

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102518.D
 Acq On : 27 Jan 2018 18:53
 Operator : SJ/JU
 Sample : J1009-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-012518-B

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-BF012218.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	5.328	546	551	567	rBV	3434927	3717448	9.93%	3.946%
2	6.363	722	727	739	rBV	3324637	3827037	10.22%	4.063%
3	6.487	743	748	751	rBV	3857776	3717312	9.93%	3.946%
4	6.716	783	787	793	rBV	910305	863380	2.31%	0.917%
5	6.869	809	813	816	rBV	3135835	2745291	7.33%	2.914%
6	7.281	879	883	886	rBV	2209531	2329444	6.22%	2.473%
7	7.998	1002	1005	1008	rVB	1152772	1030334	2.75%	1.094%
8	9.075	1183	1188	1191	rBV	3627293	3327445	8.89%	3.532%
9	9.463	1250	1254	1257	rBV	488598	483396	1.29%	0.513%
10	9.751	1299	1303	1306	rVB	1165414	1077115	2.88%	1.143%
11	10.175	1372	1375	1378	rVB	537053	386807	1.03%	0.411%
12	10.539	1434	1437	1447	rVB	1777775	1723192	4.60%	1.829%
13	10.645	1452	1455	1464	rVB	424355	533894	1.43%	0.567%
14	11.233	1552	1555	1558	rBV	985468	842635	2.25%	0.894%
15	11.651	1618	1626	1628	rBV	7928330	13960190	37.29%	14.819%
16	11.945	1669	1676	1679	rBV3	344284	470675	1.26%	0.500%
17	12.374	1734	1749	1751	rBV5	8498926	37437989	100.00%	39.742%
18	12.439	1755	1760	1763	rVB	6336220	7259413	19.39%	7.706%
19	12.657	1792	1797	1801	rBV2	253130	378534	1.01%	0.402%
20	12.833	1820	1827	1829	rBV	2126684	2212869	5.91%	2.349%
21	13.004	1850	1856	1857	rBV3	290544	422194	1.13%	0.448%
22	13.063	1861	1866	1873	rVB2	754877	1125072	3.01%	1.194%
23	13.145	1876	1880	1882	rBV	1015971	934064	2.49%	0.992%
24	13.568	1949	1952	1957	rVB2	305592	440796	1.18%	0.468%
25	13.710	1973	1976	1986	rVB	347953	480755	1.28%	0.510%
26	13.810	1989	1993	1997	rBV	1070713	908534	2.43%	0.964%
27	13.880	2001	2005	2008	rBV	765463	733737	1.96%	0.779%
28	15.304	2243	2247	2253	rBV	633199	832633	2.22%	0.884%

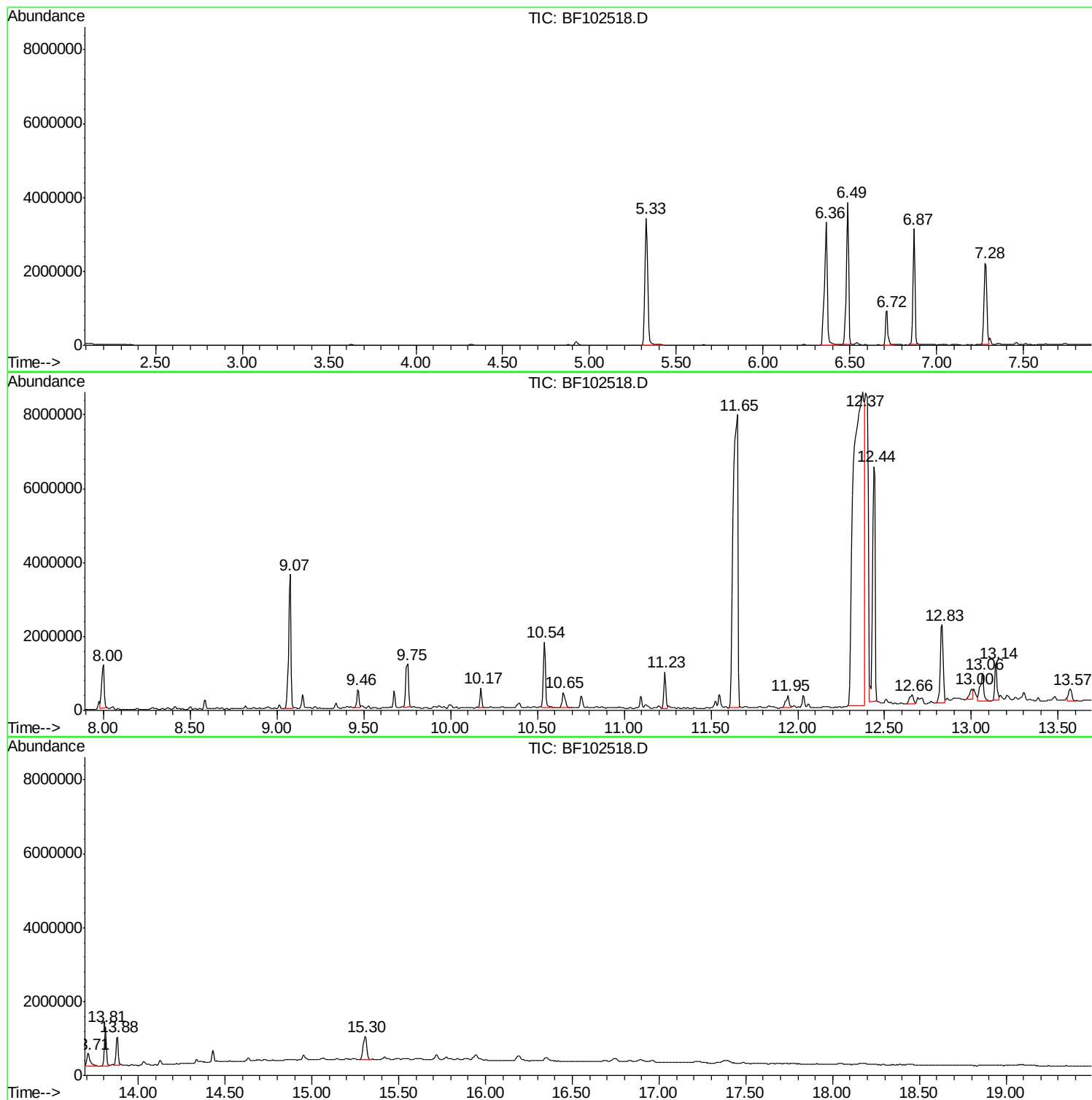
