The transition probability can typically be written as

$$P(A \to B) = P_{\text{attempt}}^A(B) P_{\text{accept}}^A(B)$$

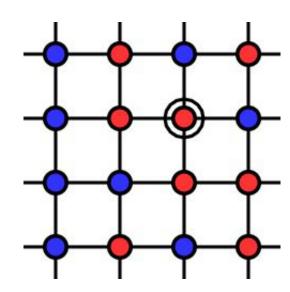
where the two factors have the following meaning:

- $P_{
 m attempt}^A(B)$ The probability of selecting B as a candidate among a number of possible new configurations
 - $P_{
 m accept}^A(B)$ The probability of actually making the transition to B after the selection of B has been done

If B has been selected but is not accepted (rejected); stay with A

For an Ising model

- ➤ Select a spin at random as a candidate to be flipped (attempt)
- ➤ Actually flip the spin with a probability to be determined (accept)
- ➤ Stay in the old configuration if the flip is not done (reject)



 $P_{\text{attempt}}^{A}(B) = 1/N_{\text{spin}}$ uniform, independent of A, B

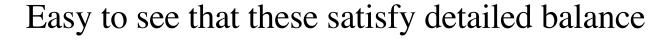
 $P_{\text{accept}}^A(B)$ constructed to satisfy detailed balance condition

$$\frac{P_{\text{accept}}^{A}(B)}{P_{\text{accept}}^{B}(A)} = \frac{W(B)}{W(A)}$$

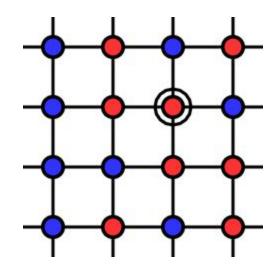
Two commonly used acceptance probabilities

Metropolis:
$$P_{\text{accept}}^A(B) = \min \left[\frac{W(B)}{W(A)}, 1 \right]$$

Heat bath:
$$P_{\text{accept}}^A(B) = \frac{W(B)}{W(A) + W(B)}$$



The ratios involve the change in energy when a spin has been flipped (or, more generally, when the state has been updated in some way)



Metropolis algorithm for the Ising model

Spin update

$$E = J \sum_{\langle i,j \rangle} \sigma_i \sigma_j, \quad \sigma_i = \pm 1$$

- Select a spin at random
- Calculate the change in energy if the spin is flipped
- Use the energy change to calculate the acceptance probability P
- Flip the spin with probability P; stay in old state with 1-P
- Repat from spin selection

Current configuration: S

Configuration after flipping spin j: S_j

Acceptance probability: $P(S \to \tilde{S}_j) = \min \left[\frac{W(\tilde{S}_j)}{W(S)}, 1 \right]$ $W(S) = e^{-E(S)/T} = e^{-\frac{J}{T} \sum \sigma_i \sigma_j} = \prod e^{-\frac{J}{T} \sigma_i \sigma_j}$

Only factors containing spin j survive in W-ratio

$$\frac{W(\tilde{S}_j)}{W(S)} = \exp\left[+\frac{2J}{T} \sigma_j \sum_{\delta(j)} \sigma_{\delta(j)} \right], \quad \delta(j) = \text{neighbor of } j$$

We want a simulation "time" unit which is normalized by the system size N (probability of a given spin having been selected after a time unit should be N independent).

1 Monte Carlo step (MC steps): N random spin-flip attempts

Measurements of physical observables done after equilibration

➤ the correct Boltzmann distribution is approached after some time that depends on the initial configuration

