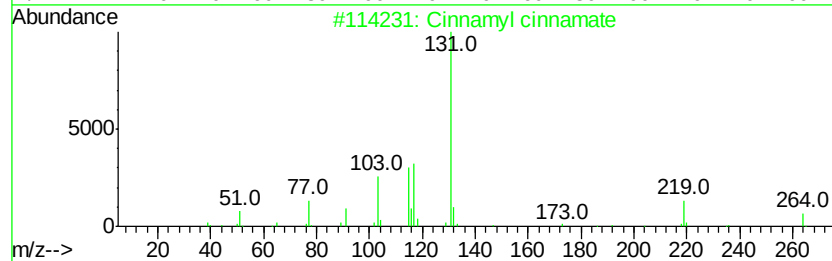
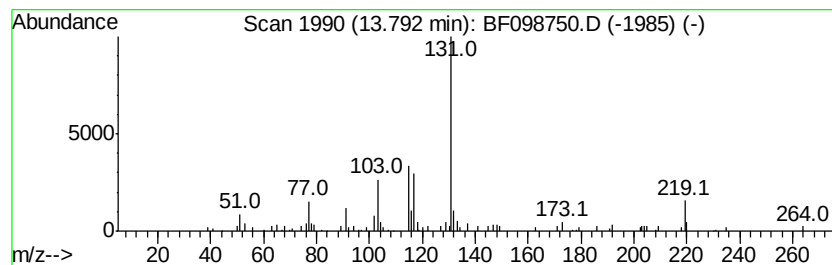
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Cinnamyl cinnamate Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	3.51 ng	152192	Chrysene-d12	14.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	87
2		2-Propenamide, N,N-diethyl-3-phe...	203	C13H17NO	003680-04-4	52
3		2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	52
4		2-Iodo-benzoic acid N'-(3-phenyl...	392	C16H13IN2O2	315672-62-9	50
5		1H-Indene, 2,3-dihydro-4-propyl-	160	C12H16	092013-16-6	50



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LT-TS-012

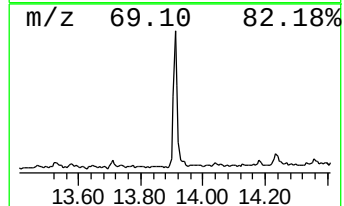
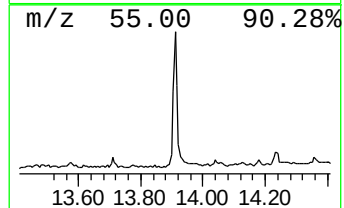
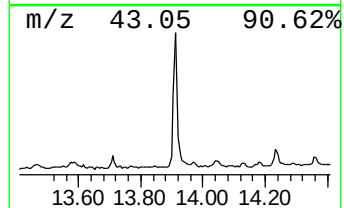
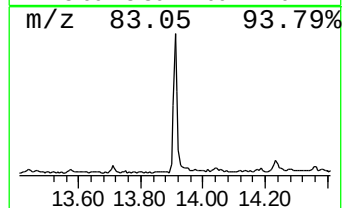
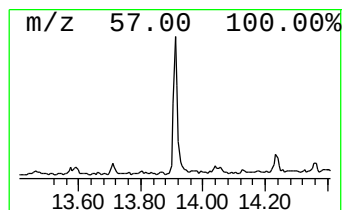
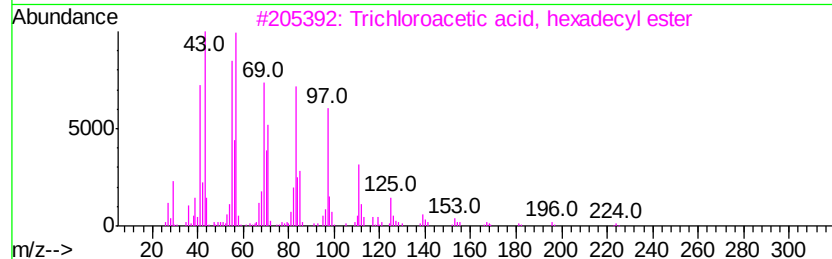
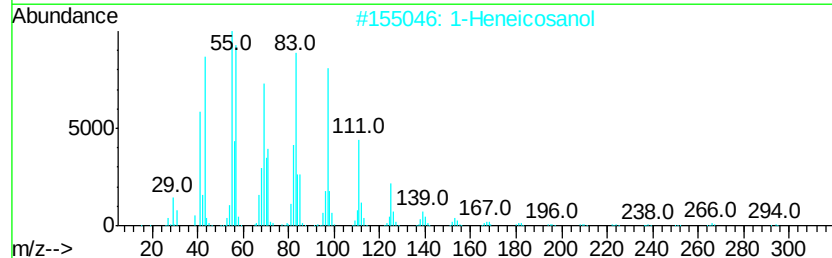
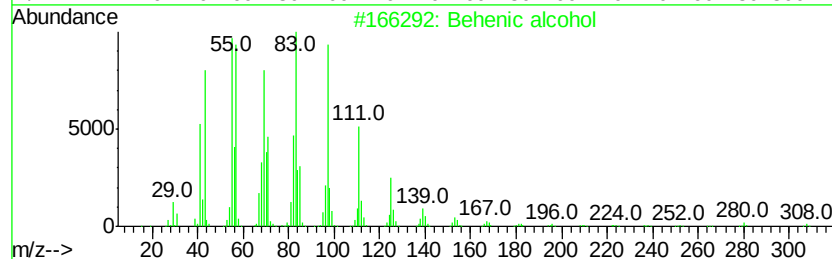
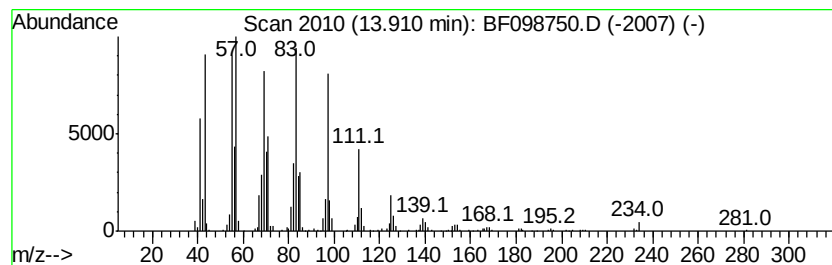
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Behenic alcohol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	9.61 ng	416796	Chrysene-d12	14.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Behenic alcohol	326	C22H46O	000661-19-8	95
