

## Lecture of Tensor Products and Heisenberg Spin Chain

### Multi-Qubit Systems:

If two independent systems are represented by the states  $|\psi_1\rangle, |\psi_2\rangle \in C^2$ , then the state of the system made of these two qubits would be  $|\psi_1, \psi_2\rangle = |\psi_1\rangle|\psi_2\rangle = |\psi_1\rangle \otimes |\psi_2\rangle$

The product state of the states  $|\psi_1\rangle, |\psi_2\rangle \in C^2$  is the pure tensor  $= |\psi_1\rangle \otimes |\psi_2\rangle$  it represents the state of a system of two independent qubits in the states  $|\psi_1\rangle$  and  $|\psi_2\rangle$

From our prior definition of tensors, we know that pure tensors are only a generating set of  $V \otimes W$ , but certain elements of  $V \otimes W$  are not pure tensors.

When the state of a 2-qubit system is not in a product state, we call that system to be in an entangled state. This definition will become clearer with the example below.

**Example:** We can start by looking the state

$$|\psi\rangle = \frac{1}{\sqrt{2}} (|0,0\rangle + |1,1\rangle) = \frac{1}{\sqrt{2}} (|0\rangle|0\rangle + |1\rangle|1\rangle) \in C^4$$

In this case, we can prove that  $|\psi\rangle$  is an entangled state. We will start by assuming that  $|\psi\rangle$  is in a product state, We let  $a_0, a_1, b_0, b_1 \in C$  s.t.

$$|\psi\rangle = |\psi_1\rangle \otimes |\psi_2\rangle$$

$$|\psi_1\rangle = a_0|0\rangle + a_1|1\rangle$$

$$|\psi_2\rangle = b_0|0\rangle + b_1|1\rangle$$

As a result, we will yield the following,

$$|\psi_1\rangle \otimes |\psi_2\rangle = a_0 b_0 |0\rangle|0\rangle + a_1 b_1 |1\rangle|1\rangle + a_0 b_1 |0\rangle|1\rangle + a_1 b_0 |1\rangle|0\rangle$$

This result will imply the below constraints:

$$a_0 b_0 = a_1 b_1 = \frac{1}{\sqrt{2}} \quad \text{and} \quad a_0 b_1 = a_1 b_0 = 0$$

The first condition contradicts the second condition. Hence, we can conclude that the state is not a product state, but rather, the state is entangled.

## Partial Measurements:

**Definition:** Given  $|\psi\rangle = a|0,0\rangle + b|0,1\rangle + c|1,0\rangle + d|1,1\rangle \in C^2 \otimes C^2$

*Measure of the 1st qubit:*

1 with probability  $a^2 + b^2$   $\rightarrow$  with posterior state  $= \frac{a|0,0\rangle + b|0,1\rangle}{\sqrt{a^2 + b^2}}$

0 with probability  $c^2 + d^2$   $\rightarrow$  with posterior state  $= \frac{c|1,0\rangle + d|1,1\rangle}{\sqrt{c^2 + d^2}}$

*Measure of the 2nd qubit:*

1 with probability  $a^2 + c^2$   $\rightarrow$  with posterior state  $= \frac{a|0,0\rangle + c|1,0\rangle}{\sqrt{a^2 + c^2}}$

0 with probability  $b^2 + d^2$   $\rightarrow$  with posterior state  $= \frac{b|0,1\rangle + d|1,1\rangle}{\sqrt{b^2 + d^2}}$

**Example :** We will use the same example state as before in order to practice partial measurements.

$$|\psi\rangle = \frac{1}{\sqrt{2}} (|0,0\rangle + |1,1\rangle) = \frac{1}{\sqrt{2}} (|0\rangle|0\rangle + |1\rangle|1\rangle) \in C^4$$

If we measure the first qubit, we obtain 0 with probability 1/2, and the system is left in the state  $|0\rangle|0\rangle$ .

Then, a measurement of the second qubit yields 0 with probability 1.

However, if we decided to measure the second qubit initially, we would get 0 with probability 1/2. As a result, the measurement of the first qubit of the system has impacted the subsequent measurements of the second qubit.

## Spins:

In physics, spin is a fundamental quantum concept, and for our purposes, we'll pragmatically consider it as a linear space representing the  $su(2)$  algebra. The discussion won't delve into the detailed physical origin of spin but emphasizes its representation through the specified  $su(2)$  algebra commutation relation.

