Robust Learning Control Design for Quantum Unitary Transformations
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Abstract—Robust control design for quantum unitary transformations has been recognized as a fundamental and challenging task in the development of quantum information processing due to unavoidable decoherence or operational errors in the experimental implementation of quantum operations. In this paper, we extend the systematic methodology of sampling-based learning control (SLC) approach with a gradient flow algorithm for the design of robust quantum unitary transformations. The SLC approach first uses a “training” process to find an optimal control strategy robust against certain ranges of uncertainties. Then a number of randomly selected samples are tested and the performance is evaluated according to their average fidelity. The approach is applied to three typical examples of robust quantum transformation problems including robust quantum transformations in a three-level quantum system, in a superconducting quantum circuit, and in a spin chain system. Numerical results demonstrate the effectiveness of the SLC approach and show its potential applications in various implementation of quantum unitary transformations.

Index Terms—Quantum learning control, quantum unitary transformation, robustness, sampling-based learning control (SLC).

I. INTRODUCTION

Quantum information and quantum computation provide the possibility to run algorithms and protocols superior to those of its classical counterparts [1]–[3]. The methodology of optimal control theory has been applied in various quantum systems to achieve different goals, such as attaining a target state, or implementing a desired quantum gate. As a basic yet influential part in quantum technology, the generation of unitary transformations is indispensable to the quantum information processing because the computation carried out in the quantum logic gates are represented by unitary transformations. Since the remarkable framework of quantum Turing machine was proposed by Deutsch [4] in 1985 and the astonishing quantum order-finding algorithm was announced by Shor [5] in 1994, the research concerning quantum information processing has blossomed. Numerous promising candidates for physical implementation of quantum systems have been proposed in recent decades, such as trapped ions [6], [7], cavity quantum electrodynamics (QED) [8], [9], nuclear magnetic resonance (NMR) [10], superconducting qubits based on Josephson junctions [11]–[13], and quantum dot in the semiconductor nanostructures [14].

Due to the unavoidable existence of imperfection and uncertainties in the construction of these quantum computation architectures, it is imperative to design robust implementation strategies. For example, in NMR, a spin ensemble consisting of around $10^{23}$ particles are utilized to perform quantum information processing. The chemical shift of their spectrometers may not be known exactly [10]. In the application of QED, when exciting atoms with lasers to high-lying Rydberg states or exploiting the long-range dipole-dipole interaction between Rydberg states, noises are unavoidable in the microwave coplanar waveguide resonators [3]. In trapped ions, the bi-chromatic laser beams may slightly interfere with each other [6]. It is also common that the temperature may influence the polarization control achieved by using liquid crystal variable retarders in the semiconductor quantum dots [14]. Moreover, operations with multiple superconducting qubits may also confront with the possible fluctuations in the coupling energy of a Josephson junction [11], [12]. Hence, it is both theoretically and practically important to develop systematic approaches for robust control design for these quantum systems. Lots of work has been done concerning this problem. For example, a noise filtering method has been presented to enhance robustness in quantum control [15]. A comprehensive approach of modulation schemes has been introduced and applied to deal with the amplitude or phase noise arising from a thermal bath in two-level systems [16]. Zhang et al. [17] used the idea of sampling uncertainty parameters to design robust control pulses for electron shuttling. A sliding model control scheme has been adopted to deal with uncertainties in two-level quantum systems [18]–[20]. In addition, stimulated Raman adiabatic passage has been extensively studied for its independence of the pulse shape, which makes it robust against moderate fluctuations in the experimental parameters [21]–[24].

On the other hand, several methods have been proposed to deal with the robust transformation problems in quantum systems. In [25], the effect of field noise upon target unitary transformations has been analyzed by investigating the spectral relationship between the Hessian and the noise in a general manner. The relationship between control time and robustness...
of a quantum system under the influence of additive white noise has been studied in [26], and the tradeoff between the fidelity and its control time has been explored. A robust optimal control landscape for the generation of quantum unitary transformations has been proposed by studying the topology of the critical regions and Grassmannian submanifolds [27].

In classical (non-quantum) engineering, feedback control is usually the preferred choice for robust control, which is also utilized in quantum engineering [28]-[30]. For instance, feedback control with quantum measurements has been used in experiments on single photons [31] and spin ensembles [32]. Nevertheless, considering the small time scales and the issues resulting from the measurement backaction in physical quantum systems, open-loop control is a more practical choice with the current level of quantum technology. For example, Khaneja et al. [33] designed optimal control pulse sequences for NMR by applying gradient algorithms. Kosut et al. [34] proposed a sequential convex programming method for designing robust quantum manipulations. However, it is still a challenging task to provide a practical approach for the implementation of general quantum unitary transformations.

In this paper, we employ the sampling-based learning control (SLC) method for designing robust quantum unitary transformations. The SLC approach was originally presented for control design of inhomogeneous quantum ensembles [35] and robust control of quantum states [36], [37]. The approach has also been applied for the robust manipulation of superconducting qubits and a set of quantum gates in the presence of fluctuations [38] and quantum ensemble classification [39]. The SLC method includes two steps of “training” and “testing”. In the training step, several samples are selected according to the distribution of Hamiltonian uncertainties to construct an augmented system, then a gradient flow based learning and optimization algorithm is applied on the augmented system to find a robust control law for a desired quantum unitary transformation. In the process of testing step, we apply the control law obtained in the “training” step to a number of samples whose parameters are selected according to uniform distribution or truncated Gaussian distribution, and evaluate the performance according to their average fidelity.

