

One-dimensional disordered Ising models by replica and cavity methods

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Using a formalism based on the spectral decomposition of the replicated transfer matrix for disordered Ising models, we obtain several results that apply both to isolated one-dimensional systems and to locally tree-like graph and factor graph (p-spin) ensembles. We present exact analytical expressions, which can be efficiently approximated numerically, for many types of correlation functions and for the average free energies of open and closed finite chains. All the results achieved, with the exception of those involving closed chains, are then rigorously derived without replicas, using a probabilistic approach with the same flavour of cavity method.

I. INTRODUCTION

The investigation of one-dimensional disordered Ising systems has a relatively long tradition among the scholars^{1–3}, since they are analytically tractable and display a non-trivial phenomenology. Moreover some of their properties also apply to diluted mean field models, where with high probability there is a unique finite path connecting any two spins and where the leading finite-size correction to the free energy is given by short and isolated loops⁴.

A pure one-dimensional system can be easily analyzed with the transfer matrix formalism. The spectral properties of this 2×2 matrix completely characterize the model, since the partition function of a chain of length ℓ is simply obtained from its ℓ -th power. This operation becomes effortless once the transfer matrix is expressed as a linear combination of orthonormal left and right eigenvectors.

The situation is much more complicated in disordered systems, where the (random) partition function is given by the product of non-commuting random matrices⁵. This situation arises when random interactions or random magnetic fields are present, as in spin glasses or random field Ising models (RFIM) respectively.

As long as the sources of disorder are independently distributed, one can express the integer moments of the partition function through the power of the $2^n \times 2^n$ transfer matrix of a system of n replicated spins. Then, as usual with replica calculations, the analytic continuation to small values of n is performed to compute relevant physical quantities. This approach was pioneered by Weigt and Monasson². Taking advantage of the symmetry group of the replicated transfer matrix (RTM) at integer n , they were able to associate the decay rate of different connected correlation functions to different eigenvalues of the RTM. They also computed the free energy of an infinite chain, rederiving a result earlier obtained by means of a recursion of a distributional equation¹.

In recent years the spectral formalism of the RTM has been the subject of further investigations⁶ and has been used to determine the DAT line in many mean field spin glasses⁷. Nevertheless its capabilities have not been pushed to full extent: a manageable decomposition in terms of left and right eigenvectors and a in-depth analysis of the $n \downarrow 0$ limit is lacking; little is known about the behaviour at finite length of the free energies and about the differences between open and closed chains; a complete characterization of the connected correlation functions is lacking, and the analysis of thermally disconnected functions, relevant to the RFIM transition, has not been taken in consideration yet. In the present paper we address all these issues, giving exact analytical expressions valid at any finite length of the aforementioned quantities.

Inspired by some recent observations linking the RTM formalism with a cavity approach⁸, we give an alternative derivation of the expressions for the connected and disconnected correlation functions and for the free energy of open chains, using a rigorous probabilistic analysis. This approach avoids the use of replicas and of uncontrolled analytic continuations and relies on iterative distributional equations involving some physically sensible observables. The integral operators arising with this method are the same we encountered in the RTM analysis, therefore closing the bridge between the two formalism.

The paper is organized as follows: in Section II we define the model we are considering and expose the main results of this paper; in Section III we develop all the spectral formalism of the RTM and we apply it in Section IV to the computation of free energies and correlation functions. Most of the results obtained with the RTM are the rederived with a purely probabilistic approach in Section V.

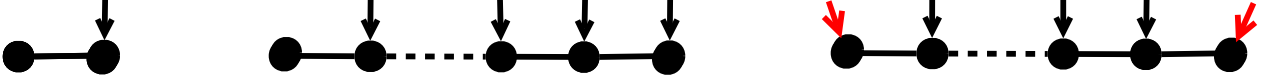


Fig. 1: Pictorial representation of the matrix T_n (left), its powers T_n^ℓ (center) and the matrix $\tilde{T}_n^{(\ell)}$ (right).

II. DEFINITIONS AND MAIN RESULTS

In this paper we consider one-dimensional Ising spin system with i.i.d. random fields and couplings, e.g. an isolated chain or a chain embedded in a locally tree-like graph, therefore described by the product of uncorrelated 2×2 random transfer matrices M_i defined by

$$M_i(\sigma_{i+1}, \sigma_i) = e^{\beta J_i \sigma_{i+1} \sigma_i + h_i \sigma_i}. \quad (1)$$

The partition function of a closed chain of length ℓ is then a random variable given by

$$Z_{\ell, c} = \text{Tr} \prod_{i=0}^{\ell-1} M_i. \quad (2)$$

A powerful technique to compute the statistical properties of this kind of objects is the well known replica method⁹. As we shall see, as long as the system stays in a replica symmetric phase, its statistical properties are encoded in the (replica symmetric) replicated transfer matrix T_n , the $2^n \times 2^n$ matrix defined by

$$T_n(\sigma, \tau) = \mathbb{E}_{J, h} e^{\beta J \sum_{a=1}^n \sigma^a \tau^a + \beta h \sum_{a=1}^n \tau^a}. \quad (3)$$

Here and in the following we denote with σ the vector $(\sigma^1, \dots, \sigma^n)$, with the n replicated spins σ^a taking values in $Z_2 = \{-1, 1\}$. A similar definition holds for τ . As usual in the replica method⁹ we shall work at integer value of n and perform the analytic continuation for $n \downarrow 0$ at the end of the computations. We shall assume in the following that the field h is an arbitrary distributed external random field, if we are considering an isolated chain, or, if we are considering a chain embedded in a locally tree-like graph, to be a random cavity field conditioned to act on a spin that is already connected to two other spins (its neighbours on the chain). See Figure 1 for a representation of T_n and its powers T_n^ℓ .

A first spectral analysis of T_n was conducted by Weigt and Monasson². Following their lead we take advantage of the replica index permutation symmetry of T_n to choose an appropriate bases to express its right eigenvectors. There are $n+1$ non-equivalent irreducible representation of the permutation group, which can be glued together to form the sectors $D^{(q)}$, $q = 0, 1, \dots, \lfloor \frac{n}{2} \rfloor$, partitioning $Z_2^{\otimes n}$. In the following, with some abuse of notation, we will denote with $D^{(q)}$ the set of eigenvalues of T_n with eigenvector in that sector. The eigenvectors of T_n in the sector $D^{(q)}$ can be parametrized by functions $g_q^\lambda(u)$ that, in the limit $n \downarrow 0$, satisfy the eigenvalue equation

$$\lambda g_q^\lambda(u) = \mathbb{E}_{J, h} \int dv g_q^\lambda(v) \delta(u - \hat{u}(J, v + h)) \left(\frac{1}{\beta} \frac{\partial \hat{u}}{\partial v} \right)^q, \quad (4)$$

where $\hat{u}(J, h) = \frac{1}{\beta} \text{atanh}(\tanh(\beta J) \tanh(\beta h))$ is the cavity iteration rule.

In this paper we extend the analysis of the spectral properties of T_n to achieve a complete description of the $n \downarrow 0$ limit, derive exact expressions for correlation functions and free energies of chains. Since T_n is the product of two non-singular symmetric matrices, it possess a complete orthonormal (in the left-right sense) basis of left and right eigenvectors with real eigenvalue. The left eigenvector corresponding to a certain right ψ_R is simply $\mathbb{E}_h e^{\beta h \sum_a \sigma^a} \psi_R(\sigma) \equiv \rho_h(\sigma) \psi_R(\sigma)$. Therefore the spectral decomposition of T_n into the subspaces $D^{(q)}$ is given by

$$T_n(\sigma, \tau) = \sum_{q=0}^{\lfloor \frac{n}{2} \rfloor} \sum_{\lambda \in D^{(q)}} \lambda \rho_q^\lambda(\sigma) \rho_h(\tau) \rho_q^\lambda(\tau) \sum_{\substack{a_1 < \dots < a_q \\ b_1 < \dots < b_q}} Q_{a_1 \dots a_q; b_1 \dots b_q} \sigma^{a_1} \dots \sigma^{a_q} \tau^{b_1} \dots \tau^{b_q}. \quad (5)$$

Here we have denoted with $\rho_q^\lambda(\tau)$ the replica symmetric part of the eigenvector in the sector $D^{(q)}$ with eigenvalue λ . The second sum is over all the eigenvalues of T_n in the sector $D^{(q)}$, given in the $n \downarrow 0$ limit by the solutions of Eq.

(4). The coefficients $Q_{a_1 \dots a_q; b_1 \dots b_q}$ have simple algebraic expressions in each sector (see Eqs. (27) and (28)) and are invariant under permutations of any of their two sets of indices. While a different left-right decomposition of T_n has already been attempted⁶, an unfortunate choice in the parametrization of the eigenvectors in terms of function of two variable led to an unmanageable formalism. Thanks to the spectral representation (5) we can easily take the powers of T_n and contract the matrix with the quantities we want to average. In Section III we derive Eqs. (4) and (5), and discuss the non-trivial aspects of the small n limit.

One of the applications of the spectral formalism is the computation of the average free energy of open and closed chains, as exposed in Section IV A. Recently it has been shown⁴ that the first finite size correction to thermodynamic free energy of systems on diluted graphs can be expressed as a linear combination of the free energies of closed and open chains. It has also been argued^{4,10} that a perturbative expansion around the Bethe approximation towards finite dimensional lattices, shall account for the presence of loops (closed chains) and will contain the free energies and the correlation function of one-dimensional objects, motivating the importance of exact and easily approximable expressions for their free energies.

Taking the trace of T_n^ℓ and performing the $n \downarrow 0$ limit one obtains the average free energy of a closed chain of size ℓ :

$$-\beta f_\ell^c = -\beta \ell f_0 + \sum_{\lambda \in D^{(1)}} \Delta_\lambda \ell \lambda^{\ell-1} + \sum_{q=1}^{\infty} \hat{d}_q \sum_{\lambda \in D^{(q)}} \lambda^\ell. \quad (6)$$

The terms Δ_λ , due to a correspondence between the eigenvalues of the sectors $D^{(0)}$ and $D^{(1)}$ at $n = 0$, are expressed in Eq. (48). The coefficients \hat{d}_q are the analytic continuation of the degeneracies of the eigenvalues, and are given in Eq. (41). We note that the correction to the intensive free energy f_0 (expressed in Eq. (43)) is given by a linear combination of exponential and ℓ times exponential terms. The decaying part of f_ℓ^c is dominated by the largest eigenvalue among the various sectors.

