

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
 Data File : BF104761.D
 Acq On : 24 Apr 2018 17:27
 Operator : JU/SJ
 Sample : J2151-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-042318-A

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-BF040618.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	5.428	564	568	580	rBV	2025084	1901061	20.02%	4.460%
2	6.451	737	742	745	rBV	1946996	1886201	19.86%	4.425%
3	6.580	760	764	767	rBV	2340763	2008227	21.14%	4.711%
4	6.804	799	802	806	rBV	708817	603385	6.35%	1.416%
5	6.963	825	829	832	rBV	1980267	1609997	16.95%	3.777%
6	7.375	894	899	902	rBV	1404295	1196256	12.59%	2.806%
7	7.969	994	1000	1008	rBV	56881	97174	1.02%	0.228%
8	8.092	1018	1021	1026	rBV	946719	772285	8.13%	1.812%
9	9.169	1200	1204	1207	rBV	2716838	2317133	24.40%	5.436%
10	9.563	1266	1271	1274	rBV	230194	237204	2.50%	0.556%
11	9.851	1316	1320	1323	rBV	890561	810462	8.53%	1.901%
12	10.245	1383	1387	1389	rBV	108241	118411	1.25%	0.278%
13	10.268	1389	1391	1394	rVB	154010	115713	1.22%	0.271%
14	10.639	1451	1454	1463	rVB	1227661	1206097	12.70%	2.829%
15	10.745	1469	1472	1481	rVB	141661	203248	2.14%	0.477%
16	11.192	1545	1548	1551	rBV	134573	112613	1.19%	0.264%
17	11.339	1569	1573	1576	rBV	856639	741403	7.81%	1.739%
18	11.727	1634	1639	1642	rBV	4205201	4017156	42.29%	9.424%
19	11.868	1659	1663	1669	rVB2	136927	140076	1.47%	0.329%
20	12.427	1751	1758	1760	rBV	5316771	9498119	100.00%	22.282%
21	12.451	1760	1762	1770	rVV2	5535772	6699373	70.53%	15.716%
22	12.521	1770	1774	1777	rVV	2564012	2264181	23.84%	5.312%
23	12.557	1778	1780	1785	rVB	143003	167034	1.76%	0.392%
24	12.927	1839	1843	1847	rVB	1955966	1763083	18.56%	4.136%
25	13.162	1881	1883	1890	rVB4	157557	188630	1.99%	0.443%
26	13.239	1894	1896	1901	rVB	209361	202940	2.14%	0.476%
27	13.656	1964	1967	1968	rBV2	89620	96682	1.02%	0.227%
28	13.815	1991	1994	2000	rVB2	227487	233932	2.46%	0.549%
29	13.915	2007	2011	2014	rBV	234110	245950	2.59%	0.577%
30	13.986	2019	2023	2026	rBV	692931	694424	7.31%	1.629%
31	15.462	2271	2274	2280	rVB	384885	478175	5.03%	1.122%

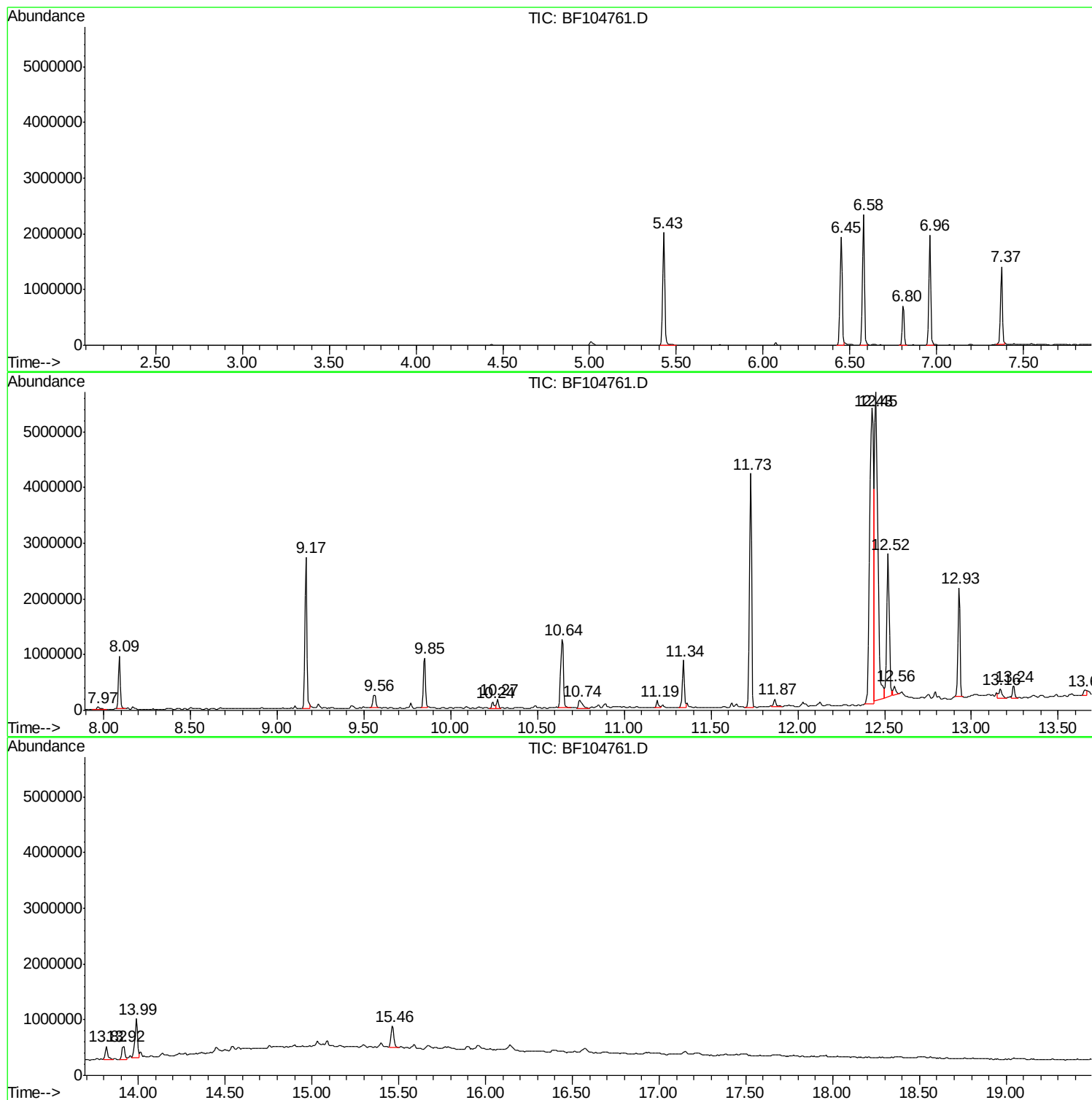
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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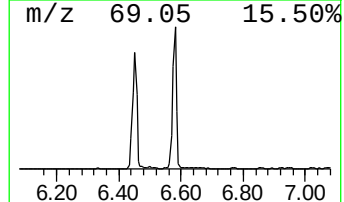
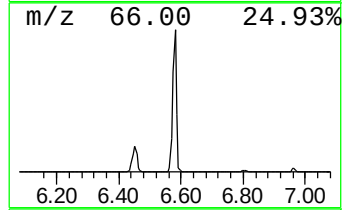
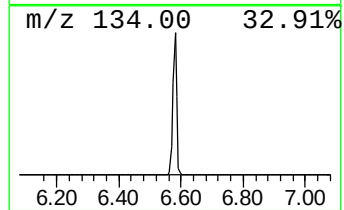
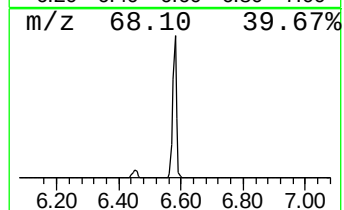
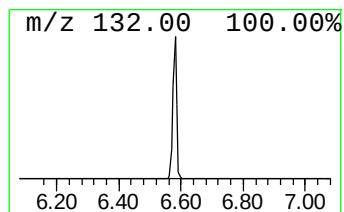
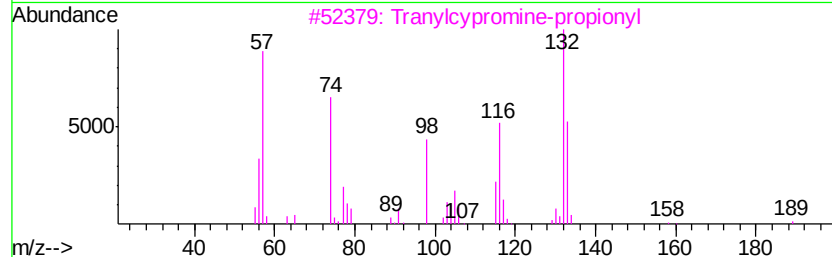
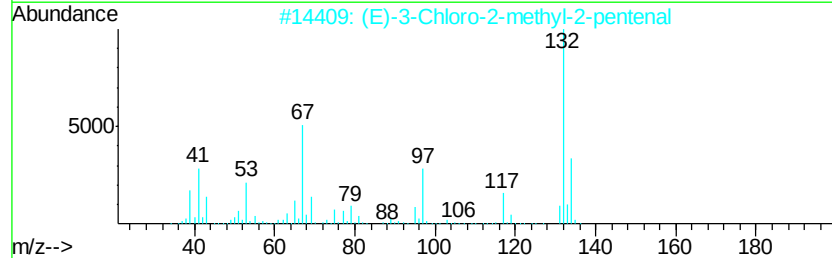
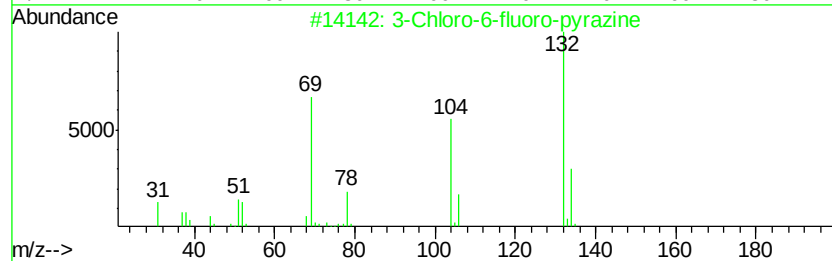
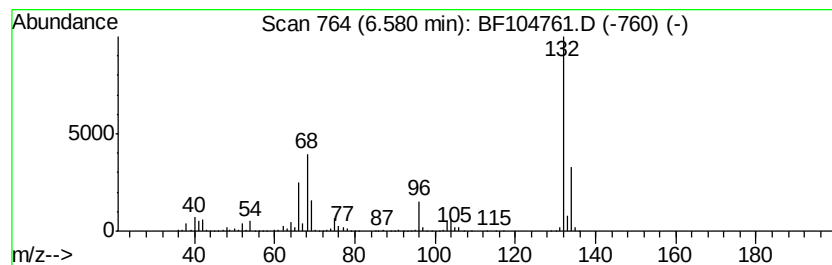
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.58 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	66.57 ng	2008230	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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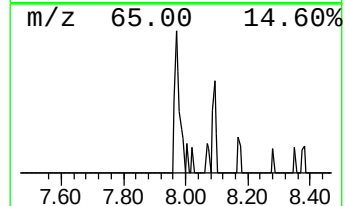
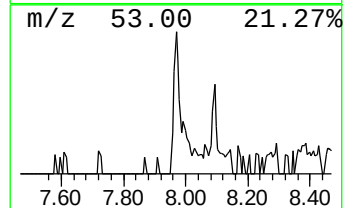
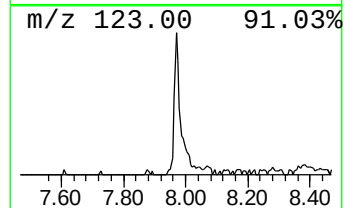
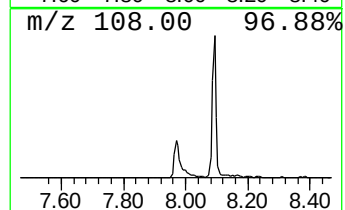
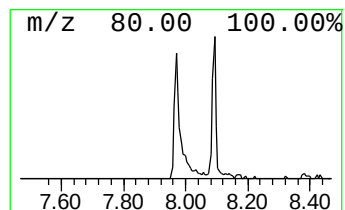
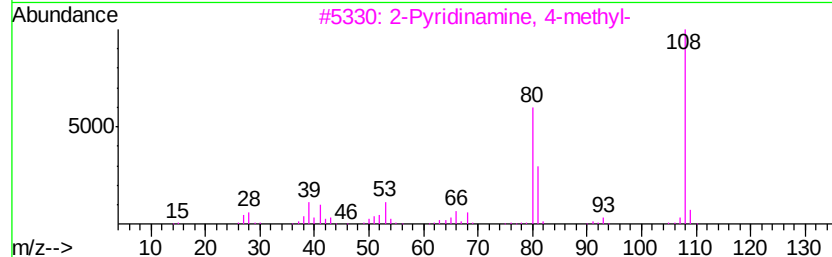
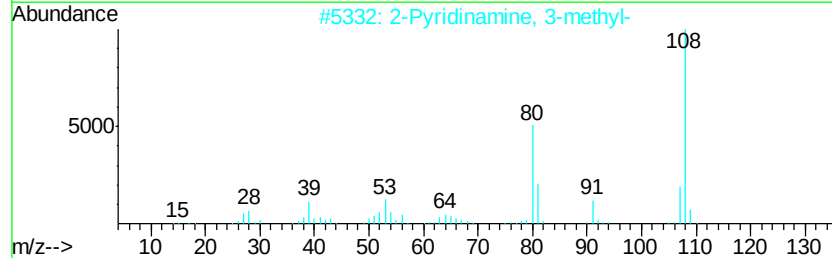
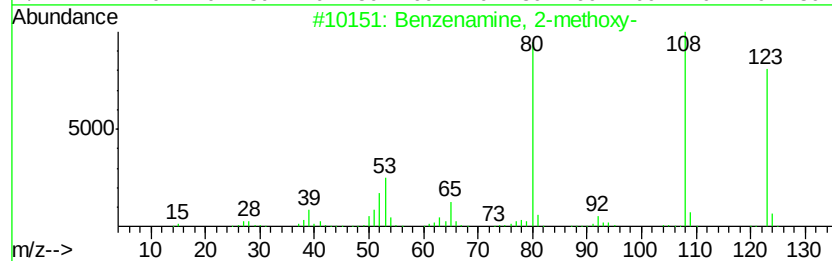
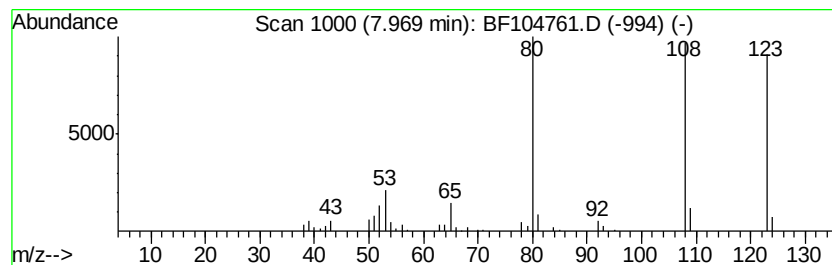
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Peak Number 3 Benzenamine, 2-methoxy- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	2.52 ng	97174	Napthalene-d8	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 2-methoxy-	123	C7H9NO	000090-04-0	95
2		2-Pyridinamine, 3-methyl-	108	C6H8N2	001603-40-3	64
3		2-Pyridinamine, 4-methyl-	108	C6H8N2	000695-34-1	64
4		Bicyclo[2.2.2]oct-2-ene	108	C8H12	000931-64-6	64
5		1,2-Benzenediamine	108	C6H8N2	000095-54-5	64



