Physical Principles in Biology Biology 3550 Spring 2024

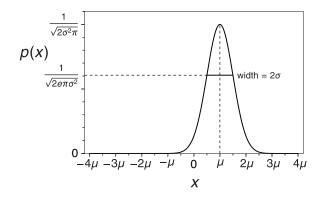
Lecture 15:

# The Gaussian Probability Distribution Function and

Introduction to Diffusion

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#### The Gaussian Probability Distribution Function

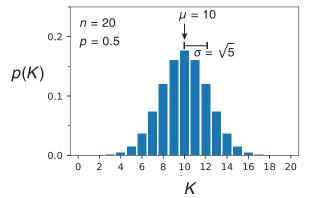


$$p(x)=rac{1}{\sqrt{2\pi\sigma^2}}e^{-rac{(x-\mu)^2}{2\sigma^2}}$$

- Also called the normal probability distribution function.
- Mean  $= \mu$
- Variance =  $\sigma^2$
- Standard deviation  $= \sigma$

#### Mean and Variance for The Binomial Distribution

I



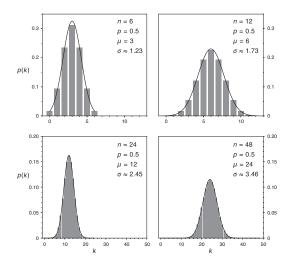
Mean  

$$\mu = np_s$$
  
Variance  
 $\sigma^2 = np_s(1 - p_s)$ 

Gaussian probability function to approximate the binomial distribution function with the same mean and variance:

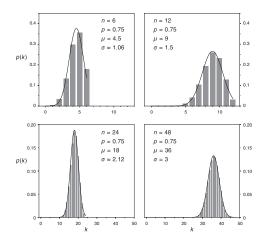
$$p(k) = rac{1}{\sqrt{2\pi\sigma^2}}e^{-rac{(k-\mu)^2}{2\sigma^2}} = rac{1}{\sqrt{2\pi n p_{\mathsf{s}}(1-p_{\mathsf{s}})}}e^{-rac{(k-np_{\mathsf{s}})^2}{2np_{\mathsf{s}}(1-p_{\mathsf{s}})}}$$

#### Approximation of Binomial Distributions by Gaussian Distributions



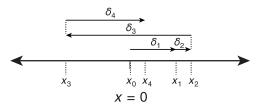
n doesn't have to be very large for a pretty good approximation!

#### Approximation of Binomial Distributions by Gaussian Distributions



- It doesn't work so well if the binomial distribution is biased, with *p*<sub>s</sub> ≠ 0.5.
- The Gaussian distribution is always symmetrical, but the binomial distribution only is if p<sub>s</sub> = 0.5.
- If n is large enough, the Gaussian distribution is a good approximation, even if p<sub>s</sub> ≠ 0.5.

The One-dimensional Random Walk with Continuously Variable Step Length



Mean displacement:

$$\langle \delta 
angle = E(\delta) = \int_{\delta_{\mathsf{min}}}^{\delta_{\mathsf{max}}} p(\delta) \delta d\delta$$

Mean-squared displacement:

$$\langle \delta^2 
angle = E(\delta^2) = \int_{\delta_{\mathsf{min}}}^{\delta_{\mathsf{max}}} p(\delta) \delta^2 d\delta$$

If 
$$\langle \delta \rangle = 0$$
:  
 $\langle x_n \rangle = 0$   
 $\langle x_n^2 \rangle = n \langle \delta^2$ 

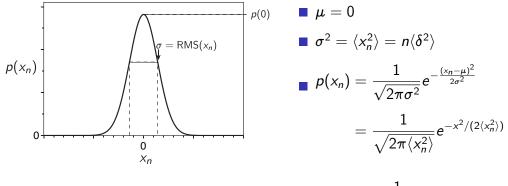
• The mean and variance of  $x_n$ :

$$\mu = \int_{x_{n,\min}}^{x_{n,\max}} p(x_n) x_n = \langle x_n \rangle = 0$$

$$\sigma^2 = \int_{x_{n,\min}}^{x_{n,\max}} p(x_n)(x_n-\mu)^2$$

$$=\int_{x_{n,\min}}^{x_{n,\max}}p(x_n)x_n^2=\langle x_n^2\rangle$$

#### The Gaussian Distribution of End Points for a 1-dimensional Random Walk with Variable Displacements



$$p(0) = \frac{1}{\sqrt{2\pi \langle x_n^2 \rangle}}$$

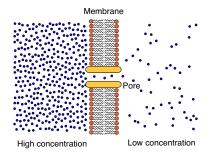
### Warning!



## **Direction Change**

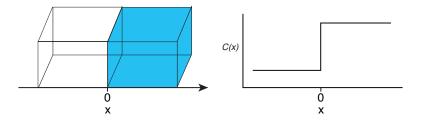
Diffusion

#### Diffusion of Molecules Across a Cellular Membrane



- How do the concentrations change with time?
- WHY do the concentrations change?
- How fast do the concentrations change?
- How fast do the molecules move?

#### An Idealized Macroscopic Diffusion Experiment



• How will plot of C(x) versus x change with time?