2		1-Heneicosanol	312	C21H44O	015594-90-8	95
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4		Cyclopentadecane	210	C15H30	000295-48-7	93
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
Data File : BF098750.D
Acq On : 23 Sep 2017 20:29
Operator : SJ/JU
Sample : I5340-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
LT-TS-012

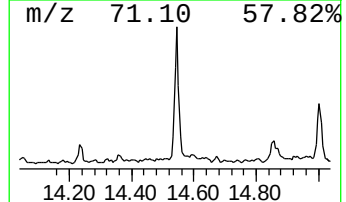
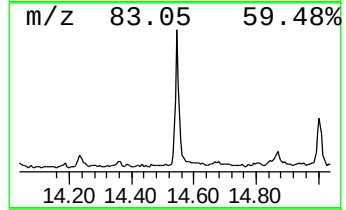
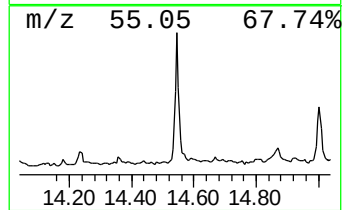
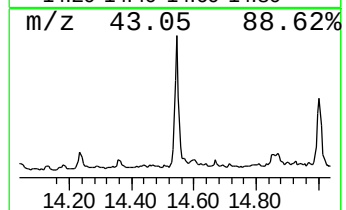
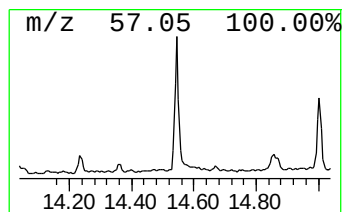
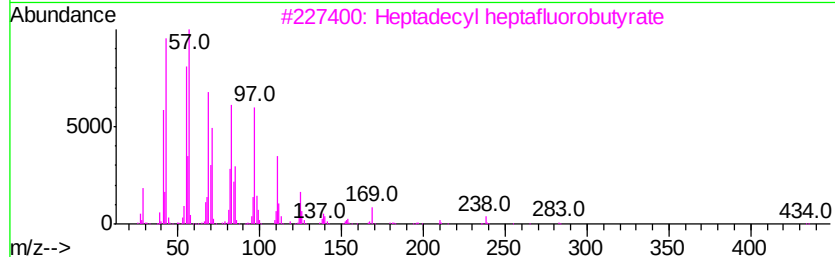
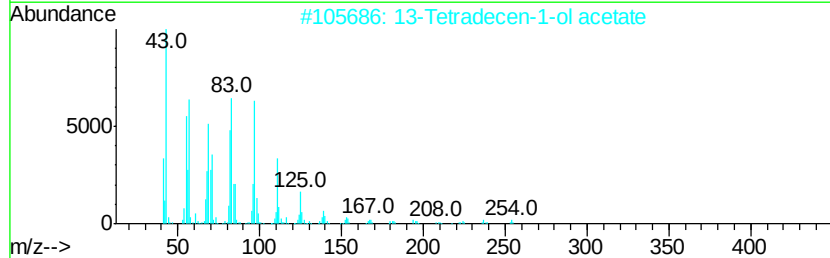
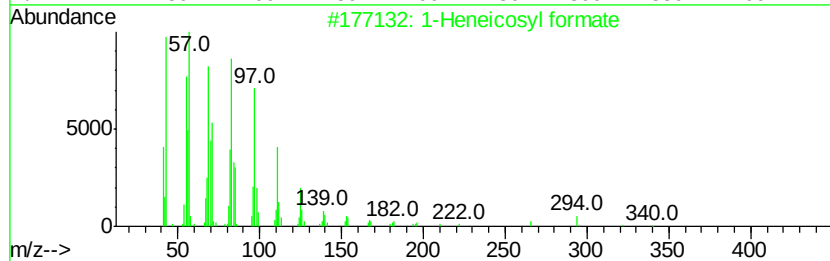
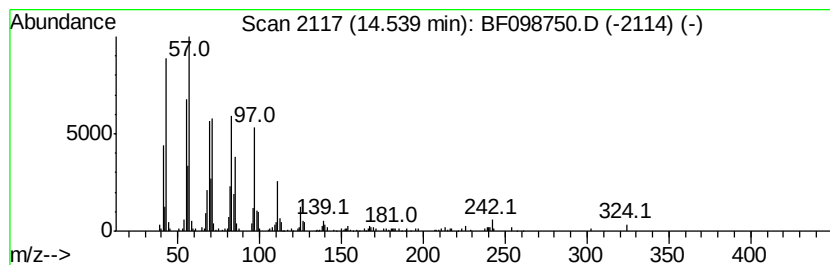
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Heneicosyl formate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	9.75 ng	422892	Chrysene-d12	14.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl	formate	340	C22H44O2	077899-03-7	95
2	13-Tetradecen-1-ol	acetate	254	C16H30O2	056221-91-1	94
3	Heptadecyl	heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4	Cyclopentadecane		210	C15H30	000295-48-7	93
5	Nonadecyl	pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
Data File : BF098750.D
Acq On : 23 Sep 2017 20:29
