Approximating ground states by neural network quantum states

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Abstract: The many-body problem in quantum physics originates from the difficulty of describing the non-trivial correlations encoded in the exponential complexity of the many-body wave function. Motivated by the Giuseppe Carleo’s work titled solving the quantum many-body problem with artificial neural networks [Science, 2017, 355: 602], we focus on finding the NNQS approximation of the unknown ground state of a given Hamiltonian \( H \) in terms of the best relative error and explore the influences of sum, tensor product, local unitary of Hamiltonians on the best relative error. Besides, we illustrate our method with some examples.

Keywords: Approximation; ground state; neural network quantum state

An artificial neural network (ANN) is an information processing paradigm that is inspired by the way biological nervous systems, such as the brain, process information. Up to now, there is a lot of research on the approximation ability of neural network architectures, such as Kolmogorov [1], Hornik [2], Cybenko [3], Funahashi [4], Hornik [5], Roux [6].

The many-body problem is a general name for a vast category of physical problems pertaining to the properties of microscopic systems made of a large number of interacting particles. In such a quantum system, the repeated interactions between particles create quantum correlations, or entanglement. As a consequence, the wave function of the system is a complicated object holding a large amount of information, which usually makes exact or analytical calculations impractical or even impossible. Thus, many-body theoretical physics most often relies on a set of approximations specific to the problem at hand, and ranks among the most computationally intensive fields of science. Recently, an idea that received a lot of attention from the scientific community consists in using neural networks as variational wave functions to approximate ground states of many-body quantum systems. In this direction, the networks are trained or optimized by the standard variational Monte Carlo (VMC) method while a few different neural-network architectures were tested [7–10], and the most promising results so far have been achieved with Boltzmann machines [10]. In particular, state-of-the-art numerical results have been obtained on popular models with restricted Boltzmann machines (RBM), and recent effort has demonstrated the power of deep Boltzmann machines to represent ground states of many-body Hamiltonians with polynomial-size gap and quantum states generated by any polynomial size quantum circuits [11,12]. Carleo and Troyer in [7] demonstrated the remarkable power of a reinforcement learning approach in calculating the ground state or simulating the unitary time evolution of complex quantum systems with strong interactions. Deng et al [13,14] show that this representation can be used to describe topological states. Besides, they have constructed exact representations for SPT states and intrinsic topologically ordered states. Very recently, Glasser et al [15] show that there are strong connections between neural network quantum states in the form of RBM and some classes of tensor-network states in arbitrary dimensions and obtain that neural network quantum states and their string-bond-state extension can describe a lattice fractional quantum Hall state exactly. In addition, there are a lot of related studies, such as [16–19].

Despite such exciting development, but it is unknown whether a general state can be expressed by neural networks efficiently. Generalizing the idea of [7], we introduced in [20] first neural networks...
quantum states (NNQSs) based on general input observables and explored some related properties about NNQSs, such as tensor product, local unitary operation and so on. Secondly, based on the construction of neural network representations for the cluster state in 1D, we proved necessary and sufficient conditions for a general graph state to be represented by an NNQS. We illustrated our method with some examples and observed that some $N$-qubit states can be represented by a normalized NNQS, such as any separable pure state, every Bell state, GHZ states and so on.

In this paper, based on the NNQSs introduced in [20], we focus on finding the NNQS approximation of the unknown ground state of a given Hamiltonian $H$. The remainder part of this paper are organized as follows. In Section 2, we recall the concept and the related properties of NNQSs introduced in [20]. In Section 3, we explore the NNQS approximation of the unknown ground state of a given Hamiltonian $H$ in terms of the best relative error and consider the influence of sum, tensor product, local unitary of Hamiltonian on the best relative error. Besides, we illustrate our method with some examples.

2. Neural-network quantum states

To start with, let us recall the concept and the related properties of NNQSs introduced in [20]. Let $Q_1, Q_2, \ldots, Q_N$ be $n$ quantum systems with state spaces $H_1, H_2, \ldots, H_N$ of dimensions $d_1, d_2, \ldots, d_N$, respectively. We consider the composite system $Q$ of $Q_1, Q_2, \ldots, Q_N$ with state space $H := H_1 \otimes H_2 \otimes \ldots \otimes H_N$.

Let $S_1, S_2, \ldots, S_N$ are non-degenerate observables of systems $Q_1, Q_2, \ldots, Q_N$, respectively. Then $S = S_1 \otimes S_2 \otimes \ldots \otimes S_N$ is an observable of the composite system $Q$. Use $\{|\psi_{k_j}\rangle\}_{k_j=0}^{d_j-1}$ to denote the eigenbasis of $S_j$ corresponding to eigenvalues $\{\lambda_{k_j}^{d_j} \}$.

