Specific-heat exponent for the three-dimensional Ising model from a 24th-order high-temperature series

Gyaő Bhanot
Thinking Machines Corporation, 245 First Street, Cambridge, Massachusetts 02142
and Institute for Advanced Study, Princeton, New Jersey 08540

Michael Creutz
Brookhaven National Laboratory, Upton, New York 11973

Uwe Glässner and Klaus Schilling
Physics Department, University of Wuppertal, Gaussstrasse 20, 42097 Wuppertal, Germany
(Received 3 December 1993)

We compute high temperature expansions of the three-dimensional Ising model using a recursive transfer-matrix algorithm and extend the expansion of the free energy to 24th order. Using inhomogeneous-differential Padé and ratio methods, we extract the critical exponent of the specific heat to be \( \alpha = 0.104(8) \).

I. INTRODUCTION

High- and low-temperature expansions constitute major tools for the calculation of critical properties in statistical systems. The Ising and Potts model low-temperature expansions were recently extended\(^1\)-\(^4\) using a technique based on the method of recursive counting.\(^5\) In a separate development, Vohwinkel\(^6\) implemented the shadow-lattice technique of Domb\(^7\) in a very clever way and added many new terms to the series. However, the extraction of critical parameters from low-temperature series is hampered by the presence of unphysical singularities. This is especially true of the three-dimensional (3D) Ising model. For this reason, low-temperature analytic methods are very often inferior to Monte Carlo methods for computing critical exponents.

High-temperature (HT) expansions, on the other hand, generally have better analytic behavior and yield more accurate exponents. Very recently, two variants of the recursive counting technique for HT expansions have been pursued. While Guttmann and Enting\(^8\) keep track of spin configurations on a set of rectangular finite lattices, Ref. 3 counts HT graphs on finite, helical lattices. Such computer based series expansions have very large memory requirements. This makes them ideal candidates for large parallel computers if communication issues can be handled efficiently. In this paper we will present the results of a HT expansion of the 3D Ising model to 24th order, obtained on a 32 node 1 GByte Connection Machine CM-5. The implementation is based on a bookkeeping algorithm of binary coded spin configurations in helical geometry.

II. COMPUTATION OF THE SERIES

We start with a discussion of the HT algorithm to compute the partition functions on finite 3D Ising lattices.

Starting from the action

\[
E(s) = -\sum_{(i,j)} s_is_j, \tag{1}
\]

the partition function is

\[
Z = \sum_{(s)} \exp(-\beta E) = \sum_{(s)} \prod_{(i,j)} \exp(\beta s_is_j) \tag{2}
\]

and is expanded in a HT series

\[
Z = (\cosh^3 \beta)^V \sum_{(s)} \prod_{(i,j)} (1 + s_is_j t) = (2 \cosh^3 \beta)^V \sum_k p(k) t^k, \tag{3}
\]

with the HT expansion parameter \( t = \tanh \beta \). \( V \) is the volume of the system. The free energy per spin is defined as

\[
f = -\frac{1}{\beta V} \ln Z = -\frac{2 \cosh^3 \beta}{\beta} = -\frac{1}{\beta} \sum_k f_k t^k. \tag{4}
\]

For simplicity, consider a finite simple cubic lattice which, in the recursion algorithm, is built up by adding one site after the other, layer by layer. This procedure defines the recursion step, which requires knowledge only of those spin states that are contained in the exposed two-dimensional surface layer. To minimize finite-size effects, it is best to use helical boundary conditions.\(^2\)\(^3\) One can visualize helical boundary conditions by imagining all spins in the layer laid out along a straight line. In this picture, the nearest neighbors to a given site in the sequence in the ith direction can be chosen to be \( h_i \) sites away, with \( i = x, y, z \). It is convenient to assume \( h_x < h_y < h_z \). It is easy to see that, as spins are added, one needs only to keep track of the states of spins on the
TABLE I. Structures and weights $w$ of the lattices used.