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ClientSampleId :
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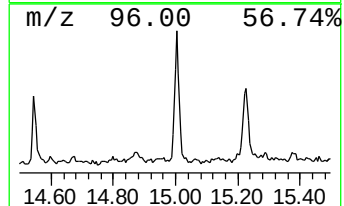
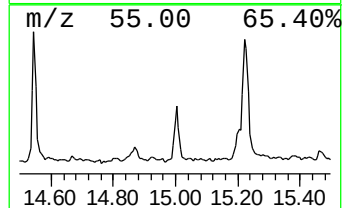
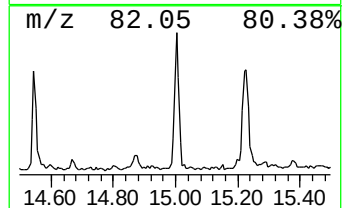
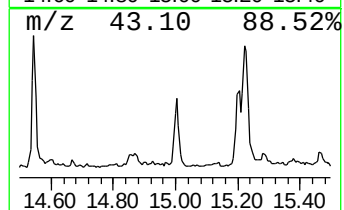
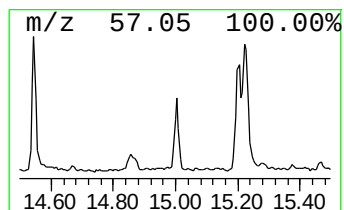
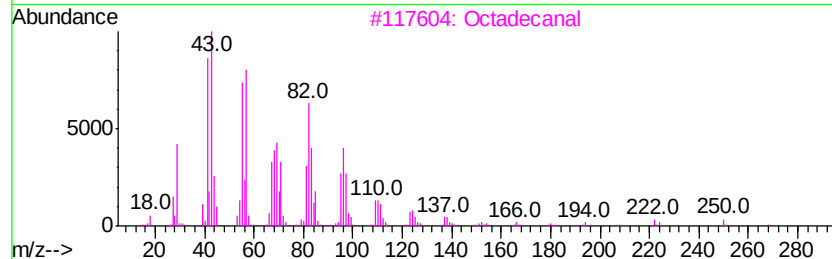
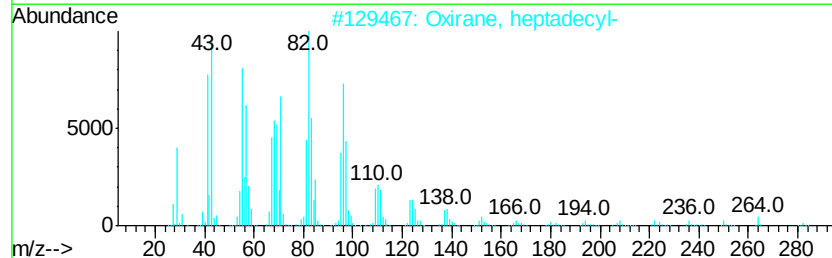
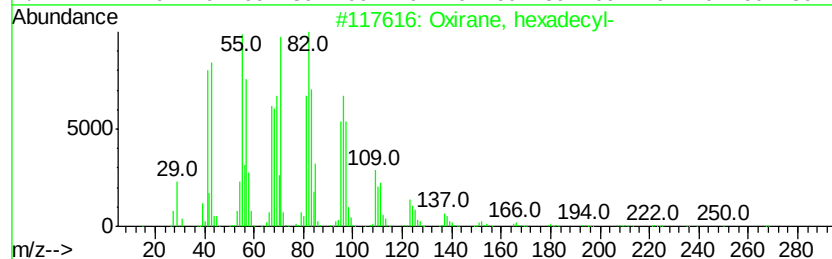
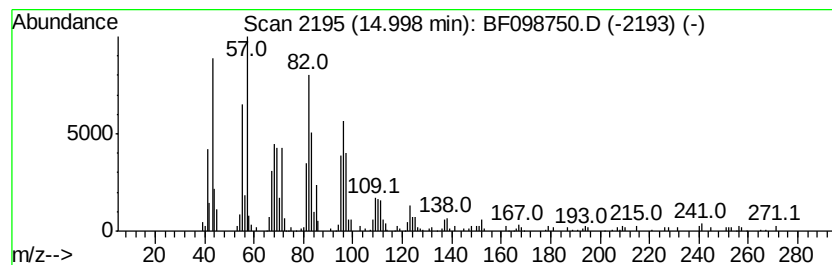
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Oxirane, hexadecyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	5.93 ng	233279	Perylene-d12	15.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	90
3			Octadecanal	268	C18H36O	000638-66-4	83
4			1,19-Eicosadiene	278	C20H38	014811-95-1	76
5			Heptadecanal	254	C17H34O	1000376-70-0	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
Data File : BF098750.D
Acq On : 23 Sep 2017 20:29
Operator : SJ/JU
Sample : I5340-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

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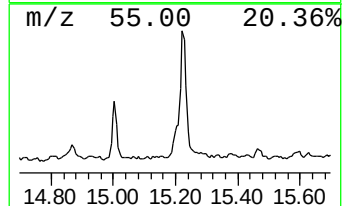
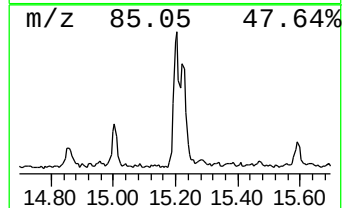
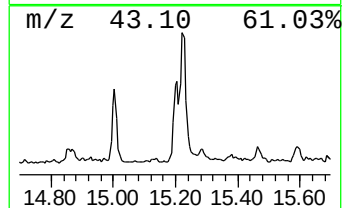
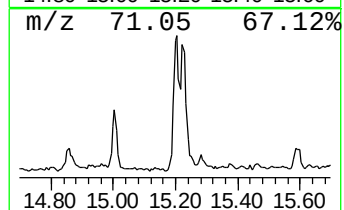
