

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101218\
 Data File : BF109846.D
 Acq On : 12 Oct 2018 20:56
 Operator : JU/SJ
 Sample : J5205-03 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-101118

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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	5.310	545	548	561	rBV	79417	159680	2.07%	0.565%
2	6.345	720	724	739	rBV	88281	237903	3.08%	0.842%
3	6.463	741	744	758	rVV	211990	291201	3.77%	1.031%
4	6.692	780	783	806	rBV	393842	504606	6.53%	1.786%
5	6.845	806	809	832	rVB	209256	271687	3.51%	0.962%
6	7.263	877	880	902	rBV	43637	161684	2.09%	0.572%
7	7.975	997	1001	1031	rBV	585615	693169	8.96%	2.453%
8	9.045	1180	1183	1197	rBV	290423	351084	4.54%	1.242%
9	9.722	1294	1298	1305	rBV	737046	701618	9.07%	2.483%
10	10.510	1429	1432	1440	rBV	103946	137522	1.78%	0.487%
11	10.733	1467	1470	1479	rBV	323256	264040	3.41%	0.934%
12	11.198	1547	1549	1552	rBV	465014	477185	6.17%	1.689%
13	11.222	1552	1553	1561	rVB	179187	164094	2.12%	0.581%
14	11.522	1598	1604	1609	rBV	461710	462218	5.98%	1.636%
15	11.610	1614	1619	1622	rBV	4192163	4094260	52.95%	14.490%
16	11.739	1637	1641	1647	rBV2	70327	83330	1.08%	0.295%
17	12.004	1682	1686	1689	rBV	106046	95543	1.24%	0.338%
18	12.292	1729	1735	1737	rBV	4812145	6431495	83.18%	22.761%
19	12.322	1737	1740	1745	rVV	6478840	7732350	100.00%	27.365%
20	12.357	1745	1746	1749	rVV	185720	134910	1.74%	0.477%
21	12.392	1749	1752	1759	rVV2	2336828	2196495	28.41%	7.773%
22	12.633	1788	1793	1797	rBV	206555	299722	3.88%	1.061%
23	12.780	1814	1818	1824	rBV	234945	240088	3.10%	0.850%
24	12.969	1844	1850	1855	rBV3	44135	111078	1.44%	0.393%
25	13.033	1855	1861	1866	rBV3	179794	260435	3.37%	0.922%