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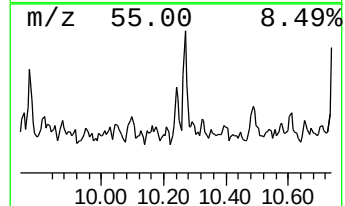
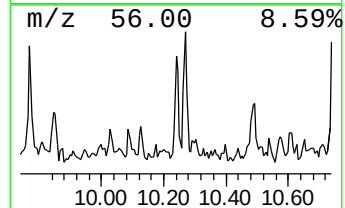
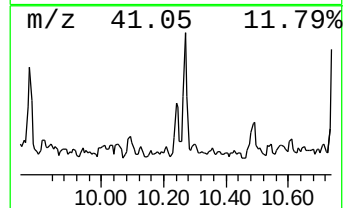
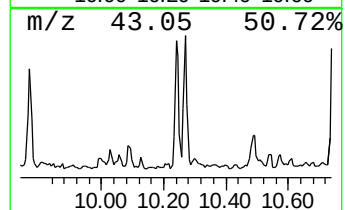
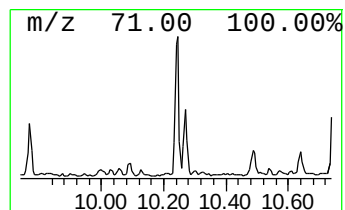
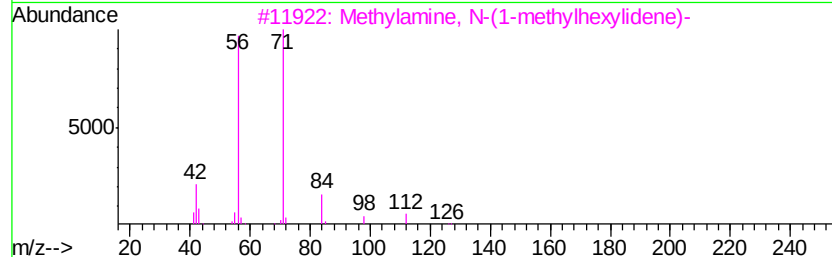
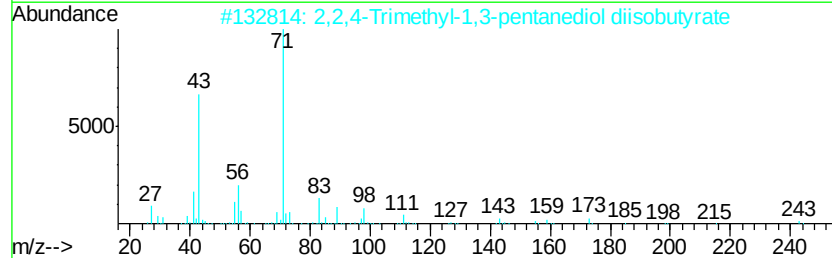
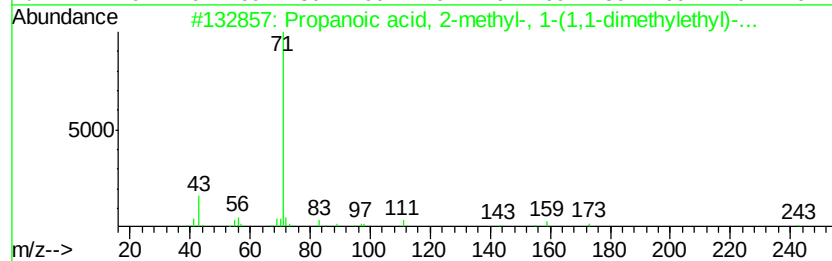
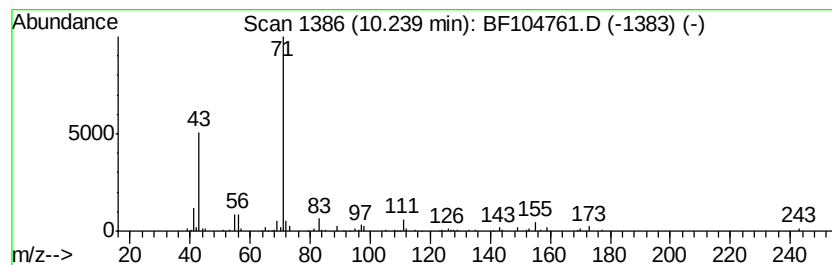
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Peak Number 4 Propanoic acid, 2-methyl-, ... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	2.92 ng	118411	Acenaphthene-d10	9.85
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Propanoic acid, 2-methyl-, 1-(1,...	286 C16H30O4	074381-40-1	72
2	2,2,4-Trimethyl-1,3-pentanediol ...	286 C16H30O4	006846-50-0	72
3	Methylamine, N-(1-methylhexylidene)...	127 C8H17N	022058-71-5	59
4	diethyl 2-hydroxy-3-(tetrahydrof...	260 C12H20O6	1000365-67-1	50
5	3,3-Dimethylheptadecane	268 C19H40	1000360-43-3	45



