Nonequilibrium scaling explorations on a two-dimensional Z(5)-symmetric model

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(Received 1 May 2014; revised manuscript received 12 July 2014; published 1 October 2014)

We have investigated the dynamic critical behavior of the two-dimensional Z(5)-symmetric spin model by using short-time Monte Carlo (MC) simulations. We have obtained estimates of some critical points in its rich phase diagram and included, among the usual critical lines the study of first-order (weak) transition by looking into the order-disorder phase transition. In addition, we also investigated the soft-disorder phase transition by considering empiric methods. A study of the behavior of $\beta/\nu z$ along the self-dual critical line has been performed and special attention has been devoted to the critical bifurcation point, or Fateev-Zamolodchikov (FZ) point. First, by using a refinement method and taking into account simulations out of equilibrium, we were able to localize parameters of this point. In a second part of our study, we turned our attention to the behavior of the model at the early stage of its time evolution in order to find the dynamic critical exponent $z$ as well as the static critical exponents $\beta$ and $\nu$ of the FZ point on square lattices. The values of the static critical exponents and parameters are in good agreement with the exact results, and the dynamic critical exponent $z \approx 2.28$ very close to the four-state Potts model ($z \approx 2.29$).

DOI: 10.1103/PhysRevE.90.042101 PACS number(s): 05.20.–y, 05.10.Ln, 05.70.Jk

I. INTRODUCTION

In statistical mechanics, nontrivial models have been extensively studied after the exact solution of the two-dimensional Ising model [1]. A lot of authors have devoted an extensive use of several methods to describe the theory of magnetic systems by studying generalizations of such models, with more complex and richer phase diagrams. Among these models, one that deserves special attention is the Z($N$) model whereas, differently than the Ising model whose spin variable can assume only two values, each spin can assume $N$ values and more than one coupling constant for $N > 4$. This leads to more delicate aspects with a phase diagram that is not completely understood yet, even, for example, for small values of $N$ such as $N = 5$.

The two-dimensional Z($N$) model contains several known systems as particular cases, for instance, the Ising ($N = 2$) and XY ($N = \infty$) models, as well as, the $N$-state scalar and vector Potts (clock) models, and the Ashkin-Teller model ($N = 4$). For $N \leq 4$, the phase diagram possesses a traditional second-order phase transition, and for $N = \infty$, it exhibits a Kosterlitz-Thouless (KT)-type phase transition [2]. But, for what $N$ value does this last phase transition appear? Several works report that the KT phase transition appears at $N = 5$ [3–7]. The Z(5) model exhibits a rich phase diagram with first-order transitions, including the five-state Potts point [8], two second-order transitions of the Ising type at Fateev-Zamolodchikov (FZ) integrability points [5], and two lines of infinite-order transitions (dual to each other) of the KT type [3,4,7,9–11] (see dashed lines in Fig. 1). Several works assert that the FZ points, henceforth named the “bifurcation points,” coincide with the points where the KT transitions are originated [5,6,10,12,13].

So, this interesting model and, especially the bifurcation points (for $N = 5$), deserves further explorations and nonequilibrium analysis can be an interesting alternative to obtain not only the static critical exponents but also the dynamical ones which have not yet been obtained in previous contributions. Moreover, this approach has proved to be efficient in determining the critical parameters of several models as shown in recent works (see, for example, Refs. [14–16]).

In this paper, we present results from the study of the critical properties of the isotropic ferromagnetic two-dimensional spin model with Z(5) symmetry, hereafter denoted as the Z(5) model, by using time-dependent MC simulations. As we are dealing with a symmetric model, the two bifurcation points are also symmetric and possess the same set of critical exponents. Hence, we concentrated on only one of them. Our contributions are divided in four parts as follows.

1. We estimated the critical parameters $x_1$ and $x_2$ of the bifurcation point [5] by using a simple refinement method, in the context of time-dependent MC simulations which searches the best power-law time decay of magnetization, as proposed in Ref. [16].

2. We obtained the dynamic critical exponent $z$ and the static critical exponents $\nu$ and $\beta$ of the two independent order parameters of the model for the bifurcation point.
Finally, in Sec. VI we summarize and conclude our work. Along the self-dual line with special attention to the FZ point, we specifically showed some estimates of critical exponents and parameters. We also show how to simulate such behaviors via time-dependent Monte Carlo simulations to calculate the critical parameters $x_1$ and $x_2$, as well as the critical exponents $z$, $\beta$, and $\nu$, we present in the next section the finite size scaling developed to describe nonequilibrium spin systems, the time-dependent power laws obtained from this approach, and some details about time-dependent MC simulations to be applied.

\[ x_1 = \exp \left[ \frac{\sqrt{5}(k_1 - k_2) - 5(k_1 + k_2)}{4} \right], \]

and

\[ x_2 = \exp \left[ \frac{\sqrt{5}(k_2 - k_1) - 5(k_1 + k_2)}{4} \right]. \]

In the particular case $k_2 = k_1$ we recover the scalar five-state Potts model and for $k_2 = 0$ the clock model. It is interesting to observe that the five-state Potts point corresponds to the meeting between the self-dual line defined by $x_1 + x_2 = (\sqrt{5} - 1)/2$ and the Potts physical line $x_1 = x_2$, this last one being a symmetry line of the diagram.

In this work we are more concerned with the bifurcation points on the self-dual line. Actually, as can be seen in Fig. 1 the model has two bifurcation points (FZ points) localized on the self-dual line. The phase transition line between the FZ points (which includes the Potts point) is of weak first order, and on the right(left) of the rightmost(leftmost) FZ point there are continuous transition lines between ordered-soft and disordered-soft phases. However the two bifurcation points are symmetric to each other and have the same set of critical exponents. For this reason, we took into account only one of them. The ratio of the coupling constants for the bifurcation point is given by $k_2/k_1 = (\sqrt{5} - 1)/2 \approx 0.618034$. Moreover, there are four order parameters but only two of them are independent ones [18], namely

\[ M_1 = \langle \delta_{n_1} - \delta_{n_2} \rangle, \]

and

\[ M_2 = \langle \delta_{n_1} - \delta_{n_3} \rangle, \]

where $\delta_{i,j}$ is the Kronecker’s delta.

Since we established the main details of the model in order to calculate the critical parameters $x_1$ and $x_2$, as well as the critical exponents $z$, $\beta$, and $\nu$, we present in the next section the finite size scaling developed to describe nonequilibrium spin systems, the time-dependent power laws obtained from this approach, and some details about time-dependent MC simulations to be applied.

\[ FIG. 1. \text{(Color online)} \text{ Phase diagram of the ZS model according to Ref. [17], (Phase I) Disordered phase; (Phase II) ordered phase; and (Phase III) soft phase. The five-state Potts and FZ points are specifically indicated on the self-dual line. The diagram is symmetric with respect to the Potts physical line.} \]

(3) We explored several points on the self-dual line of the model by estimating the exponents of its two order parameters. We showed that the exponents are different along this line but respect a peculiar symmetry. However, for the particular point corresponding to the five-state Potts model the critical exponents assume the same value.

