Dynamics of an Ising spin glass on the Bethe lattice

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Abstract

We study the dynamical low-temperature behaviour of the Ising spin glass on the Bethe lattice. Starting from Glauber dynamics we propose a cavity such as an ansatz that allows for the treatment of the slow (low temperature) part of dynamics. Assuming a continuous phase transition and ultrametricity with respect to long time scales we approach the problem perturbatively near the critical temperature. The theory is formulated in terms of correlation–response functions of arbitrary order. They can, however, be broken down completely to products of pair functions depending on two time arguments only. For binary couplings $J = \pm I$ a spin glass solution is found which approaches the corresponding solution for the SK model in the limit of high connectivity. For more general distributions $P(J)$ no stable or marginal solution of this type appears to exist. The nature of the low-temperature phase in this more general case is unclear.

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This paper is dedicated to David Sherrington on the occasion of his 65th birthday.

1. Introduction

Most of the work on spin glasses and other systems with frozen-in disorder is based on the evaluation of the free energy or the ground-state energy. This comprises replica theory, the cavity method or the TAP equations [1]. As an alternative stochastic dynamics [2] has been employed for systems with continuous freezing transition [3, 4], e.g. the Sherrington–Kirkpatrick (SK) model [5], and also for systems with discontinuous transition, e.g. the spherical spin glass with $p$-spin interactions [6–8]. The essential difference between the two

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approaches shows up in the thermodynamic limit. The computation of the free energy does not rely on any kind of dynamics and the question of how equilibrium states can be reached starting from non-equilibrium initial conditions is not addressed. In fact diverging barriers build up in the thermodynamic limit \( N \to \infty \). As a consequence the approach via dynamics in the non-ergodic low-temperature phase requires regularization by some long time scale, for instance the waiting time \( t_w \) after a quench from high temperature, by introducing slowly varying bonds [4] or slow cooling [9]. Typically the thermodynamic and long time limit are taken in the order \( \lim_{t_w \to \infty} \lim_{N \to \infty} \). This means that diverging barriers cannot be overcome or the system might be stuck in metastable states [8, 10]. The importance of the order in which the two limits are performed has been demonstrated in a different context (learning in a perceptron with binary synapses) [11]. For systems with discontinuous transition the dynamic freezing transition is higher than the temperature where replica symmetry breaking sets in, whereas both temperatures are identical for continuous transitions. In this latter case the coefficients of an expansion of the energy or the Edwards–Anderson order parameter near \( T_c \) are identical at least up to the fourth order [9]. At zero temperature, however, the energy obtained within dynamics is expected to be higher than the ground-state energy determined within replica theory. Furthermore, the energy found within dynamics may depend on details of the cooling schedule [12].

Spin glasses on diluted graphs, for instance on the Bethe lattice, share finite connectivity with spin glasses in finite dimensions. Nevertheless they are of mean field character as the fully connected SK model [1, 5]. Following earlier attempts [14, 15], the Ising spin glass on the Bethe lattice has been solved by Mézard and Parisi [13] using the cavity method on the level of one-step replica symmetry breaking. The cavity method benefits from the local tree-like structure of the Bethe lattice and deals on the level of 1RSB with distributions of local fields and distributions of those distributions. The resulting functional equations have been solved numerically with a population dynamics algorithm for binary couplings \( J = \pm I \).

Investigating the dynamics of systems of Ising spins one has to specify not only the Hamiltonian (energy) but also the kind of dynamics. In the present approach Glauber dynamics, i.e. stochastic single spin flip dynamics with transition probabilities depending on temperature, is used. In the non-ergodic low-temperature phase a separation of short and long time scales is assumed. The longest time scale is realized by a waiting time \( t_w \) or by some other means [4, 9]. On the short time scale the validity of fluctuation–dissipation theorems are assumed to hold. They are characteristics for equilibrium and in this context they describe equilibrium within a single valley of the energy landscape. On the long time scale in analogy to the cavity method [13] a functional equation for the distribution of histories of local fields is derived, investigating the iterative assembly of local subtrees of the Bethe lattice to new subtrees. Probability distributions for histories have previously been used by Hatchett et al [16] in a similar context. In contrast to the present investigation this paper deals, however, with parallel updates and short time dynamics.

Based on this functional equation, hierarchically connected relations are derived for correlation–response functions on the long time scale, connecting functions of arbitrary order. In order to truncate this system, an expansion in powers of the departure of the temperature from the freezing temperature is employed. Such an expansion is known for the dynamics of the SK model. It turns out that correlation–response functions of higher order can be broken down completely to expressions involving pair functions depending on two time arguments only. For binary couplings \( J = \pm I \) a marginal solution resembling the one for the SK model is found [3, 4].

The situation for more general distributions \( P(J) \) is unclear. There exists a critical temperature where the ergodic high temperature solution becomes unstable. The above
solution does, however, not apply and no other stable or marginal solution could be found. The expansion contains contributions corresponding to the one-step replica symmetry breaking solution. The original static 1RSB solution [13] has been evaluated for \( J = \pm I \) only. It would be interesting to see whether this solution persists for general \( P(J) \).

The paper is organized as follows. Section 2 contains a brief survey of Glauber dynamics and fixes some notations. In section 3 the dynamics of the Ising spin glass on the Bethe lattice is formulated, an effective single-site evolution is introduced and equilibrium properties are investigated. Dynamics on long time scales is dealt with in section 4 including a formulation in terms of distributions of slow components of local fields. In view of an expansion valid in the neighbourhood of the freezing temperature, this is reformulated in terms of correlation–response functions. Section 5 contains the expansion around the critical temperature in leading order for general \( P(J) \) and in next to leading order for \( J = \pm I \). The paper concludes with a stability analysis in section 6 and a brief discussion.

2. Glauber dynamics of a single spin

The present investigation is based on Glauber dynamics for Ising spins. We start with a brief outline of the formulation [11] describing the dynamics of a single spin. The Hamiltonian of the spin in an external field \( h \) is

\[
H = -h \sigma_z. \tag{2.1}
\]

The actual state of the spin at time \( t \) is described by a two-component vector

\[
|\rho(t)\rangle = \left( \begin{array}{cc} p_+(t) \\ p_-(t) \end{array} \right) = \left( \begin{array}{c} 1/2(1+m(t)) \\ 1/2(1-m(t)) \end{array} \right). \tag{2.2}
\]

Introducing the Pauli matrix

\[
\sigma_z = \left( \begin{array}{cc} 1 & 0 \\ 0 & -1 \end{array} \right) \tag{2.3}
\]

the expectation value of the spin is given by

\[
\langle \sigma_z \rangle = \frac{1}{2} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix} \begin{pmatrix} p_+(t) \\ p_-(t) \end{pmatrix} = \langle 1|\sigma_z|\rho(t)\rangle = m(t). \tag{2.4}
\]

It is convenient to use a notation resembling quantum mechanics, although Glauber dynamics is purely classical. The time evolution is described by a Liouville operator, a two-by-two matrix, acting on the state

\[
\dot{\rho} = L(t)|\rho(t)\rangle \tag{2.5}
\]

with Liouvillian

\[
L(t) = -\frac{1}{2}\gamma \begin{pmatrix} 1 - \tanh(\beta h(t)) & -1 - \tanh(\beta h(t)) \\ -1 + \tanh(\beta h(t)) & 1 + \tanh(\beta h(t)) \end{pmatrix}. \tag{2.6}
\]

