Exact Solution of a Simple Adsorption Model of **De-naturating DNA**

[†]Aleks Owczarek, [‡]Andrew Rechnitzer, and [†]Rami Tabbara

[†]MASCOS and Department of Mathematics and Statistics, The University of Melbourne

[‡] Department of Mathematics and Statistics, University of British Columbia

June, 2014



2014 SIAM Conference on Discrete Mathematics



• □ > • □ > • □ > • □ > • □ > •

Exact Solution of a Simple Adsorption Model of De-naturating DNA

AN RESEARCH COUNCI

Owczarek

- DNA is a polymer consisting of four repeating nucleic bases A,C,G,T.
- Two strands entwined with a helix structure
- Denaturation: At high *T*, strands pulled apart



The other physical motivation is the adsorption phase transition where polymers in solution can stick to the surface of a container

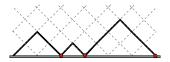
- Second order phase transition
- Order parameter is coverage of the surface by the polymer

Let use models in two dimensions and lattice models at that — Proven to be insightful and integrable

・ロト ・ 同ト ・ ヨト ・ ヨト

ADSORPTION: VERY SIMPLE ONE DIRECTED WALK MODEL

- Single Dyck path in a half space
- Energy $-\varepsilon_a$ for each time (number m_a) it visits the surface
- Boltzmann weight $a = e^{\varepsilon_a/k_B T}$



Consider the coverage, our order parameter (indicator for a phase):

$$\mathcal{A}(a) = \lim_{n \to \infty} \frac{\langle m_a \rangle}{n}$$

There exists two phases, desorbed and adsorbed, and a phase transition at a temperature T_a given by a = 2 between these:

- For $T > T_a$ (small *a*) the walk moves away entropically and A = 0
- For $T < T_a$ (large *a*) the walk is adsorbed onto the surface and A > 0

・ロト ・ 同 ト ・ ヨ ト ・ ヨ ト

ADSORPTION AND UNZIPPING

(Our system — Adsorption and Unzipping)

- double DNA strand in a solvent
- near attractive surface
- assume aligned base sequence
- so expect both adsorption and denaturation (unzipping)

E

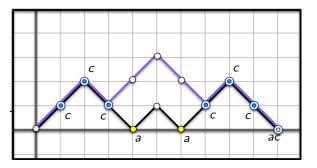
・ロト ・ 日 ・ ・ ヨ ト ・ ヨ ト

Conclusion

ALLOWED WALKS

Consider two directed walks along the square lattice. Let our model contain the class of allowed configs. with *n* steps as described:

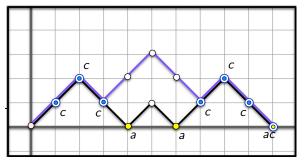
- both walks begin at (0, 0), end at (2*n*, 0).
- directed: can only take steps in the $(\pm 1, 0)$ directions.
- (∞) friendly: walks can share sites, but cannot cross



UNZIPPING ADSORPTION MODEL

Let *T* be the system temperature, k_B the Boltzmann constant.

- surface visit sites: $a \equiv e^{\varepsilon_a/k_BT}$
- shared site contacts: $c \equiv e^{\varepsilon_c/k_B T}$
- trivial walk consisting of zero steps has weight 1.



An allowed configuration of length 10. The overall weight is a^3c^7

E:

• • २ २ २ Owczarek

・ロト ・日 ・ ・ ヨ ・ ・ ヨ ・

Conclusion

GENERATING FUNCTION

- Partition function: $Z_n(a,c) = \sum_{\widehat{\varphi} \ni |\widehat{\varphi}|=n} a^{m_a(\widehat{\varphi})} c^{m_c(\widehat{\varphi})}$
- Generating function: $G(a,c) \equiv G(a,c;z) = \sum_{n \ge 1} Z_n(a,c) z^n$
- Reduced free energy:

$$\kappa(a,c) = \lim_{n \to \infty} n^{-1} \log Z_n(a,c) = \log z_s(a,c)$$

where $z_s(a, c)$ is dominant singularity of *G* w.r.t. *z*

Two order parameters:

$$\mathcal{A}(a,c) = \lim_{n \to \infty} \frac{\langle m_a \rangle}{n}$$
 and $\mathcal{C}(a,c) = \lim_{n \to \infty} \frac{\langle m_c \rangle}{n}$,

