Studies of Phase Transitions with the Ising Model and the Metropolis Algorithm

Andreas Fagerheim*

Department of Physics, University of Oslo, Norway

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Abstract

This article set forth to examine phase shifts and the critical temperature of materials. This will be done by using the well studied Ising model with an implemented algorithm called Metropolis. Using this model we will examine how the thermal energy affect a magnetic materials. At a given critical temperature, this model exhbits a phase transition from a magnetic phase (a system with a finite magnetic moment) to a phase with zero magnetization. All programs developed and used are available in the github link in the footnote.

I. Introduction

In the help of understanding phase transition we make use of fluctuations. In physics we make use of statistical models and we will here look closer at the Ising model. The behaviour of said fluctuations can be closer examined using the Isng model and will serve as a tool to better understand the underlying particle interactions. The system of interest can said to be binary where the objects in the lattice can only take on two different values. These could be -1 and 1 or other values.

We can solve the system analytically for certain expectation values in one and two dimensions and it gives a qualitatively good understanding of several types of phase transitions. The system we will look closer to can be assumed to be a ferromagnetic ordering, viz $J > 0$. We will use periodic boundary conditions and the Metropolis algorithm only. The article first examines the case of 2 x 2 lattice. Then a study of bigger surfaces $(20x20, 40x40, 80x80, 100x100)$ will closer examined.

II. Theory

i. Ising Model

The total energy of the Ising Model can in the simplest form be expressed as

$$E = -J \sum_{<kl>}^{N} s_k s_l - B \sum_{k}^{N} s_k$$

with $s_k = \pm 1$. The quantity N represents the total number of spins and J is a coupling constant expressing the strength of the interaction between neighboring spins. The symbol $< kl >$ indicates that we sum over nearest neighbors only. B is an external magnetic field interacting with the magnetic moment set up by the spins [Hjorth-Jensen, 2015]. This article will look closer at the case of witch there is no interacting external magnetic field and the energy can the be expressed as:

$$E = -J \sum_{<kl>}^{N} s_k s_l \tag{1}$$

The expectation value, the mean energy, can be calculated given a probability distribution P_i as

$$\langle E \rangle = \sum_{i=1}^{M} E_i P_i(\beta) \tag{2}$$

hvor the probability distribution is given by the Boltzmann distribution

$$P_i(\beta) = \frac{e^{-\beta E_i}}{Z}$$

where $\beta = 1/k_b T$ is the inverse temperature, k_b is the Boltzmann constant, E_i is is the energy of a microstate i. The partition function is Z and is given for

*https://github.com/AndreasFagerheim/Project-4

the canonical ensemble as a sum over all microstates, M;

$$Z = \sum_{i=1}^{M} e^{-\beta E_i}.$$

Further the the expectation values for magnetic moment $|M|$

$$\langle |M| \rangle = \sum_{i=1}^{M} M_i P_i(\beta). \tag{3}$$

The variance of E and $|M|$ is given receptively as

$$\sigma_E^2 = \langle E^2 \rangle - \langle E \rangle^2$$

$$\sigma_M^2 = \langle M^2 \rangle - \langle M \rangle^2$$

This lets us express the specific heat capacity constant volume as

$$C_v = \frac{1}{k_b T^2}(\langle E^2 \rangle - \langle E \rangle^2) \tag{4}$$

$$\chi = \frac{1}{k_b T}(\langle M^2 \rangle - \langle M \rangle^2) \tag{5}$$

ii. Analytic solutions for the Ising Model of 2x2 lattice

Looking closer at a 2x2 lattice to find analytical solutions which will serve as a bechmark for the method implemented later with T = 1. **Table 1** shows the different energies and magnetization states. Using the known states we can find analytical solutions:

$$Z = \sum_{i=1}^{M} e^{-\beta E_i} = 2e^{\beta 8J} + 2e^{-8\beta J} + 12e^0.$$

with $\beta = 1/kT$ and $T = 1$ with unit kT/J we get:

$$Z = 2e^8 + 2e^{-8} + 12e^0.$$

This gives us the energy

$$\langle E \rangle = \sum_{i=1}^{M} E_i P_i(\beta) = \frac{1}{Z} \sum_{i=1}^{M} E_i e^{-\beta E_i} \tag{6}$$

$$\langle E \rangle = \frac{1}{Z}(-8e^8 + 8e^{-8} + 8e^{-8} - 8e^8) = -7.9839$$

$$\frac{\langle E \rangle}{N} = -1.99598$$

Table 1: *The different energy states of the 2x2 Ising Model*

No. of spins	Degeneracy	Energy	Magnetization
4	1	-8J	4
3	4	0	2
2	4	0	0
2	2	8J	0
1	4	0	-2
0	1	-8J	-4

