

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
 Data File : BF088579.D
 Acq On : 1 Jul 2016 17:17
 Operator : UM/SJ
 Sample : H3808-21 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11-COMP

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-BF063016.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	1.735	12	13	23	rVB	149512	348229	3.75%	0.701%
2	1.873	23	25	88	rVB	3070338	8099475	87.29%	16.294%
3	5.393	331	333	335	rBV	516151	520735	5.61%	1.048%
4	6.433	421	424	426	rBV	382746	523285	5.64%	1.053%
5	6.570	434	436	438	rBV	618958	566604	6.11%	1.140%
6	6.810	454	457	459	rBV	857285	828204	8.93%	1.666%
7	6.959	468	470	473	rVB	464703	419150	4.52%	0.843%
8	7.359	503	505	508	rBV	230105	313002	3.37%	0.630%
9	8.090	567	569	572	rBV	1289656	1121246	12.08%	2.256%
10	9.165	661	663	665	rVB	730232	622235	6.71%	1.252%
11	9.839	720	722	724	rBV2	1044658	1128274	12.16%	2.270%
12	10.628	788	791	793	rVB2	226641	310239	3.34%	0.624%
13	10.856	809	811	814	rVV	202272	187544	2.02%	0.377%
14	11.325	849	852	855	rVB	901442	1349393	14.54%	2.715%
15	11.656	876	881	883	rVB2	254686	437320	4.71%	0.880%
16	11.736	885	888	890	rVB	4839722	5365777	57.83%	10.795%
17	11.931	903	905	909	rVV2	82779	117518	1.27%	0.236%
18	12.045	909	915	917	rVV2	85434	134460	1.45%	0.271%
19	12.137	919	923	925	rVV2	112161	154724	1.67%	0.311%
20	12.422	944	948	949	rBV	5624979	9278615	100.00%	18.667%
21	12.445	949	950	955	rVV2	5778404	7912833	85.28%	15.919%
22	12.525	955	957	959	rVV	2700236	2750727	29.65%	5.534%
23	12.605	959	964	968	rVV4	42938	119790	1.29%	0.241%
24	12.754	974	977	979	rBV	677961	677127	7.30%	1.362%
25	12.902	988	990	993	rVV	576669	556935	6.00%	1.120%
26	12.982	993	997	1001	rVV5	68292	263938	2.84%	0.531%
27	13.108	1003	1008	1010	rVV2	123302	349783	3.77%	0.704%
28	13.165	1010	1013	1018	rVV3	256917	442896	4.77%	0.891%
29	13.257	1018	1021	1024	rVV3	282570	634905	6.84%	1.277%
