

# LAB XI

## Simulating Adiabatic QC

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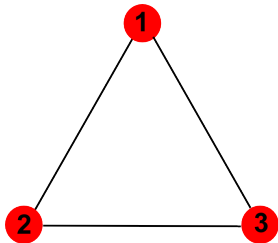
# Task

Prepare the parameter dependent Hamiltonian matrix

$$H(\lambda) = (1 - \lambda)H_0 + \lambda H_1,$$

where

$$H_0 = -\sum_i S_i^x \text{ and } H_1 = -\sum_{ij} J_{ij} S_i^z S_j^z - \sum_i h_i S_i^z.$$



Take the above system with  $h_1 = 0.6$ ,  $h_2 = h_3 = 0$ ,  $J_{12} = -1.1$ ,  $J_{13} = -2.1$ ,  $J_{23} = -3.8$  (all couplings are antiferromagnetic).

Note that  $S_i^z = S^z \otimes \mathbf{1}_{2 \times 2} \otimes \mathbf{1}_{2 \times 2}$  etc. The tensor product  $\otimes$  is just the Kronecker product of matrices, available in `Numpy`.

## Task (continued)

Calculate and plot  $\Delta E(\lambda) = E_1(\lambda) - E_0(\lambda)$ , which is the difference between the first excited and the ground state energy.

Calculate the optimal running time of the adiabatic evolution  $T_{\text{AQC}}$

$$T_{\text{AQC}} = \int_0^1 \frac{d\lambda}{[\Delta E(\lambda)]^2}.$$

It is enough to approximate the integral by a discrete sum of small intervals.

Calculate and plot  $\langle S_i^z \rangle$  for  $i = 1, 2, 3$  as a function of  $\lambda$ , label the curves. Here  $\langle S_i^z \rangle = \langle \psi | S_i^z | \psi \rangle$  is the expectation value of the operator  $S_i^z$  on the ground state eigenvector  $|\psi(\lambda)\rangle$ .

# Hints

We need the following  $2 \times 2$  spin matrices:

$$S^z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}, \quad S^x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}.$$

Spin operator  $S_i^z$  is a Kronecker product  $S_i^z = \mathbf{1} \otimes \dots \otimes S^z \otimes \dots \otimes \mathbf{1}$  with  $2 \times 2$  matrix  $S^z$  at position  $i$ . Implement  $\otimes$  directly from `numpy`:

```
from numpy import kron

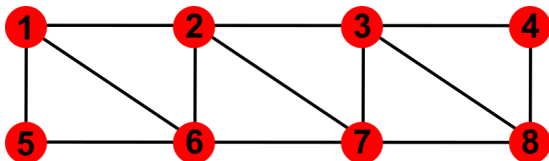
sz = np.array( [[1,0.],[0.,-1]] )
one = np.eye(2)

sz_1 = kron( sz, kron(one, one) )
```

# Extra Task

The purpose of this exercise is to write a more abstract code, which can store connectivity of the graph in a proper data structure. Introduce and use some data structure to encode the system graph. You can use e.g. a Python dictionary or tools from the package `networkx`.

Prepare the Hamiltonians  $H_0$  and  $H_1$  for the system:



Calculate and plot  $\Delta E(\lambda)$  for the test case  $J_{ij} = -1$  for all connections, and  $h_i = 0.9$  for all magnetic fields.

Calculate  $\Delta E(\lambda)$  for a few random realizations of  $J_{ij}$  and  $h_i$  and plot on a single picture. Draw the random values uniformly from the interval  $[-1, 1]$ .

# Hints

Graph in Python can be represented e.g. by dictionary of lists

```
Graph = { 'Q1': ['Q2', 'Q4'], 'Q2': ['Q3'], ... }  
Graph['Q1']
```

where a node key points to the nodes connected by edges.

One may also store the properties of a node in a dictionary

```
Node = { 'Q1': (sz1, sx1, h1), 'Q2': ... }.
```

## Extra – open

There are many more things one can compute:

- entanglement entropy (as discussed in the **reading material**)
- qubism plots e.g. from <https://arxiv.org/abs/1112.3560>
- scaling the worst case gap with  $N$  for a given lattice  
(use `scipy.sparse` for sparse matrices).

If you are interested – one of the above could be your topic for the final presentation!