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Percolation and Schramm-Loewner evolution in the 2D random-field Ising model

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ABSTRACT

The presence of random fields is well known to destroy ferromagnetic order in Ising systems in two dimensions. When the system is placed in a sufficiently strong external field, however, the size of clusters of like spins diverges. There is evidence that this percolation transition is in the universality class of standard site percolation. It has been claimed that, for small disorder, a similar percolation phenomenon also occurs in zero external field. Using exact algorithms, we study ground states of large samples and find little evidence for a transition at zero external field. Nevertheless, for sufficiently small random-field strengths, there is an extended region of the phase diagram, where finite samples are indistinguishable from a critical percolating system. In this regime we examine ground-state domain walls, finding strong evidence that they are conformally invariant and satisfy Schramm-Loewner evolution (SLE_k) with parameter $\kappa = 6$. These results add support to the hope that at least some aspects of systems with quenched disorder might be ultimately studied with the techniques of SLE and conformal field theory.

The random-field Ising model (RFIM) is one of the earliest studied and simplest disordered systems showing non-trivial and glassy behavior [1,2]. It has a number of important realizations in nature, including diluted antiferromagnets in a field and binary liquids in porous media [2]. Through its long history, researchers have managed to gain a reasonable understanding of the critical behavior, although this progress has been neither straight nor smooth, and many questions remain unanswered [2]. It is known, for example, that the RFIM in two dimensions (2D) lacks ferromagnetic order [3,4]. Even at zero temperature it remains in the paramagnetic state for non-zero disorder. Numerical ground-state calculations have shown, however, that even in the absence of a thermodynamic transition there exists a geometric transition at which the size of the spin clusters diverges in a manner bearing many similarities to classical site percolation [5,6]. While this transition is rather clearly established in the presence of an external field, it has been argued that a similar percolation phenomenon can also be observed in the absence of an external field for sufficiently small disorder [6,7]. Here, we re-investigate the zero-field behavior with large scale ground-state calculations, focusing on the possible percolation phenomenon.

The observed relations to classical site percolation at this (nonzero or zero field) geometrical transition motivate further questions of how far the similarities go. Interfaces in two-dimensional percolation satisfy Schramm-Loewner evolution (SLE) [8,9], but

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this property relies on conformal invariance which is not conserved in the presence of disorder. Recently, however, there have been suggestions that the domain walls of certain other disordered systems satisfy SLE. This intriguing possibility implies that conformal invariance, broken by disorder, is restored, at least at criticality where relevant length scales diverge.

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Schramm–Loewner evolution is a method for constructing a statistical ensemble of curves in the plane from one-dimensional Brownian motion, thus classifying curves with only one parameter, the diffusion constant κ [10]. Characteristic interfaces in many physical systems have been shown (in some cases rigorously) to satisfy SLE_{κ}. These include percolation ($\kappa = 6$), self avoiding walks ($\kappa = 4/3$), as well as spin cluster boundaries ($\kappa = 3$) and Fortuin–Kasteleyn cluster boundaries ($\kappa = 16/3$) in the Ising model. A number of numerical studies have found interfaces in certain disordered systems consistent with SLE, in particular the 2D Ising spin glass [11,12], the Potts model on dynamical triangulations [13], the random bond Potts model [14], and the disordered solid-on-solid model [15]. Here we extend this list to include the 2D random-field Ising model.

In Section 1 we introduce the RFIM and Schramm–Loewner evolution in more detail, and discuss the approach used here for determining ground states of large samples. Section 2 is devoted to an investigation of the critical behavior of the RFIM near the geometric transition, focusing on the behavior in zero external field for small disorder. In Section 3, we report the results of our tests of the correspondence between interfaces in the RFIM and Brownian motion implied by SLE.

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1. The model and method

We consider the random-field Ising model in two dimensions with Hamiltonian

$$\mathcal{H} = -J \sum_{\langle i,j \rangle} s_i s_j - \sum_i h_i s_i.$$
⁽¹⁾

Here, the spins $s_i = \pm 1$ are located on the sites of a square lattice and interact ferromagnetically with nearest neighbors. The local fields h_i are quenched random variables drawn from a normal distribution with mean H and standard deviation Δ . Since, at zero temperature, only the ratio J/Δ is relevant, we take J = 1 for simplicity. The spin-spin interaction J induces a correlation between the spins resulting in spin clusters which are compact up to a length ξ_b . Above this scale the clusters are fractal objects, the magnetization is zero and the system is paramagnetic. As the randomness Δ is decreased, the breakup length ξ_b increases. At and above three dimensions ξ_b diverges at the thermodynamic phase transition, below which the system is ferromagnetic. In two dimensions no thermodynamic phase transition exists, and ξ_b diverges only at $\Delta = 0$. It has been argued that, for H = 0, the breakup length scale ξ_b increases with decreasing Δ as [3]

