[NO NAME ARTICLE]

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This is an abstract and that's really abstract.

I. INTRODUCTION

This is cite-checking.[1]

II. MODELS AND METHODS

In the paper we consider several models: the first one is Ising model on interacting self-avoiding walk from the [REF KAMILLA], on three different lattices: 2D-square lattice, 3D-square lattice and 2D-triangle lattice. The main difference between square and triangle lattice in defining two additional diagonal monomers on lattice as nearest too, see [FIGURE FOR 2D-SQUARE AND 2D-TRIANGLE LATTICE]. Considering the case of lack of outer magnetic field in this work, the Hamiltonian of the model of fixed conformation u with length N and strength of nearest-neighbors interaction J reads:

$$H_{u,N,\{\sigma\}} = -\sum_{\langle i,j\rangle} J\sigma_i\sigma_j, \quad i,j \in u, \ |u| = N$$
 (1)

The summation runs through spins involved in conformation and only with the nearest neighbors.

The second model considered in this paper is the Ising model on the rectangular lattice from the [REF SELKE]. Simulated lattices has $L \times rL$ spins and the Hamiltonian is calculated through interaction between all spins and their nearest neighbors respectively:

$$H_{L,r,\{\sigma\}} = -\sum_{\langle i,j\rangle} J\sigma_i\sigma_j, \quad i,j \in [1..L] \times [1..rL]$$
 (2)

For comparing magnetic properties of models with similar geometric ones we also define shape factors, such as gyration tensor [REF PELLISETTO]:

$$Q_{N,\alpha\beta} = \frac{1}{N+1} \sum_{i=0}^{N} (w_{i,\alpha} - w_{c,\alpha})(w_{i,\beta} - w_{c,\beta})$$
 (3)

where N is length of the system (number of monomers in conformations of Ising-ISAW models and number of spins in the lattice in rectangular Ising), and α , β are coordinates (so, $Q_{N,xx}$ and $Q_{N,yy}$ can be defined as mean squares of coordinates of the points of the model in the cartesian coordinate system with the center in the center of model). Eugen values q1, q2 of given tensor are used

in enumerating another important shape factor - mean asphericity[REF PELLISETTO]:

$$\mathcal{A} = \left\langle \frac{(q_1 - q_2)^2}{(q_1 + q_2)^2} \right\rangle_N \tag{4}$$

The compared magnetic property of our models is the fourth order cumulant of the magnetization of the Binder cumulant, defined as [REF SELKE]:

$$U_4 = 1 - \frac{\langle m^4 \rangle}{3\langle m^2 \rangle^2} \tag{5}$$

Where $\langle m^4 \rangle$ and $\langle m^2 \rangle$ are mean fourth and second order of mean magnetization per spin respectively.

We also need to define mean proportion of monomers with fixed number i of nearest neighbors $\langle n_i \rangle$, which is counted directly for every monomer in every simulated conformation of walk.

We are interested in comparing models in their respective critical regions. For each structure, critical temperatures of Ising models are known as [REF FOSTER] [REF SELKE]:

Type of lattice	T_c
2D-square	1.199 ± 0.003
3D-square	1.90 ± 0.02
Rectangular	2.26918

TABLE I. Known values of critical temperature of different modifications of Ising-ISAW model and normal Ising on the rectangular lattice

Method. All methods of simulating for Ising model on a SAW conformation are remained the same as in [REF KAMILLA]. For simulating Rectangular Ising, however, we used cluster update of Wolff Algorithm [REF NEW-MAN BARKEMA] [REF SCHRODINGER CAT].

III. RESULTS

A. Mean Asphericity and Critical Cumulant

B. Bulk

IV. DISCUSSION

V. ACKNOWLEDGMENTS

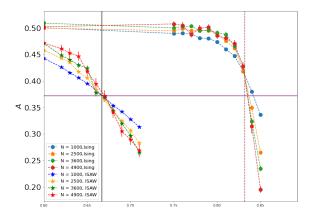


FIG. 1. Caption

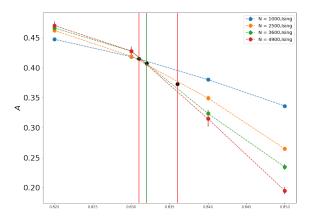


FIG. 2. Caption

PolIsing			
J	\mathcal{A}	r	U_4 Rectangular
0.831	0.415	0.465	0.340 ± 0.006
0.832	0.4072	0.47	0.343 ± 0.006
0.836	0.373	0.490 ± 0.002	0.348 ± 0.006
ISAW			
0.667	0.375	0.49	0.349 ± 0.006

TABLE II. Values of critical cumulant for Ising model on rectangular lattice with mean asphericity related to models ISAW and Ising-ISAW in their critical regions

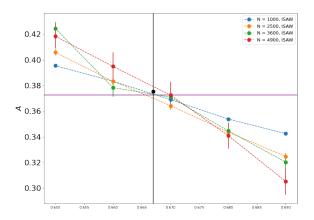


FIG. 3. Caption