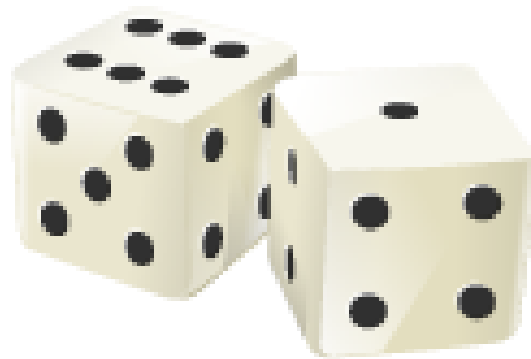


Classical Monte Carlo Simulations



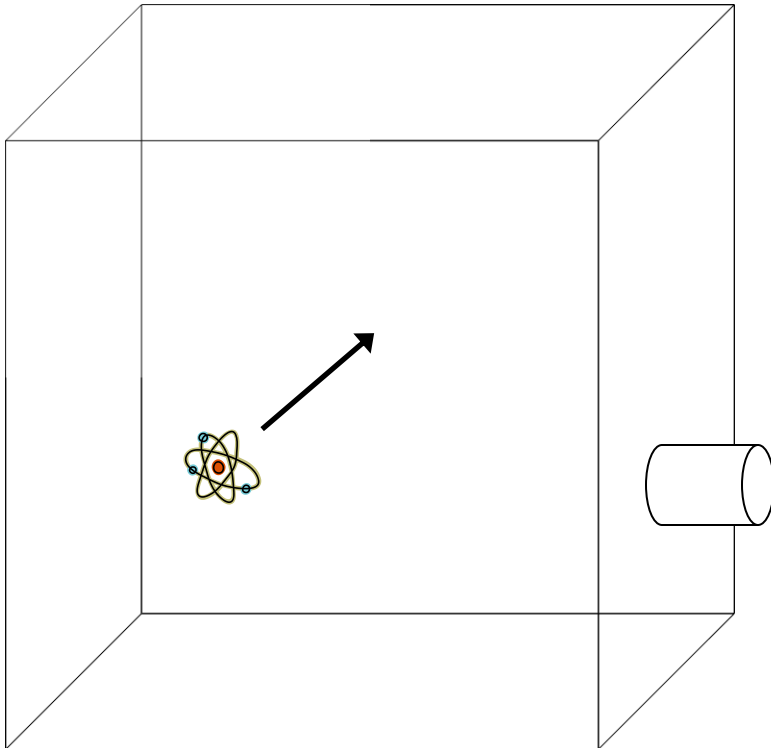
... you can calculate anything with dice.

Introduction

Calculate ...

- the number of **bounces** (mean and variance)
- and **time** (mean and variance)

for a gas molecule at temperature T to leave this box:

