What makes slow samples slow in the Sherrington–Kirkpatrick model?

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Abstract

Using results of a Monte Carlo simulation of the Sherrington–Kirkpatrick model, we try to characterize the slow disorder samples; namely we visually analyse the correlation between the relaxation time for a given disorder sample $J$ with several observables of the system for the same disorder sample. For temperatures below $T_c$ but not too low, fast samples (small relaxation times) are clearly correlated with a small value of the largest eigenvalue of the coupling matrix, a large value of the site-averaged local field probability distribution at the origin, or a small value of the squared overlap $\langle q^2 \rangle$. Within our limited data, the correlation remains as the system size increases but becomes less clear as the temperature is decreased (the correlation with $\langle q^2 \rangle$ is more robust). There is a strong correlation between the values of the relaxation time for two distinct values of the temperature, but this correlation decreases as the system size is increased. This may indicate the onset of temperature chaos.

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(Some figures in this article are in colour only in the electronic version)
mean and $\psi = 1/3$, although it has been argued [4, 5] that $\psi$ may be slightly less than 1/3 (in such a case, ln $\tau$ would be weakly self-averaging).

Whether $\psi$ is equal to 1/3 or slightly less than 1/3, there are definitely disorder samples with extremely slow dynamics. Our aim in this paper is to try to characterize these ‘slow samples’. We have two motivations in mind, one theoretical and the other more down-to-earth. The first is the question of the behaviour of $F(x)$ for large values of the argument $x$. It has been argued in [4] that $F(x)$ has an exponential behaviour with $\ln F(x) \propto -x^0$ for large $x$. If furthermore the tail of $F(x)$ is dominated by rare disorder samples with small probability $\propto \exp(-AN^{\alpha})$ and an anomalously large free-energy barrier $\propto N^\beta$ (with $\beta > 1/3$), the exponents $\alpha$, $\beta$, $\psi$ and $\eta$ fulfil [4] the consistency relation $(\beta - \psi)\eta = \alpha$. For example, a dominance by the rare samples with all exchange coupling $J_{i,j}$ positive (up to a gauge transformation, see later) would correspond to $\alpha = 2$, and $\beta = 3/2$. The value $\psi = 1/3$ would then imply that $\eta = 12/7$. This value for $\eta$ is however not compatible with the numerical results of [4] for the distribution $P(\log \tau)$. In this reference, arguments are given for the values $\alpha = \beta = 1$, $\psi = 1/4$ (and accordingly $\eta = 4/3$) instead. The second motivation for the characterization of the slow samples is of practical matter for Monte Carlo simulations. The vast majority of Monte Carlo simulations of disordered systems in the literature use the same number of iterations for all disorder samples [1]. There is accordingly a danger that some rare slow disorder samples are not thermalized. One is led naturally to the idea of concentrating the computational effort on the hard samples that may require orders of magnitude more iterations than the mainstream disorder samples. Since the measurement of the relaxation time for every disorder sample is very time consuming, any heuristic method to pinpoint the slow samples can be valuable.

The numerical method used is similar to the one used in [3]. We consider the SK model with binary exchange couplings, and the Hamiltonian

$$\mathcal{H} = -\frac{1}{2} \sum_{i \neq j} J_{i,j} \sigma_i \sigma_j$$

with $\sigma_i = \pm 1$ and $J_{i,j} = \pm 1/\sqrt{N}$. We measure the dynamic overlap

$$q_d(t) = \frac{1}{N} \sum_{i=1}^{N} \sigma_i(t_0) \sigma_i(t + t_0)$$

averaged along a very long trajectory (namely averaging over many values of $t_0$), with the Metropolis dynamics, starting from a well-equilibrated spin configuration (obtained with the parallel tempering algorithm). Obviously, $q_d(0) = 1$, and $q_d(t)$ decreases continuously towards zero, as $t$ grows. We define the relaxation time $\tau$ by the equation $q_d(t = \tau) = 1/2\sqrt{E(\langle q^2 \rangle)}$, where $q$ is the usual ($J$-dependent) overlap between two replica. Here, we depart from [3] where a definition involving $\langle q^2 \rangle$, with no disorder average, was used. Indeed, since we are looking for correlations between the value of ln $\tau$ and other $J$-dependent quantities (including $\langle q^2 \rangle$ itself), it seems more appropriate to define $\tau$ using a sample-independent condition. The parameters of the simulations are such that our estimates of ln $\tau$ have negligible thermal noise, as compared to the disorder sample to disorder sample fluctuations (see [3] for details). That the thermal noise is tamed is an essential condition for our analysis. We have data for $N = 64$–512, 1024 disorder samples, and temperatures $T = 0.4, 0.5, \ldots$ (The critical temperature is $T_c = 1$.)

