

# Computational Many-Body Physics

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SS 2020

**Sheet 3** - please submit your solutions via e-mail to Chae-Yeun Park until Monday, June 8, 2020, 12:00.

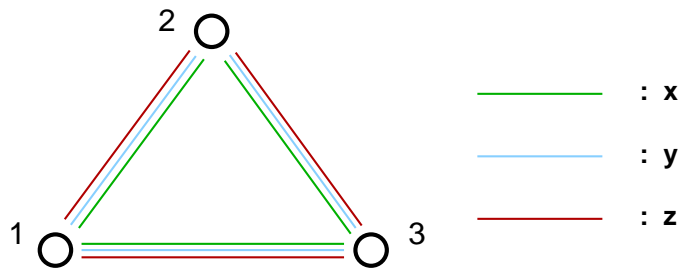
## Exercise 1: Spin-models on a three-site cluster

(7 points)

Consider the following (general) Hamiltonian for a spin-model on a three-site cluster:

$$H = - \sum_{ij\alpha} J_{ij}^\alpha S_i^\alpha S_j^\alpha ,$$

with  $i, j = 1, 2, 3$  ( $i < j$  in  $\sum_{ij}$ ) and  $\alpha = x, y, z$ .



To visualize the model, a colour code for the  $x, y$  and  $z$  components of the spin-couplings turns out to be useful, see the figure.

- a) Rewrite the Hamiltonian using the operators

$$S_i^\pm = S_i^x \pm iS_i^y , \text{ and } S_i^z .$$

(1 point)

Now set up (by hand!) the  $8 \times 8$  Hamilton matrices  $\bar{H}$  for the following three special cases:

- b) the Ising model, i.e.  $J_{ij}^\alpha = J\delta_{\alpha z}$ , (2 points)  
c) the isotropic Heisenberg model, i.e.  $J_{ij}^\alpha = J$ , and (2 points)  
d) a model with  $J_{12}^x = J_{23}^y = J_{31}^z = J$  and all other  $J_{ij}^\alpha = 0$ . (2 points)

## Exercise 2: Hamilton matrices for spin models

(13 points)

Consider a spin model of the form

$$H = - \sum_{ij\alpha} J_{ij}^{\alpha} S_i^{\alpha} S_j^{\alpha} ,$$

with  $i, j = 1, \dots, N$  and  $\alpha = x, y, z$ .

- a) Write a program which sets up the Hamilton matrix for a model with arbitrary  $N$  and an arbitrary list of couplings  $\{J_{ij}^{\alpha}\}$ . As discussed in the lecture, the  $x$ -,  $y$ - and  $z$ -links should be treated separately. For each link of the form  $-J_{ij}^{\alpha} S_i^{\alpha} S_j^{\alpha}$ , the action of  $S_i^{\alpha} S_j^{\alpha}$  on the basis state  $|n\rangle$  gives another basis state  $|l\rangle$  (times a prefactor). With the equation for  $l$  given in the lecture, the Hamiltonian matrix can be set up very efficiently. (10 points)
- b) Set up the Hamilton matrices for the three models discussed in exercise 1 (the three-site clusters) and calculate the eigenenergies and corresponding eigenstates for each model. (3 points)