

Ising model (cont'd)
PHYS 250 (Autumn 2024) – Lecture 7

David Miller

Department of Physics and the Enrico Fermi Institute
University of Chicago

October 22, 2024

Outline

Outline of the Ising model discussion

We've already discussed a lot, and I want to remind you of those topics, their progression, and where we're going from here to make sure that we're all on the same page.

- **Lecture 5: the model itself**

- The general concept of the model: lattice of spins
- Basis in thermodynamics and quantum mechanics
- Importance of simulations methods

- **Lecture 6: computational approaches**

- History of computational simulation methods: Monte Carlo
- The Metropolis Monte Carlo algorithm and its assumptions (deeply related to thermodynamics)

- **Lecture 7 (Today): analytical and computational evaluation**

- Analytical Ising model and the key numerical results
- Concept of equilibrium and how we can define it
- Observables in the Ising model and their calculation

- **“Lecture 8” (Tomorrow): Hands-on lab+lecture**

- Hands-on session in CSIL 2 for developing our Ising model simulation

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Recall our discussion of statistical mechanics

We discussed that the probability distribution of an observable such as the mean energy, $\langle E \rangle$, of a system in a particular microstate μ is given by

$$\langle E \rangle = \frac{\sum_{\mu} E_{\mu} e^{-\beta E_{\mu}}}{\sum_{\mu} e^{-\beta E_{\mu}}} = \frac{\sum_{\mu} e^{-\beta E_{\mu}}}{Z} \quad (1)$$

where Z is the **partition function** of the system ($\beta = (k_B T)^{-1}$).

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Similarly, the variance is given by:

$$\text{Var}(E) = \langle E^2 \rangle - \langle E \rangle^2 = - \frac{\partial \langle E \rangle}{\partial \beta} = \frac{\partial^2 \ln Z}{\partial \beta^2} \quad (4)$$

Ising's analytic solution for a 1D spin lattice (I)

For a single dimension, and just $N = 2$ spins (the smallest possible Ising chain model) we can use the analytic form on the previous slide to write down Z exactly, and thus the mean energy (recall $E = -J \sum_{\langle ij \rangle} s_i s_j$) and its variance.

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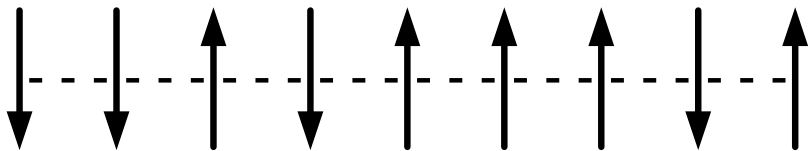
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$$= (2 \cosh \beta J)^N \quad (15)$$

$$(\text{Since } Z_1 = \sum_{s_1=\pm 1} 1 = 2)$$

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Note: at constant T , Helmholtz free energy (F) is minimized at equilibrium.

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The last expression for magnetization M , or magnetization per unit particle, $m = M/N$ shows a very important property

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It has already been pointed out by Ising himself⁴ that a linear chain of spins is not ferromagnetic. This can easily be verified by calculating the total magnetization with the help of (5) and (8):

$$M = mN \sinh C / (\sinh^2 C + e^{-4K})^{\frac{1}{2}}, \quad (10a)$$

Kramers & Wannier, 1941

Properties of the 2D Ising model

Lars Onsager, *Phys. Rev.* 65, 117 – Published 1 February 1944

Onsager showed in 1943 that one can solve the 2D Ising model analytically.

Solution for $\ln Z/2 = -\beta F - \ln 2$ (and allowing different J in each dimension)

$$\log(\lambda/2) = \frac{1}{2}\pi^{-2} \int_0^\pi \int_0^\pi \log(\cosh 2H \cosh 2H' - \sinh 2H \cos \omega - \sinh 2H' \cos \omega') d\omega d\omega'. \quad (108)$$

This leads to a relationship for the magnetization which **does allow for** $M \neq 0$ at $T \neq 0$, namely $T = T_c$:

$$\sinh\left(\frac{2J_1}{k_B T_c}\right) \sinh\left(\frac{2J_2}{k_B T_c}\right) = 1 \quad (24)$$

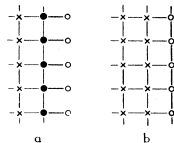


FIG. 3. Two-step extension of a two-dimensional crystal. (a) A new tier of atoms \circ is added (V_1); their configuration depends on that of atoms \bullet in previous marginal position. (b) Interaction energy between marginal atoms \circ is introduced (V_2), which modifies the distribution of configurations in this tier of atoms.