26	13.110	1871	1874	1881	rVB	104249	102089	1.32%	0.361%
27	13.774	1984	1987	1989	rBV	79871	79089	1.02%	0.280%
28	13.827	1991	1996	1999	rBV2	501267	639445	8.27%	2.263%
29	14.833	2164	2167	2176	rBV	92127	248372	3.21%	0.879%
30	15.174	2222	2225	2229	rBV2	64487	97744	1.26%	0.346%
31	15.227	2231	2234	2247	rVB	296247	532209	6.88%	1.884%

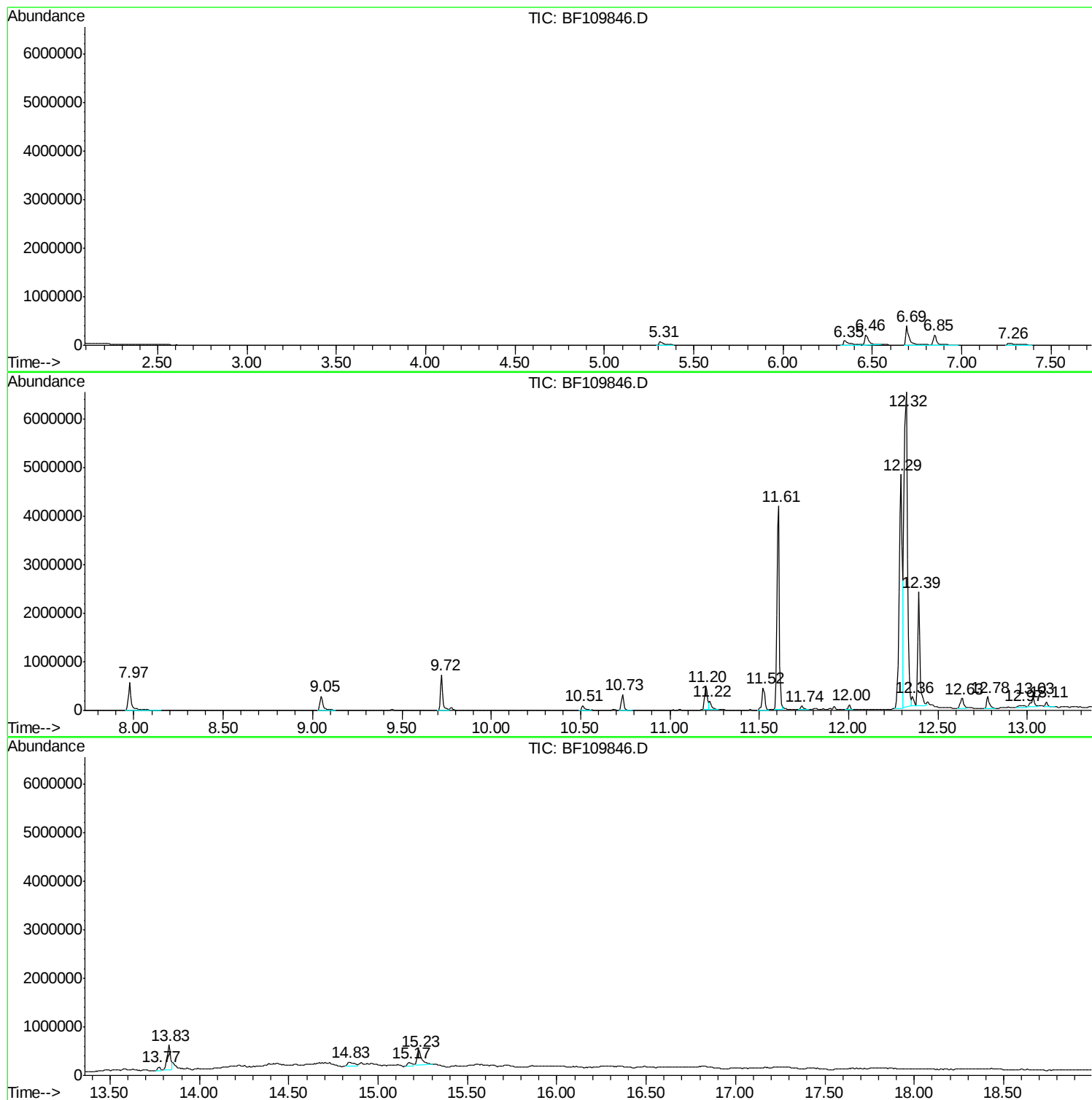
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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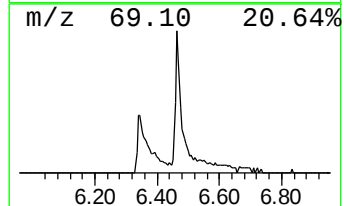
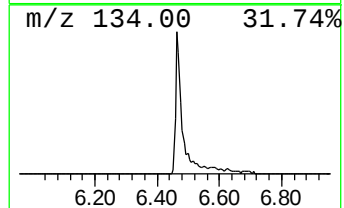
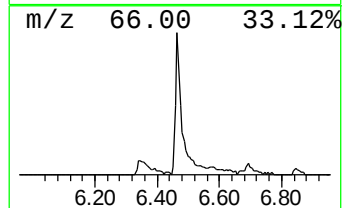
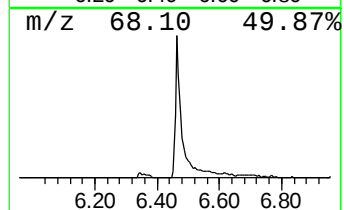
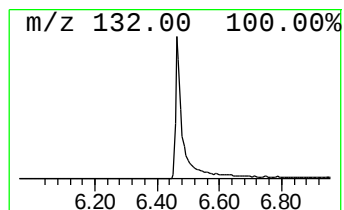
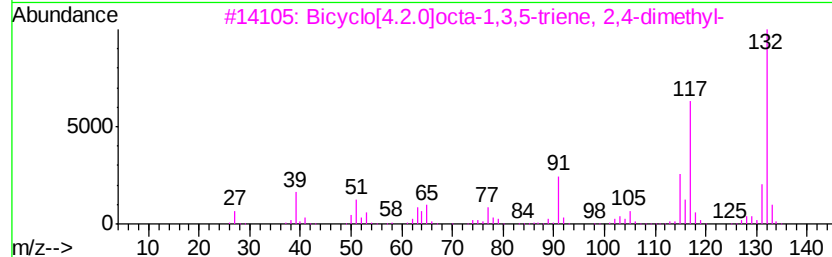
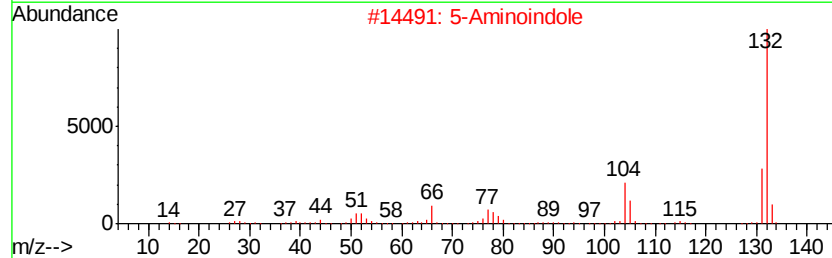
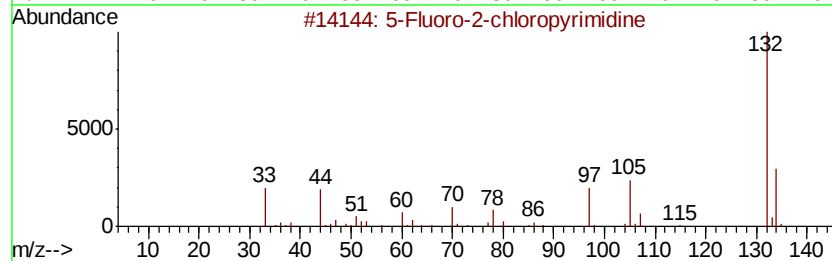
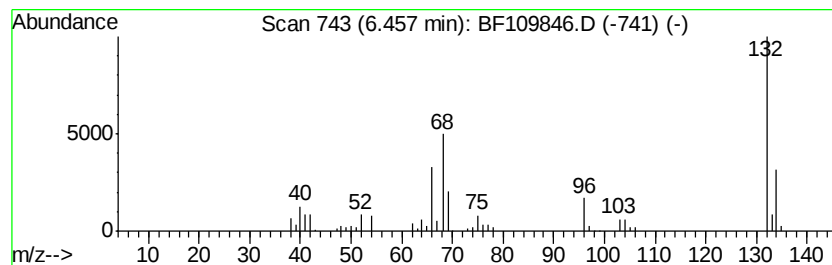
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Peak Number 1 unknown6.46 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	11.54 ng	291201	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Xenon	132	Xe	007440-63-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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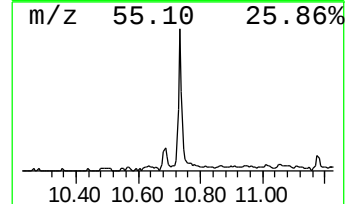
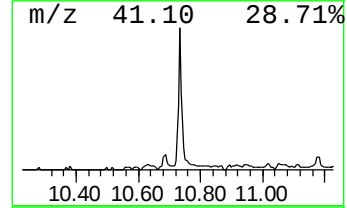
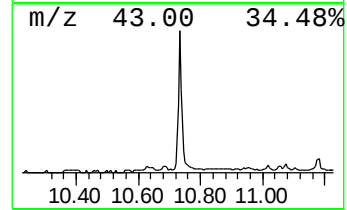
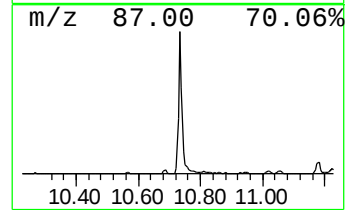
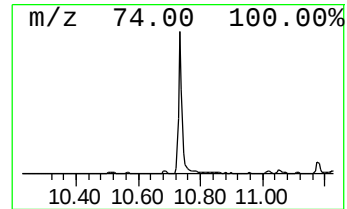
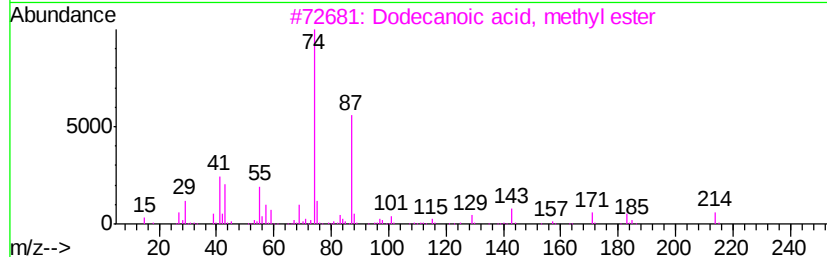
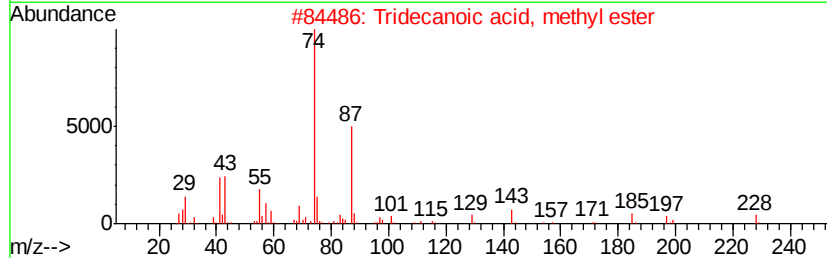
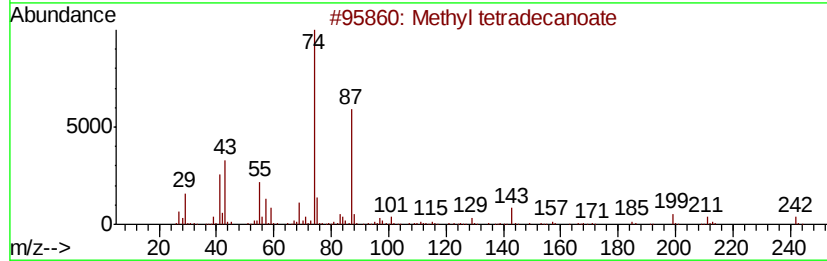
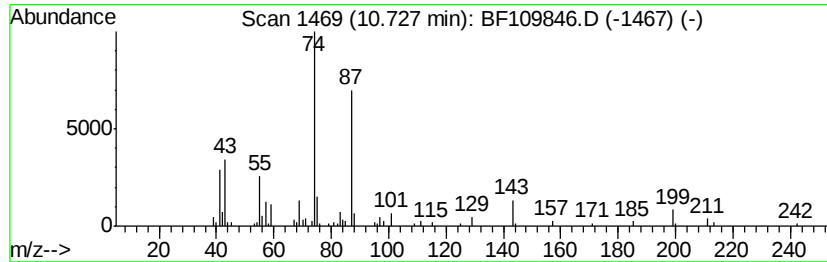