Thus,

$$S_j |\psi_{k_j}\rangle = \lambda_{k_j} |\psi_{k_j}\rangle (k_j = 0, 1, \ldots, d_j - 1).$$

It is easy to check that the eigenvalues and corresponding eigenbasis of $S = S_1 \otimes S_2 \otimes \ldots \otimes S_N$ are

$$\lambda_{k_1} \lambda_{k_2} \ldots \lambda_{k_N} \quad \text{and} \quad |\psi_{k_1}\rangle \otimes |\psi_{k_2}\rangle \otimes \ldots \otimes |\psi_{k_N}\rangle (k_j = 0, 1, \ldots, d_j - 1),$$

respectively. Put

$$V(S) = \{ \lambda_{k_1} \lambda_{k_2} \ldots \lambda_{k_N} \equiv (\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N})^T : k_j = 0, 1, \ldots, d_j - 1 \},$$

called an input space. For parameters

$$a = (a_1, a_2, \ldots, a_N)^T \in \mathbb{C}^N, \quad b = (b_1, b_2, \ldots, b_M)^T \in \mathbb{C}^M, \quad W = [W_{ij}] \in \mathbb{C}^{M \times N},$$

write $\Omega = (a, b, W)$ and put

$$\Psi_{\Sigma, \Omega}(\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N}) = \sum_{h_i = \pm 1} \exp \left( \sum_{j=1}^{N} a_j \lambda_{k_j} + \sum_{i=1}^{M} b_j h_i + \sum_{i=1}^{M} \sum_{j=1}^{N} W_{ij} h_i \lambda_{k_j} \right).$$

Then we obtain a complex-valued function $\Psi_{\Sigma, \Omega}(\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N})$ of the input variable $\Lambda_{k_1 k_2 \ldots k_N}$. We call it a neural network quantum wave function (NNQWF). It may be identically zero. In what follows, we assume that this is not the case, that is, assume that $\Psi_{\Sigma, \Omega}(\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N}) \neq 0$ for some input variable $\Lambda_{k_1 k_2 \ldots k_N}$. Then we define

$$|\Psi_{\Sigma, \Omega}\rangle = \sum_{\Lambda_{k_1 k_2 \ldots k_N} \in V(S)} \Psi_{\Sigma, \Omega}(\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N}) |\psi_{k_1}\rangle \otimes |\psi_{k_2}\rangle \otimes \ldots \otimes |\psi_{k_N}\rangle,$$
which is a nonzero vector (not necessarily normalized) of the Hilbert space \( \mathcal{H} \). We call it a \textit{neural network quantum state} (NNQS) induced by the parameter \( \Omega = (a, b, W) \) and the input observable \( S = S_1 \otimes S_2 \otimes \ldots \otimes S_N \) (Figure 1).

**Figure 1.** Artificial neural network encoding an NNQS. It is a restricted Boltzmann machine architecture that features a set of \( N \) visible artificial neurons (blue disks) and a set of \( M \) hidden neurons (yellow disks). For each value \( \Lambda_{k_1 k_2 \ldots k_N} \) of the input observable \( S \), the neural network computes the value of the NNQWF can be reduced to

\[
\Psi_{S, \Omega}(\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N}) = \prod_{j=1}^{N} e^{a_j \lambda_j} \prod_{i=1}^{M} 2 \cosh \left( b_i + \sum_{j=1}^{N} W_{ij} \lambda_j \right).
\]

(5)

It is can be described by the following “quantum artificial neural network”, see Figure 2 where \( a = 0 \) and

\[
\sum_{i=1}^{N} (x_1, x_2, \ldots, x_N) = b_i + \sum_{j=1}^{N} x_j, 2 \cosh(\lambda) = e^\lambda + e^{-\lambda}, \quad \Pi(y_1, y_2, \ldots, y_M) = \prod_{i=1}^{M} y_i,
\]

and the final outcome \( \Psi_{S, \Omega}(\lambda_{k_1}, \lambda_{k_2}, \ldots, \lambda_{k_N}) \) is given by (5).

**Figure 2.** Quantum artificial neural network with parameter \( \Omega = (0, b, W) \).

We call this network a quantum artificial neural network because that its inputs eigenvalues of quantum observables and the outcomes are values of an NNQWF, while it has a network structure similar to a usual artificial neural network.

Next, let us consider the tensor product of the two NNQSs. We have proved the following.