In the computation of the free energy of an open chain, we allow for the incoming fields at the extremities of the chain to be distributed differently from the fields h acting on the internal spins, and denote them by \tilde{h} . This is in fact what happens in general when considering an open chain embedded in a sparse graph. The expression we derived for the average free energy of an open chain of length ℓ is

$$\begin{aligned} -\beta f_\ell^o &= -\ell \beta f_0 + \mathbb{E}_{\tilde{h}} \int du P(u) 2 \log \cosh(\beta(u + \tilde{h})) - \mathbb{E}_h \int dudv P(u)P(v) \log \cosh(\beta(u + v + h)) \\ &+ \log 2 + \sum_{\lambda \in D^{(1)}} a_{\lambda,0}^2 \lambda^\ell, \end{aligned} \quad (7)$$

where $P(u)$ is the distribution of cavity messages along the chain and the coefficients $a_{\lambda,0}$ are related to the left eigenvectors of the sector $D^{(0)}$ and given in Eq. (59).

Another achievement of this paper, obtained through the spectral decomposition of the RTM, is the expression of the connected correlation functions of two spin at distance ℓ , in a form that is both analytically exact and easy to approximate numerically with high precision. In Section IV B we derive the formula

$$\overline{\langle \sigma_0 \sigma_\ell \rangle_c^q} = \sum_{\lambda \in D^{(k)}} a_{\lambda,q}^2 \lambda^\ell, \quad (8)$$

where $a_{\lambda,k}$ can be computed through the eigenfunction g_q^λ using Eq. (64). We indicate with $\bar{\bullet}$ the average over all kinds of disorder in the model considered. For Ising model on sparse random graphs with mean residual degree z , the susceptibility $\chi_q = \sum_{i < j} \frac{1}{N} \mathbb{E} \langle \sigma_i \sigma_j \rangle_c^q$ diverges when the greatest eigenvalue of $D^{(q)}$ reaches the value $1/z$. Therefore the sectors $D^{(1)}$ and $D^{(2)}$ are the relevant ones to the ferromagnetic and the spin-glass transitions respectively (see Figures 2 and 3).

The computation of the thermally disconnected correlation function, $\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} - \overline{\langle \sigma_0 \rangle} \overline{\langle \sigma_\ell \rangle}$, particularly relevant to the RFIM¹¹, requires a careful treatment of the analytic continuation to $n = 0$. The final expression we obtained, Eq. (71), is not a linear combination of terms involving only one eigenvalue, as in the previous formulas. The leading term for large ℓ is easily extracted though: let λ_1 be the greatest eigenvalue of the sector $D^{(1)}$, then

$$\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} - \overline{\langle \sigma_0 \rangle} \overline{\langle \sigma_\ell \rangle} = \Delta_{\lambda_1} a_{\lambda_1,1}^2 \ell \lambda_1^{\ell-1} + O(\lambda_1^\ell) \quad \text{for } \ell \rightarrow +\infty, \quad (9)$$

with Δ_λ and $a_{\lambda,1}$ given in Eq. (48) and Eq. (64) respectively. Therefore, on one-dimensional chains and sparse graphs, the susceptibility corresponding to the thermally disconnected correlation function present the characteristic double pole behaviour near the transition point, whose prefactor can also be computed by Eq. (9).

The expressions we found for free energies of chains, Eqs. (6) and (7), and the correlation function Eqs. (8) and (9), are exact for every value of the length ℓ of the chain but involve the computation of infinitely many terms. Fortunately it turns out from our numerical simulations that the spectrum of the integral operator in Eq. (4) is discrete and the eigenvalues are well spaced. Therefore considering only the first few highest eigenvalues one obtains very good approximations already at small values of ℓ . They can be computed numerically, discretizing the kernel of the integral operator of Eq. (4) and directly computing the eigenvalues of the associated matrix. Moreover the leading eigenvector and eigenvalue of each sector can be efficiently selected with multiple applications of the discretized operator on an arbitrarily chosen vector (as it was done to obtain Figures 2 and 3).

All the results we obtained using the replicated partition function formalism, with the noticeable exception of the formula for the average free energies of closed chains Eq. (6), can be recovered using a purely probabilistic approach in the same spirit of the usual cavity method^{9,12}.

In Section V A we devise two alternative probabilistic derivation for the average free energies of open chains. The first is based on a recursive equation involving the moments of the partition function, which leads to an expression for the moment of the random partition function Z_ℓ^n of an asymmetric open chain in terms of the left and right eigenvector of an integral operator we also encountered in the RTM formalism:

$$\overline{Z_\ell^n(u; x)} = \sum_{\lambda \in D^{(0)}} \lambda^\ell(n) g_0^\lambda(u; n) S_0^\lambda(x; n) [2 \cosh(\beta x)]^n. \quad (10)$$

Here n is not related to the number of replicas, since replicas are not present in this approach, but is an arbitrarily chosen real positive number. The other method presented in Section V A the iteration of the average free energy itself during the construction of the chain, which requires to keep track of the message of u_ℓ the cavity message propagating through a chain at distance ℓ from one of the extremities, at each iteration. The two approaches are deeply related and obviously lead to the same result.

Crucial to the probabilistic computation of the connected correlation functions, as has been noted recently⁸ is the random variable X_ℓ defined by $X_\ell \equiv \frac{\partial u_\ell}{\partial H_0}$, where H_0 field acting on the same extremity. It turns out that the connected correlation function of Eq. (8) is encoded in the q -the moments of the joint law of u_ℓ and X_ℓ at fixed u_ℓ . This object, the function

$$G_q^{(\ell)}(u) = \int dX P_\ell(u, X) X^q, \quad (11)$$

obeys a recursion rule, Eq. (87), containing the integral operator of Eq. (4). Expressing $G_q^{(\ell)}(u)$ in the basis of the eigenvalues of $D^{(q)}$ leads then straightforwardly to the expression (8) we obtained using replicas. Moreover in Section V B a more general result is presented in Eq. (90).

The thermally disconnected correlation function $\overline{\sigma_0 \sigma_\ell}_c$ is computed in Section V C, using some results we obtained for the connected correlation function and for the moment of the partition function of an open chain, thanks to the relation

$$\frac{\partial}{\partial H_0} \frac{\partial}{\partial H_\ell} Z_{\ell,o}^n = n \langle \sigma_0 \sigma_\ell \rangle_c Z_{\ell,o}^n + n^2 \langle \sigma_0 \rangle \langle \sigma_\ell \rangle Z_{\ell,o}^n. \quad (12)$$

An alternative approach, technically more difficult, outlined in Section V C involves the resolution of an iterative equation for the function $R^{(\ell)}(u) = \delta(u - u_\ell) \langle \sigma_0 \rangle^{(\ell)}$, which takes into account the shift in the magnetization of the first spin during the growth of the chain.

In the following Sections we fill-in all the technical details associated to the previous claims.

III. SPECTRAL DECOMPOSITION

We present an in-depth tractation of the spectral theory of the replica symmetric RTM. In Section III A we discuss the spectral decomposition of the matrix for integer values of the number of replicas n . We introduce in Section III B an integral representations of the eigenvectors, in order to discuss the main features of the analytic continuation to small values of n in Section III C. In Section III D we discuss some technicalities related to a peculiar aspect of the $n \downarrow 0$ limit, the degeneracy between the Longitudinal and the Anomalous sectors.

A. The Permutation Group

The $2^n \times 2^n$ matrix T_n defined by Eq. (3) is invariant under the action of the group of permutations among the replicated spins: for each permutation π acting on the n spin, we have the equivalence $T_n(\pi(\sigma), \pi(\tau)) = T_n(\sigma, \tau)$. This

symmetry allows us to block-diagonalize T_n according to the irreducible representations of the permutation group. This idea has been first introduced by Weigt and Monasson² in order to compute the eigenvalue spectrum of T_n .

For the sake of completeness we now review Weigt and Monasson's method, then we extend it further, in order to achieve the decomposition of the transfer matrix in terms of left and right eigenvectors. The replicated space is $Z_2^{\otimes n}$. Let's call Δ_m , with $m = 0, \dots, n$, the subspace of configurations having exactly m spins up. These subspaces are clearly invariant under any permutation of the replicas, therefore we can consider the representation of the permutation group in the $n + 1$ subspaces Δ_m and look for the irreducible ones. The complete decomposition of Δ_m into irreducible subspaces $D^{(m,q)}$ has been done by Wigner¹³. It reads:

$$\begin{aligned} \Delta_0 &= D^{(0,0)} , \\ \Delta_1 &= D^{(1,0)} \oplus D^{(1,1)} , \\ &\dots \\ \Delta_m &= D^{(m,0)} \oplus \dots \oplus D^{(m, \min(m, n-m))} , \\ &\dots \\ \Delta_{n-1} &= D^{(n-1,0)} \oplus D^{(n-1,1)} , \\ \Delta_n &= D^{(n,0)} . \end{aligned} \tag{13}$$

Representations $D^{(m,q)}$, at fixed q , are isomorphic and have dimension

$$d_q \equiv \dim \left(D^{(m,q)} \right) = \binom{n}{q} - \binom{n}{q-1} \quad q = 0, \dots, \lfloor n/2 \rfloor , \tag{14}$$

where $\lfloor x \rfloor$ is the smallest integer part of x . Notice that by definition $d_0 = 0$. As we have $(n + 1 - 2q)$ subspaces $D^{(m,q)}$, the q -sector of our matrix T_n will contain $(n + 1 - 2q)$ eigenvalues with degeneracy d_q . One can check that $\sum_{q=0}^{\lfloor n/2 \rfloor} d_q (n + 1 - 2q) = 2^n$.

A vector of the space $D^{(m,q)}$ can be constructed using Young tableaux⁷, and has the form

$$|m, q\rangle = (|+\rangle|-\rangle - |-\rangle|+\rangle)^q \text{SYM} (|+\rangle^{m-q} |-\rangle^{n-m-q}) , \tag{15}$$

where the operation SYM means a complete symmetrization with respect to the $n - 2q$ last entries (the first $2q$ entries are, instead, antisymmetrized). A basis of the subspace $D^{(m,q)}$ can be constructed by applying all the transformations of the permutation group to the vector $|m, q\rangle$ in Eq. (15) and choosing a maximal linearly independent subset.

