

# 1 Canonical treatment of paramagnetism (6 points)

As a classical model for paramagnetism one can consider a system of  $N$  particles with the Hamiltonian

$$\mathcal{H} = -hM \quad \text{with} \quad M = \mu \cdot \sum_{i=1}^N \cos \theta_i \quad (1)$$

where  $h$  is an external homogeneous magnetic field,  $\mu$  the magnetic moment of a single particle and  $\theta_i$  the angle between the magnetic field  $h$  and the magnetic moment  $\mu$  of particle  $i$ .

**1. Use the canonical distribution to calculate the average magnetization,  $\langle M \rangle$ , as a function of  $h$  and temperature  $T$ . (3 points)**

Partition function:

$$Z = \frac{1}{h^{3N}} \cdot \int d^{3N}r \int d^{3N}p \int d^N\Omega e^{-\beta\mathcal{H}} \quad (2)$$

$$= Z_0 \cdot \left( \int_0^{2\pi} d\varphi \int_0^\pi d\theta \cdot \sin \theta \cdot e^{h\beta\mu \sum_i \cos \theta_i} \right) \quad (3)$$

$$= Z_0 \cdot \left( \int_0^{2\pi} d\varphi \int_0^\pi d\theta \cdot \sin \theta \cdot e^{h\beta\mu \cos \theta} \right)^N \quad (4)$$

$$= Z_0 \cdot \left( \frac{4\pi \cdot \sinh(\beta\mu h)}{\beta\mu h} \right)^N \quad (5)$$

Here,  $Z_0$  is the part of the partition function that does not include the integration over orientations.

Free energy:

$$F = -k_B T \ln(Z) \quad (6)$$

$$= -k_B T \cdot \left( \ln(Z_0) + N \ln(4\pi) \right) - N k_B T \ln \left( \frac{\sinh(\beta\mu h)}{\beta\mu h} \right) \quad (7)$$

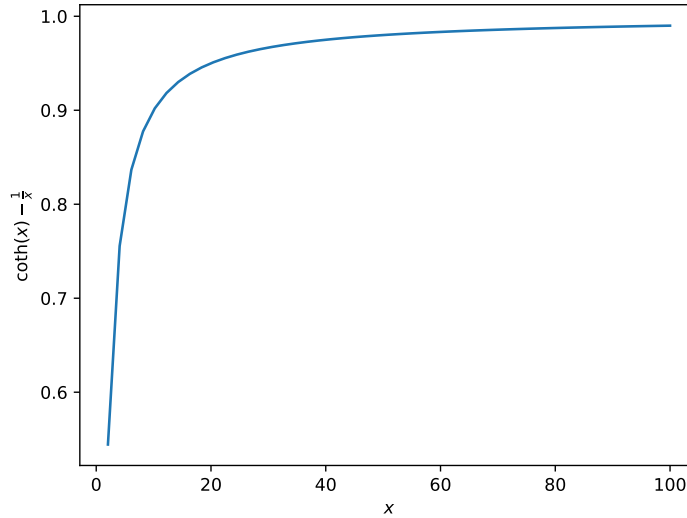
Magnetization:

$$M = -\frac{\partial F}{\partial h} \Big|_{T,V,N} = N\mu \cdot \left( \coth(x) - \frac{1}{x} \right) \quad (8)$$

with  $x := \beta\mu h$ .

**2. The ratio of which quantities determines the average magnetization? Sketch the functional dependence of the average magnetization on this ratio. (1 point)**

The average magnetization is determined by the ratio  $\beta\mu h = \mu h/k_B T$ .



**3. Discuss the two limiting cases: high temperature/weak field vs. low temperature/strong field. (2 points)**

For  $T \rightarrow 0$ , i.e.  $x \rightarrow \infty$ :

$$\coth(x) = \frac{e^{2x} + 1}{e^{2x} - 1} \approx \frac{e^{2x}}{e^{2x}} = 1 \quad \text{and} \quad 1/x \rightarrow 0 \quad (9)$$

$$\Rightarrow M \approx N\mu \quad (10)$$

As expected, all spins are aligned to the external magnetic field.

