

# General Exam

## Modeling Cooperation Between Molecular Motors Polymer Growth Against a Force

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February 9, 2006

## Introduction

- Conventional Molecular Motors
- Polymerization as a Molecular Motor
- Why Do We Care?

## Current Research

- Preliminary Model Set-Up
- Polymer Growth Simulations
- More Complicated Systems

## Future Research

# What are Molecular Motors?

## Introduction

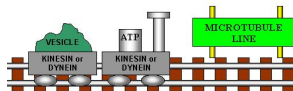
Conventional  
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### Protein molecules in the cell that:

- ▶ generate force
- ▶ cause transport



# Where are Molecular Motors?

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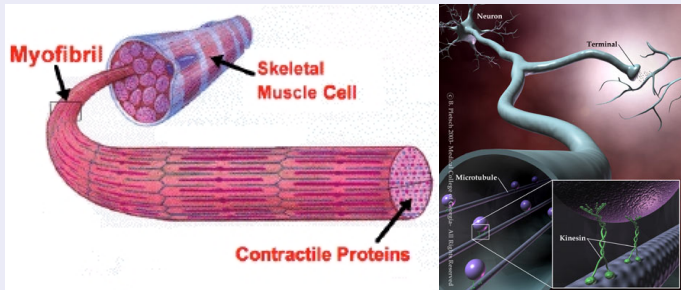
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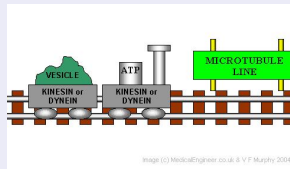
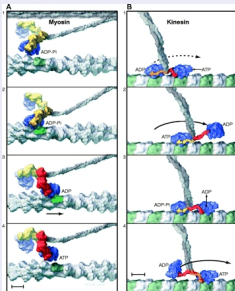
## Muscle Cells & Neural Cells



## Conventional Molecular Motors

move along polymer tracks

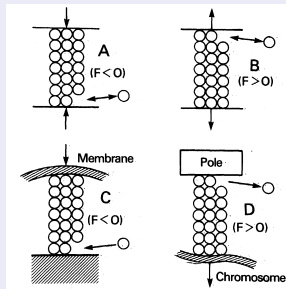
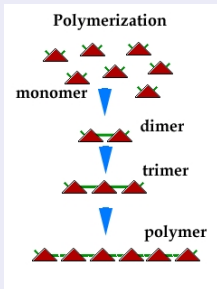
- ▶ myosin - actin microfilaments
- ▶ kinesin - tubulin microtubules



# Polymerization as a Motor

## Another way to cause motion/transport

- ▶ POLYMERIZATION-or-DEPOLYMERIZATION !
- ▶ (adding or subtracting monomers)



# Polymerization as a Motor - Biological Examples

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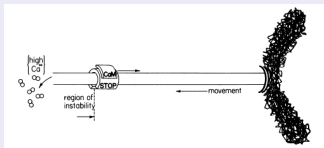
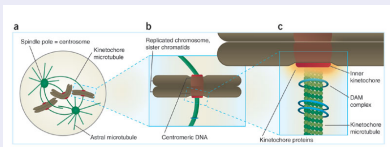
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## Chromosome Transport During Anaphase

Depolymerization of Spindle Pulls Sister Chromatids Apart:



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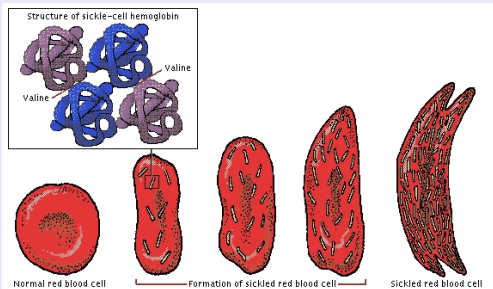
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## Cell Membrane Deformation

Sickle Hemoglobin Polymerization creates Sickle Cells:





# Why Do We Care About Molecular Motors?

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## Molecular Motors are Special Because:

- ▶ Chemical Energy  $\Rightarrow$  Mechanical Energy
  - ▶ DIRECTLY! (not via heat or electrical energy)
- ▶ Highly Efficient
  - ▶ 6 times more efficient than a car
- ▶ Models for Molecular Motors  $\Rightarrow$  Nano-Engineering of Future
  - ▶ Nano-mechano-chemical Machines
  - ▶ Tiny Robots!



