

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102521.D
 Acq On : 27 Jan 2018 20:15
 Operator : SJ/JU
 Sample : J1009-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-012518-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-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	4.916	476	481	491	rVB	76652	111226	1.19%	0.211%
2	5.328	547	551	565	rBV	3258957	3286651	35.15%	6.228%
3	6.363	722	727	744	rBV	3075639	3419381	36.57%	6.480%
4	6.487	744	748	751	rBV	3606951	3319220	35.50%	6.290%
5	6.716	783	787	794	rBV	1005333	900343	9.63%	1.706%
6	6.869	809	813	816	rBV	2830071	2447553	26.18%	4.638%
7	7.281	879	883	886	rBV	2011942	2035286	21.77%	3.857%
8	7.998	1001	1005	1013	rVB	1317337	1139009	12.18%	2.158%
9	9.075	1183	1188	1191	rBV	3406394	3005434	32.14%	5.695%
10	9.145	1197	1200	1203	rVB	110285	105240	1.13%	0.199%
11	9.469	1250	1255	1257	rBV	313595	279729	2.99%	0.530%
12	9.675	1288	1290	1294	rVB	216240	178894	1.91%	0.339%
13	9.751	1299	1303	1306	rVV	1391448	1185140	12.68%	2.246%
14	10.175	1372	1375	1378	rVB	279814	218620	2.34%	0.414%
15	10.392	1408	1412	1418	rVB	83207	117914	1.26%	0.223%
16	10.545	1434	1438	1447	rVB	1573729	1628937	17.42%	3.087%
17	10.651	1453	1456	1465	rVB2	223479	324446	3.47%	0.615%
18	11.098	1528	1532	1534	rBV	126655	112592	1.20%	0.213%
19	11.127	1534	1537	1543	rVB3	78975	101086	1.08%	0.192%
20	11.233	1552	1555	1558	rBV	982439	959318	10.26%	1.818%
21	11.257	1558	1559	1566	rVV	393537	323406	3.46%	0.613%
22	11.627	1618	1622	1625	rBV	4046385	3691764	39.49%	6.996%
23	12.327	1734	1741	1742	rBV	6463560	9349651	100.00%	17.718%
24	12.345	1742	1744	1750	rVV2	6293574	6357661	68.00%	12.048%
25	12.421	1753	1757	1759	rVV	1837567	1602611	17.14%	3.037%
26	12.451	1759	1762	1766	rVB	584499	468377	5.01%	0.888%
27	12.680	1791	1801	1808	rBV	513322	686696	7.34%	1.301%
28	12.821	1818	1825	1828	rBV	2167567	2014393	21.55%	3.817%
29	13.010	1849	1857	1860	rBV2	145590	258490	2.76%	0.490%
30	13.063	1864	1866	1871	rVB2	115564	161021	1.72%	0.305%
31	13.139	1876	1879	1882	rVB2	168462	154355	1.65%	0.293%
32	13.716	1974	1977	1986	rVB2	208246	295220	3.16%	0.559%
33	13.810	1990	1993	1996	rVV2	154309	134944	1.44%	0.256%
34	13.874	2000	2004	2007	rVV2	874553	1049500	11.23%	1.989%

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Integration Parameters: rteint.p
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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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35	13.904	2007	2009	2015	rVB	221753	218971	2.34%	0.415%
36	14.898	2175	2178	2181	rBV	168383	225153	2.41%	0.427%
37	15.245	2234	2237	2243	rVB	142978	185876	1.99%	0.352%
38	15.310	2244	2248	2257	rVB	540058	715629	7.65%	1.356%

