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Movassagh, R. and Ouyang, Y. orcid.org/0000-0003-1115-0074 (Submitted: 2020) Constructing quantum codes from any classical code and their embedding in ground space of local Hamiltonians. [Preprint - arXiv] (Submitted)

https://doi.org/10.48550/arXiv.2012.01453

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## Constructing quantum codes from any classical code and their embedding in ground space of local Hamiltonians

Ramis Movassagh \* Yingkai Ouyang †

#### Abstract

We introduce a framework for constructing a quantum error correcting code from any classical error correcting code. This includes CSS codes [CS96, Ste96b] and goes beyond the stabilizer formalism [Got96] to allow quantum codes to be constructed from classical codes that are not necessarily linear or self-orthogonal (Fig. 1). We give an algorithm that explicitly constructs quantum codes with linear distance and constant rate from classical codes with a linear distance and rate. As illustrations for small size codes, we obtain Steane's 7-qubit code [Ste96a] uniquely from Hamming's [7,4,3] code [MS77], and obtain other error detecting quantum codes from other explicit classical codes of length 4 and 6. Motivated by quantum LDPC codes [BBA+15] and the use of physics to protect quantum information, we introduce a new 2-local frustration free quantum spin chain Hamiltonian whose ground space we analytically characterize completely. By mapping classical codewords to basis states of the ground space, we utilize our framework to demonstrate that the ground space contains explicit quantum codes with linear distance. This side-steps the Bravyi-Terhal no-go theorem [BT09] because our work allows for more general quantum codes beyond the stabilizer and/or linear codes. We hesitate to call this an example of subspace quantum LDPC code with linear distance.

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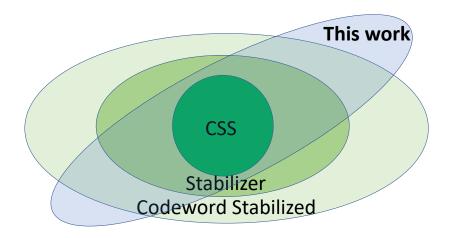


Figure 1: Comparison of the codes attainable within this work with Calderbank-Shor-Steane (CSS) [CS96, Ste96b], quantum stabilizer [Got96, CRSS97], and codeword-stabilized (CWS) [CSSZ08] codes. Inclusion of stabilizers and CWS is not strict as our codewords are supported on disjoint sets.

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## 1 Overview

Error correction is a necessary part of any reliable computation. On one hand, one can protect against errors by designing error correcting codes that allow for reliable recovery of the encoded information. On the other hand, nature herself provides innate resources for the protection and correction of information. These beg the questions:

- 1. Since quantum computers generalize classical computers, to what extent can one import the remarkable discoveries in classical coding [MS77] to the quantum realm?
- 2. Since the bulk of matter often resides in its ground state, are there physical quantum systems (e.g., 2-local Hamiltonians) whose ground states can sustain good quantum codes?

In this paper we address both of these questions in two parts:

Part 1 introduces a framework that takes any classical code and algorithmically constructs

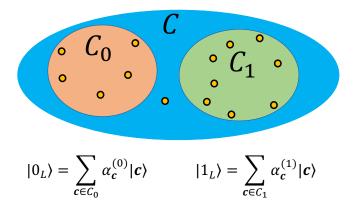


Figure 2: The encoding of one logical qubit from two disjoint subsets of a classical code C.

an explicit quantum code. The challenge in designing quantum codes is to not only correct the bit-flip errors, but also to correct phase-flip errors. We construct quantum codes with asymmetric distances given by  $d_X$  and  $d_Z$  for the bit-flip and phase-flip errors respectively. Similar to CSS codes, the logical codewords that are supported on codewords of length nare labeled by a *classical* code C. In the simplest case, we encode a logical qubit as follows

$$|0_L\rangle = \sum_{\mathbf{c} \in C_0} \alpha_{\mathbf{c}}^{(0)} |\mathbf{c}\rangle, \quad |1_L\rangle = \sum_{\mathbf{c} \in C_1} \alpha_{\mathbf{c}}^{(1)} |\mathbf{c}\rangle,$$
 (1)

where  $C_0$  and  $C_1$  are disjoint subsets of C, and each c is a product state. A priori, the subsets  $C_0$  and  $C_1$  and the coefficients are not known. By mapping the Knill-Laflamme (KL) quantum error correction criteria [KL97] to a linear algebra problem, we construct an explicit algorithm that can determine these unknowns. To design a quantum code that encodes more than one logical qubit, we provide a recursive algorithm. This algorithm finds the disjoint subsets of C and enforces the KL criteria by solving for feasible solutions of a linear program. The algorithm outputs the logical states:

$$|j_L\rangle = \sum_{\mathbf{c}\in C_i} \alpha_{\mathbf{c}}^{(j)} |\mathbf{c}\rangle, \quad j \in \{0, 1, \dots, M-1\},$$
 (2)

where  $M \leq q^{cn}$  with 0 < c < 1 resulting in quantum codes with constant rates. We prove (see Theorem 1):

**Theorem.** Take a classical code C of length n on a q-ary alphabet with the distance of  $d_X$ . Let  $V_q(r)$  be the Hamming ball of radius r. If  $|C| \geq 2V_q(d_Z - 1)$ , then Alg. 1 and Alg. 2 in the paper explicitly derive  $2 \leq M \leq q^{cn}$ , with 0 < c < 1 quantum logical states in (2) with a bit- and phase-flip distances of  $d_X$  and  $d_Z$  respectively. The overall distance is  $\min(d_X, d_Z)$ .

Our algorithm has a recursive structure. When the inequality in the theorem is satisfied, the algorithm always constructs two logical codewords. The third and subsequent logical codewords are found recursively by determining the feasibility of a sequence of linear programs.

This recursive algorithm succeeds with high probability and almost surely over random codes. In the rare case that the set of linear constraints have rows that are all non-zero

and are of the same sign, the recursion becomes infeasible and algorithm halts. We find that even if this happens we can always construct *Approximate* Quantum Error Correcting Codes (AQECC) [LNCY97] with linear distance and constant expected rate, provided that the underlying classical code has these parameters. The trade-off is between the expected number of logical states and the approximation error as given by Eq. (25).

In addition to giving asymptotic results, this formalism works equally well to construct finite size codes. To illustrate, we show examples of quantum codes that can be constructed from classic codes in the classical literature. Among the various, we find that the Steane code is the unique solution of our formalism when the input is Hamming's [7,4,3] code. The formalism introduced here is shown in relation to other known formalism in Fig. 1.

Part 2 focuses on the physics of information. Since topological models of quantum computation, it has been recognized that quantum codes may naturally appear in the ground space of physical systems [Kit03]. Most physical are 2-local interactions, and many have investigated the encoding of quantum codes in their ground spaces.

The most celebrated example is Kitaev's toric code that resides in the ground space of a 4-local Hamiltonian, which is an effective Hamiltonian of a perturbed 2-local Hamiltonian [Kit06]. Kitaev's toric code is an example of topological order and paved the way for the topological model of quantum computation. The compass model [DBM05] is 2-local on a lattice, and has recently been proposed as a candidate to encode quantum codes in the eigenbasis of the Hamiltonian [LMN+19]. However, the performance of these quantum codes are not well understood and have mainly been numerically investigated. The advantage of our work over the aforementioned works is that we construct a 2-local Hamiltonian, whose quantum error correcting properties we analytically prove.

In a nice recent result, Brandao et al. [BaCimcbuB19] gave a non-constructive proof of the existence, with high probability, of AQECCs within the low-energy sector for a multitude of translation invariant quantum spin chains including ferromagnetic Heisenberg model and spin-1 Motzkin spin chain [BaCimcbuB19]. The challenges that remained were that the codes were not explicit, the quantum error correction criteria was only approximately satisfied (hence AQECCs), errors had to be on consecutive set of spins, and the codes were in a low-energy sector of the local Hamiltonian (i.e., not the ground space). Moreover, the distance of the code grows logarithmically with the number of spins.

Our work overcomes these challenges. We construct explicit codes with linear distance that encode one logical qubit (we could have easily encoded a qudit as well). We write down a new and explicit 2-local quantum integer spin-s chain parent Hamiltonian,  $H_n$ , on n qudits. We analytically prove that its ground space can be spanned by product states. By mapping these product states to classical codewords, we reduce the problem of finding quantum codes in the ground space of our Hamiltonian to that of finding classical codes that must obey some constraints that are induced by the Hamiltonian. The classical coding problem becomes that of finding q-ary codes with forbidden sub-strings. By leveraging on existing constructions of binary codes, we construct candidate classical codes for our algorithm to run. See Fig. 3 for a comparison.

Properties:	Brandao et al (PRL 2018)	This work
QECC	Approximate with $\varepsilon = O(N^{-1/8})$	Exact
Distance d	$d = \Omega(\log(N))$	$d = \Theta(N)$
Rate	Vanishes	Vanishes
Error restriction	Consecutive spins	None
Code space	Low-energy eigenstates	Exact ground state
Translation invariance required?	Yes	No

Figure 3: Comparison of the explicit codes constructed here with previous work of Brandao et al [BaCimcbuB19]

#### The Hamiltonian and its ground space

Let us consider a spin chain of length n with open boundary conditions and the local Hilbert space dimension of 2s+1, where  $s \geq 1$  is a positive integer. We take a representation in which  $|j\rangle$  denotes the  $s_z = j$  state of a spin-s particle, such that  $\hat{S}_z|j\rangle = j$   $|j\rangle$  where  $j \in \{0, \pm 1, \pm 2, \cdots, \pm s\}$ .

The local Hamiltonian whose ground space contains the quantum code is  $H_n = H_n^J + H_n^s$ , where  $H_n^J = J \sum_{k=1}^n (|0\rangle\langle 0|)_k$ . The Hamiltonian  $H_n^s$ , is defined by

$$H_n^s = \sum_{k=1}^{n-1} \left\{ \sum_{m=-s}^s P_{k,k+1}^m + \sum_{m=1}^s Q_{k,k+1}^m \right\} , \qquad (3)$$

and the local terms are projectors acting on two neighboring spins k, k+1 are

$$P^m = |0 \leftrightarrow m\rangle\langle 0 \leftrightarrow m| , \qquad Q^m = |00 \leftrightarrow \pm m\rangle\langle 00 \leftrightarrow \pm m| ,$$
 (4)

where  $|0 \leftrightarrow m\rangle \equiv \frac{1}{\sqrt{2}}[|0, m\rangle - |m, 0\rangle]$ ,  $|00 \leftrightarrow \pm m\rangle \equiv \frac{1}{\sqrt{2}}[|0, 0\rangle - |m, -m\rangle]$ , and we denoted by  $|j, k\rangle$  the spin state  $|s_k^z = j, s_{k+1}^z = k\rangle$ . We will be mostly interested in s > 1.