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Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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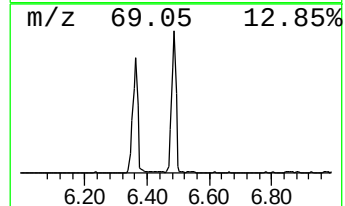
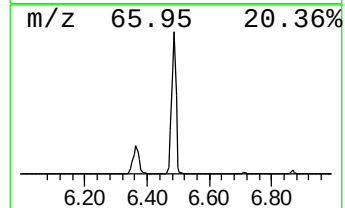
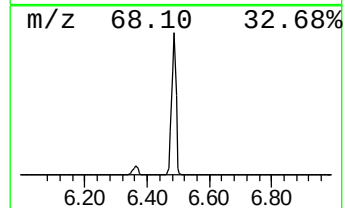
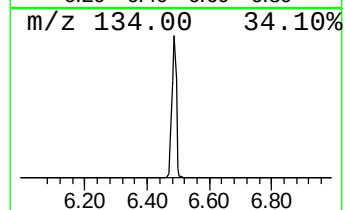
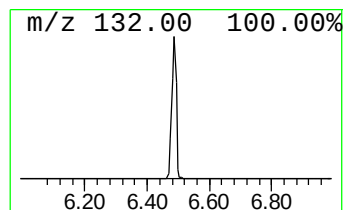
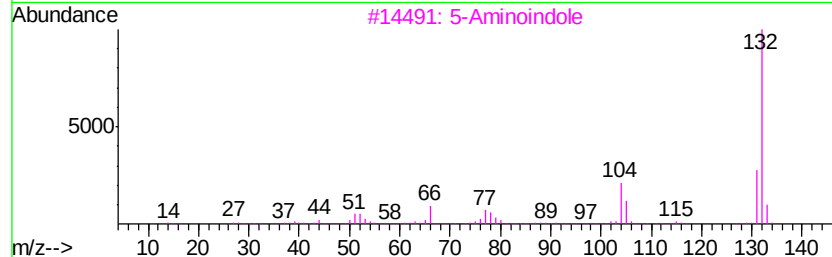
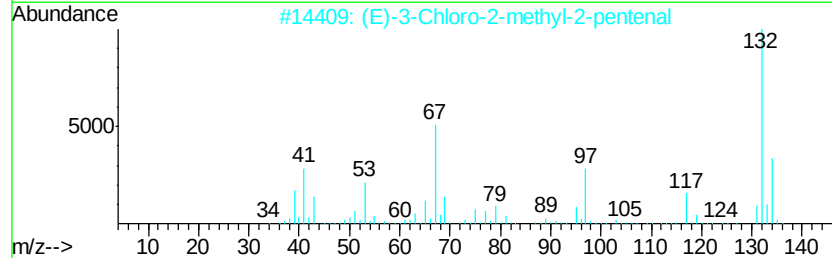
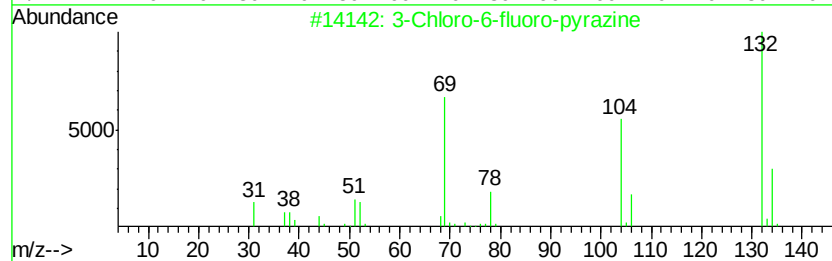
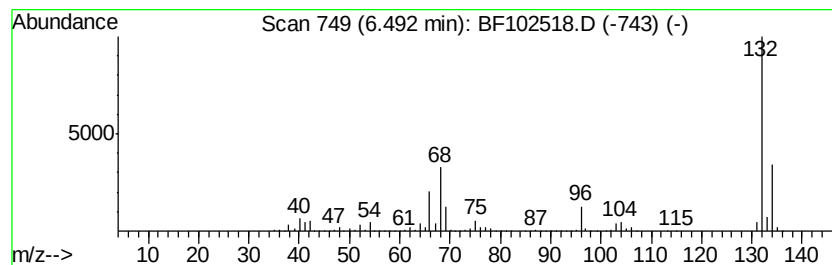
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.49 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	86.11 ng	3717310	1,4-Dichlorobenzene-d4	6.72

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	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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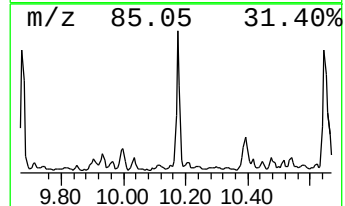
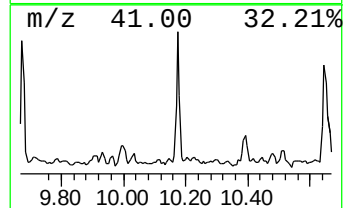
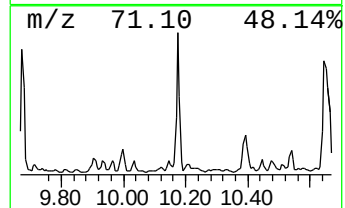
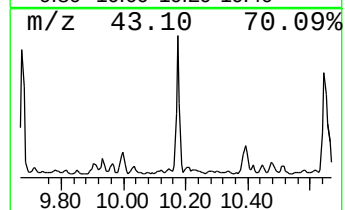
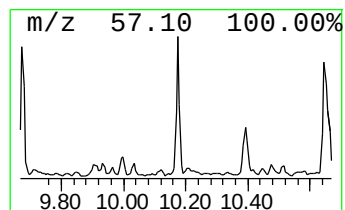
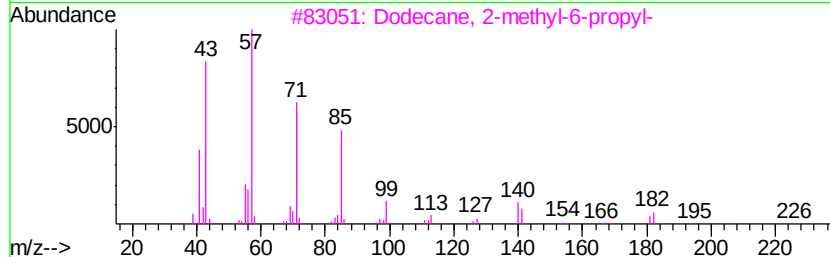
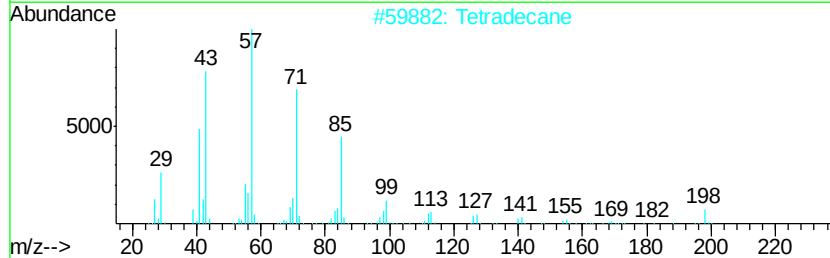
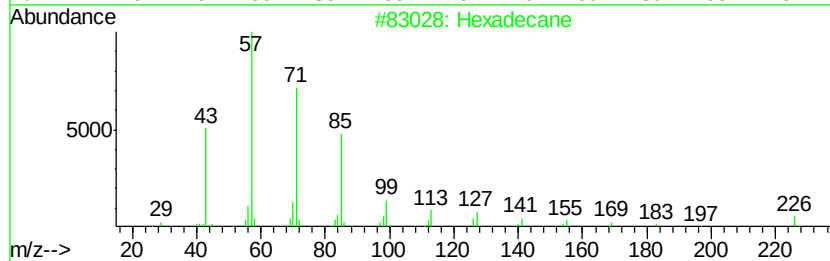
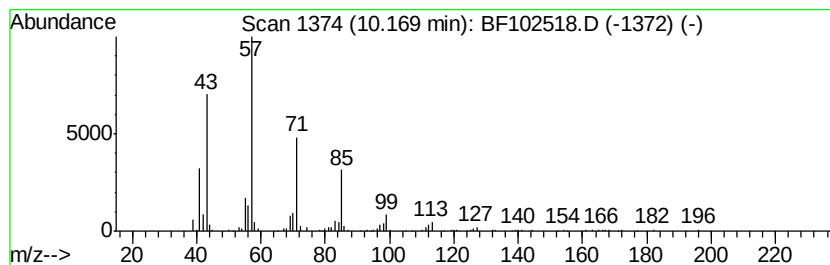
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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Peak Number 3 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.17	7.18 ng	386807	Acenaphthene-d10	9.75
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	94
2	Tetradecane	198 C14H30	000629-59-4	91
3	Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4	80
4	Pentadecane, 7-methyl-	226 C16H34	006165-40-8	72
5	Nonane, 5-butyl-	184 C13H28	017312-63-9	72



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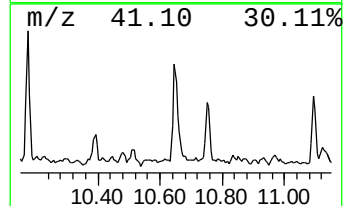
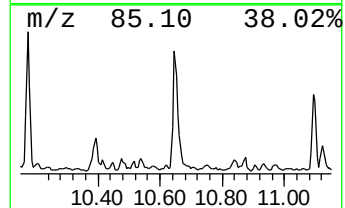
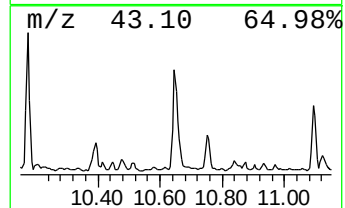
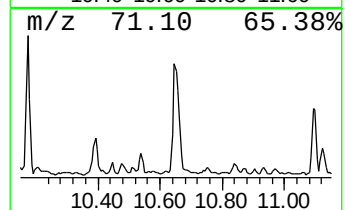
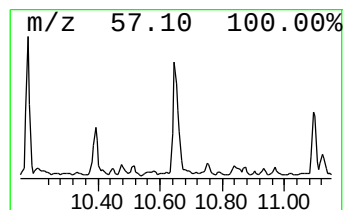
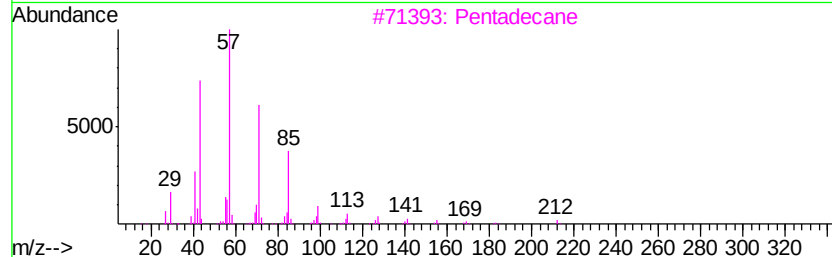
