Fault Tolerant Quantum Computation with Very High Threshold for Loss Errors

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Many proposals for fault tolerant quantum computation (FTQC) suffer detectable loss processes. Here we show that topological FTQC schemes, which are known to have high error thresholds, are also extremely robust against losses. We demonstrate that these schemes tolerate loss rates up to 24.9%, determined by bond percolation on a cubic lattice. Our numerical results show that these schemes retain good performance when loss and computational errors are simultaneously present.

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One of the most important achievements of quantum information theory was the discovery of fault tolerant quantum computation (FTQC): arbitrarily precise, scalable quantum computations can be performed using error-prone components, as long as the error rate is below a certain threshold [1,2]. Current thresholds, around $p_{\rm thres} \sim 10^{-2}$ per elementary operation [3,4], come tantalizingly close to the experimental state of the art. Of particular interest is the beautiful proposal developed by Raussendorf, Harrington, and co-workers [4–7] using ideas from topological quantum computing [8] and with only nearest neighbor gates.

Many proposals for quantum computing suffer qubit loss, such as photon loss, atom or ions escaping from traps, or, more generally, the leakage of a qubit out of the computational basis in a multilevel system. However, such errors can be detected and located without affecting the state of the remaining qubits. It is therefore expected that appropriately tailored FTQC schemes can tolerate higher rates for loss errors than for unlocatable errors [9]. Certain schemes have loss thresholds of $p_{\rm loss} < 0.5$ [10,11], although it is unclear how they perform in the presence of computational errors (unlocated errors acting within the computational subspace). Other FTQC proposals, tolerant to both error types, have thresholds of $p_{\rm loss} \lesssim 3 \times 10^{-3}$ and $p_{\rm comp} \lesssim 10^{-4}$ [12,13].

In this Letter, we describe a FTQC scheme which tolerates both loss and computational errors with very high thresholds. The thresholds are characterized by a contour in (p_{loss}, p_{comp}) parameter space passing through the points (0.249, 0) and (0, 0.0063). This represents an improvement of almost 2 orders of magnitude over earlier results [12,14]. Our approach requires only that losses are detected at the final readout stage.

Our scheme combines methods from Raussendorf's topological scheme [4–6] and our previous work on loss tolerance in surface codes [15,16]. As in the latter, the loss tolerance threshold follows from the bond percolation threshold on the relevant lattice, here the cubic lattice [17].

To this end, we develop and implement a new classical algorithm to analyze the syndrome in the presence of both loss and logical errors. Monte Carlo simulations of the resulting FTQC scheme yield an estimate of the ($p_{\rm loss}$, $p_{\rm comp}$) threshold contour. This contour exhibits a moderate tradeoff between $p_{\rm loss}$ and $p_{\rm comp}$, indicating that the code is robust against both kinds of error.

We first give an overview of Raussendorf's scheme for FTOC. The scheme is based on the one-way quantum computer: that is, a *cluster state* $|C\rangle_L$, of many physical qubits located on the faces and edges of a cubic primal lattice \mathcal{L} , is prepared by first preparing every physical qubit in the state $|+\rangle$, and then applying controlled-PHASE (CPHASE) gates between neighboring qubits on the lattice [see Fig. 1(a)]. It is convenient to introduce the dual lattice \mathcal{L}^* whose vertices, edges, faces, and cubes correspond, respectively, to the cubes, faces, edges, and vertices of \mathcal{L} . Physical qubits reside only on the faces and edges of these lattices. Computation proceeds by making a sequence of single qubit measurements on $|C\rangle_{\mathcal{L}}$, aided by classical processing of the measurement outcomes. Correlations between the measurement outcomes give rise to fault tolerance.

