

## Quantum error correction for continuously detected errors

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We show that quantum feedback control can be used as a quantum-error-correction process for errors induced by a weak continuous measurement. In particular, when the error model is restricted to one, perfectly measured, error channel per physical qubit, quantum feedback can act to perfectly protect a stabilizer codespace. Using the stabilizer formalism we derive an explicit scheme, involving feedback and an additional constant Hamiltonian, to protect an  $(n-1)$ -qubit logical state encoded in  $n$  physical qubits. This works for both Poisson (jump) and white-noise (diffusion) measurement processes. Universal quantum computation is also possible in this scheme. As an example, we show that detected-spontaneous emission error correction with a driving Hamiltonian can greatly reduce the amount of redundancy required to protect a state from that which has been previously postulated [e.g., Alber *et al.*, Phys. Rev. Lett. **86**, 4402 (2001)].

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### I. INTRODUCTION

Many of the applications of quantum-information science, such as quantum computation [1,2] and quantum cryptography [3], rely on preserving the coherence of quantum states. However, these states are typically short-lived because of unavoidable interactions with the environment. Combatting this decoherence has been the subject of much study.

Two important tools that have been developed for this task are quantum error correction [4–7] and quantum feedback [8–10]. In the usual protocol for quantum error correction, projective measurements are performed to acquire an error syndrome. A unitary operation chosen based on the results of the projective measurements is then applied to correct for the error. Quantum feedback control, on the other hand, uses the tools of continuous measurements and Hamiltonian feedback. The parameter to be controlled is typically the strength of the feedback Hamiltonian, which is conditioned on the result of the continuous measurements.

Quantum error correction and quantum feedback both rely on performing operations that are conditioned on the result of some measurements on the system, which suggests that exploring the links between these two techniques adds to our understanding of both processes and may lead to insights into future protocols and experimental implementations. In particular, this work provides an alternate avenue for examining the situation considered in Refs. [11–14] of correcting for a specific error process, such as spontaneous emission, at the expense of correcting fewer general errors. Practically, as these authors point out, it makes sense to pursue the trade-off between general correction ability and redundancy of coding, as smaller codes are more likely to be in the range of what can be experimentally realized in the near future. We shall

see that combining the pictures of quantum feedback and error correction provides a convenient framework in which to investigate this situation.

An additional motivation for considering the union of these techniques, as in Ref. [15], is to examine what is possible with different physical tools. In particular, we consider continuous measurements and Hamiltonians instead of the projective measurements and fast unitary gates generally assumed by discrete quantum error correction. Continuous error correction might well be useful even in a scenario in which near-projective measurements are possible (e.g., ion traps [16] and superconducting qubits [17]); it could be modified to provide bounds on how strong interactions in such systems would have to be to perform operations such as error correction and stay within a certain error threshold.

Reference [15] presupposed that classical processing of currents could be done arbitrarily and quickly, so the feedback was allowed to be an extremely complicated function of the entire measurement record. This can be modeled only by numerical simulations. In this paper, by contrast, we will restrict our feedback to be directly proportional to measured currents, thus removing any need for classical postprocessing. In the Markovian limit, this allows an analytical treatment. This simplification is possible because, in this paper, we assume that the errors are *detected*; that is, the experimentalist knows precisely what sort of error has occurred because the environment that caused the errors is being continuously measured. Since the environment is thus acting as a part of the measurement apparatus, the errors it produces could be considered as measurement-induced errors.

There are a number of implementations in which measurement-induced errors of this sort may be significant. In the efficient linear optics scheme of Knill *et al.* [18], gates are implemented by nondeterministic teleportation. Failure of the teleportation corresponds to a gate error in which one of the qubits is measured in the computational basis with a known result. In a number of solid-state schemes, the readout device is always present and might make an accidental mea-

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surement of a qubit, even if the readout apparatus is in a quiescent state. An example is the use of rf single-electron transistors (SET) to readout a charge transfer event in the Kane proposal. Such a measurement is modeled as a weak continuous measurement [19]. While one supposes that the SET is biased in its low conductance state during qubit processing, it is useful to know that even if the device does accidentally make a measurement, the resulting error can be corrected.

In this paper, we show that for certain error models and codes, Markovian feedback plus an additional constant Hamiltonian (a “driving Hamiltonian”) can protect an unknown quantum state encoded in a particular codespace. Using the stabilizer formalism, we show that if there is one sort of error per physical qubit, and the error is detected perfectly, then it is always possible to store  $n-1$  logical qubits in  $n$  physical qubits. This works whether the detector record consists of discrete spikes (Poisson noise) or a continuous current (white noise). This suggests that if the dominant decoherence process can be monitored then using that information to control the system Hamiltonian may be the key to preventing such decoherence (see also the example in Ref. [20]).

As a salient application of this formalism, we consider the special case of spontaneous emission. Stabilizing states against spontaneous emission by using error-correcting codes has been studied by several groups [11–14]. Here we demonstrate that a simple  $n$ -qubit error-correcting code, Markovian quantum feedback, and a driving Hamiltonian are sufficient to correct spontaneous emissions on  $n-1$  qubits. The result of encoding  $n-1$  logical qubits in  $n$  physical qubits has been recently independently derived in Ref. [21] for the special case of spontaneous emission; however, our scheme differs in a number of respects. We also show that spontaneous emission error correction by feedback can be incorporated within the framework of canonical quantum error correction, which can correct arbitrary errors.

The paper is organized as follows. We review some useful results in quantum-error-correction and quantum-feedback theory in Sec. II. In Sec. III, we present the example of detected spontaneous-emission errors, first for two qubits and then for  $n$  qubits. In Sec. IV, we generalize this for protecting an unknown state subject to any single-qubit measurements. We show how to perform universal quantum computation using our protocol in Sec. V. Section VI concludes.

## II. BACKGROUND

### A. Quantum error correction: Stabilizer formalism

Quantum error correction has specifically been designed for protecting unknown quantum states [6,7,22]. An important class of quantum error-correcting codes is that of the stabilizer codes. An elegant and simple formalism [7] exists for understanding these codes; in this paper, we will restrict ourselves to this class of codes in order to take advantage of this formalism.

In the remainder of this paper, we will use the notation of Ref. [7] in which  $X$ ,  $Y$ , and  $Z$  denote the Pauli matrices  $\sigma_x$ ,

$\sigma_y$ , and  $\sigma_z$ , respectively, and juxtaposition denotes a tensor product; hence any element of the Pauli group,

$$P_n = \{\pm 1, \pm i\} \otimes \{I, X, Y, Z\}^{\otimes n}, \quad (2.1)$$

may be denoted as a concatenation of letters (e.g.,  $ZZI = \sigma_z \otimes \sigma_z \otimes I$ ).