30	13.302	1024	1025	1028	rVV	310259	359461	3.87%	0.723%
31	13.382	1030	1032	1035	rVV	147962	214972	2.32%	0.432%
32	13.451	1035	1038	1040	rVB2	110338	160948	1.73%	0.324%
33	13.657	1053	1056	1058	rBV2	114442	176105	1.90%	0.354%
34	13.828	1069	1071	1074	rVV2	111761	142180	1.53%	0.286%

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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-BF063016.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.908	1077	1078	1079	rVV	132256	113335	1.22%	0.228%
36	13.942	1079	1081	1087	rVB2	922238	1433920	15.45%	2.885%
37	14.525	1130	1132	1135	rBV2	102849	131118	1.41%	0.264%
38	14.948	1167	1169	1174	rVV	189039	398750	4.30%	0.802%
39	15.234	1191	1194	1196	rVB	92674	123725	1.33%	0.249%
40	15.291	1196	1199	1201	rVB	110597	157045	1.69%	0.316%
41	15.348	1201	1204	1208	rBV	624139	760781	8.20%	1.531%

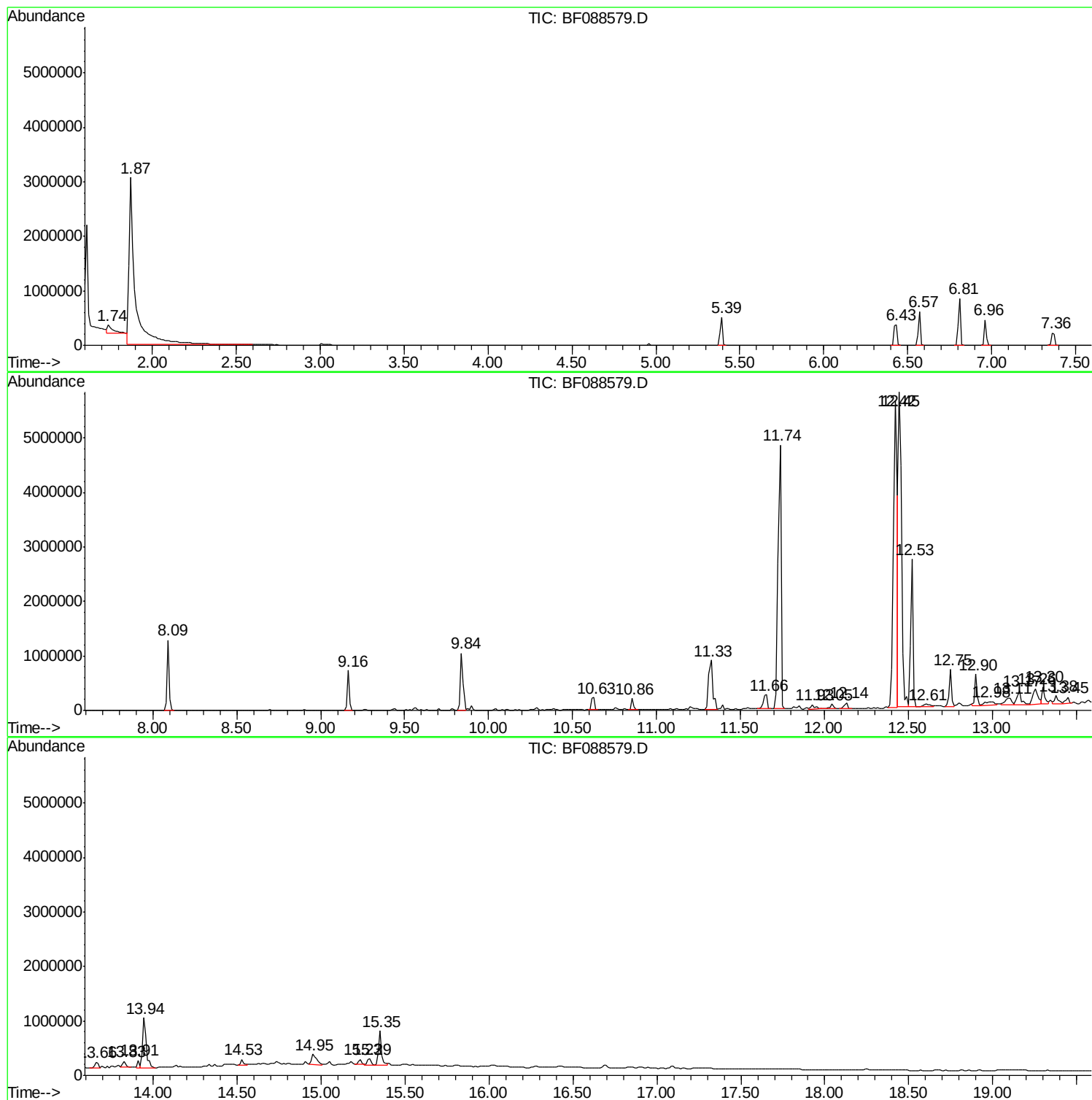
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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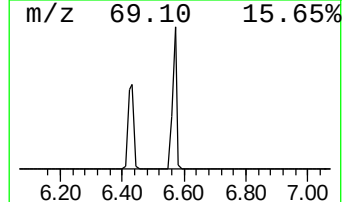
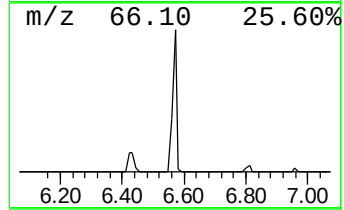
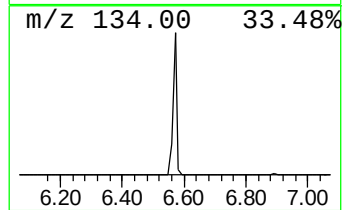
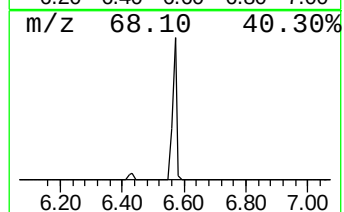
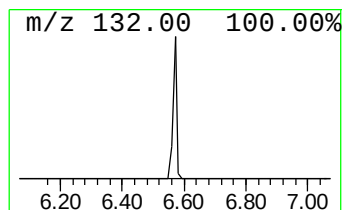
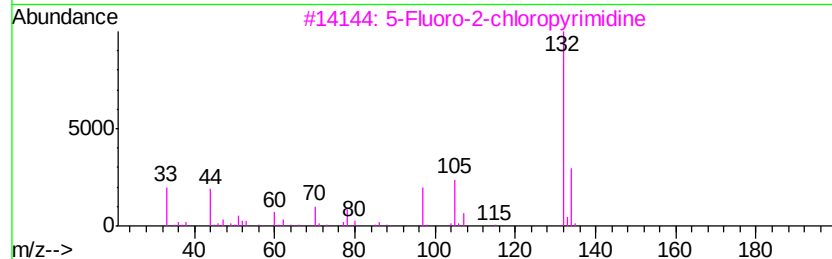
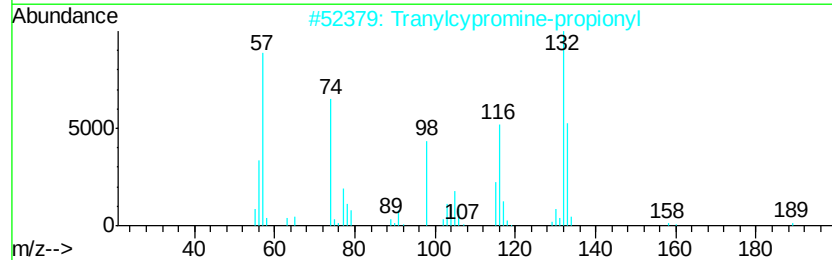
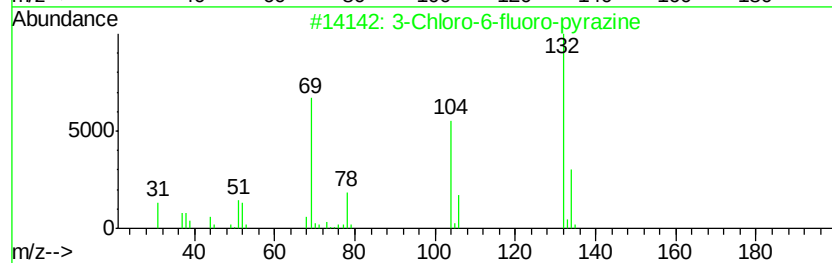
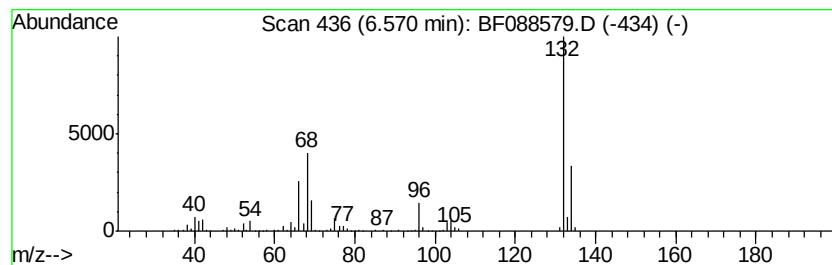
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.57 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	13.68 ng	566604	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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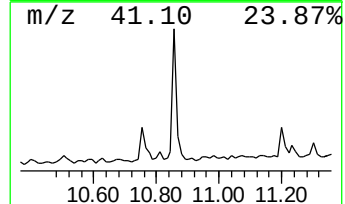
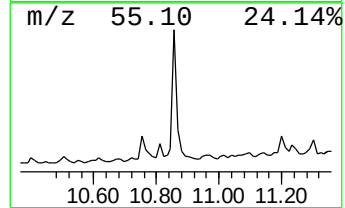
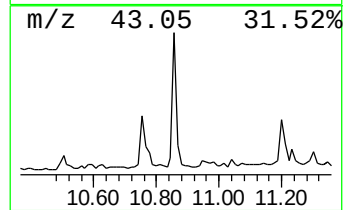
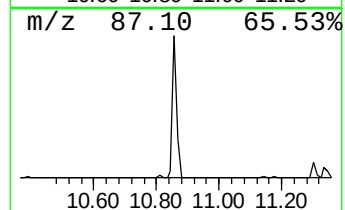
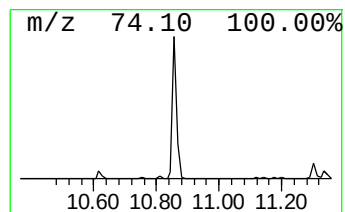
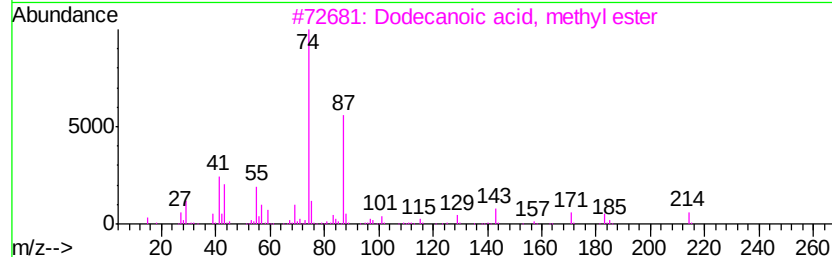
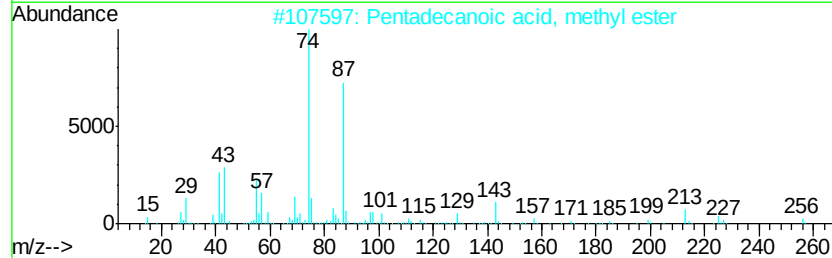
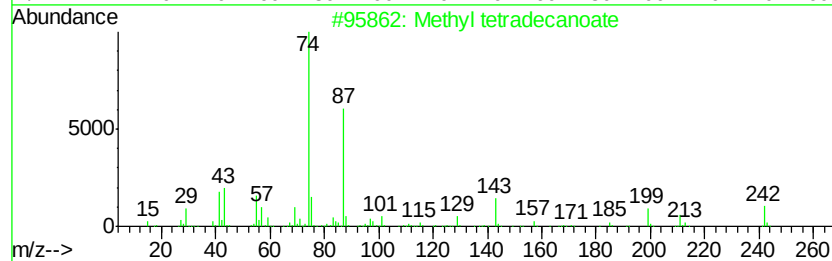