We look for the eigenvectors of T_n in the subspaces

$$D^{(q)} = \bigoplus_{m=q}^{n-q} D^{(m,q)} \quad q = 0, \dots, \left\lfloor \frac{n}{2} \right\rfloor \tag{16}$$

of dimension $d_q(n + 1 - 2q)$. Since T_n has no symmetries beside the replicas permutation one, it has $n + 1 - 2q$ different eigenvalues in $D^{(q)}$, each with multiplicity d_q . In the following we will refer to the subspaces $D^{(q)}$ as to *sectors*. Moreover, with some abuse of notation, we shall use the symbol $D^{(q)}$ for the set of eigenvalues corresponding to eigenvectors in that sector.

Of particular relevance are the sectors $D^{(0)}$, $D^{(1)}$ and $D_0^{(2)}$ since they are associated to the Longitudinal, Anomalous and Replicon modes respectively from meanfield spin-glass theory¹⁴, as we will later show when discussing correlation functions in Section IV B.

By Eqs. (15) and (16) it is possible to factorize the replica symmetric part in the eigenvectors $\psi_q^\lambda(\sigma)$ of the transfer matrix in the sector $D^{(q)}$, that is we can write

$$\psi_q^\lambda(\sigma) = \rho_q^\lambda \left(\sum_a \sigma^a \right) \sum_{a_1 < \dots < a_q} C_{a_1 \dots a_q} \sigma^{a_1} \dots \sigma^{a_q} , \tag{17}$$

where the replica symmetric part ρ_q^λ of the eigenvectors is the one relevant to the computation of the eigenvalues. By last equation the eigenvectors of the sector $D^{(0)}$ are completely replica symmetric. The coefficients $C_{a_1 \dots a_q}$ are

invariant for any permutation of the indices and are equal to zero if any two of the indices are equal. Moreover they have to satisfy the constraint

$$\sum_{a_1=1}^n C_{a_1 \dots a_q} = 0, \quad (18)$$

which is a necessary and sufficient condition for any vector of the form of Eq. (17) to belong to the subspace $D^{(q)}$. Any set of d_q linearly independent coefficient vectors C can be chosen as an appropriate basis for the subspace. It is easy to prove that the product of two non-singular symmetric matrices possess a complete orthonormal (in the left-right sense) basis of left and right eigenvectors with real eigenvalues, and this is indeed case for T_n . In fact if we define, with a little abuse of notation, the vector

$$\rho_h(\sigma) \equiv \mathbb{E}_h e^{\beta h \sum_a \sigma^a}, \quad (19)$$

than $T_n(\sigma, \tau) = \sum_{\sigma'} \mathbb{E}_J e^{\beta J \sigma \sigma'} \times \delta_{\sigma' \tau} \rho_h(\tau)$. Moreover the left eigenvector ψ_L corresponding to a certain right $\psi_R(\sigma; \lambda, k)$ is simply given by

$$\psi_L(\sigma; \lambda, k) = \rho_h(\sigma) \psi_R(\sigma; \lambda, k), \quad (20)$$

where k denotes one choice of the coefficients $C_{a_1 \dots a_q}$ among the d_q possible. Imposing the orthonormality condition

$$\sum_{\sigma} \psi_L(\sigma; \lambda, k) \psi_R(\sigma; \lambda', k') = \delta_{\lambda \lambda'} \delta_{kk'}, \quad (21)$$

with the sum ranging over all the 2^n configuration of the replicated spin, and after successive application of Eq. (18), we obtain

$$\sum_{\sigma} \rho_q^{\lambda}(\sigma) \rho_h(\sigma) \rho_q^{\lambda'}(\sigma) \prod_{a=1}^q (1 - \sigma^{2a-1} \sigma^{2a}) = \delta_{\lambda \lambda'}, \quad (22)$$

along with

$$\sum_{a_1 < \dots < a_q} C_{a_1 \dots a_q}^k C_{a_1 \dots a_q}^{k'} = \delta_{kk'}. \quad (23)$$

We are now able to write down the transfer matrix in the spectral form

$$T_n(\sigma, \tau) = \sum_{q=0}^{\lfloor \frac{n}{2} \rfloor} T_{n,q}(\sigma, \tau) \quad (24)$$

where $T_{n,q}$ is the restriction of T_n to the subspace $D^{(q)}$, defined by

$$T_{n,q}(\sigma, \tau) = \sum_{\lambda \in D^{(q)}} \lambda \rho_q^{\lambda}(\sigma) \rho_h(\tau) \rho_q^{\lambda}(\tau) \sum_{\substack{a_1 < \dots < a_q \\ b_1 < \dots < b_q}} Q_{a_1 \dots a_q; b_1 \dots b_q} \sigma^{a_1} \dots \sigma^{a_q} \tau^{b_1} \dots \tau^{b_q}. \quad (25)$$

The coefficients Q appearing in last expression are invariant for any permutation of the set of indices a or b , therefore they depended only on the number of equal indexes in the sets $\{a_1, \dots, a_q\}$ and $\{b_1, \dots, b_q\}$. They are defined by

$$Q_{a_1 \dots a_q; b_1 \dots b_q} = \sum_{k=1}^{d_q} C_{a_1 \dots a_q}^k C_{b_1 \dots b_q}^k, \quad (26)$$

and their $(q+1)$ different values can be computed applying recursively Eqs. 18 and 23. If we denote $Q_p^{(q)}$ the coefficient in the sector $D^{(q)}$ with p pairs of different indexes, for the first sectors we have

$$Q_0^{(1)} = \frac{n-1}{n} \quad Q_1^{(1)} = -\frac{1}{n} \quad (27)$$

$$Q_0^{(2)} = \frac{n-3}{2(n-1)} \quad Q_1^{(2)} = -\frac{Q_0^{(2)}}{n-2} \quad Q_2^{(2)} = -\frac{2Q_1^{(2)}}{n-3} \quad (28)$$

B. Integral representations

In order to perform the limit $n \downarrow 0$ it is convenient to find a suitable parametrization for the eigenvectors of the form (17). For the replica symmetric part of the eigenvectors ψ_q^λ , see Eq. (17), we employ the standard parametrization

$$\rho_q^\lambda(\sigma) = \int du g_q^\lambda(u; n) \frac{e^{\beta u \sum_a \sigma^a}}{[2 \cosh(\beta u)]^n}, \quad (29)$$

in terms of the functions $g_q^\lambda(u; n)$. Turns out that all the functions g_q^λ parametrizing the eigenvectors of the sector $D^{(0)}$, are by themselves the eigenfunctions of an integral operator associated to that sector. In fact, expressing the linear terms in Eq. (17) through the identity

$$\sigma_{a_1} \dots \sigma_{a_q} = \frac{\partial}{\partial \epsilon_{a_1}} \dots \frac{\partial}{\partial \epsilon_{a_q}} \Big|_{\epsilon=0} e^{\sum_a \epsilon_a \sigma^a} \quad (30)$$

and plugging Eq. (29) into the eigenvalue equation $T_n \psi_q = \lambda \psi_q$, we obtain, after some manipulations, the new eigenvalue equation

$$\lambda g_q^\lambda(u; n) = \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, h + v)) \left(\frac{1}{\beta} \frac{\partial \hat{u}}{\partial v} \right)^q Z^n(J, h, v) g_q^\lambda(v; n). \quad (31)$$

The function $\hat{u}(J, x)$, defined by

$$\hat{u}(J, x) = \frac{1}{\beta} \operatorname{atanh}(\tanh(\beta J) \tanh(\beta x)), \quad (32)$$

will be recognized by the learned reader as the update rule for cavity messages. As we shall see, the function

$$Z(J, h, v) = \frac{2 \cosh(\beta J) \cosh(\beta(v + h))}{\cosh(\beta v)} \quad (33)$$

is related to the intensive free energy of an chain. Notice that in writing down Eq. (31) we have shifted the problem of finding a complete bases of eigenvectors for the matrix T_n to the equivalent problem of the spectral decomposition of the integral operators of Eq. (31), for $q = 0, \dots, \lfloor \frac{n}{2} \rfloor$. Turns out that, for a given sector $D^{(q)}$, the integral operator has a set of left eigenfunctions in the form

$$S_q^\lambda(v; n) = \mathbb{E}_h \int du g_q^\lambda(u; n) \left[\frac{\cosh(\beta(u + v + h))}{2 \cosh(\beta u) \cosh(\beta v)} \right]^n [1 - \tanh^2(\beta(u + v + h))]^q, \quad (34)$$

as can be inferred from Eq. (20) and can be directly verified. In the rest of the paper we will assume that the left and right eigenfunctions of the sector $D^{(q)}$ satisfy the normalization condition

$$\int du S_q^\lambda(u; n) g_q^{\lambda'}(u; n) = \delta_{\lambda\lambda'} \quad (35)$$

derived from Eq. (22). We are now ready to take the $n \downarrow 0$ limit and discuss its non trivial aspects.

C. The small n limit

In the limit $n \downarrow 0$ we obtain an infinite number of sectors $D^{(q)}$, $q = 0, 1, \dots$, in a fashion that is characteristic to replicas computations. Setting $n = 0$ in Eq. (31) we obtain Eq. (4), which we rewrite for convenience:

$$\lambda g_q^\lambda(u) = \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, h + v)) \left(\frac{1}{\beta} \frac{\partial \hat{u}}{\partial v} \right)^q g_q^\lambda(v). \quad (36)$$

From now on we shall refer to $g_q^\lambda(v)$ as a solution of last equation and shall explicitly express the n dependence for the solutions of (31) at finite n . In Figure 2 and Figure 3 we show two examples of eigenvalues and eigenfunctions in the sector $D^{(1)}$ and $D^{(2)}$ respectively.

