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Quantum simulation of Pauli channels and dynamical maps: algorithm and implementation

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12 Abstract

Pauli channels are fundamental in the context of quantum computing as they model the simplest kind of noise in quantum devices. We propose a quantum algorithm for simulating Pauli channels and extend it to encompass Pauli dynamical maps (parametrized Pauli channels). A parametrized quantum circuit is employed to accommodate for dynamical maps. We also establish the mathematical conditions for an N-qubit transformation to be achievable using a parametrized circuit where only one single-qubit operation depends on the parameter. The implementation of the proposed circuit is demonstrated using IBM's quantum computers for the case of one qubit, and the fidelity of this implementation is reported.

²³ Since their inception, quantum computers were proposed as powerful tools for the ²⁴ simulation of quantum systems [1]. Being open quantum systems of fundamental [2,3]

22 1 Introduction

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25 and practical [4] interest, there has been efforts towards the simulation of the evolution
26 of open quantum systems [5-7] and specifically for quantum channels [8-10], which can
27 be used to study and model decoherence. Such quantum algorithms can be represented
28 using what is known as a quantum circuit [11], which we will study in section 3.
     Such systems have been simulated because of their many applications, such as
30 studying the emergence of multipartite entanglement [12, 13], studying dissipative
31 processes [14] and modeling non-Markovian dynamics [15]. Among quantum systems,
32 the simplest case is that of a qubit [11], and withing within them, the simplest class of
33 channels that produce decoherence are Pauli channels [16–18]. Indeed, they serve as
<sup>34</sup> effective models for the noise affecting quantum devices [19].
     In this article we propose a quantum algorithm for simulating Pauli channels and
36 implement it on one of IBM's quantum computers. The algorithm proposed is
37 straightforward and can be used for any Pauli channel by only changing a couple of
38 parameters in the operations it performs. To represent the algorithms implemented in
39 quantum computers, either to simulate a physical system or for some other purpose,
40 one often uses a quantum circuit algorithm, we use a quantum circuit, which is a
41 common way of representing algorithms intended for quantum computers. [11]. In this
42 work, we shall also work with parametrized quantum circuits, that is, quantum
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43 circuits in which some of the operations depend on variable parameters [20]. These 44 circuits play an important partin applications Furthermore, we will also simulate Pauli 45 dynamical maps, which are continuous parametrized curves of Pauli channels. These 46 maps can be used to model a continuous change of a qubit instead of only discrete 47 jumps. The generality of the algorithm proposed for Pauli channels will be very useful 48 in this part, since then using for dynamical maps will come very naturally. 49 Implementing these maps will lead us to study quantum algorithms with free 50 parameters, something that is common in areas such as quantum machine 51 learning [21] and describing general quantum transformations dependent on parameters. 52 Substantial research has been devoted to enhancing the efficiency of these circuits 53 [22]. This can be done using parametrized quantum circuits, which consist of 54 quantum circuits in which one or more operations depend on a free parameter. We start by providing the definition of quantum channels, the general framework 56 used here, and multi-qubit Pauli channels in section 2. Our first objective is to present 57 a general and straightforward quantum algorithm capable of simulating Pauli channels 58 on quantum computers; we do this in section 3, where we also demonstrate its 59 implementation using IBM's quantum computers for the particular case of single-qubit 60 several single and two-qubit Pauli channels. Expanding beyond discrete Pauli channels, 61 we introduce the concept of Pauli dynamical maps, defined as a continuous 62 parametrization of multi-qubit Pauli channels. Therefore, in section 4 we shift our 63 focus to In fact, in order to implement them, we study parametrized quantum channels, 64 aiming to adapt the algorithm developed for channels to dynamical mapscircuits in 65 chapter 4. Furthermore, we contribute to the body of work related to parametrized 66 quantum circuits by establishing a theorem, which sets the mathematical conditions for 67 the transformations that can be done using a parametrized circuit with the condition 68 restriction that only a controlled single-qubit rotation in the circuit may depend on the 69 parameter. Finally, in section 5, we conclude about analyzing the Pauli dynamical 70 maps that fulfill the conditions of theorem 2.

$_{\scriptscriptstyle{7}}\,2$ Pauli channels and dynamical maps

72 In this section we introduce the concept of quantum channels, focusing on a specific 72 type called Pauli channels. Furthermore, we define Pauli dynamical maps, which are 74 curves of Pauli channels parametrized by a variable.

75 2.1 Quantum channels

⁷⁶ In quantum mechanics, a closed system's state is represented by a vector in a Hilbert ⁷⁷ space \mathcal{H} . The state's evolution is unitary and given by Schrodinger's equation [23]. ⁷⁸ However, in real-world situations, quantum systems are usually open, which means that ⁷⁹ they interact with their environment [4]. For instance, the system's state may become ⁸⁰ entangled with the environment, leading to a loss of information about the system's ⁸¹ state over time.

To describe open systems, instead of state vectors, we use matrices ρ that act on \mathcal{H} . These matrices are called density matrices, and they include information about the system's interaction with its environment. For a density matrix ρ to be physically valid, it must satisfy two conditions: $\operatorname{tr}(\rho)=1$ and it must be positive semi-definite, which is denoted as $\rho\geq 0$ [11].

Knowing this, we can now define quantum channels. Quantum channels are 98 operators $\mathcal E$ that can describe the evolution of open quantum systems, such that 99 $\rho \to \mathcal E(\rho)$. Quantum channels are the most general linear operations that a quantum 90 system can undergo independently of its past [24,25]. These channels are constructed

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91 based on three fundamental properties: linearity, trace preservation, and complete 92 positivity.

Linearity ensures that a quantum channel \mathcal{E} maps any ensemble of density matrices into the corresponding ensemble of their evolution. The trace preserving property is given by $\operatorname{tr}(\mathcal{E}[\rho]) = \operatorname{tr}(\rho) = 1$ and guarantees that the quantum channel does not change the condition that $\operatorname{tr}(\rho) = 1$. Finally, the channel should also preserve the condition $\rho \geq 0$, and a map that does this is called a positive map. However, positivity of \mathcal{E} is not enough, and we actually require the more restrictive condition of complete positivity. Complete positivity means that $\mathcal{E} \otimes \mathbb{I}_n$ is positive for any positive integer n (where \mathbb{I}_n is entangled with another system, applying \mathcal{E} to the principal main system while doing nothing to the other one still results in a positive semidefinite state for the principal main system [11].

Given a quantum channel \mathcal{E} , the condition of trace preservation is straightforward to verify but complete positivity is not as simple. To test complete positivity of a quantum channel, Jamiołkowski and Choi [26,27] developed a simple algorithm that exploits the isomorphism between a channel \mathcal{E} and the state $\mathcal{D} = (\mathbb{I} \otimes \mathcal{E})[|\Omega\rangle\langle\Omega|]$, where $|\Omega\rangle = 1/\dim(\mathcal{H}) \sum_i^{\dim(\mathcal{H})} |i\rangle|i\rangle$ is a maximally entangled state between the original system and an ancilla. Remarkably, the map \mathcal{E} is completely positive if and only if \mathcal{D} (also known as the Choi or dynamical matrix of \mathcal{E}) is positive semidefinite.