We demonstrate the application of the SLC method to three typical examples of quantum unitary transformations. In these examples, we assume that there exist some uncertainties and aim at designing robust control fields that can achieve high-fidelity quantum unitary transformations. For simplicity, we consider time-invariant uncertainties in these numerical examples. These results are straightforwardly applicable to time-varying uncertainties. The first example is a quantum unitary transformation problem in a three-level quantum system, which is a widely used model in natural and artificial ions and atoms. The second example is concerning the superconducting quantum circuits, which have been proved to be one of the promising alternatives for quantum information processing [40]-[44]. In particular, different robust quantum transformations including SWAP, CPhase and CHadamard operations are implemented using SLC. In the third example, we investigate the application of SLC in a spin chain system, which has been widely studied and applied due to its long decoherence and relaxation time [45], [46]. Numerical results show that the SLC method is effective for robust control design of quantum unitary transformations.

The rest of the paper is organized as follows. Section II formulates the control problem of quantum unitary transformations. Section III introduce the approach of sampling-based learning control and a gradient flow based learning and optimization algorithm. The results of robust control design for quantum unitary transformations in a three-level quantum system is presented in Section IV. Section V demonstrates the application of the SLC approach to achieve three robust quantum transformations (i.e., SWAP, CPhase and CHadamard operations) in quantum superconducting circuits. In Section VI the SLC approach is utilized to learn a robust control law for a spin chain of Heisenberg XXX model. Conclusions are presented in Section VII.

II. PROBLEM FORMULATION

For a quantum system, if it can be approximated as a closed system, its state may be described by a complex vector \((\alpha_1, \alpha_2, \cdots)^T\), where \(\alpha_j (j = 1, 2, \cdots)\) are complex numbers satisfying \(\sum_j |\alpha_j|^2 = 1\). In this paper, we consider finite-dimensional systems and assume the dimension as \(D\). An operation on a quantum system can be described by a unitary transformation \(U\) which turns the system from an initial state \((\alpha_1, \alpha_2, \cdots, \alpha_D)^T\) to another state \((\beta_1, \beta_2, \cdots, \beta_D)^T\), i.e.,

\[
(\beta_1, \beta_2, \cdots, \beta_D)^T = U(\alpha_1, \alpha_2, \cdots, \alpha_D)^T,
\]

where the unitary transformation \(U\) is a \(D \times D\) complex matrix satisfying \(U^TU = I\), and \(U^\dagger\) is the conjugate transpose of \(U\).

Denoted in the matrix form, the controlled evolution of a unitary transformation \(U\) on a quantum system satisfies

\[
\begin{aligned}
\frac{d}{dt} U(t) &= -iH(t)U(t) \\
&= -i\{H_0 + \sum_{m=1}^M u_m(t)H_m\}U(t),
\end{aligned}
\]

(2)

where \(H_0\) is the free Hamiltonian of the system, \(H_m (m = 1, 2, \cdots, M)\) are the control Hamiltonians, and \(u_m(t)\) are the corresponding control pulse sequences. Here we use atomic units (a.u.) by setting the reduced Planck constant \(\hbar = 1\). At \(t = 0\), the initial unitary transformation \(U(t) = U(0) = I\).

Then we have the propagator at the final time \(T\)

\[
U(T) = T\exp(-i \int_0^T H(t)dt),
\]

(3)

where \(T\) is the time-ordering operator. In practical applications, the transfer time \(T\) is usually discretized into \(N\) equal intervals, and the control amplitudes keep constant during each interval. Let \(U_j\) denote the unitary transformation during the \(j\)th interval. Thus this quantum transformation can be divided into \(N\) smaller pieces \(U_1, U_2, \ldots, U_N\), which satisfy
where $f_0(\varepsilon_0)$ and $f_m(\varepsilon_m)$ characterize these uncertainties. We assume that $\varepsilon_0 \in [1 - E_0, 1 + E_0]$ and $\varepsilon_m \in [1 - E_m, 1 + E_m]$ are time independent. The constants $E_0 \in [0, 1]$ and $E_m \in [0, 1]$ represent the normalized bounds of the uncertainty parameters. The aim is to design a control strategy \{u_m(t), m = 1, 2, \ldots, M\} to achieve robust manipulation, i.e., to construct the unitary $U(t)$ starting from the initial transformation $U(0) = I$ to the target transformation $U_F$ with a high level of fidelity in the presence of uncertainties. In this paper, the fidelity is defined as

$$F(U_F, U(T)) = \frac{1}{D} |\langle U_F | U(T) \rangle|$$

This fidelity is often used to measure the difference between two unitary transformations \[^{47},^{48}\].

The control performance can be described by a performance function $J(u)$ for each control strategy $u = \{u_m(j), m = 1, 2, \ldots, M, j = 1, 2, \ldots, N\}$ and the control problem can be formulated as a maximization problem as follows:

$$\max_u J(u) := \max_u \{\frac{1}{D} |\langle U_F | U(T) \rangle|\}$$

s.t. $J(t) = -iH(t)U(t)$, $H(t) = f_0(\varepsilon_0)H_0 + \sum_{m=1}^{M} f_m(\varepsilon_m)u_m(t)H_m$, $t \in [0, T]$, $U(0) = I$, $\varepsilon_0 \in [1 - E_0, 1 + E_0]$, $\varepsilon_m \in [1 - E_m, 1 + E_m]$, where $J(u)$ is dependent on the control fields $u$ through the Schrödinger equation and $A[\cdot]$ denotes the average expectation fidelity with uncertainty parameters $E = \{\varepsilon_0, \varepsilon_m, m = 1, 2, \ldots, M\}$.

