

Equilibrium spin-glass transition of magnetic dipoles with random anisotropy axes on a site-diluted lattice

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We study partially occupied lattice systems of classical magnetic dipoles which point along randomly oriented axes. Only dipolar interactions are taken into account. The aim of the model is to mimic collective effects in disordered assemblies of magnetic nanoparticles. From tempered Monte Carlo simulations, we obtain the following equilibrium results. The zero-temperature entropy approximately vanishes. Below a temperature T_c , given by $k_B T_c = (0.95 \pm 0.1)x\varepsilon_d$, where ε_d is a nearest-neighbor dipole-dipole interaction energy and x is the site occupancy rate, we find a spin-glass phase. In it, (1) the mean value $\langle|q|\rangle$, where q is the spin overlap, decreases algebraically with system size N as N increases, and (2) $\delta|q| \approx 0.5\langle|q|\rangle\sqrt{T/x}$, independently of N , where $\delta|q|$ is the root-mean-square deviation of $|q|$.

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I. INTRODUCTION

Magnetic dipole-dipole interactions play a fundamental role in magnetic phenomena. Their long-range nature gives rise to magnetic domains in ferromagnets. On the other hand, because of their weak strength, they usually play an insignificant role in determining the Curie temperatures and in critical phenomena of atomic crystals. In assemblies of magnetic nanoparticles (NPs), things can be very different. Because ferromagnetic NPs of up to thousands of Bohr magnetons have a single magnetic domain,¹ dipole-dipole interactions among such NPs can be very large and can thus dominate their collective behavior.² Often, both the orientation of the crystallites of which NPs are made as well as their positions are disordered in the assembly, and therefore behave much as a system of interacting magnetic dipoles with randomly oriented magnetic easy axes. These systems of random axes dipoles (RADs) are clearly frustrated, since two different dipoles give rise to magnetic fields at any given point which are not in general collinear. The sort of time-dependent behavior that is expected of spin glasses has been observed in experiments³⁻⁵ as well as in simulations⁶⁻¹⁰ of disordered assemblies of magnetic NPs. Because these systems evolve in time very slowly (exhibiting aging^{3,6} and other memory¹¹ effects), evidence for a *thermodynamic* spin-glass phase in them is more difficult to come by. One of us has recently given numerical evidence for the existence of an equilibrium spin-glass phase for a *fully* occupied lattice of dipoles with randomly oriented axes.¹² On the other hand, numerical evidence against such a phase has been given for *site-diluted* lattice systems of magnetic dipoles with *parallel* axes.¹³

One might expect site-diluted systems of RADs to behave as fully occupied ones, and therefore to have equilibrium spin-glass phases at low temperatures. For support of this expectation, consider rescaling distances in a dilute system of RADs. Because dipole-dipole interactions decay with distance r as r^{-3} , letting $r \rightarrow br$ merely redefines dipole-dipole interaction energies as $\varepsilon_d \rightarrow b^{-3}\varepsilon_d$. This would imply all physical quantities for systems of RADs in three dimensions

(3D) with different values of site concentration x collapse onto the same curve when plotted versus T/x , where T is the temperature. This argument holds for randomly located RADs on a *continuous* space in 3D. It is therefore expected to hold approximately for a site-diluted *lattice* if $x \ll 1$, but not necessarily for lattices with higher concentrations of RADs. This is why dipoles with parallel axes can have an antiferromagnetic phase at low temperature on a fully occupied simple cubic lattice,¹⁴ but appear to have no condensed phase of any sort on a very dilute lattice.¹³ Similarly, the existence of a spin-glass phase for RADs for $x < 1$ does not follow from its existence for $x = 1$.