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Peak Number 3 Methyl tetradecanoate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	11.07 ng	264040	Phenanthrene-d10	11.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
2		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	86
3		Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	86
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	86
5		Undecanoic acid, 11-bromo-, meth...	278	C12H23BrO2	006287-90-7	78



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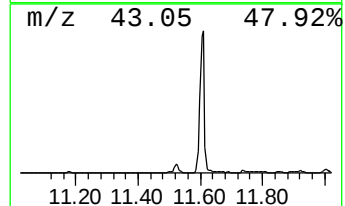
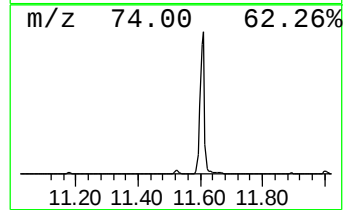
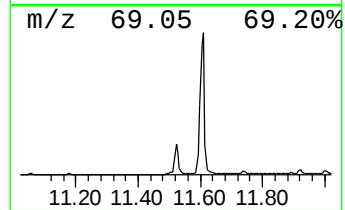
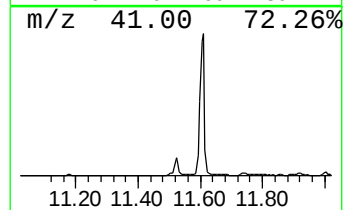
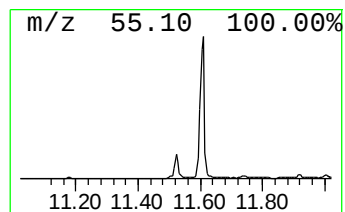
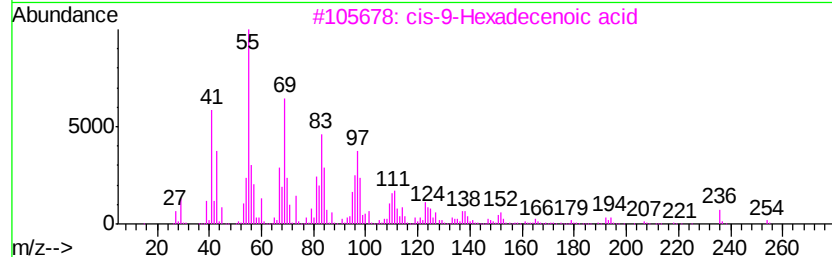
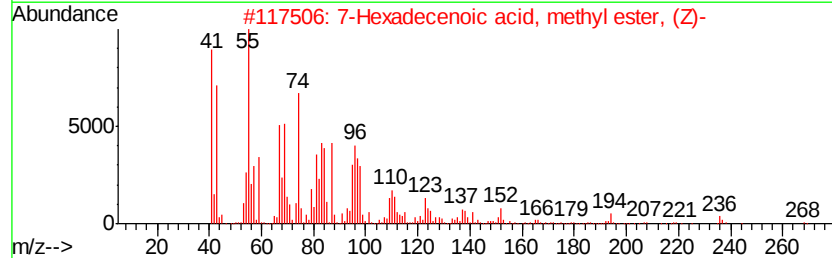
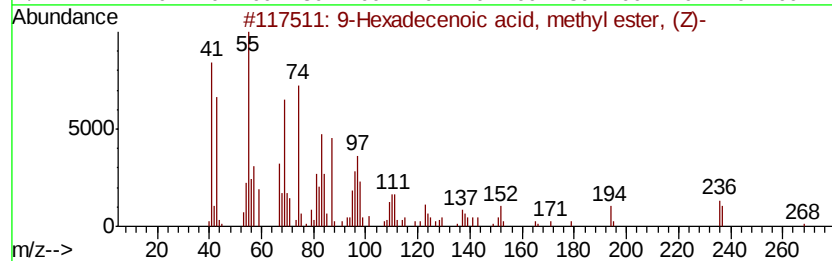
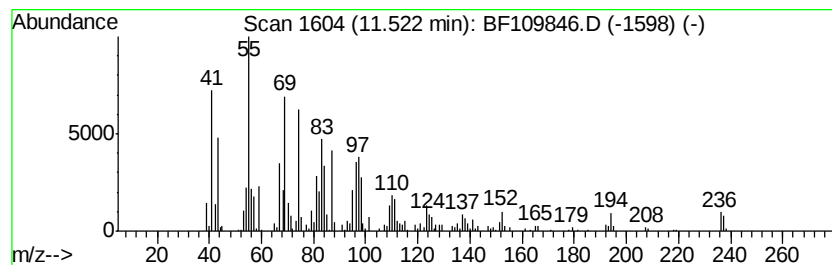
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Peak Number 4 9-Hexadecenoic acid, methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.52	19.37 ng	462218	Phenanthrene-d10	11.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	91
2		7-Hexadecenoic acid, methyl este...	268	C17H32O2	056875-67-3	87
3		cis-9-Hexadecenoic acid	254	C16H30O2	1000333-19-5	53
4		Chloromethyl 7-chlorododecanoate	282	C13H24Cl2O2	080419-03-0	45
5		13-Borabicyclo[7.3.0]tridecane, ...	236	C15H29BO	1000156-41-7	30



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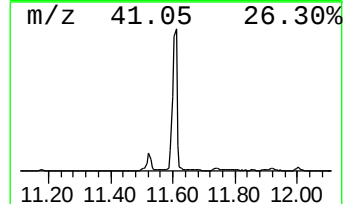
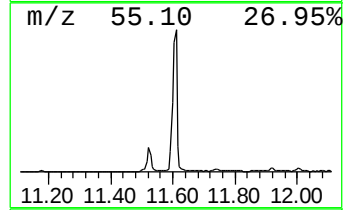
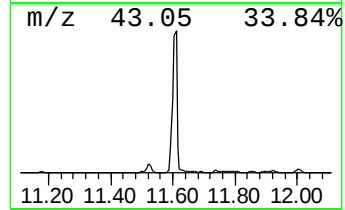
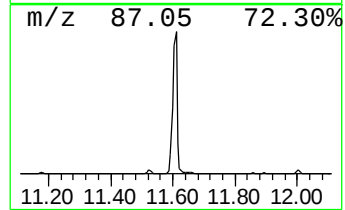
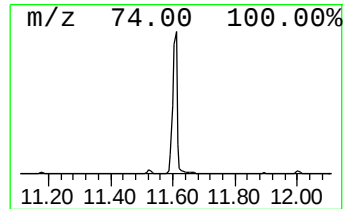
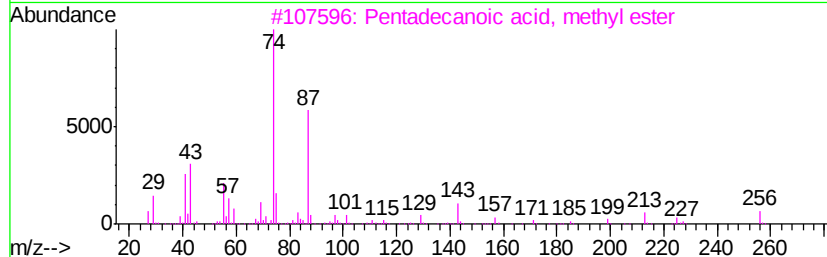
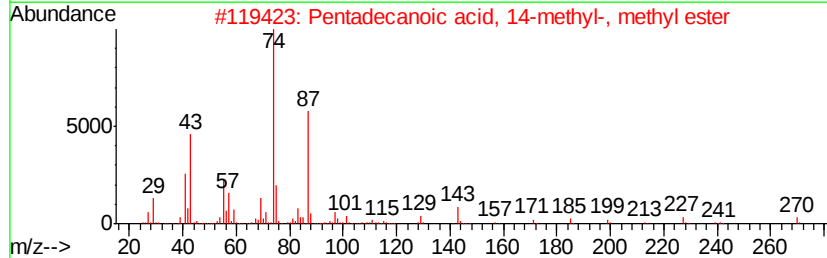
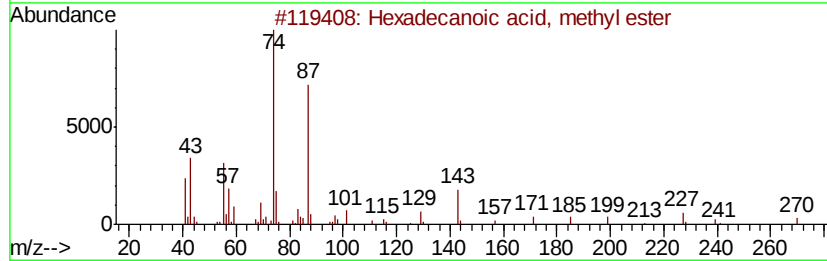