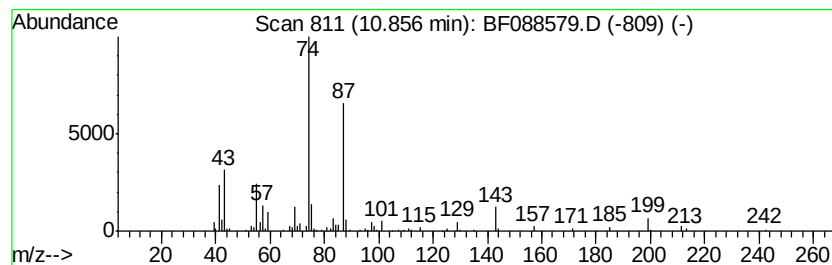
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Methyl tetradecanoate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.86	2.78 ng	187544	Phenanthrene-d10		11.32	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl tetradecanoate	242	C15H30O2	000124-10-7	94
2		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
3		Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	90
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	90
5		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90



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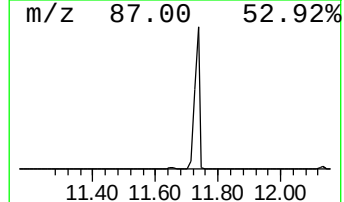
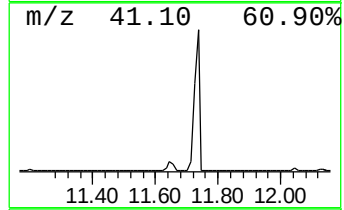
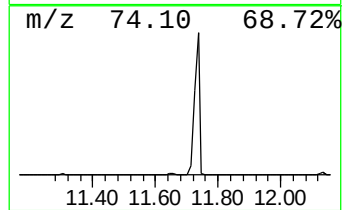
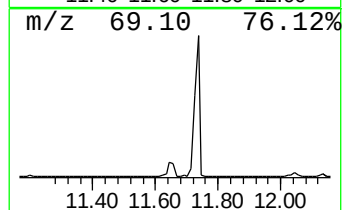
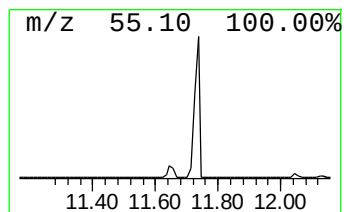
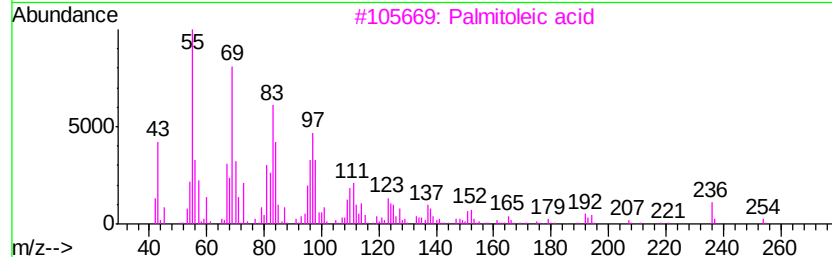
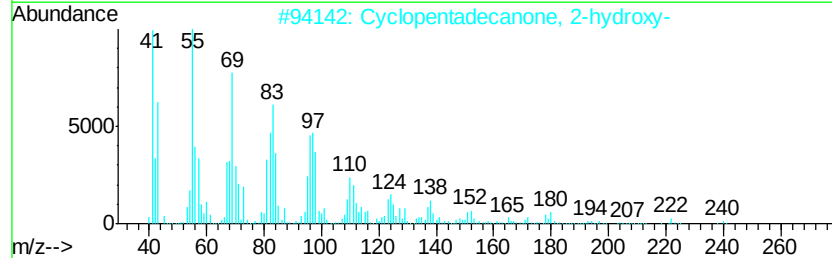
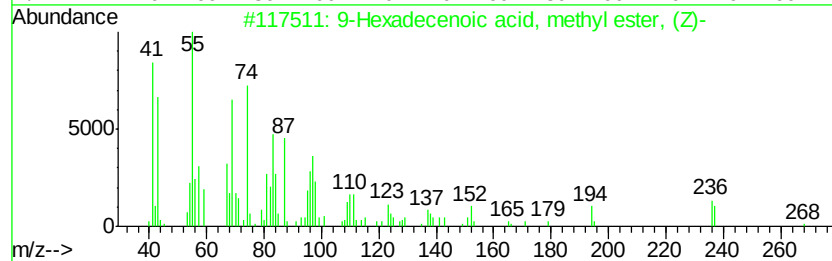
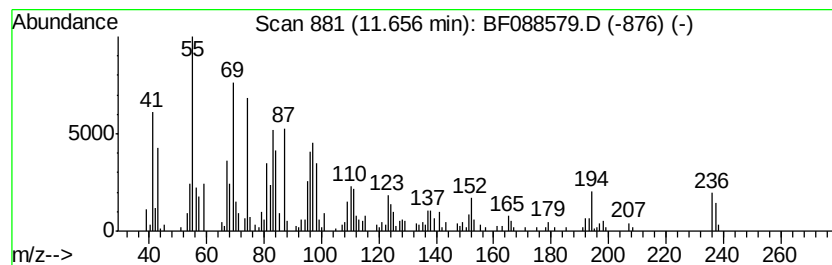
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	6.48 ng	437320	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	70
2		Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	70
3		Palmitoleic acid	254	C16H30O2	000373-49-9	53
4		7-Hexadecenoic acid, methyl este...	268	C17H32O2	056875-67-3	49
5		E-9-Tetradecenoic acid	226	C14H26O2	1000131-35-8	38



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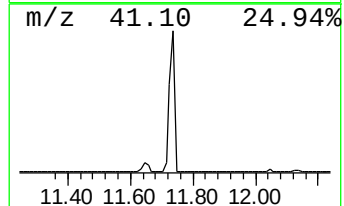
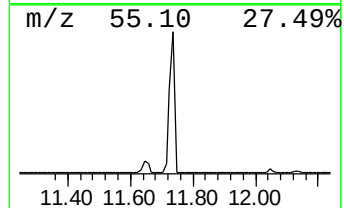
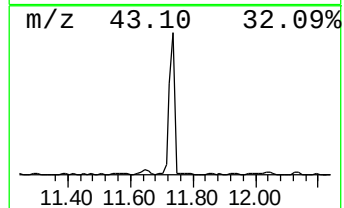
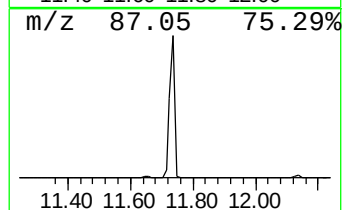
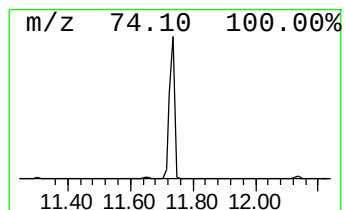
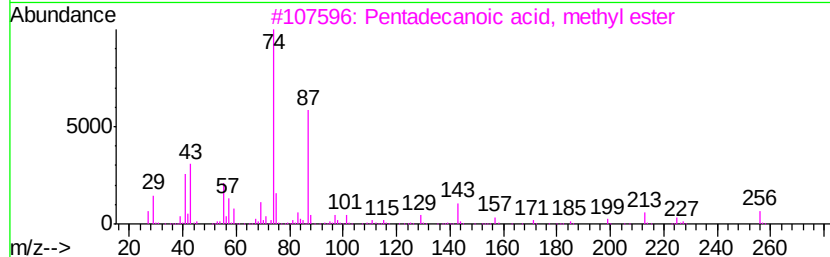
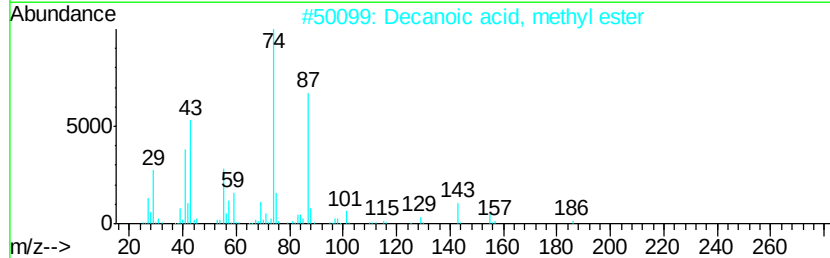
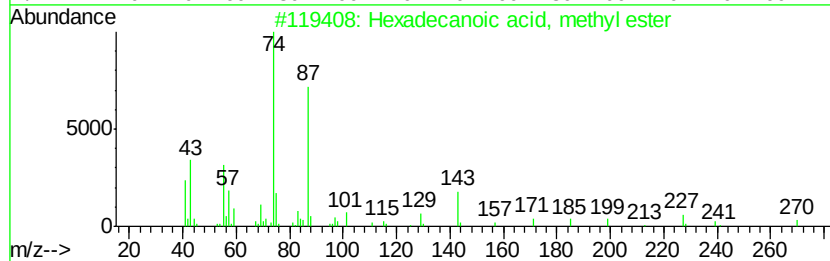
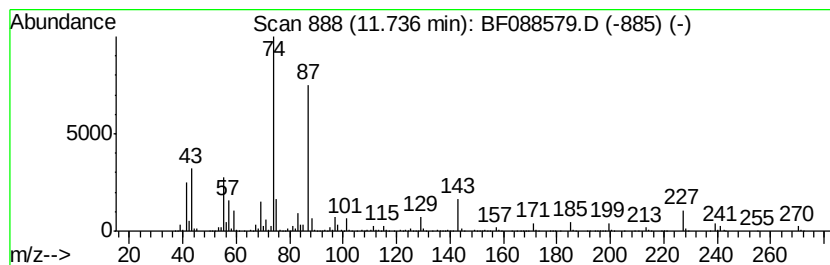
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexadecanoic acid, methyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	79.53 ng	5365780	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4		Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	80