A stabilizer code may be defined simply as follows: Consider a  $2^n$ -dimensional ( $n$ -qubit) Hilbert space and a subgroup of  $2^{n-k}$  commuting Pauli operators  $\mathcal{S} \in P_n$ . This group of operators is the *stabilizer* of the code; the codespace  $\mathcal{C}(\mathcal{S})$  is the simultaneous  $+1$  eigenspace of all the operators in  $\mathcal{S}$ . It can be shown that if  $-I$  is not an element of  $\mathcal{S}$ , the subspace stabilized is nontrivial and the dimension of  $\mathcal{C}(\mathcal{S})$  is  $2^k$ ; hence, we regard this system as encoding  $k$  qubits in  $n$ . The generators of such a group are a subset of this group such that any element of the stabilizer can be described as a product of generators. It is not hard to show that  $n-k$  generators suffice to describe the stabilizer group  $\mathcal{S}$ .

When considering universal quantum computation, it is also useful to define the *normalizer* of a code. Given a stabilizer group  $\mathcal{S}$ , the normalizer  $N(\mathcal{S})$  is the group of elements in  $P_n$  that commute with all the elements of  $\mathcal{S}$ , and it can be shown that the number of elements in  $N(\mathcal{S})$  is  $2^{n+k}$ .

Now,  $n+k$  generators suffice to describe  $N(\mathcal{S})$ . Of these,  $n-k$  can be chosen to be the generators of  $\mathcal{S}$ . It can be shown that the remaining  $2k$  generators can be chosen to be the *encoded operators*  $\bar{Z}_\mu, \bar{X}_\mu, \mu = 1, 2, \dots, k$ , where  $\bar{Z}_\mu, \bar{X}_\mu$  denote the Pauli operators  $X$  and  $Z$  acting on encoded qubit  $\mu$ , tensored with the identity acting on all other encoded qubits. These encoded operators act, as their name implies, to take states in  $\mathcal{C}(\mathcal{S})$  to other states in  $\mathcal{C}(\mathcal{S})$ .

The usual protocol for stabilizer codes, which will be modified in what follows, starts with measuring the stabilizer generators. This projection discretizes whatever error has occurred into one of  $2^{n-k}$  error syndromes labeled by the  $2^{n-k}$  possible outcomes of the stabilizer generator measurements. The information given by the stabilizer measurements about what error syndrome has occurred is then used to apply a unitary recovery operator that returns the state to the codespace.

A diagram of how such a protocol would be implemented in a physical system is given by the top portion of Fig. 1. After a state  $|\psi\rangle$  is encoded into a stabilizer subspace, as described above, an error may occur due to interaction with the environment. The encoded qubits are then entangled with a meter so that the error syndrome may be diagnosed through a direct measurement of the meter; the stabilizer formalism ensures that such a direct measurement does not disturb the state of the encoded qubits. Once the error syndrome has been diagnosed, a correcting unitary based on the particular error syndrome is applied to the encoded qubits to return the state to the codespace.

In this paper, we will use a modified version of this protocol. In particular, we will not entangle the encoded qubits with a meter in order to diagnose the error syndrome. Instead, we will assume that a limited class of errors occurs on the system and that these errors are detectable by measuring the environment directly; thus, we know when an error has

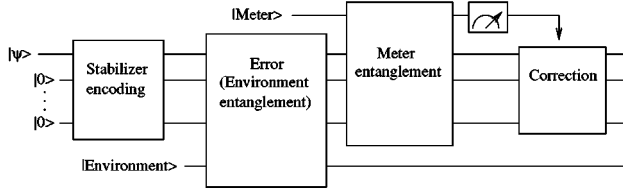
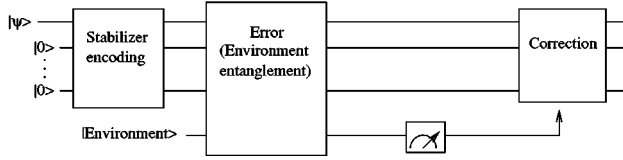
**Conventional protocol**

**Modified protocol**


FIG. 1. The top diagram shows the conventional stabilizer error-correction protocol. After the state is encoded, an error occurs through coupling with the environment. To correct this error, the encoded state is entangled with a meter in order to measure the stabilizer generators, and then feedback is applied on the basis of those measurements. The bottom diagram shows our modified protocol, in which the error and measurement steps are the same. To correct the error in this protocol, the environment qubits are measured, and we feedback on the results of the environment measurement.

happened and what the error is. The correction back to the codespace can still be performed by a unitary recovery operator based on the information from the error measurement. Figure 1 shows the difference between the conventional protocol and our modified protocol. This protocol is the same as that used in a less explicit form in Refs. [13,14].

In this paper, we will also consider operators of the form

$$T = T_1 \otimes \cdots \otimes T_n, \quad (2.2)$$

where  $T_i$  is an arbitrary traceless one-qubit operator, normalized such that its eigenvalues are  $\{-1, 1\}$ . Operators of this form are not generally Pauli-group stabilizers as presented in Ref. [7], as  $T$  is not, in general, a member of  $P_n$ . However, because of the special form of  $T$ ,  $T$  is equivalent to a Pauli operator up to conjugation by a unitary that is a product of one-qubit unitaries, i.e., there exists some  $U = \otimes_{i=1}^n U_i$  such that  $UTU^\dagger$  is a member of  $P_n$ . Therefore, choosing  $T$  as the sole stabilizer generator for a code is equivalent, up to conjugation by a unitary, to choosing a member of the Pauli group as the stabilizer generator. (Note that additional constraints are necessary if  $T$  is not the only stabilizer generator.)

**B. Quantum feedback**

Continuous quantum feedback can be defined, for the present purposes, as the process of monitoring a quantum system and using the continuous (in time) measurement record to control its dynamics. It can be analyzed by considering the dynamics of the measured system, conditioned on the continuous measurement record; this process is referred to as *unraveling*. The reduced dynamics of a system subject

to a weak continuous measurement is described by a Markov master equation, which determines the dynamics of the system averaged over all possible measurement records. However, if the time-continuous measurement record (a classical stochastic process) is known then it is possible to describe the conditional state of the measured system by a stochastic conditional evolution equation. A given master equation does not uniquely determine the conditional evolution equation, as there are many ways in which information about the system may be collected from the environment to which it is coupled as a result of the measurement; that is to say, a given master equation admits many unravelings.

