

$$\beta = \frac{1}{kT} \quad T = \frac{1}{k_B\beta}$$

$$①.) \quad F = U - TS$$

$$U = \sum E_i P(E_i) = \sum E_i \frac{e^{-\beta E_i}}{Z} = -\partial_\beta \ln Z$$

$$S = -\sum P(E_i) \ln P(E_i) = -\sum \frac{e^{-\beta E_i}}{Z} \ln \frac{e^{-\beta E_i}}{Z}$$

$$d\ln Z = \frac{\partial \ln Z}{\partial \beta} d\beta + \frac{\partial \ln Z}{\partial \mu} dm = -U d\beta + \beta \mu dm$$

$$m = \sum P(E_i)m_i = \frac{1}{Z} \frac{\partial \ln Z}{\partial \mu}$$


$$d(\mu V) = d\mu V + \mu dm \rightarrow dm = \ell(\mu V) - \mu d\beta$$

$$d\ln Z = -d(\mu V) + d\mu V + \beta \mu dm$$

$$\Rightarrow dV = \frac{1}{\mu} d\ln Z + \frac{1}{\mu} d(\mu V) - \beta dm$$

$$= \frac{1}{\mu} d(\ln Z + \mu V) - \beta dm$$

$$\text{but in } N \rightarrow \infty \quad dU = TdS + \beta dm$$

$$\Rightarrow S = \omega_B \ln Z + \frac{U}{T} \rightarrow F = U - TS$$

$$\Rightarrow F = \cancel{\lambda_+} - k_B T \ln Z + \cancel{\lambda_-}$$

```

| ex1q1.py
import numpy as np
import matplotlib.pyplot as plt
from matplotlib import cm
from matplotlib import colors
plt.rcParams["font.size"] = 15

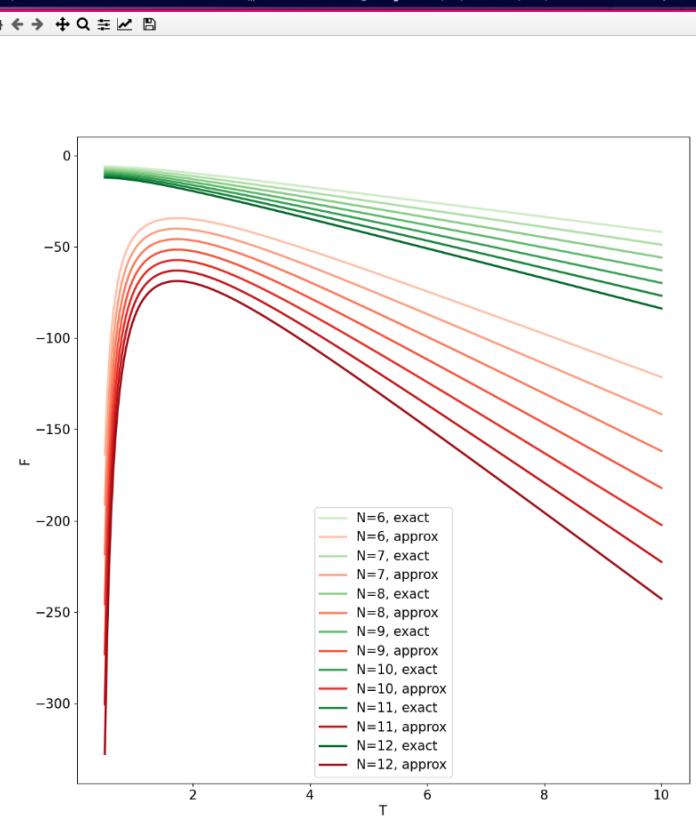
T = np.linspace(5e-01, 10, 1000)
norm = colors.Normalize(vmin=0, vmax=9)
cmap1 = cm.get_cmap('Greens', 10)
cmap2 = cm.get_cmap('Reds', 10)

for N in range(6, 13):
    Fexact = -T * N * np.log(2 * np.cosh(1 / T))
    L = np.exp(1 / T) * np.cosh(1 / T) + np.sqrt(
        np.exp(2 / T) * (np.cosh(1 / T)) ** 2 - 2 * np.sinh(2 / T)
    )
    Fapprox = -T * N * L

    plt.plot(T, Fexact, lw=2.5, label=f"N={N}, exact", color=cmap1(norm(N-4)))
    plt.plot(T, Fapprox, lw=2.5, label=f"N={N}, approx", color=cmap2(norm(N-4)))

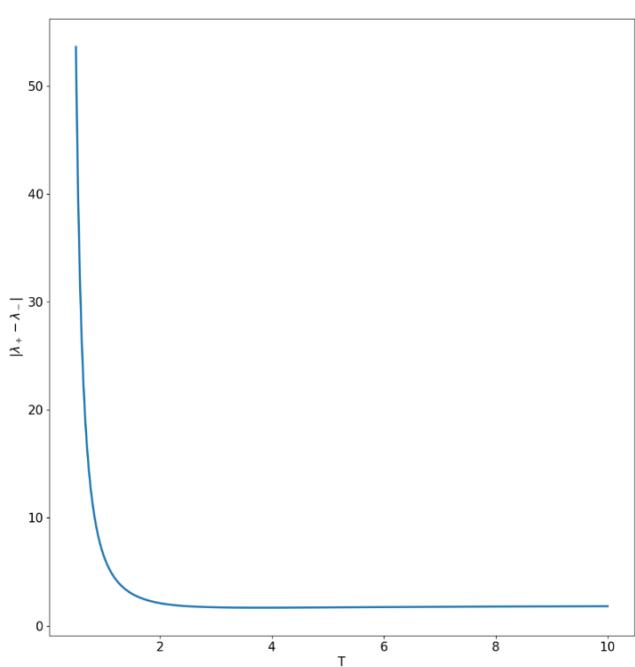
plt.ylabel("F")
plt.xlabel("T")
plt.legend(loc="best")
plt.show()

