

PHY604 Lecture 19

November 9, 2023

Review: MFT of the Poisson equation

- So, for the transformed Poisson equation:

$$\left[\exp\left(\frac{-i2\pi m}{N}\right) + \exp\left(\frac{i2\pi m}{N}\right) + \exp\left(\frac{-i2\pi n}{N}\right) + \exp\left(\frac{i2\pi n}{N}\right) - 4 \right] F_{m,n} = -\frac{h^2}{\epsilon_0} R_{m,n}$$

- Solving for the **F** matrix:

$$F_{m,n} = -\frac{h^2}{2\epsilon_0(\cos(2\pi m/N) + \cos(2\pi n/N) - 2)} R_{m,n}$$

- To get the potential, we just need to take the inverse FFT:

$$\Phi_{j,k} = \frac{1}{N^2} \sum_{m=0}^{N-1} \sum_{n=0}^{N-1} F_{m,n} \exp\left(\frac{i2\pi jm}{N}\right) \exp\left(\frac{i2\pi kn}{N}\right)$$

Review: von Neumann stability analysis

- In discretized form:

$$a_j^n = A^n e^{ikjh}$$

- Advancing the solution by one step:

$$a_j^{n+1} = A^{n+1} e^{ikjh} = \xi A^n e^{ikjh}$$

- ξ is the **amplification factor**

- **von Neumann stability analysis:** Insert this trial solution into the numerical scheme and solve for amplification factor given h and τ

- Unstable if $|\xi| > 1$

Review: Stability of FTCS for advection equation

- FTCS scheme:
$$a_i^{n+1} = a_i^n - \frac{c\tau}{2h}(a_{i+1}^n - a_{i-1}^n)$$
- Insert trial solutions: $a_j^n = A^n e^{ikjh}$ $a_j^{n+1} = \xi A^n e^{ikjh}$
$$\xi A^n e^{ikjh} = A^n e^{ikjh} \left[1 - i \frac{c\tau}{h} \sin(kh) \right]$$
- Therefore:
$$|\xi| = \left| 1 - i \frac{c\tau}{h} \sin(kh) \right| = \sqrt{1 + \left(\frac{c\tau}{h} \right)^2 \sin(kh)^2}$$
- So, the solution in general grows with each timestep, and therefore unstable

Review: Stability of the Lax scheme

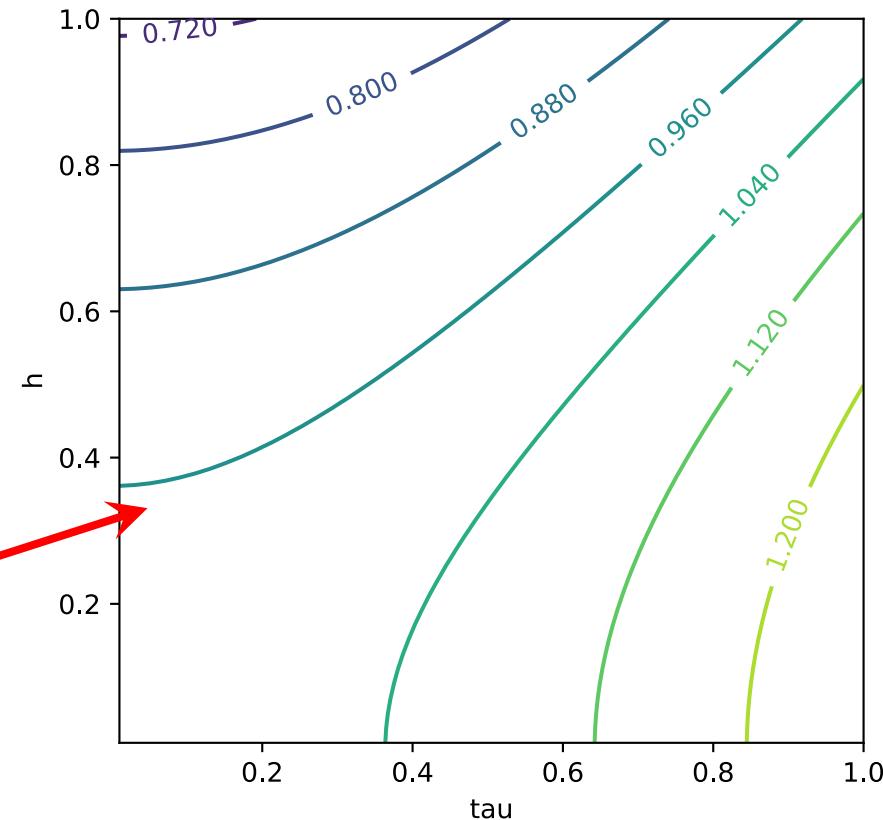
- So, we have: $|\xi| = \sqrt{\cos^2(kh) + \left(\frac{c\tau}{h}\right)^2 \sin^2(kh)}$

- Example: take $k=\pi/4$, $c=1$:

- In general: $\left|\frac{c\tau}{h}\right| \leq 1$

- Same as the Courant-Friedrichs-Lowy stability criterion

τ must be less than or equal to h



Review: Stability condition on eigenvalues

$$\mathbf{T}^{n+1} = \sum_{k=0}^{N-1} c_k (\lambda_k)^n \mathbf{v}_k$$

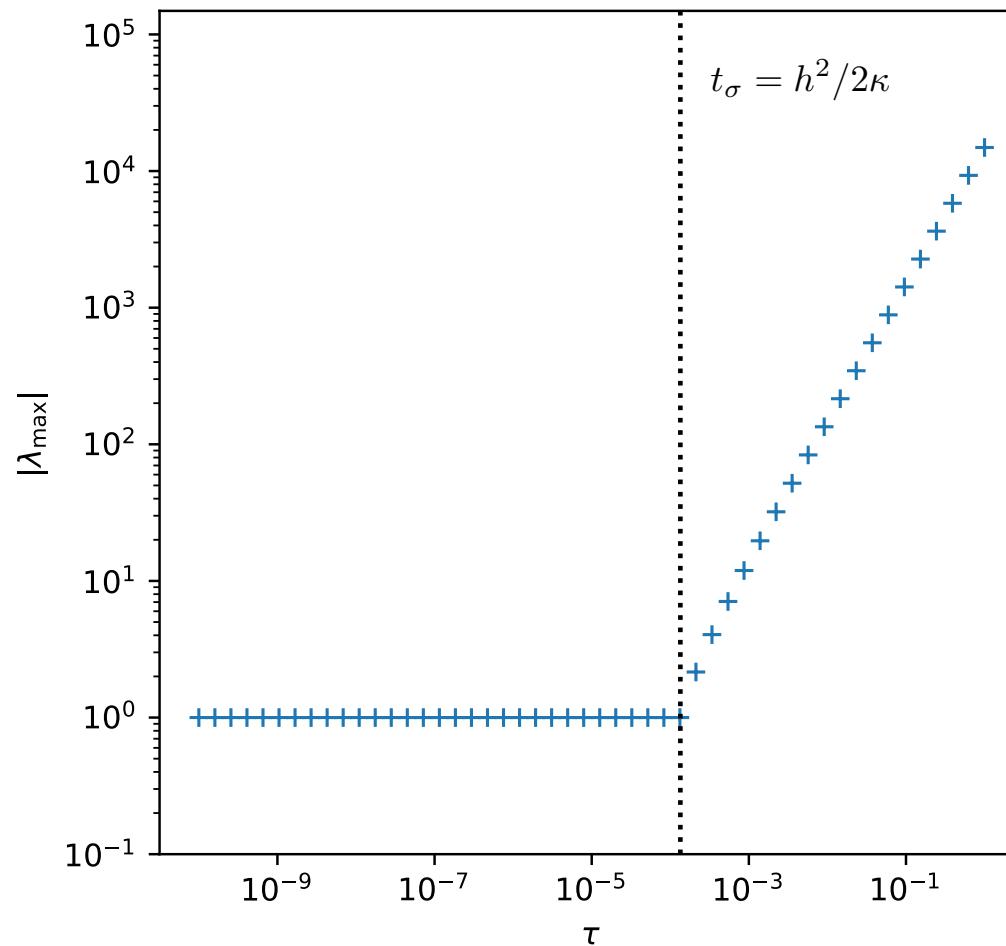
- We see that we will have divergence if we have any eigenvalues that are: $|\lambda_k| > 1$
- **Spectral radius of \mathbf{A} :** Magnitude of the largest eigenvalue

$$\rho(\mathbf{A}) = |\lambda_{\max}|$$

- Scheme is matrix stable if the spectral radius is less than or equal to unity

Review: Stability of FTCS for diffusion equation with timestep

- 61 spatial grid points with unit length, $\kappa = 1$:



Today's lecture:

Implicit schemes for PDEs and Random numbers

- Implicit schemes for PDEs: The Schrodinger equation
- Introduction to stochastic methods: Random numbers

Example for implicit schemes: Schrödinger equation

$$i\hbar \frac{\partial}{\partial t} \psi(x, t) = -\frac{\hbar^2}{2m} \frac{\partial^2}{\partial x^2} \psi(x, t) + V(x) \psi(x, t)$$

- Or:

$$i\hbar \frac{\partial \psi}{\partial t} = \mathcal{H}\psi$$

- Formal solution:

$$\psi(x, t) = \exp \left[-\frac{i}{\hbar} \mathcal{H}t \right] \psi(x, 0)$$

Discretizing the Schrödinger equation

- FTCS for the Schrödinger equation:

$$i\hbar \frac{\psi_j^{n+1} - \psi_j^n}{\tau} = -\frac{\hbar^2}{2m} \frac{\psi_{j+1}^n + \psi_{j-1}^n - 2\psi_j^n}{h^2} + V_j \psi_j^n$$

- Since the Hamiltonian is a linear operator:

$$i\hbar \frac{\psi_j^{n+1} - \psi_j^n}{\tau} = \sum_{k=0}^{N-1} H_{jk} \psi_k^n$$

- Where:

$$H_{jk} = -\frac{\hbar^2}{2m} \frac{\delta_{j+1,k} + \delta_{j-1,k} - 2\delta_{jk}}{h^2} + V_j \delta_{jk}$$

FTCS steps for Schrödinger equation

- Final FTCS scheme in matrix notation:

$$\Psi^{n+1} = \left(\mathbf{I} - \frac{i\tau}{\hbar} \mathbf{H} \right) \Psi^n$$

- First term in Taylor expansion of the formal solution for one time step:

$$\psi(x, t) = \exp \left[-\frac{i}{\hbar} \mathcal{H}t \right] \psi(x, 0)$$

Implicit schemes for the Schrödinger equation

- We have seen that the FTCS is numerically unstable for time steps that are too large
- Alternative approach: Apply the Hamiltonian to the future value of ψ

$$i\hbar \frac{\psi_j^{n+1} - \psi_j^n}{\tau} = \sum_{k=0}^{N-1} H_{jk} \psi_k^{n+1}$$

- Or:
- Solving for Ψ^{n+1} :

$$\Psi^{n+1} = \left(\mathbf{I} + \frac{i\tau}{\hbar} \mathbf{H} \right)^{-1} \Psi^n$$

Implicit FTCS scheme

- Implicit FTCS:

$$\Psi^{n+1} = \left(\mathbf{I} + \frac{i\tau}{\hbar} \mathbf{H} \right)^{-1} \Psi^n$$

- Compare with explicit FTCS:

$$\Psi^{n+1} = \left(\mathbf{I} - \frac{i\tau}{\hbar} \mathbf{H} \right) \Psi^n$$

- Equivalent as τ goes to 0 since for small ε :

$$\frac{1}{1 + \epsilon} \rightarrow (1 - \epsilon)$$

- Con: Implicit method requires evaluation of matrix inverse, which can be costly
- Pro: Unconditionally stable!

