OPTIMIZATION AND NOISE ANALYSIS OF THE QUANTUM ALGORITHM FOR SOLVING ONE-DIMENSIONAL POISSON EQUATION

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Solving differential equations is one of the most promising applications of quantum computing. Recently we proposed an efficient quantum algorithm for solving one-dimensional Poisson equation avoiding the need to perform quantum arithmetic or Hamiltonian simulation. In this paper, we further develop this algorithm to make it closer to the real application on the noisy intermediate-scale quantum (NISQ) devices. To this end, we first optimize the quantum 1D-Poisson solver by developing a new way of performing the sine transformation. The circuit depth for implementing the sine transform is reduced from n^2 to n without increasing the total qubit cost of the whole circuit, which is achieved by neatly reusing the additional ancillary quits. Then, we analyse the effect of common noise existing in the real quantum devices on our algorithm using the IBM Qiskit toolkit. We find that the phase damping noise has little effect on our algorithm, while the bit flip noise has the greatest impact. In addition, threshold errors of the quantum gates are obtained to make the fidelity of the circuit output being greater than 90%. The results of noise analysis will provide a good guidance for the subsequent work of error correction for our algorithm. The noise-analysis method developed in this work can be used for other algorithms to be executed on the NISQ devices.

Keywords: Quantum algorithm, Poisson equation, Quantum noise

1. Introduction

Quantum computer executing quantum algorithms can efficiently solve many problems that are extremely difficult for classical computer [1]. One of the most valuable applications is to solve differential equations, which is the main task in classical high-performance scientific computing [2]. A series of quantum algorithms with exponential speed-up over the classical counterparts have been developed for solving both ordinary and partial differential equations [3-7].

The main idea of most quantum algorithms for solving differential equations is to discretize the differential equations into systems of linear equations, which are then solved by the quantum linear system algorithm (QLSA) [8-9] or Hamiltonian simulation [10-12]. These algorithms aim to solve general differential equations; however, they are too expensive to be implemented using the near-term quantum computers [13-15].

Most recently, we proposed a concise quantum algorithm for solving one-dimensional (1D) Poisson equation (hereafter this algorithm will be called as quantum 1D-Possion solver) [16]. In the spirit of QLSA, our algorithm can encode the solutions of Poisson equation as a quantum state using only 3n qubits and $5/3n^3$ elementary gates, where n is the logarithmic of the number of discrete points. This algorithm aims to be implementable on the near-term noisy intermediate-scale quantum (NISQ) devices for real applications [17-18].

The prominent features of NISQ devices are the short coherent time and the noisy quantum gates. The short coherent time limits the depth of the quantum circuit, namely the total number of quantum gates; the operational error of quantum gates will change the amplitude distribution of output states. For a specific circuit representation of a quantum algorithm, it is necessary to study the influence of various noises existing in a real quantum computer on the circuit to evaluate the practicality of actually implementing the algorithm on a quantum computer.

In this paper, we first optimize the quantum 1D-Poisson solver by reducing the circuit depth to meet the requirement of limited coherent time. The depth of the circuit for implementing the sine transform is reduced from n^2 to n. Secondly, we perform noise analysis for our quantum circuit utilizing both the IBM real quantum computer, IBM SANTIAGO, and the IBM circuit simulator, Qiskit. The main contribution of the present work is the development of a particular noise-analysis method for quantum algorithms to be executed on the NISQ devices.

The paper is organized as follows. First, we provide a high-level overview of the algorithm of quantum 1D-Poisson solver. Next, we optimize the quantum 1D-Poisson solver by redesigning the way of performing the sine transform, and demonstrate the algorithm on the IBMQ_QASM_SIMULATOR. Then, noise analysis of our circuits is performed by combining the IBM real quantum computer and the circuit simulator. Finally, are the conclusions.

2. Overview of the quantum 1D-Poisson solver

The one-dimensional Poisson equation with Dirichlet boundary conditions can be expressed as follows,

$$-\frac{d^2v(x)}{dx^2} = b(x), \quad x \in (0,1),$$

$$v(0) = v(1) = 0,$$
(1)

where b(x) given as input is a smooth function and v(x) is the solution of the equation. Using central difference approximation, Eq. (1) can be discretized into a linear system of equations as follows,

$$h^{-2}(-v_{i-1} + 2v_i - v_{i+1}) = b_i + \varepsilon_i, \quad i = 1, 2, ..., N - 1$$

$$v_0 = v_N = 0,$$
(2)

where h=1/N is mesh size and the number of discrete points is N+1. Ignore the truncation error ε_i , then Eq. (2) turns to be

$$\vec{Av} = \vec{b} \iff h^{-2} \begin{pmatrix} 2 & -1 & 0 \\ -1 & \ddots & \ddots & \\ & \ddots & \ddots & -1 \\ & & -1 & 2 \end{pmatrix} \begin{pmatrix} v_1 \\ v_2 \\ \vdots \\ v_{N-1} \end{pmatrix} = \begin{pmatrix} b_1 \\ b_2 \\ \vdots \\ b_{N-1} \end{pmatrix}. \tag{3}$$

The coefficient matrix A is a tridiagonal Toeplitz matrix whose eigenvectors and corresponding eigenvalues are $u_j(k) = \sqrt{2/N} \sin(j\pi k/N)$ and $\lambda_j = 4N^2 \sin^2(j\pi/2N)$.

The quantum version of solving Eq. (3) is to produce such a quantum state that

$$|v\rangle = A^{-1}|b\rangle = \left(\sum_{j} \frac{1}{\lambda_{j}} |u_{j}\rangle\langle u_{j}|\right) \cdot \left(\sum_{j'} \beta_{j'} |u_{j'}\rangle\right) = \sum_{j} C \frac{\beta_{j}}{\lambda_{j}} |u_{j}\rangle,\tag{4}$$

where C is the normalizing constant. In Ref. [16] we show that such a state can be prepared in a very simple way, which avoid the need of invoking the high-cost subroutines of QLSA or Hamiltonian simulation. The main point is that the reciprocal of eigenvalues of matrix A can be computed according to the following equation,

$$\frac{8}{\lambda_{j}} = \left[\left(\sin \frac{\pi}{6} \right)^{m} \prod_{k=2}^{n-m} \sin \left(\frac{\left| 2^{k} - \left(2^{-m} j \right) \bmod 2^{k+1} \right|}{2^{k+1}} \pi \right) \right]^{2}. \tag{5}$$

That is, each $1/\lambda_j$ in Eq. (4) can be calculated using n-1 terms of square of sine values, where n equals $\log(N)$. Furthermore, the sine-square terms can be prepared easily by a series of single-qubit R_V rotations.