In equilibrium, for constant \( h \), \( L|\bar{\rho}\rangle = 0 \). This determines the equilibrium state

\[
|\bar{\rho}\rangle = \frac{1}{2} \begin{pmatrix} 1 + \tanh(\beta h) \\ 1 - \tanh(\beta h) \end{pmatrix} = \frac{e^{-\beta h}}{2 \cosh(\beta h)} |1\rangle \tag{2.7}
\]

with \(|1\rangle\) being the two-dimensional unit vector. The temporal evolution of the spin is

\[
\dot{\sigma}_z(t) = \sigma_z L(t) = L(t)\sigma_z. \tag{2.8}
\]
In addition to the spin operator a response operator \( \hat{\sigma}_z(t) \) is introduced. It describes the action of a variation \( \delta h(t) \) of the external field at time \( t \)

\[
\hat{\sigma}_z = \frac{\partial L(t)}{\partial h(t)}.
\]

(2.9)

Acting on an equilibrium state these operators obey a fluctuation–dissipation theorem

\[
\hat{\sigma}_z \rho = \beta \hat{\sigma}_z \rho.
\]

(2.10)

For time-dependent field \( h(t) \) the Liouvillians for different times do not commute in general. This can formally be overcome by introducing a time ordering operator \( T \) rearranging products of objects at different time such that they are ordered from left to right according to the decreasing time. This allows us to write for \( t > t_0 \)

\[
|\rho(t)\rangle = T \left[ e^{\int_{t_0}^{t} \sum_i L_i(k_i(s)) \, ds} \right] |\rho(t_0)\rangle.
\]

(2.11)

3. Ising model on the Bethe lattice

3.1. Dynamics

We consider Ising spins \( \sigma_i \) on the Bethe lattice with couplings \( J_{ij} \). The Hamiltonian (energy) is

\[
H = -\frac{1}{2} \sum_{i,j} J_{ij} \sigma_i \sigma_j - \sum_i h_i \sigma_i.
\]

(3.1)

The couplings \( J_{ij} \) are independent stochastic variables taken from some distribution \( P(J) \) such that \( J_{ij} = 0 \) and \( J_{ij}^2 = \mathcal{I}^2 \). The underlying vector space for the complete system is spanned by the direct product of the two-dimensional vectors associated with each of the spins \( \sigma_i \).

The dynamics is ruled by the Liouvilian

\[
L(t) = \sum_i L_i(k_i)
\]

(3.2)

where \( L_i(k_i) \) given by (2.6) acts on the corresponding subspace of site \( i \) and depends on the effective field

\[
k_i = \frac{\partial H}{\partial \sigma_i} = h_i + \sum_j J_{ij} \sigma_j.
\]

(3.3)

Expectation values of time-dependent observables with initial condition at \( t_0 \) are written as

\[
\langle A(t) B(t') \cdots \rangle = \langle 1_N A(t) T \left[ e^{\int_{t_0}^{t} \sum_i L_i(k_i(s)) \, ds} \right] B(t') T \left[ e^{\int_{t'}^{\cdots} \sum_i L_i(k_i(s)) \, ds} \right] \cdots \rangle_N
\]

(3.4)

for \( t > t' > \cdots > t_0 \).

Adopting in the following Ito-calculus the effective fields (3.3) are retarded, i.e.

\[
k_i(t) = h + \sum_j J_{ij} \sigma_j(t^-).
\]

(3.5)

It is convenient to integrate over the effective fields \( k_i(t) \) in (3.4) and to take (3.5) into account by introducing appropriate \( \delta \)-functions. Those are written in their Fourier representation
integrating in addition over imaginary auxiliary fields \( \hat{k}_i(t) \). This leads to a path integral representation

\[
\langle A(t) B(t') \cdots \rangle = \prod_i \int \mathcal{D}[\hat{k}_i, k_i] \ e^{\sum_i \int_0^T d\hat{k}_i(t) [\hat{h} - \hat{k}_i(t)]}
\]

\[
\times (1|N \mathcal{T}[A(t) B(t') \cdots \sum_i \int_0^T d\hat{k}_i(t) \sum_j J_{ij} \sigma_i(t) + \mathcal{L}_s(\hat{k}_i(t), t)]]) \rho(t_0) \rangle_N.
\]

A state \(|\rho\rangle_N\) of the system with \( N \) sites is in general a superposition of direct products of two-dimensional vectors each representing a single spin. The unit vector for \( N \) sites \(|1\rangle_N\) is a direct product of single-site unit vectors. \(|1\rangle_N\) is its adjoint.

The equilibrium state, for given \( J \), can be expressed as

\[
|\bar{\rho}\rangle_N = Z^{-1} e^{\sum_i h_i \sigma_i + \frac{1}{2} \sum_i J_{ij} \sigma_i \sigma_j} |1\rangle_N.
\]

It obeys

\[
\mathcal{L}_s(k_i) |\bar{\rho}\rangle_N = 0.
\]

### 3.2. Effective single-site evolution

For time-dependent expectation values involving the spin at a single site, say site \( o \) only, (3.6) can be rewritten introducing an effective retarded time evolution. It is obtained by performing the expectation values with respect to all spins except \( \sigma_o \). Assuming equilibrium initial conditions expectation values of quantities at a single site, say \( o \), can be written as (3.6),

\[
\langle \sigma_o(t) \cdots \rangle = \int \mathcal{D}[\hat{k}, k] \ e^{\sum_i \int_0^T d\hat{k}_i(t) [\hat{h} - \hat{k}_i(t)]}
\]

\[
\times (1|o \mathcal{T}[\sigma_o(t) \cdots \sigma_o(k)] |\bar{\rho}_o(k(t_0))\rangle_o).
\]

The action of the remaining part of the system is contained in

\[
\mathcal{Y}_o([\hat{k}]|\sigma_o); t_0) = \prod_i \int \mathcal{D}[\hat{k}_i, k_i] \ e^{\sum_i \int_0^T d\hat{k}_i(t) [\hat{h} - \hat{k}_i(t)]}
\]

\[
\times (1|N-o \mathcal{T}[\sum_i \int_0^T d\hat{k}_i(t) \sigma_i(t) + \mathcal{L}_s(\hat{k}_i(t), t)]]) \bar{\rho}_o(k(t_0))\rangle_{N-o}.
\]

and the initial condition in (3.9) is

\[
|\bar{\rho}_o(k)\rangle_o = \frac{1}{2 \cosh(\beta k)} e^{\beta \sigma_o} |1\rangle_o.
\]

In the above expression the primed products and sums run over \( i \neq o \). The expectation values in (3.10) refer to the system without the spin at site \( o \).

It should be pointed out that the above expression holds for given values of the couplings \( J_{ij} \) and the average over \( J_{ij} \) has still to be evaluated.