 \mathcal{L} is subdivided into three regions, denoted V, D, and S. Qubits in V and D are measured in the X and Z bases, respectively. Qubits in S are measured in either the Y or $(X+Y)/\sqrt{2}$ basis. D comprises a collection of thick structures with some characteristic transverse diameter (de-fects), embedded in V [Fig. 1(b)]. Each pair of defects encodes one logical qubit. As the D qubits are measured in the Z basis, one can omit them altogether, which we assume for the remainder of this Letter.

The topology of the braiding of defect regions effect certain Clifford gates between the logical qubits, while measurement results in *V* provide topologically protected fault tolerant error correction of the logical Clifford gates to arbitrarily high accuracy. *S* comprises a collection of well-separated single qubits, spread out among the defects.

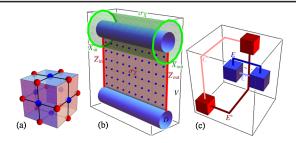


FIG. 1 (color online). (a) Unit cell showing qubits centered on faces and edges of the primal lattice. Heavy lines indicateCPHASE gates. (b) A primal identity gate showing logical input operators $\bar{X}_{\rm in}$ and $\bar{Z}_{\rm in}$ and output operators $\bar{X}_{\rm out}$ and $\bar{Z}_{\rm out}$, and the correlation surfaces $\sigma_{\bar{X}}$ and $\sigma_{\bar{Z}}$, that interpolate between them. Also shown are physical qubits (spheres) living on the correlation surface between the \bar{Z} operators, and the defect regions D embedded in the bulk vacuum V. (c) Failed syndromes for error chains on the primal E and dual E^* lattices, and candidate correction chains C and C^* .

The S qubits are used to introduce noisy encoded *magic* states into the circuit, which can then be distilled using the very accurate logical Clifford gates in V and D, implementing a universal gate set [18].

Magic state distillation tolerates a rather large amount of noise, and so it is found that the thresholds for both loss and computational errors are set by the corresponding thresholds in the bulk topological region, V and D. We therefore discuss error correction in V and D in more detail, focusing on the implementation of a fault tolerant identity gate. The details of nontrivial Clifford gates are described in [4,19].

The layout of the identity gate is show in Fig. 1(b). The boundaries of D are positioned such that faces of L lie just inside the defect regions. The alignment of the defects specify a "simulated time" axis, such that the identity gate maps the logical qubit from the input (rear) plane I to the output (front) plane, separated by an integer number of unit cells. Within these planes, the qubit is encoded in a surface code [8], containing two holes, which coincide with the intersection of the defect regions with the corresponding plane. A single logical qubit is associated to the pair of holes, with encoded logical operators \bar{X} and \bar{Z} as indicated on Fig. 1(b).

To demonstrate the operation of the noiseless gate, it suffices to show that the input logical operators $\bar{X}_{\rm in}$, $\bar{Z}_{\rm in}$ are mapped to the output operators $\bar{X}_{\rm out}$, $\bar{Z}_{\rm out}$, under the action of the CPHASE gates and single qubit measurements [20]. To this end, we introduce the cluster stabilizer operators, $K_f = X_f \bigotimes_{e \in \partial f} Z_e$, associated with each face f of \mathcal{L} or \mathcal{L}^* , where ∂f denotes the qubits at the edges of f. The initial state of the cluster satisfies $K_f |C\rangle_{\mathcal{L}} = |C\rangle_{\mathcal{L}}$. Consider the surface $\sigma_{\bar{Z}}$ illustrated in Fig. 1(b), consisting of faces in \mathcal{L} . The product of the cluster stabilizers on $\sigma_{\bar{Z}}$ yields $K(\sigma_{\bar{Z}}) = \bigotimes_{f \in \sigma_{\bar{Z}}} X_f \bigotimes_{e \in \partial \sigma_{\bar{Z}}} Z_e$, where $\partial \sigma_{\bar{Z}}$ denotes the qubits at the boundary of $\sigma_{\bar{Z}}$. Since $\bigotimes_{e \in \partial \sigma_{\bar{Z}}} Z_e = \bar{Z}_{\rm in} \bigotimes_{\bar{Z}_{\rm out}}$, it follows that measurements in the bulk effects the operator mapping $\bar{Z}_{\rm in} \to \pm \bar{Z}_{\rm in} \otimes \bar{Z}_{\rm out}$, where the sign is

given by the parity of the measurement outcomes $\Pi_{f \in \sigma_{\bar{Z}}} x_f$.