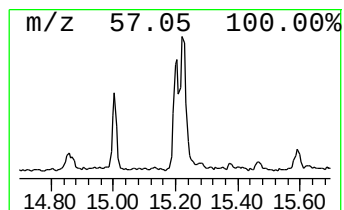
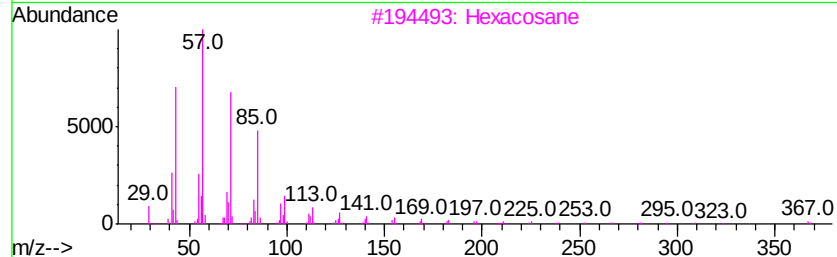
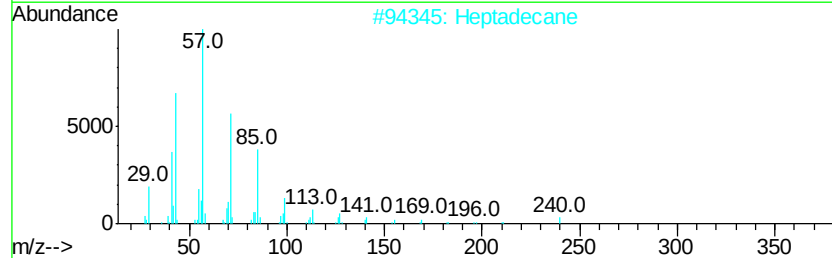
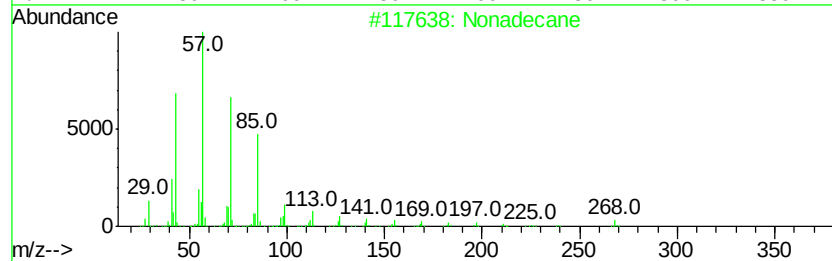
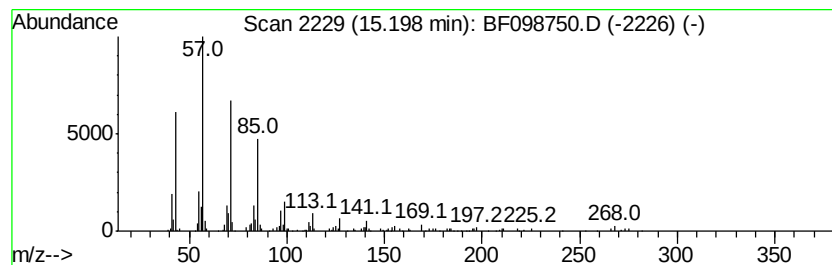
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.20	5.02 ng	197569	Perylene-d12	15.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	96
2			Heptadecane	240	C17H36	000629-78-7	91
3			Hexacosane	366	C26H54	000630-01-3	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092317\
Data File : BF098750.D
Acq On : 23 Sep 2017 20:29
Operator : SJ/JU
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Misc :
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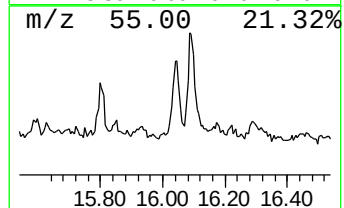
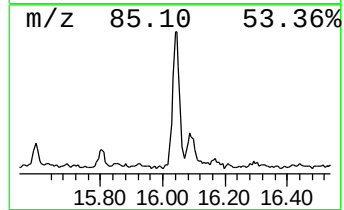
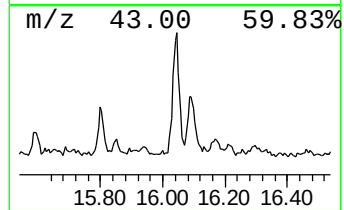
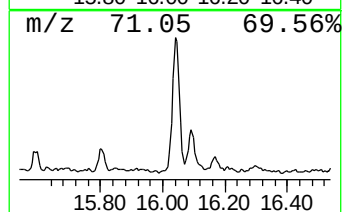
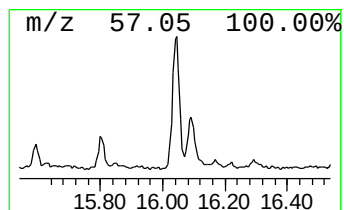
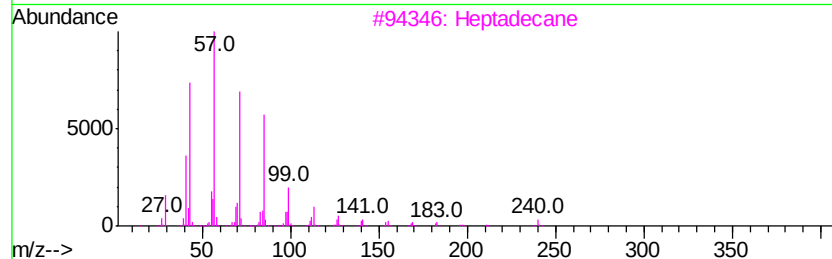
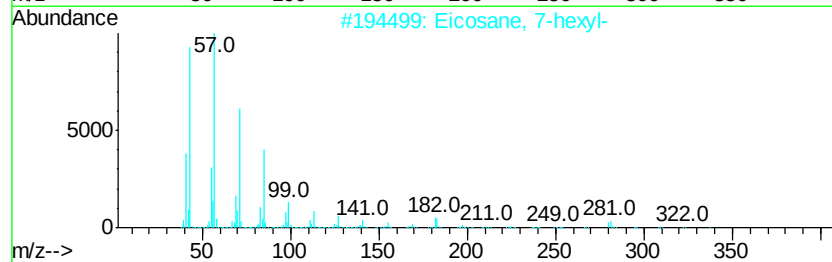
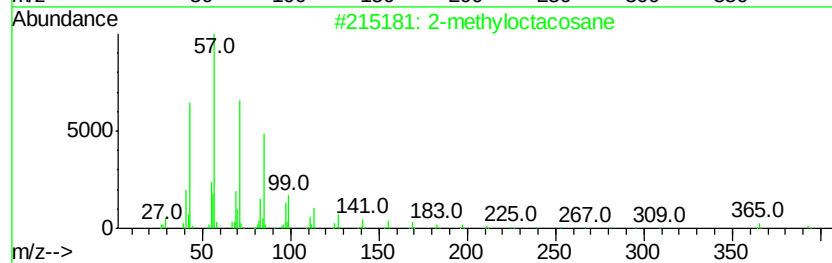
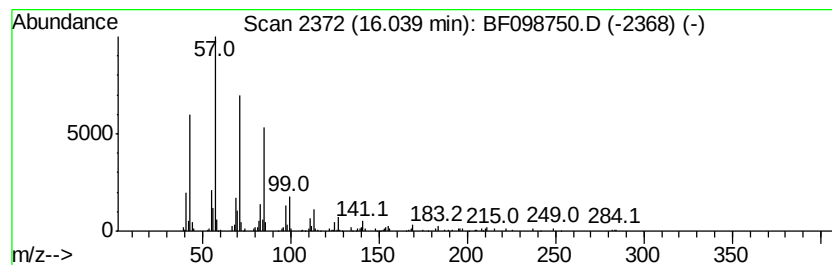
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-methyloctacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	6.26 ng	246099	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	87
2		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86