## Introduction

Conventional Molecular Motors

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# Basic Polymer Model

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## How does Polymerization Work?

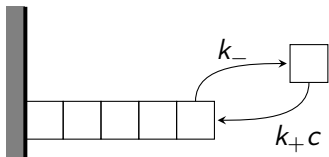
Rate Constants:

$k_+$ : second order rate constant of adding a monomer

$k_-$ : first order rate constant of subtracting a monomer

$c$ : concentration of monomers in surrounding solution

Position of the end of the polymer can be modeled as a 1-D biased random walk.



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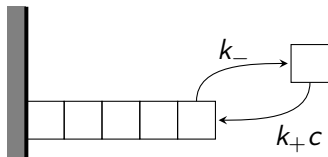
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## How does Polymerization Work?

$P_n(t)$ : probability length is  $n$  at time  $t$

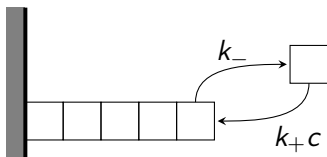
Differential Equation for Polymer Length:

$$\frac{dP_n(t)}{dt} = k_+ c P_{n-1}(t) + k_- P_{n+1}(t) - (k_+ c + k_-) P_n(t)$$



## How does Polymerization Work?

- ▶ Deterministic System:
  - ▶ Motion is continuous in Space, Time
  - ▶ Initial Condition  $\Rightarrow$  one possible trajectory
- ▶ Stochastic System:
  - ▶ Direction of motion, Time motion occurs - Random
  - ▶ Initial Condition  $\Rightarrow$  many possible trajectories



# Basic Polymer Model

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## Stochastic System: Continuous-Time Random Walk

Number of Events in time  $t$  is modeled as a Poisson Process with rate:

$$\lambda = k_+c + k_-$$

Times between Events have an Exponential Distribution with rate  $\lambda$ .

Probability of subtracting or adding a monomer:

$$P(-) = \frac{k_-}{k_- + k_+c} = \frac{k_-}{\lambda}$$

$$P(+) = \frac{k_+c}{k_- + k_+c} = \frac{k_+c}{\lambda}$$

⇒ Use this idea to create simulations!

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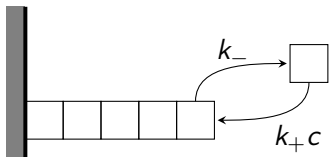
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# Simulation with One Polymer

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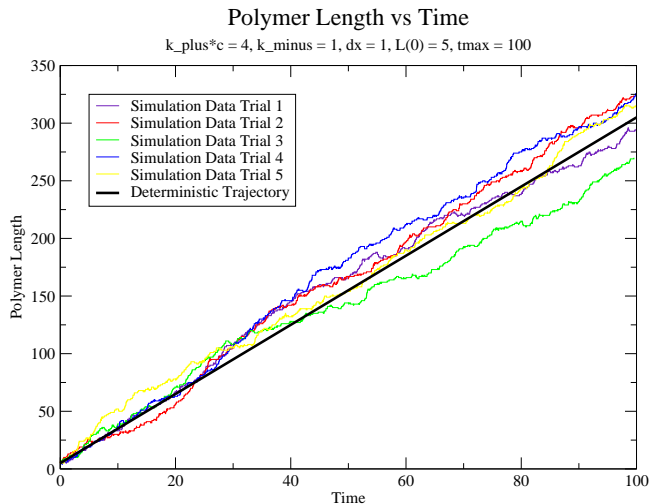
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# Polymer Model with a Moving Wall

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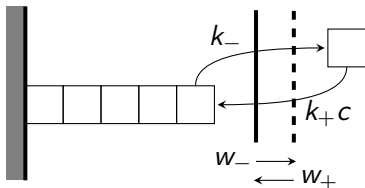
## Polymer Interacting with a Moving Wall

Let  $w$  be the position of the moving wall.