Showing the different energy and magnetization for the two-dimensional Ising model with periodic boundary conditions.

hvor $N = 4$. In the same way the magnetization turns out to be

$$\langle |M| \rangle = \frac{1}{Z} \sum_{i=1}^{M} M_i e^{E_i} = 3.9946 \tag{7}$$

$$\frac{\langle |M| \rangle}{N} = 0.9986$$

The heat capacity then becomes:

$$\langle E^2 \rangle = \frac{1}{Z} \sum_{i=1}^{M} E_i^2 e^{E_i} = \frac{128e^8 + 128e^{-8}}{5973.917} \tag{8}$$

$$\frac{C_V}{N} = \frac{\langle E^2 \rangle - \langle E \rangle^2}{4} = 0.12832 = 0.03208$$

The magnetic susceptibility can in the same way be calculated to $\frac{\chi}{N} = 0.00401$.

iii. Studies of phase transitions.

For many materials there exist a critical temperature, T_C, at which the material changes its characteristics and undergoes a phase transition. In this case the Ising model can give us a tool for examination of the characteristics. The mean magnetization can for the Ising model be expressed: As an example, for the Ising class of models, the mean magnetization is given by

$$\langle M(T) \rangle \sim (T - T_C)^{\beta},$$

where $\beta = 1/8$ is a so-called critical exponent. For the heat capacity and the magnetic susceptibility it can in the same way be expressed as

$$C_V(T) \sim |T_C - T|^{\alpha},$$

$$\chi(T) \sim |T_C - T|^{\gamma}, \tag{9}$$

with $\alpha = 0$ and $\gamma = 7/4$.

A second-order phase transition is characterized by a correlation length. The behavior at finite lattice can be related the results for an infinite lattice. Scaling of the critical temperature take the form

$$T_C(L) - T_C(L = \infty) = aL^{-1/\nu}, \tag{10}$$

with a a constant and ν defined in Eq. (11). We set $T = T_C$ and obtain a mean magnetisation

$$\langle \mathcal{M}(T) \rangle \sim (T - T_C)^{\beta} \rightarrow L^{-\beta/\nu}, \tag{11}$$

a heat capacity

$$C_V(T) \sim |T_C - T|^{-\gamma} \rightarrow L^{\alpha/\nu}, \tag{12}$$

and susceptibility

$$\chi(T) \sim |T_C - T|^{-\alpha} \rightarrow L^{\gamma/\nu}. \tag{13}$$

III. Metropolis algorithm

The Metropolis algorithm builds on the principle of natures desire to reach equlibrum towards lower energy levels. Simply explained it starts of by picking a random place in the spin matrix. It then try to flip the spin if the calculated change in energy is making the total energy lower in the system. Flipping the spin is then thougt to make the system closer to equilibrium. Here we make use off what can be said to be Monte Carlo method to repeatedly call on the Metropolis algorithm to make allowed changes to the spin. In C++ this will look like the figure below in some way.

```
for(int i=1;i<MC_cycles;i++){

    Metropolis(spin_matrix, L, energy, magnetization, w, idum, confi_accept, neighbour);
    // update expectation values
    values[0] += energy;    values[1] += energy*energy;
    values[2] += magnetization;    values[3] += magnetization*magnetization;
    values[4] += fabs(magnetization);
```

Figure 1: *Out take of code implemented where the Metropolis function get called each cycle andthe the new expectation values are saved.*

IV. Results

From **Table 2** we see that the numerical value of E compared to the analytical value for the 2 x 2 lattice, and some other values. **Figure 2** shows how the Energy develops with the increasing number of Monte Carlo cycles. Here we see that after 40000 cycles the values fluctuates less and is close to the analytical solution. The same can be said about the mean magnetization shown in the lower plot in **Figure 2**.

