

CORRECTION

Correction: DNA Sequence Determinants Controlling Affinity, Stability and Shape of DNA Complexes Bound by the Nucleoid Protein Fis

The *PLOS ONE* Staff

There is an error in Tables 2, 3 and 4. The numbering in the first row of the second column does not line up with the corresponding DNA base sequence below. For example, in the sequence aaatttGTTTGAATTTGAGCaaattt, the fourth capital “T” should be directly below the number 0. The publisher apologizes for the errors.



OPEN ACCESS

Citation: The *PLOS ONE* Staff (2016) Correction: DNA Sequence Determinants Controlling Affinity, Stability and Shape of DNA Complexes Bound by the Nucleoid Protein Fis. *PLoS ONE* 11(6): e0157224. doi:10.1371/journal.pone.0157224

Published: June 21, 2016

Copyright: © 2016 The PLOS ONE Staff. This is an open access article distributed under the terms of the [Creative Commons Attribution License](#), which permits unrestricted use, distribution, and reproduction in any medium, provided the original author and source are credited.

Table 2. Effects of flanking DNA substitutions on Fis-DNA binding affinity and complex stability¹

	-10 -9 -8 -7 -6 -5 -4 -3 -2 -1 0 1 2 3 4 5 6 7 8 9 10	K _d (nM)	Fold-difference ²	t _{1/2} (min)
F1(\pm 8T)	aaatttGTTTGAATTTGAGCaaat	0.2 ± 0.07	1	41 ± 4
F1±8A	aaatt <u>a</u> GTTTGAATTTGAGC <u>t</u> aat	2.1 ± 0.2	10	5 ± 0.5
F1±8C	aaatt <u>c</u> GTTTGAATTTGAGC <u>g</u> aat	0.7 ± 0.2	3	14 ± 1
F1±8G	aaatt <u>g</u> GTTTGAATTTGAGC <u>a</u> at	30 ± 6	150	< 0.25
F1±9A	aaat <u>a</u> GTTTGAATTTGAGC <u>t</u> at	0.5 ± 0.2	3	22 ± 2
F1±9C	aaat <u>c</u> tGTTTGAATTTGAGC <u>g</u> at	0.3 ± 0.1	2	30 ± 2
F1±9G	aaat <u>g</u> GTTTGAATTTGAGC <u>c</u> at	0.6 ± 0.2	3	25 ± 2
F1±10A	aaa <u>t</u> tGTTTGAATTTGAGCaa <u>t</u> tt	0.5 ± 0.1	3	36 ± 5
F1±10C	aaa <u>c</u> tGTTTGAATTTGAGCaa <u>g</u> tt	0.3 ± 0.1	2	44 ± 1
F1±10G	aa <u>g</u> tGTTTGAATTTGAGCaa <u>c</u> tt	0.5 ± 0.1	3	38 ± 5
F14	<u>ggg</u> tttGTTTGAATTTGAGCaa <u>ccc</u>	0.5 ± 0.1	3	ND ³
F15	<u>ccc</u> tttGTTTGAATTTGAGCaa <u>ggg</u>	0.4 ± 0.2	2	ND
F16	<u>gcgg</u> tttGTTTGAATTTGAGCaa <u>ccgc</u>	0.5 ± 0.2	3	ND
F33	aa <u>gg</u> tGTTTGAATTTGAGCaa <u>cc</u> tt	0.6 ± 0.1	6	ND
F34	a <u>ag</u> ttGTTTGAATTTGAGCaa <u>cg</u> tt	0.2 ± 0.1	1	8 ± 1
INV	<u>ttaaa</u> GTTTGAATTTGAGC <u>ttaaa</u>	33 ± 3	165	< 0.25
INV±8G	<u>ttaag</u> GTTTGAATTTGAGC <u>cttaaa</u>	250 ± 20	1250	ND
INV±8C	<u>ttaac</u> GTTTGAATTTGAGC <u>gttaaa</u>	6.0 ± 2	30	ND
INV±8T	<u>ttaa</u> tGTTTGAATTTGAGC <u>ttaaa</u>	2.3 ± 0.7	12	4 ± 0.7
INV-CAT	<u>tttca</u> tGTTTGAATTTGAGC <u>tgaaa</u>	2.9 ± 1.0	15	ND
INV-GAT	<u>tttga</u> tGTTTGAATTTGAGC <u>tcaa</u>	2.0 ± 0.6	10	ND
INV±9-10T	<u>tttatt</u> GTTTGAATTTGAGCaa <u>taaa</u>	0.4 ± 0.1	2	ND

¹Upper case letters represent the 15 bp core Fis binding site sequence and those in lower case represent flanking DNA. Underlined and bold nucleotides highlight those that differ from F1.

²Fold-difference relative to the apparent equilibrium dissociation constant (K_d) for WT Fis with F1 DNA.