Binning: Accumulate data over bins consisting of M MC steps

> Averages and statistical errors calculated from bin averages

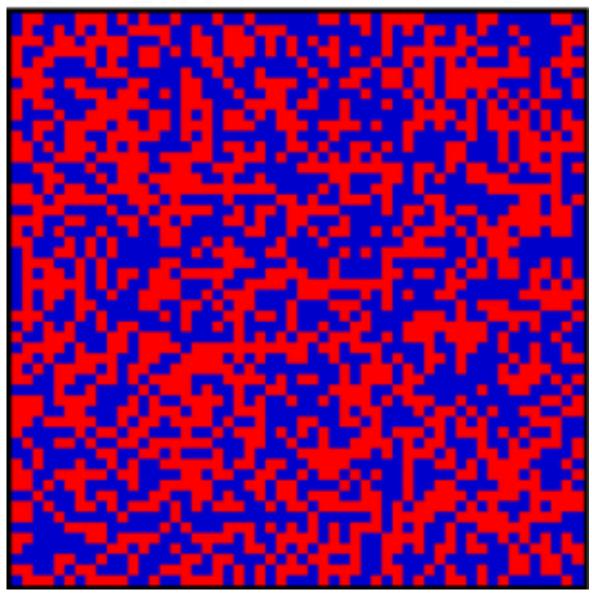
Flow of a complete simulation

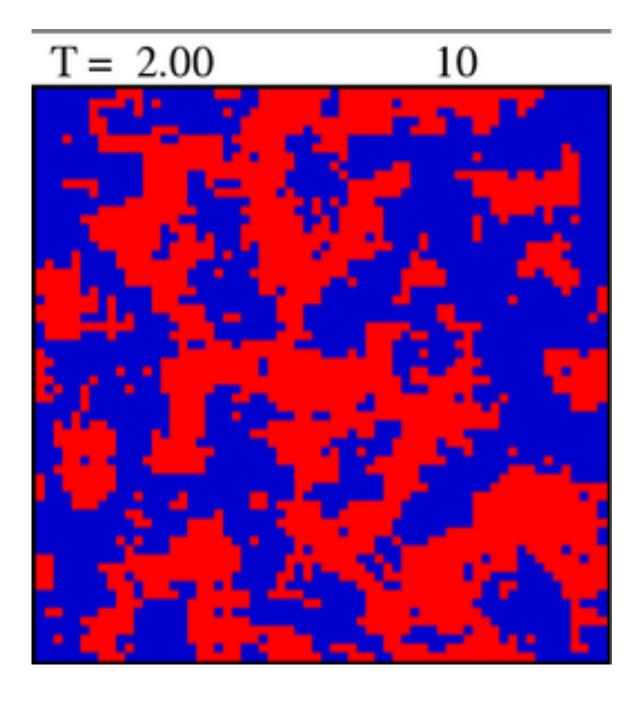
- Generate arbitrary starting state
- Carry out a number of MC steps for equilibration
- Carry out a number of bins
 - each bin consists of M MC steps
 - measurements done after every (or every few) MC step
 - save bin averages in a file after each bin (or save internally in program)
- Calculate averages and statistical errors