MC cycles	E/N	C_V/N	$\lvert M\rvert/N$	χ/N
Analytisk	-1.9959	0.03208	0.9986	0.00401
20000	-1.9951	0.03910	0.9983	0.005288
40000	-1.9966	0.02755	0.9988	0.003669
60000	-1.9966	0.02715	0.9988	0.003345
80000	-1.9964	0.02875	0.9987	0.003707
100000	-1.9966	0.02699	0.9989	0.003455

Table 2: *Comparison of the analytical values and the numerical values with increasing number of Monte Carlo cycles.*

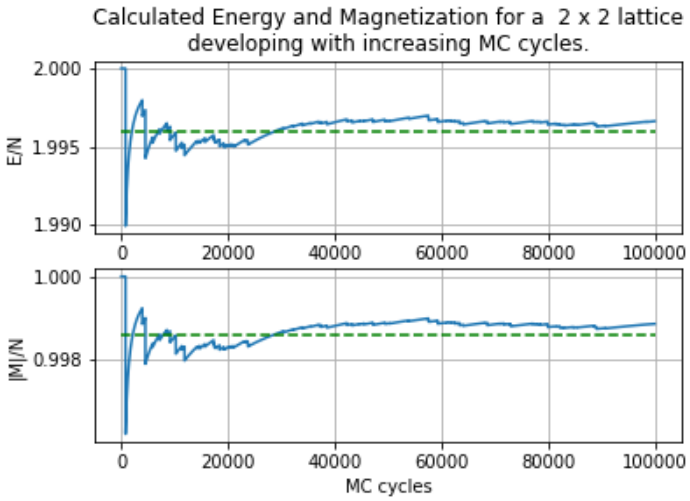

Figure 2: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 2 x 2 lattice and with the initial temperature set to $k_bT/J = 1$. The green dotted lines is the analytical value to the normalized Energy and magnetization.*

For a lattice of dimensions 20 X 20 the development of energy and mean magnetisation as a function of Monte Carlo cycles is plotted in **FIgure 3,4,5,6**.

It is evident that the system reaches a state of equilibrium already around 20 000 cycles.

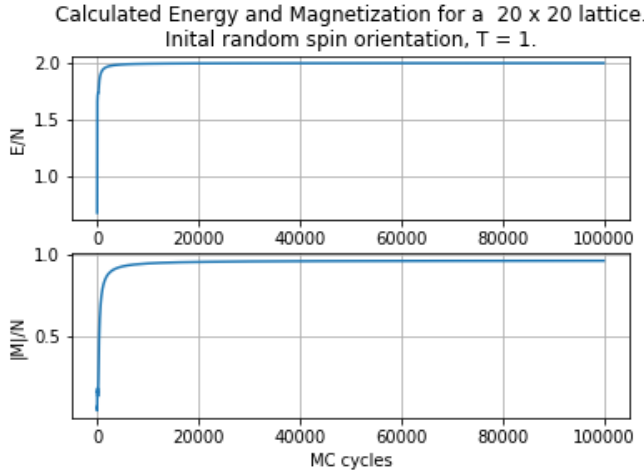

Figure 3: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 1$ and an initial random spin orientation.*

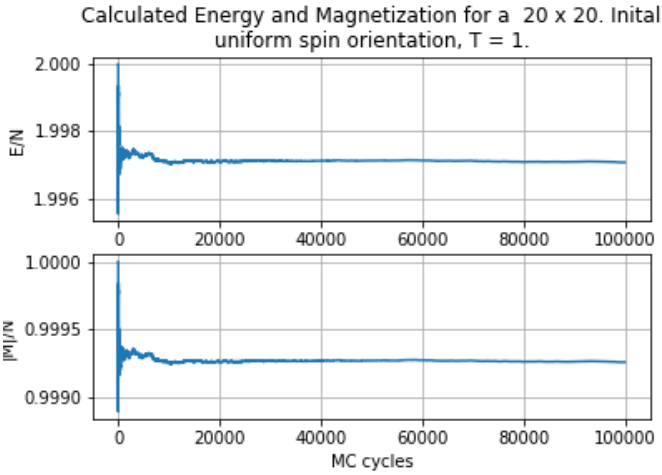

Figure 4: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 1$ and an initial uniform spin orientation.*

