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Andrew M. Childs, Andrew M. Childs, Debbie Leung, Michael A. Nielsen ...+1 more authors

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Published on: 17 Mar 2005 - Physical Review A (American Physical Society)

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Unified derivations of measurement-based schemes for quantum computation

Andrew M. Childs,^{1,2,*} Debbie W. Leung,^{1,†} and Michael A. Nielsen^{1,3,‡}

¹*Institute for Quantum Information, California Institute of Technology, Pasadena, California 91125, USA*

²*Center for Theoretical Physics, Massachusetts Institute of Technology, Cambridge, Massachusetts 02139, USA*

³*School of Physical Sciences and School of Information Technology & Electrical Engineering, The University of Queensland, QLD 4072, Australia*

(Received 1 July 2004; published 17 March 2005)

We present unified, systematic derivations of schemes in the two known measurement-based models of quantum computation. The first model (introduced by Raussendorf and Briegel, [Phys. Rev. Lett. **86**, 5188 (2001)]) uses a fixed entangled state, adaptive measurements on single qubits, and feedforward of the measurement results. The second model (proposed by Nielsen, [Phys. Lett. A **308**, 96 (2003)] and further simplified by Leung, [Int. J. Quant. Inf. **2**, 33 (2004)]) uses adaptive two-qubit measurements that can be applied to arbitrary pairs of qubits, and feedforward of the measurement results. The underlying principle of our derivations is a variant of teleportation introduced by Zhou, Leung, and Chuang, [Phys. Rev. A **62**, 052316 (2000)]. Our derivations unify these two measurement-based models of quantum computation and provide significantly simpler schemes.

DOI: 10.1103/PhysRevA.71.032318

PACS number(s): 03.67.Lx, 03.65.Ta

I. INTRODUCTION

What physical resources are needed to simulate the evolution of an arbitrary physical system? In the context of information processing, the ability to perform *universal* computation is equivalent to the ability to simulate an arbitrary evolution—any computation is performed by evolving a computing machine, and conversely, a universal computing machine can be used to simulate the evolution of any system.

In the standard quantum circuit model of quantum computation [1–3], a quantum computation involves initializing quantum systems—typically *qubits* or two-level quantum systems—that are then acted on by a sequence of *quantum gates*, followed by some measurements. In this model, simple quantum gates (say, acting on one or two qubits at a time) can be used to build up an *arbitrary* unitary transformation. Nonunitary evolution such as dissipation can also be simulated in this model by introducing and later discarding ancilla qubits.

Since measurement is generally irreversible, until recently the conventional wisdom has held that the processing of quantum information should be kept coherent and measurements should be delayed until the final readout of computation results. A notable exception to this rule of thumb is quantum teleportation [4], in which a measurement by one party determines the correction that a remote party should apply to recover a quantum state. Another notable exception is the use of syndrome measurements in quantum error correction [5]. Syndrome measurements reveal the error that has occurred without measuring the encoded quantum state, thereby preserving its coherence. Indeed, there are many approaches to fault-tolerant quantum computation in which

measurements and simple quantum gates are used to implement other quantum gates that are difficult to apply directly [6–12].

Raussendorf and Briegel [13] overturned this conventional wisdom, showing that it is possible to perform universal quantum computation using a *sequence of single-qubit measurements alone*, acting on some fixed entangled state called a *cluster state* [14]. Once the cluster state is prepared, no further interactions are required, and the only aspect of the computation that must remain coherent is the storage of quantum information. More precisely, *any* quantum circuit up to depth d and breadth b may be simulated using a single, fixed cluster state of $O(bd)$ qubits. Each simulation of a quantum gate is successful up to an additional known Pauli error. Since the act of measuring the cluster state is irreversible, this model is referred to as the *one-way quantum computer* (1WQC) model.

After the 1WQC was introduced, a very different measurement-based scheme for quantum computation was introduced by one of us [15], following the line of thought developed in [16,9,11]. We will refer to this model as a *teleportation-based model of quantum computation* (TQC), since it is conceptually derived from teleportation. The TQC uses similar physical resources to the 1WQC: (multiple-qubit) measurements, quantum memory, and feedforward. The initial TQC scheme proposed in [15] uses four-qubit measurements. It also requires a nondeterministic number of steps to perform each quantum gate. Simpler TQC schemes were later proposed [17–19], with the simplest using only two-qubit measurements and performing each gate deterministically (up to a known Pauli error).

The TQC is easy to understand since it is similar to the standard model of quantum computation. In comparison, the conceptual basis for the 1WQC is less clear. The prescriptions given in [13,20] for using a 1WQC can be easily verified, but there is no clear underlying principle. This makes it nontrivial to modify or optimize the existing 1WQC

*Email address: amchilds@cs.caltech.edu

†Email address: wcleung@cs.caltech.edu

‡Email address: nielsen@physics.uq.edu.au

schemes. It is also unclear what makes the cluster state a good substrate for quantum computation, and more generally, what makes a good or bad substrate. Finally, the 1WQC formalism is heavily based on the stabilizer language developed by Gottesman [21]. Although this language is powerful, it is also rather specialized, being limited primarily to the analysis of situations in which operations from a special set—the Clifford group—are being applied. Furthermore, the connection between the stabilizer language and the corresponding picture in terms of state vectors is not always straightforward.

On the other hand, the 1WQC has important advantages over the TQC. First, no quantum interactions are required after the initial preparation of the cluster state. Second, the cluster state is independent of the computation to be performed, except for its breadth and depth: arbitrary interactions can be extracted from the fixed cluster state. Third, there may be physical systems in which a cluster state offers experimental advantages over more conventional approaches [14,22].

Our initial goal in undertaking the research reported here was to identify simple underlying principles for the 1WQC and to systematically derive schemes similar to the proposed 1WQC. We eventually found such a systematic derivation using teleportation as an underlying principle [23], in accord with the conjecture that the 1WQC and the TQC are closely related. Our improved conceptual understanding of the 1WQC proved valuable, for we subsequently found much simpler 1WQC-like schemes [24] by choosing a simpler underlying principle, known as “one-bit teleportation” [10]. Such simplification is reminiscent of the work in [10], which simplifies the systematic fault-tolerant gate construction proposed in [9]. We then realized that one-bit teleportation also simplifies schemes in the TQC model [25].