```



The approximation fails at $T=0$.

AT $T=0$ both λ_{\pm} should have the same value cuz there's only one state for the system.



But clearly this is not the case. AT $T=0$

$$\lambda_+ - \lambda_- \xrightarrow{T \rightarrow 0} \infty$$

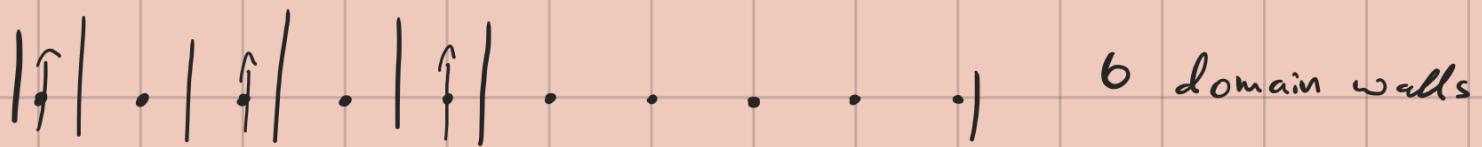
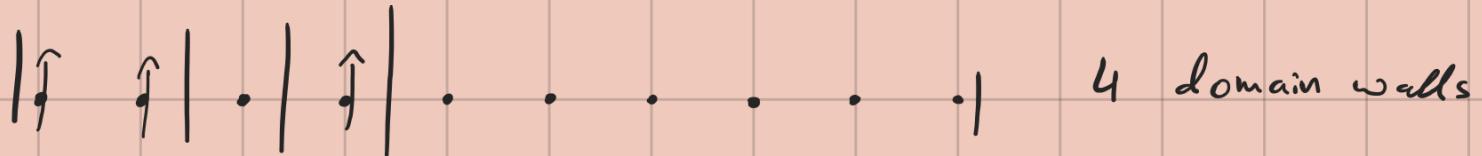
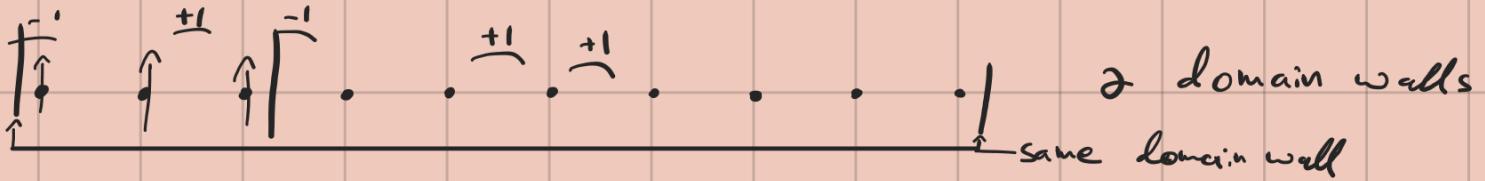
②

$$Z = \sum_{\text{m-states}} e^{-\beta H}$$

$$H = -J \sum_i C_i C_{i+1}$$

(periodic boundary condition $C_{N+1} \equiv C_1$)

It's easier to count domain walls than options.



number of options: $\binom{\text{number of domain walls}}{\text{number of positions}} = \binom{2}{10}, \binom{4}{10}, \binom{6}{10}$

$T_{\text{tot}} = 10J$ each domain wall = $-J$

$Z = (u_5 e^{8\beta J} + 210 e^{6\beta J} + 210^6 e^{4\beta J}) \cdot J^6$ for spin inversion

