

# Nucleation in the Ising Model

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# Chapter

Ising Model

History

Lattice & lattice sites

Energy & Monte Carlo

Temperature

Summary

Nucleation

Nucleation in the Ising Model

# History

- ▶ 1924: Ernst Ising - *Beitrag zur Theorie des Ferromagnetismus*<sup>1</sup>  
"Es entsteht ... [durch] ... die Beschränkung der Wechselwirkung auf diejenige benachbarter Elemente [...] kein Ferromagnetismus."
- ▶ 1936: Rudolph Peierls - *On Ising's model of ferromagnetism*<sup>2</sup>  
"[...] for sufficiently low temperatures the Ising model in two [or more] dimensions shows ferromagnetism [...]."

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<sup>1</sup>Zeitschrift für Physik Februar–April 1925, Volume 31, Issue 1, pp 253-258

<sup>2</sup>Cambridge Philosophical Society 1936 / Volume 32 / Issue 03 / October

# Lattice

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Figure: Lattice in 1 dimension

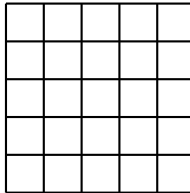


Figure: Lattice in 2 dimensions

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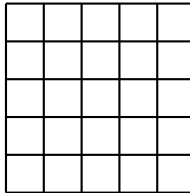


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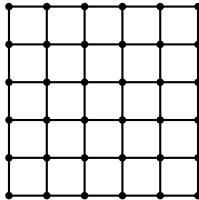


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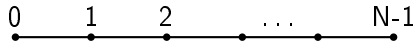


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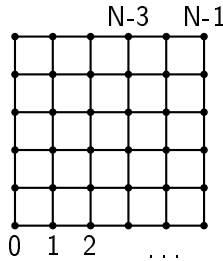


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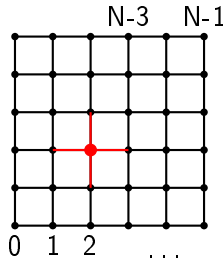


Figure: Lattice in 2 dimensions



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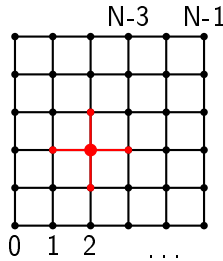


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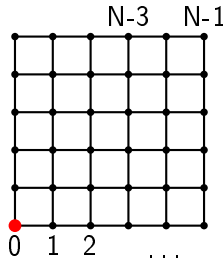


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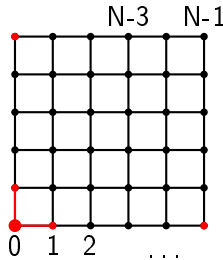


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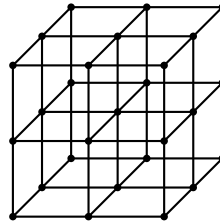
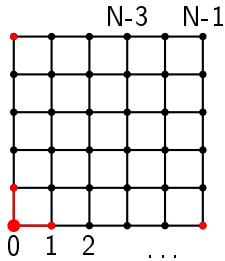


Figure: Lattice in 2 and 3 dimensions

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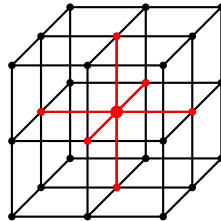
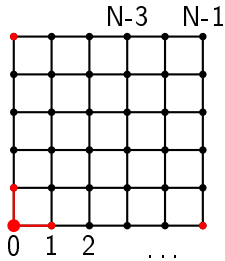


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# Lattice Sites

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- ▶ Assignment of states  $s = (s_0, s_1, s_2, \dots, s_{N-1})$  to the lattice sites is called a configuration
- ▶ Therefore  $2^N$  unique configurations for a lattice with N lattice sites.

# Energy

- ▶ Each configuration has a corresponding energy - the Hamiltonian  $H$

$$H(s) = H(s_0, s_1, \dots, s_{N-1}) = -J \sum_{\langle i,j \rangle} s_i \cdot s_j - h \sum_i s_i$$