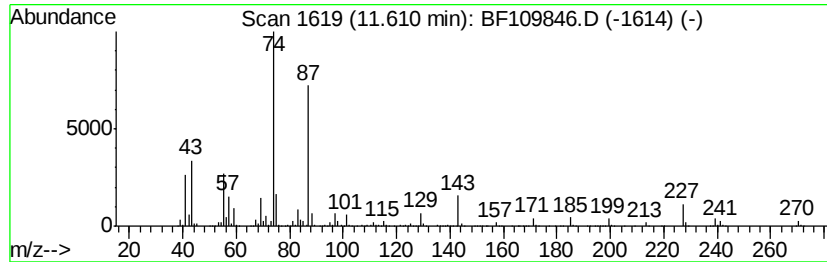
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Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	171.60 ng	4094260	Phenanthrene-d10	11.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4		Methyl tetradecanoate	242	C15H30O2	000124-10-7	83
5		Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	80



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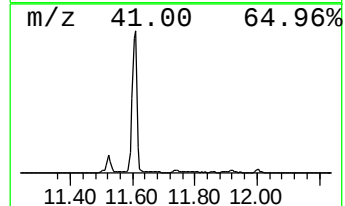
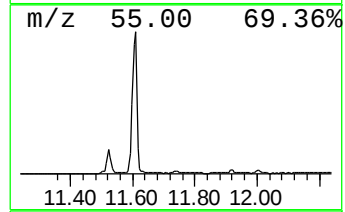
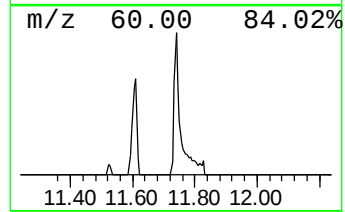
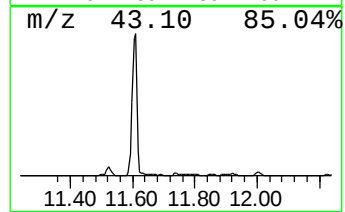
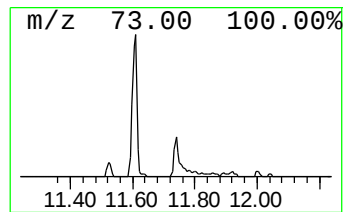
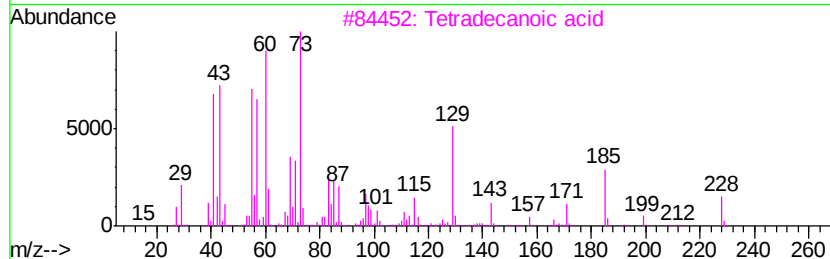
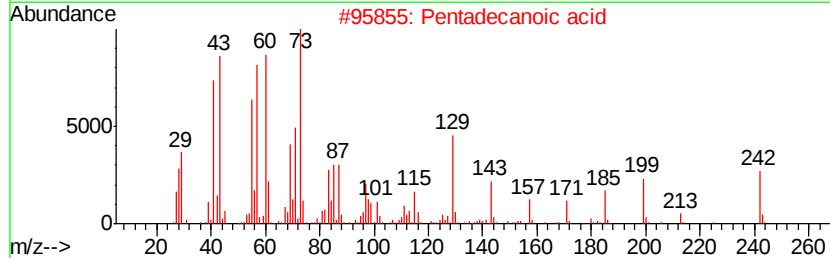
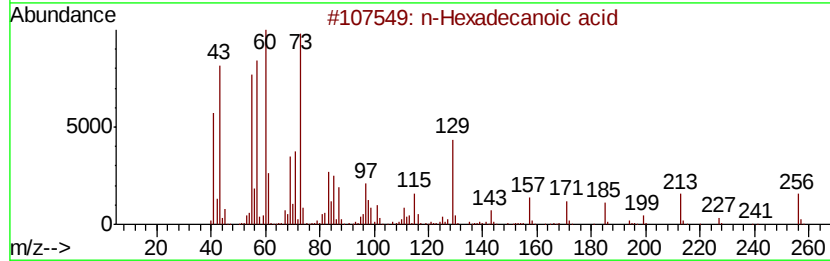
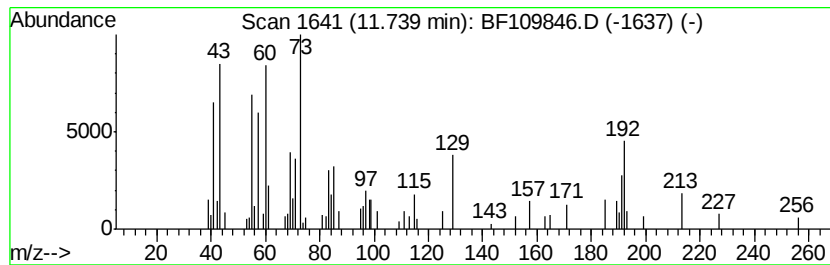
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.74	3.49 ng	83330	Phenanthrene-d10	11.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
4	Tridecanoic acid	214	C13H26O2	000638-53-9	50
5	n-Decanoic acid	172	C10H20O2	000334-48-5	43



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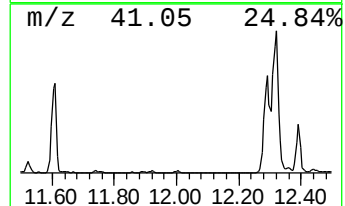
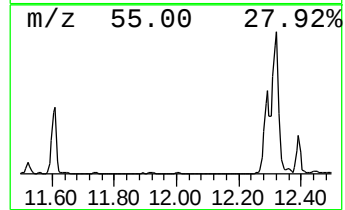
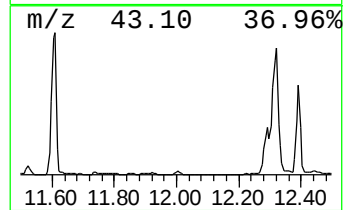
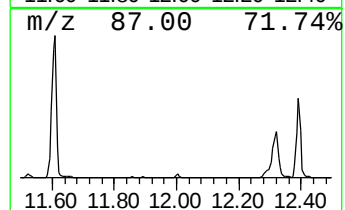
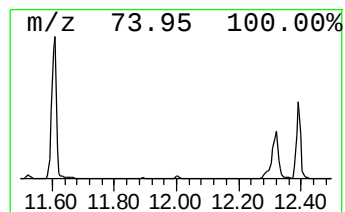
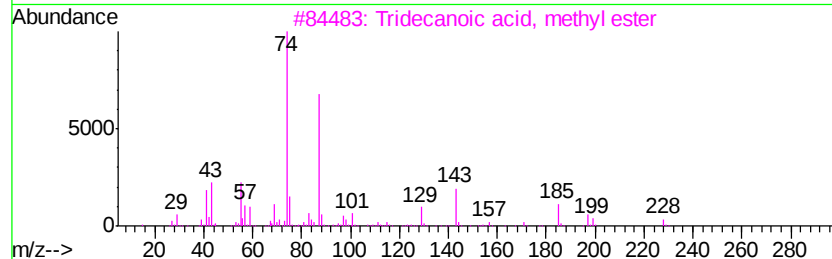
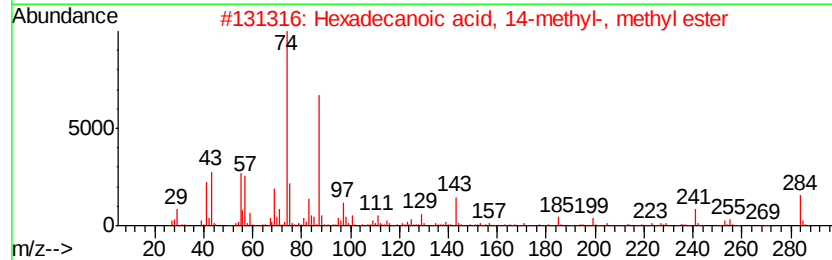
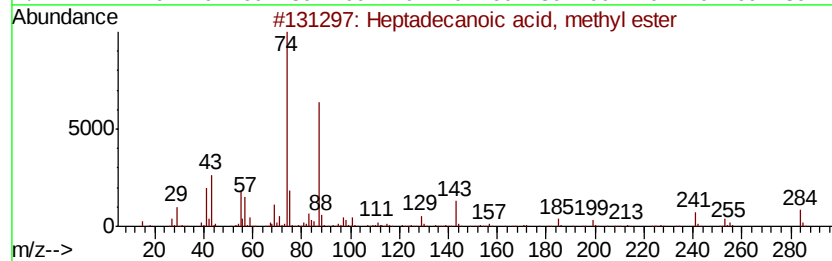
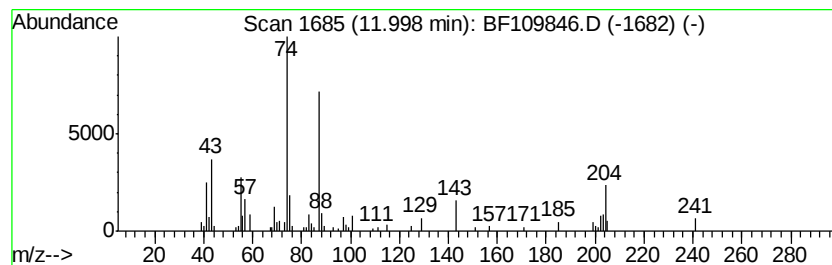
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecanoic acid, methyl ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.00	4.00 ng	95543	Phenanthrene-d10	11.20		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	94
