Welcome to

CHE 384T: Computational Methods in Materials Science

Random Walk Diffusion

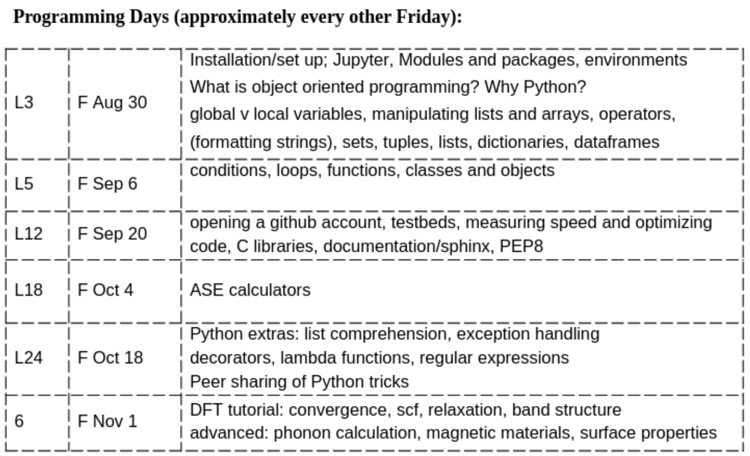
LeSar Ch. 2, App. B7, C5, I2-I3

The University of Texas at Austin
McKetta Department
of Chemical Engineering
Cockrell School of Engineering

Announcements

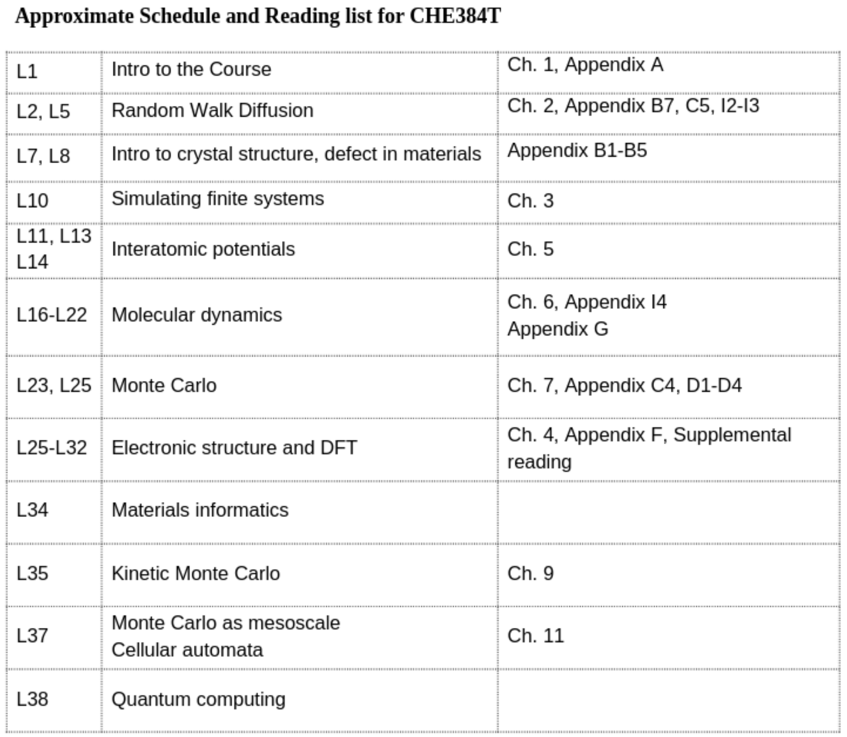
Python Pre-test
Due 08/30 11:59pm

Bring a computer to Friday Lecture



Programming Days (approximately every other Friday):

L3	F Aug 30	Installation/set up; Jupyter, Modules and packages, environments What is object oriented programming? Why Python? global v local variables, manipulating lists and arrays, operators, (formatting strings), sets, tuples, lists, dictionaries, dataframes
L5	F Sep 6	conditions, loops, functions, classes and objects
L12	F Sep 20	opening a github account, testbeds, measuring speed and optimizing code, C libraries, documentation/sphinx, PEP8
L18	F Oct 4	ASE calculators
L24	F Oct 18	Python extras: list comprehension, exception handling decorators, lambda functions, regular expressions Peer sharing of Python tricks
6	F Nov 1	DFT tutorial: convergence, scf, relaxation, band structure advanced: phonon calculation, magnetic materials, surface properties



Approximate Schedule and Reading list for CHE384T

L1	Intro to the Course	Ch. 1, Appendix A
L2, L5	Random Walk Diffusion	Ch. 2, Appendix B7, C5, I2-I3
L7, L8	Intro to crystal structure, defect in materials	Appendix B1-B5
L10	Simulating finite systems	Ch. 3
L11, L13 L14	Interatomic potentials	Ch. 5
L16-L22	Molecular dynamics	Ch. 6, Appendix I4 Appendix G
L23, L25	Monte Carlo	Ch. 7, Appendix C4, D1-D4
L25-L32	Electronic structure and DFT	Ch. 4, Appendix F, Supplemental reading
L34	Materials informatics	
L35	Kinetic Monte Carlo	Ch. 9
L37	Monte Carlo as mesoscale Cellular automata	Ch. 11
L38	Quantum computing	

Lecture Outline

What is diffusion
Examples of diffusion in materials science

Connection with continuum description

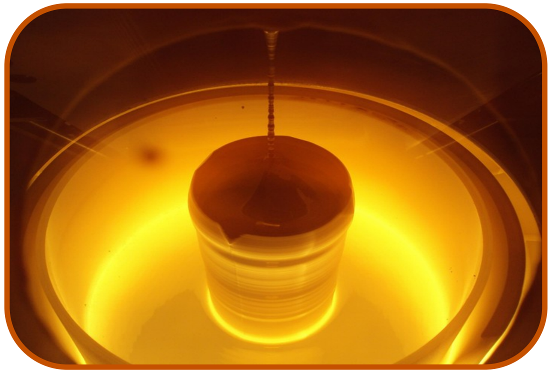
Random Walk model for Diffusion

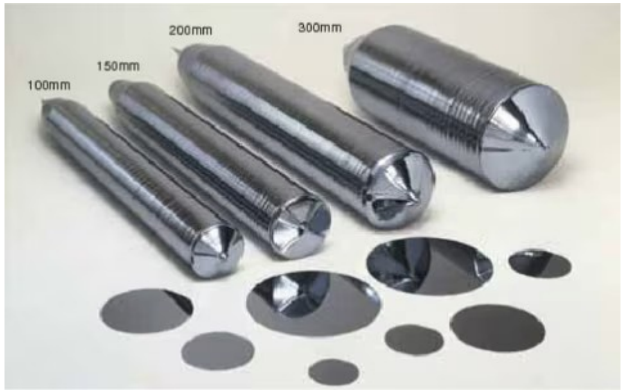
Coding considerations:
- Random number generators
- Binning probability distributions

What is diffusion?