Our main aim here is to find, by means of the parallel tempered Monte Carlo (TMC) algorithm,^{15,16} whether an equilibrium spin-glass phase exists in a site-diluted system of RADs in 3D. We also aim to establish, if the spin-glass phase does exist, whether in the condensed phase (1) there is a single extended state,¹⁷ as in a ferromagnet or in the droplet model of spin glasses,¹⁸ or (2) there are multiple extended states, as in the XY model in two dimensions (2D) (Ref. 19) or in the replica symmetry breaking theory of spin glasses.²⁰

The paper is planned as follows. In Sec. II we specify the RAD model and describe how we apply the parallel TMC algorithm.^{15,16} We do this for $x=0.35$ and $x=0.5$. Results are given in Sec. III. The entropy S follows from our data for the specific heat by numerical integration. It approximately vanishes at zero temperature.²¹ More precisely, $S < 0.01k_B$ at $T=0$. We provide numerical evidence for the existence of an equilibrium spin-glass phase below a transition temperature T_c . The evidence comes from the behavior of the distribution of the spin overlap parameter q .^{22,23} In the spin-glass phase, $\langle q^2 \rangle$ seems to vanish in the macroscopic limit, but only as a power of system size, in accordance with the existence of quasi-long-range order. For both $x=0.35$ and $x=0.5$, $k_B T_c = (0.95 \pm 0.1)x\varepsilon_d$, where ε_d is a nearest-neighbor dipole-dipole interaction energy which is defined in Sec. II. (Within errors, the value of T_c for $x=1$ we have previously obtained¹² also satisfies this expression.) For $T \leq 0.9T_c$, $\delta|q|$, that is, the root-mean-square deviation of $|q|$, fulfills $\delta|q| \approx 0.5\langle|q|\rangle\sqrt{T/T_c}$, independently of system size. This result is

compared with its counterpart for the XY model in 2D. Results are discussed in Sec. IV.

II. MODEL AND METHOD

We treat systems of magnetic dipoles on simple cubic lattices. We place a dipole on each lattice site with probability x and leave the site unoccupied with probability $1-x$. All dipoles point along randomly oriented anisotropy axes. With this model we aim to mimic assemblies of NPs in which uniaxial anisotropy energies are an order of magnitude larger than the largest dipole-dipole energy. Then, anisotropy energy barriers are not so large as to freeze spin reorientations near the spin-glass temperature, but are sufficiently large to restrict spins to point “up” or “down” approximately along the easy magnetization axes. We term this the random-axes-dipolar (RAD) model. The Hamiltonian is given by

$$\mathcal{H} = \frac{1}{2} \sum_{ij} \sum_{\alpha\beta} T_{ij}^{\alpha\beta} S_i^\alpha S_j^\beta, \quad (1)$$

where the first sum is over all occupied sites i and j of a simple cubic lattice, S_i^α is the α component of a classical three-component spin on site i ,

$$T_{ij}^{\alpha\beta} = \varepsilon_d (a/r_{ij})^3 (\delta_{\alpha\beta} - 3r_{ij}^\alpha r_{ij}^\beta / r_{ij}^2), \quad (2)$$

r_{ij} is the distance between i and j , ε_d is an energy, and a is a nearest-neighbor distance. Each spin points along a randomly chosen direction. These equations can be cast into a form that is manifestly Ising type by letting $\hat{\mathbf{u}}_j$ be (1) a null vector if the j site is unoccupied and (2) a three-component unit vector chosen randomly from a spherically uniform distribution if the j site is occupied, and defining a pseudospin $\sigma_j = \pm 1$ for each site, such that $\mathbf{S}_j = \hat{\mathbf{u}}_j \sigma_j$. We can then write

$$\mathcal{H} = -\frac{1}{2} \sum_{ij} J_{ij} \sigma_i \sigma_j, \quad (3)$$

where $J_{ij} = -\sum_{\alpha\beta} T_{ij}^{\alpha\beta} u_i^\alpha u_j^\beta$. Thus, the RAD model is an Ising model whose bonds J_{ij} are determined by the dipole-dipole terms $T_{ij}^{\alpha\beta}$ and the set of three-component randomly oriented unit vectors $\{\hat{\mathbf{u}}_j\}$.

From here on, unless we state otherwise, we let $k_B=1$, where k_B is Boltzmann’s constant and give all temperatures in terms of ε_d . We use periodic boundary conditions. Details and justification are given in Refs. 12, 14, 16, and 24.