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



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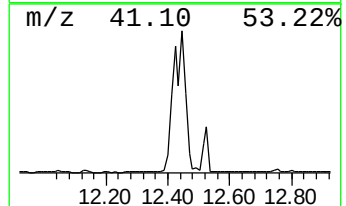
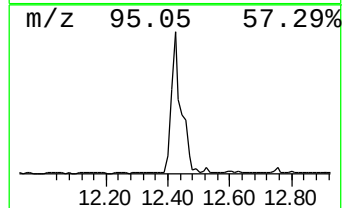
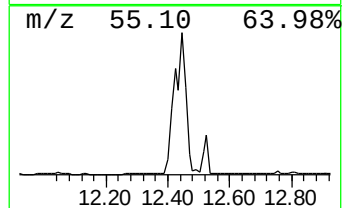
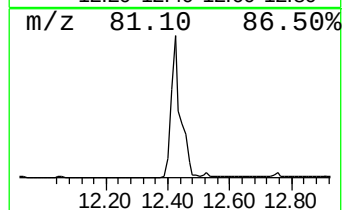
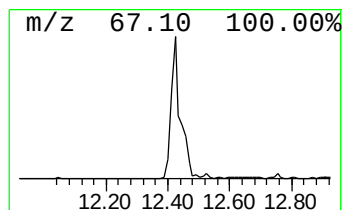
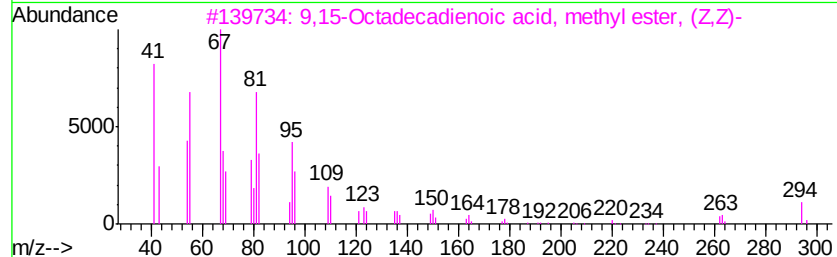
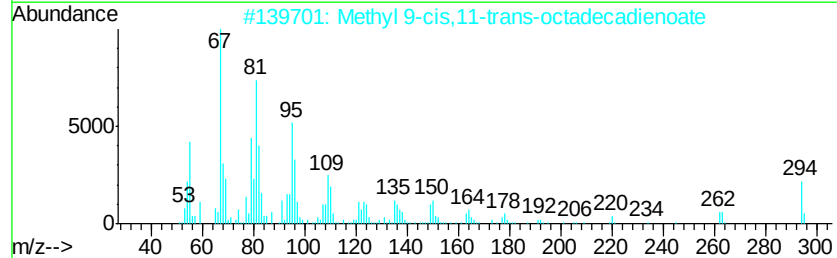
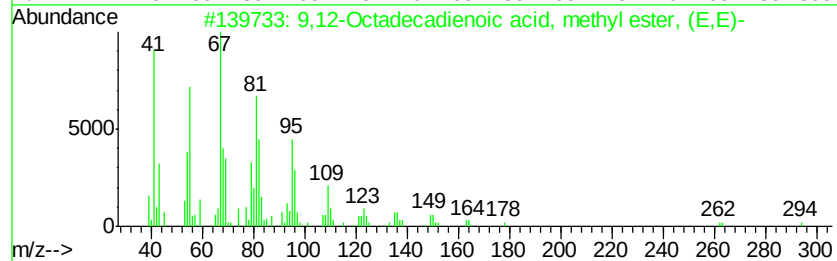
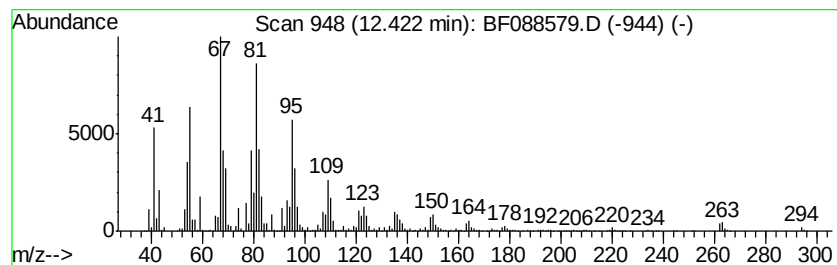
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Peak Number 9 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	137.52 ng	9278620	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99
3		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
Data File : BF088579.D
Acq On : 1 Jul 2016 17:17
Operator : UM/SJ
Sample : H3808-21 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-11-COMP

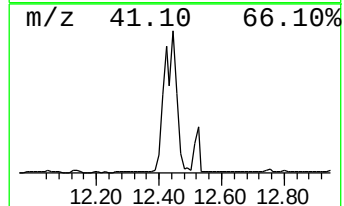
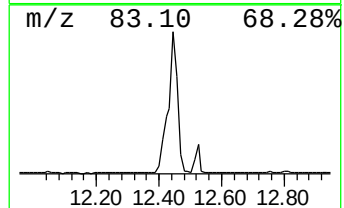
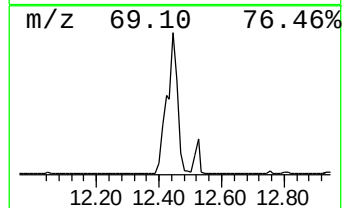
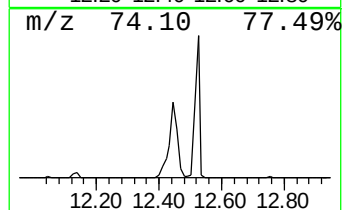
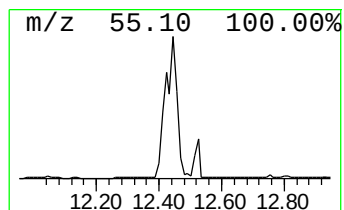
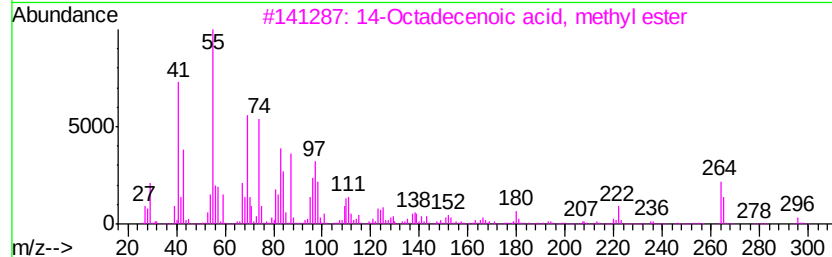
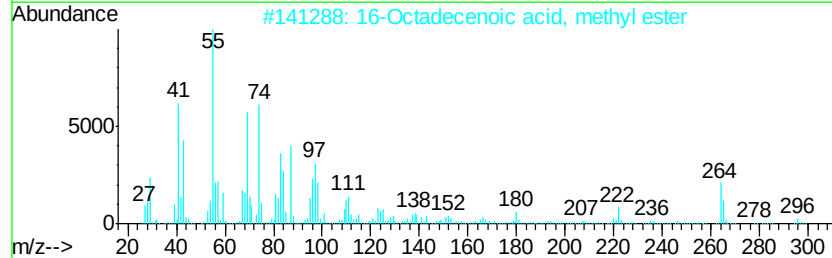
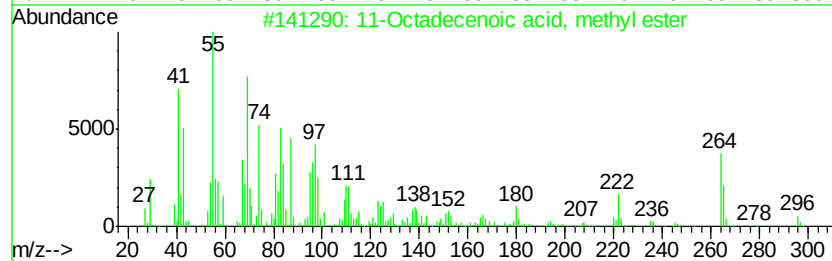
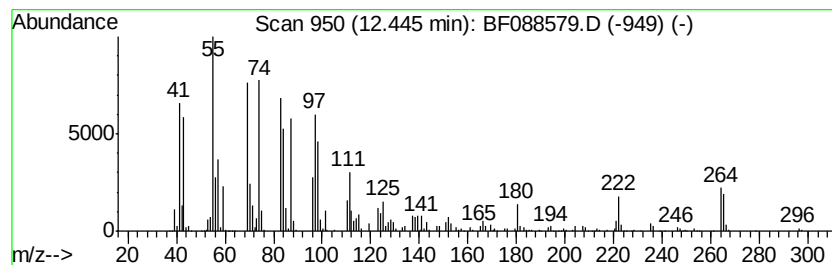
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11-Octadecenoic acid, methyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	117.28 ng	7912830	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	93
2		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	89
3		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	89
4		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	89
5		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
Data File : BF088579.D
Acq On : 1 Jul 2016 17:17
Operator : UM/SJ
Sample : H3808-21 5X
