

Lipkin Model

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I INTRODUCTION

Consider a two level system with the lower energy state ($\sigma = -1$) having an energy of $-\epsilon/2$ and the upper energy state ($\sigma = 1$) having an energy of $\epsilon/2$. Now, consider a another system with $\Omega \in \mathbb{Z}^+$ copies of this two level system.

This gives a system with two quantum numbers, $\sigma \in \{-1, 1\}$, and $m \in \{1, 2, \dots, \Omega\}$. In Fock space, the Hamiltonian for this system is

$$\hat{H} = \frac{1}{2}\epsilon \sum_{m\sigma} \sigma a_{m\sigma}^\dagger a_{m\sigma} - \frac{1}{2}V \sum_{mm'\sigma} a_{m\sigma}^\dagger a_{m'\sigma}^\dagger a_{m'-\sigma} a_{m-\sigma}. \quad (\text{I.1})$$

The V term here is the two particle interaction term between adjacent particles.

II QUASI-SPIN OPERATORS

One can consider the operators

$$\hat{K}_3 = \frac{1}{2} \sum_{m=1}^{\Omega} (a_{m+}^\dagger a_{m+} - a_{m-}^\dagger a_{m-}), \quad (\text{II.1})$$

$$\hat{K}_+ = \sum_{m=1}^{\Omega} a_{m+}^\dagger a_{m-}, \quad (\text{II.2})$$

$$\hat{K}_- = (\hat{K}_+)^{\dagger}, \quad (\text{II.3})$$

$$= \sum_{m=1}^{\Omega} a_{m-}^\dagger a_{m+}. \quad (\text{II.4})$$

Using these "quasi-spin" operators, one can write I.1 as

$$\hat{H} = \epsilon \hat{K}_3 - \frac{1}{2}V(\hat{K}_+^2 + \hat{K}_-^2). \quad (\text{II.5})$$

Furthermore, there are two additional "quasi-spin" operators given by the equations:

$$\hat{K}_1 = \frac{1}{2}(\hat{K}_+ + \hat{K}_-), \quad (\text{II.6})$$

$$\hat{K}_2 = \frac{1}{2i}(\hat{K}_+ - \hat{K}_-). \quad (\text{II.7})$$

These create a "quasi-spin" space, similar to the angular momentum space with three components representing the three spatial dimensions. One can show that these operators are $SU(2)$ generators, just like the angular momentum operators, by showing that the following commutation relations hold (see the appendix for the algebra):

$$[\hat{K}_+, \hat{K}_-] = 2\hat{K}_3, \quad (\text{II.8})$$

$$[\hat{K}_3, \hat{K}_{\pm}] = \pm \hat{K}_{\pm}. \quad (\text{II.9})$$

A Signature Operator

Consider the quasi-spin vector, \mathbf{K} , where

$$\mathbf{K} \cdot \mathbf{k}_i = \langle \hat{K}_i \rangle, \quad (\text{II.10})$$

where \mathbf{k}_i are basis vectors in the quasi-spin space. One can consider a rotation in this space by π in some plane of this space. Let the signature operator be given by

$$\hat{R} = e^{i\pi \hat{K}_3}. \quad (\text{II.11})$$

This operator has the effect of rotating by π in the k_{12} plane.

The signature operator commutes with the Hamiltonian ($[\hat{R}, \hat{H}] = 0$), so an energy eigenstate is also an eigenstate of the signature operator, with eigenvalue $r \in \{-1, 1, -i, i\}$. If there are an even number of particles, $r \in \{-1, 1\}$. If there are an odd number of particles, $r \in \{-i, i\}$.

III EXAMPLE

Consider a system with 12 particles ($N = 12$), $\epsilon = 1$, $\Omega = 12$. If one imagines "turning on" the interaction term by increasing V from zero, then the energy eigenvalues of the Hamiltonian will begin to diverge, as shown in figure III.1.

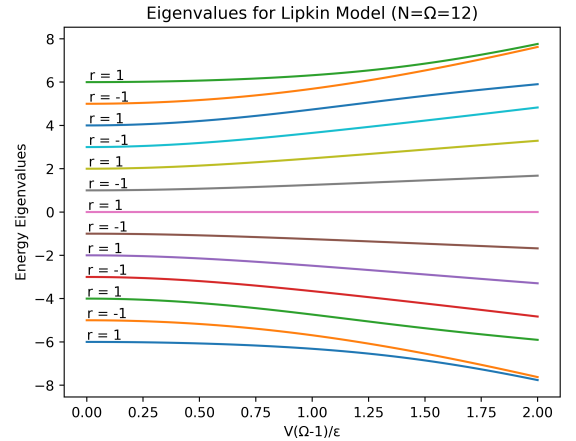


Figure III.1: Each line in this figure corresponds to an eigenvalue of the Hamiltonian, as the interaction term is "switched on". Each line is labeled with the signature eigenvalue, r . As the interaction term is increased, the eigenvalues begin to diverge. The eigenvalues start ($V = 0$) evenly spaced by a value of one, but separate as V increases.