Now, we take the Hadamard operation on a two-level quantum system in \[^{49}\] as an example to demonstrate the demand of robust control design. The Hadamard operation is a special unitary transformation denoted as

$$\text{Hadamard} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}.$$
The task of the training step is to find an optimal control field \( u^* = (u_1, u_2, \ldots, u_M)^* \) to maximize the performance index \( F_X(u) \).

In order to obtain good performance, it is necessary to choose representative samples for the uncertainties. If the distributions of parameters are uniform, the intervals \([-E_0, E_0]\) and \([-E_m, E_m]\) can be divided into \( N_0 + 1 \) and \( N_m + 1 \) subintervals, then we select \( N_0 \) samples for \( \varepsilon_0 \) and \( N_m \) samples for \( \varepsilon_m \), i.e., \( X = N_0 \prod_{m=1}^{M} N_m \) samples in total. Usually, larger \( N_0 \) and \( N_m \) could lead to better performance. However, they require longer computational time. In this paper, we select 5 samples for each uncertainty parameter, which are randomly selected according to the parameter distributions of parameters are uniform, the intervals \([-E_0, E_0]\) and \([-E_m, E_m]\) can be divided into \( N_0 + 1 \) and \( N_m + 1 \) subintervals, then we select \( N_0 \) samples for \( \varepsilon_0 \) and \( N_m \) samples for \( \varepsilon_m \), i.e., \( X = N_0 \prod_{m=1}^{M} N_m \) samples in total. Usually, larger \( N_0 \) and \( N_m \) could lead to better performance. However, they require longer computational time. In this paper, we select 5 samples for each uncertainty parameter, which are randomly selected according to the parameter uncertainties. Each sample will be evaluated to get its performance.

In the testing step, we apply the control fields \( u^* \) obtained in the training step to a large number of additional samples, which are randomly selected according to the parameter uncertainties. If the average fidelity of all the tested samples is satisfactory, the learned control strategy is acceptable and the quantum unitary transformation we achieved is robust. In this paper, we use 1000 samples to test the learned control strategy in the testing step.

Within the SLC framework, a key task is to develop an appropriate algorithm to solve the optimization problem in the training step. The gradient flow method has shown to be one of the most efficient methods to solve optimal control problems. In this paper, we consider the problem of realizing a desired quantum unitary transformation with a high level of fidelity within a given time \( T \). The gradient flow algorithm is described as follows.

Let \( U_F \) denote the desired target quantum transformation, the performance index can be defined by the Hilbert-Schmidt norm of the difference between the target unitary transformation \( U_F \) and the controlled unitary transformation \( U(T) \) as

\[
\|U_F - U(T)\|^2 = \|U_F\|^2 - 2\text{Re}(U_F^* U(T)) + \|U(T)\|^2. \tag{12}
\]

For practical applications, considering the existence of an arbitrary global phase factor \( e^{i\varepsilon} \), the problem is changed to minimize

\[
\|U_F - e^{i\varepsilon} U(T)\|^2 = \|U_F\|^2 - 2\text{Re}(U_F^* e^{i\varepsilon} U(T)) + ||e^{i\varepsilon} U(T)||^2, \tag{13}
\]

which is equivalent to maximize \( \text{Re}(U_F^* e^{i\varepsilon} U(T)) \). It can be verified that this problem is equivalent to maximize

\[
\Phi = |\langle U_F | e^{i\varepsilon} U(T) \rangle|^2. \tag{14}
\]

With operators \( A_j \) and \( B_j \) defined as \( A_j = U_j \cdots U_1 \) and \( B_j = U_{j+1}^* \cdots U_N^* U_F = A_j U(T)^\dagger U_F \), we can derive the performance function as

\[
\Phi = |\langle U_F | e^{i\varepsilon} U(T) \rangle|^2 = \langle U_F | e^{i\varepsilon} U_N \cdots U_1 \rangle e^{i\varepsilon} U_N \cdots U_1 |U_F \rangle = \langle U_F^M \rangle_{1}^M \langle U_F^M \rangle_{1}^M = \langle B_j | A_j \rangle \langle A_j | B_j \rangle. \tag{15}
\]

Let us see how the performance \( \Phi \) changes when we perturb the control amplitude from \( u_m(j) \) to \( u_m(j) + \delta u_m(j) \) at step \( j \). According to Eq. (2), we have

\[
\frac{d}{dt} (U_j(t) + \delta U_j(t)) = -i \{ H_0 + \sum_{m=1}^{M} u_m(t) H_m + \delta u_m(j) H_m \} (U_j(t) + \delta U_j(t)). \tag{16}
\]

Compared with Eq. (2), Eq. (16) becomes

\[
\frac{d}{dt} \delta U_j(t) = -i \{ H_0 + \sum_{m=1}^{M} u_m(t) H_m \} \delta U_j(t) - i \delta u_m(j) H_m U_j(t). \tag{17}
\]