We have therefore unified the 1WQC and the TQC models and obtained simplified measurement-based quantum computation schemes. The 1WQC schemes we derive combine the conceptual simplicity of the TQC with the practical advantages of the 1WQC. We have also identified one-bit teleportation as a single principle underlying both approaches to measurement-based quantum computation.

During the course of our investigation and preparation of the manuscript, several related results have been reported. A different explanation of the 1WQC model in terms of valence bond solids was reported by Verstraete and Cirac [26]. Whereas our 1WQC-like schemes differ from the original schemes of [13,20], an exact explanation of the latter in terms of teleportation was given by Aliferis and one of us [27]. A partial explanation of the 1WQC model in terms of one-bit teleportation was reported very recently by Jorrand and Perdrix [28], while schemes similar to our simplified TQC schemes were independently reported by Perdrix [29]. One of us has combined the 1WQC model with linear optics [22]. Several results announced after our initial posting may also be of interest. These include a model of measurement-based universal quantum Turing machines [30], a more efficient method to combine the 1WQC with linear optics [31], and a new fault-tolerance study in the 1WQC [32].

The structure of the paper is as follows. We begin in Sec. II by briefly reviewing the circuit model, introducing a no-

tion of circuit simulation with Pauli errors, and describing the TQC and 1WQC models of quantum computation in more detail. We emphasize some conceptual ideas that may be useful elsewhere. Section III reviews one-bit teleportation and presents useful techniques obtained from it. The techniques are used to derive a simplified TQC scheme in Sec. IV A. Section IV C explains how one-bit teleportation can be used to derive a scheme that simulates arbitrary circuits using a circuit-dependent entangled initial state, single-qubit measurements, and feedforward. Section IV D explains several techniques to remove the dependence of the initial state on the circuit being simulated (except for its breadth and depth). We conclude the paper in Sec. V. Our initial systematic derivation of 1WQC-like schemes based on teleportation can be found in [23].

II. THE QUANTUM CIRCUIT MODEL, THE TQC, AND THE 1WQC

In this section, we summarize the circuit model of quantum computation as a way of introducing the notation used throughout the paper. We describe a notion of circuit simulation that is crucial to our discussion. Then, we outline the main features of the TQC and the 1WQC that motivate our derivation. Since we derive simplified TQC and 1WQC schemes, a full summary of the existing schemes will be omitted. Finally, we describe a diagrammatic representation of the 1WQC, which we call the *substrate representation*.

A. The quantum circuit model

Any unitary evolution can be built from simple quantum gates (say, acting on one or two qubits at a time). A circuit diagram represents a partially ordered set of unitary evolutions and measurements. The input states and measurement outcomes may be included. In a circuit diagram, time runs from left to right. Each horizontal line represents quantum information propagating forward in time, or equivalently, quantum storage. Often, each line represents one qubit of quantum information. A unitary gate is represented by a box on the line(s), and a symbol for the gate is written inside the box. Thus, the circuit symbol for a single-qubit gate U is given by

$$\text{---} \boxed{U} \text{---} \tag{1}$$

In the basis $\{|0\rangle, |1\rangle\}$, the matrix representations of some useful single-qubit gates are given by

$$X_\theta = e^{-i\theta X}, \quad Z_\theta = e^{-i\theta Z}, \quad H = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}, \tag{2}$$

where I, X, Y, Z are used to represent the Pauli operators

$$I = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, \quad X = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \\ Y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad Z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}. \tag{3}$$

is not a hindrance to the computation, provided subsequent measurements are adapted accordingly. Since our schemes share the same feature, it is useful to introduce some conventions that simplify later discussions of simulation.

A quantum circuit consists of ancilla preparations, gates, and measurements endowed with a partial time-ordering. We can group together operations that can be performed in parallel in a time step, although the grouping may not be unique. The minimal number of time steps is called the *logical depth* of the circuit. For each grouping of \mathcal{C} , an input state $|\psi_0\rangle$ specifies a sequence of quantum states $\{|\psi_i\rangle\}$ where $|\psi_i\rangle$ is the quantum state at the end of the i th time step of the computation. We say that a circuit \mathcal{C}' *simulates* \mathcal{C} with Pauli errors if there is a grouping of \mathcal{C} so that, for any input state $|\psi_0\rangle$ and any given Pauli error P , applying \mathcal{C}' on the input $P|\psi_0\rangle$ produces a sequence of states $P_i|\psi_i\rangle$ where P_i are known Pauli errors. These Pauli errors redefine the intermediate states and the final measurement outcomes, but do not affect the intended computation. From now on, we will simply say that \mathcal{C}' *simulates* \mathcal{C} to mean that \mathcal{C}' simulates \mathcal{C} with Pauli errors. As a simple example, a unitary gate simulates the identity if and only if it is in the Pauli group. As a second example, a unitary gate simulates itself if and only if it is in the Clifford group.

Our notion of simulation is defined to ensure an important property—circuit simulation is preserved under the composition of circuits. Therefore, to simulate a circuit, it suffices to compose simulations of individual circuit elements. Universality can then be proved by showing how to efficiently simulate all possible circuit elements required for universal quantum computation. Furthermore, even though a circuit element may act on part of an entangled state, it is sufficient to verify the simulation of a circuit element for all possible pure state inputs.

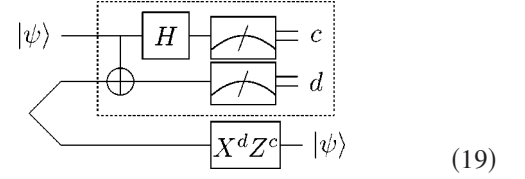
In the context of measurement-based quantum computation, initial (or ancilla) state preparation and read out of computation results can be simulated as follows. We will only need initial states that can be prepared by a simple measurement, up to known Pauli errors. For example, measuring a single qubit in the computational basis with outcome $d \in \{0, 1\}$ results in the state $X^d|0\rangle$. Similarly, a measurement of two qubits in the Bell basis produces a singlet state up to a known Pauli error. Throughout the paper, we suppress the known Pauli errors in the ancillas whenever their effects are straightforward, so as to keep the discussion and the simulation circuits simple. We also omit physically irrelevant global phase factors that arise from the composition of Pauli errors. We restrict our attention to measurements that are deterministically affected by known Pauli errors, so that the actual outcomes and the knowledge of the Pauli errors can be used to determine the Pauli-error-free measurement outcomes. For example, the result of a measurement in the computational basis is simply flipped by an X error and unaffected by a Z error. Now, it suffices to focus on simulating a universal set of unitary gates in the measurement-based model of quantum computation.