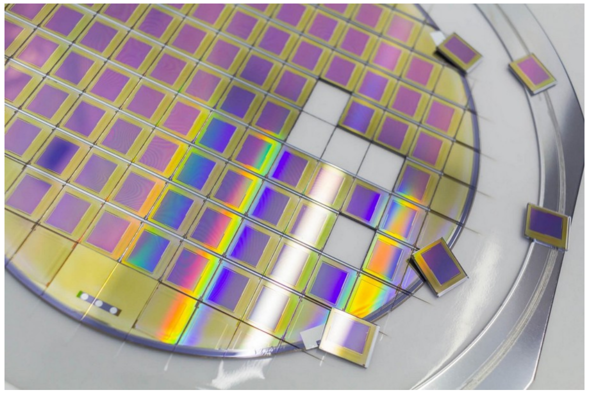
https://www.youtube.com/watch?v=UlnubVuJvZM

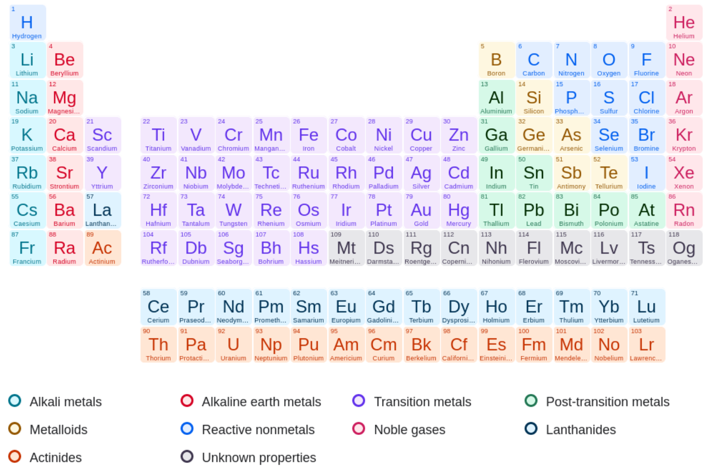
Silicon Wafer Processing: Example of Diffusion in Materials Science





99.999999999% pure Silicon

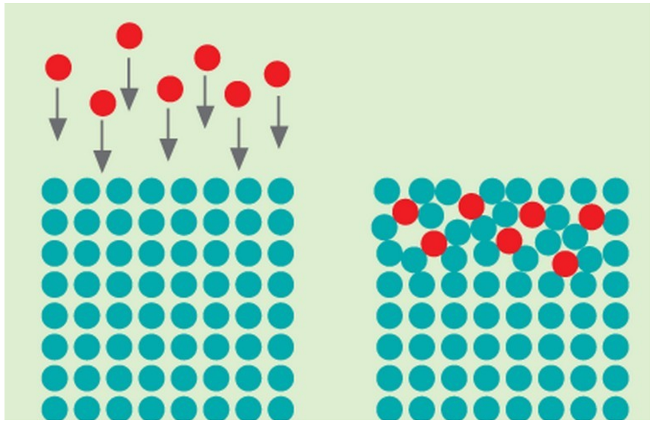




Silicon Wafer Processing:
Example of Diffusion in Materials Science

Intentional incorporation of impurities (e.g., boron, phosphorous)

Step 1: Steady-state gas diffusion
or ion implantation



Step 2: Drive-in process (higher temperature)
→ uniform distribution of impurities

Diffusion of Lithium ions in a battery

Fick’s First and Second Law

See also Appendix B7

Fick’s Second Law: Derivation

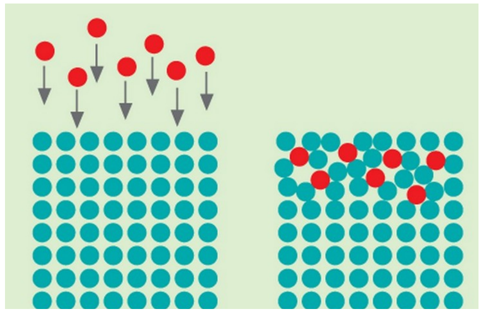
Based on conservation of mass

Fick’s First and Second Law: Interpretation

Example: Silicon wafer processing

Intentional incorporation of impurities (e.g., boron, phosphorous)

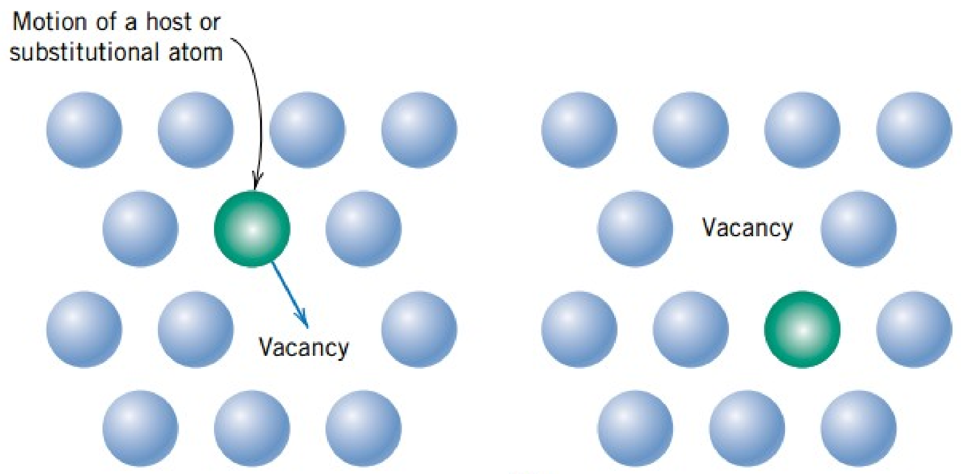
Step 1: Steady-state gas diffusion or ion implantation



Step 2: Drive-in process (higher temperature)

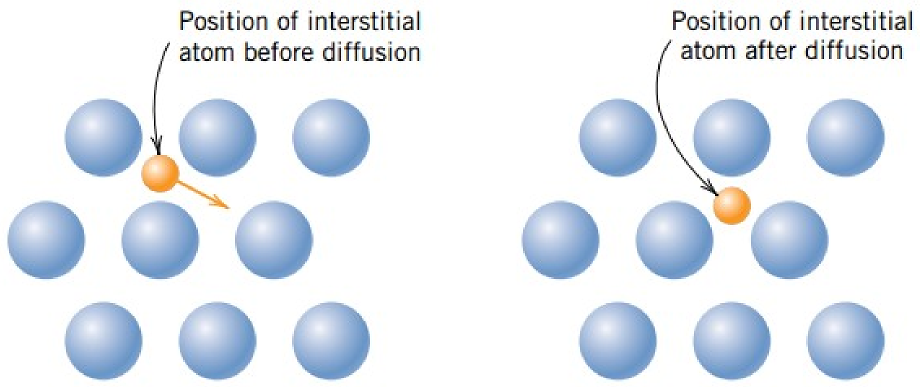
Types of Diffusion

Vacancy diffusion



(a)