(4) We also explored and obtained some estimates of weak first-order points on the self-dual line and other second-order points on the soft-disorder transition line using a heuristic method, developed in this paper, that takes into account the second moment of the order parameters. This article is organized as follows. In the next section we define the model and briefly discuss some peculiarities of its phase diagram. In Sec. III we present some finite size scaling relations in nonequilibrium spin systems theory and describe the power laws which are considered in this work to measure the required exponents and parameters. We also show how to simulate such behaviors via time-dependent Monte Carlo simulations. Our results are divided in two sections: In Sec. IV, we determined estimates of the phase transition points in the phase diagram by using a nonequilibrium approach and in Sec. V we specifically showed some estimates of critical exponents along the self-dual line with special attention to the FZ point. Finally, in Sec. VI we summarize and conclude our work.

II. THE MODEL AND ITS PHASE DIAGRAM

In this article we have studied the dynamic critical behavior of the Z(5) model by using short-time Monte Carlo simulations. The most general Hamiltonian of this model is given by

\[ -\beta \mathcal{H} = \sum_{i,j} k_1 \left[ \cos \left( \frac{2\pi}{5} (n_i - n_j) \right) - 1 \right] + k_2 \left[ \cos \left( \frac{4\pi}{5} (n_i - n_j) \right) - 1 \right], \]

where $(i,j)$ indicates that the spin variables interact only with their nearest neighbors, $i$ and $j$ label the sites of a two-dimensional lattice of size $L \times L$, $k_1$ and $k_2$ are the two positive coupling constants, and $n_i = 0, 1, 2, 3, 4$ label the degrees of freedom of each site of the lattice.

In Fig. 1 (according to Ref. [17]) we can observe the phase diagram of this model translated to the suitable variables:

\[ x_1 = \exp \left[ \frac{\sqrt{5}(k_1 - k_2) - 5(k_1 + k_2)}{4} \right], \]

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III. NONEQUILIBRIUM DYNAMICS AND TIME-DEPENDENT MC SIMULATIONS

Until a few years ago, the numerical calculation of critical exponents was carried out only in equilibrium. Unfortunately, in this stage, the measurements of such exponents are very difficult due to severe critical slowing down which takes place in the vicinity of the critical temperature. To circumvent this difficulty, some algorithms were proposed, for instance, the cluster algorithm [19,20] that, although it is very efficient in...
the study of static properties, violates the dynamic universality class of the specific local dynamics, such as Model A.

Another way to avoid problems with the critical slowing down was proposed by Janssen, Schaub, and Schmittmann [21] and Huse [22], both in 1989. They discovered that when using renormalization group techniques and numerical calculations, respectively, that there is universality and scaling behavior far from equilibrium. Since then, the so-called short-time regime has become an important method for the study of phase transitions and critical phenomena.

The dynamic scaling relation obtained by Janssen et al. for the kth moment of the order parameter, extended to systems of finite size [23], is written as

$$\langle M^k \rangle(t,\tau, L, m_0) = b^{-k\nu z} \langle M^k \rangle(b^{-1/t \cdot \xi \cdot \nu \cdot z \cdot t, b^{-1/L \cdot m_0})$$

(4)

where t is the time evolution, b is an arbitrary spatial rescaling factor, $\tau = (T - T_c)/T_c$ is the reduced temperature, and L is the linear size of the lattice. The exponents $\beta$ and $\nu$ are the equilibrium critical exponents associated with the order parameter and the correlation length, and z is the dynamic exponent characterizing temporal correlations in equilibrium. Here, the operator $\langle \ldots \rangle$ denotes averages over different configurations due to different possible time evolution from each initial condition of a given initial magnetization $m_0$. For a large lattice size $L$ and small initial magnetization $m_0$ at the critical temperature ($\tau = 0$), Eq. (4) is governed by the new dynamic exponent $\theta$, according to

$$\langle M \rangle_{m_0} \sim m_0^\theta,$$

(5)

if we choose the scaling factor $b = t^{1/z}$. This new exponent characterizes the so-called critical initial slip, the anomalous behavior of the order parameter when the system is quenched to the critical temperature $T_c$.

In addition, a new critical exponent $x_0$, which represents the anomalous dimension of the initial magnetization $m_0$, is introduced to describe the dependence of the scaling behavior on the initial conditions. This exponent is related to $\theta$ as $x_0 = \theta z + \beta/\nu$. Actually the relaxation of spin systems is determined by two different behaviors: this initial slip and a second behavior corresponding to a power-law decay. This can be derived from Eq. (4). After the scaling $b^{-1}L = 1$ at the critical temperature $T_c$, the first ($k = 1$) moment of the order parameter is $\langle M \rangle(t,L,m_0) = b^{-\beta/\nu} \langle M \rangle(b^{-1/t \cdot \xi \cdot \nu \cdot z \cdot t, L_0 \cdot m_0)$. Denoting $u = tL^{-z}$ and $w = L^{\nu z}m_0$, one has $\langle M \rangle(u,w) = L^{-\beta/\nu} \langle M \rangle(L^{-z}t, L_0 \cdot m_0)$. The derivative with respect to $L$ is given by

$$\partial_L \langle M \rangle = (\beta/\nu) L^{-\beta/\nu - 1} \langle M \rangle(u,w) + L^{-\beta/\nu} [\partial_u \langle M \rangle \partial_L u + \partial_w \langle M \rangle \partial_L w]$$

where explicitly we have $\partial_u \langle M \rangle = -\beta z L^{-z-1}$ and $\partial_L w = x_0 m_0 L^{-z-1}$. In the limit $L \rightarrow \infty$, which implicates in $\partial_1 \langle M \rangle \rightarrow 0$, one has $x_0 w \partial_u \langle M \rangle - zu \partial_u \langle M \rangle - \beta/\nu \langle M \rangle = 0$. The separability of the variables u and w, i.e., $\langle M \rangle(u,w) = M_u(u) M_w(w)$ leads to

$$x_0 w M_u/M_w = \beta/\nu + zu M_u/M_u,$$

where the prime means the derivative with respect to the argument. Since the left-hand side of this equation depends only on $w$ and the right-hand side depends only on $u$, both sides must be equal to a constant $c$. Thus, $M_u(u) = w^{L_z} - \beta/\nu v < \nu v >$ and $M_w(w) = w^{L_z} - \beta/\nu v < \nu v >$, resulting in $M(t,w) = m_0^{L_z} L^{\nu z} (t^{L_z} - \beta/\nu v < \nu v > )$. Returning to the original variables, one has $\langle M \rangle(t,L,m_0) = m_0^{L_z} L^{\nu z} (t^{L_z} - \beta/\nu v < \nu v > )$.