$[S^\alpha, S^\beta] = i\epsilon^{\alpha\beta\gamma}S^\gamma$ ,  $\alpha, \beta, \gamma = 1, 2, 3$  where  $\epsilon^{\alpha\beta\gamma}$  is the Levi-Civita symbol. The fundamental representation in which the spin operators are given by the Pauli matrices show below are the simplest representation of this algebra:

$$S^1 = \frac{1}{2}\sigma^x \quad S^2 = \frac{1}{2}\sigma^y \quad S^3 = \frac{1}{2}\sigma^z$$

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

The two basis vectors are below; they are called “spin up” and “spin down”.

$$|\uparrow\rangle \equiv \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \quad |\downarrow\rangle \equiv \begin{pmatrix} 0 \\ 1 \end{pmatrix}$$

There exists two useful operators in the fundamental representation:

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

These operators act on the basis vectors as below:

$$\begin{aligned} S^+|\uparrow\rangle &= 0, & S^-|\uparrow\rangle &= |\downarrow\rangle, & S^z|\uparrow\rangle &= \frac{1}{2}|\uparrow\rangle, \\ S^+|\downarrow\rangle &= |\uparrow\rangle, & S^-|\downarrow\rangle &= 0, & S^z|\downarrow\rangle &= -\frac{1}{2}|\downarrow\rangle. \end{aligned}$$

### Heisenberg Spin Chain:

A spin chain is essentially a series of spins arranged on a 1-dimensional lattice with  $L$  sites. Each site contains a spin, and the interactions between these spins are determined by a specified quantum Hamiltonian.

### Hilbert Space:

The Hilbert space of a spin chain is constructed by taking the direct product of linear spaces corresponding to individual spins. Denoted as  $V$ , it is expressed as the tensor product of  $V_1 \otimes V_2 \otimes \dots \otimes V_L$ , where  $V_k$  represents the linear space at site  $k$ . This Hilbert space,  $V$ , has a dimension of  $2^L$ , and a convenient basis is formed by states such as  $|\uparrow\rangle_1 \otimes |\uparrow\rangle_2 \otimes \dots \otimes |\uparrow\rangle_L$  and  $|\downarrow\rangle_1 \otimes |\uparrow\rangle_2 \otimes \dots \otimes |\downarrow\rangle_L$ . The dimension corresponds to the 2 possible choices (spin-up or spin-down) at each site, and common notation condenses the tensor product symbol for brevity.

Hamiltonian: we've previously discussed a chain of  $L$  spins, where the interactions are governed by a Hamiltonian. Specifically, the Hamiltonian for the Heisenberg spin chain is now introduced.

$$\hat{H} = \sum_{n=1}^L (J_x S_n^x S_{n+1}^x + J_y S_n^y S_{n+1}^y + J_z S_n^z S_{n+1}^z) .$$

### Special cases:

Firstly,  $J_x, J_y, J_z$  are three parameters which specify how strong the spins interact in each direction. We have the following important special cases

1.  $J_x = J_y = 0, J_z \neq 0$ . This is the Ising Spin Chain.
2.  $J_z = 0, J_x = J_y \neq 0$ . This is the XX spin chain, which is equivalent to a free lattice fermion by Jordan-Wigner transformation.
3.  $J_x = J_y = J_z \neq 0$ . This is the isotropic case, which is called the XXX spin chain.
4.  $J_x = J_y \neq J_z \neq 0$ . This is the anisotropic case called XXZ spin chain.
5.  $J_x \neq J_y \neq J_z \neq 0$ . This is the completely anisotropic case, which is called the XYZ spin chain.

**Interacting range:** The interaction in the Heisenberg spin chain is characterized by nearest neighbor interactions, where each spin  $S_n^\alpha$  ( $\alpha = x, y, z$ ) only interacts with its immediate neighbor  $S_{n+1}$ . This is referred to as nearest neighboring interaction, and the term "interacting range" is defined as the number of sites involved in the Hamiltonian; commonly studied cases have a range of  $k = 2$ . However, there's a growing interest in exploring integrable spin chains with larger ranges ( $k > 2$ ), known as medium or long-range interacting spin chains. It's important

to note that the spin operators, denoted as  $S_n^\alpha$ , act locally on the spins at site-n without affecting other sites, earning them the label of local spin operators

### The XXX Spin Chain:

For the XXX Spin chain, we will modify the Hamiltonian as below:

$$\begin{aligned}
 H_{\text{XXX}} &= -J \sum_{n=1}^L (S_n^x S_{n+1}^x + S_n^y S_{n+1}^y + S_n^z S_{n+1}^z) \\
 &= -\frac{J}{2} \sum_{n=1}^L (S_n^- S_{n+1}^+ + S_n^+ S_{n+1}^- + 2S_n^z S_{n+1}^z) \quad \text{and} \\
 & \quad S_n^a = \frac{1}{2} \sigma_n^a, \quad a = x, y, z
 \end{aligned}$$

Structure of Hilbert space to simplify calculations, the Hilbert space is divided into smaller subspaces based on the number of spin-downs, considering spin-ups as the 'vacuum' and spin-downs as 'excitations.' This perspective becomes clearer in the context of the Bethe ansatz. For a spin chain of length L, the Hilbert space is decomposed into sectors with 0, 1, 2, and so on, spin-downs. For instance, for L = 3, various sectors are formed based on the number of spin-downs.

- Vacuum:  $|\uparrow\uparrow\uparrow\rangle$
- One spin-down:  $|\downarrow\uparrow\uparrow\rangle, |\uparrow\downarrow\uparrow\rangle, |\uparrow\uparrow\downarrow\rangle$
- Two spin-downs:  $|\uparrow\downarrow\downarrow\rangle, |\downarrow\uparrow\downarrow\rangle, |\downarrow\downarrow\uparrow\rangle$
- Three spin-down  $|\downarrow\downarrow\downarrow\rangle$