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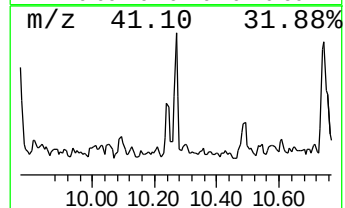
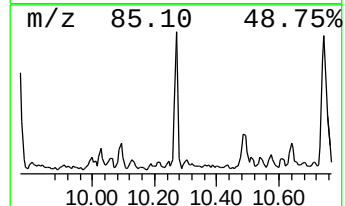
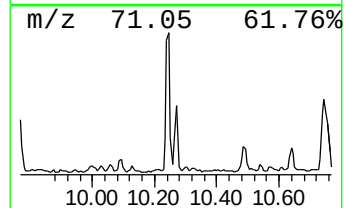
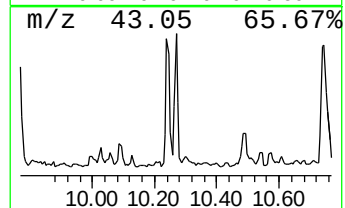
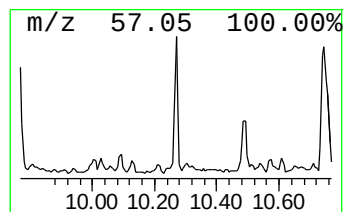
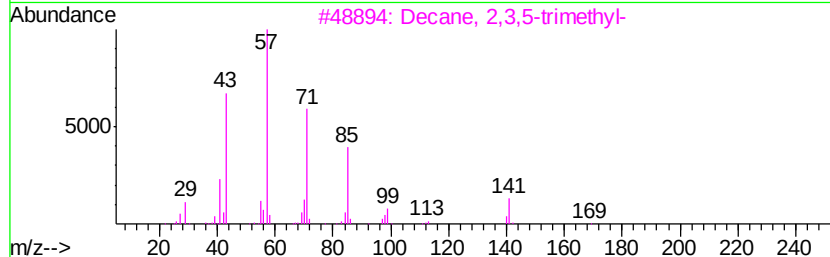
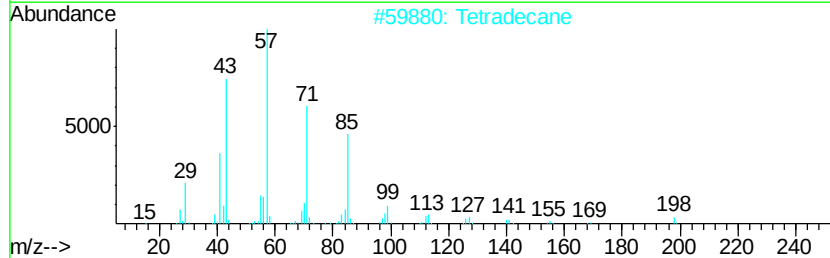
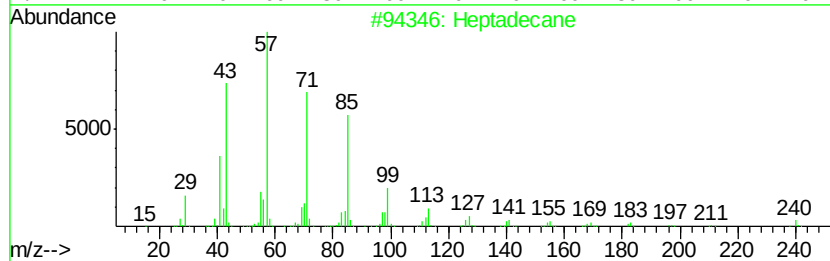
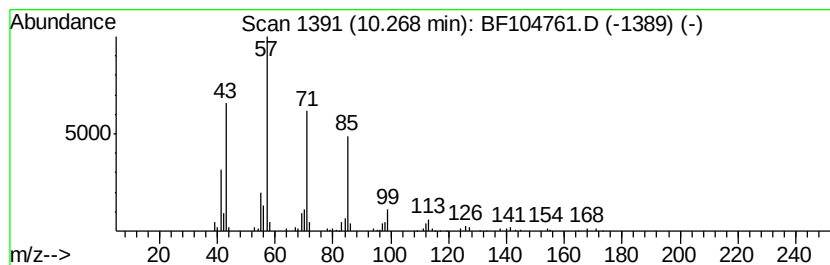
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Peak Number 5 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.27	2.86 ng	115713	Acenaphthene-d10	9.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	90
2	Tetradecane	198	C14H30	000629-59-4	86
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86
4	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1000309-12-5	86
5	Hexadecane	226	C16H34	000544-76-3	83



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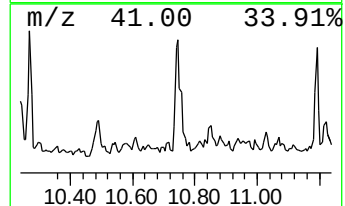
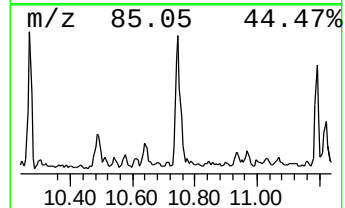
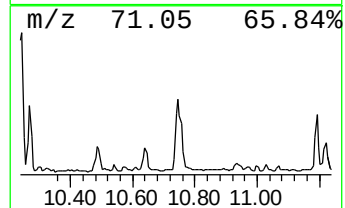
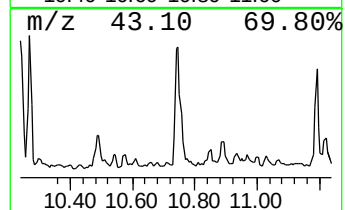
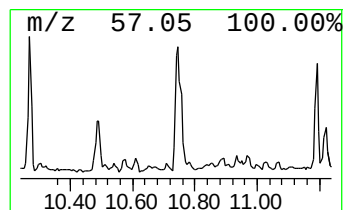
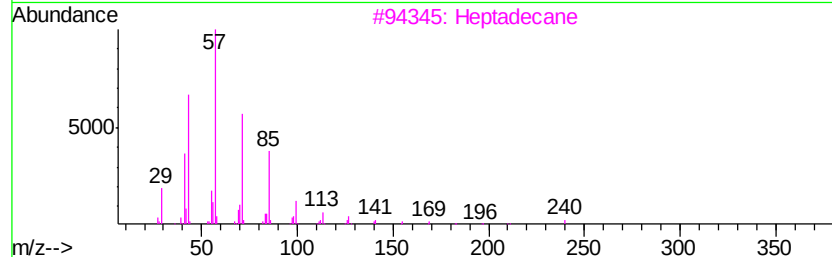
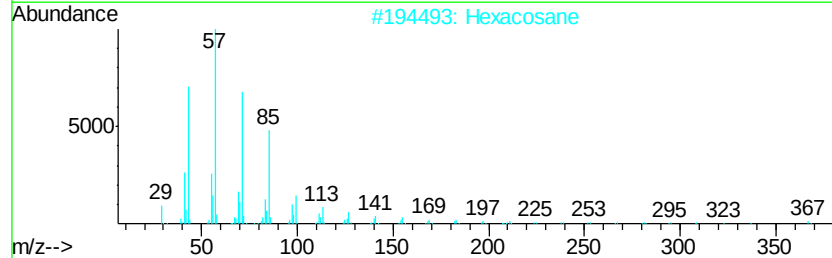
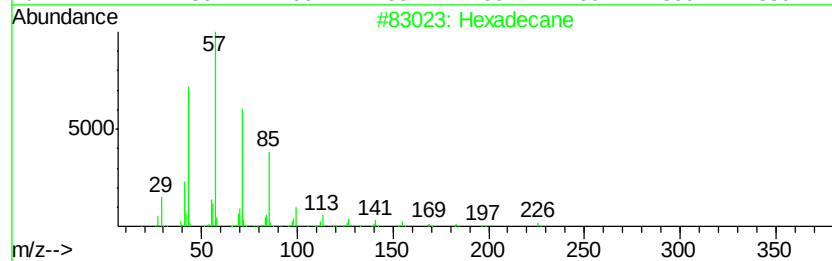
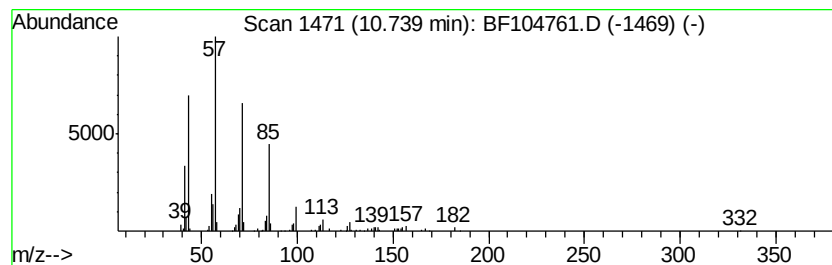
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Peak Number 6 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.74	5.48 ng	203248	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Hexacosane	366	C26H54	000630-01-3	90
3	Heptadecane	240	C17H36	000629-78-7	90
4	Pentadecane	212	C15H32	000629-62-9	87
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86