In order to arrive at equilibrium results, we make use of the parallel TMC algorithm.¹⁵ This enables one to circumvent large energy barriers that can trap a system’s state. We apply the TMC algorithm as follows. We run (in parallel) several identical systems at different temperatures: a system at temperature T_0 , a second one at $T_0+\Delta T$, and so on, at equally spaced temperatures, up to T_m . We choose T_m to be at least twice as large as what we expect to be the transition temperature between the paramagnetic and spin-glass phases. We let each system evolve under the Metropolis MC algorithm for 10 MC sweeps before pairs of systems are given a chance to exchange their states. More specifically, pairs of systems at temperatures $(T_0$ and $T_0+\Delta T)$, $(T_0+2\Delta T$ and T_0

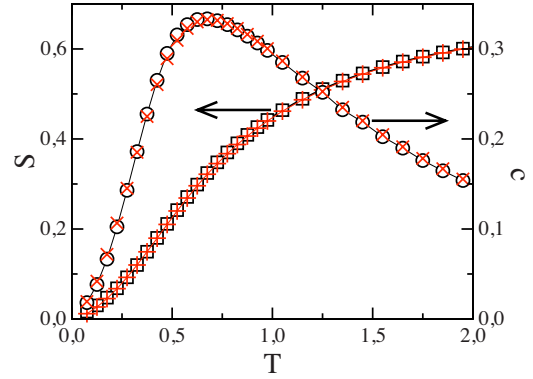


FIG. 1. (a) (Color online) Plots of c and of S vs T , for $x=0.35$. Data points (black) \circ (for $N=605$) and (red) \times (for $N=179$) are for the specific heat, while (black) \square (for $N=605$) and (red) $+$ (for $N=179$) are for the entropy. Errors, δS , for S increase as T decreases, from $\delta S \approx 0.002$ for $T=4$ to $\delta S \approx 0.004$ for $T < 0.1$.

+ $3\Delta T$), and so on, are given a chance to exchange states every 10 MC sweeps, and pairs of systems at temperatures $(T_0+\Delta T$ and $T_0+2\Delta T)$, $(T_0+3\Delta T$ and $T_0+4\Delta T)$, and so on, are given a chance to exchange states at times in between. Under TMC rules, these chances are as follows. A system at temperature T that is in state 2, and another one at $T+\Delta T$ in state 1, exchange states with probability $P=1$ if $E_1 < E_2$ and $\Delta T > 0$, but $P = \exp(-\Delta\beta\delta E)$, where $\Delta\beta = 1/T - 1/(T+\Delta T)$ and $\delta E = E_1 - E_2$, if $E_1 > E_2$ and $\Delta T > 0$.¹⁶

A sufficiently small value of ΔT must be chosen in order to keep $\exp(-\Delta\beta\delta E)$ from becoming too small. This will often be fulfilled if $\Delta\beta\Delta E \leq 1$, where ΔE is the mean energy difference between two systems at temperatures T and $T+\Delta T$. The required condition, $\Delta T \leq T/\sqrt{Nc}$, where N is the number of dipoles in the system, follows for ΔT if we replace ΔE by $cN\Delta T$ (we thus define c as the specific heat per spin). From plots of the specific heat vs T , such as the one shown in Fig. 1 for systems of 179 and of 605 magnetic dipoles on lattices with 0.35 of their sites occupied, one can get upper bounds for ΔT . (How the data points shown in Fig. 1 were obtained is explained in Sec. III.) Values of ΔT as well as of other parameters for all TMC runs are given in Table I.

In order to probe for spin-glass behavior, we define, as is usual, the spin overlap parameter.^{22,23} First, let

$$\phi_j = \sigma_j^{(1)} \sigma_j^{(2)}, \quad (4)$$

where $\sigma_j^{(1)}$ and $\sigma_j^{(2)}$ are the pseudospins [defined above Eq. (3)] on site j of identical twin replicas 1 and 2 of the system. Clearly,

$$q = N^{-1} \sum_j \phi_j \quad (5)$$

is a measure of the spin configuration overlap between replicas 1 and 2. Thus, $|q|=1$ if either $\sigma_j^{(1)} = \sigma_j^{(2)}$ for all j or $\sigma_j^{(1)} = -\sigma_j^{(2)}$ for all j . We also define the moments of q , $q_k = \langle |q|^k \rangle$, for $k=1, 2$, and 4, where $\langle \dots \rangle$ stands for an average which we next specify.

