

16.5/20

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Answer 1

For any two sites i, j the interaction is termed as J .

The role of J is as follows:

$J > 0 \rightarrow$ Interaction is ferromagnetic

$J < 0 \rightarrow$ Interaction is anti-ferromagnetic

$J = 0 \rightarrow$ Non-interacting

✓ $2\frac{1}{2}/3$

good job !

Answer 2

When we consider periodic boundary conditions, the last spins are considered to be their nearest neighbour of each other.

✓ $1/2/1$

(1-D chain becomes a ring
How did you apply it in code ?)

Answer 3

Defining the necessary modules

```
import matplotlib.pyplot as plt
import numpy as np
from numpy.random import random
from random import choice
```

The 1-D Ising Model

```
def H(J, h, s): # Hamiltonian
    ret_value = 0.0
    n = len(s)
    for x in range(n):
        # We start by iterating over all spin states
        ret_value += -J*(s[(x-1)%n]*s[x]) - h*s[x] # for periodic
    boundary condition
    return ret_value
```

```
def S_random(n):
    # Defining array of length n filled with +1 or -1 random spin
    # states and appending it to S
    s = []
    for x in range(N):
        s.append(choice((+1, -1)))
    return s
```

it should be n here so that it doesn't need a global variable to run.

We have considered $J = 1$ and $k_B = 1$ as suggested in the question
 $J=1$

After differentiating the expression for $\langle m \rangle$ analytically, we get

the following expression:

```
def m_exact(n,h,T):  
    # ret_m = (np.tanh(h/T))/n  
    # return float(ret_m)  
    sin = np.sinh(h)  
    cos = np.cosh(h)  
    exp = np.exp(-4*J)  
    ratio = (cos-np.sqrt(sin**2+exp))/(cos+np.sqrt(sin**2+exp))  
    ret_value = (1-ratio**N)/(1+ratio**N)  
    ret_value *= sin/np.sqrt(sin**2+exp)  
    return ret_value
```

$$\frac{9}{10}$$

(points deducted only for error analysis
exact fn and calculating standard deviation would
have been better)

here its supposed to
be h/T and J/T

There isn't any temperature
dependence in your exact function
thats why your plots don't
agree.

```
def m_aprox(n,h,T,I):  
    num = 0.0  
    den = 0.0  
    for i in range(I):  
        S = S_random(N) # create random  
        num += (sum(S)/N)*np.exp(-H(J,h,S)/T) # use the observable  
        den += np.exp(-H(J,h,S)/T)  
    return num/den
```

The relevant dimensionless ratios are as follows:

(i) J/T and,

(ii) h/T

$m > i$ vs h for fixed N

```
plt.figure(figsize=(32,18))  
plt.rc('font', size=50)
```

```
N = 10 # Number of lattice sites <= ~20  
I = int(2**(N+1)) # Number of generated spin configurations
```

The range of h is given as [-1,1]

```
hRange = np.arange(-1,1,0.001)
```

```
exactVal=[m_exact(N,h,0.5)for h in hRange]  
aproxVal=[m_aprox(N,h,0.5,I)for h in hRange]
```

Use
these
in m-exact

$$1/1$$

```

aproxError=[val*I**-0.5 for val in aproxVal]

plt.plot(hRange ,exactVal ,'r',color="green")
# We have chosen the value of T as 0.5
plt.legend(("T = 0.5"))

plt.errorbar(hRange, aproxVal, yerr=aproxError, fmt='ro',
markersize=2)

plt.ylabel('<m>')
plt.xlabel('h')
plt.title("Graph for <m> vs h for fixed N")
plt.show()

```

```

/tmp/ipykernel_7392/3058133422.py:16: UserWarning: color is
redundantly defined by the 'color' keyword argument and the fmt string
"r" (-> color=(1.0, 0.0, 0.0, 1)). The keyword argument will take
precedence.