<table>
<thead>
<tr>
<th>$h_x$</th>
<th>9</th>
<th>1</th>
<th>9</th>
<th>5</th>
<th>7</th>
<th>10</th>
<th>5</th>
<th>14</th>
<th>11</th>
<th>14</th>
<th>9</th>
<th>9</th>
<th>5</th>
<th>5</th>
<th>16</th>
<th>10</th>
<th>16</th>
<th>1</th>
<th>17</th>
</tr>
</thead>
<tbody>
<tr>
<td>$h_y$</td>
<td>11</td>
<td>12</td>
<td>14</td>
<td>15</td>
<td>13</td>
<td>15</td>
<td>15</td>
<td>16</td>
<td>17</td>
<td>16</td>
<td>17</td>
<td>19</td>
<td>20</td>
<td>20</td>
<td>21</td>
<td>21</td>
<td>22</td>
<td>22</td>
<td>22</td>
</tr>
<tr>
<td>$w$</td>
<td>-3</td>
<td>3</td>
<td>-3</td>
<td>3</td>
<td>-3</td>
<td>3</td>
<td>-3</td>
<td>3</td>
<td>-1</td>
<td>-2</td>
<td>-1</td>
<td>1</td>
<td>-2</td>
<td>5</td>
<td>2</td>
<td>-2</td>
<td>2</td>
<td></td>
<td></td>
</tr>
</tbody>
</table>

The topmost $h_z$ sites. Let these spins be denoted $s_1, \ldots, s_{h_z}$. Then the partition function can be rewritten as

$$Z = (2 \cosh^3 \beta)^V \sum_{k} \sum_{s_1, \ldots, s_{h_z}} p(k; s_1, \ldots, s_{h_z}) t^k. \quad (5)$$

The recursion step, which consists of adding another spin $s_0$ to the system, changes the partition function into

$$Z = 2^V (\cosh^3 \beta)^{V+1} \sum_{s_0} \sum_{k} \sum_{s_1, \ldots, s_{h_z-1}} p(k; s_1, \ldots, s_{h_z-1}, s_0) t^k (1 + s_0 s_{h_z} t) (1 + s_0 s_{h_z} t) (1 + t) \times \left[ x^2 + p(k; s_1, \ldots, s_{h_z-1}, s_0) t^k (1 + s_0 s_{h_z} t) (1 + t) \right]. \quad (7)$$

The contribution in the second (third) line of this equation contains the part with $s_{h_z}$ being parallel (antiparallel), denoted by $s_0$ to $s_0$. Comparing this expression with the HT series (5) for the new system yields the recursion relation induced for the coefficients $p$,

$$2p(k; s_0, s_1, \ldots, s_{h_z-1}) = p(k - 0; s_1, \ldots, s_{h_z-1}, s_0) + p(k - 0; s_1, \ldots, s_{h_z-1}, s_0)$$
$$+ p(k - 1; s_1, \ldots, s_{h_z-1}, s_0) (s_0 s_{h_z} + s_0 s_{h_z} + 1)$$
$$+ p(k - 1; s_1, \ldots, s_{h_z-1}, s_0) (s_0 s_{h_z} + s_0 s_{h_z} - 1)$$
$$+ p(k - 2; s_1, \ldots, s_{h_z-1}, s_0) (s_0 s_{h_z} + s_0 s_{h_z} + s_0 s_{h_z})$$
$$+ p(k - 2; s_1, \ldots, s_{h_z-1}, s_0) (s_0 s_{h_z} - s_0 s_{h_z} - s_0 s_{h_z})$$
$$+ p(k - 3; s_1, \ldots, s_{h_z-1}, s_0) (s_0 s_{h_z} + s_0 s_{h_z} + s_0 s_{h_z}). \quad (8)$$

It is crucial to remove finite-size errors by combining the results of different lattice structures as described in Refs. 2 and 3. We use the set of lattices listed in Table I and obtain the free energy coefficients up to 24th order as given in Table II. In order to eliminate the contribution from (unphysical) loops with an odd number of links in any direction, we use the cancellation technique of Ref. 3. This amounts to inserting additional signature factors into Eq. (6) for each of the three link factors

$$(1 + s_0 s_{h_z} t) \to \sigma_i (1 + s_0 s_{h_z} t), \quad i = x, y, z \quad (9)$$

