

Numerical study of the branching tree of states in spin glasses

G. Parisi, F. Ricci-Tersenghi, **D. Yllanes**

Dipartimento di Fisica
La Sapienza Università di Roma

CompPhys13, Leipzig University, November 2013



The Sherrington-Kirkpatrick model

- **Spin glasses:** Random, mixed-interacting magnetic systems that experience a random, yet cooperative, freezing of spins below some critical temperature T_c .

The Sherrington-Kirkpatrick model

- **Spin glasses:** Random, mixed-interacting magnetic systems that experience a random, yet cooperative, freezing of spins below some critical temperature T_c .

Edwards-Anderson model

$$\mathcal{H} = - \sum_{\langle i,j \rangle} J_{ij} s_i s_j, \quad s_i = \pm 1$$

- $J_{ij} = \pm 1$ with 50% probability.
- Disorder and **frustration**
- Order parameter from the overlap

$$q = \lim_{t \rightarrow \infty} \frac{1}{N} \sum_x \langle s_x(0) s_x(t) \rangle_t \quad \rightsquigarrow \quad q = \frac{1}{N} \sum_x \langle s_x \rangle^2.$$

The Sherrington-Kirkpatrick model

The Sherrington-Kirkpatrick model

- The EA model is too difficult to handle analytically.
- We consider its mean-field version:

$$\mathcal{H}_J = - \sum_{i,j} J_{ij} s_i s_j, \quad s_i = \pm 1 \quad J_{ij} = \pm \frac{1}{\sqrt{N}}.$$

The Sherrington-Kirkpatrick model

- The EA model is too difficult to handle analytically.
- We consider its mean-field version:

$$\mathcal{H}_J = - \sum_{i,j} J_{ij} s_i s_j, \quad s_i = \pm 1 \quad J_{ij} = \pm \frac{1}{\sqrt{N}}.$$

- Solution by G. Parisi (Replica Symmetry Breaking, RSB).
- The overlap can take any value in $[0, q_M]$ with non-zero probability density $p(q)$.
- $p(q)$ is smooth in $[0, q_M)$, but has a δ function at q_M

The Sherrington-Kirkpatrick model

- The EA model is too difficult to handle analytically.
- We consider its mean-field version:

$$\mathcal{H}_J = - \sum_{i,j} J_{ij} s_i s_j, \quad s_i = \pm 1 \quad J_{ij} = \pm \frac{1}{\sqrt{N}}.$$

- Solution by G. Parisi (Replica Symmetry Breaking, RSB).
- The overlap can take any value in $[0, q_M]$ with non-zero probability density $p(q)$.
- $p(q)$ is smooth in $[0, q_M)$, but has a δ function at q_M
- The order parameter is not a number, but a function $q(x)$.
- $x(q)$ is the cumulative probability of q : $x(q) = \int_0^q dq' p(q')$.

Infinitely many states

- From the previous discussion, we know that there are infinitely many relevant **pure states**:

$$F_\alpha - F_\beta = \mathcal{O}(1), \quad \text{even if } N \rightarrow \infty.$$

Infinitely many states

- From the previous discussion, we know that there are infinitely many relevant **pure states**:

$$F_\alpha - F_\beta = \mathcal{O}(1), \quad \text{even if } N \rightarrow \infty.$$

- It makes sense to consider restricted averages: $\langle A \rangle_\alpha$, so that intensive quantities do not fluctuate in $\langle \dots \rangle_\alpha$

Infinitely many states

- From the previous discussion, we know that there are infinitely many relevant **pure states**:

$$F_\alpha - F_\beta = \mathcal{O}(1), \quad \text{even if } N \rightarrow \infty.$$

- It makes sense to consider restricted averages: $\langle A \rangle_\alpha$, so that intensive quantities do not fluctuate in $\langle \dots \rangle_\alpha$
- Each state will have a weight w_α

$$\langle A \rangle = \sum_{\alpha} w_{\alpha} \langle A \rangle_{\alpha}.$$

(ferromagnet: consider $\langle A \rangle = \frac{1}{2} \langle A \rangle_+ + \frac{1}{2} \langle A \rangle_-$).