- ▶  $w_+$  - rate that the wall moves towards the polymers
- ▶  $w_-$  - rate the wall moves away from the polymers

Additional Constraint, if  $w - x < \Delta x$ :

- ▶ monomer cannot be added.
- ▶ wall cannot move towards the polymer.





# Simulation with One Polymer and a Moving Wall

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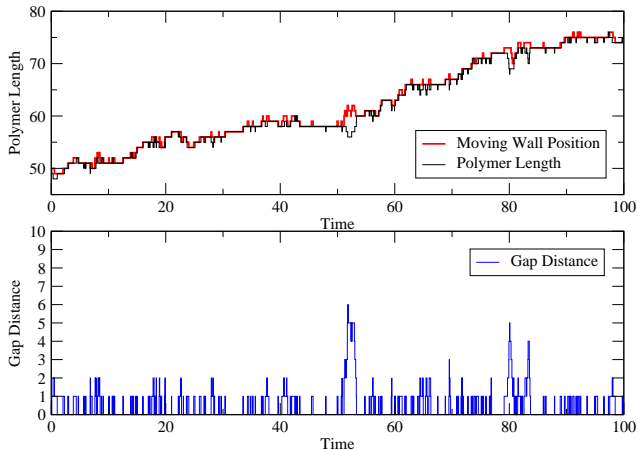
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## Polymer Length and Gap Distance vs Time

$k_{\text{plus}} * c = 4$ ,  $k_{\text{minus}} = 1$ ,  $w_{\text{plus}} = 2$ ,  $w_{\text{minus}} = 1$ ,  $t_{\text{max}} = 10,000$



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## Modeling the Gap Distance

- ▶  $\alpha$  - rate gap distance shrinks

$\beta$  - rate gap distance grows

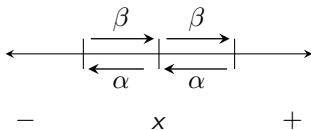
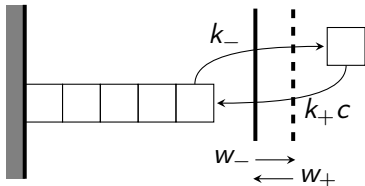
$$\alpha = k_+c + w_+ \quad \beta = k_- + w_-$$

- ▶  $p(x, t)$ : probability that the gap distance is  $x$  at time  $t$

$$\beta p_t(x, t) = \alpha p(x + \Delta x, t) + \beta p(x - \Delta x, t) - (\alpha + \beta)p(x, t)$$

- ▶ Solve for Steady-State:

$$\Rightarrow p = p(x), \quad p_t = 0$$



# Polymer Model with a Moving Wall

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## Model the Steady-State Gap Distance

- ▶ Discrete Space  $\Leftrightarrow$  Random Walk

$$p_i = \frac{\alpha}{\alpha+\beta} p_{i+1} + \frac{\beta}{\alpha+\beta} p_{i-1} \quad i > 0$$

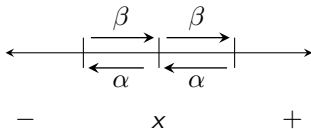
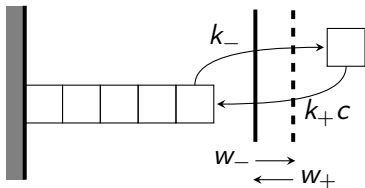
$$\frac{\beta}{\alpha+\beta} p_0 = \frac{\alpha}{\alpha+\beta} p_1 \quad (\text{B.C. for } i=0)$$

- ▶ Continuous Space Limit  $\Leftrightarrow$  Brownian Motion

$$p_t = D p_{xx} + V p_x = 0 \quad x > 0$$

$$D p_x + V p = 0 \quad (\text{B.C. for } x=0)$$

$$D = \lim_{\Delta x \rightarrow 0} (\alpha + \beta) \frac{(\Delta x)^2}{2} \quad V = \lim_{\Delta x \rightarrow 0} (\alpha - \beta) \Delta x$$