with $\{\sigma_x, \sigma_y, \sigma_z\} = \{\pm, \pm, \pm\}$. By performing eight separate runs corresponding to all possible values of $\sigma$ and adding the results, one achieves a complete elimination of the unwanted loops. Possible contributions of higher-order finite-size loops are at least of order 25 for this set of lattices. Since we use open boundary conditions, the coefficients $p$ are invariant under the global transformation $s_i \to -s_i$. This $Z(2)$ symmetry enables us to reduce memory requirements by a factor of 2. Unlike Refs. 2,3,4 we use multiple-word arithmetic to account for the size of the coefficients. This implementation needs about 100% more memory but leads to a doubling in performance. Since the number of words can be adjusted separately for every order, the computational effect can be reduced.

TABLE II. Free energy up to 24th order.

<table>
<thead>
<tr>
<th>Order $k$</th>
<th>Free energy $f_n$</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>0</td>
</tr>
<tr>
<td>2</td>
<td>0</td>
</tr>
<tr>
<td>4</td>
<td>3</td>
</tr>
<tr>
<td>6</td>
<td>22</td>
</tr>
<tr>
<td>8</td>
<td>375/2</td>
</tr>
<tr>
<td>10</td>
<td>1980</td>
</tr>
<tr>
<td>12</td>
<td>24044</td>
</tr>
<tr>
<td>14</td>
<td>319170</td>
</tr>
<tr>
<td>16</td>
<td>18059031/4</td>
</tr>
<tr>
<td>18</td>
<td>201010408/3</td>
</tr>
<tr>
<td>20</td>
<td>5162283633/5</td>
</tr>
<tr>
<td>22</td>
<td>16397040750</td>
</tr>
<tr>
<td>24</td>
<td>260938787382</td>
</tr>
</tbody>
</table>
acCORDINGLY. ON THE 32 NODE CM-5 THE TOTAL TIME FOR ALL COMPUTATIONS WAS ABOUT 50 h.

COMParED TO THE finite-LATtICE appRACh OF GUTTMANN and ENTing, 4 our method appears to require more CPU time since we need to cancel unphysical loops. IT should be noted, however, that helical lattices ARE truly implemented in data parallel software environments and thus lead to better performance. IN THE usual finite-LATtICE method, 6 the HT EXPANsion can only be extended in fairly coarse steps, using lATTICES with (4 × 5) cross section for 22nd order and (5 × 5) cross section for 26th order, respectively. FOR this reason, a 24th-order computation would not have been feasible using that method with our computer resources.

III. eRITical eXponent

THE specific heat is defined as

\[ c|_{\beta=0} = \beta^2 \frac{\partial^2}{\partial^2 T} \ln Z = \sum c_k \beta^2 \]

and is expected to behave near \( T_C \) as

\[ c|_{\beta=0} = A(T)|T - T_C|^{-\alpha} [1 + B(T)|T - T_C|^\theta + \ldots] \]

with \( A \) and \( B \) being analytic near \( T_C \). 9,10 We analyze the series using unbiased and biased inhomogeneous-differential Padé approximants (IDP's) 11 as well as ratio tests.

A. Padé Analysis

IN Fig. 1 we plot \( \alpha \) against \( t_c^2 \) for each IDP approximant \([J/L;M]\). FITting the linear dependence of \( \alpha \) on \( t_c^2 \), we find

\[ \alpha = 0.102 \pm 0.008 \] at the value \( t_c = 0.218092 \) as obtained in Monte Carlo simulations. 12 A direct, biased-IDP analysis was also performed. We obtained \( \alpha = 0.109 \pm 0.016 \).

IDP's can also be used to predict the most significant digits of the next term in the specific-heat series. 4 The estimate of the 24th-order term as obtained in Ref. 4 agrees perfectly with our exact result. Using the same method we can estimate the 26th-order term in the expansion to be

\[ f_{26} = 443762(4) \times 10^7 \]

where the errors quoted are two standard deviations.

B. ratio test

THE main problem in the determination of critical exponents in the low-temperature case is the presence of unphysical singularities nearer to the origin than the physical one. Since the expansion coefficients \( c_n \) are dominated by these unphysical singularities, ratio methods cannot be applied.