```

```

    plt.plot(hRange ,exactVal ,'r',color="green")

```

```

-----
-----
ValueError                                Traceback (most recent call
last)
Cell In [6], line 20
    17 # We have chosen the value of T as 0.5
    18 plt.legend(("T = 0.5"))
--> 20 plt.errorbar(hRange, aproxVal, yerr=aproxError, fmt='ro',
markersize=2)
    22 plt.ylabel('<m>')
    23 plt.xlabel('h')

```

```

File ~/.local/lib/python3.8/site-packages/matplotlib/pyplot.py:2482,
in errorbar(x, y, yerr, xerr, fmt, ecolor, elinewidth, capsize,
barsabove, lolims, uplims, xlolims, xuplims, errorevery, capthick,
data, **kwargs)

```

```

    2476 @_copy_docstring_and_deprecators(Axes.errorbar)
    2477 def errorbar(
    2478     x, y, yerr=None, xerr=None, fmt='', ecolor=None,
    2479     elinewidth=None, capsize=None, barsabove=False,
lolims=False,
    2480     uplims=False, xlolims=False, xuplims=False,
errorevery=1,
    2481     capthick=None, *, data=None, **kwargs):
-> 2482     return gca().errorbar(
    2483         x, y, yerr=yerr, xerr=xerr, fmt=fmt, ecolor=ecolor,
    2484         elinewidth=elinewidth, capsize=capsize,
barsabove=barsabove,

```

```

2485         lolims=lolims, uplims=uplims, xlolims=xlolims,
2486         xuplims=xuplims, errorevery=errorevery,
capthick=capthick,
2487         **({"data": data} if data is not None else {}),
**kwargs)

```

File ~/.local/lib/python3.8/site-packages/matplotlib/__init__.py:1423, in _preprocess_data.<locals>.inner(ax, data, *args, **kwargs)

```

1420 @functools.wraps(func)
1421 def inner(ax, *args, data=None, **kwargs):
1422     if data is None:
-> 1423         return func(ax, *map(sanitize_sequence, args),
**kwargs)
1425     bound = new_sig.bind(ax, *args, **kwargs)
1426     auto_label = (bound.arguments.get(label_namer)
1427                  or bound.kwargs.get(label_namer))

```

File

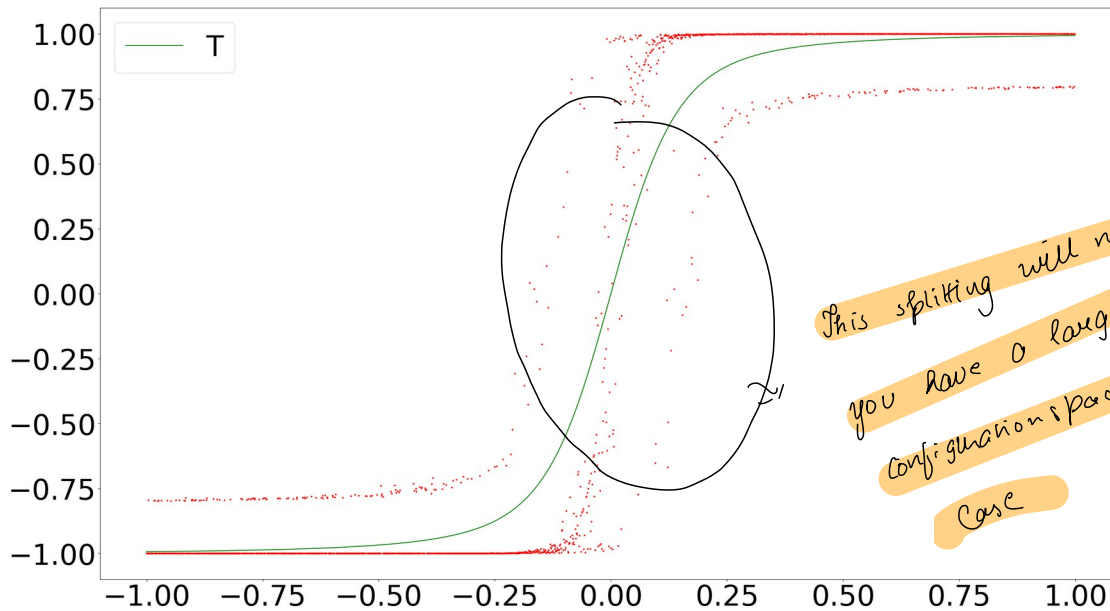
~/.local/lib/python3.8/site-packages/matplotlib/axes/_axes.py:3587, in Axes.errorbar(self, x, y, yerr, xerr, fmt, ecolor, elinewidth, capsize, barsabove, lolims, uplims, xlolims, xuplims, errorevery, capthick, **kwargs)

```

3584 res = np.zeros(err.shape, dtype=bool) # Default in case of
nan
3585 if np.any(np.less(err, -err, out=res, where=(err == err))):
3586     # like err<0, but also works for timedelta and nan.
-> 3587     raise ValueError(
3588         f"'{dep_axis}err' must not contain negative values")
3589 # This is like
3590 #     elow, ehigh = np.broadcast_to(...)
3591 #     return dep - elow * ~lolims, dep + ehigh * ~uplims
3592 # except that broadcast_to would strip units.
3593 low, high = dep + np.row_stack([-(1 - lolims), 1 - uplims]) *
err

