

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062516\
 Data File : BF088407.D
 Acq On : 26 Jun 2016 7:58
 Operator : UM/SJ
 Sample : H3602-13
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q8-B5(0-16)C

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-BF060716.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	1.701	9	10	37	rVB	202738	387791	1.88%	0.800%
2	5.142	309	311	335	rBV	1130274	1398823	6.79%	2.884%
3	6.227	403	406	413	rBV	1126146	1332926	6.47%	2.748%
4	6.330	413	415	430	rVB	1466312	1514891	7.35%	3.123%
5	6.559	433	435	442	rBV	356745	376338	1.83%	0.776%
6	6.719	447	449	451	rBV	977565	1193813	5.79%	2.461%
7	7.142	483	486	488	rBV	578721	854065	4.14%	1.761%
8	7.850	546	548	550	rBV	397893	491644	2.38%	1.014%
9	8.925	640	642	644	rBV	1764993	1643185	7.97%	3.388%
10	9.325	675	677	681	rBV	317184	279217	1.35%	0.576%
11	9.588	698	700	703	rVV	558967	552315	2.68%	1.139%
12	10.376	767	769	772	rBV	684267	742202	3.60%	1.530%
13	10.616	787	790	792	rBV	360213	397876	1.93%	0.820%
14	11.062	827	829	834	rBV	423561	497095	2.41%	1.025%
15	11.405	855	859	861	rBV2	508612	793087	3.85%	1.635%
16	11.496	863	867	870	rBV2	3786235	6529404	31.67%	13.462%
17	12.228	922	931	934	rBV4	4162846	20616158	100.00%	42.504%
18	12.285	934	936	938	rVV	2989143	3109972	15.09%	6.412%
19	12.354	938	942	946	rVB2	81231	218483	1.06%	0.450%
20	12.502	953	955	957	rBV2	310974	326505	1.58%	0.673%
21	12.662	966	969	971	rBV	1013432	1177807	5.71%	2.428%
22	12.765	971	978	979	rVV4	92895	365368	1.77%	0.753%
23	12.845	983	985	988	rVV3	152189	351593	1.71%	0.725%
24	12.902	988	990	995	rVB2	296831	619898	3.01%	1.278%
25	12.994	995	998	1002	rBV2	247859	593082	2.88%	1.223%
26	13.062	1002	1004	1006	rVV	121523	248222	1.20%	0.512%
27	13.199	1013	1016	1022	rVB5	142334	279077	1.35%	0.575%
28	13.417	1031	1035	1039	rVB2	266494	585345	2.84%	1.207%
29	13.645	1053	1055	1057	rVV2	140908	213283	1.03%	0.440%
30	13.691	1057	1059	1066	rVB	292967	464140	2.25%	0.957%
31	15.051	1176	1178	1182	rVB	232937	350088	1.70%	0.722%

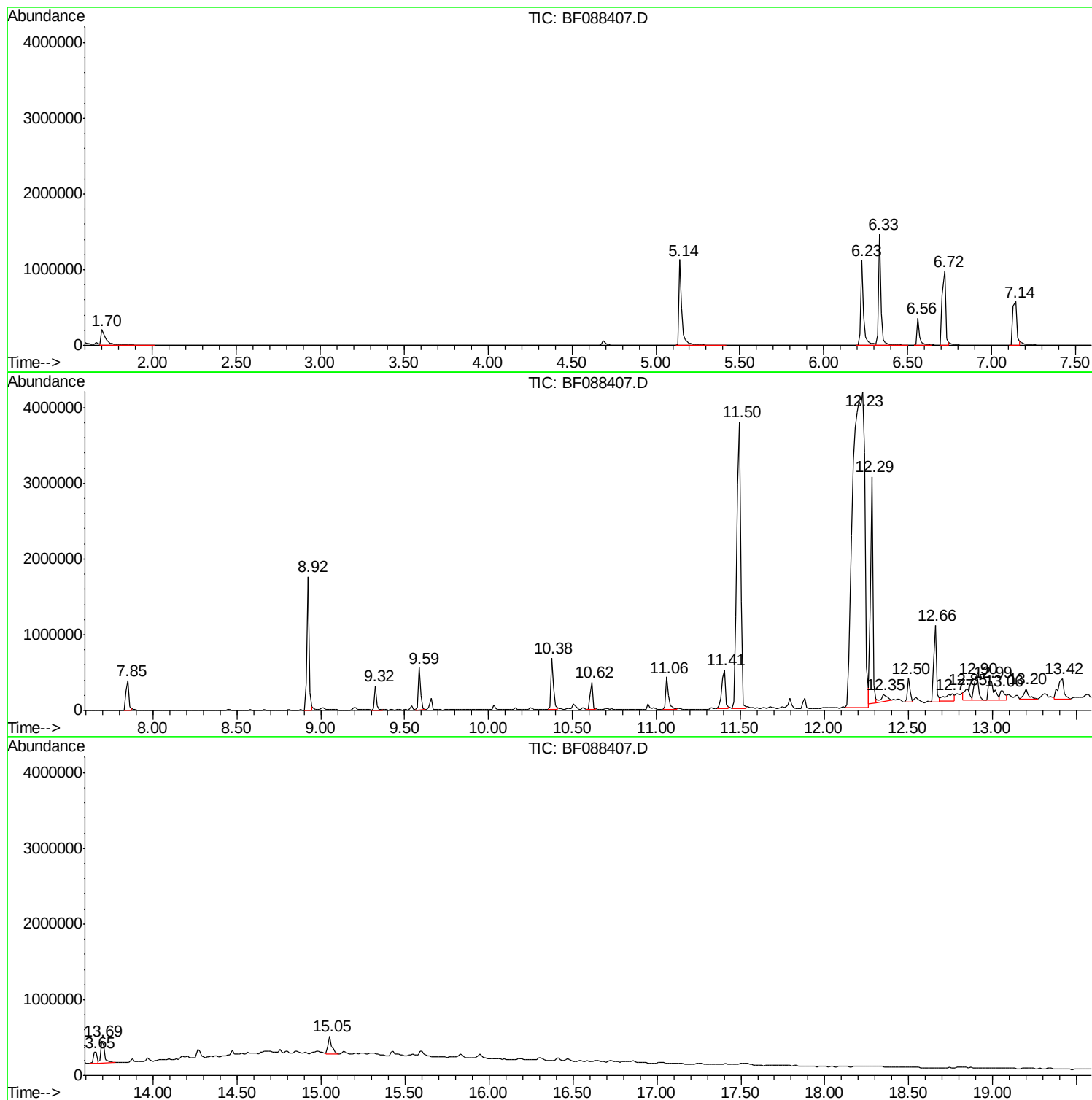
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ClientSampleId :