2		Hexadecanoic acid, 14-methyl-, m...	284	C18H36O2	002490-49-5	87
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	74
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	74
5		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
Data File : BF109846.D
Acq On : 12 Oct 2018 20:56
Operator : JU/SJ
Sample : J5205-03 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-04-101118

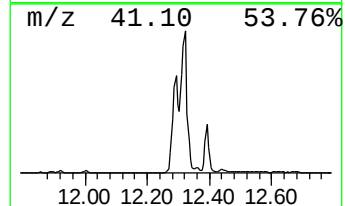
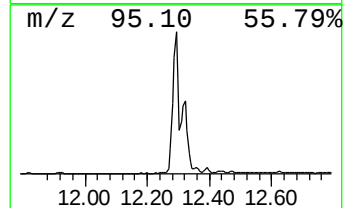
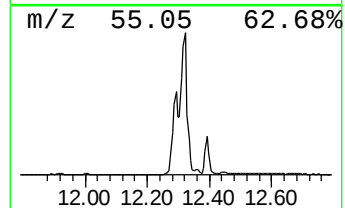
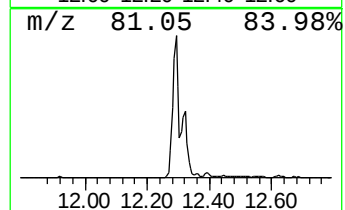
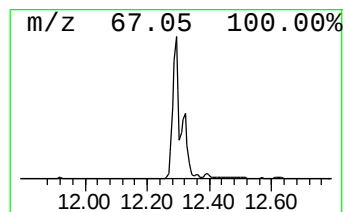
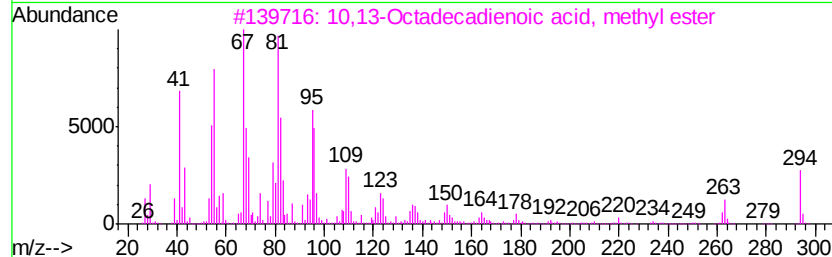
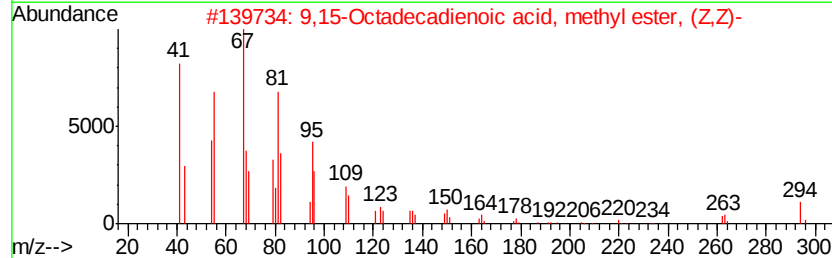
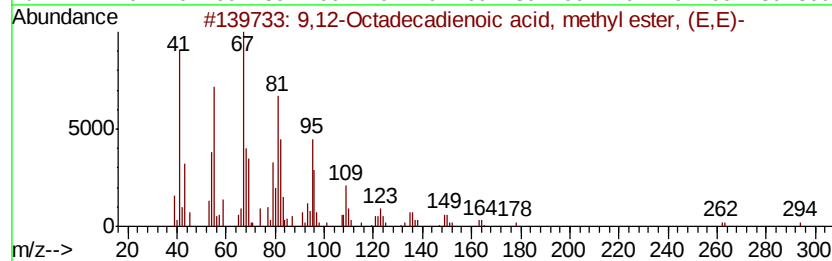
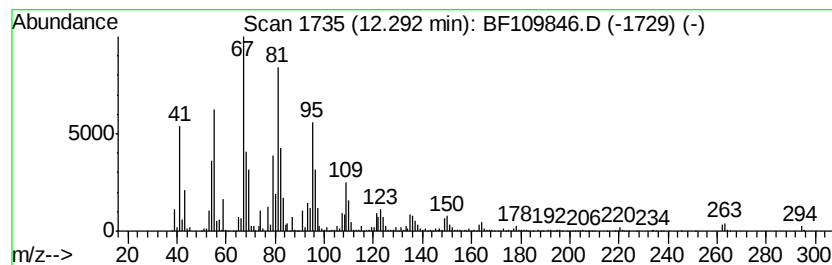
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	269.56 ng	6431500	Phenanthrene-d10	11.20

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	98
4		11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	98
5		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
Data File : BF109846.D
Acq On : 12 Oct 2018 20:56
Operator : JU/SJ
Sample : J5205-03 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

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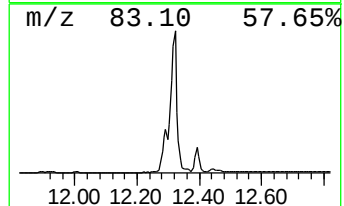
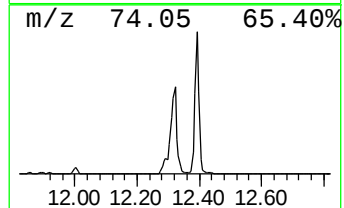
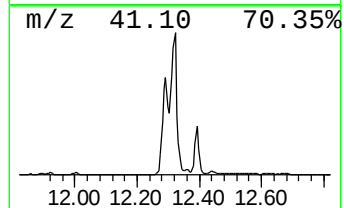
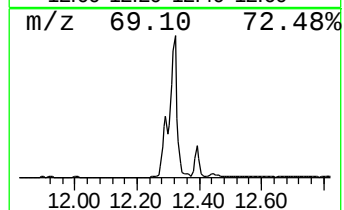
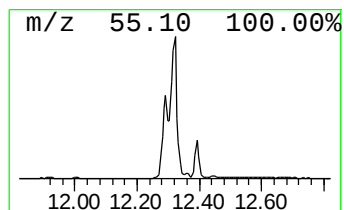
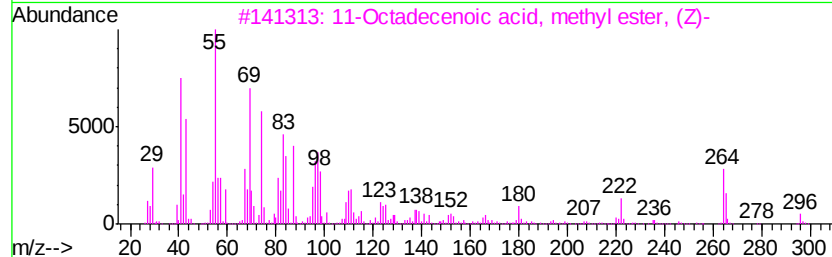
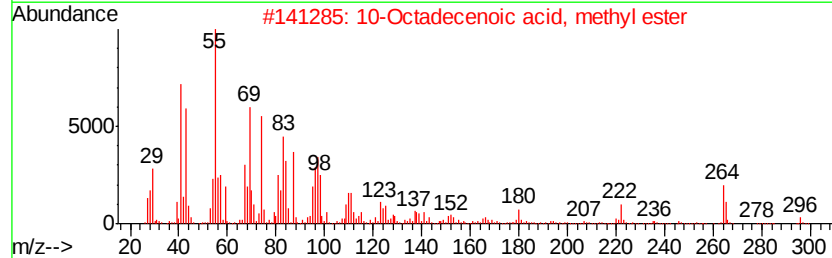
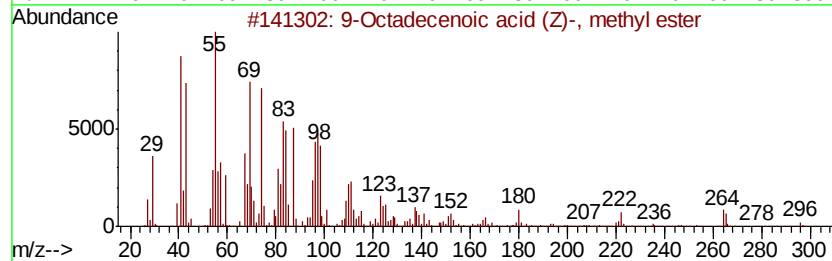
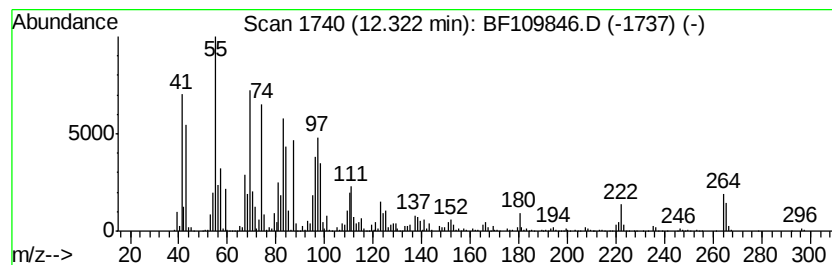
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	324.08 ng	7732350	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