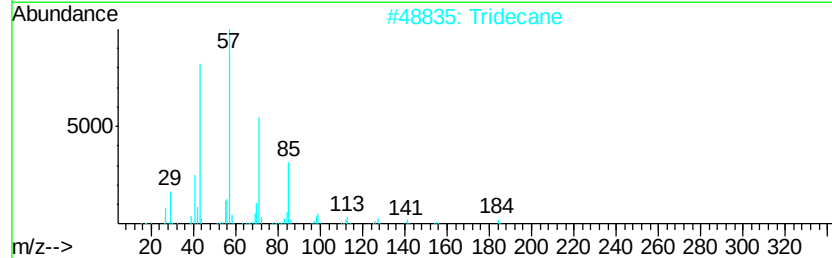
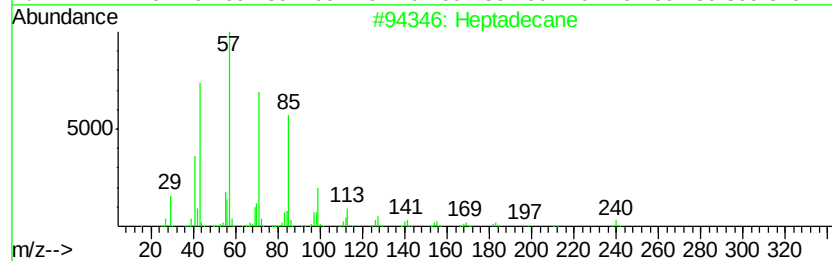
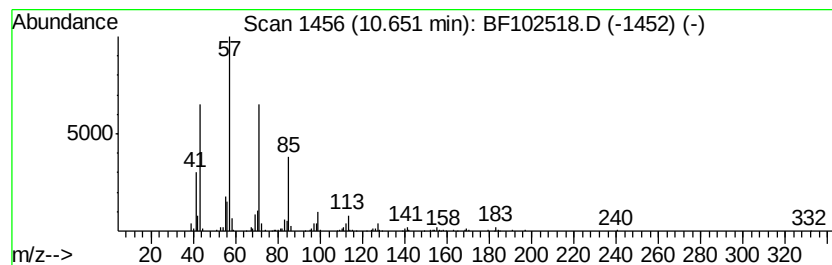
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	12.67 ng	533894	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Tridecane		184	C13H28	000629-50-5	90
3	Pentadecane		212	C15H32	000629-62-9	90
4	Hexadecane		226	C16H34	000544-76-3	90
5	Decane, 3-methyl-		156	C11H24	013151-34-3	87



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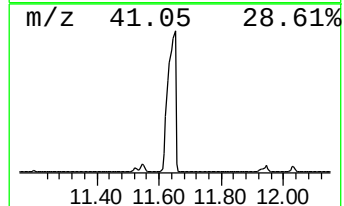
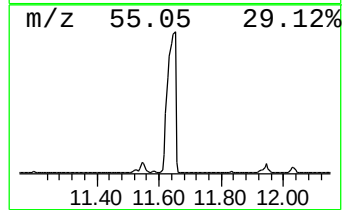
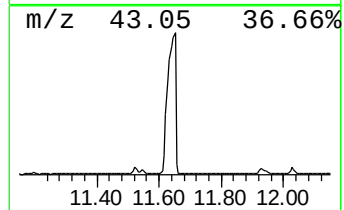
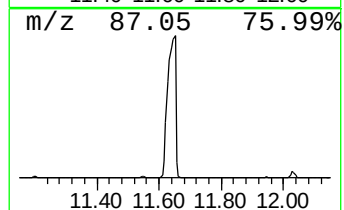
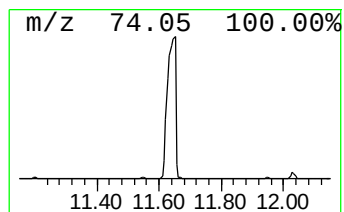
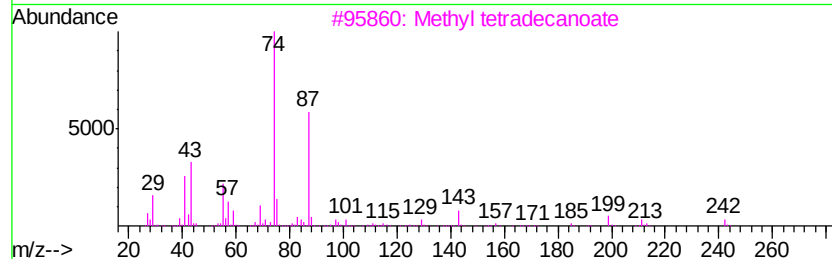
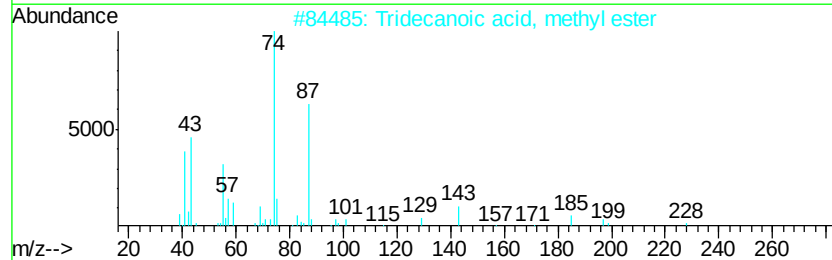
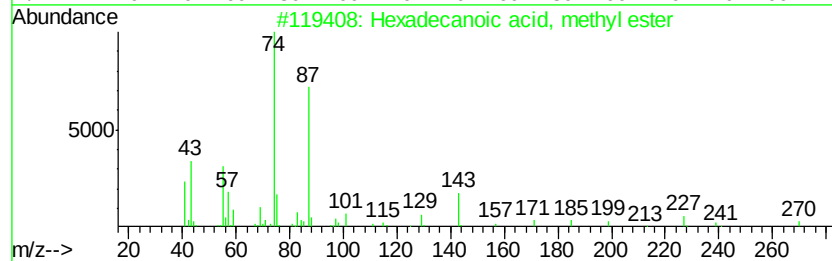
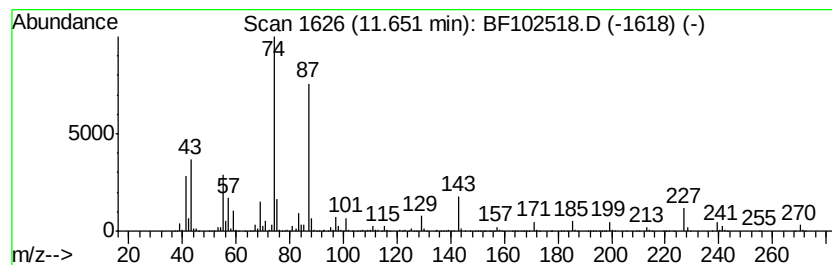
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Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	331.35 ng	13960200	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	95
3		Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
5		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	87



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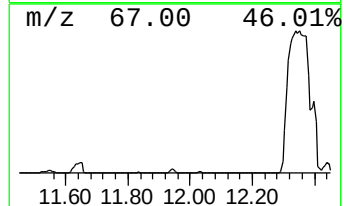
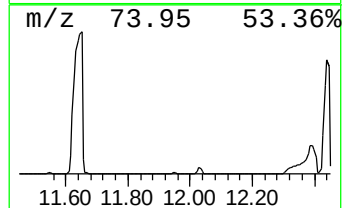
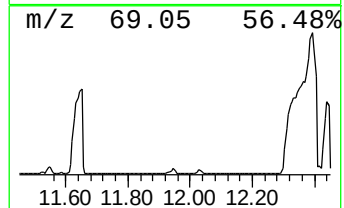
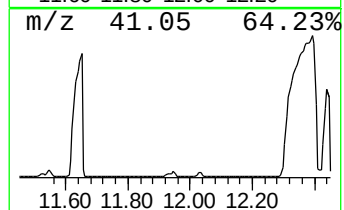
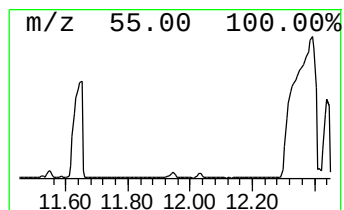
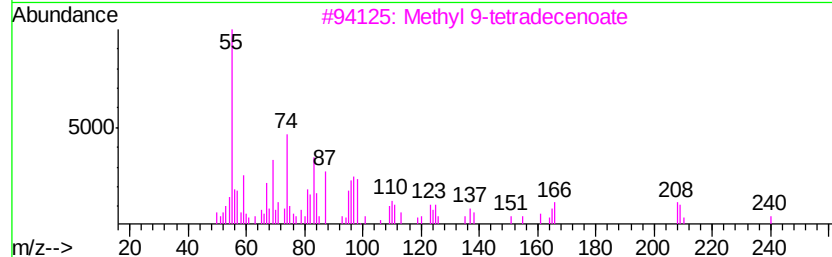
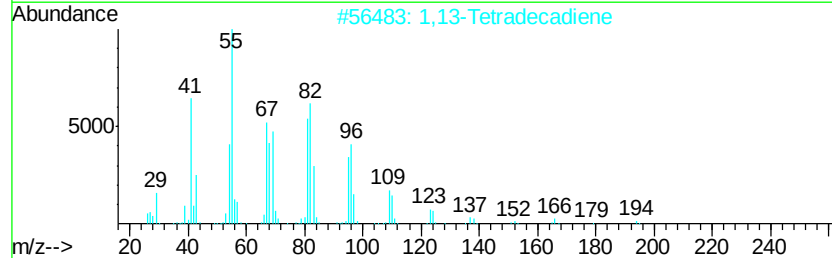
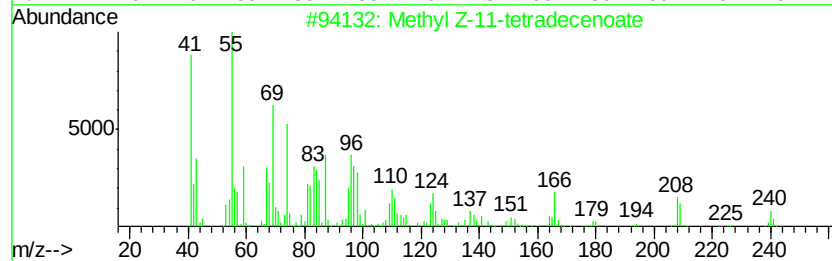
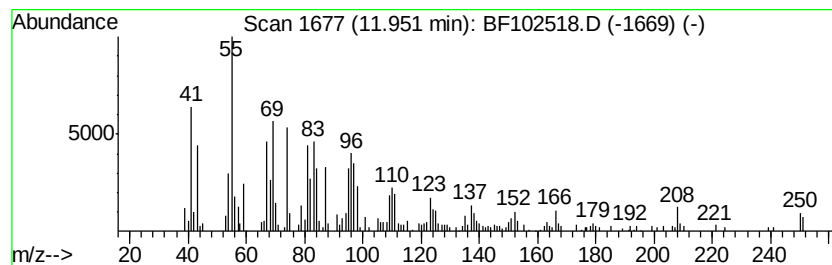
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Peak Number 6 Methyl Z-11-tetradecenoate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.95	11.17 ng	470675	Phenanthrene-d10	11.23	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl Z-11-tetradecenoate	240	C15H28O2	1000130-82-8	70
2	1,13-Tetradecadiene	194	C14H26	021964-49-8	60
3	Methyl 9-tetradecenoate	240	C15H28O2	1000336-47-3	50