Exact Solution of a Simple Adsorption Model of De-naturating DNA

Owczarek

CONTEXT

No wall/interaction:

- Vicious dir. walks: Lindström-Gessel-Viennot thm. ('85), Essam & Guttmann ('95), Guttmann, Owczarek & Viennot ('98)
- Friendly walks & Osculating walks: Brak ('97), Guttmann & Vöge ('02), Bousquet-Mélou ('06)

With wall (but no interaction)

• Vicious: Krattenhaler, Guttmann & Viennot ('00)

With wall (interaction)

- Vicious: Brak, Essam & Owczarek ('99, '01)
- Friendly: Owczarek, Rechnitzer & Wong ('12)

3

・ロト ・ 同 ト ・ ヨ ト ・ ヨ ト

GENERALISED GENERATING FUNCTION

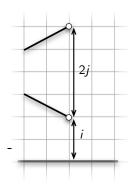
We consider walks φ in the larger set, where each walk can end at any possible height.

- To find G(a, c), consider larger class of configs.
- Generalised generating function:

F

$$\begin{aligned} (\mathbf{r},\mathbf{s}) &\equiv F(\mathbf{r},\mathbf{s},a,c;z) \\ &= \sum_{\varphi \in \Omega} a^{m_a(\varphi)} c^{m_c(\varphi)} \mathbf{r}^i \mathbf{s}^j z^n \end{aligned}$$

• G(a,c) = F(0,0)



・ロト ・ 同ト ・ ヨト ・ ヨト

ntroduction and motivations	ntroductio	on and r	motivations
-----------------------------	------------	----------	-------------

Obstinate. kernel method

ESTABLISHING A FUNCTIONAL EQUATION

- By considering the addition of a single column onto a configuration, and the types of walks obtained, we can find a decomposition of all configurations
- Translating back to generating functions we end up with

$$\begin{aligned} K(r,s)F(r,s) &= \frac{1}{ac} + \left(C(c) - \frac{zr}{s}\right)F(r,0) \\ &+ \left[A(a) - \frac{z}{r}\left(s+1\right)\right]F(0,s) - A(a)C(c)F(0,0) \end{aligned}$$

where

$$A(a) = \frac{a-1}{a}, \quad C(c) = \frac{c-1}{c}$$

and the kernel K(r, s) is

$$K(r,s) \equiv K(r,s;z) = \left(1 - z\left[r + \frac{s}{r} + \frac{r}{s} + \frac{1}{r}\right]\right).$$

・ロト ・日 ・ ・ 日 ・ ・ 日 ・

Mishna Contemp. Math. 520 (2010)

OBSTINATE KERNEL METHOD

- Equation is written as "bulk = boundary terms" where bulk term is
 - product of kernel and bulk generating function

Obstinate, kernel method

- Answer needed is one of the boundary generating functions so try to remove bulk by setting the value of a catalytic variable to a value that makes the kernel vanish
- Standard kernel method due to Knuth (1968): use values of "catalytic variable' to "kill" kernel
- From \approx early '00's applied to a number of dir. walk problems
- Sometimes need multiple values of catalytic variable(s): obstinate kernel method
- More than one catalytic variable requires this
- Earliest combinatorial application of the obstinate kernel method due to Bousquet-Mélou ('02).
- See Bousquet-Mélou Math. and Comp. Sci 2 (2002)), Bousquet-Mélou,