The overall quantum circuit to solve the one-dimensional Poisson equation is shown in Figure 1. The sine transform (ST) is used to change the basis of Register B from eigenstates $|u_j\rangle$ to computational basis $|j\rangle$, namely $\sum_j \beta_j |u_j\rangle \rightarrow \sum_j \beta_j |j\rangle$. Then the reciprocals of eigenvalues are calculated under the control of computational basis $|j\rangle$ through a series of R_y rotations according to Eq. (5). Details of this part can be found in Ref. [16]. After the multi-controlled CNOT operation and the inverse sine transform (ST[†]), the state evolves to

$$\sum_{j=1}^{2^{n}-1} \beta_{j} \left[\sqrt{1 - \left(\frac{8}{\lambda_{j}} \right)^{2}} \left| \gamma \right\rangle_{E} \left| 0 \right\rangle + \frac{8}{\lambda_{j}} \left| 1 \right\rangle_{E}^{\otimes 2n-2} \left| 1 \right\rangle \right] \left| \mu_{j} \right\rangle_{B}, \tag{6}$$

where $|\gamma\rangle_E$ represents all the states with at least one qubit being $|0\rangle$ in register E. Finally, the ancilla qubit is measured, and if the result is $|1\rangle$, the solution state of the one-dimensional Poisson equation is created successfully in register B. Before the measurement, amplitude amplification is usually implemented to increase the success probability of obtaining the expected state.

The complexity of the circuit in Figure 1 is 3n in qubits and $5/3n^3$ in elementary gates, which is rather low comparing with the algorithms in Refs. [19,24]. Here we find that the cost can be reduced further by changing the way of implementing the sine transform and combining it properly with the

following series of R_y rotations. In the next section, we will show the optimized quantum 1D-Poisson solver.

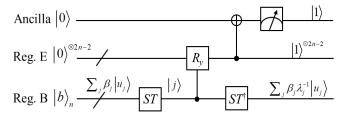


Figure 1 The overall circuit for solving one-dimensional Poisson equation. Register B encodes the input states $|b\rangle = \sum_{j} \beta_{j} |u_{j}\rangle$, and ST represents the sine transform. Register E is used to compute the reciprocals of eigenvalue through a series of R_{y} rotations. Finally, Ancilla qubit will be measured with the success flag being $|1\rangle$.

3. Optimization and demonstration of the quantum 1D-Poisson solver

3.1. Optimize the quantum 1D-Poisson solver

We optimize the quantum 1D-Poisson solver by redesigning the way of performing the sine transform (ST). The sine transform corresponds to a matrix $\left\{ST_{i,j}\right\}_{i,j=1,\dots,N-1}$ with $ST_{i,j} = \sqrt{2/N}\sin(\pi ij/N)$. Generally, the sine transform can be implemented through the Fourier transform (FT) as follows,

$$T_{2N}^{\dagger} F T_{2N} T_{2N} = \begin{pmatrix} C T_{N+1} & 0 \\ 0 & -i \cdot S T_{N-1} \end{pmatrix}, \tag{7}$$

where CT_{N+I} (ST_{N-I}) represents the cosine (sine) transform and the subscript denotes the size of the corresponding matrix [20]. As shown in Figure 2, the sine transform can be extracted easily from the matrix of $T_{2N}^{\dagger}FT_{2N}T_{2N}$ by utilizing one ancilla qubit. The quantum Fourier transform can be implemented efficiently in an exact way [21] or an approximated way [22] (with n circuit depth). Below we discuss the T_{2N} transformation, which is the starting point we improve the way of implementing the sine transform.

Ancilla
$$|1\rangle$$
 T_{2N} FT_{2N} T_{2N}^{\dagger}

Figure 2 The frame to implement the sine transform. The ancilla qubit initialized to $|1\rangle$ is used to pick out the sine transform from the transformation matrix $T_{2N}^{\dagger}FT_{2N}T_{2N}$.

The general quantum circuits for implementing the unitary T_{2N} are discussed in Refs. [23]. The T_{2N} transformation mainly consists of an add-one operation, which is usually implemented in a way as shown in Figure 3(a) [24]. The advantage of this way is that it does not need any ancillary qubits,

while the disadvantage is the heavy cost of operations and circuit depth mainly resulting from the multi-controlled NOT gates.

In the present work, we adapt the quantum ripple carry adder [25] to implement the add-one operation as shown in Figure 3(b). The carry adder can reduce the depth of the add-one circuit greatly, that is, the n n-controlled NOT gates are replaced by n TOFFOLI gates as shown in the figure. Specifically, using the new way of performing addition-one operation, the sine transform requires at most $3n^2+6n$ basic gates and the depth of the circuit is 4n. As comparison, the previous method needs at least $4n^2+n$ gates, and the depth is at least n^2+5n according to Ref. [26]. This improvement brings the quantum 1D-Poisson solver closer to the real application on quantum computer in the near future.

It is worth noting that although the carry adder in Figure 3(b) needs n-1 additional ancillary qubits comparing with the previous method in Figure 3(a), these ancillary qubits (i.e. the carry register C in Figure 3(b)) are reversed to zero and will be employed in the following R_y rotation operations as discussed below. Therefore, we reduce the circuit depth for implementing the sine transform from n^2 to n without increasing the total qubit cost of the whole circuit for solving the Poisson equations.

