# Cluster algorithms for the Ising model

Monte Carlo methods which overcome the problem of critical slowing down close to second order phase transitions

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- Autocorrelation, Binning Analysis
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Phase transitions

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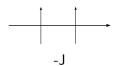
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Figure : Ernst Ising (1900 - 1998)

$$\mathscr{H}(\sigma) = -J \sum_{\langle i,j \rangle} \sigma_i \sigma_j - H \sum_i \sigma_i$$

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### Ising model - Canonical ensemble

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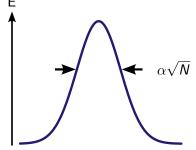
$$p(\sigma, T) = \frac{e^{-\beta \mathscr{H}(\sigma)}}{\mathscr{Z}(T)}, \quad \beta = \frac{1}{k_B T}$$

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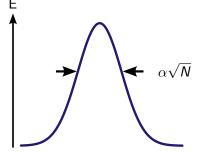
$$p(\sigma, T) = \frac{e^{-\beta \mathcal{H}(\sigma)}}{\mathcal{Z}(T)}, \quad \beta = \frac{1}{k_B T}$$
$$\langle M \rangle_T = \sum_{\sigma} M(\sigma) p(\sigma, T)$$

• We can't compute all configurations  $(2^N)$ 

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- ► We can't sample uniformely distributed over energy



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► Solution: Importance sampling using Metropolis algorithm

1. Choose one site (uniformly randomly)

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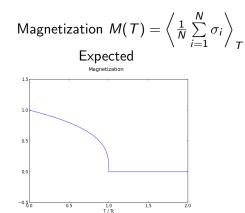
Magnetization 
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Expected

Magnetization

1.0 T / Tc

-0.5



# Obtained Magnetization 1.5 1.0 0.5 0.0 1.2 3 4 5 6

# Ising model - 2 spins

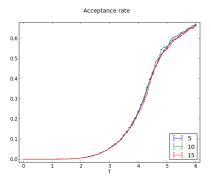
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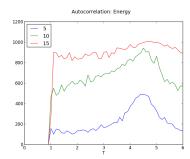
$$\mathcal{H}(\sigma) = -J \sum_{\langle i,j \rangle} \sigma_i \sigma_j$$



$$\phi_{\mathcal{A}}(t) = rac{\langle \mathcal{A}(t_0)\mathcal{A}(t)
angle - \langle \mathcal{A}
angle^2}{\langle \mathcal{A}(t_0)^2
angle - \langle \mathcal{A}
angle^2}$$

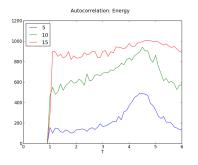
$$\phi_{\mathcal{A}}(t) = rac{\langle \mathcal{A}(t_0)\mathcal{A}(t) \rangle - \langle \mathcal{A} \rangle^2}{\langle \mathcal{A}(t_0)^2 \rangle - \langle \mathcal{A} \rangle^2} \propto \mathrm{e}^{-\frac{t}{\tau_{\mathcal{A}}}}$$

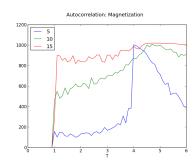
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### Autocorrelation time

$$\phi_{A}(t) = rac{\langle A(t_0)A(t)
angle - \langle A
angle^2}{\langle A(t_0)^2
angle - \langle A
angle^2} \propto e^{-rac{t}{ au_A}}$$





$$Var X := E\left[X^2\right] - E\left[X\right]^2$$

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  $(\Delta X)^2 = rac{\mathsf{Var} X}{N} \left(1 + 2 au_X
ight)$ 

$$A_i^{(I)} = \frac{1}{2} \left( A_{2i-1}^{(I-1)} + A_{2i}^{(I-1)} \right)$$

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$$\tau_A = \lim_{I \to \infty} \left( \frac{2^I \mathsf{Var} A^{(I)}}{\mathsf{Var} A^{(0)}} - 1 \right)$$

# What really happens in the System

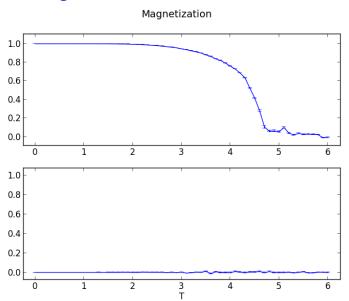
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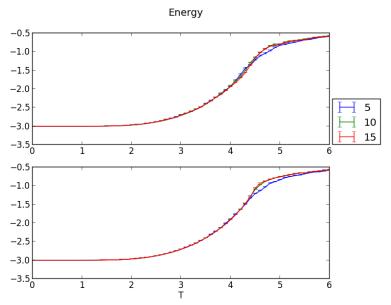
$$P(\sigma_x, \sigma_y) = 1 - \min(1, e^{-2\beta\sigma_x\sigma_y})$$

- 1. Choose one site (uniformly randomly)
- 2. Flip its spin and add it to the cluster
- 3. For all sites in the cluster:
  - 3.1 Visit every unknown neighbour, flip its spin and add it to the cluster with probability given above

### Why cluster algorithms are better

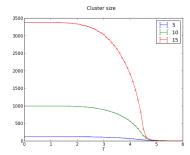


### Energy



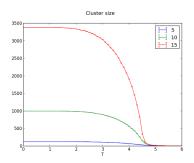
### Cluster size

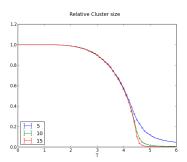
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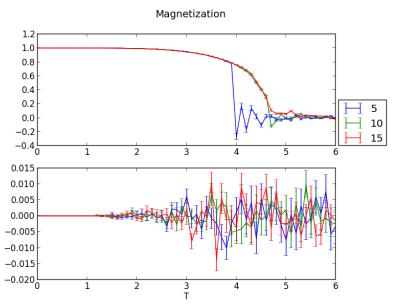
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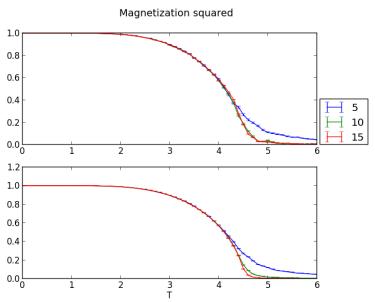




### Magnetization



### Magnetization squared



# Computation time per spin flip

Similar to Wolff, but:

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- Always touches every spin

# Questions

# Binning analysis in detail

$$VarX := E[X^{2}] - E[X]^{2}$$

$$(\Delta X)^{2} = \frac{1}{N^{2}} \sum_{i,j=1}^{N} \left( E[X_{i}X_{j}] - E[X]^{2} \right)$$

$$= \frac{VarX}{N} + \frac{1}{N^{2}} \sum_{i \neq j} \left( E[X_{i}X_{j}] - E[X]^{2} \right)$$

$$= \frac{VarX}{N} + \frac{2}{N^{2}} \sum_{i=1}^{N} \sum_{t} \left( E[X_{i}X_{i+t}] - E[X]^{2} \right)$$

$$:= \frac{VarX}{N} (1 + 2\tau_{X})$$

### Wolff or Swendsen-Wang?

Swendsen-Wang better for parallelization because it touches the whole lattice.

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