This is a rank 4 tensor, each index taking on one of two values. It is however, natural to group the left two indices and the right two indices to write this as a 4×4 matrix. That is, we create a matrix whose rows correspond to st = 00,01,10,11. One usually just does this in one's head, but one can also formally add graphical "combiners"

where the rank 3 tensors have non-zero elements:

$$\Gamma_{v_{00}}^{s_0 t_0} = \Gamma_{v_{10}}^{s_1 t_0} = \Gamma_{v_{01}}^{s_0 t_1} = \Gamma_{v_{11}}^{s_1 t_1} \tag{1.24}$$

It is just a trivial relabeling of two indices as one composite index. In fact, when we write computer programs, we will make functions which exactly do that.

The reason for calculating E_I is that if we use periodic boundary conditions the norm is just $\langle \psi | \psi \rangle =$ Tr $(E_I)^N$. That means if λ is the largest eigenvalue of E_I , the norm is $\langle \psi | \psi \rangle = \lambda^N$. The state can then be normalized by dividing each M by $1/\sqrt{\lambda}$.

I. HW 1 – Due : Jan 29

Problem 1. (For Credit) Consider the "spin singlet": $\uparrow \downarrow - \downarrow \uparrow$. Write this as a matrix product state. Hint the matrices are 1×2 and 2×1 (so I guess this could be called a "vector product state.") Solution 1.1.

$$|\uparrow\downarrow-\downarrow\uparrow\rangle = \left(|\uparrow\rangle \quad -|\downarrow\rangle\right) \left(\begin{array}{c}|\downarrow\rangle\\|\uparrow\rangle\end{array}\right)$$
(1.25)

Problem 2. (For Credit) The "GHZ" or "cat" state of five spins is $|\uparrow\uparrow\uparrow\uparrow\uparrow\rangle + |\downarrow\downarrow\downarrow\downarrow\downarrow\rangle$. Write this as a matrix product state.

Solution 2.1.

$$|\psi\rangle = \left(|\uparrow\rangle |\downarrow\rangle \right) (1.26)$$

Problem 3. (For Credit) The "W"-state is $|\uparrow\downarrow\downarrow\downarrow\downarrow\rangle + |\downarrow\uparrow\downarrow\downarrow\downarrow\rangle + |\downarrow\downarrow\uparrow\downarrow\downarrow\rangle + |\downarrow\downarrow\downarrow\uparrow\downarrow\rangle + |\downarrow\downarrow\downarrow\downarrow\uparrow\downarrow\rangle$, Write this as a matrix product state. Hint: It is the same as what we used for the single-particle state. **Solution 3.1.**

$$|\psi\rangle = \left(|\downarrow\rangle_1 |\uparrow\rangle_1 \right) \left(|\downarrow\rangle_2 |\uparrow\rangle_2 \\ 0 |\downarrow\rangle_2 \right) \left(|\downarrow\rangle_3 |\uparrow\rangle_3 \\ 0 |\downarrow\rangle_3 \right) \left(|\downarrow\rangle_4 |\uparrow\rangle_4 \\ 0 |\downarrow\rangle_4 \right) \left(|\uparrow\rangle_5 \\ |\downarrow\rangle_5 \right)$$
(1.27)

Problem 4. (For Credit) The ferromagnetic 1D transverse field Ising Model is a spin model, defined by a Hamiltonian

$$H = \sum_{j} \left[-J\sigma_z^j \sigma_z^{j+1} - h\sigma_x^j \right].$$
(1.28)

Here σ_z and σ_x are the regular Pauli matrices. We can get a simple understanding of how this model works through a variational calculation. The simplest variational wavefunction we can use is a product: $|\psi\rangle = \otimes_j |\psi_j\rangle$ – where $|\psi_j\rangle$ is a two-component spinor. In the \hat{z} basis we can parametrize $|\psi_j\rangle$ as

$$|\psi_j\rangle = \cos(\theta/2)|\uparrow\rangle + \sin(\theta/2)|\downarrow\rangle, \qquad (1.29)$$

so that

$$\langle \psi_j | \sigma_z^j | \psi_j \rangle = \cos(\theta)$$
 (1.30)

$$\langle \psi_j | \sigma_x^j | \psi_j \rangle = \sin(\theta).$$
 (1.31)

4.1. Show that up to boundary terms (which are irrelevant in the thermodynamic limit),

$$\langle \psi | H | \psi \rangle = -JN \cos^2 \theta - hN \sin(\theta), \qquad (1.32)$$

where N is the total number of sites.

Solution 4.1. This is trivial. Each spin is independent, and there are N terms.

Problem 4. cont...

4.2. Write $x = \sin(\theta)$. Minimize $\langle \psi | H | \psi \rangle$ with respect to x (with the constraint that -1 < x < 1). Make a plot of the magnetization $m = \langle \sigma_z^j \rangle = \sqrt{1 - x^2}$ as a function of the ratio h/J.

You should see two phases – a "ferromagnetic" phase where $m \neq 0$, and a "paramagnetic" one where m = 0. This is the simplest example of what is referred to as a "quantum phase transition." Note, the mean-field theory over-estimates the stability of the ordered phase, so you should not take the numbers too seriously.

Solution 4.2. The we scale the energy, and write

$$\bar{E} = \frac{E}{JN} = -(1 - x^2) - \frac{h}{J}x.$$
(1.33)

The slope $d\bar{E}/dx = 2x - h/J$ vanishes at x = h/(2J). If h/2J < 1 the energy is minimized at x = h/2J, otherwise it is minimized at x = 1. Thus the magnetization is

$$m = \begin{cases} \sqrt{1 - \frac{h^2}{4J^2}} & h < 2J \\ 0 & h > 2J \end{cases}$$
(1.34)



Problem 5. Challenge – not for Credit

If I have N hard core bosons on M sites, and use the algorithm in Sec. D for parameterizing the states, how big is the m'th matrix? As we will see later, this is related to the entanglement entropy of a generic state. Hint: the m'th matrix in this product has d_1 rows and d_2 columns. Express d_1 and d_2 as sums over binomial coefficients. Separately consider the cases m < N and m > N. For m < N you should be able to do the sums. For m > N there is no closed form. **Solution 5.1.** The matrix has d_1 rows and d_2 columns.

Case 1: If m < N then

$$d_1 = \sum_{n=0}^{N} \begin{pmatrix} m-1\\ n \end{pmatrix} \tag{1.35}$$

and

$$d_2 = \sum_{n=0}^{N} \begin{pmatrix} m \\ n \end{pmatrix}. \tag{1.36}$$

The logic is that each row corresponds to a different configuration of the previous m-1 sites, while each column corresponds to a configuration of the first m sites. We therefore just count the number of ways of putting n particles on m-1 sites, then sum over n.

These sums are elementary, and $d_1 = 2^{m-1}$ and $d_2 = 2^m$. You can verify that this works for the case N = 2 that we explicitly did in class.

Case 2: If m > N, then we have to modify this slightly, since the configuration with N particles just gives a single row. Therefore

$$d_1 = 1 + \sum_{n=0}^{N-1} \begin{pmatrix} m-1\\ n \end{pmatrix}$$
(1.37)

and

$$d_2 = 1 + \sum_{n=0}^{N-1} \begin{pmatrix} m \\ n \end{pmatrix}.$$
 (1.38)

As stated in the question there is no closed form for these sums.