Misc :
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ClientSampleId :
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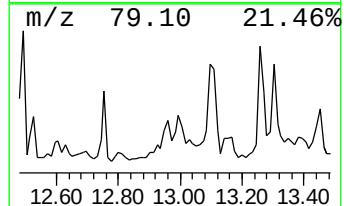
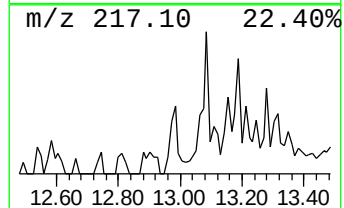
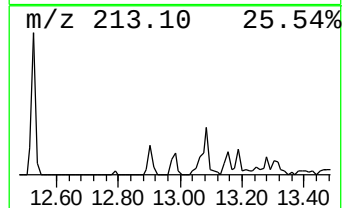
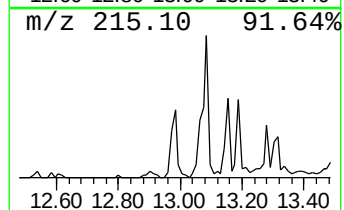
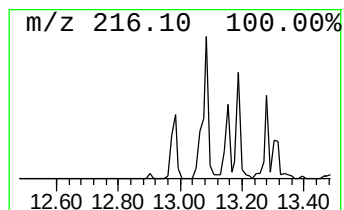
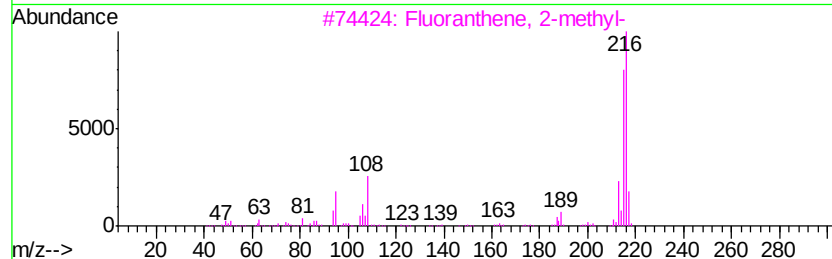
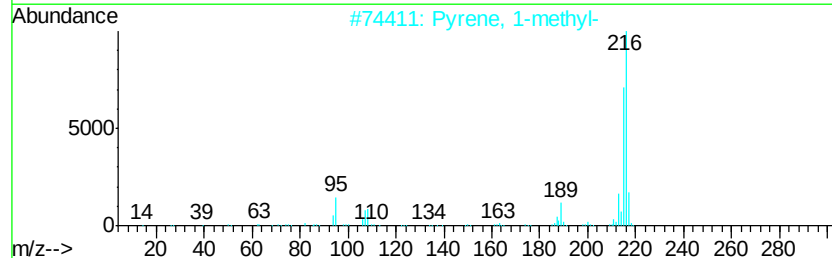
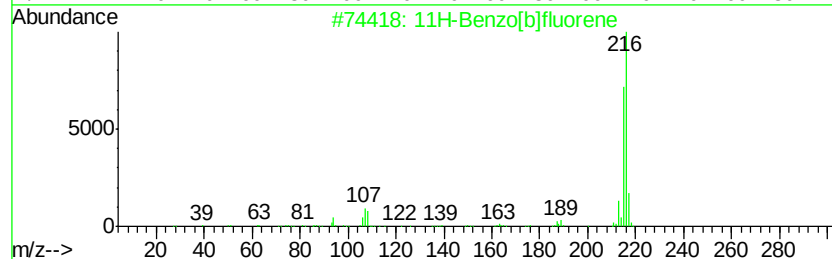
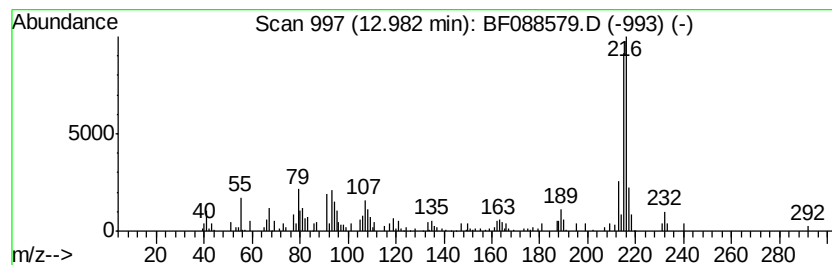
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	3.68 ng	263938	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
Data File : BF088579.D
Acq On : 1 Jul 2016 17:17
Operator : UM/SJ
Sample : H3808-21 5X
Misc :
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ClientSampleId :
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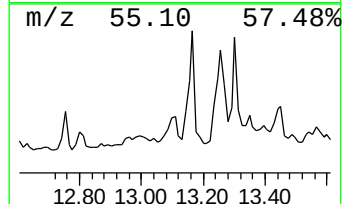
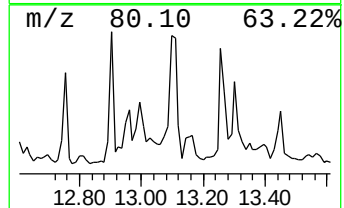
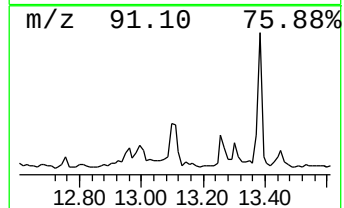
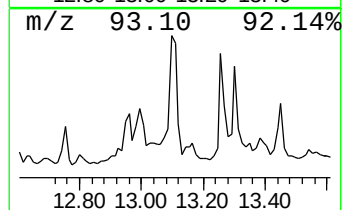
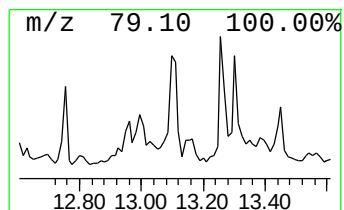
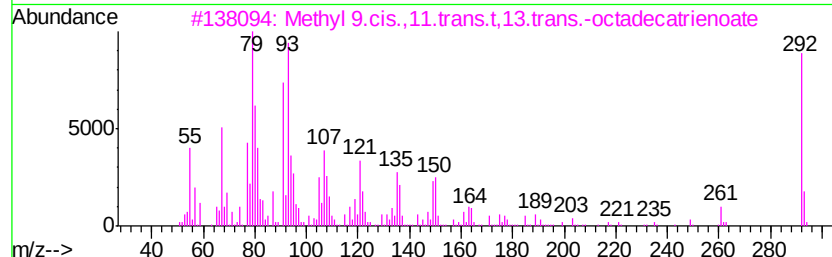
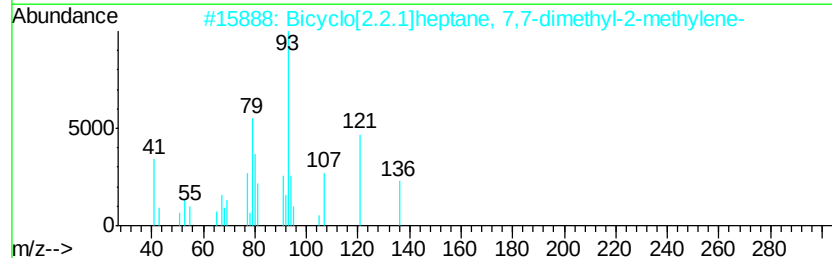
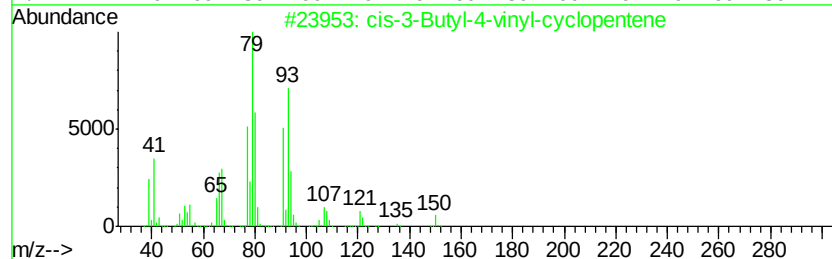
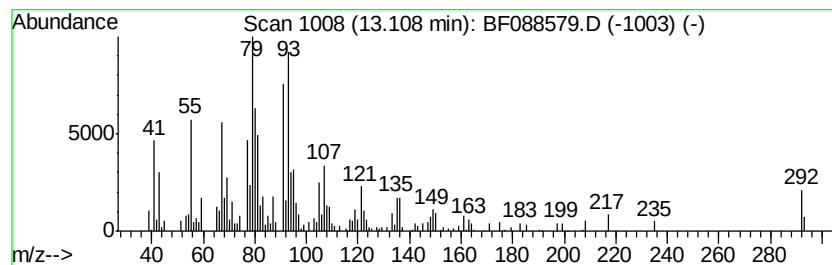
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 cis-3-Butyl-4-vinyl-cyclope... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	4.88 ng	349783	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-3-Butyl-4-vinyl-cyclopentene	150	C11H18	093779-52-3	81
2			Bicyclo[2.2.1]heptane, 7,7-dimet...	136	C10H16	000471-84-1	74
3			Methyl 9.cis.,11.trans.t,13.tran...	292	C19H32O2	1000336-42-6	64
4			Methyl 6-cis,9-cis,11-trans-octa...	292	C19H32O2	1000336-37-7	64
5			Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-03-6	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
Data File : BF088579.D
Acq On : 1 Jul 2016 17:17
Operator : UM/SJ