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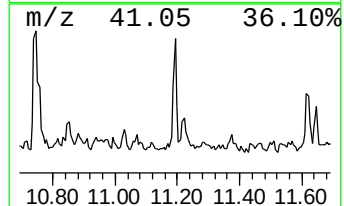
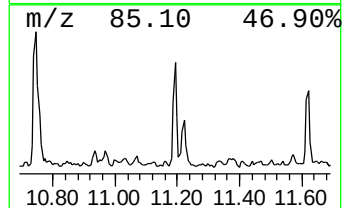
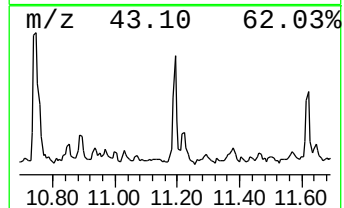
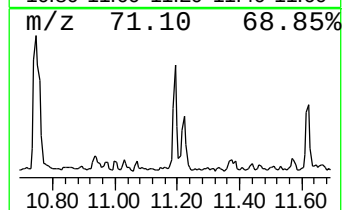
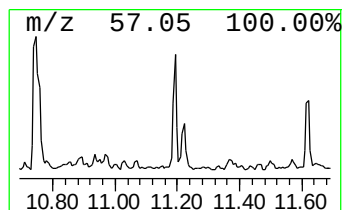
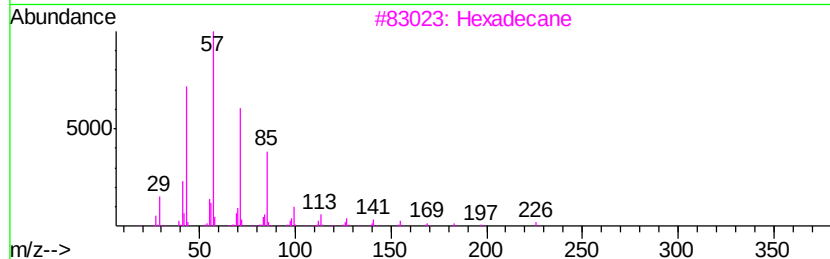
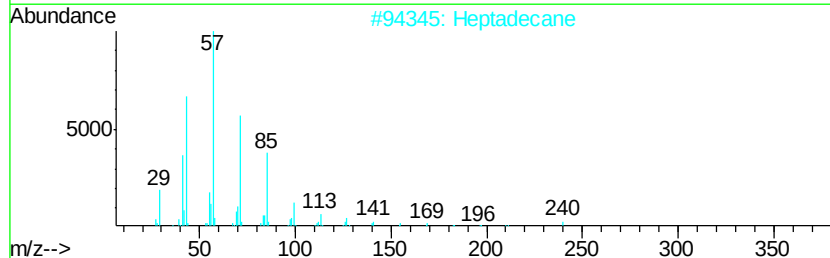
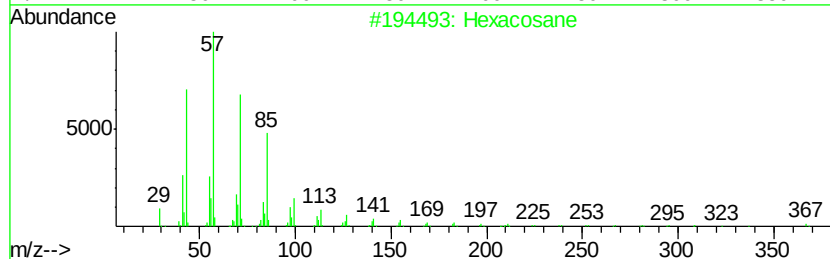
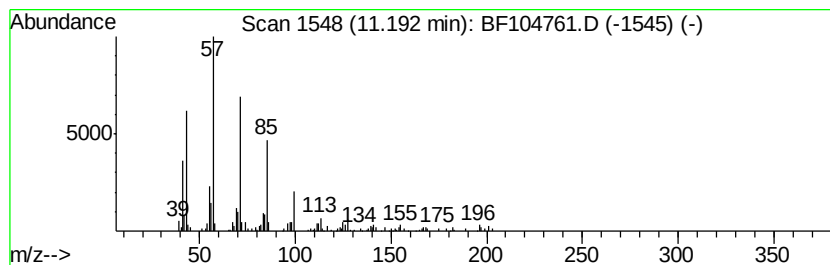
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	3.04 ng	112613	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	83
2		Heptadecane	240	C17H36	000629-78-7	80
3		Hexadecane	226	C16H34	000544-76-3	80
4		Pentacosane	352	C25H52	000629-99-2	80
5		Pentadecane	212	C15H32	000629-62-9	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104761.D
Acq On : 24 Apr 2018 17:27
Operator : JU/SJ
Sample : J2151-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-042318-A

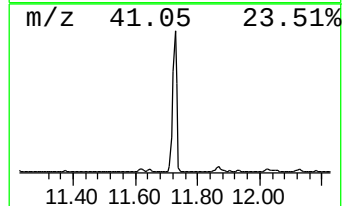
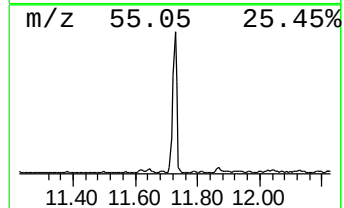
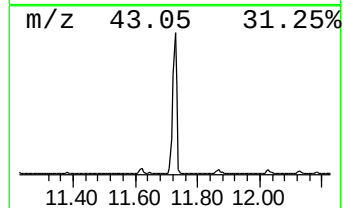
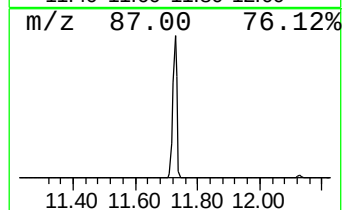
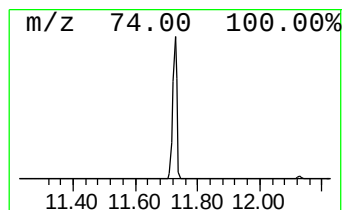
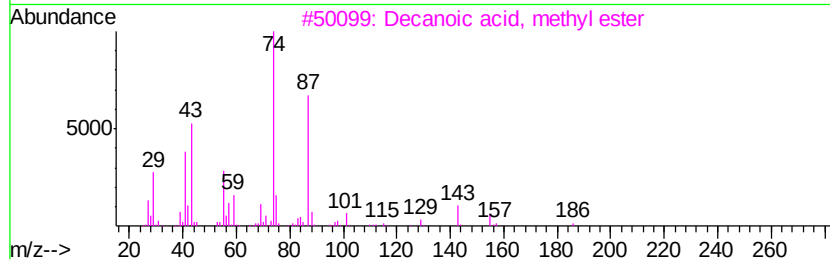
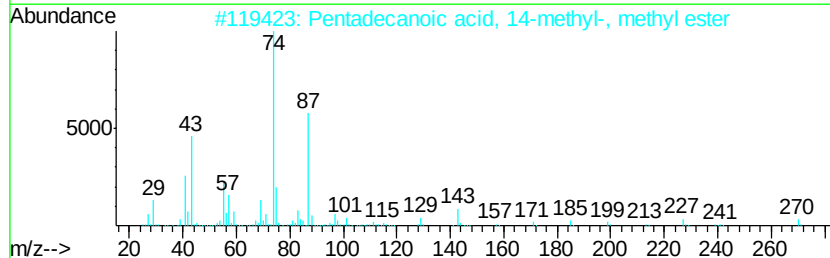
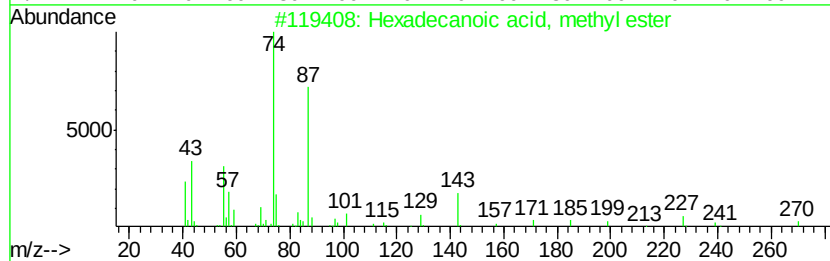
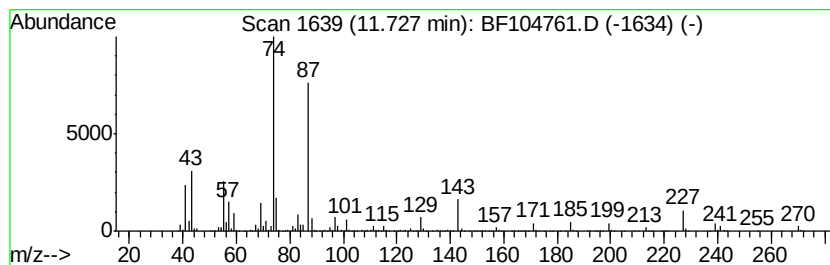
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	108.37 ng	4017160	Phenanthrene-d10	11.34