Since  $H_n^s$  is free of the sign problem (i.e., stoquastic), the local projectors define an effective Markov chain, which have the following correspondence:

Local Projector	Local moves	Interpretation
$P^m$	$0m \longleftrightarrow m0$	Spin transport: local exchange of spin $m$ with 0
$Q^m$	$00 \longleftrightarrow m, -m$	Spin interaction: local creation/annihilation of $m, -m$

We prove that the ground state degeneracy is exponentially large in the number of spins:

$$\dim(\ker(H_n^s)) = \frac{\left(-2 + r_+^{n+1} + r_-^{n+1}\right)}{2(s-1)} \approx \frac{r_+^{n+1}}{2(s-1)} \quad , \quad n \gg 1$$

where  $r_{\pm} \equiv (1 \pm \sqrt{1 - 1/s})$ , and  $\ker(M)$  the kernel of the operator M. Let us denote by  $\operatorname{Ent}_q : [0,1] \to [0,1]$  the q-ary entropy function defined by  $\operatorname{Ent}_q(x) = -x \log_q x - (1 - x) \log_q (1-x) + x \log_q (q-1)$ . Our main theorem is (this is Theorem 2 in the paper):

**Theorem.** Let  $0 < \tau \le 1/2$  be a real and positive constant. There exist quantum codes in  $\ker(H_n)$  that encode one logical qubit and have the distance of  $2\tau n$  whenever

$$\operatorname{Ent}_2(2\tau) + \operatorname{Ent}_{2s+1}(2\tau) \log_2(2s+1) + o(1) \le 1.$$

Second, there are explicit quantum codes which encode one logical qubit with a distance of  $2\tau n$  whenever

$$1/2 - \tau/0.11 \ge \log_2(2s+1)\operatorname{Ent}_{2s+1}(2\tau).$$

**Remark.** We call the constructions given by optimizing (51) and (52) as the Gilbert-Varshamov (GV) [MS77, Chpt. 1] and Justesen construct [MS77, Chpt. 10, Thm. 11] respectively. The GV construct arises from choosing a random C, while the Justesen construct uses the classical Justesen code to define C.

This side-steps the Bravyi-Terhal no-go theorem [BT09] because our work allows for more general quantum codes beyond the stabilizer and/or linear codes. This model may be called an example of *subspace* quantum LDPC codes with linear distance.

This work could pave the way for constructing the first Quantum LDPC codes with linear distance in the ground space of translation invariant local spin chains. We note that had we used *all* of the ground space to construct the codes, this Hamiltonian could have made the case for the first example of topological order in one-dimension, which has been conjectured to be impossible. The practical advantage of our work is that such explicit Hamiltonians are easily constructed in the laboratory in the near term, especially in atomic or ion trap architectures. Lastly, the Hamiltonian is a generalization of the highly entangled colored Motzkin spin chain [MS16], which may be of independent interest.

## 1.1 Discussions and open problems

This paper provides a rigorous framework for the systematic construction of a quantum codes from any classical code. We illustrate the theory through a series of examples and proved that new quantum codes with linear distance and constant rate can be constructed using this work. Our formalism encapsulates the CSS formalism, and has an intersection with stabilizer and codeword-stabilized (CWS) formalisms (see subsection 2.4.4). However, there are codes inside the stabilizer and CWS that our formalism does not capture. For example, the five-qubit code is not covered by our formalism because its logical codewords cannot be written as superpositions over disjoint computational bases. Alg. 2 can construct logical states beyond a logical qubit. The relation of our work to the previous is faithfully depicted in the Venn diagram (Fig. 1) shows.

An open problem is whether starting from a classical linear, self-orthogonal, and binary code, do we always get a CSS code? Another open problem would be to find an alternative way of getting at our logical qudit construction by directly using the (high-dimensional) kernel of the matrix A. It would also be interesting to see an extension of our formalism to

encompass all stabilizer codes, and indeed, any arbitrary quantum code. With the exception of permutation invariant codes (subsection 2.4.5), most of the analysis herein takes the classical codewords and uses them to define a product basis over which the logical quantum states are defined. The extension of our results to include non-product basis for the logical codewords calls for further investigation.

Motivated by the physical implementation of quantum error correcting codes, we gave a local, frustration-free, Hamiltonians whose ground space has a quantum code with linear distance as a subspace. The model we have is translation-invariant but that is not essential and can easily be relaxed. A major open problem has been to construct 'good' quantum LDPC (QLDPC) codes, where 'good' means that the quantum code must have a linear distance. To the best of our knowledge, there is no unique definition for QLDPC codes besides that they should be in the ground space of local Hamiltonians. To this end, one may call our work a 'good' subspace QLDPC code. However, it would have been more satisfactory if the code is all of the ground space. It is our hope that this work might help in eventually realizing a good QLDPC code.

Another interesting problem is to find a Hamiltonian whose ground space is quantum code with a macroscopic distance, and the local and global ground states satisfy a consistency criterion as defined in [BH11]. This would then serve as the first example of topological quantum order in one-dimension. We would have found it easy to prove a gap above the degenerate ground space of the local Hamiltonian herein; however, since the code occupies a subspace of the ground space one would need to prove a 'local gap' lower bound. This means that in order to move from a subspace of the ground space to another subspace an operator with a large support needs to be applied.

## 2 Part 1: Explicit quantum codes from classical codes

## 2.1 Constructing a logical qubit with linear distance

In this section, we want to design q-ary quantum codes with bit-flip distance of  $d_X$  and a phase-flip distance of  $d_Z$  using a q-ary classical code  $C \subset \{0, 1, \dots, q-1\}^n$  as an input. The minimum distance of the quantum code is then  $d = \min(d_X, d_Z)$ .

To correct errors on q-ary quantum codes, we consider errors in the generalized Pauli basis. Let us denote by  $\omega$  the primitive root of unity  $\omega \equiv \exp(2\pi i/q)$ . The Z and X type Pauli matrices are respectively given by

$$Z = \sum_{j=0}^{q-1} \omega^j |j\rangle\langle j| , \quad X = \sum_{j\in\mathbb{Z}_q} |j\rangle\langle j+1| .$$
 (5)

And any Pauli operator is equivalent to  $X^aZ^b$  up to a phase for some  $a, b = 0, \ldots, q - 1$ . We consider the set Pauli operators on n qudits  $\mathcal{P}_n = \{X^aZ^b : a, b = 0, \ldots, q - 1\}^{\otimes n}$  that span the space of linear operators on n qudits. Given any Pauli in  $\mathcal{P}_n$  that has the form  $P = X^{a_1}Z^{b_1} \otimes \cdots \otimes X^{a_n}Z^{b_n}$ , we denote  $\operatorname{wt}_X(P) = \operatorname{wt}(\mathbf{a})$  and  $\operatorname{wt}_Z(P) = \operatorname{wt}(\mathbf{b})$ , where  $\mathbf{a} = (a_1, \ldots, a_n)$ ,  $\mathbf{b} = (b_1, \ldots, b_n)$ , and  $\operatorname{wt}(\cdot)$  denotes the Hamming weight of the vector. It

is easy to see that for any non-negative integer  $r \leq n$ , we have  $\sum_{w=0}^{r} |\mathcal{Z}_w| = V_q(r)$ , where

$$V_q(r) = \sum_{w=0}^{r} \binom{n}{w} (q-1)^w$$

denotes the volume of the q-ary Hamming ball of radius r.

For the KL criteria to hold for a quantum code with logical codewords  $|0_L\rangle$  and  $|1_L\rangle$  on n qubits with a bit-flip distance of  $d_X$  and a phase-flip distance of  $d_Z$ , it suffices to require that for all generalized Pauli matrices P such that  $\operatorname{wt}_X(P) \leq d_X - 1$  and  $\operatorname{wt}_Z(P) \leq d_Z - 1$ , these hold:

$$\langle 0_L | P | 0_L \rangle = \langle 1_L | P | 1_L \rangle \tag{6}$$

$$\langle 0_L | P | 1_L \rangle = 0. (7)$$

Eqs. (6) and (7) are the *non-deformation* and *orthogonality* conditions respectively. If we demand that

- 1.  $\operatorname{dist}(C) \geq d_X$ : the minimum distance of C is at least  $d_X$
- 2. Supp $(0_L) \cap \text{Supp}(1_L) = \emptyset$ : (the logical codewords  $|0_L\rangle$  and  $|1_L\rangle$  are supported on distinct codewords in C),

then the orthogonality condition (Eq. (7)) trivially holds. To verify the non-deformation condition (Eq. (6)), we note that  $\langle 0_L|P|0_L\rangle = \langle 1_L|P|1_L\rangle = 0$  whenever P is not diagonal. Hence the only non-trivial cases to be verified are the diagonal generalized Pauli operators, where the set of diagonal Pauli operators of weight w is

$$\mathcal{Z}_w = \{ Z^{z_1} \otimes \cdots \otimes Z^{z_n} : \operatorname{wt}(\mathbf{z}) = w \}.$$
 (8)

In general, for any diagonal Pauli operator P, the expectations  $\langle 0_L|P|0_L\rangle$  and  $\langle 1_L|P|1_L\rangle$  are complex numbers. This is in contrast to the case where P are Kraus operators of the amplitude damping channel, in which all such expectations are real, or when P are diagonal operators and q=2. For Eq. (6) to hold, the following has to hold for all Paulis P with a weight at most d-1:

$$\operatorname{Re}(\langle 0_L | P | 0_L \rangle) - \operatorname{Re}(\langle 1_L | P | 1_L \rangle) = 0$$
  
$$\operatorname{Im}(\langle 0_L | P | 0_L \rangle) - \operatorname{Im}(\langle 1_L | P | 1_L \rangle) = 0,$$

The quantum code we define depends on a "balanced" real non-zero column vector  $\mathbf{x} = (x_1, x_2, \dots, x_m)^T$  in the sense that

$$\sum_{i=1}^{m} x_i = 0.$$

Let  $x_k^+ = \max\{x_k, 0\}$ ,  $x_k^- = \min\{-x_k, 0\}$ , and  $x = x_1^+ + \dots + x_m^+$ . We can decompose the vector  $\mathbf{x}$  into its positive and negative components  $\mathbf{x} = \mathbf{x}^+ - \mathbf{x}^-$  where  $\mathbf{x}^+ = (x_1^+, \dots, x_m^+)^T$  and  $\mathbf{x}^- = (x_1^-, \dots, x_m^-)^T$  respectively. We also have  $x = \|\mathbf{x}^+\|_1 = \|\mathbf{x}^-\|_1 = \|\mathbf{x}\|_1/2$ . For example,

using this notation  $\mathbf{x} = (1, 2, -1, -1, -1)^T$  gives  $\mathbf{x}^+ = (1, 2, 0, 0, 0)^T$  and  $\mathbf{x}^- = (0, 0, 1, 1, 1)^T$ , and x = 3.

In our construction of quantum codes using the classical code C, we only consider logical codewords that are linear combinations over labels in C with only real coefficients. Hence, we define the two logical codewords of our quantum code as

$$|0_L\rangle = \frac{1}{\sqrt{x}} \left( \sqrt{x_1^+} |\mathbf{c}_1\rangle + \dots + \sqrt{x_m^+} |\mathbf{c}_m\rangle \right) ,$$
 (9)

$$|1_L\rangle = \frac{1}{\sqrt{x}} \left( \sqrt{x_1^-} |\mathbf{c}_1\rangle + \dots + \sqrt{x_m^-} |\mathbf{c}_m\rangle \right) .$$
 (10)

Since that  $x_j^+ x_j^- = 0$  for all  $j \in [m]$ , the logical states  $|0_L\rangle$  and  $|1_L\rangle$  have disjoint supports.

We next clarify the connection between  $\mathbf{x}$  and the non-deformation conditions by constructing a real matrix A that enforces these conditions. Roughly speaking, this matrix has rows labeled by diagonal Pauli errors of weight at most  $d_Z - 1$  and columns labeled by the states  $|\mathbf{c}_1\rangle, \ldots, |\mathbf{c}_m\rangle$ . While the ordering of the rows of A is unimportant, we will collect the rows in groups corresponding to the weights of P. The matrix A is defined by

$$A = \sum_{P \in \mathcal{Z}_0} \sum_{k=1}^m |P\rangle\langle k| + \sum_{w=1}^{d-1} \sum_{P \in \mathcal{Z}_w} \sum_{k=1}^m \left\{ \operatorname{Re}\left(\langle \mathbf{c}_k | P | \mathbf{c}_k \rangle\right) |P, 0\rangle\langle k| + \operatorname{Im}\left(\langle \mathbf{c}_k | P | \mathbf{c}_k \rangle\right) |P, 1\rangle\langle k| \right\}.$$
(11)

In matrix representation A is a wide rectangular matrix and writes

$$A = \begin{bmatrix} 1 & \cdots & 1 \\ a_{2,1} & \cdots & a_{2,m} \\ \vdots & \cdots & \vdots \\ a_{2V_q(2t)-1,1} & \cdots & a_{2V_q(2t)-1,m} \end{bmatrix} \equiv \begin{bmatrix} -\mathbf{a}_1^T - \\ -\mathbf{a}_2^T - \\ \vdots \\ -\mathbf{a}_{2V_q(2t)-1}^T - \end{bmatrix}$$
(12)

where  $\mathbf{a}_r = (a_{r,1}, \dots, a_{r,m})^T$  are column vectors. The reason for introducing the matrix A is that the non-deformation condition for correcting t errors using the vector  $\mathbf{x}$  is enforced by the constraint (see Fig.4)

$$A\mathbf{x} = 0$$
.

