Anti-ferromagnetic Ising Model in Hierarchical Networks

Xiang Cheng and Stefan Boettcher Department of Physics, Emory University, Atlanta, GA 30322, USA

The Ising antiferromagnet is a convenient model of glassy dynamics. It can introduce geometric frustrations and may give rise to a spin glass phase and glassy relaxation at low temperatures. We apply the antiferromagnetic Ising model to 3 hierarchical networks which share features of both small world networks and regular lattices. Their recursive and fixed structures make them suitable for exact renormalization group analysis as well as numerical simulations. We first explore the dynamical behaviors using simulated annealing and discover an extremely slow relaxation at low temperatures. Then we employ the Wang-Landau algorithm to investigate the energy landscape and the corresponding equilibrium behaviors for different system sizes. Besides the Monte Carlo methods, renormalization group is used to study the equilibrium properties in the thermodynamic limit and to compare with the results from simulated annealing and Wang-Landau sampling.

I. INTRODUCTION

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II. RG CALCULATIONS

The setup of renormalization is started by separating the anti-ferromagnetic (AFM) Ising Hamiltonian into hierarchies

$$-\beta \mathcal{H} = \sum_{n=1}^{k-2} (-\beta \mathcal{H}_n) + \mathcal{R}(K_2, K_3, \cdots)$$
 (1)

where \mathcal{R} is the coupling beyond \mathcal{H}_n of levels k > 2. \mathcal{H}_n depends on the interactions K_0 on the backbone and L_0 , L_1, K_1, \cdots among the long range couplings. We describe the RG procedure network by network.

A. RG Analysis of HNNP and HN6 without magnetic field

1. RG setup and procedures

For HNNP and HN6, the \mathcal{H}_n for each hierarchy is

$$-\beta \mathcal{H}_n = K_0 \left(x_{n-2} x_{n-1} + x_{n-1} x_n + x_n x_{n+1} + x_{n+1} x_{n+2} \right)$$

$$+ K_1 \left(x_{n-2} x_{n+1} + x_{n-1} x_{n+2} \right) + y L_1 \left(x_{n-2} x_{n+2} \right)$$

$$+ L_0 \left(x_{n-2} x_n + x_n x_{n+2} \right) + 4I$$
 (2)

I is a constant parameter emerging after the first step of RG, and its initial value is 0. L_0 also stands for the coupling term emerged in the RG process, and its initial value is also 0. All the initials values for these activity parameters are

$$I = 0$$

 $K_0 < 0$
 $K_1 < 0$
 $L_0 = 0$ (3)

where K_1 is not changing in the RG flow because it is introduced 'freshly' at every RG step. Therefore, it can

be used as a reference of temperature. High temperatures $T \to \infty$ stands for $K_1 \to -0$; while low temperatures $T \to 0$ corresponds to large $K_1 \to -\infty$.

The renormalization trace the sites x_n and x_{n+1} to get the renormalized form

$$-\beta \mathcal{H}_n = 2I' + K_0' (x_{n-2}x_n + x_n x_{n+2}) + L_0' (x_{n-2}x_{n+2})$$
(4)

which is the same as that in HN3 and HN5.

$$C = e^{-4I} = e^{-A[0]}$$

$$\kappa = e^{-4K_0} = A[1]$$

$$\lambda = e^{-4L_0} = A[2]$$

$$\mu = e^{-2K_1} = e^{-2L_1} = A[3]$$
(5)

where A[i] are the variables used in Mathematica calculations. The corresponding initial values are

$$C = 1, A[0] = 0$$

$$\kappa = A[1] = \mu^{2} > 1$$

$$\lambda = A[2] = \mu^{2y}$$

$$\mu = A[3] > 1$$
(6)

In the calculation, we only need to specify the temperature μ , then other initial values are determined accordingly. The relationship between μ and T is,

$$T \to 0 \Leftrightarrow K_1 \to -\infty \Leftrightarrow \mu \to \infty \Leftrightarrow 1/\mu \to 0$$

$$T \to \infty \Leftrightarrow K_1 \to -0 \Leftrightarrow \mu \to 1 \Leftrightarrow 1/\mu \to 1$$
(7)

Eqs. 2 and 4 yields the RG recursions

$$\kappa' = \frac{\kappa \lambda (1+\mu)^2}{(1+\mu\kappa)^2}$$

$$\lambda' = \mu^{2y} \frac{(\kappa+\mu)^2}{(1+\mu\kappa)^2}$$
(8)

$$C' = \frac{C^2 \kappa \mu^2}{(1 + \mu^2)(\kappa + \mu)(1 + \mu \kappa)}$$

These solutions are also the same as those in Ref. [1]; we also reproduced the phase diagram of HNNP Ferromagnetic model