**Proposition 1.** [20] Suppose that \( |\Psi^{s'}_{S', \Omega'}\rangle \) and \( |\Psi^{s''}_{S'', \Omega''}\rangle \) are two NNQSs with parameters

\[
S' = S'_1 \otimes \ldots \otimes S'_{N'}, S'' = S''_1 \otimes \ldots \otimes S''_{N''}, \Omega' = (a', b', W'), \Omega'' = (a'', b'', W''),
\]

respectively. Then \( |\Psi^{s'}_{S', \Omega'} \otimes |\Psi^{s''}_{S'', \Omega''}\rangle \) is also an NNQS \( |\Phi_{S, \Omega}\rangle \) with parameters

\[
S = S' \otimes S'', \Omega = (a, b, W), N = N' + N'', M = M' + M'',
\]
\[ a = \begin{pmatrix} a' \\ a'' \end{pmatrix}, \quad b = \begin{pmatrix} b' \\ b'' \end{pmatrix}, \quad W = [W_{ij}] = \begin{pmatrix} W_{M' \times N'} & 0 \\ 0 & W_{M'' \times N''} \end{pmatrix}. \]

Now, we discuss the influence of local unitary operation (Luo) on an NNQS. We conclude this conclusion as following.

**Proposition 2.** [20] Suppose that \(|\Psi_{S,\Omega}\rangle\) is an NNQS and \(U = U_1 \otimes U_2 \otimes \cdots \otimes U_N\) is a local unitary operator on \(H\). Then \(U|\Psi_{S,\Omega}\rangle = |\Psi_{USU^\dagger,\Omega}\rangle\), which is also an NNQS with the input observable USU\(^\dagger\) and the parameter \(\Omega\), and has the same NNQWF as \(|\Psi_{S,\Omega}\rangle\).

**Remark 1.** It can be seen from Proposition 2 that if two pure states are LU-equivalent and an NNQS representation of one of the two states is easily given, then that of another state can be obtained from that of the former.

As the end of this section, we discuss a special classes of NNQSs. When \(S = \sigma_1^2 \otimes \sigma_2^2 \otimes \cdots \otimes \sigma_N^2\), we have

\[ \lambda_k = \begin{cases} 1, & k_j = 0 \\ -1, & k_j = 1 \end{cases}, \quad |\psi_{k_j}\rangle = \begin{cases} |0\rangle, & k_j = 0 \\ |1\rangle, & k_j = 1 \end{cases} \quad (1 \leq j \leq N), \]

and \(V(S) = \{1, -1\}^N\).

In this case, the NNQS (4) becomes

\[ |\Psi_{S,\Omega}\rangle = \sum_{\lambda_{k_1},\lambda_{k_2},\ldots,\lambda_{k_N} \in \{1, -1\}^N} \Psi_{S,\Omega}(\lambda_{k_1},\lambda_{k_2},\ldots,\lambda_{k_N})|\psi_{k_1}\rangle \otimes |\psi_{k_2}\rangle \otimes \cdots \otimes |\psi_{k_N}\rangle. \tag{6} \]

This leads to the NNQS induced in [7] and discussed in [13]. We call such an NNQS a spin-z NNQS.

**3. Approximating ground states by neural network quantum states**

In this section, we try to find approximate solution to the static Schrödinger equation \(H|\psi\rangle = E|\psi\rangle\) for a given Hamiltonian \(H\). For example, to find approximation of ground states by neural network quantum states.

Let \(|\Psi_{S,\Omega}\rangle\) be an NNQS given by Eq.(4) and \(H\) be a Hamiltonian whose smallest eigenvalue \(E_{\text{exact}}\) is not zero. Put

\[ E_H(S, \Omega) = \frac{\langle \Psi_{S,\Omega}|H|\Psi_{S,\Omega}\rangle}{\langle \Psi_{S,\Omega}|\Psi_{S,\Omega}\rangle}. \]

We seek the minimum relative error between \(E_H(S, \Omega)\) and \(E_{\text{exact}}\) over \(\Omega\),

\[ \epsilon = \min_\Omega \frac{|E_H(S, \Omega) - E_{\text{exact}}|}{|E_{\text{exact}}|}. \tag{7} \]

We call \(\epsilon\) the best relative error between \(E_H(S, \Omega)\) and \(E_{\text{exact}}\). The neural network quantum state \(|\Psi_{S,\Omega}\rangle\) corresponding to the minimum of \(\epsilon\) is the best neural network representation of the ground state of \(H\).