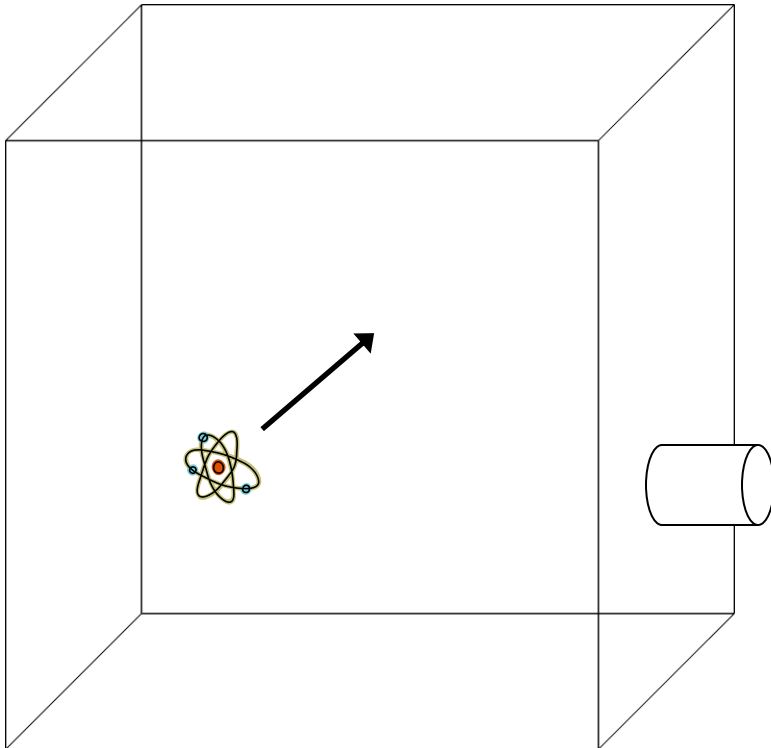


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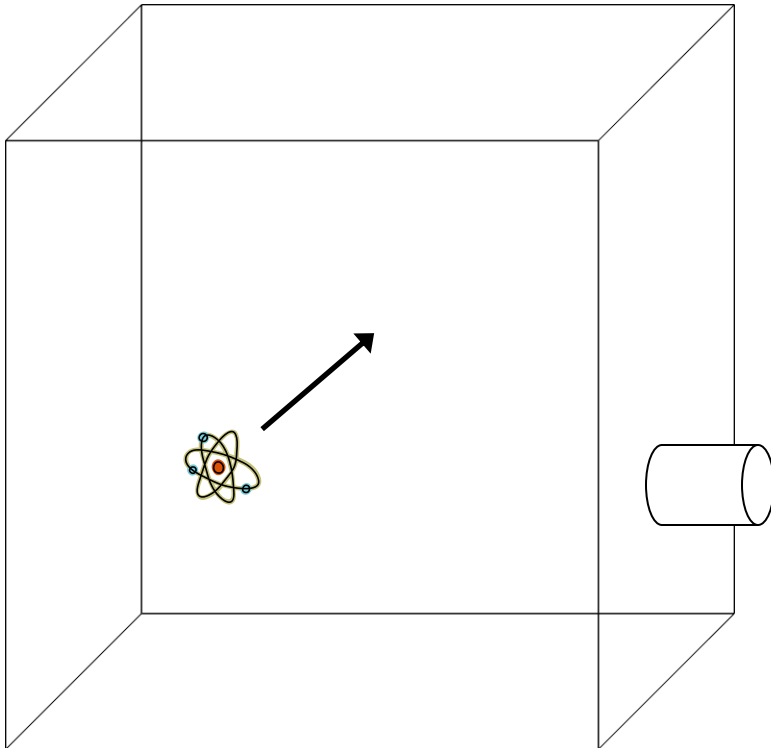
- molecule re-thermalize on each wall bounce.
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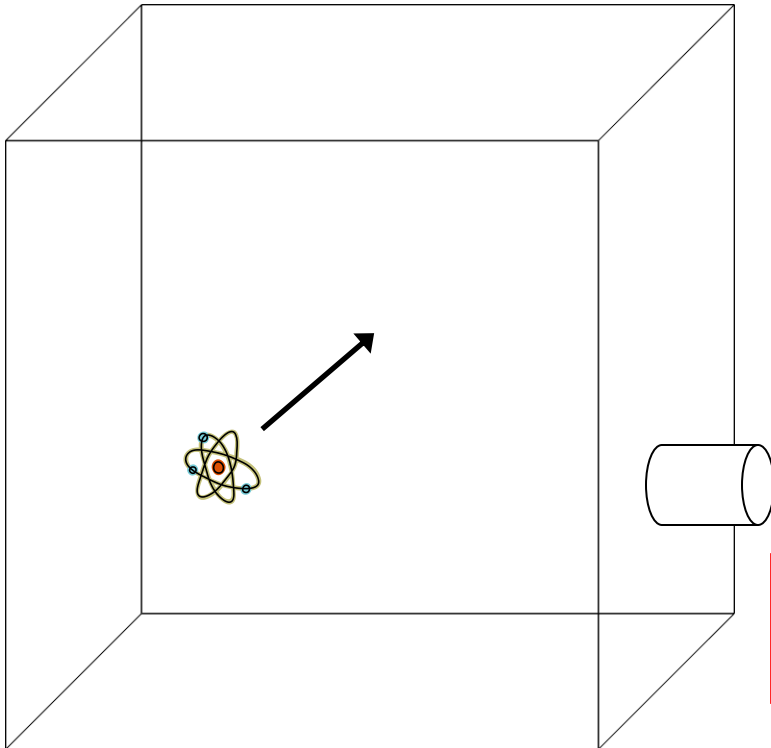
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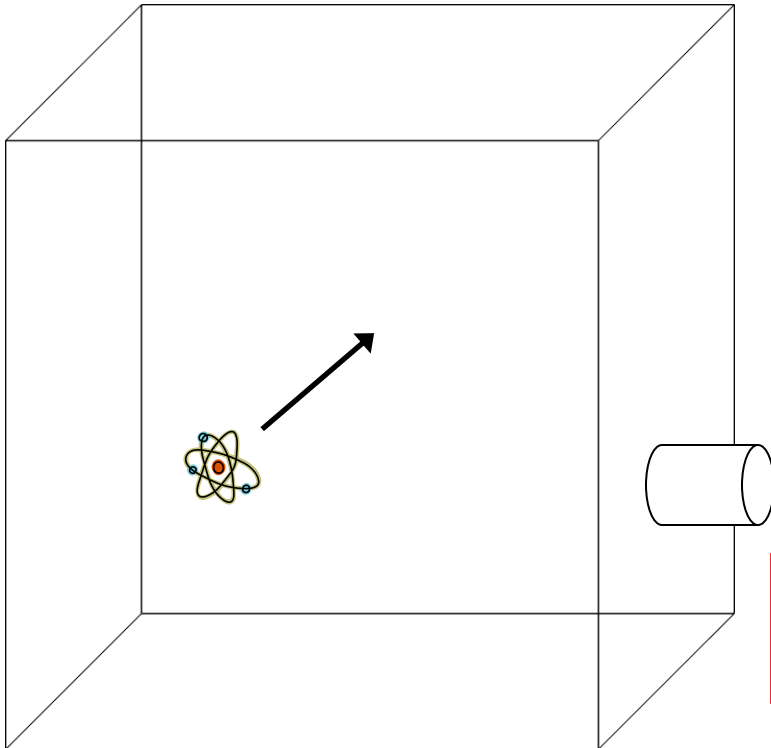
Solution: simulate many individual molecular trajectories and look at statistics ($\langle n_{\text{exit}} \rangle$, σ_n)