111 2.2 Pauli channels

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¹¹² We have discussed the main features of quantum channels and now we turn our ¹¹³ attention to a specific type of channels for N-qubit systems called Pauli channels. First ¹¹⁴ we will define these channels for single-qubit systems, whose most general density ¹¹⁵ matrix can be written as [11]:

$$\rho = \frac{1}{2} \sum_{\alpha=0}^{3} r_{\alpha} \sigma_{\alpha}, \tag{1}$$

117 with $\sigma_0 = \mathbb{I}$, and $\sigma_{1,2,3}$ the usual Pauli matrices. The condition $\operatorname{tr}(\rho) = 1$ requires that 118 $r_0 = 1$ while $\rho \geq 0$ implies that the remaining $r_{1,2,3}$ form a vector $\vec{r} = (r_1, r_2, r_3)$ inside 119 a unit sphere known as the Bloch sphere [28]. That is, every possible density matrix for 120 a one-qubit system is uniquely associated with a point in a unit sphere.

Given a one-qubit system described by ρ , a Pauli channel is defined as an operation that with probability k_{γ} applies the Pauli matrix σ_{γ} to the system, for $\gamma = 0, 1, 2, 3$ [16]. Mathematically, the Pauli channel is written in the following way:

$$\mathcal{E}(\rho) = \sum_{\gamma=0}^{3} k_{\gamma} \sigma_{\gamma} \rho \sigma_{\gamma}, \tag{2}$$

where the probabilities k_{γ} of applying σ_{γ} are non-negative real numbers such that $\sum_{\gamma} k_{\gamma} = 1$ (these conditions also ensure that the channel is trace preserving and completely positive).

Pauli channels are some of the most fundamental noise models in quantum information science [29]. Some notable examples of Pauli channels are the following:

• Bit Flip Channel: This is a channel that with probability 1-p leaves the qubit as it is and with probability p applies the σ_1 matrix (which flips the basis states $|0\rangle$ and $|1\rangle$ of the qubit), and so it is given by:

$$\mathcal{E}(\rho) = (1-p)\rho + p\sigma_1\rho\sigma_1.$$

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Analogous channels exist using σ_3 (called the bit flip channel, which has a probability p of adding a relative phase π to the state) or using σ_2 (called the phase flip channel, which has a probability p of flipping the base states and also add a relative phase π).

• **Depolarizing channel:** This channel has a probability 1-p of doing nothing to the qubit and a probability p of converting it into the maximally mixed state $\frac{1}{2}\mathbb{I}$ and it can be written as:

$$\mathcal{E}(\rho) = (1-p)\rho + p\frac{1}{2}\mathbb{I} = \left(1 - \frac{3p}{4}\right)\sigma + \frac{p}{4}\sigma_1\rho\sigma_1 + \frac{p}{4}\sigma_2\rho\sigma_2 + \frac{p}{4}\sigma_3\rho\sigma_3. \tag{3}$$

We can also see how an arbitrary Pauli channel acts on an arbitrary density matrix.

To do it, we substitute Eq (1) in Eq (2):

$$\mathcal{E}(\rho) = \frac{1}{2} \sum_{\gamma,\alpha=0}^{3} k_{\gamma} r_{\alpha} \sigma_{\gamma} \sigma_{\alpha} \sigma_{\gamma}. \tag{4}$$

141 This can be simplified by using the following property noting that

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143 which was obtained following the commutation relation of Pauli matrices:

which leads to $[\sigma_{\alpha}, \sigma_{\beta}] = 2i\epsilon_{\alpha\beta\gamma}\sigma_{\gamma}$ and $\epsilon_{\alpha\beta\gamma}$ is the Levi-Civita tensor. Then, Eq. (4) takes the form

$$\mathcal{E}(\rho) = \frac{1}{2} \sum_{\alpha} \left(\sum_{\gamma} A_{\alpha,\gamma} k_{\gamma} \right) r_{\alpha} \sigma_{\alpha}. \tag{6}$$

¹⁴⁸ Eq (6) once again has the form of Eq (1) but with components $\left(\sum_{\gamma} A_{\alpha,\gamma} k_{\gamma}\right) r_{\alpha}$. This ¹⁴⁹ gives us another way of understanding Pauli channels as operations that take each ¹⁵⁰ component r_{α} of the density matrix and multiplies them by $\sum_{\gamma} A_{\alpha,\gamma} k_{\gamma}$, that is:

$$r_{\alpha} \xrightarrow{\text{Pauli}} \tau_{\alpha} r_{\alpha}, \quad \tau_{\alpha} := \sum_{\gamma} A_{\alpha, \gamma} k_{\gamma}. \tag{7}$$

Notice that $\tau_0=1$, which is a consequence of $\sum_{\gamma}k_{\gamma}=1$ and ensures that after the channel, the resulting density matrix still has trace one. Furthermore, reverting the definition of τ_{α} by using that $A^{-1}=\frac{1}{4}A$, we get that $k_{\gamma}=\frac{1}{4}\sum_{\alpha}A_{\alpha,\gamma}\tau_{\alpha}$. Then, using that $k_{\gamma}\geq 0$ we get the following conditions on the multipliers τ_{α} :

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$$1 + \tau_i - \tau_j - \tau_k \ge 0, \quad \text{for } i, j, k \text{ different numbers in } \{1, 2, 3\}, \tag{8}$$

$$1 + \tau_1 + \tau_2 + \tau_3 \ge 0. \tag{9}$$

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These conditions imply that (τ_1, τ_2, τ_3) has to be inside a tetrahedron with vertices (1,1,1), (1,-1,-1), (-1,1,-1) and (-1,-1,1). Therefore, the $\tau_{1,2,3}$ are numbers between -1 and 1, which means that the components r_{α} of the density matrix are always attenuated and possibly sign flipped.

Having defined the one qubit case, we can now generalize to N qubits. In order to do it, we need to introduce the so-called *Pauli strings*, defined as

$$\sigma_{\vec{\alpha}} = \sigma_{\alpha_1} \otimes \sigma_{\alpha_2} \otimes \cdots \otimes \sigma_{\alpha_N}, \tag{10}$$

165 where $\vec{\alpha}$ denotes a multi-index $(\alpha_1, \cdots, \alpha_N)$ and $\alpha_i \in \{0, 1, 2, 3\}$. These operators form 166 an orthogonal basis in the space of operators acting on N qubits. Similarly to the 167 single-qubit case, the density matrix ρ of a system of N qubits can be written using 168 Pauli strings as:

$$\rho = \frac{1}{2^N} \sum_{\vec{\sigma}} r_{\vec{\alpha}} \sigma_{\vec{\alpha}}. \tag{11}$$

Then, just as before, we define a Pauli channel as a transformation that applies the 171 operator $\sigma_{\vec{\gamma}}$ to ρ with probability $k_{\vec{\gamma}}$ and is therefore described mathematically by:

$$\mathcal{E}(\rho) = \sum_{\vec{\gamma}} k_{\vec{\gamma}} \sigma_{\vec{\gamma}} \rho \sigma_{\vec{\gamma}}, \tag{12}$$

where just as before, $k_{\vec{\gamma}}$ are non-negative real numbers such that $\sum_{\vec{\gamma}} k_{\vec{\gamma}} = 1$.