In this section, we will introduce some of the results of this formalism; for more details, see Ref. [23]. We will assume that the change in the state of the system over a time interval  $dt$  due to its interaction with the environment can be described by a single jump operator  $c$ . By this we mean that jumps are represented by a Kraus operator  $\Omega_1 = c\sqrt{dt}$ , so that they occur with probability  $\langle c^\dagger c \rangle dt$ . Normalization requires another Kraus operator  $\Omega_0 = 1 - c^\dagger c dt/2 - iHdt$ , where  $H$  is Hermitian; then the unconditional master equation without feedback is just the familiar Lindblad form [24]

$$\begin{aligned} d\rho &= \Omega_0 \rho \Omega_0 + \Omega_1 \rho \Omega_1 - \rho \\ &= -i[H, \rho] dt + c \rho c^\dagger dt - \frac{1}{2} (c^\dagger c \rho + \rho c^\dagger c) dt \\ &\equiv -i[H, \rho] dt + \mathcal{D}[c] \rho dt. \end{aligned} \quad (2.3)$$

A bosonic example is given in Ref. [25], while a fermionic example is given in Ref. [26].

**1. Jump unravelings**

One way to unravel this master equation is to assume that the environment is measured so that the time of each jump event is determined. If the measured number of jumps up to time  $t$  is denoted  $N(t)$ , then the increment  $dN(t)$  is defined by

$$dN_c(t)^2 = dN_c(t), \quad (2.4)$$

$$E[dN_c(t)] = \langle c^\dagger c \rangle_c dt. \quad (2.5)$$

Here  $E[\ ]$  defines a classical ensemble average and the subscript  $c$  on the quantum average reminds us that the rate of the process at time  $t$  depends on the conditional state of the quantum system up to that time; that is to say, it depends on the state of the quantum system conditioned on the entire previous history of the current  $dN/dt$ . This conditional state is determined by a stochastic Schrödinger equation

$$\begin{aligned} d|\psi_c(t)\rangle &= \left[ dN_c(t) \left( \frac{c}{\sqrt{\langle c^\dagger c \rangle_c(t)}} - 1 \right) + dt \right. \\ &\quad \left. \times \left( \frac{\langle c^\dagger c \rangle_c(t)}{2} - \frac{c^\dagger c}{2} - iH \right) \right] |\psi_c(t)\rangle. \end{aligned} \quad (2.6)$$

We will refer to this as a *jump unraveling*. If we average over the measurement record to form  $\rho(t) = E[|\psi_c(t)\rangle\langle\psi_c(t)|]$ , it is easy to show using Eqs. (2.4) and (2.5) that  $\rho(t)$  obeys the unconditional master equation given in Eq. (2.3).

Now consider Markovian Hamiltonian feedback, where the strength of the feedback is a linear function of the measurement current,

$$H_{fb}(t) = \frac{dN(t)}{dt} V, \quad (2.7)$$

with  $V$  an Hermitian operator. Taking into account that the feedback must act after the measurement, it can be shown [8] that the feedback modifies the conditional evolution by changing the  $c$  in the numerator of the first term into  $e^{-iV}c$ . Since likewise changing all of the other occurrences of  $c$  has no effect, the ensemble average behavior is the same as before, with  $c$  changed to  $e^{-iV}c$ ; that is, to say, the feedback-modified master equation is

$$\dot{\rho} = -i[H, \rho] + \mathcal{D}[e^{-iV}c]\rho. \quad (2.8)$$

## 2. Diffusive unravelings

A very different unraveling may be defined by first noting that given some complex number  $\gamma = |\gamma|e^{i\phi}$ , we may make the transformation

$$c \rightarrow c + \gamma, \quad (2.9)$$

$$H \rightarrow H - \frac{i|\gamma|}{2}(e^{-i\phi}c - e^{i\phi}c^\dagger),$$

and obtain the same master equation. In the limit as  $|\gamma|$  becomes very large, the rate of the Poisson process is dominated by the term  $|\gamma|^2$ . In this case it may become impossible to monitor every jump process, and a better strategy is to approximate the Poisson stochastic process by a Gaussian white-noise process.

For large  $\gamma$ , we can consider the system for a time  $\delta t$  in which the system changes negligibly but the number of detections  $\delta N(t) \approx |\gamma|^2 \delta t$  is very large; then we can approximate  $\delta N(t)$  as [27]

$$\delta N(t) \approx |\gamma|^2 \delta t + |\gamma| \langle e^{-i\phi}c + c^\dagger e^{i\phi} \rangle_c \delta t + |\gamma| \delta W(t), \quad (2.10)$$

where  $\delta W(t)$  is normally distributed with mean zero and variance  $\delta t$ .

We now define the stochastic measurement record as the current

$$\frac{dQ(t)}{dt} = \lim_{\gamma \rightarrow \infty} \frac{\delta N(t) - |\gamma|^2 \delta t}{|\gamma| \delta t} \quad (2.11)$$

$$= \langle e^{-i\phi}c + e^{i\phi}c^\dagger \rangle_c + dW(t)/dt. \quad (2.12)$$

Given this stochastic measurement record, we can determine the conditional state of the quantum system by a stochastic Schrödinger equation analogous to Eq. (2.6). The equiva-

lence (in the ensemble average) to the master equation (2.3) is, in this case, easier to see by considering  $\rho_c = |\psi_c\rangle\langle\psi_c|$ , which obeys the stochastic master equation

$$d\rho_c(t) = -i[H, \rho_c(t)]dt + \mathcal{D}[e^{-i\phi}c]\rho_c(t)dt + \mathcal{H}[e^{-i\phi}c]\rho_c(t)dW(t). \quad (2.13)$$

In the above equations, the expectation  $\langle a \rangle_c$  denotes  $\text{tr}(\rho_c a)$ ,  $dW$  is a normally distributed infinitesimal random variable with mean zero and variance  $dt$  (a *Wiener increment* [28]), and  $\mathcal{H}$  is a superoperator that takes a jump operator as an argument and acts on density matrices as

$$\mathcal{H}[c]\rho = c\rho + \rho c^\dagger - \rho \text{tr}[c\rho + \rho c^\dagger]. \quad (2.14)$$

We thus have a different unraveling of the original master equation (2.3). Because of the white noise in the stochastic master equation (2.13), we call this a *diffusive unraveling*. It applies, for example, when one performs a continuous weak homodyne measurement of a field  $c$  by first mixing it with a classical local oscillator in a beam splitter and then measuring the output beams with photodetectors [27]. In this case, the measurement process  $dQ(t)$  determines the observed photocurrent. Another measurement model in which it may be appropriate to approximate the Poisson measurement process as a white-noise measurement process is the electronic point contact model for monitoring a single quantum dot [29,30]. In this case, the form of the master equation itself determines a large background jump rate, rather than an imposed classical field prior to detection.