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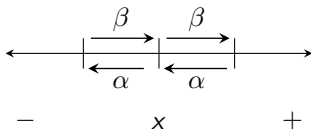
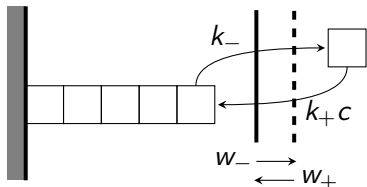
## Steady-State Gap Distance

- ▶ Discrete Space  $\Leftrightarrow$  Random Walk

$$\Rightarrow p_i = \frac{\alpha - \beta}{\alpha} \left( \frac{\beta}{\alpha} \right)^i$$

- ▶ Continuous Space Limit  $\Leftrightarrow$  Brownian Motion

$$\Rightarrow p(x) = \frac{V}{D} e^{-\frac{V}{D}x}$$



# Moving Wall Steady State Gap Distance

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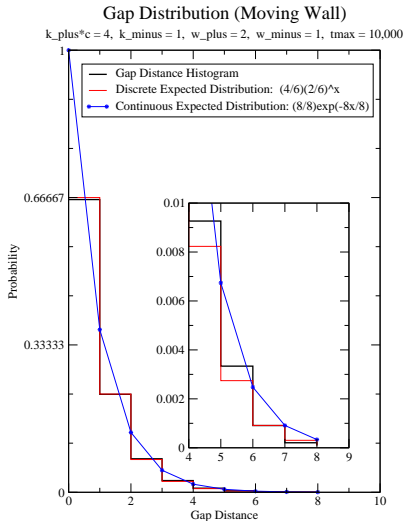
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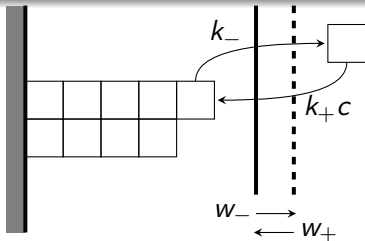
## Multiple Polymers Interacting with a Moving Wall

Build upon the basic model and simulations to study a more interesting system:

- ▶  $N$  Polymers

Gap distances between each polymer and the wall can be modeled as an N-D biased random walk.

- ▶ Special Case:  $N=2$



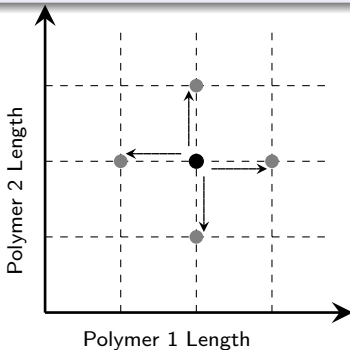
# More Complicated Polymer Model

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## Multiple Polymers Interacting with a Moving Wall

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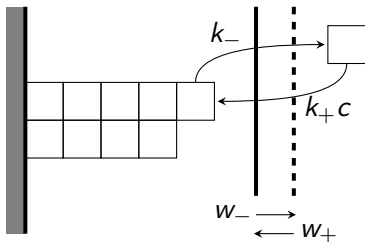
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## Multiple Polymers Interacting with a Moving Wall

- ▶ Gillespie-type Algorithm Generates:
  - ▶ Position of each Polymer Tip
  - ▶ Position of the Moving Wall
- ▶ Simulations can be used to:
  - ▶ Investigate the System
  - ▶ Compare with Theoretical Results





# Simulation with Many Polymers

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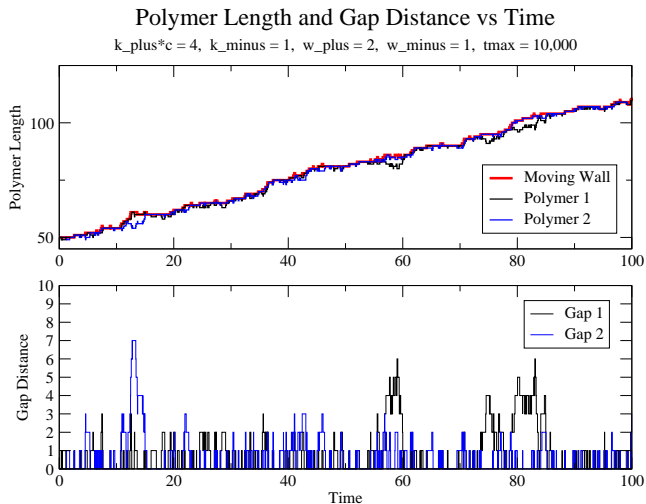
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# More Complicated Polymer Model