Q8-B5(0-16)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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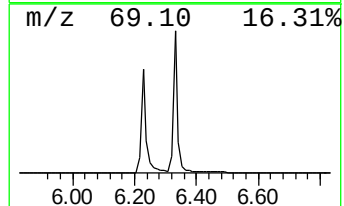
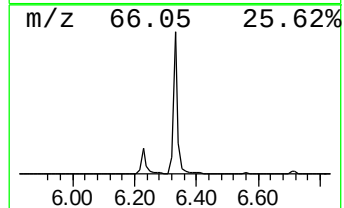
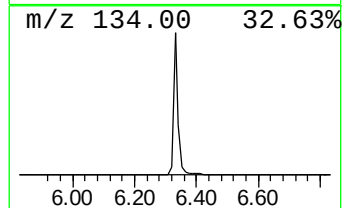
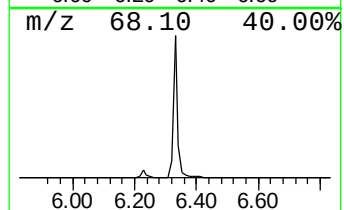
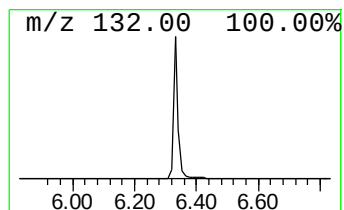
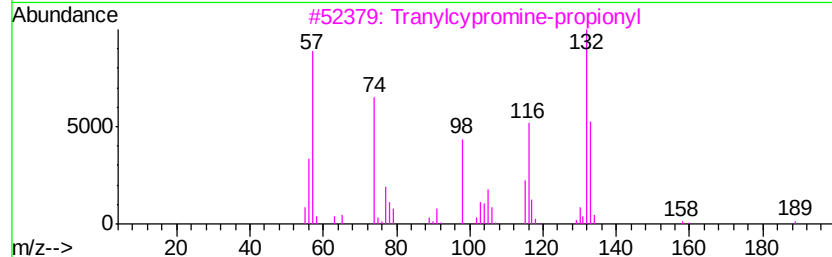
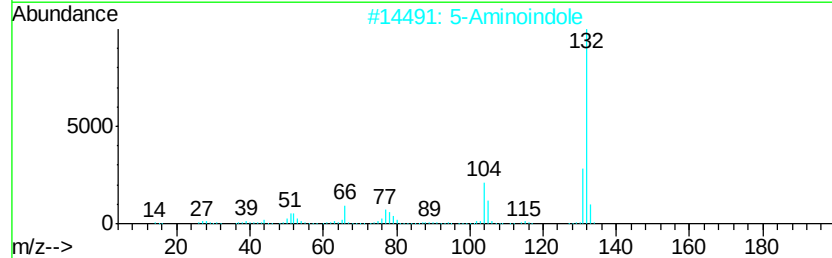
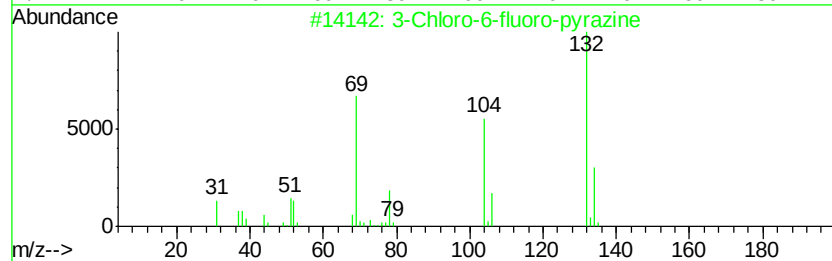
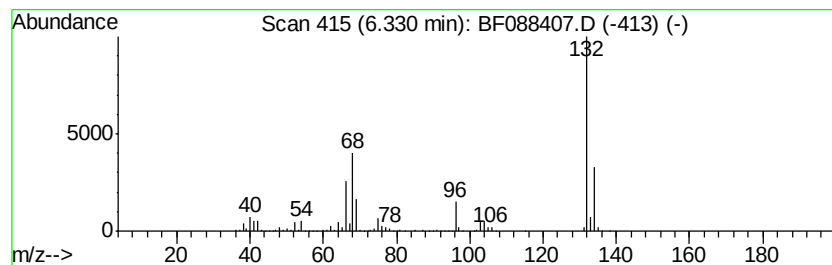
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.33 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	80.51 ng	1514890	1,4-Dichlorobenzene-d4	6.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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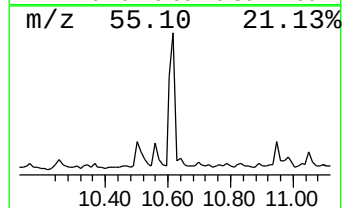
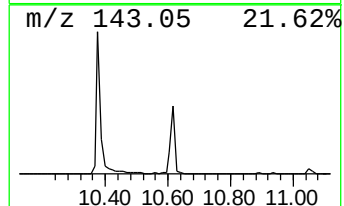
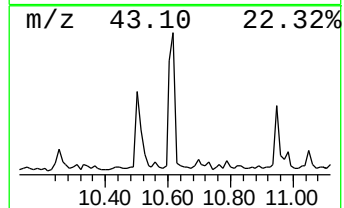
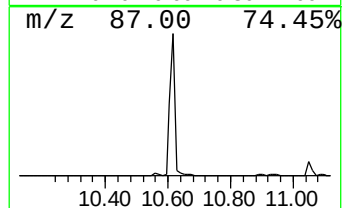
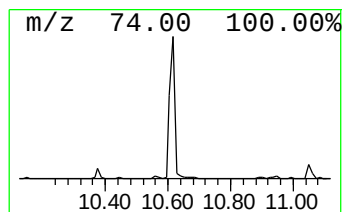
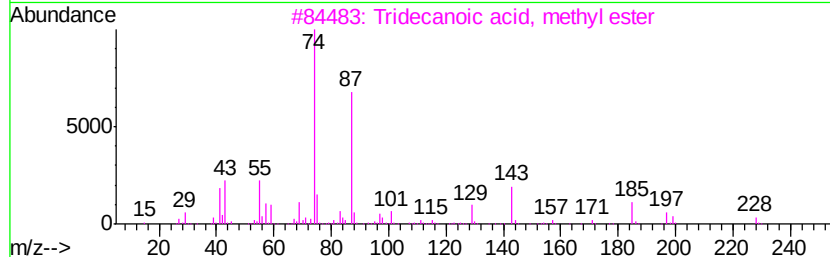
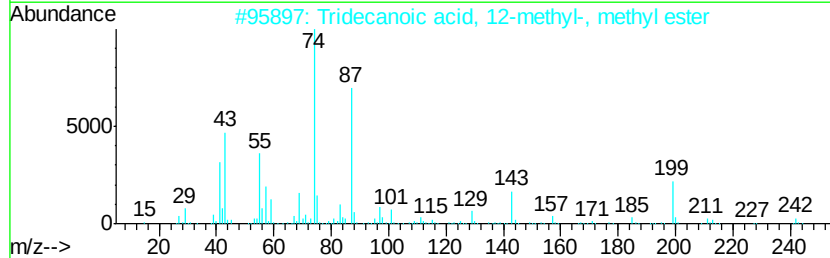
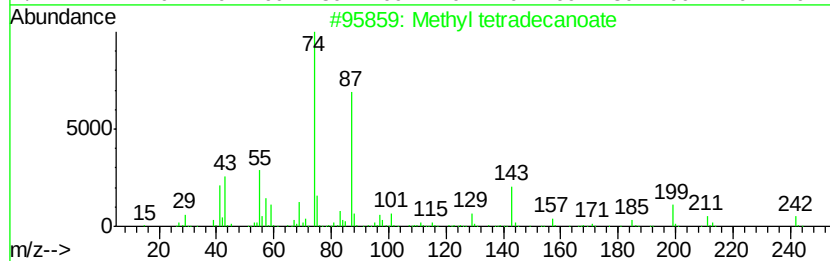
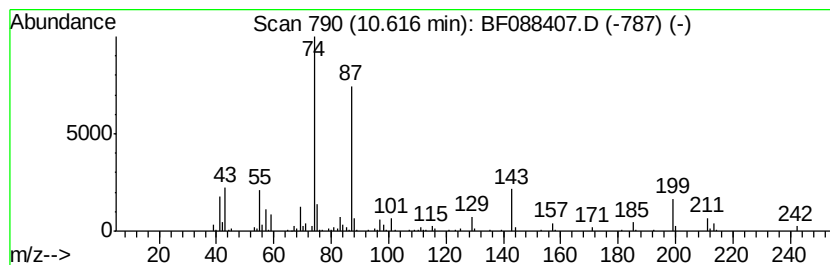
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Peak Number 4 Methyl tetradecanoate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.62	16.01 ng	397876	Phenanthrene-d10	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
2		Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	89
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	80
4		Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	80
5		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	72



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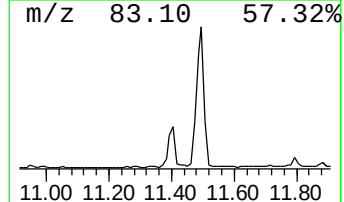
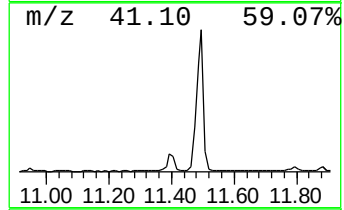
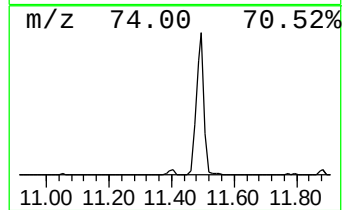
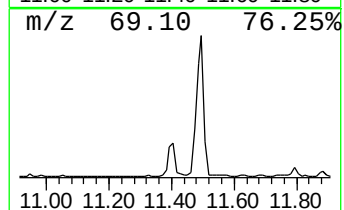
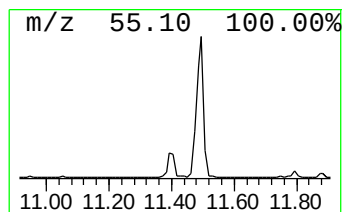
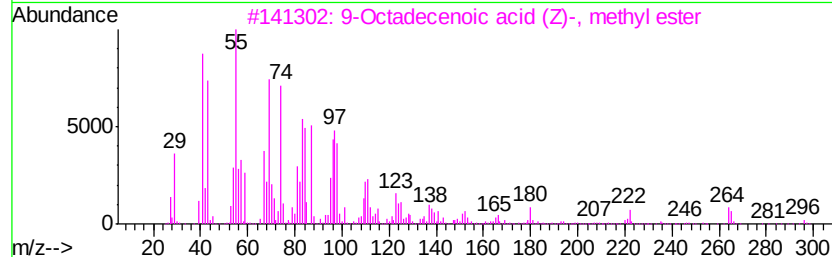
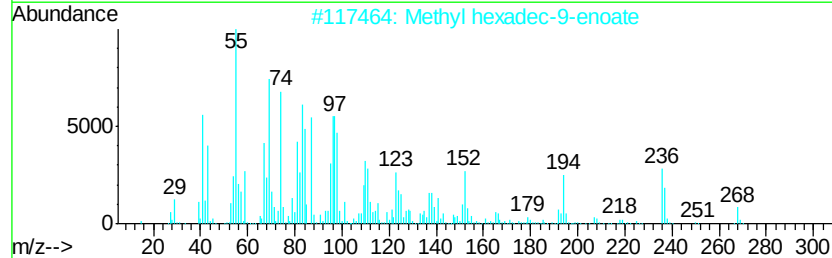
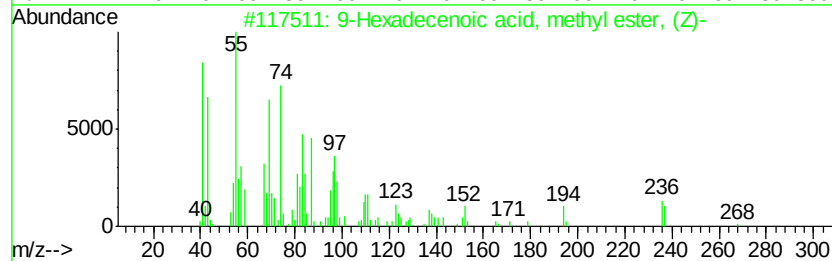
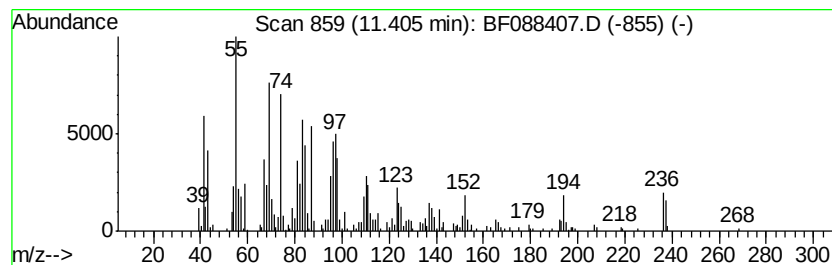
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Peak Number 5 9-Hexadecenoic acid, methyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.41	31.91 ng	793087	Phenanthrene-d10	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	99