For  $T \rightarrow \infty$ , i.e.  $x \rightarrow 0$ :

$$\coth x = \frac{e^{2x} + 1}{e^{2x} - 1} \approx \frac{x}{2} \quad (\text{Laurent power series}) \quad (11)$$

$$\Rightarrow M \approx \frac{1}{2} N \beta \mu^2 h \rightarrow 0 \quad (12)$$

For high temperatures, the mean magnetization is zero, due to the non-uniform distribution of  $\theta_i$ .

## 2 One-dimensional lattice gas (6 points)

Consider a one-dimensional lattice model for a non-ideal gas with  $N$  lattice sites and periodic boundary conditions. Each lattice site  $i$  is either empty (occupancy  $n_i = 0$ ) or occupied by at most one atom (occupancy  $n_i = 1$ ). There is an attractive energy  $J$  between atoms occupying neighbouring sites. The chemical potential of the atoms is  $\mu$ . The Hamiltonian of this lattice gas is

$$H = -J \cdot \sum_{\langle ij \rangle} n_i n_j - \mu \cdot \sum_i n_i, \quad (13)$$

where  $\sum_{\langle ij \rangle}$  is the sum over all pairs of neighbouring sites.

**1. Express the partition sum of the one-dimensional lattice gas in terms of the transfer matrix  $T$ . Calculate the transfer matrix  $T$  and its eigenvalues. (2 points)**

The partition sum is given by

$$Z_N = \sum_{\{n_i\}} e^{-\beta H} = \sum_{n_1=0,1} \dots \sum_{n_N=0,1} e^{-\beta H}. \quad (14)$$

Let us rewrite the Hamiltonian in the following manner:

$$\begin{aligned} H &= -J \sum_{\langle ij \rangle} n_i n_j - \mu \sum_{i=1}^N n_i \\ &= -J \sum_{i=1}^N n_i n_{i+1} - \frac{1}{2} \mu \sum_{i=1}^N (n_i + n_{i+1}). \end{aligned} \quad (15)$$

This is equivalent since we have periodic boundary conditions ( $n_{N+1} = n_1$ ) and the lattice is one-dimensional. Now, we can define

$$-\beta H = K \sum_{i=1}^N n_i n_{i+1} + \frac{1}{2} L \sum_{i=1}^N (n_i + n_{i+1}), \quad (16)$$

where  $K = \beta J$  and  $L = \beta \mu$ . This allows us to write the argument of the partition sum as

$$e^{-\beta H} = T_{1,2} \cdot T_{2,3} \cdot \dots \cdot T_{N,1}, \quad (17)$$

where

$$T_{i,i+1} = e^{K n_i n_{i+1} + \frac{1}{2} L (n_i + n_{i+1})} \quad (18)$$

is the transfer matrix from lattice site  $i$  to site  $i + 1$ . From the fact that there are two possible values for  $n_i$  and  $n_{i+1}$  (unoccupied or occupied) which are independent of the index  $i$ , one can conclude that the  $2 \times 2$  transfer matrix is independent of the index  $i$ . Thus, one can write

$$T = \begin{bmatrix} 1 & e^{\frac{1}{2}L} \\ e^{\frac{1}{2}L} & e^{K+L} \end{bmatrix}, \quad (19)$$

where  $T_{00}$  corresponds to  $n_i = n_{i+1} = 0$ ,  $T_{01}$  corresponds to  $n_i = 0$  and  $n_{i+1} = 1$ ,  $T_{10}$  corresponds to  $n_i = 1$  and  $n_{i+1} = 0$  and  $T_{11}$  corresponds to  $n_i = n_{i+1} = 1$ . Since the transfer matrix is independent of

the index  $i$ , one can simplify the partition sum further by defining

$$|n_i = 0\rangle = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, \quad (20)$$

$$|n_i = 1\rangle = \begin{bmatrix} 0 \\ 1 \end{bmatrix}. \quad (21)$$

With this definition it follows that  $\sum_{n_i=0,1} |n_i\rangle\langle n_i| = 1$  and

$$Z_N = \sum_{n_1=0,1} \dots \sum_{n_N=0,1} T_{1,2} \cdot T_{2,3} \cdot \dots \cdot T_{N,1} \quad (22)$$

$$= \sum_{n_1=0,1} \dots \sum_{n_N=0,1} \langle n_1|T|n_2\rangle \langle n_2|T|n_3\rangle \dots \langle n_N|T|n_1\rangle \quad (23)$$

$$= \sum_{n_1=0,1} \langle n_1|T^N|n_1\rangle \quad (24)$$

$$= \lambda_1^N(K, L) + \lambda_2^N(K, L), \quad (25)$$

where  $\lambda_1$  and  $\lambda_2$  are the eigenvalues of  $T$  which are given by

$$\lambda_{1/2}(K, L) = \frac{1 + e^{K+L} \pm \sqrt{(1 + e^{K+L})^2 - 4e^L(e^K - 1)}}{2}. \quad (26)$$