More accurate approximations: Crank-Nicholson

- As we saw before, numerically stable does not mean accurate
- More accurate scheme: **Crank-Nicholson**
 - Average of implicit and explicit FTCS:

$$i\hbar \frac{\psi_j^{n+1} - \psi_j^n}{\tau} = \frac{1}{2} \sum_{k=0}^{N-1} H_{jk} (\psi_k^n + \psi_k^{n+1})$$

- In matrix form:

$$\Psi^{n+1} = \Psi^n - \frac{i\tau}{2\hbar} \mathbf{H} (\Psi^n + \Psi^{n+1})$$

- Isolating the $n+1$ term:

$$\Psi^{n+1} = \left(\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right)^{-1} \left(\mathbf{I} - \frac{i\tau}{2\hbar} \mathbf{H} \right) \Psi^n$$

Properties of Crank-Nicolson

$$\Psi^{n+1} = \left(\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right)^{-1} \left(\mathbf{I} - \frac{i\tau}{2\hbar} \mathbf{H} \right) \Psi^n$$

- Unconditionally stable
- Centered in both space and time
- “Páde” approximation for exponential is
 - See (https://en.wikipedia.org/wiki/Pad%C3%A9_approximant)

$$e^{-z} \simeq \frac{1 - z/2}{1 + z/2}$$

- CN can be interpreted as Páde for the formal solution
- Preserves the unitarity of e^{-z}

Example: Numerical solution of the Schrödinger equation

- Initial conditions: Gaussian wave packet

- Localized around x_0
- Width of σ_0
- Average momentum of: $p_0 = \hbar k_0$

$$\psi(x, t = 0) = \frac{1}{\sqrt{\sigma_0 \sqrt{\pi}}} \exp(ik_0 x) \exp\left[-\frac{(x - x_0)^2}{2\sigma_0^2}\right]$$

- Which is normalized so that:

$$\int_{-\infty}^{\infty} |\psi|^2 dx = 1$$

- Also, has the special property that uncertainty produce $\Delta x \Delta p$ is minimized ($\hbar/2$)

Propagation of wave packet in free space

- Wavefunction evolves like:

$$x \rightarrow x - \frac{p_0 t}{2m}, \quad \sigma_0^2 \rightarrow \alpha^2 \equiv \sigma_0^2 + \frac{i\hbar t}{m}$$

- So we have:

$$\psi(x, t) = \frac{1}{\sqrt{\sigma_0 \sqrt{\pi}}} \frac{\sigma_0}{\alpha} \exp \left[ik_0 \left(x - \frac{p_0 t}{2m} \right) \right] \exp \left[-\frac{(x - x_0 - \frac{p_0 t}{2m})^2}{2\alpha^2} \right]$$

- And for the probability density:

Remains a Gaussian in
time

→ $P(x, t) = |\psi(x, t)|^2 = \frac{\sigma_0}{|\alpha|^2 \sqrt{\pi}} \exp \left[- \left(\frac{\sigma_0}{|\alpha|} \right)^4 \frac{(x - x_0 - \frac{p_0 t}{m})^2}{\sigma_0^2} \right]$

Propagation of wave packet in free space

- By symmetry, max of Gaussian equals its expectation value:

$$\langle x \rangle = \int_{-\infty}^{\infty} x P(x, t) dx$$

- In time, it moves as: $\langle x \rangle = x_0 + \frac{p_0 t}{m}$
- And the wave packet spreads as:

$$\sigma(t) = \sigma_0 \sqrt{1 + \frac{\hbar^2 t^2}{m^2 \sigma_0^4}}$$

Why does the rough spatial discretization give errors?

- The reason is a poor representation of the initial conditions
- Rough discretization suppresses the higher wave number modes
 - Difficult to represent those modes on a coarse grid
- Because of this suppression, the discretized version has a lower momentum than $\psi(x,t)$

Can we avoid the taking the inverse of the matrix?

- As usual, we can trade taking the matrix inverse for solving a linear system of equations:

$$\begin{aligned}\Psi^{n+1} &= \left(\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right)^{-1} \left(\mathbf{I} - \frac{i\tau}{2\hbar} \mathbf{H} \right) \Psi^n \\ &= \left(\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right)^{-1} \left[2\mathbf{I} - \left(\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right) \right] \Psi^n \\ &= \left[2 \left(\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right)^{-1} - \mathbf{I} \right] \Psi^n\end{aligned}$$

- Or:

$$\Psi^{n+1} = \mathbf{Q}^{-1} \Psi^n - \Psi^n, \quad \mathbf{Q} = \frac{1}{2} \left[\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right]$$

Crank-Nicolson for tridiagonal matrices

$$\Psi^{n+1} = \mathbf{Q}^{-1}\Psi^n - \Psi^n, \quad \mathbf{Q} = \frac{1}{2} \left[\mathbf{I} + \frac{i\tau}{2\hbar} \mathbf{H} \right]$$

- Now we can solve for the next timestep by solving the linear system:

$$\mathbf{Q}\chi = \Psi^n$$

- And then:

$$\Psi^{n+1} = \chi - \Psi^n$$

- Recall that for banded matrices, solving linear systems via, e.g., Gaussian elimination, is particularly efficient