2		Methyl hexadec-9-enoate	268	C17H32O2	010030-74-7	89
3		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	87
4		Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	70
5		7-Hexadecenoic acid, methyl este...	268	C17H32O2	056875-67-3	60



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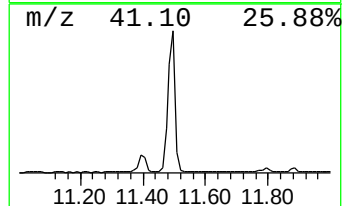
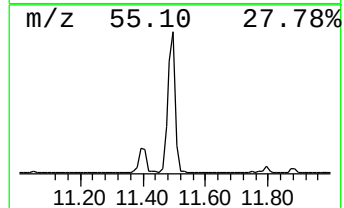
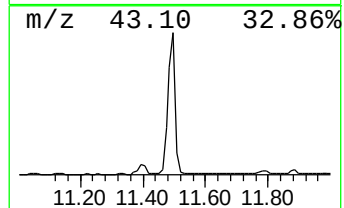
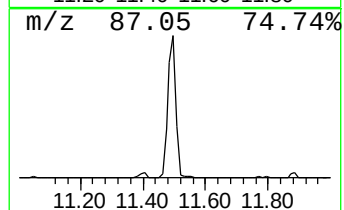
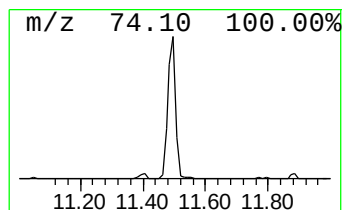
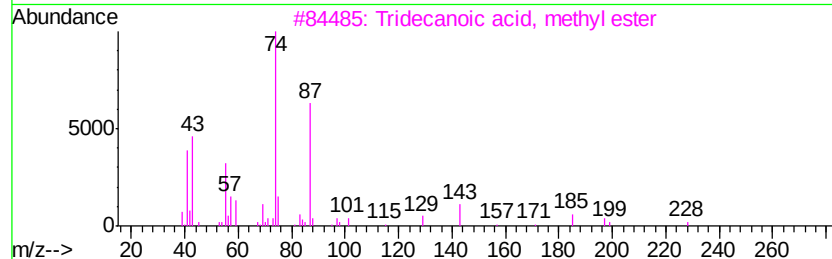
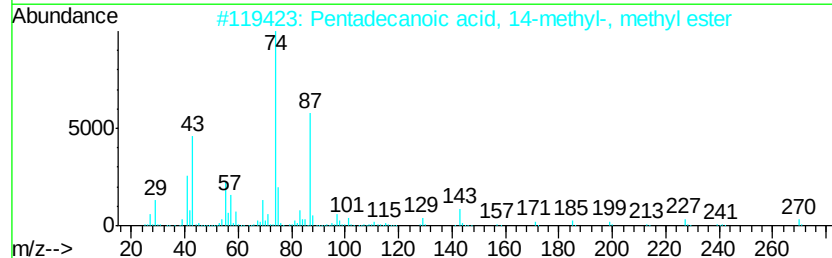
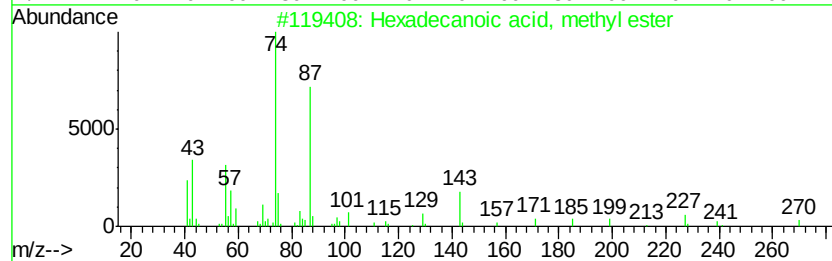
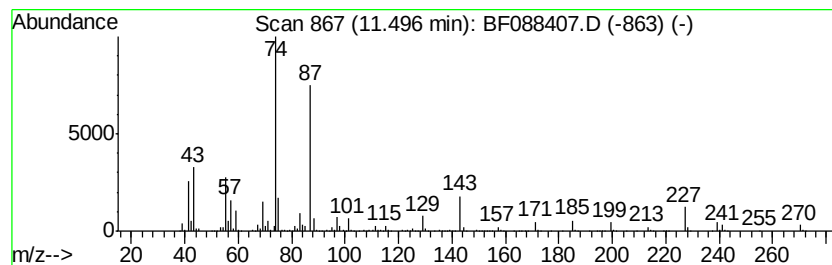
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Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.50	262.70 ng	6529400	Phenanthrene-d10	11.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87



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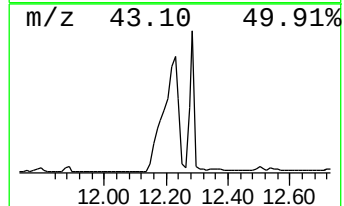
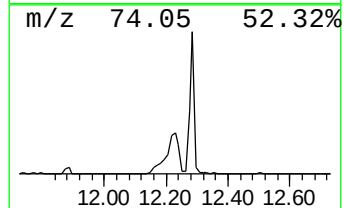
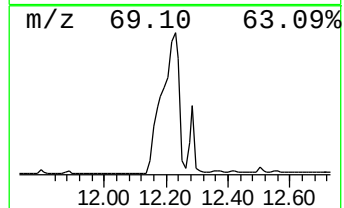
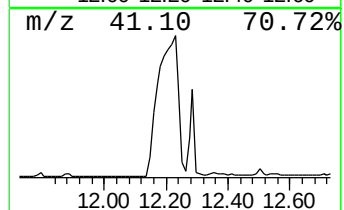
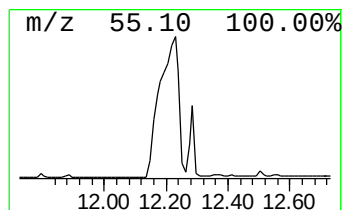
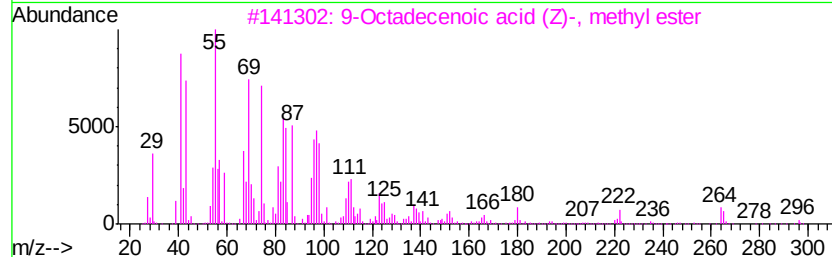
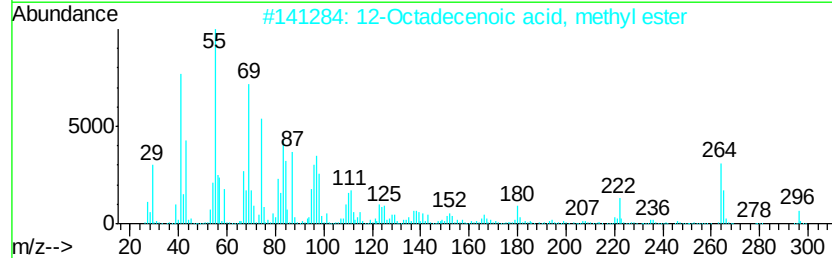
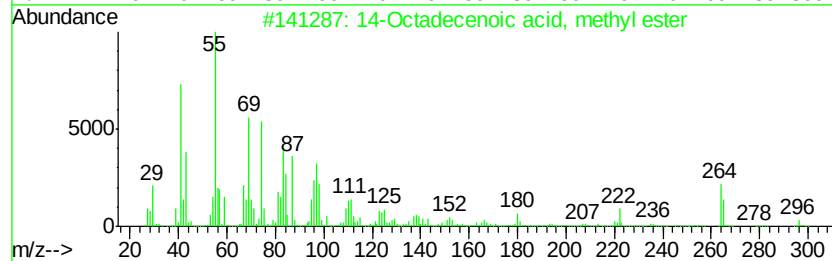
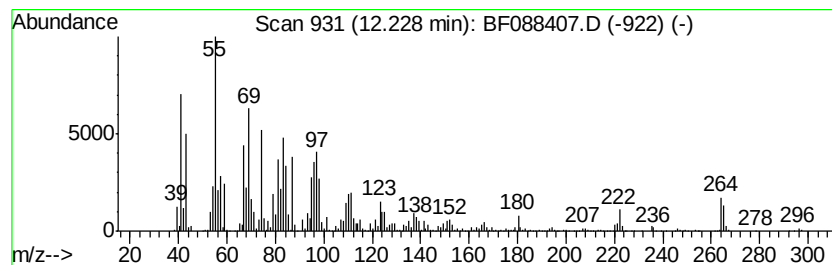
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Peak Number 7 14-Octadecenoic acid, methyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	829.47 ng	20616200	Phenanthrene-d10	11.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
2		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
3		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
4		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
5		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	98



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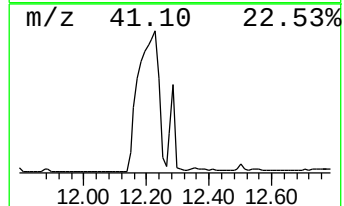
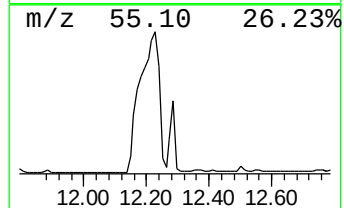
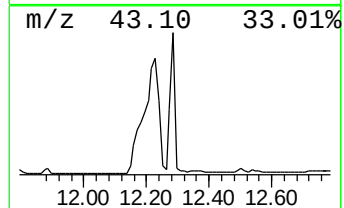
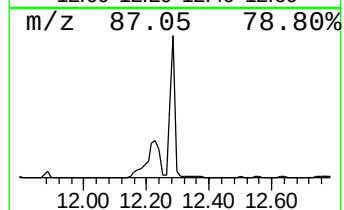
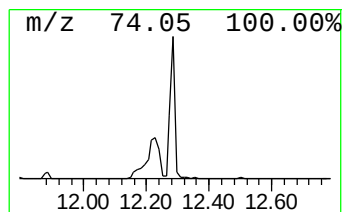
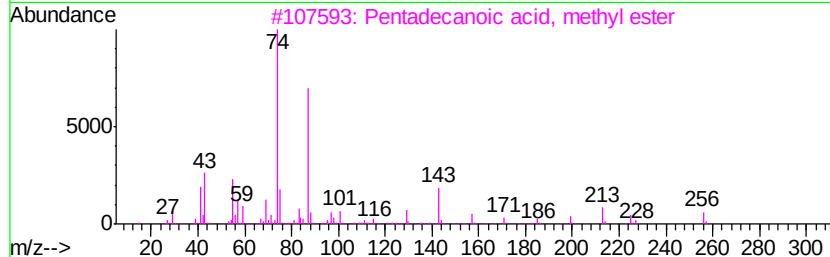
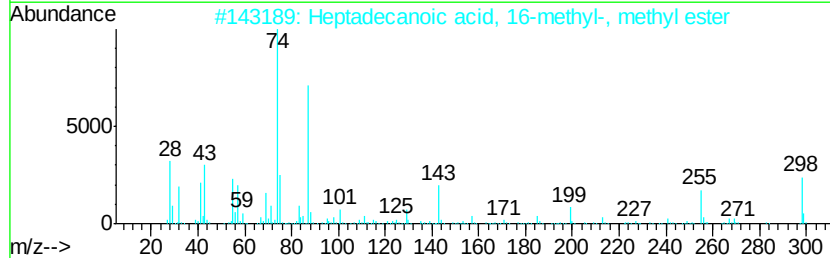
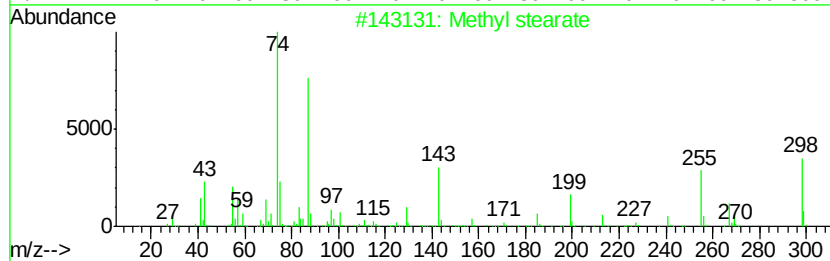
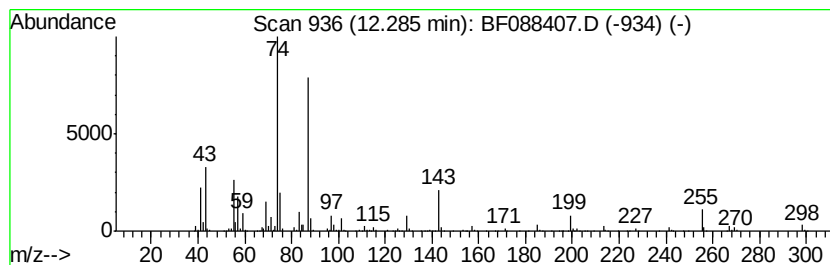