Sample : H3808-21 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

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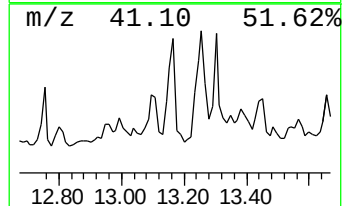
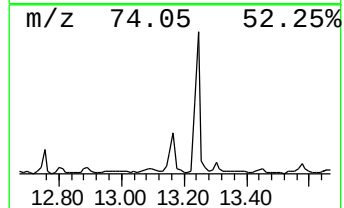
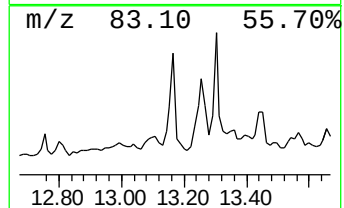
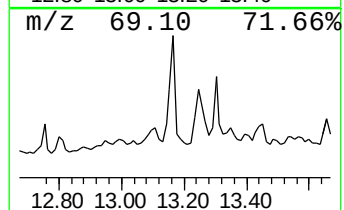
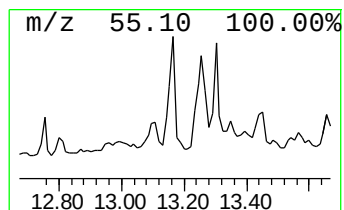
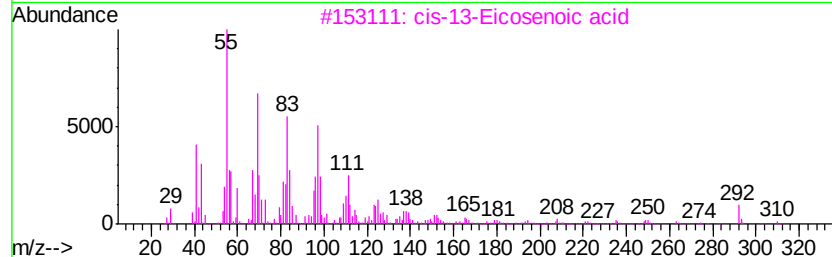
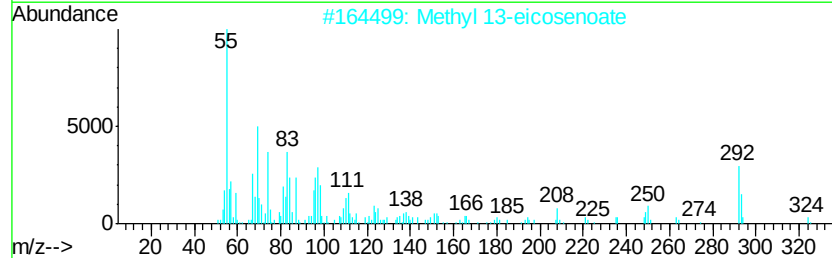
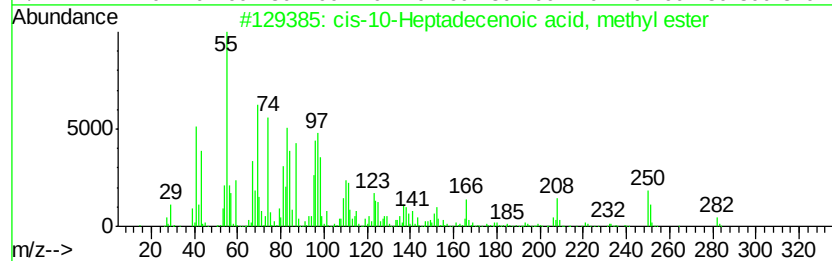
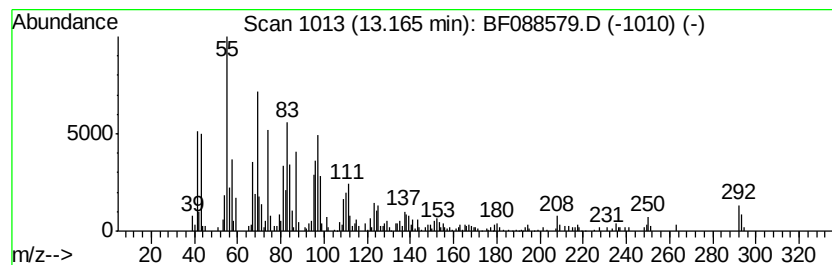
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.17	6.18 ng	442896	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-10-Heptadecenoic acid, methy...	282	C18H34O2	1000333-62-1	91
2			Methyl 13-eicosenoate	324	C21H40O2	1000336-48-4	90
3			cis-13-Eicosenoic acid	310	C20H38O2	017735-94-3	74
4			cis-Vaccenic acid	282	C18H34O2	000506-17-2	72
5			cis-9-Hexadecenoic acid	254	C16H30O2	1000333-19-5	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
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Operator : UM/SJ
Sample : H3808-21 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

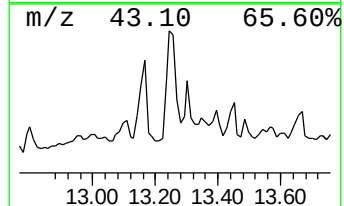
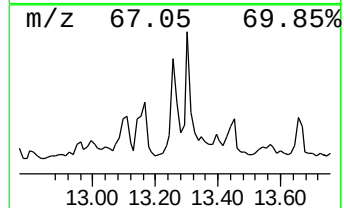
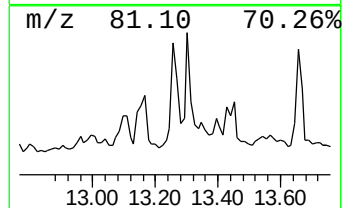
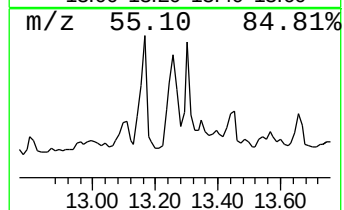
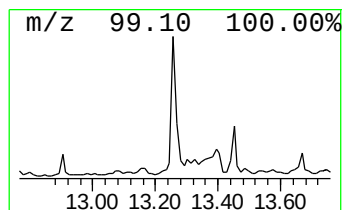
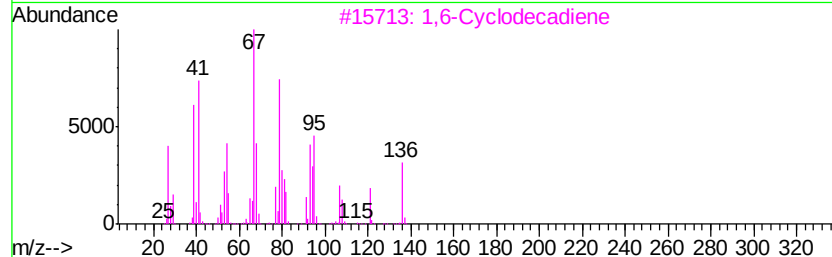
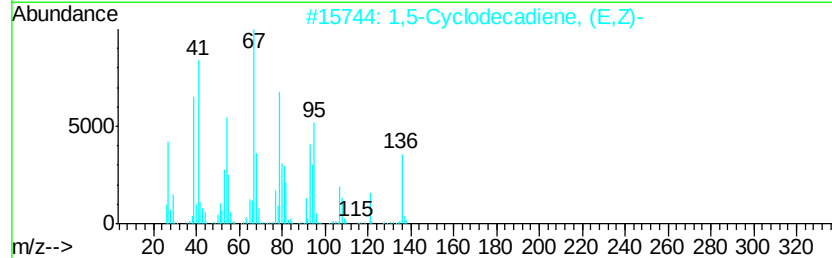
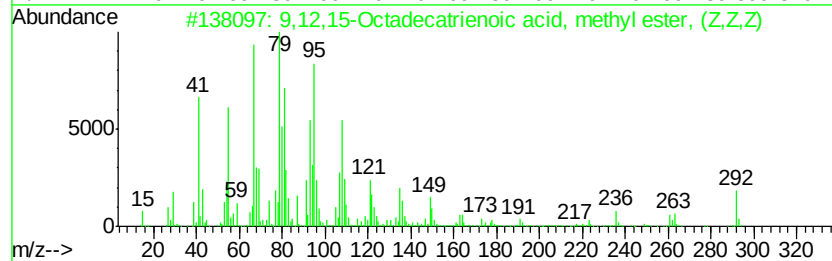
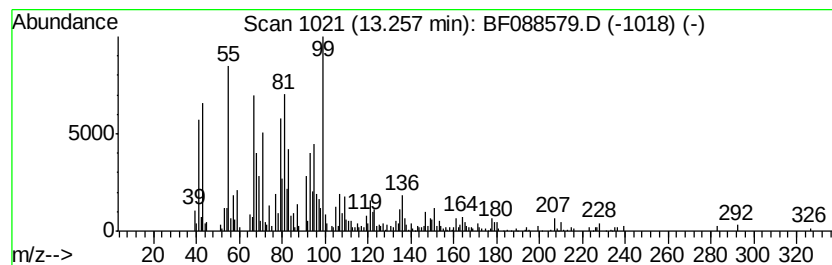
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 9,12,15-Octadecatrienoic ac... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.26	8.86 ng	634905	Chrysene-d12	13.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12,15-Octadecatrienoic acid, m...	292	C19H32O2	000301-00-8	64
2	1,5-Cyclodecadiene, (E,Z)-	136	C10H16	001124-78-3	55
3	1,6-Cyclodecadiene	136	C10H16	007049-13-0	43
4	9-Octadecen-12-ynoic acid, methy...	292	C19H32O2	056847-05-3	42
