

# Phys730 Mid-Term-Project Proposal

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February 26, 2016

## Numerical Solution of Quantum Heisenberg Chain

### 1 Introduction

The  $S = \frac{1}{2}$  isotropic quantum Heisenberg chain is a model of mutually interacting quantum spins arranged with regular spacing along a line. The interaction is limited to nearest-neighbour spins and the nature of interaction is antiferromagnetic ( $J > 0$ ), meaning that they favor the opposite alignment of spins.

$$\hat{H} = J \sum_i^L \left[ \frac{1}{2} (\hat{S}_i^+ \hat{S}_{i+1}^- + \hat{S}_i^- \hat{S}_{i+1}^+) + \hat{S}_i^z \hat{S}_{i+1}^z \right]$$

Where  $L$  is the length of the chain.

In 1931, the model was solved analytically by Bethe and the exact ground state energy per site is  $E_0 = \frac{1}{4} - \log_e(2)$ .

### 2 Computational Method

Python code will be written to generate a Hamiltonian matrix and to calculate its eigenvalues. The lowest eigenvalue is the ground state energy. The Calculation will be carried out on finite-size system of increasing size, and the result will be extrapolated to the thermodynamic limit. Numerical result will be compared to the known result.

### 3 Tool

numpy  
scipy  
matplotlib