2. Fixed point analysis

In the AFM scenario, for HNNP (y = 0), there are only 2 possible fixed points

$$\kappa^* = 0, \ \lambda^* = \mu^2$$

$$\kappa^* = \lambda^* = 1$$
(9)

For convenience, we set the the interaction constant J = -1 and Boltzmann constant $k_B = 1$. Then the variables of interest become

$$J^{(0)} = -1,$$

 $T = 2/\log(\mu)$ (10)
 $J^{(n)} = -\frac{T}{4}\log\kappa^{(n)}$

The Jocobian matrix

$$J_{11} = \frac{\partial \kappa'(\kappa, \lambda)}{\partial \kappa} = \frac{\lambda(\mu + 1)^2}{(\kappa \mu + 1)^2} - \frac{2\kappa \lambda \mu(\mu + 1)^2}{(\kappa \mu + 1)^3}$$
(11)

$$J_{12} = \frac{\partial \kappa'(\kappa, \lambda)}{\partial \lambda} = \frac{\kappa(\mu + 1)^2}{(\kappa \mu + 1)^2}$$
 (12)

$$J_{21} = \frac{\partial \lambda'(\kappa)}{\partial \kappa} = \frac{2(\kappa + \mu)\mu^{2y}}{(\kappa \mu + 1)^2} - \frac{2(\kappa + \mu)^2 \mu^{2y+1}}{(\kappa \mu + 1)^3} \quad (13)$$

$$J_{22} = \frac{\partial \lambda'(\kappa)}{\partial \lambda} = 0 \tag{14}$$

Fixed Point Stability Done—

In this AFM model, the fixed point analysis may be messy. Some figures may be not easy to make sense. In order to address the 'messy' fixed point analysis, we show the analysis from 3 perspectives

1. Consistency

At low temperatures for certain values of y, κ is oscillating with a certain pattern along the RG steps. The questions is whether the oscillation pattern is changing for different number of RG steps. Our analysis shows that the pattern does not change after ~ 1000 RG steps. In other words, for example, the oscillation pattern of of steps (1000 \sim 5000) is the same as that of steps (5000 \sim 9000), (9000 \sim 14000), etc. Fig. 2 is an example of this 'consistency analysis'. Note for different y and μ , the histogram curves may be very different. It is easy to see the 'disorderness' of the fixed points in the following two perspectives.

The 2 Eigenvalues for HNNP (y=0)

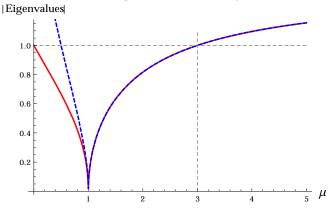


FIG. 1. The absolute values (1st norm) of the 2 complex eigenvalues for HNNP (y=0). For $0<\mu<1$, that is for the ferromagnetic case; while the antiferromagnetic region is $\mu>1$. You see the transition temperature is $\mu=3.0$ because after that the |eigenvalues|>1.

- 2. Histogram of κ at different μ and y Generally, at high temperatures (μ is small), there is usually one stable fixed point; but, at low temperatures (μ is big), κ is usually oscillating. The lower the temperature is, the wider κ spreads. Fig. 3 (positive y) and Fig. 4 (negative y) show the distribution of different κ . The next figures (Figs. 5 and 6) can help explain.
- 3. All values of κ at different μ and y. Inspired by Ref. [2] Fig 2, we plot all the κ in the RG steps (steps 18000 \sim 20000). Each $(\kappa, 1/\mu)$ is a data point in Figs. 5 and 6. You can see why the histogram plots are messy. The values of κ covers a large range and may be not very discrete.

When y < -2.0, the whole system became similar to ferromagnetic model as shown in Fig. 7.

The transition temperature can be solved by setting an equation of abs(eigenvalue) == 1 because

- 1. the eigenvalues are complex numbers, and the abs(eigenvalue) is the 1st norm;
- as what I read from mathematical literatures, for complex eigenvalues, if the 1st norm is smaller than
 the recursive equations can reach a stable fixed point; otherwise, it won't. Also, complex eigenvalues usually mean oscillating flow of the parameters;
- 3. The actually equation to get the transition is very long, complex, and not solvable analytically. We won't show it here, but we can find the transition for HNNP \sim HN6 for different y values. (An example of the eigenvalues are shown in Fig 1.)