### 3.3. Equilibrium

The above path integral comprises an integration at the initial time \( t_0 \). This is investigated in the following. The single time (static) expectation value of \( \sigma_o(t_o) \) is

\[
\langle \sigma_o(t_o) \rangle = Z^{-1} \int \frac{d\hat{k} dk}{2\pi} e^{\{\hat{h} - \hat{k}\}} (1|\sigma_o(\hat{k}) \bar{\rho}(\hat{k}) e^{\beta \sigma_o} |1\rangle.
\]
and the effective action (3.10) reduces to
\[ Y_o(h) = \left\{ e^{\sum J_{io} \sigma_i} \right\}_{N-1}. \] (3.13)

In analogy to the cavity method we define \( c_o \) and \( k_o \) writing
\[ e^{\beta(c_o + k_o \sigma)} = \int \frac{dk}{2\pi} e^{\hat{k}(h-k)} Y_o(\hat{k}) e^{\beta \sigma}. \] (3.14)

This yields for the partition function and the local magnetization
\[ Z_o = 2 e^{\beta c_o} \cosh(\beta k_o) \quad \text{and} \quad \langle \sigma_o \rangle = \tanh(\beta k_o). \] (3.15)

The two quantities defined above result in
\[ c_o = \frac{1}{\beta} \left\{ \ln \left( \left\{ e^{\beta \sum J_{io} \sigma_i} \right\}_{N-1} \right) + \ln \left( \left\{ e^{-\beta \sum J_{io} \sigma_i} \right\}_{N-1} \right) \right\}, \] (3.16)
\[ k_o = h + \frac{1}{2\beta} \left\{ \ln \left( \left\{ e^{\beta \sum J_{io} \sigma_i} \right\}_{N-1} \right) - \ln \left( \left\{ e^{-\beta \sum J_{io} \sigma_i} \right\}_{N-1} \right) \right\}. \] (3.17)

For the SK model and the Bethe lattice the expectation values for sites \( i \) can be assumed to factorize. For Ising spins and using (3.15) we find the identity
\[ \langle e^{\pm \beta J_{io} \sigma_i} \rangle_{N-1} = \cosh(\beta J_{io}) \pm \tanh(\beta k_i) \sinh(\beta J_{io}) \] (3.18)
which allows us to express \( c_o \) and \( k_o \) in terms of the fields of the adjacent sites:
\[ c_o = \sum_i c(J_{io}, k_i), \quad k_o = h + \sum_i u(J_{io}, k_i), \] (3.19)
\[ c(J, k) = \frac{1}{2\beta} \ln(\cosh^2(\beta J) - \tanh^2(\beta k)) \sinh(\beta J)), \] (3.20)
\[ u(J, k) = \frac{1}{\beta} \text{artanh}(\tanh(\beta k) \tanh(\beta J)). \] (3.21)

In particular for \( J \to 0 \), i.e. for the SK model,
\[ u(J, k) = J \tanh(\beta k), \quad c(J, k) = \frac{\beta J^2}{2 \cosh^2(\beta k)}. \] (3.22)

In view of this limit correlation functions are later defined with rescaled quantities
\[ U(J, \kappa) = \frac{1}{I} u(J, \kappa), \] (3.23)
where \( I \) is the typical size of the couplings. This ensures, e.g., that the Edwards–Anderson order parameter stays finite in the SK limit.

4. Slow dynamics

4.1. Distribution of slow fields

Dynamics on short time scales takes place within a single valley of the energy landscape; dynamics on long time scales is supposed to be due to transitions among different valleys. Assume there exists a time scale \( t^* \gg 1 \) separating fast and slow motions. This time scale might be realized by some waiting time, slowly changing couplings or other means [4, 9]. The local field \( \hat{k}(t) \) and the conjugate field \( \dot{\hat{k}}(t) \) are split into fast and slow contributions
\[ k(t) \to k(t) + \kappa(t/t^*), \quad \dot{\hat{k}}(t) \to \dot{\hat{k}}(t) + \frac{1}{t^*} \hat{k}(t/t^*). \] (4.1)
The fast parts are due to the fluctuations around the quasi-equilibrium state within a single valley. The slow part $\kappa(\tau)$ acts like an external field and the quasi-equilibrium state follows this field adiabatically. Evaluating time-dependent correlation functions of the spin $\sigma_o$ on this long time scale, this spin can be viewed as being in equilibrium in the slow field resulting in

$$\langle \sigma_o(t^* \tau) \sigma_o(t^* \tau') \cdots \rangle = \tanh(\beta \kappa_o(\tau)) \tanh(\beta \kappa_o(\tau')) \cdots. \quad (4.2)$$

The effective field is given by (3.19) with $k$ replaced by $\kappa(\tau)$, i.e.

$$\kappa_o(\tau) = h + \sum_i u(J_{oi}, \kappa_i(\tau^-)). \quad (4.3)$$

In the low-temperature regime of a spin glass this field is, however, distributed and (4.2) has to be averaged over some distribution $P_o(k_0)$. With (4.3) this distribution can be calculated from the distribution of slow fields at the neighbouring sites $i$. Continuing this mapping to outer shells of the lattice an iteration scheme is set up similar to the course of action in the cavity approach [13]. This program involves the following steps.

Performing the average over the couplings $J_{ij}$ and defining

$$Q_o([\hat{k}, \kappa]) = \left\langle e^{\int dt \sum [h(J, \kappa_i(\tau^-)) + h(J, \kappa_o(\tau^-))]} \right\rangle^{N-1}_{\hat{k}} \quad (4.4)$$

this distribution is obtained by integration over $\hat{k}$

$$P_o([\kappa]) = \int D[\hat{k}] e^{\int dt \langle [h(\hat{k}(\tau)) + h(\hat{k}(\tau^-))] \rangle_{\hat{k}}^2} Q_o([\hat{k}, \kappa]). \quad (4.5)$$

The average on the right-hand side of (4.4) involves a corresponding joint distribution $Q_{N-1}$ of the fields $\hat{k}_i$ and $\kappa_i$ at the $K + 1$ neighbouring sites. The second term in the exponent is the contribution of site $o$ to the local field at site $i$, corresponding to the Onsager reaction field.

Assuming factorization of the expectation values on the right-hand side of (4.4) leads to

$$Q_o([\hat{k}, \kappa]) = \prod_i \int D[\hat{k}_i, \kappa_i] e^{\int dt \langle [h(\hat{k}_i(\tau)) + h(\hat{k}_i(\tau^-))] \rangle_{\hat{k}_i}^2} Q_o([\hat{k}_i, \kappa_i]) \times e^{\int dt \langle [h(J_{oi}, \kappa_o(\tau^-)) + h(J_{oi}, \kappa_i(\tau^-))] \rangle_{J_{oi}}^2}. \quad (4.6)$$

The functional $Q_o([\hat{k}, \kappa])$ contains the action of all $K + 1$ spins surrounding site $o$. $Q_i([\hat{k}_i, \kappa_i])$, on the other hand, contains only the action of the $K$ spins on the outgoing branches originating at site $i$. The action of the spin $\sigma_o$ is taken into account by the bond-averaged exponential in (4.6). The functional $Q_i$ is given by a similar average over the distributions on the next shell of sites on the tree. Iterating this process, a fixed point distribution is assumed to exist. It is given by

$$Q([\hat{\kappa}, \kappa]) = \left[ \int D[\hat{\lambda}, \lambda] e^{\int dt \langle [h(\hat{\lambda}(\tau)) + h(\hat{\lambda}(\tau^-))] \rangle_{\hat{\lambda}}^2} Q([\hat{\lambda}, \lambda]) e^{\int dt \langle [h(J, \lambda_o(\tau^-)) + h(J, \lambda_o(\tau^-))] \rangle_{J_o}} \right]^K. \quad (4.7)$$

The functional $Q_o([\hat{k}, \kappa])$ for the central spin at site $o$ is given by the same expression with $K$ replaced by $K + 1$.