By ignoring higher order terms, we have

\[
\delta U_j(\Delta t) = \int_0^{\Delta t} -i \delta u_m(j) \delta U_j(\Delta t - \tau) H_m U_j(\tau) d\tau. \tag{18}
\]

Employing the standard formula

\[
e^{-SA} B e^{SA} = B - S [A, B] + \frac{s^2}{2!} [A, [A, B]] + \cdots, \tag{19}
\]

Eq. (18) becomes

\[
\delta U_j(\Delta t) = -i \delta u_m(j) \left( \int_0^{\Delta t} U_j(\Delta t - \tau) H_m U_j(\tau) - \frac{(\Delta t)^2}{2} i [H(j), H_m] + \cdots \right) U_j(\Delta t). \tag{20}
\]

Considering the first order approximation, we can obtain the change in \( U_j \) as

\[
\delta U_j(\Delta t) = -i \Delta t \delta u_m(j) H_m U_j(\Delta t). \tag{21}
\]

Therefore, the corresponding gradient \( \delta \Phi / \delta u_m(j) \) to first order in \( \Delta t \) is given by

\[
\frac{\delta \Phi}{\delta u_m(j)} = -\langle B_j | i \Delta t H_m A_j \rangle \langle A_j | B_j \rangle - \langle B_j | A_j \rangle \langle i \Delta t H_m A_j | B_j \rangle = -2 Re \{ \langle B_j | i \Delta t H_m A_j \rangle \langle A_j | B_j \rangle \}, \tag{22}
\]

where \( u_m(j) \) is the amplitude of \( u_m \) at the \( j \)th step. During the iteration, the control fields can be updated according to the law:

\[
u_m^{k+1}(j) = u_m^k(j) + \alpha_s \frac{\delta \Phi}{\delta u_m(j)}, \tag{23}
\]

where \( \alpha_s \) is the step size.

Remark 1: It is clear that in the iterative updating process, we try to maximize the performance function \( \Phi \) by a forward Euler method. Although as an explicit method, the forward Euler method may be not able to ensure numerical stability and there is a limitation on the step size \( \alpha_s \), it is very easy to implement and has low computation cost compared with the implicit backward Euler method. A large number of numerical
results of quantum optimization problems show that it is easy to choose a suitable step size to achieve the numerical stability for the forward Euler method.

Algorithm 1 Gradient flow method with sampling-based learning control

1: Set the index of iterations $k = 0$
2: Choose a set of arbitrary controls $u^k = \{u^0_m(j), m = 1, 2, \ldots, M, j = 1, 2, \ldots, N\}$
3: repeat (for each iterative process)
4: repeat (for each training samples $i = 1, 2, \ldots, X$)
5: Compute the propagator $U^k_i(j)$ with the control strategy $u^k$
6: $U_i(T, u^k) = U^k_i(N) \cdots U^k_i(1) U_0$
7: until $i = X$
8: Compute the fidelity $F^k$ of these $X$ samples
9: $F^k = \frac{1}{X} \sum_{i=1}^{X} \frac{1}{2} |\langle U_F | U_i(T, u^k) \rangle|$
10: repeat (for each control $u_m (m = 1, 2, \ldots, M)$ of the control field $u$)
11: $\delta^k_m(j) = -\frac{\alpha}{2} \sum_{i=1}^{X} \text{Re} \{ \langle B_i | i \Delta t H_m A_i^\dagger | A_i^\dagger | B_j^\dagger \rangle \}$
12: $u^{k+1}_m(j) = u^{k}_m(j) + \alpha \delta^k_m(j)$
13: until $m = M$
14: $k = k + 1$
15: until the learning process ends
16: The optimal control strategy $u^* = \{u^*_m\} = \{u^k_m\}, m = 1, 2, \ldots, M$

Algorithm 1 gives the algorithm of gradient flow with the SLC approach in the training step. In Algorithm 1, the updating rule (line 10-13) is formalized as a batch gradient descent method to maximize the averaged performance index from the traditional gradient ascent/descent algorithm, there are lots of other algorithms in the optimal control theory, such as the conjugate gradient method, and the biconjugate gradient method. There are also some second-order algorithms, like the Newton method or the quasi-Newton method, by using the Hessian. In this paper, the gradient ascent pulse engineering (GRAPE) method [33] is used due to its simplicity and effectiveness.

IV. ROBUST QUANTUM TRANSFORMATION IN THREE-LEVEL QUANTUM SYSTEMS

A. The system

Considering a V-type quantum system, assume that the initial state is $|\psi_1\rangle = (\alpha_1, \beta_1, \gamma_1)^T$ and the target state is $|\psi_2\rangle = (\alpha_2, \beta_2, \gamma_2)^T$, and the target quantum transformation is $U_F$ that satisfies $|\psi_2\rangle = U_F |\psi_1\rangle$. The evolution equation of the quantum transformation is given as

$$\frac{d}{dt} U(t) = -i [u_0 H_0 + u_1(t) H_1 + u_2(t) H_2 + u_3(t) H_3] U(t).$$

Fig. 2. The learning performance (a) the fidelity; (b) the infidelity versus the number of iterations for the quantum transformation $U_{3 \times 3}$ with uncertainty parameters of $\varepsilon_0$ and $\varepsilon_1$.