**Lemma 1.** Let  $\mathbf{x}$  be a non-zero real vector such that  $A\mathbf{x} = 0$ . Let  $|0_L\rangle$  and  $|1_L\rangle$  be logical codewords that depend on  $\mathbf{x}$  as in Eqs. (9) and (10). Then  $\langle 0_L|P|0_L\rangle = \langle 1_L|P|1_L\rangle$  for any diagonal Pauli of weight at most  $d_Z - 1$ .

*Proof.* Recall that  $x = x_1^+ + \cdots + x_m^+$ . Since each  $|\mathbf{c}_k\rangle$  is a product state, we have  $\langle \mathbf{c}_j | P | \mathbf{c}_k \rangle = 0$  for all distinct j and k, and for all diagonal Pauli of weight at most  $d_Z - 1$ . We can use the definitions of logical codewords (Eqs. (9) and (10)) to write

$$x \left( \langle 0_L | P | 0_L \rangle - \langle 1_L | P | 1_L \rangle \right) = \sum_{k=1}^m x_k^+ \langle \mathbf{c}_k | P | \mathbf{c}_k \rangle - \sum_{k=1}^m x_k^- \langle \mathbf{c}_k | P | \mathbf{c}_k \rangle$$
$$= \sum_{k=1}^m x_k \langle \mathbf{c}_k | P | \mathbf{c}_k \rangle$$

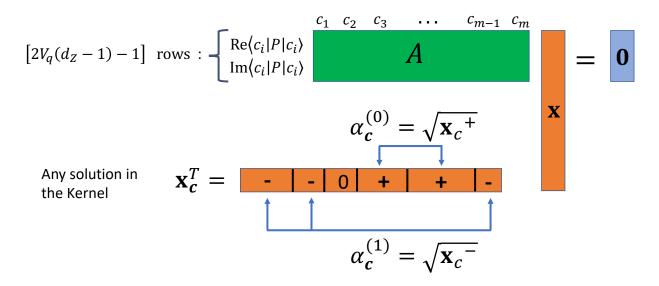


Figure 4: Illustration of solution of  $A\mathbf{x} = 0$  via Lemma 1.

where on the second line we used  $x_k = x_k^+ - x_k^-$ . Using Eq. (11) we see that  $\langle \mathbf{c}_k | P | \mathbf{c}_k \rangle = \langle P, 0 | A | k \rangle + i \langle P, 1 | A | k \rangle$ . Therefore

$$x(\langle 0_L | P | 0_L \rangle - \langle 1_L | P | 1_L \rangle) = \sum_{k=1}^m x_k (\langle P, 0 | A | k \rangle + i \langle P, 1 | A | k \rangle)$$
$$= \sum_{k=1}^m (\langle P, 0 | A x_k | k \rangle + i \langle P, 1 | A x_k | k \rangle)$$
$$= \langle P, 0 | A \mathbf{x} + i \langle P, 1 | A \mathbf{x} ,$$

because  $\mathbf{x} = \sum_{k=1}^{m} x_k | k \rangle$ . Since the first row of A is all ones,  $A\mathbf{x} = 0$  implies  $\langle 0_L | 0_L \rangle = \langle 1_L | 1_L \rangle$ . Moreover, in the above, P is an arbitrary diagonal Pauli of weight at most  $d_Z - 1$ , and the requirement  $A\mathbf{x} = 0$  implies  $\langle 0_L | P | 0_L \rangle = \langle 1_L | P | 1_L \rangle$  for any diagonal Pauli of weight at most  $d_Z - 1$ .

The condition  $A\mathbf{x} = 0$  is satisfied for any  $\mathbf{x} \in \ker(A)$ . It is then important to understand the structure of the kernel. Since A is real, any vector in its kernel must be real, which we then use to design an explicit quantum code that obeys the generalized orthogonality conditions above. Whenever  $|C| \geq 2V_q(d_Z - 1)$ , that is the number of codewords in C is strictly greater than the number of rows in A, by the rank-nullity theorem, A must have a non-trivial kernel. We thus have the following existence theorem for quantum codes that reside within the ground space of  $H_n$ .

**Lemma 2.** (Existence) Let C have a minimum distance at least  $d_X$ . If  $|C| \ge 2V_q(d_Z - 1)$ , then Alg. 1 constructs a quantum code with one logical qubit and with a bit-flip and phase-flip distance of  $d_X$  and  $d_Z$  respectively. Moreover, this quantum code can be constructed from any nonzero  $\mathbf{x} \in \ker(A)$ .

This lemma shows that we can derive the logical codewords of our quantum code from a classical code, and only requires C to have a minimum distance of  $d_X$  and for |C| to be at least twice the size of the q-ary Hamming ball of radius  $d_Z - 1$ .

We proceed to derive a more explicit expression for the entries of the matrix A in Eqs. (11) and (12). We will show that these entries are elements of a finite-sized set given by

$$\{\cos(2\pi k/q) , \sin(2\pi k/q) : k = 0, \dots, q-1\}.$$
 (13)

For any classical string  $\mathbf{c} = (c_1, \dots, c_n)^T \in C$  with  $c_i = 0, 1, \dots, q-1$ , the quantum state is  $|\mathbf{c}\rangle = |c_1, c_2, \dots, c_n\rangle$ , and when  $P = Z^{z_1} \otimes \cdots \otimes Z^{z_n}$ , we have

$$\langle \mathbf{c}|P|\mathbf{c}\rangle = \prod_{j=1}^{n} \langle c_j|Z^{z_j}|c_j\rangle = \prod_{j=1}^{n} \langle c_j|\sum_{k\in\Sigma_c} (\omega^k)^{z_j}|k\rangle\langle k|c_j\rangle = \prod_{j=1}^{n} \omega^{z_j c_j} = \omega^{\mathbf{z}^T \mathbf{c}}$$
(14)

$$= \cos\left(\frac{2\pi \,\mathbf{z}^T \mathbf{c}}{q}\right) + i \sin\left(\frac{2\pi \,\mathbf{z}^T \mathbf{c}}{q}\right). \tag{15}$$

The first row of A is all ones, and all other entries are given by

$$A' = \sum_{w=1}^{d_z - 1} \sum_{\text{wt}(\mathbf{z}) = w} \sum_{\mathbf{c} \in C} \left[ \cos \left( \frac{2\pi \, \mathbf{z}^T \mathbf{c}}{q} \right) | \mathbf{z}, 0 \rangle \langle \mathbf{c} | + \sin \left( \frac{2\pi \, \mathbf{z}^T \mathbf{c}}{q} \right) | \mathbf{z}, 1 \rangle \langle \mathbf{c} | \right]. \tag{16}$$

Since  $\mathbf{z}^T \mathbf{c}$  is always an integer, it follows that the entries of A must take values from the set in (13).

**Remark 1.** In Eq. (16) finding a solution in the kernel amounts to finding linear combination of roots of unity that vanish. There is a vast literature on this topic and properties of the underlying code that controls the values of the integers  $\mathbf{z}^T \mathbf{c}$  can in principle be utilized to give analytic solutions.

**Remark 2.** When q = 2, all rows in A' that are labeled by  $|\mathbf{z}, 1\rangle$  are equal to zero, because the argument of the sine is always an integer multiple of  $\pi$ .

This section is summarized in the following algorithm:

**Algorithm 1. Input:** A classical code  $C \subset \{0, 1, \dots, q-1\}^n$  with  $m \equiv |C|$ .

- Form the matrix A defined by Eq.(11).
- Solve  $A\mathbf{x} = 0$  to find  $\mathbf{x} \neq 0$ . Define  $\mathbf{x}^+, \mathbf{x}^- \geq 0$  such that  $\mathbf{x} = \mathbf{x}^+ \mathbf{x}^-$  as in Lemma 1
- Let  $C_0$  be the set of codewords  $\mathbf{c}_k$  with  $k \in supp(\mathbf{x}^+)$ . And let  $C_1$  be the set of codewords  $\mathbf{c}_k$  with  $k \in supp(\mathbf{x}^-)$ .
- For all  $\mathbf{c} \in C_0$ , assign  $\alpha_{\mathbf{c}}^{(0)} = \sqrt{x_c^+}$ , and for all  $\mathbf{c} \in C_1$ , assign  $\alpha_{\mathbf{c}}^{(1)} = \sqrt{x_c^-}$ .

**Output:** A logical quantum bit with codewords  $|o_L\rangle$  and  $|1_L\rangle$  as defined in Eqs. (9) and (10). The code distances are  $d_X = dist(C)$  and  $d_Z$  that satisfies  $m > 2V_q(d_Z - 1) - 1$ .

## 2.2 Constructing logical states with linear distance and constant rate

In building a single logical qubit we used any one non-zero solution in the kernel of A to identify two disjoint subsets  $C_0$  and  $C_1$ . Since the number of rows is  $2V_q(d_Z-1)-1$  and the number of columns by Gilbert-Varshamov bound satisfies  $m \geq q^n/V_q(d_Z-1)$ , we find that the ratio of the number of columns to the number of rows is asymptotically exponentially large

$$\frac{1}{n}\log_q\left(\frac{m}{\# \text{ rows}}\right) \ge 1 - 2\text{Ent}_q\left(\frac{d_Z}{n}\right) , \qquad \frac{d_Z}{n} \in \left[0, \frac{q-1}{q}\right]. \tag{17}$$

Here we exploit this to derive roughly  $q^{n(1-2\operatorname{Ent}_q(d_Z/n))}$  logical quantum states, and hence obtaining a linear rate  $r \approx (1-2\operatorname{Ent}_q(d_Z/n))$ . Below we think of  $M \propto m/(\# \text{ rows})$ .