$$\xi_b \sim e^{A/\Delta^2}.\tag{2}$$

Though there is no thermodynamic transition in 2D, the linear extent of the largest clusters diverges for sufficiently large *H*. In most aspects, this divergence appears to be consistent with standard site percolation [5–7]. It has been suggested that the divergence occurs even at H = 0 if Δ is below a critical value [5,7]. It is the characteristics of clusters of aligned spins and their boundaries at this geometric transition which we focus on in this study.

We restrict our investigation to ground-state spin configurations, which can be efficiently constructed through a mapping to the well-known minimum cut (or maximum flow) problem in graph theory [16,17]. Consider a directed graph with N + 2 vertices, and edges (i, j) furnished with weights c_{ij} . The minimum (s, t) cut is given by a subset of the edges of the graph, whose removal disconnects vertices *s* and *t*, such that the sum of the weights of the cut edges is minimal. Using the variables x_i , which are 1 if vertex *i* is connected to *s* and 0 otherwise, the total weight of a cut can be represented as

$$C(\{\mathbf{x}\}) = \sum_{i,j=1}^{N} x_i (1 - x_j) c_{ij}.$$
(3)

With the identification of x_i with spin variables (excepting the vertices s and t), and an appropriate choice of the weights c_{ij} , this function can be made to precisely match the RFIM Hamiltonian (1). The minimum cut separating the graph into vertices connected to s and vertices connected to t gives the minimum-energy way of cutting the RFIM lattice into clusters of up and down spins, and thus corresponds to the ground state of Eq. (1). The choice for the edge weights is, for $i \notin (s, t)$

$$c_{ij} = \begin{cases} 0, & i \ge j, \\ 4J, & \text{else.} \end{cases}$$
(4)

For the edges connecting the "spin" vertices to *s* and *t*, the result is given in terms of the quantity $u_i = -h_i - \frac{1}{2} \sum_j (c_{ij} - c_{ji})$ as

$$c_{si} = \begin{cases} 0, & u_i > 0, \\ -u_i, & \text{else}, \end{cases} \qquad c_{it} = \begin{cases} u_i, & u_i > 0, \\ 0, & \text{else}. \end{cases}$$
(5)

 c_{is} and c_{ti} are taken to be zero for all *i*. Here, we use a fast algorithm for solving the minimum cut problem based on the idea of

"augmenting paths" [18,19]. The worst case scenario for the running time of this algorithm is an unimpressive $O(N^3)$ (or more generally V^2E where V is the number of vertices and E is the number of edges), however, the algorithm was designed to optimize the typical case. The optimization was carried out for vision and image analysis problems, but even for the graph structure corresponding to the RFIM ground-state calculation the running time is proportional to N for the samples considered here. In practice, the maximum system size is limited more by computer memory constraints than by time.

Schramm–Loewner evolutions are defined in terms of a family g_t of conformal maps which take, formally, the upper half plane minus the curve γ_t (parametrized by "time" t) to the upper half plane, $g_t : \mathbb{H} \setminus \gamma_t \to \mathbb{H}$. This map can be defined (using complex notation and suitable normalization) in terms of the differential equation

$$\frac{\partial g_t(z)}{\partial t} = \frac{2}{g_t(z) - \xi_t},\tag{6}$$

where ξ_t is the driving function for the curve γ_t . If the curve (or the process generating the curve) satisfies SLE_{κ} , then ξ_t will be a Brownian motion with zero mean and variance κt . Numerically, rather than solving the differential equation, the map g_t is instead pictured as a series of maps g_i , which iteratively remove a small section (Δt_i) from the beginning of the curve. To calculate the driving function from a given curve, the incremental map g_i is approximated using a vertical slit map [20]

$$g_i(z) = i\sqrt{-(z-\xi_i)^2 - 4\Delta t_i + \xi_i}.$$
 (7)

The parameters ξ_i and Δt_i are determined from the coordinates of the curve segment to be removed through the relations $\xi_i = x_{i,i-1}$ and $\Delta t_i = y_{i,i-1}^2/4$. More specifically, $x_{i,i-1}$ and $y_{i,i-1}$ are the coordinates of the i'th segment of the curve after undergoing the i-1 successive maps $g_{i-1} \circ \cdots \circ g_1$. The parameter ξ_i is the value of the driving function ξ_t sampled at time $t_i = \sum_{j \leq i} \Delta t_j$. The complex square root in Eq. (7) is calculated, as usual, with the branch cut along the negative real axis.