In this model, we use the Gell-Mann matrices to represent the free and the control Hamiltonians. Let $H_0 = \sigma_3$, $H_1 = \sigma_1$, $H_2 = \sigma_4$, and $H_3 = \sigma_6$, where

$$\sigma_1 = \begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}, \sigma_3 = \begin{pmatrix} 1 & 0 & 0 \\ 0 & -1 & 0 \\ 0 & 0 & 0 \end{pmatrix},$$

$$\sigma_4 = \begin{pmatrix} 0 & 0 & 1 \\ 0 & 0 & 0 \\ 1 & 0 & 0 \end{pmatrix}, \sigma_6 = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & 1 \\ 0 & 1 & 0 \end{pmatrix},$$

are the Gell-Mann matrices. Suppose that uncertainties exist in both of the free Hamiltonian and control Hamiltonians, i.e.,

$$H(t) = f_0(\varepsilon_0) u_0 H_0 + \sum_{m=1}^{3} f_m(\varepsilon_m) u_m(t) H_m.$$  \hspace{1cm} (26)

For simplicity, we assume that $f_0(\varepsilon_0) = \varepsilon_0$ and $f_m(\varepsilon_m) = f(\varepsilon_1) = \varepsilon_1$ for all $m = 1, 2, 3$ and $E_0 = E_1 = E$, i.e., $\varepsilon_0, \varepsilon_1 \in [1 - E, 1 + E]$, where $E \in [0, 1]$ is a given constant. $\varepsilon_0$ and $\varepsilon_1$ are assumed to have uniform distributions in their ranges of fluctuations, respectively. In the training step of SLC, an augmented system is constructed by selecting $N_0 = 5$ for $\varepsilon_0$ and $N_1 = 5$ for $\varepsilon_1$. The samples are selected as:

$$\begin{cases} 
\varepsilon_0 = 1 - E + \frac{2 \text{fix}((n - 1)/5) + 1) E}{5}, \\
\varepsilon_1 = 1 - E + \frac{2 \text{mod}((n - 1)/5) + 1) E}{5},
\end{cases}$$

where $n = 1, 2, \ldots, 25$, fix($X$) rounds $X$ to the nearest integer towards zero, and mod($X, Y$) returns the remainder of the division of $X$ by $Y$. In the testing step of SLC, we use 1000 samples for testing.
Now we use the proposed method to find a robust control sequence to achieve the quantum transformation $U_{3\times 3}$. The infidelity is adopted as the performance index (Infidelity = 1 – Fidelity). We assume $T = 8$, $u_0 = 1$, $u_m \in [-5, 5]$ and approximate each control field using piece-wise pulses that may be easy to implement in some practical quantum systems. For example, the manipulation time $T$ can be divided into 40 intervals where a constant pulse is applied during each interval. Here we use 40 piece-wise subpulses to consider the tradeoff between the computational cost and the performance. Usually, more subpulses could achieve better performance while more computational cost is required. The default initial control field is $u_m = \sin t$. The boundary of the fluctuations is set as $E = 0.2$. The iteration step size is set as $\alpha_s = 0.1$.

The training performance is illustrated in Fig. 2, which shows that the average fidelity of the augmented system converges to 0.9999 after 1,000,000 iterations. The learned control strategy is shown as in Fig. 3. Then in the testing step, the learned fields are applied to 1000 randomly selected samples whose parameters are chosen according to the uniform distribution. As shown in Fig. 4, the average fidelity reaches 0.99998 and shows that the realization of the quantum transformation under the learned control is of great robustness.

Fig. 3. The learned control strategy for the quantum transformation $U_{3\times 3}$.

![Fig. 3](image)

Fig. 4. The testing performance of the quantum transformation $U_{3\times 3}$: (a) 1000 randomly selected testing samples; (b) the testing performances regarding fidelity.

![Fig. 4](image)

### B. Numerical results

The quantum unitary transformation on a three-level quantum system can be denoted by a $3 \times 3$ unitary matrix $U_{3\times 3}$. In this section, we randomly select one as the target unitary transformation, e.g.,

$$U_{3\times 3} = \begin{pmatrix} -\frac{1}{\sqrt{3}} & \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{6}} \\ -\frac{1}{\sqrt{3}} & 0 & -\frac{1}{\sqrt{6}} \\ -\frac{1}{\sqrt{3}} & \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{6}} \end{pmatrix}. \quad (28)$$

In this section, the proposed approach is applied to some physical systems of superconducting circuits. The presented method is very flexible in the selection of the initial control $u_m(t)$ and the operation time $T$, as well as the target unitary transformation. It is also robust against fluctuations in different parameters.

### V. Robust Unitary Transformation in Superconducting Circuits

In this section, the proposed approach is applied to some physical systems of superconducting circuits. The presented method is very flexible in the selection of the initial control $u_m(t)$ and the operation time $T$, as well as the target unitary transformation. It is also robust against fluctuations in different parameters.

#### A. The physical setup

Considering the physical realization of quantum computers, solid-state devices may be promising candidates. Among them, superconducting quantum circuits have been widely studied. Based on Josephson junctions, these superconducting circuits can behave quantum mechanically like artificial atoms, offering a promising way for quantum information processing. Furthermore, superconducting quantum circuits provide efficient solutions for quantum computer architectures when extended to a large number of qubits due to their exceptional ability of scalability, tunability and design flexibility.

