

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
 Data File : BF093939.D
 Acq On : 27 Mar 2017 12:54
 Operator : SJ/MA
 Sample : I2367-08
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-BB10-NSW

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-BF032217.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	3.599	286	291	298	rBV	1120109	1288563	1.31%	0.654%
2	4.481	433	441	445	rBV	3265823	4359910	4.44%	2.213%
3	5.551	611	623	626	rVV	3288243	4445391	4.52%	2.256%
4	5.693	642	647	650	rVV	4043943	5104892	5.19%	2.591%
5	5.951	687	691	694	rBV	1046195	1002465	1.02%	0.509%
6	6.128	717	721	725	rVB	3171543	3223300	3.28%	1.636%
7	6.598	794	801	804	rBV	2497698	2822200	2.87%	1.433%
8	7.451	942	946	949	rVV	1398503	1460854	1.49%	0.742%
9	7.969	1019	1034	1037	rVV2	751619	1835288	1.87%	0.932%
10	8.781	1158	1172	1175	rVV	4060664	5192315	5.28%	2.636%
11	9.433	1277	1283	1285	rVB	848701	1324510	1.35%	0.672%
12	9.592	1307	1310	1315	rVV2	1272184	1393881	1.42%	0.708%
13	9.928	1359	1367	1371	rBV4	658280	1214628	1.24%	0.617%
14	10.569	1472	1476	1480	rVB	2001379	2128561	2.17%	1.080%
15	11.175	1560	1579	1582	rBV	4170668	15391637	15.66%	7.813%
16	11.422	1615	1621	1623	rBV	1126221	1243741	1.27%	0.631%
17	11.557	1641	1644	1648	rBV2	1002801	1383741	1.41%	0.702%
18	11.651	1653	1660	1662	rBV2	813880	1211380	1.23%	0.615%
19	12.433	1739	1793	1794	rBV8	7295601	98281329	100.00%	49.887%
20	12.633	1825	1827	1831	rVB	1347045	1332481	1.36%	0.676%
21	12.792	1850	1854	1857	rBV2	1563507	1678545	1.71%	0.852%
22	13.027	1889	1894	1897	rVB	2909178	3584124	3.65%	1.819%
23	13.180	1904	1920	1928	rBV3	4575292	25387299	25.83%	12.886%
24	13.374	1950	1953	1959	rVB3	1738540	2284374	2.32%	1.160%
25	13.433	1959	1963	1969	rVB	1679887	2039533	2.08%	1.035%
26	18.962	2893	2903	2918	rVB	1444814	3928853	4.00%	1.994%
27	19.245	2942	2951	2964	rVB2	910416	2463566	2.51%	1.250%

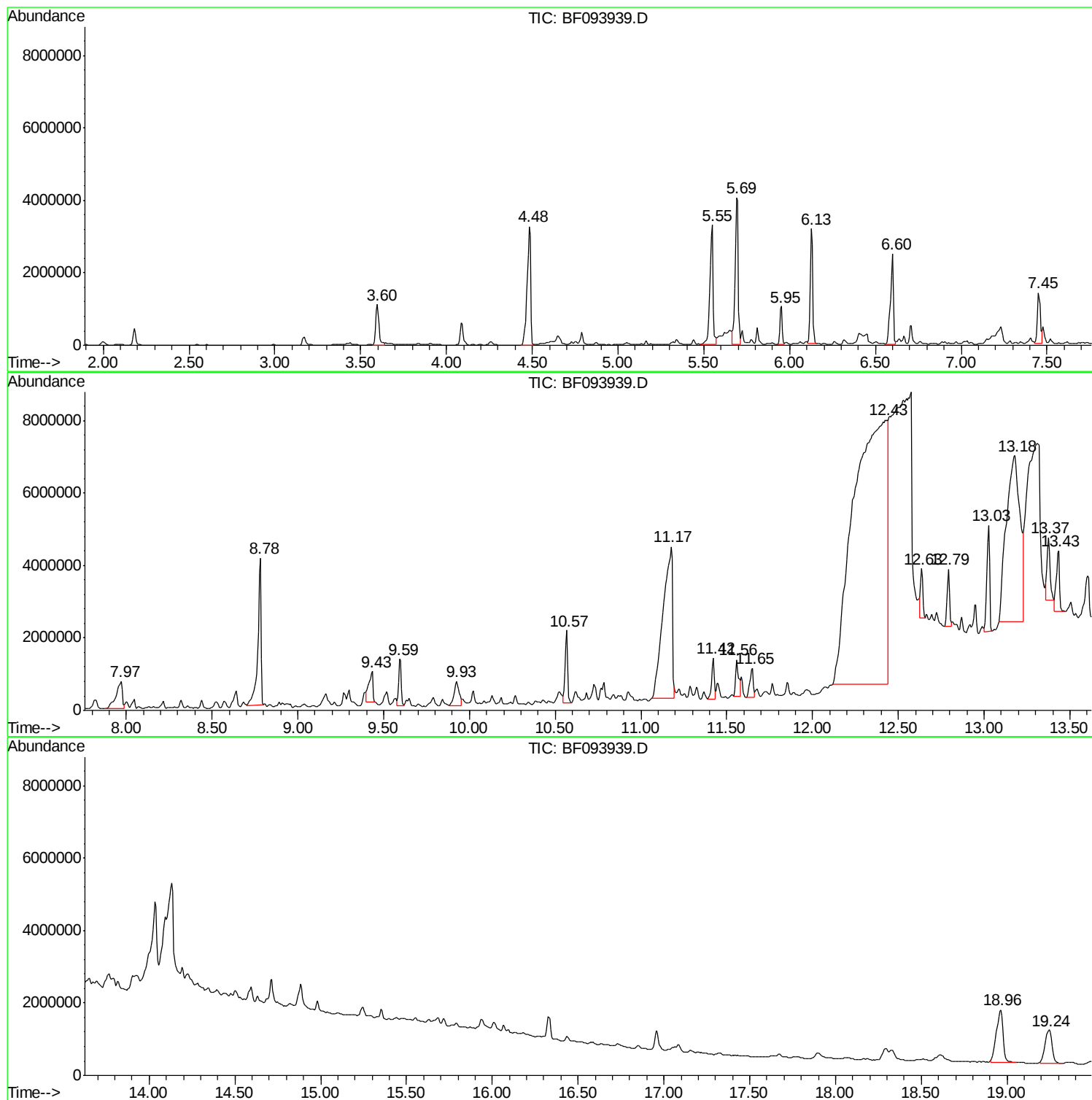
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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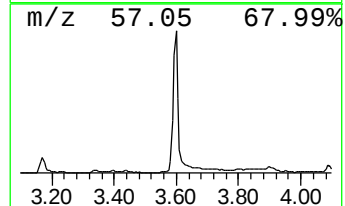
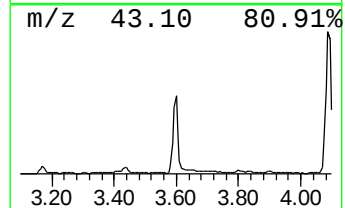
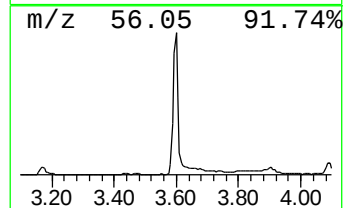
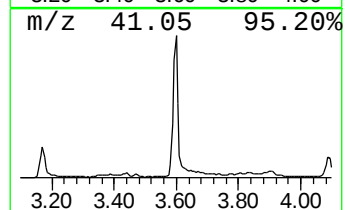
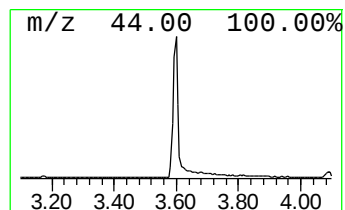
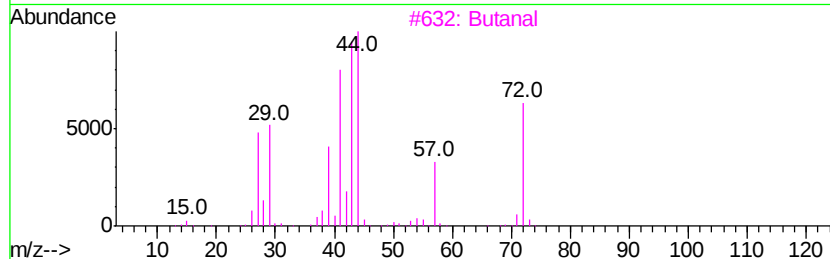
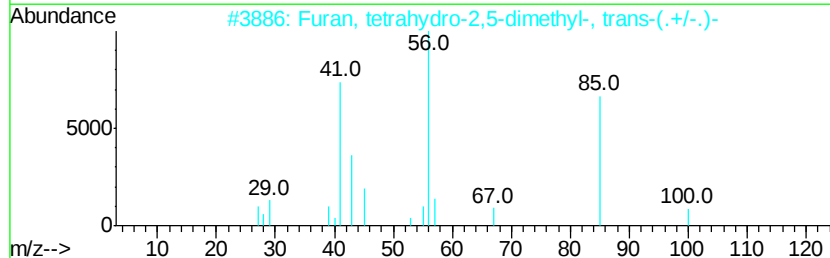
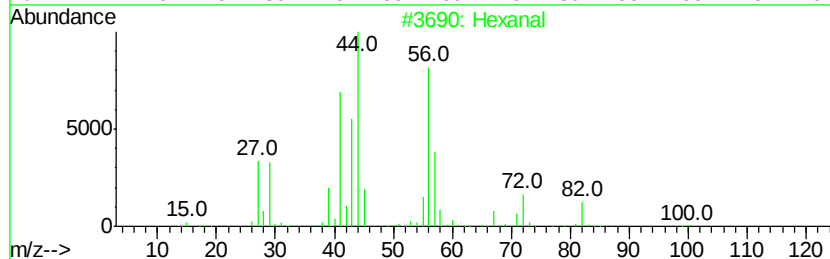
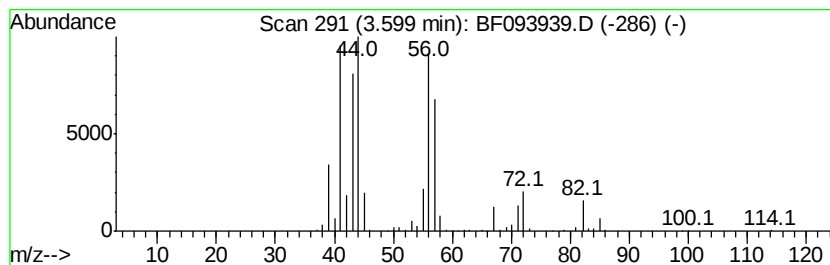
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.60	25.71 ng	1288560	1,4-Dichlorobenzene-d4	5.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanal	100	C6H12O	000066-25-1	83
2		Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	038484-59-2	38
3		Butanal	72	C4H8O	000123-72-8	35
4		Cyanoic acid, butyl ester	99	C5H9NO	001768-24-7	32
5		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	25



