The speed of time in an interacting particle system

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1 Comparison of the speed of two models

Let (Λ, E) be a transitive graph. For each $i \in \Lambda$ we let \mathcal{N}_i denote the set of vertices adjacent to i. We will be interested in an interating particle system on the lattice (Λ, E) whose local state space is $S = \{-1, +1\}^2$. We denote an element of S^{Λ} as (x, y) where $x = (x(i))_{i \in \Lambda} \in \{-1, +1\}^{\Lambda}$ and $y = (y(i))_{i \in \Lambda} \in \{-1, +1\}^{\Lambda}$. For $x \in \{0, 1\}^{\Lambda}$ and $i \in \Lambda$, we let

$$F_i(x) := \frac{\#\{j \in \mathcal{N}_i : x(j) = x(i)\}}{\#\mathcal{N}_i}.$$

denote the fraction of neighbors of i where x has the same value as in i. We also write

$$1_{\{x(i)=y(i)\}} := \left\{ \begin{array}{ll} 1 & \text{if } x(i)=x(j), \\ 0 & \text{otherwise,} \end{array} \right.$$

and we define $1_{\{x(i)\neq y(i)\}}$ similarly. There are two parameters $\alpha, \beta \geq 0$. Each vertex $i \in \Lambda$ becomes active at the times of independent Poisson point set with rate one. Once a vertex i is active, we either flip x(i), that is, we change its value from whatever it was to the other possible value, or we flip y(i), or we do nothing. The probabilities are as follows:

- x(i) flips with probability $\frac{1}{2}e^{-\beta F_i(x)} \alpha 1_{\{x(i)=y(i)\}}$,
- y(i) flips with probability $\frac{1}{2}e^{-\beta F_i(y)} \alpha 1_{\{x(i)\neq y(i)\}}$.

In the note aper.pdf a model is considered such that for each $i \in \Lambda$:

- x(i) flips with rate $e^{-x(i)[\beta'M_i(x) + \alpha'y(i)]}$,
- y(i) flips with rate $e^{-y(i)[\beta' M_i(y) \alpha' x(i)]}$,

where for $x \in \{0,1\}^{\Lambda}$ and $i \in \Lambda$, we call

$$M_i(x) := \frac{1}{\# \mathcal{N}_i} \sum_{j \in \mathcal{N}_i} x(j)$$

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the local magnetisation of x around i. We observe that

$$M_i(x) = x(i) [2F_i(x) - 1]$$
 and $x(i)y(i) = 21_{\{x(i) = y(i)\}} - 1 = 1 - 21_{\{x(i) \neq y(i)\}}$.

Using the fact that $x(i)^2 = 1$ always, it follows that

$$e^{-x(i)[\beta' M_i(x) + \alpha' y(i)]} = e^{-\beta' x(i) M_i(x) - \alpha' x(i) y(i)]}$$

$$= e^{-\beta' [2F_i(x) - 1] - \alpha' [21_{\{x(i) = y(i)\}} - 1]} = e^{\beta' + \alpha'} e^{-2\beta' F_i(x) - 2\alpha' 1_{\{x(i) = y(i)\}}}.$$

and similarly

$$e^{-y(i)[\beta' M_i(y) - \alpha' x(i)]} = e^{-\beta' y(i) M_i(y) + \alpha' x(i) y(i)]}$$

= $e^{-\beta' [2F_i(y) - 1] + \alpha' [1 - 21_{\{x(i) \neq y(i)\}}]} = e^{\beta' + \alpha'} e^{-2\beta' F_i(y) - 2\alpha' 1_{\{x(i) \neq y(i)\}}}.$

Thus, if we set $\alpha' := \frac{1}{2}\alpha$ and $\beta' := \frac{1}{2}\beta$, then this is the same model as before, except that all rates are a factor

$$_{2e}(\beta+\alpha)/2$$

larger. So for example, the model with $\beta=3$ and $\alpha=1$ corresponds to the model in aper.pdf with $\beta'=3/2$ and $\alpha'=1/2$, except for a factor

$$2e^{(\beta+\alpha)/2} = 2e^2 \approx 14.778$$

in the speed of time.

2 Faster simulation

In numerical simulations, the transitive graph (Λ, E) is finite. In each step, we choose a point (i, σ) uniformly from $\Lambda \times \{1, 2\}$. Then we apply the following rules:

- If $\sigma = 1$, then we flip x(i) with probability $e^{-\beta F_i(x) \alpha \mathbb{1}_{\{x(i) = y(i)\}}}$.
- If $\sigma = 2$, then we flip y(i) with probability $e^{-\beta F_i(y)} \alpha 1_{\{x(i) \neq y(i)\}}$.

We now measure time in steps of size

$$\frac{1}{\#(\Lambda \times \{1,2\})} = \frac{1}{2\#\Lambda}.$$

When α and β are large, these probabilities are quite small for a lot of the sites, which means that most of the time our program does nothing. We can speed the program up as follows. We define:

$$R(i,\sigma) := \begin{cases} e^{-\beta F_i(x) - \alpha \mathbb{1}_{\{x(i) = y(i)\}}} & \text{if } \sigma = 1, \\ e^{-\beta F_i(y) - \alpha \mathbb{1}_{\{x(i) \neq y(i)\}}} & \text{if } \sigma = 2. \end{cases}$$

and set

$$R := \sum_{\sigma=1}^{2} \sum_{i \in \Lambda} R(i, \sigma).$$

Instead of choosing (i, σ) uniformly from $\Lambda \times \{1, 2\}$, we now choose (i, σ) with the probability

$$\frac{R(i,\sigma)}{R},$$

and after we have chosen (i, σ) , we always flip, i.e., if $\sigma = 1$ we flip x(i) and if $\sigma = 2$ we flip y(i). This yields roughly the same as the previous model, provided we measure time in steps of size

 $\frac{1}{R}$.

Note that R can change during our simulation, so not all time steps have the same size!