Following the early superconducting qubits scheme proposed by Shnirman et al. [42], a series of results of superconducting qubits have been conducted to explore the properties. In superconducting quantum circuits, the Josephson qubit can be achieved in a Cooper-pair box, which is a small superconducting island weakly coupled to a bulk superconductor through a Josephson-Junction and driven by a voltage source through a gate capacitance. In a superconducting qubit, two significant quantities are the Josephson coupling energy $E_J$ and the charging energy $E_C$, whose ratio determines that the behaviour of qubit is dominated by phase or charge [43, 44].

- **Fig. 2**: The training performance of the quantum transformation.
- **Fig. 3**: The learned control strategy for the quantum transformation $U_{3\times 3}$.
- **Fig. 4**: The testing performance of the quantum transformation $U_{3\times 3}$.
(see Fig. 5(a)). When $E_C \gg E_J$, a superconducting charge qubit is constructed, whose Hamiltonian can be described as

$$H = E_C (n - n_g)^2 - E_J \cos \phi,$$  

(29)

where the phase drop $\phi$ across the Josephson-Junction is conjugate to the number $n$ of extra Cooper pairs in the box. $n_g = C_g V_g / 2e$ is controlled by the external gate voltage $V_g$, $C_g$ is the gate capacitance and $2e$ is the charge of each Cooper pair. In most experiments, in order to get a better control over the qubit, physicists often use a dc superconducting interference device (SQUID) loop instead, which is constructed by two Josephson-Junctions, as shown in Fig. 5(b). The Hamiltonian of the system can be approximated as

$$H = f(V_g) \sigma_z - g(\Phi) \sigma_x,$$  

(30)

where $f(V_g)$ is relevant to the charging energy $E_C$ and can be adjusted by $V_g$, $g(\Phi)$ is relevant to the coupling energy $E_J$ and can be controlled by the magnetic flux $\Phi$ applied through the SQUID loop. The Pauli matrices $\sigma = (\sigma_x, \sigma_y, \sigma_z)$ are

$$\sigma_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \sigma_y = \begin{pmatrix} 0 & i \\ i & 0 \end{pmatrix}, \sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}.$$  

(31)

Here, we consider an example of two coupled superconducting qubits [12]. Each qubit is a nonlinear resonator built from an Al/AlO$_x$/Al Josephson junction, and the two qubits are coupled via a modular four-terminal device. This four-terminal device is constructed using two nontunable inductors, a fixed mutual inductance and a tunable inductance. The equivalent Hamiltonian can be described as [11]

$$H = \frac{\hbar \omega_1(t)}{2} \sigma_x^{(1)} + \frac{\hbar \omega_2(t)}{2} \sigma_z^{(2)} + \frac{\hbar \omega_3(t)}{2} \sigma_x^{(1)} + \frac{\hbar \omega_4(t)}{2} \sigma_x^{(2)} + \frac{\hbar \Omega_c(t)}{2} (\sigma_x^{(1)} \sigma_x^{(2)} + \frac{1}{6N_{q1}N_{q2}} \sigma_z^{(1)} \sigma_z^{(2)}),$$  

(32)

where $N_{q1}$ and $N_{q2}$ are the number of levels in the potentials of qubits 1 and 2. The typical values for $N_{q1}$ and $N_{q2}$ are $N_{q1} = N_{q2} = 5$.

B. Numerical results

We assume that the frequencies $\omega_1(t), \omega_2(t) \in [-5, 5]$ GHz can be adjusted by changing the bias currents of two phase qubits, and $\Omega_c(t) \in [-800, 800]$ MHz can be adjusted by changing the bias current in the coupler. Let $\omega_3 = \omega_4 = 1$ GHz, the operation time $T = 8$ ns and $T$ is divided into 40 smaller time intervals. The iteration step size is set as $\alpha = 0.1$. The default initial control fields are $\omega_1 = \omega_2 = \sin t$ GHz, $\Omega_c = 0.05 \sin t$ GHz.

Due to possible fluctuations, we assume that the practical Hamiltonian has the following form

$$H = \frac{\hbar \epsilon_1 \omega_1(t)}{2} \sigma_x^{(1)} + \frac{\hbar \epsilon_2 \omega_2(t)}{2} \sigma_z^{(2)} + \frac{\hbar \epsilon_3 \omega_3(t)}{2} \sigma_x^{(1)} + \frac{\hbar \epsilon_4 \omega_4(t)}{2} \sigma_x^{(2)} + \frac{\hbar \epsilon_5 \Omega_c(t)}{2} (\sigma_x^{(1)} \sigma_x^{(2)} + \frac{1}{30} \sigma_z^{(1)} \sigma_z^{(2)}),$$  

(33)

with $\epsilon_j \in [1 - E, 1 + E]$ ($j = 1, 2, 3$). We assume $E = 0.1$.