On one hand, by choosing $c = \xi_0$ at criticality ($\tau = 0$), one obtains $\langle M \rangle_{m_0} \sim m_0^\theta$, where $\theta = (x_0 - \beta/\nu)/v$ that corresponds to a regime of small initial magnetization soon after a finite time scaling $b = t^{1/z}$ in Eq. (4). This leads to $\langle M \rangle(t,m_0) = t^{-\beta/\nu \langle \nu \rangle} \langle M \rangle(1,m_0)$. By calling $x = t^{L_z} m_0$, an expansion of the averaged magnetization around $x = 0$ results in $\langle M \rangle(1,x) = \langle M \rangle(1,0) + \partial_x \langle M \rangle|_{x=0} + O(x^2)$. By construction $\langle M \rangle(1,0) = 0$ and, since $u = t^{L_z} m_0 \ll 1$, we can discard quadratic terms resulting in $\langle M \rangle_{m_0} \sim m_0^\theta$. This anomalous behavior of initial magnetization is valid only for a characteristic time scale $t_{\text{max}} \sim m_0^{-z/\nu}$.

On the other hand, the choice $c = 0$ corresponds to the case where the system does not depend on the initial trace and $m_0 = 1$ leads to simple power law:

$$\langle M \rangle_{m_0 = 1} \sim t^{-\beta/\nu \langle \nu \rangle},$$

(6)

that similarly corresponds to the decay of magnetization (for $t > t_{\text{max}}$) of a system previously evolved from an initial small magnetization ($m_0$), and that had its magnetization increased according to Eq. (5) up to a peak.

For $m_0 = 0$, it is not difficult to show that the second moment of the magnetization is given by

$$\langle M^2 \rangle_{m_0 = 0} \sim t^\xi,$$

(7)

with $\xi = (d - 2\beta/\nu)/z$, where $d$ is the dimension of the system. By using short-time MC simulations, when lattices are suitably prepared with a fixed initial magnetization, many authors have obtained the dynamic exponent $z$ as well as the static ones $\beta$ and $\nu$, for many different models (see, for example, two good reviews found in Refs. [24, 25]).

In order to estimate independently the critical exponents, we can, first, determine $\xi$ by using a power law that mixes initial conditions [26] as follows:

$$F_c(t) = \langle M^2 \rangle_{m_0 = 0} \sim t^\xi,$$

(8)

where $\xi = d/\nu$. With the estimate of $\xi$, denoted here by $\tilde{\xi}$, we are able to obtain an estimate of $z$ (given by $\tilde{\xi} = d/\nu$) independent of other parameters. In order to obtain $\nu$, we use an alternative power law. When considering $m_0 = 1$ in Eq. (4), one can see that there is no dependence on the initial conditions. Therefore, when $L \rightarrow \infty$, one can have $\langle M \rangle(t,\tau) = b^{-\beta/\nu \langle \nu \rangle} (b^{-1/t \cdot \xi \cdot \nu \cdot z \cdot t, b^{-1/L \cdot m_0})$. By scaling $b^{-z} = 1$, we have $\langle M \rangle(t,\tau) = t^{-\beta/\nu \langle \nu \rangle} f(t^{1/\nu \langle \nu \rangle})$ where $f(x) = \langle M \rangle(1,x)$ and so $\partial \ln \langle M \rangle(1,\tau)/\partial \tau = \xi \partial \langle M \rangle(1,\tau)/\partial \tau$. Therefore we have

$$D(t) = \frac{\partial \ln \langle M \rangle}{\partial \tau} \bigg|_{\tau=0} = f_0 \cdot t^{1/\nu \langle \nu \rangle} \sim t^\phi,$$

(9)

where $f_0 = f(0)$ is a constant and $\phi = 1/\nu \langle \nu \rangle$. Since we have already obtained the exponent $\xi$, we are able to obtain $\nu$. With these two exponents in hand, we can obtain $\beta$ by estimating the exponent $\mu = \beta/\nu \langle \nu \rangle$ from Eq. (6).
In order to simulate numerically the theoretical moments of the magnetization of the spin systems as functions of time, we used a local dynamic evolution of the spins which are updated by the heat-bath algorithm. In our simulations we used two different initial states: to obtain the power laws given by Eqs. (6) and (9), we used the initial ordered state, i.e., $m_0 = 1 (\sigma_i = 1, i = 1, ..., N = L^d)$. On the other hand, when considering Eq. (7) we used an initial state with $m_0 = 0$, i.e., the spins of each site were chosen at random on the sites but keeping the same proportion-$L^d/5$ spins of each type: $\sigma_i = 0,1,2,3,4$. Here it is important to mention that $m_0 = 0$ for any order parameter proposed in our analysis [Eqs. (2) and (3)].

In the context of time-dependent MC simulations, the magnetization ($k = 1$) and its higher moments ($k > 1$) have statistical estimators for the theoretical moments (4) given by

$$\langle M^k \rangle(t) = \frac{1}{N_{\text{run}} L^d} \sum_{j=1}^{N_{\text{run}}} \sum_{j=1}^{L^d} \sigma_{i,j}(t) \right)^k,$$

where $\sigma_{i,j}(t)$ denotes the $i$th spin variable on the lattice at the $r$th MC step of the $j$th run. Here $N_{\text{run}}$ denotes the number of different repetitions (runs) or different time series used to compute the averages.

IV. RESULTS I: EXPLORING THE PHASE DIAGRAM VIA NONEQUILIBRIUM MC SIMULATIONS

Our initial plan was to study the phase transition points of the $Z(5)$ model via time-dependent MC simulations by estimating the best $x_2$ given as input the parameter $x_1$ according to the phase diagram (see Fig. 1). We performed this task for several points in this diagram and the analysis was carried out by using an approach developed in [16] in the context of generalized statistics. This tool had also been applied successfully to study multicritical points, for example, tricritical points [15,27] and Lifshitz point of the ANNNI model [14].

Since at criticality it is expected that the order parameter obeys the power-law behavior of Eq. (6), we fixed the value of $x_1$ and changed the value of $x_2$ according to a resolution $\Delta x_2$. Then, we calculated the known coefficient of determination [28] that, for our case, is given by

$$r = \sum_{l=1}^{N_{\text{MC}}} \frac{(\ln(M) - a - b \ln t)^2}{\sum_{l=1}^{N_{\text{MC}}} (\ln(M(t)/l)^2},$$

with $\ln(M) = (1/N_{\text{MC}}) \sum_{l=1}^{N_{\text{MC}}} \ln(M(t)$, for each value $x_2 = x_2^{(\text{min})} + i \Delta x_2$, with $i = 1,...,n$, where $n = \lfloor (x_2^{(\text{max})} - x_2^{(\text{min})})/\Delta x_2 \rfloor$, and the critical value corresponds to $x_2^{(\text{opt})} = \arg \max_{x_2 \in [x_2^{(\text{min})}, x_2^{(\text{max})}]} r$. The coefficient $r$ has a very simple explanation: It measures the ratio: (expected variation)/(total variation). The bigger the $r$, the better the linear fit in log scale, and therefore, the better the power law which corresponds to the critical parameter except for an error $O(\Delta x_2)$.