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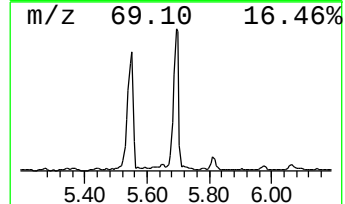
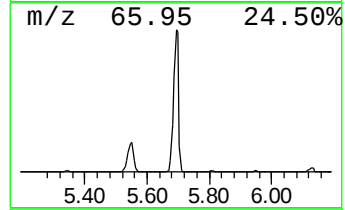
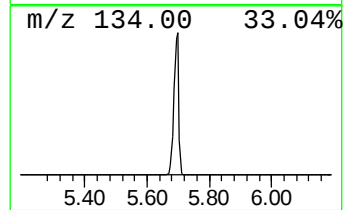
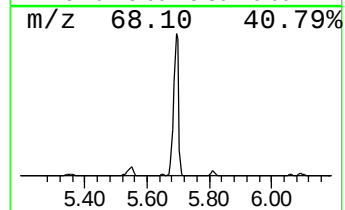
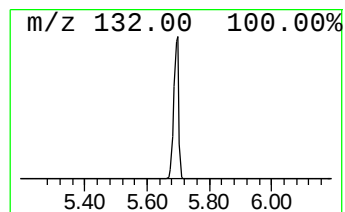
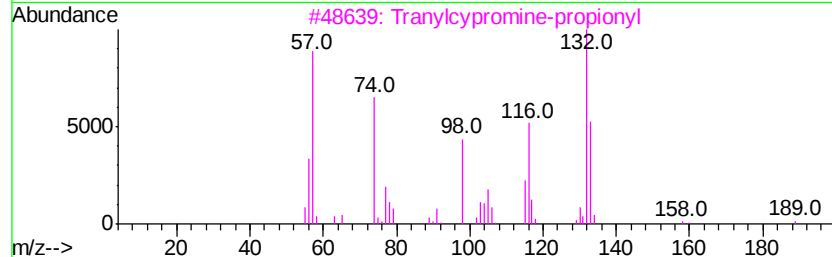
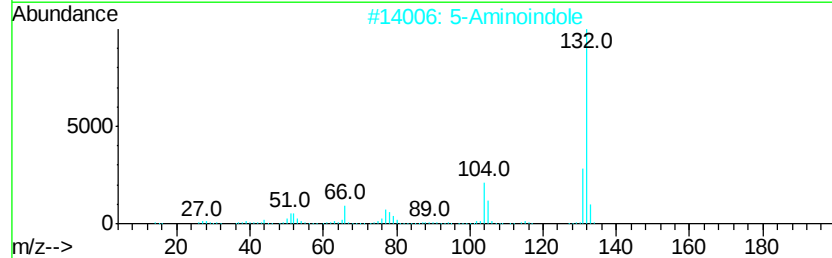
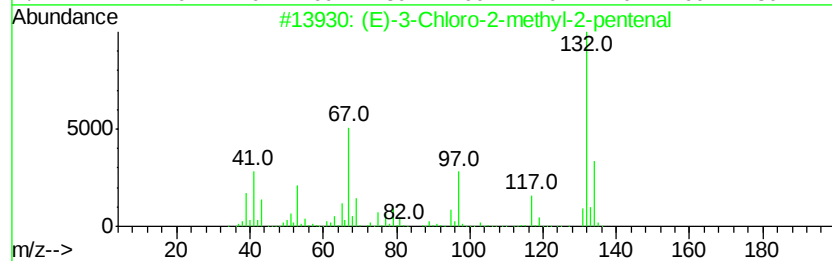
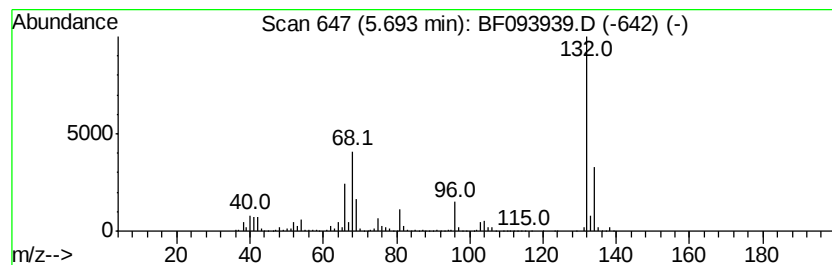
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Peak Number 2 unknown5.69 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.69	101.85 ng	5104890	1,4-Dichlorobenzene-d4	5.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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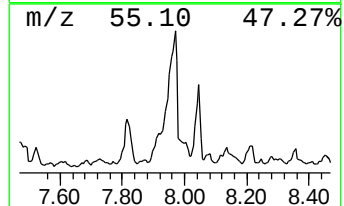
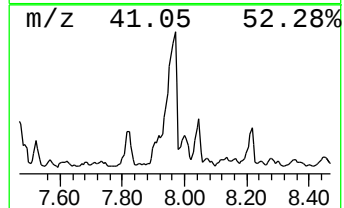
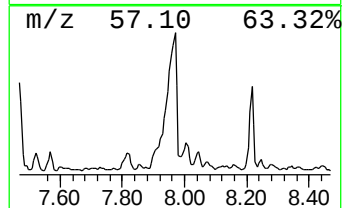
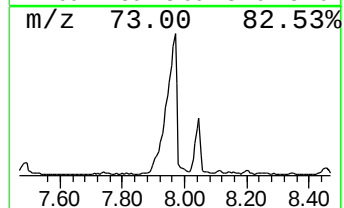
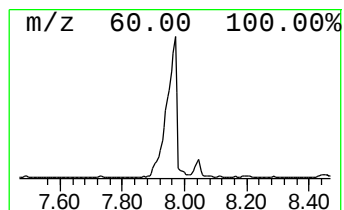
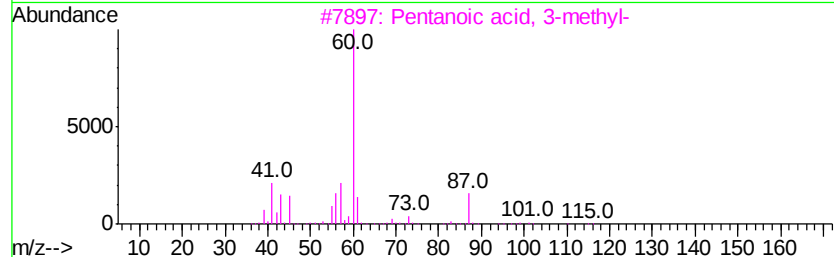
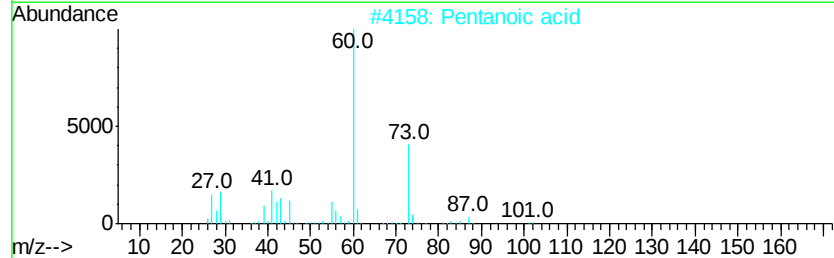
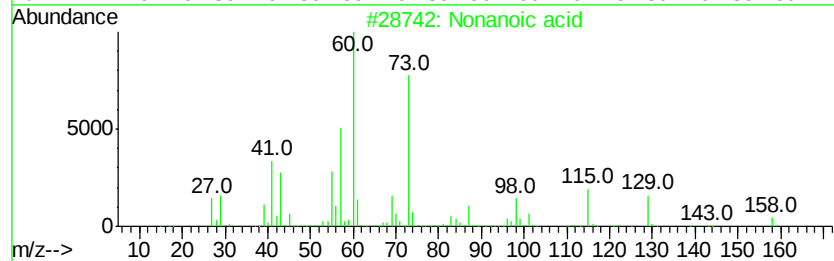
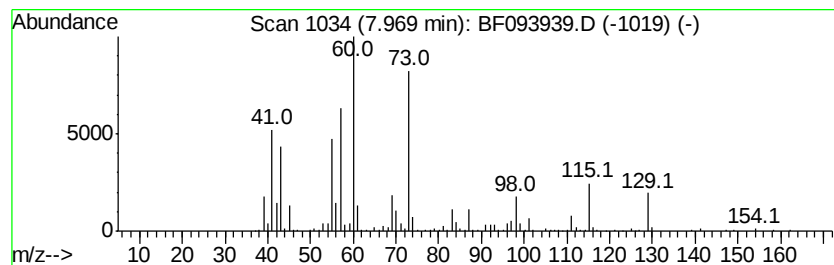
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Peak Number 4 Nonanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	25.13 ng	1835290	Naphthalene-d8	7.45