Now we use the proposed method to achieve the SWAP, CPhase and CHadamard gates, respectively. The SWAP gate

![Image](59x491 to 290x736)
Fig. 7. The learned control strategy for the quantum operations in the superconducting qubits, the first, second, third column is the control fields of the SWAP, CPhase and CHadamard operation respectively.

is a quantum gate which swaps the states of two qubits, and it can be represented by the matrix

$$\text{SWAP} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix}. \quad (34)$$

Controlled gates act on two or more qubits, where one or more qubits act as a control for some operation. For example, the controlled-Phase/Hadamard gate (or CPhase/CHadamard) acts on two qubits, and performs the Phase/Hadamard operation on the second qubit only when the first qubit is $|1\rangle$, otherwise leaves it unchanged. They are represented by the matrices

$$\text{CPhase} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & -1 \end{pmatrix},$$

$$\text{CHadamard} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \\ 0 & 0 & -\frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \end{pmatrix}. \quad (35)$$

The result is shown in Fig. 6. In the training step, the precision of the SWAP operation can achieve slightly above 99.0%, and the precision of the CPhase operation and the CHadamard operation can achieve around 99.90%. The learning process takes 4.2 hours using a computer with CPU@3.00GHz, Windows 7, Matlab R2013a. The learned control strategies are shown in Fig. 7 where the first, second and third columns are the control fields of the SWAP, CPhase and CHadamard operations, respectively. In the testing step, the learned fields are applied to 1000 samples which are generated randomly by selecting values of fluctuation parameters according to the uniform distribution. The average fidelity of the SWAP, CPhase and CHadamard operation reaches 0.9934, 0.9987 and 0.9991, respectively (as shown in Fig. 8), which verifies the robust realization of these operations. These results are listed in Table I. It is clear that the higher fidelity of one operation achieved in the training step, the higher its average fidelity of the tested samples can be reached in the testing step.

| Operation   | SWAP    | CPhase  | CHadamard |
|-------------|---------|---------|-----------|
| Training fidelity | 0.9935  | 0.9988  | 0.9991    |
| Testing fidelity  | 0.9934  | 0.9987  | 0.9991    |

VI. ROBUST UNITARY TRANSFORMATION IN QUANTUM SPIN CHAIN

One of the admirable features of quantum technologies is its ability to establish amazing correlations between a pair of
particles. In particular, spin chains with nearest-neighbor interaction have been recognized as prototypical quantum models, since they provide a wide range of solutions for quantum transformation [50], [51]. In addition to their simple theoretical descriptions, they can be efficiently implemented by using arrays of trapped ions [52] or optical lattices with cold atoms [53]. In this section, the presented SLC approach is further tested in a quantum spin chain system with uncertainties.

A. The physical system

As shown in Fig. 9 a theoretical model of a spin chain with nearest-neighbor interactions is demonstrated with the spin up/down in each particle representing the quantum state. The Hamiltonian of an isotropic Heisenberg spin 1/2 chain with nearest-neighbor interactions is given by

\[
H_0 = J \sum_{n=1}^{N_s-1} (S_n^x S_{n+1}^x + S_n^y S_{n+1}^y + S_n^z S_{n+1}^z),
\]

where \( H_0 \) is the Heisenberg Hamiltonian and

\[
H_c(t) = u_1^x(t) S_1^x + u_1^y(t) S_1^y + u_2^x(t) S_2^x + u_2^y(t) S_2^y
\]

is the control Hamiltonian. The time-dependent control fields may be applied only on the first two spins. \( N_s \) is the length of the spin chain, \( S_\alpha = \sigma_\alpha/2 \) (\( \alpha = x, y, z \)) are spin 1/2 operators, \( \sigma_x, \sigma_y, \sigma_z \) are Pauli operators, and \( J > 0 \) is the antiferromagnetic exchange interaction between spins. All frequencies and control field amplitudes can be expressed in units of the coupling strength \( J \), and all times in units of \( 1/J \). For convenience, throughout this section we set \( J = 1 \). The above spin chain model is called the Heisenberg XXX model.

We assume that the Hamiltonian with uncertainties can be written as

\[
H(t) = \varepsilon_0 H_0 + \varepsilon_c H_c(t),
\]

where \( \varepsilon_0 \) and \( \varepsilon_c \) represent uncertainty parameters in the free Hamiltonian and control Hamiltonian, respectively. We assume
that the uncertainty parameters satisfy \( \varepsilon_0 \in [1 - E, 1 + E] \) and \( \varepsilon_c \in [1 - E, 1 + E] \).

One role of the Heisenberg spin chain in quantum computation is to be used to perform quantum transformations. Taking the Toffoli gate, or the Controlled-Controlled-NOT (CCNOT) gate as an example, the action of the \( N_s \)-qubit gate CCNOT\(_{N_s}\), which performs the CCNOT operation on the last three qubits in the chain, can be defined as

\[
\text{CCNOT}_{N_s} := I \otimes I \otimes \ldots \otimes I \otimes \text{CCNOT}. \tag{40}
\]

The Toffoli gate (or CCNOT) acts on three qubits, and flips the third qubit if and only if the first and the second qubits are both at state \( |1 \rangle \), otherwise leaves it unchanged. This unitary transformation can be represented by the matrix

\[
\text{CCNOT} = 
\begin{pmatrix}
1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 1 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 1 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 1 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 1 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 1 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 1
\end{pmatrix}. \tag{41}
\]

The fidelity between the quantum unitary transformation \( U(T) \) and the target transformation \( U_F \) can be defined as follows

\[
F(T) = \frac{1}{2N_s} \text{tr}\{U_F^T U(T)\}, \tag{42}
\]

where \( N_s \) denotes the number of qubits under consideration in the spin chain. For simplicity, we only consider the evolution of the last three qubits in the chain in this section, i.e., \( N_s = 3 \). During the learning step of SLC, this performance function is used to measure the fidelity of the system under a given control law. An optimal control law can be found by maximizing \( F(T) \). In this section, we set transformation time \( T = 20 \). The iteration step size is set as \( \alpha_s = 0.01 \).