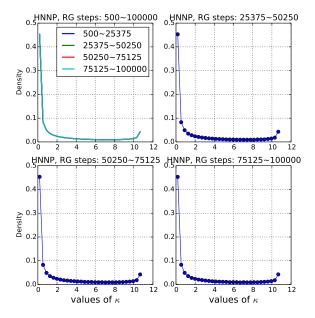


FIG. 2. Consistency of the pattern for different RG steps $(y=0; \mu=4)$. The first figure is the histogram of all the RG steps; while the rest three figures are for different step-range. They are the same, which is important for further analysis.

B. RG Analysis of HN3 and HN5 without magnetic field

1. RG setup and procedures

For HN3 and HN5, the \mathcal{H}_n for each hierarchy is

$$-\beta \mathcal{H}_{n} = K_{0} \left(x_{n-2} x_{n-1} + x_{n-1} x_{n} + x_{n} x_{n+1} + x_{n+1} x_{n+2} \right)$$

$$+ K_{1} \left(x_{n-1} x_{n+1} \right) + L_{0} \left(x_{n-2} x_{n} + x_{n} x_{n+2} \right)$$

$$+ y L_{1} \left(x_{n-2} x_{n+2} \right) + 4I$$

$$(15)$$

where y is 0 for HN3 and 1 for HN5. I is a constant emerging since the first step of RG, and its initial value is 0. L_0 also stands for the coupling term emerged in the RG flow, and its initial value is also 0. The <u>initial values</u> for these parameters are

$$I = 0$$

 $K_0 < 0$
 $K_1 < 0$
 $L_0 = 0$
 $L_1 < 0$ (16)

where K_1 is not changing in the RG flow because it is introduced again at every RG step. Therefore, it can be used as a reference of temperature. High temperatures $T \to \infty$ stands for $K_1 \to -0$; while low temperatures $T \to 0$ corresponds to large $K_1 \to -\infty$.

After tracing the sites x_{n-1} and x_{n+1} , in order to use

RG, we have to get this form of equation

$$-\beta \mathcal{H}_n = 2I' + K_0' (x_{n-2}x_n + x_n x_{n+2}) + L_0' (x_{n-2}x_{n+2})$$
(17)

which is 'half' of original RG setup. For convenience of the partition function and Mathematica calculation, we rewrite these parameters as

$$C = e^{-4I} = e^{-A[0]}$$

$$\kappa = e^{-4K_0} = A[1]$$

$$\lambda = e^{-4L_0} = A[2]$$

$$\mu = e^{-2K_1} = e^{-2L_1} = A[3]$$
(18)

The array $\{A[i]\}$ is used in the Mathematica calculations. The pre-factors of K_0, K_1, I , and L_1 is for the convenience of the RG calculations. The reason to use exponential form for C and A[0] is to utilize more precision in Mathematica. Thus, the <u>initial values</u> for them are

$$C = 1, A[0] = 0$$

$$\kappa = A[1] = \mu^{2} > 1$$

$$\lambda = A[2] = \mu^{2y} = 1$$

$$\mu = A[3] > 1$$
(19)

Generally, we only specify μ in the first step of RG; the initial values for HN3(y=0) and HN5(y=1) are determined by μ and y.

The relationship between μ and T is the opposite as between temperature T and K_1 ,

$$T \to 0 \Leftrightarrow K_1 \to -\infty \Leftrightarrow \mu \to \infty \Leftrightarrow 1/\mu \to 0$$

$$T \to \infty \Leftrightarrow K_1 \to -0 \Leftrightarrow \mu \to 1 \Leftrightarrow 1/\mu \to 1$$
(20)

By tracing over all the $\{x_n\}$ in Eq. 15 and Eq. 17, we can get equations and the following solutions

$$\kappa' = \frac{2\kappa\lambda(1+\mu)}{\kappa^2 + 2\mu\kappa + 1}$$

$$\lambda' = \mu^{2y} \frac{(1+\mu)(1+\kappa)^2}{2(\kappa^2 + 2\mu\kappa + 1)}$$

$$C' = \frac{C^2\kappa\mu}{\sqrt{2}(1+\kappa)(1+\mu)^{3/2}\sqrt{\kappa^2 + 2\mu\kappa + 1}}$$
(21)

In Mathematica, what we get equivalently is actually

$$A'[1] = \frac{2A[1]A[2](1+\mu)}{1+2\mu A[1] + A[1]^2}$$

$$A'[2] = \mu^{2y} \frac{(1+\mu)(1+A[1])^2}{2(1+2\mu A[1]+A[1])}$$

$$A'[0] = \frac{1}{2} \log \left(\frac{e^{-4A[0]} \mu^2 A[1]^2}{2(1+\mu)^3 (1+A[1])^2 (1+2\mu A[1]+A[1]^2)} \right)$$
(22)