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid		158	C9H18O2	000112-05-0	91
2	Pentanoic acid		102	C5H10O2	000109-52-4	46
3	Pentanoic acid, 3-methyl-		116	C6H12O2	000105-43-1	43
4	Butanoic acid, 3-methyl-		102	C5H10O2	000503-74-2	43
5	Heptanoic acid		130	C7H14O2	000111-14-8	43



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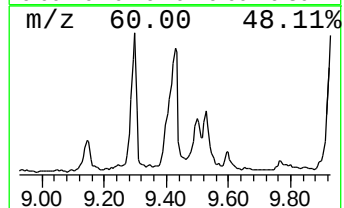
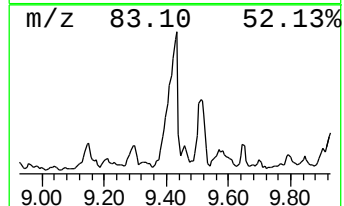
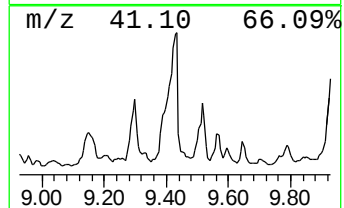
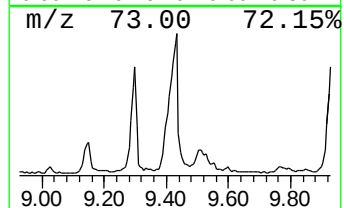
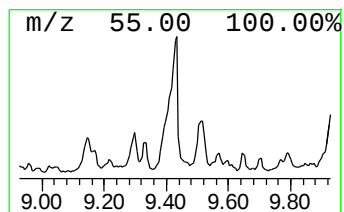
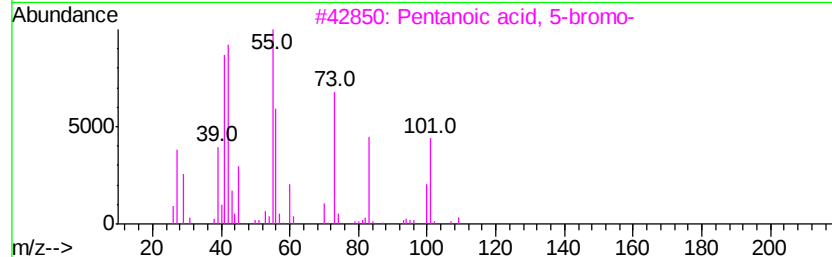
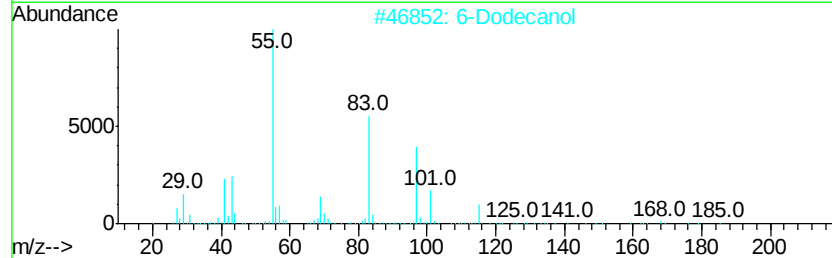
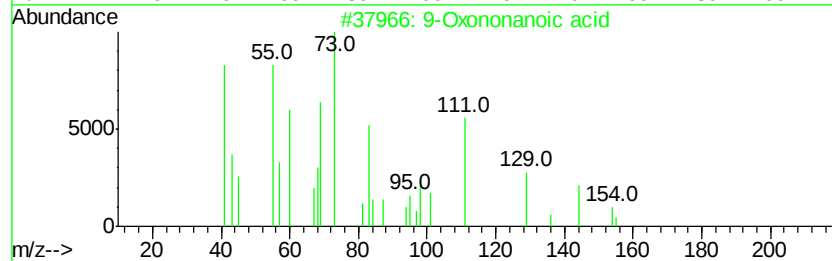
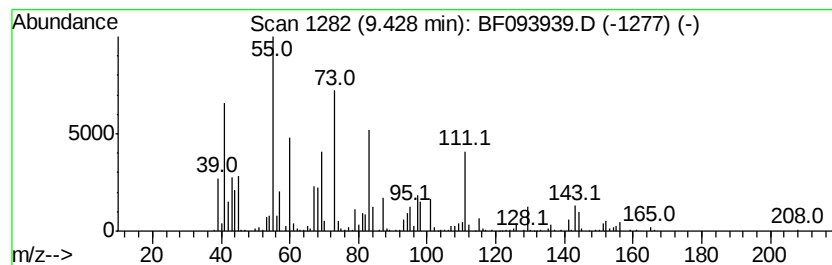
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Peak Number 5 9-Oxononanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.43	19.00 ng	1324510	Acenaphthene-d10	9.59

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Oxononanoic acid	172	C9H16O3	002553-17-5	59
2		6-Dodecanol	186	C12H26O	006836-38-0	22
3		Pentanoic acid, 5-bromo-	180	C5H9BrO2	002067-33-6	16
4		Undecane, 4-cyclohexyl-	238	C17H34	013151-79-6	14
5		4-Nonanol	144	C9H20O	005932-79-6	14



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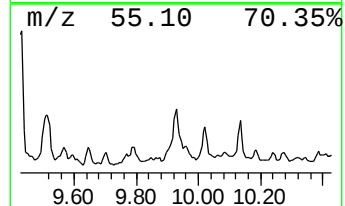
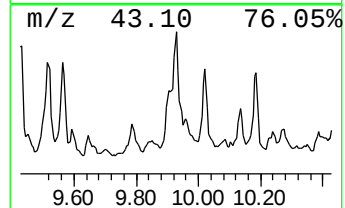
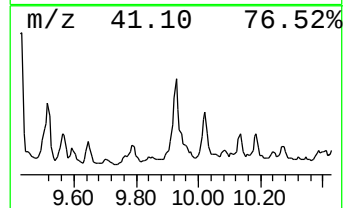
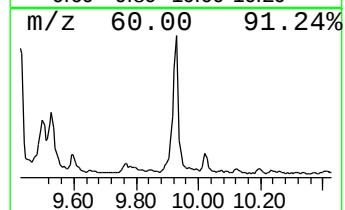
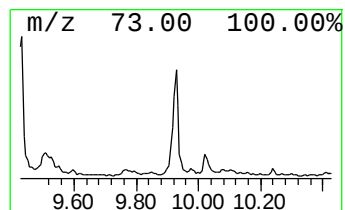
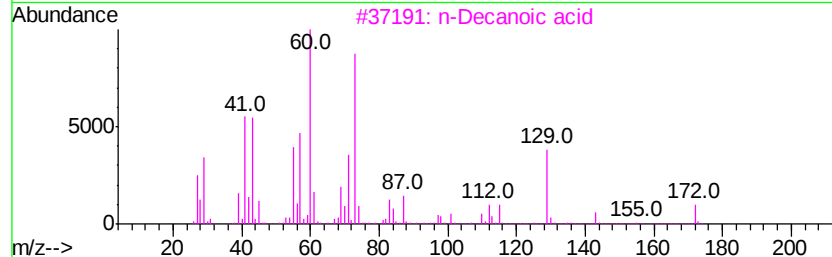
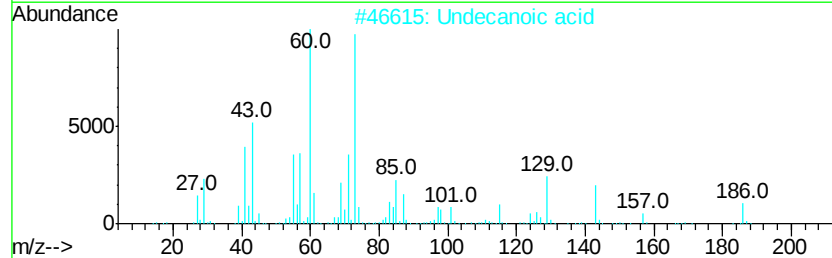
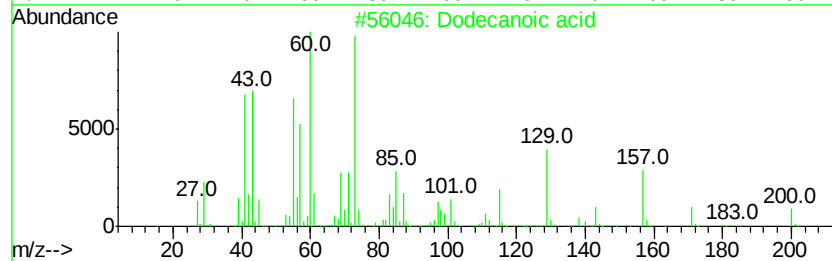
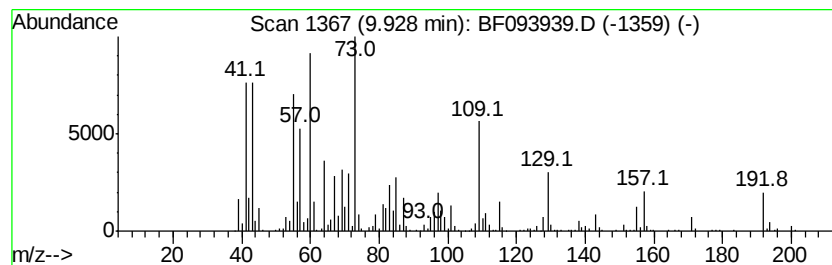
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Dodecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	17.43 ng	1214630	Acenaphthene-d10	9.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanoic acid	200	C12H24O2	000143-07-7	64
2		Undecanoic acid	186	C11H22O2	000112-37-8	46
3		n-Decanoic acid	172	C10H20O2	000334-48-5	35
4		Nonanoic acid	158	C9H18O2	000112-05-0	30
5		d-Glucohexodialdose	178	C6H10O6	1000149-19-1	27