As in the one qubit case, Pauli channels for N qubits attenuate the components $r_{\vec{\alpha}}$ of the density matrix. This can be seen by substituting Eq (11) in Eq (12) and using the property of Eq (5):

$$\mathcal{E}(\rho) = \frac{1}{2^N} \sum_{\vec{\gamma}, \vec{\alpha}} k_{\vec{\gamma}} r_{\vec{\alpha}} \sigma_{\vec{\gamma}} \sigma_{\vec{\alpha}} \sigma_{\vec{\gamma}}$$
$$= \frac{1}{2^N} \sum_{\vec{\alpha}} \left(\sum_{\vec{\gamma}} (A^{\otimes^N})_{\vec{\alpha}, \vec{\gamma}} k_{\vec{\gamma}} \right) r_{\vec{\alpha}} \sigma_{\vec{\alpha}},$$

which means that applying the Pauli channel multiplies the components $r_{\vec{\alpha}}$ by $\tau_{\vec{\alpha}} := \sum_{\vec{\gamma}} (A^{\otimes^N})_{\vec{\alpha},\vec{\gamma}} k_{\vec{\gamma}}$.

176 2.3 Pauli dynamical maps

¹⁷⁷ As seen in the last section, Pauli channels and in general quantum channels are discrete maps that transform a density matrix ρ into $\mathcal{E}(\rho)$. However, we could also define a continuous set of channels ε_p with p a real parameter.

For the special case of Pauli channels, we define a Pauli dynamical map as a continuous parametrized curve drawn inside the set of Pauli channels and starting at the identity channel. Therefore, a Pauli dynamical map can be written as

$$\mathcal{E}_p(\rho) = \sum_{\vec{\gamma}} k_{\vec{\gamma}}(p) \sigma_{\vec{\gamma}} \rho \sigma_{\vec{\gamma}}, \tag{13}$$

where p is a parameter in an interval [a, b] and \mathcal{E}_p is a Pauli channel for every p, with \mathcal{E}_a being the identity channel.

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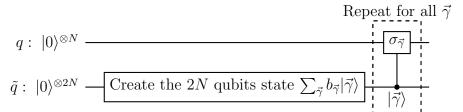


Fig 1. Circuit for an N-qubit Pauli channel. The circuit creates the state $\sum_{\vec{\gamma}} b_{\vec{\gamma}} |\vec{\gamma}\rangle$ on the 2N ancilla qubits denoted by \tilde{q} .

3 Circuit for a Pauli channel

187 In this section we propose a quantum circuit that simulates N-qubit Pauli channels. A 188 quantum circuit is a model for representing computation on a quantum system. This 189 computation may include preparing the initial state of the system, applying unitary 190 operations and measuring the whole system or parts of it [23]. Our systems always 191 consists of some number of qubits, that can be divided into the ones to which we will 192 apply the quantum channel, or main qubits and the ones that help us complete said 193 operation, or ancilla qubits. The circuit implements a unitary operator U on the 194 system consisting of all the qubits and then the effect on the main qubits is measured. 195 Moreover, we implement the circuit for N=1 on a quantum computer and analyze the 196 results using the diamond norm. We find that close to the depolarizing channel, the 197 general circuit simulator can implement such channels with the highest fidelity.

3.1 Description of the circuit for a Pauli channel

¹⁹⁹ To design the circuit that implements Eq (12), we construct a state that includes the ²⁰⁰ probabilities k_{γ} on the ancilla qubits and subsequently apply controlled Pauli operations ²⁰¹ on the main qubits. The circuit that does this is presented in Fig 1.

The first part of the circuit involves the creation of the state

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Ancilla state =
$$\sum_{\vec{\gamma}} b_{\vec{\gamma}} |\vec{\gamma}\rangle$$
 (14)

on the ancilla qubits, where $b_{\vec{\gamma}}$ are numbers such that $|b_{\vec{\gamma}}|^2 = k_{\vec{\gamma}}$ and the 2N-qubit state $|\vec{\gamma}\rangle$ is defined as $|\gamma_1\rangle\cdots|\gamma_N\rangle$. When measured in the computational basis, the state given in Eq (14) collapses to $|\vec{\gamma}\rangle$ with a probability $|b_{\vec{\gamma}}|^2 = k_{\vec{\gamma}}$. The circuit in Fig 1 uses this fact to apply $\sigma_{\vec{\gamma}}$ on the main qubits with a probability $k_{\vec{\gamma}}$ by using controlled operations conditioned on the state of the system being $|\vec{\gamma}\rangle$, just as the Pauli channel is supposed to do.

210 3.2 Simulation for one-qubit Pauli channels

For the particular case of a Pauli channel on one qubit, the circuit that simulates it can be constructed as in Fig 2, which is a special case of Fig 1 but with all details explicitly shown. In said figure, the ancilla state of Eq (14) can be taken to be $\sqrt{k_0}|00\rangle + \sqrt{k_1}|01\rangle + \sqrt{k_2}|10\rangle + \sqrt{k_3}|11\rangle$ and it is created on the ancilla qubits with the

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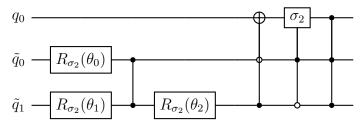


Fig 2. One-qubit Pauli channel circuit. Circuit for a one-qubit Pauli channel, which is a particular case of Fig 1. Here we have two ancilla qubits and we use three rotations of angles given by Eq (15) to create the Ancilla State of Eq (14) for the two ancilla qubits.

help of three rotations of angles defined by the following equations:

$$\cos\left(\frac{\theta_0}{2}\right) = \sqrt{k_0 + k_1},$$

$$\tan\left(\frac{\theta_1 + \theta_2}{2}\right) = \sqrt{k_1/k_0},$$

$$\tan\left(\frac{\theta_2 - \theta_1}{2}\right) = \sqrt{k_3/k_2}.$$
(15)

We took a sample of one-qubit Pauli channels and evaluated their implementation ²¹² on IBM's ibmq-lima quantum computer [30], as shown in Fig 3. For each of the ²¹³ channels sampled, we used quantum process tomography [30,31] to obtain the operator ²¹⁴ ξ_I corresponding to the implementation of the circuit in the quantum computer. Then, ²¹⁵ we compared ξ_I with the theoretical operator ξ_T of the Pauli channel we wanted to ²¹⁶ implement. To see how close the operators ξ_I and ξ_T are, we shall use the diamond ²¹⁷ distance [32], which is defined by

$$||\xi_I - \xi_T||_{\diamond} = \max_{\rho} ||(\xi_I \otimes I)\rho - (\xi_T \otimes I)\rho||_1, \tag{16}$$

with I the identity map, $||\cdot||_1$ the trace norm and the maximization done over all density matrices ρ . The calculation of this norm is done using the semi-definite program from reference [33]. When the two channels are the same, the diamond distance has a value of 0, while in the case that the channels are completely distinguishable, the distance reaches its maximum value of 2 [34]. For the analysis done in Fig 3, we define a value of "diamond fidelity" as:

$$f = 1 - \frac{1}{2} ||\varepsilon_I - \varepsilon_T||_{\diamond}, \tag{17}$$

²²⁶ which ranges from 0, when the channels have a maximum distance, to 1, when they are ²²⁷ exactly equal.