## 2. Find a transformation of the occupancies $n_i$ to map the lattice gas model to the Ising model with spins $s_i$ . (2 points)

A map from the lattice gas model,

$$Z_L(K, L) = \sum_{\{n_i\}} e^{K \sum_i n_i n_{i+1} + L \sum_i n_i}, \quad (27)$$

with occupancies  $n_i = 0, 1$  to the Ising model,

$$Z_I(K', L') = \sum_{\{s_i\}} e^{K' \sum_i s_i s_{i+1} + L' \sum_i s_i}, \quad (28)$$

with spins  $s_i = \pm 1$  can be given by

$$n_i = \frac{s_i + 1}{2}. \quad (29)$$

Inserting Eq. (29) into Eq. (27) yields

$$Z_L(K, L) = e^{\frac{N}{2}(\frac{K}{2}+L)} \cdot Z_I(K'(K, L), L'(K, L)), \quad (30)$$

with  $K'(K, L) = K/4$  and  $L'(K, L) = (K + L)/2$ . Thus, the lattice gas model and the Ising model are in fact equivalent. Expressing the eigenvalues of the lattice model  $\lambda_{1/2}(K, L)$  with the new variables  $K'$  and  $L'$ ,

$$\lambda_{1/2}(K', L') = e^{L'} \left( \cosh(L') \pm \sqrt{\cosh^2(L') - 2e^{-2K' \sinh(2K')}} \right), \quad (31)$$

and calculating the partition sum of the Ising model,

$$\begin{aligned} Z_I(K', L') &= e^{-N(L'-K')} Z_L(K', L') = e^{-N(L'-K')} (\lambda_1^N(K', L') + \lambda_2^N(K', L')) \\ &= \tilde{\lambda}_1^N(K', L') + \tilde{\lambda}_2^N(K', L'), \end{aligned} \quad (32)$$

gives in fact the correct eigenvalues of the Ising model:

$$\tilde{\lambda}_{1/2}(K', L') = e^{K'} \left( \cosh(L') \pm \sqrt{\cosh^2(L') - 2e^{-2K' \sinh(2K')}} \right). \quad (33)$$

**3. Derive an expression for the average  $\langle n_i \rangle$  in the limit of  $N \rightarrow \infty$  in terms of the eigenvalues of the transfer matrix. (2 points)**

Since  $\lambda_1 > \lambda_2$  the partition sum can be simplified for  $N \rightarrow \infty$ :

$$Z = \lambda_1^N + \lambda_2^N = \lambda_1^N \left[ 1 + \left( \frac{\lambda_2}{\lambda_1} \right)^N \right] \longrightarrow \lambda_1^N. \quad (34)$$

We see that only  $\lambda_1$  matters in the limit  $N \rightarrow \infty$ . Thus, we will neglect  $\lambda_2$  in the following. An expression for  $\langle n_i \rangle$  can be given as:

$$\begin{aligned} \langle n_i \rangle &= \frac{1}{Z} \sum_{n_1=0,1} \dots \sum_{n_N=0,1} n_i e^{-\beta H} \\ &= \frac{1}{Z} \sum_{n_1=0,1} \dots \sum_{n_N=0,1} n_i T_{1,2} \cdot T_{2,3} \cdot \dots \cdot T_{N,1} \\ &= \frac{1}{Z} \sum_{n_1=0,1} \dots \sum_{n_N=0,1} \langle n_1 | T | n_2 \rangle \dots n_i \langle n_i | T | n_{i+1} \rangle \dots \langle n_N | T | n_1 \rangle \\ &= \frac{1}{Z} \sum_{n_1=0,1} \langle n_1 | T^{i-1} | n_i = 1 \rangle \langle n_i = 1 | T^{N-(i-1)} | n_1 \rangle \\ &\approx \frac{1}{Z} \lambda_1^N \langle n_1 = 0 | n_i = 1 \rangle \langle n_i = 1 | n_1 = 0 \rangle \\ &= \langle n_1 = 0 | n_i = 1 \rangle \langle n_i = 1 | n_1 = 0 \rangle. \end{aligned} \quad (35)$$

Expressing the states in the eigenvector basis of  $T$  and calculating the scalar products yields the wanted expression for  $\langle n_i \rangle$  in the limit of  $N \rightarrow \infty$ .