TABLE I. Simulation parameters. x is the probability that any given site is occupied by a magnetic dipole; N is the mean number of magnetic dipoles in the system; ΔT is the temperature step in the TMC runs; T_m is the highest temperature of all systems; N_s is the number of disordered system pairs; MCS is the number of tempered Monte Carlo sweeps.

x	0.35	0.35	0.35	0.35	0.5	0.5	0.5	0.5
N	22.4	75.6	179.2	604.8	32	108	256	864
ΔT	0.05	0.05	0.05	0.05	0.05	0.05	0.05	0.05
T_m	1.0	1.0	1.0	1.0	1.3	1.3	1.3	1.3
N_s	10000	3000	1250	900	5000	3000	1500	150, 600
MCS	50000	10^5	5×10^6	10^6	50000	50000	10^6	$10^6, 10^5$

Suppose our TMC runs last for a time t , and assume that, within the temperature range we are interested in, equilibration takes place in a time less than $t/2$. During the TMC runs, we write, at equally spaced time intervals within the $(t/2, t)$ range, the spin configuration for each temperature. We later draw from these written configurations the numbers for $\sigma_j^{(1)}$ that go into Eq. (4). This procedure is repeated for the same system (that is, for a system with the same anisotropy axes), but starting from different initial conditions. From this second set of configurations, we draw the numbers for $\sigma_j^{(2)}$ that go into Eq. (4). A value of q for each time t and temperature T is thus obtained from Eq. (5). We first average all quantities of interest over t and finally repeat these pairs of TMC runs a number N_s of times for different realizations of axes orientations, from which the average values we are reporting are obtained. Values we have used in our TMC runs for N_s and for t (labeled MCS therein) are given in Table I.

Finally, in order to check that equilibration actually takes place as assumed, we also calculate the spin overlap \tilde{q} , not between identical twin replicas, but between spin configurations at two different times t_0 and t_1 of the same TMC run. We do this for several values of t_0 and t_1 in the neighborhoods of $t/2$ and t , respectively, and average over different random axes realizations, in order to obtain \tilde{q}_1 . For t sufficiently large, $\tilde{q}_1 = q_1$. We find this is fulfilled for all parameter values given in Table I, except for (1) $x=0.35$, $N=604.8$, and $T \leq 0.2$, and (2) for $x=0.5$, $N=864$, and $T \leq 0.35$.

III. RESULTS

Plots of the specific heat and of the entropy vs T for the RAD model are shown in Fig. 1 for $x=0.35$. The data for the specific heat follow from numerical derivatives of the energy with respect to T . We obtain S from $S(T) = \ln 2 + \int_{\infty}^T c(T')/T' dT'$. Our data cover the temperature range (not all of it shown in Fig. 1) $0.05 < T < 4$. For a numerical integration, we must extrapolate our data for $c(T)$ beyond $T=4$. To this end, we use the leading term of an energy expansion in powers of $1/T$, which gives $c(T) \rightarrow A/T^2$. A fit of the value of A to our data for $c(T)$ at $T=4$ leads to a $\Delta S \approx 0.027$ contribution to the entropy from the $T > 4$ range. We thus obtain $S(T)$. It is exhibited in Fig. 1. Errors for $S(T)$ come in two approximately equal pieces: (1) an error of roughly 0.002 from the $4 < T < \infty$ range and (2) an error of roughly 0.002, from errors in the data for the specific, which enter the integral $\int_{\infty}^T c(T')/T' dT'$. We finally obtain $S = 0.015 \pm 0.004$ for

$T=0.05$, and extrapolations below $T=0.05$ yield $S < 0.01$ at $T=0$. Similarly, we find $S < 0.01$ at $T=0$ for $x=0.50$.