SYMMETRIES OF THE KERNEL

The kernel is symmetric under the following two transformations, which are involutions:

$$(r,s)\mapsto \left(r,\frac{r^2}{s}\right),$$
 $(r,s)\mapsto \left(\frac{s}{r},s\right)$

Transformations generate a family of 8 symmetries ('group of the walk')

$$(r,s), \left(r, \frac{r^2}{s}\right), \left(\frac{s}{r}, \frac{s}{r^2}\right), \left(\frac{r}{s}, \frac{1}{s}\right), \left(\frac{1}{r}, \frac{1}{s}\right), \left(\frac{1}{r}, \frac{s}{r^2}\right), \left(\frac{r}{s}, \frac{r^2}{s}\right), \text{ and } \left(\frac{s}{r}, s\right)$$

- We make use of four of these which only involve positive powers of r.
- This gives us four equations.
- One can eliminate many of the unknown generating functions by a clever choice of adding these equations

ROOTS OF THE KERNEL

- The kernel has two roots as function of either *r* or *s*
- choose the one which gives a positive term power series expansion in *z*
- with Laurent polynomial coefficients in *s* (*r*):

$$\hat{r}(s;z) \equiv \hat{r} = \frac{s\left(1 - \sqrt{1 - 4\frac{(1+s)^2 z^2}{s}}\right)}{2(1+s)z} = \sum_{n>0} C_n \frac{(1+s)^{2n+1} z^{2n+1}}{s^n},$$

where $C_n = \frac{1}{n+1} \binom{2n}{n}$ is a Catalan number.

• Make the substitution $r \mapsto \hat{r}$

Exact Solution of a Simple Adsorption Model of De-naturating DNA

Owczarek

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □

FINDING THE SOLUTION

Key idea

- Treat *K* as fn. of *r* or *s* to get roots \hat{r} and \hat{s}
- Then use subset of \mathcal{F} to get system of eqns. E.g. Using \hat{r} :

(\hat{r},s)	$F(\hat{r},0)$	F(0,s)	F(0, 0)
$(\hat{r},\hat{r}^2/s)$	$F(\hat{r},0)$	$F(0,\hat{r}^2/s)$	F(0, 0)
$(\hat{r}/s,\hat{r}^2/s)$	$F(\hat{r}/s,0)$	$F(0,\hat{r}^2/s)$	F(0, 0)
$(\hat{r}/s, 1/s)$	$F(\hat{r}/s,0)$	F(0, 1/s)	F(0, 0)

• Combine these eqns. to get new fn. eqn

$$N_1^{\star}(s;z)F(0,1/s) + N_2^{\star}(s;z)F(0,s) = \left[M^{\star}(s) - c^2 H^{\star}(s;z)\right] \left(\frac{1}{ac} - ACF(0,0)\right),$$

- Can do the same using $\hat{s}!$
- Nice things happen when a = 1 or c = 1 to $N_1^{\star}(s; z)$ etc

SOLUTION FOR G(a, 1)

Exact solution for G(a, 1) is known and can be found using the method described

- Brak, Essam & Owczarek (1998, 2001): Partition fn. using Lindström-Gessel-Viennot Thm.
- Owczarek, Rechnitzer & Wong (2012): Gen. fn calculated by employing same kernel method techniques.

Specifically:

$$G(a,1) = \sum_{n\geq 0} z^{2n} \sum_{k=0}^{n} a^k \frac{k(k+1)(k+2)}{(2n-k)(n+1)^2(n+2)} {2n-k \choose n} {2n \choose n}.$$

E

・ロト ・ 日 ・ ・ ヨ ト ・ ヨ ト

SOLUTION FOR G(1, c)

• No known previous solution for G(1, c)