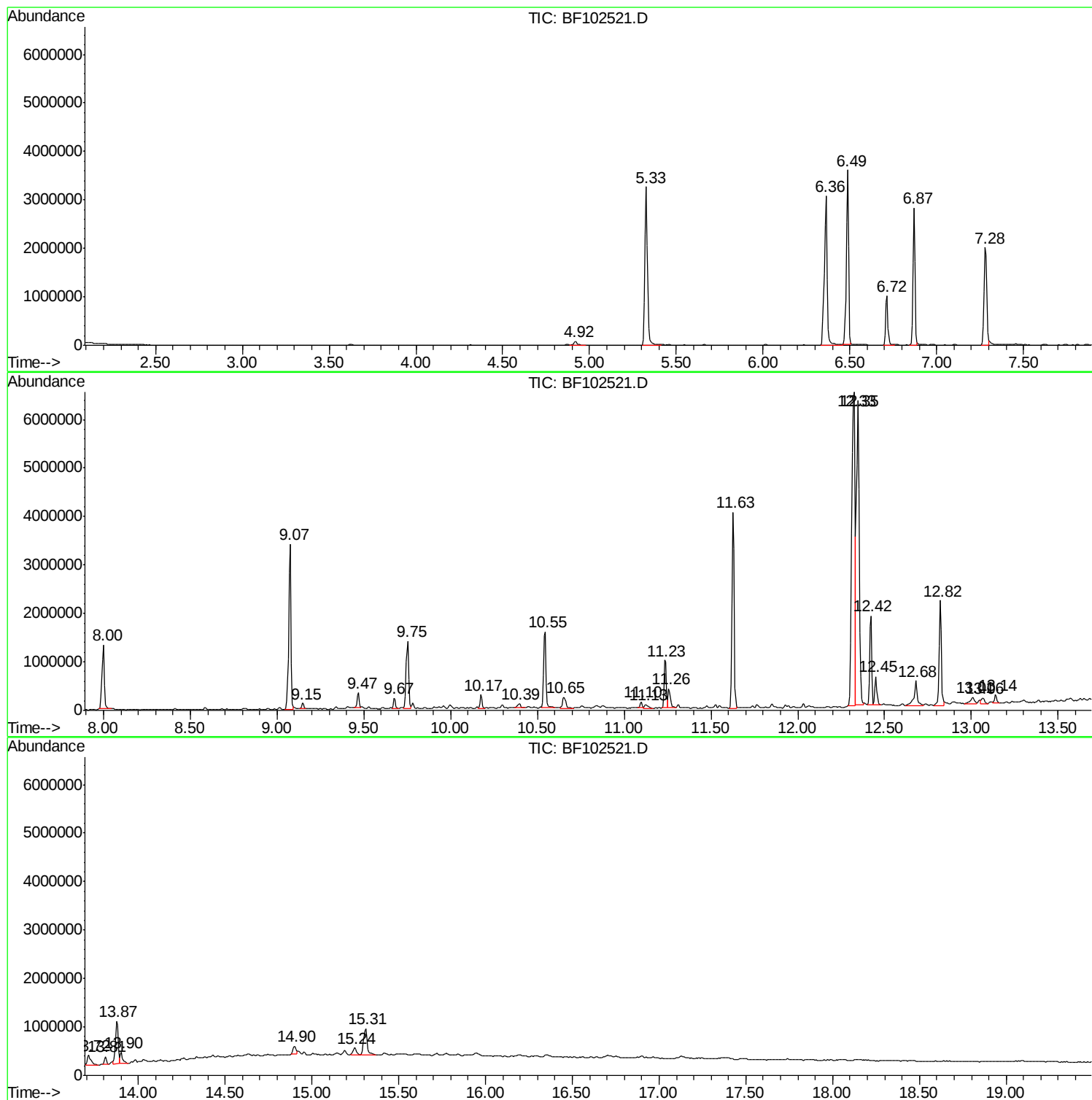
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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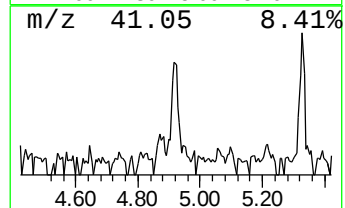
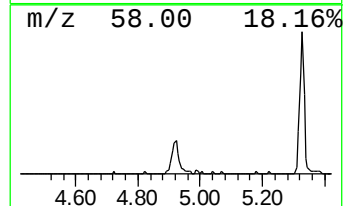
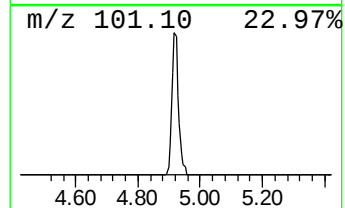
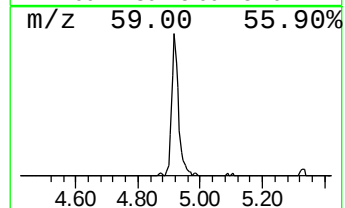
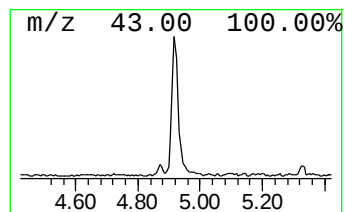
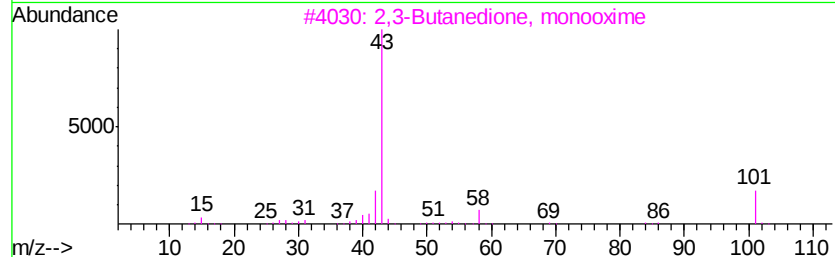
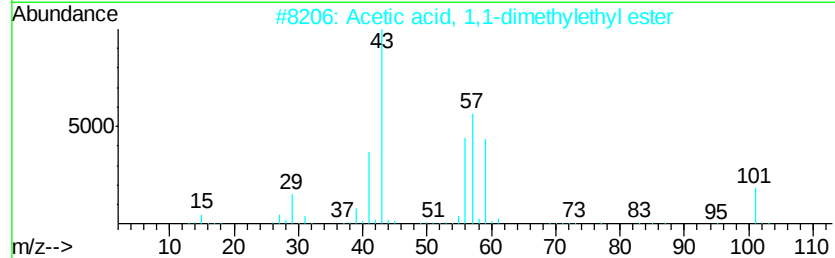
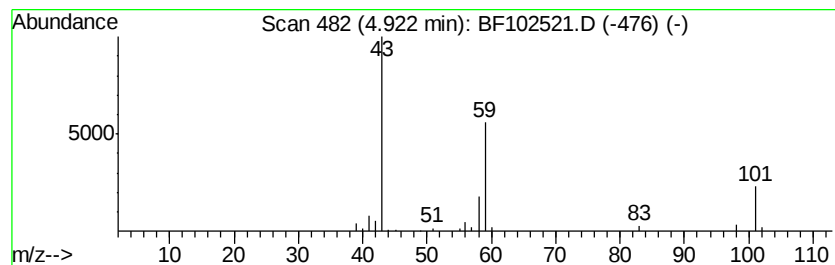
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	2.47 ng	111226	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4			3-Hexanol	102	C6H14O	000623-37-0	28
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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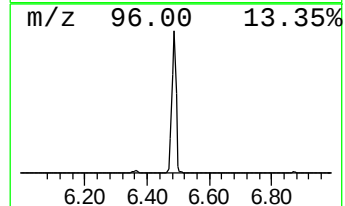
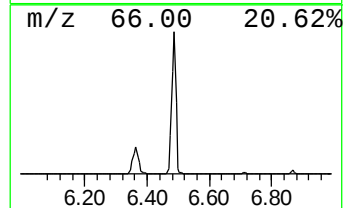
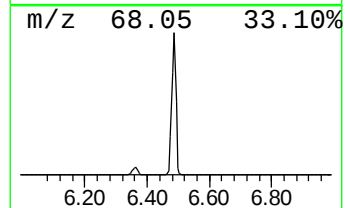
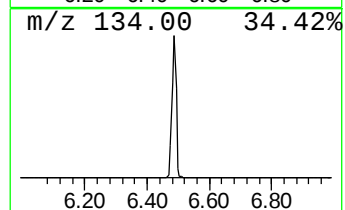
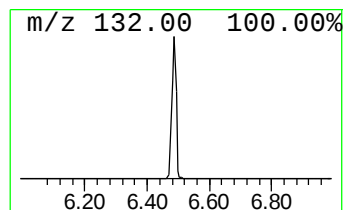
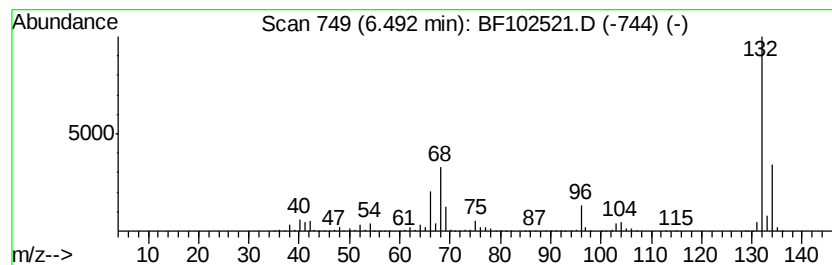
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.49 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	73.73 ng	3319220	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	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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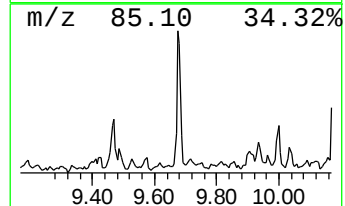
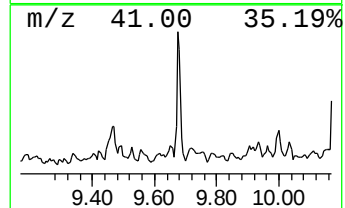
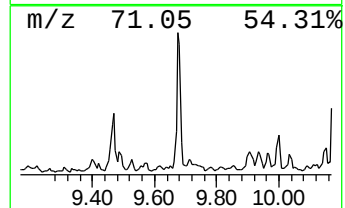
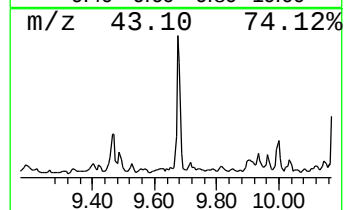
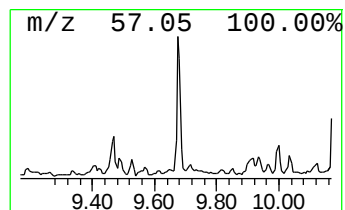
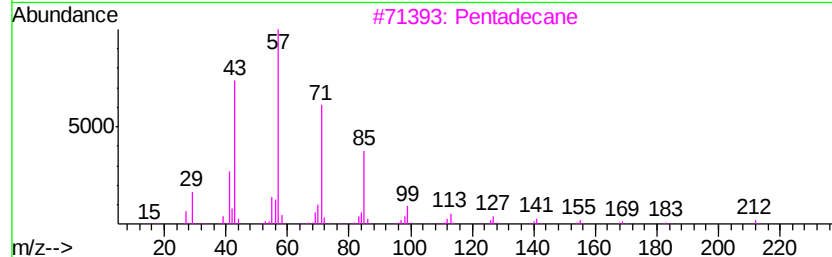
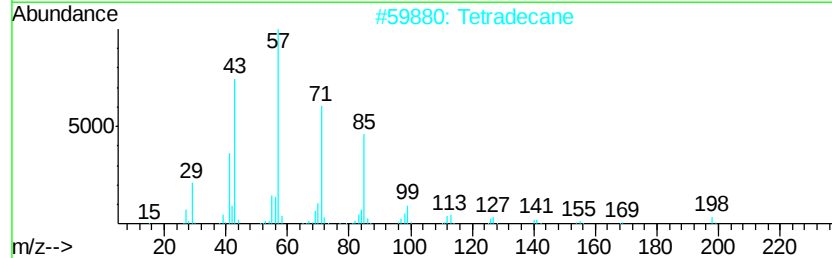
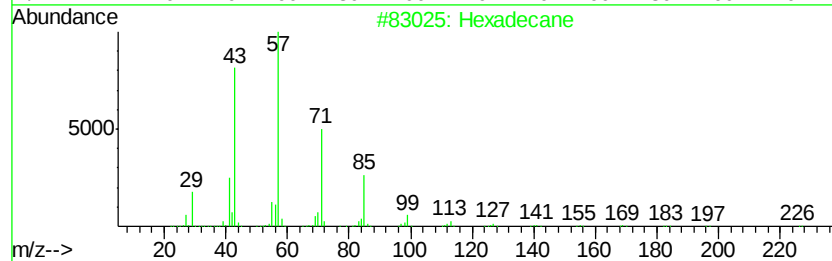
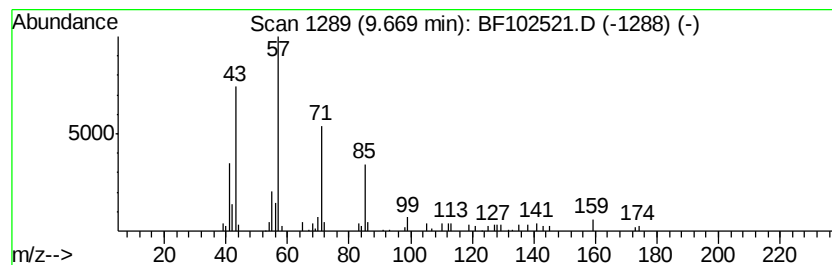
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.67	3.02 ng	178894	Acenaphthene-d10	9.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetradecane		198	C14H30	000629-59-4	90
3	Pentadecane		212	C15H32	000629-62-9	86
4	Tridecane		184	C13H28	000629-50-5	86
5	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	78