Our results for q_1 and q_2 follow. Note we use an absolute value in the definition of q_1 . Recall that $Nq_2 = O(1)$ in the paramagnetic phase and diverges at the transition temperature T_c . Above the lower critical dimension d_c , $q_2 = O(1)$ in the droplet model of the spin-glass phase.¹⁸ Our data show that q_2 decreases as N increases, for all nonzero temperatures. Log-log plots of q_2 vs N are exhibited in Figs. 2(a) and 2(b) for various values of T . The behavior of q_2 for $T \leq 0.3$ and $T \leq 0.45$ for $x=0.35$ and $x=0.5$, respectively, is consistent with $q_2 \sim N^{-y}$, where y is some positive parameter that depends on T . This behavior is reminiscent of the XY model in 2D.¹⁹ For higher values of T , q_2 vs N clearly curves downward, in accordance with a faster than algebraic in N decay, as one expects for the paramagnetic phase. Log-log plots of q_1^2 are also shown in Figs. 2(a) and 2(b) for various values of T . Note that $q_2/q_1^2 > 1$ for all N and nonzero T , which implies a nonvanishing uncertainty in $|q_r|$, where $q_r = q/q_1$, for $T > 0$.

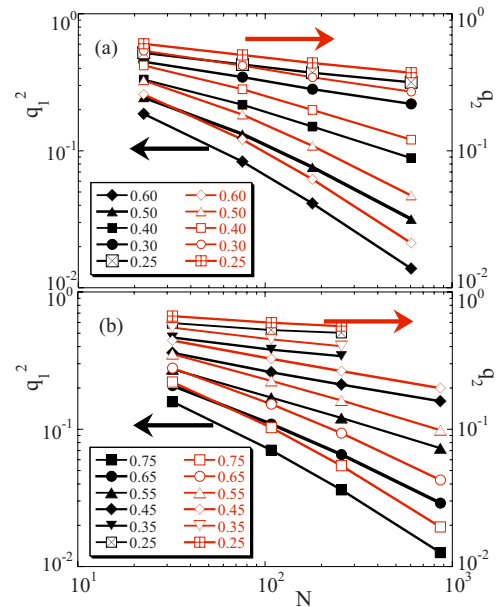


FIG. 2. (Color online) (a) Log-log plots of q_1 and of q_2 vs N for $x=0.35$ and the shown values of T . Lines are guides for the eyes. (b) Same as in (a) but for $x=0.5$. Some data points for small T and large N are missing because our TMC runs did not reach equilibration for them.

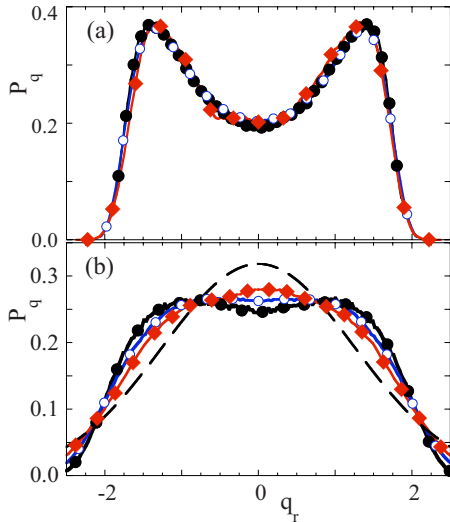


FIG. 3. (Color online) (a) Plots of the probability P_q vs q_r , for $x=0.50$ and $T=0.45$. \bullet , \circ , and \blacklozenge are for $N=108, 256$, and 864 , respectively. Lines are guides for the eyes. For clarity, only 30% of the data points are shown. (b) Same as in (a) but for $T=0.60$. For comparison, a dashed line is shown for $(1/\pi)\exp(-q_r^2/\pi)$, which ensues for a macroscopic paramagnetic phase.