T = 4.00

T = 2.30

T = 2.3010 T = 2.00





T = 1.00

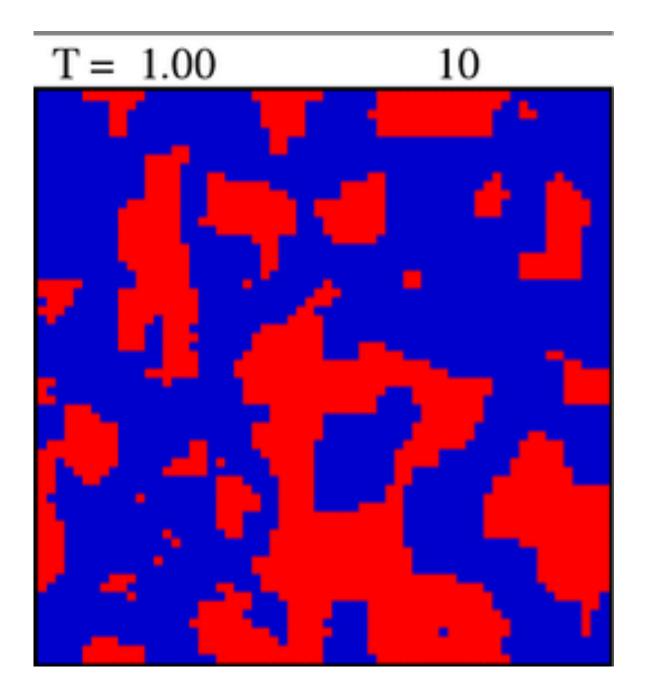
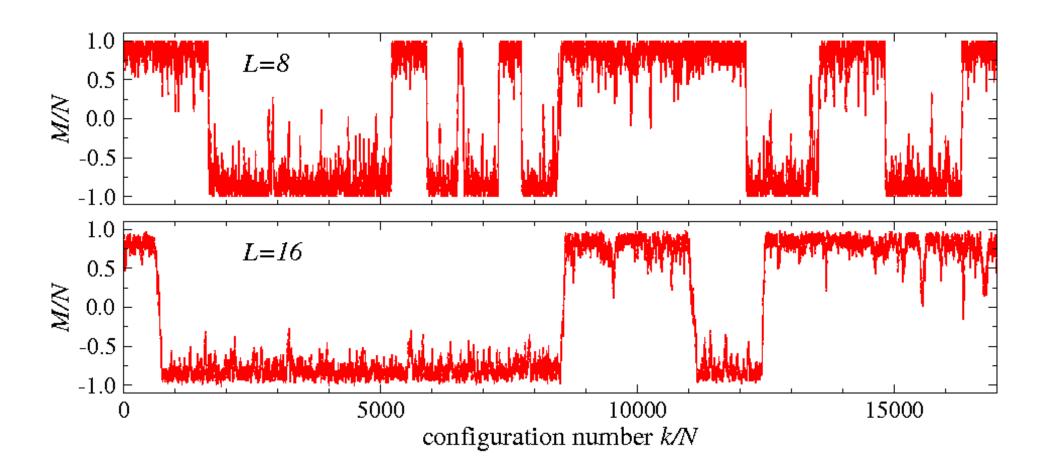


Illustration of simulation

Evolution of the magnetization, 2D Ising model, T/J=2.2 (below Tc)

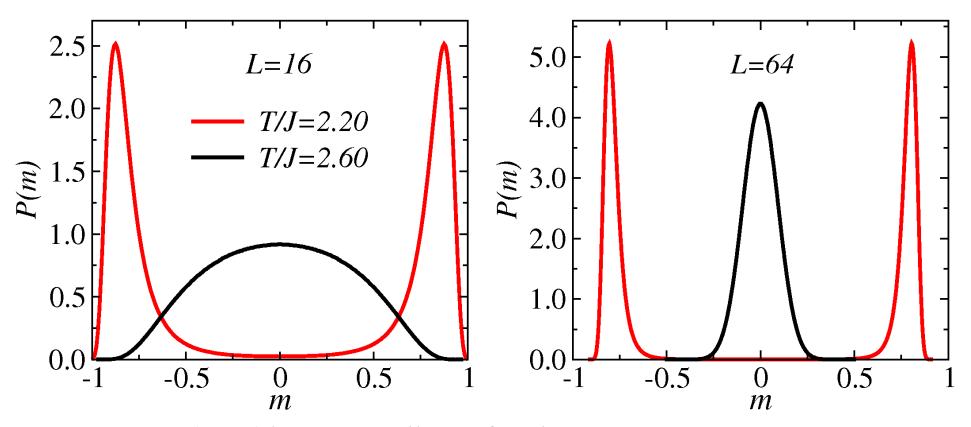
- <M>=0, but time scale for M-reversal increases with L
- Symmetry-breaking occurs in practice for large L



Magnetization distribution P(m)

The distribution depends on T and L

- single peak around m=0 for T>Tc
- double peak around +<m> and-<m> for T<Tc



Symmetry breaking (sampling of only m>0 or m<0 states) occurs in practice for large L

- because extremely small probability to go between them