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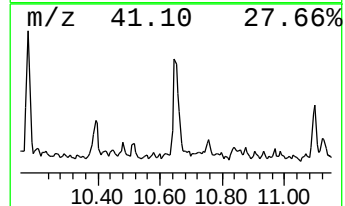
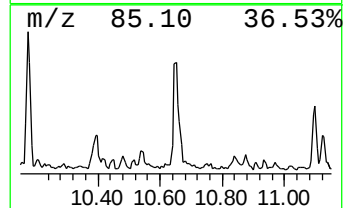
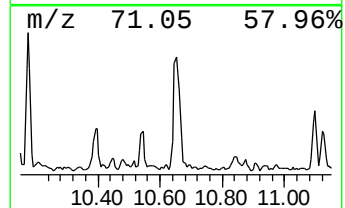
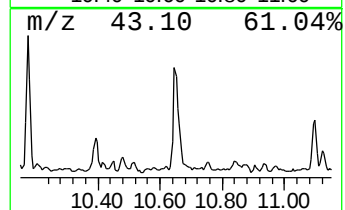
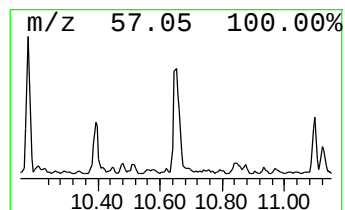
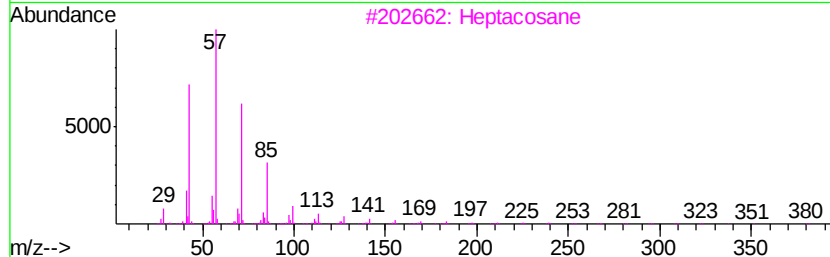
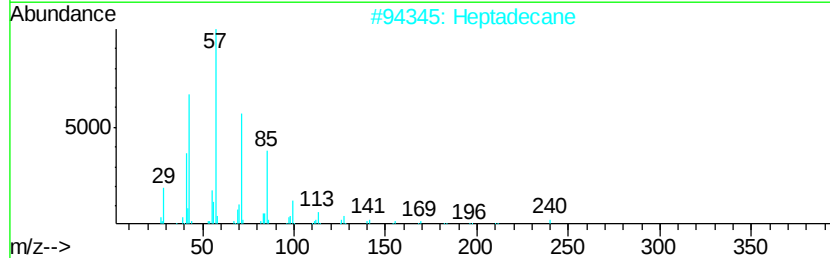
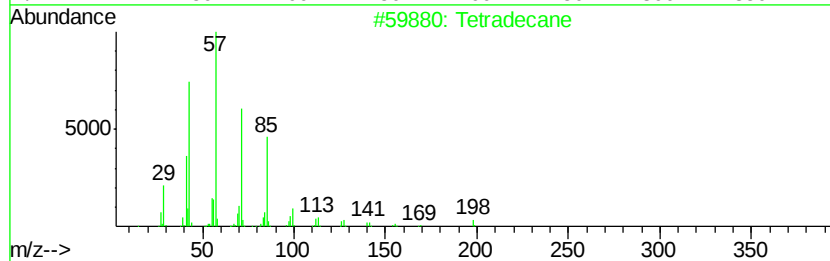
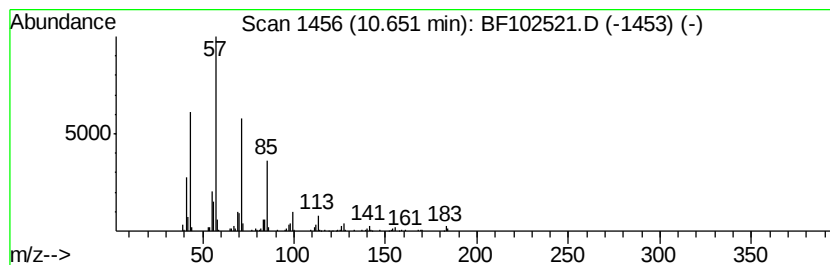
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	6.76 ng	324446	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Heptacosane	380	C27H56	000593-49-7	90
4		Hexadecane	226	C16H34	000544-76-3	90
5		Tridecane	184	C13H28	000629-50-5	87



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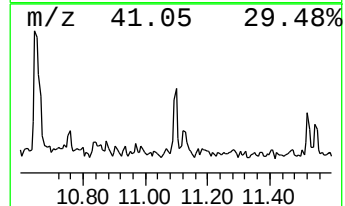
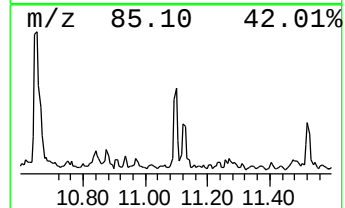
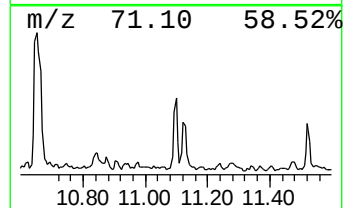
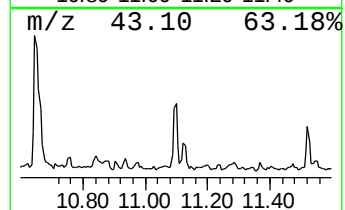
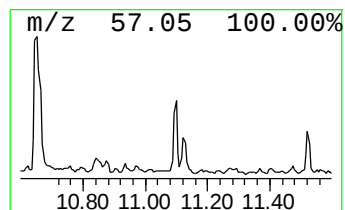
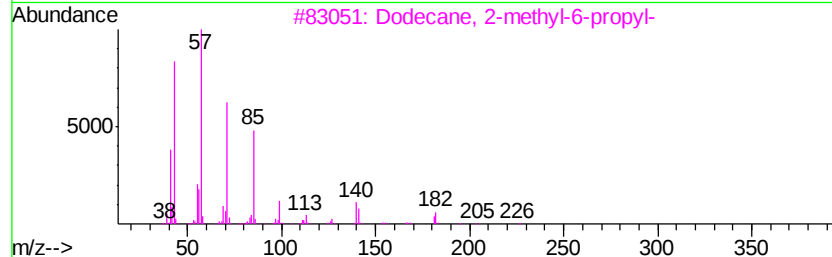
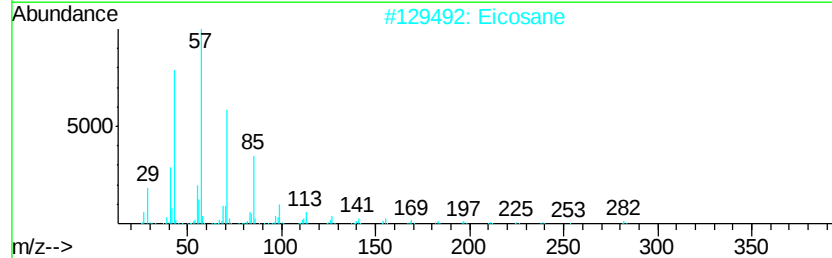
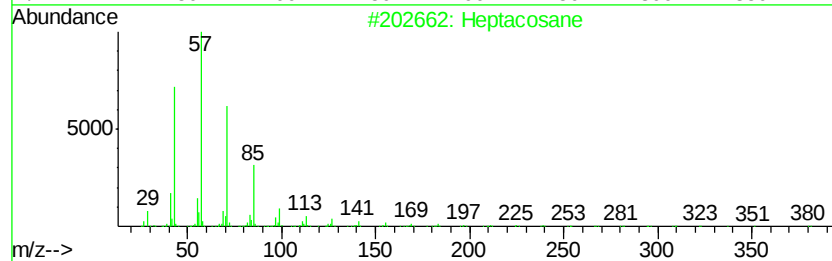
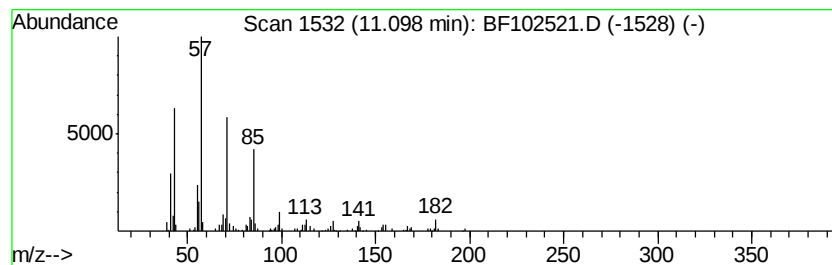
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Heptacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.35 ng	112592	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Eicosane		282	C20H42	000112-95-8	86
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86
4	Hexadecane		226	C16H34	000544-76-3	86
5	Docosane		310	C22H46	000629-97-0	86