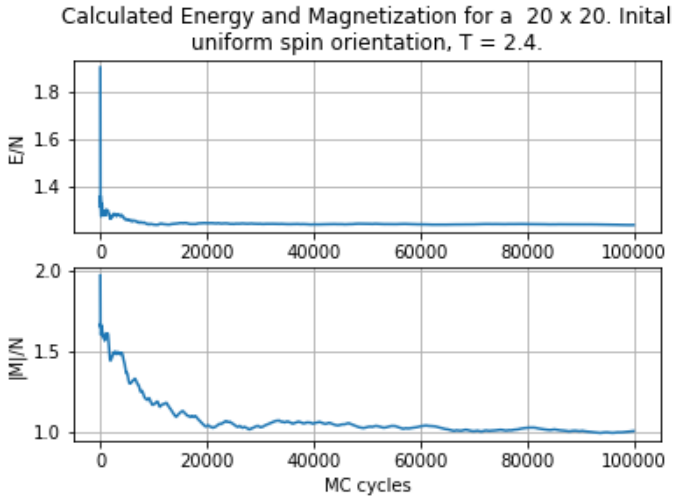

Figure 5: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 2.4$ and an initial random spin orientation.*

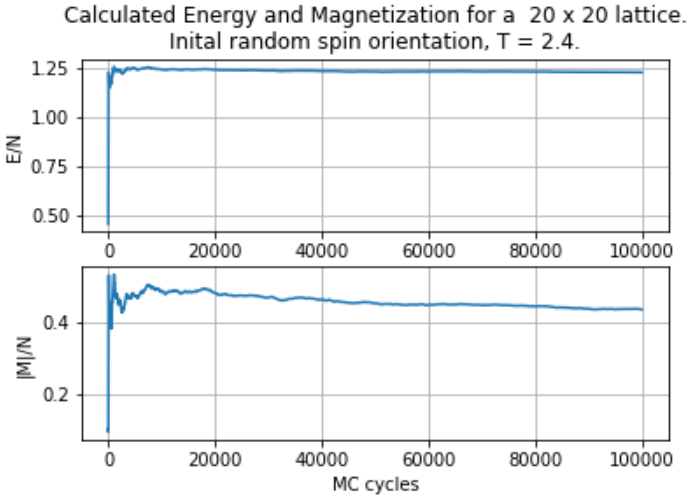

Figure 6: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 2.4$ and an uniform random spin orientation.*

Looking at **Figure 7** and **8** should give and indication on how the number of accepted states should develop with the increasing number of cycles. Number of accepted configurations seems to increase with the temperature.There should although be said that the numbers in **Fig. 8** don't makes to much sense i

the magnitude nesseceraly, but the number of states should behave in the same way as the number of cycles increase. The initial spin orientation does not affect the number of possible configurations after some time has passed. Temperature does seem to increase the number of acceptable states per cycle, but i don't know if by as much as the results indicates. This aligns with the principal that with higher temperature, the probability for accepting a state increases.

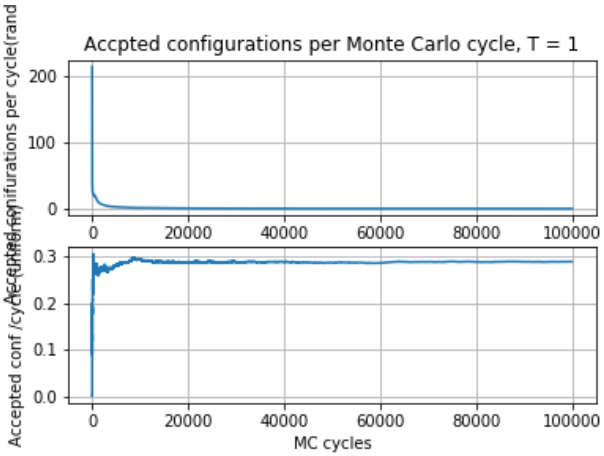

Figure 7: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 2.4$ and an uniform random spin orientation.*

