

If there any hope to find divergence it will be in the thermodynamic limit. Fluctuations with respect to the order parameter. Remind that:

We can see that:

$$\chi_T = \frac{1}{k_B T} \left[\langle M^2 \rangle - \langle M \rangle^2 \right] \quad (1)$$

Idea: replace the M with an integral.

$$M(\mathbb{C}) = \int d^3 \vec{r} m(\vec{r}) \quad (2)$$

$$\chi_T k_B T = \int d\vec{r} d\vec{r}' \left[\langle m(\vec{r}) m(\vec{r}') \rangle - \langle m(\vec{r}) \rangle \langle m(\vec{r}') \rangle \right] \quad (3)$$

Translational symmetry implies:

$$\langle m(\vec{r}) \rangle = m \quad (4)$$

and the 2-points correlation function

$$\langle m(\vec{r}) m(\vec{r}') \rangle \equiv G(\vec{r} - \vec{r}') \quad (5)$$

I could see:

$$G_c(\vec{r} - \vec{r}') = \left\langle \underbrace{(m(\vec{r}) - \langle m(\vec{r}') \rangle)}_{\substack{\delta m(\vec{r}') \\ \text{fluctuations}}} (m(\vec{r}') - \langle m(\vec{r}') \rangle) \right\rangle \quad (6)$$

$$\chi_T k_B T = \int d\vec{r} d\vec{r}' \left[G(\vec{r} - \vec{r}') - \langle m(\vec{r}) \rangle \langle m(\vec{r}') \rangle \right] \quad (7)$$

it means that \vec{r}' is not no more a parameter. You can put it tends to zero. If in the last equation we put $\vec{r}' = 0$, we obtain:

$$\Rightarrow \chi_T k_B T = V(\Omega) \int d\vec{r} G_c(\vec{r} - \vec{r}_0) \quad (8)$$

that is the *fluctuations and dissipation relation*. In general the correlation function $G_c(\vec{r})$ what gives? Usually it behaves in an exponential way. In principle we expect that:

$$G_c(\vec{r}) \simeq e^{-|\vec{r}|/\xi} \quad (9)$$

where ξ is called the *correlation length*. It is related to the correlation function. In general is finite but if you approach T_c diverges. In fact, at the critical point this correlation will expand in the whole space and reaches the size of all the system, it goes to infinity. When ξ will diverge, there is not anymore the exponential and the integrall cannot be kept finite. From the thermodynamic divergence you obtain that the quantity ξ goes to infinity. You are looking a critical phenomena.

Therefore, as $T \rightarrow T_c^\pm$ we have

$$\xi \stackrel{t \rightarrow 0^\pm}{\sim} t^{-\nu_\pm} \quad (10)$$

Supposing $T = T_c$:

$$G_c(r) \sim \frac{1}{r^{d-2-\eta}} \quad (11)$$

Lecture 5.

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If we have a system with finite size L , the temperature is not $T = T_c$, because we are in a finite system.

Consider a finite system L , for sure we have $\xi \leq L$. We have also a small scale as a , the length of the Bravais lattice. For instance, suppose $L = 1 \text{ cm}$ and $T \equiv \frac{T-T_c}{T_c}$, we can see that in general

$$\xi = \xi_0 t^{-\nu} \rightarrow t = \left(\frac{\xi}{\xi_0} \right)^{-1/\nu} \simeq \left(\frac{L}{\xi_0} \right)^{-1/\nu} = \left(\frac{10^{-2} m}{10^{-9} m} \right)^{-1/\nu} \sim (10^7)^{-2} \sim 10^{-14} \quad (12)$$

this explain why at the critical point we do not see this effect.

We can find the critical point by doing Montecarlo simulation. Supposing a Montecarlo simulation of a Ising model, for which there is no an analitic solution and compute the energy

1 fig

Try to estrapolate for example the position of the peak as N increases. If you start to see this behaviour it shown that something is happening.

Nevertheless, there are two approaches: I want to study the system by looking for all the details: example a protein, that interact with other proteins. I can look at all the electrons (or atoms). Even if you think at the simple protein that exist there are a lot of degree of freedom. For doing a simulation, if you are interested in long time behaviour and in large scale behaviour, in that case details are not important. What it is important are symmetry, range of interaction, therefore you can forget about all the details. You introduce the effective potentials as wander-walls or Lenard Jones potential. You study collective effect. This is the second approach.

0.1 Ising Model D=1

This model has a critical point that is $T_c = 0$, it means that there is no phase transition. Therefore, Lenz (?) thought that the ising model was unuselles for studing the paramagnetic system. This is not true, because at dimension 2, the critical temperature is $T_c \neq 0$. This is the result of Onsager, Remember

2fig

Think a square lattice, in which in each point you have a spin $s_i = \pm 1$. Given $N(\Omega)$ sites in your lattice. The number of configuration is $\#\{\mathcal{C}\} = 2^N$. Remember that a term as $\sum_i S_i$ is not correlated. We need an interaction. The simpler one is putting the interaction

$$\sum_i S_i H_i + \frac{1}{2} \sum_{i \neq j} J_{ij} S_i S_j \quad (13)$$

$$J_{ij} = \frac{J}{|\vec{r}_i - \vec{r}_j|} \quad (14)$$