We now generalize the construction of two quantum states (a logical qubit) to more quantum states (a logical qu dit). Suppose we identify M subsets  $\{C_0, C_1, \ldots, C_{M-1}\}$  such that  $C_i \subset C$  for all  $i \in \{0, \ldots, M-1\}$  and that the subsets are pairwise disjoint  $C_i \cap C_j = \emptyset$  for all  $i \neq j$ . Define the logical qudit as

$$|0_L\rangle = \sum_{\boldsymbol{c}\in C_0} \alpha_{\boldsymbol{c}}^{(0)} |\boldsymbol{c}\rangle, |1_L\rangle = \sum_{\boldsymbol{c}\in C_1} \alpha_{\boldsymbol{c}}^{(1)} |\boldsymbol{c}\rangle, \dots, |(M-1)_L\rangle = \sum_{\boldsymbol{c}\in C_{M-1}} \alpha_{\boldsymbol{c}}^{(M-1)} |\boldsymbol{c}\rangle.$$
 (18)

where , 
$$\sum_{c \in C_0} (\alpha_c^{(0)})^2 = \sum_{c \in C_1} (\alpha_c^{(1)})^2 = \dots = \sum_{c \in C_{M-1}} (\alpha_c^{(M-1)})^2 = 1$$
: normalization.(19)

Clearly these states are orthonormal. The KL criteria then writes

$$\Pi P \Pi = c_P \Pi \tag{20}$$

where  $\Pi = |0_L\rangle\langle 0_L| + |1_L\rangle\langle 1_L| + \cdots + |(M-1)_L\rangle\langle (M-1)_L|$  is a projector of (potentially exponentially large) rank M. The orthogonality and non-deformation conditions now write:

$$\langle i_L | P | j_L \rangle = c_P \, \delta_{i,j} \,, \quad i, j \in \{0, 1, 2, \dots, M - 1\} \,.$$

We first check orthogonality. Since the subsets are pairwise disjoint and futhermore C has a minimum distance of  $d_X$ , we have  $\langle i_L|P|j_L\rangle=0$  for all  $i\neq j$  for and all diagonal Paulis. Moreover,  $\langle i_L|P|j_L\rangle=0$  if  $1\leq \operatorname{wt}_X(P)\leq d_X-1$ , which is inherited from the classical code's distance  $d_X$ . We now turn our attention to the non-deformation condition. As before it is clear that  $\langle i_L|P|i_L\rangle=0$  for all  $i\in\{0,1,\ldots,M-1\}$  and Paulis with  $1\leq \operatorname{wt}_X(P)\leq d_X-1$ . This follows from the distance of the code as before. Therefore, it is again sufficient to prove the non-deformation condition for diagonal Paulis. We need to find the subsets  $C_1,C_2,\ldots,C_{M-1}$  such that

$$\langle 0_L | P | 0_L \rangle = \langle 1_L | P | 1_L \rangle = \dots = \langle (M-1)_L | P | (M-1)_L \rangle. \tag{21}$$

Our method constructs the sets  $C_0, \ldots, C_{M-1}$  such that they are disjoint and satisfy the foregoing equation. Unlike the qubit case, our proof technique will not be seeking a solution in the kernel of A anymore, rather we recursively build the logical quantum states.

Comment: It would be exciting to see an alternative construction that uses the exponentially large kernel of the matrix A.

For every new logical state we need to call the following algorithm once:

**Algorithm 2. Input:** A set of  $2V_q(d_Z-1)$  columns of A and a vector **b** of size  $2V_q(d_Z-1)-1$ .

- Check that the augmented homogeneous linear system as described in Section 5 of [Din26] does not have a row that is all of the same sign.
- If no row has non-zero entries of the same sign, apply the algorithm of [Din26] to find a point in the feasible set, i.e., a solution  $\mathbf{x} \geq 0$ . End the algorithm.
- If there is a row that is all of same sign, augment the set of columns by one and repeat.

  If all the columns are exhausted, then output fail.

**Output:** If succeeded, output a solution  $\mathbf{x} \geq 0$  in the feasible set.

If recursively successful, then the algorithm can at least be called  $m/2V_q(d_Z-1)=q^{n(1-2\mathrm{Ent}_q(d_Z/n))}$  times. The success is guaranteed if at any step of recursion no row of all the same sign is encountered.

We demonstrate the recursive construction of logical quantum states by first building a qutrit (three logical states). Recall that A has  $2V_q(d_Z-1)-1$  rows whose first row is all ones, we now proceed to find a natural partitioning of the columns of A. To construct a logical qubit, we use the first  $2V_q(d_Z-1)$  columns of A as follows. Let the matrix  $A'_1$  be defined by the first  $2V_q(d_Z-1)$  columns of A. Now solve for the solution  $\mathbf{x_1}$  in  $A'_1\mathbf{x}=\mathbf{0}$ , where  $\mathbf{x_1}$  has  $2V_q(d_Z-1)$  components. This matrix equation certainly has a non-trivial kernel because it has more columns than rows. Just as in the one logical qubit construction in Alg. 1 we will find an  $\mathbf{x}=\mathbf{x}^+-\mathbf{x}^-$  where  $\mathbf{x}^+\geq 0$  and  $\mathbf{x}^-\geq 0$ ,  $\sum_i x_i^+=\sum_i x_i^-$ , and that  $supp(\mathbf{x}^+)\cap supp(\mathbf{x}^-)=\emptyset$ . We build the logical qubit exactly as in Eqs. (9) and (10).

We then rearrange the columns of  $A'_1$  according to the supports of  $\mathbf{x}^+$  and  $\mathbf{x}^-$  and write the concatenated matrix  $[A_1A_2]$ , where  $A_1$  has a set of columns labeled by codewords  $c_s$  such that  $s \in supp(\mathbf{x}^+)$  and  $A_2$  has columns labeled by codewords  $c_s$  such that  $s \in supp(\mathbf{x}^-)$ . The new equation being satisfied is

$$[A_1 A_2] \left[ \begin{array}{c} \boldsymbol{x_1^+} \\ -\boldsymbol{x_1^-} \end{array} \right] = 0 ,$$

where  $x_1^+, x_1^- \ge 0$ . To build the third logical state, we select a set of  $2V_q(d_Z - 1)$  new columns out of A and solve

$$[A_2A_3]$$
 $\begin{bmatrix} -x_1^- \\ x_2 \end{bmatrix} = 0, \quad x_2 \ge 0,$ 

where  $-x_1^-$  is treated fixed from the previous step and we solve for a solution  $x_2 \ge 0$ . A solution exists as  $A_3x_2 = A_2x_1^-$  is under-constrained. Although this can be formally thought of as a linear programming problem, where the objective function is just zero and a point in the feasible set is sought, we proceed differently and give an explicit algorithm based on Dines' Annals of Math (1926) [Din26].

Once  $x_2 \geq 0$  is found this way, we proceed to solve

$$[A_3A_4] \left[ \begin{array}{c} \boldsymbol{x_2} \\ -\boldsymbol{x_3} \end{array} \right] = 0 , \quad \boldsymbol{x_3} \ge 0,$$

where now  $x_3$  is the unknown. Just as before this amounts to solving  $A_4x_3 = -A_3x_2$  where the right-hand side is a known vector. Continuing this way we can build  $2 \le M \le q^{n(1-2\operatorname{Ent}_q(\frac{d_Z}{n}))}$  logical states.

The results of this section prove the following main theorem of part 1 of this work:

**Theorem 1.** Take a classical code C of length n on a q-ary alphabet with a minimum distance of  $d_X$ . If  $|C| \geq 2V_q(d_Z - 1)$ , then Alg. 1 and multiple calls to Alg. 2 explicitly derives the M quantum logical states in (18) with bit- and phase-flip distances of  $d_X$  and  $d_Z$  respectively. The overall distance is  $\min(d_X, d_Z)$ .

When A matrix is wide (under-constrained) Alg. 1 always constructs two logical codewords (a logical qubit). The third and subsequent logical codewords are found recursively by calling Alg. 2 multiple times. Alg. 2 succeeds with high probability and almost surely over random codes. In the rare case that the set of linear constraints have rows that are all non-zero and are of the same sign, the recursion becomes infeasible and algorithm halts. As shown in [Din26] this is the only way that the algorithm can fail. We define a random classical code as one whose codewords have entries over  $\Sigma = \{0, 1, 2, ..., q - 1\}$  such that each entry is independently and randomly drawn from the uniform distribution over  $\Sigma$ .

**Lemma 3.** M quantum logical states can be constructed as long as in each step of the recursion, no row other than the first has entries that are all non-zero and with the same sign. When C is a random classical code, then with high probability, Alg. 2 succeeds in providing a quantum code with constant rate.

*Proof.* The first part of the Lemma follows from the proof of Dines [Din26] applied to every step of the recursion. To prove the second part, first recall that at each step we are solving a linear system with  $\rho \equiv 2V_q(d_Z-1)$  columns and  $\rho-1$  rows. To find positive solutions of this linear system, we employ Dines algorithm which is itself recursive. Hence we will prove that at every step of Dines recursive algorithm, it fails with low probability. We prove this by induction, first starting with the base case.

Recall that  $\omega \equiv \exp(2\pi i/q)$ . For every diagonal P of weight at least one,  $\langle c|P|c\rangle$  is a random variable that takes the values  $\{1,\omega,\ldots\omega^{q-1}\}$  with a uniform probability. From Eq. (15) we know that each  $\langle c|P|c\rangle$  is equal to a  $\omega^j$  for some  $j \in \{0,1,\ldots,q-1\}$ . That is,  $\operatorname{Re}(\langle c|P|c\rangle)$  and  $\operatorname{Im}(\langle c|P|c\rangle)$  are random variables that take values in [-1,1]. Since the code words are random and uniformly distributed over the symbols, by symmetry the probability of an entry having a positive (or negative) sign is a half when q is even and is at most 2/3 when q is odd. Moreover the entries are independent. Hence the probability that a given row has all the same sign is  $(3/2)^{-\rho+1}$ . And by a union bound, the probability that any of the  $\rho-1$  rows have entries whose all entries have the same sign is  $O(\rho(3/2)^{-\rho})$ .

Now we prove the induction step. We take q to be even for now, which ensure that  $a_{ij}$  are symmetric random variables with mean zero. Note that Dines algorithm takes a matrix with matrix elements  $a_{ij}$ , and constructs a new matrix  $a_{r,ij} = a_{1i}a_{rj} - a_{1j}a_{ri}$ . In the new matrix, the indices i and j belong to disjoint sets I and J. The number of columns in the new matrix is |I||J|. For the induction hypothesis, we assume that the matrix elements  $a_{ij}$  are identical and symmetric random variables with zero mean, which are furthermore independent with respect to the column index j. We will show that  $a_{r,ij}$  will then also be

a symmetric random variable with zero mean and also furthermore be independent with respect to the new column indices ij.

We now show that the  $a_{r,ij}$  has zero mean. For that we see  $\mathbb{E}(a_{r,ij}) = \mathbb{E}(a_{1i}a_{rj}) - \mathbb{E}(a_{1j}a_{ri})$  because the expectation is linear. Next the independence of i and j imply that  $\mathbb{E}(a_{1i}a_{rj}) = \mathbb{E}(a_{1i})\mathbb{E}(a_{rj})$  and  $\mathbb{E}(a_{1j}a_{ri}) = \mathbb{E}(a_{1j})\mathbb{E}(a_{ri})$ . Substituting this shows that  $\mathbb{E}(a_{r,ij}) = 0$  because  $a_{ij}$  are independent random variables with mean zero.

We now note that the random variables  $a_{r,ij}$  are independent with respect to the column labels ij. This follows readily from the independence of  $a_{ij}$  with respect to j. For instance, treat i to be fixed and consider  $a_{r,ij}$  and  $a_{r,ij'}$  for  $j \neq j'$ .

We next show that the random variables  $a_{r,ij}$  are symmetric. For this, we use the fact that the product of non-degenerate symmetric random variables is a symmetric random variable [HW85].

In the base case, we have shown that  $a_{ij}$  satisfies the induction hypothesis with high probability. We have just shown that  $a_{r,ij}$  is symmetric, independent, and has zero mean with respect to ij. This proves that its entries have equal probability of being positive or negative.

It remains to bound the probabilities of failure of our algorithm under our recursion and Dines recursion. With high probability, the product |I||J| is going to be greater than the number of columns in  $a_{ij}$ . As shrinking the number of columns will only be due to at most one entry of a different sign, this happens with very low probability. The probability of this happening is at most  $2\rho(1/2)^{\rho-1}$  for large  $\rho$ . The probability of a single row with all the same sign is at most  $2(1/2)^{\rho}$ . Hence the total probability of a given row being pathological is at most  $2(1/2)^{\rho} + 2\rho(1/2)^{\rho-1} = O(\rho(1/2)^{\rho})$ . The probability of a matrix at any step of Dines recursion to have a pathological row is therefore at most  $O(\rho^2(1/2)^{\rho})$ . Since our algorithm has  $O(m/\rho)$  steps, our algorithm's failure probability is at most  $O(m\rho(1/2)^{\rho})$ . Since  $m2^{-\rho}$  is at most  $(2^n2^{-2^{cn}})$  for some constant  $c \in [0, 1]$ , our algorithm will succeed with overwhelming probability. Although this proof is specialized for the case where q is even, a similar argument will work for q odd.