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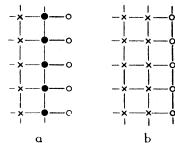


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Which corresponds to

$$\frac{k_B T_c}{J} = \frac{2}{\ln(1 + \sqrt{2})} \quad (25)$$

$$= 2.2692 \quad (26)$$

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Equilibrium

In order to measure properties of the system using a simulation, we have to run our simulation for a suitably long period of time until it has come to equilibrium at the temperature (or whatever state-related property) we are interested in. This period is called the equilibration time τ_{eq} .

Only then can we obtain a reliable measurement of the quantity we are interested in. And then, in order to measure it's average values we must wait another suitably long period of time and average it, to evaluate the estimator of that quantity.

So many questions!

Questions you should ask yourself

- What exactly do we mean by “allowing the system to come to equilibrium”?
- How long is a “suitably long time” for it to happen?
- How do we go about measuring our quantity of interest once that happens?
- How long do we have to average over to get a result of a desired degree of accuracy?

These are very general questions which we need to consider every time we do a Monte Carlo simulation.

Discussion of equilibrium

“**Equilibrium**” means that the average probability of finding our system in any particular state μ is proportional to the Boltzmann weight $e^{-\beta E_\mu}$ of that state.

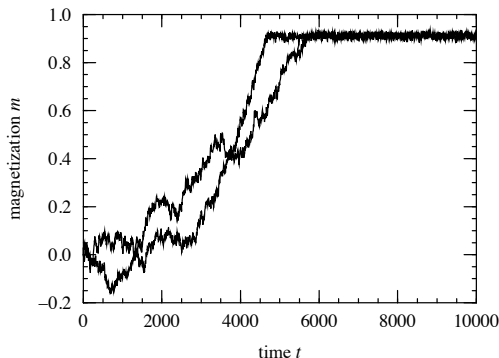
If we start our system off in a state such as with $T = 0$ (e.g. all spins aligned), and we want to perform a simulation at some finite non-zero temperature, $T \neq 0$, it will take a little time before we reach equilibrium.

A system at equilibrium spends the overwhelming majority of its time in a small subset of states in which its internal energy and other properties take a narrow range of values. This is what is “mapped out” by the temperature. To determine if the system has reached equilibrium, we could just “look at it” but this isn’t very reliable.

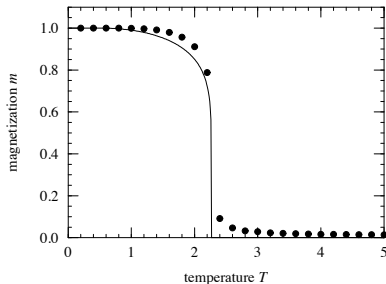
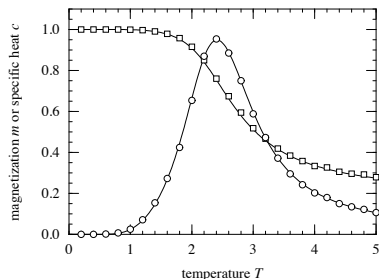
Equilibrium in simulations: magnetization

A standard approach is to plot a graph of some quantity of interest, like the magnetization per spin $m = M/N$ of the system or the energy of the system E , as a function of “time” (iterations or steps) from the start of the simulation.

Up until equilibrium, the energy and the magnetization are changing, but after equilibrium is reached, they just fluctuate around a steady average value. This is what we should measure.



Simulation vs. calculations for the 2D Ising model



These results come from **two simulations (markers)**, a 5×5 lattice (left) and a 100×100 lattice (right). These are compared to **analytic calculations (solid lines)**. Many iterations are run for each simulation, and averaged after equilibrium is reached to measure the magnetization and the energy. $m = M/N$ and $c = C/N$ are obtained from Eqs **??,??**, and $J = k_B = 1$.