Finally, using the representation of Pauli channels in a tetrahedron as in Eq (8), we show in Fig 3 the diamond fidelity defined by Eq (17) for the channels analyzed. We can see that channels close to the completely depolarizing channel (that is close to the tetrahedron) have a high f, while those close to unitary channels have much lower f. This is reasonable because quantum computers are prone to errors that depolarize qubits, which isn't very problematic when trying to simulate depolarization but it is when simulating unitary processes. Moreover, Moreover, as can be seen in Fig 3, the algorithm of Fig 2 is not optimal for unitary channels (that is, the channels corresponding to the vertices channels in the edges of the tetrahedron). These

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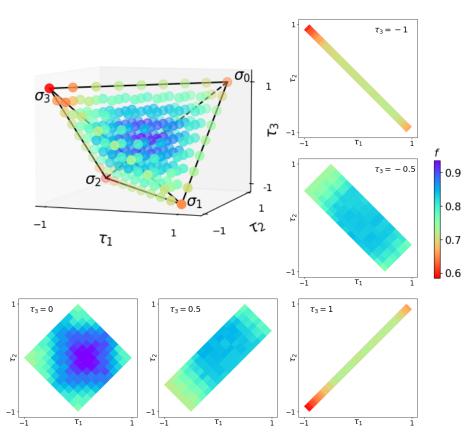


Fig 3. Results of the diamond fidelities as defined in Eq (17) for a sample of Pauli channels in the tetrahedron. Notice that channels close to the center of the tetrahedron have high fidelities, while those close to the borders do not. Moreover, we show the results for cuts of the tetrahedron at different values of τ_3 .

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straightforward, which corresponds to unitary channels. These channels could be accomplished more efficiently by simply applying the corresponding Pauli operation directly. Nevertheless, due to its general design to accommodate any unitary operator directly. In this case we would be trading generality for fidelity. The main advantage of the algorithm is that we can simulate with a unitary circuit any Pauli channel, only varying a couple of angles in the circuit. In fact, as we will see, sometimes only changing one angle in the circuit will be enough to simulate many non-unitary dynamical maps of physical interest.

245 3.3 Simulation of a two-qubit Pauli channel

For the general case of a two-qubit Pauli channel, the algorithm employs numerous quantum gates even in such scenarios circuit of Fig 1 requires two main qubits and four ancilla qubits. However, the quantum computer which we are using has only 5 qubits. We could still simulate a two qubit Pauli channel that applies at most 8 Pauli operations, instead of 16 which would be the most general case, using the three ancilla qubits. For example, consider the following two-qubit dephasing channel:

$$\varepsilon(\rho) = (1 - p)\rho + \frac{p}{3}(I \otimes \sigma_z)\rho(I \otimes \sigma_z) + \frac{p}{3}(\sigma_z \otimes I)\rho(\sigma_z \otimes I) + \frac{p}{3}(\sigma_z \otimes \sigma_z)\rho(\sigma_z \otimes \sigma_z),$$
(18)

where $p \in [0,1]$ is a parameter that defines the channel. Since this channel has four Pauli operations, only 2 ancilla qubits will be necessary, and thus it can be implemented using the circuit of Fig 4b.

Just as for the one-qubit Pauli channels, this channel was implemented on IBM's ibmq-lima quantum computer. We obtained the Choi matrix by doing quantum process tomography on the main qubits. Then, we calculated the diamond fidelity for different values of p and obtained the results shown in Fig 4a. As can be seen in the figure, the fidelity doesn't vary much as we change the value of p. The average fidelity over all values of p is 0.758 and it is biggest when p=0, corresponding to the identity channel, where it reaches a fidelity of 0.874.

55 4 One parameter circuits

²⁵⁷ Just Finally, we consider the simulation of Pauli dynamical maps. This is pretty much already solved, since just as Pauli channels, Pauli dynamical maps can be implemented ²⁵⁸ using the circuit of Fig 1. However, there is one difference: the state to be created on ²⁶⁰ the ancilla qubits now depends on a parameter p, and it is represented by the expression:

$$\sum_{\vec{\gamma}} b_{\vec{\gamma}}(p) |\vec{\gamma}\rangle. \tag{19}$$

²⁶² Thus, we temporarily shift our focus from Pauli channels and dynamical maps to the general problem of creating a circuit to generate a curve of states like the one described ²⁶⁴ in Eq (19).

In general, producing this curve of states for N qubits will require many rotations parametrized by p, such as the three rotations used for the ancilla qubits in Fig 2. However, it would be preferable to achieve the same effect using only one parametrized rotation. This would allow us to interpret said rotation as a knob that smoothly traverses the curve of states. Consequently, we are faced with the question of which curves of states, such as the one described in Eq (19), can be produced using just a rotation are produced using just a rotation. To clarify this, we provide the following definition for a circuit with one parametrized rotation.

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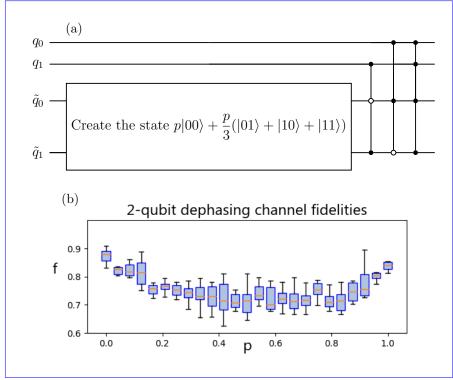


Fig 4. Circuit for 2-qubit dephasing channel. Circuit implementation for the 2-qubit dephasing channel of Eq (18) and fidelity results of its implementation using ibmq-lima. The circuit is shown in figure a), where the state on the two ancilla qubits can be constructed just as it was done for Fig 2. In figure b) we show the results of fidelities after implementing this circuit on ibmq-lima for 25 values of p between 0 and 1, each of which is simulated 10 times.

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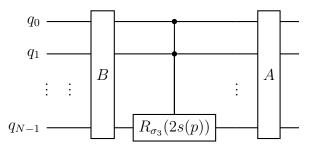


Fig 5. General form of a 1PR circuit. Any 1PR circuit can be transformed into this form, where the rotation on the last qubit can be controlled or not by any of the other qubits. A and B are N-qubit gates that do not depend on the parameter p and s = s(p) is a function of the parameter.