Some comments in implicit schemes

- Recall that the killer app of implicit methods was that they are unconditionally stable
- Major downside is that for higher-dimensional problems, matrices become very large and difficult to manipulate
 - Can use approaches to separately perform implicit steps in different dimensions

Today's lecture:

Implicit schemes for PDEs and Random numbers

- Implicit schemes for PDEs: The Schrodinger equation
- Introduction to stochastic methods: Random numbers

Monte Carlo and stochastic methods

- Randomness is an important part of physics
 - E.g., radioactive decay, Brownian motion
 - In standard interpretations of quantum mechanics, microscopic phenomena are random
- Random sampling can be a useful tool for integration
 - Whole family of techniques based on this idea



Wikipedia

How can we model randomness on the computer

- In order to implement stochastic methods, we need random (or pseudorandom) numbers
- What do we need from a random number generator? (according to Pang)
 - Long “period” before sequences of numbers are repeated
 - Small correlation between numbers generated in sequence
 - Very fast, so we can get many random numbers to accumulate statistics
- Typical random number generators return a number in $[0,1)$
 - Should uniformly fill that space
 - Seeds can be used to allow for reproducibility (from one run to the next)

Example of a simple random number generator

- Simplest generator made using the linear congruent scheme

- Random numbers are generated in sequence from the linear relation:

$$x_{i+1} = (ax_i + b) \mod c$$

- a , b , and c are “magic numbers” which determine the quality of the generator

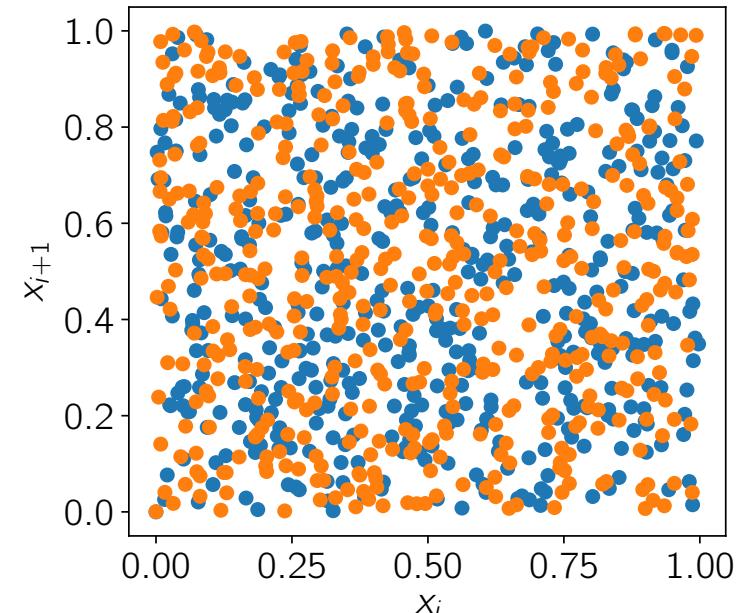
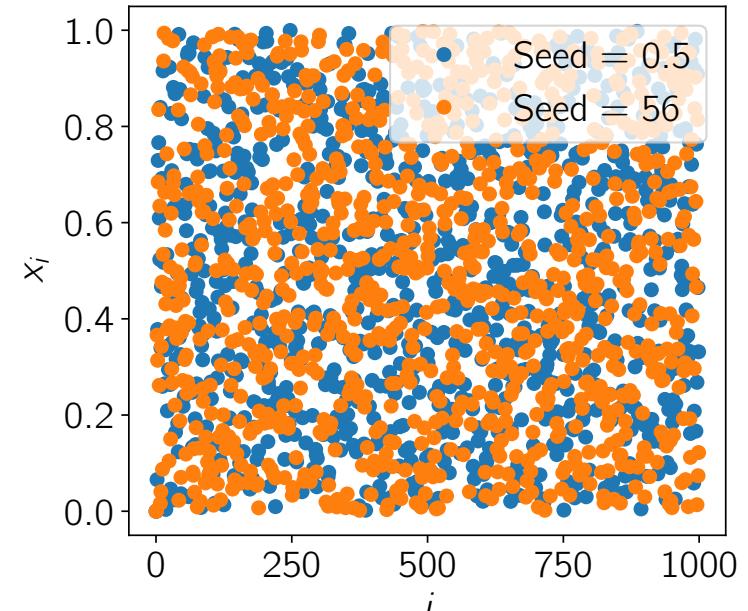
- Typical choices: $a = 7^5$, $b = 0$, $c = 2^{31} - 1$
 - x_0 is the seed, allows for reproducibility

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Random versus pseudo random

- The numbers generated on the previous slide are actually **pseudorandom**
 - Wikipedia: Appears to be statistically random despite having been produced by a completely deterministic and repeatable process
 - Usually, good enough for most applications (except if you are doing cryptography)
 - Good for testing code, since you get the same values every time
 - Can randomize the seed (e.g., clock time)
- “True” random numbers can be generated by physically random processes
 - Some noise or random process of the computer hardware (e.g., clock time)
 - Thermal noise from a resistor
 - Quantum shot noise
 - Atmospheric noise: <https://www.random.org/>
 - Lava lamps: <https://patents.google.com/patent/US5732138>