Interstitial diffusion


Position of interstitial atom before diffusion
Position of interstitial atom after diffusion

(b)

Factors that affect diffusion: Diffusion coefficient

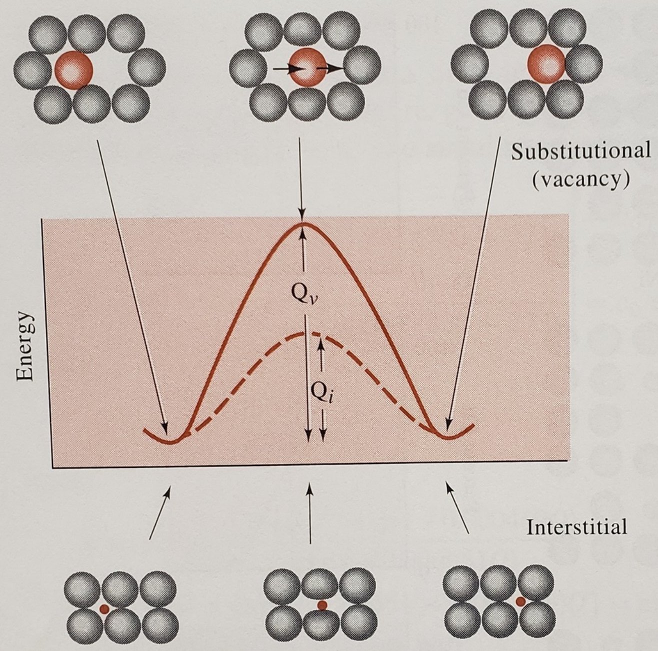
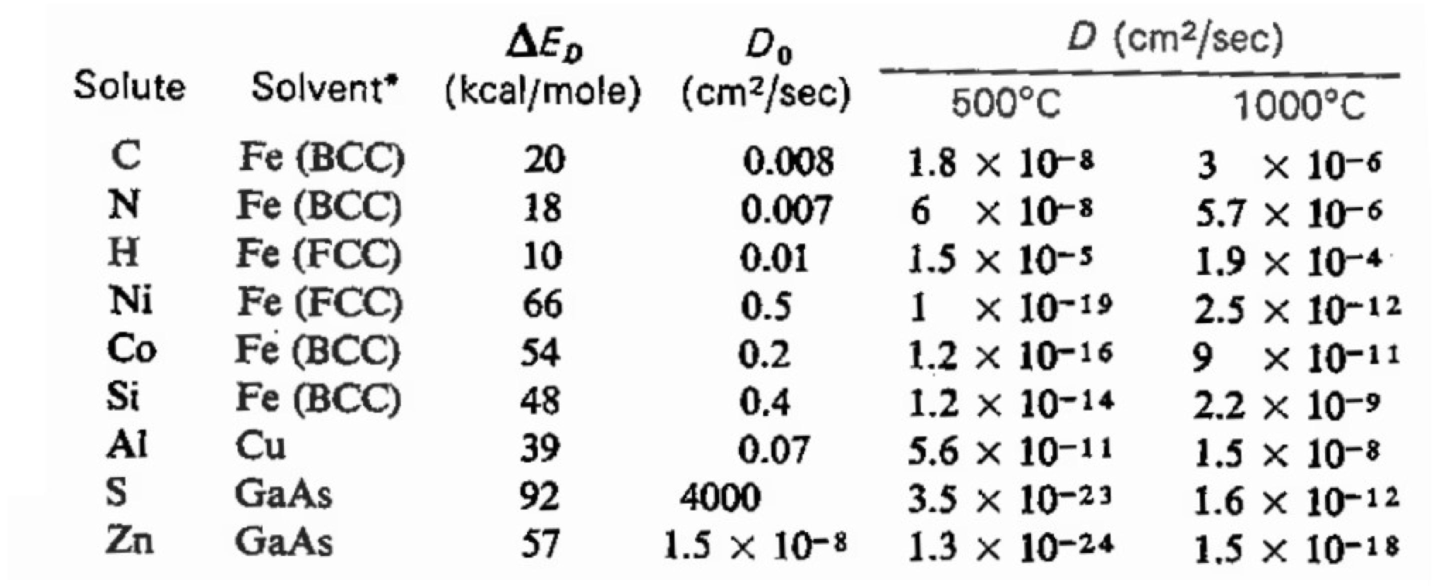


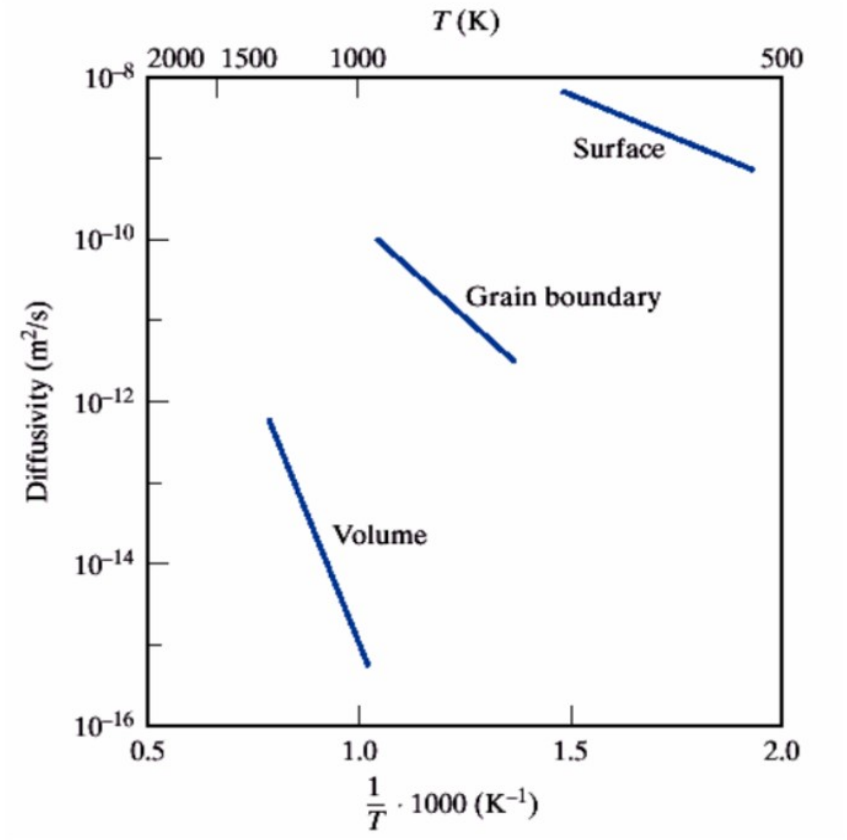
FIGURE 5-4 A high energy is required to squeeze atoms past one another during diffusion. This energy is the activation energy *Q*. Generally more energy is required for a substitutional atom than for an interstitial atom.