273 **Definition 1** 1-Parameter Rotation Circuit: A 1-Parameter Rotation (1PR) 274 circuit is a parametrized quantum circuit that includes only one gate dependent on a 275 parameter p. Moreover, the parametrized gate is a one-qubit rotation about any axis, 276 whether controlled or not.

Therefore, a 1PR circuit implements a unitary operator U(p) on some number N of gubits, such that U depends on the parameter p only locally on one specific gubit.

Based on this definition, we aim to determine which curves of states can be generated using 1PR circuits. To accomplish this, we begin by proving that all 1PR circuits have the form depicted in Fig 5, where the parametrized rotation is around σ_3 and is applied to the last qubit.

Theorem 1 An N-qubit 1PR circuit can always be transformed into the form shown 284 in Fig 5.

Proof: First, we observe that according to the definition, a 1PR circuit always consists of an operation B followed by the parametrized rotation on a specific qubit and then another operation A, where A and B are not parametrized.

Next, we note that it is not necessary to consider rotations about an arbitrary axis, 289 as a rotation about any axis \hat{n} parameterized by p can be transformed into a rotation 290 about σ_3 without introducing gates that depend on p. To see this, consider the rotation 291 $R_{\hat{n}}(2s)$, where 2s is a function of p (the factor of 2 is for convenience later on) and 292 $\hat{n} = (n_1, n_2, n_3)$ represents the rotation axis. We can express \hat{n} as 293 $(\sin\theta\cos\phi, \sin\theta\sin\phi, \cos\theta)$, where θ and ϕ are fixed angles dependent on \hat{n} . The 294 rotation can then be rewritten as follows:

$$R_{\hat{n}}(2s) = R_{\sigma_3}(\phi) R_{\sigma_2}(\theta) R_{\sigma_3}(2s) R_{\sigma_2}(-\theta) R_{\sigma_3}(-\phi). \tag{20}$$

²⁹⁶ Since the angles θ and ϕ do not depend on the parameter p, any 1PR circuit can be ²⁹⁷ transformed into a circuit where the parametrized rotation is around σ_3 instead of an ²⁹⁸ arbitrary axis. Moreover, without loss of generality, we can choose the last qubit as the ²⁹⁹ target qubit for the rotation, since if it weren'tis not, we could use swap gates to move ³⁰⁰ the rotation to the first qubit without adding gates that depend on p.

Therefore, a 1PR circuit can be transformed such that the rotation is around σ_3 and σ_3 is applied to the last qubit (possibly controlled by other qubits), resulting in the form depicted in Fig 5. \blacksquare

With the aid of this theorem, we can now determine the curves of states of N qubits that can be generated using a 1PR circuit. This result is stated in the following theorem.

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Theorem 2 Consider a 1PR circuit of N qubits parametrized by p and denote by U(p) the operator it implements on this system. Then, for every $j \in \{0, 1, \dots, 2^N - 1\}$, we have that:

$$U|j\rangle = e^{is(p)}|a^j\rangle + e^{-is(p)}|b^j\rangle + |c^j\rangle$$

$$U(p)|j\rangle = e^{is(p)}|a^j\rangle + e^{-is(p)}|b^j\rangle + |c^j\rangle$$
(21)

307 with s(p) some function of p, $|a^j\rangle, |b^j\rangle, |c^j\rangle$ orthogonal states and 308 $\langle a^j|a^j\rangle + \langle b^j|b^j\rangle + \langle c^j|c^j\rangle = 1$. Here $|j\rangle$ represents an element of the N-qubit 309 computational basis.

Proof: We can conclude from theorem 1 that U=ARB, where A and B are unitary matrices. As before, from the definition of a 1PR circuit, we know that U(p) can be decomposed as ARB, where A, B are unitary operators acting on the N-qubit system and B are rotation applied to one qubit. From theorem 1, we conclude that B can be taken in general as a B3 rotation of angle B3. Prove this theorem, we will apply B4 to B5. First, applying B5, then B6, B7 results in B8, where B8, where B9 results in B9. This can be rewritten by separating the last qubit and First, applying B8 to B9. This can be rewritten by separating the last qubit from the other B9. This can be rewritten by separating the last qubit from the other B9.

$$B|j\rangle = \sum_{k=0}^{2^{N-1}-1} (B_{2k,j}|k\rangle|0\rangle + B_{2k+1,j}|k\rangle|1\rangle).$$
 (22)

After the operator B, the circuit applies the controlled rotation $R_{C_3}(2s(p))$. To simplify the analysis, we separate the states of the first N-1 qubits into those that fulfill the control conditions of the rotation (which we denote as the set C) and those that do not, and write it as

$$B|j\rangle = \sum_{k \in \mathcal{C}} \left(B_{2k,j} |k\rangle |0\rangle + B_{2k+1,j} |k\rangle |1\rangle \right) + \sum_{k \notin \mathcal{C}} \left(B_{2k,j} |k\rangle |0\rangle + B_{2k+1,j} |k\rangle |1\rangle \right). \tag{23}$$

Then, the rotation R will only affect the states on the first sum (since they fulfill the control conditions) and not the others. Therefore, remembering that a σ_3 using that a $R_{\sigma_3}(2s(p))$ rotation acts by adding a phase $e^{-is(p)}$ to $|0\rangle$ and a phase $e^{is(p)}$ to $|1\rangle$, we have that,

$$\underbrace{RBR_{\sigma_3}(2s(p))B}_{k\in\mathcal{C}}|j\rangle = e^{-is(p)} \sum_{k\in\mathcal{C}} B_{2k,j}|k\rangle|0\rangle + e^{is(p)} \sum_{k\in\mathcal{C}} B_{2k+1,j}|k\rangle|1\rangle + \sum_{k\notin\mathcal{C}} (B_{2k,j}|k\rangle|0\rangle + B_{2k+1,j}|k\rangle|1\rangle$$

$$= e^{-is(p)} |\tilde{b}^j\rangle + e^{is(p)} |\tilde{a}^j\rangle + |\tilde{c}^j\rangle, \tag{24}$$

where we defined

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$$\begin{split} |\tilde{a}^{j}\rangle &= \sum_{k \in \mathcal{C}} B_{2k,j} |k\rangle |\underline{01}\rangle, \\ |\tilde{b}^{j}\rangle &= \sum_{k \in \mathcal{C}} B_{2k+1,j} |k\rangle |0\rangle, \\ |\tilde{c}^{j}\rangle &= \sum_{k \notin \mathcal{C}} \left(B_{2k,j} |k\rangle |0\rangle + B_{2k+1,j} |k\rangle |1\rangle\right). \end{split}$$

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321 These states are clearly orthogonal because they are each linear combinations of 322 different orthogonal states of the computational basis. Moreover, they satisfy 323 $\langle a^j|a^j\rangle + \langle b^j|b^j\rangle + \langle c^j|c^j\rangle = 1$ because this quantity is the squared norm of the jth 324 column of B, which is unitary.