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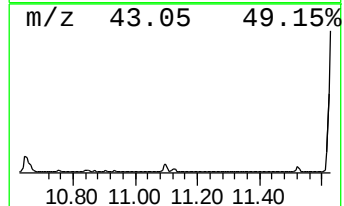
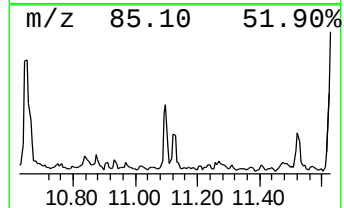
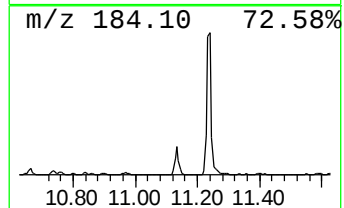
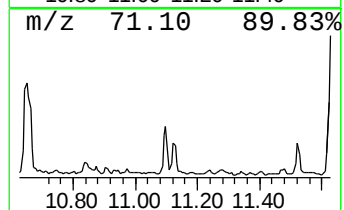
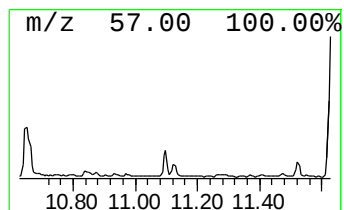
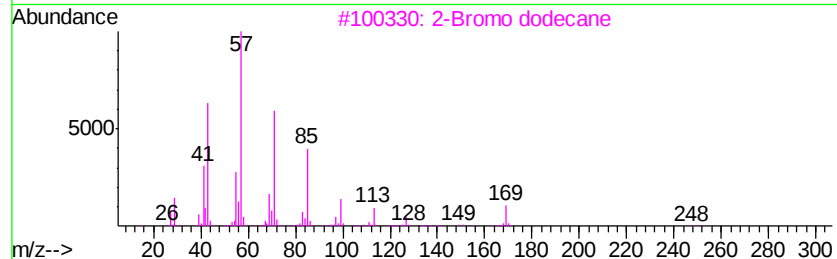
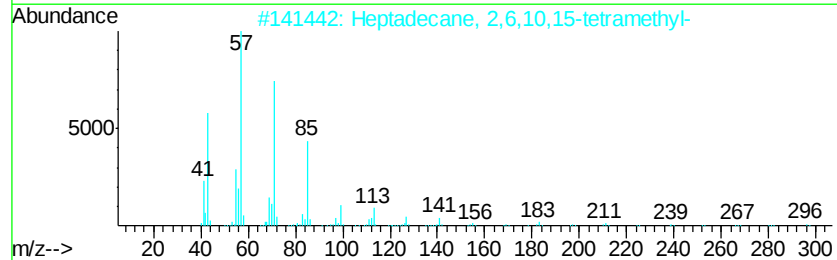
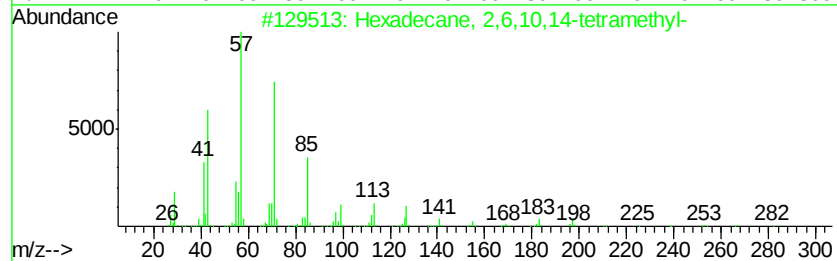
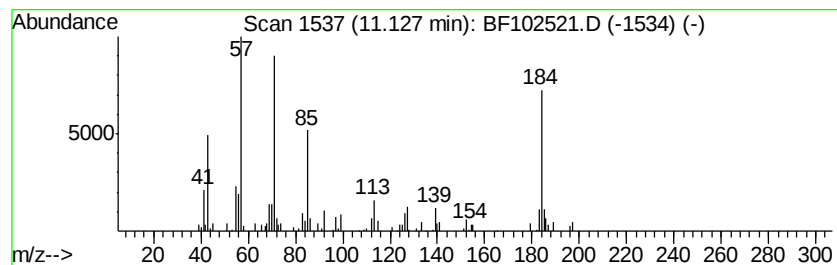
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ClientSampleId :
AU-06-012518-A

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 Hexadecane, 2,6,10,14-tetra... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.13	2.11 ng	101086	Phenanthrene-d10	11.23		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	58	
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	49	
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	47	
4	Tridecane, 3-methyl-	198	C14H30	006418-41-3	47	
5	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	43	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-A

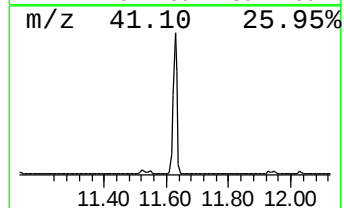
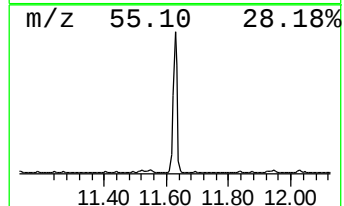
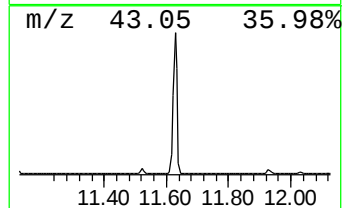
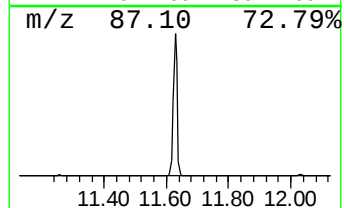
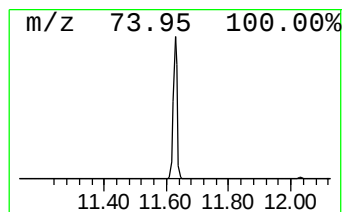
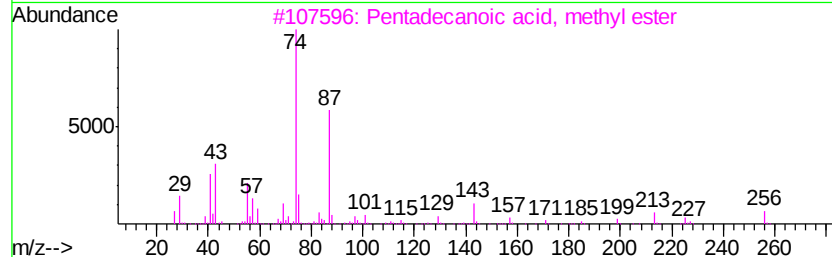
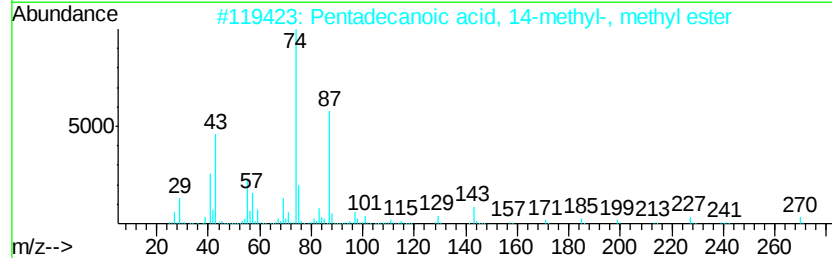
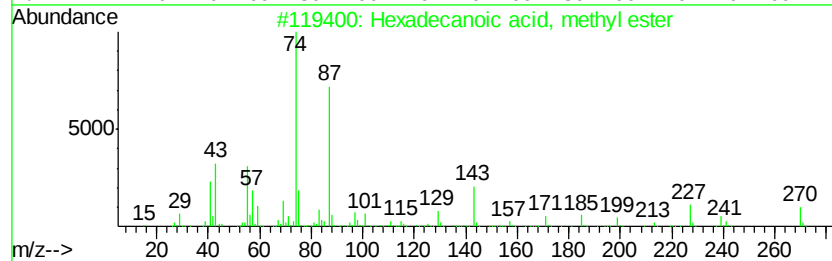
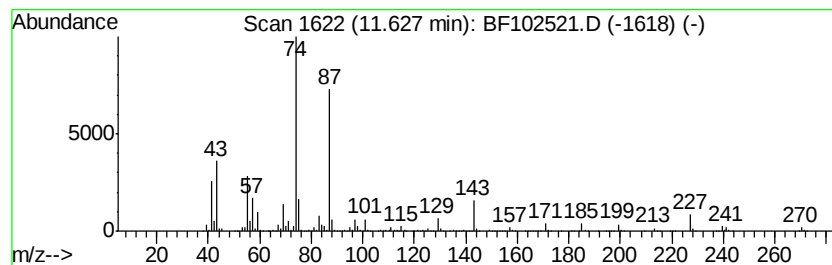
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.63	76.97 ng	3691760	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
4		Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	90
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-A