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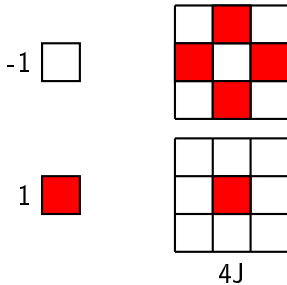
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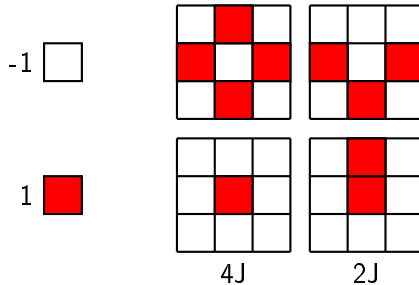
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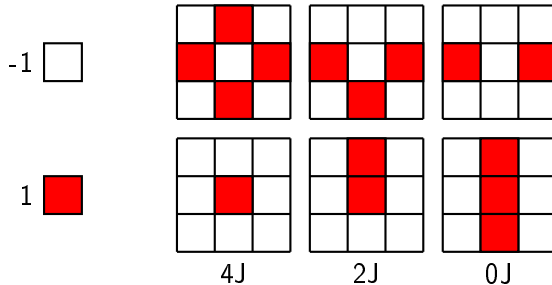
**Figure:** Energy contribution (from nearest neighbor interaction) of the central lattice site

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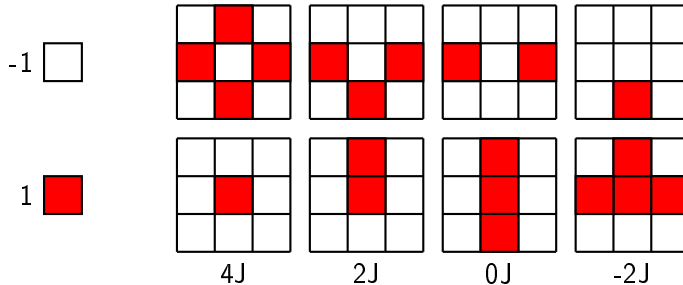
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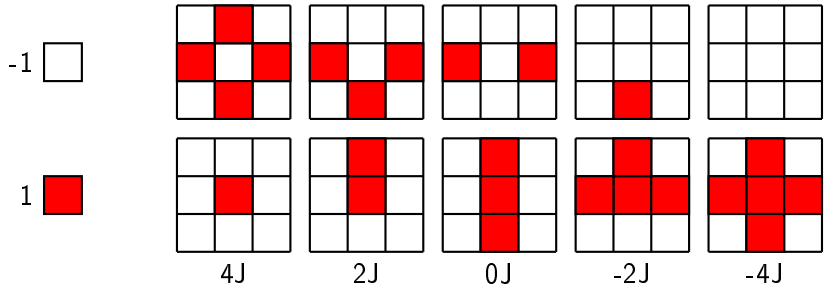
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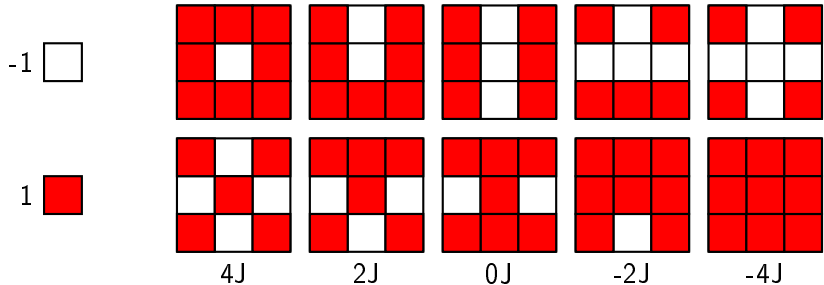
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## Single Flip Monte Carlo (Metropolis- Hastings Algorithm)

- ▶ Current configuration is  $s_t$ 
  - ▶ first configuration arbitrary
- ▶ Flip one lattice site  $\rightarrow s_f$ 
  - ▶ has to be chosen randomly - suitable RNG necessary
- ▶ Calculate energy difference  $\Delta H = H(s_f) - H(s)$
- ▶ Calculate acceptance probability  $P$

$$P = \min \left( 1, e^{-\beta \cdot \Delta H} \right), \quad \beta = 1/kT > 0$$

- ▶ Generate random number  $r \in [0, 1[$ 
  - ▶  $r < P \rightarrow s_{t+1} = s_f$
  - ▶  $r > P \rightarrow s_{t+1} = s_t$

## Temperature in the Ising Model

$$P = \min \left( 1, e^{-\beta \cdot \Delta H} \right), \quad \beta = 1/kT > 0$$

- ▶  $\Delta H < 0 \rightarrow P = 1$
- ▶ high temperature leads to higher acceptance probability  $\rightarrow$  unordered (low magnetization, Curie Temperature  $T_c$ )
- ▶ critical temperature  $T_c$  when  $\langle \sum_i^N s_i \rangle / N \approx 0$

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  - ▶  $kT_c/J = 2.269$

## Summary - Ising Model

- ▶ molecules on a lattice - each with with one of two possible states
- ▶ (magnetic) moments prefer to align
- ▶ low temperatures: ordered
- ▶ high temperatures: disordered

# Chapter

Ising Model

Nucleation

What is Nucleation?