IN the HT expansion, the physical singularity dominates the asymptotic behavior, so that the ratio \( r_n = c_n/c_{n-1} \) of successive coefficients of the series is expected to behave as 9

\[ r_n = \frac{1}{t_c^{2 \nu}} \left[ 1 + \frac{\nu - 1}{n \nu} + \frac{c}{n^{1+\nu}} + \frac{d}{n^{1+2\nu}} + O \left( \frac{1}{n^{1+3\nu}} \right) \right]. \]

Assuming that the correction-to-scaling exponent \( \theta \) is close to 0.5, 12,13 the following sequence \( s_n \) is expected to converge toward \( \alpha \) like

\[ s_n := (t_c^2 r_n - 1) n + 1 = \alpha + \frac{c}{n^{1/2}} + \frac{d}{n} + O \left( \frac{1}{n^{3/2}} \right). \]
A plot of this sequence against $n$ is shown in Fig. 2. Obviously the first four values are dominated by higher-order corrections. To obtain estimates for $\alpha$ we therefore use only the values $\{s_6, \ldots, s_{13}\}$. A three-parameter least-square-fit using the ansatz of Eq. (15) yields the values shown as diamonds in Fig. 3. The value of $\alpha = 0.113$ obtained by the fit to the points $\{s_6, \ldots, s_{11}\}$ is in perfect agreement with the result of Ref. 4. Their estimate of $\alpha = 0.110$ using the extrapolated term $s_{12}$ appears to be slightly above our value of $\alpha = 0.108$ using the exact term. Including our value for $s_{13}$ of the ID Padé extrapolation Eq. (13) we obtain $\alpha = 0.105(2)$. The error represents the uncertainty of the extrapolation. However, from Fig. 3 it is quite suggestive that the $\alpha$ values might converge to a value below 0.105.

To get an estimate of the uncertainties of our results, we investigate the stability of the fits. For this purpose, we repeat the analysis after eliminating the point $s_6$ from the data. As a result we obtain sizable changes for $\alpha$. The new data are shown as crosses in Fig. 3.

In Fig. 4 we present the results for the first correction-to-scaling coefficient $c$ from our three-parameter fits. In contrast to Ref. 4, our values suggest that $c$ changes sign with increasing $n_{\text{max}}$. Because of the sensitivity of the
fits to the number of terms we keep, it is difficult to determine the value of $c$ very precisely. Our best estimate is $c = 0.01(4)$. Since $c$ vanishes within error, it seems reasonable to also try a two-parameter ansatz with $c = 0$ to fit the data. The results of these fits are shown in Fig. 5. We now find that the fits are much more stable and the $\alpha$ estimates show much more of a convergence to their asymptotic values. The best value (from the largest $n_{\text{max}}$) is $\alpha = 0.1045(3)$. This value supports the impression of the three-parameter fits, which suggested that $\alpha$ was slightly below 0.105. Finally we investigate the influence of the uncertainty in the correction-to-scaling exponent $\theta$ on our results. Repeating the analysis with $\theta = 0.53^{12}$ we find a change on $\alpha$ of less than 0.0005. Taking into account the fact that neglecting $c$ causes an additional systematic error, our final estimate for the critical exponent is,

$$\alpha = 0.104(4).$$

(16)

### IV. DISCUSSION AND OUTLOOK

The crucial element in the estimate of the error in $\alpha$ [Eq. (16)] is our neglect of the correction-to-scaling coef-
The resulting systematic error is rather large. From Fig. 4 one might speculate that the estimates for $c$ begin to exhibit asymptotic behavior at the 26th order. Therefore an exact calculation of the 26th term of the expansion might reduce the uncertainty of $c$ significantly. If the magnitude of $c$ turns out to be really negligible, one could adopt the errors of the linear fits, and $\alpha$ would be obtained accurate to the fourth significant digit.

ACKNOWLEDGMENTS

This work was partly funded under Contracts No. DE-AC02-76CH00016 and DE-FG02-90ER40542 of the U.S. Department of Energy. The work of G.B. was also partly supported by a grant from the Ambrose Monell Foundation. U.G. and K.S. are grateful to Deutsche Forschungsgemeinschaft for its support to the Wuppertal CM-Project.