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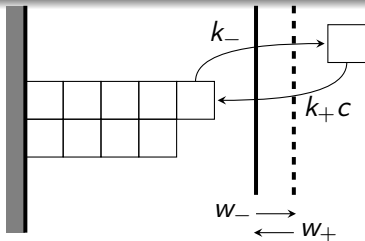
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## Multiple Polymers Interacting with a Moving Wall

- ▶ The Polymers are *Identical*:
  - ▶  $k_+ c_{p1} = k_+ c_{p2} = k_+ c$
  - ▶  $k_- p1 = k_- p2 = k_-$
- ▶ Polymers do not *Explicitly* Interact
- ▶ Are the Polymers *Independent*?
  - ▶ *Independent*  $\Rightarrow p_i = \frac{\alpha - \beta}{\alpha} \left( \frac{\beta}{\alpha} \right)^i$



# Steady State Gap Distribution for 2-Polymer System

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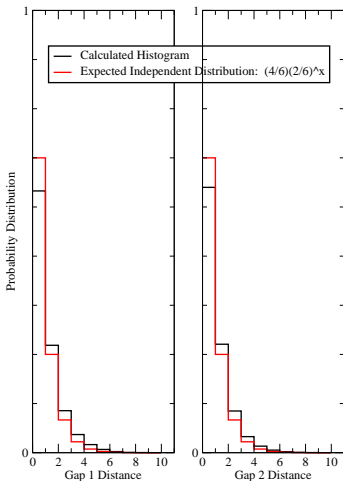
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Steady State Distribution for Gaps 1 & 2  
 $k_{plus} * c = 4$ ,  $k_{minus} = 1$ ,  $w_{plus} = 2$ ,  $w_{minus} = 1$ ,  $t_{max} = 10,000$



# Polymer Cooperation - 2D Random Walk

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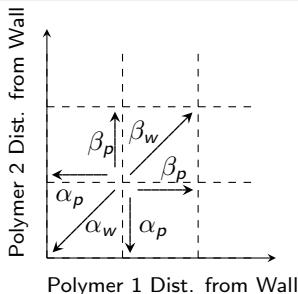
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## Polymer Motion from Wall's POV - Gap Distances

Rates of motion are given by:

polymer moves	wall moves	gaps
$\alpha_p$ - towards wall	$\alpha_w$ - towards polymer	shrink
$\beta_p$ - away from wall	$\beta_w$ - away from polymer	grow

(Origin represents both polymers touching the wall)



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# Polymer Cooperation - 2D Random Walk

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$p(x, y, t)$  - gap 1 distance is  $x$ , gap 2 distance is  $y$ , at time  $t$

- ▶ PDE for Gap Distance Probability

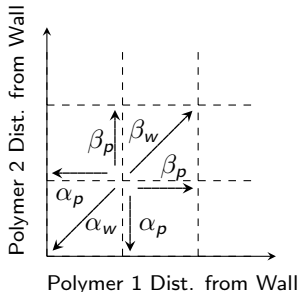
- ▶  $p_t = D_1(p_{xx} + p_{yy}) + 2D_2p_{xy} + V(p_x + p_y)$

- ▶ No-Flux Boundary Conditions:

- ▶  $J_1(0, y) = D_1p_x + D_2p_y + Vp = 0$

- ▶  $J_2(x, 0) = D_1p_y + D_2p_x + Vp = 0$

- ▶ Solve for Steady-State Solution



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## Research to be Done

- ▶ Finish analysis of 2-Polymer Model
- ▶ Generalize to N-Polymer Model
  - ▶ Gap Distances  $\Rightarrow$  N+1-D PDE for  $p(x_1, x_2, \dots, x_N, t)$
  - ▶ N spatial, 1 time
- ▶ Mathematical & Biophysical Analysis of N-Polymer Model
- ▶ Study Density Function for an N-Polymer Model
  - ▶ Gap Distances  $\Rightarrow$  2-D PDE for  $c(x, t)$
  - ▶ 1 space, 1 time

# Questions?

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