Homogeneous Nucleation

Heterogeneous Nucleation

Nucleation in the Ising Model

Outlook

# Nucleation

- ▶ is a phase transformation process
- ▶ phase transformation grows from small nucleus

## Examples

- ▶ cloud chamber
- ▶ supercooled water

# Nucleation

- ▶ Homogeneous nucleation
  - ▶ in a uniform substance
  - ▶ no nucleation until nucleus with critical size "appears" (due to stochastic processes)
  - ▶ higher supersaturation leads to smaller critical radius.
  - ▶ rarely occurs in nature
- ▶ Heterogeneous Nucleation
  - ▶ begins at some preferable interface and grows from there
  - ▶ much (!) more likely
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# Chapter

Ising Model

Nucleation

Nucleation in the Ising Model

Homogeneous Nucleation

Transition Path Sampling

Heterogeneous Nucleation

Forward Flux Sampling

Outlook

# Homogeneous Nucleation in the Ising Model

- ▶ Necessary modifications:
  - ▶ none
- ▶ Problems
  - ▶ long time until nucleus of critical size appears
  - ▶ inefficient to simulate billions of cycles until phase change takes place

## Cluster size

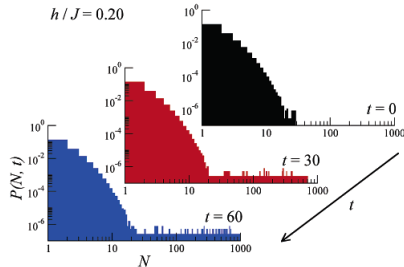


Figure: Propability of finding a cluster of size  $N$  at different times<sup>3</sup>

<sup>3</sup>Pan, Rappl, Chandler, Balsara: J. Phys. Chem. B 2006

## Transition Path Sampling (TPS) - "shooting" method<sup>4</sup>

- ▶ needs two stable states (A & B)
- ▶ path through configuration space connecting these
- ▶ change the path a little at a random point between A and B
- ▶ sample new path and accept if it connects A with B

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<sup>4</sup>aaDellago, Bolhuis, Chandler: Advances in Chemical Physics 123 (1998)

## Transition Path Sampling

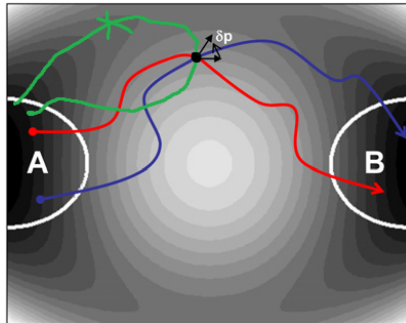


Figure: First path (red), slightly changed and accepted path (blue), rejected path (green)<sup>5</sup>

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<sup>5</sup>Esobedo, Borrero, Araque - J. Phys.: Condens. Matter 21 (2009)

# Heterogeneous Nucleation

- ▶ Necessary modifications:
  - ▶ handle boundaries in heterogeneous nucleation

$$H(s) = -J \sum_{\langle i,j \rangle} s_i \cdot s_j - h \sum_i s_i - J_s \sum_{\langle i,j \rangle} s_i \cdot s_j$$

- ▶ implement walls/surfaces with fixed spins

## Nucleation in and out of Pores<sup>6</sup>

$$H(s) = -0.8 \sum_{\langle i,j \rangle} s_i \cdot s_j - 0.05 \sum_i s_i - 0 \sum_{\langle i,j \rangle} s_i \cdot s_j, \quad kT = 1$$

- ▶ nucleation near surfaces  $10^{12}$  times faster
- ▶ fastest in pores
- ▶ nucleation in 2 steps
- ▶ diversified pore sizes lead to fastest reaction as probability of existence optimal pore size is higher

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<sup>6</sup>Page, Sear - Heterogeneous Nucleation in and out of Pores PRL 97, 065701 (2006)

## Nucleation in and out of Pores<sup>6</sup>

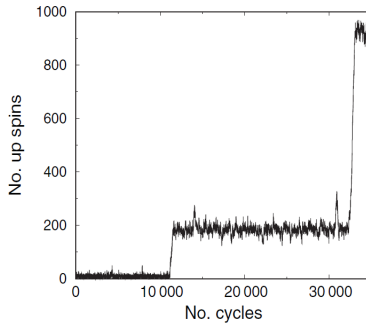


Figure: 2 phases of nucleation

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<sup>6</sup>Page, Sear - Heterogeneous Nucleation in and out of Pores PRL 97, 065701 (2006)



# Problems

- ▶ phase transitions are rare events (with realistic values for the coupling constant, ...)
- ▶ nonequilibrium systems therefore TPS (transition path sampling) not applicable.