Finally, after having applied the rotation, the circuit applies gate A, so that the result is given by:

$$U|j\rangle = ARB|j\rangle = e^{-is(p)}A|\tilde{a}^j\rangle + e^{is(p)}A|\tilde{b}^j\rangle + A|\tilde{c}^j\rangle = e^{-is(p)}|a^j\rangle + e^{is(p)}|b^j\rangle + |c^j\rangle,$$

$$U[j] = AR_{\sigma_3}(2s(p))B[j]$$

$$= e^{-is(p)}A[\tilde{a}^j] + e^{is(p)}A[\tilde{b}^j] + A[\tilde{c}^j]$$

$$= e^{-is(p)}|a^j| + e^{is(p)}|b^j| + |c^j|, \qquad (25)$$

where $|a^j\rangle=A|\tilde{a}^j\rangle, |b^j\rangle=A|\tilde{b}^j\rangle, |c^j\rangle=A|\tilde{c}^j\rangle$ are still orthogonal states that satisfy $\langle a|a\rangle+\langle b|b\rangle+\langle c|c\rangle=1$ because A is unitary.

This theorem implies that when starting from the state $|0\rangle$ or any other initial state, the only possible curves of states that can be created using a 1PR circuit are of the following form:

$$|\eta(p)\rangle = |c\rangle + e^{is(p)}|a\rangle + e^{-is(p)}|b\rangle,$$
 (26)

332 with conditions defined by the equations:

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$$\langle a|a\rangle + \langle b|b\rangle + \langle c|c\rangle = 1, \quad \langle a|b\rangle = \langle a|c\rangle = \langle b|c\rangle = 0.$$
 (27)

Moreover, it is possible to construct a 1PR circuit to generate any given curve of states described by Eq (26). One approach to achieve this is by utilizing the circuit depicted in Fig 5, with the parametrized rotation R applied to the last qubit controlled by all the other qubits. The There is some freedom when choosing the operators A and B can be defined as follows:

$$\begin{split} & \underline{B|0\rangle = \sqrt{\langle a|a\rangle}|2^N - 1\rangle + \sqrt{\langle b|b\rangle}|2^N - 2\rangle + \sqrt{\langle c|c\rangle}|2^N - 3\rangle,} \\ & A|2^N - 1\rangle = \frac{1}{\sqrt{\langle a|a\rangle}}|a\rangle, \ A|2^N - 2\rangle = \frac{1}{\sqrt{\langle b|b\rangle}}|b\rangle, \ A|2^N - 3\rangle = \frac{1}{\sqrt{\langle c|c\rangle}}|c\rangle. \end{split}$$

, we only need to make sure that:

$$B|0\rangle = \sqrt{\langle a|a\rangle}|2^{N} - 1\rangle + \sqrt{\langle b|b\rangle}|2^{N} - 2\rangle + \sqrt{\langle c|c\rangle}|2^{N} - 3\rangle,$$

$$A|2^{N} - 1\rangle = \frac{1}{\sqrt{\langle a|a\rangle}}|a\rangle, \ A|2^{N} - 2\rangle = \frac{1}{\sqrt{\langle b|b\rangle}}|b\rangle, \ A|2^{N} - 3\rangle = \frac{1}{\sqrt{\langle c|c\rangle}}|c\rangle. \tag{28}$$

The remaining part of the operators A and B can be defined chosen in any arbitrary manner as long as they are unitary. By starting from the initial state $|0\rangle$ and applying the circuit shown in Fig 5, straightforward calculations lead us to obtain the resulting Examples of these matrices are shown for particular dynamical maps in the next section.

To see that $AR_{\sigma_3}(s(p))B|0\rangle$ in fact creates the curve of states described in Eq (26).

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To see it, we can rewrite the expression of $B|0\rangle$ by separating the first N-1 qubits from the last one:

$$B|0\rangle = \sqrt{\langle a|a\rangle}|2^{N-1}-1\rangle|1\rangle + \sqrt{\langle b|b\rangle}|2^{N-1}-1\rangle|0\rangle + \sqrt{\langle c|c\rangle}|2^{N-1}-2\rangle|1\rangle.$$

Since the parametrized rotation $R = R_{\mathcal{O}_3}(2s(p))$ is controlled by all the first N-1 qubits, and it only applies to the first two terms of $B|0\rangle$. As a result, we obtain:

$$\underline{RB}R_{\sigma_3}(2s(p))B|0\rangle = \sqrt{\langle a|a\rangle}e^{is(p)}|2^{N-1}-1\rangle|1\rangle + \sqrt{\langle b|b\rangle}e^{-is(p)}|2^{N-1}-1\rangle|0\rangle + \sqrt{\langle c|c\rangle}|2^{N-1}-2\rangle|1\rangle.$$

Finally, applying the $\frac{\text{defined}}{\text{operator}}$ operator A to this state yields $\frac{\text{desired result}}{\text{constant}}$.

$$AR_{\sigma_3}(2s(p))B = |c\rangle + e^{is(p)}|a\rangle + e^{-is(p)}|b\rangle.$$

With this, we prove that any curve of states as that in Eq (26) can be constructed by the circuit in Fig 5 with the parametrized rotation controlled by all the other qubits by correctly choosing matrices A and B.

345 5 1PR circuit for a Pauli map

³⁴⁶ We can now use the previous results to conclude directly which Pauli dynamical maps ³⁴⁷ can be implemented with a 1PR circuit. For this, the curve of states of Eq (19) has to ³⁴⁸ be constructed with only one parametrized rotation, so it has to satisfy the conditions ³⁴⁹ of theorem 2. Therefore, this implies that the map

$$\varepsilon_p(\rho) = \sum_{\vec{\gamma}} k_{\vec{\gamma}}(p) \sigma_{\vec{\gamma}} \rho \sigma_{\vec{\gamma}}, \tag{29}$$

 $\beta_{\vec{\gamma}}(p)$ such that $|\beta_{\vec{\gamma}}(p)|^2 = k_{\vec{\gamma}}(p)$ and

$$\sum_{\vec{\gamma}} \beta_{\vec{\gamma}}(p) |\vec{\gamma}\rangle = |c\rangle + e^{is(p)} |a\rangle + e^{-is(p)} |b\rangle, \tag{30}$$

353 where $|a\rangle, |b\rangle, |c\rangle$ fulfill the conditions of Eq. (27).

For the particular case of one qubit, we can show some examples of Pauli dynamical maps implementable with a 1PR circuit, which are plotted in Fig 6. The examples we show include some of the most common maps: the bit flip, phase flip, bit-phase flip and depolarizing. However, we also include the parabolic dynamical map, defined in Eq (32) and shown in Fig 6. This map traces a parabola inside the tetrahedron connecting two of its vertices and it describes a frontier in the tetrahedron between Pauli channels that are reachable by Lindbladian dynamics and those that are not [18].