### B. Numerical results

In physical realization, in order to avoid the error arising from the interfere of two control pulses, these two control pulses may be applied to the system alternately. During the first half of the pulse duration \((\Delta t = T/N)\) in the \( j \)-th interval, we apply one \( x \) control to the first spin of the chain with amplitude \( u_x^j(t) \) and another \( x \) control to the second spin with amplitude \( u_x^j(t) \). That is, the system is governed by the Hamiltonian \( H^x_j = H_0 + u_x^j(t)S_z^1 + u_y^j(t)S_z^2 \). Subsequently we apply one \( y \) control with amplitude \( u_y^j(t) \) and another \( y \) control with amplitude \( u_y^j(t) \) to the first and the second spin of the chain, respectively, in the second half of the \( j \)-th time interval. Therefore the system evolves under the Hamiltonian \( H^y_j = H_0 + u_y^j(t)S_y^1 + u_y^j(t)S_y^2 \). The whole quantum unitary transformation during the operation time \( T \) may be described as

\[
U(T) = U_N^y U_N^x \cdots U_3^y U_3^x \cdots U_1^y U_1^x U_0,
\]

where \( U_j^x = e^{-iH_j^x \Delta t/2} \) and \( U_j^y = e^{-iH_j^y \Delta t/2} \) are the half-interval unitary transformation, respectively.

In this subsection, the proposed method is applied to find a robust control for a CCNOT operation of three interacting qubits in a spin chain system. The Hamiltonian and its parameters are presented in the previous subsection. The training performance is shown in Fig. 10. The average fidelity of the augmented system of the CCNOT operation converges to 0.9808 after 200,000 iterations. The learned control strategy is shown in Fig. 11. Then in the testing step, the learned fields are applied to 1000 randomly selected samples whose uncertainty parameters have truncated Gaussian distribution (with mean 1 and standard deviation \( E/3 \)), and the average fidelity of the CCNOT operation reaches 0.9924 shown in Fig. 13(b), which demonstrates the robustness of our proposed method.

In practical applications, it would be more convenient if we use less number of controls. But at the same time, the fidelity of one unitary transformation we achieved may be slightly decreased with less number of controls, as well as larger control amplitudes are usually needed. In the particular quantum model presented in the previous subsection, we may use up to six control pulse sequences; that is, we may apply both \( x \) and \( y \) direction controls on all of the three qubits. Under this circumstance, the control Hamiltonian may be rewritten as

\[
H_x(t) = \sum_{i=1}^{3} (u_x^i(t)S_z^i + u_y^i(t)S_y^i), \tag{44}
\]

and the half-interval system Hamiltonian may be rewritten as

\[
H^x_j = H_0 + \sum_{i=1}^{3} u_x^i(j)S_z^i, \tag{45}
\]

\[
H^y_j = H_0 + \sum_{i=1}^{3} u_y^i(j)S_y^i.
\]
The whole quantum unitary transformation during the operation time $T$ is the same as Eq. (43). Besides, we assume the settings of the uncertainty parameters are still the same as Eq. (39).

The training performance with six control pulse sequences is shown in Fig. 10 (dashed line). The average fidelity of the augmented system of the CCNOT operation converges to 0.9927 after 200,000 iterations, which is better than that with four controls. The learned control strategy is shown in Fig. 12. It can be found that the control amplitudes are smaller than that with four controls. Usually, it is easier to achieve the control objective using more control pulse sequences because we have more flexibility for control design. The maximum amplitude and the mean amplitude under these two circumstances are all listed in Table II. Then in the testing step, the learned fields are applied to 1000 randomly selected samples whose uncertainty parameters have truncated Gaussian distribution (with mean $\varepsilon_0$ and standard deviation $E/3$), and the average fidelity of the CCNOT operation reaches 0.9973 (see Fig. 13(c)), which further demonstrates the robustness of the proposed method. More detailed results are summarized in Table II.

All these results further prove that the higher fidelity of a unitary transformation achieved in the training step, the higher its average fidelity of the tested samples can achieve in the testing step. By comparing Table II and Table III it is clear that if the uncertainty parameters of the tested samples are selected according to the truncated Gaussian distribution other than the uniform distribution, better fidelities in the testing step may be achieved.

### VII. Conclusion

In this paper, an SLC approach was employed to achieve quantum optimal control laws for robust unitary transformations. The proposed method has been applied to three typical examples of robust control problems including robust quantum transformations in a three-level quantum system, achieving robust SWAP, CPhase and CHadamard operations in a superconducting quantum circuit, and CCNOT gate in a spin chain system. Several groups of numerical results demonstrate that even when the uncertainty parameters have quite large fluctuations, the proposed SLC method is still effective for the creation of robust unitary transformations. If we could estimate...
the fluctuation bound better (i.e., a smaller bound), we may achieve better performance using the proposed method. Our future work will focus on expanding the proposed method to other types of uncertainties and other tasks in quantum systems (e.g., synchronization and switching control), as well as developing more efficient and practical methods by combing other learning algorithms, such as reinforcement learning and approximate dynamic programming algorithms genetic algorithms or differential evolution.

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