BirminghamInterviewExercise

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1 Birmingham Interview Exercise

The spin-1 matrices have the following 3x3 representations

In [167]: Sz

Out[167]:

Quantum object: dims = [[3], [3]], shape = [3, 3], type = oper, isherm = True

$$\left(\begin{array}{cccc}
1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & -1.0
\end{array}\right)$$

In [168]: Sy

Out[168]:

Quantum object: dims = [[3], [3]], shape = [3, 3], type = oper, isherm = True

$$\begin{pmatrix}
0.0 & -0.707j & 0.0 \\
0.707j & 0.0 & -0.707j \\
0.0 & 0.707j & 0.0
\end{pmatrix}$$

In [169]: Sx

Out[169]:

Quantum object: dims = [[3], [3]], shape = [3, 3], type = oper, isherm = True

$$\left(\begin{array}{cccc}
0.0 & 0.707 & 0.0 \\
0.707 & 0.0 & 0.707 \\
0.0 & 0.707 & 0.0
\end{array}\right)$$

The Hamiltonian

$$H = J_x S_1^x S_2^x + J_y S_1^y S_2^y + J_z S_1^z S_2^z$$

is the weighted sum of tensor products of spin operators, which in the spin basis can be represented as a 9x9 matrix

$$2 ext{ Jx} = ext{Jy} = ext{Jz} = 1$$

Out [309]:

Quantum object: $\dim s = [[3, 3], [3, 3]], \text{ shape} = [9, 9], \text{ type} = \text{oper, isherm} = \text{True}$

The matrix has dimension 9x9 and is not rank-deficient, and thus has 9 eigenvalues and eigenvectors. The groundstate is the eigenvector corresponding to the lowest eigenvalue (this uses a procedure from QuTiP which uses the Cholesky solver from the Numpy Library and returns the lowest energy eigenvalue and vector)

Out[342]: True

By inspection, the ground state eigenvector can be expressed in terms of the tensor products of N=3 spin states, where the expansion (schmidt) coefficients are

$$p_1 = -\sqrt{\frac{1}{3}}, p_2 = \sqrt{\frac{1}{3}}, p_3 = -\sqrt{\frac{1}{3}}$$

and the vectors are

$$v_1 = |0, 2\rangle, v_2 = |1, 1\rangle, v_3 = |2, 0\rangle$$

In [316]: ground_state_ket_1

Out[316]:

Quantum object: dims = [[3, 3], [1, 1]], shape = [9, 1], type = ket

$$\begin{pmatrix}
0.0 \\
0.0 \\
-0.577 \\
0.0 \\
0.577 \\
0.0 \\
-0.577 \\
0.0 \\
0.0
\end{pmatrix}$$

3 Jx = 0.1, Jy = 0.2, Jz = 1

The non-maximally entangled case

Out [410]:

Quantum object: dims = [[3, 3], [3, 3]], shape = [9, 9], type = oper, isherm = True

$$\begin{pmatrix} 1.0 & 0.0 & 0.0 & 0.0 & -0.050 & 0.0 & 0.0 & 0.0 & 0.0 \\ 0.0 & 0.0 & 0.0 & 0.150 & 0.0 & -0.050 & 0.0 & 0.0 & 0.0 \\ 0.0 & 0.0 & -1.0 & 0.0 & 0.150 & 0.0 & 0.0 & 0.0 & 0.0 \\ 0.0 & 0.150 & 0.0 & 0.0 & 0.0 & 0.0 & 0.0 & -0.050 & 0.0 \\ -0.050 & 0.0 & 0.150 & 0.0 & 0.0 & 0.0 & 0.150 & 0.0 & -0.050 \\ 0.0 & -0.050 & 0.0 & 0.0 & 0.0 & 0.0 & 0.150 & 0.0 & -0.050 \\ 0.0 & 0.0 & 0.0 & 0.0 & 0.150 & 0.0 & -1.0 & 0.0 & 0.0 \\ 0.0 & 0.0 & 0.0 & -0.050 & 0.0 & 0.150 & 0.0 & 0.0 & 0.0 \\ 0.0 & 0.0 & 0.0 & 0.0 & -0.050 & 0.0 & 0.0 & 0.0 & 1.0 \end{pmatrix}$$