We now consider Markovian feedback of the white-noise measurement record via a Hamiltonian, where the strength of the feedback is a linear function of the measurement current,

$$H_{fb}(t) = \frac{dQ(t)}{dt} F, \quad (2.15)$$

where  $F$  is a Hermitian operator. It can be shown that the addition of such feedback leads to the conditioned master equation [8,31]

$$\dot{\rho} = -i[(e^{i\phi}c^\dagger F + e^{-i\phi}F c)/2 + H, \rho] + \mathcal{D}[e^{-i\phi}c - iF]\rho + dW(t)\mathcal{H}[e^{-i\phi}c - iF]\rho. \quad (2.16)$$

In order to derive analytic results given such feedback, it is convenient to consider the average over many such a evolution trajectories. Since the expectation value of  $dW$  is zero, averaging yields an unconditioned master equation

$$\dot{\rho} = -i[(e^{i\phi}c^\dagger F + e^{-i\phi}F c)/2 + H, \rho] + \mathcal{D}[e^{-i\phi}c - iF]\rho. \quad (2.17)$$

Note that these equations are only valid for perfect (unit-efficiency) detection; the correspondences between error correction and feedback are more readily seen in this case, and we discuss the case of imperfect detection in Sec. IV D.

These feedback equations are easily generalized in the following way: Given  $n$  qubits, denote a set of measurement operators by  $\{c_1, c_2, \dots, c_n\}$ , where  $c_j$  acts on the  $j$ th qubit,

and a set of feedback operators by  $\{F_1, \dots, F_n\}$ , where the action of  $F_j$  is conditioned on the measurement of the  $j$ th qubit. Then the unconditional master equation (2.17), for example, generalizes to

$$\begin{aligned} \dot{\rho} = & \sum_{j=1}^n \{ -i[(e^{i\phi_j} c_j^\dagger F_j + e^{-i\phi_j} F_j c_j)/2 + H, \rho] \\ & + \mathcal{D}[e^{i\phi_j} c_j - iF_j] \rho \}. \end{aligned} \quad (2.18)$$

### III. EXAMPLE: SPONTANEOUS-EMISSION CORRECTION

A particular example of a Poisson process error is spontaneous emission, in which the jump operator is proportional to  $|0\rangle\langle 1|$ , so that the state simply decays from  $|1\rangle$  to  $|0\rangle$  at random times. Indeed, if the decay is observed (say by emitting a photon that is then detected), this may be regarded as a destructive measurement of the operator  $|1\rangle\langle 1|$ .

The process of stabilizing states against the important decay process of spontaneous emission through application of error-correcting codes has been studied by several groups [11–14]. In Ref. [12] Plenio, Vedral, and Knight considered the structure of quantum-error-correction codes and addressed the problem that spontaneous emission implies continuous evolution of the state even when no emission has occurred. They developed an eight-qubit code that corrects both one general error and the no-emission evolution to arbitrary order.

More recently, in several papers, Alber *et al.* [13,14] have addressed a somewhat more specific problem relating to spontaneous emission from statistically independent reservoirs. In this formulation, the only errors possible are spontaneous-emission errors, and the time and position of a particular spontaneous emission are known. They showed that given these constraints, a reduction of the redundancy in Ref. [12] was possible and constructed a four-qubit code that corrects for one spontaneous emission error.

Here we show that for the case considered in Refs. [13,14], a very simple error-correcting code consisting of just two qubits with feedback is sufficient to correct spontaneous emissions for a single logical qubit. A crucial difference from Refs. [13,14] is that we call for a constant driving Hamiltonian in addition to the feedback Hamiltonian. This driving Hamiltonian enables us to *correct* the no-emission evolution, rather than using an encoding that ensures that the no-emission evolution has no effect, as in Refs. [13,14]. This makes our encoding much more efficient. Indeed, a simple code of  $n$  qubits, with the appropriate feedback and driving Hamiltonians, can encode  $n-1$  qubits and correct for spontaneous emissions when the position (i.e., which qubit) and time of the jump are known. We also show that an equally effective protocol can be found for a diffusive unraveling of the spontaneous emission (as in homodyne detection).

#### A. Two-qubit code: Jump unraveling

The simplest system for which we can protect against detected spontaneous emissions is a system of two qubits.

We consider the model in which the only decoherence process is due to spontaneous emission from statistically independent reservoirs. We will show that a simple code, used in conjunction with a driving Hamiltonian, protects the codespace when the time and location of a spontaneous emission are known and a correcting unitary is applied instantaneously; the codespace suffers no decoherence.

The codewords of the code are given by the following:

$$\begin{aligned} |\bar{0}\rangle & \equiv (|00\rangle + |11\rangle)/\sqrt{2}, \\ |\bar{1}\rangle & \equiv (|01\rangle + |10\rangle)/\sqrt{2}. \end{aligned} \quad (3.1)$$

In the stabilizer notation, this is a stabilizer code with stabilizer generator  $XX$ . Both codewords are  $+1$  eigenstates of  $XX$ .

Following the presentation in Sec. II B, the jump operators for spontaneous emission of the  $j$ th qubit are

$$\Omega_j = \sqrt{\kappa_j dt} (X_j - iY_j) \equiv \sqrt{\kappa_j} dt a_j, \quad (3.2)$$

where  $4\kappa_j$  is the decay rate for this qubit. In the absence of any feedback, the master equation is

$$\dot{\rho} = \sum_{j=1,2} \kappa_j \mathcal{D}[X_j - iY_j] \rho - i[H, \rho]. \quad (3.3)$$

If the emission is detected, such that the qubit  $j$  from which it originated is known, it is possible to correct back to the codespace without knowing the state. This is because the code and error fulfill the necessary and sufficient conditions for appropriate recovery operations [6]:

$$\langle \psi_\mu | E^\dagger E | \psi_\nu \rangle = \Lambda_E \delta_{\mu\nu}. \quad (3.4)$$

Here,  $E$  is the operator for the measurement (error) that has occurred and  $\Lambda_E$  is a constant. The states  $|\psi_\mu\rangle, |\psi_\nu\rangle$  are the encoded states in Eq. (3.1) with  $\langle \psi_\mu | \psi_\nu \rangle = \delta_{\mu\nu}$ . These conditions differ from the usual condition only by taking into account that we *know* a particular error  $E = \Omega_j$  has occurred.