Similarly, the product of cluster stabilizers acting on the surface $\sigma_{\bar{X}}$, consisting of faces in \mathcal{L}^* , yields $K(\sigma_{\bar{X}}) = \bigotimes_{f^* \in \sigma_{\bar{X}}} X_{f^*} \bigotimes_{e^* \in \partial \sigma_{\bar{X}}} X_{e^*}$, and since $\bigotimes_{e^* \in \partial \sigma_{\bar{X}}} X_{e^*} = \bar{X}_{\mathrm{in}} \otimes \bar{X}_{\mathrm{out}}$ it follows that measurements on $\sigma_{\bar{X}}$ effect the mapping $\bar{X}_{\mathrm{in}} \to \pm \bar{X}_{\mathrm{in}} \otimes \bar{X}_{\mathrm{out}}$. Finally, measuring the qubits in I leads to the desired transformation (up to an unimportant Pauli frame update) $\bar{Z}_{\mathrm{in}} \to \pm \bar{Z}_{\mathrm{out}}$ and $\bar{X}_{\mathrm{in}} \to \pm \bar{X}_{\mathrm{out}}$.

Error syndromes are revealed by correlations of measurement outcomes. In particular, considering the product of K_f 's centered on the faces of a unit cube c of \mathcal{L} leads to the parity check $P_c = \prod_{f \in \partial c} x_f$. Absent any errors, the parity checks satisfy $P_c = +1$. Failed parity checks, i.e., cubes which instead satisfy $P_c = -1$, reveal the end points of *chains*, E, of E errors [see Fig. 1(c)]. These chains reside on edges of E, and are corrected by inverting the recorded measurement outcomes on some correction chain E (i.e., E in E for E in E sharing the same end points as E (i.e., E in E in

Finding a suitable correction chain *C* requires an algorithm for pairing failed parity checks, then finding a suitable path in the lattice between each pair of syndromes. Choosing *C* optimally is computationally difficult, since it is equivalent to minimizing the free energy of a particular spin-glass model [6,21]. However, efficient heuristic methods for decoding error syndromes have been developed, including those using Edmonds' perfect matching algorithm [22,23], and belief propagation [24].

Realistic error models depend on the particular implementation under consideration. As in [7] we assume single qubit depolarizing errors take place during preparation, storage, and measurement, with rates p_P , p_S , and p_M respectively, while CPHASE gates suffer two-qubit depolarizing errors with rate p_2 . For numerical calculations, we assume these errors are independent, identically distributed (iid) and occur with equal probability, p_{comp} , i.e., $p_P =$ $p_S = p_M = p_2 = p_{\text{comp}}$. These parameters may vary, with modest changes in the threshold [7]. For this error model, previous results established an error threshold of p_{comp} < $p_t = 0.0058$ [7]. Note that without losses, exploiting error correlations within each sublattice improves the threshold to $p_t = 0.0075$ [5]. A further improvement may be possible by exploiting correlations between \mathcal{L} and \mathcal{L}^* [24]. For simplicity, we do not exploit correlations, at the cost of a marginal reduction in the threshold.