3		Heptadecane	240	C17H36	000629-78-7	86
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
5		Triacontane	422	C30H62	000638-68-6	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092317\
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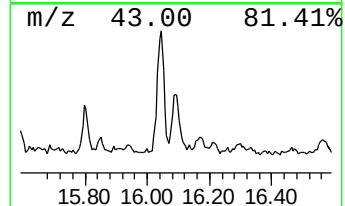
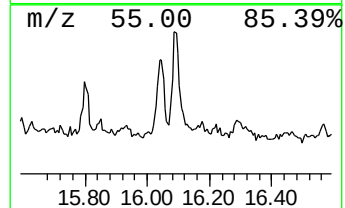
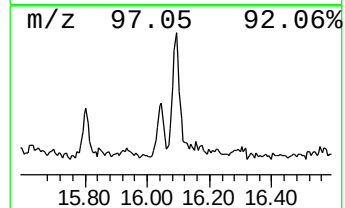
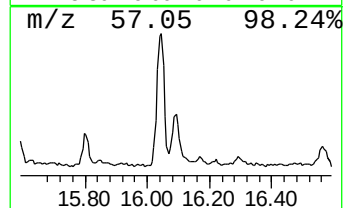
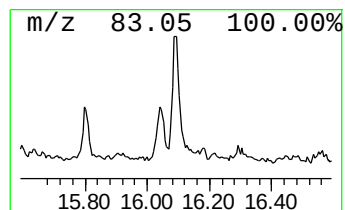
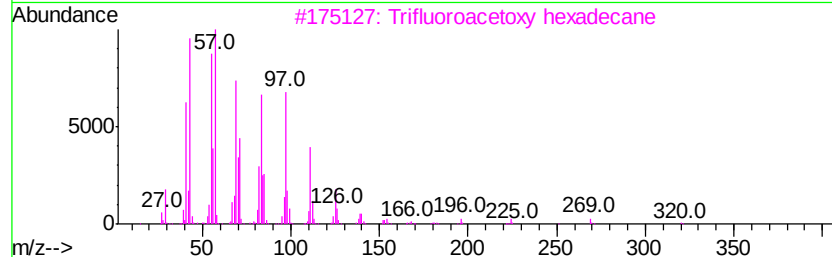
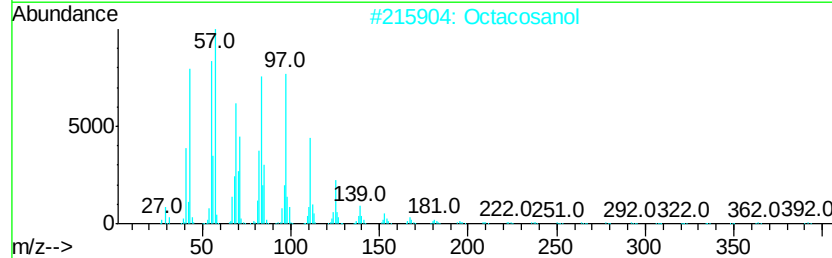
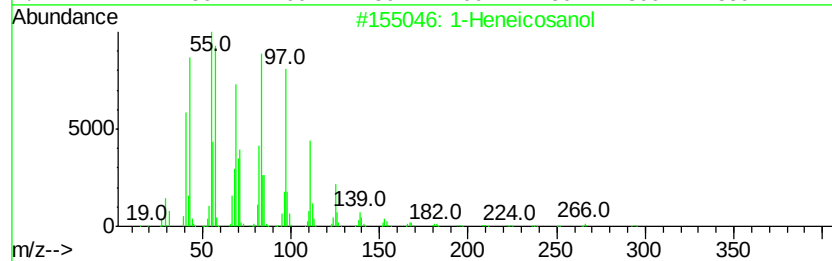
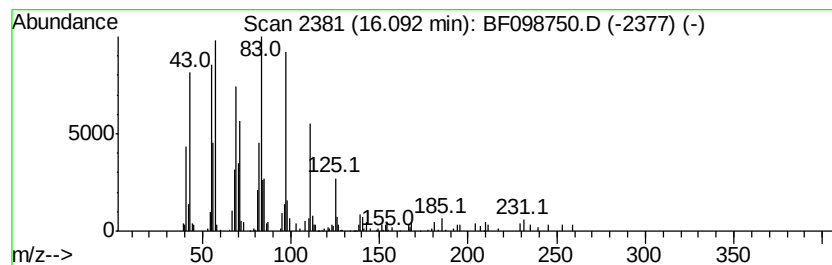
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Heneicosanol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	4.81 ng	189371	Perylene-d12	15.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Heneicosanol	312 C21H44O	015594-90-8 94
2		Octacosanol	410 C28H58O	000557-61-9 81
3		Trifluoroacetoxy hexadecane	338 C18H33F3O2	006222-03-3 62
4		13-Tetradecen-1-ol acetate	254 C16H30O2	056221-91-1 58
5		Heptadecyl heptafluorobutyrate	452 C21H35F7O2	959085-66-6 58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
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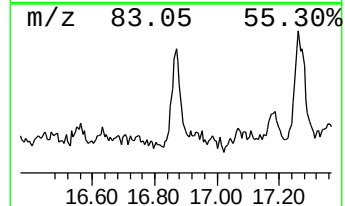