More explicitly, if a spontaneous emission on the first qubit occurs,  $|\bar{0}\rangle \rightarrow |01\rangle$  and  $|\bar{1}\rangle \rightarrow |00\rangle$ , and similarly for a spontaneous emission on the second qubit. Since these are orthogonal states, this fulfills the condition given in Eq. (3.4), so a unitary exists that will correct this spontaneous-emission error. One choice for the correcting unitary is

$$\begin{aligned} U_1 & = (XI - ZX)/\sqrt{2}, \\ U_2 & = (IX - XZ)/\sqrt{2}. \end{aligned} \quad (3.5)$$

As pointed out in Ref. [12], a further complication is the nontrivial evolution of the state in the time between spontaneous emissions. From Sec. II B, this is described by the measurement operator

$$\Omega_0 = II(1 - (\kappa_1 + \kappa_2)dt) - \kappa_1 dt ZI - \kappa_2 dt IZ - iHdt. \quad (3.6)$$

The nonunitary part of this evolution can be corrected by assuming a driving Hamiltonian of the form

$$H = -(\kappa_1 YX + \kappa_2 XY). \quad (3.7)$$

This result can easily be seen by plugging Eq. (3.7) into Eq. (3.6) with a suitable rearrangement of terms:

$$\begin{aligned} \dot{\Omega}_0 = & II(1 - (\kappa_1 + \kappa_2)dt) - \kappa_1 dt ZI(II - XX) \\ & - \kappa_2 dt IZ(II - XX), \end{aligned} \quad (3.8)$$

and since  $II - XX$  acts to annihilate the codespace,  $\Omega_0$  acts trivially on the codespace.

We then have the following master equation for the evolution of the system:

$$d\rho = \Omega_0 \rho \Omega_0^\dagger - \rho + dt \sum_{j=\{1,2\}} \kappa_j U_j a_j \rho a_j^\dagger U_j^\dagger, \quad (3.9)$$

where  $U_j$  is the recovery operator for a spontaneous emission from qubit  $j$ . From Sec. II B, these unitaries can be achieved by the feedback Hamiltonian

$$H_{fb} = \sum_{j=1,2} \frac{dN_j(t)}{dt} V_j, \quad (3.10)$$

where  $N_j(t)$  is the spontaneous emission count for qubit  $j$  and  $U_j = \exp(-iV_j)$ . Here, we can see from the simple form of Eq. (3.5) that  $V_j$  can be chosen as proportional to  $U_j$ . Since  $U_j a_j \rho a_j^\dagger U_j^\dagger$  acts as the identity on the codespace by definition, and since we have shown that  $\Omega_0 \rho \Omega_0^\dagger$  preserves the codespace, Eq. (3.9) must preserve the codespace.

Such a code is optimal in the sense that it uses the smallest possible number of qubits required to perform the task of correcting a spontaneous emission error, as we know that the information stored in one unencoded qubit is destroyed by a spontaneous emission.

### B. Two-qubit code: Diffusive unraveling

A similar situation applies for feedback of a continuous measurement record with white noise, as from homodyne detection of the emission. We use the same codewords, and choose  $\phi_j = -\pi/2$  for the measurement. Then Eq. (3.5) suggests using the following feedback operators:

$$\begin{aligned} F_1 &= \sqrt{\kappa_1} (XI - ZX), \\ F_2 &= \sqrt{\kappa_2} (IX - XZ). \end{aligned} \quad (3.11)$$

If we use these feedback Hamiltonians with the same driving Hamiltonian (3.7) as in the jump case, the resulting master equation is, using Eq. (2.18),

$$\dot{\rho} = \kappa_1 \mathcal{D}[YI - iZX]\rho + \kappa_2 \mathcal{D}[IY - iXZ]\rho. \quad (3.12)$$

We can see that this master equation preserves the codespace by again noting that  $YI - iZX = YI(II - XX)$ , and similarly for  $IY - iXZ$ . The operator  $II - XX$ , of course, acts to annihilate the codespace. This insight will be used in the following section to derive a feedback procedure for a more general measurement operator.

### C. Generalizations to $n$ qubits

We will now demonstrate a simple  $n$ -qubit code that corrects for spontaneous emission errors only, while encoding  $n-1$  qubits. Both of the above calculations (jump and diffusion) generalize. The master equation is the same as Eq. (3.3), but now the sum runs from 1 to  $n$ . Again we need only a single stabilizer generator, namely,  $X^{\otimes n}$ . The number of codewords is thus  $2^{n-1}$ , enabling  $n-1$  logical qubits to be encoded. Since it uses only one physical qubit in excess of the number of logical qubits, this is again obviously an optimal code.

First, we consider the jump case. As in Sec. III A, a spontaneous emission jump fulfills the error-correction condition (3.4) (see Sec. IV A below). Therefore, there exists a unitary that will correct for the spontaneous-emission jump. Additionally, it is easy to see by analogy with Eq. (3.8) that

$$H = \kappa_j \sum_j X^{\otimes j-1} Y X^{\otimes n-j} \quad (3.13)$$

protects against the nontrivial no-emission evolution. Therefore, the codespace is protected.

Next, for a diffusive unraveling, we again choose  $\phi_j = -\pi/2$ , as in Sec. III B. The same driving Hamiltonian (3.13) is again required, and the feedback operators generalize to

$$F_j = \sqrt{\kappa_j} (I^{\otimes j-1} X I^{\otimes n-j} + X^{\otimes j-1} Z X^{\otimes n-j}). \quad (3.14)$$

The master equation becomes

$$\dot{\rho} = \sum_j \kappa_j \mathcal{D}[I^{\otimes j-1} Y I^{\otimes n-j} (I^{\otimes n} - X^{\otimes n})]. \quad (3.15)$$

These schemes with a driving Hamiltonian do not have the admittedly desirable property of the codes given in Refs. [12–14] that if there is a time delay between the occurrence of the error and the application of the correction, the effective no-emission evolution does not lead to additional errors. Nevertheless, as pointed out in Ref. [14], the time delay for these codes must still be short so as to prevent two successive spontaneous emissions between corrections; they numerically show that the fidelity decays roughly exponentially as a function of delay time. Therefore, we believe that this drawback of our protocol is not significant.

### IV. ONE-QUBIT GENERAL MEASUREMENT OPERATORS

The form of the above example strongly indicates that there is a nice generalization to be obtained by considering stabilizer generators in more detail. In this section, we consider an arbitrary measurement operator operating on each qubit. We find the condition that the stabilizers of the codespace must satisfy. We show that it is always possible to find an optimal codespace (that is, one with a single stabilizer group generator). We work out the case of diffusive feedback in detail and derive it as the limit of a jump process.

### A. General unraveling

Different unravelings of the master equation (2.3) may be usefully parametrized by  $\gamma$ . In Sec. II B, we have seen that a simple jump unraveling has  $\gamma=0$ , while the diffusive unraveling is characterized by  $|\gamma|\rightarrow\infty$ . We will now address the question of when a unitary correction operator exists for an arbitrary  $\gamma$ , i.e., when a measurement scheme with a given  $\gamma$  works to correct the error.