As we are dealing with a rich phase diagram, a careful analysis of the order of the phase transition is necessary, mainly when taking into account first-order “critical” points. As pointed out earlier, the phase diagram of the $Z(5)$ model possesses two second-order phase transition points which coincide with the FZ integrability points, as well as two lines of infinite-order transition (dual to each other) also known as self-dual lines. The phase transitions of the points on these lines which extend from the five-state Potts point to the FZ points are expected to be of first order. Although a power-law behavior of the order parameter at strong first-order points is not expected, it is possible to obtain this behavior for weak first-order ones, whereas for $k > k_c$ a disorder metastable state vanishes at a certain $k^*$ and, for $k < k_c$, there is an ordered metastable state which disappears at $k^*$. Both parameter values look like critical points if the system remains in the disordered or ordered metastable states, and so in both points a power-law behavior must be observed as studied by Schulke and Zheng [29] through the analysis of the weakness of the first-order phase transition in the $q$-state Potts model. In that case a good estimate for $k_c$ would be $(k^* + k^*)/2$. For the five-state Potts model, for example, the difference between the pseudocritical points $k^*$ and $k^*$ and $k_c$ is in the fourth decimal digit. Moreover, the difference between power laws obtained from the pseudo-critical points and $k_c$ is observed for $t \sim 1000$ MC steps.

Since the self-dual line of the $Z(5)$ model is analytically described by $x_2 = (1/\sqrt{5} - 1/2) - x_1$ and the points extending from $x_1 = (\sqrt{5} - 1)/4 \approx 0.30901...$ (but not including) the FZ point (which corresponds to $x_1 \approx 0.3473834...$) are points of weak first-order transition, we determined the corresponding $x_2$ via the method previously described. In this case, by looking into the difference between $x_2$ (exact) and $x_2$ (simulation), it was possible to have a measure of weakness of the considered points.

In Table I, column 3, we show our results for $x_2 (x_2^{(\text{opt})})$ for five points along the self-dual line whose transitions are expected to be of first order, as well as for the FZ point (sixth line). In order to obtain these results, we used resolution of $\Delta x_2 = 0.002$ and applied a simple algorithm that makes a process of refinement of the parameter in order to localize the best $x_2$ along the simulations. These values must be compared to the exact predictions of the self-dual

<table>
<thead>
<tr>
<th>$x_1$</th>
<th>$x_2^{(\text{exact})}$</th>
<th>$x_2^{(\text{opt})}$ (simulation)</th>
<th>$r(x_2^{(\text{opt})} - \Delta x_2)$</th>
<th>$r(x_2^{(\text{opt})})$</th>
<th>$r(x_2^{(\text{opt})} + \Delta x_2)$</th>
<th>$x_2^{(\text{opt})}^{(2)}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Potts</td>
<td>0.30901...</td>
<td>0.308(2)</td>
<td>0.994 251</td>
<td>0.999 605</td>
<td>0.999 557</td>
<td>0.3094(1)</td>
</tr>
<tr>
<td>0.31</td>
<td>0.30803...</td>
<td>0.308(2)</td>
<td>0.997 386</td>
<td>0.999 514</td>
<td>0.998 977</td>
<td>0.3083(1)</td>
</tr>
<tr>
<td>0.32</td>
<td>0.29803...</td>
<td>0.298(2)</td>
<td>0.997 535</td>
<td>0.999 696</td>
<td>0.999 920</td>
<td>0.2979(1)</td>
</tr>
<tr>
<td>0.33</td>
<td>0.28803...</td>
<td>0.288(2)</td>
<td>0.998 707</td>
<td>0.999 715</td>
<td>0.998 626</td>
<td>0.2873(1)</td>
</tr>
<tr>
<td>0.34</td>
<td>0.27803...</td>
<td>0.278(2)</td>
<td>0.998 385</td>
<td>0.999 572</td>
<td>0.998 690</td>
<td>0.2781(1)</td>
</tr>
<tr>
<td>FZ</td>
<td>0.27065...</td>
<td>0.270(2)</td>
<td>0.999 401</td>
<td>0.999 701</td>
<td>0.999 168</td>
<td>0.2702(1)</td>
</tr>
</tbody>
</table>
line (column 2). It is important to notice that columns 4, 5, and 6 represent, respectively, the values of \( r \) obtained for the fits with respective values of \( x_2: x_2^{\text{opt}} - \Delta x_2, x_2^{\text{opt}}, \) and \( x_2^{\text{opt}} + \Delta x_2 \). For instance, we observe that, for the Potts point \( r(x_2^{\text{opt}} - \Delta x_2) = 0.994251 \), \( r(x_2^{\text{opt}}) = 0.999605 \), and \( r(x_2^{\text{opt}} + \Delta x_2) = 0.999557 \). From that, we applied a second refinement for the interval \([x_2^{\text{opt}} - \Delta x_2, x_2^{\text{opt}} + \Delta x_2]\) by using \( \Delta x_2 = 10^{-4} \) and we found 0.30941(1) (column 7). When compared to the exact value 0.30901..., we observed an error only in the fourth decimal place which is reasonable according to lattice used in our MC simulations for this optimization, \( L = 160 \).

Now, since we analyzed the first-order (weak) transition up to the bifurcation point, we turned our attention to points after it via time-dependent MC simulations. According to these phase diagrams (Fig. 1), after the bifurcation point, \( x_1 > 0.3473834... \), there are two second-order lines separating the ordered and disordered phases and the soft one.

For example, by applying our refinement process for \( x_1 = 0.42 \), the method produces a clear point where \( r \) is maximum \( x_1^{\text{opt}} = 0.198(2) \) [see Fig. 2(a)]. This value is in complete agreement with the exact value of the self-dual line, \( x_2 = \left(\frac{\sqrt{5} - 1}{2}\right) - 0.42 = 0.19803... \). However, it is important to notice that we did not find the two points which we would expect by looking into the phase diagram corresponding to the two critical lines. In order to better exploit such specificities, we simulated our method for two other inputs: \( x_1 = 0.44 \) and \( x_1 = 0.46 \); the first one corresponds to the end of order-disorder transition and the second one was chosen because there is no ordered phase at this point [see Figs. 2(b) and 2(c)].