LONGITUDINAL

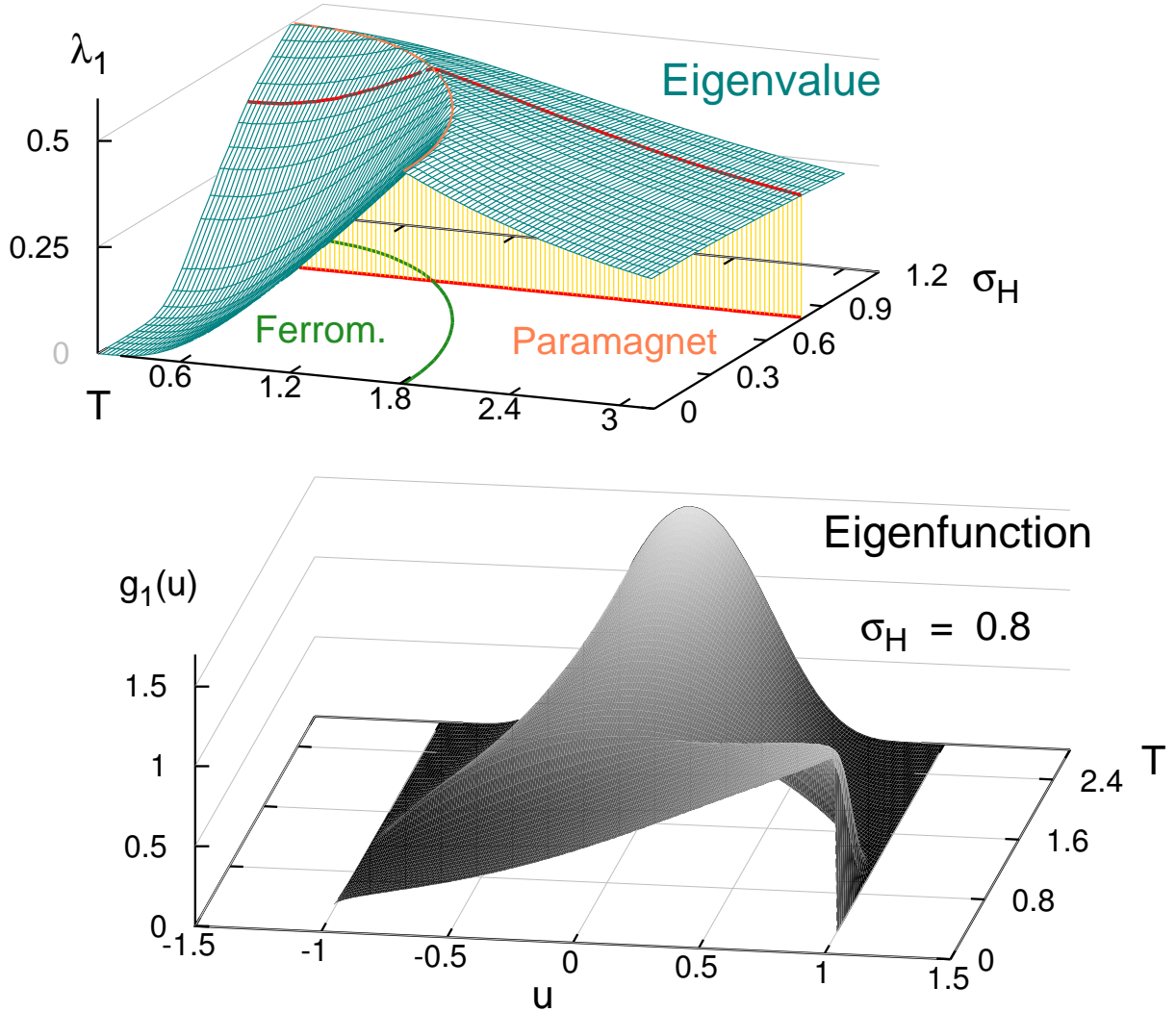


Fig. 2: (Top) The leading eigenvalue λ_1 of the sector $D^{(1)}$ in the RFIM, as a function of the temperature and of the gaussian external field with variance σ_H^2 . (Bottom) The corresponding right eigenfunction $g_1(u)$ at $\sigma_H = 0.8$. The random fields h and \tilde{h} are distributed as the cavity fields arriving on a chain embedded in a RRG with connectivity $z = 3$, therefore the transition point is localized at $\lambda_1 = \frac{1}{2}$.

For $q = 0$, i.e. in the sector $D^{(0)}$, Eq. (36) admits a unique maximum eigenvalue $\lambda = 1$ by Perron-Frobenius theorem. The corresponding eigenfunction is the probability distribution of cavity biases, which we call $P(u)$ ¹². We have thus established a first connection between the cavity method and the RTM formalism, and we shall enforce this connection in Section V. The other eigenfunctions of $D^{(0)}$ are characterized by $\int du g_0^\lambda(u) = 0$ at $n = 0$. It is convenient, to held compatibility with the normalization condition Eq. (35) as we will see, to impose a diverging scaling for all the eigenfunctions of $D^{(0)}$ except for the first one:

$$g_0^\lambda(u; n) \sim \frac{1}{\sqrt{n}} (g_0^\lambda(u) + n \tilde{g}_0^\lambda(u)). \quad (37)$$

The symbol \sim denotes equivalence between the r.h.s. and l.h.s. up to higher order correction in n , and \tilde{g}_0^λ is the

first correction to the leading order of the eigenfunction in $D^{(0)}$. Using Eq. (37) for the right eigenfunctions and considering also the correction in n to the eigenvalues, we can compute the left eigenfunctions of $D^{(0)}$ from Eq. (34). In fact we obtain at the leading order

$$S_0^\lambda(v; n) \sim \sqrt{n} S_0^\lambda(v) = \sqrt{n} \left[c_\lambda + \mathbb{E}_h \int du g_0^\lambda(u) \log \left(\frac{\cosh(\beta(u+v+h))}{\cosh(\beta u)} \right) \right], \quad (38)$$

where c_λ is the normalization of the first order correction to the eigenfunction g_0^λ , that is

$$c_\lambda \equiv \int du \tilde{g}_0^\lambda(u) = \frac{1}{\lambda - 1} \mathbb{E}_h \int du g_0^\lambda(u) \log \left(\frac{\cosh(\beta(u+h))}{\cosh(\beta u)} \right). \quad (39)$$

In all calculations involving the sector $D^{(0)}$ we will express the eigenvectors using Eqs. (37) and (38), then proceed carefully to take the $n \downarrow 0$ limit.

To find an expression for the left eigenfunctions in the other sectors no such care is needed to take the $n \downarrow 0$ limit in Eq. (34), therefore we straightly obtain

$$S_q^\lambda(v) = \mathbb{E}_h \int du g_q^\lambda(u) [1 - \tanh^2(\beta(u+v+h))]^q \quad \text{for } q \geq 1. \quad (40)$$

The degeneracy between $D^{(0)}$ and $D^{(1)}$ corresponds to the degeneracy between the Longitudinal and Anomalous eigenvalues in the Hessian of the Sherrington-Kirkpatrick model^{14,15}. The multiplicity of the eigenvalues in the two sectors, $d_0 = 1$ and $d_1 = n - 1$, sum up to give an $O(n)$ contribution as should be expected, while from Eq. (14) the other sectors have degeneracies of order $O(n)$ without the need of further elisions. Therefore it is convenient to define

$$\hat{d}_q = \begin{cases} 1 & \text{for } q = 1, \\ \lim_{n \rightarrow 0} \frac{d_q}{n} = (-1)^{q+1} \frac{2q-1}{q(q-1)} & \text{for } q \geq 2. \end{cases} \quad (41)$$

The first eigenvalue of $D^{(0)}$ requires separate considerations. We define the coefficient f_0 from its n expansion:

$$\lambda(n) \sim 1 - \beta f_0 n. \quad (42)$$

As we already noted, the cavity messages distribution $P(u)$ is the eigenvector associated to the largest eigenvalue of the sector $D^{(0)}$ for $n = 0$. The corresponding left eigenvalue is $S(u) \equiv 1$. In Section IV A we shall see that f_0 is the intensive free energy of a chain. From Eq. (47) we obtain

$$-\beta f_0 = \mathbb{E}_{J,h} \int dv \log [Z(J, h, v)] P(v). \quad (43)$$

D. The degeneracy between the Longitudinal and the Anomalous sector

A close inspection of the eigenvalue equation (36) reveals a surprising relation between the sectors $D^{(0)}$ and $D^{(1)}$ at $n = 0$. It can be shown, respectively deriving or integrating both members of Eq. (36) for $q = 1$ and $q = 0$, that all the eigenfunctions of $D^{(1)}$ have a corresponding eigenfunction in $D^{(0)}$ with the same eigenvalue. On the other hand, all the eigenfunctions of $D^{(0)}$, except for the first one, i.e. the ones having zero sum, have a corresponding eigenfunction in $D^{(1)}$ with the same eigenvalue. We have thus established a degeneracy between the Longitudinal and the Anomalous sectors. The following relations hold:

$$g_0^\lambda(u) = \frac{1}{\beta} \partial_u g_1^\lambda(u); \quad \frac{1}{\beta} \partial_u S_0^\lambda(u) = -S_1^\lambda(u). \quad (44)$$

Particular attention has to be taken in the limits involving these two sectors, keeping track of the $O(n)$ corrections both to eigenvalues and eigenvectors. A double pole contribution to some observables, as we shall later see, stems from the first correction in n to the paired eigenvalues in $D^{(0)}$ and $D^{(1)}$. In fact if we define the eigenvalue shifts $\delta\lambda_0$ and $\delta\lambda_1$ by