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Methyl stearate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	125.13 ng	3109970	Phenanthrene-d10	11.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl stearate	298	C19H38O2	000112-61-8	99
2		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	90
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
4		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
Data File : BF088407.D
Acq On : 26 Jun 2016 7:58
Operator : UM/SJ
Sample : H3602-13
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
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ClientSampleId :
Q8-B5(0-16)C

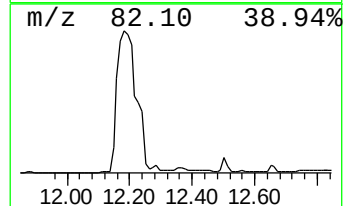
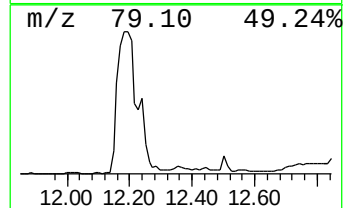
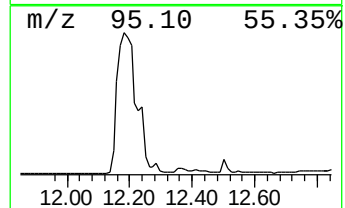
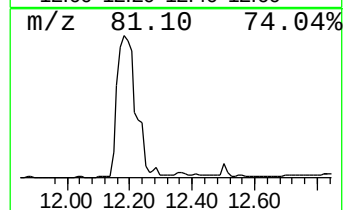
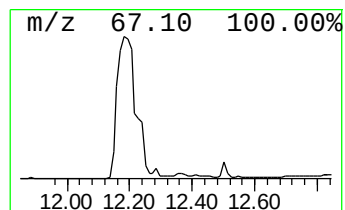
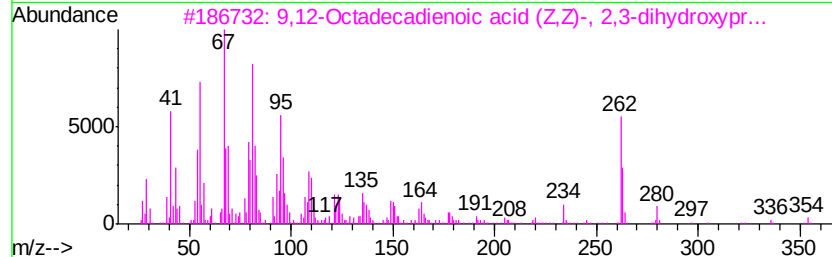
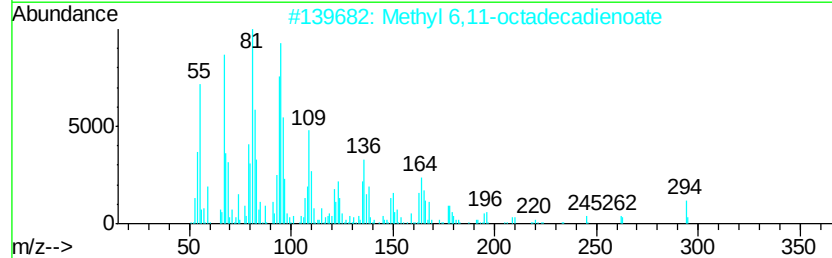
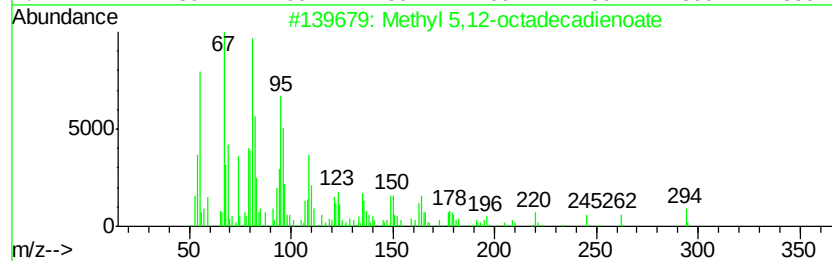
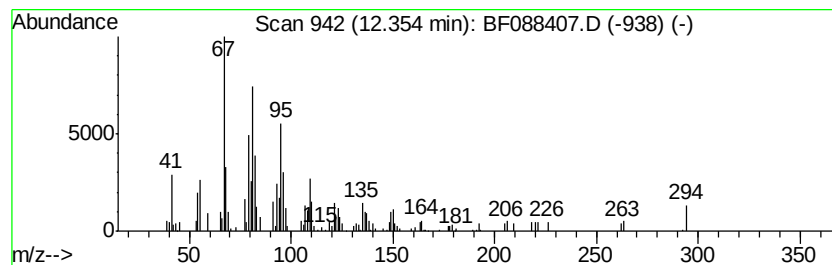
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl 5,12-octadecadienoate Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	8.79 ng	218483	Phenanthrene-d10	11.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 5,12-octadecadienoate	294	C19H34O2	1000336-43-1	95
2		Methyl 6,11-octadecadienoate	294	C19H34O2	1000336-43-2	93
3		9,12-Octadecadienoic acid (Z,Z)-...	354	C21H38O4	002277-28-3	91
4		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	86
5		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
Data File : BF088407.D
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Operator : UM/SJ
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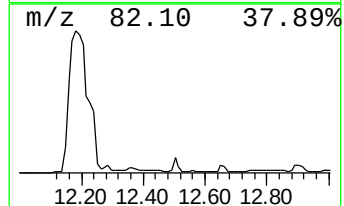
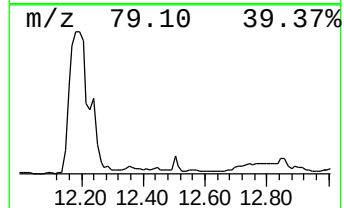
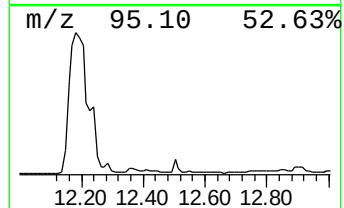
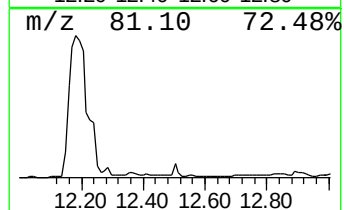
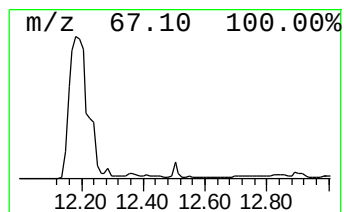
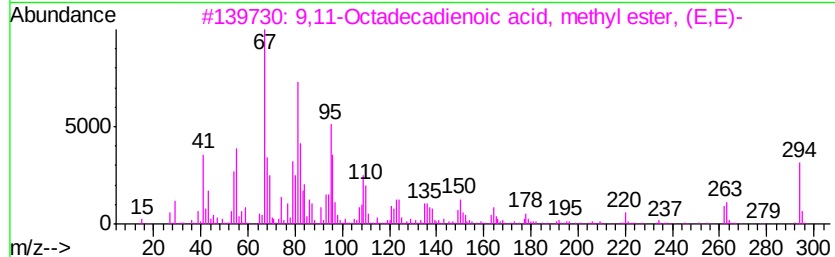
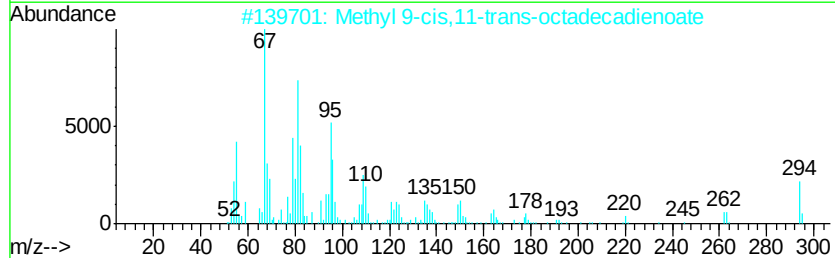
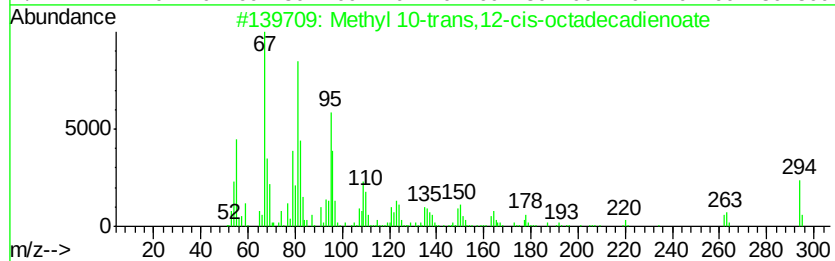
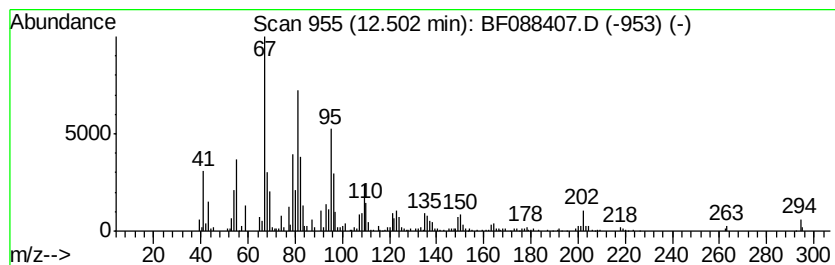
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Methyl 10-trans,12-cis-octa... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	14.07 ng	326505	Chrysene-d12	13.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99