Factors that affect diffusion: Diffusion coefficient

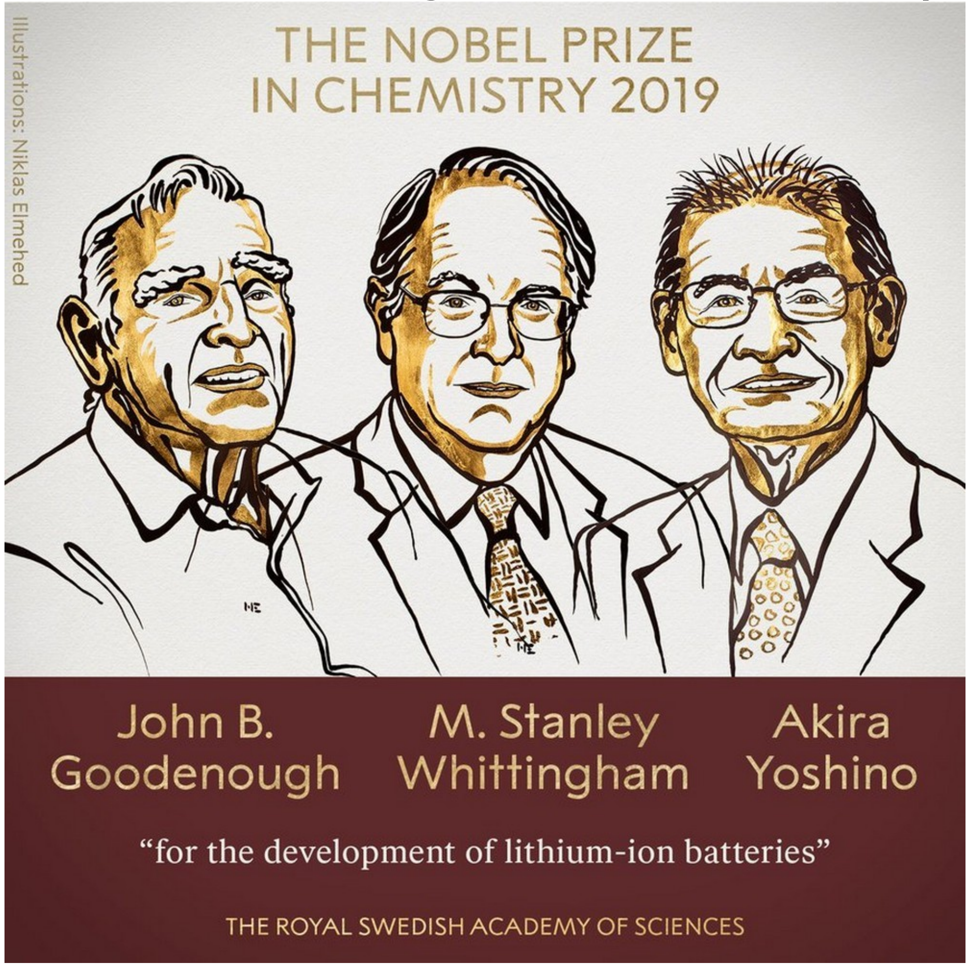


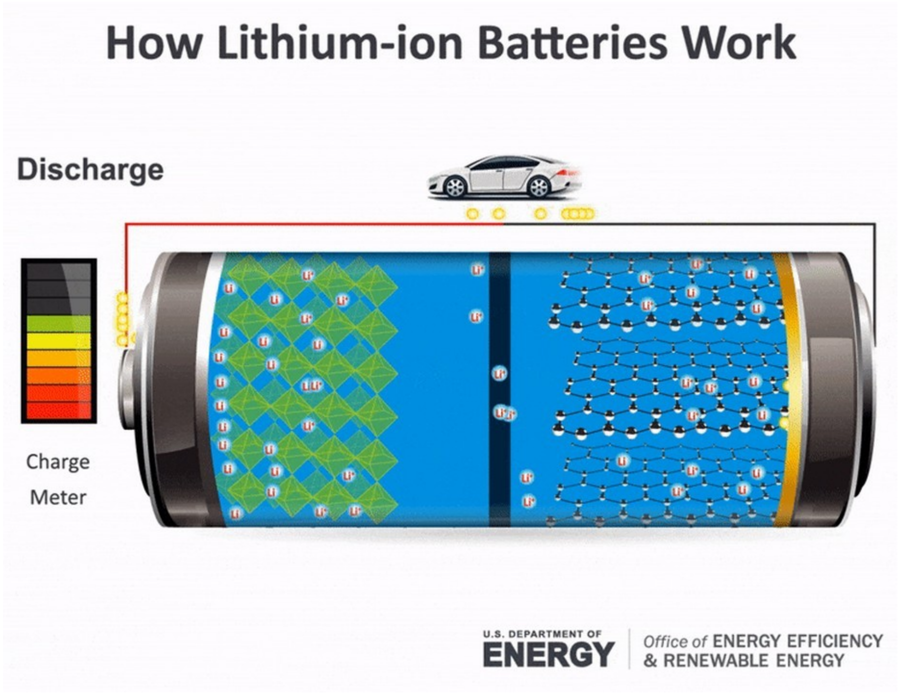
Solute	Solvent*	ΔE_D (kcal/mole)	D_0 (cm²/sec)	D (cm²/sec) 500°C	1000°C
C	Fe (BCC)	20	0.008	1.8×10^{-8}	3×10^{-6}
N	Fe (BCC)	18	0.007	6×10^{-8}	5.7×10^{-6}
H	Fe (FCC)	10	0.01	1.5×10^{-5}	1.9×10^{-4}
Ni	Fe (FCC)	66	0.5	1×10^{-19}	2.5×10^{-12}
Co	Fe (BCC)	54	0.2	1.2×10^{-16}	9×10^{-11}
Si	Fe (BCC)	48	0.4	1.2×10^{-14}	2.2×10^{-9}
Al	Cu	39	0.07	5.6×10^{-11}	1.5×10^{-8}
S	GaAs	92	4000	3.5×10^{-23}	1.6×10^{-12}
Zn	GaAs	57	1.5×10^{-8}	1.3×10^{-24}	1.5×10^{-18}

Factors that affect diffusion: Diffusion coefficient



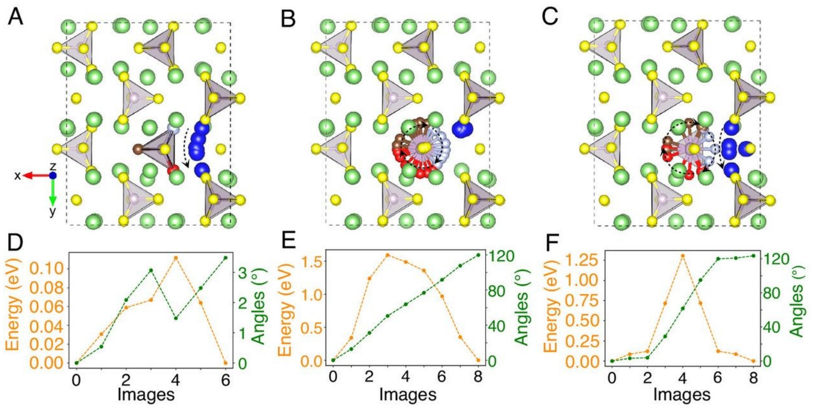
What is the materials science behind the Li-ion battery winning the 2019 Nobel prize in Chemistry?