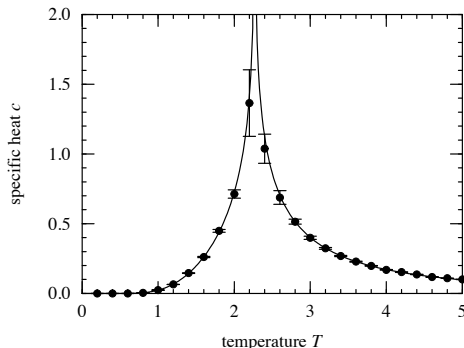
The features at the Curie temperature T_c are an indication of a **phase transition**, for which the magnetization is the **order parameter**

Specific heat: a measure of the variance of the energy

As we wrote down in Eq. ??:

$$C = \frac{\partial \langle E \rangle}{\partial T} = -k_B \beta^2 \frac{\partial^2 \ln Z}{\partial \beta^2} = \frac{N(\beta J)^2}{\cosh^2 \beta J} = k_B \beta^2 \text{Var}(E) \quad (27)$$

we can calculate this in terms of the energies of the system.



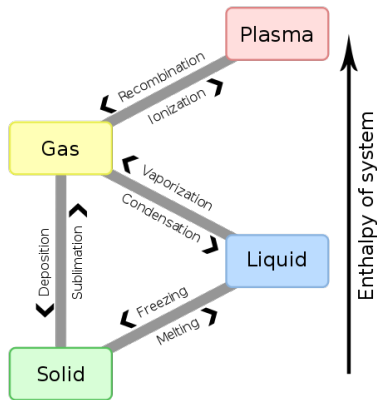
The divergence at T_c is another indication of a **phase transition**

Phase transitions

We have all heard of, talked about, played with, and even tasted phase transitions since shortly after we were born:

Basic idea of a phase transition (Wikipedia)

A phase of a thermodynamic system and the states of matter have uniform physical properties. During a phase transition of a given medium, certain properties of the medium change, often discontinuously, as a result of the change of some external condition, such as temperature, pressure, or others.

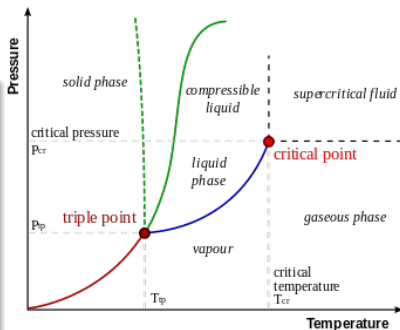


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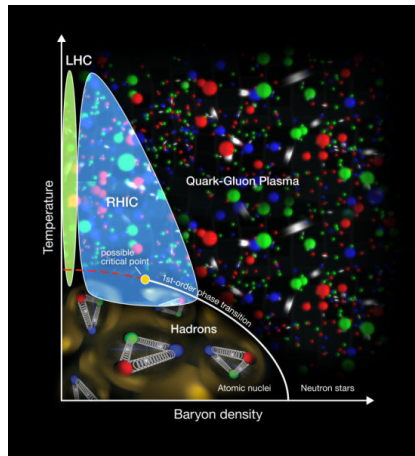
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More exotic phase transitions

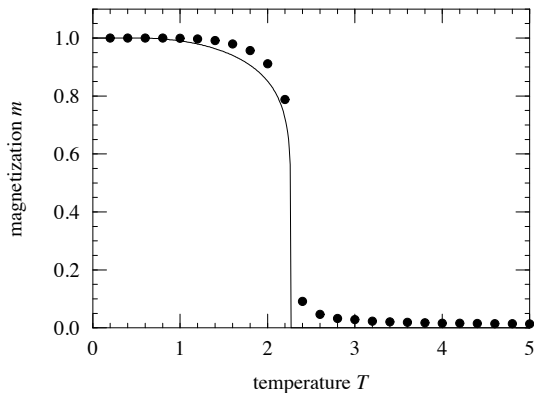
What you have likely **not heard of or talked about**, certainly have not “**played with**”, and **definitely have not tasted phase transitions in the early universe** (or at the LHC):



In each case, the description of the system has gross characteristics that depend on certain parameters, and very often, some notion of **temperature**.

Let's look at that plot of magnetization again

Pretty clear that there is a phase transition occurring!



Outline

Some hints for HW3

We'll work on this together tomorrow (remember, CSIL at 2:30pm!) but here are some hints in the mean time

Approach

Make sure you have short snippets that allow you to test each function you write and make sure it's doing what you think it should be doing.

Spin sums

```
# Can be used instead of my example code for  
# less readable, but much faster sums  
np.roll(lattice)
```

Energy variance

```
# Useful for specific heat  
# Modularize your manipulations of the lattice!  
np.var(energy_density/(T**2))
```

Equilibrium

Test the “time” required to reach equilibrium **before** doing many simulations with many iterations each.