$$\lambda_0(n) \sim \lambda + n \delta\lambda_0, \quad (45)$$

$$\lambda_1(n) \sim \lambda + n \delta\lambda_1, \quad (46)$$

REPLICON

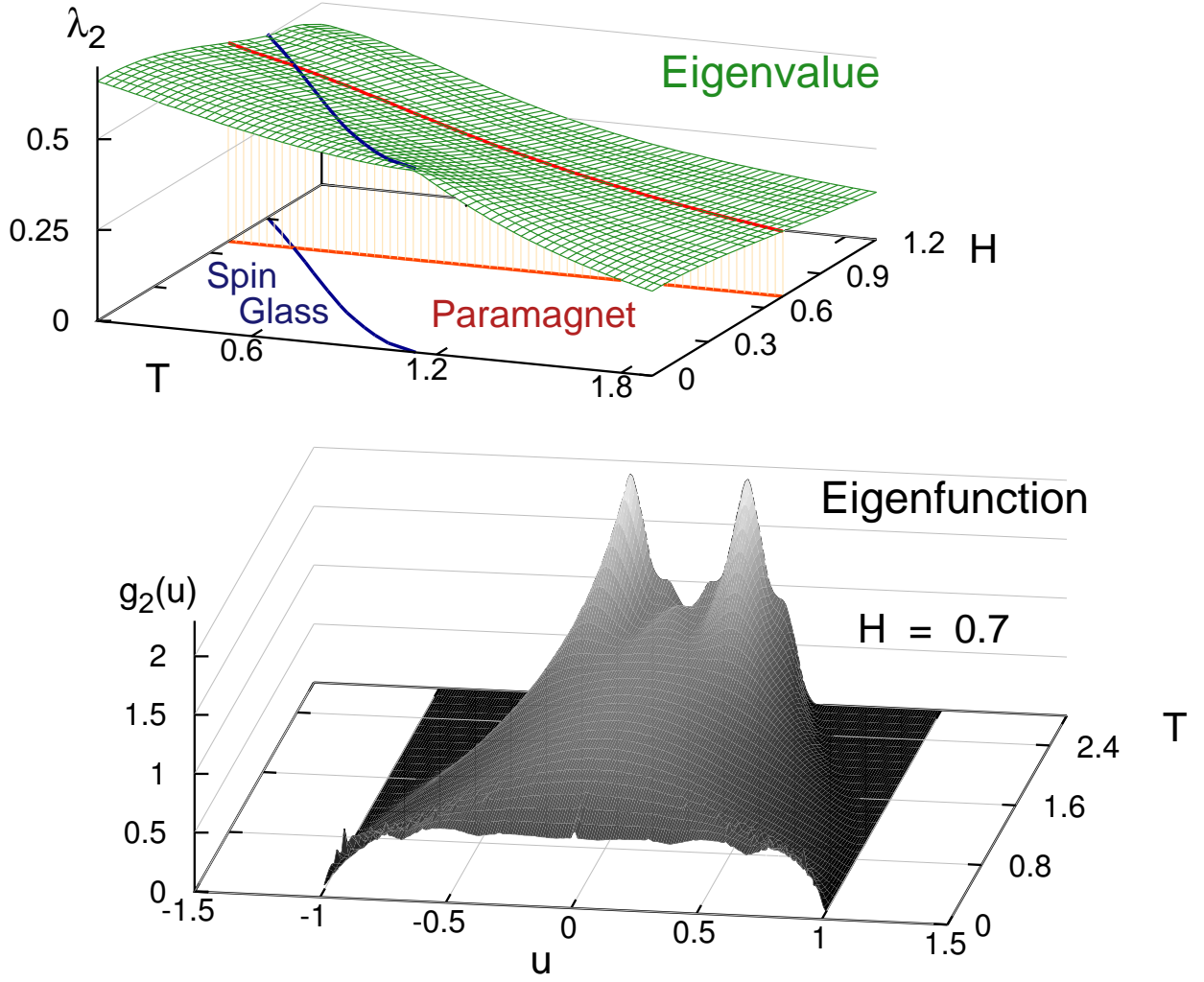


Fig. 3: (*Top*) The leading eigenvalue λ_2 of the sector $D^{(2)}$ in a $J = \pm 1$ spin-glass, as a function of the temperature and of the uniform external field H . The phase diagram is also shown in the $H - T$ plane. (*Bottom*) The corresponding right eigenfunctions $g_2(u)$ along the orange line of the top picture. The random fields h and \tilde{h} are distributed as the cavity fields arriving on a chain embedded in a RRG with connectivity $z = 3$, therefore the transition point is localized at $\lambda_2 = \frac{1}{2}$.

and consider the expansion to the first order in n of the eigenvalue equation (31) for $q =$, from standard perturbation theory we have

$$\delta\lambda_q = \mathbb{E}_{J,h} \int du dv S_q^\lambda(u) \log[Z(J, h, v)] \delta(u - \hat{u}(J, h + v)) \left(\frac{1}{\beta} \frac{\partial \hat{u}}{\partial v} \right)^q g_q^\lambda(v). \quad (47)$$

The shift difference $\Delta_\lambda = \delta\lambda_0 - \delta\lambda_1$ is the relevant quantity we are looking after, since it arises in the calculation of the free energies of closed chains and of the thermally disconnecter correlation function, see Section IV. Using Eq.

(47) and the relation (44) between the eigenfunctions in the two sectors, we obtain the expression

$$\Delta_\lambda = -\mathbb{E}_{J,h} \int du dv S_0^\lambda(u) \delta(u - \hat{u}(J, v + h)) [\tanh(\beta(v + h)) - \tanh(\beta v)] g_1^\lambda(v) \quad (48)$$

If we call (\bullet, \bullet) the scalar product in L^2 and define the kernel

$$Q(u, v) = \mathbb{E}_{J,h} \delta(u - \hat{u}(J, v + h)) [\tanh(\beta(v + h)) - \tanh(\beta v)], \quad (49)$$

then Eq. (48) can be rewritten as

$$\Delta_\lambda = -(S_0^\lambda, Q g_1^\lambda). \quad (50)$$

In the next Section we shall apply the formalism we have developed to the computation of some physically relevant quantities.

IV. SOME APPLICATIONS OF THE FORMALISM

A. Free energy of chains

Let us first consider the average free energy of a closed chain of length ℓ , each node receiving i.i.d. random fields h , and call it f_ℓ^c . If the chain considered is embedded in a locally tree-like graph, the random fields h are distributed according to the cavity messages distribution on that graph ensemble. Since $\text{Tr } T_n^\ell$ is the replicated partition function of this system, the free energy is given by

$$-\beta f_\ell^c = \lim_{n \rightarrow 0} \partial_n \text{Tr } T_n^\ell, \quad (51)$$

where, thanks the orthonormal decomposition of T_n , the trace can be written in the form

$$\text{Tr } T_n^\ell = \sum_{q=0}^{\frac{n}{2}} d_q \sum_{\lambda \in D^{(q)}} \lambda^\ell. \quad (52)$$

In last equation the eigenvalue degeneracies d_q are given in Eq. (14), and the eigenvalues λ depends implicitly on n . In the small n limit the sum over q can be extended to infinity. The considerations over the eigenvalues' shifts and degeneracies of last Section lead to the final expression

$$-\beta f_\ell^c = -\beta \ell f_0 + \sum_{\lambda \in D^{(1)}} \Delta_\lambda \ell \lambda^{\ell-1} + \sum_{q=1}^{\infty} \hat{d}_q \sum_{\lambda \in D^{(q)}} \lambda^\ell. \quad (53)$$

The coefficients \hat{d}_q are given by Eq. (41), the shift differences Δ_λ given by Eq. (48) and an expression for the intensive free energy f_0 is found in Eq. (43). We notice that all the quantities entering Eq. (53) can be expressed in terms of the eigenvalues and eigenfunctions of Eq. (36).

The computation of the average free energy of open chains is a little more involved. In the definition of open chains, we allow the spins at the extremities to receive a random field \tilde{h} that could have a distribution different from the one of the fields acting on the internal spins of the chain. We introduce this relaxation of the model in order to apply our formalism to the case of open chains embedded in a generic tree-like random graph.

It is convenient to define the replicated partition function of an open chain of length ℓ , conditioned on the configuration of the replicated spins at its extrema in the following way: starting from T_n^ℓ , we remove the field h on the right and substitute it with a field \tilde{h} , then we add the other field \tilde{h} on the left (see Figure 1). Therefore we define

$$\tilde{T}_n^{(\ell)}(\sigma, \tau) \equiv \rho_{\tilde{h}}(\sigma) T_n^\ell(\sigma, \tau) \rho_h^{-1}(\tau) \rho_{\tilde{h}}(\tau), \quad (54)$$

where, with a little abuse of notation, the vector $\rho_{\tilde{h}}$ is defined by

$$\rho_{\tilde{h}}(\sigma) \equiv \mathbb{E}_{\tilde{h}} e^{\beta \tilde{h} \sum_a \sigma^a}. \quad (55)$$

By definition the matrix $\tilde{T}_n^{(\ell)}$ is symmetric (see Figure 1 for a pictorial representation). From Eq. (54) and Eq. (24) we obtain the spectral decomposition

$$\tilde{T}_n^{(\ell)}(\sigma, \tau) = \sum_{q=0}^{\lfloor \frac{n}{2} \rfloor} \sum_{\lambda \in D^{(q)}} \lambda^\ell \rho_{\tilde{h}}(\sigma) \rho_q^\lambda(\sigma) \rho_{\tilde{h}}(\tau) \rho_q^\lambda(\tau) \sum_{\substack{a_1 < \dots < a_q \\ b_1 < \dots < b_q}} Q_{a_1 \dots a_q; b_1 \dots b_q} \sigma^{a_1} \dots \sigma^{a_q} \tau^{b_1} \dots \tau^{b_q}. \quad (56)$$

The average free energy of an open chain of length ℓ is then given by

$$-\beta f_\ell^o = \lim_{n \rightarrow 0} \partial_n \sum_{\sigma, \tau} \tilde{T}_n^{(\ell)}(\sigma, \tau). \quad (57)$$

From Eq. (56) it is easy to see that only the $D^{(0)}$ sector of $\tilde{T}_n^{(\ell)}$ contributes to last equation.

A different behaviour characterizes the terms corresponding to the leading eigenvalue at $n = 0$ (the cavity one) from the others. As in the case of the closed chain, the extensive contribution to the free energy comes from the leading eigenvalue of $D^{(0)}$, $\lambda \sim 1 - n\beta f_0$. An $O(1)$ contribution comes from the leading eigenfunction $g_0^\lambda(u; n) = P(u) + O(n)$, while each other eigenvalue, the ones degenerate with $D^{(0)}$, gives an exponential term. Therefore, after a careful treatment of the small n limit, we arrive to the expression

$$\begin{aligned} -\beta f_\ell^o &= -\ell\beta f_0 + \mathbb{E}_{\tilde{h}} \int du P(u) 2 \log \cosh(\beta(u + \tilde{h})) - \mathbb{E}_h \int dudv P(u)P(v) \log \cosh(\beta(u + v + h)) \\ &\quad + \log 2 + \sum_{\lambda \in D^{(1)}} a_{\lambda,0}^2 \lambda^\ell, \end{aligned} \quad (58)$$

with

$$a_{\lambda,0} = \frac{1}{\lambda - 1} \mathbb{E}_h \int du g_0^\lambda(u) \log \left[\frac{\cosh(\beta(u + h))}{\cosh(\beta u)} \right] + \mathbb{E}_{\tilde{h}} \int du g_0^\lambda(u) \log \left[\frac{\cosh(\beta(u + \tilde{h}))}{\cosh(\beta u)} \right] \quad (59)$$

In Eq. (58) it is clearly expressed at the order $O(1)$ in ℓ the free energy shift, with respect to the free energy of a closed chain, due to the addition of two extremal spins and the removal of an internal one.