THE NOBEL PRIZE IN CHEMISTRY 2019
Illustrations: Niklas Elmehed
John B. Goodenough
M. Stanley Whittingham
Akira Yoshino
"for the development of lithium-ion batteries"
THE ROYAL SWEDISH ACADEMY OF SCIENCES

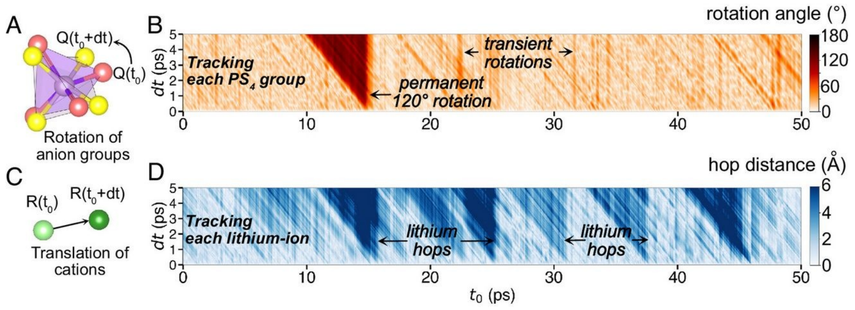

How Lithium-ion Batteries Work
Discharge
Charge Meter
U.S. DEPARTMENT OF ENERGY
Office of ENERGY EFFICIENCY & RENEWABLE ENERGY

https://www.energy.gov/science/doe-explainsbatteries

Collective motion of lithium with crystal lattice

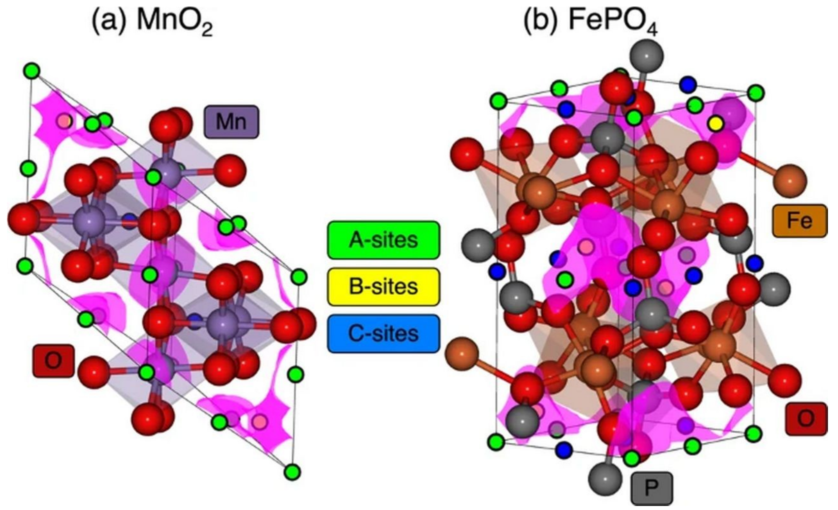


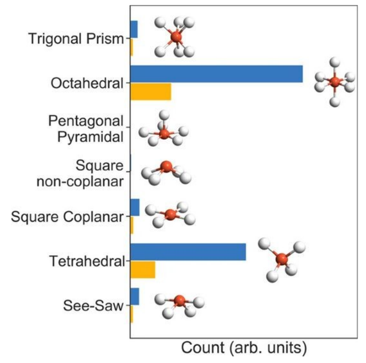
β-Li_3PS_4



PNAS. 121 (18) e2316493121 (2024).

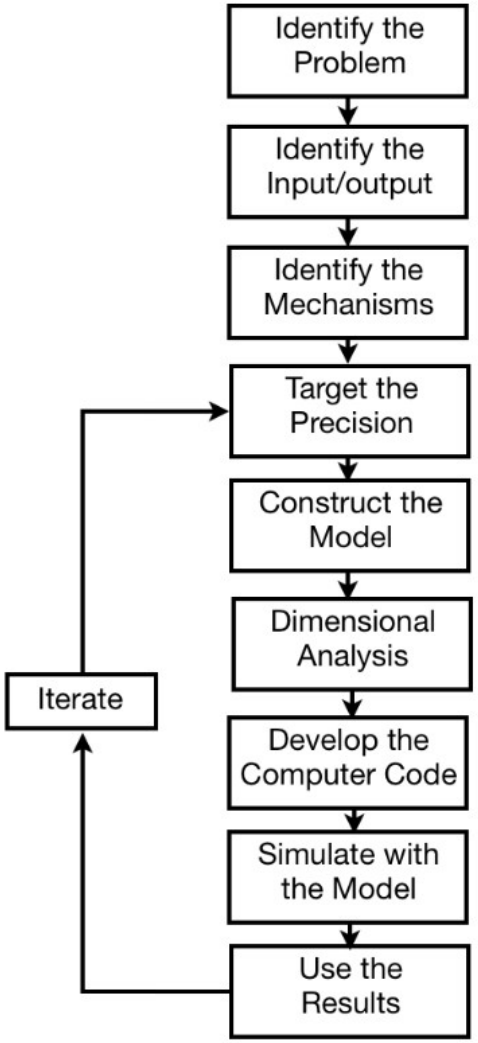
Intercalation of lithium in novel cathode materials





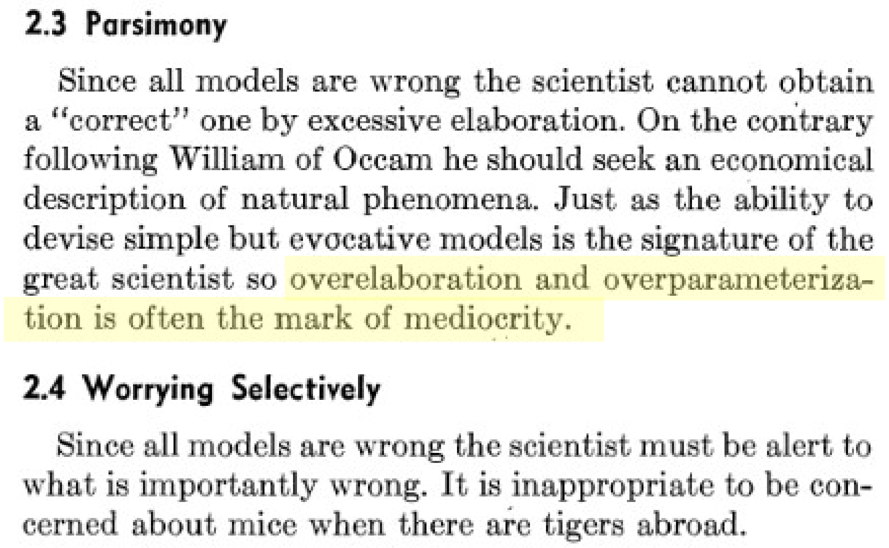
npj Comput Mater 6, 161 (2020).