Infinitely many states

- From the previous discussion, we know that there are infinitely many relevant **pure states**:

$$F_\alpha - F_\beta = \mathcal{O}(1), \quad \text{even if } N \rightarrow \infty.$$

- It makes sense to consider restricted averages: $\langle A \rangle_\alpha$, so that intensive quantities do not fluctuate in $\langle \dots \rangle_\alpha$
- Each state will have a weight w_α

$$\langle A \rangle = \sum_{\alpha} w_{\alpha} \langle A \rangle_{\alpha}.$$

- Example: the overlap:

$$q_{\alpha\beta} = \frac{1}{N} \sum_i \langle s_i \rangle_{\alpha} \langle s_i \rangle_{\beta} \implies p(q) = \overline{\sum_{\alpha\beta} w_{\alpha} w_{\beta} \delta(q - q_{\alpha\beta})}.$$

Infinitely many states

- From the previous discussion, we know that there are infinitely many relevant **pure states**:

$$F_\alpha - F_\beta = \mathcal{O}(1), \quad \text{even if } N \rightarrow \infty.$$

- It makes sense to consider restricted averages: $\langle A \rangle_\alpha$, so that intensive quantities do not fluctuate in $\langle \dots \rangle_\alpha$
- Each state will have a weight w_α

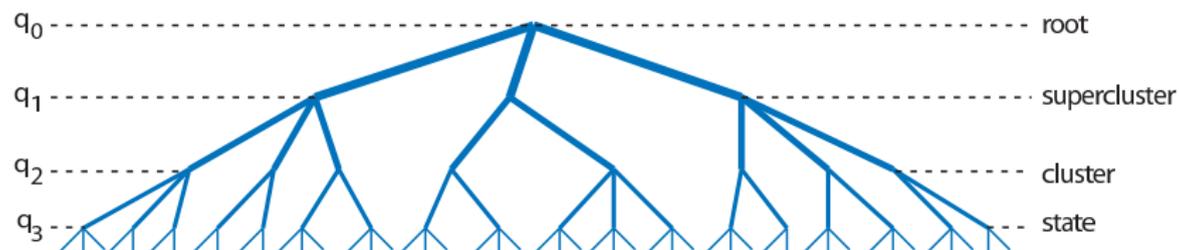
$$\langle A \rangle = \sum_{\alpha} w_{\alpha} \langle A \rangle_{\alpha}.$$

- Example: the overlap:

$$q_{\alpha\beta} = \frac{1}{N} \sum_i \langle s_i \rangle_{\alpha} \langle s_i \rangle_{\beta} \implies p(q) = \overline{\sum_{\alpha\beta} w_{\alpha} w_{\beta} \delta(q - q_{\alpha\beta})}.$$

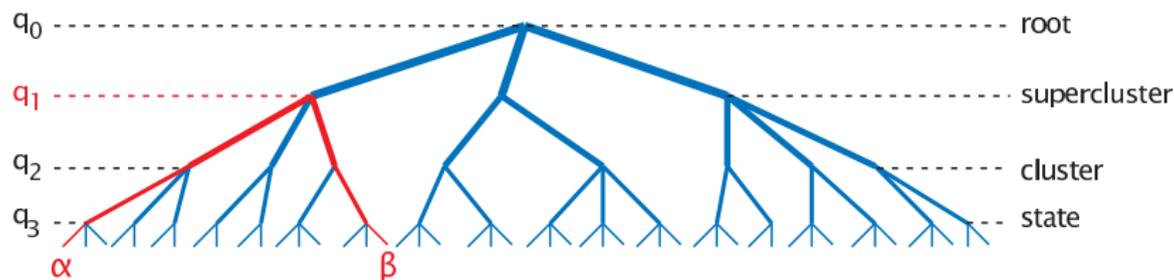
- It turns out that the states live in an **ultrametric space**, using $q_{\alpha\beta}$ to define a distance
- We can classify the states in a taxonomic tree, which branches out as we break the replica symmetry.

The tree of states (I)



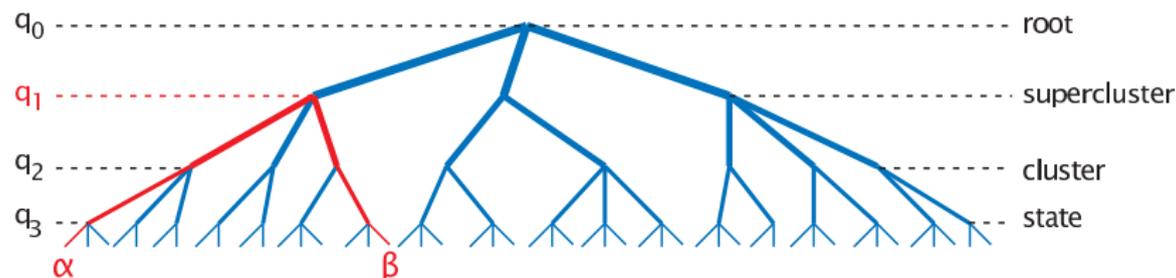
- Consider a simplified case where q can only take four discrete values (equivalently $K = 3$ steps of RSB).
- We can group the **configurations** (organisms) \rightarrow **states** (species) \rightarrow **clusters** (geni) \rightarrow **superclusters** (families).