Note that as in [3], we do not measure relaxation times larger than some $t_{\text{window}} \approx 3.7 \times 10^6$ (and $t_{\text{window}} \approx 1.7 \times 10^8$ for $N = 512$.). For a given couple $N$ and $T$, the disorder samples

1 For counter examples where the CPU time is adjusted to the disorder sample sluggishness see [6] or [7].
with relaxation time larger than $t_{\text{window}}$ are thus 'censored'. By convention, $\ln \tau = -1$ for such samples, namely $-1$ means overflow. Note that the real $\ln \tau$ is never equal to $-1$ (our relaxation times are integers).

As previously explained, we found it more proper to define the relaxation time using a condition involving the disorder averaged $E(\langle q^2 \rangle)$ rather than $\langle q^2 \rangle$ as was done in [3]. It turns out however that this makes little difference, as shown in figure 1, where we compare the two definitions in the case $N = 128$ and $T = 0.4$. The same conclusion holds for other couples of $N$ and $T$. On closer inspection, one notes that for low values of $\ln \tau$, the definition used here gives systematically lower results than the one used in [3]. This is explained by the fact that low values of $\ln \tau$ are strongly correlated to low values of $\langle q^2 \rangle$ (see later in the text and figure 5). This small systematic difference between the two definitions of the relaxation time disappears when $\ln \tau$ grows.

The first question we would like to address is whether the tail of the distribution $P(\ln \tau)$ is dominated by ferromagnetic disorder configurations. In order to proceed, we need a measure of the ferromagnetic character of a disorder configuration. The sum $\sum_{i \neq j} J_{i,j}$ is not a suitable indicator since it is not invariant under the local gauge symmetry of the model $\sigma_i \rightarrow \epsilon_i \sigma_i$, $J_{i,j} \rightarrow \epsilon_i \epsilon_j J_{i,j}$, with $\epsilon_i = \pm 1$, whereas the dynamics is invariant under this symmetry. Said another way, for every disorder configuration with all $J_{i,j} > 0$, there is a huge number of $2^N - 1$ other gauge-transformed configurations with the same value of $\ln \tau$ (and the same weight) but a different value of $\sum_{i \neq j} J_{i,j}$, and any correlation is washed out. (We nevertheless checked that there is indeed no sign of correlations between $\ln \tau$ and $\sum_{i \neq j} J_{i,j}$ in our data.) A better indicator of the ferromagnetic character of a disorder configuration, which has been proposed in [5], is the largest eigenvalue $\lambda_N$ of the matrix $\{J_{i,j}\}$. In the SK model, the diagonal elements of this matrix are not used, but they are obviously needed in order to compute the eigenvalues of the matrix, and we have set them equal to zero. Our results for the correlation between $\ln \tau$ and $\lambda_N$ can be found in figure 2 for $N = 64$ and $T = 0.8$. The points on the $x$-axis are concentrated around the value 2, in agreement with random matrix results for the Gaussian orthogonal ensemble (GOE) with the normalization $E(J_{i,i}^2) = 1/N$ for $i \neq j$. (In the GOE the diagonal elements are random with $E(J_{i,i}^2) = 2/N$, and not identically zero, but...
Figure 2. The logarithm of the relaxation time as a function of the largest eigenvalue $\lambda_N$ of the $\{J_{i,j}\}$ matrix for $N = 64$ and $T = 0.8$. There are no disorder samples with $\tau > t_{\text{window}}$ in the data.

This should not change the asymptotic behaviour.) There is a clear correlation between low values of $\lambda_N$ and low values for $\ln \tau$, but this correlation becomes fuzzy as $\lambda_N$ grows. We were looking for a characterization of the slow disorder samples but we found a characterization of the fast disorder samples instead.