Developing a model



"Science and Statistics." George E.P. Box (1976)

"All models are wrong, but some are useful."



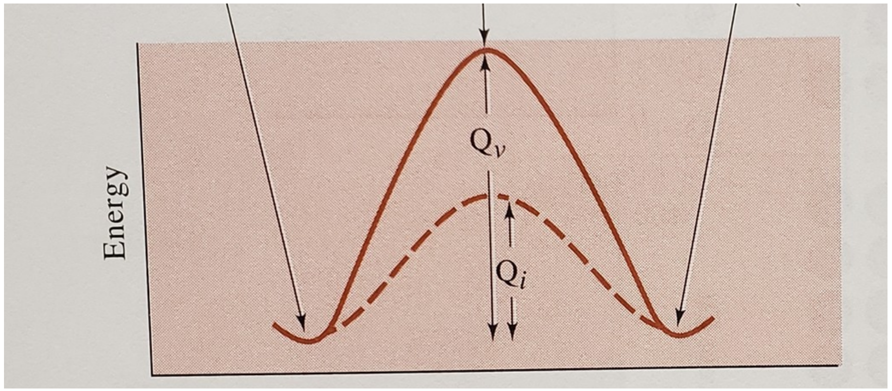
2.3 Parsimony

Since all models are wrong the scientist cannot obtain a "correct" one by excessive elaboration. On the contrary following William of Occam he should seek an economical description of natural phenomena. Just as the ability to devise simple but evocative models is the signature of the great scientist so overelaboration and overparameterization is often the mark of mediocrity.

2.4 Worrying Selectively

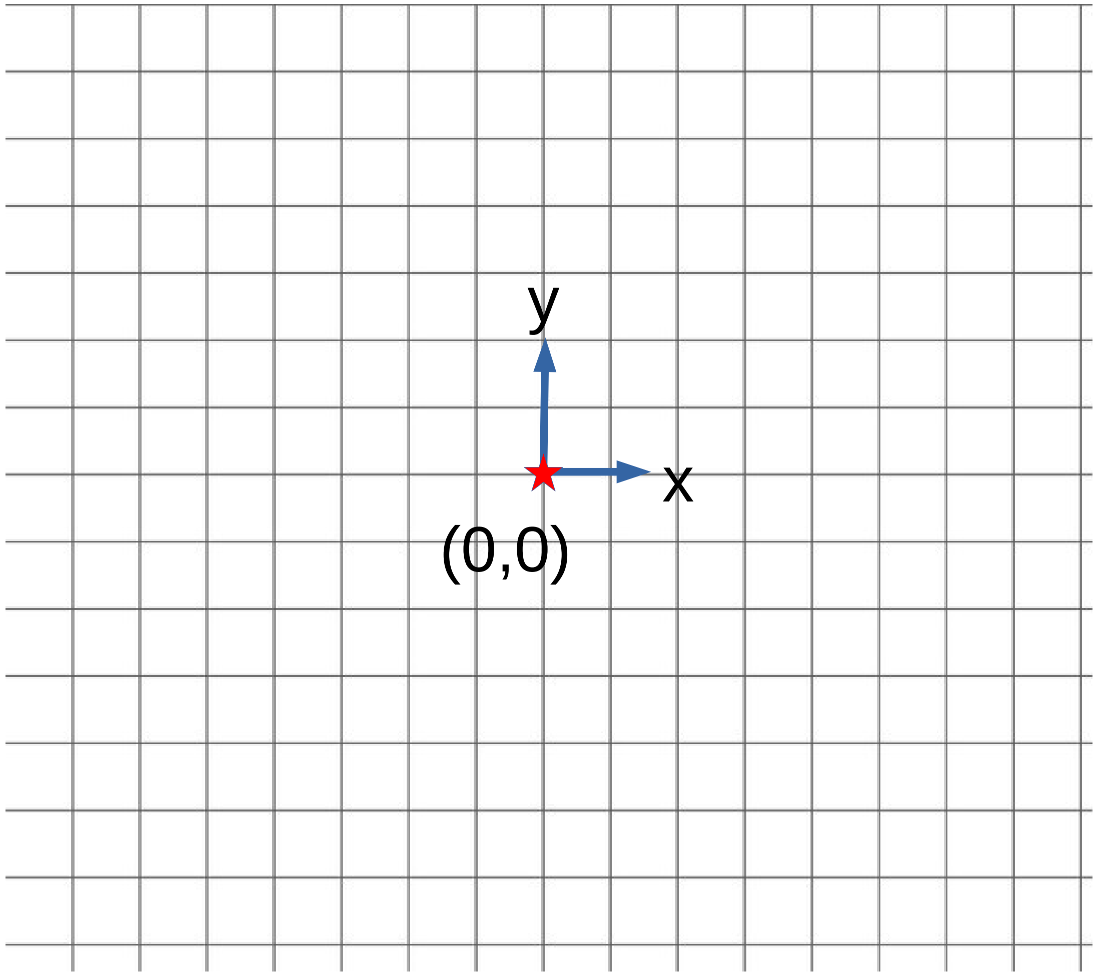
Since all models are wrong the scientist must be alert to what is importantly wrong. It is inappropriate to be concerned about mice when there are tigers abroad.