Generally, \(E_H(S, \Omega) \geq E_{\text{exact}}\). Hence, \(\epsilon\) can also be expressed as

\[ \epsilon = \min_\Omega \frac{E_H(S, \Omega) - E_{\text{exact}}}{|E_{\text{exact}}|}. \]

Next, we discuss the influence of the sum of Hamiltonians on the best relative error. We obtain the following conclusion.
Proposition 3. Suppose that \( H_1 \) and \( H_2 \) are two Hamiltonians, \( E_{1,\text{exact}}' \), \( E_{1,\text{exact}}'' \) and \( E_{\text{exact}} \) are the smallest eigenvalue of \( H_1 \), \( H_2 \) and \( H_1 + H_2 \), respectively, \(|\Psi_{S,\Omega}\rangle\) is an NNQS. Then

\[
E_{H_1+H_2}(S,\Omega) = E_{H_1}(S,\Omega) + E_{H_2}(S,\Omega).
\]

Furthermore, if \( \min_{\Omega}(E_{H_1}(S,\Omega) + E_{H_2}(S,\Omega)) = \min_{\Omega} E_{H_1}(S,\Omega) + \min_{\Omega} E_{H_2}(S,\Omega) \), then

\[0 \leq \epsilon \leq \epsilon_1 + \epsilon_2,
\]

where

\[
\epsilon_1 = \min_{\Omega} \frac{|E_{H_1}(S,\Omega) - E_{\text{exact}}'|}{|E_{\text{exact}}'|}, \quad \epsilon_2 = \min_{\Omega} \frac{|E_{H_2}(S,\Omega) - E_{\text{exact}}''|}{|E_{\text{exact}}''|}, \quad \epsilon = \min_{\Omega} \frac{|E_{H_1+H_2}(S,\Omega) - E_{\text{exact}}|}{|E_{\text{exact}}|}.
\]

Proof. We can easily compute that

\[
E_{H_1+H_2}(S,\Omega) = \langle \Psi_{S,\Omega}|H_1 + H_2|\Psi_{S,\Omega}\rangle = \langle \Psi_{S,\Omega}|H_1|\Psi_{S,\Omega}\rangle + \langle \Psi_{S,\Omega}|H_2|\Psi_{S,\Omega}\rangle = E_{H_1}(S,\Omega) + E_{H_2}(S,\Omega).
\]

It is easily seen that \( \epsilon \geq 0 \). Generally,

\[
\min_{\Omega} E_{H_1}(S,\Omega) \geq E_{\text{exact}}', \quad \min_{\Omega} E_{H_2}(S,\Omega) \geq E_{\text{exact}}'', \quad \min_{\Omega} E_{H_1+H_2}(S,\Omega) \geq E_{\text{exact}}.
\]

Besides, when \( \min_{\Omega}(E_{H_1}(S,\Omega) + E_{H_2}(S,\Omega)) = \min_{\Omega} E_{H_1}(S,\Omega) + \min_{\Omega} E_{H_2}(S,\Omega) \), we see from \( E_{\text{exact}} \geq E_{\text{exact}}' + E_{\text{exact}}'' \) that

\[
\epsilon = \frac{|\min_{\Omega} E_{H_1+H_2}(S,\Omega) - E_{\text{exact}}|}{|E_{\text{exact}}|} = \frac{|\min_{\Omega}(E_{H_1}(S,\Omega) + E_{H_2}(S,\Omega)) - E_{\text{exact}}|}{|E_{\text{exact}}|} \leq \frac{|\min_{\Omega} E_{H_1}(S,\Omega) + \min_{\Omega} E_{H_2}(S,\Omega) - E_{\text{exact}}'|}{|E_{\text{exact}}'|} + \frac{|\min_{\Omega} E_{H_2}(S,\Omega) - E_{\text{exact}}''|}{|E_{\text{exact}}''|} = \epsilon_1 + \epsilon_2.
\]

Now, we discuss the influence of tensor product of Hamiltonians on the best relative error. We get this conclusion as following.

Proposition 4. Suppose that \( H_1 \) and \( H_2 \) are two Hamiltonians, \( E_{\text{exact}}', E_{\text{exact}}'' \) and \( E_{\text{exact}} \) are the smallest eigenvalue of \( H_1 \), \( H_2 \) and \( H_1 \otimes H_2 \), respectively, \(|\Psi'_{S',\Omega'}\rangle\) and \(|\Psi''_{S'',\Omega''}\rangle\) are two NNQSs with parameters

\[
S' = S'_1 \otimes \ldots \otimes S'_{N'}, S'' = S''_1 \otimes \ldots \otimes S''_{N''}, \Omega' = (a', b', W'), \Omega'' = (a'', b'', W''),
\]

respectively. Let

\[
S_0 = S' \otimes S'', \quad \Omega_0 = (a_0, b_0, W_0), \quad N = N' + N'', \quad M_0 = M' + M'',
\]
Thus, we have

\[ E_{H_1 \otimes H_2}(S_0, \Omega_0) = E_{H_1}(S', \Omega') \cdot E_{H_2}(S'', \Omega''). \]