The tree of states (I)



- Consider a simplified case where q can only take four discrete values (equivalently $K = 3$ steps of RSB).
- We can group the configurations \rightarrow states \rightarrow clusters \rightarrow superclusters .
- It makes sense, because the **overlap** (distance) depends only on the first **common ancestor**: $q_{\alpha\beta} = q_1$.

The tree of states (I)



- Consider a simplified case where q can only take four discrete values (equivalently $K = 3$ steps of RSB).
- We can group the configurations \rightarrow states \rightarrow clusters \rightarrow superclusters .
- It makes sense, because the **overlap** (distance) depends only on the first **common ancestor**: $q_{\alpha\beta} = q_1$.
- Notice that the self-overlap is the same for all states:

$$q_{\alpha\alpha} = q_M, \quad \forall \alpha$$

The tree of states (II)

- The real tree of states of a mean-field spin glass is more complicated:
 - Infinitely many levels: the tree branches out for any value of $q < q_M$.
 - Infinite number of branches at any level.

The tree of states (II)

- The real tree of states of a mean-field spin glass is more complicated:
 - Infinitely many levels: the tree branches out for any value of $q < q_M$.
 - Infinite number of branches at any level.
- Amusing analogy with QED: number of emitted photons diverges at low energy \leftrightarrow infinite number of branches as we consider small values of w .

The tree of states (II)

- The real tree of states of a mean-field spin glass is more complicated:
 - Infinitely many levels: the tree branches out for any value of $q < q_M$.
 - Infinite number of branches at any level.
- Amusing analogy with QED: number of emitted photons diverges at low energy \leftrightarrow infinite number of branches as we consider small values of w .
- However, it is possible to compute analytically the distribution of the W_i , at any level in q :

$$P(W; q) = \frac{W^{x(q)-1}(1-W)^{x(q)-1}}{\Gamma(1-x(q))\Gamma(x(q))} .$$

The tree of states (II)

- The real tree of states of a mean-field spin glass is more complicated:
 - Infinitely many levels: the tree branches out for any value of $q < q_M$.
 - Infinite number of branches at any level.
- Amusing analogy with QED: number of emitted photons diverges at low energy \leftrightarrow infinite number of branches as we consider small values of w .
- However, it is possible to compute analytically the distribution of the W_i , at any level in q :

$$P(W; q) = \frac{W^{x(q)-1}(1-W)^{x(q)-1}}{\Gamma(1-x(q))\Gamma(x(q))}.$$

- $\sum_i W_i = 1 \implies$ the weights are not independent:

$$P(W, W'; q) = x(q) \frac{\Theta(1-W-W')(WW')^{x(q)-1}(1-W-W')^{2x(q)-1}}{\Gamma(1-x(q))\Gamma(1-x(q))\Gamma(2x(q))}.$$

The tree of states (II)

- The real tree of states of a mean-field spin glass is more complicated:
 - Infinitely many levels: the tree branches out for any value of $q < q_M$.
 - Infinite number of branches at any level.
- Amusing analogy with QED: number of emitted photons diverges at low energy \leftrightarrow infinite number of branches as we consider small values of w .
- However, it is possible to compute analytically the distribution of the W_i , at any level in q :

$$P(W; q) = \frac{W^{x(q)-1}(1-W)^{x(q)-1}}{\Gamma(1-x(q))\Gamma(x(q))}.$$

- $\sum_i W_i = 1 \implies$ the weights are not independent:

$$P(W, W'; q) = x(q) \frac{\Theta(1-W-W')(WW')^{x(q)-1}(1-W-W')^{2x(q)-1}}{\Gamma(1-x(q))\Gamma(1-x(q))\Gamma(2x(q))}.$$

- We want to generate explicit realisations of this tree.

Free-energy fluctuations

- The W_i are cumbersome to handle.
- We consider instead the **free-energy fluctuations**.

Free-energy fluctuations

- The W_i are cumbersome to handle.
- We consider instead the **free-energy fluctuations**.
- The free energy per spin of each state has fluctuations $f_\alpha/N = \mathcal{O}(1/N)$.