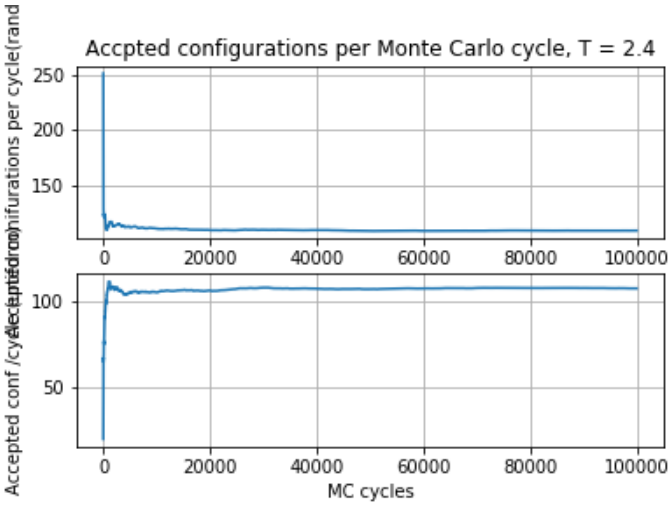

Figure 8: *Energy and magnetization as a function of Monte Carlo cycles. The plot is for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 2.4$ and an uniform random spin orientation.*

i. Probability

Looking at **Figure 9** and **Figure 10**. For the case $T = 2.4$ the distribution favours towards the right, which is towards the origin. This coincides with the Boltzmann distribution.

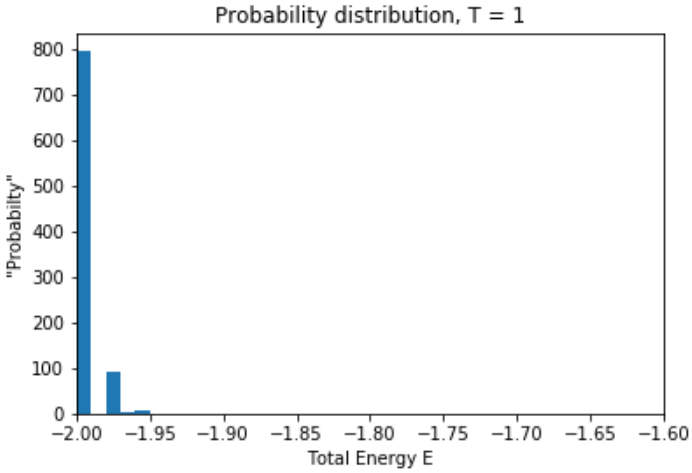

Figure 9: *Probability distribution for a 20 x 20 lattice and with the initial temperature set to $k_bT/J = 1$.*

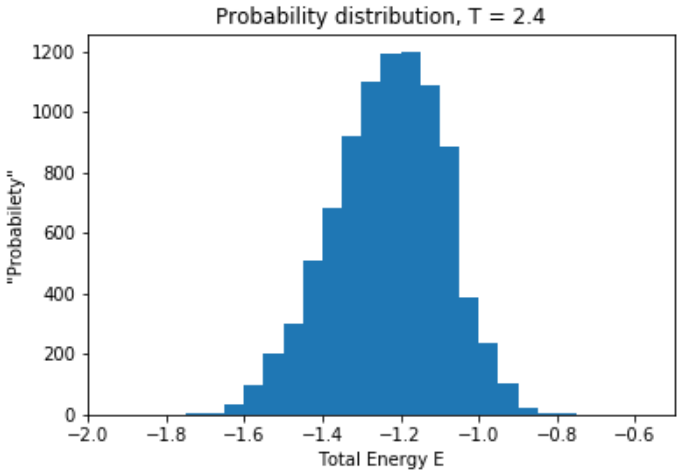

Figure 10: *Probability distribution for a 20 x 20 lattice and with the initial temperature set to* $k_bT/J = 2.4$.

V. Conclusion

The Ising model has given us a taste of some of the characteristics the system undergoes. The system clearly reaches an equlibrum and the temperature can bee seen to impact the system. With a higher initial temperature of the system there will be a higher number of possible microstates. The analytical solutions of the 2x2 lattice serves as a way of validating the method and insight to how it works. This study lack a closer look in to the study of the critical temperature T_C due to problems in parallelizing the code for. This would have made it possible to closer examine how the magnetic susceptibility, ,χ and the heat capacity, C_V, for the material changes with the temperature. These characteristics drastically changes when the material reaches its critical temperature.

References

[Hjorth-Jensen, 2015] Hjort-Jensen, M. (2015). Computational Physics.

[Hjorth-Jensen] Hjort-Jensen, M. https://github.com/CompPhysics/ComputationalPhysics