In those cases we can clearly see that there is no unique point where \( r \) assumes a maximum value. Finally in Fig. 2(d) we show the behavior of this same coefficient for some important points just for an appropriated comparison: the five-state Potts model (weak first-order transition point), \( x_1 = 0.4 \) (crossing two second-order lines), \( x_1 = 0.5 \), and especially the FZ point whose critical exponents are estimated in this paper. Now we would like to consider alternatives to determine (localize) points after the bifurcation point that are localized on the soft-disorder transition line. From now on, we will be much more empirical in our techniques. As we reported above, our optimization method captures the points on the self-dual line but the points corresponding to soft-disorder and soft-order transitions seem to be neglected by the method and this deserves a better investigation.

Since we used the power laws for ordered initial spin systems, this can be the reason whereas such transitions are not order-disorder-like. In order to localize such points we prepared a second algorithm similar to the previous method. However, instead of optimizing Eq. (6), by performing several time-dependent MC simulations starting from \( m_0 = 1 \), we monitored simulations starting from \( m_0 = M_1(0) = 0 \) and, in this case, we expected that the second moment of the order parameter has the power law given by Eq. (7) (see [30]). Moreover, we also monitored the value of \( \zeta \)
and evaluation of the coefficient \( \varsigma \) different power-law fits:
two input values:

\[ \langle \frac{M}{m_0} \rangle_m \sim \frac{r_{\beta/\nu z}}{x^2} \quad \text{and} \quad \langle M^2 \rangle_m \sim \frac{r_{\beta/\nu z}}{x^2} \]

evaluation of the coefficient \( \varsigma = \frac{(d - 2\beta/\nu z)}{x^2} \) for the different values of \( x_2 \) considering as input: \( x_1 = 0.36 \) (a) and \( x_1 = 0.46 \) (b).

whereas it can be estimated, even without significance, when the coefficient of determination is not satisfactory.

Figure 3 shows the behavior of the coefficient of determination when one takes into account the power laws for \( \langle M \rangle_{m=1} \) and \( \langle M^2 \rangle_{m=0} \) along with the numerical estimates of \( \varsigma \), for two input values: \( x_1 = 0.36 \) (a) and \( x_1 = 0.46 \) (b). We can see that determination for \( \langle M \rangle_{m=0} \) for both values decreases abruptly for a value of \( x_2 \) followed by a subsequent abrupt increase. Such behavior was found for several other studied points ranging from the five-state Potts model to \( x_1 = 0.6 \). We also can see that the peak of the curves of the determination coefficient corresponds to the points where the numerical estimates of \( \varsigma \) change their signal. For instance, for \( x_1 = 0.36 \), we found a clear maximum of the determination coefficient for \( x_2 = 0.258(2) \) when we considered fits for \( \langle M \rangle \) [Eq. (6)]. On the other hand, when one considers fits for \( \langle M^2 \rangle \) [Eq. (7)] the value of \( x_2 \) at the peak of the determination coefficient \( (x_2 = 0.278(2)) \) does not coincide with the previous one.

In order to establish some relationships between the estimates of the points where there is an abrupt decreasing of coefficient \( r \) for \( \langle M^2 \rangle \) and values of the soft-disorder transition, we decided to digitize the phase diagram of the model (Fig. 1, Ref. [17]) in order to localize (by using a pointer on the bitmap figure) and compare some points of the soft-disorder phase transition to the values obtained in our simulations.

We can observe that after \( x_1 = 0.40 \) (see Table II) there is excellent agreement between unofficial estimates (Ref. [17]) and our empirical method (EM). It is important to mention that before \( x_1 = 0.44 \) our method for optimization of the power law for \( \langle M \rangle_{m=1} \) has already localized very well the considered points on the self-dual transition line. So from this analysis we have two important conclusions.

1. By taking into account points with \( (\sqrt{5} - 1)/4 < x_1 < 0.44 \), we are able to estimate the best values of \( x_2 \) which corresponds to the self-dual line by optimizing Eq. (6).

2. For \( x_1 \geq 0.40 \) we estimated some values of \( x_2 \) through Eq. (7) by using an empirical approach and analyzed the soft-disorder transition, the only transition above the self-dual line, in this region predicted by the phase diagram (see [17]).

Finally, it is important to mention a technical detail in our simulations. Here, our initial condition for obtaining \( m_0 = 0 \) for \( \langle M^2 \rangle \) was built only with spins related to the first-order parameter \( \psi_0 \) [Eq. (2)], i.e., \( n_1 = 1 \) or 2. This case does not correspond to the correct critical values of \( \beta \) and \( \nu \), whereas the correct way to vanish the initial configuration is to put \( n_1 = 0,1,2,3,4 \) in the proportion of \( 1/5 \) for each one, as used in this paper to compute the critical exponents. However, when considering the empirical method presented above this initial condition \( (n_2 = 1 \text{ or } 2) \) brings a change of signal of \( \varsigma \) which was not observed when considering the initial natural condition (proportion of \( 1/5 \)).

V. RESULTS II: ESTIMATING THE CRITICAL EXPOUNTS (STATIC AND DYNAMIC ONES) OF THE BIFURCATION POINT

Now we explored the critical exponents of the \( Z(5) \) model with special attention to the bifurcation point. Before showing the estimates for this point, we presented some estimates of the exponent \( \mu_1 = -\beta/\nu z \) from Eq. (6), with \( i = 1 \) or 2, along the self-dual line by using the two order parameters \( M_i \) [Eqs. (2) and (3)]. Our main idea here is to study the symmetry between these two order parameters via nonequilibrium MC simulations and to explore if there is some pair \( (x_1,x_2) \) for which \( \mu_1 = \mu_2 \). It is important to mention that \( \mu \) is a sort of effective exponent since it was used to analyze first weak and second-order points.

A. Exploring the self-dual line

We prepared an algorithm that measures \( \mu \) for each \( (x_1,x_2) \) pair in the self-dual line, \( x_2 = (\sqrt{5} - 1)/2 - x_1 \), and performed time-dependent MC simulation to obtain averages of the order parameter [Eqs. (2) and (3)] and, consequently, the
exponents $\mu_1$ and $\mu_2$ from the power-law decay [Eq. (6)]. For these simulations, we considered $x_1$ ranging from $x_1^{(\text{min})} = 0.2$ to $x_1^{(\text{max})} = 0.4$, with $\Delta x_1 = 5 \times 10^{-3}$. For each input pair $(x_1, x_2)$ we used $N_{\text{run}} = 1200$ runs, $N_{\text{MC}} = 150$, and $L = 160$ (enough after a fast finite size scaling study as shown in the next subsection).

In Fig. 4 we show the behavior of $\mu_1$ and $\mu_2$ as a function of $x_1$. We can observe that the curves meet each other at the point $x_1 = 0.310(5)$ which corresponds to the five-state Potts point. The inset plot represents the same time evolution of $F_2$ as a function of $t$. In order to verify the finite-size effects, we have used lattice of linear sizes, $L = 10, 20, 40, 80, 160$, and 240. In Fig. 5 we can observe robust power laws for the time evolution of the ratio $F_2$. As can be seen in the figure, the power-law behavior of the first-order parameter $M_1$ is shown as points while the second one, $M_2$, is represented by lines. Then, it is possible to notice in this figure that both order parameters share the same exponent $z$.