Free-energy fluctuations

- The W_I are cumbersome to handle.
- We consider instead the **free-energy fluctuations**.
- The free energy per spin of each state has fluctuations $f_\alpha/N = \mathcal{O}(1/N)$.
- The f_α are related to the w_α :

$$w_\alpha = \frac{e^{-\beta f_\alpha}}{\sum_\gamma e^{-\beta f_\gamma}} .$$

Free-energy fluctuations

- The W_I are cumbersome to handle.
- We consider instead the **free-energy fluctuations**.
- The free energy per spin of each state has fluctuations $f_\alpha/N = \mathcal{O}(1/N)$.
- The f_α are related to the w_α :

$$w_\alpha = \frac{e^{-\beta f_\alpha}}{\sum_\gamma e^{-\beta f_\gamma}} .$$

- However, it turns out that the f_α are **independent**:

$$\mathcal{P}(f_\alpha) \propto e^{-\beta x(q_M) f}$$

Free-energy fluctuations

- The W_I are cumbersome to handle.
- We consider instead the **free-energy fluctuations**.
- The free energy per spin of each state has fluctuations $f_\alpha/N = \mathcal{O}(1/N)$.
- The f_α are related to the w_α :

$$w_\alpha = \frac{e^{-\beta f_\alpha}}{\sum_\gamma e^{-\beta f_\gamma}} .$$

- However, it turns out that the f_α are **independent**:

$$\mathcal{P}(f_\alpha) \propto e^{-\beta x(q_M) f}$$

- We can do this at any level of q :

$$W_I = \frac{e^{-\beta f_I}}{\sum_J e^{-\beta f_J}}, \quad \mathcal{P}_q(f) \propto e^{-\beta x(q) f} .$$

Free-energy fluctuations

- The W_I are cumbersome to handle.
- We consider instead the **free-energy fluctuations**.
- The free energy per spin of each state has fluctuations $f_\alpha/N = \mathcal{O}(1/N)$.
- The f_α are related to the w_α :

$$w_\alpha = \frac{e^{-\beta f_\alpha}}{\sum_\gamma e^{-\beta f_\gamma}} .$$

- However, it turns out that the f_α are **independent**:

$$\mathcal{P}(f_\alpha) \propto e^{-\beta x(q_M) f}$$

- We can do this at any level of q :

$$W_I = \frac{e^{-\beta f_I}}{\sum_J e^{-\beta f_J}}, \quad \mathcal{P}_q(f) \propto e^{-\beta x(q) f} .$$

- Universality: everything is encoded in $x(q)$.

Generating the tree from the ground up (I)

- The formulation of the f_l considering a single q level in isolation is convenient for analytical computations.
- Here, we need a different approach.
- We will generate the tree step by step, beginning by the root and down to the states, at each step computing the weights.

Generating the tree from the ground up (I)

- The formulation of the f_l considering a single q level in isolation is convenient for analytical computations.
- Here, we need a different approach.
- We will generate the tree step by step, beginning by the root and down to the states, at each step computing the weights.
- We start considering a discretised $q(x)$ (equivalently, a finite number K of RSB steps).

Generating the tree from the ground up (I)

- The formulation of the f_l considering a single q level in isolation is convenient for analytical computations.
- Here, we need a different approach.
- We will generate the tree step by step, beginning by the root and down to the states, at each step computing the weights.
- We start considering a discretised $q(x)$ (equivalently, a finite number K of RSB steps).
- We have, therefore, $K + 1$ levels.

Generating the tree from the ground up (I)

- The formulation of the f_l considering a single q level in isolation is convenient for analytical computations.
- Here, we need a different approach.
- We will generate the tree step by step, beginning by the root and down to the states, at each step computing the weights.
- We start considering a discretised $q(x)$ (equivalently, a finite number K of RSB steps).
- We have, therefore, $K + 1$ levels.
- In addition, since we cannot handle an infinite number of states, we will 'prune' the tree.

Generating the tree from the ground up (I)

- The formulation of the f_l considering a single q level in isolation is convenient for analytical computations.
- Here, we need a different approach.
- We will generate the tree step by step, beginning by the root and down to the states, at each step computing the weights.
- We start considering a discretised $q(x)$ (equivalently, a finite number K of RSB steps).
- We have, therefore, $K + 1$ levels.
- In addition, since we cannot handle an infinite number of states, we will 'prune' the tree.
- We eliminate at each level all the clusters with $W_l < \epsilon$ (equivalent to neglecting all the states with $w_\alpha < \epsilon$).
- We are losing a total probability of $\sim \epsilon^{1-x(q_M)}$.

Generating the tree from the ground up (II)

- Given a cluster at level q_i , we want its subclusters at level q_{i+1} .