In the next section, We find that even if this happens we can always construct *Approximate* Quantum Error Correcting Codes (AQECC) [LNCY97] with linear distance and constant rate, provided that, the underlying classical code also has a linear distance and a sufficiently large rate.

## 2.3 AQECCs with designed rates

In the unlikely case, where building logical quantum states using Alg. 2 of the previous section fails, we can always build an AQECC as we now show. We introduce an algorithm that produces a quantum code with M logical codewords satisfying the KL criteria approximately. Here M can be strictly larger than 2 at the expense of an approximation error, which is equal to the infidelity of quantum code.

Suppose  $C_1, \ldots, C_{M/2}$  are disjoint subsets of the classical code C whose minimum distance is  $d_X$ . Hence each  $C_j$ ,  $j \in [M/2]$  inherits the distance  $d_X$ , where we take M to be even for simplicity. Suppose that for every  $j \in [M/2]$ , it holds that

$$|C_j| \ge 2V_q(d_Z - 1) . (22)$$

For each classical code  $C_j$ , we use Alg. 1 to construct a corresponding matrix  $A_j$  from which we derive logical codewords  $|(2j)_L\rangle$ ,  $|(2j-1)_L\rangle$  that satisfy the KL criteria for quantum codes with a bit-flip and phase-flip distance of  $d_X$  and  $d_Z$  respectively.

It is clear from our construction that for every diagonal Pauli P of weight at most  $d_Z - 1$ ,

$$\langle (2j)_L | P | (2j)_L \rangle = \langle (2j-1)_L | P | (2j-1)_L \rangle = \gamma_{j,P}$$
 (23)

where  $\gamma_{j,P}$  is a complex number of norm at most one. Now for each  $j \in [M/2]$ , let  $\Gamma_j = (\gamma_{j,P})$  be a row vector of length  $V_q(d_Z - 1)$  with components that correspond to diagonal Paulis of weight at most  $d_Z - 1$ . Then it follows that each  $\Gamma_j$  lies in a  $|V_q(d_Z - 1)|$ -dimensional complex unit ball corresponding to a hyper-cube of length two centered at the origin with respect to the infinity norm.

Suppose  $\max_{j,k\in[M/2]} \|\Gamma_j - \Gamma_k\|_{\infty} = \delta$ . Then because of Eq. (23) we have

$$\max_{j,k=1,\dots,M} \max_{P} |\langle j_L | P | j_L \rangle - \langle k_L | P | k_L \rangle| = \delta.$$
 (24)

It remains to find a suitable upper bound for  $\delta$ . To relate the error  $\delta$  in satisfying the non-deformation to the size of the code |C| and the number of logical qubits we can construct, we rely on the following fact.

**Fact.** Consider the complex hyper-cube of side length two in N dimensions. Let x be the number of points distributed randomly inside it. Then there exists a ball of radius  $\delta$  in the infinity norm that contains at least  $x(\delta/2)^N$  points in expectation.

The number of points inside the unit hyper-cube is  $x = \lfloor |C|/(2V_q(d_Z - 1))\rfloor$ . The radius of the ball is  $\delta$ . The dimension of the hyper-cube is  $V_q(d_Z - 1)$ . Hence from the above fact, we have that the expected number of logical codeword pairs is  $\mathbb{E}M = 2x(\delta/2)^N$  which writes

$$\mathbb{E}M \ge 2 \left| \frac{|C|}{2V_q(d_Z - 1)} \right| \left( \frac{\delta}{2} \right)^{V_q(d_Z - 1)}. \tag{25}$$

The infidelity  $\epsilon$  defined by one minus the worst case entanglement fidelity of the quantum code can be upper bounded as shown in [Ouy14], to be

$$\epsilon \le O\left(\delta V_a^4(d_Z - 1)\right),\tag{26}$$

when the noisy quantum channel introduces bit-flip and phase-flip weights at most  $d_X - 1$  and  $d_Z - 1$  uniformly at random.

#### 2.4 Illustrations

In this section we illustrate our framework through a series of examples. It is noteworthy that the first example is a one-to-one correspondence between the celebrated classical and quantum results of Hamming's and Steane's respectively.

#### 2.4.1 C as a [7,4,3] Hamming code gives the Steane code

In classical coding theory, we use the notation (n, m, d) to denote a binary code with codewords of length n that has m codewords, a distance of d. We use  $[n, \log_2 m, d]$  to denote a binary code with codewords of length n that has m codewords, a distance of d, and is furthermore a linear code. Consider the case when C is generated from the codewords 1000110, 0100101, 0010011, and 0001111. This classical code is the celebrated [7,4,3] Hamming code, and has been used previously by Steane to obtain the [[7,1,3]] Steane code. Applying our framework to C, we get the unique solution

$$|0_{L}\rangle = \frac{1}{\sqrt{8}} (|0,0,0,0,0,0,0\rangle + |0,0,0,1,1,1,1\rangle + |0,1,1,0,1,1,0\rangle + |0,1,1,1,0,0,1\rangle + |1,0,1,0,1,0,1\rangle + |1,0,1,1,0,1,0\rangle + |1,1,0,0,0,1,1\rangle + |1,1,0,1,1,0,0\rangle)$$
(27)  
$$|1_{L}\rangle = \frac{1}{\sqrt{8}} (|0,0,1,0,0,1,1\rangle + |0,0,1,1,1,0,0\rangle + |0,1,0,0,1,0,1\rangle + |0,1,0,1,0,1,0\rangle + |1,0,0,0,1,1,0\rangle + |1,0,0,1,0,0,1\rangle + |1,1,1,0,0,0,0\rangle + |1,1,1,1,1,1,1\rangle).$$
(28)

This is in fact equivalent to the Steane code.

#### **2.4.2** *C* as a [6,3,3] code

The set of all three-bit strings comprises of 000, 001, 010, 011, 100, 101, 110, and 111. Now append each string with another three-bit string so that C = [6, 3, 3] comprises of the codewords 000000, 001110, 010101, 011011, 100011, 101101, 110110, and 111000. When the designed minimum distance of the quantum code is 3, our algorithm finds no quantum code. However if the designed distance is reduced to 2, then we derive the following error detecting quantum code

$$|0_L\rangle = \frac{1}{2} (|0,0,0,0,0,0\rangle + |0,1,1,0,1,1\rangle + |1,0,1,1,0,1\rangle + |1,1,0,1,1,0\rangle)$$
 (29)

$$|1_L\rangle = \frac{1}{2} (|0,0,1,1,1,0\rangle + |0,1,0,1,0,1\rangle + |1,0,0,0,1,1\rangle + |1,1,1,0,0,0\rangle)$$
(30)

## 2.4.3 C as a [4,2,2] code.

When the designed minimum distance of our quantum code is 2, we find no quantum code when our algorithm uses the classical code C = [4, 2, 2] with codewords 0000, 1010, 1101, and 0111.

#### 2.4.4 C as a nonlinear cyclic code

Now consider the nonlinear (4,8,2) code with codewords that are cyclic permutations of 0001 and 1110. The corresponding kernel of the matrix A has dimension 3, and one solution to this gives an error detecting quantum code with logical codewords

$$|0_L\rangle = \frac{1}{\sqrt{2}}(|0,0,0,1\rangle + |1,1,1,0\rangle)$$
 (31)

$$|1_L\rangle = \frac{1}{\sqrt{2}}(|0,0,1,0\rangle + |1,1,0,1\rangle).$$
 (32)

In fact, we can also have the additional logical codewords

$$|2_L\rangle = \frac{1}{\sqrt{2}}(|0, 1, 0, 0\rangle + |1, 0, 1, 1\rangle)$$
 (33)

$$|3_L\rangle = \frac{1}{\sqrt{2}}(|1,0,0,0\rangle + |0,1,1,1\rangle).$$
 (34)

This gives a quantum code of dimension 4 and a minimum distance of 2. This quantum code is also an example of a CWS code. This is because for every j=0,1,2,3, there is a Pauli operator that takes the stabilizer state  $\frac{1}{\sqrt{2}}(|0,0,0,0\rangle+|1,1,1,1\rangle)$  to  $|j_L\rangle$ . Since this quantum code is CWS, quantum error correction can proceed using formalism developed for CWS codes [CSSZ08]. This quantum code also has some other attractive properties. First, by inducing cyclic shifts in the underlying qubits, we can move from one logical codeword to another. Second, this quantum code is stabilized by  $X^{\otimes 4}$  and is affected uniformly by  $Z^{\otimes 4}$ . Together, his implies that the quantum code is invariant under transversal X and Y and Z operations.

#### 2.4.5 Permutation-invariant quantum codes

Our quantum code construction formalism can also extend to quantum codes with logical codewords that are supported on non-product basis states. One example of such codes are permutation-invariant quantum codes, which are invariant under any permutation of the underlying particles. Permutation-invariant quantum codes have been explicitly constructed using a variety of different techniques [Rus00, PR04, Ouy14, OF16, Ouy17, OC19]. Recently, permutation-invariant quantum codes have been consider for applications such as for quantum storage [Ouy19], or for robust quantum metrology [OSM19], and they can also be prepared in physically realistic scenarios [WWG+19].

When permutation-invariant quantum codes are constructed on n qubits, they must be superpositions over Dicke states

$$|D_w^n\rangle = \frac{1}{\sqrt{\binom{n}{w}}} \sum_{\substack{x_1,\dots,x_N \in \{0,1\}\\x_1+\dots+x_n=w}} |x_1\rangle \otimes \dots \otimes |x_n\rangle.$$
(35)

Here w is the weight of the Dicke state, and counts the Hamming weights of its constituent computation basis states' labels. The Dicke states for qubit states are labeled by only their weights, of which there are only n+1 possibilities. For our quantum code construction, we can choose the logical states to be supported on  $|D_{w_1}^n\rangle, \ldots, |D^n w_m\rangle$  where  $w_{j+1} - w_j \ge d$  for any  $j = 1, \ldots, m-1$ , and d is the desired minimum distance of the quantum code.

When a quantum code is permutation-invariant, we only need to consider equivalence classes of Pauli operators up to a permutation. Since because for Dicke states,  $\langle D_w^n|P|D_w^n\rangle$  are not necessarily zero even when the Pauli P is non-diagonal [OSM19], we need to count the number of all Paulis of weight at most d-1 up to a permutation. The number of unique qubit-Paulis up to a permutation having a weight of at most w is equal to the number of ways to order a w-tuple in  $\{1,2,3\}^w$  in a non-decreasing sequence, and this number is just  $\binom{n+w-1}{w}$ . Hence to total number of Paulis that we need to consider for the non-deformation conditions is at most  $\sum_{w=0}^{d-1} 3^w$ .

Now we consider a variation of the A-matrix from Alg. 1 with the matrix elements

$$\langle D_{w_i}^n | P | D_{w_i}^n \rangle, \tag{36}$$

where P labels the rows and j labels the columns.

From this, we can get permutation-invariant quantum codes with a minimum distance of d whenever

$$(\lfloor n/d \rfloor + 1) \ge 1 + \sum_{w=0}^{d-1} 3^w.$$
 (37)

For instance, when d = 3 this inequality becomes

$$|n/3| \ge (1+13),\tag{38}$$

and this formalism show that we can get gives permutation-invariant codes with a distance of d=3 when  $n\geq 42$ . This bound is however loose, because there are permutation-invariant quantum codes with d=3 on 9 qubits [Rus00, Ouy14], and even on 7 qubits [PR04]. This suggests that rather than using loose bounds on the nullity of the A matrix in Alg. 1, we need to exploit additional structure about the kernel of A to realize the full potential of our formalism.

#### 2.5 Optimality

If we take the metrric of optimality to be maximization of the distance d for fixed length n and number of encoded qubits k, then our construction is not optimum. This is because the five-qubit code is the unique [[5,1,3]] code, and our framework does not encompass it.