We can write functional equation as

$$G(1,c) = F(0,0,1,c;z) = [r^1] \frac{\hat{s} (r^2 - 1) [r - cr + cz (1 + r^2 - \hat{s})]}{(c-1) (\hat{s} - c\hat{s} + crz)},$$

expanding RHS as power series in *c* and so obtain, after some work:

$$\begin{split} &G(1,c;z) = 1 + c^2 z^2 + c^3 \left(1 + 2z\right) z^4 \\ &+ \sum_{i=3}^{\infty} z^{2i} \sum_{m=3}^{2i} c^m \sum_{k=3}^m (-1)^{k+1} \frac{k(k-1)(k-2)(2i-k+1)(i-k+2)}{i^2(i-1)^2(i+1)(i-2)} \binom{m}{k} \binom{2i-k}{i-2} \binom{2i-k-1}{i-3}. \end{split}$$

Solution for G(1, c)

- While we have an explicit solution for *G*(1, *c*) it is advantageous for analysis to directly read off the singularities
- Alternative find differential equation satisfied by generating function
- Use Zeilberger-Gosper algorithm: Maple: DETools package, Zeilberger hyperexp. implementation
- Result: DE for G(1, c) is order 6 with poly. coeff of deg_z = 12

FORTUNATE DECOMPOSITION OF G(a, c)

Using various combinatorial relationships between the generating functions we can re-write G(a, c) in terms of G(a, 1) and G(1, c):

$$G(a,c) = \frac{1}{(a-1)(c-1)} + \frac{p_1(a,c,z)}{p_2(a,c,z) + p_3(a,c,z)G(a,1) + p_4(a,c,z)G(1,c)}$$

where p_i are polynomials in a, c and z: quadratics in z^2 .

Key point: With solutions to G(a, 1) and G(1, c) we additionally have solved for G(a, c).

SINGULARITIES OF G(a, 1) & G(1, c)

- Recall, free energy $\kappa(a, c) = \log z_s(a, c)$
- For G(a, 1), prev. known:

$$z_s(a,1) = \begin{cases} z_b \equiv 1/4, & a \le 2\\ z_a \equiv \frac{\sqrt{a-1}}{2a}, & a > 2 \end{cases}$$

• For *G*(1, *c*), we use the DE (roots of leading poly. coeff.):

$$z_s(1,c) = \begin{cases} z_b \equiv 1/4, & c \le 4/3\\ z_c \equiv \frac{1-c+\sqrt{c^2-c}}{c}, & c > 4/3 \end{cases}$$

Exact Solution of a Simple Adsorption Model of De-naturating DNA

(ロ) (同) (E) (E) (E) (E)

Analysis of solution

Conclusion

TRANSITIONS OF G(a, 1) & G(1, c)

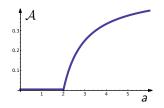
• For *G*(*a*, 1): the order parameter associated with the phase transition is the surface coverage

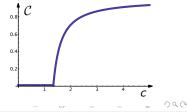
$$\mathcal{A}(a,1) = \begin{cases} 0, & a \le 2\\ \frac{a-2}{2(a-1)}, & a > 2 \end{cases}$$

• For *G*(1, *c*): the order parameter associated with the phase transition is the shared site density

$$\mathcal{C}(1,c) = \begin{cases} 0, & c \leq 4/3 \\ \frac{c-2+\sqrt{c(c-1)}}{2(c-1)}, & c > 4/3 \end{cases}$$

• Second-order adsorption and zipping phase trans. resp.







SINGULARITIES AND PHASES

This leads us to associate the singularities of G(a, 1) and G(1, c) with the phases as

• $z_b = 1/4$ with a desorbed phase where $\mathcal{A} = 0$ and $\mathcal{C} = 0$

•
$$z_a = \frac{\sqrt{a-1}}{2a}$$
 with an adsorbed phase where $A > 0$

•
$$z_c = \frac{1-c+\sqrt{c^2-c}}{c}$$
 with a zipped phase where $C > 0$

ORDER PARAMETERS FOR THE FULL MODEL

Four possible phases:

- Desorbed: A = C = 0
- Adsorbed: (a-rich) A > 0, C = 0
- Zipped: (c-rich) $\mathcal{A} = 0, \mathcal{C} > 0$
- Zipped & Adsorbed: (ac-rich) A > 0, C > 0