Furthermore, if \( H_1 \) and \( H_2 \) are positive definite, then \( e' e'' \leq e^0 \) where

\[ e' = \min_{\Omega'} E_{H_1}(S', \Omega') - E_{\text{exact}}', \quad e'' = \min_{\Omega''} E_{H_2}(S'', \Omega'') - E_{\text{exact}}'', \]

\[ e^0 = \min_{\Omega_0} \frac{|E_{H_1}(S_0, \Omega_0) - E_{\text{exact}}|}{|E_{\text{exact}}|}. \]

**Proof.** Since \( |\Psi_{S', \Omega'}\rangle \) and \( |\Psi_{S'', \Omega''}\rangle \) are two NNQSs, we know from Proposition 1 that \( |\Psi_{S', \Omega'}\rangle \otimes |\Psi_{S'', \Omega''}\rangle = |\Phi_{S_0, \Omega_0}\rangle \) is also an NNQS. Furthermore, we can compute

\[
E_{H_1 \otimes H_2}(S_0, \Omega_0) = \frac{\langle \Psi_{S_0, \Omega_0} | H_1 \otimes H_2 | \Psi_{S_0, \Omega_0} \rangle}{\langle \Psi_{S_0, \Omega_0} | \Psi_{S_0, \Omega_0} \rangle} = E_{H_1}(S', \Omega') \cdot E_{H_2}(S'', \Omega'').
\]

Since \( H_1 \) and \( H_2 \) are positive, \( E_{\text{exact}} = E_{\text{exact}}' E_{\text{exact}}'' \). Observe that

\[
\min_{\Omega'} E_{H_1}(S', \Omega') \geq E_{\text{exact}}' > 0, \quad \min_{\Omega''} E_{H_2}(S'', \Omega'') \geq E_{\text{exact}}'' > 0, \quad \min_{\Omega_0} E_{H_1 \otimes H_2}(S_0, \Omega_0) \geq E_{\text{exact}} > 0.
\]

Thus, we have

\[
e^0 = \frac{|\min_{\Omega_0} E_{H_1 \otimes H_2}(S_0, \Omega_0) - E_{\text{exact}}|}{|E_{\text{exact}}|} = \frac{|\min_{\Omega'} E_{H_1}(S', \Omega') \cdot \min_{\Omega''} E_{H_2}(S'', \Omega'') - E_{\text{exact}}'|}{|E_{\text{exact}}'| \cdot |E_{\text{exact}}''|} = \frac{|\min_{\Omega} E_{H_1}(S', \Omega') - E_{\text{exact}}'|}{|E_{\text{exact}}'|} \cdot \frac{|\min_{\Omega''} E_{H_2}(S'', \Omega'') - E_{\text{exact}}''|}{|E_{\text{exact}}''|} = e' e''.
\]

Now, we discuss the influence of local unitary operation on the best relative error. We conclude this conclusion as following.

**Proposition 5.** Suppose that \( H \) is an Hamiltonian, \( |\Psi_{S, \Omega}\rangle \) is an NNQS and \( U = U_1 \otimes U_2 \otimes \ldots \otimes U_N \) is a local unitary operator on \( H \). \( E_{\text{exact}}', E_{\text{exact}}'' \) are the smallest eigenvalue of \( H \) and \( UHU^\dagger \), respectively. Then

\[ E_{UHU^\dagger}(S, \Omega) = E_H(U^\dagger SU, \Omega), \]

and \( e = e' \) where

\[
e = \min_{\Omega} \frac{|E_H(U^\dagger SU, \Omega) - E_{\text{exact}}|}{|E_{\text{exact}}|}, \quad e' = \min_{\Omega} \frac{|E_{UHU^\dagger}(S, \Omega) - E_{\text{exact}}'|}{|E_{\text{exact}}'|}.
\]
We can easily calculate that $E_{\hat{U}S\Omega}(S, \Omega) = \langle \Psi_{S\Omega} | \hat{U}S\Omega \rangle$, which is also an NNQS. Therefore

\[
E_{\hat{U}S\Omega}(S, \Omega) = \frac{\langle \Psi_{S\Omega} | \hat{U}S\Omega \rangle}{\langle \Psi_{S\Omega} | \Psi_{S\Omega} \rangle} = \frac{\langle \Psi_{U^*S\Omega} | H | \Psi_{U^*S\Omega} \rangle}{\langle \Psi_{U^*S\Omega} | \Psi_{U^*S\Omega} \rangle} = E_H(U^*S\Omega).
\]

Since $U$ is a local unitary operator, $E_{\text{exact}} = E'_{\text{exact}}$. We can easily obtain that $\epsilon = \epsilon'$.