In our experiments we used $N_{\text{MC}} = 150$ MC steps and calculated the exponents for different time windows of size $\Delta N = 10$ MC steps with respective goodness of fit $q$. In Table III (column 3) we show the different values obtained for $z$. All intervals presented excellent goodness of fit (column 6), with $q_z > 0.73$.

Similarly, the plots in Figs. 6 and 7 show the time evolution of $D(t)$ and $M(t)$, for the two different order parameters. Here, $D(t)$ was numerically estimated according to

$$D(t) \approx \frac{1}{2\delta} \ln \left[ \frac{\langle M(t, T_c + \delta) \rangle}{\langle M(t, T_c - \delta) \rangle} \right].$$

### TABLE III. Estimates of exponents for different time windows by using the order parameter $M_1$

<table>
<thead>
<tr>
<th>Interval</th>
<th>$\phi = 1/\nu z$</th>
<th>$z$</th>
<th>$\mu = \beta/\nu z$</th>
<th>$q_{\phi/\nu z}$</th>
<th>$q_z$</th>
<th>$q_{\beta/\nu z}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>[30, 40]</td>
<td>0.666(6)</td>
<td>2.38(3)</td>
<td>0.064(2)</td>
<td>0.994</td>
<td>0.998</td>
<td>0.989</td>
</tr>
<tr>
<td>[40, 50]</td>
<td>0.649(5)</td>
<td>2.43(5)</td>
<td>0.0650(4)</td>
<td>0.998</td>
<td>1.000</td>
<td>1.000</td>
</tr>
<tr>
<td>[50, 60]</td>
<td>0.667(6)</td>
<td>2.34(6)</td>
<td>0.0650(7)</td>
<td>0.999</td>
<td>1.000</td>
<td>1.000</td>
</tr>
<tr>
<td>[60, 70]</td>
<td>0.659(6)</td>
<td>2.40(5)</td>
<td>0.065(1)</td>
<td>0.995</td>
<td>1.000</td>
<td>1.000</td>
</tr>
<tr>
<td>[70, 80]</td>
<td>0.64(1)</td>
<td>2.28(6)</td>
<td>0.066(1)</td>
<td>1.000</td>
<td>1.000</td>
<td>1.000</td>
</tr>
<tr>
<td>[80, 90]</td>
<td>0.66(2)</td>
<td>2.24(6)</td>
<td>0.066(1)</td>
<td>1.000</td>
<td>1.000</td>
<td>1.000</td>
</tr>
<tr>
<td>[90, 100]</td>
<td>0.65(2)</td>
<td>2.34(6)</td>
<td>0.067(2)</td>
<td>0.998</td>
<td>1.000</td>
<td>1.000</td>
</tr>
<tr>
<td>[100, 110]</td>
<td>0.63(2)</td>
<td>2.35(5)</td>
<td>0.065(1)</td>
<td>1.000</td>
<td>0.993</td>
<td>1.000</td>
</tr>
<tr>
<td>[110, 120]</td>
<td>0.66(1)</td>
<td>2.32(3)</td>
<td>0.067(2)</td>
<td>0.999</td>
<td>0.917</td>
<td>1.000</td>
</tr>
<tr>
<td>[120, 130]</td>
<td>0.64(2)</td>
<td>2.32(4)</td>
<td>0.066(3)</td>
<td>1.000</td>
<td>0.968</td>
<td>1.000</td>
</tr>
<tr>
<td>[130, 140]</td>
<td>0.68(2)</td>
<td>2.33(5)</td>
<td>0.066(2)</td>
<td>1.000</td>
<td>0.986</td>
<td>0.999</td>
</tr>
<tr>
<td>[140, 150]</td>
<td>0.66(1)</td>
<td>2.29(4)</td>
<td>0.067(3)</td>
<td>0.999</td>
<td>0.737</td>
<td>1.000</td>
</tr>
</tbody>
</table>

### B. The exponents $z, \nu$, and $\beta$ of the FZ point

Initially we performed simulations to obtain $F_2$ as a function of $t$. In order to verify the finite-size effects, we have used lattice of linear sizes, $L = 10, 20, 40, 80, 160$, and 240. From these simulations we considered $x_1$ ranging from $x_1^{(\text{min})} = 0.2$ to $x_1^{(\text{max})} = 0.4$, with $\Delta x_1 = 5 \times 10^{-3}$. For each input pair $(x_1, x_2)$ we used $N_{\text{run}} = 1200$ runs, $N_{\text{MC}} = 150$, and $L = 160$ (enough after a fast finite size scaling study as shown in the next subsection).

In Fig. 4 we show the estimates of the exponent $\mu = -\beta/\nu z$ (a sort of effective exponent) along the self-dual line. We can observe that curves assume the same value in $x_1 = 0.310(5)$ which corresponds to the five-state Potts point.
where \( \langle M \rangle (t, T) \pm \delta \) means the magnetizations above (below) critical temperature of a quantity \( \delta \), starting from an ordered initial state. Since our parameters are \( k_1 = J_1/k_BT \) and \( k_2 = J_2/k_BT \) a perturbation of \( \delta \) in \( T \) corresponds to \( k_1' = J_1/k_BT (\pm \delta) = k_1/(1 \pm \delta') \) and \( k_2' = k_2/(1 \pm \delta') \), where \( \delta' = \delta/T \).

In Table III we also present our results for \( \phi \) and \( \mu \) exactly as previously reported for \( z \). We can observe again good fits in all time windows. All the analysis and estimates presented above for \( M_1 \) were also performed for the second-order parameter \( M_2 \). However, for economy they were not reported here whereas a compilation of our main estimates, including \( M_1 \) and \( M_2 \), are presented in Table IV. The results from columns 2–7 are estimated by using the regular method to obtain the error bars in the context of short-time critical MC simulations, via error propagation (see first part of the appendix).

In this table, the term “best” means the best value found which reproduces the most similar conjectured values for the static parameters \( \nu \) and \( \beta \) (columns 10 and 11, respectively). The term “prop” refers to uncertainty which was calculated by error propagation. The term “aver” means the average of exponents performed from larger time windows taking the estimates from [70,80] up to [140,150].

We used an alternative method to obtain better estimates, considering the bootstrap re-sampling method for the uncertainty calculation (see second part of the appendix for a detailed description). The idea is to overcome possible statistical correlation among the exponents. The results are presented in columns 8 and 9. Our estimates by using bootstrap re-sampling (boot in Table IV) corroborate the exact values for \( \nu \) and \( \beta \).