³Not determined.

doi:10.1371/journal.pone.0157224.t002

Table 3. Interplay between Fis residues contacting the flanking sequences and binding site variants

	-10 -9 -8 -7 -6 -5 -4 -3 -2 -1 0 1 2 3 4 5 6 7 8 9 10	Fis protein	K _d (nM)	Fold-difference ¹
F1	aaatttGTTGAATTTGAGCaaattt	WT	0.2 ± 0.05	1
		R71A	0.5 ± 0.2	2.5
		T75A	0.3 ± 0.8	1.5
		N73A	29 ± 0.2	140
F27	aaatttGTTGA <u>A</u> CTTTGAGCaaattt	WT	0.2 ± 0.1	1
		R71A	2.8 ± 0.5	14
		T75A	3.5 ± 0.9	18
F28	aaatttGTTGA <u>GCG</u> TTGAGCaaattt	WT	28 ± 4	140
		R71A	470 ± 100	2300
		T75A	> 1000	> 5000
F32	aaattt <u>G</u> <u>GAG</u> AA <u>TTT</u> <u>TC</u> Caaattt	WT	28 ± 5	140
		R71A	73 ± 11	370
		T75A	76 ± 10	380
F1±8G	aaatt <u>g</u> GTTTGAATTTGAGC <u>c</u> aattt	WT	30 ± 0.8	150
		R71A	450 ± 20	2270
		T75A	48 ± 5	240

¹Fold-difference relative to the apparent equilibrium dissociation constant (K_d) for WT Fis with F1 DNA.

doi:10.1371/journal.pone.0157224.t002

Table 4. Effects of flanking and core substitutions on Fis-induced DNA bending

	-10 -9 -8 -7 -6 -5 -4 -3 -2 -1 0 1 2 3 4 5 6 7 8 9 10	Gel mobility assay		In-gel FRET assay		Distance (Å) (complex) ¹	Angle (°) (complex) ²		
		(Bend angle°)		(FRET efficiency)					
		Free	Complex	Free	Complex				
F1	aaatttGTTGAATTTGAGCaaattt	51 ± 3	119 ± 3	0.09 ± 0.01	0.27 ± 0.01	82.6	68		
F1±8A	aaatt <u>a</u> GTTGAATTTGAGC <u>t</u> aattt	48 ± 3	109 ± 3	-	-	-	-		
F1±8C	aaatt <u>c</u> GTTGAATTTGAGC <u>g</u> aattt	47 ± 3	126 ± 2	0.13 ± 0.01	0.26 ± 0.01	83.3	66		
F1±8G	aaatt <u>g</u> GTTGAATTTGAGC <u>c</u> aattt	41 ± 1	99 ± 1	0.09 ± 0.01	0.23 ± 0.01	85.6	61		
F1±9A	aaat <u>a</u> tGTTGAATTTGAGC <u>a</u> tattt	46 ± 3	104 ± 2	-	-	-	-		
F1±9C	aaat <u>c</u> tGTTGAATTTGAGC <u>g</u> attt	45 ± 1	99 ± 1	-	-	-	-		
F1±9G	aaat <u>g</u> tGTTGAATTTGAGC <u>c</u> attt	42 ± 2	89 ± 2	0.08 ± 0.03	0.18 ± 0.04	90.1	50		
INV	<u>t</u> <u>t</u> <u>a</u> <u>a</u> GTTGAATTTGAGC <u>t</u> <u>t</u> <u>a</u> <u>a</u>	≤ 36	62 ± 1	0.06 ± 0.04	0.14 ± 0.01	94.7	36		
F34	a <u>g</u> <u>ttt</u> GTTGAATTTGAGCaa <u>cg</u> tt	~ 0	62 ± 3	0.12 ± 0.02	0.26 ± 0.02	83.3	66		
F18	aaatttGTT <u>G</u> GAATTT <u>C</u> AGCaaattt	51 ± 2	116 ± 1	-	-	-	-		
F31	aaatttG <u>T</u> <u>AG</u> GAATTT <u>C</u> TGCAAattt	51 ± 2	108 ± 2	-	-	-	-		
F32	aaatttG <u>G</u> <u>AG</u> GAATTT <u>C</u> TCCAAattt	52 ± 1	106 ± 2	-	-	-	-		

¹Inter-fluorophore distance calculated from FRET efficiency as detailed in the Methods.

²Angle calculated assuming a single central bend in the Fis-bound DNA as detailed in the Methods.

doi:10.1371/journal.pone.0157224.t003

Reference

- Hancock SP, Stella S, Cascio D, Johnson RC (2016) DNA Sequence Determinants Controlling Affinity, Stability and Shape of DNA Complexes Bound by the Nucleoid Protein Fis. PLoS ONE 11(3): e0150189. doi:[10.1371/journal.pone.0150189](https://doi.org/10.1371/journal.pone.0150189) PMID: [26959646](https://pubmed.ncbi.nlm.nih.gov/26959646/)