Lastly, we give two examples in order to illustrate our method.

**Example 1.** Suppose that $H = |00\rangle\langle 00| + 2|01\rangle\langle 01| + 3|10\rangle\langle 10| + 4|11\rangle\langle 11|$. Then $H$ can be represented under the basis $\{00\}, \{01\}, \{10\}, \{11\}$ by $H = \text{diag}(1,2,3,4)$. It is easy to see that the minimum eigenvalue of $H$ is 1, the ground state is $|00\rangle$.

Next we use spin-$z$ NNQSs

\[
|\Psi_{S\Omega}(\lambda_1, \lambda_2) = \sum_{\lambda_1, \lambda_2 \in (1,-1)^2} \Psi_{S\Omega}(\lambda_1, \lambda_2)|\psi_k\rangle \otimes |\psi_k\rangle.
\]

to approximate the ground state $|00\rangle$ of $H$, where

\[
|\Psi_{S\Omega}(\lambda_1, \lambda_2) = \prod_{j=1}^{M} e^{a_j \lambda_j} \cdot \prod_{i=1}^{M} 2 \cosh \left( b_i + \sum_{j=1}^{M} W_{ij} \lambda_j \right).
\]

When $N = M = 2$, we have

\[
|\Psi_{S\Omega} = 4e^{a_1}e^{a_2} \cosh(b_1 + W_{11} + W_{12}) \cosh(b_2 + W_{21} + W_{22})|00\rangle + 4e^{a_1}e^{-a_2} \cosh(b_1 + W_{11} - W_{12}) \cosh(b_2 + W_{21} - W_{22})|01\rangle + 4e^{-a_1}e^{a_2} \cosh(b_1 - W_{11} + W_{12}) \cosh(b_2 - W_{21} + W_{22})|10\rangle + 4e^{-a_1}e^{-a_2} \cosh(b_1 - W_{11} - W_{12}) \cosh(b_2 - W_{21} - W_{22})|11\rangle.
\]

We can easily calculate that

\[
E_H(S, \Omega) = \left( e^{a_1}e^{a_2} \cosh(b_1 + W_{11} + W_{12}) \cosh(b_2 + W_{21} + W_{22}) \right)^2 + \left( e^{a_1}e^{-a_2} \cosh(b_1 + W_{11} - W_{12}) \cosh(b_2 + W_{21} - W_{22}) \right)^2 + \left( e^{-a_1}e^{a_2} \cosh(b_1 - W_{11} + W_{12}) \cosh(b_2 - W_{21} + W_{22}) \right)^2 + \left( e^{-a_1}e^{-a_2} \cosh(b_1 - W_{11} - W_{12}) \cosh(b_2 - W_{21} - W_{22}) \right)^2.
\]

Next we seek the minimum value of $E_H(S, \Omega)$ over $\Omega$. By letting

\[
b_1 = x_1, b_2 = x_2, W_{11} = x_3, W_{12} = x_4, W_{21} = x_5, W_{22} = x_6, a_1 = x_7, a_2 = x_8,
\]
we define a function $g$ by
\[
g(x_1, x_2, \ldots, x_8) = (e^{x_1+x_5} \cdot \cosh(x_1 + x_3 + x_4) \cdot \cosh(x_2 + x_5 + x_6))^2 \\
+ 2(e^{x_1-x_5} \cdot \cosh(x_1 + x_3 - x_4) \cdot \cosh(x_2 + x_5 - x_6))^2 \\
+ 3(e^{-x_1+x_5} \cdot \cosh(x_1 - x_3 + x_4) \cdot \cosh(x_2 - x_5 + x_6))^2 \\
+ 4(e^{-x_1-x_5} \cdot \cosh(x_1 - x_3 - x_4) \cdot \cosh(x_2 - x_5 - x_6))^2)
\]

and then numerically minimize $g$ over $x_1, x_2, \ldots, x_8$ (see Figure 3).

\[\text{Figure 3. Numerically minimize } g \text{ over } x_1, x_2, \ldots, x_8 \text{ by optimization.}\]

By using Matlab, we find
\[
\min_{x_i} g(x_1, x_2, \ldots, x_8) = g(0.743, 5.788, 2.843, 4.274, 5.501, 5.148, 3.312, 1.916) = 1.
\]