The scatter plots become progressively harder to interpret as $N$ grows and/or $T$ decreases, since we are missing more and more points that correspond to censored values of $\ln \tau$. One can nevertheless conclude from our data that the correlation remains as the system size increases, but becomes more fuzzy as temperature is decreased. (The neat correlation seen on the left of figure 2 fades away as the temperature is decreased.) The net conclusion is that, in our data, there is no visible dominance of the large relaxation region by ferromagnetic disorder samples. This is in agreement with the findings of [4] for the tail of the distribution of $\ln \tau$.

In the spherical SK model [8], the relaxation time is fixed by the difference between the two largest eigenvalues of the $\{J_{i,j}\}$ matrix, through the equation

$$\ln \tau = N/(2T)(1 - T)(\lambda_N - \lambda_{N-1}).$$

For the SK model with binary couplings that we consider here, the correlation between $\ln \tau$ and $\lambda_N - \lambda_{N-1}$ is more fuzzy, as one can see in figure 3. It is in fact fuzzier than the relation between $\ln \tau$ and $\lambda_N$ alone. This is true for all the considered values of $N$ and $T$, and for all values of $\ln \tau$ (both large and small). Note that in the spherical SK model, equation (3) is used to prove that $E(\ln \tau) \propto N^{1/3}$. It turns out [2] that this scaling also holds in the usual SK model, although equation (3) does not hold.

In the spirit of [4], we have also looked at the correlation between $\ln \tau$ and the site-averaged local field probability distribution $P(h_{\text{local}})$:

$$P(h_{\text{local}}) = \frac{1}{N} \left\langle \sum_i \delta(\bar{h} - \sum_{j \neq i} J_{i,j} \sigma_j) \right\rangle.$$  

Although one may always argue that truly ferromagnetic disorder samples correspond to $\lambda_N$ of order $\sqrt{N}$, deep in the tail of the distribution that we do not see due to limited statistics.

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Figure 3. The logarithm of the relaxation time as a function of the gap between the two largest eigenvalues of the $\{J_{i,j}\}$ matrix for $N = 64$ and $T = 0.8$. There are no disorder samples with $\tau > t_{\text{window}}$ in the data.

Figure 4. The logarithm of the relaxation time as a function of $P(h_{\text{local}} = 0)$ for $N = 64$ and $T = 0.8$. The normalization is such that $\int_0^\infty dh_{\text{local}} P(h_{\text{local}}) = 1$. There are no disorder samples with $\tau > t_{\text{window}}$ in the data.

Specifically, we looked at the correlation between $\ln \tau$ and the value of $P(h_{\text{local}})$ at the origin. Our results can be found in figure 4 for $N = 64$ and $T = 0.8$. In [4], it was argued that the disorder samples with large relaxation times are the ones with all local fields (on every site $i$) large. This is indicative of a correlation between large values of $\ln \tau$ and a distribution $P(h_{\text{local}})$ that is depleted at the origin. (Indeed in [10], $\ln \tau$ increases as $T$ decreases, while $P(h_{\text{local}} = 0)$ decreases as $T$ decreases, with $P(h_{\text{local}} = 0) = 0$ at zero temperature in the $N = \infty$ limit.) Our data indeed show a neat correlation between large values of $P(h_{\text{local}} = 0)$, and small $\ln \tau$, ...
Figure 5. The logarithm of the relaxation time as a function of \( \langle q^2 \rangle \) for \( N = 64 \) and \( T = 0.8 \). There are no disorder samples with \( \tau > t_{\text{window}} \) in the data.

but this correlation becomes fuzzy for lower values of \( P(h_{\text{local}} = 0) \). Again we were looking for a characterization of the slow disorder samples but we found a characterization of the fast disorder samples instead. The observed correlation seems to persist as \( N \) grows, but fades as \( T \) is decreased. There is a similar correlation between \( \ln \tau \) and the average value of \( h_{\text{local}} \) within the tail of \( P(h_{\text{local}}) \), for example within the last 5% of the distribution. A small value of \( h_{\text{local}} \) inside the tail is correlated to a small relaxation time.