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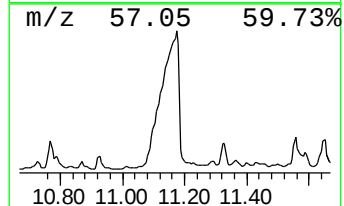
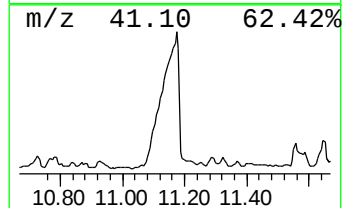
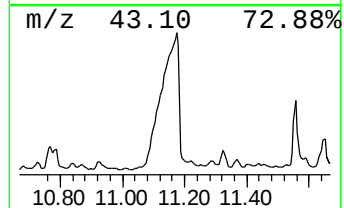
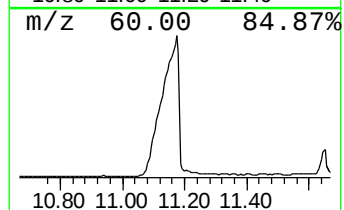
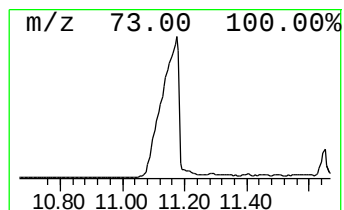
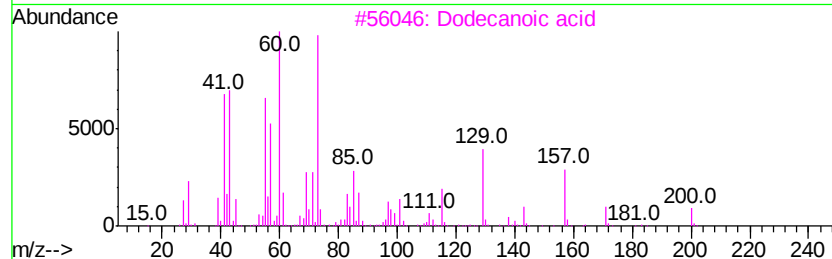
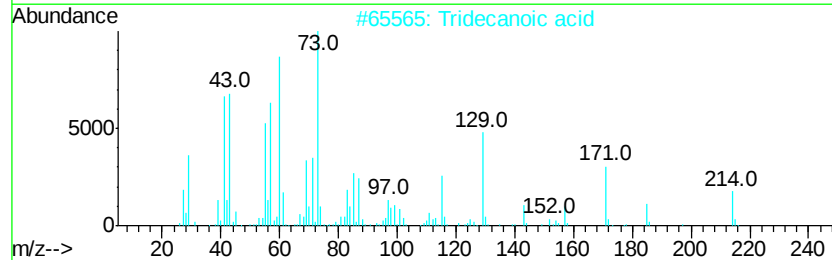
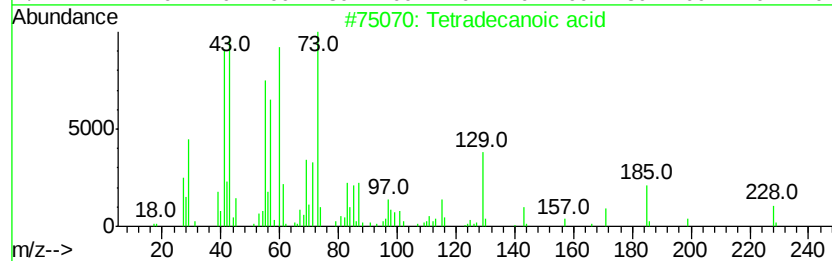
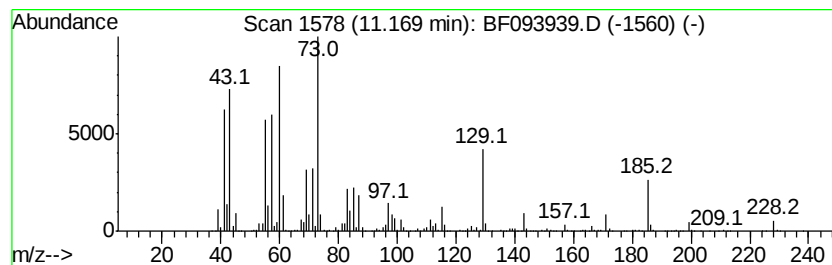
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Tetradecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.17	247.51 ng	15391600	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanoic acid	228	C14H28O2	000544-63-8	97
2		Tridecanoic acid	214	C13H26O2	000638-53-9	80
3		Dodecanoic acid	200	C12H24O2	000143-07-7	74
4		n-Decanoic acid	172	C10H20O2	000334-48-5	64
5		Undecanoic acid	186	C11H22O2	000112-37-8	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
Data File : BF093939.D
Acq On : 27 Mar 2017 12:54
Operator : SJ/MA
Sample : I2367-08
Misc :
ALS Vial : 3 Sample Multiplier: 1