The above line of arguments follows pretty much the steps in the cavity method calculation of Mézard and Parisi [13]. The functional fixed point equation (4.7) is, however, more general. The only assumptions made are the separation of time scales and the factorization of expectation values on different subtrees. In particular, it contains in principle contributions corresponding to the full replica symmetry breaking solutions. Within this formulation the replica symmetric solution is obtained by leaving out the reaction term in (4.7) as shown later.
4.2. Correlation–response functions

There is little hope to find solutions without relying on approximations or expansions. In the following we investigate an expansion in powers of the difference $\epsilon = \frac{T_c - T}{T_c}$ between transition temperature $T_c$ and the actual temperature $T$. Such an expansion is expected to apply in the neighbourhood of a continuous freezing transition where the Edwards–Anderson order parameter $q_{EA} \sim \epsilon$ near $T_c$. This is the case for the SK model.

Rather than working with the full functional $Q(\{\hat{k}, \kappa\})$, it is sufficient to investigate its moments, the correlation–response functions on the long time scale. With the rescaled quantities $U(J, \kappa)$, (3.23), they are defined as

$$C_{nm}(\tau_1, \ldots, \tau_n; \tau'_1, \ldots, \tau'_m; J) = \int \mathcal{D}[\hat{k}, \kappa] e^{\int d\kappa(\sigma) \hat{h}_c(\sigma)} Q(\{\hat{k}, \kappa\})$$

$$\times \prod_{\nu=1}^n U(J, \kappa(\tau_\nu)) \prod_{\mu=1}^m \hat{k}(\tau'_\mu).$$  (4.8)

Inserting (4.7) and expanding the exponential, a hierarchy of equations for the correlation–response functions is obtained which is derived along the following lines: evaluating a function $C_{nm}(\tau_1, \ldots, \tau_n; \tau'_1, \ldots, \tau'_m; J)$ there is the product over $U(J, \kappa(\tau_\nu))$ involving $\kappa$ at the times $\tau_\nu$. The factors $\delta / \delta \kappa(\tau_\nu)$ can be replaced under the path integral by functional derivatives $\delta / \delta \kappa(\tau_\nu)$ acting on $Q(\{\hat{k}, \kappa\})$, at least as long as all time arguments are different (otherwise the derivatives could act on the products of $U(J, \kappa(\tau_\nu))$ as well). The derivatives acting on $Q$ create additional terms involving $\kappa(\tau'_\mu)$. Expanding the second term in the exponential of (4.7), internal integrations over times $\sigma_\rho$ are generated. They contribute additional factors now depending on $\kappa(\sigma_\rho)$. In order $l$ of this expansion there are altogether functions of $n + m + l$ discrete times. For $\tau \neq \tau_\nu, \tau'_\mu, \sigma_\rho$ the path integral over $\kappa(\tau)$ can be performed resulting in $\hat{k}(\tau) = 0$. This leaves integrations over $\hat{k}$ and $\kappa$ at those discrete times only. With (4.7) they reduce to functions of correlation–response functions. This general strategy is applied in the following to correlation–response functions of lowest order.

5. Expansion around $T_c$

5.1. Leading order for general $P(J)$

We investigate an expansion in powers of $\epsilon = \frac{T_c - T}{T_c}$. For the SK model $C_{nm} \sim \epsilon^{(n+3m)/2}$ is found [3, 4] and the same scaling with $\epsilon$ is proposed for the present investigation as well. Later it is shown that this scaling is fulfilled in a consistent manner. For simplicity the following discussion will be restricted to $h = 0$.

We may start with an expansion of (4.7)

$$Q(\{\hat{k}, \kappa\}) \approx \left[1 + \frac{1}{2} I^2 \int \! d\sigma \! d\sigma' \hat{k}(\sigma) \hat{k}(\sigma') C_{2,0}(\sigma, \sigma'; J)''ight.$$

$$+ I^2 \int \! d\sigma \! d\sigma' \hat{k}(\sigma) \hat{k}(\sigma') U(J', \kappa(\sigma')) C_{1,1}(\sigma, \sigma'; J)''$$

$$+ \frac{1}{4!} I^4 \int \! d\sigma_1 \! d\sigma_2 \! d\sigma_3 \! d\sigma_4 \hat{k}(\sigma_1) \hat{k}(\sigma_2) \hat{k}(\sigma_3) \hat{k}(\sigma_4) C_{4,0}(\sigma_1, \sigma_2, \sigma_3, \sigma_4; J)''$$

$$+ \frac{1}{3!} I^4 \int \! d\sigma_1 \! d\sigma_2 \! d\sigma_3 \! d\sigma_4 \hat{k}(\sigma_1) \hat{k}(\sigma_2) \hat{k}(\sigma_3) \hat{k}(\sigma_4) \hat{U}(J', \kappa(\sigma_4)) C_{3,1}(\sigma_1, \sigma_2, \sigma_3; J)''$$

$$+ \cdots \right]^{\hat{k}}.$$  (5.1)
Evaluating $C_{2,0}(\tau, \tau'; J)$ and keeping terms $\sim \epsilon^2$

$$C_{2,0}(\tau, \tau'; J) = \int \mathcal{D}[\hat{\kappa}, \kappa] \exp \left[ i \kappa(\sigma) - \kappa(\sigma') \right] U(J, \kappa(\tau)) U(J, \kappa(\tau'))$$

$$\times \left\{ 1 + \frac{1}{2} I^2 K \int \sigma \sigma' \kappa(\sigma) \kappa(\sigma') C_{2,0}(\sigma, \sigma'; J') \right.$$ 

$$\left. + \frac{1}{4!} I^4 K \int \sigma_1 \sigma_2 \sigma_3 \sigma_4 \kappa(\sigma_1) \kappa(\sigma_2) \kappa(\sigma_3) \kappa(\sigma_4) \left[ C_{4,0}(\sigma_1, \sigma_2, \sigma_3, \sigma_4; J') \right] \frac{1}{C_{2,0}(\sigma_1, \sigma_2; J') C_{2,0}(\sigma_3, \sigma_4; J')} \right\}. \quad (5.2)$$

Performing the path integration over $\kappa(\sigma)$ for $\sigma \neq \tau, \tau'$ the integrals over $\sigma$ reduce to sums over $\sigma = \tau, \tau'$ and the path integrals become ordinary integrations over variables $\hat{\kappa} = \hat{\kappa}(\tau), \kappa = \kappa(\tau), \hat{\kappa}' = \hat{\kappa}(\tau')$ and $\kappa' = \kappa(\tau')$.