The strongest correlations we found are between \( \ln \tau \) and the average overlap squared \( \langle q^2 \rangle \), as can be seen in figure 5. The observed correlation seems to persist as \( N \) grows, and as \( T \) is decreased. This correlation is an empirical finding and we have no dynamical explanation for it. On the other hand, \( \langle q^2 \rangle \) is fairly easy to estimate with Monte Carlo, even with little statistics, and this correlation could be used to flag slow samples in numerical simulations. We finally remark that if there is a correlation between \( \ln \tau \) and the average overlap squared \( \langle q^2 \rangle \), then there is no correlation between \( \ln \tau \) and the number of peaks in the order parameter distribution \( P(q) \). We have looked for such a correlation using the data of [9], where the number of peaks was estimated for a subset of the disorder samples considered here. The presence of thermal noise in the measured distribution makes it difficult to count the number of peaks. In this reference, the authors did their best to count the number of peaks by visual inspection of the plots of \( P(q) \) for 192 disorder samples, with \( T = 0.4 \). We find no correlation between the value of \( \ln \tau \) and the number of peaks for \( N = 64 \) and 256.

We have finally found that, disorder sample by disorder sample, the relaxation time measured at two temperatures (both in the spin glass phase) are strongly correlated. This can be seen in figure 6 where we show \( \ln \tau(T = 0.5) \) as a function of \( \ln \tau(T = 0.6) \) both for \( N = 64 \). This figure shows a very strong correlation. One notes a couple of samples for which \( \ln \tau(T = 0.5) = -1 \) (censored data). Figure 7 shows the same figure but with all systems’ sizes together. Obviously, there are more and more disorder samples censored as \( N \) grows, now with a whole horizontal segment with \( \ln \tau = -1 \). Interestingly, all points in the scatter plot scale on the same \( N \)-independent thick line. This thick line would extend further towards large values had we used a larger observation window. (But it would be quite CPU time consuming to obtain a large extension.) Note that the \( N = 512 \) data appear
to be more scattered than the other data; an optimistic interpretation is that we are seeing some onset of a temperature chaotic behaviour. In order to be more quantitative, we have analysed the data as follows. First, we made a linear least-squares fit of the $N = 64$ data to the form $\ln \tau(T = 0.5) = a + b \ln \tau(T = 0.6)$, with parameters $a$ and $b$ (we obtain the values $a = 0.195$ and $b = 1.297$). Then we computed, for every system size $N$, the deviation $\delta^2_N = 1/(N \sum_j (a + b \ln \tau(T = 0.6) - \ln \tau(T = 0.5))^2$, with the values of $a$ and $b$ obtained from the fit of the $N = 64$ data, a sum over those disorder samples such that both

Figure 6. Logarithm of the relaxation time for $T = 0.5$ as a function of the logarithm of the relaxation time for the same disorder sample but $T = 0.6$. The number of sites is $N = 64$. There are a couple of disorder samples with $\tau > t_{\text{window}}$ for $T = 0.5$ in this figure. Those are the points with $\ln \tau = -1$ (by convention).

Figure 7. Same as in figure 6 but with $N = 64, 128, 256$ and $512$ together. There are a couple of disorder samples with $\tau > t_{\text{window}}$ for both $T = 0.5$ and $T = 0.6$ in this figure. Those are the points with $\ln \tau = -1$ (by convention).
relaxation times are less than $t_{\text{window}}(N/512)^{1/3}$, with $t_{\text{window}}$ the cutoff used for the $N = 512$ data, and $N_J$ the number of disorder samples that satisfy the constraint. In other words, $\delta_N^2$ is the mean squared deviation from the linear-squares fit, properly taking into account the relaxation time observational cutoff. We find that $\delta_N/N^{1/3} = 0.0423, 0.0485, 0.0509$ and $0.0568$ for $N = 64, 128, 256$ and 512, respectively. The correlation between the values of $\ln \tau$ at two temperatures becomes looser as $N$ grows, confirming quantitatively the indication of the onset of temperature chaos.

In conclusion, we have measured the equilibrium relaxation time $\tau$ of the Sherrington–Kirkpatrick (SK) model with binary couplings for many samples of the quenched disorder, and several values of the temperature, with system sizes $N = 64–512$, taking great care that the thermal noise is negligible. We confirm the result of [4] that the slow samples are not correlated to ‘ferromagnetic’ disorder configurations, but we did not find evidence for a dominance by configurations with a small value at the origin of the site-averaged local field probability distribution. We find a strong correlation between the relaxation times measured at two distinct values of the temperature (with the same disorder sample). A closer look shows a broadening that is possibly an indication of the onset of temperature chaos.

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