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	92
3		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
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Acq On : 24 Apr 2018 17:27
Operator : JU/SJ
Sample : J2151-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

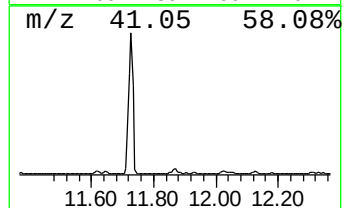
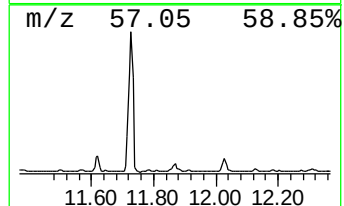
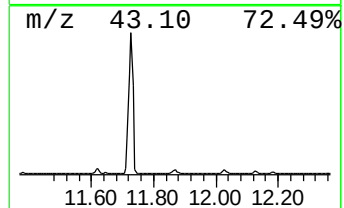
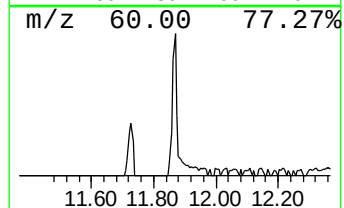
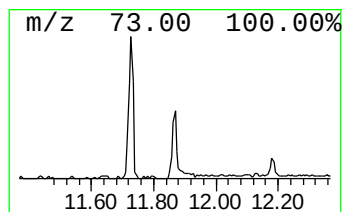
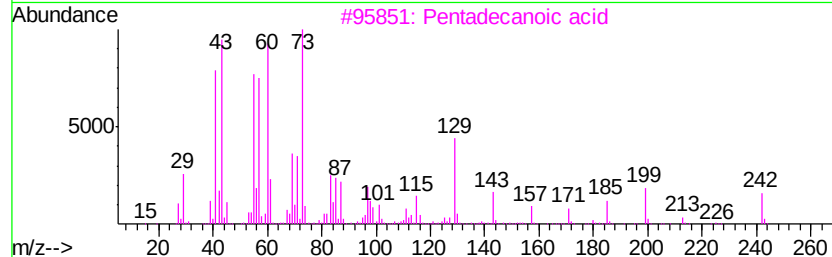
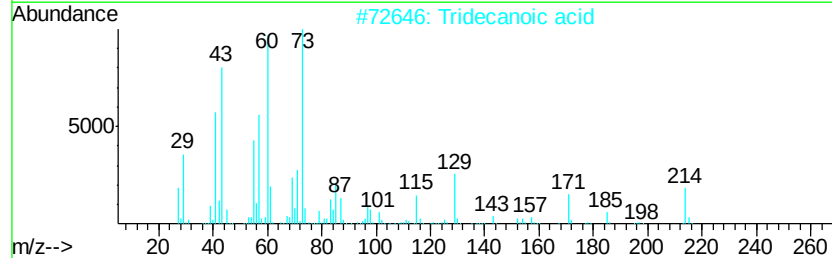
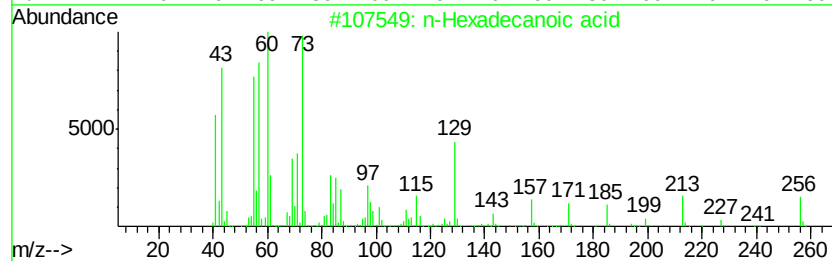
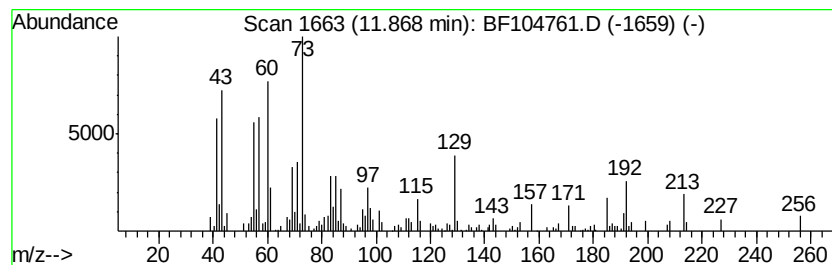
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.87	3.78 ng	140076	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	89
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	n-Decanoic acid	172	C10H20O2	000334-48-5	60
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
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Operator : JU/SJ
Sample : J2151-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-042318-A