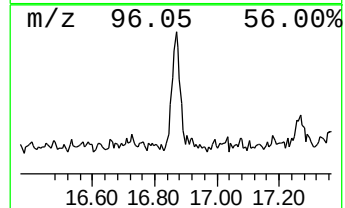
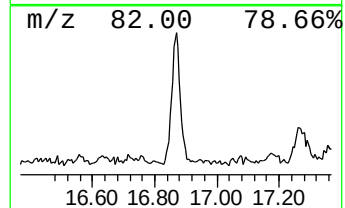
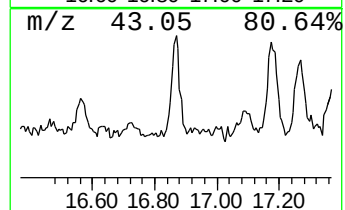
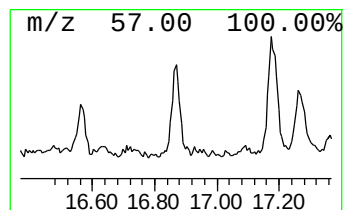
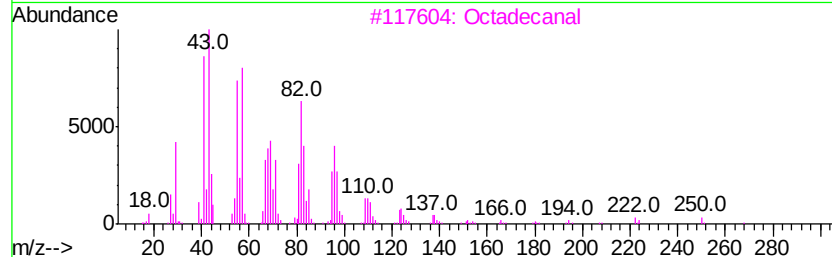
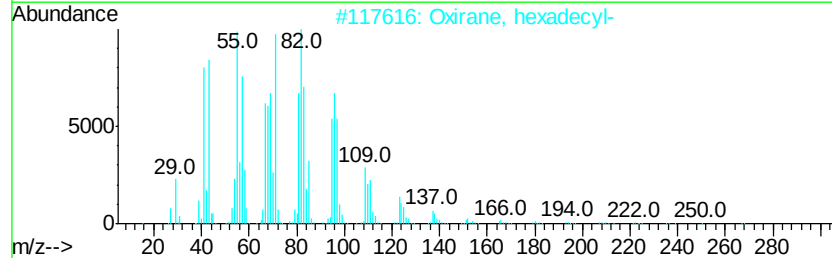
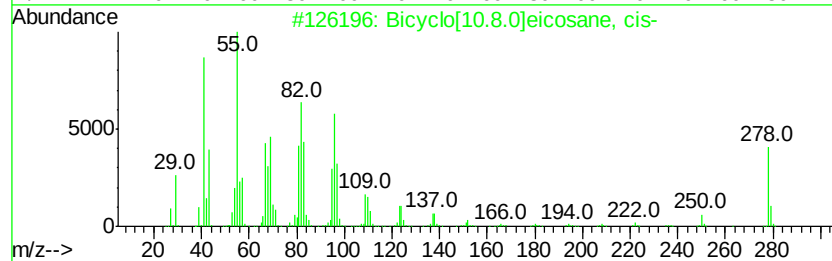
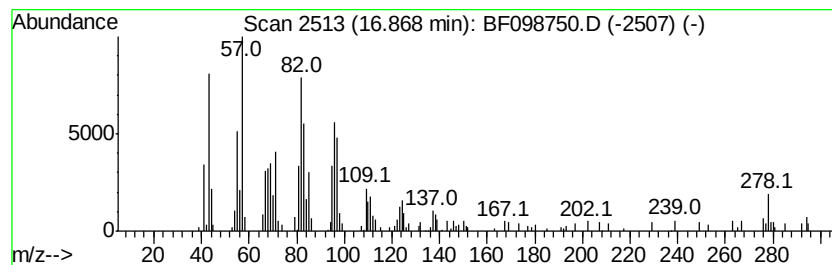
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Bicyclo[10.8.0]eicosane, cis- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.87	5.12 ng	201412	Perylene-d12		15.57	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	89
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	86
3		Octadecanal	268	C18H36O	000638-66-4	72
4		11-Dodecen-1-ol trifluoroacetate	280	C14H23F3O2	128792-46-1	70
5		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092317\
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ClientSampleId :
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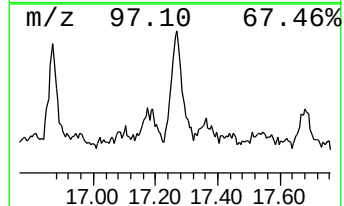
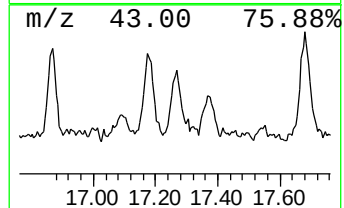
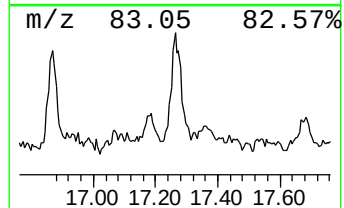
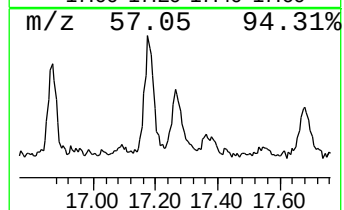
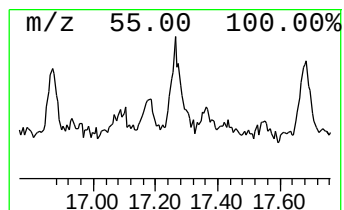
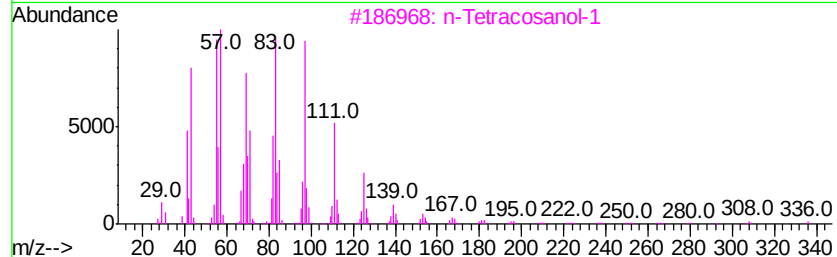
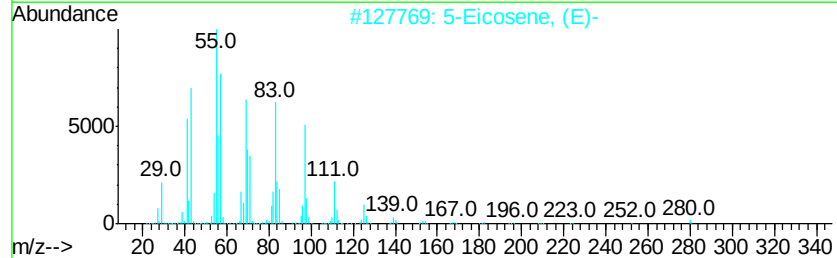
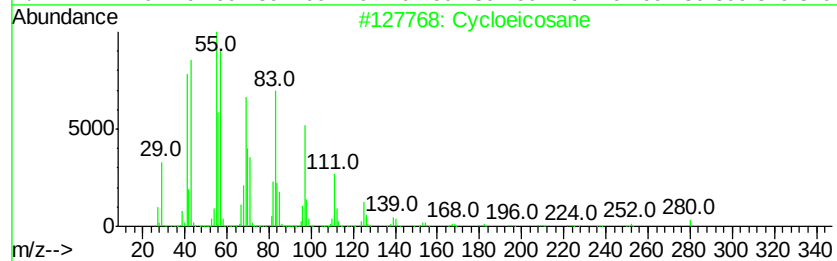