C. The TQC model

In this section, we review some elements of the TQC. Our review follows the simplified approach of [19], which allows

the computation to proceed with a deterministic number of steps. The universality of the TQC model follows from the ability to simulate any single-qubit gate U and a two-qubit gate such as $\Lambda(X)$.

The crucial ingredient of the simulation is teleportation [4], which transmits a qubit $|\psi\rangle$ using the following circuit:



When two qubits are connected on the left side of a circuit diagram, as the bottom two qubits are in Eq. (19), they denote a two-qubit maximally entangled state $|\Phi_{00}\rangle = (|00\rangle + |11\rangle)/\sqrt{2}$. The dashed box performs a *Bell measurement* along the basis

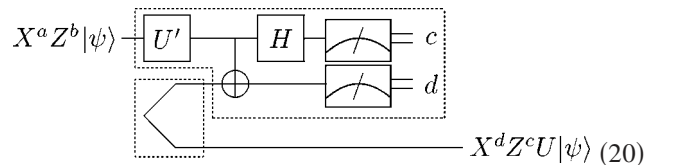
$$|\Phi_{00}\rangle = \frac{1}{\sqrt{2}}(|00\rangle + |11\rangle), \quad |\Phi_{10}\rangle = \frac{1}{\sqrt{2}}(|00\rangle - |11\rangle),$$

$$|\Phi_{01}\rangle = \frac{1}{\sqrt{2}}(|01\rangle + |10\rangle), \quad |\Phi_{11}\rangle = \frac{1}{\sqrt{2}}(|01\rangle - |10\rangle).$$

The teleportation circuit can be verified by rewriting the initial state $|\psi\rangle|\Phi_{00}\rangle$ as $\frac{1}{2}\sum_{c,d}|\Phi_{cd}\rangle \otimes (Z^c X^d|\psi\rangle)$.

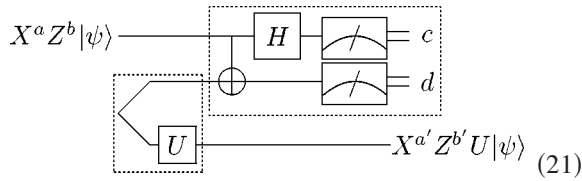
The teleportation circuit, Eq. (19), simulates the identity gate (in the sense described in Sec. II B). In fact, it does so even when the postmeasurement correction $X^d Z^c$ is omitted: for any input state $X^a Z^b|\psi\rangle$, the output from the teleportation circuit without correction is simply $X^{a+d} Z^{b+c}|\psi\rangle$ (up to a global phase, as is our convention throughout the paper). From now on, Pauli corrections are always omitted in what we call teleportation.

Suppose we want to simulate a single-qubit gate U . Consider a circuit in which we apply the gate $U' = UZ^b X^a$ to the input state $X^a Z^b|\psi\rangle$ and perform teleportation on the resulting state $U|\psi\rangle$. Following Eq. (19), the output quantum state of the circuit is $X^d Z^c U|\psi\rangle$. In other words, the following circuit simulates U :



The circuit in Eq. (20) can be divided into subcircuits, each of which can be simulated by a single two-qubit measurement. The first subcircuit (the dashed box in the lower left corner) is just the preparation of the state $|\Phi_{00}\rangle$. It can be simulated by a single Bell measurement, up to a Pauli error. The rest of the circuit, $(U' \otimes I)$ followed by a Bell measurement, is just a two-qubit measurement along a rotated Bell basis $\{(U'^{\dagger} \otimes I)|\Phi_{cd}\rangle\}$. Composing these two measurements provides a simulation of Eq. (20), and thus a simulation of U , in the TQC model.

An alternative simulation is applicable to gates in the Clifford group [9]:



The above is simply a simulation of the identity by teleportation followed by a simulation of U by itself. More specifically, the above circuit teleports the input state $X^a Z^b |\psi\rangle$ to produce the state $X^{a+d} Z^{b+c} |\psi\rangle$, and then applies the gate U . The output state is thus $U X^{a+d} Z^{b+c} |\psi\rangle$. Using the fact U is in the Clifford group, the output state can be rewritten as $X^{a'} Z^{b'} U |\psi\rangle$ for known a', b' .

The simulation of two-qubit gates is similar to that depicted in Eqs. (20) and (21). We omit the details of the existing simulation schemes, and we refer the interested reader to [19]. Instead, we will present a simplified simulation in Sec. IV A.

Comparing Eqs. (20) and (21), the latter simulation has a simpler teleportation measurement but a more complicated initial state $(I \otimes U) |\Phi_{00}\rangle$. This tradeoff is useful in the simulation of two-qubit gates in the Clifford group.

Note that in the TQC model we assume the ability to apply measurements to any subset of the qubits, without worrying about the respective locations of those qubits, just as in the circuit model. The TQC is simply a method for implementing each gate in the circuit model by a sequence of measurements.

D. The one-way quantum computer model

Since we will present a simple systematic derivation of 1WQC-like schemes, we omit the details of the existing 1WQC schemes, and refer interested readers to the ingenious schemes discussed in [13,20,33,34]. Instead, we focus on features of the 1WQC that differ from the TQC model.

The 1WQC is based on an input state known as the *cluster state* [14]. The cluster state consists of a two-dimensional square lattice of qubits. To simulate a computation of n qubits with logical depth m , the lattice is chosen to be of size $O(n) \times O(m)$. One way of defining the cluster state is as the result of a two-stage preparation procedure: (i) prepare all lattice sites in the state $|+\rangle = (|0\rangle + |1\rangle) / \sqrt{2}$, and (ii) apply $\Lambda(Z)$ between each pair of qubits that are adjacent in the lattice. Since the $\Lambda(Z)$ operations all commute, it does not matter in what order they are applied. Note that this prescription is merely a convenient way of defining the cluster state, and there is no need to actually prepare it by following these steps.