Generating the tree from the ground up (II)

- Given a cluster at level q_i , we want its subclusters at level q_{i+1} .
- We introduce M variables f_1, \dots, f_M . They are **not independent**:

$$\mathcal{P}_{q_i \rightarrow q_{i+1}}(f_1, \dots, f_M) \propto \exp \left[-\beta x(q_{i+1}) \sum_{i=1}^M f_i \right] \left[\sum_{i=1}^M \exp(-\beta f_i) \right]^{x(q_i)} .$$

Generating the tree from the ground up (II)

- Given a cluster at level q_i , we want its subclusters at level q_{i+1} .
- We introduce M variables f_1, \dots, f_M . They are **not independent**:

$$\mathcal{P}_{q_i \rightarrow q_{i+1}}(f_1, \dots, f_M) \propto \exp\left[-\beta x(q_{i+1}) \sum_{i=1}^M f_i\right] \left[\sum_{i=1}^M \exp(-\beta f_i)\right]^{x(q_i)}.$$

- Now the weights of the subclusters are (W = weight of the cluster at q_i)

$$w_i = W \frac{\exp(-\beta f_i)}{\sum_{i=1}^M \exp(-\beta f_i)}.$$

Generating the tree from the ground up (II)

- Given a cluster at level q_i , we want its subclusters at level q_{i+1} .
- We introduce M variables f_1, \dots, f_M . They are **not independent**:

$$\mathcal{P}_{q_i \rightarrow q_{i+1}}(f_1, \dots, f_M) \propto \exp\left[-\beta x(q_{i+1}) \sum_{i=1}^M f_i\right] \left[\sum_{i=1}^M \exp(-\beta f_i)\right]^{x(q_i)}.$$

- Now the weights of the subclusters are (W = weight of the cluster at q_i)

$$w_i = W \frac{\exp(-\beta f_i)}{\sum_{i=1}^M \exp(-\beta f_i)}.$$

- It is not immediately obvious, but this method generates the same probability distributions for the f_α .
- This is because the correlation in the f_i of the subclusters at level q_{i+1} compensates the correlations of the weights of the clusters at q_i .

Generating the tree from the ground up (II)

- Given a cluster at level q_i , we want its subclusters at level q_{i+1} .
- We introduce M variables f_1, \dots, f_M . They are **not independent**:

$$\mathcal{P}_{q_i \rightarrow q_{i+1}}(f_1, \dots, f_M) \propto \exp\left[-\beta x(q_{i+1}) \sum_{i=1}^M f_i\right] \left[\sum_{i=1}^M \exp(-\beta f_i)\right]^{x(q_i)}.$$

- Now the weights of the subclusters are (W = weight of the cluster at q_i)

$$w_i = W \frac{\exp(-\beta f_i)}{\sum_{i=1}^M \exp(-\beta f_i)}.$$

- It is not immediately obvious, but this method generates the same probability distributions for the f_α .
- This is because the correlation in the f_i of the subclusters at level q_{i+1} compensates the correlations of the weights of the clusters at q_i .
- Notice that the first step, the root, is going from $q = 0 \rightarrow q_0$. Since $x(q = 0) = 0$, in the first step we have independent f_i .

The cavity method

- We have seen how to generate the tree, we now want to compute physical quantities from it. We consider a **cavity approach**.
- We add an $N + 1$ spin s_0 to a system with N spins s_j .
- The average properties of s_0 are the same as those of the s_j .
- Consider the N -spin system in equilibrium and compute the cavity field

$$h = \sum_{k=1}^N J_{0k} s_k$$

The cavity method

- We have seen how to generate the tree, we now want to compute physical quantities from it. We consider a **cavity approach**.
- We add an $N + 1$ spin s_0 to a system with N spins s_j .
- The average properties of s_0 are the same as those of the s_j .
- Consider the N -spin system in equilibrium and compute the cavity field

$$h = \sum_{k=1}^N J_{0k} s_k$$

- The cavity field h_α for each state is the sum of $K + 1$ Gaussian random variables:

$$h_\alpha = h_\alpha^0 + h_\alpha^1 + \dots + h_\alpha^K,$$

- h_α^0 has variance βq_0 and is common to the whole tree.
- h_α^i has variance $\beta(q_i - q_{i-1})$ and is common to all the states down the same branch.

The cavity method

- We have seen how to generate the tree, we now want to compute physical quantities from it. We consider a **cavity approach**.
- We add an $N + 1$ spin s_0 to a system with N spins s_j .
- The average properties of s_0 are the same as those of the s_j .
- Consider the N -spin system in equilibrium and compute the cavity field

$$h = \sum_{k=1}^N J_{0k} s_k$$

- The cavity field h_α for each state is the sum of $K + 1$ Gaussian random variables:

$$h_\alpha = h_\alpha^0 + h_\alpha^1 + \dots + h_\alpha^K,$$

- Example: $m_\alpha = \tanh(\beta h_\alpha)$.