It is convenient to introduce vertices

$$V_n(J) = \left. \frac{\partial^{n-1}}{\partial \kappa^{n-1}} U(J, \kappa) \right|_{\kappa = 0}$$

with $V_n = 0$ for odd $n$. In the following we use the notation $\overline{A(J)B(J)} \rightarrow AB$. Then (5.2) yields

$$C_{2,0}(\tau, \tau'; J) = \left\{ I^2 K V^2(J) + \frac{1}{2} I^2 K (K - 1) V_2(J) V_4(J) \left[ \overline{C_{2,0}(\tau, \tau)} + \overline{C_{2,0}(\tau', \tau')} \right] \right.$$ 

$$\left. \times \overline{C_{2,0}(\tau, \tau')} + \frac{1}{4!} I^4 K V_2(J) V_4(J) \left[ \overline{C_{4,0}(\tau, \tau, \tau', \tau)} + \overline{C_{4,0}(\tau', \tau, \tau, \tau')} \right] \right\}. \quad (5.4)$$

Evaluation of $C_{4,0}$, again with $h = 0$, in order $\epsilon^2$ results in

$$C_{4,0}(\tau_1, \tau_2, \tau_3, \tau_4; J) = I^4 K V^2(J) \overline{C_{2,0}(\tau_1, \tau_2, \tau_3, \tau_4)} + I^4 K (K - 1) V_2(J)$$

$$\times \left[ \overline{C_{2,0}(\tau_1, \tau_2) C_{2,0}(\tau_3, \tau_4)} + \overline{C_{2,0}(\tau_1, \tau_3) C_{2,0}(\tau_2, \tau_4)} + \overline{C_{2,0}(\tau_1, \tau_4) C_{2,0}(\tau_2, \tau_3)} \right] \quad (5.5)$$

and performing the average over $J$ we get

$$\overline{C_{4,0}(\tau_1, \tau_2, \tau_3, \tau_4)} = \frac{I^4 K (K - 1)}{1 - I^2 K V^2} \left[ \overline{C_{2,0}(\tau_1, \tau_2) C_{2,0}(\tau_3, \tau_4)} + \overline{C_{2,0}(\tau_1, \tau_3) C_{2,0}(\tau_2, \tau_4)} + \overline{C_{2,0}(\tau_1, \tau_4) C_{2,0}(\tau_2, \tau_3)} \right]. \quad (5.6)$$

With this $C_{2,0}$ becomes

$$C_{2,0}(\tau, \tau'; J) = \left\{ I^2 K V^2(J) + \frac{1}{2} I^2 K (K - 1) \frac{V_2(J) V_4(J)}{1 - I^4 K V^2} \left[ \overline{C_{2,0}(\tau, \tau)} + \overline{C_{2,0}(\tau', \tau')} \right] \right.$$ 

$$\left. \times \overline{C_{2,0}(\tau, \tau')} \right\}. \quad (5.7)$$

It should be noted that the correlation function $\overline{C_{4,0}}$ has been broken down completely to products of pair functions $\overline{C_{2,0}}$. This can be done for correlation–response functions $\overline{C_{n,m}}$ with general $n$ and $m$ in higher orders as well. In this case the products also contain response functions $\overline{C_{1,1}}$.

Averaging over $J$ a non-trivial solution with $\overline{C_{2,0}(\tau, \tau')} \neq 0$ requires that the expression in the curly bracket on the right-hand side equals 1. Including terms of higher orders in $\epsilon$ shows that a corresponding expression holds only for the time derivative. Assuming time translational invariance we can write

$$\partial_\tau \overline{C_{2,0}(\tau)} = Y(\tau) \partial_\tau \overline{C_{2,0}(\tau)} \quad (5.8)$$
with

\[ Y(\tau) = I^2 K \overline{V_2^2} + I^4 K (K - 1) \frac{\overline{V_2 V_4}}{1 - I^2 K \overline{V_2^4}} \overline{c_{2,0}(0)}. \] (5.9)

In the first order \( Y(\tau) \) is constant and (5.8) would hold for \( \overline{c_{2,0}(\tau)} \) as well. Including higher orders, however, it depends explicitly on \( \tau \) and (5.8) holds for the derivative only. A non-trivial solution requires \( Y(\tau) = 1 \) for all \( \tau \) and this eventually determines the full time dependence of \( \overline{c_{2,0}(\tau)} \).

The evaluation of \( C_{1,1} \) follows similar lines. Since \( C_{1,1} \sim \epsilon^2 \) contributions \( \sim \epsilon^3 \) are kept, and the response function becomes

\[ C_{1,1}(\tau, \tau'; J) = I^2 K V_2(J) \overline{V_2 C_{1,1}(\tau, \tau')} + \frac{1}{2} I^4 K (K - 1) V_4(J) \overline{c_{2,0}(\tau, \tau)} \overline{V_2 C_{1,1}(\tau, \tau')} + \frac{1}{2} I^4 K (K - 1) V_2(J) \overline{V_2 C_{1,1}(\tau; \tau') \overline{c_{2,0}(\tau', \tau')}} + \frac{1}{3!} I^4 K V_4(J) \overline{V_2 C_{3,1}(\tau, \tau, \tau'; \tau')} + \frac{1}{2} I^4 K V_2(J) \overline{V_2 C_{3,1}(\tau, \tau', \tau'; \tau')} \] (5.10)

Likewise the four-point function on the right-hand side obeys

\[ C_{3,1}(\tau_1, \tau_2, \tau_3; \tau_4; J) = I^4 K V_2^2(J) \overline{V_2 C_{3,1}(\tau_1, \tau_2, \tau_3; \tau_4) + I^4 K (K - 1) V_2^3(J) \overline{V_2 C_{1,1}(\tau_1; \tau_4) \overline{c_{2,0}(\tau_2, \tau_3)}} + \overline{V_2 C_{1,1}(\tau_2; \tau_4) \overline{c_{2,0}(\tau_1, \tau_3)}} + \overline{V_2 C_{1,1}(\tau_3; \tau_4) \overline{c_{2,0}(\tau_1, \tau_2)}}]. \] (5.11)

The average over \( J \) requires some attention because on the right-hand side of (5.10) and (5.11) the averages contain different vertices \( V_\delta(J) \). Nevertheless they can be evaluated successively. Multiplying (5.11) with \( V_2(J) \) and performing the \( J \)-average, \( \overline{V_2 C_{3,1}} \) can be expressed in terms of \( \overline{c_{2,0}} \) and \( \overline{V_2 C_{1,1}} \). Multiplying (5.11) with \( V_4(J) \), averaging and using the result for \( \overline{V_2 C_{3,1}}, \overline{V_4 C_{3,1}} \) is again written in terms of pair functions. This is inserted into (5.10) resulting in

\[ C_{1,1}(\tau; \tau'; J) = \left[ I^2 K V_2(J) + \frac{1}{2} I^4 K (K - 1) V_4(J) \overline{c_{2,0}(\tau, \tau)} + I^4 K V_2(J) \overline{V_2 C_{2,0}(\tau', \tau')} \right] \overline{V_2 C_{1,1}(\tau; \tau')} + \frac{1}{2} I^4 K (K - 1) V_2(J) \overline{V_2 C_{1,1}(\tau; \tau')} \overline{V_2 C_{1,1}(\tau; \tau')} \] (5.12)

This yields in lowest order

\[ \overline{V_2 C_{1,1}(\tau; \tau')} = I^2 K V_2 V_4 V_4 \overline{V_2 C_{1,1}(\tau; \tau')} \] (5.13)