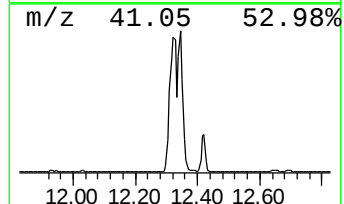
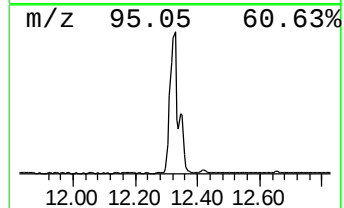
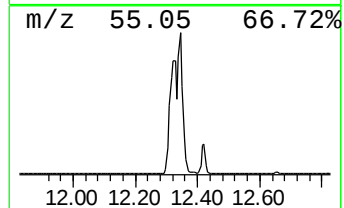
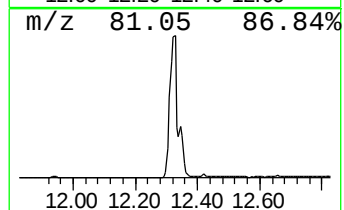
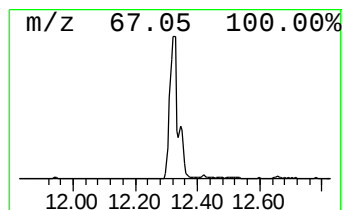
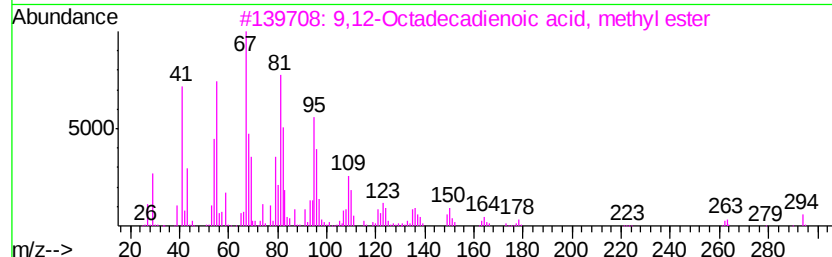
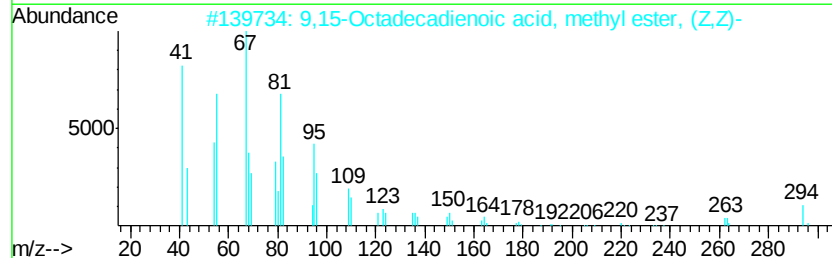
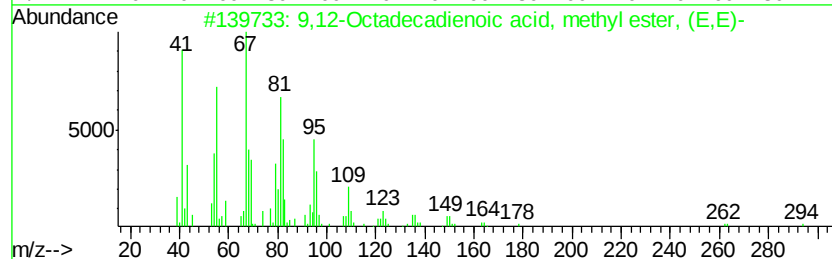
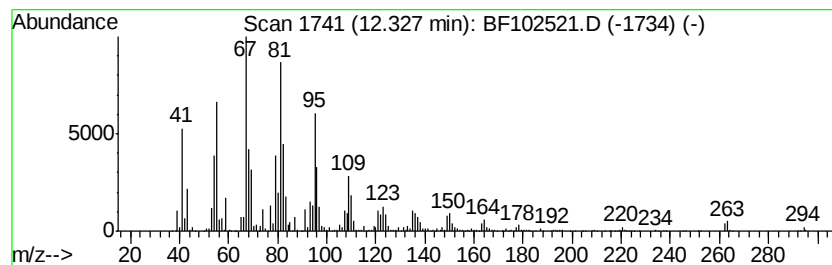
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
12.33	194.92 ng	9349650	Phenanthrene-d10	11.23

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		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
4		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-A

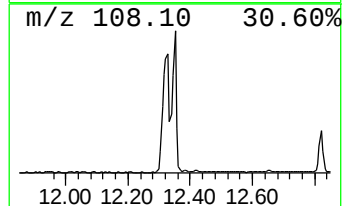
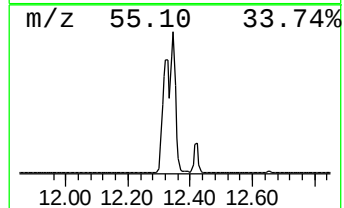
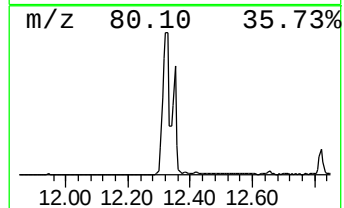
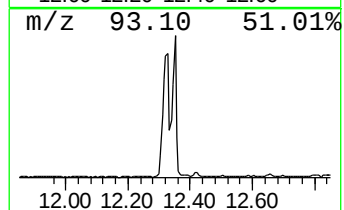
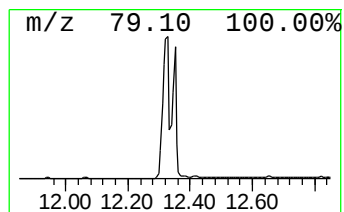
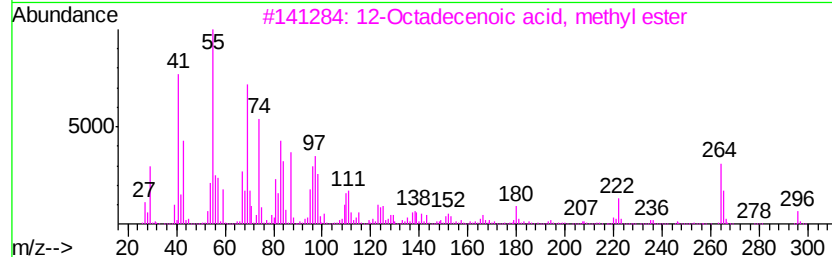
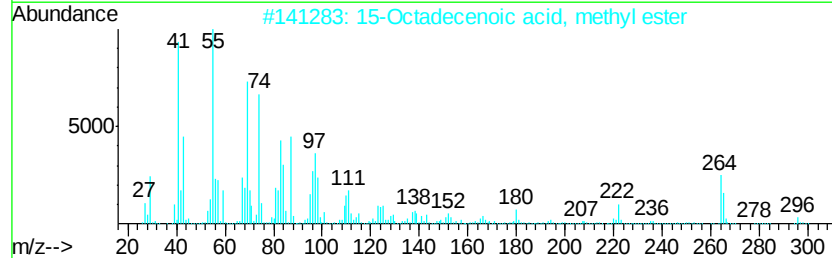
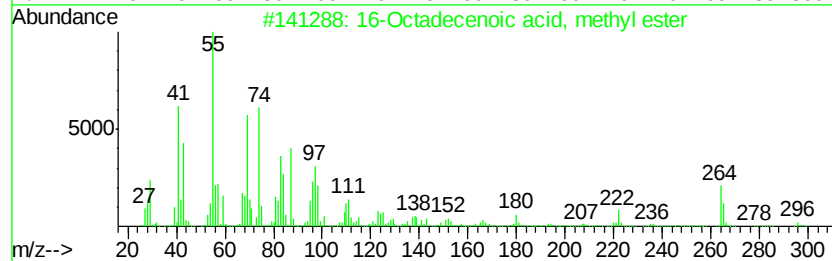
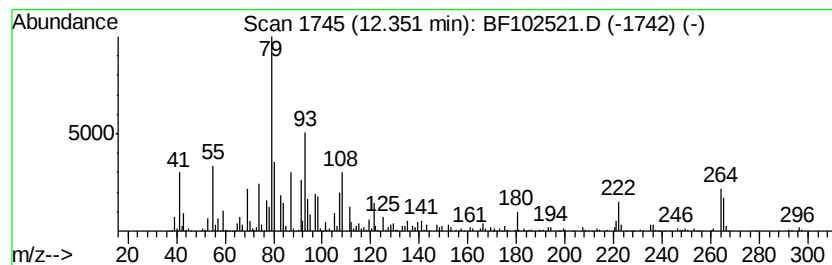
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 16-Octadecenoic acid, methyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	132.55 ng	6357660	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	94
2		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	94
3		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	93
4		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	93
5		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