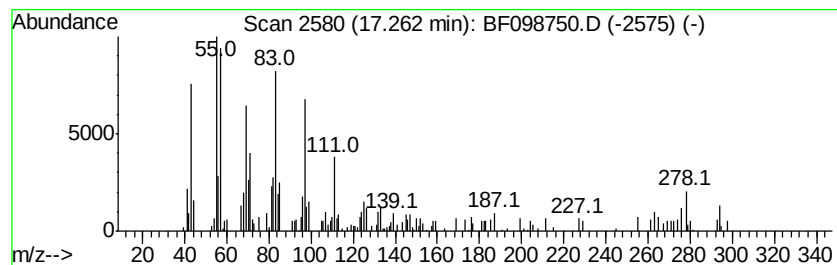
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Cycloeicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.26	3.95 ng	155301	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cycloeicosane	280	C20H40	000296-56-0	89
2		5-Eicosene, (E)-	280	C20H40	074685-30-6	78
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	76
4		1-Docosene	308	C22H44	001599-67-3	76
5		10-Heneicosene (c,t)	294	C21H42	095008-11-0	68



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
Data File : BF098750.D
Acq On : 23 Sep 2017 20:29
Operator : SJ/JU
Sample : I5340-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LT-TS-012

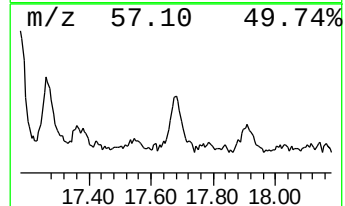
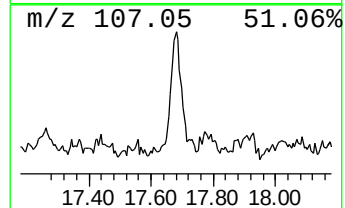
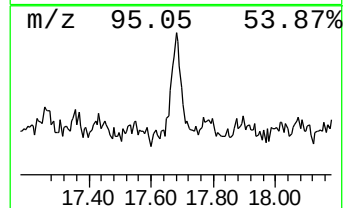
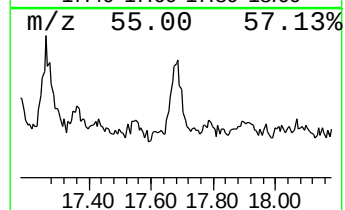
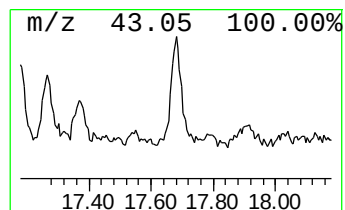
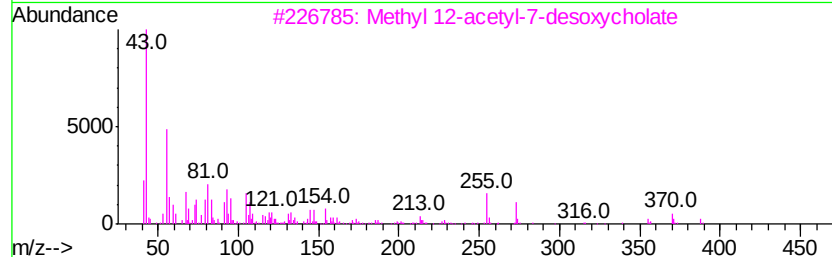
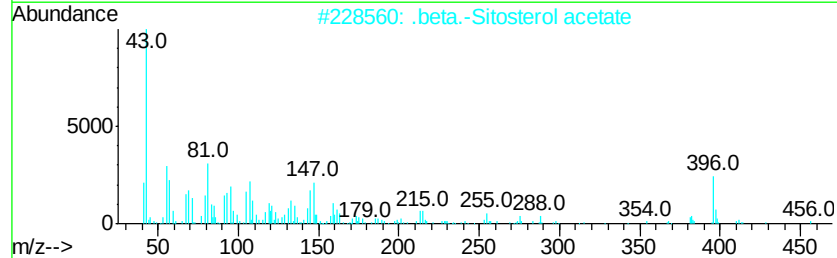
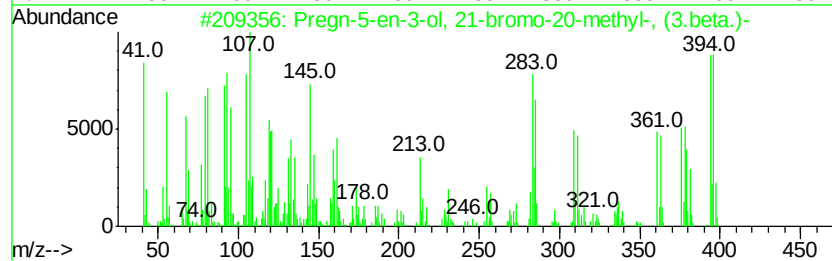
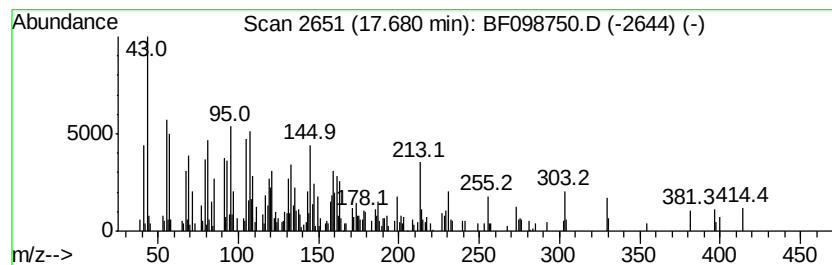
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown17.68 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	10.04 ng	394866	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	42