Consider a Hilbert space of  $n$  qubits with a stabilizer group  $\{S_j\}$ . Let us consider a single jump operator  $c$  acting on a single qubit. We may then write  $c$  in terms of Hermitian operators  $A$  and  $B$  as

$$e^{-i\phi}c = \chi I + A + iB \quad (4.1)$$

$$\equiv \chi I + \vec{a} \cdot \vec{\sigma} + i\vec{b} \cdot \vec{\sigma}, \quad (4.2)$$

where  $\chi$  is a complex number,  $\vec{a}$  and  $\vec{b}$  are real vectors, and  $\vec{\sigma} = (X, Y, Z)^T$ .

We now use the standard condition (3.4), where we take  $E = c + \gamma$ . Henceforth,  $\gamma$  is to be understood as real and positive, since the relevant phase  $\phi$  has been taken into account in definition (4.1). The relevant term is

$$\begin{aligned} E^\dagger E &= (|\chi + \gamma|^2 + \vec{a}^2 + \vec{b}^2)I + \text{Re}(\chi + \gamma)A + \text{Im}(\chi + \gamma)*iB \\ &\quad + (\vec{a} \times \vec{b}) \cdot \vec{\sigma} \\ &\equiv (|\chi + \gamma|^2 + \vec{a}^2 + \vec{b}^2)I + D, \end{aligned} \quad (4.3)$$

where  $D$  is Hermitian.

Now we can use the familiar sufficient condition for a stabilizer code [7]: the stabilizer should anticommute with the traceless part of  $E^\dagger E$ . This condition becomes explicitly

$$0 = \{S, D\}. \quad (4.4)$$

As long as this is satisfied, there is some feedback unitary  $e^{-iV}$  that will correct the error.

Normalization implies that when  $E$  does not occur, there may still be nontrivial evolution. In the continuous time paradigm, where one Kraus operator is given by  $E\sqrt{dt}$ , the transform (2.9) tells us that the no-jump normalization Kraus operator is given by

$$\Omega_0 = 1 - \frac{1}{2}E^\dagger E dt - \frac{\gamma}{2}(e^{-i\phi}c - e^{i\phi}c^\dagger)dt - iHdt. \quad (4.5)$$

Now we choose the driving Hamiltonian

$$H = \frac{i}{2}DS + \frac{i\gamma}{2}(e^{-i\phi}c - e^{i\phi}c^\dagger). \quad (4.6)$$

This is a Hermitian operator because of (4.4). Then the total evolution due to  $\Omega_0$  is just the identity, apart from a term proportional to  $D(1-S)$ , which annihilates the codespace. Thus, for a state initially in the codespace, the condition (4.4) suffices for correction of both the jump and the no-jump evolution.

A nice generalization may now be found for a set  $\{c_j\}$  of errors such that  $c_j$  [with associated operator  $D_j$  as defined in Eq. (4.3)] acts on the  $j$ th qubit alone. Since  $D_j$  is traceless, it is always possible to find some other Hermitian traceless one-qubit operator  $s_j$  such that  $\{s_j, D_j\} = 0$ . Then we may pick the stabilizer group by choosing the single stabilizer generator

$$S = s_1 \otimes \cdots \otimes s_n, \quad (4.7)$$

so that the stabilizer group is  $\{1, S\}$ . As noted in Sec. II A, this is not strictly a stabilizer group, as  $S$  may not be in the Pauli group, but this does not change the analysis. Choosing  $H$  according to this  $S$ , such that

$$H = \sum_j \frac{i}{2}D_j S + \frac{i\gamma_j}{2}(e^{-i\phi_j}c_j - e^{i\phi_j}c_j^\dagger) \quad (4.8)$$

will, by our analysis above, provide a total evolution that protects the codespace, and the errors will be correctable; furthermore, this codespace encodes  $n-1$  qubits in  $n$ .

Note that we can now easily understand the  $n$ -qubit jump process error of spontaneous emission considered in Sec. III. Here,  $\gamma=0$ ,  $S = X^{\otimes n}$ , and  $D_j = 2\kappa_j Z_j$ . Thus, Eq. (4.4) is satisfied, and the Hamiltonian (3.13) is derived directly from Eq. (4.8).

Moreover, one is not restricted to the case of one stabilizer; it is possible to choose a different  $S_j$  for each individual error  $c_j$ . For example, for the spontaneous emission errors  $c_j = X_j - iY_j$ , we could choose  $S_j$  as different stabilizers of the five-qubit code. This choice is easily made, as the usual generators of the five-qubit code are  $\{XXXXI, IXZZX, XIXZZ, ZXIXZ\}$  [2]. For each qubit  $j$ , we may pick a stabilizer  $S_j$  from this set which acts as  $X$  on that qubit, and  $X$  anticommutes with  $D_j = Z_j$ . This procedure would be useful in a system where spontaneous emission is the dominant error process; it would have the virtue of both correcting spontaneous emission errors by means of feedback as well as correcting other (rarer) errors by using canonical error correction in addition.

We note that the work in this section can very easily be modified to generalize the results of Ref. [21]. This work has the same error model as ours: known jumps occurring on separate qubits so that the time and location of each jump are known; but Ref. [21] postulates fast unitary pulses instead of a driving Hamiltonian. Their scheme for spontaneous emission depends on applying the unitary  $X^{\otimes n}$  at intervals  $T_c/2$  that are small compared to the rate of spontaneous emission jumps. They show that after a full  $T_c$  period, the no-jump evolution becomes

$$U = e^{-iT_c H_c/2} X^{\otimes n} e^{-iT_c H_c/2} X^{\otimes n} = e^{-T_c/2} \sum_{i=1}^N \kappa_i 1. \quad (4.9)$$

Thus, the application of these pulses acts, as does our driving Hamiltonian, to correct the no-jump evolution. The generalization from spontaneous emission to general jump operator  $c_j$  for their case is simple: the code is the same as in the

above one-stabilizer protocol, with single stabilizer equal to Eq. (4.7). The fast unitary pulses are, in this case, also simply equal to Eq. (4.7).

### B. Diffusive unraveling

The case of white-noise feedback, where  $\gamma \rightarrow \infty$ , is easily treated by recalling the master equation (2.17) for white-noise measurement and feedback. It is clear that the first term in Eq. (2.17) can be eliminated by choosing the constant driving Hamiltonian

$$H = -(e^{i\phi}c^\dagger F + e^{-i\phi}Fc)/2, \quad (4.10)$$

which is automatically Hermitian. The problem then becomes choosing a feedback Hamiltonian  $F$  such that  $c - iF$  annihilates the codespace. The choice for  $F$  can be made simply by noting that if the codespace is stabilized by some stabilizer  $S$ , we can choose

$$F = B - iAS. \quad (4.11)$$

Now, note that the decoherence superoperator  $\mathcal{D}$  acts such that

$$\mathcal{D}[\chi I + L]\rho = \mathcal{D}[L]. \quad (4.12)$$

Then we know that  $\mathcal{D}[c - iF] = \mathcal{D}[\chi I + A(I - S)]$  annihilates the codespace.