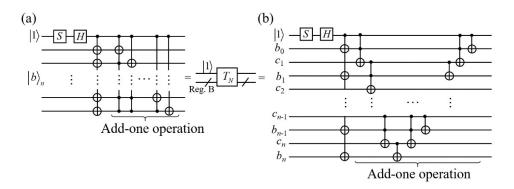


Figure 3 The quantum circuits for implementing T_{2N} transformation before (a) and after (b) optimization. In the optimized circuit (b), the ripple carry adder is used to perform add one operation, and an additional register C (i.e. $c_1, c_2...c_n$) is used to stores the carry.

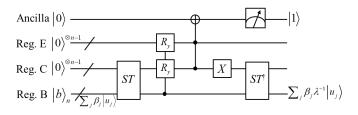


Figure 4 The optimized quantum circuit for solving one-dimensional Poisson equation. Register B and Ancilla are the same as those in Figure 1. Register C is the carry register in Figure 3. Register C and E is combined to implement the series of R_y rotations.

Combine the new sine transform with the following R_y rotations, and then we obtain the optimized quantum 1D-Poisson solver as shown in Figure 4. Comparing with the previous algorithm as shown in Figure 1, the main difference is that the Register E in Figure 1 is divided into two registers in Figure 4. One of the registers, namely Register C, is first used to implement the sine transformation before reversing the eignvalues.

Now let us sketch how the quantum states evolve through the circuit in Figure 4. As before, first the sine transform change the basis in Register B. After the implement of the R_y rotations and controlled NOT operations, the state evolves into

$$\sum_{j=1}^{2^{n}-1} \beta_{j} \left[\sqrt{1 - \left(\frac{8}{\lambda_{j}} \right)^{2}} \left| \gamma \right\rangle_{E \& C} \left| 0 \right\rangle + \frac{8}{\lambda_{j}} \left| 1 \right\rangle_{E}^{\otimes n-1} \left| 1 \right\rangle_{C}^{\otimes n-1} \left| 1 \right\rangle \right] \left| j \right\rangle, \tag{8}$$

where $|\gamma\rangle_{E\&C}$ represents all the states with at least one qubit being $|0\rangle$ in register E and C. Next, the X module flip the states of Register C qubits, i.e. the second term in equation (8) from $|1\rangle$ to $|0\rangle$ as follows,

$$\sum_{j=1}^{2^{n}-1} \beta_{j} \left[\sqrt{1 - \left(\frac{8}{\lambda_{j}} \right)^{2}} \left| \gamma \right\rangle_{E \& C} \left| 0 \right\rangle + \frac{8}{\lambda_{j}} \left| 1 \right\rangle_{E}^{\otimes n-1} \left| 0 \right\rangle_{C}^{\otimes n-1} \left| 1 \right\rangle \right] \left| j \right\rangle. \tag{9}$$

Next step is to inverse the sine transform, which transforms the state into

$$\sum_{j=1}^{2^{n}-1} \beta_{j} \sqrt{1 - \left(\frac{8}{\lambda_{j}}\right)^{2}} ST^{\dagger} |j\rangle |\gamma\rangle_{regE\&C} |0\rangle + \sum_{j=1}^{2^{n}-1} \beta_{j} \frac{8}{\lambda_{j}} |u_{j}\rangle |1\rangle_{regE}^{\otimes n-1} |0\rangle_{regC}^{\otimes n-1} |1\rangle. \tag{10}$$

Before measurement, we can perform amplitude amplification to increase the success probability of obtaining the targeted state. Finally, we measure the ancilla qubit; if the measurement result is $|1\rangle$, then the solution state to the one-dimensional Poisson equation is created successfully in Register B.

Here we remark that our algorithm solves the Poisson equations in a probabilistic way as shown by Eq. (10) that is actually an implementation of non-unitary operation. In quantum computing, the non-unitary transformation can be implemented by the method of linear combination of unitaries (LCU), which was proposed in Ref. [27]. In fact, the well-known HHL algorithm [2] is a LCU algorithm as shown in Ref. [28]; this is so for the present algorithm. The LCU method has become one of the five major techniques for designing quantum algorithms as discussed in Ref. [29], and has been extensively used, for instance, in full quantum eigensolver [30] and quantum optimization [31].

3.2. Demonstration of the optimized quantum 1D-Poisson solver

We take the cases of n=2 and n=3 for example to demonstrate the optimized quantum 1D-Poisson solver. In the cases of n=2, the discretized matrix of the Poisson equation is 3×3 corresponding to 5 discretized points, while for n=3 the matrix is 7×7 with 9 points. Figure 5 shows the circuit for n=2.

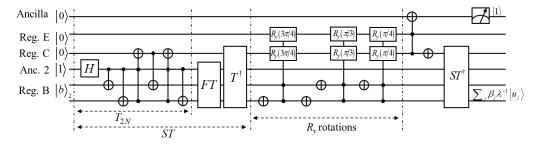


Figure 5 The demonstrated circuit for quantum 1D-Poisson equation with n=2. The Anc. 2 qubit initialized in $|1\rangle$ corresponds to the ancillary qubit in Figure 3. The NOT gate before the ST[†] transform is the X module in Figure 4. The circuit for n=3 can be obtained accordingly.

We use the IBM quantum-circuit simulator, QASM_SIMULATORv0.1.547 [17], to execute the circuits of n=2 and n=3. The initial state $|b\rangle_2$ for n=2 and $|b\rangle_3$ for n=3 are prepared as $1/\sqrt{2}|01\rangle+1/2|10\rangle+1/2|11\rangle$

and $1/2|001\rangle + \sqrt{2}/4|010\rangle + \sqrt{2}/4|011\rangle + \sqrt{2}/4|100\rangle + \sqrt{2}/4|101\rangle$ $\sqrt{2}/4|110\rangle + \sqrt{2}/4|111\rangle$ respectively. In theory, probabilities of the computational basis $\{01, 10, 11\}$ are $\{0.205, 0.304, 0.161\}$ for n=2 circuit and $\{0.029, 0.078, 0.118, 0.132, 0.114, 0.073, 0.025\}$ for basis $\{001, 010, 011, 100, 101, 110, 111\}$ of n=3 circuit. Note that the above probabilities of the computational basis are the ones when the state of Ancilla in Figure 5 is measured as $|1\rangle$, so the summation of them would not be 1.