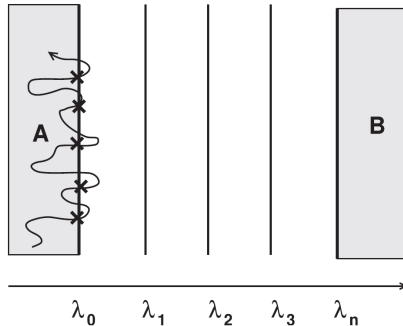
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- ▶ → Forward Flux Sampling

## Forward Flux Sampling

- ▶ Similar to TIS (transition interface sampling - a modified TPS)
  - ▶ initial state A:  $\lambda < \lambda_A = \lambda_0$
  - ▶ final state B:  $\lambda > \lambda_B = \lambda_n$
  - ▶ path has to pass every  $\lambda_i$  in increasing order (can go backwards in between too) until it reaches  $\lambda_n$  (B)
- ▶ after reaching a new interface ( $\lambda_{i+1}$ ) configuration is stored
- ▶ stored configurations used as starting point for new trial runs
- ▶ trial runs continued until path reaches A ( $\rightarrow$  failure) or a new interface  $\lambda_{i+1}$  ( $\rightarrow$  success)

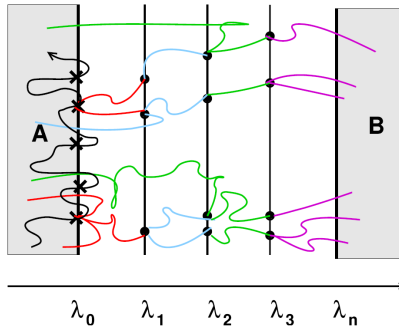
## (Direct) Forward Flux Sampling<sup>7</sup>



**Figure:** Sampling path starting in A - store configurations where the path leaves A (X)

<sup>7</sup>Allen, Valeriani, Rein ten Wolde: 2009 J. Phys.: Condens. Matter 21

## (Direct) Forward Flux Sampling<sup>7</sup>



**Figure:** Sampling new paths from every stored configuration. Discard if path goes back to A

<sup>7</sup>Allen, Valeriani, Rein ten Wolde: 2009 J. Phys.: Condens. Matter 21

# Chapter

Ising Model

Nucleation

Nucleation in the Ising Model

Outlook

Limitations of the Ising Model

Potts Model

# Limitations of the Ising Model

- ▶ only nearest neighbor interaction (can be changed to e.g. next-nearest neighbor interaction)
- ▶ no forces from the outside (can be changed with additional term in the Hamiltonian)
- ▶ → more qualitative than quantitative results

## Potts Model

- ▶ states not only  $-1 \wedge 1$  but (discrete) angles.
- ▶  $H = -J_c \sum_{i,j} \cos(\theta_i - \theta_j) + \dots$
- ▶ Applications
  - ▶ percolation (Wu: "Percolation and the Potts Model" (1978))
  - ▶ flow of foam (Sanyal, Soma: "Viscous instabilities in flowing foams" (2006))
  - ▶ cancerous tumors (Sun, Chang, Cai: "A Discrete Simulation of Tumor Growth Concerning Nutrient Concentration" (2004))



- ▶ Page, Sear - Heterogenous Nucleation in and out of Pores (2006): PRL 97, 065701
- ▶ Allen, Valeriani, Rein ten Wolde - Forward Flux Sampling for rare event simulations (2009): J. Phys.: Condens. Matter 21 (2009) 463102 (21pp)
- ▶ Allen, Frenkel, Rein ten Wolde - Forward Flux Sampling-type schemes for simulating rare events: Efficiency analysis (2008): <http://arxiv.org/abs/cond-mat/0602269v1>
- ▶ Escobedo, Borrero, Araque - Transition path sampling and forward flux sampling. Applications to biological systems 2009 J. Phys.: Condens. Matter 21 333101

- ▶ Sourcefiles and binaries on my github  
<https://github.com/oerpli/Ising2D>