The only caveat is that  $F$  is a Hamiltonian and therefore must be Hermitian. Then the choice (4.11) for  $F$  is only possible if the anticommutator of  $S$  and  $A$  is zero:

$$\{S, A\} = 0. \quad (4.13)$$

Therefore, if we are given the measurement operator  $e^{-i\phi}c = \chi + A + iB$ , we must choose a code with some stabilizer such that condition (4.13) applies; then it is possible to find a feedback and a driving Hamiltonian such that the total evolution protects the codespace.

At first glance, it may seem odd that the condition for feedback does not depend at all upon  $B$ . This independence has to do with the measurement unraveling; the diffusive measurement record (2.12) depends only upon  $e^{-i\phi}c + e^{i\phi}c^\dagger = 2(A + \chi)$ .

### C. Diffusion as the limit of jumps

It is instructive to show that the diffusive feedback process can be derived by taking the limit of a jump feedback process using the transformation (2.9). This takes several steps, and we use the treatment in Ref. [32] as a guide. But to begin, note that the condition (4.13) follows by considering Eq. (4.4) in the limit  $\gamma \rightarrow \infty$ , as the leading-order term in  $D$  is proportional to  $A$ .

Consider the jump unraveling picture with jump operator  $c + \gamma$  for large  $\gamma$  (but not infinite). Recall that in the error-correction picture, given in Sec. III, we postulated a feedback Hamiltonian  $(dN/dt)V$  that produces a unitary correction  $e^{-iV}$  that acts instantaneously after the jump. In addition, we will postulate a driving Hamiltonian  $K$  that acts when no

jump happens. In this picture, we will show that given the condition (4.13), it is possible to find asymptotic expressions for  $V$  and  $K$  so that the deterministic equation for the system preserves the stabilizer codespace. Finally, we will show that taking the limit  $\gamma \rightarrow \infty$  leads to the expressions for the feedback and driving Hamiltonians (4.10) and (4.11).

Let us consider the measurement operators for the unraveling with large  $\gamma$  and  $H = 0$ . Following Eq. (2.9), these are

$$\begin{aligned} \Omega_1 &= \sqrt{dt}(c + \gamma), \\ \Omega_0 &= 1 - \frac{dt}{2}[c\gamma - c^\dagger\gamma + (c + \gamma)^\dagger(c + \gamma)], \end{aligned} \quad (4.14)$$

where we have assumed, for simplicity, that  $\gamma$  is real. Now, including the feedback and driving Hamiltonians modifies these to

$$\begin{aligned} \Omega_1'(dt) &= \sqrt{dt}e^{-iV}(c + \gamma), \\ \Omega_0'(dt) &= e^{-iKdt}\Omega_0(dt) = 1 - iKdt - \frac{dt}{2}(c^\dagger c + 2\gamma c + \gamma^2). \end{aligned} \quad (4.15)$$

Following Ref. [32], expand  $V$  in terms of  $1/\gamma$  to second order:  $V = V_1/\gamma + V_2/\gamma^2$ , where the  $V_i$  are Hermitians. Then expanding the exponential in Eq. (4.15), we get to second order

$$\begin{aligned} \Omega_1'(dt) &= \sqrt{dt} \left[ 1 - i \left( \frac{V_1}{\gamma} + \frac{V_2}{\gamma^2} \right) - \frac{1}{2} \frac{V_1^2}{\gamma^2} \right] (A + iB + \gamma) \\ &= \sqrt{dt} \gamma \left[ 1 + \frac{\chi}{\gamma} + \frac{1}{\gamma} (A + iB - iV_1) \right. \\ &\quad \left. + \frac{1}{\gamma^2} (V_1^2/2 - iV_2 - i(A + iB)V_1) \right]. \end{aligned} \quad (4.16)$$

A reasonable choice for  $V_1$ , by analogy to Eq. (4.11), is  $B - iAS$ . Following Ref. [32], we also use Eqs. (4.10) and (4.11) to choose  $V_2$  and  $K$ ; note that Eq. (4.11) is exactly the expression we would expect for  $K$  from Eq. (4.6) in the limit as  $\gamma$  is taken to infinity. We will proceed to show that the choice for  $V$  and  $K$ ,

$$V_1 = B - iAS, \quad (4.17)$$

$$V_2 = -(c^\dagger F + Fc)/2, \quad (4.18)$$

$$K = -\gamma(B - iAS), \quad (4.19)$$

leads to the correct evolution to second order in  $\gamma$ .

Now, the deterministic evolution is given by

$$d\rho = \Omega_0'\rho\Omega_0' + \Omega_1'\rho\Omega_1' - \rho. \quad (4.20)$$

Substituting Eqs. (4.15)–(4.19) into Eq. (4.20) to second order in  $\gamma$ , after some algebra, gives the deterministic jump equation

$$d\rho = \mathcal{D}[A(1-S)]\rho, \quad (4.21)$$

which of course acts as zero on the codespace.

Now we will show that taking the limit as  $\gamma \rightarrow \infty$  leads to the feedback operators given in Eqs. (4.10) and (4.11). We saw in Eqs. (4.17) and (4.18) that the feedback Hamiltonian needed to undo the effect of the jump operator  $c + \gamma$  was just

$$H_{fb} = \frac{dN(t)}{dt} \left( \frac{B - iAS}{\gamma} - \frac{c^\dagger F + Fc}{2} \right). \quad (4.22)$$

Keeping terms of two orders in  $\gamma$  gives

$$H_{fb} = \gamma(B - iAS) - \frac{c^\dagger F + Fc}{2} + \frac{dN(t) - \gamma^2 dt}{\gamma dt} (B - iAS). \quad (4.23)$$

The last term just becomes the current  $\dot{Q}(t)$  as  $\gamma$  approaches infinity, as in Eq. (2.11). Furthermore, we have not yet added the driving Hamiltonian to the expression for the feedback. Doing so yields

$$H_{\text{total}}(t) = H_{fb} + K = \dot{Q}(t)(B - iAS) - \frac{c^\dagger F + Fc}{2}, \quad (4.24)$$

which is just what we obtained in the preceding section. Thus, we can see that this continuous current feedback can be thought of as an appropriate limit of a jump plus unitary correction process.

#### D. Imperfect detection

These results for feedback were obtained by assuming unit efficiency, i.e., perfect detection. Realistically, of course, the efficiency  $\eta$  will be less than unity. This results in extra terms in the feedback master equations we have derived [8]. In the jump case, the extra term is

$$\dot{\rho} = (1 - \eta) \sum_j (c_j \rho c_j^\dagger - U_j c_j \rho c_j^\dagger U_j^\dagger). \quad (4.25)$$

In the diffusion case, it is

$$\dot{\rho} = \frac{1 - \eta}{\eta} \sum_j \mathcal{D}[F_j]. \quad (4.26)$$

In both the cases, this results in the exponential decay of coherence in the codespace. This is because the error-correction protocol here relies absolutely upon detecting the error when it occurs. If the error is missed (jump case), or imperfectly known (diffusion case), then it cannot be corrected. This behavior is, of course, a property of any continuous-time error-correction protocol that depends on correcting each error instantaneously (e.g., Refs. [13,14,21]).