At a value of $V(\Omega - 1)/\epsilon = 1$, the two smallest and two largest eigenvalues start becoming noticeably closer together, and the difference between them continues to decrease as this V increases, seemingly becoming arbitrarily close.

IV TRANSITION OPERATOR

The \hat{K}_1 operator is sometimes called the "transition" operator. The reason for this is unknown to me.

For the ground state of the Hamiltonian, regardless of the number of particles, value of ϵ , or value of Ω , the expectation value of this operator is zero.

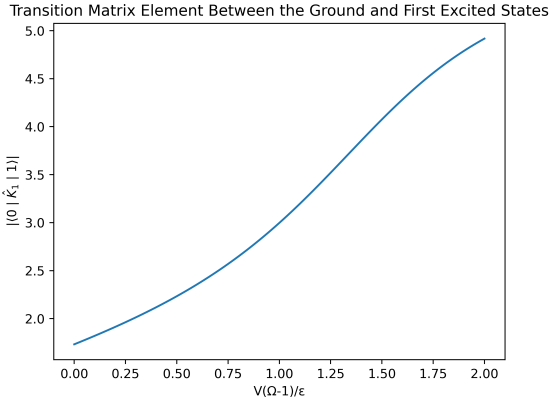


Figure IV.1: Here, the matrix element for the transition operator of the ground and first excited states is shown as a function of the $V(\Omega - 1)/\epsilon$.

V APPENDIX

A Commutation Relations

$$[\hat{K}_3, \hat{K}_\pm] = \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{m+}^\dagger a_{m+} - a_{m-}^\dagger a_{m-}) a_{n\pm}^\dagger a_{n\mp} - \frac{1}{2} \sum_{m,n=1}^{\Omega} a_{n\pm}^\dagger a_{n\mp} (a_{m+}^\dagger a_{m+} - a_{m-}^\dagger a_{m-}), \quad (\text{V.6})$$

$$= \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{m+}^\dagger a_{m+} a_{n\pm}^\dagger a_{n\mp} - a_{m-}^\dagger a_{m-} a_{n\pm}^\dagger a_{n\mp}) - \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{n\pm}^\dagger a_{n\mp} a_{m+}^\dagger a_{m+} - a_{n\pm}^\dagger a_{n\mp} a_{m-}^\dagger a_{m-}) \quad (\text{V.7})$$

$$= \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{m+}^\dagger \delta_{m+n\pm} a_{n\mp} - a_{m-}^\dagger \delta_{m-n\pm} a_{n\mp} - a_{m+}^\dagger a_{n\pm}^\dagger a_{m+} a_{n\mp} + a_{m-}^\dagger a_{n\pm}^\dagger a_{m-} a_{n\mp}) - \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{n\pm}^\dagger \delta_{n\mp m+} a_{m+} - a_{n\pm}^\dagger \delta_{n\mp m-} a_{m-} - a_{n\pm}^\dagger a_{m+}^\dagger a_{n\mp} a_{m+} + a_{n\pm}^\dagger a_{m-}^\dagger a_{n\mp} a_{m-}), \quad (\text{V.8})$$

$$= \pm \sum_{m=1}^{\Omega} a_{m\pm}^\dagger a_{m\mp} + \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{m-}^\dagger a_{n\pm}^\dagger a_{m-} a_{n\mp} - a_{m+}^\dagger a_{n\pm}^\dagger a_{m+} a_{n\mp}) + \frac{1}{2} \sum_{m,n=1}^{\Omega} (a_{n\pm}^\dagger a_{m+}^\dagger a_{n\mp} a_{m+} - a_{n\pm}^\dagger a_{m-}^\dagger a_{n\mp} a_{m-}), \quad (\text{V.9})$$

$$= \pm \hat{K}_\pm. \quad (\text{V.10})$$

$$[\hat{K}_+, \hat{K}_-] = \sum_{m,n=1}^{\Omega} a_{m+}^\dagger a_{m-} a_{n-}^\dagger a_{n+} - a_{n-}^\dagger a_{n+} a_{m+}^\dagger a_{m-}, \quad (\text{V.1})$$

$$= \sum_{m,n=1}^{\Omega} a_{m+}^\dagger (\delta_{m-n-} - a_{n-}^\dagger a_{m-}) a_{n+} - \sum_{m,n=1}^{\Omega} a_{n-}^\dagger (\delta_{m+n+} - a_{m+}^\dagger a_{n+}) a_{m-}, \quad (\text{V.2})$$

$$= \sum_{m=1}^{\Omega} a_{m+}^\dagger a_{m+} - a_{m-}^\dagger a_{m-} + \sum_{m,n=1}^{\Omega} a_{n-}^\dagger a_{m+}^\dagger a_{n+} a_{m-} - a_{m+}^\dagger a_{n-}^\dagger a_{m-} a_{n+}, \quad (\text{V.3})$$

$$= 2\hat{K}_3 + 0, \quad (\text{V.4})$$

$$= 2\hat{K}_3. \quad (\text{V.5})$$