The simulation results for n=2 and n=3 circuits are shown in Figure 6. The codes for the two circuits can be found in the Appendix A. The codes are executed on the IBM Qiskit open-source software. They are written in an intelligible quantum assembly language developed by the IBM. The simulation results are in good coincidence with the theoretical ones, and thus the optimized quantum 1D-Poisson solver is verified.

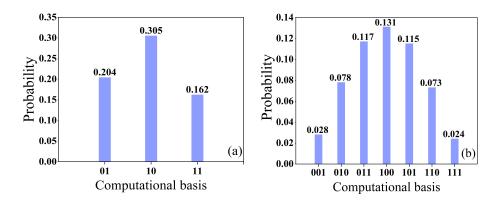


Figure 6 The simulation results of the optimized quantum 1D-Poisson solver for n=2 (a) and n=3 (b) using the IBM circuit simulator, Qiskit. The computational basis is that of Reg. B in Figure 5. The results coincidence well with the theoretical ones.

4. Noise analysis of the optimized quantum 1D-Poisson solver

4.1. Noise in the real quantum computer

Noise existing in the quantum computer is from the interaction between quantum systems and environments. Every completely positive trace-preserving map ε can be regarded as a channel which can be represented in the Kraus form [32],

$$\rho' = \varepsilon(\rho) = \sum_{k} E_{k} \rho E_{k}^{\dagger}, \tag{11}$$

where ρ is the density matrix of initial state ρ' is density matrix of final state after dynamic process and E_k represents a series of Kraus operators. It describes the dynamic evolution of the quantum system. Different noise can be described by a particular series of Kraus operators.

In general, there are four common kinds of noise, that is, amplitude damping, phase damping, bit flip and depolarizing noise. The amplitude damping noise can be used to describe the loss of energy from the quantum system to the environment, such as the process of photon emission. The Kraus operators E_k corresponding to amplitude damping are [21]

$$E_0 = \begin{bmatrix} 1 & 0 \\ 0 & \sqrt{1-p} \end{bmatrix}, \quad E_1 = \begin{bmatrix} 0 & \sqrt{p} \\ 0 & 0 \end{bmatrix}, \tag{13}$$

where $p(0 \le p \le 1)$ is the probability of a quantum state occurring error. The Kraus operator E_0 leaves the state $|0\rangle$ unchanged and reduces the amplitude of state $|1\rangle$, while E_1 changes the state $|1\rangle$ into $|0\rangle$ corresponding to the process of losing energy for the quantum system. Note that a quantum state occurring error means that there appears noise in the circuit, so the noise intensity is represented by the magnitude of the probability p. This relationship will be used below.

The phase damping noise describes the loss of quantum information of a quantum system, but without loss of energy, like the process of photon scattering [21]. Its Kraus operators are

$$E_0 = \begin{bmatrix} 1 & 0 \\ 0 & \sqrt{1-p} \end{bmatrix}, \quad E_1 = \begin{bmatrix} 0 & 0 \\ 0 & \sqrt{p} \end{bmatrix}, \tag{14}$$

where $p(0 \le p \le 1)$ is the probability of a quantum state occurring error, i.e. the probability of occurring photons scattering (without loss of energy). The Kraus operator E_0 leaves the state $|0\rangle$ unchanged and reduces the amplitude of the state $|1\rangle$, while E_1 destroys state $|0\rangle$ and reduces the amplitude of state $|1\rangle$.

The bit flip noise changes the state of a qubit from $|0\rangle$ to $|1\rangle$ and vice versa with a probability of 1-p, and leave the state unchanged with probability p. So the Kraus operators is as follows [21],

$$E_0 = \sqrt{p} \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, \quad E_1 = \sqrt{1-p} \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}.$$
 (15)

The depolarizing channel depolarizes a state of qubit into a completely mixed state I/2 with probability p, and with a probability 1-p the state is left unchanged. The Kraus operators is as follows [21],

$$E_{0} = \sqrt{1 - \frac{3p}{4}} \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, E_{1} = \frac{\sqrt{p}}{2} \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}, E_{2} = \frac{\sqrt{p}}{2} \begin{bmatrix} 0 & -i \\ i & 0 \end{bmatrix}, E_{3} = \frac{\sqrt{p}}{2} \begin{bmatrix} 1 & 0 \\ 0 & -1 \end{bmatrix}. \tag{16}$$

In order to analyze the noise effect on the quantum circuit, one needs to place the Kraus operators after each quantum gate with certain probability in the quantum circuit, and analyze how the quantum states are changed along with the probability.

4.2. Noise analysis of the quantum 1D-Poisson solver

We perform the noise analysis through three steps: first benchmarking the noise module provided by the IBM Qiskit, then analyzing the effect of the four kinds of noise discussed above respectively, finally quantifying the noise effect on the quantum 1D-Poisson solver circuit.

To start, the quantum circuits are executed in three ways, i.e. by the ideal circuit simulator without noise, by the circuit simulator with device backend noise model provided in the IBM Qiskit, and by the real quantum computer IBMQ_SANTIAGO (five qubits) [17]. The obtained results in the three scenarios for n=2 circuit are compared in Figure 7. The running codes are given in the Appendix A.

As can be seen from the Figure 7, the noise existing in the real quantum computer indeed has a great effect on the output states. In fact, the noises nearly smooth the characteristic amplitude distribution of the computational basis of the output states. The deviation would be up to 80 percent. That is, the circuit fail to encode the solutions of the 1D-Poisson equation into its output states. In addition, we can see that the device backend noise model in the IBM Qiskit can simulate, to some extent, the effect of the noise in the real quantum device. This noise model is a good approximation of the real noise, which can provide us a good reference and starting point to analyze the noise effect in more detail. In fact, the device backend noise model is generated using the calibration information of IBM's real device, which can mimic a real quantum computer approximately (here we chose IBMQ_SANTIAGO).

The backend noise model includes a single qubit depolarizing error followed by a single qubit thermal relaxation error, a two-qubit depolarizing error followed by single-qubit thermal relaxation errors, and the single-qubit readout errors when measuring individual qubit.