The cluster states naturally generalize to *graph states*, for which an arbitrary graph defines the adjacency relation [34–36]. We will use a feature of any graph state called the *deletion principle*. When one of the qubits is measured in the computational basis, the unmeasured qubits will be left in a different graph state (up to known Pauli errors) corresponding to the graph obtained by deleting the measured qubit from the original graph.

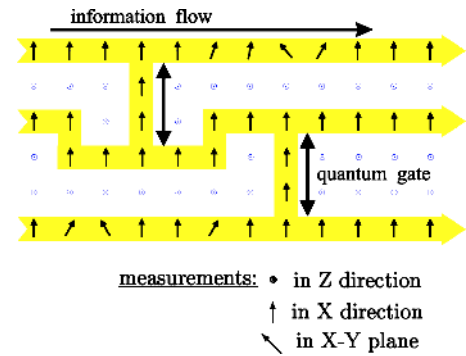


FIG. 1. A schematic diagram of a 1WQC computation. Figure courtesy H. J. Briegel and R. Raussendorf.

In the 1WQC model, once the cluster state is prepared, an *arbitrary* circuit \mathcal{C} can be simulated using only single-qubit measurements. The first step is to “imprint” the circuit \mathcal{C} onto the cluster state by deleting qubits from the lattice to obtain some graph state that depends on \mathcal{C} . Roughly speaking, this graph consists of horizontal lines of vertices, each line corresponding to a qubit in \mathcal{C} , and vertical edges connecting neighboring lines corresponding to interactions between the simulated qubits. The qubits in the graph are then measured from the left to the right. Each measurement propagates quantum information from the measured qubit to its unmeasured right neighbor. In general, the measurement basis will depend on previous measurement outcomes. Each region of the graph is used to simulate a circuit element. Qubits entering a region from the left boundary carry the input state of the circuit element, and qubits exiting at the right boundary carry the output. When composing element-wise simulation in the 1WQC model, the output registers of one region have to match the input registers of the next region. A schematic diagram for a computation in the 1WQC model is given in Fig. 1.

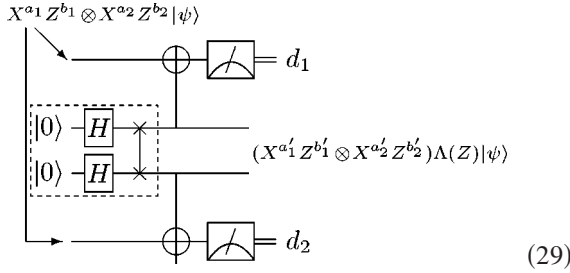
Simulation of a circuit using the 1WQC is discussed in detail in [20]. The precise protocols and their verification are beyond the scope of this paper. Interpretations in terms of valence bond solids [26] and teleportation [27,37] have been reported recently. In this paper, the derivation is based on a different underlying principle and the schemes are significantly simpler than those discussed in [20].

E. The substrate representation

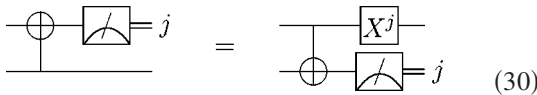
Circuit representations of a computation in the 1WQC model turn out to be rather unwieldy. Instead, we prefer to use a more concise “substrate representation” in which we label each vertex in the graph representing the cluster with the measurement to be performed. The measurement bases may depend on previous measurement outcomes, and this dependency should be indicated in the labeling. Note that the interdependence of the measurement outcomes and measurement bases specifies a partial order in which the qubits must be measured. An example of the substrate representation is Eq. (44), in Sec. IV C 2, which simulates the circuit in Eq. (39).

Thus in Eq. (28), $a'_1 = a_1 + d_1$, $b'_1 = b_1 + a_2 + d_2$, $a'_2 = a_2 + d_2$, and $b'_2 = b_2 + a_1 + d_1$.

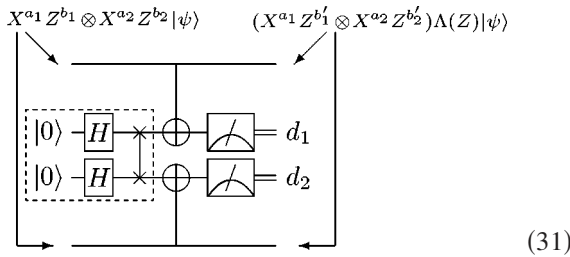
We can derive useful simulation circuits from Eq. (28). Suppose we commute $\Lambda(Z)$ to the left of the controlled-NOTs, and reorder the qubits so that the second qubit from the top becomes the last:



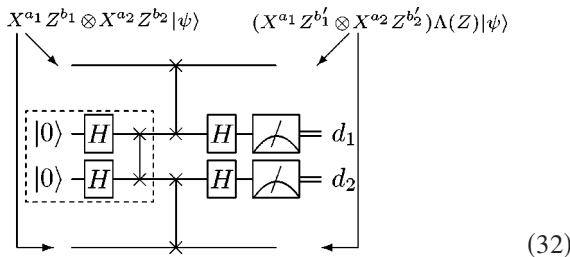
Furthermore, for the same input state, the following circuits produce the same outcomes and corresponding post-measurement states:



Thus Eq. (29) implies the following:

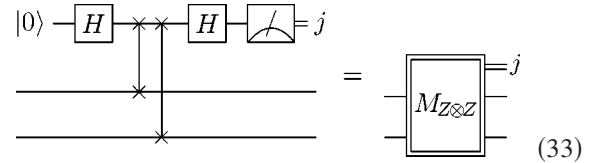


where, according to Eq. (30), the output X errors in Eq. (31) are obtained by adding d_1, d_2 to a'_1, a'_2 defined in Eq. (29). The results are simply a_1, a_2 . Finally, rewrite both controlled-NOTs using Eq. (7), and note that the state in the dashed box in Eq. (31) is stabilized by $H \otimes H$, giving a “remote $\Lambda(Z)$ ” construction:

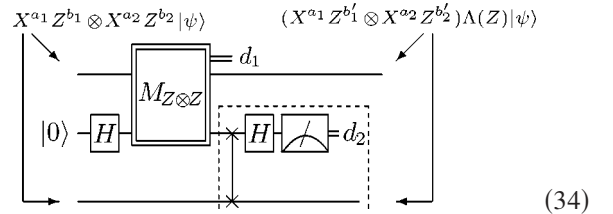


If we perform a remote controlled-NOT by performing H before and after the remote $\Lambda(Z)$ according to Eq. (7), we obtain the well-known remote $\Lambda(X)$ circuit due to Gottesman [38]. The current derivation is based only on composing the identity simulation (by one-bit teleportation) with self-simulation of the desired gate, and is different from the derivation in [10].

