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LETTER TO THE EDITOR

On the critical dynamics of the diluted Q -state Potts models

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Abstract. The results of Monte Carlo simulations of the four-state Potts model on a square lattice at the bond percolation threshold are presented. Estimates are given for the new dynamical exponents A and B . Our result for A is in clear contradiction with a recent conjecture of Nunes da Silva and Lage.

A recent [1] neutron scattering experiment on $\text{Rb}_2\text{Co}_p\text{Mg}_{1-p}\text{F}_4$, a site-diluted two-dimensional Ising antiferromagnet, near the percolation threshold, $p_c = 0.5927$, has generated considerable interest [2-8] in the critical dynamics of diluted spin systems.

In the dynamic scaling hypothesis the average relaxation time, $\tau_{\text{AV}}(T)$, scales as

$$\ln(\tau_{\text{AV}}(T)) = f(\ln \xi_T) \quad (1)$$

where ξ_T is the thermal correlation length. Conventional dynamic scaling [9] would imply that $f(\ln \xi_T) \sim Z(\ln \xi_T)$, where Z is a dynamic critical exponent. For percolating systems ($p = p_c$), however, it is found [2-8] that Z is no longer a constant but a temperature-dependent function

$$Z(T) = A(\ln \xi_T) + B. \quad (2)$$

Very recently, the breakdown of dynamic scaling has also been confirmed for two-dimensional diluted Potts models [6, 7] and the bond-diluted Ising model on hypercubic lattices [8]. The analytic work of Nunes da Silva and Lage [7, 8] is particularly interesting because it suggests that the new dynamical exponents depend very strongly on both the dimensionality, d , and the number of spin components, q .

The dimensionality and symmetry of the problem are two crucial factors governing the critical behaviour of a system. By varying just the number of spin components and keeping fixed all other features of the system (the type of dilution, the updating rate, etc) we are able to investigate the q dependence of the dynamic behaviour.

In this letter we present the results from Monte Carlo simulations of the two-dimensional four-state Potts model on a square lattice at the bond percolation threshold. We shall show that our results contradict Nunes da Silva and Lage [7].

We choose the Hamiltonian to be

$$H = -\sum_{\langle ij \rangle} J_{ij} \delta_{\alpha_i, \alpha_j} \quad (3)$$

where $\alpha_i (\alpha_i = 1, \dots, q)$ are the Potts spins situated on every site of a 64×64 square lattice ($q = 4$ in our simulations) and the nearest-neighbour ferromagnetic couplings are selected according to

$$P(J_{ij}) = \frac{1}{2}[\delta(J_{ij}) + \delta(J_{ij} - 1)]. \quad (4)$$

Imposing periodic boundary conditions, we update the spins via the Metropolis transition probability. The data were collected over the temperature range $0.44 \leq T/T_c(\text{pure}) \leq 1.10$ and have been averaged over 10-248 samples; the largest statistical error bar is less than 10%. The only difference between the system discussed here and the one simulated by Jain *et al* [6] is the number of Potts states (this should be consulted for further technical details).

We define an average relaxation time by

$$\tau_{AV} = \frac{4}{3} \int_0^{+\infty} \left(N^{-1} \sum_i \delta_{\alpha_i(t_0)\alpha_i(t+t_0)} - \frac{1}{4} \right) dt \tag{5}$$

where $N = 4096$, the number of spins, and $t = t_0$ indicates an equilibrium state of the system. The asymptotic behaviour of the spatial correlation function,

$$\Gamma(n) = \frac{4}{3} \left(N^{-1} \sum_i \langle \delta_{\alpha_i, \alpha_{i+n}} \rangle_T - \frac{1}{4} \right) \tag{6}$$

where $\langle \dots \rangle_T$ implies a thermal average and n ($n = 0, 1, \dots, 10$) is the displacement in the x direction, is given by

$$\Gamma(n) \sim \exp(-n/\xi_T) \quad \text{for} \quad n \gg \xi_T. \tag{7}$$

Equation (7) enables us to extract the thermal correlation length for any temperature. Now, one expects [10]

$$\xi_T(q) = \xi_0(q) \exp(\beta\nu) \tag{8}$$

where $\beta = 1/T$, $\nu (= \nu_T)$ is the (universal) thermal exponent and $\xi_0(q)$ is the non-universal amplitude.