First of all, it is important to mention that we obtained estimates of exponent \( \nu \) for both order parameters which, to our knowledge, have never been calculated. We can see values greater than estimates for the Ising model, for example, \((2.14 \lesssim \nu \lesssim 2.16)\) and the three-state Potts model \((\nu \approx 2.19)\) [26], but similar to results obtained for the four-state Potts model \((\nu \approx 2.29)\) [32]. The exponents \( \nu \), for both order parameters, are in complete agreement according to error bars. By using error propagation, our estimates for \( \beta \) (\( \beta_{\text{prop}} \)) over any criteria are rigorously according to conjecture value \( \beta = 0.08 \) for the order parameter \( M_2 \). On the contrary, although we have reasonable results for the order parameter \( M_1 \), \( \beta_{\text{prop}} = 0.107(4) \) and \( \beta_{\text{aver}} = 0.105(3) \), the error bars are not enough to cover the conjectured value \( \beta = 0.12 \).

Alternatively, with the procedure described in the second part of the appendix that combines bootstrap and selection, we have as best estimate \( \beta_{\text{best}} = 0.119(3) \) satisfying the conjecture.

We finally found \( \nu_{\text{best}} = 0.70(2) \) and \( 0.70(3) \) for \( M_1 \) and \( M_2 \), respectively, which corroborates the conjecture \( \nu = 0.7 \).

### VI. DISCUSSION AND CONCLUSIONS

In this paper we studied the phase diagram of the Z(5) model through the nonequilibrium finite size scaling study in the context of time-dependent MC simulations. We determined some critical values and weak first-order transition values along the self-dual line with special attention to the FZ point that, to our knowledge, have never been analyzed using this approach. We also determined some transition points along the soft-disorder transition line by using a nonconventional way that looks for an abrupt “depression” on the second moment for both order parameters which, to our knowledge, have never been calculated. We can see values greater than estimates for the Ising model, for example, \((2.14 \lesssim \nu \lesssim 2.16)\) and the three-state Potts model \((\nu \approx 2.19)\) [26], but similar to results obtained for the four-state Potts model \((\nu \approx 2.29)\) [32]. The exponents \( \nu \), for both order parameters, are in complete agreement according to error bars. By using error propagation, our estimates for \( \beta \) (\( \beta_{\text{prop}} \)) over any criteria are rigorously according to conjecture value \( \beta = 0.08 \) for the order parameter \( M_2 \). On the contrary, although we have reasonable results for the order parameter \( M_1 \), \( \beta_{\text{prop}} = 0.107(4) \) and \( \beta_{\text{aver}} = 0.105(3) \), the error bars are not enough to cover the conjectured value \( \beta = 0.12 \).

Alternatively, with the procedure described in the second part of the appendix that combines bootstrap and selection, we have as best estimate \( \beta_{\text{best}} = 0.119(3) \) satisfying the conjecture.

We finally found \( \nu_{\text{best}} = 0.70(2) \) and \( 0.70(3) \) for \( M_1 \) and \( M_2 \), respectively, which corroborates the conjecture \( \nu = 0.7 \).

### ACKNOWLEDGMENTS

This research was partially supported by the Conselho Nacional de Desenvolvimento Científico e Tecnológico (CNPq), by the grant 11862/2012-8. The authors thank CESUP (Super Computer Center of Federal University of Rio Grande do Sul) as well as Professor Leonardo G. Brunet (IF-UFRGS) for the available computational resources. We are grateful for support from Clustered Computing (ada.if.ufrgs.br). We also would like to thank the anonymous referees of Physical Review E for helpful suggestions.
APPENDIX

In this section we present our methods to estimate uncertainties. In this paper we used two approaches: (1) error propagation, generally used in short-time dynamics literature and (2) alternative error analysis by using bootstrap estimate.

1. Error propagation

In this paper, we used \( N_{\text{run}} = 4 \times 10^3 \) runs for the computation of averaged time series of the second moment of the order parameters, Eq. (7), in which are required disordered initial configurations, and \( N_{\text{run}} = 10^4 \) runs for experiments that demand ordered initial configurations, such as those which take into account the power laws given by Eqs. (6), (8), and (9).

The error bars were obtained from \( N_b = 5 \) different bins. Our results, presented in the following plots, correspond to five different time series, \( \hat{\beta} \) which corresponds to five different time series, \( \hat{\nu} \), and with a sampling distribution of the respective uncertainty: 

\[
\hat{\sigma}/\sqrt{N_b} = \left( \sum_{i=1}^{N_b} (M^i(t))^{(i)} - \langle M(t) \rangle \right)^{1/2},
\]

where \( (M^i(t))^{(i)} \) denotes the average of the \( k \)th moment of magnetization of the \( i \)th bin.

The exponent \( \nu \) was estimated from Eq. (8) as \( \hat{\beta} = 2/\hat{\nu} \) (by setting \( d = 2 \)) and its error \( \sigma_{\nu} \) was obtained through the equation \( \sigma_{\nu} = (2/\hat{\nu}^2)\sigma_t \), where \( \sigma_t \) is the error obtained from the power-law fit. With the estimate of \( \nu \) and its respective uncertainty in hand, we were able to obtain an estimate of \( \nu \) through the fitting of Eq. (9), i.e., \( \hat{\nu} = \hat{\phi}^{-1} \nu^{-1} \), with its respective uncertainty:

\[
\sigma_{\nu} = \left[ \frac{\hat{\phi}^{-2}\hat{\nu}^{-2}\hat{\phi}^{2}\sigma_t^2 + \hat{\phi}^{-4}\hat{\nu}^{-4}\sigma_t^2}{(\hat{\phi}^{2}\sigma_t^2)} \right]^{1/2}.\]

Now, we can estimate \( \beta \). Whereas we have in hand an estimate of \( \hat{\phi} \), we can estimate \( \beta \), where by fitting Eq. (6) \( \hat{\beta} = \hat{\mu}/\hat{\phi} \), with respective uncertainty,

\[
\sigma_{\beta} = \left[ \frac{\hat{\phi}^{-4}\hat{\mu}^{2}\sigma_t^2 + \hat{\phi}^{-4}\hat{\mu}^{2}\sigma_t^2}{(\hat{\phi}^{4}\sigma_t^2)} \right]^{1/2}.
\]

2. Alternative approach with bootstrap estimates

Now we describe an alternative analysis for estimating exponents with uncertainties calculated by the bootstrap method. Let us start by the independent exponent \( \beta \). So, instead of determining this exponent by combining five seeds which corresponds to five different time series, \( t \times F_2(t) \), and obtaining the error bars over these five seeds for each point of averaged time series, we used a different procedure. Since we have five seeds for \( \langle M^i(t) \rangle_{i=1} \) and five seeds for \( \langle M^2 \rangle_{i=1} \), we can obtain \( N_{\text{bin}} = 25 \) different time series \( t \times F_2(t) \) by crossing the seeds. So, we obtain \( N_{\text{sample}} \) different re-sampled data sets obtained with replacement. For each data set, each time series \( t \times F_2(t) \), corresponds to a specific bin \( i = 1,...,N_{\text{bin}} \), and an exponent \( \nu_i \) is calculated. Then, for every re-sampled data set would be, for example: 

\[
sample_1 = (z_1(1), z_1(2),...,z_2(1), z_2(2),...,z_{25}(2),...,z_{25}(25)),
\]

\[
sample_2 = (z_1(1), z_2(1),...,z_2(2),...,z_{25}(1), z_{25}(2),...,z_{25}(25)).
\]