2			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	96
3			9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	93
4			7,10-Octadecadienoic acid, methy...	294	C19H34O2	056554-24-6	93
5			9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
Data File : BF088407.D
Acq On : 26 Jun 2016 7:58
Operator : UM/SJ
Sample : H3602-13
Misc :
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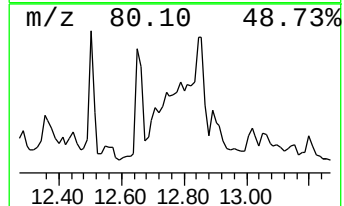
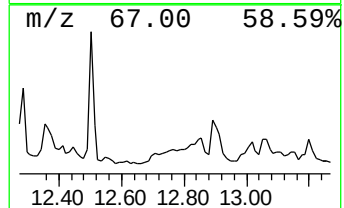
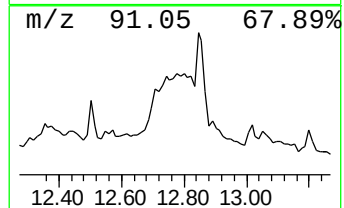
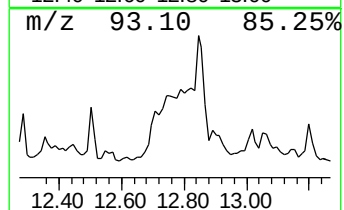
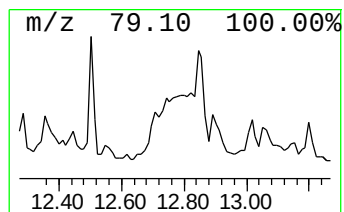
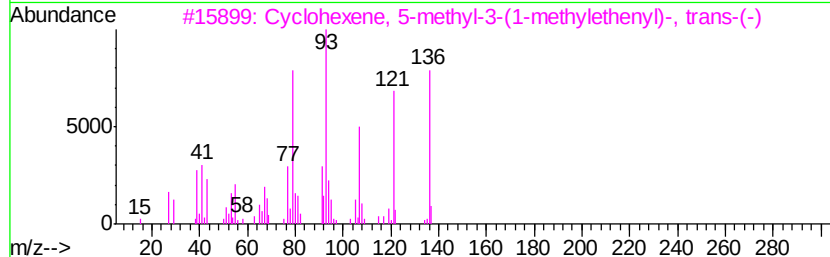
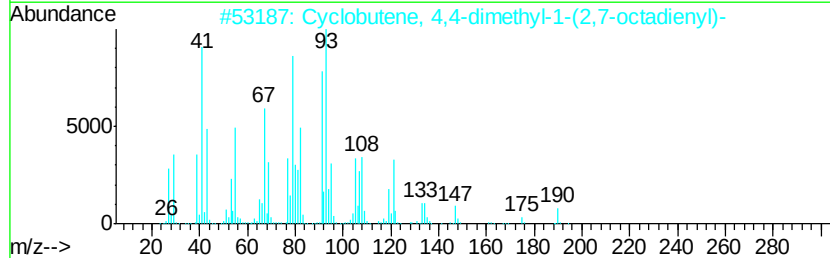
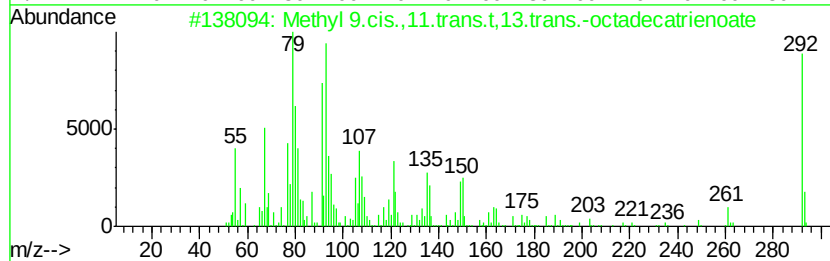
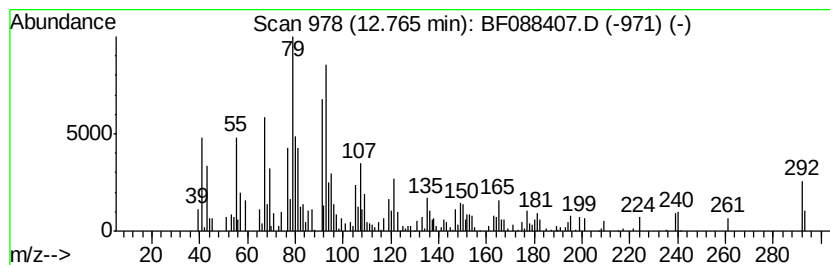
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methyl 9.cis.,11.trans.t,13... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	15.74 ng	365368	Chrysene-d12	13.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 9.cis.,11.trans.t,13.tran...	292	C19H32O2	1000336-42-6	93
2		Cyclobutene, 4,4-dimethyl-1-(2,7...	190	C14H22	062338-42-5	76
3		Cyclohexene, 5-methyl-3-(1-methv...	136	C10H16	056816-08-1	76
4		Methyl 6-cis,9-cis,11-trans-octa...	292	C19H32O2	1000336-37-7	72
5		6-Butyl-1,4-cycloheptadiene	150	C11H18	022735-58-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062516\
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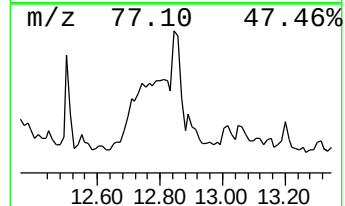
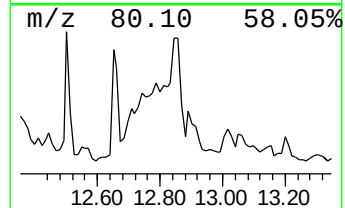
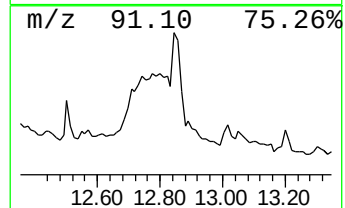
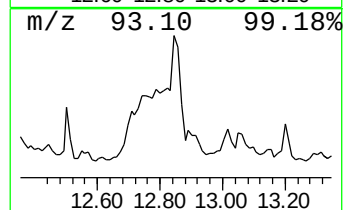
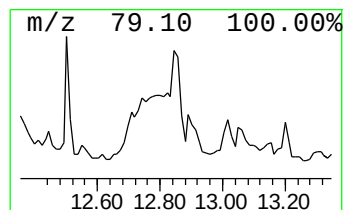
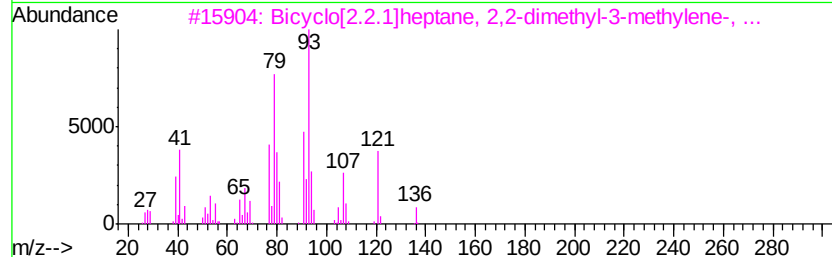
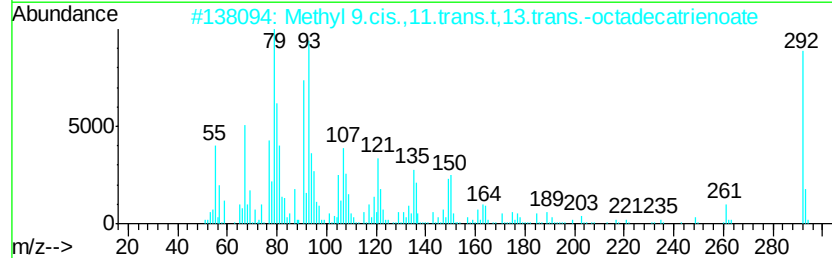
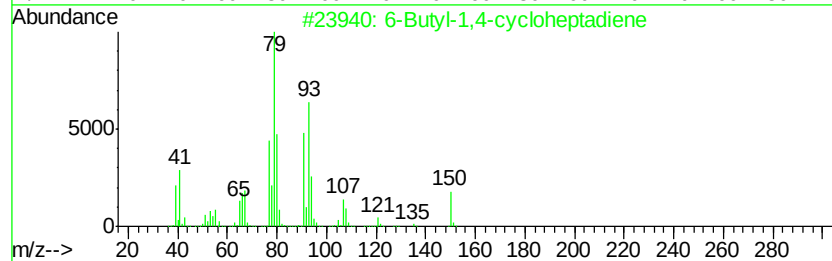
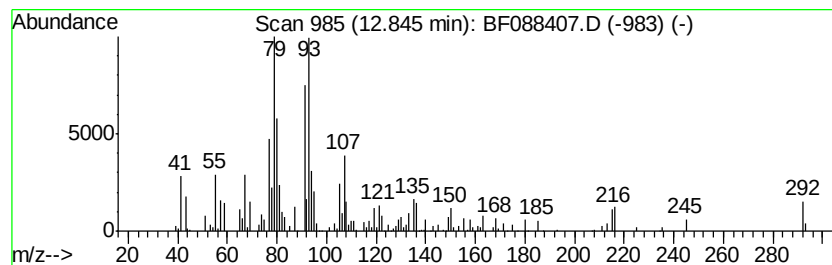
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 6-Butyl-1,4-cycloheptadiene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.85	15.15 ng	351593	Chrysene-d12	13.70	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Butyl-1,4-cycloheptadiene	150	C11H18	022735-58-6	94
2	Methyl 9.cis.,11.trans.t,13.trans...	292	C19H32O2	1000336-42-6	90