4 Schmidt Decomposition

The schmidt decomposition is related to the SVD decomposition (for a finite dimensional Hilbert space) as follows

$$|\psi\rangle = \sum_{i} \sum_{j} C_{ij} |a\rangle_{i} \otimes |b\rangle_{j} = \sum_{i} \sqrt{p_{i}} |\alpha\rangle \otimes |\beta\rangle_{i}$$

where the SVD of the matrix C_{ij} gives vectors U, Σ, V such that

$$C = U\Sigma V^{\dagger}$$

and

$$|\alpha\rangle_i = U|a\rangle_i$$
$$|\beta\rangle_i = V^*|b\rangle_i$$
$$\Sigma_{ii} = \sqrt{p_i}$$

```
# System 2
          0.00488704 * qt.tensor(
              qt.basis(3, 0), qt.basis(3, 0)) - 0.69284525 * qt.tensor(
              qt.basis(3, 2), qt.basis(3, 0)) + 0.19970764 * qt.tensor(
              qt.basis(3, 1), qt.basis(3, 1)) - 0.69284525 * qt.tensor(
              qt.basis(3, 0), qt.basis(3, 2)) + 0.00488704 * qt.tensor(
              qt.basis(3, 2), qt.basis(3, 2));
          # canonical basis vectors for n dim system
          dim = 3
          basis = [qt.basis(dim, i).full() for i in range(dim)]
In [414]: # from inspection of system 1
          C_1 = np.array(
              [[0, 0, -np.sqrt(1/3)],
               [0, np.sqrt(1/3), 0],
               [-np.sqrt(1/3), 0, 0]])
          res(C_1).T
          # perform svd
          U_1, sig_1, Vst_1 = LA.svd(C_1)
          beta_1 = [Vst_1.T.dot(vec) for vec in basis]
          alpha_1 = [U_1.dot(vec) for vec in basis]
          schmidts_1 = [l for l in sig_1]
          ## expansion in terms of the schmidt coefficients
          ## reproduces the original state vector (to machine precision)
          assert not (sum(
                  [schmidts_1[i]*
                   qt.tensor(qt.Qobj(alpha_1[i]), qt.Qobj(beta_1[i])) for i in range(dim)]) -
                      ground_state_ket_1).full().any()
          ## new vector is normalised
          assert np.sum([schmidt**2 for schmidt in schmidts_1]) == 1
          schmidts_1
Out[414]: [0.57735026918962573, 0.57735026918962573, 0.57735026918962573]
In [418]: # from inspection of system 2
          C_2 = np.array([[0.00488704, 0, -0.69284525],
                          [0, 0.19970764, 0],
                          [-0.69284525, 0, 0.00488704]])
          # perform decomposition, produce new basis vectors and schmidt coefficients
          prec = 8
          U_2, sig2, Vst_2 = LA.svd(C_2)
          beta_2 = [np.around(Vst_2.T.dot(vec), decs) for vec in basis]
          alpha_2 = [np.around(U_2.dot(vec), decs) for vec in basis]
          schmidts_2 = [np.around(1, decs) for 1 in sig_2]
          ## The expansion in terms of the schmidt coefficients
          ## reproduces the original state vector to within 10**8 tolerance
          assert not (sum(
                  [schmidts_2[i]*qt.tensor(
                          qt.Qobj(alpha_2[i]), qt.Qobj(beta_2[i])) for i in range(dim)]) -
```

```
ground_state_ket_2).tidyup(10**-prec).full().any()
          ## new vector is normalised
          assert np.isclose(np.sum([schmidt**2 for schmidt in schmidts_2]), 1)
          schmidts_2
Out [418]: [0.69773229000000003, 0.68795821000000001, 0.1997076399999999]
```

5 Von Neumann entropies

def res(array): ret_list = [] for row in array:

def vec(array): ret_list = []

ret_list+=list(row) return np.asarray([ret_list])

ret_list+=list(col.T) return np.asarray([ret_list])

for col in array.T:

The von neumann entropy in terms of the Schmidt coefficients

```
S(p) = -\sum \rho_i ln \rho_i
```

```
In [419]: def von_neumann_entropy(nums):
              '''takes schmidt coefficients and returns von-neumann entropy'''
              return -np.sum([num**2*np.log(num**2) for num in nums])
In [420]: system_1_von_neumann_entropy = von_neumann_entropy(schmidts_1)
          system_2_von_neumann_entropy = von_neumann_entropy(schmidts_2)
In [421]: system_1_von_neumann_entropy, np.log(3)
Out [421]: (1.0986122886681096, 1.0986122886681098)
In [422]: system_2_von_neumann_entropy
Out [422]: 0.83297935252154198
  System 1 is maximally entangled, and has the corresponding entanglement entropy log(N), w/ N=3 the
number of schmidt coefficients
In [423]: # Imports, setup, tools
          import numpy as np
          import qutip as qt
          import numpy.linalg as LA
          import cmath
          Sx = qt.Qobj((1/np.sqrt(2))*np.array([[0, 1, 0], [1, 0, 1], [0, 1, 0]]))
          Sy = qt.Qobj((1/(np.sqrt(2)*1j))*np.array([[0, 1, 0], [-1, 0, 1], [0, -1, 0]]))
          Sz = qt.Qobj(np.array([[1, 0, 0], [0, 0, 0], [0, 0, -1]]))
          def re_shuffle(array, m, n):
              return np.asarray([[np.trace(np.outer(basis[i], basis[j]).T.dot(array)) for i in range(m)
          def res_shuffle(array, m, n):
              return np.asarray([[res(np.outer(basis[i], basis[j])).dot(res(array)) for i in range(m)]
```