We obtain
\[
\epsilon = \min_{\Omega} \frac{|E_H(S, \Omega) - E_{\text{exact}}|}{|E_{\text{exact}}|} = 0
\]

Meanwhile, the corresponding NNQS is
\[
|\Psi_{S, \Omega}⟩ = 6.6458 \times 10^{12} |00⟩ + 4.6761 \times 10^3 |01⟩ + 505.6622 |10⟩ + 406.2882 |11⟩,
\]

the normalized NNQS is
\[
|\Psi_{S, \Omega}'⟩ = \frac{|\Psi_{S, \Omega}⟩}{\sqrt{⟨\Psi_{S, \Omega} | \Psi_{S, \Omega}⟩}} \approx |00⟩.
\]

Besides, we can also calculate the distance between the actual ground state $|00⟩$ and the approximate state $|\Psi_{S, \Omega}'⟩$ to be
\[
dist(|00⟩, |\Psi_{S, \Omega}'⟩) = \| |00⟩ - |\Psi_{S, \Omega}'⟩ \| \approx 0.
\]

Example 2. Suppose that
\[
H^{(\text{cluster})}_N = -\sum_{i=1}^{N} \sigma_{i-1}^x \sigma_i^x \sigma_{i+1}^x,
\]
where \( \sigma_0^z = 1, \sigma_{N+1}^z = 1 \). It is easily to see that the minimum eigenvalue of \( H_{N}^{\text{cluster}} \) is \(-N\), the ground state is cluster state \( |C_N\rangle \). Hence, \( E_{\text{exact}} = -N \).

Next we use spin-z NNQSs

\[
|\Psi_{S,\Omega}\rangle = \sum_{\Lambda_{k1,2, \cdots , k_N} \in \{1,-1\}^N} \Psi_{S,\Omega}(\Lambda_{k1}, \Lambda_{k2}, \cdots , \Lambda_{k_N}) |\psi_{k1}\rangle \otimes |\psi_{k2}\rangle \otimes \cdots \otimes |\psi_{k_N}\rangle,
\]

to approximate the ground state \( |C_N\rangle \) of \( H_{N}^{\text{cluster}} \), where

\[
\Psi_{S,\Omega}(\Lambda_{k1}, \Lambda_{k2}, \cdots , \Lambda_{k_N}) = \prod_{j=1}^{N} e^{a_j \Lambda_{kj}} \cdot \prod_{i=1}^{M} 2 \cosh \left( b_i + \sum_{j=1}^{N} W_{ij} \Lambda_{kj} \right).
\]

(i) When \( N = M = 2 \). By letting

\[
a_1 = x_1 + x_2 i, \quad a_2 = x_3 + x_4 i, \quad b_1 = x_5 + x_6 i, \quad b_2 = x_7 + x_8 i,
\]
\[
W_{11} = x_9 + x_{10} i, \quad W_{12} = x_{11} + x_{12} i, \quad W_{21} = x_{13} + x_{14} i, \quad W_{22} = x_{15} + x_{16} i,
\]

using Matlab (see Figure 4), we find

\[
\epsilon = 1.438 \times 10^{-6},
\]

where

\[
a = \begin{pmatrix} 0.065 + 0.194 i \\ 0.008 + 0.37 i \end{pmatrix}, \quad b = \begin{pmatrix} 0.022 + 0.693 i \\ -0.431 - 0.056 i \end{pmatrix}, \quad W = \begin{pmatrix} 0.437 + 0.909 i & 0.018 + 0.733 i \\ -0.272 + 0.952 i & 0.2 + 0.771 i \end{pmatrix}.
\]

![Figure 4](https://example.com/figure4.png)

**Figure 4.** Numerically minimize \( \epsilon \) by optimization.

Meanwhile, the corresponding NNQS is

\[
|\Psi_{S,\Omega}\rangle = (3.5877 + 0.4407 i)|00\rangle + (3.5755 + 0.5038 i)|01\rangle + (3.5805 + 0.4372 i)|10\rangle - (3.5698 + 0.5169 i)|11\rangle,
\]

then normalized NNQS is

\[
|\Psi'_{S,\Omega}\rangle = \frac{|\Psi_{S,\Omega}\rangle}{\sqrt{\langle \Psi_{S,\Omega} | \Psi_{S,\Omega} \rangle}} = (0.4969 + 0.0610 i)|00\rangle + (0.4952 + 0.0704 i)|01\rangle
\]
\[
+ (0.4959 + 0.0606 i)|10\rangle - (0.4944 - 0.0716 i)|11\rangle.
\]

Besides, we can also calculate the fidelity between the actual ground state

\[
|C_2\rangle = \frac{1}{2} |00\rangle + \frac{1}{2} |01\rangle + \frac{1}{2} |10\rangle - \frac{1}{2} |11\rangle
\]
and the approximate state $|\Psi_{S,\Omega}^{'}\rangle$ to be

$$F(|C_2\rangle,|\Psi_{S,\Omega}^{'}\rangle) = |\langle C_2 | \Psi_{S,\Omega}^{'} \rangle| = 0.9999 \approx 1.$$ 

Hence, $|C_2\rangle \approx |\Psi_{S,\Omega}^{'}\rangle$.