However, the device backend noise model provided by the IBM Qiskit can only be used to simulate the total effect of all kinds of noise. Moreover, the maximum number of bits it supports is equal to that of real device it mimics. In order to get more detail about the noise effect on our quantum 1D-Poisson solver circuit, we individually add the four common noises in section 4.1 into the circuit with reasonable error value (i.e. the noise intensity). When adding a noise to a quantum gate, the corresponding Kraus operators of the noise is placed before the logic gate.

We add the four kinds of noise into the n=2 and n=3 circuits of the quantum 1D-Poisson solver. The IBM Qiskit toolkit is employed to accomplish this task. It contains classes and functions to help building desired noise model to simulate a quantum circuit. The simulation results for the n=2 circuit is shown in Figure 8. The corresponding code is given in Appendix A. The noise intensity, namely

the magnitude of the probability p in Kraus operators, is set to be 1.8×10^{-2} according to the real parameters of the 5-qubit IBMQ_SANTIAGO quantum computer.

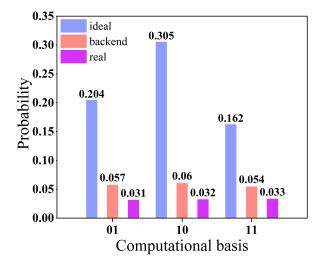


Figure 7 The execution results of the quantum 1D-Poisson solver circuit (with n=2) by ideal circuit simulator without noise, circuit simulator with device backend noise model, and real quantum computer. The real quantum computer is the IBMQ_SANTIAGO. All the probability values in the figure are the average over three sets of data and the same is true for the following figures.

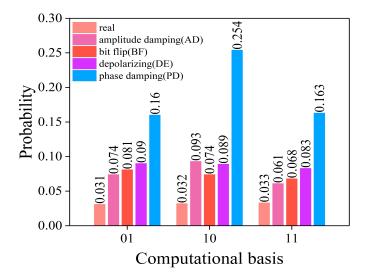


Figure 8 The running results of the n=2 quantum 1D-Poisson solver on the IBM real quantum computer, as well as the numerical simulation results under the effect of four types of noise. For each computational basis, the order of the data from left to right is the same as that in the legend from up to down.

It is interesting to see from Figure 8 that the noise of phase damping has small effect on the circuit output. This is due to the fact that our algorithm is designed to evolve the quantum states on the amplitudes that the phase damping would has little effect on the output states. The other three kinds of noise, the amplitude damping, bit flip and depolarizing channel, has similar effect on the circuit. This conclusion is useful, which tell us that when performing error correction, the phase damping error can be ignored with small loss of accuracy. Furthermore, it shows that when considering the four kinds of noise, we can properly mimic the effect of the noise existing in the real quantum device.

In the last step, we quantify the noise effect on our quantum 1D-Poisson solver. The four kinds of noise are added into n=2 and n=3 circuits with different order of magnitudes of noise intensity. Specifically, the probability p in Kraus operators for each kind of noise is set to be changed from 10^{-4} to 10^{-2} step-by-step according to the following equation [33],

$$p_i = 10^{-4} \times 700^{0.1 \times i}, \quad i = 1, 2, ..., 9.$$
 (17)

In addition, the deviation of the probability of each computational basis of the output states is taken as a measure to quantify the noise effect. It is defined as

$$\overline{D} = E_{basis}(D), \quad D = \left| \frac{p^{noise} - p^{\text{theory}}}{p^{\text{theory}}} \right|,$$
 (18)

where p^{noise} denotes the measured probability of the target computational basis with a given noise type and level, and p^{theory} is the theoretical one without noise effect. The $E_{basis}(D)$ represents the average of the deviation D over all targeted computational basis.

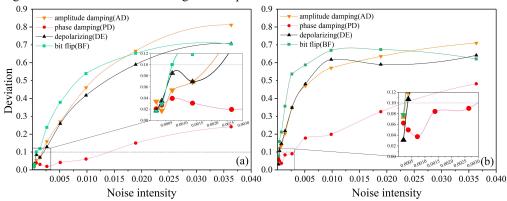


Figure 9 The deviation of the probability of targeted computational basis along with the increasing of noise intensity for the n=2 (a), and n=3 (b) 1D-Poisson solver circuit. The noise intensity is set according to Eq. (17). The specific data can be found in Table 1 and 2 in the Appendix B.

The simulation results of the deviation \overline{D} for n=2 and n=3 circuit are shown in Figure 9. (The running codes are given in the Appendix A.) Obviously, as can be seen from the figure, along with the error decreasing, the deviation become smaller and smaller. In order to guarantee that the deviation of the amplitude distribution is below 10%, the error of the quantum gates should be lower than 7.1×10^{-4} for the n=2 circuit with 131 basic quantum gates; while for the n=3 circuit with 361

basic gates, the error should lower than 1.9×10^{-4} . This result also provides a good quantitative reference for the noise effect on NISQ algorithms whose number of gates are similar or proportional to the above two circuits. In addition, we find that when the noise level is moderate (neither too imperceptible nor too noisy, which is also the typical noise level of the present NISQ devices), the bit flip noise will cause the largest deviation for the probability distribution of the output states. Furthermore, the four types of noise can be added to the quantum circuit simultaneously to approximate the noise situation in the real quantum computers.

5. Conclusion

In the present work, we optimize the quantum algorithm for solving the one-dimensional Poisson equation by reducing the circuit depth. To achieve the optimization, we develop a new way of implementing the sine transformation, and its circuit depth is reduced from n^2 to n. This new method could be a useful tool for other quantum algorithms. The optimized quantum 1D-Poisson solver is demonstrated using the IBM circuit simulator.

In order to make the quantum 1D-Poisson solver closer to the real applications on the NISQ devices, we analyze the noise effect of four common kinds of noise on the circuits. We first execute the circuits successfully on a 5-qubits real quantum computer, the IBM_SANTIAGO, and show the great effect of the noise on the circuit output states. Then we quantify the noise effect of the four kinds of noise, including the amplitude damping, phase damping, bit flip and depolarizing noise, using the IBM Qiskit toolkit. We find that the phase damping noise has little effect on our algorithm, while the bit flip noise causes the greatest effect in the concerned cases. In addition, for n=2 and n=3 quantum 1D-Poisson solver circuits with 131 and 361 basic gates, the threshold error of the quantum gates should be lower than 7.1×10^{-4} and 1.9×10^{-4} , respectively, to make the deviation of the output probability distribution below 10%. These results provide a good guidance for our subsequent work to implement error correction on our quantum 1D-Poisson solver.