In figure 1 we show a plot of $\xi_T(q=3)/\xi_T(q=4)$ against T . The weighted line of best fit indicates that $\xi_0(q=3)/\xi_0(q=4) = 1.20 \pm 0.11$. Since [6] $\xi_0(q=3) = 0.16 \pm 0.02$, we have that $\xi_0(q=4) = 0.14 \pm 0.03$. Further, if we assume that $\nu = \frac{4}{3}$, the conjectured theoretical value [9], then we have that for all temperatures simulated $\xi_T \ll$ linear size

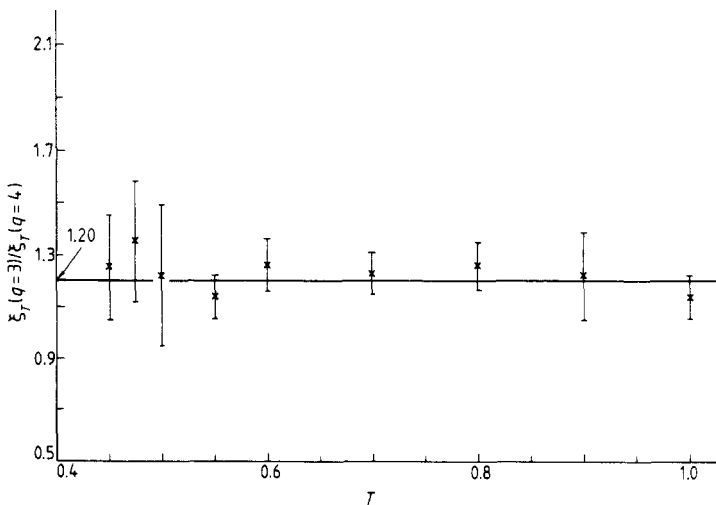


Figure 1. A plot of $\ln \xi_T(q=3)/\ln \xi_T(q=4)$ against T . The intercept on the y axis implies that $\xi_0(q=3)/\xi_0(q=4) = 1.20 \pm 0.11$.

of the lattice and $\xi_T \gg 1$ for $T \leq 0.625$. So our results are not expected to be influenced by finite-size effects [4].

Writing

$$\ln \tau_{AV} = Y_1/T^2 + Y_2/T + Y_3 \tag{9}$$

we have that $A(q) = Y_1/\nu^2$, $B(q) = \nu^{-1}(Y_2 - 2Y_1\nu^{-1} \ln \xi_0(q))$ and $Y_3 = \text{constant}$. In figure 2 we show $\ln \tau_{AV}$ against $1/T$. The best quadratic fit to the data for $T \leq 0.625$ gives $A(q=4) = 0.56 \pm 0.04$ and $B(q=4) = 4.76 \pm 0.80$. Table 1 contains the various estimates which have been made for $A(q)$ and $B(q)$ in two dimensions. We note that the value of $A(q=4)$ is not consistent with the conjecture of [7] who suggest that $A(q=4)/A(q=2) = \frac{3}{2}$.

We can make a more direct comparison with the $q=3$ case by considering

$$\ln[\tau_{AV}(q=3)/\tau_{AV}(q=4)] = \alpha_1/T^2 + \alpha_2/T + \alpha_3 \tag{10}$$

where $\alpha_1 = \nu^2[A(q=3) - A(q=4)]$, $\alpha_2 = \nu\{2[A(q=3)\xi_0(q=3) - A(q=4)\xi_0(q=4)] + (B(q=3) - B(q=4))\}$ and $\alpha_3 = \text{constant}$. Thus, $A(q=3) = A(q=4)$ would imply that a plot of $\ln[\tau_{AV}(q=3)/\tau_{AV}(q=4)]$ plotted against $1/T$ should be linear. Such a plot is shown in figure 3. The line of best fit for $T \leq 0.625$ has slope 1.00 ± 0.22 and intercept -1.75 ± 0.30 . The data presented in table 1 are consistent with a slope in the range