In addition, we find that when $N = 2$, $\epsilon$ gets smaller and smaller as $M$ changes, see Table 1.

**Table 1.** The numerical simulation results of $N, M$.

<table>
<thead>
<tr>
<th>$N$</th>
<th>$M$</th>
<th>$\epsilon$</th>
</tr>
</thead>
<tbody>
<tr>
<td>2</td>
<td>2</td>
<td>$1.438 \times 10^{-6}$</td>
</tr>
<tr>
<td>2</td>
<td>4</td>
<td>$1.0716 \times 10^{-6}$</td>
</tr>
<tr>
<td>2</td>
<td>6</td>
<td>$6.7887 \times 10^{-7}$</td>
</tr>
<tr>
<td>2</td>
<td>8</td>
<td>$4.987 \times 10^{-7}$</td>
</tr>
</tbody>
</table>

(ii) When $N = 3, M = 3$. By using Matlab (see Figure 5), we find

$$\epsilon = 2.981 \times 10^{-4}.$$ 

where

$$a = \begin{pmatrix} 0.956 + 1.669i \\ 1.309 - 0.255i \\ -0.148 - 0.152i \end{pmatrix}, \quad b = \begin{pmatrix} 0.653 + 0.863i \\ 0.569 + 0.706i \\ -0.613 + 0.894i \end{pmatrix}, \quad W = \begin{pmatrix} -0.066 + 0.969i & -1.213 + 2.029i & -0.354 - 0.647i \\ -0.233 + 3.12i & 0.986 + 0.198i & 0.438 + 0.16i \\ 0.74 + 1.206i & 0.749 - 0.985i & -0.445 + 0.8i \end{pmatrix}.$$ 

**Figure 5.** Numerically minimize $\epsilon$ by optimization.

Meanwhile, the corresponding NNQS is

$$|\Psi_{S,\Omega}\rangle = (-4.5329 - 9.8797i)|000\rangle + (-4.4661 - 9.6734i)|001\rangle + (-4.5709 - 9.9717i)|010\rangle + (4.3557 + 9.7498i)|011\rangle + (-4.4258 - 9.8957i)|100\rangle + (-4.4603 - 9.6152i)|101\rangle + (4.6706 + 9.8781i)|110\rangle + (-4.1489 - 9.7979i)|111\rangle.$$
then normalized NNQS is
\[ |\Psi'_{S,\Omega}\rangle = \frac{|\Psi_{S,\Omega}\rangle}{\sqrt{\langle \Psi_{S,\Omega} | \Psi_{S,\Omega} \rangle}} = (-0.1488 - 0.3242i)|000\rangle + (-0.1466 - 0.3175i)|001\rangle \\
+ (-0.1500 - 0.3272i)|010\rangle + (0.1429 + 0.32i)|011\rangle + (-0.1452 - 0.3248i)|100\rangle \\
+ (-0.1464 - 0.3156i)|101\rangle + (0.1533 + 0.3242i)|110\rangle + (-0.1362 - 0.3215i)|111\rangle. \]

Besides, we can also calculate the fidelity between the actual ground state
\[ |C_3\rangle = \frac{1}{\sqrt{2}}(|000\rangle + |001\rangle + |010\rangle - |011\rangle + |100\rangle + |101\rangle - |110\rangle + |111\rangle) \]
and the approximate state \( |\Psi'_{S,\Omega}\rangle \) to be
\[ F(|C_3\rangle, |\Psi'_{S,\Omega}\rangle) = |\langle C_3 | \Psi'_{S,\Omega} \rangle| = 0.9999 \approx 1. \]

Hence, \( |C_3\rangle \approx |\Psi'_{S,\Omega}\rangle \).

### 4. Conclusions

In this paper, the question of approximating ground states by neural network quantum states has been discussed in terms of the best relative error (BRE), some properties of the BREs have been obtained, including the BREs of sums, tensor products, local unitary transformations of Hamiltonians. Besides, our method have been illustrated with two examples.
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References