• Depolarizing: This dynamical map is given by

$$\varepsilon_p(\rho) = (1 - 3p/4)\rho + (p/4)\sigma_1\rho\sigma_1 + (p/4)\sigma_2\rho\sigma_2 + (p/4)\sigma_3\rho\sigma_3,$$

$$\varepsilon_p(\rho) = (1 - 3p/4)\rho + (p/4)\sigma_1\rho\sigma_1 + (p/4)\sigma_2\rho\sigma_2 + (p/4)\sigma_3\rho\sigma_3,$$
 (31)

with $p \in [0,1]$. Therefore, the curve of states $|\beta(p)\rangle$ needed on the ancilla qubits is such that $|\beta_0(p)|^2 = (1-3p/4), \ |\beta_1(p)|^2 = |\beta_2(p)|^2 = |\beta_3(p)|^2 = p/4$. Then, taking the β_j to be real, the curve of states can be

$$|\beta(p)\rangle = \sqrt{1-3p/4}|\underline{000}\rangle + \sqrt{p/4}|\underline{011}\rangle + \sqrt{p/4}|\underline{102}\rangle + \sqrt{p/4}|\underline{113}\rangle.$$

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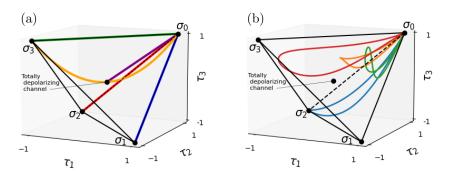


Fig 6. Some Pauli dynamical maps that can be implemented with a 1PR circuit. The curves painted in these tetrahedrons represent Pauli dynamical maps that can be implemented with 1PR circuits. (a) shows the dynamical maps mentioned in the main text, which are: depolarizing (purple), bit flip (blue), phase flip (green), bit-phase flip (red) and parabolic (orange). (b) shows dynamical maps selected at random that can be implemented with 1PR circuits.

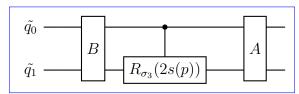


Fig 7. 1PR Circuit for one qubit curve of states. This is a 1PR circuit for creating a curve of states that satisfies the conditions of theorem 2. given the curve of states, operations A and B can be defined as in Eq (28).

This state can be rewritten as:

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$$\begin{split} |\beta(p)\rangle &= e^{is} \left(\frac{1}{2}|\underline{000}\rangle - \frac{i}{2\sqrt{3}}|\underline{011}\rangle - \frac{i}{2\sqrt{3}}|\underline{102}\rangle - \frac{i}{2\sqrt{3}}|\underline{113}\rangle\right) \\ &+ e^{-is} \left(\frac{1}{2}|\underline{000}\rangle + \frac{i}{2\sqrt{3}}|\underline{011}\rangle + \frac{i}{2\sqrt{3}}|\underline{102}\rangle + \frac{i}{2\sqrt{3}}|\underline{113}\rangle\right), \\ &\stackrel{\equiv}{=} e^{is}|a\rangle + e^{-is}|b\rangle \end{split}$$

with $\sin s = \sqrt{3p/4}$. We can see that this curve satisfies the conditions of Eq (27), meaning that it can be created with a 1PR circuit. Explicitly, following the discussion after theorem 2, we know that this curve of states can be constructed by the circuit of Fig 7.

The circuit requires defining A and B as in Eq (28), and for the particular vectors $|a\rangle, |b\rangle, |c\rangle$ of this dynamical map, we have that

$$A = \begin{pmatrix} 0 & 0 & 1/\sqrt{2} & 1/\sqrt{2} \\ 1/\sqrt{6} & -1/\sqrt{2} & i/\sqrt{6} & -i/\sqrt{6} \\ 1/\sqrt{6} & 1/\sqrt{2} & i/\sqrt{6} & -i/\sqrt{6} \\ -\sqrt{2/3} & 0 & i/\sqrt{6} & -i/\sqrt{6} \end{pmatrix}, B = \begin{pmatrix} 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 \\ 1/\sqrt{2} & -1/\sqrt{2} & 0 & 0 \\ 1/\sqrt{2} & 1/\sqrt{2} & 0 & 0 \end{pmatrix},$$

where, as mentioned before, the first column of A and the last three of B can be chosen arbitrarily as long as the resultant matrices are unitary. We ran this

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circuit in IBM's quantum computer ibmq-lima for 25 different values of p between 0 and 1, with 20 repetitions for each value of p. The diamond fidelity was calculated in each case and is shown in Fig 8.

• Parabolic dynamical map: We define the parabolic dynamical map as:

$$\epsilon(\rho) = (1-p)^2 \rho + (1-pp-p^2)\sigma_1 \rho \sigma_1 + (1-pp-p^2)\sigma_2 \rho \sigma_2 + (1+p)^2 \sigma_3 \rho \sigma_3, (32)$$

with $p \in [-1,1]$ $p \in [0,1]$. If we take the β_j to be real, the curve of states needed on the ancilla qubits can be:

$$|\beta(p)\rangle = (1-p)|\underline{000}\rangle + \sqrt{1-p^2}\sqrt{p-p^2}|\underline{011}\rangle + \sqrt{1-p^2}\sqrt{p-p^2}|\underline{102}\rangle + (1+p)|\underline{113}\rangle. (33)$$

This can be rewritten as

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$$\begin{split} |\beta(p)\rangle &= \left(\frac{1}{2}|\underline{000}\rangle + \frac{1}{2}|\underline{113}\rangle\right) + e^{is}\left(\frac{i}{4}|\underline{000}\rangle + \frac{1}{4}|\underline{011}\rangle + \frac{1}{4}|\underline{102}\rangle - \frac{i}{4}|\underline{113}\rangle\right) \\ &+ e^{-is}\left(\frac{-i}{4}|\underline{000}\rangle + \frac{1}{4}|\underline{011}\rangle + \frac{1}{4}|\underline{102}\rangle + \frac{i}{4}|\underline{113}\rangle\right) \\ &\equiv |c\rangle + e^{is}|a\rangle + e^{-is}|b\rangle, \end{split}$$

with $\sin s = p \sin s = 2p - 1$, so that this map curve fulfills the conditions of Eq (27). This means that it can be created with the 1PR circuit of figure Fig 7. In this case, matrices A and B can be taken to be:

$$A = \begin{pmatrix} 0 & 1/\sqrt{2} & -i/2 & i/2 \\ -1/\sqrt{2} & 0 & 1/2 & 1/2 \\ 1/\sqrt{2} & 0 & 1/2 & 1/2 \\ 0 & 1/\sqrt{2} & i/2 & -i/2 \end{pmatrix}, B = \begin{pmatrix} 0 & 1 & 0 & 0 \\ 1/\sqrt{2} & 0 & 0 & -1/\sqrt{2} \\ 1/2 & 0 & 1/\sqrt{2} & 1/2 \\ 1/2 & 0 & -1/\sqrt{2} & 1/2 \end{pmatrix},$$

where again, the first column of A and the last three of B can be chosen arbitrarily. As before, we ran this circuit for 25 values of p between 0 and 1, with 20 repetitions for each, obtaining as a result the fidelities shown in Fig 8.