Inserted into (5.12) a closed equation for \( \overline{V_2 C_{1,1}} \) is obtained

\[ \overline{V_2 C_{1,1}(\tau; \tau')} = \left\{ I^2 K \overline{V_2^2} + \frac{1}{2} I^4 K (K - 1) \overline{V_2 V_4} \overline{c_{2,0}(\tau, \tau)} + \frac{1}{2} I^4 K (K - 1) \right\} \overline{V_2 C_{1,1}(\tau; \tau')} + \overline{I^4 K V_2^2 V_4} \overline{V_2 C_{2,0}(\tau', \tau')} \] (5.14)

which is again of the form

\[ \overline{V_2 C_{1,1}(\tau)} = Y'(\tau) \overline{V_2 C_{1,1}(\tau)} \] (5.15)

assuming time translational invariance.
Pointing out the difference between \( Y(\tau), (5.9) \) and \( Y'(\tau) \) we write

\[
Y'(\tau) = Y(\tau) + \frac{1}{2} I^4 K (K - 1) \left\{ \frac{I^4 K (V_2^2 V_3 V_4 - V_2^2 V_2 V_4)}{1 - I^4 K V_2^2} \right\}
- \left[ 1 - I^2 K V_2^2 \right] V_2 V_4 \right\} C_{2,0}(0).
\]

(5.16)

This allows for the following non-trivial solutions:

(a) \( Y(\tau) = 1 \), \( Y'(\tau) \neq 1 : \overline{C_{2,0}(\tau)} \neq 0, \overline{C_{1,1}(\tau)} = 0. \)

(b) \( Y(\tau) \neq 1, \ Y'(\tau) = 1 : \overline{C_{2,0}(\tau)} = 0, \overline{C_{1,1}(\tau)} \neq 0. \)

(c) \( Y(\tau) = Y'(\tau) = 1 : \overline{C_{2,0}(\tau)} \neq 0, \overline{C_{1,1}(\tau)} \neq 0. \)

A solution of type (a) has been found for instance for the spherical version of the SK model [17]. It does not describe glassy behaviour since the response function vanishes on the long time scale, although the correlation function shows non-trivial properties. A solution of type (b) does actually not exist as shown later.

Only the last solution is supposed to be characteristic for a spin glass phase. Since \( I^2 K V_2^2 = 1 + O(\epsilon) \) the last term in (5.16) does not contribute in order \( \epsilon \) and only the first term in the bracket has to be taken into account. This means, however, that solution (c) requires

\[
\overline{V_2^2 V_2 V_4} = \overline{V_2^2 V_2 V_4}.
\]

(5.17)

With (3.21), (3.23) and (5.3)

\[
V_2(J) = I^{-1} \tanh(\beta J) \quad V_4(J) = -2\beta^2 I^{-1} \tanh(\beta J)(1 - \tanh^2(\beta J))
\]

(5.18)

and

\[
V_2^2(J) V_2^2(J') V_4(J') - V_2^2(J) V_2(J') V_2(J') V_4(J')
= 2\beta^2 I^{-6} \tanh^2(\beta J) \tanh^2(\beta J') (\tanh^2(\beta J) - \tanh^2(\beta J')).
\]

(5.19)

This shows that (5.17) can be fulfilled only with \( J = \pm I \) and that this kind of solution does not exist for more general \( P(J) \). This result is quite remarkable, indicating the special role of the binary distribution of couplings.

5.2. Next to leading order for \( J = \pm I \)

The correlation–response functions (4.8) have the symmetry

\[
C_{nm}(\cdots; J) = (-1)^n C_{nm}(\cdots; -J).
\]

(5.20)

For binary couplings \( J = \pm I \) the bond average therefore becomes almost trivial. For simplicity we set \( h = 0 \) in the following and assume time translational invariance on the long time scale. The strategy to evaluate contributions of higher orders in \( \epsilon \) is analogous to what has been described in section 5.1. Evaluating \( C_{2,0} \) in order \( \epsilon^3 \) requires to compute \( C_{4,0} \) and \( C_{6,0} \) in the same order and to break them down successively to products of pair functions \( C_{2,0} \). In addition, in this order an internal integration shows up taking into account the third line on the right-hand side of (5.1) leading to a contribution of the form

\[
C_{2,0}(\tau, \tau') = \cdots + V_2 \int d\sigma C_{3,1}(\tau, \tau', \sigma; \sigma^*).
\]

(5.21)

The four-point function \( C_{3,1} \) can again be broken down to products of pair functions according to (5.11).
Using the notation
\[ \tilde{q}(\tau - \tau') = C_{2,0}(\tau, \tau'; I), \quad \tilde{r}(\tau - \tau') = C_{1,1}(\tau, \tau'; I), \quad V_n = V_n(I) \] (5.22)
and collecting all contributions up to order \( \epsilon^3 \) one obtains
\[
\tilde{q}(\tau) = I^2 K V_2^4 \tilde{g}(\tau) + I^4 K (K - 1) \frac{V_2 V_4}{1 - I^4 K V_2^2} \tilde{q}(0) \tilde{q}(\tau)
+ I^6 K (K - 1) \frac{1}{1 - I^4 K V_2^6} \left\{ K - 2 + 3I^4 K (K - 1) \frac{V_4^2 V_2^4}{1 - I^4 K V_2^2} \right\} \tilde{q}^2(0)
\times \left\{ \frac{1}{4} (V_2 V_6 + V_4^2) + 3I^4 K V_2^8 \frac{V_4^4 V_2^4}{1 - I^4 K V_2^2} \right\} \tilde{q}^2(\tau) \tilde{q}(\tau)
+ I^4 K (K - 1) \frac{V_2^3}{1 - I^4 K V_2^2} \int d\sigma [\tilde{r}(\tau + \sigma) \tilde{q}(\sigma) + \tilde{q}(\tau - \sigma) \tilde{r}(\sigma)].
\] (5.23)

The corresponding calculation for the response function yields
\[
\tilde{r}(\tau) = I^2 K V_2^4 \tilde{r}(\tau) + I^4 K (K - 1) \frac{V_2 V_4}{1 - I^4 K V_2^2} \tilde{q}(0) \tilde{r}(\tau)
+ I^6 K (K - 1) \frac{1}{1 - I^4 K V_2^6} \left\{ K - 2 + 3I^4 K (K - 1) \frac{V_4^2 V_2^4}{1 - I^4 K V_2^2} \right\} \tilde{q}^2(0)
\times \left\{ \frac{1}{4} (V_2 V_6 + V_4^2) + 3I^4 K V_2^8 \frac{V_4^4 V_2^4}{1 - I^4 K V_2^2} \right\} \tilde{q}^2(\tau) \tilde{r}(\tau)
+ I^4 K (K - 1) \frac{V_2^3}{1 - I^4 K V_2^2} \int d\sigma [\tilde{r}(\tau + \sigma) \tilde{r}(\sigma) + \tilde{q}(\tau - \sigma) \tilde{r}(\sigma)].
\] (5.24)

5.3. Ultrametric time parametrization

The low-temperature phase of the SK model is characterized by a hierarchy of long time scales ranging from \( t^* \) to some longest time scale \( t_w \). Eventually the limit \( t_w \to \infty \) is taken. In order to keep track of the long time scales it is convenient to introduce the parametrization [4]
\[
t = t_w^{1 - x(t)}, \quad x(t) = 1 - \frac{\ln(t)}{\ln(t_w)},
\] (5.25)
and write
\[
\tilde{q}(\tau) = \tilde{Q}(x(t^* \tau)), \quad \tilde{r}(\tau) = -t^* \dot{x}(t^* \tau) \tilde{R}(x(t^* \tau)).
\] (5.26)