The coefficients $a_{\lambda,0}$ are strictly related to the left eigenfunctions S_0^λ defined in Eq. (38). In fact if the random field at the extremities of the chain are distributed as the one on the internal spins, i.e. $\tilde{h} \stackrel{d}{=} h$ as in the case of a chain embedded in a Poissonian random graph, then $a_{\lambda,0} = S_0^\lambda(0)$. More generally if a probability distribution $G^{(0)}(u)$ exists such that

$$\tilde{P}(\tilde{h}) = \mathbb{E}_h \int du G^{(0)}(u) \delta(\tilde{h} - (u + h)) \quad (60)$$

holds, then Eq. (59) can be written in the compact form $a_{\lambda,0} = (S_0^\lambda, G^{(0)})$. Obviously if $\tilde{h} \stackrel{d}{=} h$ we have $G^{(0)}(u) = \delta(u)$. For a chain embedded in a random regular graphs ensemble instead, $G^{(0)}(u)$ is given by the distribution of cavity biases $P_{cav}(u)$ ¹².

B. Correlation functions

We take advantage of the spectral representation of the RTM to find some analytical expressions for the two-point correlation functions. We consider two spins, σ_0 and σ_ℓ , at distance ℓ along a chain. As in the previous paragraph, we admit the possibility for the chain to be embedded in a locally tree-like graph, therefore the random fields \tilde{h} acting on σ_0 and σ_ℓ can be distributed differently from the fields h on the internal spin of the chain. The decomposition of $\tilde{T}_n^{(\ell)}(\sigma, \tau)$ in Eq. (56) can be exploited to obtain the correlation functions. In fact contracting $\tilde{T}_n^{(\ell)}(\sigma, \tau)$ with two spins having the same replica index constrains them to be in the same thermal state, as in $\overline{\langle \sigma_0 \sigma_\ell \rangle} = \lim_{n \rightarrow 0} \sum_{\sigma, \tau} \sigma^1 \tilde{T}_n^{(\ell)}(\sigma, \tau) \tau^1$. Choosing different replica indexes instead corresponds to choosing different thermal states, e.g. $\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} = \lim_{n \rightarrow 0} \sum_{\sigma, \tau} \sigma^1 \tilde{T}_n^{(\ell)}(\sigma, \tau) \tau^2$. Generalizing this considerations is easy to obtain

$$\overline{\langle \sigma_0 \sigma_\ell \rangle^k} = \lim_{n \rightarrow 0} \sum_{\sigma, \tau} \sigma^1 \dots \sigma^k \tilde{T}_n^{(\ell)}(\sigma, \tau) \tau^1 \dots \tau^k. \quad (61)$$

Since vectors of the form $\sigma^1 \dots \sigma^k$ have non-zero projections in $D^{(q)}$ only for $q \leq k$, only these sectors of the spectral representation of $\tilde{T}_n^{(\ell)}$ contribute to Eq. (61). The expression for $\overline{\langle \sigma_0 \sigma_\ell \rangle_c^k}$ is quite complicated and it involves also the correction for small n to the eigenfunction of $D^{(0)}$ and $D^{(1)}$, as in the case of the thermally disconnected correlation function we shall later see. Therefore, since this kind of correlation function has little physical relevance, we won't report its expression in terms of the transfer matrix eigenvalues and eigenfunctions.

Far more interesting from the physical viewpoint are the connected correlation functions. The ferromagnetic connected correlation functions can be expressed as $\overline{\langle \sigma_0 \sigma_\ell \rangle_c} = \lim_{n \rightarrow 0} \frac{1}{2} \sum_{\sigma, \tau} (\sigma^1 - \sigma^2) \tilde{T}_n^{(\ell)}(\sigma, \tau) (\tau^1 - \tau^2)$, as one can rapidly check, and this expression can be easily generalized to

$$\overline{\langle \sigma_0 \sigma_\ell \rangle_c^k} = \lim_{n \rightarrow 0} \frac{1}{2^k} \sum_{\sigma, \tau} (\sigma^1 - \sigma^2) \dots (\sigma^{2k-1} - \sigma^{2k}) \tilde{T}_n^{(\ell)}(\sigma, \tau) (\tau^1 - \tau^2) \dots (\tau^{2k-1} - \tau^{2k}) . \quad (62)$$

It is worth noticing that the vector $v = (\sigma^1 - \sigma^2) \dots (\sigma^{2k-1} - \sigma^{2k})$ belongs to the subspace $D^{(k)}$, therefore we can choose a basis for the spectral representation of $\tilde{T}_n^{(\ell)}$ such that all but one vectors are orthogonal to v . This leads to the following compact expression for the connected correlation functions:

$$\overline{\langle \sigma_0 \sigma_\ell \rangle_c^k} = \sum_{\lambda \in D^{(k)}} a_{\lambda, k}^2 \lambda^\ell , \quad (63)$$

with the coefficients $a_{\lambda, k}$ given by

$$a_{\lambda, k} = \mathbb{E}_{\tilde{h}} \int du g_k^\lambda(u) \left[1 - \tanh^2 \left(\beta(u + \tilde{h}) \right) \right]^k . \quad (64)$$

As in the case of the coefficient $a_{\lambda, 0}$ defined in Eq. (59), if a solution $G^{(0)}$ of (60) exist then $a_{\lambda, k}$ is simply given by the projection of $G^{(0)}$ on S_k^λ , that is $a_{\lambda, k} = (S_k^\lambda, G^{(0)})$.

Equation (63) allows us to easily compute the susceptibilities $\chi_k = \lim_{N \rightarrow \infty} \frac{1}{N} \sum_{i, j} \overline{\langle \sigma_i \sigma_j \rangle_c^k}$ in a random graph with mean degree and mean residual degree z_0 and z respectively, in fact in thermodynamic limit we have

$$\begin{aligned} \chi_k &= \overline{(1 - m^2)^k} + \sum_{\ell=1}^{\infty} z_0 z^{\ell-1} \overline{\langle \sigma_0 \sigma_\ell \rangle_c^k} \\ &= \overline{(1 - m^2)^k} + z_0 \sum_{\lambda \in D^{(k)}} a_{\lambda, k}^2 \frac{\lambda}{1 - z\lambda} \end{aligned} \quad (65)$$

At a transition point the largest eigenvalue of one of the sectors $D^{(q)}$ reaches the value $\frac{1}{z}$ and the corresponding susceptibility diverges. Assuming a smooth behaviour for the eigenvalue in the high temperature region before the transition, $\lambda(T) = \frac{1}{z} + O(T - T_c)$ for $T \rightarrow T_c^+$, we obtain the mean-field critical exponent $\gamma = 1$.

The computation of the thermal disconnected correlation function $\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle}$, relevant to the RFIM transition, is more complicated, since it involves the subleading corrections in n to the eigenvectors of T_n . Great care has to be taken in the limit $\lim_{n \rightarrow 0} \sum_{\sigma, \tau} \sigma^1 \tilde{T}_n^{(\ell)}(\sigma, \tau) \tau^2 = \overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle}$. As in Eq. (37), let us call $\tilde{g}_0^\lambda(u)$ the correction to the eigenfunction $g_0^\lambda(u)$. We denote with $\langle \bullet \rangle_q$ the expectation over $\tilde{T}_n^{(\ell)}(\sigma, \tau)$ restricted to the sector $D^{(q)}$. Then in $D^{(0)}$ we obtain

$$\begin{aligned} \langle \sigma^1 \tau^2 \rangle_0 &\sim \overline{\langle \sigma_\infty \rangle^2} + \sum_{\lambda \in D^{(1)}} \frac{1}{n} a_{\lambda, 1}^2 \lambda^\ell + a_{\lambda, 1}^2 \ell \delta \lambda_0 \lambda^{\ell-1} \\ &\quad - 2a_{\lambda, 1} \lambda^\ell \left[\int du \tilde{g}_0^\lambda(u) \tanh(\beta(u + \tilde{h})) + \int du g_0^\lambda(u) \tanh(\beta(u + \tilde{h})) \log \frac{\cosh(\beta(u + \tilde{h}))}{\cosh(\beta u)} \right] , \end{aligned} \quad (66)$$

where the contribution $\overline{\langle \sigma_\infty \rangle}$ comes from the cavity eigenvector and is the average magnetization of a spin at the end of an infinite chain.

Similarly, if we define $\tilde{g}_1^\lambda(u)$ by $g_1^\lambda(u; n) \sim g_1^\lambda(u) + n \tilde{g}_1^\lambda(u)$, in the sector $D^{(1)}$ we have

$$\begin{aligned} \langle \sigma^1 \tau^2 \rangle_1 \sim & \sum_{\lambda \in D^{(1)}} \frac{-1}{n} a_{\lambda,1}^2 \lambda^\ell - a_{\lambda,1}^2 \ell \delta \lambda_1 \lambda^{\ell-1} \\ & - 2a_{\lambda,1} \lambda^\ell \left[\int du \tilde{g}_1^\lambda(u) \left(1 - \tanh^2(\beta(u + \tilde{h})) \right) + \int du g_1^\lambda(u) \left(1 - \tanh^2(\beta(u + \tilde{h})) \right) \log \frac{\cosh(\beta(u + \tilde{h}))}{\cosh(\beta u)} \right] \end{aligned} \quad (67)$$

Summing the two contributions, the final result for the disconnected correlation function is

$$\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} = \overline{\langle \sigma_\infty \rangle}^2 + \sum_{\lambda \in D^{(1)}} \Delta_\lambda a_{\lambda,1}^2 \ell \lambda^{\ell-1} + \alpha_\lambda \lambda^\ell. \quad (68)$$

Therefore each eigenvalue of the Anomalous sector contributes to $\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle}$ with a simple exponential term and with a term that leads to a double pole behaviour in the associated susceptibility, with coefficients Δ_λ given in Eq. (48). The coefficients α_λ of the exponential decays instead are given by

$$\begin{aligned} \alpha_\lambda = 2a_{\lambda,1} \left[\int du g_1^\lambda(u) \tanh(\beta(u + \tilde{h})) \left(\tanh(\beta(u + \tilde{h})) - \tanh(\beta u) \right) \right. \\ \left. - \int du \tilde{g}_1^\lambda(u) \left(1 - \tanh^2(\beta(u + \tilde{h})) \right) - \int du \tilde{g}_0^\lambda(u) \tanh(\beta(u + \tilde{h})) \right] \end{aligned} \quad (69)$$