The cavity method

- We have seen how to generate the tree, we now want to compute physical quantities from it. We consider a **cavity approach**.
- We add an $N + 1$ spin s_0 to a system with N spins s_j .
- The average properties of s_0 are the same as those of the s_j .
- Consider the N -spin system in equilibrium and compute the cavity field

$$h = \sum_{k=1}^N J_{0k} s_k$$

- The cavity field h_α for each state is the sum of $K + 1$ Gaussian random variables:

$$h_\alpha = h_\alpha^0 + h_\alpha^1 + \dots + h_\alpha^K,$$

- Example: $m_\alpha = \tanh(\beta h_\alpha)$.
- The cavity step shifts the free energies:

$$w'_\alpha \sim w_\alpha \cosh(\beta h_\alpha)$$

⇒ we have an iterative method to refine the tree

Testing the program

- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).

Testing the program

- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).
- As a consistency check of the program, we can try to find $q(x)$ at a given T , starting from an educated guess.

Testing the program

- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).
- As a consistency check of the program, we can try to find $q(x)$ at a given T , starting from an educated guess.
- We consider $T = 0.85$, close to the critical point ($T_c = 1$),

Testing the program

- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).
- As a consistency check of the program, we can try to find $q(x)$ at a given T , starting from an educated guess.
- We consider $T = 0.85$, close to the critical point ($T_c = 1$),
- In these conditions, $q(x)$ is linear with a very good approximation.

Testing the program

- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).
- As a consistency check of the program, we can try to find $q(x)$ at a given T , starting from an educated guess.
- We consider $T = 0.85$, close to the critical point ($T_c = 1$),
- In these conditions, $q(x)$ is linear with a very good approximation.
- We only need to find two parameters: q_M and x_M .

Testing the program

- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).
- As a consistency check of the program, we can try to find $q(x)$ at a given T , starting from an educated guess.
- We consider $T = 0.85$, close to the critical point ($T_c = 1$),
- In these conditions, $q(x)$ is linear with a very good approximation.
- We only need to find two parameters: q_M and x_M .
 - 1 Find the correct q_M for a fixed x_M and compute $F(x_M)$.
 - 2 Minimize $F(x_m)$ to find the correct x_M .

Testing the program

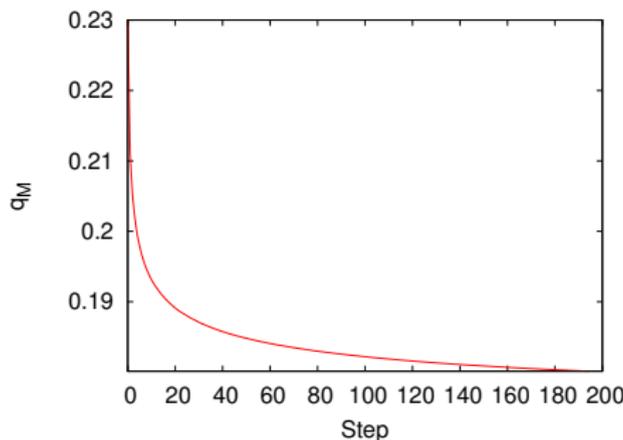
- We have described how to generate the tree given a $q(x)$.
- Also, how to refine the w_α with a cavity step (and with them get a new $q'(x)$).
- As a consistency check of the program, we can try to find $q(x)$ at a given T , starting from an educated guess.
- We consider $T = 0.85$, close to the critical point ($T_c = 1$),
- In these conditions, $q(x)$ is linear with a very good approximation.
- We only need to find two parameters: q_M and x_M .
 - 1 Find the correct q_M for a fixed x_M and compute $F(x_M)$.
 - 2 Minimize $F(x_m)$ to find the correct x_M .
- Let us study the solution at x_M fixed to the known correct value $x_M \approx 0.233122$.

$q(x)$ at $T = 0.85$

- We start with $q_M = 0.23 \approx x_M$ and iterate.
- We consider $\epsilon = 10^{-5} \implies \sim 1 - \epsilon^{1-x_M} = 99.99\%$ of the probability.
- $K = 20$ is enough, since $q(x)$ is so simple.
- We generate 10^6 samples (trees) and iterate
- Each iteration takes ≈ 2 min in a single CPU.