Simplifications of the Random Walk Model



- Reduced dimensionality (to 2D or 1D)
- Simplification of the crystal structure
- Model the hops the atoms take as random (for a particular atom)

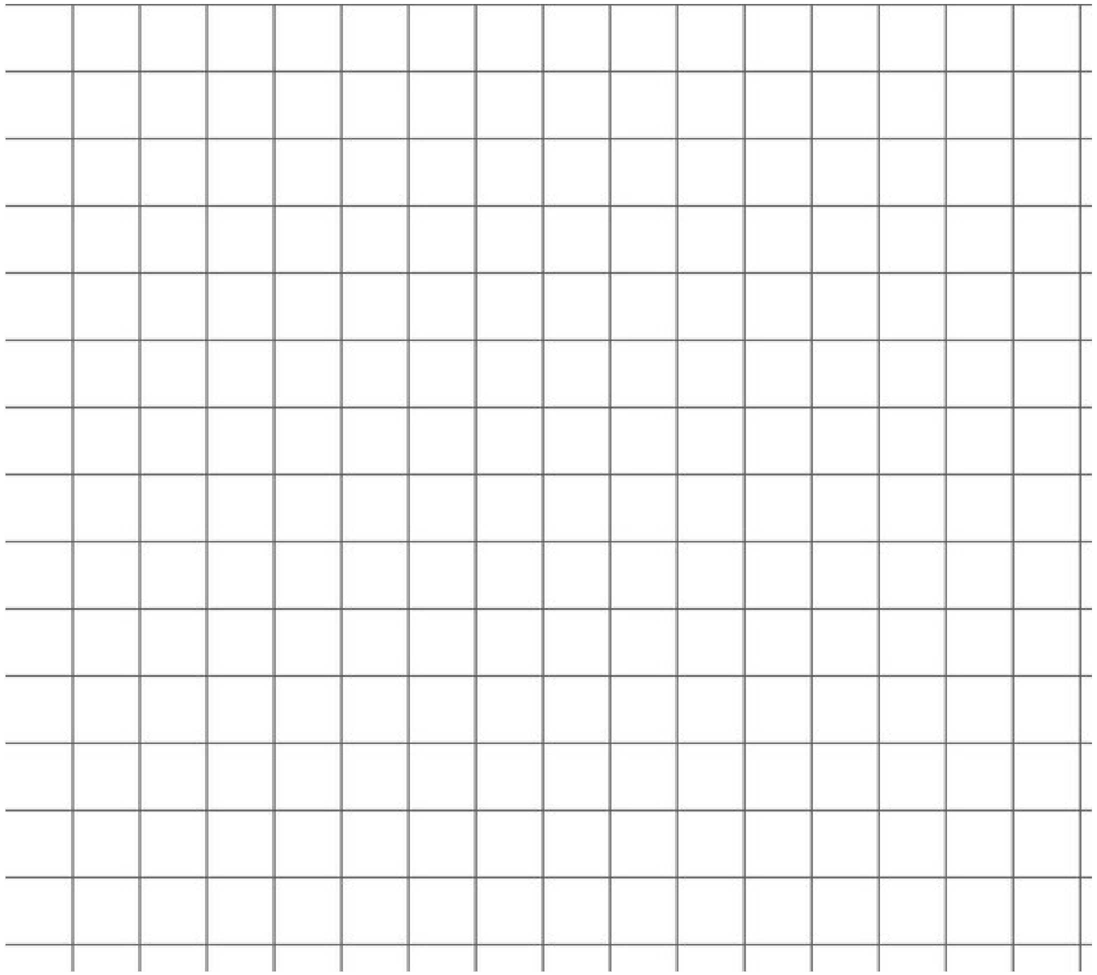
Random walk diffusion: an atomic model for diffusion



Rules for the random walker:

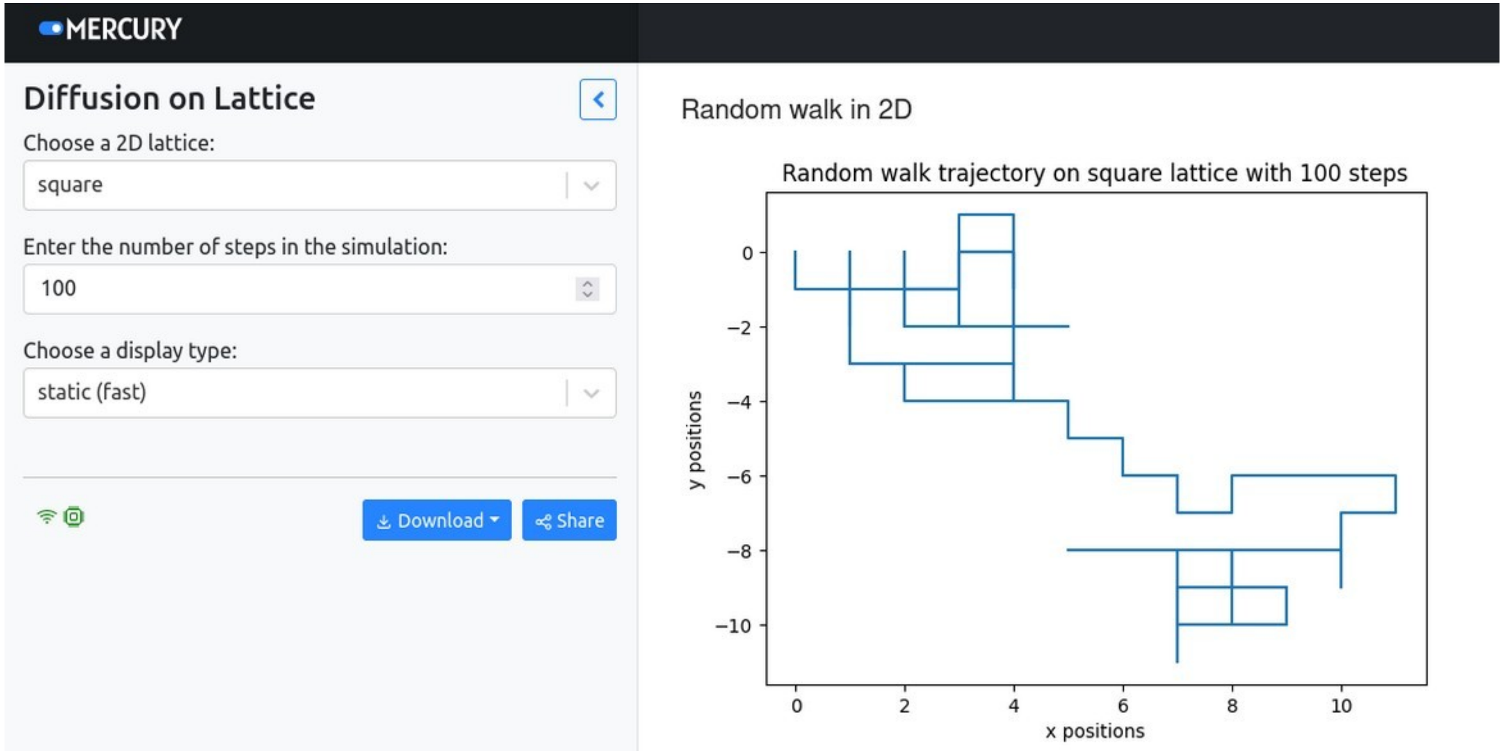
- divide time into *nt* discrete steps spaced by Δt time, where *nt* is an integer and Δt is a number
- can only move 1 space at each time step
- equal and random probability of moving up, down, left, right