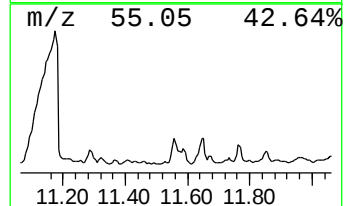
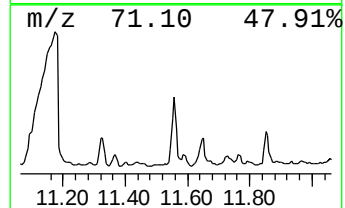
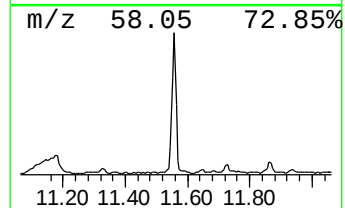
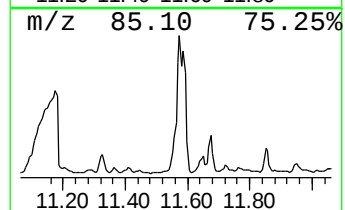
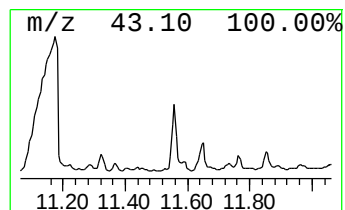
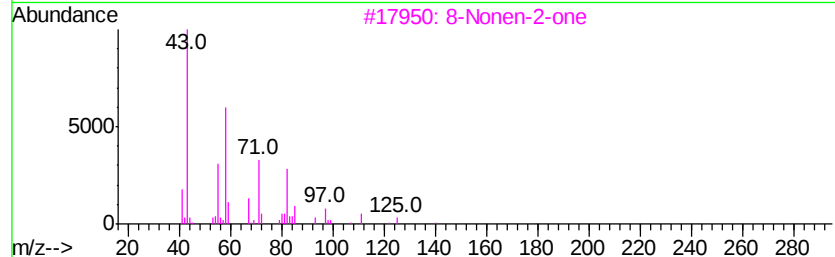
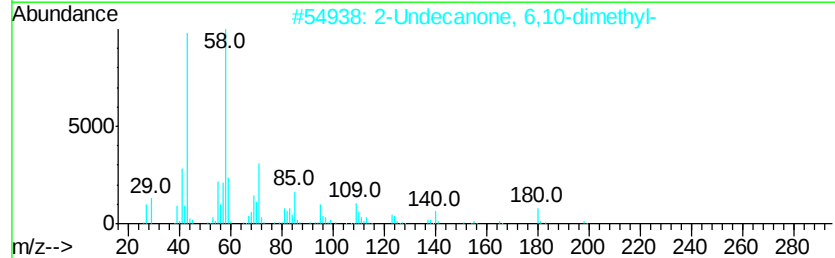
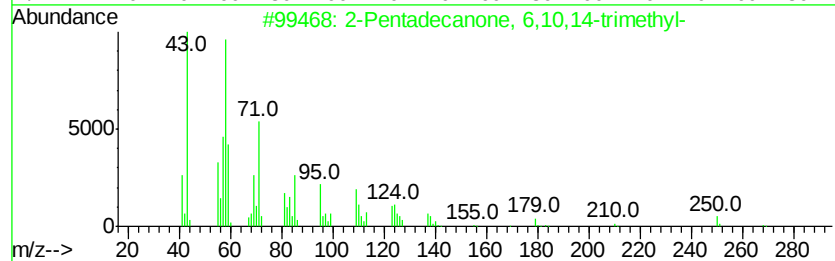
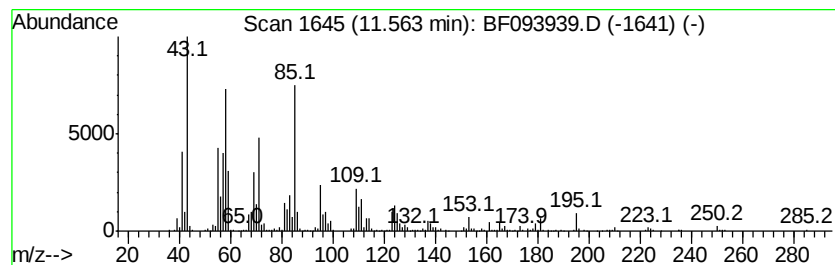
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2-Pentadecanone, 6,10,14-tr... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.56	22.25 ng	1383740	Phenanthrene-d10	11.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentadecanone, 6,10,14-trimethyl-	268	C18H36O	000502-69-2	64
2	2-Undecanone, 6,10-dimethyl-	198	C13H26O	001604-34-8	59
3	8-Nonen-2-one	140	C9H16O	005009-32-5	58
4	2-Heptanone, 4-methyl-	128	C8H16O	006137-06-0	47
5	5-Azulenemethanol, 1,2,3,3a,4,5,...	222	C15H26O	055255-90-8	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032717\
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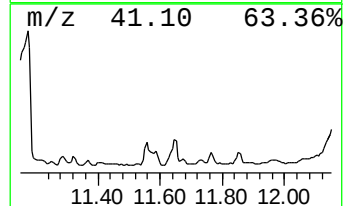
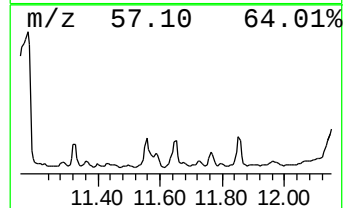
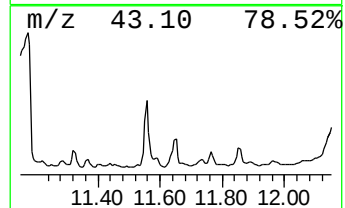
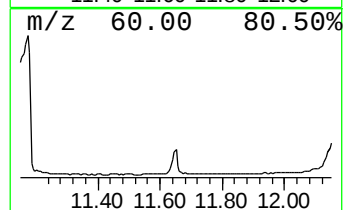
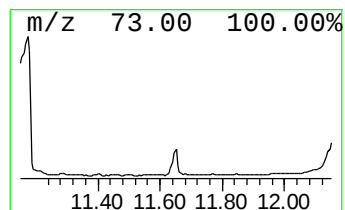
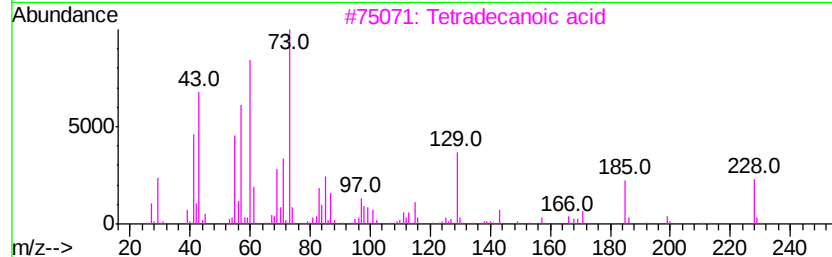
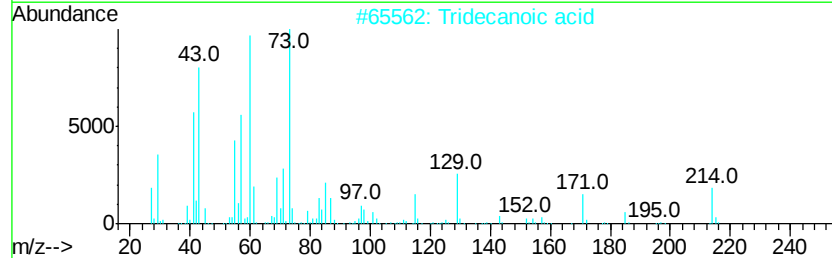
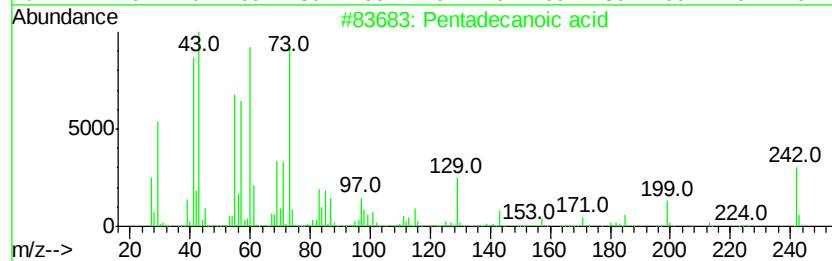
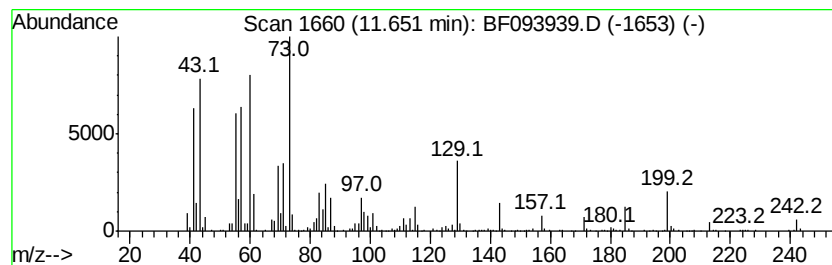
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Pentadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	19.48 ng	1211380	Phenanthrene-d10	11.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecanoic acid	242	C15H30O2	001002-84-2	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Undecanoic acid	186	C11H22O2	000112-37-8	68
5		n-Decanoic acid	172	C10H20O2	000334-48-5	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032717\
Data File : BF093939.D
Acq On : 27 Mar 2017 12:54
Operator : SJ/MA
Sample : I2367-08
Misc :
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Instrument :
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ClientSampleId :
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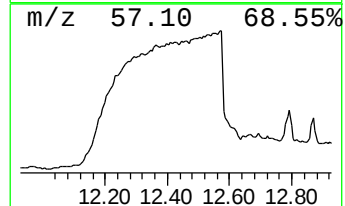
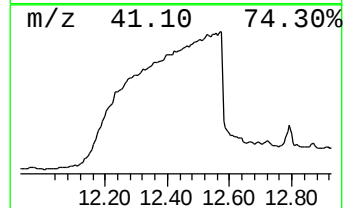
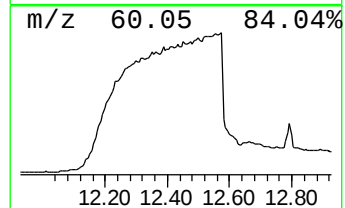
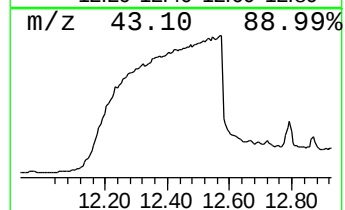
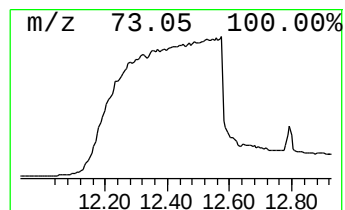
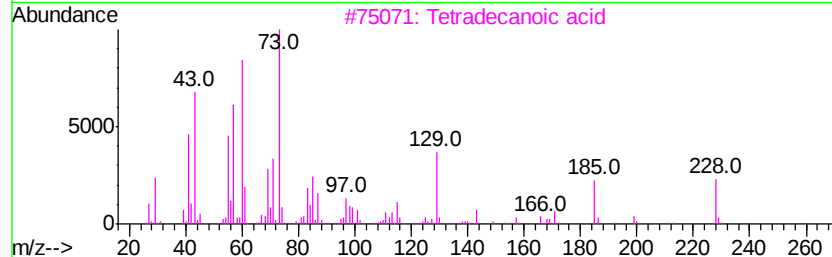
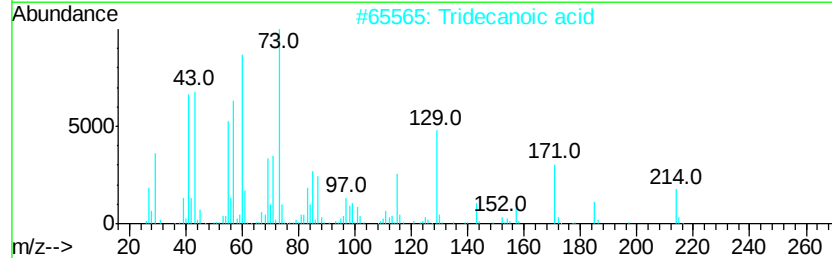
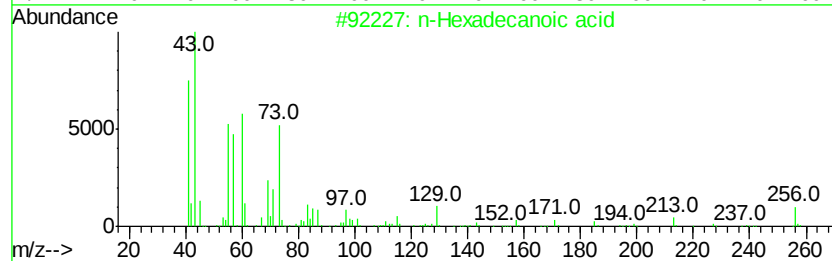
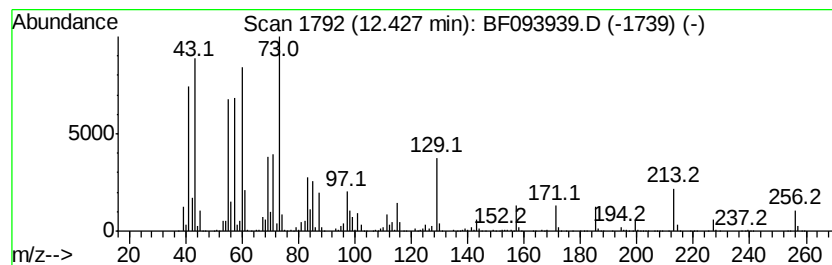
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	1580.41 ng	98281300	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
5		Undecane	156	C11H24	001120-21-4	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
Data File : BF093939.D
Acq On : 27 Mar 2017 12:54
Operator : SJ/MA
Sample : I2367-08
Misc :
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Instrument :
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ClientSampleId :
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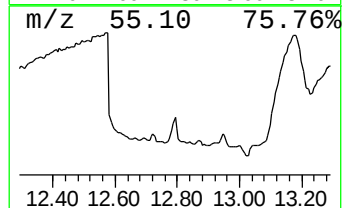
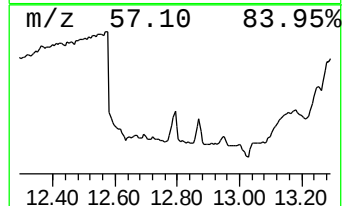
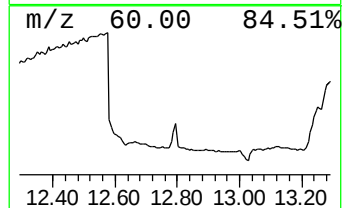
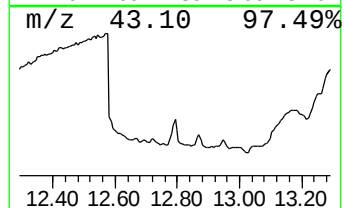
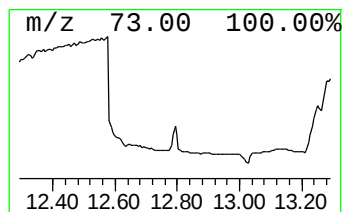
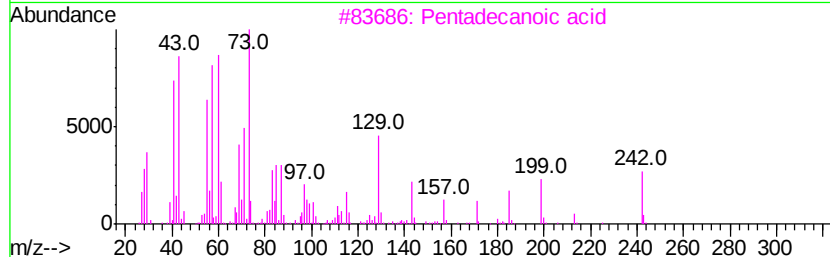
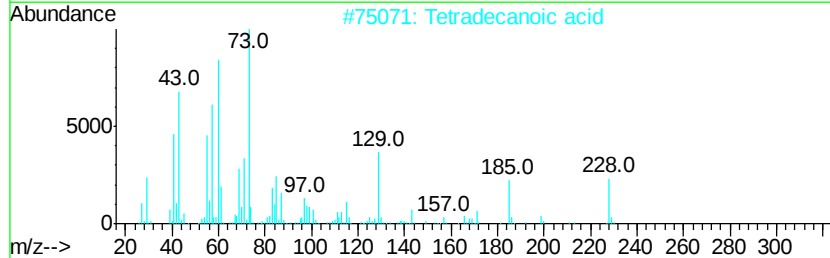
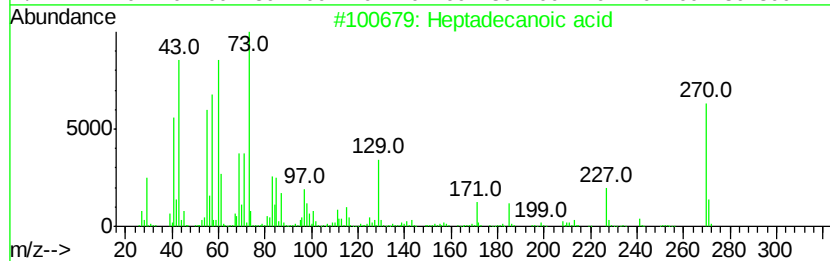
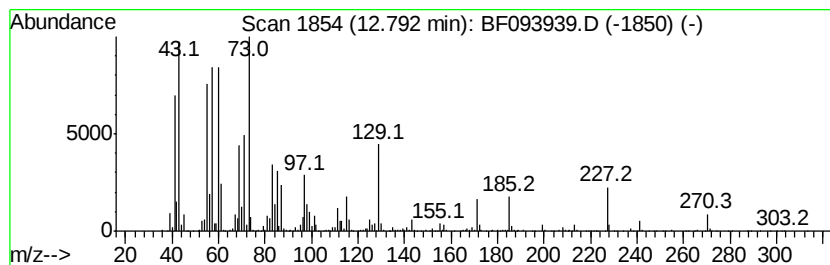
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	26.99 ng	1678550	Phenanthrene-d10	11.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecanoic acid	270	C17H34O2	000506-12-7	95
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4		n-Decanoic acid	172	C10H20O2	000334-48-5	68
5		Tridecanoic acid	214	C13H26O2	000638-53-9	68