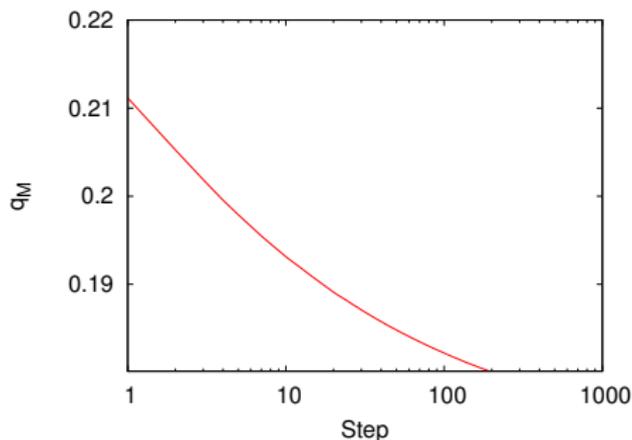
$q(x)$ at $T = 0.85$

- We start with $q_M = 0.23 \approx x_M$ and iterate.
- We consider $\epsilon = 10^{-5} \implies \sim 1 - \epsilon^{1-x_M} = 99.99\%$ of the probability.
- $K = 20$ is enough, since $q(x)$ is so simple.
- We generate 10^6 samples (trees) and iterate
- Each iteration takes ≈ 2 min in a single CPU.



$q(x)$ at $T = 0.85$

- We start with $q_M = 0.23 \approx x_M$ and iterate.
- We consider $\epsilon = 10^{-5} \implies \sim 1 - \epsilon^{1-x_M} = 99.99\%$ of the probability.
- $K = 20$ is enough, since $q(x)$ is so simple.
- We generate 10^6 samples (trees) and iterate
- Each iteration takes ≈ 2 min in a single CPU.



The convergence is very slow (logarithmic).

$q(x)$ at $T = 0.85$

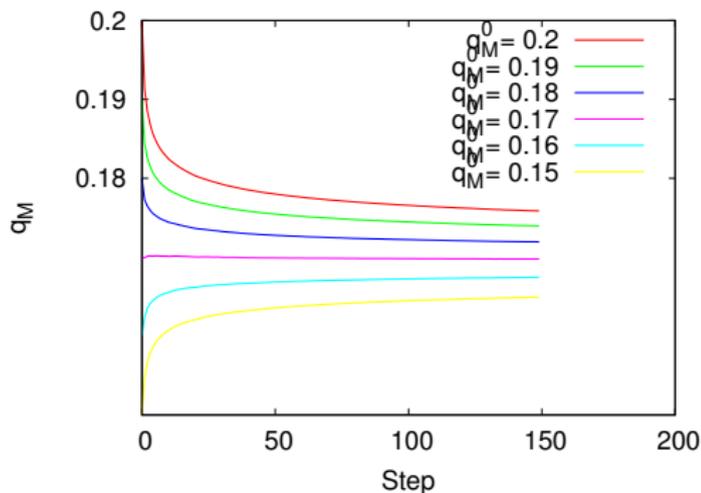
- We need a different approach.

$q(x)$ at $T = 0.85$

- We need a different approach.
- Start several computations, sweeping a wide range of starting q_M :

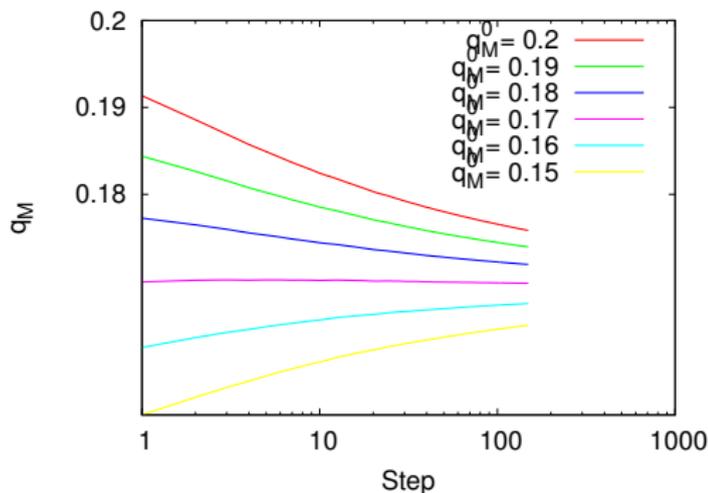
$q(x)$ at $T = 0.85$

- We need a different approach.
- Start several computations, sweeping a wide range of starting q_M :