With the above definition
\[
x(t + t') = 1 - \frac{\ln \left( t_w^{1 - x(t)} + t_w^{1 - x(t')} \right)}{\ln (t_w)} = x(t) - \frac{\ln \left( 1 + t_w^{x(t) - x(t')} \right)}{\ln (t_w)}
\] (5.27)
and for \( t > t' \) and with it \( x(t) < x(t') \)
\[
x(t + t') = x(t) - \frac{\ln \left( 1 + t_w^{x(t) - x(t')} \right)}{\ln (t_w)} \approx x(t) - \frac{t_w^{x(t) - x(t')}}{\ln (t_w)} \approx x(t).
\] (5.28)
The corresponding result for \( t < t' \) is \( x(t + t') \approx x(t') \). This yields the ultrametric relation
\[
x(t + t') = x(\max(t, t')).
\] (5.29)

With it the integrals in (5.23) and (5.24) become
\[
\int_0^\tau d\sigma \tilde{r}(\tau - \sigma)\tilde{q}(\sigma) + \int_0^\infty d\sigma [\tilde{r}(\tau + \sigma)\tilde{q}(\sigma) + \tilde{q}(\tau + \sigma)\tilde{r}(\sigma)]
\to 2\tilde{Q}(x) \int_0^1 dx'\tilde{R}(x') + 2 \int_0^x dx'\tilde{Q}(x')\tilde{R}(x')
\] (5.30)

and
\[
\int_0^\tau d\sigma \tilde{r}(\tau - \sigma)\tilde{r}(\sigma) \to -2r^*\dot{x}(t + \tau)\tilde{R}(x) \int_0^1 dx'\tilde{R}(x').
\] (5.31)

This allows us to write (5.24) in the form
\[
\tilde{R}(x) = \tilde{Y}(x)\tilde{R}(x)
\] (5.32)

with
\[
\tilde{Y}(x) = Y_0 + Y_1 Y_1(1) + \frac{1}{2} Y_2 Y_2^2(1) + \frac{1}{2} Y_2 Y_2^2(x) + Y_r \int_1^x dx'\tilde{R}(x')
\] (5.33)

and the coefficients
\[
Y_0 = I^2K V_2^2,
\] (5.34)
\[
Y_1 = I^3(K - 1) \frac{V_2 V_4}{1 - I^4K V_2^4},
\] (5.35)
\[
Y_2 = I^5(K - 1) \frac{1}{1 - I^4K V_2^4} \left\{ K - 2 + 3I^4K(K - 1) \frac{V_2^4}{1 - I^4K V_2^4} \right\}
\times \left\{ \frac{1}{2}(V_2 V_6 + V_2^2) + 6I^4K \frac{V_2^4 V_2^2}{1 - I^4K V_2^4} \right\},
\] (5.36)
\[
Y_{2v} = I^6K(K - 1) \frac{1}{1 - I^4K V_2^4} \left\{ K - 2 + 3I^4K(K - 1) \frac{V_2^4}{1 - I^4K V_2^4} \right\}
\times \left\{ V_2^2 + 2I^3K \frac{V_2^4 V_2^2}{1 - I^4K V_2^4} \right\},
\] (5.37)

and
\[
Y_r = 2I^4K(K - 1) \frac{V_2^3}{1 - I^4K V_2^4}.
\] (5.38)

The correlation function (5.23) becomes
\[
\tilde{Q}(x) = \left\{ Y_0 + Y_1 Y_1(1) + \frac{1}{2} Y_2 Y_2^2(1) + \frac{1}{3!} Y_{2v} Y_2^2(x) \right\} \tilde{Q}(x)
\] \[+ Y_r \left\{ \tilde{Q}(x) \int_0^1 dx'\tilde{R}(x') + \int_0^x dx'\tilde{Q}(x')\tilde{R}(x') \right\}.
\] (5.39)

Differentiation with respect to \( x \) results in
\[
\partial_x \tilde{Q}(x) = \tilde{Y}(x)\partial_x \tilde{Q}(x)
\] (5.40)

with the same \( \tilde{Y}(x) \) as above. This allows a non-trivial solution for \( \tilde{Y}(x) = 1 \).
5.4. Results in next to leading order

In particular for \( x = 1 \) and in the second order we get

\[
\bar{Q}(1) = q_{EA} = 1 - \frac{Y_0}{Y_1} - \frac{1}{2} \frac{(Y_2 + Y_{2x})(1 - Y_0)^2}{Y_1^3}.
\] (5.41)

Lowering the temperature a non-zero Edwards–Anderson order parameter \( q_{EA} > 0 \) shows up first at a critical temperature where

\[
Y_0 = I^2KV_2^2 = K \tanh^2(\beta I) = 1.
\] (5.42)

This determines the critical temperature

\[
T_c = \frac{I}{\text{artanh}(1/\sqrt{K})},
\] (5.43)

and choosing

\[
I = \text{artanh}(1/\sqrt{K})
\] (5.44)

the critical temperature is \( T_c = 1 \).

Expanding the vertices and coefficients (5.36)–(5.38) in powers of \( \epsilon = 1 - T \)

\[
Y_0 = 1 + 2I \frac{K - 1}{\sqrt{K}} \epsilon + I^2(K - 3I + 2\sqrt{K}) \frac{K - 1}{K} \epsilon^2,
\]

\[
Y_1 = -2I^2(K - 1) - 4I^2(I\sqrt{K} + 1)(K - 1) \epsilon,
\]

\[
Y_2 = 10I^4(K^2 - 1), \quad Y_{2x} = 4I^4(K^2 - 1), \quad Y_r = 2I\sqrt{K}
\] (5.45)

and (5.41) becomes

\[
q_{EA} = \frac{1}{I\sqrt{K}} \epsilon + \left[ 2 \frac{K - 1}{K} - \frac{1}{I\sqrt{K}} \right] \epsilon^2.
\] (5.46)

Differentiating (5.33) with respect to \( x \) gives

\[
\bar{R}(x) = \frac{Y_{2x}}{Y_r} \bar{Q}(x) \partial_x \bar{Q}(x) = 2I^3 \frac{K^2 - 1}{\sqrt{K}} \bar{Q}(x) \partial_x \bar{Q}(x).
\] (5.47)

This verifies that \( \bar{R}(x) \sim \epsilon^2 \).

In equilibrium a fluctuation–dissipation theorem (FDT) holds, i.e. \( r(t) = -\beta \delta_t q(t) \). As pointed out earlier this is also expected to hold in the glassy state for \( t < t^* \). On the long time scale, however, the FDT is violated. We may introduce a measure \( X \) for the violation of the FDT

\[
r(\tau) = -\beta X(\tau) \partial_\tau \bar{q}(\tau), \quad \bar{R}(x) = \beta \bar{X}(\bar{Q}(x)) \partial_x \bar{Q}(x)
\] (5.48)

with

\[
\bar{X}(q) = \frac{Y_{2x}}{Y_r} q = 2I^3 \frac{K^2 - 1}{\sqrt{K}} q \quad \text{for} \quad 0 < q < q_{EA}.
\] (5.49)

in lowest order.