Since the magnetization of a spin conditioned to be to be the extremity of a chain of size ℓ is given by

$$\overline{\langle \sigma_\ell \rangle} = \lim_{n \rightarrow 0} \sum_{\sigma, \tau} \sigma^1 \tilde{T}_n^{(\ell)}(\sigma, \tau) = \overline{\langle \sigma_\infty \rangle} - \sum_{\lambda \in D^{(1)}} a_{\lambda,1} a_{\lambda,0} \lambda^\ell, \quad (70)$$

if we call Λ the highest eigenvalue of the sector $D^{(1)}$, the most relevant contributions to the thermally-disconnected disorder-connected correlation function is given by

$$\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} - \overline{\langle \sigma_0 \rangle} \overline{\langle \sigma_\ell \rangle} = \Delta_\Lambda a_{\Lambda,1}^2 \ell \Lambda^{\ell-1} + \left(\alpha_\Lambda - 2\overline{\langle \sigma_\infty \rangle} a_{\Lambda,1} a_{\Lambda,0} \right) \Lambda^\ell + o(\Lambda^\ell) \quad \text{for } \ell \rightarrow +\infty. \quad (71)$$

We notice that, while the coefficient of the exponential term is quite hard to compute, the coefficient $a_{\Lambda,1}^2 \Delta_\Lambda$, which regulates the leading behaviour, has a much simpler expression given in Eq. (48) and Eq. (64). From Eq. 71 turns out that near a ferromagnetic transition point, i.e. $\Lambda = \frac{1}{z}$, as long as Δ_Λ is not zero, the leading behaviour of the disconnected susceptibility $\chi_{disc} = \sum_{i,j} \overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} - \overline{\langle \sigma_0 \rangle} \overline{\langle \sigma_\ell \rangle}$ reads

$$\chi_{disc} \simeq z_0 \Delta_\Lambda a_{\Lambda,1}^2 \frac{1}{(1 - z\Lambda)^2}. \quad (72)$$

The expected double-pole behaviour of the disconnected susceptibility is thus recovered.

V. CAVITY DERIVATION

In this section we present the derivation of several of the results of last Section using a probabilistic approach, in the same spirit of the usual cavity method calculations^{9,12}. While being this approach more physically intuitive with respect to the RTM formalism, it requires the set up of an ad-hoc recursion rule for each observable. Noticeably we could not recover Eq. (53) for the free energy of closed chains.

A. Open chains

We want to study the statistical properties of a random Ising open chain without the use of replicas. We start with an asymmetric chain of length ℓ , whose random partition function we denote with Z_ℓ , constructed iteratively according to the following procedure: Z_0 is the partition function of a single spin receiving a random field u_0 , i.e. $Z_0 = 2 \cosh(\beta u_0)$; at the i -th step of the construction we add a spin σ_i , a random coupling J_i between σ_i and σ_{i-1}

and a random field h_{i-1} on σ_{i-1} ; the random variable Z_ℓ is the partition function of the system obtained after the ℓ -th step of the procedure. Note that the last spin added to the chain has no external fields acting on it. The following distributional identity can be easily derived:

$$Z_{\ell+1} = \frac{2 \cosh(\beta J_\ell) \cosh(\beta(u_\ell + h_\ell))}{\cosh(\beta u_\ell)} \times Z_\ell \equiv Z(J_\ell, h_\ell, u_\ell) \times Z_\ell. \quad (73)$$

It is convenient to introduce the quantity $\overline{Z_\ell^n(u)} \equiv \overline{\delta(u - u_\ell) Z_\ell^n}$, which corresponds to the expectation of Z_ℓ^n along with the indicator function of the event $u_\ell = u$. Here n is an arbitrary chosen positive real number, the symbol being chosen to stress the analogy with the replica formalism where the quantity n (integer in this case) is the number of replicated systems. Using this definition from Eq. (73) follows readily

$$\overline{Z_{\ell+1}^n(u)} = \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, v + h)) Z^n(J, h, v) \overline{Z_\ell^n(v)}, \quad (74)$$

where $\tilde{u}(J, x) = \frac{1}{\beta} \operatorname{atanh}(\tanh(\beta J) \tanh(\beta x))$ is the usual message passing rule. The integral operator of Eq. (74) is the same we found in the RTM formalism in Eq. (31) for the sector $D^{(0)}$, therefore we can make use of the spectral analysis result from those paragraphs, in particular of the completeness relation

$$\mathbb{E}_{J,h} \delta(u - \hat{u}(J, v + h)) Z^n(J, h, v) = \sum_{\lambda \in D^{(0)}} \lambda(n) g_0^\lambda(u; n) S_0^\lambda(v; n), \quad (75)$$

between left and right eigenvectors. The definition of the left eigenfunctions of $D^{(0)}$ was already given in Eq. (34), but we rewrite it for convenience:

$$S_0^\lambda(v; n) = \mathbb{E}_h \int du g_0^\lambda(u; n) \left[\frac{\cosh(\beta(u + v + h))}{2 \cosh(\beta u) \cosh(\beta v)} \right]^n. \quad (76)$$

Let us define another random partition function, $Z_\ell(u; x)$, obtained from $Z_\ell(u)$ conditioning on the value of the message u_0 on the first spin, that is $Z_\ell(u; x) = Z_\ell(u) | (u_0 = x)$. Since also $\overline{Z_\ell^n(u; x)}$ as a function of u obeys equation (74), using the decomposition Eq. (75) and the initial condition $Z_0 = 2 \cosh(\beta u_0)$ we arrive to the important result

$$\overline{Z_\ell^n(u; x)} = \sum_{\lambda \in D^{(0)}} \lambda^\ell(n) g_0^\lambda(u; n) S_0^\lambda(x; n) [2 \cosh(\beta x)]^n. \quad (77)$$

Using last equation it is easy to compute any moment $\overline{Z_\ell^n}$, n not necessarily integer, of the partition function of a random asymmetric Ising chain of length ℓ . More interesting is the computation of the properties of a symmetric Ising open chain, the one considered in Section IV A, which receives on each extremity an external field distributed according to a certain probability distribution $\tilde{P}(\tilde{h})$. As already stated, this is definition stems from the need to cover the important case of a chain embedded in a locally tree-like graph. Let us call $Z_{\ell,o}$ the random partition function of this open chain. It is related to the random partition function Z_ℓ of the asymmetric open chain by

$$Z_{\ell,o} = \frac{\cosh(\beta(u_\ell + \tilde{h}_\ell))}{\cosh(\beta u_\ell)} \times Z_{\ell-1}(u_\ell; u_1) \times \frac{2 \cosh(\beta J_0) \cosh(\beta \tilde{h}_0)}{\cosh(\beta u_1)} \quad (78)$$

where u_1 is distributed as $\tilde{u}(J_0, \tilde{h}_0)$. From Eq. (78) along with Eq. (77) and Eq. (76), we derive the main result of this paragraph:

$$\overline{Z_{\ell,o}^n} = \sum_{\lambda \in D^{(0)}} \lambda^\ell(n) a_{\lambda,0}^2(n). \quad (79)$$

where $a_{\lambda,0}(n)$ is defined by

$$a_{\lambda,0}(n) \equiv \mathbb{E}_{\tilde{h}} \int du \left[\frac{\cosh(\beta(u + \tilde{h}))}{\cosh(\beta u)} \right]^n g_0^\lambda(u; n). \quad (80)$$

In the RTM formalism of Section III and IV, last expression could be derived from $\tilde{T}_n^{(\ell)}$ defined in Eq. (56) by analytic continuation of $\overline{Z_{\ell,o}^n} = \sum_{\sigma,\tau} \tilde{T}_n^{(\ell)}(\sigma, \tau)$ to non-integer n .

The average free energy of an open chain of length ℓ can then be obtained by

$$-\beta f_\ell^o = \lim_{n \rightarrow 0} \partial_n \overline{Z_{\ell,o}^n}. \quad (81)$$

The computation involves computing the order n of all the quantities present in Eq. (79), as it was done in Section (IV A). In this paragraph however, without any use of replicas, we gave a purely probabilistic argument valid for any real value of n . We refer therefore to Section (IV A) for the successive step of the computation of f_ℓ^o , leading to the final result Eq. (58). Notice that in the notation of that paragraph $a_{\lambda,0}$ is related to $a_{\lambda,0}(n)$ defined in Eq. (80) by $a_{\lambda,0}(n) \sim \sqrt{n} a_{\lambda,0}$.

The expression (58) for f_ℓ^o could also be obtained by a different approach that does not involve any limit $n \downarrow 0$ but is technically more difficult. We define the function $\varphi^{(\ell)}(u)$ by

$$\varphi^{(\ell)}(u) \equiv \overline{\delta(u - u_\ell) \log Z_\ell}, \quad (82)$$

and observe that given the distribution of the cavity message at distance ℓ along the chain, u_ℓ , which we call $G_0^{(\ell)}(u)$, it obeys the iterative rule

$$\begin{aligned} \varphi^{(\ell+1)}(u) &= \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, h + v)) \varphi^{(\ell)}(v) \\ &\quad + \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, h + v)) \log \left[\frac{2 \cosh(\beta J) \cosh(\beta(v + h))}{\cosh(\beta v)} \right] G_0^{(\ell)}(v) \end{aligned} \quad (83)$$

Last equation can be solved decomposing $\varphi^{(\ell)}(u)$ and $G_0^{(\ell)}(v)$ along the eigenfunctions of $D^{(0)}$ at $n = 0$, then $\varphi^{(\ell)}(u)$ can be used to obtain f_ℓ^o .