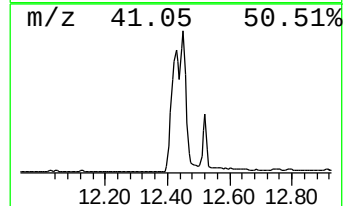
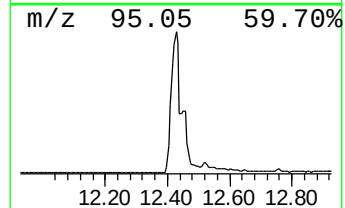
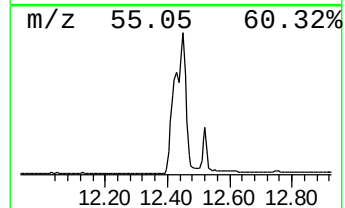
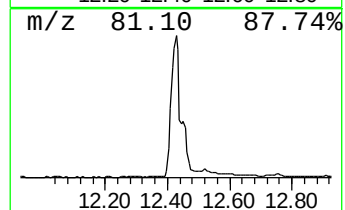
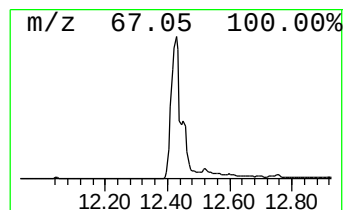
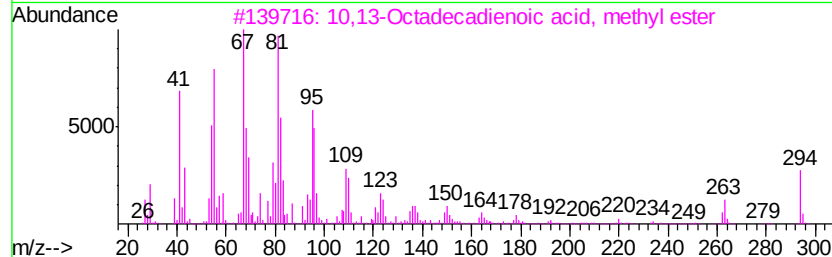
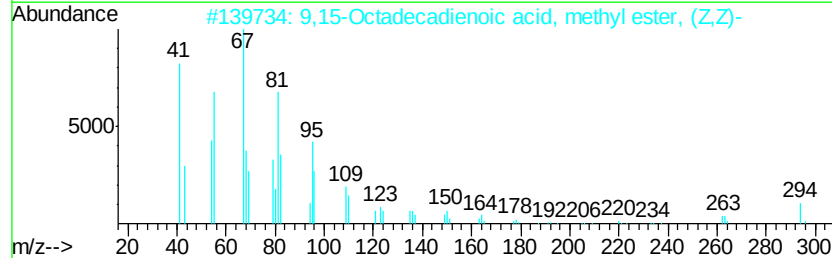
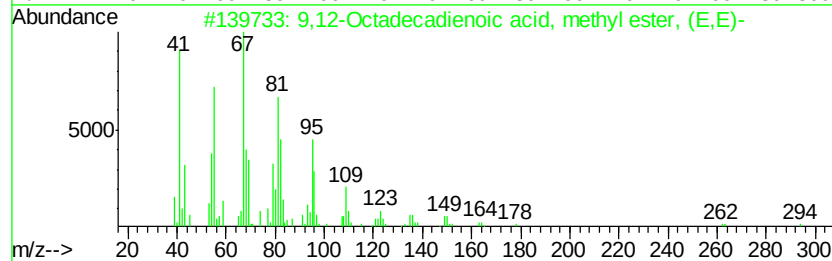
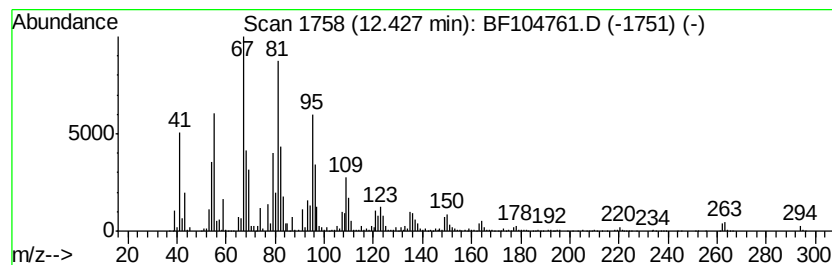
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	256.22 ng	9498120	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
4		11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	98
5		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
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Acq On : 24 Apr 2018 17:27
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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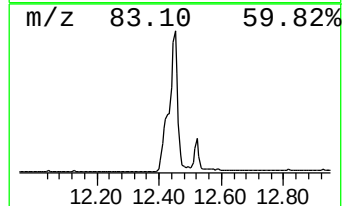
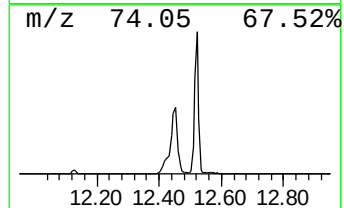
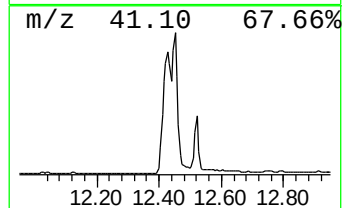
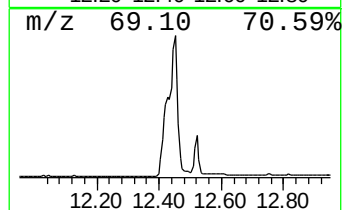
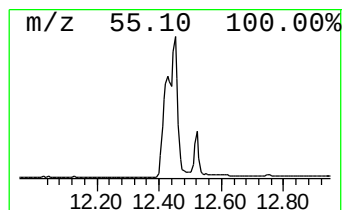
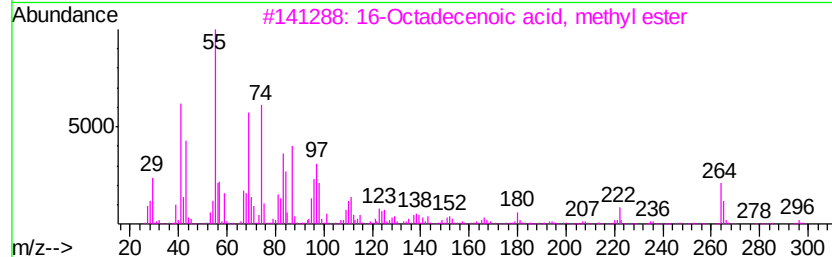
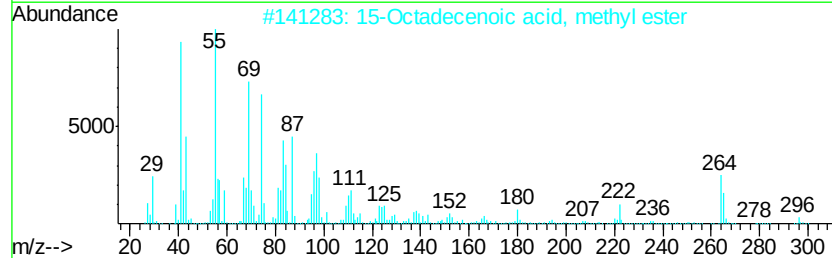
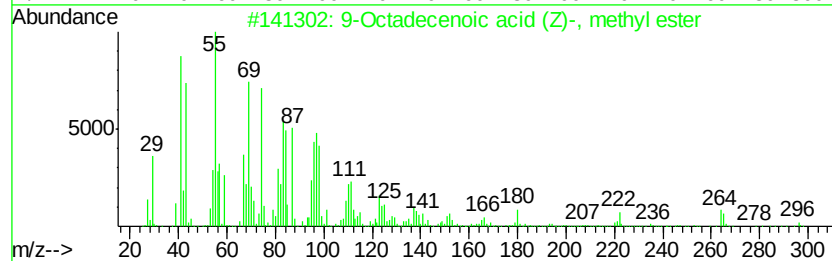
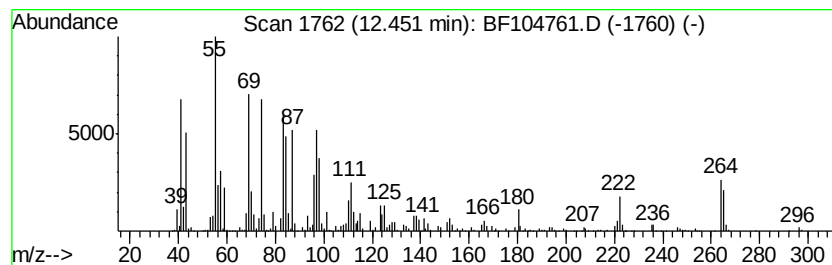
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	180.72 ng	6699370	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	90
2		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	90
3		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	90
4		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	87
5		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104761.D
Acq On : 24 Apr 2018 17:27
Operator : JU/SJ
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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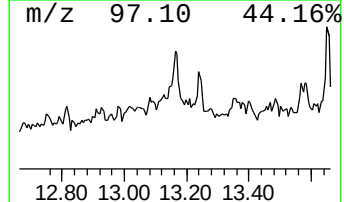
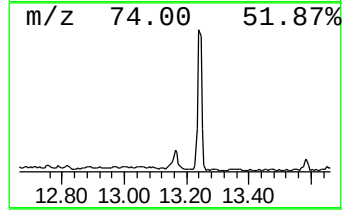
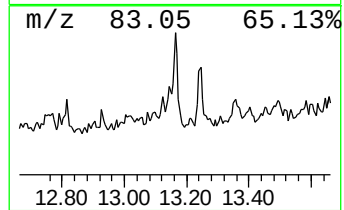
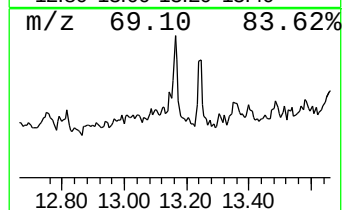
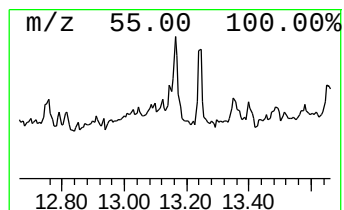
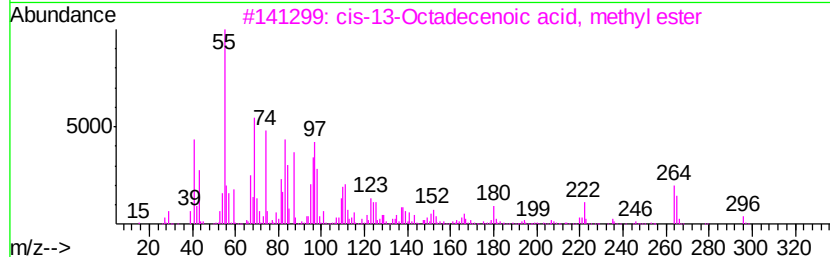
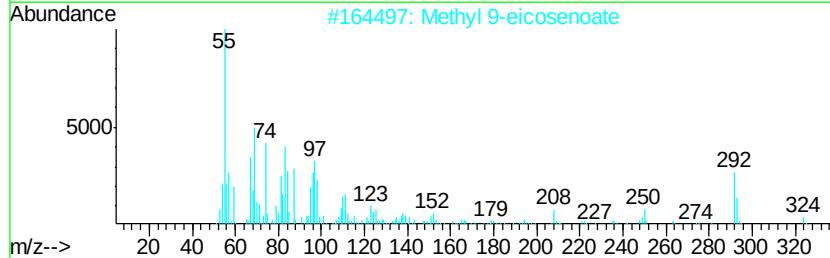
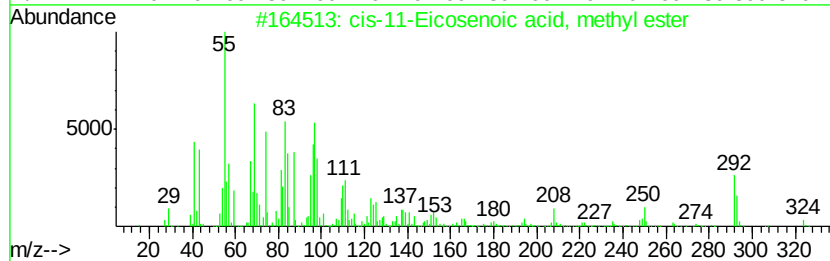
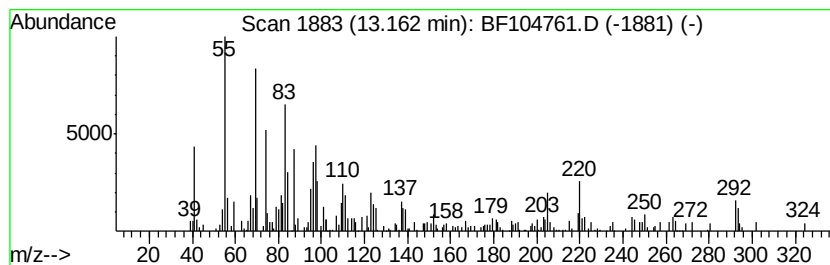
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 cis-11-Eicosenoic acid, met... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	5.43 ng	188630	Chrysene-d12	13.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-11-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-63-8	56
2			Methyl 9-eicosenoate	324	C21H40O2	1000336-50-5	53
3			cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	46
4			11-Eicosenoic acid, methyl ester	324	C21H40O2	003946-08-5	45
5			cis-13-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-52-1	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042418\
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ALS Vial : 12 Sample Multiplier: 1