We demonstrate examples of optimal quantum codes using our framework which use the Hamming code and a nonlinear cyclic code in Sec. 2.4.1 and Sec. 2.4.3 respectively.

However, let us additionally impose the constraint (C1) that logical codewords must be supported on disjoint subsets of computational basis states. (C1) is satisfied by all CSS codes. Our framework generalizes CSS codes, but does not encompass all quantum codes. For example many stabilizer codes are not captured by our framework. Then our framework can give optimal quantum codes with (C1). For construction of a logical qubit, we can impose a second constraint, (C2), which is that the quantum code must be supported on a fixed classical code C, in addition to satisfying (C1). Then as long as the A-matrix from Alg. 1 has a kernel of dimension one, then the construction is optimal with respect to (C2).

We can also define a notion of optimality for our quantum codes that encode more than a logical qubit. Given a fixed classical code C, if the number of logical codewords in our derived quantum codeword using Alg. 1 and Alg. 2 is equal to the dimension of the nullspace of the A-matrix in Alg. 1 plus on, then we say that our quantum code is optimal. For instance, we find that the Steane code is an optimal construction, and also find an optimal construction using the nonlinear cyclic code that encodes four logical codewords.

# 3 Part 2: Linear distance codes in ground space of local Hamiltonians

### 3.1 Why introduce a Hamiltonian?

#### **Theoretical**

This model is of theoretical interest because it allows for the encoding of linear distance codes in its ground space. Encoding quantum information in the ground space of physical Hamiltonians has a long history dating back to Kitaev's toric code [Kit03]. A motivation for our work was the nice work of Brandao et al. [BaCimcbuB19] where existence of codes in low energy eigenstates of local translation invariant spin chains were found. In this work we came up with explicit and exact constructions and overcome some of the limitations of that work. These were detailed in the introduction and will be elaborated on below.

It is our hope that this line of work will prove useful in constructing QLDPC codes with linear distance, which despite recent progress [TZ13, CDZ13, BH14, KT20, EKZ20, HHO20] has been a long standing open problem.

#### Applied and engineering

Finding new quantum codes can help hasten the dream of fault-tolerant quantum computation. Our method is distinct from other widely used methods to construct quantum codes from classical codes, such as for CSS codes, stabilizer codes, and codeword-stabilized codes. A key feature of our code construction is that we can take as input a general classical code, and demand that the quantum code we construct must be supported on the computational basis vectors that are labelled by these classical codewords. The non-trivial solution of the nullspace then gives us the amplitudes over with the logical zero and logical one of our quantum code will assume.

Our proposal to engineer a two-local Hamiltonian to stabilize quantum information is in line with the ideas utilizing quantum control techniques to suppress the noise before employing quantum error correction. We differ from traditional approaches in quantum control procedures where one typically applies dynamical decoupling pulses to create an essentially a trivial identity Hamiltonian that acts on the system when no quantum gates are performed. In this situation, however, local errors are not energetically penalized. In contrast, one might envision that quantum control methods can engineer a Hamiltonian that energetically penalizes the dominant noise rates that occur in a quantum system before introducing quantum error correction [ML14].

In many practical physical systems, noise is biased and can be dominated by bit-flip or phase-flip type errors. We consider a base noise model that is dominated by bit-flip type errors. On a spin-system, we expect our Hamiltonian to energetically penalize bit-flip errors. This would allow our engineered Hamiltonian to greatly suppress the noise rates of the dominant (bit-flip) type of errors. The remaining errors can then be cleaned up using our quantum code with a linear distance. The advantage of engineering our Hamiltonian, as compared to, for instance, the surface code Hamiltonian, is that the Hamiltonian terms that we require are two-local, whereas the surface code requires many-body interactions, which

is challenging to realize in practice.

The Hamiltonian we will introduce has the form  $H_n = \sum_{k=1}^{n-1} H_{k,k+1}$ ; see Eq.(40). Its ground space has a strict subspace that is the quantum code. We could make the model sums of commuting local terms by considering a new Hamiltonian that skips all (say) even interactions and write  $H'_n = \sum_{k \text{ odd}} H_{k,k+1}$ . The ground space of  $H'_n$  contains the ground space of  $H_n$  and therefore also includes the quantum code. One could continue this way and eventually get a Hamiltonian that is trivial (i.e., no interaction) H'' = I for which any quantum code is in the "ground space". The price one pays following this crooked path is that Nature will help less and less in suppressing the rate at which errors appear.

Since our quantum code does not occupy the entire subspace of our designed 2-local Hamiltonian, one might call it a subspace quantum LDPC code. The fact that we only use a fraction of the ground space of our Hamiltonian indeed leads to our quantum code imposing no energy penalty on phase errors. However, in a noise model where phase errors are rare, this is not a problem, as we can clean up the few phase errors using subsequent quantum error correction.

#### 3.2 Local Hamiltonian and its ground space

Let us consider a spin chain of length n with open boundary conditions and the local Hilbert space dimension of 2s + 1, where  $s \ge 1$  is a positive integer. We take a representation in which  $|j\rangle$  denotes the  $s_z = j$  state of a spin-s particle:

$$\hat{S}_z|j\rangle = j |j\rangle, \qquad j \in \Sigma \quad .$$

The local Hamiltonian whose ground space will be shown to contain a nontrivial quantum error correcting code is

$$H_n = H_n^J + H_n^s \tag{39}$$

where  $H_n^J = J \sum_{k=1}^n (|0\rangle\langle 0|)_k$ . Recall that the Hamiltonian  $H_n^s$ , is defined by

$$H_n^s = \sum_{k=1}^{n-1} \left\{ \sum_{m=-s}^s P_{k,k+1}^m + \sum_{m=1}^s Q_{k,k+1}^m \right\} , \qquad (40)$$

and the local terms are projectors acting on two neighboring spins k, k+1 defined by

$$P^{m} = |0 \leftrightarrow m\rangle\langle 0 \leftrightarrow m| , \qquad Q^{m} = |00 \leftrightarrow \pm m\rangle\langle 00 \leftrightarrow \pm m|; \tag{41}$$

where

$$|0 \leftrightarrow m\rangle \equiv \frac{1}{\sqrt{2}} [|0, m\rangle - |m, 0\rangle]$$
 (42)

$$|00 \leftrightarrow \pm m\rangle \equiv \frac{1}{\sqrt{2}} [|0,0\rangle - |m,-m\rangle],$$
 (43)

and we denoted by  $|j,k\rangle$  the spin state  $|s_k^z=j,s_{k+1}^z=k\rangle$ . There are 3s local projectors as  $P_{k,k+1}^0$  automatically vanishes. See Fig. 5.

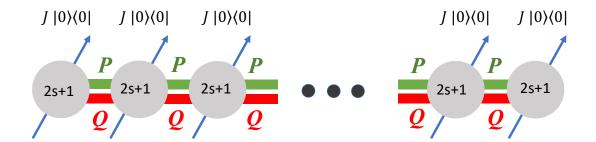


Figure 5: The new local integer spin-s Hamiltonian  $H_n$ .

In the simplest form s = 1, and we have

$$H_n^1 = \frac{1}{2} \sum_{k=1}^{n-1} \left\{ |0 \leftrightarrow 1\rangle \langle 0 \leftrightarrow 1| + |0 \leftrightarrow -1\rangle \langle 0 \leftrightarrow -1| + |00 \leftrightarrow \pm 1\rangle \langle 00 \leftrightarrow \pm 1| \right\} ,$$

where  $|1 \leftrightarrow 0\rangle \propto |0, 1\rangle - |1, 0\rangle$ ,  $|-1 \leftrightarrow 0\rangle \propto |0, -1\rangle - |-1, 0\rangle$ , and  $|\pm 1 \leftrightarrow 00\rangle \propto |0, 0\rangle - |1, -1\rangle$ . Below we will be mostly interested in s > 1.

**Lemma 4.** Suppose  $H_1 \geq 0$  and  $H_2 \geq 0$ , and  $H_1 + H_2$  is a frustration free (FF) Hamiltonian with zero energy ground state. Then the ground space of  $H_1 + H_2$  coincides with the intersection of the ground spaces of  $H_1$  and  $H_2$ .

*Proof.* Any state in the intersection of the kernels of  $H_1$  and  $H_2$  automatically vanishes on  $H_1 + H_2$ . Conversely, a state  $|\psi\rangle$  that is in the kernel of the sum  $H_1 + H_2$  obeys  $\langle \psi | (H_1 + H_2) | \psi \rangle = 0$ . Since each summand is a positive operator, so is their sum, and for  $|\psi\rangle$  to be a zero energy ground state of  $H_1 + H_2$  it has to vanish on each summand  $\langle \psi | H_i | \psi \rangle = 0$  for i = 1, 2. Therefore  $|\psi\rangle$  is a FF ground state of each  $H_i$  as well.

It is clear that

$$\operatorname{spec}(H_n^J) = J\{0, 1, 2, \dots, n\}$$
(44)

whose gap we denote by  $\Delta(H_n^J) = J$ . The kernel of  $H_n^J$  is the span of all product states  $|\mathbf{t}\rangle$  of weight n, where  $\mathbf{t} \in \Sigma_*^n$ ; note that the letter 0 is excluded in these strings.

We will obtain the ground space of  $H_n$  by taking the intersection of the ground space of  $H_n^J$  with  $H_n^s$ . We proceed to analytically derive the ground space of  $H_n^s$  after preliminary definitions. From now we assume that m is a positive integer unless stated otherwise.

Since  $H_n^s$  is free of the sign problem (i.e., stoquastic), the local projectors define an effective Markov chain, which have the following correspondence:

Local Projector	Local moves	Interpretation
$P^m$	$0m \longleftrightarrow m0$	Spin transport: local exchange of spin $m$ with 0
$Q^m$	$00 \longleftrightarrow m, -m$	Spin interaction: local creation/annihilation of $m, -m$

We say two strings  $\mathbf{t}$ ,  $\mathbf{z}$  are equivalent, denoted by  $\mathbf{t} \sim \mathbf{z}$  if  $\mathbf{z}$  can be reached from  $\mathbf{t}$  by applying a sequence of the local moves defined in the table. We define a set of equivalence classes as follows. Let k denote the number of nonzero letters in a product state (Eq. (45)). Using a consecutive set of the local moves stated above, we can take any state  $\dots m0 \dots 0(-m) \dots \dots \dots 0 \dots 0m(-m)0 \dots 0 \dots \dots \dots 0000 \dots$  We then move all the zeros to the rightmost end and ensure that all strings are of the form

$$c_{x_1,\dots,x_k} = x_1 \dots x_k \underbrace{0 \dots 0}_{n-k} \tag{45}$$

where  $x_i \in \Sigma_*$ . By assumption the string  $x_1 \dots x_k$  cannot be further reduced and n-k is the maximum number of zeros. Then it follows that if  $x_i = m$  then it must not have to its immediate right an  $x_{i+1} = -m$  for otherwise the annihilation rule  $(m, -m) \to 00$  would further reduce it.