Our last simulation circuit for $\Lambda(Z)$ uses the standard (and easily-verified) result that the following circuit implements $M_{Z \otimes Z}$ on the two input qubits:



We can apply Eq. (33) to Eq. (32), and identify the operations involving the second qubit (from the top) in Eq. (32) as a two-qubit measurement on the first and third qubits. This gives a simulation circuit for $\Lambda(Z)$:



The operations in the dashed box can be implemented by $M_{X \otimes Z}$ followed by M_Z on the first qubit [see Eq. (18) in Sec. II A]. With this argument we have rederived Gottesman’s remote controlled-NOT using a single-qubit ancilla and two two-qubit measurements [39], and shown that it is easily understood as a consequence of one-bit teleportation and the simple circuit identities (18) and (30).

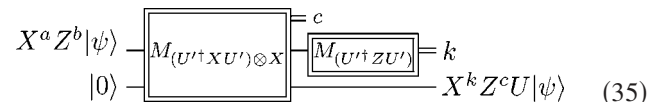
IV. MEASUREMENT-BASED UNIVERSAL QUANTUM COMPUTATION SCHEMES

In this section we derive simple variants of both the TQC and 1WQC models of computation using the principles described in earlier sections. Following the discussion in Sec. II B, it suffices to show how to perform a universal set of gates in each measurement-based model of quantum computation. We will first see that the simulation circuits derived in the previous section immediately give a universal scheme in the TQC model. This scheme is much simpler than those based on teleportation. (A similar simplified scheme was reported independently in [29].) Then we discuss a method to further reduce the required resources in the TQC model by identifying and simulating certain subunits of a circuit. We then turn to the 1WQC model and present a systematic derivation of universal quantum computation schemes using primitives discussed in the previous section.

A. Derivation of simplified TQC schemes

1. Universality

Consider the universal set consisting of the single-qubit gates and $\Lambda(Z)$. A single-qubit gate can be performed in the TQC model using either Eq. (24) or Eq. (25)—the operations in the dashed boxes are of the form of Eq. (17), with V in the Clifford group and $V^\dagger Z V$ in the Pauli group. Thus Eq. (18) without the Pauli correction $V^\dagger Z V$ can be used to implement the dashed boxes in the TQC model. More concretely, Eqs. (24) and (25) imply the following simulation circuits:



$$\begin{array}{c}
 X^a Z^b |\psi\rangle \\
 |+\rangle
 \end{array}
 \begin{array}{c}
 \boxed{M_{(U^\dagger Z U') \otimes Z}} \\
 \boxed{M_{(U^\dagger X U')}}
 \end{array}
 \begin{array}{c}
 =d \\
 =k
 \end{array}
 \begin{array}{c}
 \\
 Z^k X^d U |\psi\rangle
 \end{array}
 \quad (36)$$

In Eqs. (35) and (36) the ancillas can be prepared up to known Pauli errors that *commute* with the subsequent two-qubit measurements. Simulation circuits for $\Lambda(Z)$ can be obtained from Eqs. (32) and (34):

$$\begin{array}{c}
 X^{a_1} Z^{b_1} \otimes X^{a_2} Z^{b_2} |\psi\rangle \\
 (X^{a_1} Z^{b_1'} \otimes X^{a_2} Z^{b_2'}) \Lambda(Z) |+\rangle
 \end{array}
 \begin{array}{c}
 \boxed{M_{Z \otimes X}} \\
 \boxed{M_{X \otimes Z}}
 \end{array}
 \begin{array}{c}
 =d_1 \\
 =d_2
 \end{array}
 \begin{array}{c}
 \boxed{M_Z} \\
 \boxed{M_Z}
 \end{array}
 \begin{array}{c}
 =k_1 \\
 =k_2
 \end{array}
 \quad (37)$$

$$\begin{array}{c}
 X^{a_1} Z^{b_1} \otimes X^{a_2} Z^{b_2} |\psi\rangle \\
 |+\rangle
 \end{array}
 \begin{array}{c}
 \boxed{M_{Z \otimes Z}} \\
 \boxed{M_{X \otimes Z}}
 \end{array}
 \begin{array}{c}
 =d_1 \\
 =d_2
 \end{array}
 \begin{array}{c}
 \boxed{M_Z} \\
 \boxed{M_Z}
 \end{array}
 \begin{array}{c}
 =k_2 \\
 =k_2
 \end{array}
 \quad (38)$$

In the above, k_1 should be added to the value of b_1' from Eq. (32), and k_2 should be added to b_2' . The state $\Lambda(Z)|+\rangle|+\rangle$ in Eq. (37) can be prepared by a two-qubit measurement. In both Eqs. (37) and (38), the ancillas can be prepared up to known Z errors, which have no effect other than flipping the measurement outcomes of subsequent $M_{X \otimes Z}$ and $M_{Z \otimes X}$. The simulation (37) uses two ancillary qubits, three two-qubit measurements, and two single-qubit measurements, and its logical depth is 3. The simulation (38) uses one ancillary qubit, two two-qubit measurements, and two single-qubit measurements, but its logical depth is 4.

2. Reduced-cost combined pseudosimulations

In the TQC model, how many single- and two-qubit measurements are required to simulate a circuit \mathcal{C} consisting of single-qubit gates, m $\Lambda(Z)$ gates, and n final single-qubit measurements on the n computation qubits? We can do better than the method described above by combining some of the gates in the circuit into larger subunits, and simulating the subunits directly in the TQC model. In particular, without loss of generality, there are single-qubit gates U_i, V_i for $i = 1, \dots, m$, such that \mathcal{C} only consists of m “composite” gates $(U_i^\dagger \otimes V_i^\dagger) \Lambda(Z) (U_i \otimes V_i)$ applied in order, followed by single-qubit measurements.