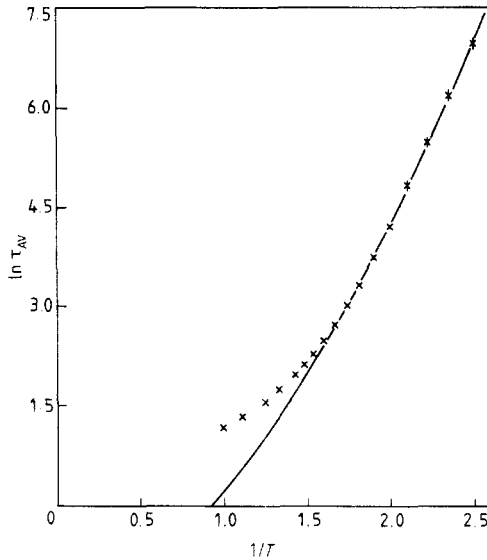


Figure 2. A plot of $\ln \tau_{AV}$ against $1/T$ for $q=4$. The quadratic fit shown yields $A(q=4) = 0.56 \pm 0.04$ and $B(q=4) = 4.76 \pm 0.80$.

Table 1. The behaviour of $A(q)$ and $B(q)$ for $d=2$.

Reference	$A(q)$	$B(q)$
Jain [4]; $q=2$	0.51 ± 0.05	3.25 ± 0.41
Jain <i>et al</i> [6]; $q=3$	0.78 ± 0.15	3.35 ± 0.88
This work; $q=4$	0.56 ± 0.04	4.76 ± 0.80
Nunes da Silva and Lage [7]	$A(q)/A(2) = 2(q-1)/q$	

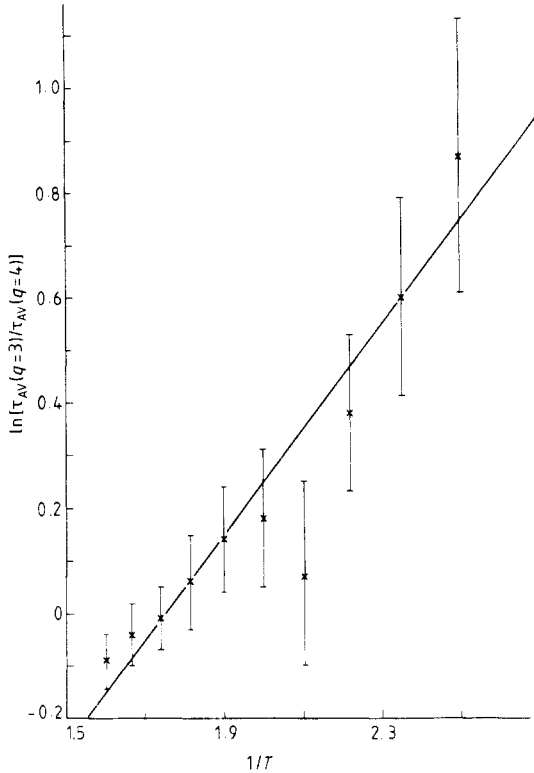


Figure 3. A plot of $\ln[\tau_{AV}(q=3)/\tau_{AV}(q=4)]$ against $1/T$. The straight line, which has slope 1.00 ± 0.22 , is a consequence of assuming that $A(q=3) = A(q=4)$.

-4.19 to 0.53 . Note that, if we also have $B(q=3) = B(q=4)$, then the slope would be expected to be 0.05 ± 0.12 . So it would appear that A is possibly independent of q . B , however, probably depends on the number of Potts states.

To conclude, we have given estimates for the new dynamical exponents A and B for the four-state Potts model at the bond percolation threshold. Our results, although in agreement with Jain *et al* [6], contradict a recent conjecture of Nunes da Silva and Lage [7].

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