• Bit flip map: This dynamical map is defined as

$$\varepsilon_n(\rho) = (1-p)\rho + p\sigma_1\rho\sigma_1,$$

for $p \in [0,1]$. In particular, if we take the β_j to be real, we need to create the curve of states:

$$|\beta(p)\rangle = \sqrt{1-p}|\underline{000}\rangle + \sqrt{p}|\underline{011}\rangle.$$

This can be rewritten as

$$|\beta(p)\rangle = e^{is}\left(\frac{1}{2}|\underline{000}\rangle - \frac{i}{2}|\underline{011}\rangle\right) + e^{-is}\left(\frac{1}{2}|\underline{000}\rangle + \frac{i}{2}|\underline{011}\rangle\right),$$

with $\sin s = \sqrt{p}$. Therefore, we can see that the curve satisfies the conditions of Eq (27) and it can be created with a 1PR circuit. Note that in this case we actually only need one ancilla qubit since the state $|\beta(p)\rangle$ is only two dimensional. Matrices A and B can be obtained as for the other dynamical maps and we ran the resulting circuit in ibmq-lima with the same specifications as for the other examples, with the results are shown in Fig 8.

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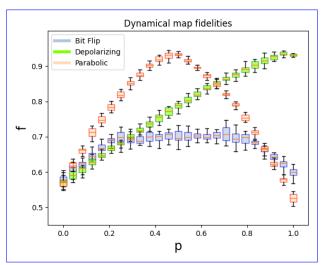


Fig 8. Fidelities for dynamical maps. Results of the fidelities obtained when simulating the three dynamical maps considered (depolarizing, parabolic and bit flip). The implementation was done on IBM's ibmq-lima quantum computer. Furthermore, we can see that the fidelities are comparable to those of Fig 3, so that using the 1PR circuit doesn't considerably affect fidelity.

The exact same thing can be done for the phase flip and bit phase flip dynamical maps by changing σ_1 to σ_3 and σ_2 respectively. For example, the bit phase flip map was implemented in [12,35] using an optical arrangement, and it was indeed done by varying only one angle that depends on the parameter p (the angle of a half waveplate).

From the results of Fig 8, we can see that just as in figure Fig 3, fidelities are high for points close to the depolarizing channel (p = 1 for depolarizing map, p = 1/2 for) parabolic map) and lower for points close to unitary channels (p = 0 for all maps and) p = 1 for parabolic and bit flip). Furthermore, we can see that the actual values of the fidelities are comparable to those of Fig 3, which means that using the 1PR circuit as opposed to implementing each channel with Fig 2 does not considerably affect fidelity.

Furthermore, we can construct other examples of Pauli dynamical maps such that they can be implemented with a 1PR circuit. To do it, we only need to choose the three states $|a\rangle$, $|b\rangle$, $|c\rangle$ that satisfy the conditions of Eq (27). For example, this can be done systematically for the case of curves of states of two qubits (that is, for Pauli dynamical maps of one qubit) with the following procedure:

1. We first choose the norms |a|, |b|, |c| such that $\langle a|a\rangle + \langle b|b\rangle + \langle c|c\rangle = 1$. This can be done by selecting two angles $\mu \in [0, \pi/2]$, $\nu \in [0, \pi/2]$ and defining:

$$|a| = \sin \nu \cos \mu$$
, $|b| = \sin \nu \sin \mu$, $|c| = \cos \nu$.

2. We define $|a'\rangle = |a||0\rangle$, $|b'\rangle = |b||1\rangle$, $|c'\rangle = |c||2\rangle$.

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3. Finally, we choose a unitary matrix V with the condition that its first row is equal to $e^{i\theta}(|a|,|b|,|c|,0)$ with θ a uniform random phase. That way, we can define $|a\rangle = V|a'\rangle$, $|b\rangle = V|b'\rangle$, $|c\rangle = V|c'\rangle$ and since V is unitary, these unprimed vectors will fulfill the conditions of Eq (27). Furthermore, the form of the first row ensures that the dynamical map begins at the identity, since it implies that when s=0, the state created in Eq (30) is $|a\rangle + |b\rangle + |c\rangle = e^{i\theta}|0\rangle$, which corresponds with applying the identity channel.

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Such a matrix V can be randomly constructed by first finding three vectors $\vec{w}_1, \vec{w}_2, \vec{w}_3$ orthogonal to the first row using the Gram-Schmidt process. Then selecting random complex numbers r_1, r_2, r_3 such that $|r_1|^2 + |r_2|^2 + |r_3|^2 = 1$ and defining the second row of V to be $r_1\vec{w}_1 + r_2\vec{w}_2 + r_3\vec{w}_3$. Once the first two rows are chosen, use Gram-Schmidt to find two vectors \vec{v}_1, \vec{v}_2 orthogonal to them and similarly define the third row as $q_1\vec{v}_1 + q_2\vec{v}_2$ with $|q_1|^2 + |q_2|^2$ and q_1, q_2 selected at random. Finally, there is only one choice for the fourth row so that it is orthonormal to the first three and a random phase can be given to it.

Following this procedure for random angles and unitary matrices V, we plot four Pauli dynamical maps selected at random that can be implemented with a 1PR circuit in Fig 6.

419 6 Conclusion

 $_{420}$ In this work, we found a quantum algorithm for simulating Pauli channels in N-qubit $_{421}$ systems and generalized it to Pauli dynamical maps by using parametrized quantum $_{422}$ circuits. Furthermore, we implemented single-qubit Pauli channels on one of IBM's $_{423}$ quantum computers and obtained their fidelities. Finally

Then, when working with Pauli dynamical maps, we searched for a way of
simplifying the parametrized minimizing the amount of parametric operations in the
circuit by requiring that only one single-qubit rotation depends on the parameter. In
theorem 2 we found the general mathematical conditions for this, applicable to any
parametrized circuit. Given this condition, we considered the space of Pauli maps that
are possible to parametrize with one parameter, and gave some explicit examples for a
system of one qubit. As the number of qubits gets bigger, the degrees of freedom for
choosing a Pauli dynamical map also gets bigger, so we would expect that the space of
maps that can be done with only one parameter becomes a relatively smaller part of
the whole space of Pauli dynamical maps. However, some paradigmatic maps as the
dephasing one we studied for 2 qubits will still be possible to do with a 1PR circuit
for this case, this can be seen by noticing that said dephasing channel has the same
form as the depolarizing channel for one qubit, which we proved to be possible to
minimizing the amount of parametric operations in the
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ThereforeIn conclusion, this work presents yet another example of the current exploration into simulating open quantum systems in quantum computers, and we observe the big effect that the error of quantum computers a quantum computer have on these simulations—, quantified by the fidelity. On the other hand, the result of theorem 2 shows what can be done with the condition of using only one parametrized rotation and can be applied to any quantum algorithm that requires parametrized circuits, such as those used for quantum machine learning [21].

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