1

ANALYSING G(a, c)

Recall

$$G(a,c) \sim \frac{p_1(a,c,z)}{p_2(a,c,z) + p_3(a,c,z)G(a,1) + p_4(a,c,z)G(1,c)}$$

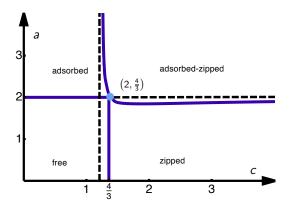
- \Rightarrow Singularities: Look at G(a, 1), G(1, c) and root of above denom.
- root of denominator is associated with the zipped-adsorbed phase

The dominant singularity $z_s(a, c)$ of the generating function G(a, c; z) is one of four types associated with the four phases

$$z_{s}(a,c) = \begin{cases} z_{b} \equiv 1/4, & a \leq 2, c \leq 4/3 \\ z_{a}(a) \equiv \frac{\sqrt{a-1}}{2a}, & a > 2, c \leq \alpha(a) \\ z_{c}(c) \equiv \frac{1-c+\sqrt{c^{2}-c}}{c}, & a \leq \gamma(c), c > 4/3 \\ z_{ac}(a,c), & a > \gamma(c), c > \alpha(a) \end{cases}$$

- $\alpha(a)$ is boundary between adsorbed and zipped-adsorbed phases
- $\gamma(c)$ is the boundary between zipped and zipped-adsorbed phases

PHASE DIAGRAM



All transitions found to be second order

Low-temp argument gives

•
$$c \to \infty, \gamma(c) \to 2$$

•
$$a \to \infty$$
, $\alpha(a) \to \sqrt{5} - 1$

Ξ

Conclusion

ASYMPTOTICS

Table : The growth rates of the coefficients $Z_n(a, c)$ modulo the amplitudes of the full generating function G(a, c; z) over the entire phase space.

phase region	$Z_n(a,c) \sim$
free	$4^{n}n^{-5}$
free to adsorbed boundary	$4^{n}n^{-3}$
free to zipped boundary	$4^{n}n^{-3}$
a = 2, c = 4/3	$4^{n}n^{-3}$
adsorbed	$z_a(a)^{-n}n^{-3/2}$
zipped	$z_c(c)^{-n}n^{-3/2}$
adsorbed to adsorbed-zipped boundary ($\alpha(a)$)	$z_a(c)^{-n}n^{-1/2}$
zipped to adsorbed-zipped boundary $(\gamma(c))$	$z_c(c)^{-n}n^{-1/2}$
adsorbed-zipped	$z_{ac}(a,c)^{-n}n^{-1}$

Ξ

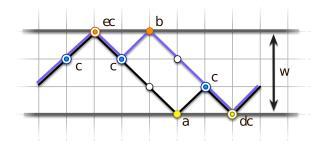
CONCLUSION

- Simple model of DNA as two friendly walks near a boundary
- Used combinatorial decomposition to obtain linear functional equation
- Used obstinate kernel method to solve functional equations (using symmetries to provide sufficient information)
- Explicit series solutions for G(a, 1) and G(1, c)
- Combined these equations to relate G(a, c) to both G(a, 1) and G(1, c)
- Also used Zeilberger-Gosper algorithm to find linear DE for G(1, c)
- Full analysis of asymptotics and phase diagram
- R. Tabbara, A. L. Owczarek and A. Rechnitzer, J. Phys. A.: Math. Theor, 47, 015202 (34pp), 2014

・ロト ・ 日 ・ ・ 日 ・ ・ 日 ・ ・

FUTURE WORKS

- Combine single, double surface and unzipping interactions: work in progress
- Consider three walks with multiple unzipping interactions: work in progress
- Also in progress is work in a slit: see Thomas Wong's talk
- Extend these to various interactions



(ロ) (同) (E) (E) (E) (E)