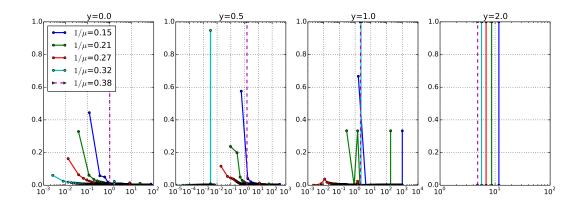


FIG. 3. Histogram of κ at different μ and positive y of steps 1000 \sim 20000 (Density vs κ). For y=0 (HNNP), the transition happens at $\mu=3$, i.e., for $\mu\leq 3.0$, there is one fixed point at $\kappa=1$, while, for $\mu>3.0$, κ is oscillating and never converge to one point. Figs. 5 and 6 can help explain.

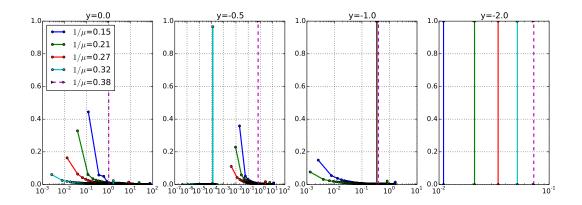


FIG. 4. Histogram of κ at different μ and negative y of steps $1000 \sim 20000$ (Density vs κ). The legend in the 1st subfigure works for all subfigures. Figs. 5 and 6 can help explain.

2. Fixed point analysis

For HN3 with y=0, we get 2 analytical fixed point solutions which are

$$\kappa = 1, \lambda = 1;$$

$$\kappa = 0, \lambda = \frac{1+\mu}{2}$$
(23)

where only the first one is a stable fixed point solution. The second one is only stable at infinitely high temperatures; any finite temperatures lead to the fixed point $\kappa = 1, \lambda = 1$.

For HN5 with y=1, the fixed points are more interesting. The fixed point solution are

$$\kappa = 0, \quad \lambda = \mu^{2y} \frac{\mu + 1}{2} \tag{24}$$

$$\kappa = \frac{1}{2} \left[\mu^2 - \mu + \sqrt{(\mu + 1)(\mu^3 - 3\mu^2 + 8\mu - 4)} \right]$$
 (25)

$$\lambda = \frac{\mu}{4} \left[\mu^2 - \mu + 2 + \sqrt{(\mu + 1)(\mu^3 - 3\mu^2 + 8\mu - 4)} \right]$$
(26)

For the parameter of interest κ , its solution Eq.25 can be written in a general form as a function of μ ($\mu > 1$)

$$\kappa = \frac{1}{2} \left[\mu^{1+y} + \mu^y - 2\mu + \sqrt{4(\mu^{1+y} + \mu^y - 1) + (\mu^{1+y} + \mu^y - 2\mu)^2} \right]$$
(27)

 κ changes with μ ; the values of μ satisfying $\kappa(\mu) = 0$ are likely to be the phase transition temperatures. After transformation of Eq.27, the equation $\kappa(\mu) = 0$ can be rewritten as

$$\mu^{1+y} + \mu^y - 1 = 0 \tag{28}$$

where, for any values of y, the real roots of $\mu > 1$ are the acceptable solutions. The examples with specific values