Random numbers from the RAND corporation

- If you want your random numbers in analog format, you can download a book of them:

https://www.rand.org/pubs/monograph_reports/MR1418.html

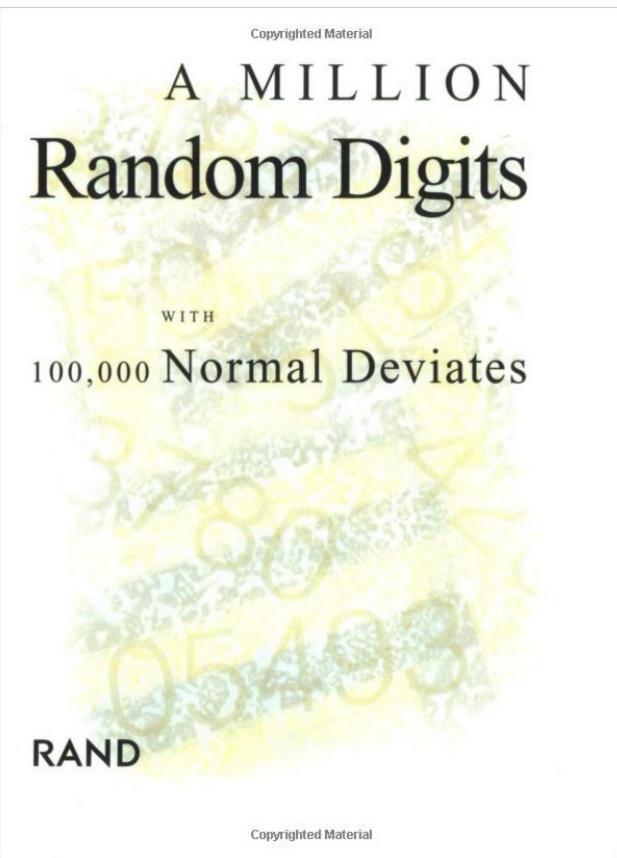


TABLE OF RANDOM DIGITS

3

00100	03991	10461	93716	16894	66083	24653	84609	58232	88618	19161
00101	38555	95554	32886	59780	06355	60860	29735	47762	71299	23853
00102	17546	73704	92052	46215	55121	29281	59076	07936	27954	58909
00103	32643	52861	95819	06831	00911	98936	76355	93779	80863	00514
00104	69572	68777	39510	35905	14060	40619	29549	69616	33564	60780
00105	24122	66591	27699	06494	14845	46672	61958	77100	90899	75754
00106	61196	30231	92962	61773	41839	55382	17267	70943	78038	70267
00107	30532	21704	10274	12202	39685	23309	10061	68829	55986	66485
00108	03788	97599	75867	20717	74416	53166	35208	33374	87539	08823
00109	48228	63379	85783	47619	53152	67433	35663	52972	16818	60311
00110	60365	94653	35075	33949	42614	29297	01918	28316	98953	73231
00111	83799	42402	56623	34442	34994	41374	70071	14736	09958	18065
00112	32960	07405	36409	83232	99385	41600	11133	07586	15917	06253
00113	19322	53845	57620	52606	66497	68646	78138	66559	19640	99413
00114	11220	94747	07399	37408	48509	23929	27482	45476	85244	35159

The 100,000 “normal deviates” cited in the title of this volume constitute a subset of random numbers whose occurrence can be plotted on a bell-shaped curve. RAND legend has it that this seemingly self-contradictory mathematical expression caused the New York Public Library to misslelve the volume in the Psychology section.

Best bet is to use previous implementations
for random number generators

- Correlations between random samples can be difficult to detect and cause errors in computations
- See: <https://docs.python.org/3/library/random.html> or <https://numpy.org/doc/stable/reference/random/index.html> for details on how python does it

Radioactive decay

(see Newman Sec. 10.1)

- One of the quintessential random processes in physics
- Parent atoms decay with characteristic half-life τ
- We will consider ^{208}TI , which decays to ^{208}Pb with $\tau = 183.18$ sec.

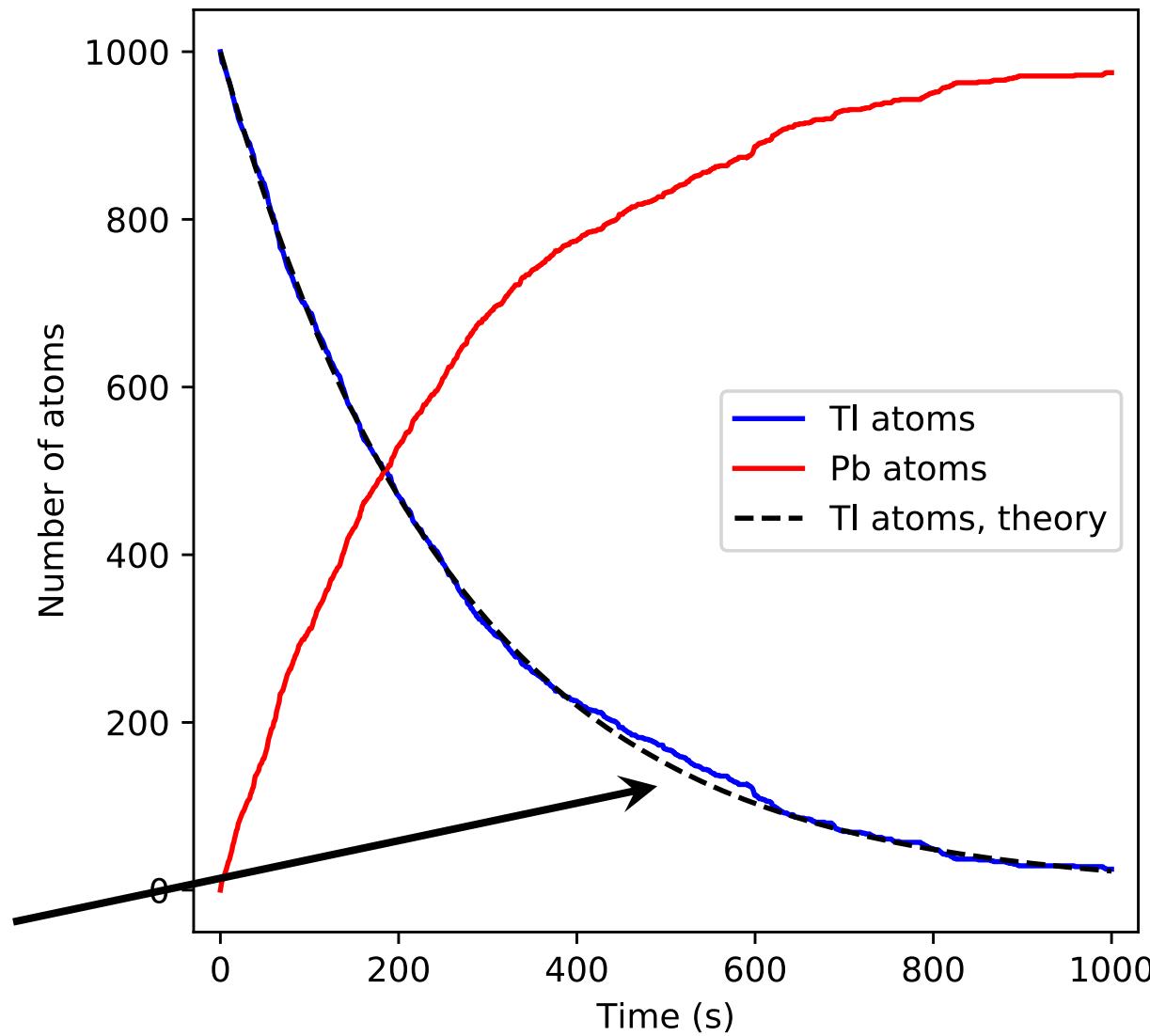
- Number of parent atoms falls off exponentially:

$$N(t) = N(0)2^{-t/\tau}$$

- Probability that a particular atom has decayed in a time interval t :

$$p(t) = 1 - 2^{-t/\tau}$$

Radioactive decay



$$N(t) = N(0)2^{-t/\tau}$$

Nonuniform distributions

- We can also select random numbers from a distribution that is not constant over the range
 - I.e., all numbers are not selected with equal probability

- Consider the radioactive decay example:

- Probability of decay in time interval dt is:

$$p(t) = 1 - 2^{-dt/\tau} = 1 - \exp\left(-\frac{dt}{\tau} \ln 2\right) \simeq \frac{\ln 2}{\tau} dt$$

- What is the probability to decay in time window $t + dt$?

- Needs to survive without decay until t (probability $2^{-t/\tau}$)
 - Then must decay in dt
 - Total probability is:

$$P(t)dt = 2^{-t/\tau} \frac{\ln 2}{\tau} dt$$

Nonuniform distribution for decay example

- Nonuniform probability distribution:

$$P(t)dt = 2^{-t/\tau} \frac{\ln 2}{\tau} dt$$

- Decay times t are distributed in proportion to $2^{-t/\tau}$
- We could calculate the decay of N atoms by drawing N random samples from this distribution
 - More efficient than previous method
 - Need to generate nonuniform distribution of random numbers
- Can generate nonuniform random numbers from a uniform distribution

Transformation method for changing distributions

- We have a source of random numbers z drawn from distribution $q(z)$
 - Probability of generating a number between z and $z+dz$ is $q(z)dz$
- Now we choose a function $x = x(z)$ whose distribution $p(x)$ is the one we want
- We know that:
$$p(x)dx = q(z)dz$$
- If our random numbers are drawn from a uniform distribution $[0,1]$,
 $q(z)=1$ from 0 to 1, zero elsewhere
- Then:
$$\int_{-\infty}^{x(z)} p(x')dx' = \int_0^z dz' = z$$
- We need to do the integral on the left and then solve for $x(z)$
 - Not always possible

Transformation method to exponential distribution

- Say we want to generate random real numbers that are > 0 with the distribution:

$$p(x) = \mu e^{-\mu x}$$

- μ is for normalization

- Then:

$$\mu \int_{-\infty}^{x(z)} e^{-\mu x'} dx' = 1 - e^{-\mu x} = z$$

- So:

$$x = -\frac{1}{\mu} \ln(1 - z)$$

Nonuniform distribution for decay example

- We can write the probability distribution for the decay example as

$$P(t)dt = 2^{-t/\tau} \frac{\ln 2}{\tau} dt = e^{-t \ln 2 / \tau} \frac{\ln 2}{\tau}$$

- So:

$$x = -\frac{\tau}{\ln 2} \ln(1 - z)$$

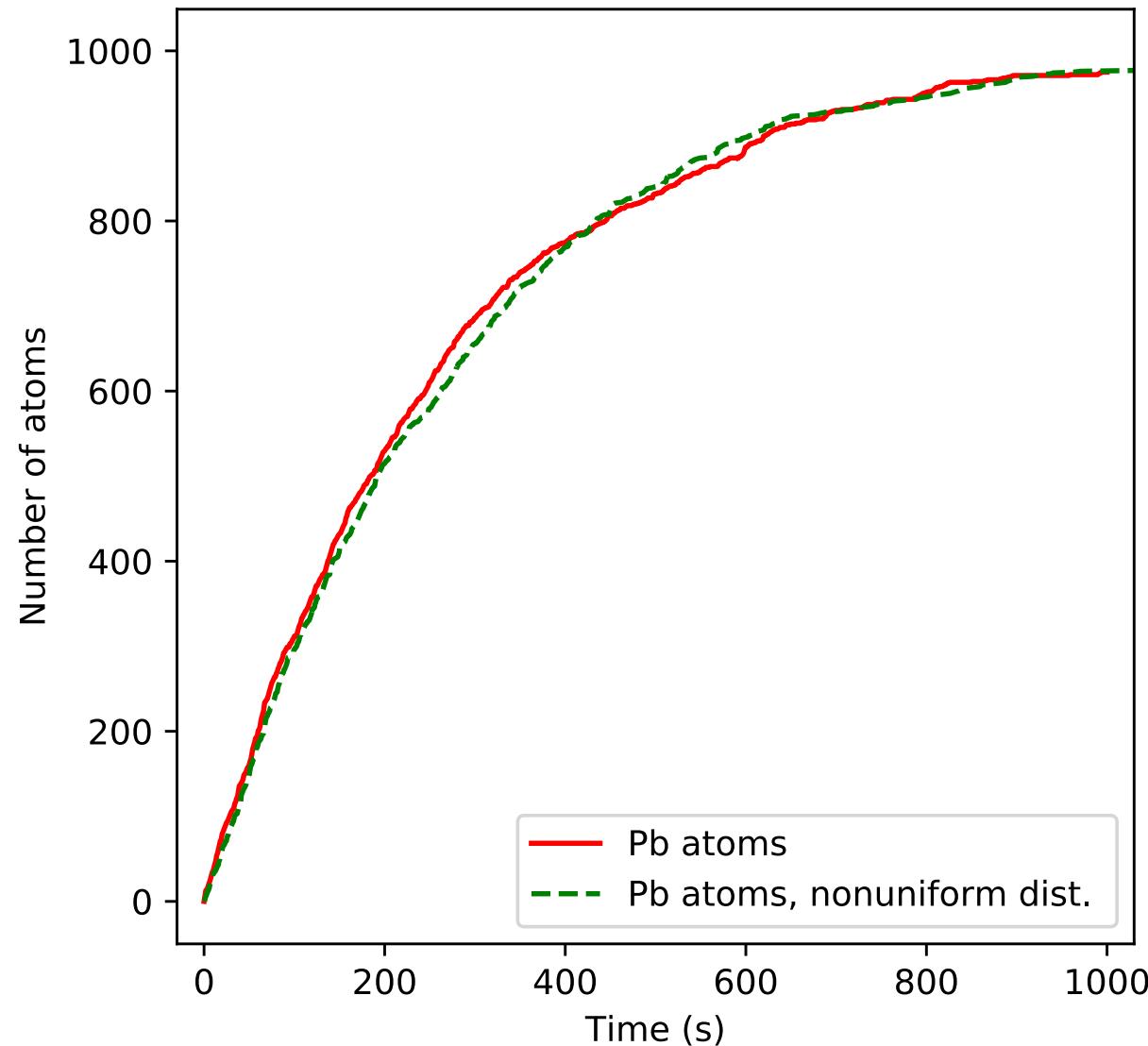
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Gaussian random numbers

- In many cases we would like to draw numbers from a Gaussian (i.e., normal) distribution:

$$p(x) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{x^2}{2\sigma^2}\right)$$

- Let's try the transformation method:

$$\frac{1}{\sqrt{2\pi\sigma^2}} \int_{-\infty}^x \exp\left(-\frac{x^2}{2\sigma^2}\right) dx = z$$

- The solution to this integral and the resulting equation is complicated

Gaussian random numbers

- Trick: consider two random numbers x and y , both drawn from Gaussian distribution with the same standard deviation
- Probability that point with position (x,y) falls in some element $dxdy$ on xy plane is:

$$\begin{aligned} p(x)dx \times (y)dy &= \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{x^2}{2\sigma^2}\right) dx \times \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{y^2}{2\sigma^2}\right) dy \\ &= \frac{1}{2\pi\sigma^2} \exp\left(-\frac{x^2 + y^2}{2\sigma^2}\right) dxdy \end{aligned}$$

- Now convert to polar coordinates:

$$p(r, \theta)drd\theta = \frac{1}{2\pi\sigma^2} \exp\left(-\frac{r^2}{2\sigma^2}\right) r dr d\theta = \frac{r}{\sigma^2} \exp\left(-\frac{r^2}{2\sigma^2}\right) dr \frac{d\theta}{2\pi}$$

2D transformation method

$$p(r)dr \times p(\theta)d\theta = \frac{r}{\sigma^2} \exp\left(-\frac{r^2}{2\sigma^2}\right) dr \frac{d\theta}{2\pi}$$

- The point in polar coordinates will have the same distribution as the original point in cartesian (x,y)
 - Solving in polar coordinates and transforming back to Cartesian gives us two random points from a Gaussian distribution
- θ part is just a uniform distribution: $p(\theta) = 1/2\pi$
- Radial part can be treated with transformation method:

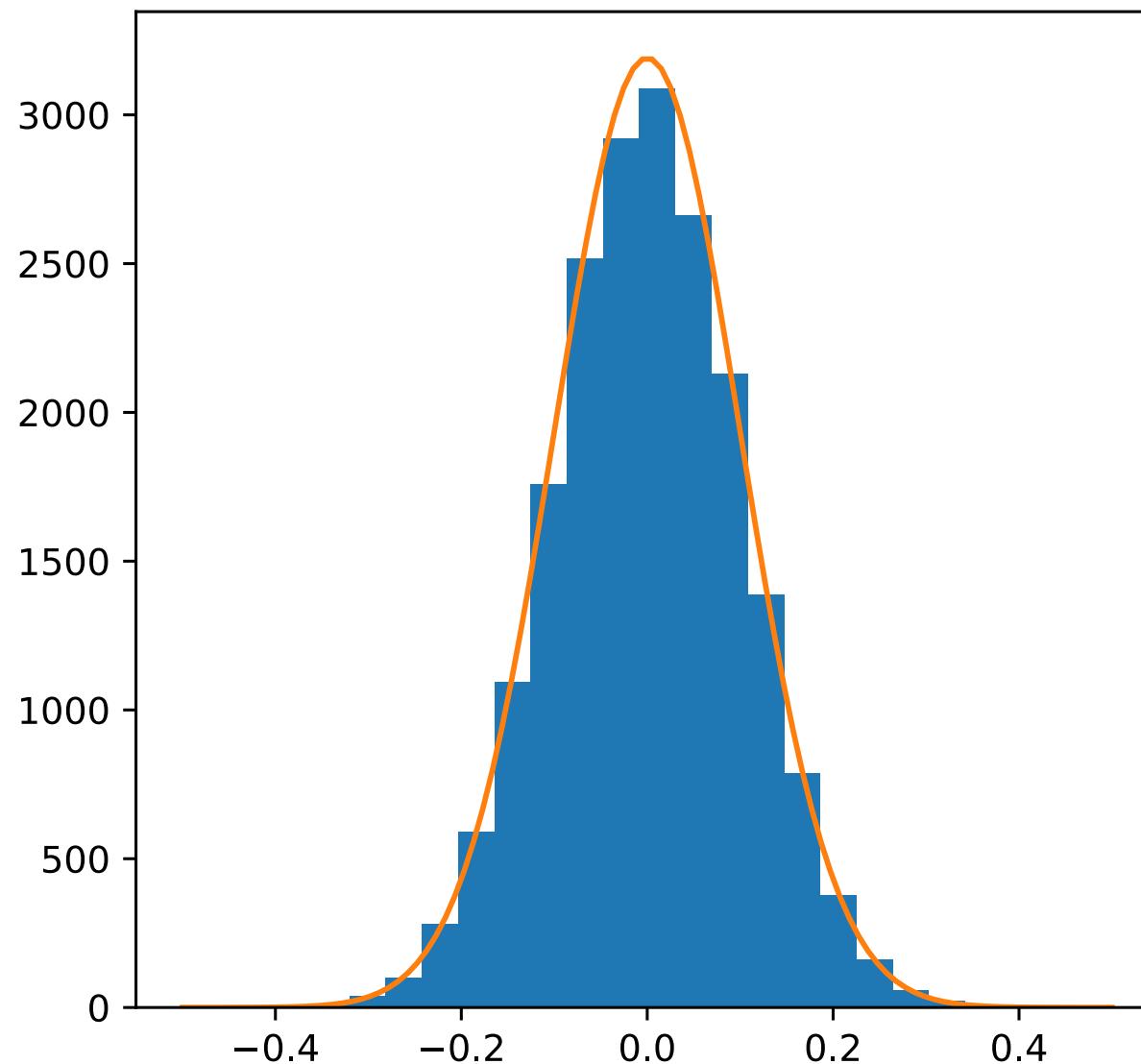
$$\frac{1}{\sigma^2} \int_0^r \exp\left(-\frac{r'^2}{2\sigma^2}\right) r' dr' = 1 - \exp\left(-\frac{r^2}{2\sigma^2}\right) = z$$

- So:

$$r = \sqrt{-2\sigma^2 \ln(1 - z)}$$

- And random numbers are: $x = r \cos \theta, \quad y = r \sin \theta$

Random numbers from Gaussian distribution



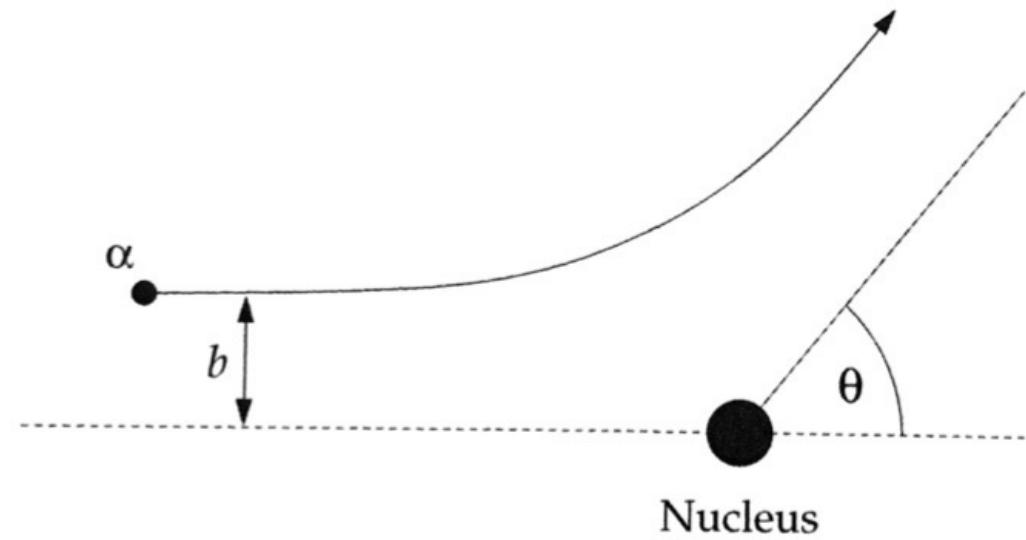
Example: Rutherford scattering

- α particles (helium nuclei) scatter when they pass close to an atom with angle:

$$\tan\left(\frac{\theta}{2}\right) = \frac{Ze^2}{2\pi\epsilon_0 Eb}$$

- E is the kinetic energy of the α particle, b is the impact parameter
- Consider Gaussian beam of particles with $\sigma=a_0/100$ and $E=7.7\text{MeV}$ fired at a gold atom
- How many “bounce back” (scattering angle > 90 degrees)?

$$b \leq \frac{Ze^2}{2\pi\epsilon_0 E}$$



After class tasks

- Homework 4 is posted, due Nov. 14, 2023
- Readings
 - Newman Sec. 10.1
 - Pang Sec. 2.5
 - Garcia Sec. 11.2