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
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Acq On : 27 Mar 2017 12:54
Operator : SJ/MA
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
E2-BB10-NSW

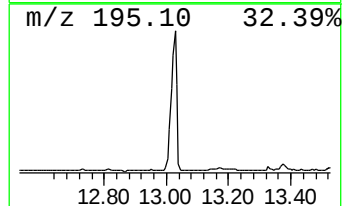
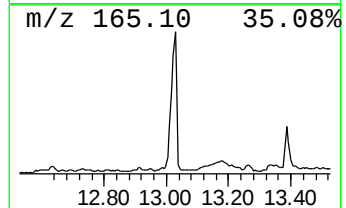
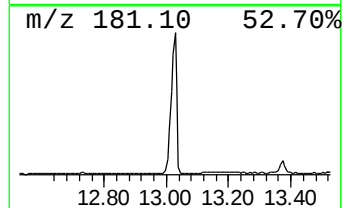
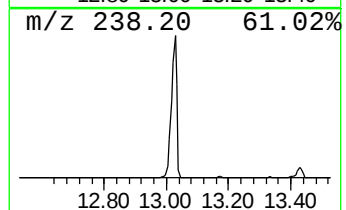
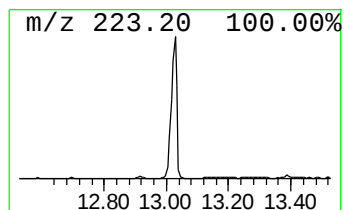
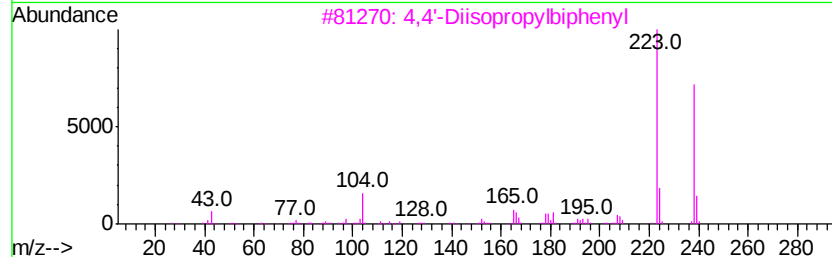
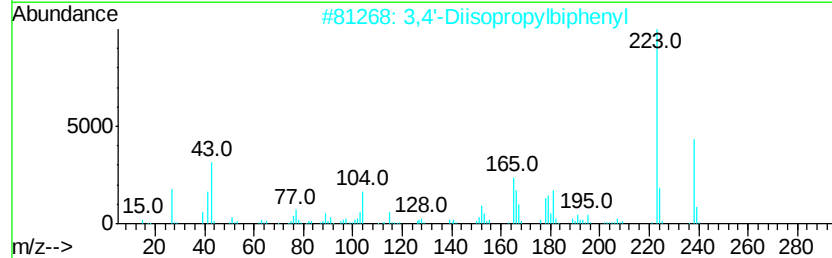
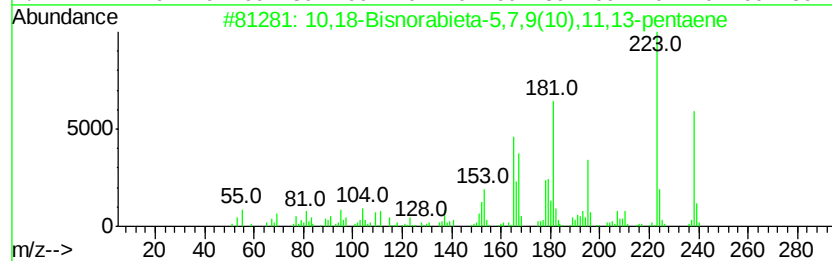
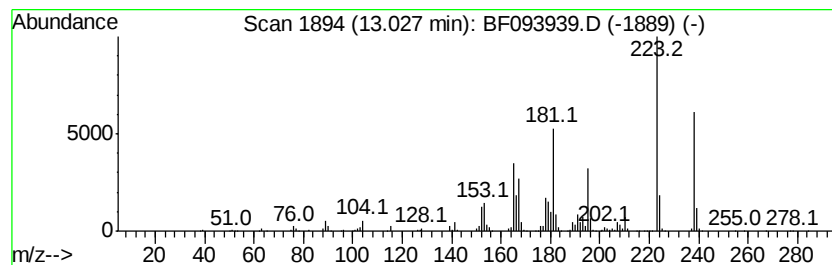
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	57.63 ng	3584120	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	86
3		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60
4		4,4'-Diacetyl biphenyl	238	C16H14O2	000787-69-9	46
5		9,10-Anthracenedione, 2-amino-	223	C14H9NO2	000117-79-3	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
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Acq On : 27 Mar 2017 12:54
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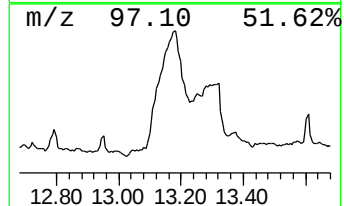
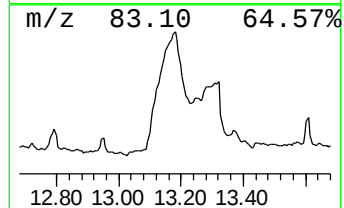
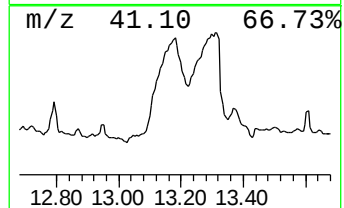
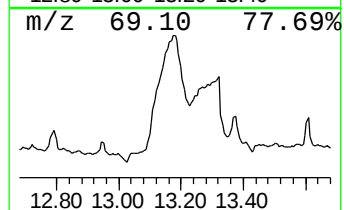
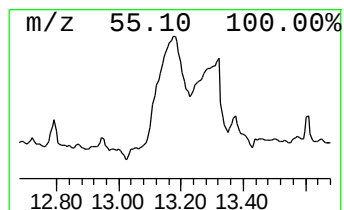
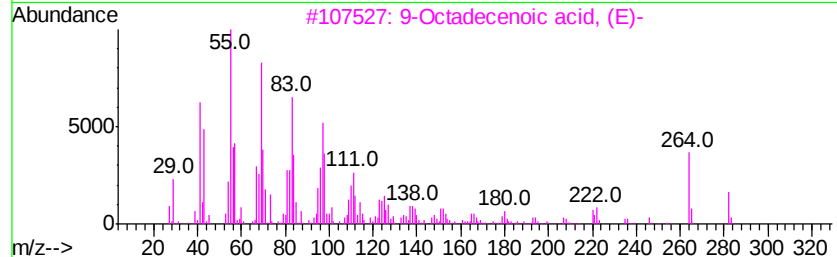
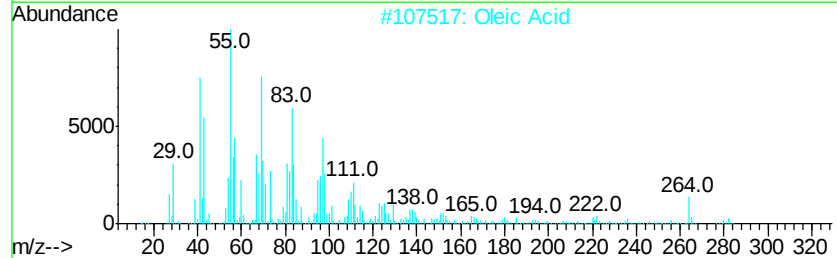
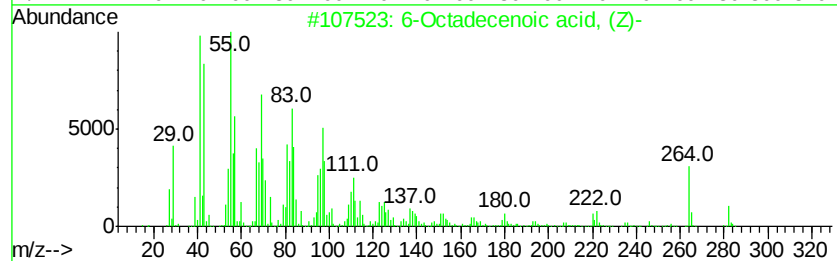
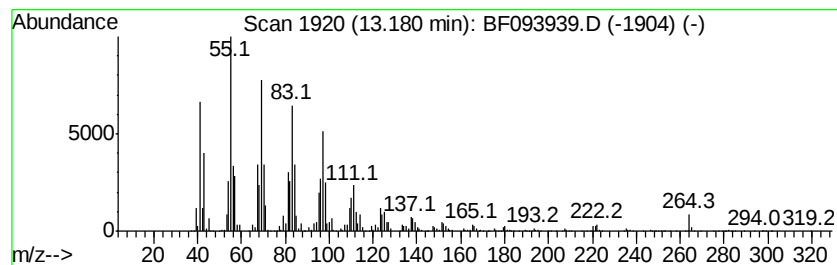
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 6-Octadecenoic acid, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	248.95 ng	25387300	Chrysene-d12	14.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	96
2		Oleic Acid	282	C18H34O2	000112-80-1	91
3		9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	90
4		15-Tetracosenoic acid, methyl es...	380	C25H48O2	002733-88-2	60
5		9-Hexadecenoic acid	254	C16H30O2	002091-29-4	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
Data File : BF093939.D
Acq On : 27 Mar 2017 12:54
Operator : SJ/MA
Sample : I2367-08
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-BB10-NSW