We next report results for the probability distribution $P_q(q_r)$. Unless stated otherwise, all data given below are for a normalized $P_q(q_r)$, that is, $\int P_q(q_r) dq_r = 1$. Recall that in the paramagnetic phase, because spin-spin correlation lengths are finite, the central limit theorem implies q_r is normally distributed for macroscopic systems. Plots of P_q vs q_r are shown in Fig. 3(a) (for $T=0.45$) and Fig. 3(b) (for $T=0.60$), both for $x=0.50$. The distribution of q appears to be size independent for $T=0.45$. For $T=0.60$, on the other hand, $P_q(q_r)$ drifts with system size. The drift seems consistent

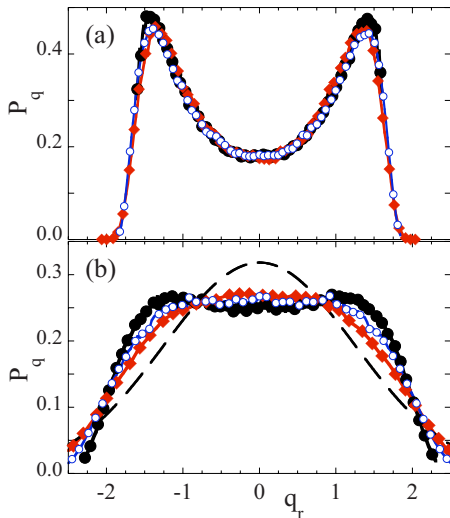


FIG. 4. (Color online) (a) Plots of the probability distribution P_q vs q_r , for $x=0.35$ and $T=0.3$. \bullet , \circ , and \blacklozenge are for $N=76, 179$, and 605 , respectively. Lines are guides for the eyes. For clarity, only 30% of the data points are shown. (b) Same as in (a) but for $T=0.45$. For comparison, a dashed line is shown for $(1/\pi)\exp(-q_r^2/\pi)$, which ensues for a macroscopic paramagnetic phase.

with $P_q(q_r) \rightarrow (1/\pi)\exp(-q_r^2/\pi)$ as $N \rightarrow \infty$, which would be in accordance with a paramagnetic phase. Similar remarks apply to the plots exhibited in Fig. 4(a) (for $T=0.30$) and Fig. 4(b) (for $T=0.45$), both for $x=0.35$. Clearly, $0.30 < T_c < 0.45$ for $x=0.35$, and $0.45 < T_c < 0.60$ for $x=0.50$. For $T < T_c$, plots of P_q vs q_r (not shown) become sharply peaked only as $T \rightarrow 0$. Now, if it turns out that $d_c=3$ for systems of RADs, then $P_q(q_r)$ ought to exhibit critical behavior for all $T \leq T_c$, more specifically, $P_q(q_r)$ ought to be size independent. This sort of behavior is best summarized by the mean square deviation of $|q_r|$, $\Delta_q^2 \equiv \langle q_r^2 \rangle - \langle |q_r| \rangle^2$.

Plots of Δ_q^2 vs T/x are shown in Fig. 5(a) (for $x=0.35$) and in Fig. 5(b) for $x=0.50$ for various values of N . Clearly, curves in Figs. 5(a) and 5(b) differ only slightly, as expected from the argument given in Sec. I, that variations in T and x have an effect only through T/x if $x \ll 1$. The data points in both figures suggest $\Delta_q^2 \rightarrow \pi/2 - 1$ as $N \rightarrow \infty$, for $T/x \gtrsim 1$. This is expected for a macroscopic paramagnet, since $P_q(q_r)$ is a normal distribution then. For $T/x \lesssim 0.8$,

$$\Delta_q \approx 0.5\sqrt{T/x} \tag{6}$$

provides the best fit to the data points shown in Figs. 5(a) and 5(b). This is in contrast with the behavior, $\Delta_q^2 \rightarrow 0$ as