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Appendix A Codes

Here we provide the Qiskit codes for the case of n=2 and 3. In the section first in order to visualize the designed circuit the code is taken the form in IBM online computing platform which can be transformed into Qiskit code easily by drop-down window on the interface [1]. Then The Qiskit code for ideal circuit simulation without noise, for device backend noise model mimicked IBMQ_SANTIAGO, for the real quantum computer IBMQ_SANTIAGO, for adding four common kinds of noise are given in turn. The execution of Qiskit code depends on the Python environment. So, first, we need install the software Anaconda, next in Anaconda prompt create a python environment. Then basing on the environment build a Qiskit package [2]. Finally typing "python" to start the simulation.

The circuit code for n=2

```
#the simulator has its basic gates of H, S, X, T, CNOT, TOFFOLI, U, U2, U3, etc
#they are converted to one & two qubit
gates to get the number of 79
#5 qubit omitted the auxiliary bits in fig.5 instead of observing Reg. E and Reg. C directly
#qubit 0 and 2 corresponds to Reg. E and C respectively
#qubit 3 and 1 Reg. B
#qubit 4 Anc.2
qreg q[5];
creg c[4];
#state b preparation
h q[3];
x q[4];
x q[3];
h q[4];
cx q[3],q[1];
x q[3];
ch q[3],q[1];
#state b preparation done
```

```
cx q[4],q[3];
cx q[4],q[1];
ccx q[4],q[1],q[2];
cx q[2],q[3];
\cos q[4],q[1],q[2];
cx q[4],q[1];
h q[4];
cu1(pi/2) q[3],q[4];
cu1(pi/4) q[1],q[4];
h q[3];
cu1(pi/2) q[1],q[3];
h q[1];
swap q[1],q[4];
cx q[4],q[1];
ccx q[4],q[1],q[2];
cx q[2],q[3];
\cos q[4],q[1],q[2];
cx q[4],q[1];
cx q[4],q[3];
#sin transform done
#controlled R_y module
x q[3];
ccx q[3],q[1],q[2];
ccx q[1],q[3],q[0];
cry(pi/8) q[1],q[2];
cry(pi/8) q[1],q[0];
cx q[3],q[1];
cry(-pi/8) q[1],q[2];
cry(-pi/8) q[1],q[0];
cx q[3],q[1];
cry(pi/8) q[3],q[2];
cry(pi/8) q[3],q[0];
x q[1];
x q[3];
cry(pi/6) q[1],q[2];
cry(pi/6) q[1],q[0];
cx q[3], q[1];
cry(-pi/6) q[1],q[2];
cry(-pi/6) q[1],q[0];
cx q[3],q[1];
cry(pi/6) q[3],q[2];
cry(pi/6) q[3],q[0];
x q[1];
cry(pi/8) q[1],q[2];
cry(pi/8) q[1],q[0];
cx q[3],q[1];
cry(-pi/8) q[1],q[2];
cry(-pi/8) q[1],q[0];
```

#sin transform

```
cx q[3],q[1];
cry(pi/8) q[3],q[2];
cry(pi/8) q[3],q[0];
#controlled Ry module done
x q[2]; #X module in Fig. 4
cx q[4],q[3];
cx q[4],q[1];
ccx q[4],q[1],q[2];
cx q[2],q[3];
ccx q[4],q[1],q[2];
cx q[4],q[1];
swap q[1],q[4];
h q[1];
cu1(-pi/2) q[1],q[3];
h q[3];
cu1(-pi/4) q[1],q[4];
cu1(-pi/2) q[3],q[4];
h q[4];
{\rm cx} \ {\rm q}[4], {\rm q}[1];
\cos q[4],q[1],q[2];
cx q[2],q[3];
ccx q[4],q[1],q[2];
cx q[4],q[1];
x q[2];
cx q[4],q[3];
h q[4];
x q[4];
measure q[3] -> c[3];
measure q[1] \rightarrow c[2];
measure q[2] -> c[1];
measure q[0] -> c[0];
The circuit code for n=3
#9 qubit and 217 one & two qubit
gates
#qubit 0 Anc
#qubit 1, 2 and 4, 6 corresponds to Reg. E and C respectively
#qubit 3, 5, 7 Reg. B
#qubit 8 Anc.2
qreg q[9];
creg c[4];
#state b preparation
h q[5];
h q[7];
x q[8];
x q[5];
x q[7];
```

```
ccx q[7],q[5],q[3];
x q[5];
x q[7];
cry(pi/4) q[5],q[3];
cx q[7],q[5];
cry(-pi/4) q[5],q[3];
cx q[7],q[5];
cry(pi/4) q[7],q[3];
x q[5];
cry(pi/4) q[5],q[3];
cx q[7],q[5];
cry(-pi/4) q[5],q[3];
cx q[7], q[5];
cry(pi/4) q[7],q[3];
x q[5];
x q[7];
cry(pi/4) q[5],q[3];
ex q[7],q[5];
cry(-pi/4) q[5],q[3];
ex q[7],q[5];
cry(pi/4) q[7],q[3];
x q[7];
#state b preparation done
#sin transform
cx q[8],q[7];
cx q[8],q[5];
cx q[8],q[3];
ccx q[3],q[8],q[4];
ccx q[4],q[5],q[6];
cx q[6],q[7];
ccx q[4],q[5],q[6];
cx q[4],q[5];
ccx q[3],q[8],q[4];
cx q[8],q[3];
h q[8];
cp(pi/2) q[7],q[8];
cp(pi/4) q[5],q[8];
cp(pi/8) q[3],q[8];
h q[7];
cp(pi/2) q[5],q[7];
cp(pi/4) q[3],q[7];
h q[5];
cp(pi/2) q[3],q[5];
h q[3];
swap q[3],q[8];
swap q[5],q[7];
cx q[8],q[3];
ccx q[3],q[8],q[4];
```