4	Methyl myristoleate	240	C15H28O2	056219-06-8	50
5	9-Methylbicyclo[3.3.1]nonane	138	C10H18	025107-01-1	46



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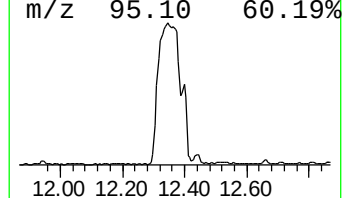
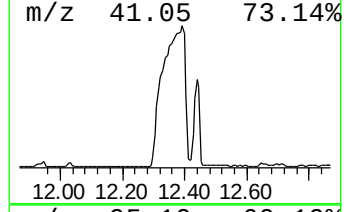
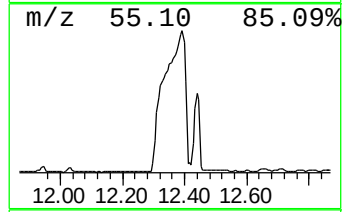
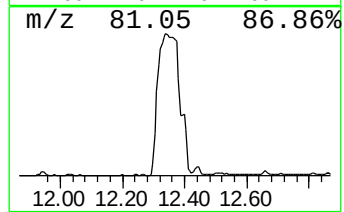
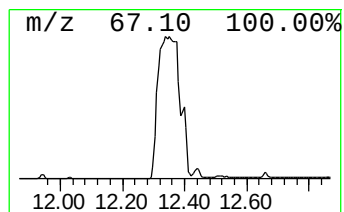
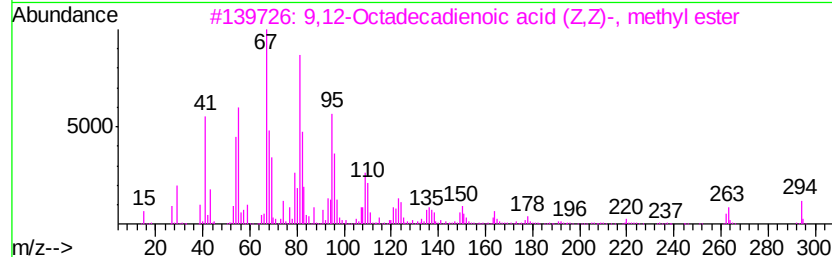
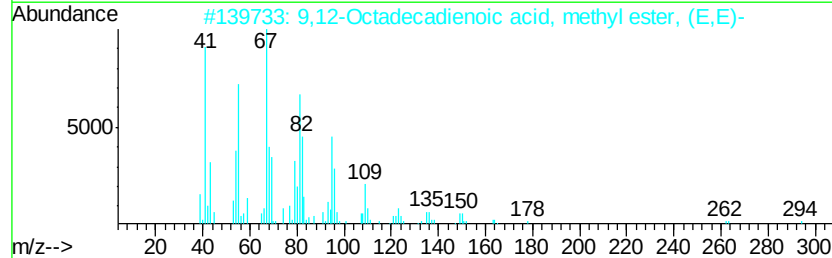
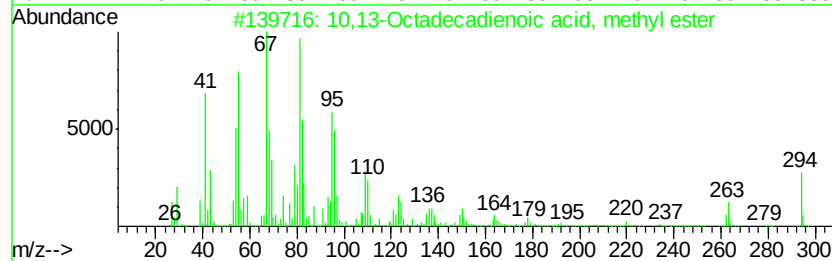
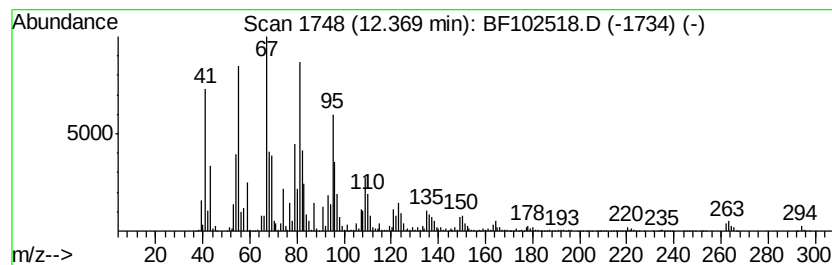
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 10,13-Octadecadienoic acid,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	888.59 ng	37438000	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
2		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	98
4		Methyl 7,12-octadecadienoate	294	C19H34O2	1000336-43-8	96
5		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102518.D
Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
Sample : J1009-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

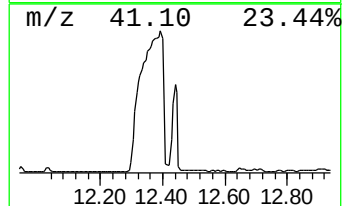
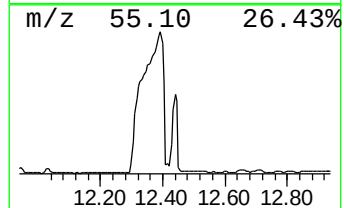
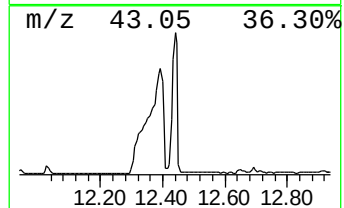
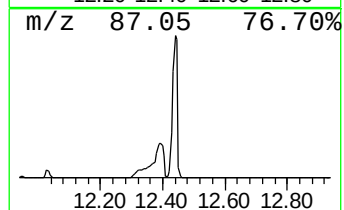
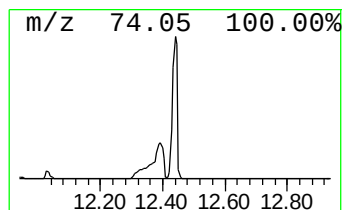
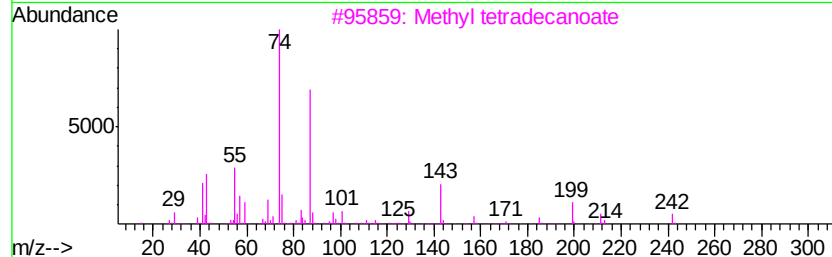
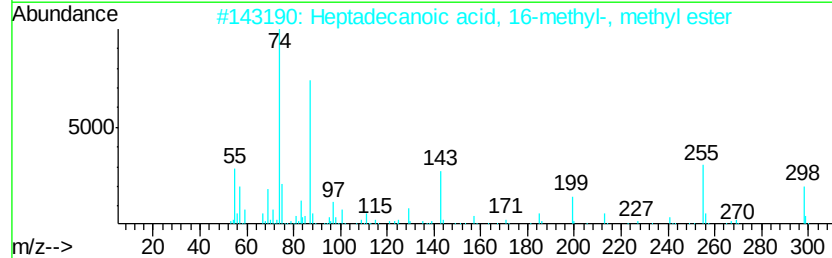
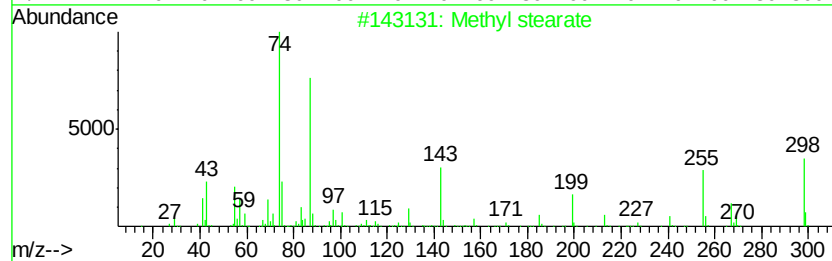
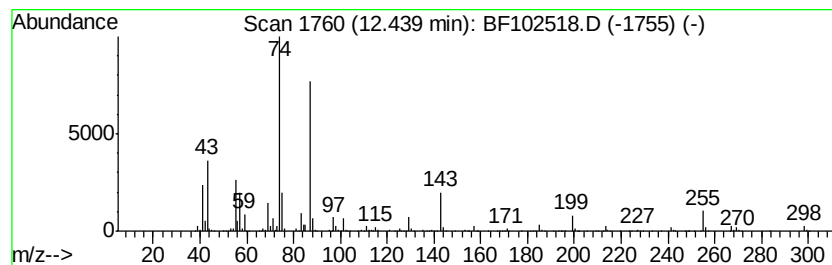
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
12.44	172.30 ng	7259410	Phenanthrene-d10	11.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 98
2		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 94
3		Methyl tetradecanoate	242 C15H30O2	000124-10-7 91
4		Hexadecanoic acid, methyl ester	270 C17H34O2	000112-39-0 87
5		Pentadecanoic acid, methyl ester	256 C16H32O2	007132-64-1 86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102518.D
Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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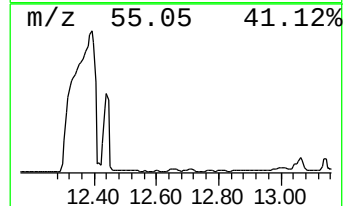