3	Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-03-6	83
4	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	016022-04-1	72
5	m-Mentha-4,8-diene, (1S,3S)-(+) -	136	C10H16	005208-51-5	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
Data File : BF088407.D
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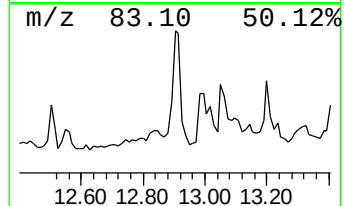
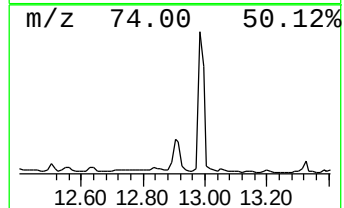
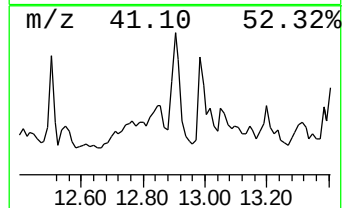
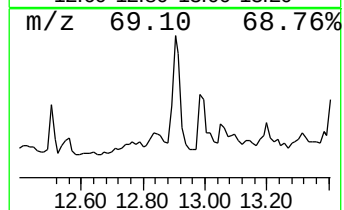
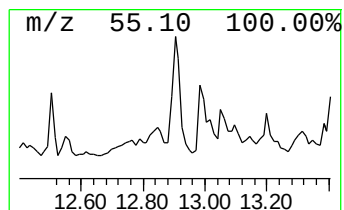
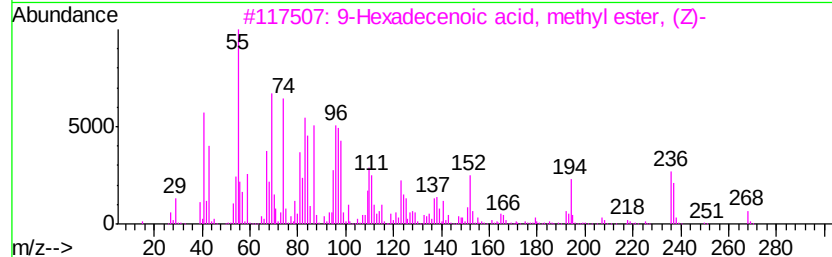
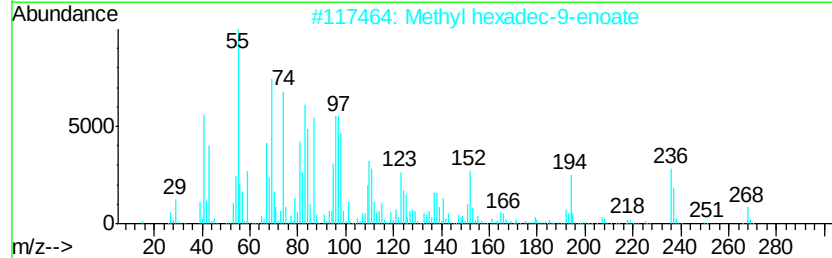
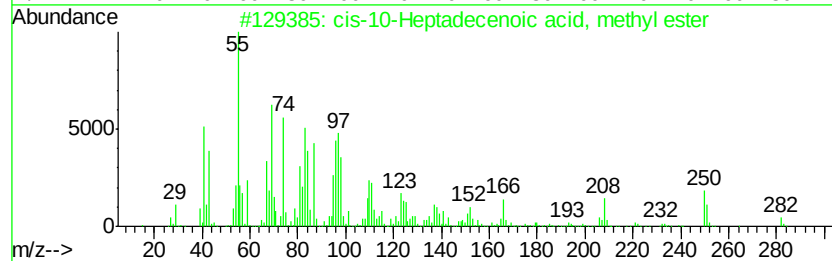
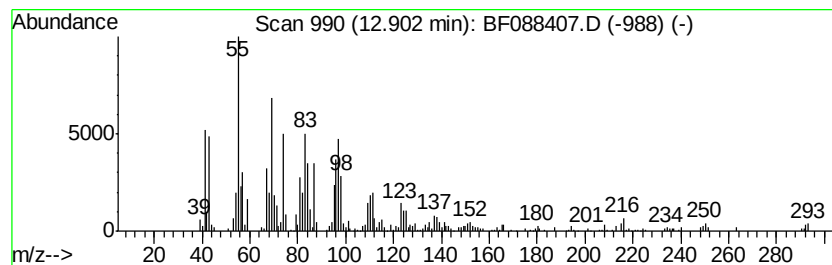
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 cis-10-Heptadecenoic acid, ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	26.71 ng	619898	Chrysene-d12	13.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-10-Heptadecenoic acid, methy...	282	C18H34O2	1000333-62-1	91
2			Methyl hexadec-9-enoate	268	C17H32O2	010030-74-7	90
3			9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	87
4			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	87
5			Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
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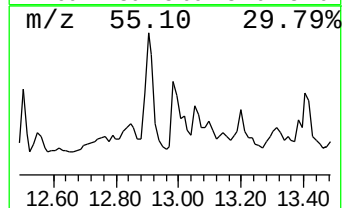
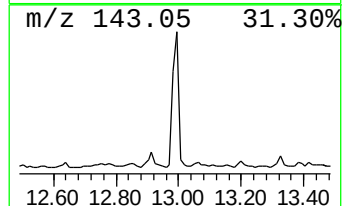
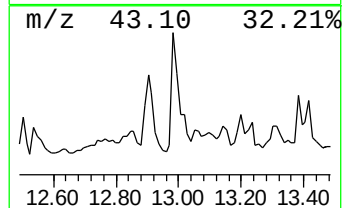
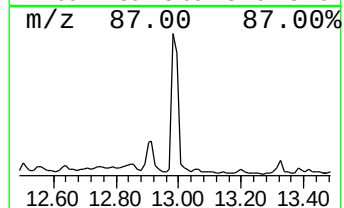
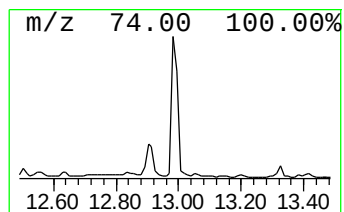
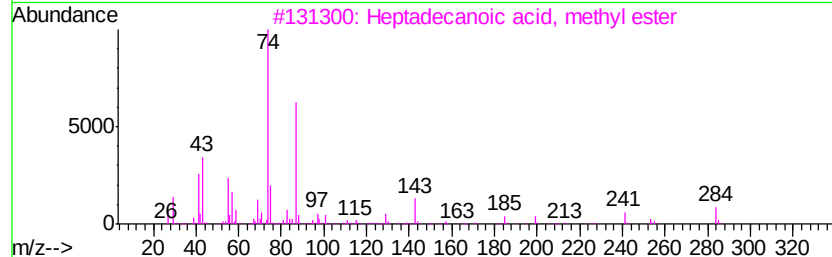
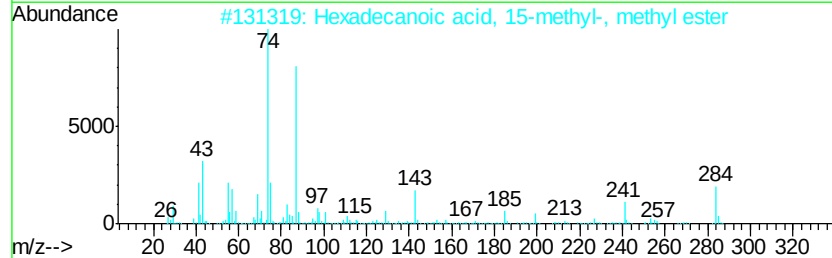
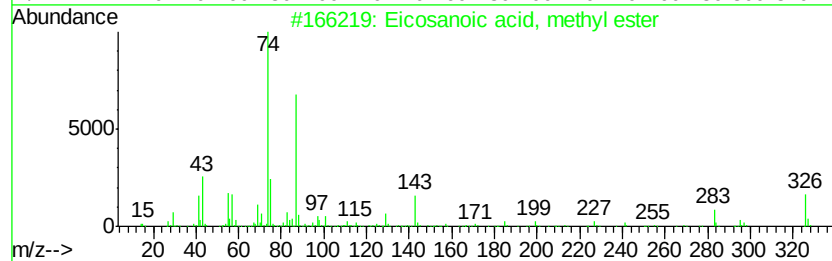
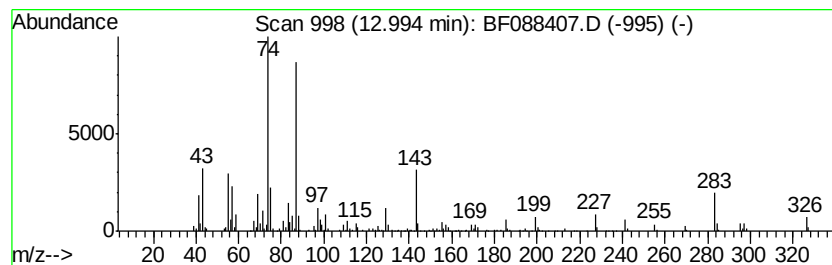
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosanoic acid, methyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	25.56 ng	593082	Chrysene-d12	13.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	97
2		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	96
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	93
4		Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	91
5		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	90



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Sample : H3602-13
