

Supplementary Figures and Tables: CSI-Tree: A  
regression tree approach for modeling binding properties  
of DNA binding molecules based on cognate site  
identification (CSI) data

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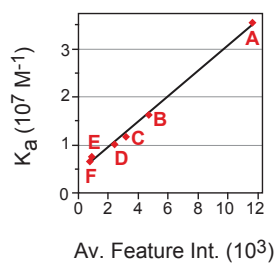
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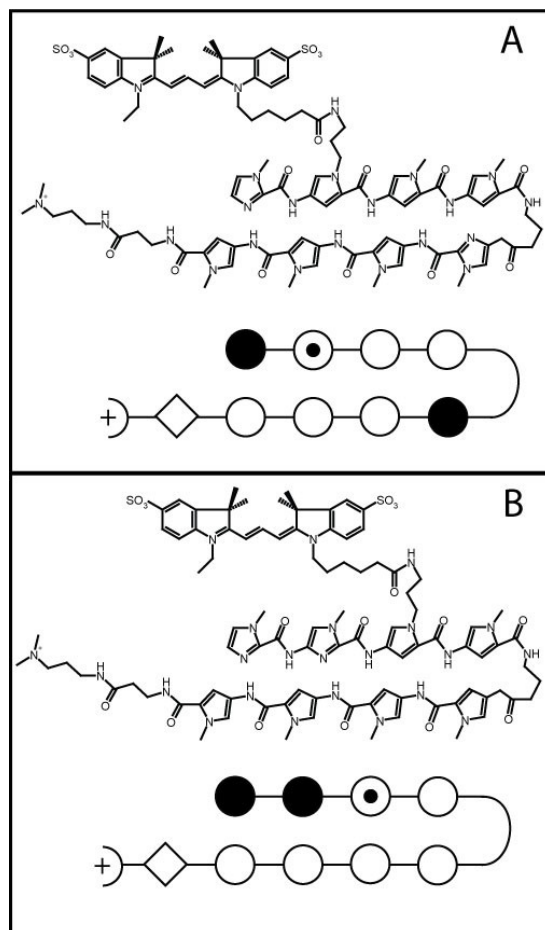
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Data Point	Sequence	$K_a$ ( $10^7 \text{ M}^{-1}$ )	Intensity ( $10^3$ )
A	AAGGTTA	3.59	11.6
B	AACGTTA	1.62	4.71
C	ATGGAAT	1.17	3.16
D	CAGGTTA	0.95	2.39
E	TACGATT	0.73	0.86
F	AAGGTCA	0.65	0.79

*Supplementary Table 1. CSI fluorescent intensities and corresponding binding affinities  $K_a$  determined from nuclease protection assays for the six sequences spanning the range of the PA2 CSI data.*



*Supplementary Figure 1. CSI fluorescent intensities versus binding affinities  $K_a$  for six sequences spanning the range of the PA2 CSI data. Similar linear correlations are observed for PA1 [1] and Nkx-2.5 (data not shown).*



*Supplementary Figure 2. Chemical structures of polyamides 1 (PA1) and 2 (PA2).*

## References

- [1] C.L. Warren, N. C. Kratochvil, K. E. Hauschild, S. Foister, M. L. Brezinski, P. B. Dervan, G. N. Phillips, and A.Z. Ansari. Defining the sequence-recognition profile of dna-binding molecules. *Proceedings of the National Academy of Sciences U.S.A*, 103(4):867–872, January 2006.