3		11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99
4		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
5		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
Data File : BF109846.D
Acq On : 12 Oct 2018 20:56
Operator : JU/SJ
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Misc :
ALS Vial : 22 Sample Multiplier: 1

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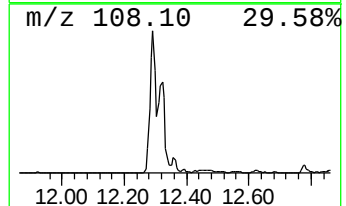
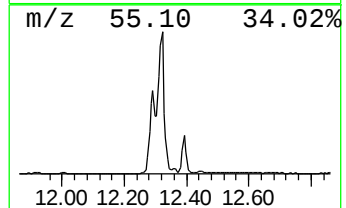
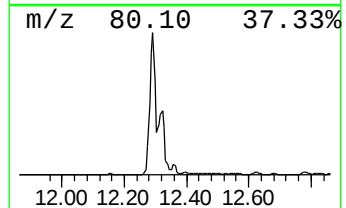
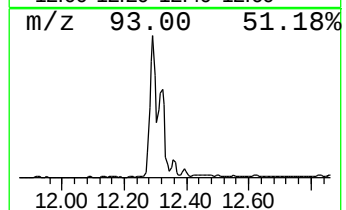
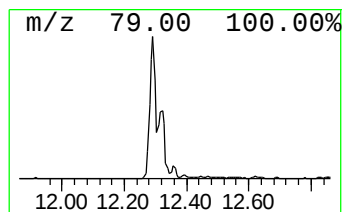
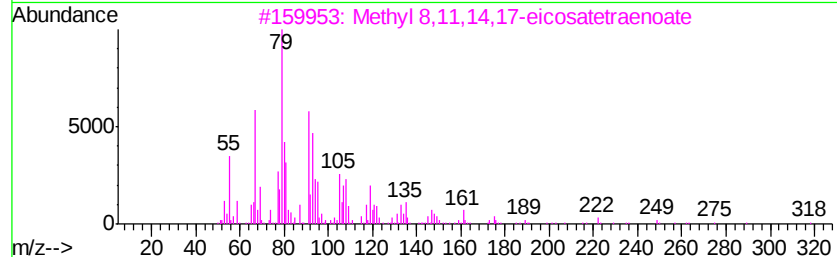
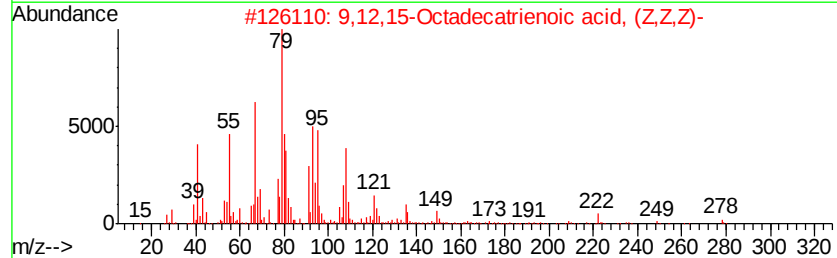
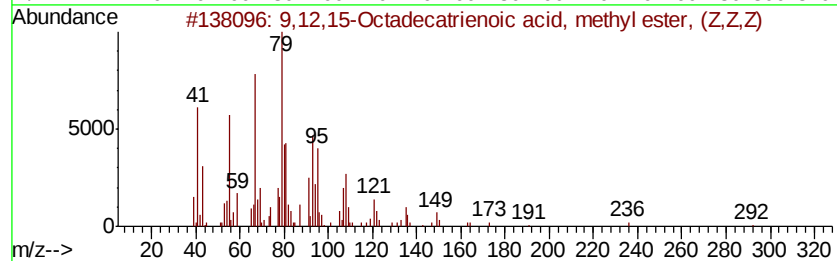
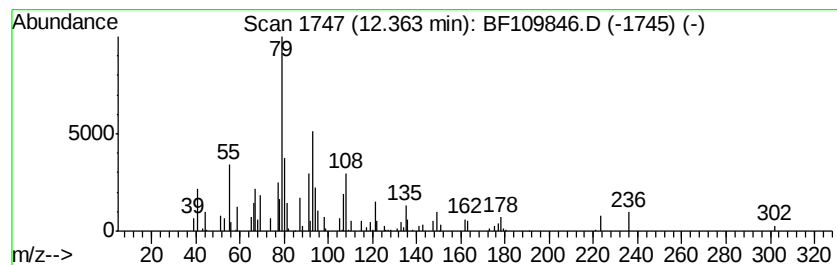
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,12,15-Octadecatrienoic ac... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	5.65 ng	134910	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12,15-Octadecatrienoic acid, m...	292	C19H32O2	000301-00-8	91		
2	9,12,15-Octadecatrienoic acid, (...	278	C18H30O2	000463-40-1	86		
3	Methyl 8,11,14,17-eicosatetraenoate	318	C21H34O2	1000336-47-0	83		
4	1,3-Cyclooctadiene	108	C8H12	001700-10-3	70		
5	3-Tetradecen-5-yne, (Z)-	192	C14H24	074663-68-6	64		



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
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Acq On : 12 Oct 2018 20:56
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Misc :
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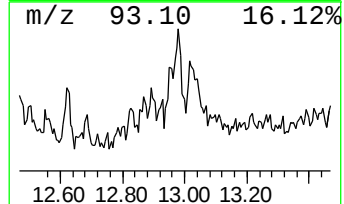
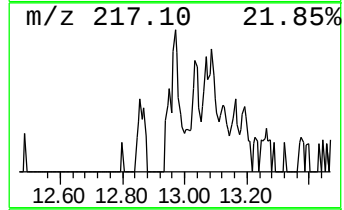
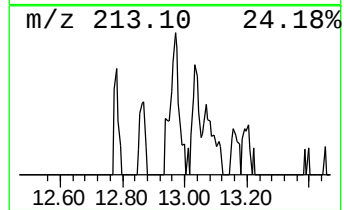
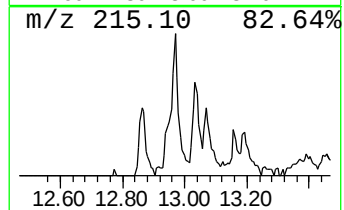