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
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ClientSampleId :
Q8-B5(0-16)C

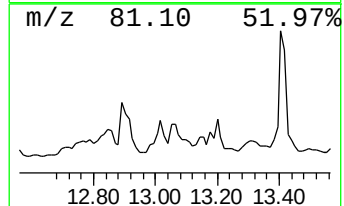
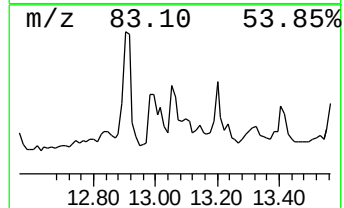
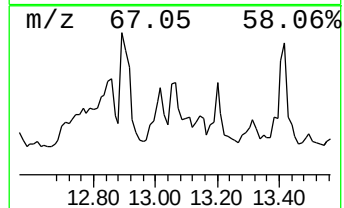
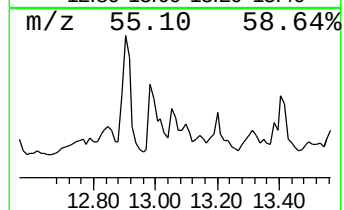
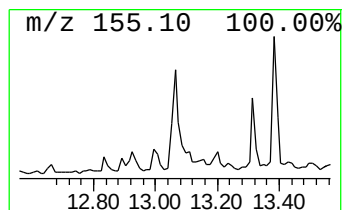
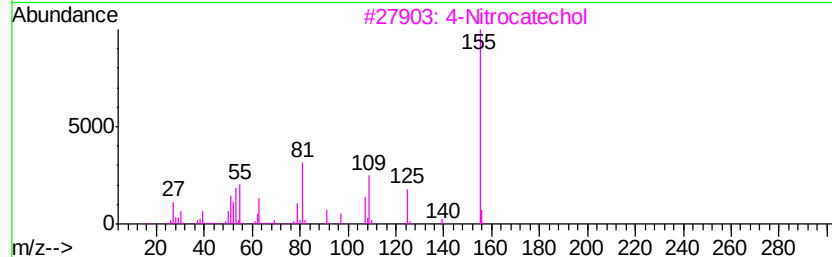
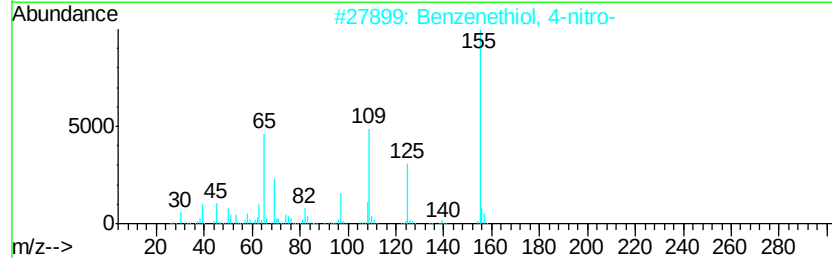
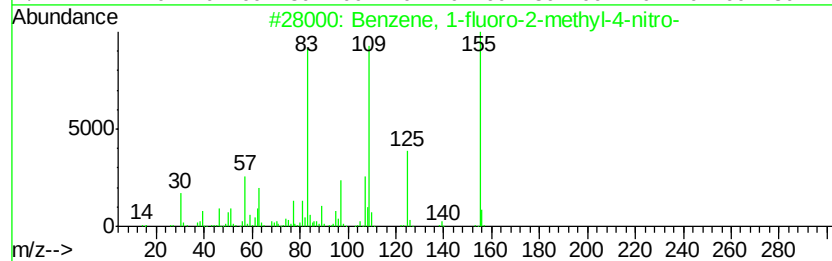
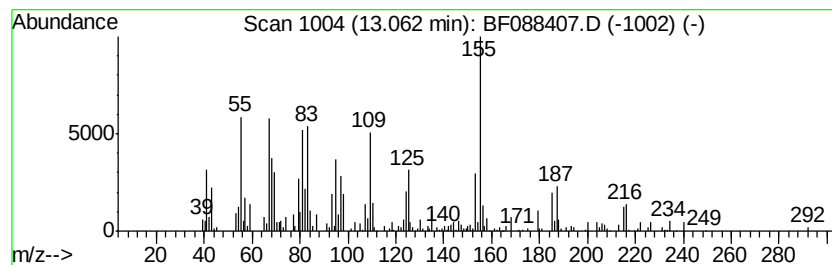
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown13.06 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	10.70 ng	248222	Chrysene-d12	13.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-fluoro-2-methyl-4-nitro-	155	C7H6FN02	000455-88-9	43
2		Benzenethiol, 4-nitro-	155	C6H5N02S	001849-36-1	43
3		4-Nitrocatechol	155	C6H5N04	003316-09-4	43
4		Imidazole-4-carboxylic acid, 2-a...	155	C6H9N3O2	149520-94-5	38
5		Benzene, 2-fluoro-1-methyl-4-nitro-	155	C7H6FN02	001427-07-2	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062516\
Data File : BF088407.D
Acq On : 26 Jun 2016 7:58
Operator : UM/SJ
Sample : H3602-13
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
Q8-B5(0-16)C

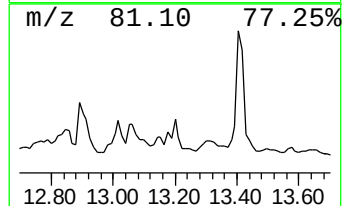
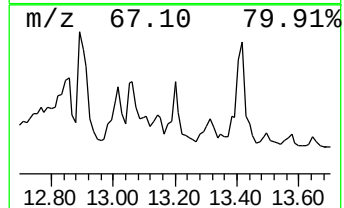
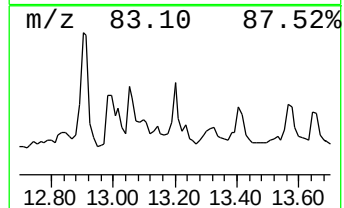
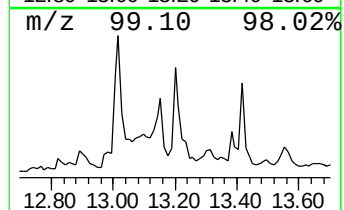
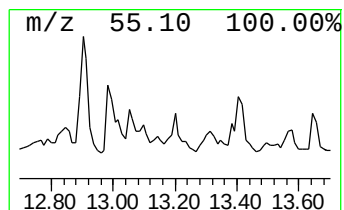
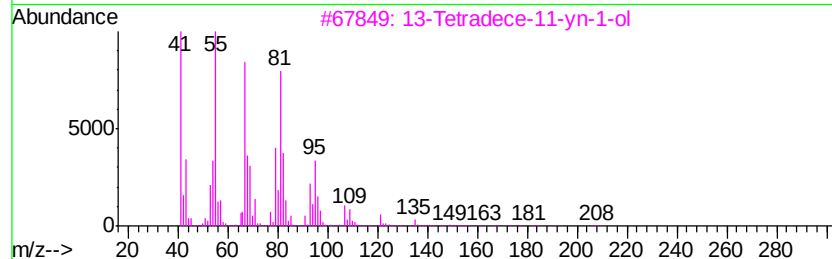
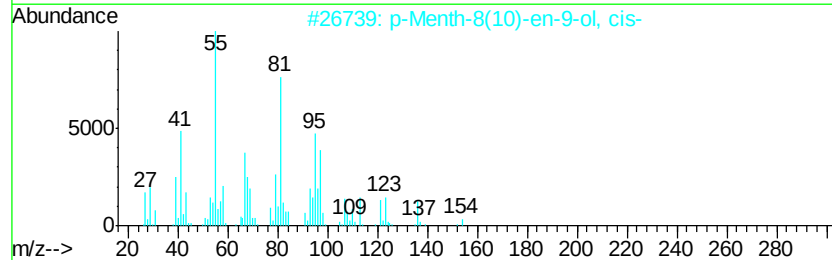
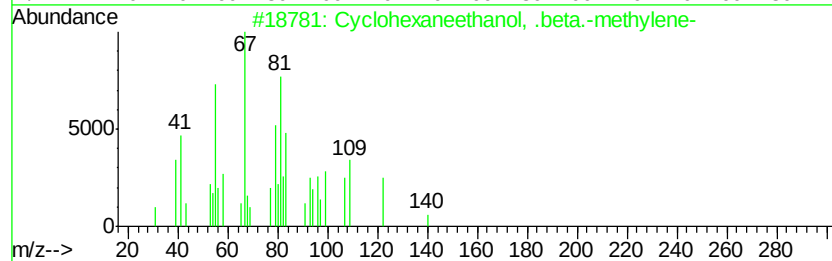
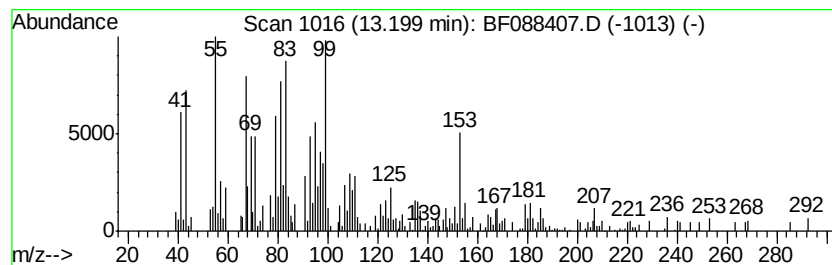
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Cyclohexaneethanol, .beta.-... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	12.03 ng	279077	Chrysene-d12	13.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexaneethanol, .beta.-methy...	140	C9H16O	007697-55-4	56
2		p-Menth-8(10)-en-9-ol, cis-	154	C10H18O	015714-13-3	47
3		13-Tetradec-11-yn-1-ol	208	C14H24O	1000131-00-4	30
4		7-Tetradecyne	194	C14H26	035216-11-6	30
5		Bicyclo[3.1.1]heptane, 2,6,6-tri...	178	C13H22	050746-55-9	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062516\
Data File : BF088407.D
Acq On : 26 Jun 2016 7:58
Operator : UM/SJ
Sample : H3602-13