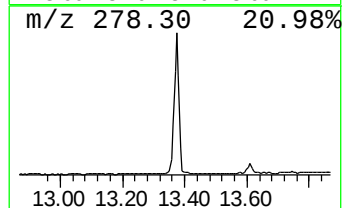
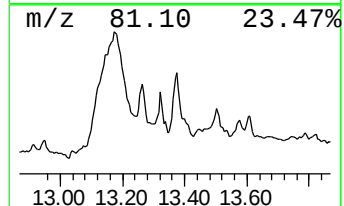
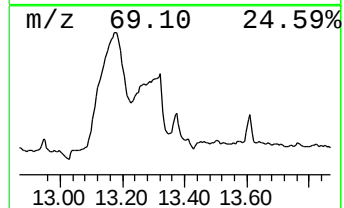
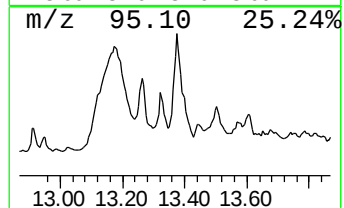
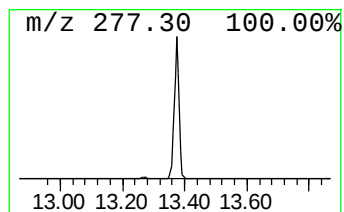
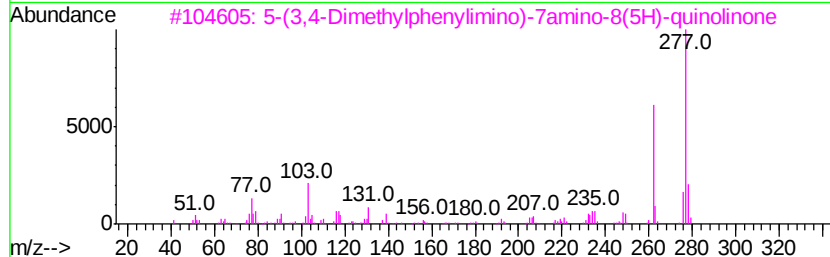
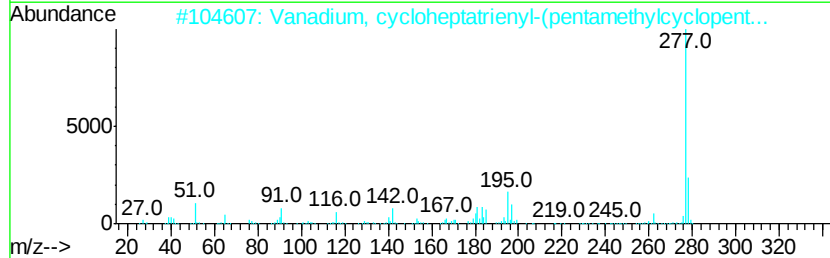
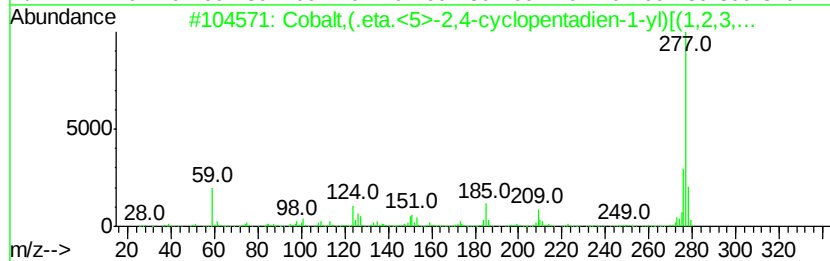
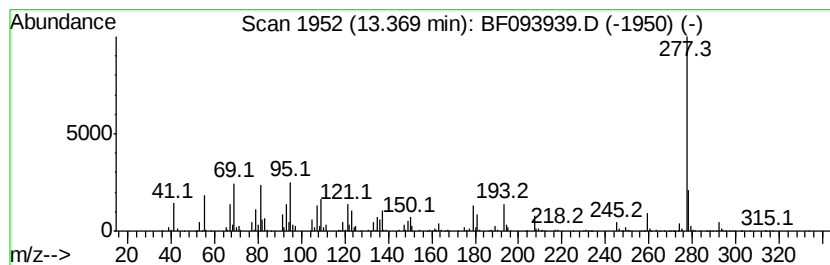
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown13.37 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	22.40 ng	2284370	Chrysene-d12	14.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cobalt,(.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	47
2		Vanadium, cycloheptatrienyl-(pen...	277	C17H22V	1000155-20-5	43
3		5-(3,4-Dimethylphenylimino)-7ami...	277	C17H15N3O	1000254-01-4	43
4		Benzonitrile, 4-[3-(3,4-methylen...	277	C17H11N03	1000273-95-6	43
5		Quinoline-2,4(1H,3H)-dione, 3-(2...	277	C16H11N3O2	1000263-14-4	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032717\
Data File : BF093939.D
Acq On : 27 Mar 2017 12:54
Operator : SJ/MA
Sample : I2367-08
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-BB10-NSW