This is an iterative process in which the coordinates of the interface are successively updated for each step, and thus the computational complexity is $O(L_I^2)$ where L_I is the length of the interface. If L_I grows like $\sim N^{d_f/2}$ (with $d_f = 7/4$ expected for percolation) in the number of spins *N*, the resulting computational complexity is $O(N^{7/4})$, which is significantly slower than the ground-state calculation which is O(N) on average. We use a fast implementation of this "zipper" algorithm [21], in which blocks of multiple slit maps are approximated using a Laurent series. Treating blocked maps in one step dramatically speeds up the calculation such that it scales on average as $L_I^{1.3}$. The loss in accuracy from the approximation is minimal.

If the driving function ξ_t can be shown to be Brownian motion then the curves satisfy SLE_{κ}. In practice, the finite size of the lattice and the zipper algorithm introduce correlations between the increments of ξ_i and in their associated time values t_i . In particular the distribution of time steps $t_i - t_{i-1}$ is non-trivial and has significant correlations. However, it should be emphasized that correlations in the times t_i at which ξ_t is sampled does not imply correlations in the underlying continuous driving function ξ_t .

2. Phase diagram and behavior at zero field

For very large disorder, $\Delta \gg Jz$, where *z* is the lattice coordination number, the interactions between the spins play no role and the system reduces, trivially, to the classical site percolation model. Each spin is determined solely by the independent random variable h_i . Identifying spin up with "site occupation", the site

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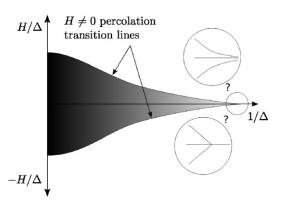


Fig. 1. Schematic representation of the dependence of the location H_c of the percolation transition on the random-field strength Δ . For $H < H_c$ the spin clusters have finite extent. For $H \ge H_c$ there exists at least one infinite cluster.

occupation probability *p* is simply the probability for h_i to be positive: $p = \frac{1}{2} \operatorname{erfc}\{-H/(\sqrt{2}\Delta)\}$. For smaller disorder the spin-spin interaction of the RFIM complicates the analogy with percolation. However, it has been demonstrated that there exists a line of critical external fields $H_c(\Delta)$, as pictured schematically in Fig. 1, for which observables like the crossing probability and the fractal dimension of the spin clusters maintain the percolation values [6].

For very large disorder the line of critical external fields is found from the critical site percolation probability p_c ($p_c \approx 0.5927$ for the square lattice)

$$H_c(\Delta \gg z) \approx -\Delta\sqrt{2} \operatorname{erfc}^{-1}\{2p_c\}.$$
(8)

For small disorder the behavior is less well understood. $H_c(\Delta)$ decreases as Δ decreases, approaching the limiting value $H_c(\Delta =$ 0) = 0. It has been claimed [6,7] that at a finite disorder strength, $\Delta_c \approx 1.65(5)$, the curve becomes identically zero $H_c(\Delta < \Delta_c) = 0$, cf. Fig. 1. We tested this claim by looking at the behavior of the system varying Δ at H = 0. We calculated the probability that a connected cluster of up spins crosses a rectangle of aspect ratio $e^{-2/5}$ in the vertical (shorter) direction (this aspect ratio is chosen to be different from 1/2, but the specific value is somewhat arbitrary). In the upper panel of Fig. 2, we present the results of these calculations. The crossing probability π for large Δ is zero (as expected for a square lattice with $p_c > 0.5$). As the disorder is decreased, π approaches the exact percolation value, which is indicated by the horizontal line in the plot. At very small disorder, when the breakup length scale ξ_b becomes comparable to the system size L, the system appears ferromagnetic and the crossing probability falls away from the plateau value.