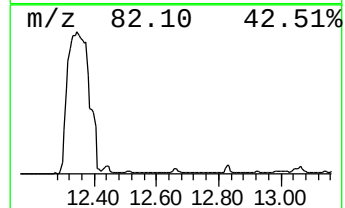
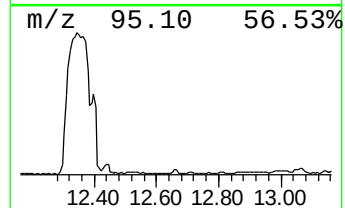
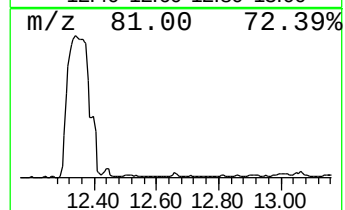
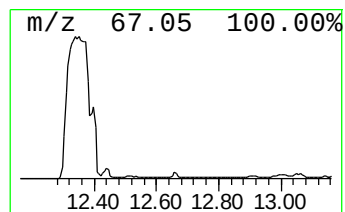
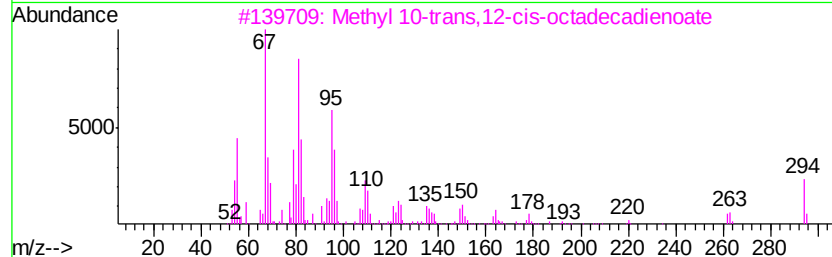
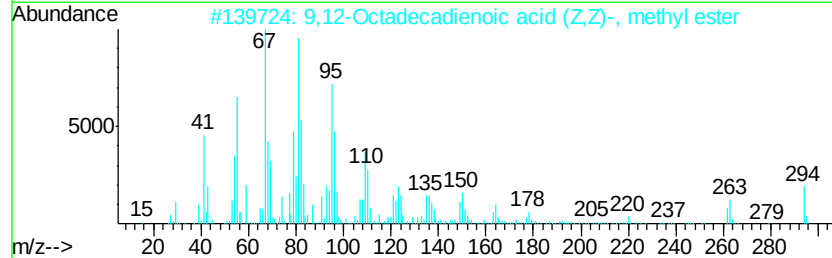
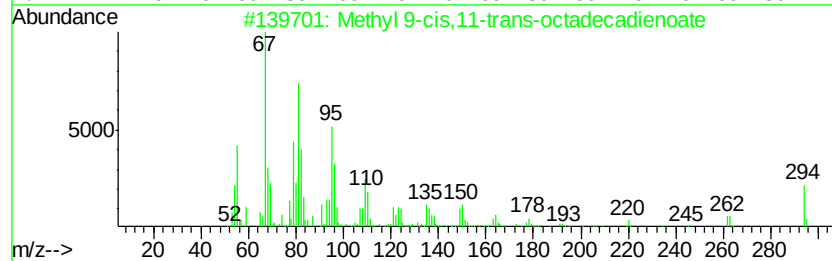
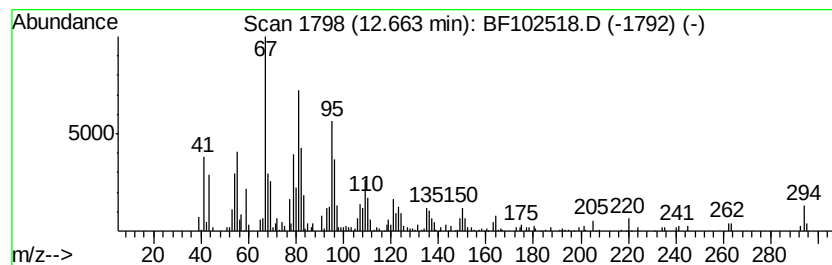
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl 9-cis,11-trans-octad... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	10.32 ng	378534	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99
2		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	97
4		8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	96
5		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102518.D
Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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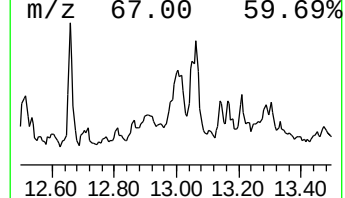
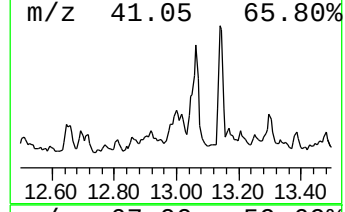
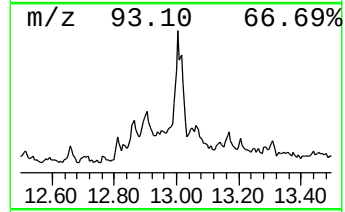
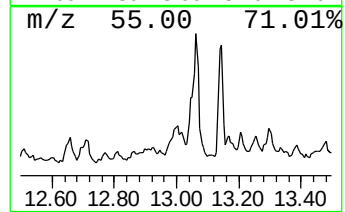
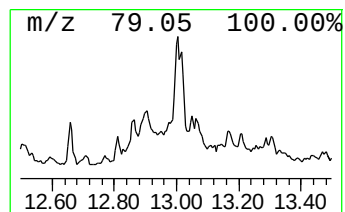
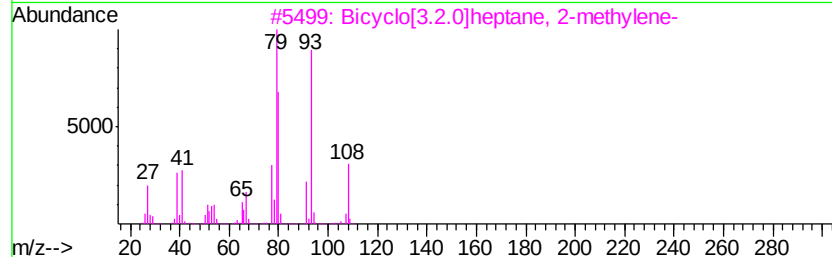
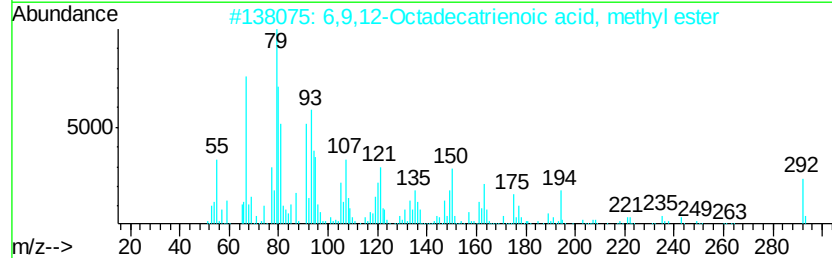
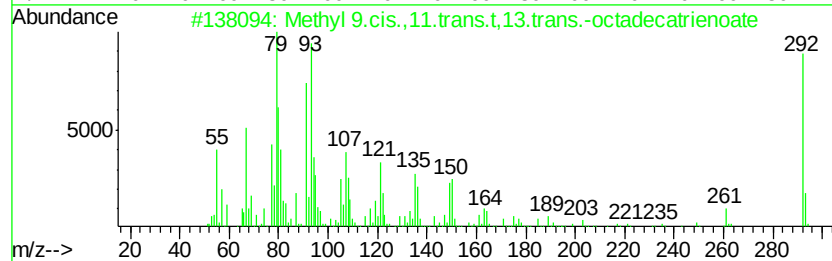
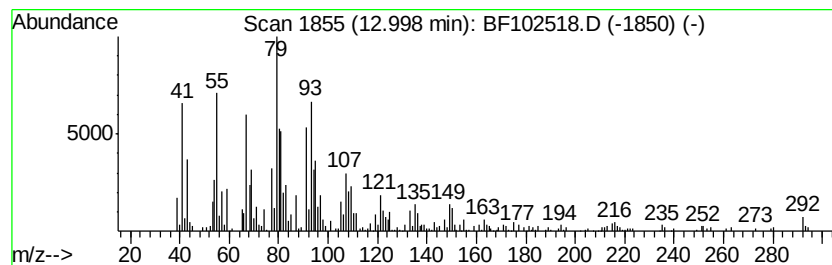
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Methyl 9.cis.,11.trans.t,13... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	11.51 ng	422194	Chrysene-d12	13.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 9.cis.,11.trans.t,13.tran...	292	C19H32O2	1000336-42-6	95
2			6,9,12-Octadecatrienoic acid, me...	292	C19H32O2	002676-41-7	78
3			Bicyclo[3.2.0]heptane, 2-methylene-	108	C8H12	089243-94-7	64
4			Cyclobutene, 4,4-dimethyl-1-(2,7...	190	C14H22	062338-42-5	62
5			Methyl 6-cis,9-cis,11-trans-octa...	292	C19H32O2	1000336-37-7	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102518.D
Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
Sample : J1009-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