Starting from Eqs. (32) and (34) and using Eq. (18), analogues of Eqs. (37) and (38) can be used to attempt the simulation of $W = (U^\dagger \otimes V^\dagger) \Lambda(Z) (U \otimes V)$ for any single-qubit gates U and V . These analogues of Eqs. (37) and (38) simply have $M_{U^\dagger Z U' \otimes X}$, $M_{X \otimes V^\dagger Z V'}$, and $M_{U^\dagger Z U' \otimes Z}$ in place of $M_{Z \otimes X}$, $M_{X \otimes Z}$, and $M_{Z \otimes Z}$ respectively. We call these analogues “pseudosimulations,” because W is simulated up to

possible left multiplications of $U^\dagger Z U$ and $V^\dagger Z V$, which can easily be compensated for in the next pseudosimulation involving the same qubit.

The reduced-cost combined pseudosimulations use measurements of complexity comparable to those required in Eqs. (35) and (36). Altogether, a computation using m $\Lambda(Z)$ gates and n computation qubits can be pseudosimulated in TQC using m ancillary qubits, $2m$ two-qubit measurements, and $2m+n$ single-qubit measurements. In comparison, a full simulation [say, using Eqs. (35) and (38)] requires $3m$ ancillary qubits, $4m$ two-qubit measurements, and $6m+n$ single-qubit measurements.

B. Derivation of schemes similar to the 1WQC starting from the TQC

In this and subsequent subsections, we present our derivation of 1WQC-like schemes using one-bit teleportation as the underlying principle, preserving the conceptual simplicity of the TQC. The derivation is motivated by the differences between the TQC and 1WQC models. The TQC and 1WQC models differ in three main respects.

(a) The TQC model is similar to the circuit model in that no action is required on a qubit unless a nonidentity gate is applied. In contrast, in the 1WQC model, it is necessary to keep measuring qubits simply to propagate quantum information forward on the lattice.

(b) In the TQC model, interactions are effected by multi-qubit measurements. In contrast, no interactions are used in the 1WQC model after the initial preparation of the cluster state. In some sense, all interactions are built into the initial state before the computation begins.

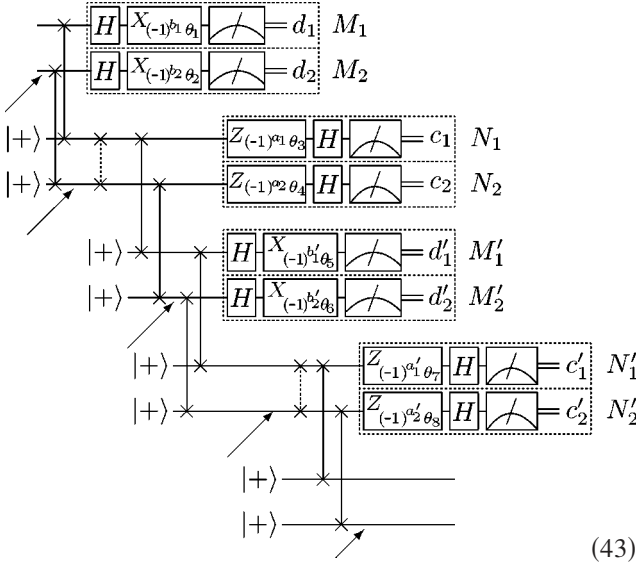
(c) In the 1WQC model, a circuit \mathcal{C} can be simulated using a \mathcal{C} -dependent graph state, which can in turn be produced from a \mathcal{C} -independent cluster state. Thus, the built-in interactions in the 1WQC model can be made independent of \mathcal{C} . In contrast, a TQC simulation has a one-to-one correspondence with \mathcal{C} .

These differences suggest a strategy to derive 1WQC-like schemes using the principles of the TQC model: every gate is performed by simulation circuits based on teleportation or one-bit teleportation [such as Eqs. (20), (21), (26), (27), and (32)]. Suppose the goal is to simulate a circuit \mathcal{C} with n qubits and m time steps.

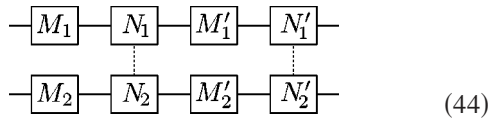
(a) Each gate in \mathcal{C} will be simulated by circuits like Eqs. (26), (27), and (32). Furthermore, in each time step, identity gates will be explicitly simulated on qubits that are not being acted on. Thus, each qubit will be “teleported” in each step. Matching the output of one gate simulation to the input of the next, we obtain a circuit \mathcal{C}' that “teleports” each of the n qubits m times, with the desired gates performed along the way. \mathcal{C}' contains initial $|0\rangle$ states, one- or two-qubit gates, and single-qubit measurements. Note that in this circuit we do not interpret a two-qubit gate followed by a single-qubit measurement as an incomplete two-qubit measurement, as we did in the TQC. The reason is that in the next step we will build the two-qubit gates into the initial state, leading to an equivalent circuit containing only single-qubit measurements.

Each arrow in Eq. (42) indicates where the output of a certain teleportation step matches the input of the subsequent teleportation. The values of a'_i and b'_i can be read from Eqs. (40) and (41). The circuit of Eq. (42) generalizes easily to n qubits with multiple optional $\Lambda(Z)$ gates.

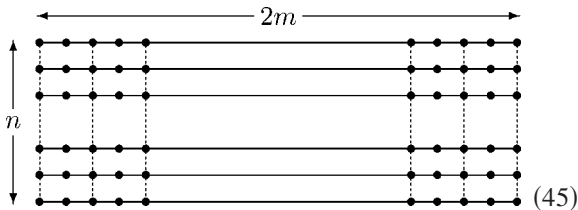
The simulation (42) can be simplified by (i) rewriting $\Lambda(X)$ as $(I \otimes H)\Lambda(Z)(I \otimes H)$, (ii) canceling out consecutive Hadamard gates (since $H^2=I$), (iii) rewriting $H|0\rangle$ as $|+\rangle$, and (iv) absorbing H before a single-qubit measurement as part of the measurement. We thus obtain a simpler simulation scheme for Eq. (39):



We can view the operations in the dashed boxes as single-qubit measurements, and the rest of the circuit as an initial state $|\psi_C\rangle$. Note that $|\psi_C\rangle$ is a graph state, as defined in Sec. II D. The circuit \mathcal{C} determines whether each $\Lambda(Z)^k=I$ or $\Lambda(Z)$ in $|\psi_C\rangle$. The substrate representation of Eq. (43) is



where the measurements are as specified in Eq. (43). The initial graph state for an n -qubit circuit \mathcal{C} with m time steps can be chosen to be