So for every re-sampled data we calculate \( z_{i(1)} \) and with a sampling distribution of \( z_{i(1)} \) we calculate \( \hat{z}_{i} = (1/N_{\text{bin}}^{(\text{boot})}) \sum_{i=1}^{N_{\text{bin}}^{(\text{boot})}} z_{i(1)} \). The standard deviation of the sampling is given by \( \sigma_{\nu} = \sqrt{(N_{\text{sample}}^{(\text{boot})} - 1)^{-1} \sum_{i=1}^{N_{\text{sample}}^{(\text{boot})}} (z_{i(1)} - \hat{z}_{i})^2} \), which is a standard error of the mean (this is the more important point).

Since we obtained previously an estimate of \( \nu \), we used it as input and we calculated \( \nu^{(\text{boot})} \) by using time series \( t \times \frac{\Delta}{s} \ln[\langle M^2(t) \rangle_{i=1} - \langle M(t) \rangle_{i=1}] \). We also crossed the seeds to obtain \( N_{\text{bin}} = 25 \) bins and for each bin, a linear fit is performed producing \( \hat{\phi}_i \rightarrow \nu_i = 1/(\phi_i \cdot z) \). We repeat the re-sampling procedure in order to obtain \( \sigma_{\nu} = \sqrt{(N_{\text{sample}}^{(\text{boot})} - 1)^{-1} \sum_{i=1}^{N_{\text{sample}}^{(\text{boot})}} (\nu_{i(1)} - \nu_{i})^2} \). Finally, since we have estimates for \( \nu \) and \( \nu \) we repeat the procedures to obtain the error estimate of \( \beta \): (a) Linear fit produces \( \beta_i = \nu_i \cdot \beta_i, i = 1,...,N_{\text{bin}} \), (b) re-sampling to obtain the bootstrap estimate of the error estimate, \( \sigma_{\beta} = \sqrt{(N_{\text{sample}}^{(\text{boot})} - 1)^{-1} \sum_{i=1}^{N_{\text{sample}}^{(\text{boot})}} (\beta_{i(1)} - \beta_{i})^2} \). The only difference here is that \( N_{\text{bin}} = 5 \) since there is no crossing of seeds for this estimate.

So, our method follows the following prescription.

(1) We obtain two estimates of the dynamic exponent \( \nu \) (minimum and maximum) estimates where the error bars were obtained with bootstrap re-sampling, under \( N_{\text{sample}}^{(\text{boot})} = 10^4 \).

(2) From these two estimates (input), we obtain a list of worst and best estimates of the static exponent \( \nu \). From these estimates we select the nearest and the farthest estimates with uncertainties calculated by the bootstrap method.

(3) Finally with best and worst values of \( \nu \), our re-sampling bootstrap results in a list of worst and best estimates of \( \beta \) and its uncertainty.

For example, for the order parameter \( M_1 \) we have the results for \( \beta \) according to column 2 in Table V for the different intervals. Taking the two more different estimates (maximum and minimum) we replicated the bootstrap method in order to obtain candidate estimates for \( \nu \) and \( \beta \), which is shown in columns 3–6 in the same table. Here \( \nu_{\text{best}} \) are the values obtained for \( \nu = 2.25 \) while the values for \( \nu_{\text{worst}} \) were obtained by using \( z = 2.36 \) as input. The columns \( \beta_{\text{best}} \) and \( \beta_{\text{worst}} \) correspond to the best and worst values by using previous input values. So we choose \( \nu = 0.70(3) \) and \( \beta = 0.119(3) \) as the better estimates among best estimates. Similar analysis was performed for \( M_2 \) which is shown in columns 8 and 9 in Table IV.

<table>
<thead>
<tr>
<th>Interval</th>
<th>( z )</th>
<th>( \nu_{\text{best}} )</th>
<th>( \nu_{\text{worst}} )</th>
<th>( \beta_{\text{best}} )</th>
<th>( \beta_{\text{worst}} )</th>
</tr>
</thead>
<tbody>
<tr>
<td>[70, 80]</td>
<td>2.28(5)</td>
<td>0.71(4)</td>
<td>0.66(4)</td>
<td>0.117(3)</td>
<td>0.104(5)</td>
</tr>
<tr>
<td>[80, 90]</td>
<td>2.25(5)</td>
<td>0.69(1)</td>
<td>0.64(1)</td>
<td>0.117(1)</td>
<td>0.104(1)</td>
</tr>
<tr>
<td>[90, 100]</td>
<td>2.35(4)</td>
<td>0.70(3)</td>
<td>0.65(2)</td>
<td>0.119(3)</td>
<td>0.106(5)</td>
</tr>
<tr>
<td>[100, 110]</td>
<td>2.36(7)</td>
<td>0.71(2)</td>
<td>0.66(2)</td>
<td>0.116(3)</td>
<td>0.103(5)</td>
</tr>
<tr>
<td>[110, 120]</td>
<td>2.32(6)</td>
<td>0.69(2)</td>
<td>0.64(2)</td>
<td>0.117(3)</td>
<td>0.104(5)</td>
</tr>
<tr>
<td>[120, 130]</td>
<td>2.32(6)</td>
<td>0.71(3)</td>
<td>0.66(3)</td>
<td>0.117(3)</td>
<td>0.104(2)</td>
</tr>
<tr>
<td>[130, 140]</td>
<td>2.30(5)</td>
<td>0.68(2)</td>
<td>0.63(2)</td>
<td>0.117(1)</td>
<td>0.104(1)</td>
</tr>
<tr>
<td>[140, 150]</td>
<td>2.27(4)</td>
<td>0.69(2)</td>
<td>0.64(2)</td>
<td>0.118(8)</td>
<td>0.105(3)</td>
</tr>
</tbody>
</table>
Here it is important to mention that the lattice was randomly vanished by considering only two spin variables, \( n_i = 1 \) and \( 2 \), different than the experiments performed to calculate the critical exponents where the lattice was vanished by putting \( 1/5 \) of spin variables of each kind. Our choice was based on numerical experiments that were shown to be appropriate for this kind of analysis. On the other hand, for the former prepared initial configurations, the exponent \( \zeta \) probably does not correspond to the correct value \( (d - 2\beta/\nu)/z \). However, this does not forbid our approach whereas in this stage of the paper, our aim was only to explore alternatives for the localization of the critical points and not to estimate critical exponents which was correctly performed in the appropriate section.