(3) 1)

$$1 - S^2 = \begin{cases} 0 & S = -1 \\ 1 & S = 0 \\ 0 & S = 1 \end{cases}$$

if the spin is in
an excited state
what so ever

$$2) e^{\beta J S_1 S_2 + \frac{\beta B}{2} (S_z + S_{\sigma}) + \frac{\beta \lambda}{2} (S_x - S_{\sigma} - S_z)}$$

symmetric

$$\begin{matrix} S_1 = -1 & S_2 = -1 \\ = -1 & = 0 \\ = -1 & = 1 \end{matrix} \quad \begin{matrix} \vdots \\ \vdots \\ \vdots \end{matrix} \quad \begin{matrix} e^{\beta J + 2\beta B - \beta \lambda} \\ e^{-\beta B} \\ e^{-\beta J - \beta \lambda} \end{matrix}$$

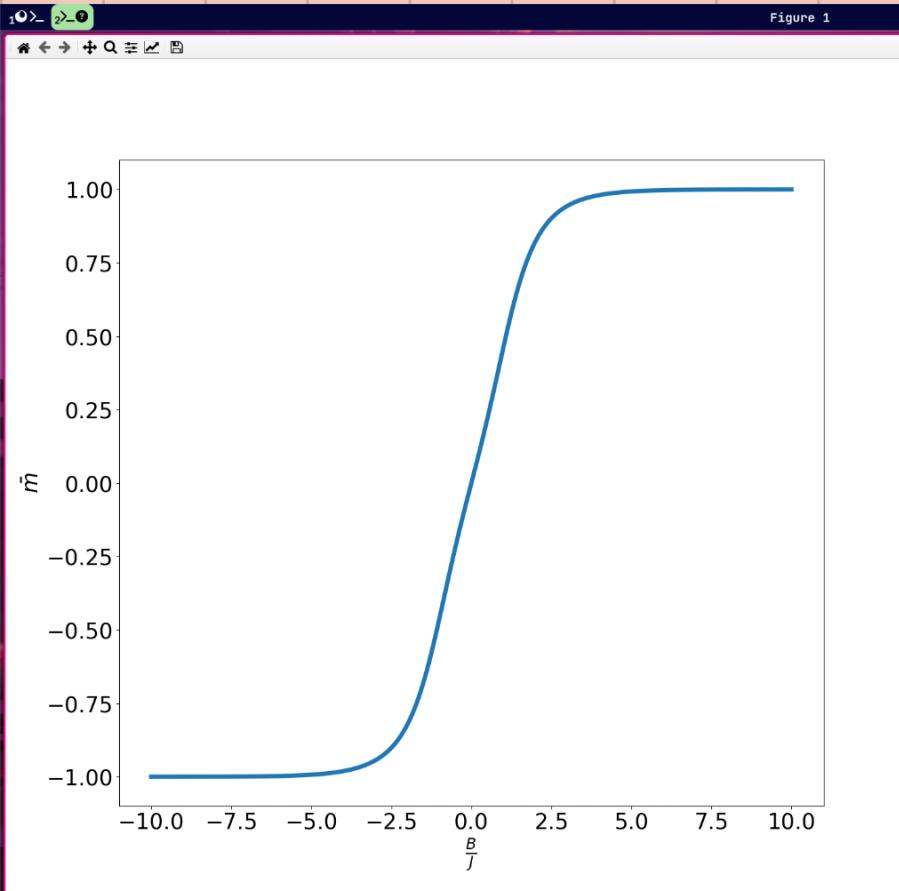
$$\begin{matrix} S_1 = -1 \\ = -1 \\ = -1 \end{matrix} \quad \begin{matrix} S_2 = -1 \\ = 0 \\ = 1 \end{matrix} \quad \begin{matrix} S_1 = 0 \\ = 0 \\ = 1 \end{matrix} \quad \begin{matrix} S_2 = 0 \\ = 1 \\ = 1 \end{matrix}$$

$$\left(\begin{matrix} e^{-\frac{\beta}{J}} & e^{-\frac{1}{2}\beta B + \frac{1}{2}\beta \lambda} & e^{-\frac{1}{2}\beta J} \\ e^{\beta J - \beta B} & e^{-\frac{1}{2}\beta B + \frac{1}{2}\beta \lambda} & e^{\frac{1}{2}\beta J + \frac{1}{2}\beta \lambda} \\ e^{-\beta J} & e^{\frac{1}{2}\beta B + \frac{1}{2}\beta \lambda} & e^{\beta J + \beta B} \end{matrix} \right)$$

$$\lambda = \alpha T \quad \beta = \frac{1}{T}$$

3) The eigen values + vectors are monstrous.
 Anyway $\bar{m} = \lambda_{\text{largest}}^N$ from wolfram