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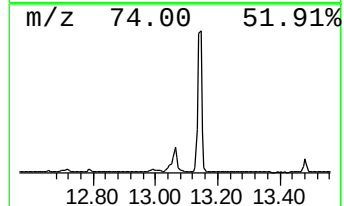
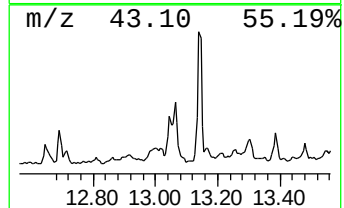
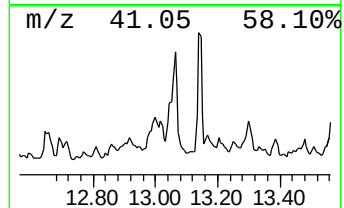
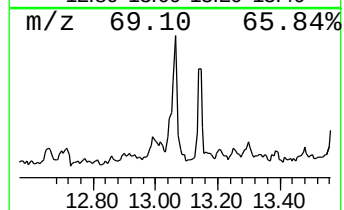
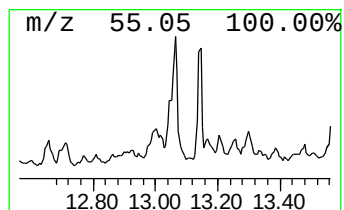
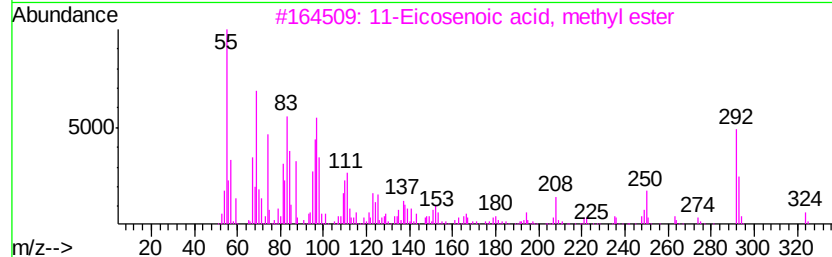
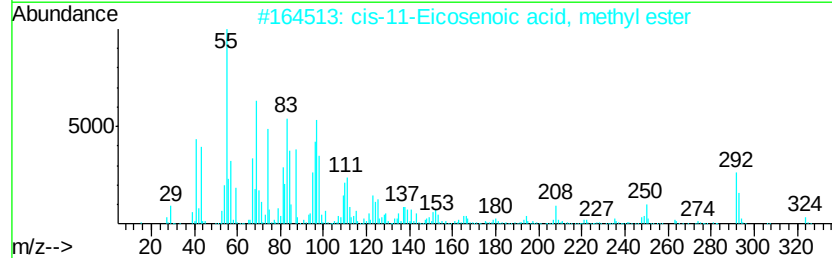
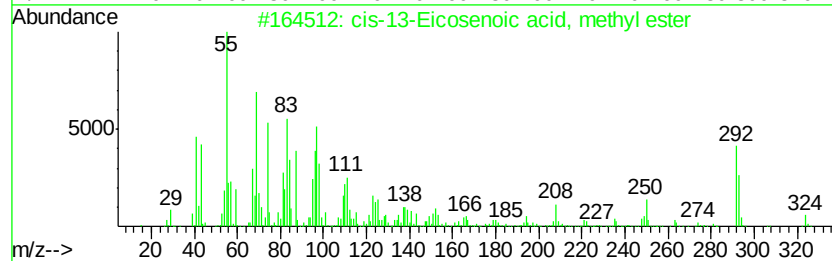
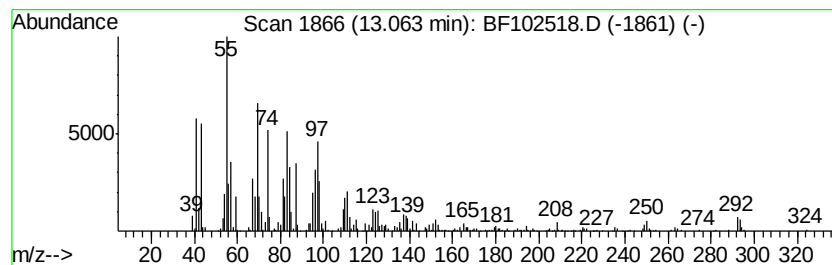
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 cis-13-Eicosenoic acid, met... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	30.67 ng	1125070	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-13-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-52-1	99
2		cis-11-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-63-8	99
3		11-Eicosenoic acid, methyl ester	324	C21H40O2	003946-08-5	98
4		11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	94
5		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102518.D
Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
Sample : J1009-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

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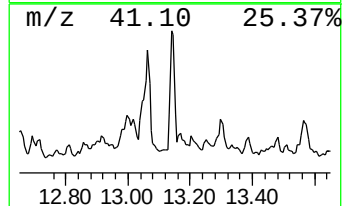
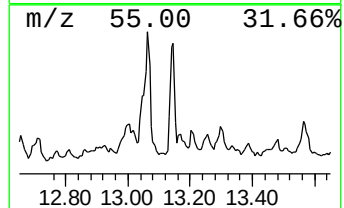
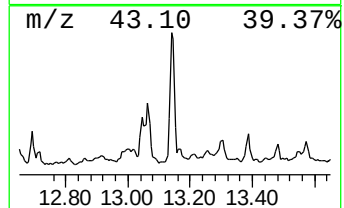
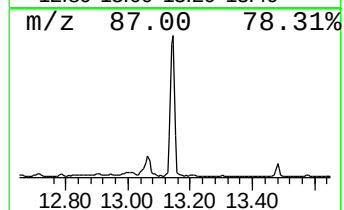
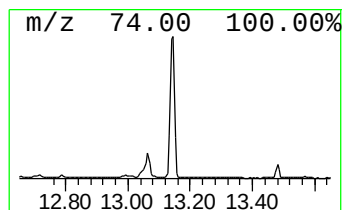
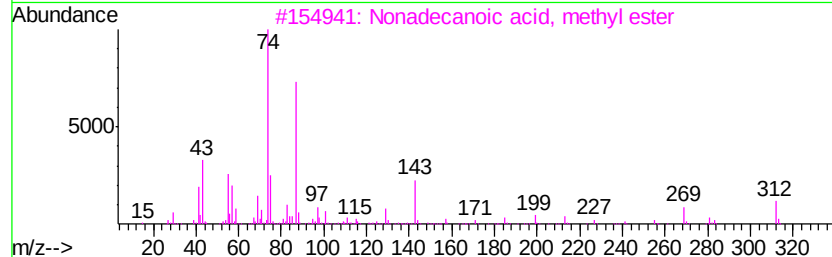
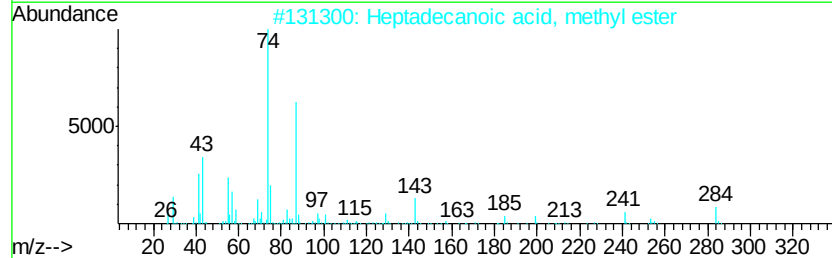
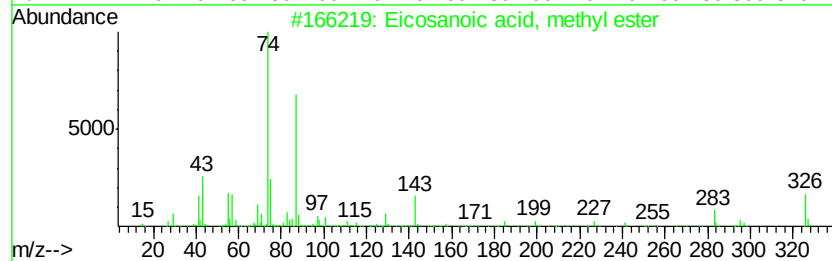
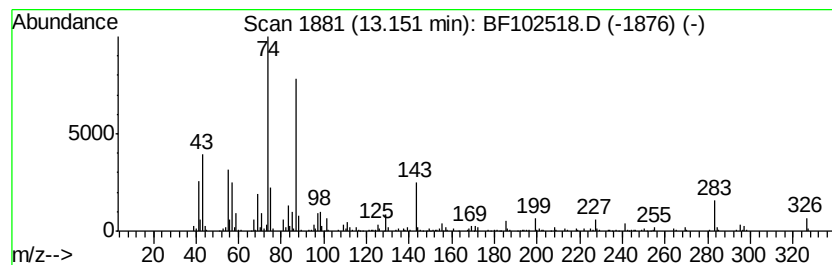
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosanoic acid, methyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	25.46 ng	934064	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	98
2		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	93
3		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	90
4		Methyl stearate	298	C19H38O2	000112-61-8	90
5		Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