Random walk diffusion: an atomic model for diffusion



Random walk diffusion: a small simulation

https://rwd2d-mercury.runmercury.com/


MERCURY

Diffusion on Lattice

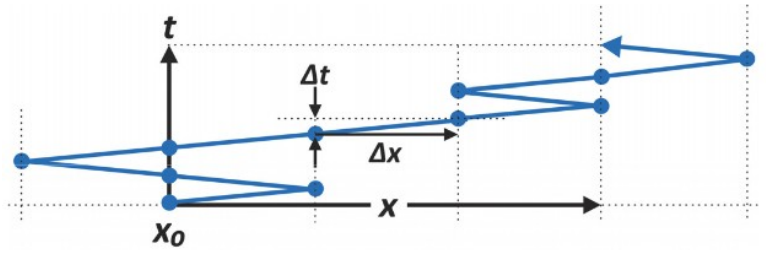
Choose a 2D lattice: square

Enter the number of steps in the simulation: 100

Choose a display type: static (fast)

Connection between Random Walk and Diffusivity

1D case



Connection between Random Walk and Diffusivity

$$D = \frac{1}{6t}\langle R^2 \rangle \quad \text{(in 3D)}$$

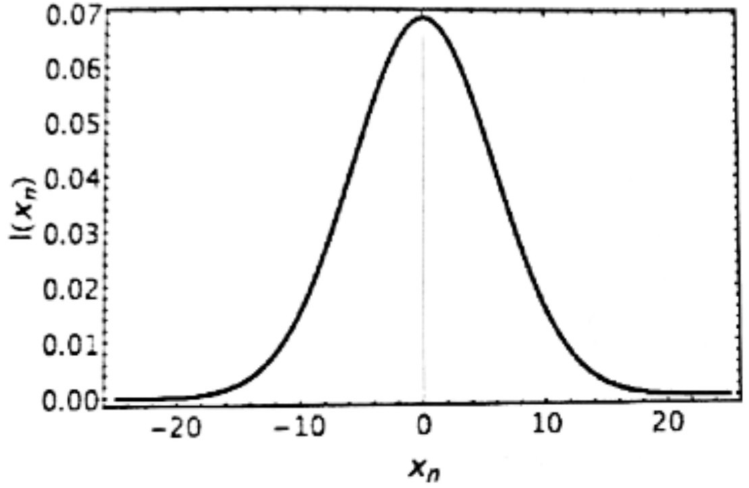
For the random walk model

Statistics of the Random Walk Model: End-to-end distribution

1D case

Statistics of the Random Walk Model: End-to-end distribution

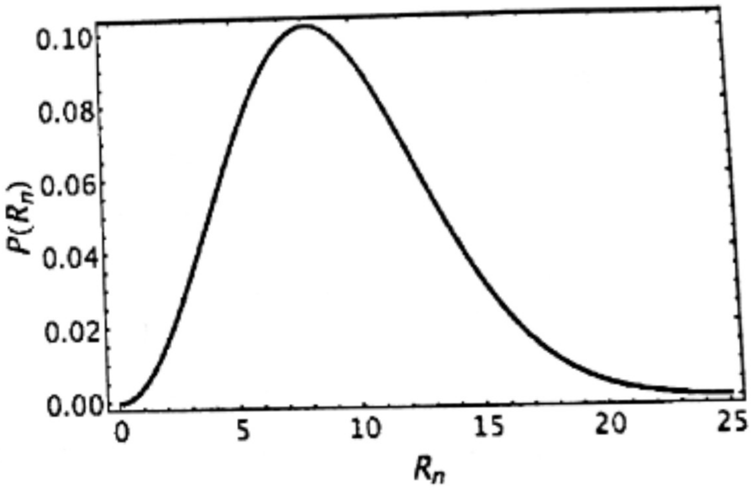
(In 1D)



$$\mathcal{P}(n_{tot}, q) = \frac{1}{2}^{n_{tot}} \frac{n_{tot}!}{(\frac{n_{tot}+q}{2})!(\frac{n_{tot}-q}{2})!}$$

Statistics of the Random Walk Model: End-to-end distribution

(In 3D)



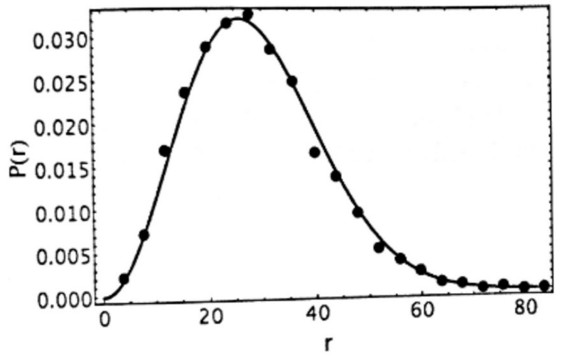
Coding considerations:

Binning distributions

Discretizing continuous functions
Appendix I.3

Choose n_{bin}

$$\Delta = \frac{R_n^{max} - R_n^{min}}{n_{bin}}$$



Random Number Generator

Discretizing continuous functions
Appendix I.2

To reproduce a “random” result,
pick a consistent *seed*

Python:

numpy.random.rand(...): generates (pseudo-)random number over [0,1)

numpy.ceil(…): round to next highest integer

Donald R. Askeland. The Science and Engineering of Materials 3rd ed.

An implementation of the 2D Random Walk Diffusion

Objective	Concept	Python Representation
User chooses *nt* = number of time steps	Variable, integer	`nt`

An implementation of the 2D Random Walk Diffusion

Objective	Concept	Python Representation
User chooses *nt* = number of time steps	Variable, integer	`nt`

Keep track of position of random walker at each time step.
Let's assume it starts at the origin.
	Array (list of items)
	e.g., [3, 4.5, 8, -1]
	
	2D → *x* and *y* coordinate for each position
		Use the library numpy, shorthand is np:
		
		`x = np.zeros(nt+1)`
		`y = np.zeros(nt+1)`

An implementation of the 2D Random Walk Diffusion

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Specify how the position changes at each time step.		`delx = np.array([?,?,?,?])`
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An implementation of the 2D Random Walk Diffusion

Objective	**Concept**	**Python Representation**
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		`dely = np.array([?,?,?,?])`

Save each new position of the diffusion path	Index the array
	i.e., access a specific element
	“zero index”
		`x = [1,2,3]`
		`x[0] = 1`
		`x[1] = 2`

An implementation of the 2D Random Walk Diffusion

Objective	Concept	Python Representation

Repeat for *nt* times	for loop
	range function
		Input:
		`for i in range(3):`
		`    print(i)`
		Output:
		`0`
		`1`
		`2`

Encode the random number to a change in position of the random walker	Generate a (pseudo)-random number	`np.floor(4* np.random.rand(nt))` Generate random number b/t 0 and 1 Random number b/t 0 and 4 Random integer: 0, 1, 2, 3