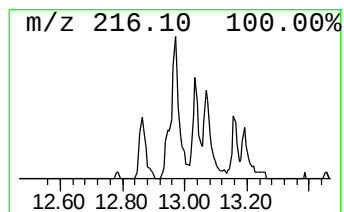
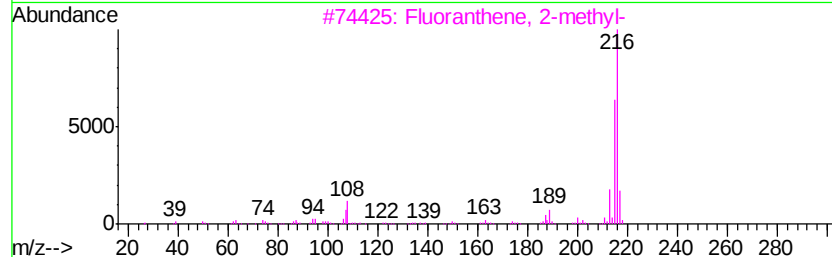
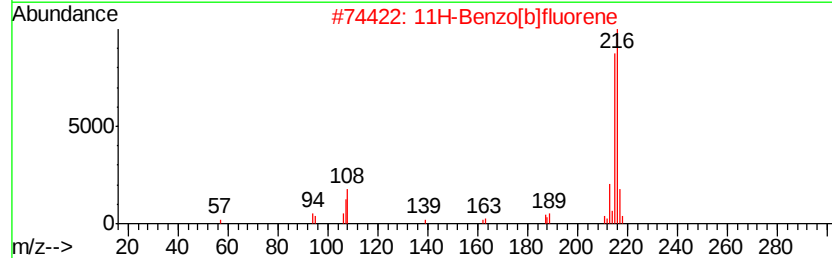
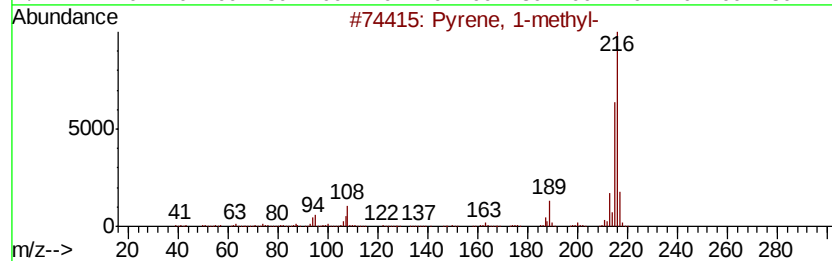
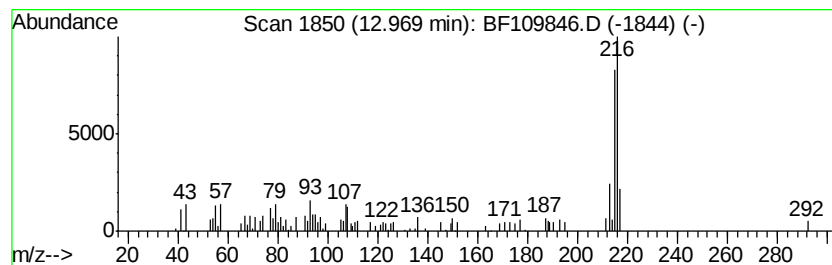
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	3.47 ng	111078	Chrysene-d12	13.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
Data File : BF109846.D
Acq On : 12 Oct 2018 20:56
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Misc :
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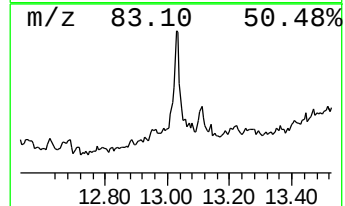
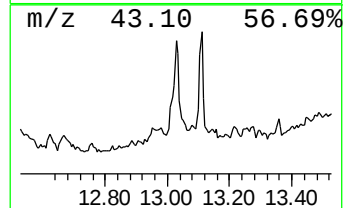
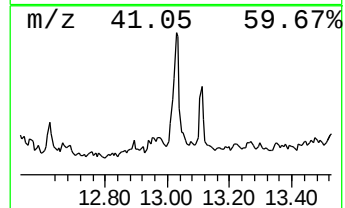
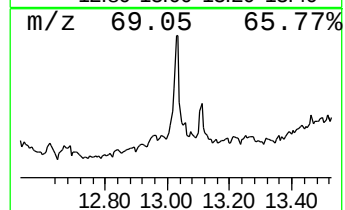
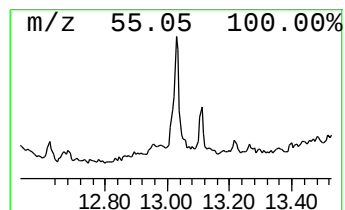
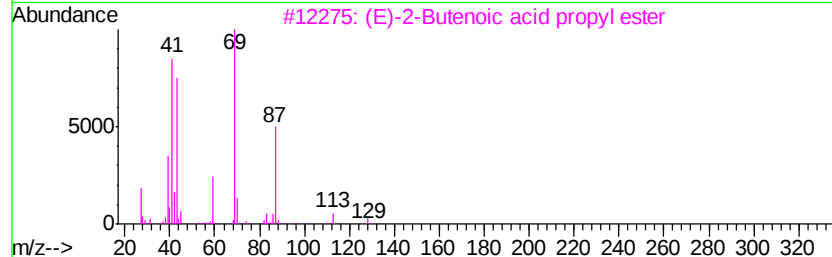
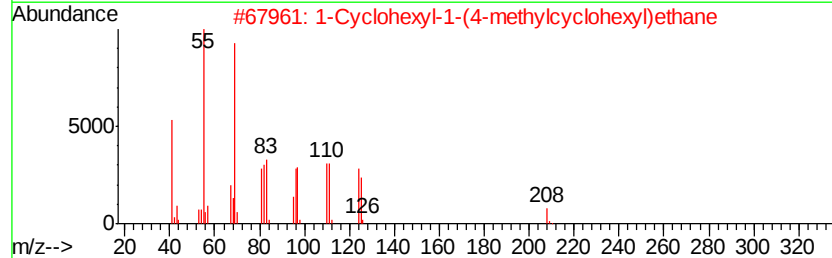
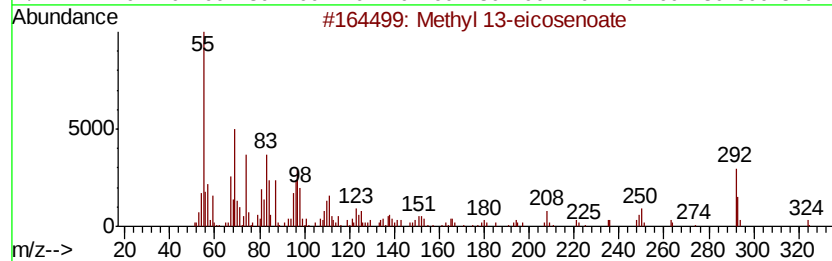
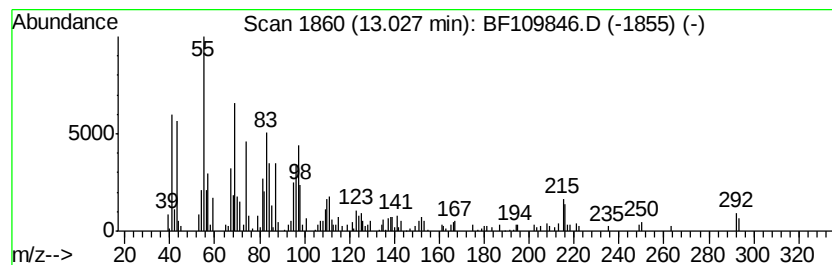
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown13.03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	8.15 ng	260435	Chrysene-d12	13.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 13-eicosenoate	324	C21H40O2	1000336-48-4	49
2		1-Cyclohexyl-1-(4-methylcyclohexyl)ethane	208	C15H28	002027-12-5	43
3		(E)-2-Butenoic acid propyl ester	128	C7H12O2	010352-87-1	42
4		Cyclohexane, 1-(cyclohexylmethyl)-	208	C15H28	054934-92-8	42
5		Methyl 9-heptadecenoate or 9-17:1	282	C18H34O2	1000336-38-0	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
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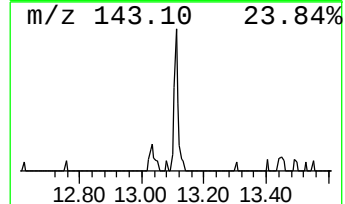
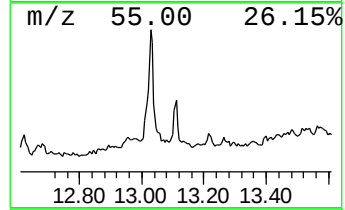
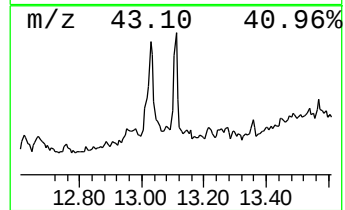
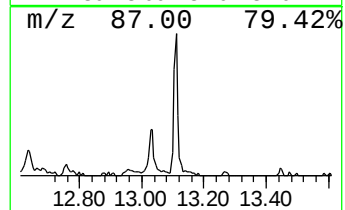