h q[8];

```
cx q[4],q[5];
ccx q[4],q[5],q[6];
cx q[6],q[7];
ccx q[4],q[5],q[6];
ccx q[3],q[8],q[4];
cx q[8],q[3];
cx q[8],q[5];
barrier q[5];
cx q[8],q[7];
#sin transform done
#controlled R_y module
x q[5];
x q[7];
cry(pi/4) q[7],q[6];
cry(pi/4) q[7],q[4];
cx q[3],q[7];
cry(-pi/4) q[7],q[6];
cry(-pi/4) q[7],q[4];
cx q[3],q[7];
cry(pi/4) q[3],q[6];
x q[7];
cry(pi/4) q[3],q[4];
cry(pi/8) q[5],q[6];
x q[7];
cry(pi/8) q[5],q[4];
cx q[3], q[5];
cry(-pi/8) q[5],q[6];
cry(-pi/8) q[5],q[4];
cx q[3],q[5];
cry(pi/8) q[3],q[6];
cry(pi/8) q[3],q[4];
x q[5];
cry(pi/8) q[3],q[6];
cry(pi/8) q[3],q[4];
cx q[7],q[5];
cry(pi/4) q[5],q[2];
cry(pi/4) q[5],q[1];
cx q[3], q[5];
cry(-pi/4) q[5],q[2];
cry(-pi/4) q[5],q[1];
cx q[3], q[5];
cry(pi/4) q[3],q[2];
cx q[7],q[5];
cry(pi/4) q[3],q[1];
cry(pi/6) q[5],q[6];
x q[7];
cry(pi/4) q[3],q[2];
cry(pi/6) q[5],q[4];
x q[7];
```

```
cry(pi/4) q[3],q[1];
x q[3];
cx q[3],q[5];
cry(-pi/6) q[5],q[6];
cry(-pi/6) q[5],q[4];
cx q[3],q[5];
cry(pi/6) q[3],q[6];
cry(pi/6) q[3],q[4];
cry(pi/4) q[7],q[2];
cry(pi/4) q[7],q[1];
ccx q[3],q[5],q[7];
cry(-pi/4) q[7],q[2];
cry(-pi/4) q[7],q[1];
ccx q[3],q[5],q[7];
cry(pi/8) q[5],q[2];
x q[7];
cry(pi/8) q[5],q[1];
cry(pi/6) q[7],q[6];
cx q[3],q[5];
cry(-pi/8) q[5],q[2];
cry(-pi/8) q[5],q[1];
cx q[3],q[5];
cry(pi/8) q[3],q[2];
cry(pi/8) q[3],q[1];
cry(pi/8) q[5],q[2];
cry(pi/8) q[5],q[1];
cx q[3], q[5];
cry(-pi/8) q[5],q[2];
cry(-pi/8) q[5],q[1];
cx q[3],q[5];
cry(pi/8) q[3],q[2];
x q[5];
cry(pi/8) q[3],q[1];
cry(pi/6) q[7],q[4];
x q[3];
x q[3];
cry(pi/6) q[7],q[2];
cry(pi/6) q[7],q[1];
ccx q[3],q[5],q[7];
cry(-pi/6) q[7],q[6];
cry(-pi/6) q[7],q[4];
cry(-pi/6) q[7],q[2];
cry(-pi/6) q[7],q[1];
ccx q[3],q[5],q[7];
cry(pi/12) q[5],q[6];
cx q[8],q[7];
cry(pi/12) q[5],q[4];
cry(pi/12) q[5],q[2];
cry(pi/12) q[5],q[1];
```

cx q[3],q[5];

```
cry(-pi/12) q[5],q[6];
cry(-pi/12) q[5],q[4];
cry(-pi/12) q[5],q[2];
cry(-pi/12) q[5],q[1];
cx q[3],q[5];
cry(pi/12) q[3],q[6];
cry(pi/12) q[3],q[4];
x q[5];
cry(pi/12) q[3],q[2];
cry(pi/12) q[3],q[1];
crz(pi) q[1],q[0];
x q[3];
#controlled R_y module done
cry(pi/2) q[1],q[0];
crz(pi) q[2],q[0];
cry(pi/2) q[2],q[0];
ccx q[4],q[6],q[0];
cry(-pi/2) q[2],q[0];
crz(-pi/2) q[2],q[0];
ccx q[4],q[6],q[0];
crz(-pi/2) q[2],q[0];
cry(-pi/2) q[1],q[0];
crz(-pi/2) q[1],q[0];
crz(pi) q[2],q[0];
cry(pi/2) q[2],q[0];
ccx q[4],q[6],q[0];
cry(-pi/2) q[2],q[0];
crz(-pi/2) q[2],q[0];
ccx q[4],q[6],q[0];
crz(-pi/2) q[2],q[0];
x q[6]; #X module
crz(-pi/2) q[1],q[0];
x q[4]; #X module
cx q[8],q[5];
rx(pi) q[0];
cx q[8],q[3];
x q[0];
ccx q[3],q[8],q[4];
ccx q[4],q[5],q[6];
cx q[6],q[7];
ccx q[4],q[5],q[6];
cx q[4], q[5];
ccx q[3],q[8],q[4];
cx q[8],q[3];
swap q[5],q[7];
swap q[3],q[8];
h q[3];
cp(-pi/2) q[3],q[5];
h q[5];
```

```
cp(-pi/4) q[3],q[7];
cp(-pi/2) q[5],q[7];
h q[7];
cp(-pi/8) q[3],q[8];
cp(-pi/4) q[5],q[8];
cp(-pi/2) q[7],q[8];
h q[8];
cx q[8],q[3];
ccx q[3],q[8],q[4];
cx q[4], q[5];
ccx q[4],q[5],q[6];
cx q[6],q[7];
ccx q[4],q[5],q[6];
ccx q[3],q[8],q[4];
cx q[8],q[3];
cx q[8],q[5];
cx q[8],q[7];
h q[8];
x q[8];
measure q[7] -> c[3];
measure q[5] \rightarrow c[2];
measure q[3] \rightarrow c[1];
measure q[0] \rightarrow c[0];
```

