14.4 Monte Carlo simulation of 2-d U(1) lattice gauge theory

José Antonio García Hernández

August 13, 2024

We call $U_{\mu}(x)$ to a U(1) variable at point x in the μ direction. For U(1), the lattice gauge action reads

$$S[U] = \beta \sum_{x} \sum_{\mu < \nu} \text{Re} \left[1 - U_{\mu\nu}(x) \right],$$
 (1)

where $\beta = \frac{2}{e^2}$ and e being the gauge coupling. The plaquettes are defined by

$$U_{\mu\nu}(x) = U_{\mu}(x)U_{\nu}(x+\hat{\mu})U_{\mu}^{\dagger}(x+\hat{\nu})U_{\nu}^{\dagger}(x)$$
 (2)

The sum of plaquette variables is computed using

$$S_{\rm P} = \sum_{x} \sum_{\mu < \nu} \text{Re } U_{\mu\nu}(x), \tag{3}$$

so the expectation value of an elementary Wilson loop or plaquette reads

$$E_{\rm P} = \frac{\langle S_{\rm P} \rangle}{DV},\tag{4}$$

where $V = L^d$ is the volume of the lattice and $D = \frac{d(d-1)}{2}$ is the number of planes of rotation in d dimensions.

We define the sum of staples as

$$\Sigma_{\mu}(x) = \sum_{\nu \neq \mu} \left[U_{\nu}(x) U_{\mu}(x+\hat{\nu}) U_{\nu}^{\dagger}(x+\hat{\mu}) + U_{\nu}^{\dagger}(x-\hat{\nu}) U_{\mu}(x-\hat{\nu}) U_{\nu}(x+\hat{\mu}-\hat{\nu}) \right]. \tag{5}$$

The change in the action by a local update when changing $U_{\mu}(x) \to U'_{\mu}(x)$ is

$$\Delta S = -\beta \operatorname{Re} \operatorname{Tr} \left[\left(U'_{\mu}(x) - U_{\mu}(x) \right) \Sigma_{\mu}^{\dagger} \right]. \tag{6}$$

We use the Metropolis algorithm for the simulation as decribed in the following steps:

- 1. Given some gauge field configuration, we go through all the lattice points in a lexicographic way. At each point x in the μ direction we generate a random U(1) variable $U'_{\mu}(x)$.
- 2. We compute the sum of the staples and compute the change of the action according to eq. (6). We generate a uniform random number $r \in [0, 1)$ and accept the change if $r \leq p$, where p is

$$p = \min(1, \exp(-\Delta S)). \tag{7}$$

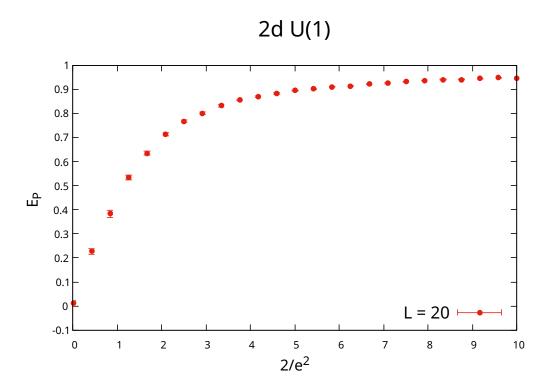


Figure 1: Expectation value of the plaquette variable vs. $\beta=2/e^2$ with a volume $V=20^2$. The errors were computed using the Jackknife error. The parameters of the simulation were: at each value of β , 5000 measurements were made separated by 20 sweeps and a thermalization of 1000 sweeps.