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... fairly simple and quick on a computer.

Definition

The Monte Carlo method is any numerical method in which the solution is obtained by ***averaging over many probabilistic simulation instances***.

Example: Numerical Integration

The Monte Carlo method is frequently used to evaluate difficult integrals (in many dimensions):

$$\text{"calculus" average: } \langle f(x) \rangle_{[a,b]; \text{calculus}} = \frac{1}{b-a} \int_a^b f(x) dx$$

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x_i = probabilistic variable

i.e. choose x_i 's randomly on $[a,b]$ with a uniform probability distribution.

Theorem

If $f(x)$ is well behaved on $[a,b]$ (i.e. does not diverge), then in the limit of $N \rightarrow \infty$, the following is true (in the probabilistic sense)

$$\langle f(x) \rangle_{[a,b]; \text{calculus}} = \langle f(x) \rangle_{[a,b]; \text{statistical}, N} \pm \frac{\sigma_{f(x), N}}{\sqrt{N}}$$

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where
$$\sigma_{f(x), N}^2 = \frac{1}{N-1} \left(\sum_{i=1}^N f(x_i)^2 - N \langle f(x) \rangle_{[a,b], N}^2 \right) = \frac{1}{N-1} \sum_{i=1}^N (f(x_i) - \langle f(x) \rangle)^2$$

= standard deviation of simulations

Advantages

- Monte Carlo simulations are generally easy to formulate and set-up.
- Monte Carlo simulations are generally faster than other numerical methods, especially for problems in a large number of dimensions.