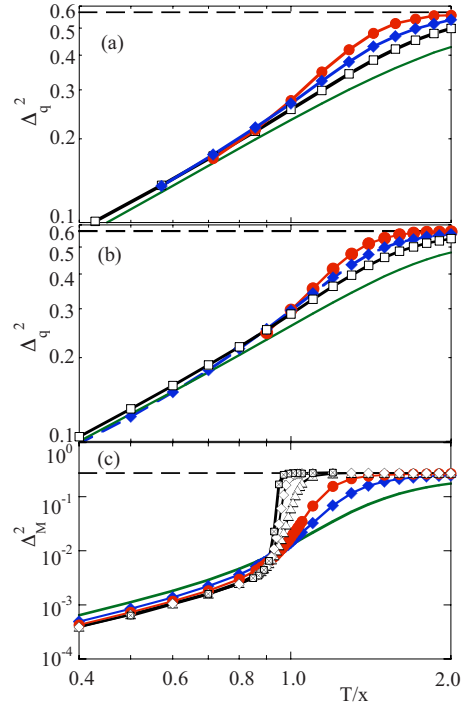


FIG. 5. (Color online) (a) Plots of Δ_q^2 vs T for $x=0.35$. (red) \bullet , (blue) \blacklozenge , and (black) \square are for $N=604.8, 179.2$, and 75.6 , respectively. The continuous (green) line is for $N=22.4$. The dashed line is for macroscopic paramagnet. (b) Same as in (a) but for $x=0.50$. \bullet (red), \blacklozenge (blue), and \square (black) are for $N=864, 256$, and 108 , respectively. The continuous (green) line is for $N=32$, and the dashed line is for $\pi/2 - 1$. (c) Same as in (a), but for the mean square relative deviation Δ_M^2 of $|M|$, where M is the magnetic moment of XY systems of $L \times L$ spins in 2D vs T/x , where $x=1$, for $L=4$ (green line), $L=8$ (blue \blacklozenge), $L=16$ (red \bullet), $L=64$ (black \triangle), $L=256$ (black \diamond), and $L=1024$ (black \boxtimes).

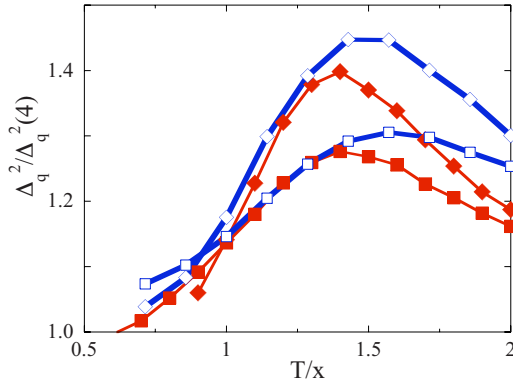


FIG. 6. (Color online) Plots of $\Delta_q^2/\Delta_q^2(L=4)$ vs T , where $\Delta_q^2(L=4)$ is Δ_q^2 for $L=4$. All (\blacklozenge) are for $x=0.35$ and $L=12$, and all (\blacksquare) are for $x=0.35$ and $L=8$. All (\diamond) are for $x=0.50$ and $L=12$, and all (\square) are for $x=0.50$ and $L=8$. Lines are guides for the eyes.

$N \rightarrow \infty$, that one expects for the ordered phase of ferromagnets such as the Ising model in two or higher dimensions. For a more relevant comparison, we show in Fig. 5(c) plots of the mean square relative deviation Δ_M^2 of $|M|$, where M is the magnetic moment, of the XY model in 2D. Notice there is no counterpart in Figs. 5(a) and 5(b) for the data points shown for rather large systems in Fig. 5(c).

Now, it is hard to see in Figs. 5(a) and 5(b) where curves merge or cross, near $T/x=0.9$. In order to enhance the differences between Δ_q^2 curves for different values of L , we plot $\Delta_q^2/\Delta_q^2(L=4)$ vs T/x . These plots are shown in Fig. 6 for $x=0.35$ and $x=0.50$. The weak dependence on x is remarkable. Notice that whereas curves for $L=12$ and $L=8$ cross near $T/x=0.95$ for $x=0.35$ and for $x=0.50$, the same curves cross the $\Delta_q^2/\Delta_q^2(L=4)=1$ horizontal line only, if at all, for much smaller values of T/x . It thus appears that curves for Δ_q^2 increasingly larger system sizes merge or cross at increasingly larger values of T/x .