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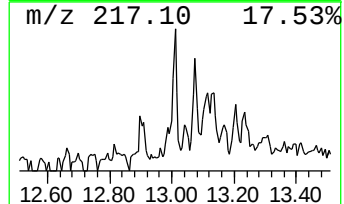
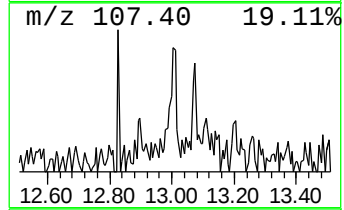
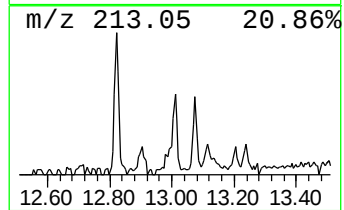
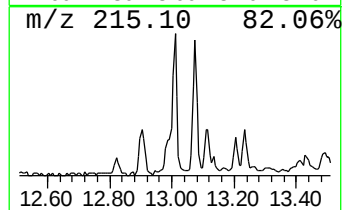
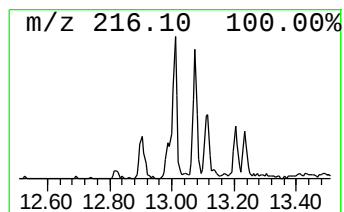
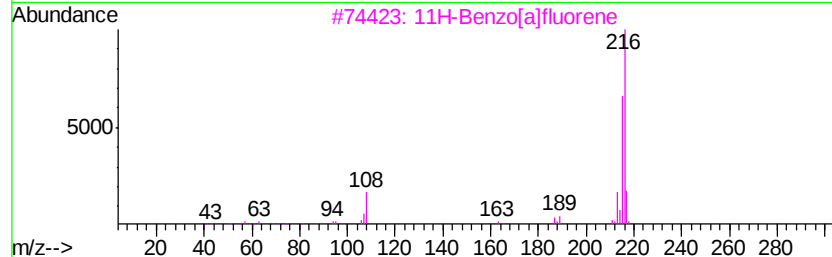
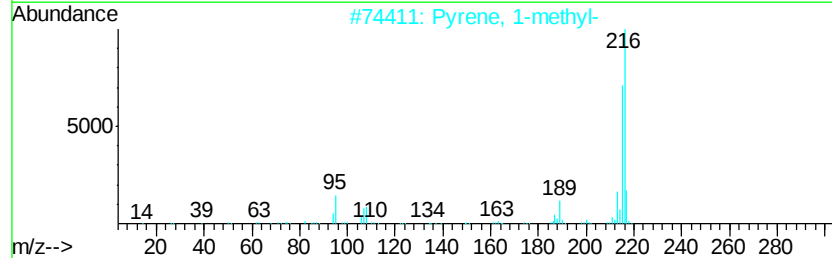
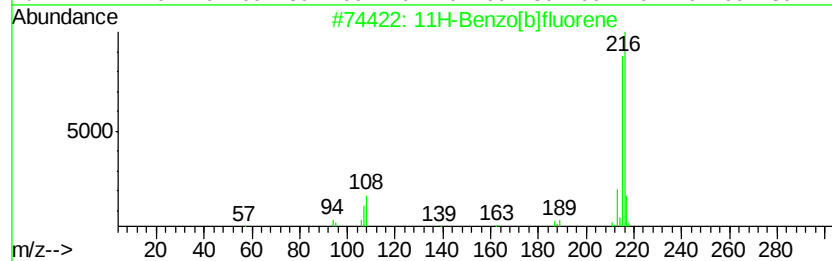
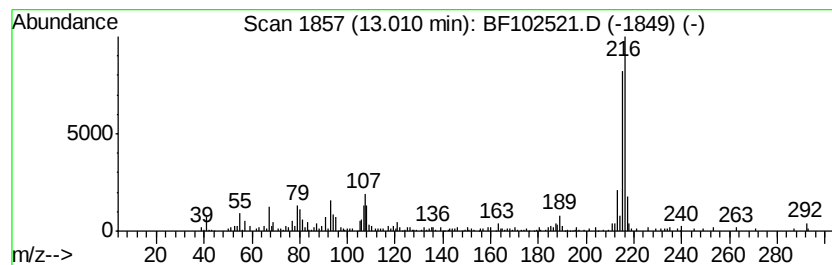
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	4.93 ng	258490	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	89
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	89
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AU-06-012518-A

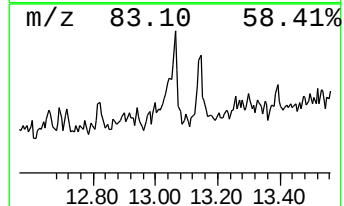
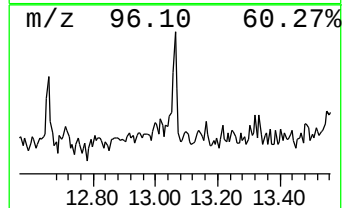
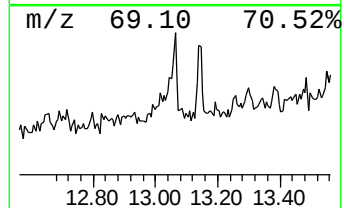
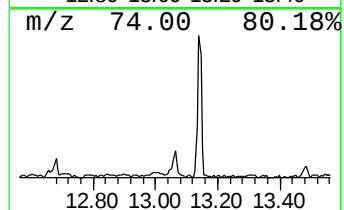
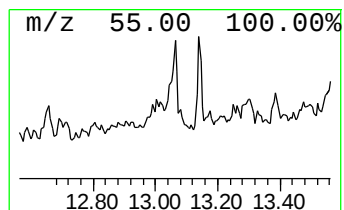
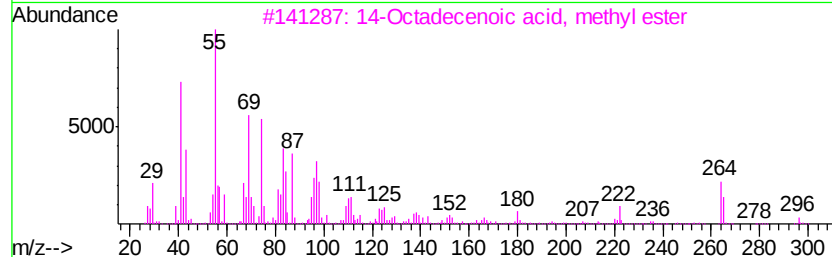
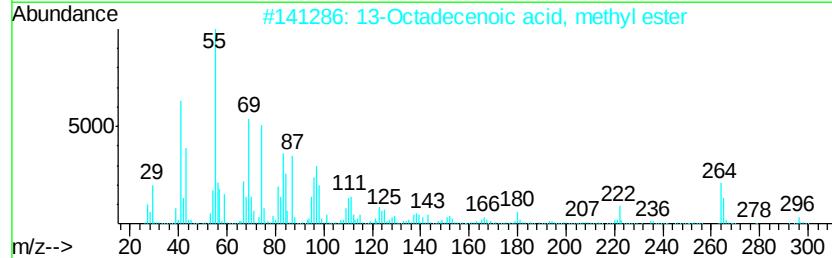
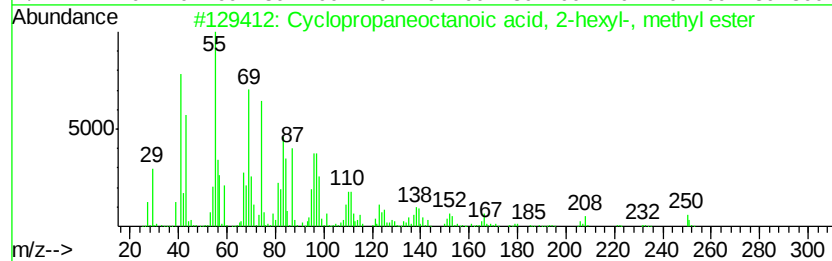
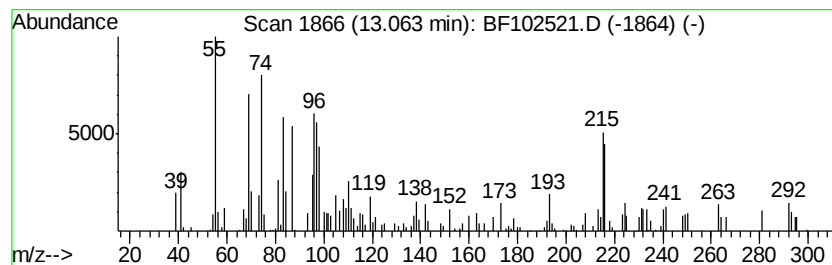
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopropaneoctanoic acid, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	3.07 ng	161021	Chrysene-d12	13.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropaneoctanoic acid, 2-hex...	282	C18H34O2	010152-61-1	52
2			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	50
3			14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	50
4			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	49
5			Cyclopropanenonanoic acid, methy...	212	C13H24O2	010152-60-0	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AU-06-012518-A