The fact that \( \bar{X} \) depends on \( \bar{Q} \) only holds in higher orders as well. For \( t < t^* \) the correlation function obeys \( q(t) > q_{EA} \) and we can set \( \bar{X}(q) = 1 \) for \( t < t^* \). This quantity is related to Parisi’s overlap distribution function \( P(q) = \partial_q \bar{X}(q) \).

Equation (5.47) does not fix the actual form of \( \bar{Q}(x) \) or \( \bar{R}(x) \). Using the so-called Parisi gauge [18], \( \beta X(\bar{Q}(x)) = x \),

\[
\bar{Q}(x) = \frac{Y_r}{Y_{2x}} x \quad \text{for} \quad x < x^*
\] (5.50)
with
\[ x^* = \frac{Y_{2k}}{Y_r} q_{\text{EA}} \] (5.51)
and \( t^* = t_{w}^* \).

For \( K \to \infty \) the known results for the SK model are recovered.

6. Stability analysis

6.1. Solutions for general distribution of couplings

For general distributions \( P(J) \) the difference \( \bar{Y}'(x) - \bar{Y}(x) \), given in (5.16) is non-zero. In order \( \varepsilon \) the last term \( \sim 1 - I_2 V_2^2 V_4^2 \) does not contribute. The remaining term contains
\[ V_2^2 V_4^2 V_4 - V_2^4 V_2 V_4 \geq 0, \] (6.1)
which is positive according to (5.19) unless \( J = \pm I \). Accordingly \( \bar{Y}'(x) > \bar{Y}(x) \) for general distributions \( P(J) \).

This still allows for solution (a) of section 5.1 with \( \bar{R}(x) = 0 \). The last term in (5.33) vanishes and \( \bar{Q}(x) = q_{\text{EA}} \), given in (5.46). This solution does not show any time dependence on long time scales. This property is shared by correlation functions of higher order, \( C_n(\tau_1 \cdots \tau_n) = C_n \). Correlation–response functions \( C_{n,m}(\cdots) \) with \( m > 0 \) vanish.

Reconstructing the full distribution \( Q(\{\hat{\kappa},\kappa\}) \) from (5.1) one finds the replica symmetric solution [13]
\[ Q(\{\hat{\kappa},\kappa\}) = \int d\hat{\kappa} P(\hat{\kappa}) e^{\int d\sigma \hat{\kappa}(\sigma)[\sigma - \hat{\kappa}]} \] (6.2)
with
\[ P(\kappa) = \prod_{i=1}^{K} \int d\kappa_i P(\kappa_i) \delta \left( h + \sum_{i} u(J_i, \kappa_i) - \kappa \right) \] (6.3)

In next to leading order
\[ \bar{Y}(x) = Y_0 + Y_1 q_{\text{EA}} + \left[ \frac{1}{2} Y_2 + \frac{1}{3!} Y_{2s} \right] q_{\text{EA}}^2 \] (6.4)
which yields
\[ q_{\text{EA}} = \frac{1}{I\sqrt{K}} \varepsilon - \left[ \frac{13}{3} \frac{K + 1}{K} \right] \frac{1}{I\sqrt{K}} \varepsilon^2. \] (6.5)

As shown below this solution is unstable.

Solution (b) of section 5.1 requires \( \bar{Q}(x) = 0 \) and with (5.33)
\[ \bar{R}(x) = Y_0 + Y_r \int_{x}^{1} dx' \bar{R}(x') = 1. \] (6.6)

This has, however, no solution with \( \bar{R}(x) \neq 0 \).

6.2. Stability analysis

Various criteria can be used to test the stability of a given solution. One may, for instance, ask whether the correlation–response functions on the short time scale approach the values determined by the dynamics on the long time scale. For example the pair correlation function
\( q(t) = \langle \sigma(t) \sigma(0) \rangle \) should approach \( q_{EA} \) for \( t \approx t^* \). This decay is expected to be algebraic, but a corresponding analysis is beyond the scope of this paper.

The general procedure used in statics as well as in the present formulation of dynamics consists in iteratively connecting subtrees to a new tree and expressing the properties at the new vertex by properties at the base vertices of the subtrees searching for fixed points of this mapping. A necessary condition for stability is the decay of small perturbations under this mapping. Equations (5.23) and (5.24) are actually such mappings.

For \( T > T_c \) the trivial solution \( \bar{Q}(x) = 0 \) and \( \bar{R}(x) = 0 \) is expected to be valid. A small perturbation \( \delta \bar{Q}_i(x) \) on site \( i \) is mapped onto

\[
\delta \bar{Q}_o(x) = I^2 K V_o^2 \delta \bar{Q}_i(x)
\]  

(6.7)

at site \( o \), and a corresponding mapping for \( \bar{R}_0(x) \). Since \( I^2 K V_o^2 < 1 \) for \( T > T_c \) the trivial solution fulfills this stability criterion. By the same token this solution is unstable for \( T < T_c \).

The stability criterion for solution (a) reads with (6.4) and \( \hat{Y}(x) = 1 \)

\[
\delta \bar{Q}_o(x) = [1 + \frac{1}{3} Y_2, q_{EA}^2] \delta \bar{Q}_i(x) > \delta \bar{Q}_i(x)
\]  

(6.8)

and the same for \( \delta \bar{R}(x) \). This solution is therefore unstable for \( T < T_c \) as well.

Finally solution (c) for \( J = \pm J \) is investigated. We may test the stability with respect to a perturbation \( \delta \bar{Q}(x) \) and \( \delta \bar{R}(x) \). This yields with \( \hat{Y}(x) = 1 \), \( \delta \bar{Q}_o(x) = \delta \bar{Q}_i(x) \) and \( \delta \bar{R}_o(x) = \delta \bar{R}_i(x) \). This means that this solution is marginal with respect to perturbations of this kind. Investigating a perturbation \( \delta \bar{R}(x) \) with \( \delta \bar{Q}_i(x) = 0 \) one obtains

\[
\delta \bar{R}_o(x) = \int_0^1 dx' \delta \bar{R}_i(x') \bar{R}(x).
\]  

(6.9)

Considering the second contribution a perturbation at some value \( x' \) creates fluctuations at \( x < x' \) only. This means that the perturbation at \( x' \) is not enhanced due to this term and this solution is marginal with respect to a perturbation of this kind as well. Such a marginal stability criterion is actually expected because of the reparametrization invariance mentioned at the end of section 5.4.

7. Discussion

In this paper, we have shown that the long time dynamics of an Ising spin glass with binary couplings \( J = \pm J \) on the Bethe lattice is of the form known from the Sherrington–Kirkpatrick model. This is remarkable in so far as the interaction is restricted to nearest neighbours. In contrast to a lattice in finite dimensions, however, the typical size of closed loops scales with \( \ln(N) \) for a system of \( N \) sites.

The second main result of this paper, the failure of a corresponding solution for general distributions \( P(J) \), is unexpected, and the nature of the low-temperature phase in this case is not known. A similar breakdown of a replica or cavity method calculation for general \( P(J) \) cannot be excluded since the relevant equations have been evaluated for binary coupling only [13]. It might be of interest to perform an expansion of the relevant equations around \( T_c \), possibly within the extended replica scheme proposed by de Dominicis et al. [18]. In this scheme two-order parameter functions corresponding to \( \bar{Q}(x) \) and \( \bar{R}(x) \) are used.

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