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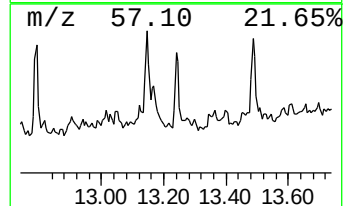
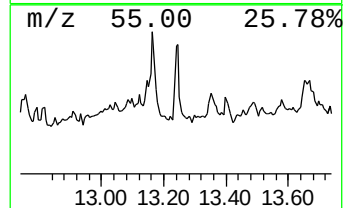
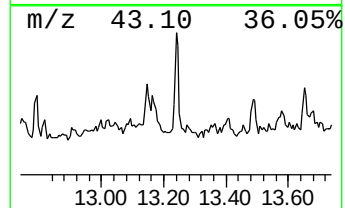
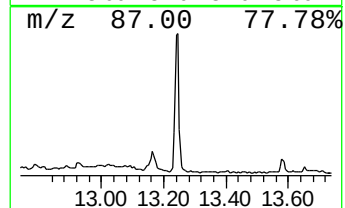
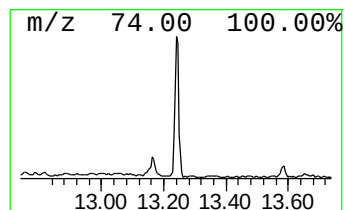
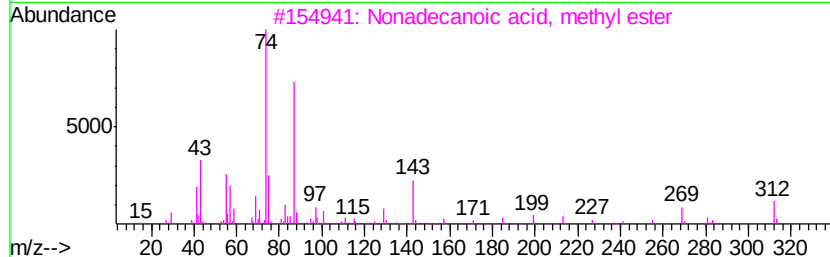
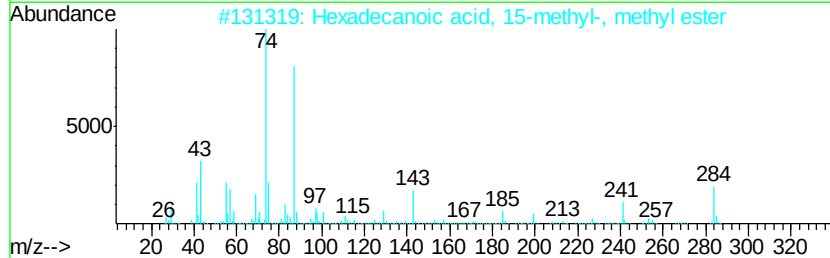
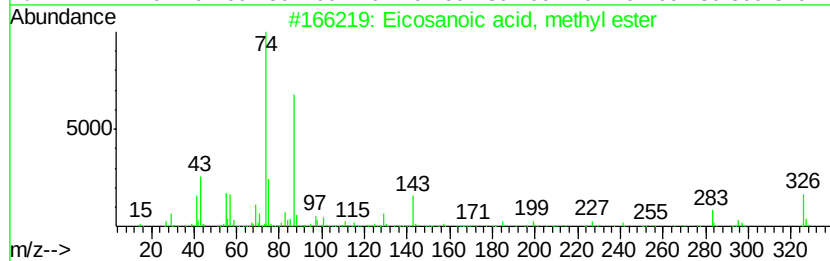
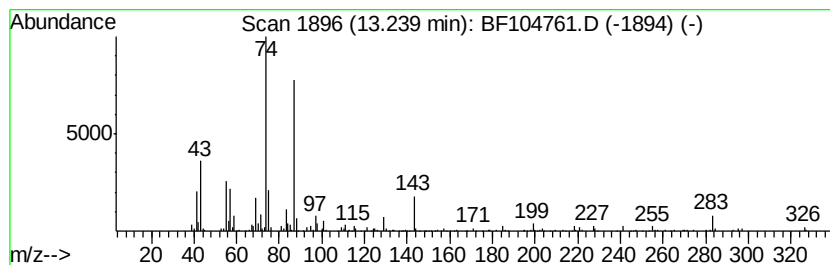
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosanoic acid, methyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	5.84 ng	202940	Chrysene-d12	13.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	97
2			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	93
3			Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	90
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90
5			Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	90



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Misc :
ALS Vial : 12 Sample Multiplier: 1

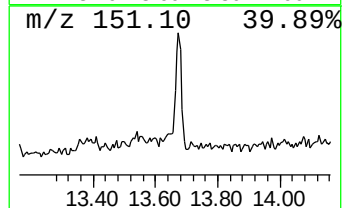
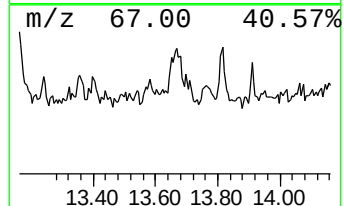
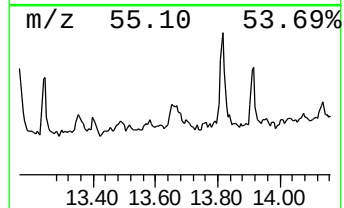
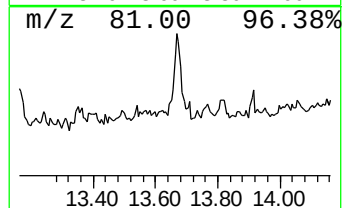
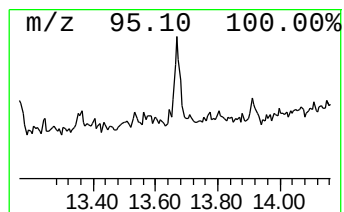
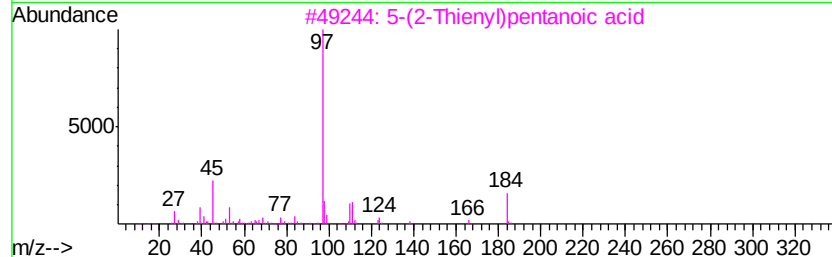
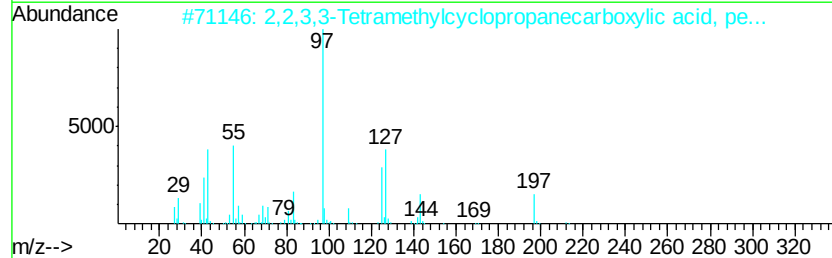
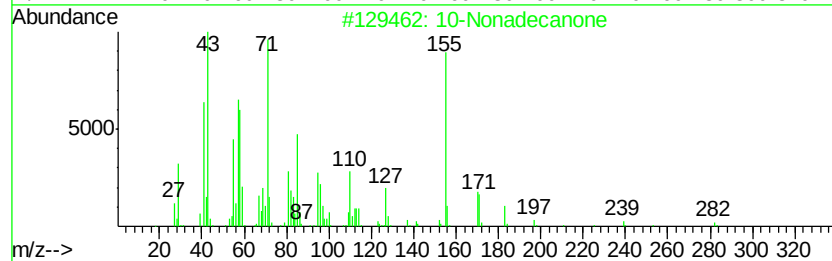
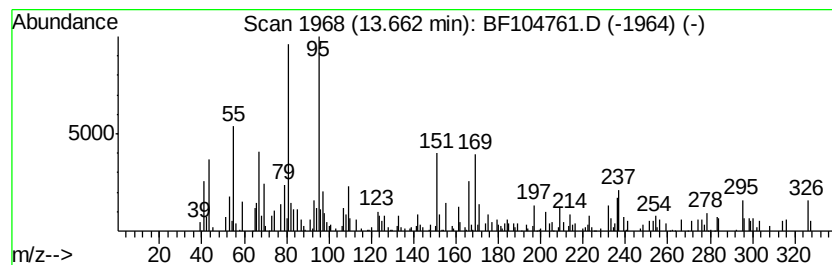
Instrument :
BNA_F
ClientSampleId :
AU-05-042318-A

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown13.66 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.66	2.78 ng	96682	Chrysene-d12		13.99		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10		Nonadecanone	282	C19H38O	000504-57-4	30
2	2		2,2,3,3-Tetramethylcyclopropanec...	212	C13H24O2	1000221-98-6	25
3	5		(2-Thienyl)pentanoic acid	184	C9H12O2S	021010-06-0	15
4	3		Thiopheneethanol	128	C6H8OS	013781-67-4	11
5	5		(7,7-Dimethyl-1-methyl)dithiome...	328	C16H24O3S2	1000210-90-4	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104761.D
Acq On : 24 Apr 2018 17:27
Operator : JU/SJ
Sample : J2151-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