### 3 Renormalization of the Ising chain (3 points)

In the lecture we have derived the following RG flow equation for the coupling constant  $K$  of the Ising chain without magnetic field: the new value  $K'$  is given by

$$K'(K) = \frac{1}{2} \ln \cosh(2K). \quad (36)$$

In addition we have derived the absolute increase in free energy per spin arising in each iteration:

$$g(K) = \frac{1}{2} \ln 2 + \frac{1}{4} \ln \cosh(2K) \quad (37)$$

**1. Write a short computer program (e.g. in Mathematica or Python) that defines the flow equation  $K'(K)$  and the free energy increase  $g(K)$  as functions. Start with a coupling constant  $K_0 = 1$  and iterate through  $K_1, K_2, K_3$  up to  $K_4$ . Also calculate the corresponding values  $g_0 = g(K_0)$  to  $g_4 = g(K_4)$ . What are the limits for these two series? (1.5 points)**

Python functions:

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1 def K_new(K):
2     return log(cosh(2*K)) / 2
3
4
5 def g(K):
6     return log(2) / 2 + log(cosh(2*K)) / 4

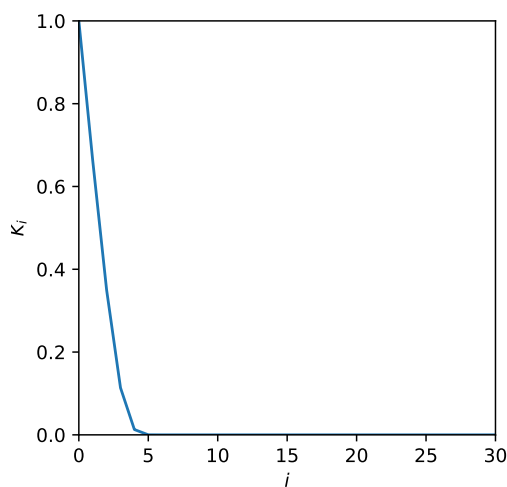
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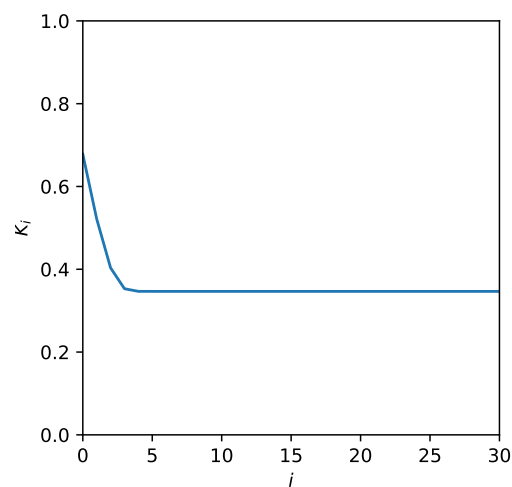
Iteration to  $i = 4$  yields

	i=0	i=1	i=2	i=3	i=4
$K_i$	1	0.66	0.35	0.11	0.01
$g_i$	0.68	0.52	0.4	0.35	0.35

The  $K_i$ -series converges to zero for large  $i$ , while the  $g_i$ -series converges to  $\frac{1}{2} \ln 2 \approx 0.35$ .



(a) coupling constant  $K$



(b) free energy increase  $g$

**2. Use these results to estimate the dimensionless free energy per spin  $f = -\beta F/N$  in fourth order (simply cut the appropriate sum after the term with  $g_4$ ; you can also include the next order term, but now by simply using the first term in  $g(K)$ ). Compare to the known exact result for the Ising chain. How good is the numerical agreement? (1.5 points)**

$$f = \sum_{i=0}^4 g_i \approx 2.3 \tag{38}$$