2		.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	37
3		Methyl 12-acetyl-7-desoxycholate	448	C27H44O5	055547-48-3	25
4		Clionasterol acetate	456	C31H52O2	004651-54-1	25
5		Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092317\
 Data File : BF098750.D
 Acq On : 23 Sep 2017 20:29
 Operator : SJ/JU
 Sample : I5340-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LT-TS-012

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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...	5.16	16.9	ng	796034	1	6.92	943227	20.0
unknown6.69	6.69	87.5	ng	4126790	1	6.92	943227	20.0
n-Hexadecanoic acid	11.96	8.0	ng	372135	4	11.45	930472	20.0
Octadecanoic acid	12.75	4.2	ng	193123	4	11.45	930472	20.0
Cinnamyl cinnamate	13.79	3.5	ng	152192	5	14.08	867794	20.0
Behenic alcohol	13.91	9.6	ng	416796	5	14.08	867794	20.0
1-Heneicosyl formate	14.54	9.8	ng	422892	5	14.08	867794	20.0
Oxirane, hexadecyl-	15.00	5.9	ng	233279	6	15.57	786658	20.0
Nonadecane	15.20	5.0	ng	197569	6	15.57	786658	20.0
2-methyloctacosane	16.04	6.3	ng	246099	6	15.57	786658	20.0
1-Heneicosanol	16.09	4.8	ng	189371	6	15.57	786658	20.0
Bicyclo[10.8.0]ei...	16.87	5.1	ng	201412	6	15.57	786658	20.0
Cycloeicosane	17.26	4.0	ng	155301	6	15.57	786658	20.0
unknown17.68	17.68	10.0	ng	394866	6	15.57	786658	20.0