The Qiskit code for ideal circuit simulation without noise

preparation for execution
from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute,Aer
from qiskit.visualization import plot_histogram
from numpy import pi

#Put Qiskit circuit code for n=2 or n=3

#execute the circuit
backend=Aer.get_backend('qasm_simulator')

backend=Aer.get_backend('qasm_simulator job=execute(circuit,backend,shots=16384) result = job.result() counts = result.get_counts(circuit) plot_histogram(counts).show()

#Put Qiskit circuit code for n=2

The *n*=2 Qiskit code for device backend noise model mimicked IBMQ_SANTIAGO

#preparation for execution from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute from qiskit.providers.aer import QasmSimulator from qiskit.visualization import plot_histogram from qiskit.test.mock import FakeSantiago from numpy import pi #execute the circuit
device_backend = FakeSantiago()
Santiago_simulator=QasmSimulator.from_backend(device_backend)
result=execute(circuit,Santiago_simulator).result()
counts =result.get_counts(circuit)
plot histogram(counts).show()

The n=2 Qiskit code for executing circuit on real quantum computer IBMQ SANTIAGO

#preparation for execution
from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute
from qiskit.visualization import plot_histogram
from numpy import pi
from qiskit import IBMQ
provider = IBMQ.load_account()
backend=provider.get_backend('ibmq_santiago')

#execute the circuit
job=execute(circuit,backend,shots=8192)
result = job.result()
counts = result.get_counts(circuit)
plot histogram(counts).show()

#Put Qiskit circuit code for n=2

The Qiskit code of adding four common kinds of noise

#preparation for execution from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute,Aer from qiskit.visualization import plot_histogram from numpy import pi

#Put Qiskit circuit code for n=2 or n=3

#execute the circuit backend=Aer.get_backend('qasm_simulator') job=execute(circuit,backend,shots=16384) result = job.result() counts = result.get_counts(circuit) plot_histogram(counts).show()

The n=2 Qiskit code for device backend noise model mimicked IBMQ SANTIAGO

#preparation for execution from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute from qiskit.providers.aer import QasmSimulator from qiskit.visualization import plot_histogram from qiskit.test.mock import FakeSantiago from numpy import pi

#Put Qiskit circuit code for n=2

#execute the circuit device backend = FakeSantiago() Santiago simulator=QasmSimulator.from backend(device backend) result=execute(circuit,Santiago simulator).result() counts =result.get counts(circuit) plot_histogram(counts).show()

The n=2 Qiskit code for executing circuit on real quantum computer IBMQ SANTIAGO

#preparation for execution from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute from qiskit.visualization import plot histogram from numpy import pi from qiskit import IBMQ provider = IBMQ.load account() backend=provider.get backend('ibmq santiago')

#Put Qiskit circuit code for n=2

#execute the circuit job=execute(circuit,backend,shots=8192) result = job.result() counts = result.get counts(circuit) plot histogram(counts).show()

#put selected noise on gates of the circuit error gatec = error.tensor(error)

The Qiskit code of adding four common kinds of noise

#preparation for execution

from qiskit import QuantumRegister, ClassicalRegister, QuantumCircuit, execute from qiskit.providers.aer import QasmSimulator from giskit.providers.aer.noise import NoiseModel from qiskit.providers.aer.noise import QuantumError from qiskit.providers.aer.noise import amplitude damping error, phase damping error, depolarizing error, pauli error from qiskit.visualization import plot histogram from numpy import pi #Put Qiskit circuit code for n=2 or n=3#select the type of noise to add and set the noise intensity p = 1e-04*pow(700,1)#set intensity noise error=amplitude damping error(p) error = phase damping error(p) error = depolarizing error(p, 1)error = pauli error([('X',p error), ('I', 1 - p error)])

```
error_gateccx=error.tensor(error).tensor(error)
noise_model = NoiseModel()
noise_model.add_all_qubit_quantum_error(error, ["h"])
noise_model.add_all_qubit_quantum_error(error, ["x"])
noise_model.add_all_qubit_quantum_error(error_gatec, ["cx"])
noise_model.add_all_qubit_quantum_error(error_gatec, ["cry"])
noise_model.add_all_qubit_quantum_error(error_gatec, ["cul"])
noise_model.add_all_qubit_quantum_error(error_gatecc, ["ccx"])
noise_model.add_all_qubit_quantum_error(error, "measure")

#execute the noisy circuit
noise_simulator=QasmSimulator(noise_model=noise_model)
job = execute(circuit, noise_simulator)
noisy_result = job.result()
noisy_counts=noisy_result.get_counts(0)
plot_histogram(noisy_counts).show()
```

Appendix B Date

The detailed data on Figure 9 in the table below.

Table.1 The deviation for n=2 under four types of noise						
i	AD	PD	DE	BF		
1	0.0341	0.0206	0.0175	0.0161		
2	0.0177	0.0275	0.0350	0.0293		
3	0.0543	0.0396	0.0848	0.1001		
4	0.0696	0.0310	0.0697	0.1184		
5	0.1588	0.0196	0.1288	0.2381		
6	0.2701	0.0412	0.2591	0.3782		
7	0.4608	0.0606	0.4171	0.5390		
8	0.6660	0.1504	0.5901	0.6528		

0.2412

0.7074

0.7047

0.8118

Table.2 The deviation for n=3 under four types of noise						
i	AD	PD	DE	BF		
1	0.0745	0.0625	0.0310	0.0778		
2	0.1205	0.0498	0.1069	0.1580		
3	0.1363	0.0372	0.1471	0.2116		
4	0.2031	0.0842	0.2183	0.3565		
5	0.3497	0.0900	0.3481	0.5365		
6	0.4702	0.1781	0.4811	0.5875		
7	0.5709	0.1993	0.6183	0.6690		
8	0.6358	0.3257	0.5904	0.6736		
9	0.7105	0.4825	0.6418	0.6212		

Further details

- See IBM Quantum. https://quantum-computing.ibm.com/ (2021)
 See https://qiskit.org/documentation/getting_started.html