5	1,13-Tetradecadiene	194	C14H26	021964-49-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
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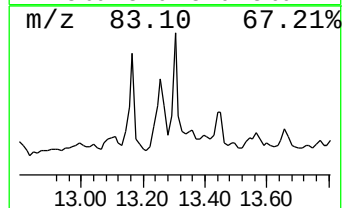
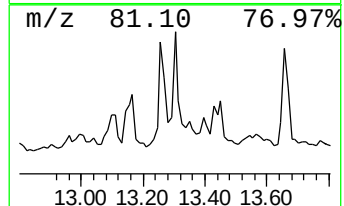
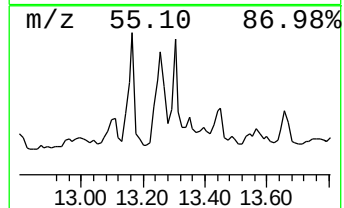
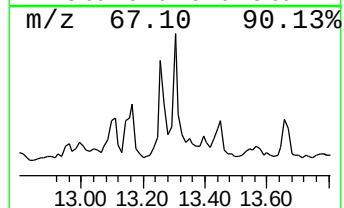
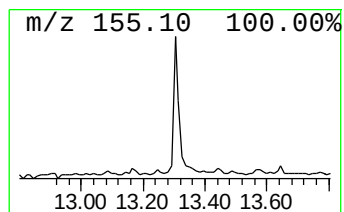
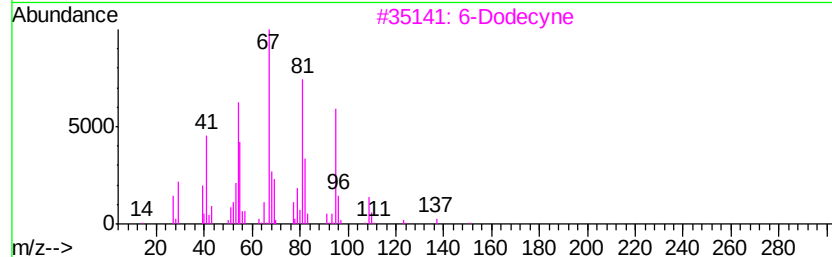
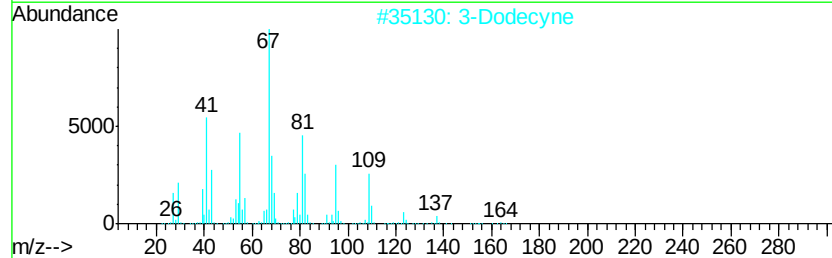
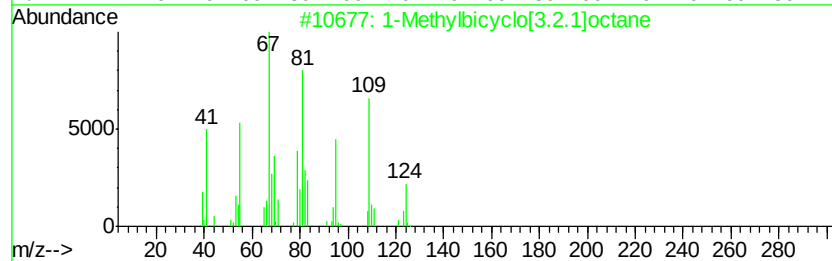
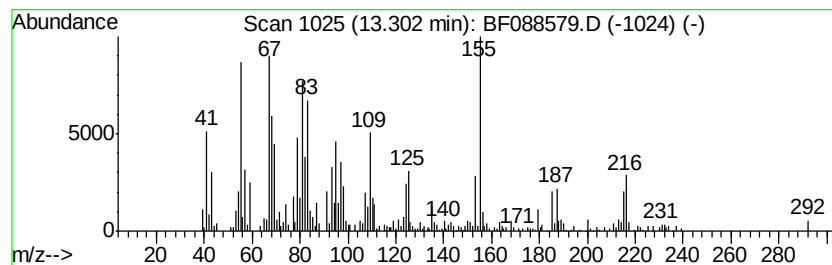
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Methylbicyclo[3.2.1]octane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.30	5.01 ng	359461	Chrysene-d12		13.95		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	55
2			3-Dodecyne	166	C12H22	006790-27-8	43
3			6-Dodecyne	166	C12H22	006975-99-1	38
4			5,10-Pentadecadien-1-ol, (Z,Z)-	224	C15H28O	064275-51-0	35
5			4-Fluoro-3-nitrotoluene	155	C7H6FN02	000446-11-7	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
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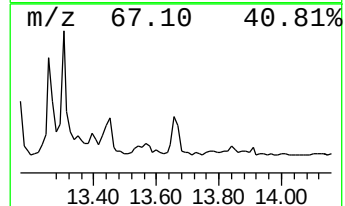
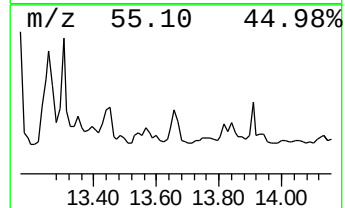
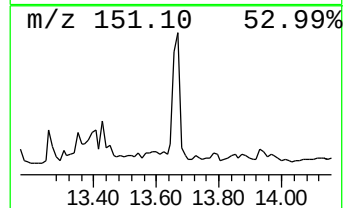
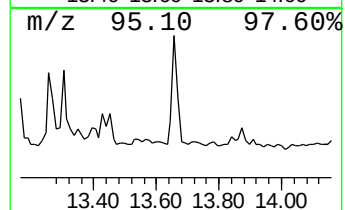
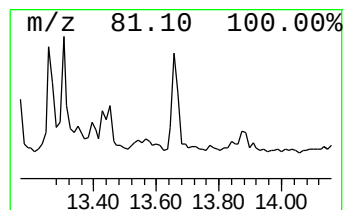
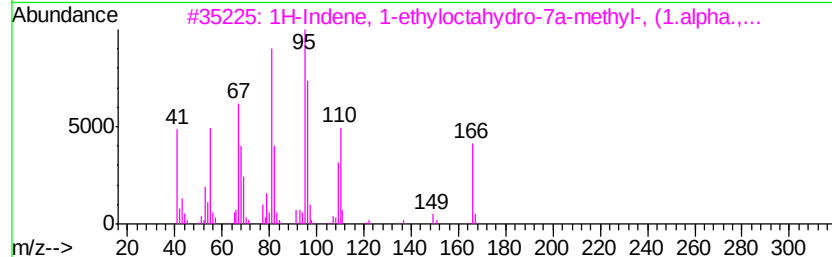
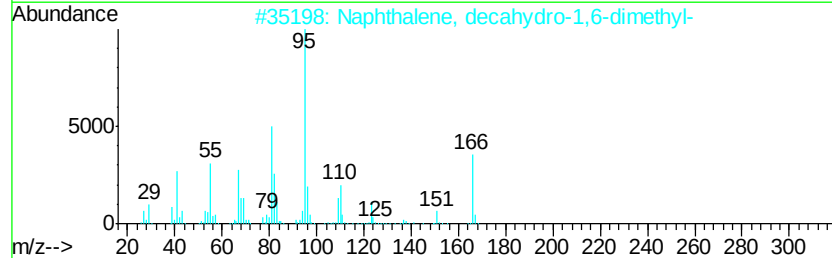
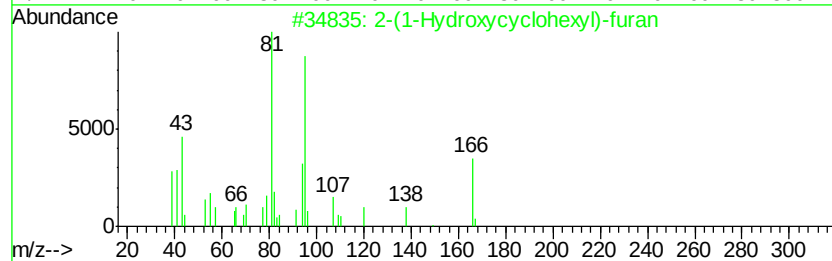
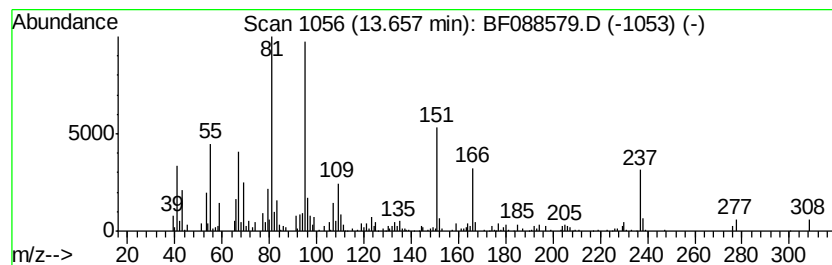
Instrument :
BNA_F
ClientSampleId :
SB-11-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown13.66 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.66	2.46 ng	176105	Chrysene-d12	13.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(1-Hydroxycyclohexyl)-furan	166	C10H14O2	036169-67-2	46
2	Naphthalene, decahydro-1,6-dimet...	166	C12H22	001750-51-2	43
3	1H-Indene, 1-ethyloctahydro-7a-m...	166	C12H22	056324-71-1	43
4	Cyclopentene, 1,4-dimethyl-5-(1-...	138	C10H18	061142-33-4	27