To separate the approach to the plateau from the finite size effects we also considered another quantity, the probability π_{tot} that there exists a connected cluster of either up or down spins crossing either horizontally or vertically. The small disorder limit of this quantity is $\pi_{tot}(\Delta \rightarrow 0) = 1$ for both the onset of percolation and the ultimate ferromagnetic (and finite size) ordering. If there is a percolation transition at finite disorder ($\Delta_c > 0$), in the thermodynamic limit $L = \infty$ the curve $\pi_{tot}(\Delta)$ will be a step function. Hence, the curves for finite systems approaching this step function will intersect at Δ_c . As is apparent from our data for π_{tot} shown in the lower panel of Fig. 2, such a crossing is not observed at least down to $\Delta \approx 1.45$, significantly below the previously conjectured value of $\Delta_c \approx 1.65$. The study of even smaller disorder strengths, while insuring $L \gg \xi_b$, is preempted by the exponential growth of the breakup length ξ_b . Although it seems very unlikely that the system undergoes a true percolation phase transition for H = 0, there exists a large plateau region in which the system appears to be at critical percolation at H = 0, even up to very large system

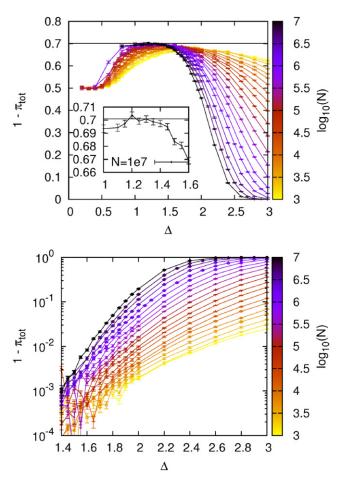


Fig. 2. Crossing probabilities for spin clusters of the 2D RFIM on rectangular domains of aspect ratio $e^{-2/5}$ for a number of system sizes $N = e^{-2/5}L^2$ at H = 0. Top panel: probability of crossing of clusters of up spins in the vertical direction. For Δ decreasing from infinity, the crossing probability approaches the exact percolation crossing value [22], indicated by the horizontal line. At very small disorder, the breakup length ξ_b is larger than the system size and the systems are effectively ferromagnetic. Inset: enlarged view of the plateau region for the largest system size. The bottom panel shows $1 - \pi_{tot}$, the total crossing probability. If there was a transition at H = 0, the lines would have to cross around the transition point.

sizes. Presumably, this holds true even for system sizes occurring in experimental realizations of the RFIM.

3. Schramm-Loewner evolution

Finally, we tested directly the conformal mapping upon which SLE is based. We studied the statistics of 10000 interfaces generated in a half disc. This geometry is used to optimally mimic the full half plane. The interface is initiated at the origin (the center of the flat edge) by two fixed spins and is considered ended when it touches the curved boundary. We find that the variance of the driving function calculated from the interfaces using the method described in Section 1 is $\hat{k} = \langle (\xi_t - \langle \xi_t \rangle)^2 \rangle / t = 6.09 \pm 0.09$, and the normalized mean is $\hat{\xi} = \langle \xi_t \rangle / \sqrt{\hat{k}t} = 0.01 \pm 0.01$. The agreement with SLE is good, but is only expected to be perfect at criticality. The difference from the expected percolation value $\kappa = 6$ which is perhaps just visible here is due to the calculations being carried out at $\Delta = 1.65$. The position distribution of the resulting stochastic process at several fixed times is shown in Fig. 3 along with curves representing the expectations for a perfect random walk.

We have focused here on the case of zero external field, where we have shown, using significantly larger system sizes than had been accessible before, that there is no percolation phase transition for H = 0, at least down to $\Delta = 1.45$. Due to the fact, however,

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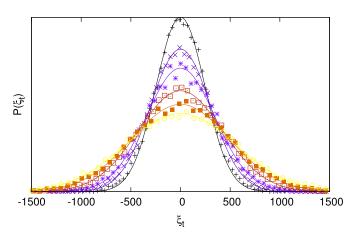


Fig. 3. Distribution of the random walk at several different "times" *t* as extracted from the RFIM interfaces at $\Delta = 1.65$. The lines are normal distributions with zero mean and variance κt with $\kappa = 6$.

that the percolation transition line comes exponentially close to H = 0 for small random-field strengths Δ (cf. Fig. 1), there exists a plateau region where even at H = 0 the behavior appears nearly indistinguishable from criticality. In this regime we have tested the applicability of Schramm–Loewner evolution to the RFIM and found good agreement. Further studies at non-zero external field and using an array of further tests for consistency with SLE confirm this result [23].

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