```

ValueError: 'yerr' must not contain negative values



$\langle m \rangle$ vs N for fixed h

```
plt.figure(figsize=(32,18))
plt.rc('font', size=50)
```

```
h = 0.25 # Fixed external magnetic field
I = int(2**(N+1)) # Number of generated spin configurations
```

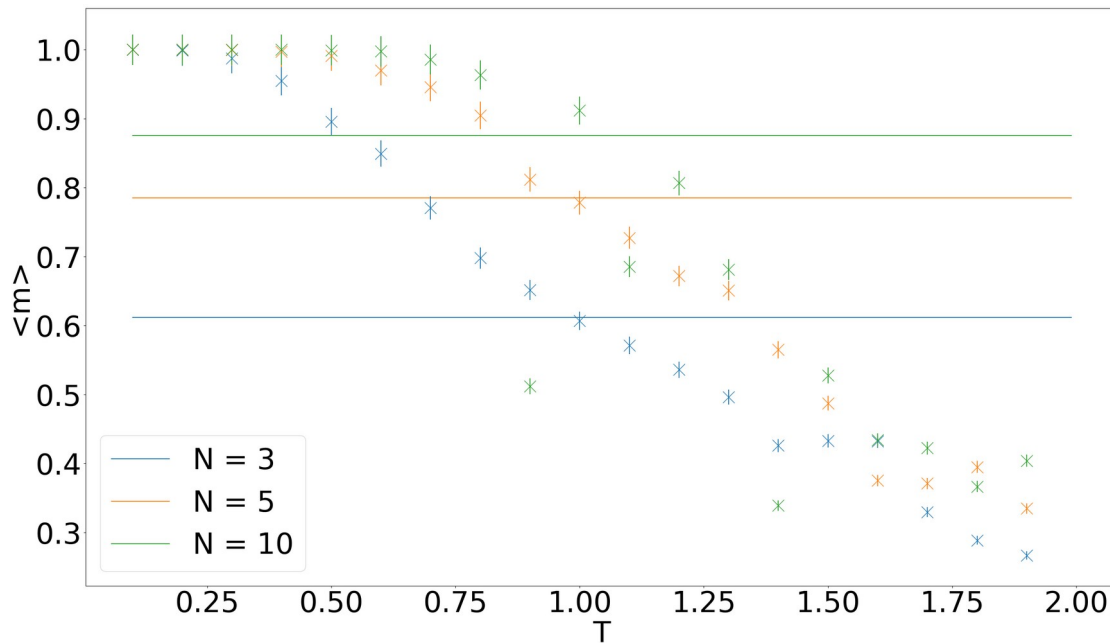
```
NRange = [3,5,10] # Number of lattice sites
```

```
TExactRange = np.arange(0.1, 2, 0.01) #For a range of Temperature
TAproxRange = np.arange(0.1, 2, 0.1)
```

```
for N in NRange:
    exactVal=[m_exact(N,h,T) for T in TExactRange]
    plt.plot(TExactRange, exactVal)
plt.legend(("N = 3", "N = 5", "N = 10"))
```

```
plt.gca().set_prop_cycle(None) #reset colour cycle
```

```
for N in NRange:
    aproxVal=[m_aprox(N,h,T,I) for T in TAproxRange]
    aproxError=[val*I**-0.5 for val in aproxVal]
    plt.errorbar(TAproxRange, aproxVal, yerr=aproxError, fmt = 'x',
markersize=20)
plt.ylabel('<math>\langle m \rangle</math>')
plt.xlabel('<math>T</math>')
plt.show()
```



(Not done comparison with thermodynamic limit plot for m -approx is wrong because doesn't have T dep)

$$\frac{3.5P}{5}$$

for thermodynamic limit: take the limit of the theoretical expression in case $N \rightarrow \infty$ you should get

$$\langle m \rangle = - \frac{\sinh(h)}{\sqrt{\sinh^2(h) + e^{-4J}}}$$

where
 $h = \frac{h}{T}$
 $J = \frac{J}{T}$