**Lemma 5.** Any string  $\mathbf{t} \in \Sigma^n$  is equivalent to one and only one  $c_{x_1,\dots,x_k}$  (Eq. (45)).

Proof. By applying the local moves in the table above to any string  $\mathbf{t}$ , one can make sure that there are no substrings m(-m) or  $m0\cdots 0(-m)$  for any  $m\in[s]$ , where  $[s]=\{1,2,\ldots,s\}$ . Now suppose we apply these moves to  $\mathbf{t}$  as much as possible to bring it as close as possible to the state of all zeros. Then if  $\mathbf{t}$  contains a single m, then the first non-zero letter to its right cannot be -m. Similarly if  $\mathbf{t}$  contains at least one -m then the first non-zero letter to its left must not be m. By applying  $0(-m) \to (-m)0$  and  $0m \to m0$  we move all the zeros to the right to obtain a string of the form given by Eq. (45). To prove that the set of all strings equivalent to  $c_{x_1,\ldots,x_k}$  is indeed an equivalence class, we need to prove that the classes are distinct. It is clear that any string is equivalent to itself (reflexive). If  $\mathbf{x} \sim \mathbf{y}$  and  $\mathbf{y} \sim c_{x_1,\ldots,x_k}$  then  $\mathbf{x} \sim c_{x_1,\ldots,x_k}$  (transitive). Lastly if  $\mathbf{x} \sim \mathbf{y}$ , then  $\mathbf{y} \sim \mathbf{x}$  because of the reversibility of the local moves (symmetric). Therefore indeed the set of strings equivalent to  $c_{x_1,\ldots,x_k}$  form an equivalence class and it is an elementary fact that equivalence classes are distinct and partition the state space (i.e., the set of all strings) into disjoint subsets.

**Lemma 6.** The uniform superposition of all strings in an equivalence class (i.e., equivalent to the irreducible string in Eq. (45)) is a (frustration free) zero energy ground state of  $H_n^s$ .

Proof. The Hamiltonian  $H_n^s$  in Eq. (40) is a sum of local projectors (Eq. (41)). If a ground state  $\psi$  vanishes on each local projector, then for any  $m \in [s]$  it must obey  $\langle \psi | 0m \rangle = \langle \psi | m0 \rangle$ ,  $\langle \psi | 0(-m) \rangle = \langle \psi | (-m)0 \rangle$  and  $\langle \psi | 00 \rangle = \langle \psi | m(-m) \rangle$ . It follows that  $\psi$  has the same amplitude on a pair of equivalent strings  $\mathbf{s} \sim \mathbf{t}$ , which means  $\langle \psi | \mathbf{s} \rangle = \langle \psi | \mathbf{t} \rangle$ . It follows that the ground subspace of  $H_n^s$  is frustration free and is spanned by the pairwise orthogonal states

$$|c_{\mathbf{x_k}}\rangle \propto \sum_{\mathbf{s} \sim c_{\mathbf{x_k}}} |\mathbf{s}\rangle$$
 (46)

where to simplify the notation, we denoted  $\mathbf{x_k} = (x_1, \dots, x_k)$ . Clearly each distinct  $\mathbf{x_k}$  results in a distinct ground state  $|c_{\mathbf{x_k}}\rangle$ .

The ground states can be highly entangled. However, among the many ground states there is a substantial subset that are all product states, i.e., k = n. We will use these to

construct quantum error correcting codes. Before doing so let us answer: How many product state ground states are there?

Let  $T_n$  be the set of all allowed 2s-ary strings  $\mathbf{t} = t_1 t_2 \dots t_n$  of length n defined by

$$T_n \equiv \{ \mathbf{t} \in \Sigma_*^n \mid \text{if } t_i = m, m \in [s], \text{ then } t_{i+1} \neq -m \}.$$
 (47)

Let  $|T_n|$  be the size of this set. Since  $T_0 = \emptyset$  and  $T_1 = \Sigma_*$ , we have that  $|T_0| = 1$  and  $|T_1| = 2s$ .

**Lemma 7.**  $|T_{n+2}| = 2s|T_{n+1}| - s|T_n|$ , with  $|T_0| = 1$  and  $|T_1| = 2s$ . We have

$$|T_n| = \frac{s^n}{2\sqrt{1 - 1/s}} \left\{ \left( 1 + \sqrt{1 - 1/s} \right)^{n+1} - \left( 1 - \sqrt{1 - 1/s} \right)^{n+1} \right\}. \tag{48}$$

Asymptotically it holds that

$$|T_n| \approx \frac{1 + \sqrt{1 - 1/s}}{2\sqrt{1 - 1/s}} \left[ s(1 + \sqrt{1 - 1/s}) \right]^n ; \qquad n \gg 1 \quad .$$
 (49)

Proof. We prove this by induction, any  $\mathbf{t} \in T_n$  is  $\mathbf{t} = t_1 \dots t_n$ , where  $t_j \in \Sigma_*$ . Since  $t_1$  can be either a m or -m for some  $m \in [s]$ ,  $\mathbf{t}$  is either  $\mathbf{t} = (-m)\mathbf{t}'$ , where the string  $\mathbf{t}' \in T_{n-1}$  or  $\mathbf{t} = m\mathbf{t}'$  where  $\mathbf{t}'$  denotes the subset of strings in  $T_{n-1}$  that do not start with the letter -m, i.e.,  $t_1' \neq -m$ . The number of strings  $\mathbf{t} = (-m)\mathbf{t}'$  with  $\mathbf{t}' \in T_{n-1}$  is clearly  $s|T_{n-1}|$ . Now the set of all string  $\mathbf{t} = m\mathbf{t}'$  with  $t_1' \neq -m$  coincides with the set that excludes the strings  $\mathbf{t} = m(-m)\mathbf{t}''$  where  $\mathbf{t}'' \in T_{n-2}$ . Since m takes on s different values, we have that the number of strings  $\mathbf{t} = m\mathbf{t}'$  with  $t_1' \neq (-m)$  is  $s(|T_{n-1}| - |T_{n-2}|)$ .

The total size of the set is then  $|T_n| = 2s|T_{n-1}| - s|T_{n-2}|$ , which is a linear recursion of second order with initial conditions  $|T_0| = 1$  and  $|T_1| = 2s$ . Shifting the indices to reflect  $|T_0|$  and  $|T_2|$  as the starting values, we have

$$|T_{n+2}| = 2s|T_{n+1}| - s|T_n|.$$

The solution is elementary and of the form  $|T_n| = Ar_+^n + Br_-^n$ , where the two roots  $r_\pm$  are  $r_\pm = s(1 \pm \sqrt{1-1/s})$  and  $A = \frac{1+\sqrt{1-1/s}}{2\sqrt{1-1/s}}$ ,  $B = \frac{-1+\sqrt{1-1/s}}{2\sqrt{1-1/s}}$  are obtained from the initial conditions  $|T_0| = 1$  and  $|T_1| = 2s$ . This proves Eq. (48) and the observation  $(1 - \sqrt{1-1/s}) < 1$  proves the asymptotic formula Eq. (49).

Comment: In the limit we have  $\lim_{s\to 1} |T_n| = n+1$ , which is the number of distinct product ground states  $|(-)_1 \dots (-)_p (+)_{p+1} \dots (+)_n \rangle$  where  $p \in \{0, 1, \dots, n\}$  with p = 0 corresponding to  $|++\dots+\rangle$ .

Corollary 1. The dimension of the Kernel of  $H_n^s$  is  $\sum_{k=0}^n |T_k| = \frac{\left(-2+r_+^{n+1}+r_-^{n+1}\right)}{2(s-1)}$ , where  $r_{\pm} \equiv s(1 \pm \sqrt{1-1/s})$ . Asymptotically we have  $\dim(\ker(H_n^s)) \approx r_+^{n+1}/[2(s-1)]$ .

*Proof.* The total number of equivalent classes is the dimension of the Kernel. For each  $k \in [n]$  there are  $|T_k|$  equivalent classes and we have

$$\dim(\ker(H_n^s)) = \sum_{k=0}^n |T_k| = \frac{r_+^{n+1} + r_-^{n+1} - 2}{2(s-1)}.$$

**Remark 3.** The fraction of product state ground states is a constant independent of n

$$\frac{|T_n|}{\sum_{k=0}^n |T_k|} = \sqrt{\frac{s-1}{s}} \left[ \frac{1 - (r_-/r_+)^{n+1}}{1 + (r_-/r_+)^{n+1} - 2r_+^{-n-1}} \right] \approx \sqrt{\frac{s-1}{s}} \quad , \quad s > 1 \quad ;$$

whereas for s = 1,  $\lim_{s \to 1} \dim(\ker(H_n^s)) = \frac{1}{2}(n+1)(n+2)$  and the fraction vanishes with the system's size as  $2/(n+2) \approx 2/n$ .

We now return to the ground space of  $H_n$ .

**Lemma 8.** Ground space of  $H = H_n^s + H_n^J$  with J > 0 coincides with the span of the equivalent classes  $|c_{\mathbf{x_n}}\rangle$ , which are all product states. The ground space dimension is  $|T_n|$ .

*Proof.* The set of FF ground states of  $H_n^s$  is given by Eq. (46) in Lemma 6. The kernel of  $J\sum_{k=1}^n(|0\rangle\langle 0|)_k$  is the span of all product states of weight n, i.e., states  $|\mathbf{t}\rangle$  where  $\mathbf{t}\in\Sigma_*^n$ . By Lemma 4 the intersection of the two is  $|c_{\mathbf{x_n}}\rangle$ , which we recall are the product states of the irreducible strings of weight n and there are  $|T_n|$  of them (Eqs. (48) and (49)).

#### 3.3 Constructing good quantum codes in the ground space

In this section we construct quantum codes that are supported on a selected subset of computational basis states that lie within the kernel of our 2-local Hamiltonian. We show that these quantum codes that encode a single logical qubit can have a linear distance.

Recall that the standard local spin states are  $|j\rangle$ , with  $j \in \Sigma = \{-s, \ldots, +s\}$ . We also define the non-zero alphabet  $\Sigma_* \equiv \{-s, \ldots, -1, +1, \ldots, s\}$ . Clearly  $|\Sigma| = 2s + 1$  and  $|\Sigma_*| = 2s$ . Define the generalized (non-Hermitian) Pauli operators in terms of these basis states as

$$X = \sum_{j \in \Sigma_c} |j\rangle\langle j + 1| , \qquad Z = \sum_{j \in \Sigma_c} \omega^j |j\rangle\langle j| ,$$

where  $\omega = \exp(2\pi i/(2s+1))$  is a root of unity, and by  $\Sigma_c$ , we mean the set  $\Sigma$  with the cyclic property that s+1=-s and -s-1=+s. We denote by  $X_k$  and  $Z_k$  the generalized Pauli operators that act on the  $k^{\text{th}}$  spin (qudit) and act trivially on the rest.

Recall that  $\ker(H_n)$  is spanned by certain product states of weight n. We denote the basis of  $\ker(H_n)$  by  $\mathcal{T}_n$  where

$$\mathcal{T}_n = \{ |\mathbf{t}\rangle : \mathbf{t} \in T_n \}$$

and  $T_n$  is defined in Eq. (47).

The quantum code that we construct will be a two-dimensional subspace of  $\mathcal{T}_n$ . In general, the logical codewords can be supported on an exponential number of basis states in  $\mathcal{T}_n$ . However, we select a subset of computational basis states labels  $C \subset T_n$  such that the minimum Hamming distance of C is at least 2t + 1, where t is the designed maximum number of correctable errors. The set of labels C has the interpretation as a classical code, and we denote its distance by  $\operatorname{dist}(C)$ :

$$dist(C) \equiv \min \left\{ dist(\mathbf{t}, \mathbf{t}') \mid \mathbf{t}, \mathbf{t}' \in T_n, \ \mathbf{t} \neq \mathbf{t}' \right\} ,$$

and

$$\operatorname{dist}(\mathbf{t}, \mathbf{t}') = |\{t_i' \neq t_i : i \in [n]\}|$$

is the usual Hamming distance between codewords  $\mathbf{t}, \mathbf{t}' \in \Sigma_*^n$ . The logical codewords of our quantum code will be supported only on the set of basis sets labeled by the classical code C.

Since the basis are product states, finding the subset C with the desired distance can be seen as a problem in classical coding theory. For example, when s = 2 one can map the elements of  $T_n$  to  $\mathbb{F}_4^n$  and use the properties of quartenary codes over  $\mathbb{F}_4^n$  to derive a classical code with the prescribed distance. For instance, one can apply the mapping

$$\varphi(1) = 0$$
,  $\varphi(-1) = 1$ ,  $\varphi(2) = a$ ,  $\varphi(-2) = b$ , (50)

where  $\mathbb{F}_4 = \{0, 1, a, b\}$ , b = a + 1 and  $a^3 - 1 = 0$ , and for every  $\mathbf{t} = t_1 \dots t_n \in T_n$ , define  $\varphi(t_1 \dots t_n) = (\varphi(t_1), \dots, \varphi(t_n))$ .