Equations (44) [with measurements specified in Eq. (43)] and (45) form a complete recipe for the simulation of a known quantum circuit by the 1WQC_{TG} scheme. This approach is generally more efficient than the circuit-independent (1WQC_T) schemes described in the following section. Indeed, the 1WQC_{TG} approach is useful in its own

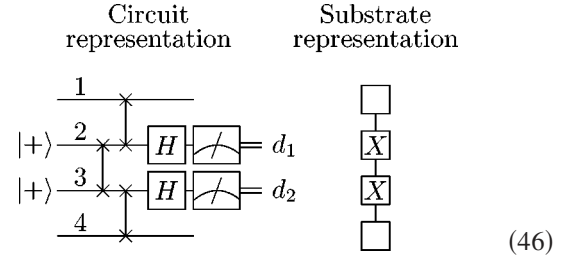
right when the simulated circuit is known at the time of initial state preparation.

D. Derivation of schemes using a universal family of initial states (1WQC_T)

We now present methods for performing an optional $\Lambda(Z)$. Composing the optional $\Lambda(Z)$ simulation with the 1WQC_{TG} scheme described in the previous section, we obtain various 1WQC_T schemes.

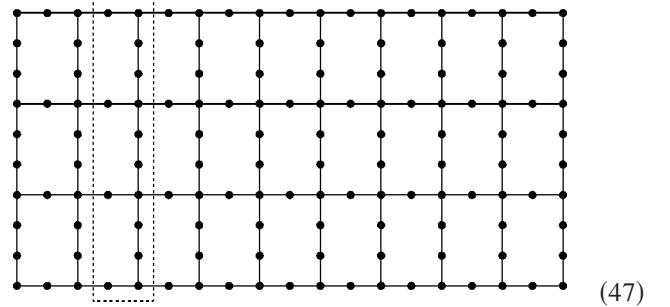
1. The remote $\Lambda(Z)$ approach (I)

Consider the circuit (32) that simulates $\Lambda(Z)$:

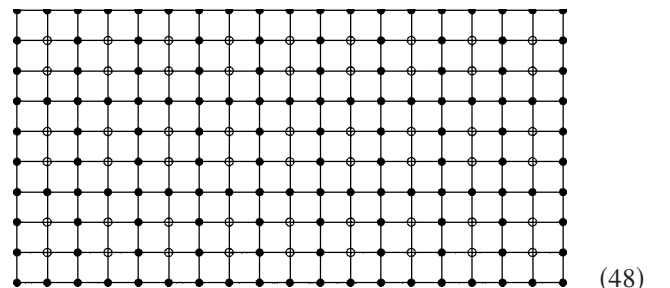


Note that we have explicitly labeled all the qubits. The circuit in Eq. (46) starts with a graph state, and applies the gate $Z^{d_2} \otimes Z^{d_1} \Lambda(Z)$ to qubits 1 and 4. On the other hand, the deletion principle from [13] (see Sec. II D) shows that, if the H gates on qubits 2 and 3 are simply omitted, and those qubits are measured along $\{|0\rangle, |1\rangle\}$, then qubits 1 and 4 are disentangled, and an identity gate is simulated instead.

Substituting the initial graph state of Eq. (46) for the optional $\Lambda(Z)$ in Eq. (45), the initial graph state for our first 1WQC_T scheme is given by



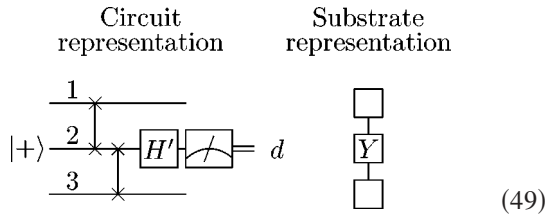
This graph state can be used to simulate a 4-qubit circuit for 9 cycles of (i) arbitrary X rotations and (ii) arbitrary Z rotations and optional nearest-neighbor $\Lambda(Z)$. The region corresponding to the simulation of one cycle is marked by a dashed box. The above state can in turn be obtained by “deleting” qubits denoted by empty circles (by measuring them in the computational basis) in the following cluster state:



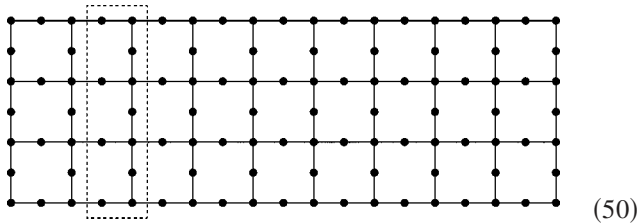
This cluster state-based simulation requires six physical qubits per logical qubit per cycle.

2. The remote $\Lambda(Z)$ approach (II)

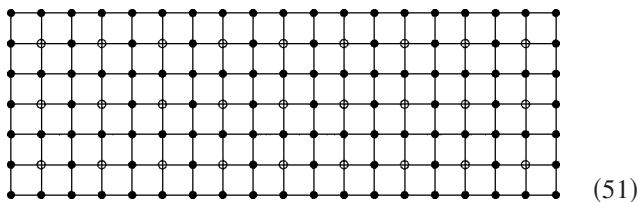
The remote $\Lambda(Z)$ described in the previous section requires two ancilla qubits per $\Lambda(Z)$. We can save resources by relaxing the exact simulation condition, and use the following circuit:



where $H' = (Z+Y)/\sqrt{2}$ and H' followed by M_Z is simply M_Y . Let R_d be the Z -rotation $Z_{(-1)^{d+1}\pi/2}$. The above circuit implements the gate $(R_d \otimes R_d)\Lambda(Z)$, and yields a $1WQC_T$ scheme from the $1WQC_{TG}$ model in Sec. IV C 2, because arbitrary Z rotations are always simulated with the optional $\Lambda(Z)$. Substituting Eq. (49) into Eq. (45) gives another universal initial graph state



which can again be obtained from the cluster state by deleting the sites marked by empty circles:



This scheme requires only four physical qubits per logical qubit per cycle.