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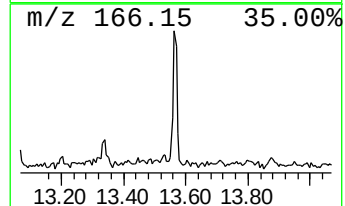
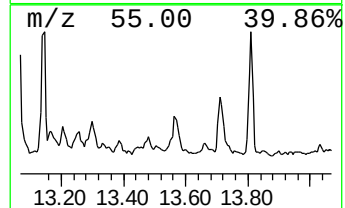
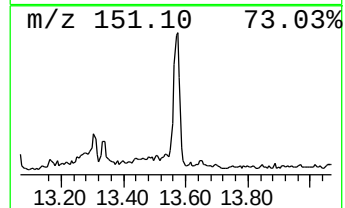
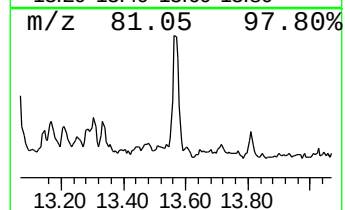
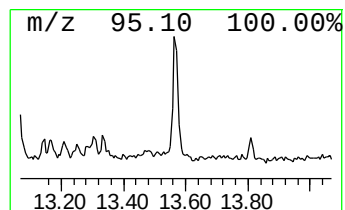
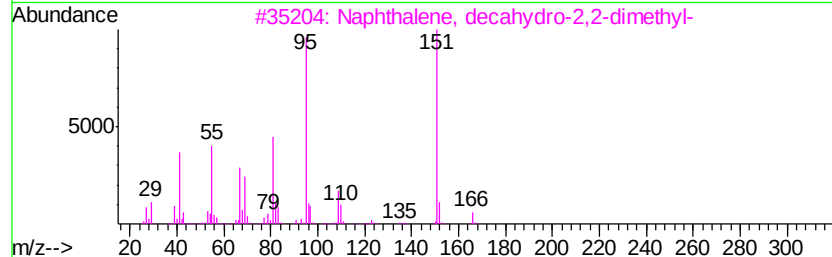
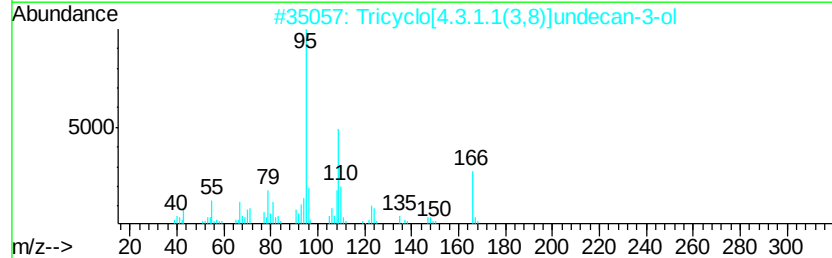
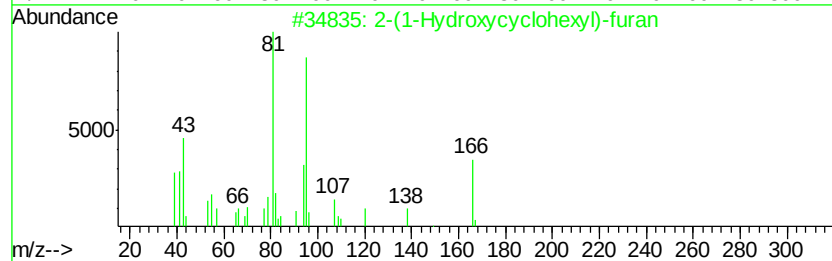
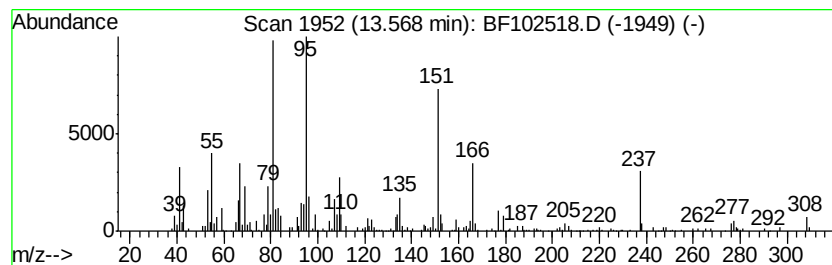
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-(1-Hydroxycyclohexyl)-furan Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.57	12.02 ng	440796	Chrysene-d12	13.88		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-(1-Hydroxycyclohexyl)-furan	166	C10H14O2	036169-67-2	58
2		Tricyclo[4.3.1.1(3,8)]undecan-3-ol	166	C11H18O	014504-80-4	42
3		Naphthalene, decahydro-2,2-dimet...	166	C12H22	031124-85-3	38
4		4-Ethyl-2-hexynal	124	C8H12O	071932-97-3	25
5		1-Chlorosulfonyl-3-methyl-1-azas...	251	C9H14ClN03S	016933-89-4	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
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Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
Sample : J1009-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AU-05-012518-B

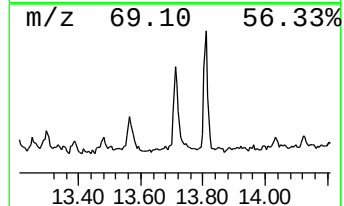
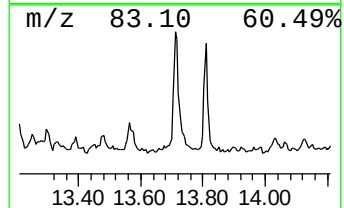
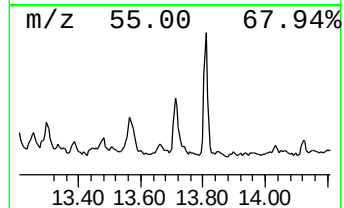
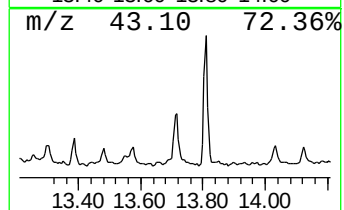
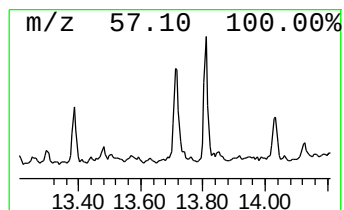
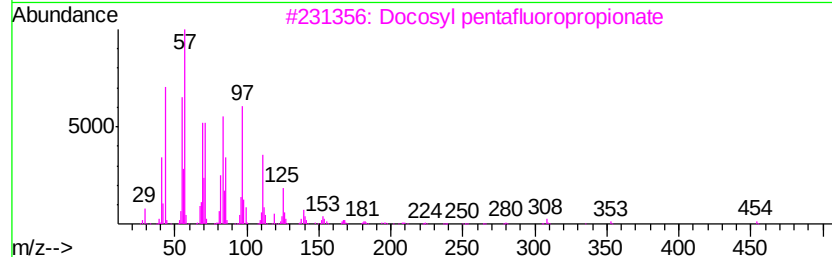
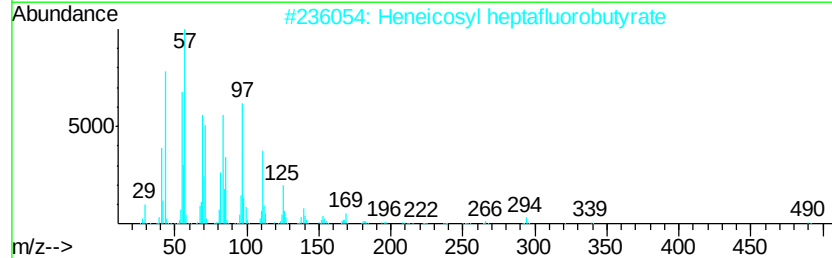
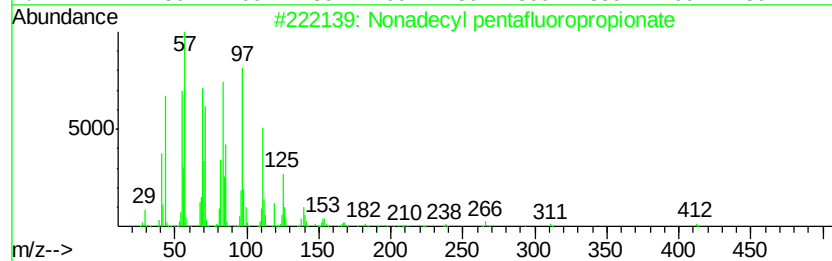
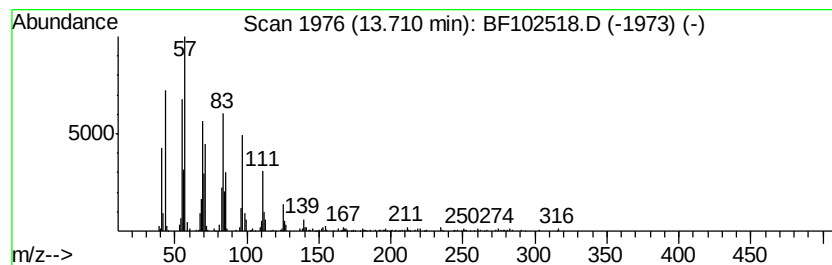
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecyl pentafluoropropio... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	13.10 ng	480755	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
2		Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	91
3		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
4		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102518.D