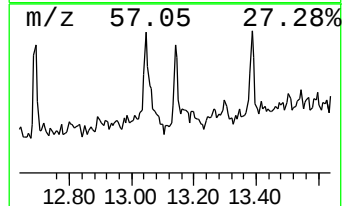
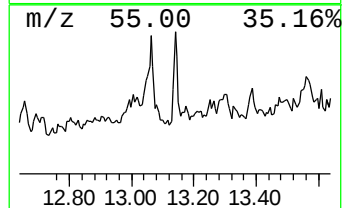
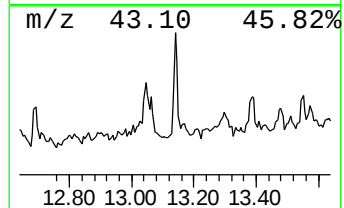
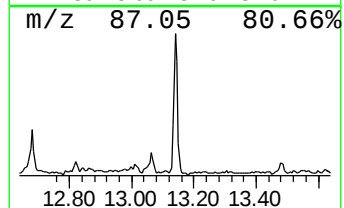
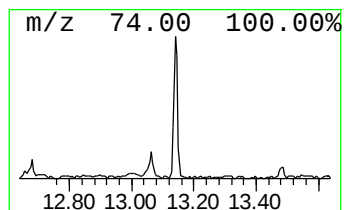
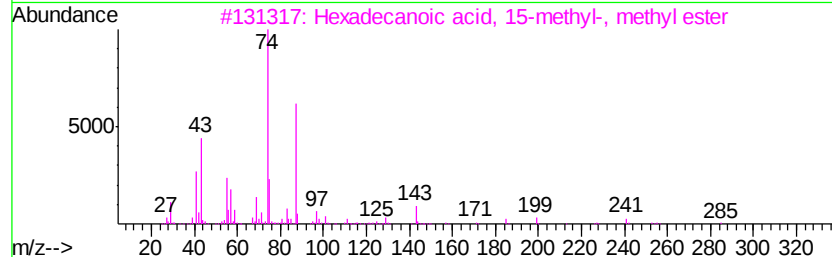
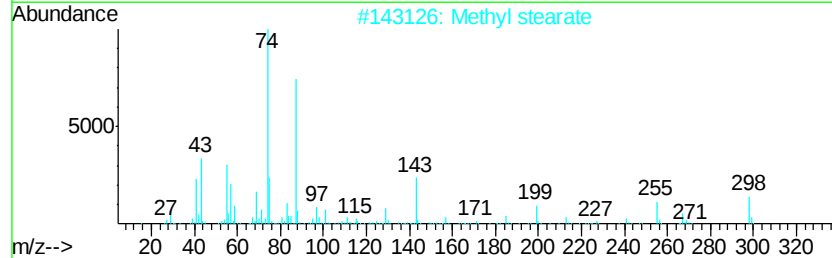
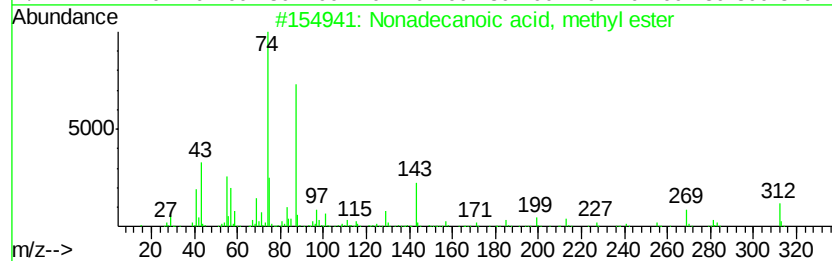
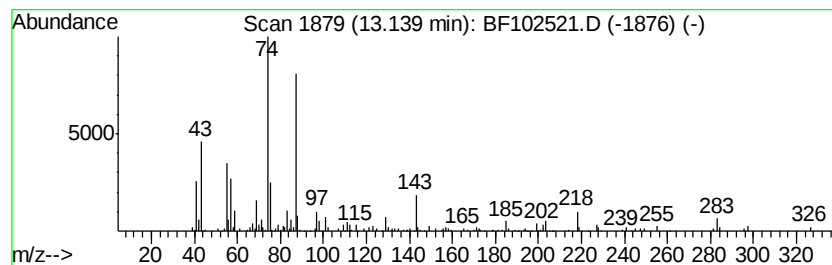
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecanoic acid, methyl e... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.14	2.94 ng	154355	Chrysene-d12	13.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	87
2			Methyl stearate	298	C19H38O2	000112-61-8	87
3			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86
4			Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	86
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-A