We assume losses are iid with total rate $p_{\rm loss}$, and only occur either before or after all CPHASE gates have acted upon each qubit, but not at intermediate times. This is consistent with, e.g., optical lattices where loading losses [25] are much higher than storage losses [26], or with a photonic implementation where source and detector inefficiencies are expected to dominate. CPHASE gates acting

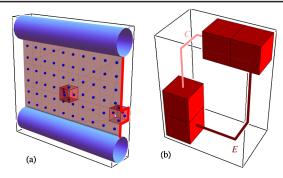


FIG. 2 (color online). (a) A correlation surface suffering loss is repaired by deforming around lost qubits (white spheres), by multiplying by the corresponding check operator. An affected logical input or output operator is similarly deformed. (b) In the bulk, supercheck operators are formed by products of elementary check operators that are affected by a loss. Error chains terminate within a supercheck operator.

on a lost input are assumed to implement the identity operation, plus depolarizing noise at rate p_2 , on the remaining qubit. Losses need only be detected at the measurement step, so loss detection is permitted to be a destructive process. Loss detection may be imperfect, but such errors are accounted for in p_M or p_{loss} .

To deal with these losses, we adapt the Edmonds' matching approach, paralleling [15,16]. Suppose qubit q is lost. The two parity checks P_{c_q} and $P_{c_q'}$ on the adjacent cubes c_q and $c_{q'}$ are incomplete and so cannot detect error syndromes. Likewise any correlation surface σ that depends on q is damaged and so cannot mediate logical gates.

We recover from this damage in two ways. First, multiplying $K(\sigma)$ by $K(\partial c_q) = \bigotimes_{f \in \partial c_q} X_f$, yields a new surface $\tilde{\sigma}$ that avoids q [see Fig. 2(a)], i.e., $K(\tilde{\sigma}) = K(\sigma)K(\partial c_q)$ is independent of q. Then $\tilde{\sigma}$ determines an equivalent mapping of input operators to output operators. This procedure can be iterated for all of the lost qubits, and succeeds as long as there is some 2D correlation surface that percolates between the logical qubit operators while completely avoiding lost qubits. In the limit where the defect dimensions and spacings between defects becomes large, such a surface can be found providing the loss rate is lower than the cubic-lattice bond percolation threshold, which is approximately 0.249 [17].

Second, the product of the two damaged check operators $\tilde{P}_q = P_{c_q} P_{c_q'}$ is independent of q, so the sign of \tilde{P}_q yields information about ∂E , as illustrated in Fig. 2(b). Thus for each instance of losses, we form new *supercheck* operators that are insensitive to the losses. The lattice of check operators is then no longer cubic, so we perform the syndrome matching on an irregular lattice.

In the context of surface codes, Edmonds' matching has typically been used to find maximum likelihood correction chains C that maximize the probability $P(C|\partial E)$ [4,22]. It is simple to show that such chains are also of minimum length [22]. For a typical pattern of loss, it turns out that

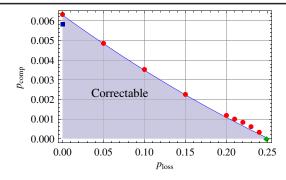


FIG. 3 (color online). Phase diagram showing correctable region of parameter space. Red circles are numerically calculated computational error thresholds, p_t at different loss rates, for an error model in which $p_P = p_S = p_M = p_2 = p_{\rm comp}$. The green diamond is the 3D bond percolation threshold [17]. The blue square is the threshold computed in [7], which ignores degeneracy in the matching algorithm.

there are very many minimum length chains. Furthermore, as discussed in [15] some minimum-distance pairings have many more matching chains than other minimum-distance pairings, and so they are *a priori* much more likely. Accounting for this degeneracy in the pairing is computationally cheap, and leads to modest improvements in the threshold. In the 3D FTQC scheme on an irregular lattice, the degeneracy of matchings may also be computed efficiently by modifying Dijkstra's algorithm for finding shortest paths on a graph.