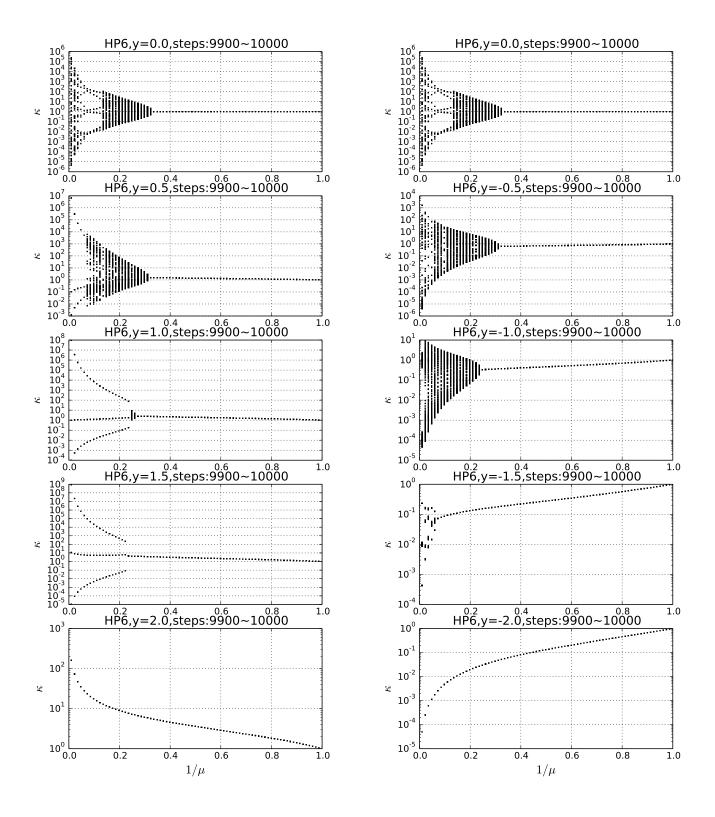


FIG. 5. All values of κ with position y for steps 18000-20000. This figure is similar to Ref [2] Fig.2. The reason of just using 2000 (20000 - 18000) is because of cercern of figure file size. The more steps included, the more data points in the figure, which makes the figure file very large and slow to open.

FIG. 6. All values of κ with negative y for steps 18000 – 20000. This figure is similar to Ref [2] Fig.2.

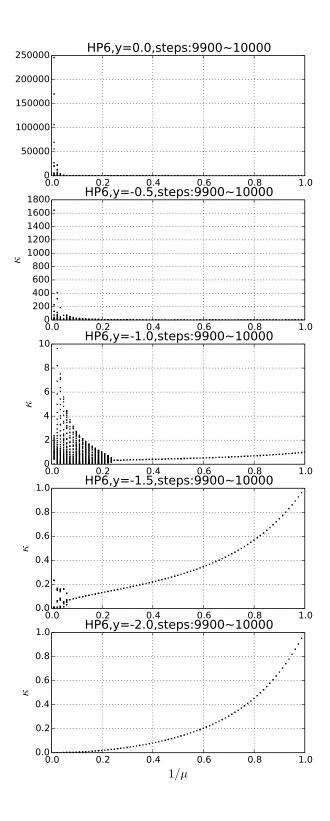


FIG. 7. All values of κ with negative y for steps 18000-20000. This figure is similar to Ref [2] Fig.2. $y\to\infty:\ \mu^{1+y}+\mu^y-1=0;\ \mu\to1(\ {\rm FM},\ T\to\infty)$

$$y \to \infty$$
: $\mu^{1+y} + \mu^y - 1 = 0$; $\mu \to 1$ (FM, $T \to \infty$)

$$y = 2: \mu^3 + \mu^2 - 1 = 0; \mu = 0.754878 \text{ (FM)}$$

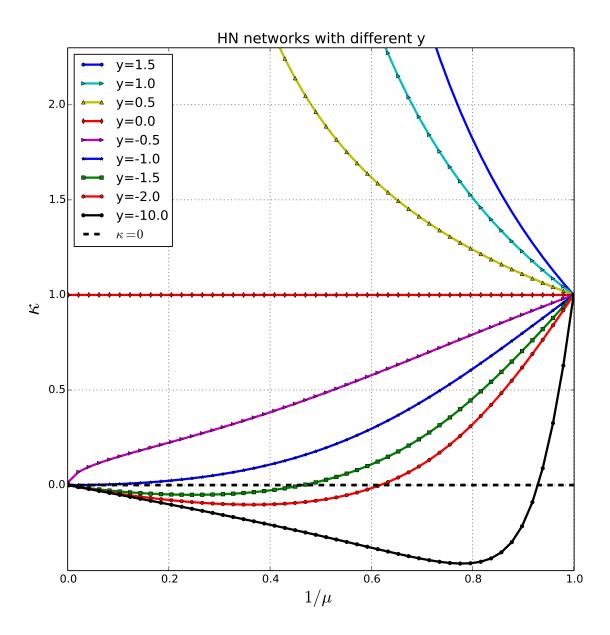


FIG. 8. κ vs $1/\mu$ calculating using Eq.27. y=0 corresponds to HNNP; y=1 corresponds to HN6; and y<0 corresponds that some bonds are ferromagnetic;. In AF model, our temperature parameter μ is greater than 1. In this case, $1/\mu=0 \Leftrightarrow T=0$; $1/\mu=1 \Leftrightarrow T \to \infty$.

^[1] S. Boettcher and C. Brunson, Physical Review E $\bf 83$, 021103 (2011).

^[2] S. R. McKay, A. N. Berker, and S. Kirkpatrick, Physical Review Letters 48, 767 (1982).