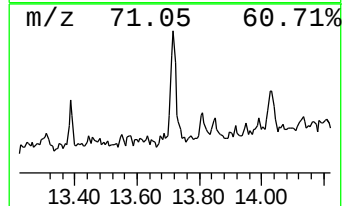
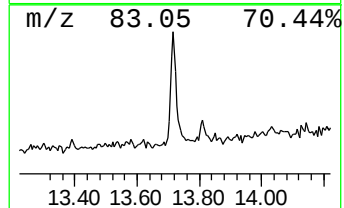
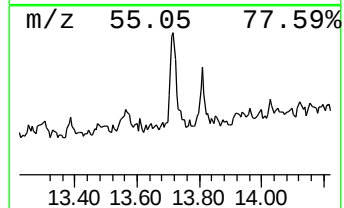
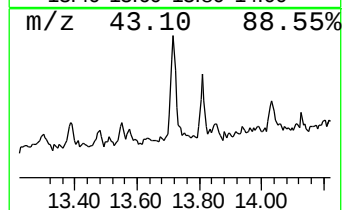
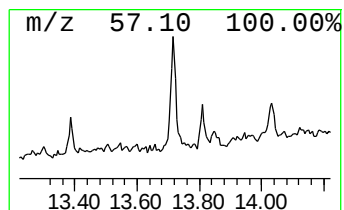
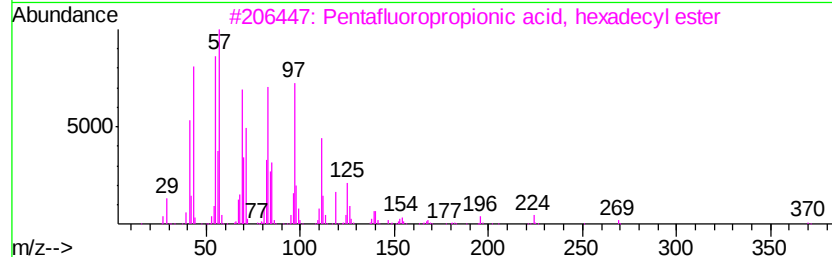
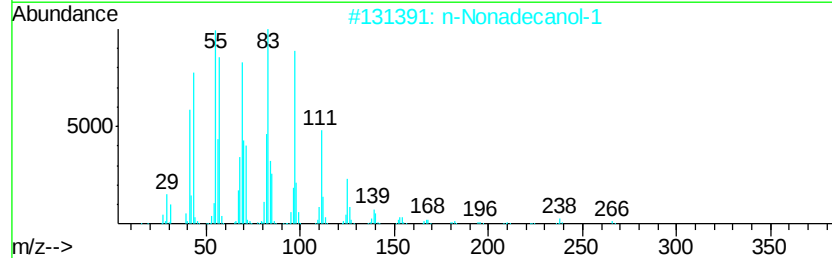
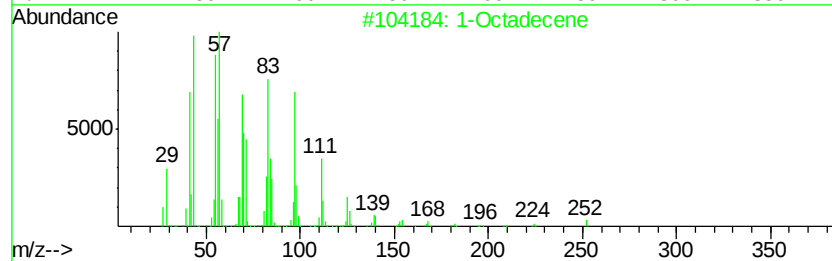
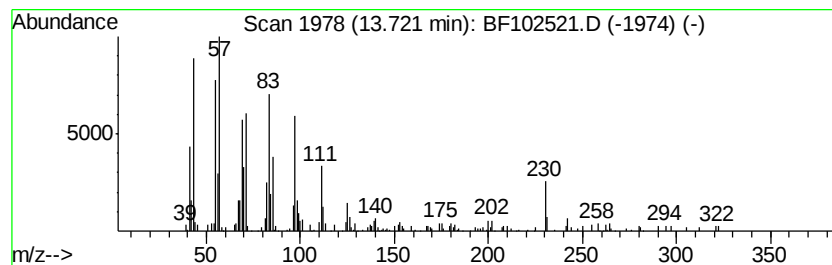
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Octadecene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	5.63 ng	295220	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	97
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	93
3		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
4		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102521.D
Acq On : 27 Jan 2018 20:15
Operator : SJ/JU
Sample : J1009-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-A

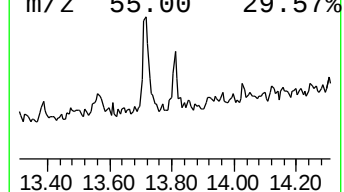
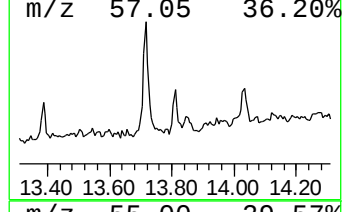
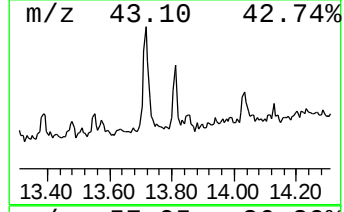
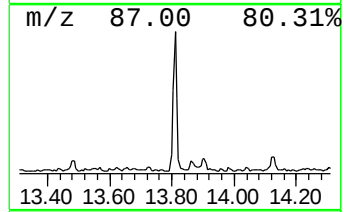
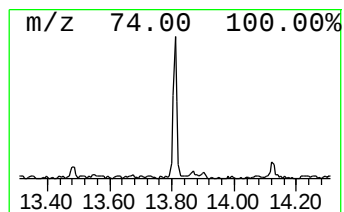
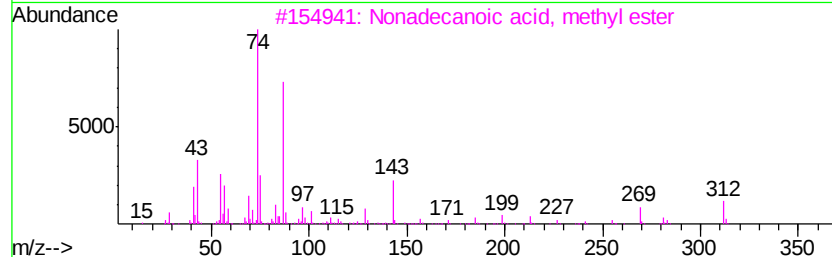
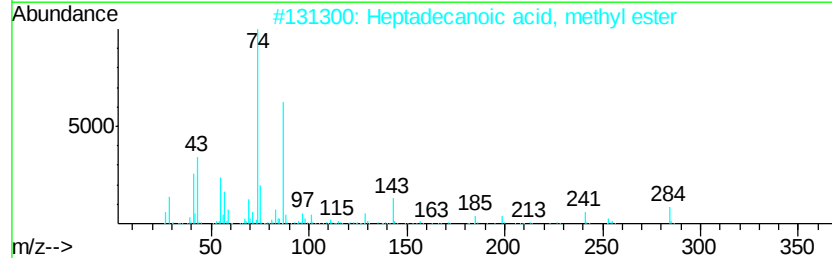
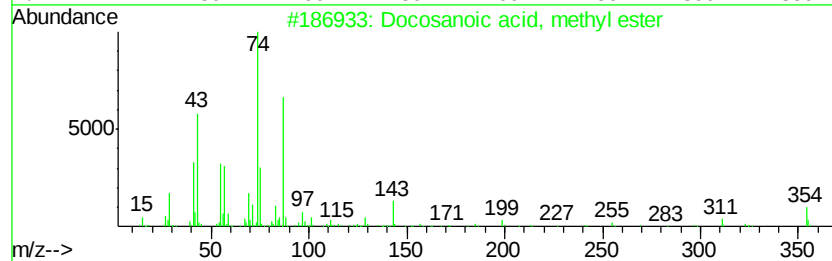
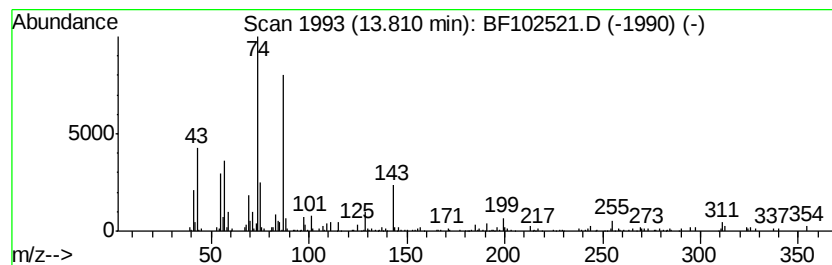
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Docosanoic acid, methyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	2.57 ng	134944	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	97
2		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
3		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	91
4		Methyl stearate	298	C19H38O2	000112-61-8	91
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102521.D
 Acq On : 27 Jan 2018 20:15
 Operator : SJ/JU
 Sample : J1009-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-012518-A

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
2-Pentanone, 4-hy...	4.92	2.5	ng	111226	1	6.72	900343	20.0
unknown6.49	6.49	73.7	ng	3319220	1	6.72	900343	20.0
Hexadecane	9.67	3.0	ng	178894	3	9.75	1185140	20.0
Tetradecane	10.65	6.8	ng	324446	4	11.23	959318	20.0
Heptacosane	11.10	2.4	ng	112592	4	11.23	959318	20.0
Hexadecane, 2,6,1...	11.13	2.1	ng	101086	4	11.23	959318	20.0
Hexadecanoic acid...	11.63	77.0	ng	3691760	4	11.23	959318	20.0
9,12-Octadecadien...	12.33	194.9	ng	9349650	4	11.23	959318	20.0
16-Octadecenoic a...	12.35	132.6	ng	6357660	4	11.23	959318	20.0
11H-Benzo[b]fluorene	13.01	4.9	ng	258490	5	13.87	1049500	20.0
Cyclopropaneoctan...	13.06	3.1	ng	161021	5	13.87	1049500	20.0
Nonadecanoic acid...	13.14	2.9	ng	154355	5	13.87	1049500	20.0
1-Octadecene	13.72	5.6	ng	295220	5	13.87	1049500	20.0
Docosanoic acid, ...	13.81	2.6	ng	134944	5	13.87	1049500	20.0