$$\therefore \bar{m} = \frac{1}{N} \frac{\partial F}{\partial B} = \frac{1}{N} \frac{\partial -\beta m z}{\partial B} = -\beta \frac{\partial \ln \lambda}{\partial B}$$



```

ex1.py
import numpy as np
import matplotlib.pyplot as plt
plt.rcParams["font.size"] = 25

l = 250      # E741 ambiguous variable name 'l'
B = np.linspace(-10, 10, l)
array = np.zeros(l)

for i, b in np.ndenumerate(B):
    TM = np.array([
        [np.e ** (1 - b), np.e ** (1 - 0.5 * b), np.e ** (-1)],
        [np.e ** (1 - 0.5 * b), np.e ** (2), np.e ** (1 + 0.5 * b)],
        [np.e ** (-1), np.e ** (1 + 0.5 * b), np.e ** (1 + b)],
    ])
    eigv, eigvec = np.linalg.eig(TM)
    array[i] = np.log(eigv.max())

grad = np.gradient(array, B)
plt.plot(B, grad, lw=5)
plt.ylabel(r"\bar{m}")
plt.xlabel(r"\frac{B}{J}")
plt.show()

```

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(u) $NN + NNN$

We don't need: $\langle G_1 G_2 G_3 | M | G_1 G_2 G_3 \rangle \langle G_1 G_2 G_3 | M | G_2 G_3 G_1 \rangle \dots$

This is 8 basis vectors, one for each combination

We can make due with 4 basis vectors:

$\langle G_1 G_2 | M | G_1 G_2 \times G_1 G_3 | M | G_2 G_3 \times G_2 G_3 | M | G_3 G_1 \rangle \dots$

but this is still more free variables than we

need. 4×4 matrix is 16 elements, we have 8 combinations

which is still less than 16 free elements for a symmetric

4×4 matrix \Rightarrow this will not be a symmetric matrix!

$$H = \frac{J_1}{2} (G_1 G_2 + G_2 G_3) + J_2 G_1 G_3 + \frac{\Omega}{3} (G_1 + G_2 + G_3)$$

The Hamiltonian knows about 3 spins at a time: $\langle G_1 G_2 | M | G_2 G_3 \rangle$

> python ex1q4.py

$$\begin{bmatrix} -\beta \cdot (-B - 2 \cdot J_1 - J_2) & -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} & -\beta \cdot (-B - J_2) & -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} - 2 \cdot J_1 + J_2 \\ e & e & e & e \\ -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} & -\beta \cdot \begin{pmatrix} B \\ - \\ + \\ 3 \end{pmatrix} + 2 \cdot J_1 + J_2 & -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} + 2 \cdot J_1 - J_2 & -\beta \cdot \begin{pmatrix} B \\ - \\ + \\ 3 \end{pmatrix} + J_2 \\ e & e & e & e \\ -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} + J_2 & -\beta \cdot \begin{pmatrix} B \\ - \\ + \\ 3 \end{pmatrix} + 2 \cdot J_1 - J_2 & -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} + 2 \cdot J_1 + J_2 & -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} + J_2 \\ e & e & e & e \\ -\beta \cdot \begin{pmatrix} B \\ - \\ - \\ 3 \end{pmatrix} - 2 \cdot J_1 + J_2 & -\beta \cdot (B - J_2) & -\beta \cdot \begin{pmatrix} B \\ - \\ + \\ 3 \end{pmatrix} & -\beta \cdot (B - 2 \cdot J_1 - J_2) \\ e & e & e & e \end{bmatrix}$$

We need to diagonalize it and then $Z = \tilde{\lambda}_{\max}$

```
| + ex1q4.py    x
import itertools
from sympy import symbols, exp, Matrix, pprint

# Define symbolic parameters
beta, J1, J2, B = symbols("beta J1 J2 B")
spin_values = [-1, 1] # Possible spin values

# Generate all configurations of three spins
configurations = list(itertools.product(spin_values, repeat=2))

# Initialize the transfer matrix
n_configs = len(configurations)
T = Matrix.zeros(n_configs, n_configs)

# Compute each element of the transfer matrix
for i, C in enumerate(configurations):
    for j, C_prime in enumerate(configurations):
        # Unpack spins for C and C_prime
        s_i, s_ip1 = C
        s_ip1p, s_ip2 = C_prime

        # Hamiltonian between these configurations
        H = (
            - J1 * (s_i * s_ip1 + s_ip1p * s_ip2)
            - J2 * (s_i * s_ip2)
            + (B / 3) * (s_i + s_ip1 + s_ip2)
        )

        # Transfer matrix element
        T[i, j] = exp(-beta * H)

# Print the symbolic transfer matrix
pprint(T)
```

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