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
Q8-B5(0-16)C

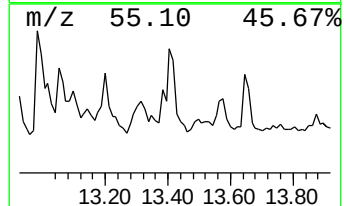
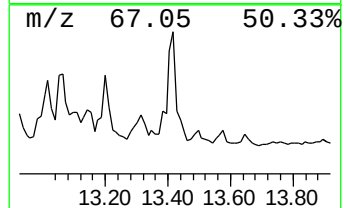
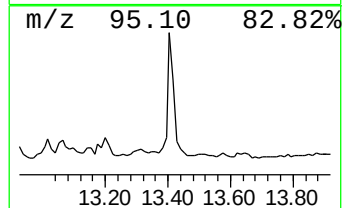
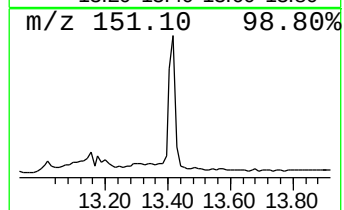
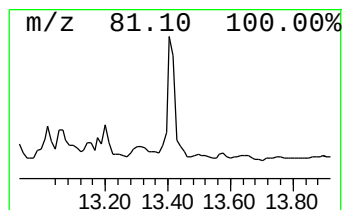
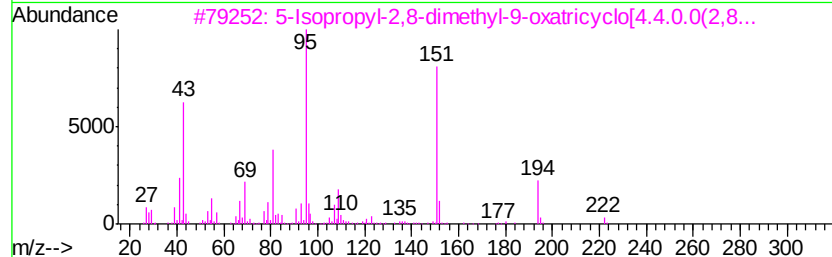
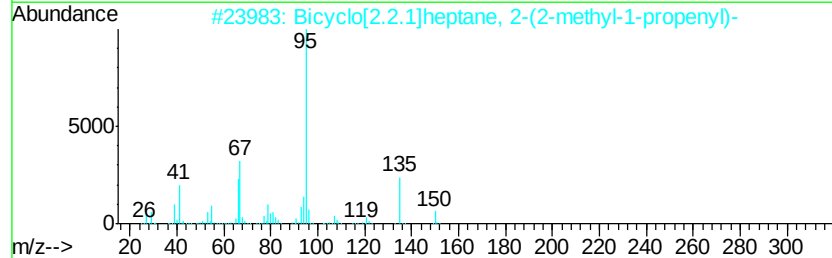
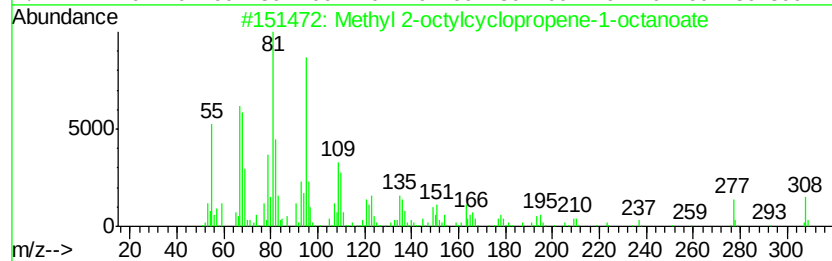
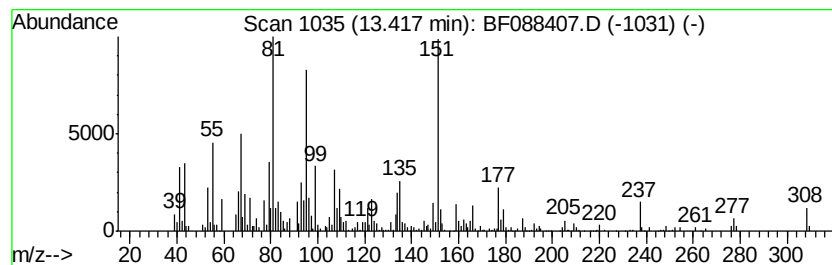
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown13.42 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	25.22 ng	585345	Chrysene-d12	13.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 2-octylcyclopropene-1-oct...	308	C20H36O2	003220-60-8	46
2		Bicyclo[2.2.1]heptane, 2-(2-meth...	150	C11H18	061142-27-6	35
3		5-Isopropyl-2,8-dimethyl-9-oxatr...	222	C14H22O2	1000185-23-8	27
4		2-Cyanomethyl-hexahydro-pyrimidi...	175	C8H9N5	1000185-88-3	15
5		Butyric acid, 2-(3-oxo-3H-benzo[...	237	C11H11NO3S	1000316-00-8	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062516\
 Data File : BF088407.D
 Acq On : 26 Jun 2016 7:58
 Operator : UM/SJ
 Sample : H3602-13
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q8-B5(0-16)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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
unknown6.33	6.33	80.5	ng	1514890	1	6.56	376338	20.0
Methyl tetradecan...	10.62	16.0	ng	397876	4	11.06	497095	20.0
9-Hexadecenoic ac...	11.41	31.9	ng	793087	4	11.06	497095	20.0
Hexadecanoic acid...	11.50	262.7	ng	6529400	4	11.06	497095	20.0
14-Octadecenoic a...	12.23	829.5	ng	20616200	4	11.06	497095	20.0
Methyl stearate	12.29	125.1	ng	3109970	4	11.06	497095	20.0
Methyl 5,12-octad...	12.35	8.8	ng	218483	4	11.06	497095	20.0
Methyl 10-trans,1...	12.50	14.1	ng	326505	5	13.70	464140	20.0
Methyl 9.cis.,11...	12.77	15.7	ng	365368	5	13.70	464140	20.0
6-Butyl-1,4-cyclo...	12.85	15.2	ng	351593	5	13.70	464140	20.0
cis-10-Heptadecen...	12.90	26.7	ng	619898	5	13.70	464140	20.0
Eicosanoic acid, ...	12.99	25.6	ng	593082	5	13.70	464140	20.0
unknown13.06	13.06	10.7	ng	248222	5	13.70	464140	20.0
Cyclohexaneethano...	13.20	12.0	ng	279077	5	13.70	464140	20.0
unknown13.42	13.42	25.2	ng	585345	5	13.70	464140	20.0