On the other hand, such behavior for the Markovian feedback protocol is in contrast to the state-estimation procedure used in Ref. [15]. The latter is much more robust under non-unit efficiency; indeed, given nonunit efficiency, it still works to protect an unknown quantum state without exponential

loss [33]. This difference in performance occurs because state estimation is a function of the entire measurement record, not just instantaneous measurement results, and thence does not propagate errors to the same extent that the Markovian feedback system does. Thus, we can see that there is a certain trade-off. Our Markovian feedback scheme relies upon calculational simplicity, but at the expense of robustness. The state-estimation procedure, conversely, is designed to be robust, but at the cost of computational complexity.

#### V. UNIVERSAL QUANTUM GATES

Given a protected code subspace, one interesting question, as in Ref. [21], is to investigate what kinds of unitary gates are possible on such a subspace. For universal quantum computation on the subspace—the ability to build up arbitrary unitary gates on  $k$  qubits—it suffices to be able to perform arbitrary one-qubit gates for all  $k$  encoded qubits and a two-qubit entangling gate such as the controlled-NOT gate for all encoded qubits  $\mu, \nu$ . Indeed, as is noted in Ref. [21], it is enough to be able to perform the Hamiltonians  $\bar{X}_\mu, \bar{Z}_\mu$ , and  $\bar{X}_\mu \bar{X}_\nu$  for all  $\mu, \nu$  [34]. We will demonstrate that performing these Hamiltonians with our protocol is possible for the spontaneous emission scheme given in Sec. III, and then we will show how that construction generalizes for an arbitrary jump operator.

Recall that the example in Sec. III has single stabilizer  $X^{\otimes n}$  and encodes  $n - 1$  logical qubits in  $n$  physical qubits. To find the  $2(n - 1)$  encoded operations, we must find operators that together with the stabilizer generate the normalizer of  $X^{\otimes n}$  [2]. In addition, if these operators are to act as encoded  $X$  and  $Z$  operations, they must satisfy the usual commutation relations for these operators:

$$\begin{aligned} \{X_\mu, Z_\mu\} &= 0, \\ [X_\mu, Z_\nu] &= 0, \quad \mu \neq \nu, \\ [X_\mu, X_\nu] &= [Z_\mu, Z_\nu] = 0. \end{aligned} \quad (5.1)$$

Operators satisfying these constraints are easily found for this code:

$$\begin{aligned} \bar{X}_\mu &= I^{\otimes \mu - 1} X I^{\otimes n - \mu}, \\ \bar{Z}_\mu &= I^{\otimes \mu - 1} Z I^{\otimes n - \mu - 1} Z, \\ \bar{X}_\mu \bar{X}_\nu &= I^{\otimes \mu - 1} X I^{\otimes \nu - \mu - 1} X I^{\otimes n - \nu}, \end{aligned} \quad (5.2)$$

where we assume  $1 \leq \mu < \nu \leq n - 1$ . If we apply a Hamiltonian  $H_{enc}$ , given by any linear combination of the operators in (5.2), the resulting evolution is encapsulated in the expression for  $\Omega_0$ , from Eq. (3.6),

$$\Omega_0 = \left( 1 - \sum_j \kappa_j dt \right) 1 - \sum_j \kappa_j Z_j (1 - X^{\otimes n}) dt - i H_{enc} dt. \quad (5.3)$$

As the first term is proportional to the identity and the second term acts as zero on the codespace, the effective evolution is

given solely by  $H_{enc}$  as long as the state remains in the codespace under that evolution. But because the encoded operations are elements of the normalizer, as we saw in Sec. II A, applying  $H_{enc}$  does not take the state out of the codespace. Furthermore, our protocol assumes that the spontaneous emission jumps are corrected immediately and perfectly, so jumps during the gate operation will also not take the state out of the codespace. Thus, we can perform universal quantum computation without having to worry about competing effects from the driving Hamiltonian.

The generalization to the scheme, given in Sec. IV A, to encode  $n-1$  logical qubits in  $n$  physical qubits is easily done. First, we note that for the stabilizer  $S$  given in the general scheme, we know that

$$S = UX^{\otimes n}U^\dagger \quad (5.4)$$

for some unitary  $U = \otimes_{i=1}^n U_i$ , so the encoded operations for that code are similarly given by

$$\begin{aligned} \bar{X}_\mu &= I^{\otimes \mu-1} U_\mu X U_\mu^\dagger I^{\otimes n-\mu}, \\ \bar{Z}_\mu &= I^{\otimes \mu-1} U_\mu Z U_\mu^\dagger I^{\otimes n-\mu-1} U_n Z U_n^\dagger, \\ \bar{X}_\mu \bar{X}_\nu &= I^{\otimes \mu-1} U_\mu X U_\mu^\dagger I^{\otimes \nu-\mu-1} U_\nu X U_\nu^\dagger I^{\otimes n-\nu}. \end{aligned} \quad (5.5)$$

Now, from Eq. (4.6), we can see that the generalization of Eq. (5.3) is

$$\Omega_0 = (1 - f dt) 1 - \sum_j g_j D_j (1 - S) dt - i H_{enc} dt, \quad (5.6)$$

for real numbers  $f, g_i$  given by expanding the expression (4.5). Again, since  $D(1-S)$  annihilates the codespace, the effective evolution is given solely by  $H_{enc}$  as long as the

state remains in the codespace under this evolution. Again,  $H_{enc}$  is made up of normalizer elements, which do not take the state out of the codespace; and again jumps that occur while the gate is being applied are immediately corrected and thus do not affect the gate operation. Therefore, universal quantum computation is possible under our general scheme.

## VI. CONCLUSION

We have shown that it is possible to understand a particular variant of quantum control as quantum error correction. This method is very general in that it can correct any single qubit detected errors, while requiring only  $n$  physical qubits to encode  $n-1$  logical qubits. As a particular example, we have shown how to correct for a spontaneous emission evolution using a feedback and a driving Hamiltonian, which allows less redundancy than has previously been obtained. We have additionally shown that universal quantum computation is possible under our method.

We expect that this work will provide a starting point for practically implementable feedback schemes to protect unknown states. The fact that only two qubits are required for a demonstration should be particularly appealing. We also expect a more complete theoretical development. Fruitful avenues for further research include applying notions of fault tolerance to this sort of quantum control.

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