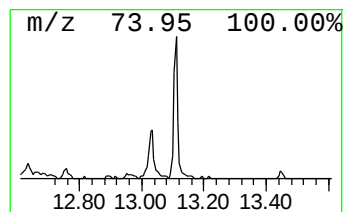
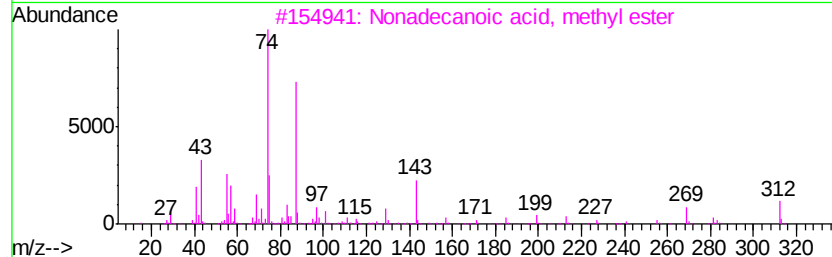
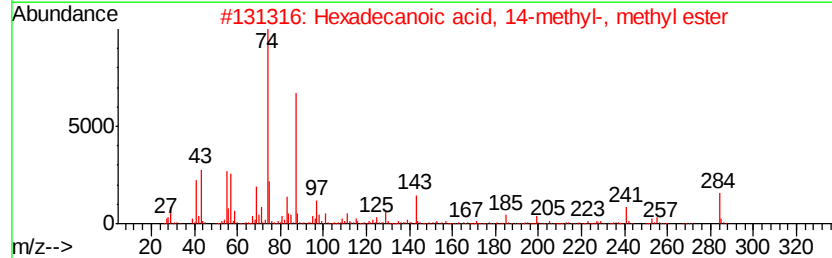
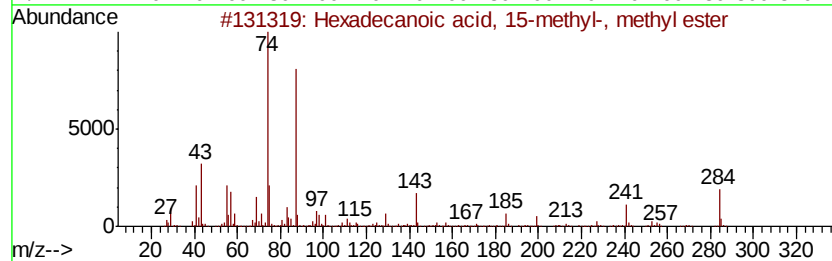
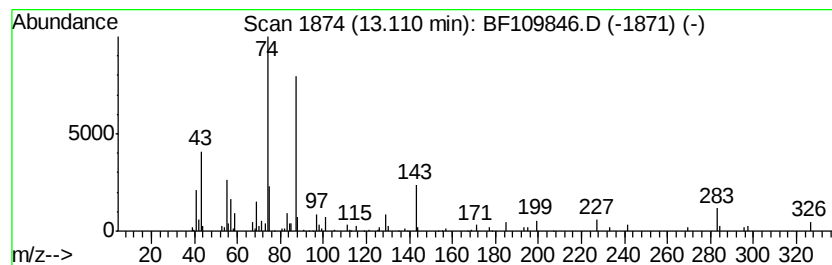
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecanoic acid, 15-methy... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.11	3.19 ng	102089	Chrysene-d12		13.83
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	91
2	Hexadecanoic acid, 14-methyl-, m...	284	C18H36O2	002490-49-5	90
3	Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	87
4	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	86
5	Methyl stearate	298	C19H38O2	000112-61-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101218\
 Data File : BF109846.D
 Acq On : 12 Oct 2018 20:56
 Operator : JU/SJ
 Sample : J5205-03 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-101118

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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
unknown6.46	6.46	11.5	ng	291201	1	6.69	504606	20.0
Methyl tetradecan...	10.73	11.1	ng	264040	4	11.20	477185	20.0
9-Hexadecenoic ac...	11.52	19.4	ng	462218	4	11.20	477185	20.0
Hexadecanoic acid...	11.61	171.6	ng	4094260	4	11.20	477185	20.0
n-Hexadecanoic acid	11.74	3.5	ng	83330	4	11.20	477185	20.0
Heptadecanoic aci...	12.00	4.0	ng	95543	4	11.20	477185	20.0
9,12-Octadecadien...	12.29	269.6	ng	6431500	4	11.20	477185	20.0
9-Octadecenoic ac...	12.32	324.1	ng	7732350	4	11.20	477185	20.0
9,12,15-Octadecat...	12.36	5.7	ng	134910	4	11.20	477185	20.0
Pyrene, 1-methyl-	12.97	3.5	ng	111078	5	13.83	639445	20.0
unknown13.03	13.03	8.2	ng	260435	5	13.83	639445	20.0
Hexadecanoic acid...	13.11	3.2	ng	102089	5	13.83	639445	20.0