$q(x)$ at $T = 0.85$

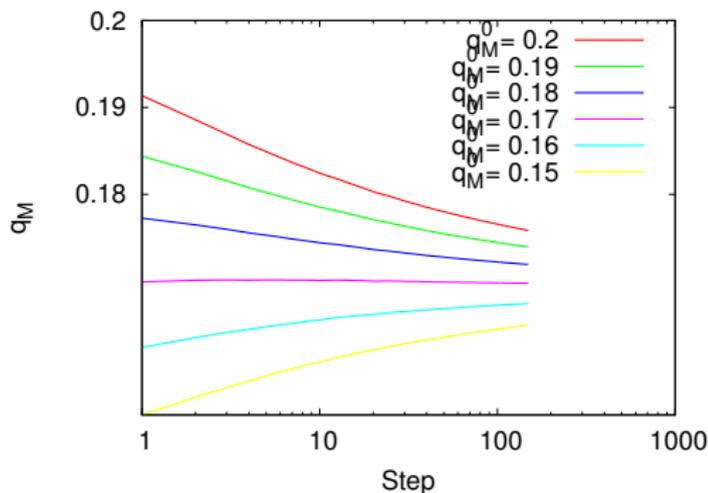
- We need a different approach.
- Start several computations, sweeping a wide range of starting q_M :



- The evolution is monotonic \implies it is easy to find the stable solution.

$q(x)$ at $T = 0.85$

- We need a different approach.
- Start several computations, sweeping a wide range of starting q_M :



- The evolution is monotonic \implies it is easy to find the stable solution.
- After 100 steps, starting with $q_M^{(0)} = 0.17$, we find $q_M^{(100)} = 0.169687(3)$ (expected value $q_M \approx 0.169691$).

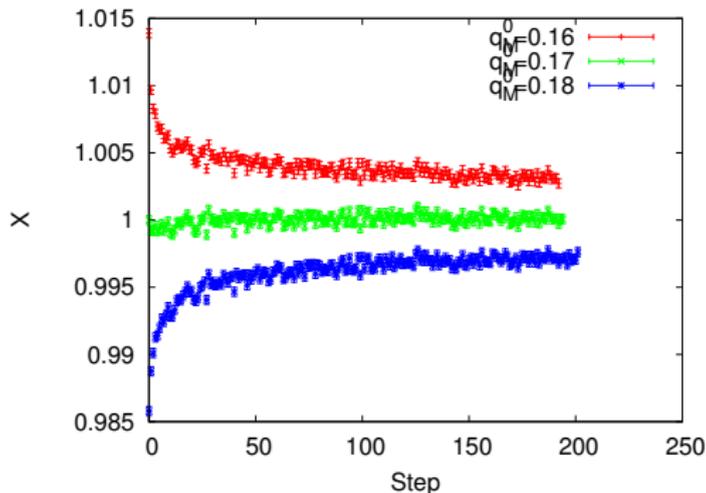
- The spin-glass susceptibility is $\chi_{\text{SG}} = \overline{((1 - m_0^2)^2)^2} / [1 - \beta^2 \overline{(1 - m_0^2)^2}]$.
- It diverges for $T < T_c$ so

$$X = \beta^2 \overline{(1 - m_0^2)^2} = \beta^2 \sum_{\alpha} \overline{w_{\alpha} (1 - m_{\alpha}^2)^2} = 1.$$

Replicon

- The spin-glass susceptibility is $\chi_{\text{SG}} = \overline{((1 - m_0^2)^2)} / [1 - \beta^2 \overline{(1 - m_0^2)^2}]$.
- It diverges for $T < T_c$ so

$$X = \beta^2 \overline{(1 - m_0^2)^2} = \beta^2 \sum_{\alpha} \overline{w_{\alpha} (1 - m_{\alpha}^2)^2} = 1.$$



- We obtain $X = 0.99972(33)$.

Conclusions

- We have reviewed the main properties of the tree of states in mean-field spin glasses.
- We have shown how to generate explicit realisations of this tree, in a self-consistent way.
- The states computed from the tree can be used to study physical quantities.
- Applications
 - Evaluate all the correlation functions of the model for fixed q .
 - Study other mean-field models (such as the full-RSB solution for the spin glass in a Bethe lattice).
 - Finite-size effects: generate the $P_{\mathcal{J}}(q)$ for single trees (single samples) and study the smoothing of the individual peaks.

Conclusions

- We have reviewed the main properties of the tree of states in mean-field spin glasses.
- We have shown how to generate explicit realisations of this tree, in a self-consistent way.
- The states computed from the tree can be used to study physical quantities.
- Applications
 - Evaluate all the correlation functions of the model for fixed q .
 - Study other mean-field models (such as the full-RSB solution for the spin glass in a Bethe lattice).
 - Finite-size effects: generate the $P_{\mathcal{J}}(q)$ for single trees (single samples) and study the smoothing of the individual peaks.

THANK YOU!