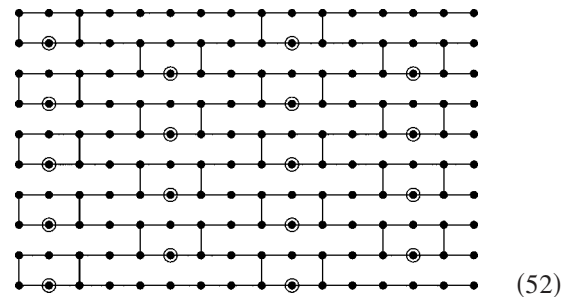
The above schemes are quite efficient in terms of the number of physical qubits per logical operation. Moreover, such efficiency is achieved with remarkably little manipulation; rather, it arises from following simple guidelines in a systematic derivation.

3. The cancellation approach

The cancellation approach uses the fact that in $1WQC_{TG}$, the angle of single-qubit rotations can be entirely determined by the measurement basis, and can be chosen online. The idea is to intersperse two consecutive $\Lambda(Z)$ gates with a single-qubit rotation so that proper choices of the angle of rotation allow the two interactions to add up or cancel out. In particular,

$$\Lambda(Z)(I \otimes X_\theta)\Lambda(Z) = \begin{cases} I & \text{if } \theta = 0, \\ (I \otimes X_{\pi/2})\Lambda(Z) & \text{if } \theta = -\frac{\pi}{2}. \end{cases}$$

The gate $\Lambda = (I \otimes X_{\pi/2})\Lambda(Z)$ is locally equivalent to $\Lambda(Z)$, and is universal given the ability to perform all single-qubit gates. Thus, we can perform optional nearest-neighbor Λ gates by introducing pairs of $\Lambda(Z)$ with a variable X rotation interspersed between each pair. This observation can be used to extend $1WQC_{TG}$ to a scheme with a fixed initial state. The idea is to represent a quantum computation as a set of repeating subunits. A subunit consists of the following: (i) arbitrary X rotations, (ii) arbitrary Z rotations and $\otimes_{i \text{ odd}} \Lambda(Z)_{i,i+1}$, (iii) interspersed X rotations, (iv) arbitrary Z rotations and $\otimes_{i \text{ odd}} \Lambda(Z)_{i,i+1}$, (v) arbitrary X rotations, (vi) arbitrary Z rotations and $\otimes_{i \text{ even}} \Lambda(Z)_{i,i+1}$, (vii) interspersed X rotations, and (viii) arbitrary Z rotations and $\otimes_{i \text{ even}} \Lambda(Z)_{i,i+1}$, where subscripts on $\Lambda(Z)$ denote the qubits being acted on. Following the discussion in Sec. IV C 2, the initial graph state is given by



In this diagram, qubits corresponding to the interspersed X rotations are marked by extra circles. The above state can be used to simulate a five-qubit circuit for two cycles.

The above graph state can also be produced from a cluster state, though the resulting scheme is not as efficient as the first approach.

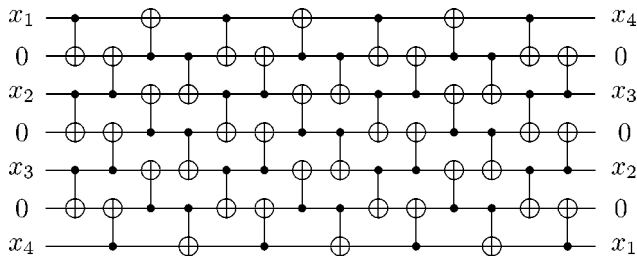
4. The routing approach

In the routing approach, a qubit can be teleported to an interaction or a non-interaction site. The interaction is always applied at the interaction site, but it may be applied to the quantum data or to some dummy state.

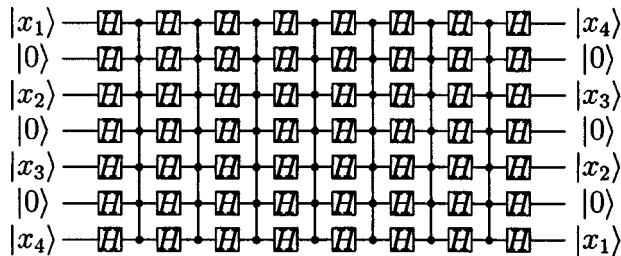
Using X teleportation, it is possible to teleport a given input state to one of several possible destinations. To see how this works, we consider the case of two destinations:

Examining this circuit identity and comparing with the circuit for X -teleportation, we see that by measuring the third qubit we can effect an X -teleportation of the first qubit to the second. Alternately, if we had decided instead to measure the second qubit, we would have been able to effect an X -teleportation of the first qubit to the third. Thus, we are able to choose to route the state $|\psi\rangle$ to one of two destina-

where the $n=4$ leftmost (rightmost) qubits carry the input (output). Every qubit in the input or the $(2n-1) \times (2n-1)$ middle block is measured in the X basis. To derive Eq. (A1), we first depict a classical circuit for the bit-reversal operation:



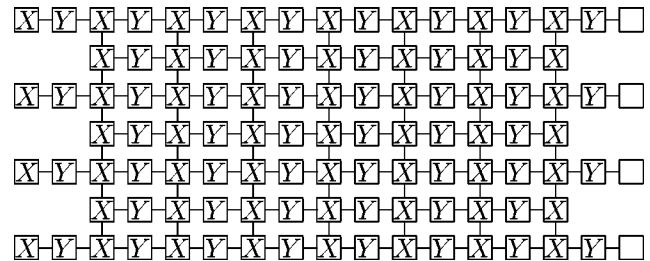
Note that the above is a valid quantum circuit for bit-reversal as well. Adding H to each of the input and output qubits in bit-reversal preserves the operation. Furthermore, rewriting $\Lambda(X)$ using Eq. (7), we obtain the following quantum circuit decomposition for bit-reversal:



Using the Euler angle decomposition

$$H = ie^{-iZ\pi/4}e^{-iX\pi/4}e^{-iZ\pi/4} \quad (A2)$$

and following our procedure in Sec. IV C, we obtain a 1WQC_{TG} scheme with substrate representation:



Finally, we can simplify the above to become Eq. (A1) by making the replacement:

$$\square - Y - \square \rightarrow \square - \square$$

due to Eq. (49). (The extra Z rotations on the two qubits always occur together, and compose to become a Pauli operation.)

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