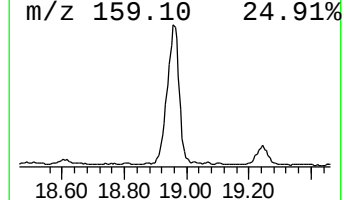
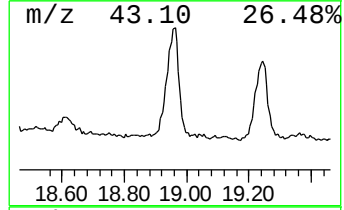
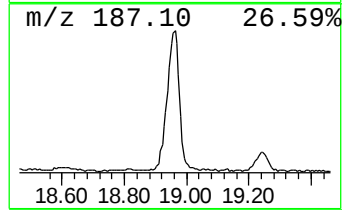
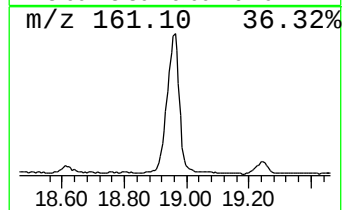
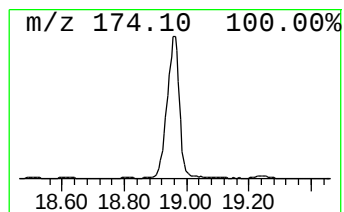
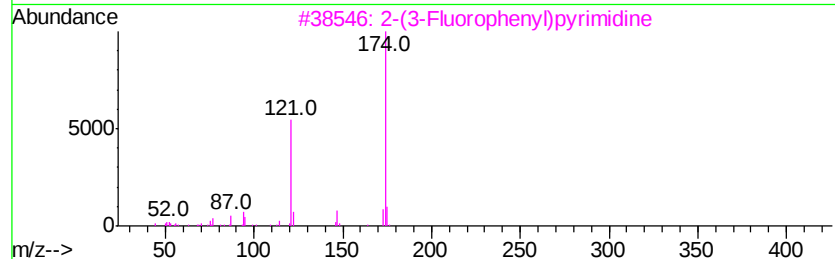
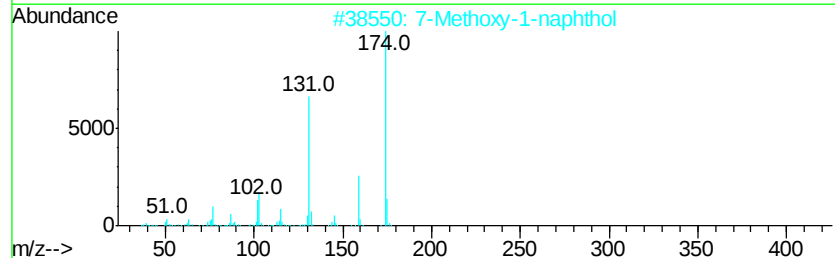
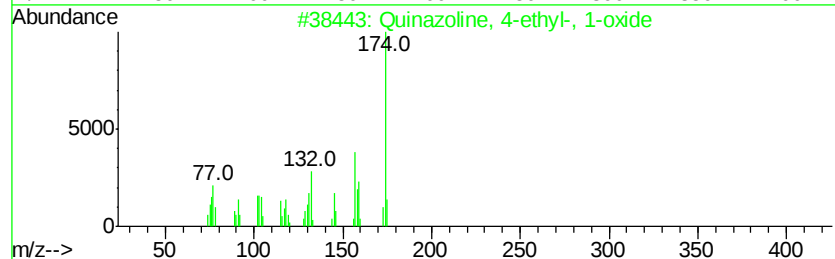
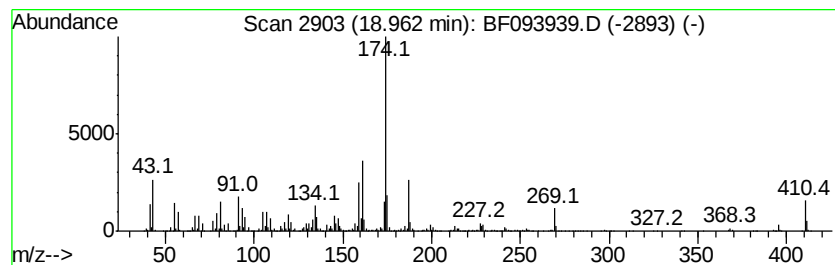
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown18.96 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.96	20.00 ng	3928850	Perylene-d12	16.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Quinazoline, 4-ethyl-, 1-oxide	174	C10H10N2O	037920-75-5	47
2		7-Methoxy-1-naphthol	174	C11H10O2	067247-13-6	43
3		2-(3-Fluorophenyl)pyrimidine	174	C10H7FN2	068049-16-1	43
4		2-(4-Fluorophenyl)pyrimidine	174	C10H7FN2	068049-17-2	38
5		1-Phenyl-4-hydroxy-5-methylpyrazole	174	C10H10N2O	089193-19-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032717\
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Operator : SJ/MA
Sample : I2367-08
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-BB10-NSW

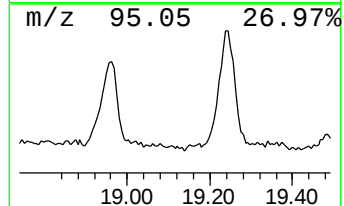
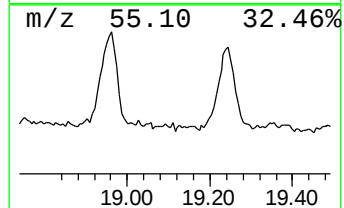
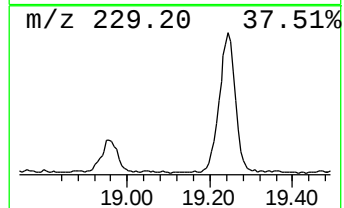
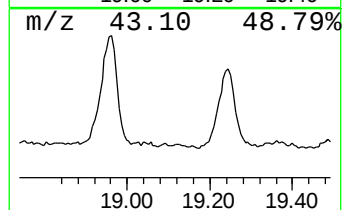
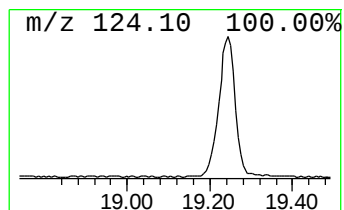
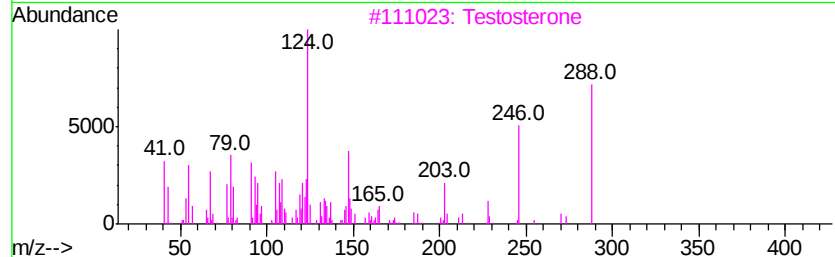
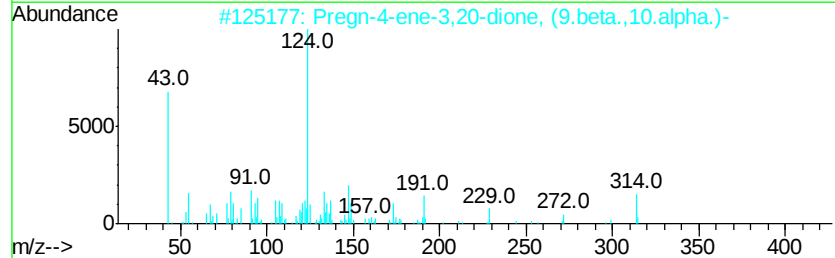
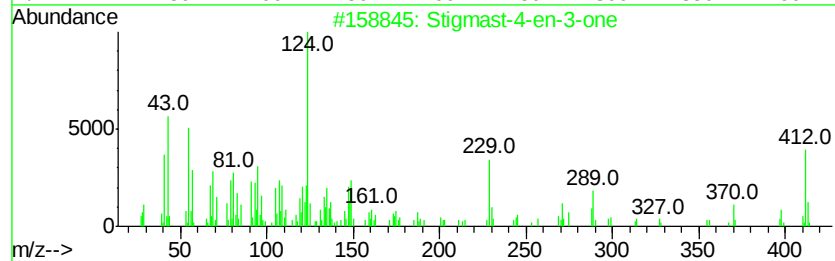
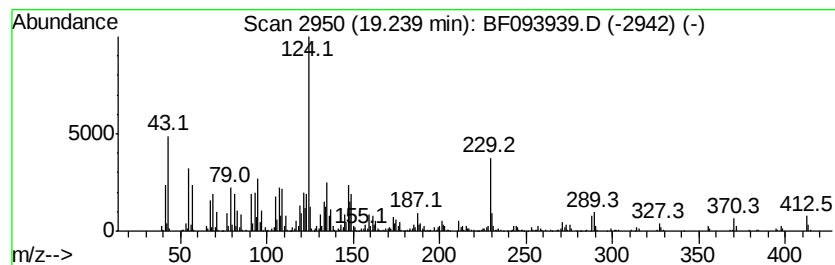
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Stigmast-4-en-3-one Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	12.54 ng	2463570	Perylene-d12	16.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	90
2		Pregn-4-ene-3,20-dione, (9.beta....	314	C21H30O2	002755-10-4	47
3		Testosterone	288	C19H28O2	000058-22-0	46
4		Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000604-39-7	45
5		2,3-Diaminophenol	124	C6H8N2O	059649-56-8	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032717\
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 Operator : SJ/MA
 Sample : I2367-08
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-BB10-NSW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexanal	3.60	25.7	ng	1288560	1	5.95	1002470	20.0
unknown5.69	5.69	101.8	ng	5104890	1	5.95	1002470	20.0
Nonanoic acid	7.97	25.1	ng	1835290	2	7.45	1460850	20.0
9-Oxononanoic acid	9.43	19.0	ng	1324510	3	9.59	1393880	20.0
Dodecanoic acid	9.93	17.4	ng	1214630	3	9.59	1393880	20.0
Tetradecanoic acid	11.17	247.5	ng	15391600	4	11.42	1243740	20.0
2-Pentadecanone, ...	11.56	22.3	ng	1383740	4	11.42	1243740	20.0
Pentadecanoic acid	11.65	19.5	ng	1211380	4	11.42	1243740	20.0
n-Hexadecanoic acid	12.43	1580.4	ng	98281300	4	11.42	1243740	20.0
4b,8-Dimethyl-2-i...	12.63	21.4	ng	1332480	4	11.42	1243740	20.0
Heptadecanoic acid	12.79	27.0	ng	1678550	4	11.42	1243740	20.0
10,18-Bisnorabiet...	13.03	57.6	ng	3584120	4	11.42	1243740	20.0
6-Octadecenoic ac...	13.18	248.9	ng	25387300	5	14.72	2039530	20.0
unknown13.37	13.37	22.4	ng	2284370	5	14.72	2039530	20.0
unknown18.96	18.96	20.0	ng	3928850	6	16.33	3928850	20.0
Stigmast-4-en-3-one	19.24	12.5	ng	2463570	6	16.33	3928850	20.0