We show plots of $P_q(0)$ vs N for $x=0.35$ and various temperatures in Fig. 7. In the paramagnetic phase, we expect $P_q(0) \rightarrow 1/\pi$ as $N \rightarrow \infty$, as follows from a normal distribution of q_r about $q_r=0$. Now, for $T/x \geq 1.0$, we see that $P_q(0) < 1/\pi$, but $P_q(0)$ increases as N increases. On the other hand, for $T/x \leq 0.857$, $P_q(0)$ remains, within errors, constant. This is again in contrast with the behavior, $P_q(0) \rightarrow 0$ as $N \rightarrow \infty$, that one expects for the ordered phase of ferromagnets such as the Ising model in two or higher dimensions, but is in accordance with $d_c \approx 3$. Finally, taking into account all our observations above, we arrive at our best estimate for T_c : $T_c/x = 0.95 \pm 0.1$.

IV. DISCUSSION

By tempered Monte Carlo calculations, we have studied systems of RADs in simple cubic lattices in which each site is occupied by a magnetic dipole with probability x . System sizes, Monte Carlo run lengths, and other details about the calculations can be found in Table I. The entropy S as a function of temperature, which follows from our data for the specific heat, is shown in Fig. 1 for $x=0.35$. S approximately vanishes at zero temperature. More precisely, $S < 0.01k_B$ at

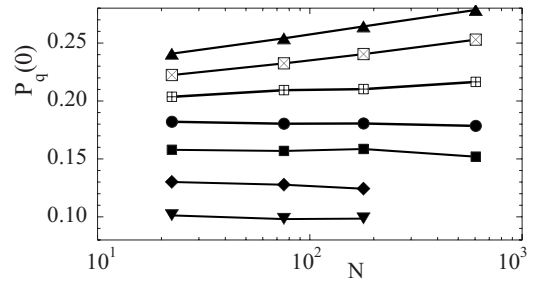


FIG. 7. Plots of $P_q(q=0)$ vs N , for $x=0.35$, for $T/x=1.286$ (\blacktriangle), $T/x=1.143$ (\boxtimes), $T/x=1.0$ (\boxplus), $T/x=0.857$ (\bullet), $T/x=0.714$ (\blacksquare), $T/x=0.57$ (\blacklozenge), and $T/x=0.429$ (\blacktriangledown).

$T=0$ for both $x=0.35$ and $x=0.50$.²¹ Note also that a vanishing zero-temperature entropy implies that trapping above the lowest energy states as $T \rightarrow 0$ occurs only very rarely in our TMC runs.

We have obtained the spin overlap parameter defined in Eq. (5). More specifically, we have obtained the mean value of $|q|$, which we term q_1 , and the mean square value of q , which is q_2 . The plots in Fig. 2(a) (for $x=0.35$) and in Fig. 2(b) (for $x=0.50$) suggest both q_1 and q_2 vanish for all non-zero T as $N \rightarrow \infty$. As can be gathered from those plots, both $q_1^2 \sim N^{-y}$ and $q_2 \sim N^{-y}$, where y is some positive parameter that depends on T , for $k_B T/x \leq 0.9\epsilon_d$. This points to quasi-long-range order. It is reminiscent of the XY model in 2D and suggests the lower critical dimension d_c of the RAD model is at or near 3D. However, we have no grounds for taking the analogy between the XY model in 2D and the RAD model in 3D any further than this. We make no claim on the relation between the universality classes these systems may belong to.

We have also studied the probability distribution $P_q(q_r)$, where $q_r = q/q_1$. Within errors, $P_q(q_r)$ is independent of system size at $T=T_c$, as is illustrated in Figs. 4(a) and 3(a) for $k_B T/x \approx 0.9\epsilon_d$. We have also found (not shown) $P_q(q_r)$ to be independent of system size at lower temperatures. The results shown in Figs. 5(a) and 5(b) and in Eq. (6), for the mean square deviation of $|q_r|$, Δ_q^2 , are consistent with a $P_q(q_r)$ that is independent of system size for all $k_B T/x \leq 0.9\epsilon_d$. Again, this points to $d_c \approx 3$ for the RAD model and is in contradistinction to results for the nearest-neighbor random bond Ising model in 3D.²⁵

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