There's a theory for that! Fick's laws of diffusion

#### Clicker Question #1

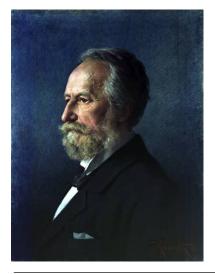
How long has it been since there was a sharp boundary?



- A) 10 min
- **B)** 1 hr
- C) 12 hr
- D) 48 hr

All answers count! (for now)

### Adolf Eugen Fick



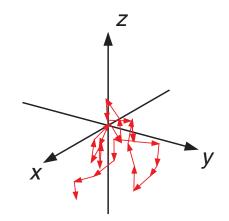
- 1829-1901
- Physician and physiologist
- Nephew, Adolf Gaston Eugen Fick, invented contact lenses in 1888.

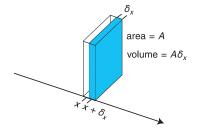
https://en.wikipedia.org/wiki/Adolf\_Eugen\_Fick Portrait by Anton Klamroth.

#### Diffusion as a Random Walk

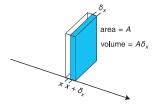
- ⟨*I*<sup>2</sup><sub>*i*</sub>⟩: Mean-square step length in three dimensions
- $\delta_x = \sqrt{\langle l_i^2 \rangle / 3}$ : RMS displacement along the *x*-direction.
- τ: Average time interval between changes in direction
- t: Time interval of interest
- n = t/τ: Average number of steps in time t

$$\ \, |\langle x_n^2\rangle = n\delta_x^2 = t\delta_x^2/\tau$$



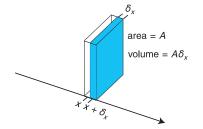


- Slices are  $\delta_x$  thick and have cross-sectional area across the *x*-axis of *A*
- Volume of each slice is  $A\delta_x$
- During time τ, all of the molecules will move (on average) the distance δ<sub>x</sub> along the x-axis, to the left or right.
- In a given slice, 1/2 of the molecules will move to the right and half to the left.



- $N_x$  = number of molecules starting in the slice centered at *x*.
- $N_{x+\delta_x}$  = number of molecules starting in the slice centered at  $x + \delta_x$
- The net number of molecules moving from slice x to slice  $x + \delta_x$ , in time  $\tau$ :

$$dN = rac{1}{2}N_x - rac{1}{2}N_{x+\delta_x}$$
 $= -rac{1}{2}(N_{x+\delta_x} - N_x)$ 



Definition: Flux, J = net number of molecules (or moles) moving past a cross section, per unit time, per unit area.

$$J = -\frac{1}{A\tau} \frac{1}{2} (N_{x+\delta_x} - N_x)$$

Movement is defined in the direction of positive *x*.

Express number of molecules in each slice in terms of the concentrations and volumes of each slice.

$$N_x = C_x \cdot V = C_x A \delta_x$$

$$N_{x+\delta_x} = C_{x+\delta_x} \cdot V = C_{x+\delta_x} A \delta_x$$

Re-write the flux equation as:

$$egin{aligned} J &= -rac{1}{A au}rac{1}{2}ig( extsf{N}_{ extsf{x}+\delta_{ extsf{x}}} - extsf{N}_{ extsf{x}}ig) \ &= -rac{1}{A au}rac{1}{2}ig( extsf{C}_{ extsf{x}+\delta_{ extsf{x}}} extsf{A}\delta_{ extsf{x}} - extsf{C}_{ extsf{x}} extsf{A}\delta_{ extsf{x}}ig) \ &= -rac{\delta_{ extsf{x}}}{ au}rac{1}{2}ig( extsf{C}_{ extsf{x}+\delta_{ extsf{x}}} - extsf{C}_{ extsf{x}}ig) \end{aligned}$$

Write the concentration difference in terms of a derivative with respect to x:

$$\frac{dC}{dx} = \lim_{\delta_x \to 0} \frac{\left(C_{x+\delta_x} - C_x\right)}{\delta_x}$$

in the limit of small  $\delta_{x}$ :

$$\left(\mathsf{C}_{x+\delta_x}-\mathsf{C}_x
ight)=\delta_xrac{dC}{dx}$$

Flux equation:

$$J = -\frac{\delta_x}{\tau} \frac{1}{2} (C_{x+\delta_x} - C_x) = -\frac{\delta_x^2}{2\tau} \frac{dC}{dx}$$

#### Fick's First Law of Diffusion



- Symbols:
  - *J* = flux of molecules per unit area per unit time:
  - $\delta_x = \text{RMS}$  step length along the *x*-direction.
  - $\tau =$  average duration of random steps.
  - $\frac{dC}{dx}$  = derivative of concentration with x, "concentration gradient."
- If concentration increases with *x*, flux is in the negative *x* direction.
- Why do molecules "move down the concentration gradient"?

#### The Diffusion Coefficient, D

Consider the term from Fick's first law:  $\frac{\delta_x^2}{2\tau}$ 

Both  $\delta_{\times}$  and  $\tau$  are parameters describing the random walk of molecules (or larger particles) undergoing diffusion, and are constant for a given type of particle under defined solution conditions.

Define a new parameter, the diffusion coefficient, *D*:

$$D = \frac{\delta_x^2}{2\tau}$$

The usual form of Fick's first law:

$$J = -D\frac{dC}{dx}$$

 D can be experimentally determined, without knowing anything about the microscopic random walk steps.