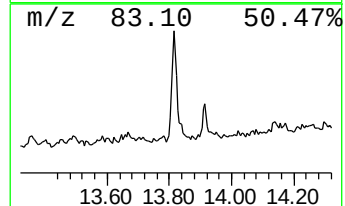
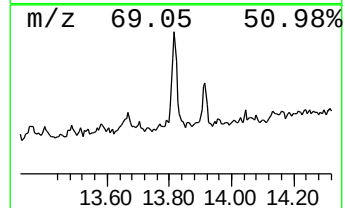
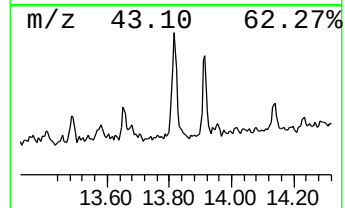
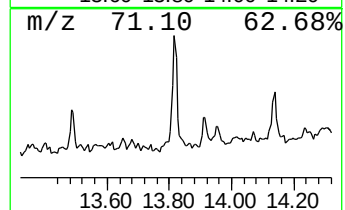
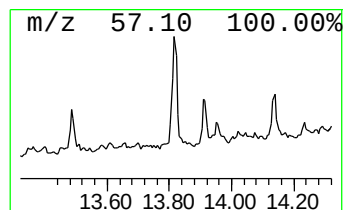
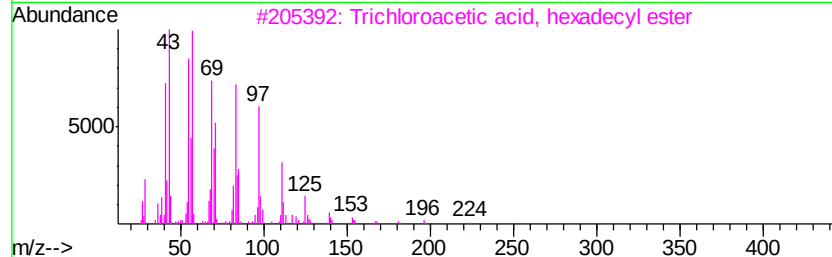
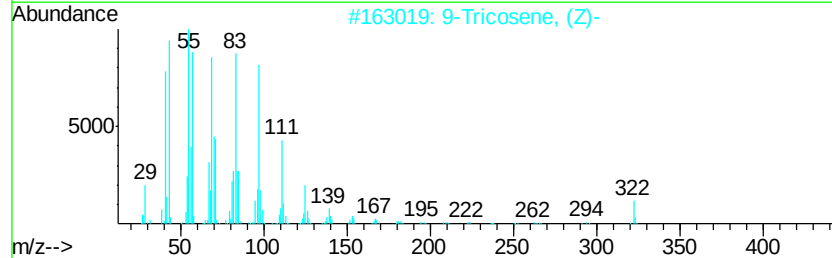
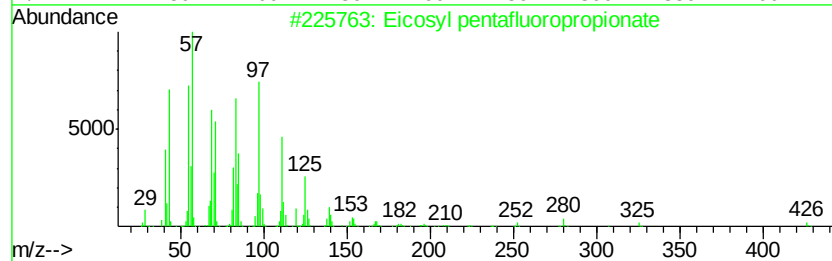
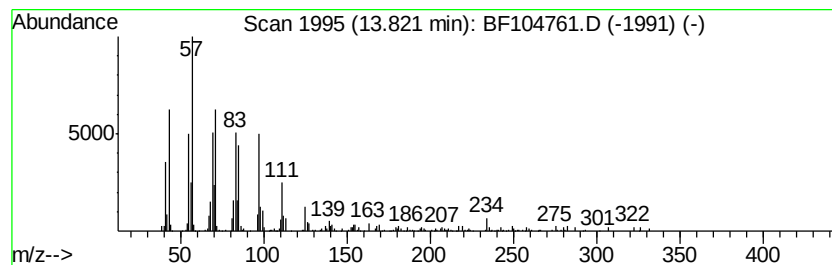
Instrument :
BNA_F
ClientSampleId :
AU-05-042318-A

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosyl pentafluoropropionate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.82	6.74 ng	233932	Chrysene-d12	13.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	90
2	9-Tricosene, (Z)-	322	C23H46	027519-02-4	89
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
4	1-Tricosene	322	C23H46	018835-32-0	87
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104761.D
Acq On : 24 Apr 2018 17:27
Operator : JU/SJ
Sample : J2151-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-042318-A

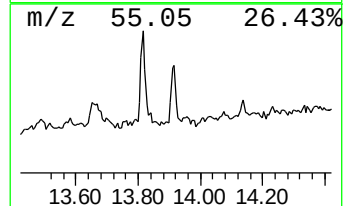
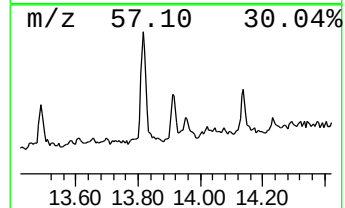
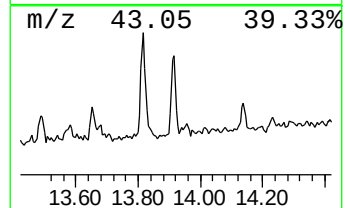
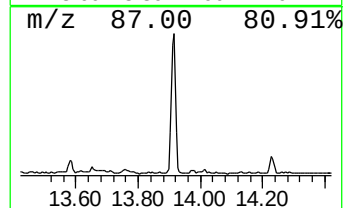
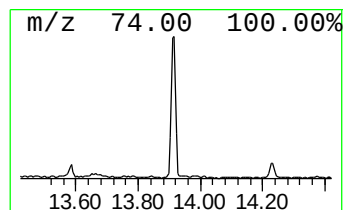
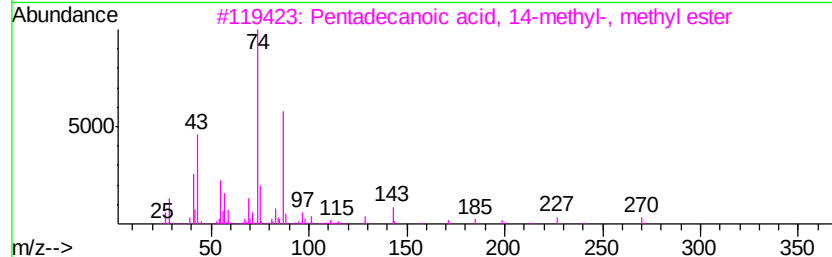
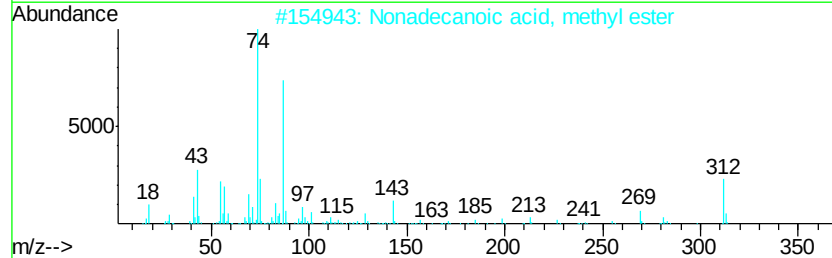
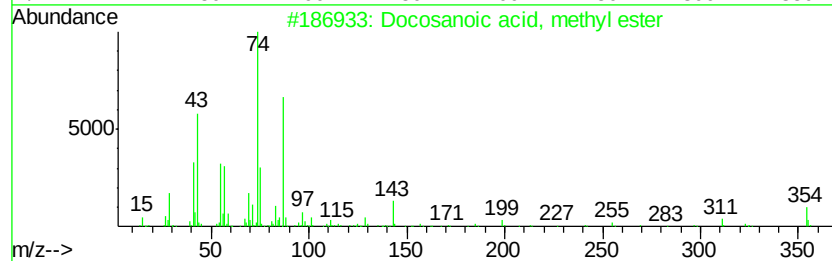
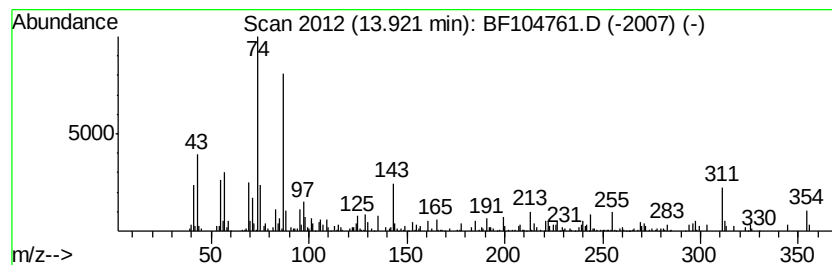
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Docosanoic acid, methyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	7.08 ng	245950	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	98
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	93
3		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90
4		Octadecanoic acid, 10-methyl-, m...	312	C20H40O2	002490-19-9	90
5		Methyl stearate	298	C19H38O2	000112-61-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042418\
 Data File : BF104761.D
 Acq On : 24 Apr 2018 17:27
 Operator : JU/SJ
 Sample : J2151-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-042318-A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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
unknown6.58	6.58	66.6	ng	2008230	1	6.80	603385	20.0
Benzenamine, 2-me...	7.97	2.5	ng	97174	2	8.09	772285	20.0
Propanoic acid, 2...	10.24	2.9	ng	118411	3	9.85	810462	20.0
Heptadecane	10.27	2.9	ng	115713	3	9.85	810462	20.0
Hexadecane	10.74	5.5	ng	203248	4	11.34	741403	20.0
Hexacosane	11.19	3.0	ng	112613	4	11.34	741403	20.0
Hexadecanoic acid...	11.73	108.4	ng	4017160	4	11.34	741403	20.0
n-Hexadecanoic acid	11.87	3.8	ng	140076	4	11.34	741403	20.0
9,12-Octadecadien...	12.43	256.2	ng	9498120	4	11.34	741403	20.0
9-Octadecenoic ac...	12.45	180.7	ng	6699370	4	11.34	741403	20.0
cis-11-Eicosenoic...	13.16	5.4	ng	188630	5	13.99	694424	20.0
Eicosanoic acid, ...	13.24	5.8	ng	202940	5	13.99	694424	20.0
unknown13.66	13.66	2.8	ng	96682	5	13.99	694424	20.0
Eicosyl pentafluo...	13.82	6.7	ng	233932	5	13.99	694424	20.0
Docosanoic acid, ...	13.92	7.1	ng	245950	5	13.99	694424	20.0