Acq On : 27 Jan 2018 18:53
Operator : SJ/JU
Sample : J1009-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-012518-B

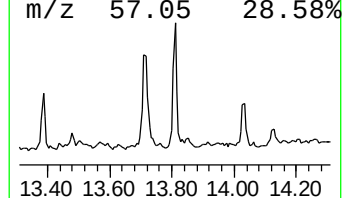
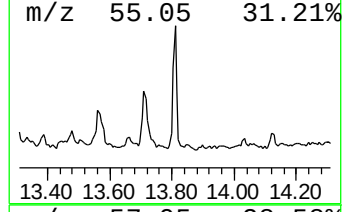
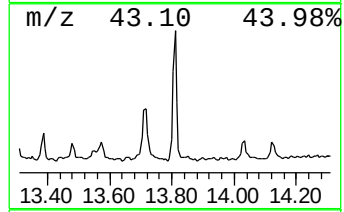
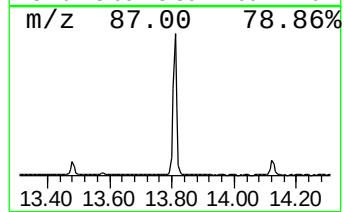
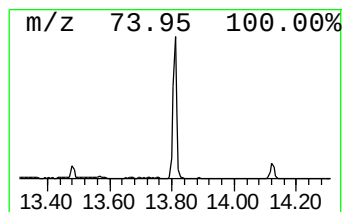
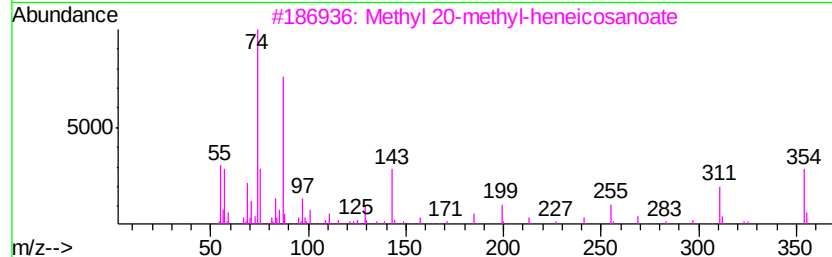
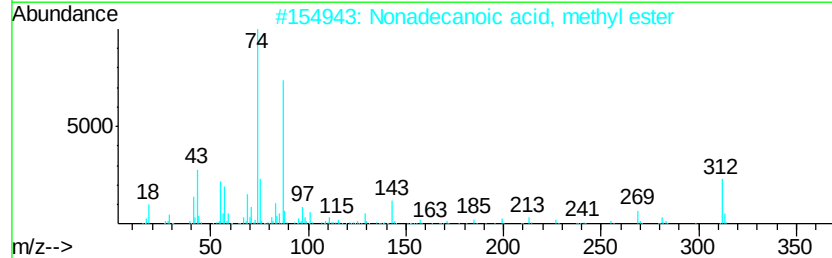
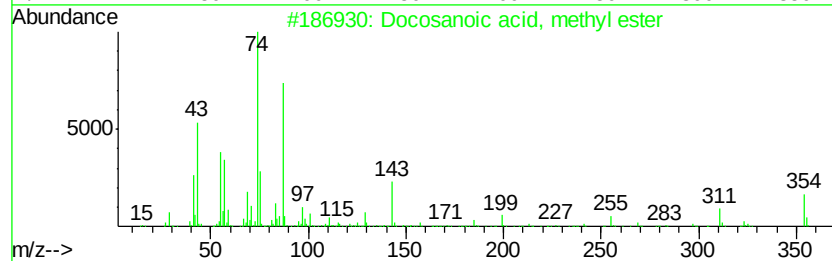
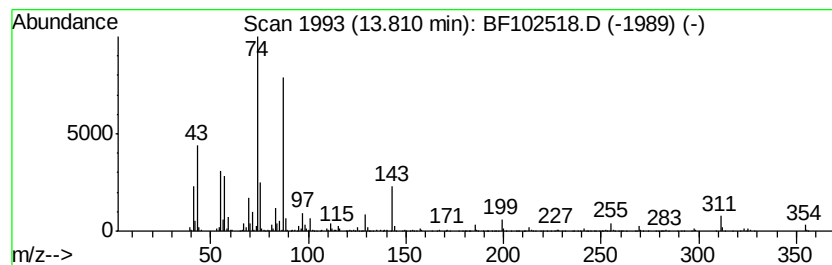
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Docosanoic acid, methyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	24.76 ng	908534	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	99
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	97
3		Methyl 20-methyl-heneicosanoate	354	C23H46O2	1000336-47-4	93
4		Octadecanoic acid, 10-methyl-, m...	312	C20H40O2	002490-19-9	90
5		Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102518.D
 Acq On : 27 Jan 2018 18:53
 Operator : SJ/JU
 Sample : J1009-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.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
unknown6.49	6.49	86.1	ng	3717310	1	6.72	863380	20.0
Hexadecane	10.17	7.2	ng	386807	3	9.75	1077120	20.0
Heptadecane	10.65	12.7	ng	533894	4	11.23	842635	20.0
Hexadecanoic acid...	11.65	331.4	ng	13960200	4	11.23	842635	20.0
Methyl Z-11-tetra...	11.95	11.2	ng	470675	4	11.23	842635	20.0
10,13-Octadecadie...	12.37	888.6	ng	37438000	4	11.23	842635	20.0
Methyl stearate	12.44	172.3	ng	7259410	4	11.23	842635	20.0
Methyl 9-cis,11-t...	12.66	10.3	ng	378534	5	13.88	733737	20.0
Methyl 9.cis.,11....	13.00	11.5	ng	422194	5	13.88	733737	20.0
cis-13-Eicosenoic...	13.06	30.7	ng	1125070	5	13.88	733737	20.0
Eicosanoic acid, ...	13.15	25.5	ng	934064	5	13.88	733737	20.0
2-(1-Hydroxycyclo...	13.57	12.0	ng	440796	5	13.88	733737	20.0
Nonadecyl pentafl...	13.71	13.1	ng	480755	5	13.88	733737	20.0
Docosanoic acid, ...	13.81	24.8	ng	908534	5	13.88	733737	20.0