B. Connected correlation functions

Let us derive the eigenvalue equation (4) and the expression for the connected correlation functions Eq. (8), without making any use of replicas. Here we consider straightly the random open chain with partition function $Z_{\ell,o}$, characterized by independent random external field distributes a h on the internal spins and as \tilde{h} on the extremities. The connected correlation function $\langle \sigma_0 \sigma_\ell \rangle_c = \frac{\partial \langle \sigma_\ell \rangle}{\partial H_0}$, where H_0 is an auxiliary field acting on σ_0 , can be expressed as a function of the message u_ℓ , coming through the chain to the spin σ_ℓ , and its derivative with respect to H_0 . In fact we have

$$\langle \sigma_0 \sigma_\ell \rangle_c = \left(1 - \tanh^2(\beta(\tilde{h}_\ell + u_\ell)) \right) \frac{\partial u_\ell}{\partial H_0}, \quad (84)$$

where \tilde{h}_ℓ , as usual, is the random effective field acting on σ_ℓ and coming eventually from the rest of the graph. Let us define the random variable X_ℓ by $X_\ell \equiv \frac{\partial u_\ell}{\partial H_0}$. The average over disorder of Eq. (84) and its moments $\langle \sigma_0 \sigma_\ell \rangle_c^q$ can then be computed once we know the joint law of the random variables u_ℓ and X_ℓ , which we call $P_\ell(u, X)$. Since X_ℓ obeys the chain rule $X_{\ell+1} = \frac{\partial u_{\ell+1}}{\partial u_\ell} X_\ell$ the recursion rule for P_ℓ reads

$$P_{\ell+1}(u, X) = \mathbb{E}_{J,h} \int dv dY \delta \left(X - \frac{\partial \hat{u}}{\partial v} Y \right) \delta(u - \hat{u}(J, h + v)) P_\ell(v, Y), \quad (85)$$

where \hat{u} is the message passing rule defined in Eq. (32). From last expression it turns out we can write an iteration rule for the momenta of X_ℓ at fixed u_ℓ ,

$$G_q^{(\ell)}(u) = \int dX P_\ell(u, X) X^q, \quad (86)$$

which reads

$$G_q^{(\ell+1)}(u) = \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, h + v)) \left(\frac{\partial \hat{u}}{\partial v} \right)^q G_q^{(\ell)}(v). \quad (87)$$

Equations (85) and (87) with $q = 2$ have been recently introduced in literature⁸ in order to derive an analytical expression for the spin-glass susceptibility.

We note that the knowledge of the maximum eigenvalue of the integral operator of Eq. (87) for a generic q allows one to reconstruct the full distribution of the connected correlation function at large distance¹⁶.

From last equation it is clear the relation of $G_q^{(\ell)}$ with the eigenfunctions g_q^λ of Eq. (4). In fact, decomposing $G_q^{(\ell)}(u)$ along the eigenfunctions of $D^{(0)}$, projecting Eq. (87) on the left eigenvectors $S_q^\lambda(u)$ and with some computations analogue to the ones leading from Eq. (79) to Eq. (80), we arrive to

$$G_q^{(\ell)}(u) = \sum_{\lambda \in D^{(q)}} a_{\lambda,q} \lambda^\ell g_q^\lambda(u), \quad (88)$$

where $a_{\lambda,q}$ is defined in Eq. (64). Equation (88), along with Eq. (84), gives the expression (63) obtained with the RTM formalism for the connected correlation functions.

Following the lead of the previous paragraph, we can extend the above derivation to compute the disorder averages $\langle \sigma_0 \sigma_\ell \rangle_c^q$ along with an arbitrary power of $Z_{\ell,o}$. The generalization of Eq. (87) in fact becomes

$$G_q^{(\ell+1)}(u; n) = \mathbb{E}_{J,h} \int dv \delta(u - \hat{u}(J, h + v)) \left(\frac{\partial \hat{u}}{\partial v} \right)^q Z(J, h, v) G_q^{(\ell)}(v; n), \quad (89)$$

and Eq. (88) generalizes trivially as well. The final result is

$$\overline{\langle \sigma_0 \sigma_\ell \rangle_c^q Z_{\ell,o}^n} = \sum_{\lambda \in D^{(q)}} \lambda^\ell(n) a_{\lambda,q}^2(n), \quad (90)$$

which extrapolates smoothly to the result we obtained for $n = 0$, i.e. Eq. (63). In last equation the coefficients $a_{\lambda,q}(n)$ are defined by

$$a_{\lambda,q}(n) \equiv \mathbb{E}_{\tilde{h}} \int du \left[\frac{\cosh(\beta(u + \tilde{h}))}{\cosh(\beta u)} \right]^n \left[1 - \tanh^2(\beta(u + \tilde{h})) \right]^q g_q^\lambda(u; n), \quad (91)$$

such that $a_{\lambda,q}(0) = a_{\lambda,q}$.

C. The disconnected correlation function

The computations of the thermally disconnected correlation function $\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle}$ is straightforward once we use the results we obtained in the two preceding paragraphs. In fact, calling H_0 and H_ℓ two auxiliary fields we add to the first and the last spin of the chain respectively and set to zero after the computation, the following relation holds:

$$\frac{\partial}{\partial H_0} \frac{\partial}{\partial H_\ell} \overline{Z_{\ell,o}^n} = n \overline{\langle \sigma_0 \sigma_\ell \rangle_c} \overline{Z_{\ell,o}^n} + n^2 \overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} \overline{Z_{\ell,o}^n}. \quad (92)$$

Using Eqs. (79) and (90) last expression leads to the main result of this paragraph, that is

$$\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle} \overline{Z_{\ell,o}^n} = \frac{1}{n} \left[\sum_{\lambda \in D^{(0)}} \lambda^\ell(n) b_{\lambda,0}^2(n) - \sum_{\lambda \in D^{(1)}} \lambda^\ell(n) a_{\lambda,1}^2(n) \right], \quad (93)$$

where the coefficient $b_{\lambda,0}(n) \equiv \frac{\partial a_{\lambda,0}(n)}{\partial H_{0/\ell}}$ reads

$$b_{\lambda,0}(n) = \mathbb{E}_{\tilde{h}} \int du \left[\frac{\cosh(\beta(u + \tilde{h}))}{\cosh(\beta u)} \right]^n \tanh(\beta(u + \tilde{h})) \sqrt{n} g_0^\lambda(u; n). \quad (94)$$

We included a factor \sqrt{n} in the definition of $b_{\lambda,0}(n)$ to facilitate the extrapolation of Eq. (93) to small n . In fact for all but the first eigenfunctions of $D^{(0)}$ the normalization condition imposes the scaling $g_0^\lambda(u; n) \sim \frac{1}{\sqrt{n}} [g_0^\lambda(u) + n \tilde{g}_0^\lambda(u)]$.

We could derive Eq. (93) also in the RTM formalism for integer values of n and then perform an analytic continuation to arbitrary real n . In the limit $n \downarrow 0$ it is easy to see that the contribution to $\overline{\langle \sigma_0 \rangle \langle \sigma_\ell \rangle}$ from the first and the second sums of Eq. (93) are given in Eqs. (66) and (67) of Section IV B respectively.

An alternative probabilistic derivation of the formula (68) for $\overline{\langle\sigma_0\rangle\langle\sigma_\ell\rangle}$, which does not require the knowledge of the momentums of the partition function and of $\overline{\langle\sigma_0\sigma_\ell\rangle_c} Z_{\ell,o}^n$, goes through the definition of

$$R^{(\ell)}(u) \equiv \overline{\delta(u - u_\ell)\langle\sigma_0\rangle^{(\ell)}}. \quad (95)$$

We used the symbol $\langle\sigma_0\rangle^{(\ell)}$ to denote the magnetization of the first spin at the ℓ -th iteration of the construction of the asymmetric chain described in Section V A. It can be easily shown that the knowledge of $R^{(\ell)}(u)$ allows the computation of $\overline{\langle\sigma_0\rangle\langle\sigma_\ell\rangle}$. Since $\langle\sigma_0\rangle^{(\ell)}$ is given by the derivative of the free energy of the chain at the step ℓ with respect to a field on the the first spin, considering the free energy difference after an iteration it is easy to arrive to the relation

$$\langle\sigma_0\rangle^{(\ell+1)} = \langle\sigma_0\rangle^{(\ell)} + [\tanh(\beta(u_\ell + h)) - \tanh(\beta u_\ell)] \frac{\partial u_\ell}{\partial h_0}. \quad (96)$$

Therefore the recursion rule for $R^{(\ell)}(u)$ is given by

$$\begin{aligned} R^{(\ell+1)}(u) &= \mathbb{E}_{J,h} \int dv R^{(\ell)}(v) \delta(u - \hat{u}(J, h + v)) \\ &+ \mathbb{E}_{J,h} \int dv G_1^{(\ell)}(v) \delta(u - \hat{u}(J, h + v)) [\tanh(\beta(v + h)) - \tanh(\beta v)], \end{aligned} \quad (97)$$

where $G_1^{(\ell)}(v)$ was defined in Eq. (88) in last paragraph. Last equation can be solved decomposing $R^{(\ell)}(u)$ along the eigenfunctions of $D^{(0)}$ at $n = 0$, and using Eq. (88) for $G_1^{(\ell)}(v)$. The computation is lengthy and not trivial, since it involves expressing \tilde{g}_0^λ and \tilde{g}_1^λ (defined in Section IV B) respectively in terms of the basis of $D^{(0)}$ and $D^{(1)}$ at $n = 0$. In the end though one arrives at the expression (68) for the disconnected correlation function.

VI. CONCLUSIONS

In the present paper we presented a thorough analysis of the spectral properties of the RTM. We have developed a formalism that is suitable to compute many different types of connected and disconnected correlation functions and can be applied both to one-dimensional systems and to locally tree-like graphs. The expressions we found are exact for any value ℓ of the spin distance and can be approximated numerically considering only the top eigenvalues of certain integral operators. Also the formalism can be trivially adapted to perform the same computations in diluted p-spin models.

We also managed to obtain exact formulas for the moments of the partition function and of the average free energies of open and closed chains of finite length. It has been recently found that short chains have an important role in the finite size corrections to disordered models on diluted graphs⁴ and in perturbative expansions around the Bethe approximation on Euclidean systems¹⁰. Therefore the analytical tools we have developed also apply to these contexts.

Most of the results have also been derived using rigorous probabilistic arguments. This approach has the merits of avoiding the complication of the decomposition of the replicated space $Z_2^{\otimes n}$ and of being more physically intuitive of the replica one. The advantage of the replica method instead is that once the spectral representation of the RTM is obtained all the observables can be computed just with opportune contraction. In the cavity analysis an ad-hoc iterative function or a computation strategy has to be devised for each observable.

Noticeably we did not manage to derive Eq. (53) for the free energy of closed chains using a cavity argument. This is the only point withstanding the proof of the complete equivalence between the two methods.

A limitation of both the RTM formalism and of its cavity counterpart, is the fact that it is applicable to the analysis of disordered Ising models only in their replica symmetric phase. This includes all isolated one-dimensional systems but not diluted models in the spin glass phase. Therefore an investigation of the spectral properties of the 1RSB replicated transfer matrix, extending Wigner's decomposition¹³ to the 1RSB symmetry group, is desirable. Another direction for the extension of our results, which should not require too much analytical effort¹⁷, is toward the investigation of Potts models.

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