Our strategy is to construct a code C over  $\mathbb{F}_4^n$  that has a guaranteed minimum distance and delete all codewords in it that have the forbidden substrings (01) and (ab) to obtain a code C'. Then we let  $C = \varphi^{-1}(C')$ , which will be the strings that define the computational (product) basis states.

The quantum code in the ground space of our Hamlitonian,  $H_n$ , with an interaction graph given by a line-graph may be called a subspace QLDPC codes. Here, we prove that there are quantum codes within  $\ker(H_n)$  that have linear distance in n. We leverage on the existence of good binary codes. The relative distance of a code is the ratio of its distance to its length. Good binary codes are defined as binary codes with positive relative distance. To use the results from binary codes, we define a map  $\beta$  from the binary symbols 0 and 1 to 2 and 1 respectively. It is then easy to see that given any binary code C,  $\beta(C)$  is guaranteed to be a feasible subset of  $T_n$ , and hence we may use  $\beta(C)$  to construct our quantum code.

Our main theorem is:

**Theorem 2.** Let  $0 < \tau \le 1/2$  be a real and positive constant. There exist quantum codes in  $\ker(H_n)$  that encode one logical qubit and have the distance of  $2\tau n$  whenever

$$\operatorname{Ent}_{2}(2\tau) + \operatorname{Ent}_{2s+1}(2\tau)\log_{2}(2s+1) + o(1) \le 1.$$
(51)

Second, there are explicit quantum codes which encode one logical qubit with a distance of  $2\tau n$  whenever

$$1/2 - \tau/0.11 \ge \log_2(2s+1)\operatorname{Ent}_{2s+1}(2\tau). \tag{52}$$

**Remark 4.** We call the constructions given by optimizing (51) and (52) as the Gilbert-Varshamov (GV) [MS77, Chpt. 1] and Justesen construct [MS77, Chpt. 10, Thm. 11] respectively. The GV construct arises from choosing a random C, while the Justesen construct uses the classical Justesen code to define C.

*Proof.* To prove the first result, we use random coding arguments. Namely, we turn our attention to random binary codes. By the Gilbert-Varshamov bound [MS77, Chapter 1], we know that such binary codes are almost surely good binary codes. Moreover we know that for any positive integer t, there exists a classical binary code C that corrects 2t+1 errors where

$$|C| \ge 2^n / V_2(2t). \tag{53}$$

Using Lemma 2, by setting  $C = \beta(C)$ , this implies that if

$$2^{n}/V_{2}(2t) \ge 2V_{2s+1}(2t),\tag{54}$$

then there exists some quantum code encoding a single qubit in  $\ker(H_n)$  that also corrects  $t = \tau n$  errors. Since the inequality (54) is equivalent to the (51), the first result of our theorem follows.

The second result follows from using Justesen's concatenated construction, which gives binary codes of with asymptotically linear distance and positive rate [MS77][Chapter 10, Theorem 11]. More specifically, a binary Justesen code  $C_{\text{Justesen}}$  is a concatenated code, with a Reed-Solomon outer code on the finite field of dimension  $2^m$  as the outer code, and distinct inner codes each encoding m bits into 2m bits. When  $C_{\text{Justesen}}$  has a length of n, the length of the Reed-Solomon outer code is n/(2m). Each inner code is a binary code and has rate 1, and is a distinct mapping from  $\mathbb{F}_{2^m}$  to  $\mathbb{F}_2^{2m}$ . The relative distance  $\delta = d/n$  of this family of Justesen codes is given by

$$\delta > 0.110(1 - 2\log_2|C_{\text{Justesen}}|/n) + o(1).$$
 (55)

Rearranging this inequality and using  $\tau = \delta/2$ , we get

$$\log_2 |C_{\text{Justesen}}| \ge n(1/2 - \tau/0.11 + o(1)). \tag{56}$$

Hence the number of codewords of a Justesen code with distance 2t + 1 is asymptotically at least  $2^{n(1/2-9.1\tau+o(1))}$ . Now we set  $C = \beta(C_{\text{Justesen}})$ , and use Lemma 2 to find that a quantum code in the ground space of  $\ker(H_n)$  that corrects asymptotically  $\tau n$  errors exists whenever  $\tau < (2s)/(2s+1)$  and the inequality (52) holds.

We plot the attainable values of  $\tau$  for different values of spins in Fig. 6.

In the next two sections we illustrate explicit quantum codes on 8 and 6 qudits respectively. These quantum code were obtained from punctured variants of the classical ternary Golay code, where by punctured we mean that the first three symbols of the code were ignored for the 8 qudit code, and five symbols were ignored for the 6 qudit code.

#### 3.3.1 A ground subspace Steane code that corrects a single error

By slightly modifying the 7-qubit Steane code, we can embed it in the ground space of  $H_7$ . Namely, on the set of computational basis vectors, we can apply the map  $|0\rangle \rightarrow |1\rangle$  and  $|1\rangle \rightarrow |2\rangle$ . This also corresponds to the quantum code with logical codewords

$$|0_{L}\rangle = \frac{1}{\sqrt{8}} (|1, 1, 1, 1, 1, 1, 1, 1, 1\rangle + |1, 1, 1, 2, 2, 2, 2\rangle + |1, 2, 2, 1, 2, 2, 1\rangle + |1, 2, 2, 2, 1, 1, 2\rangle + |2, 1, 2, 1, 2, 1, 2\rangle + |2, 1, 2, 2, 1, 2, 1\rangle + |2, 2, 1, 1, 1, 2, 2\rangle + |2, 2, 1, 2, 2, 1, 1\rangle)$$
(57)  
$$|1_{L}\rangle = \frac{1}{\sqrt{8}} (|1, 1, 2, 1, 1, 2, 2\rangle + |1, 1, 2, 2, 2, 1, 1\rangle + |1, 2, 1, 1, 2, 1, 2\rangle + |1, 2, 1, 2, 1, 2, 1\rangle + |2, 1, 1, 1, 2, 2, 1\rangle + |2, 1, 1, 2, 1, 1, 2\rangle + |2, 2, 2, 1, 1, 1, 1\rangle + |2, 2, 2, 2, 2, 2, 2\rangle).$$
(58)

Using the KL criteria, we can verify that this quantum code corrects any single error.

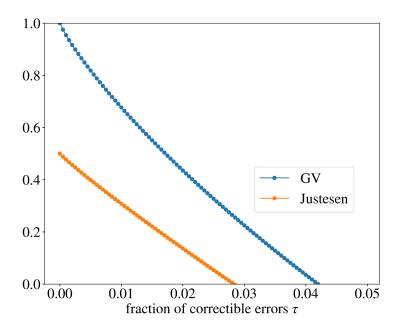


Figure 6: The vertical axis is a lower bound on the size of the kernel of the A matrix in log-scale. The horizontal axis is  $\tau = t/n$ , where t denotes the number of correctable errors for our quantum code in the ground space of  $\ker(H_n)$ . The length n is taken to be asymptotically large. This demonstrates that there are linear distance quantum codes in the ground space of the frustration-free 2-local Hamiltonian  $H_n$ .

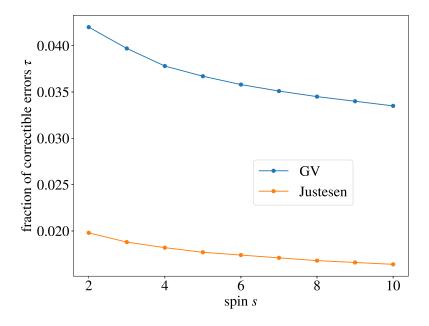


Figure 7: The vertical axis is  $\tau = t/n$ , where t denotes the number of correctable errors for our quantum code in the ground space of  $\ker(H_n)$ . The horizontal axis is the spin number our spin-chain.

#### 3.3.2 A ground subspace code on eight spins that corrects a single error

Here, we give an example of a quantum code that lies in the kernel of  $H_n^s$  for n=8 and s=2. This quantum code encodes a single logical qubit, corrects an arbitrary one-qubit error, and has logical codewords

$$|0_{L}\rangle = (|\phi_{0}\rangle|\theta_{0}\rangle + |\phi_{1}\rangle|\theta_{1}\rangle + |\phi_{2}\rangle|\theta_{2}\rangle + |\phi_{3}\rangle|\theta_{3}\rangle + |\phi_{4}\rangle|\theta_{4}\rangle + |\phi_{5}\rangle|\theta_{5}\rangle)/\sqrt{6}$$

$$|1_{L}\rangle = (|\phi_{1}\rangle|\theta_{4}\rangle + |\phi_{0}\rangle|\theta_{3}\rangle + |\phi_{3}\rangle|\theta_{0}\rangle + |\phi_{2}\rangle|\theta_{5}\rangle + |\phi_{5}\rangle|\theta_{2}\rangle + |\phi_{4}\rangle|\theta_{1}\rangle)/\sqrt{6}, \tag{59}$$

where

$$|\phi_{0}\rangle = |1, 1, 1, -2\rangle, |\phi_{1}\rangle = |1, -2, -1, -1\rangle, |\phi_{2}\rangle = |-1, -2, -2, -1\rangle, |\phi_{3}\rangle = |-1, -1, 1, 1\rangle, |\phi_{4}\rangle = |2, -1, -1, 1\rangle, |\phi_{5}\rangle = |2, 1, -2, -2\rangle,$$
(60)

and

$$|\theta_{0}\rangle = |-2, 2, 2, 1\rangle, |\theta_{1}\rangle = |1, -2, -2, -2\rangle, |\theta_{2}\rangle = |-1, 2, 2, 1\rangle, |\theta_{3}\rangle = |-2, -2, -2, -2\rangle, |\theta_{4}\rangle = |1, 2, 2, 1\rangle, |\theta_{5}\rangle = |-1, -2, -2, -2\rangle.$$
(61)

The KL criteria for correcting a single error using this quantum code are satisfied. Also it easy to see that this code is not a CWS code, and hence gives an example of a quantum code that falls outside of the CWS, stabilizer and CSS quantum coding formalisms.

Next we point out that the quantum code in (59) has a concatenated structure. The logical codewords of the outer code, given in (59) are simply maximally entangled states on two six-level systems.

From the structure of the inner codes, it is clear that to perform a logical bit-flip on our quantum code, it suffices to induce the transition  $|-2, -2, -2\rangle \leftrightarrow |2, 2, 1\rangle$  on the last three spins. Performing other logical computation operations is significantly more complicated and we leave this for future work.

#### 3.3.3 A ground subspace code that detects a single error

We also construct an error detecting quantum code (distance equal to 2) on six spins with s = 2 using the logical operators

$$|0_L\rangle = \frac{|1, 1, 2, 1, -2, 1\rangle + |-2, 1, -2, -2, 2\rangle}{\sqrt{2}}$$
 (62)

$$|1_L\rangle = \frac{|1, 1, -2, -2, 2, 1\rangle + |-2, 1, 2, 1, -2, 2\rangle}{\sqrt{2}}.$$
 (63)

Such a construction is not unique, and we have many other error detecting codes on six spins.

#### Acknowledgement

RM acknowledges funding from the MIT-IBM Watson AI Lab under the project *Machine Learning in Hilbert space*. The research was supported by the IBM Research Frontiers Institute. YO acknowledges support from the EPSRC (Grant No. EP/M024261/1) and the QCDA project (Grant No. EP/R043825/1) which has received funding from the QuantERA ERANET Cofund in Quantum Technologies implemented within the European Union's Horizon 2020 Programme.

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