5	1-(1,2-Dimethyl-cyclopent-2-enyl...	138	C9H14O	070987-82-5	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
Data File : BF088579.D
Acq On : 1 Jul 2016 17:17
Operator : UM/SJ
Sample : H3808-21 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-11-COMP

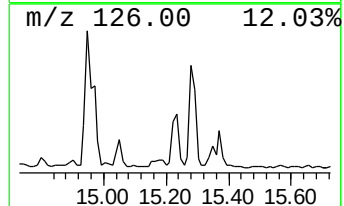
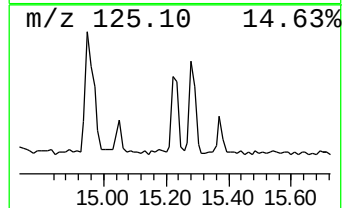
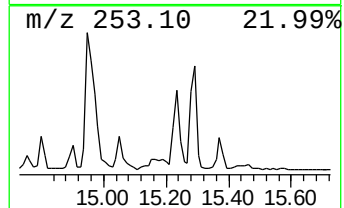
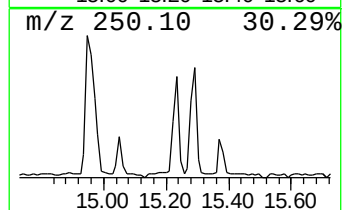
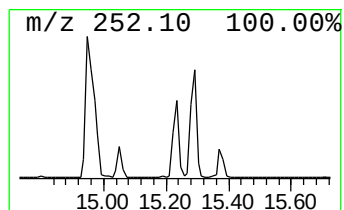
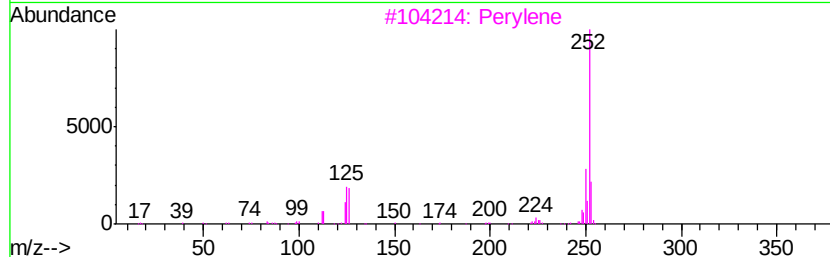
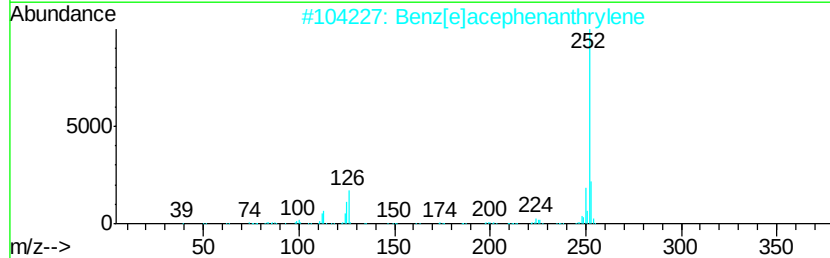
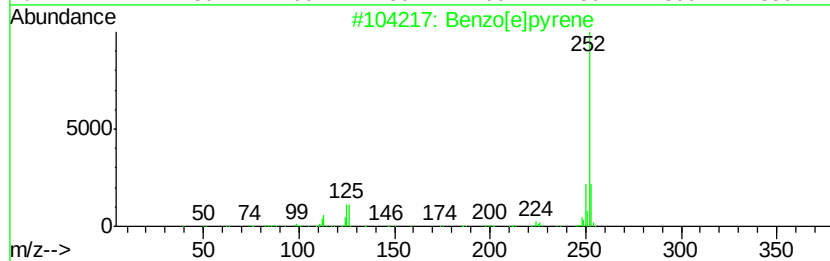
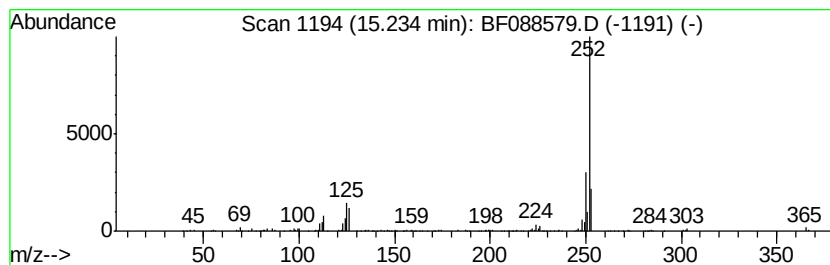
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[e]pyrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.23	3.25 ng	123725	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
3		Perylene	252	C20H12	000198-55-0	96
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
 Data File : BF088579.D
 Acq On : 1 Jul 2016 17:17
 Operator : UM/SJ
 Sample : H3808-21 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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
unknown6.57	6.57	13.7	ng	566604	1	6.81	828204	20.0
Methyl tetradecan...	10.86	2.8	ng	187544	4	11.32	1349390	20.0
9-Hexadecenoic ac...	11.66	6.5	ng	437320	4	11.32	1349390	20.0
Hexadecanoic acid...	11.74	79.5	ng	5365780	4	11.32	1349390	20.0
9,12-Octadecadien...	12.42	137.5	ng	9278620	4	11.32	1349390	20.0
11-Octadecenoic a...	12.45	117.3	ng	7912830	4	11.32	1349390	20.0
11H-Benzo[b]fluorene	12.98	3.7	ng	263938	5	13.95	1433920	20.0
cis-3-Butyl-4-vin...	13.11	4.9	ng	349783	5	13.95	1433920	20.0
cis-10-Heptadecen...	13.17	6.2	ng	442896	5	13.95	1433920	20.0
9,12,15-Octadecat...	13.26	8.9	ng	634905	5	13.95	1433920	20.0
1-Methylbicyclo[3...	13.30	5.0	ng	359461	5	13.95	1433920	20.0
unknown13.66	13.66	2.5	ng	176105	5	13.95	1433920	20.0
Benzo[e]pyrene	15.23	3.3	ng	123725	6	15.35	760781	20.0