To establish quantitative error thresholds in the presence of losses, we perform Monte Carlo simulations on L^3 cubic lattices with periodic boundary conditions. For each simulation we do the following:

- (1) Generate an instance of errors E and losses.
- (2) From losses, derive \tilde{P} and $\tilde{\sigma}$. If losses percolate the cubic lattice then FAILURE (no such $\tilde{\sigma}$ exists).
 - (3) From E, derive ∂E on the irregular lattice.
 - (4) From ∂E , compute the correction chain C.
- (5) Compute the homology class \mathcal{H}_{E+C} (count the number of intersections of E+C with $\tilde{\sigma}$, mod 2).
 - (6) If $\mathcal{H}_{E+C} = 0$ then SUCCESS. Otherwise FAILURE.

At a given loss rate, we generate 198 000 instances of errors while varying p_{comp} and $L=6,8,\ldots,16$. For each loss rate, we fit a Taylor's expansion of a universal scaling function $p_{\text{fail}}(p_{\text{loss}},p_{\text{comp}},L)\approx a+bx+cx^2$, where $x=(p_{\text{comp}}-p_t)L^{1/\nu}$, with fitting parameters p_t (the loss-dependent computational error threshold), ν (the scaling exponent), a,b, and c.

Figure 3 is the central result in this Letter, and shows p_t at different values of $p_{\rm loss}$ (red circles). The shaded region indicates the correctable region of parameter space; if $p_{\rm comp} < p_t$, then errors are almost always correctable (in the limit $L \to \infty$). The solid curve is simply a quadratic fit to the computed values of p_t for $0 \le p_{\rm loss} \le 0.15$ [27], and extrapolates to $p_{\rm loss} = 0.252 \pm 0.005$ at $p_{\rm comp} = 0$, which agrees with the percolation threshold $p_{\rm perc} = 0.249$ [17] (green diamond) on a 3D cubic lattice.

The loss-tolerant thresholds above rely on properties of the FTQC scheme in the bulk V and D regions, well away from S qubits. Special attention must be given to the effect of losses affecting S qubits (and also to nearby V and D regions), which are used to inject encoded *magic states* into the circuit. Computational errors close to the S qubits increase the effective noise on the encoded magic states to $p_S \approx 6p_{\rm comp}$, but the threshold for magic state purification is so high that the overall threshold is still set by that of the bulk [4]. This remains true with losses, and we now describe a scalable *postselection* method which, while profligate, demonstrates that losses near S qubits do not limit the thresholds.

If a loss occurs either on or near an S qubit, we discard the corresponding encoded magic state by measuring the encoded state, i.e., fusing the neighboring defect strands together. This may be performed in the bulk region, provided the error rates are below the corresponding bulk thresholds. Thus, we require that the only magic state qubits which are injected into the circuit are those for which no losses occur within a $d \times d \times d$ volume centered on the S qubit itself. Then the postselected error rate on the encoded magic states is $p_S \approx 6p_{\rm comp} + p_{\rm fail}(p_{\rm loss}, p_{\rm comp}, d)$. Below the bulk error threshold, $p_{\rm fail}$ is exponentially suppressed in d. The magic state distillation threshold $p_S < 1/\sqrt{35}$ [18] fixes d to be a (small) constant and so the additional overhead for this postselected scheme, $\sim (1-p_{\rm loss})^{d^3}$, is independent of the size of the algorithm, ensuring scalability.

In this Letter we have established high thresholds FTQC for both computational errors and loss errors. This result, which sets an important benchmark for experimental implementations, follows from the large redundancy of the surface codes that underpin Raussendorf's FTQC scheme. This redundancy serves two complementary purposes: to enable deformed correlation surfaces to mediate logical gates between encoded qubits, and to enable expanded parity check operators to identify end points of computational error chains.

A number of lines of enquiry follow from this work. We assumed deterministic CPHASE gates, so relaxing this to allow heralded, nondeterministic two-qubit gates in the cluster preparation [28–30] is open (but see [31]). Also, we expect that losses which occur *during* cluster construction will reduce thresholds further, since these induce additional *Z* errors on neighboring qubits. Since these are located errors, percolation phenomena still determine the loss threshold, so we anticipate this will remain substantially higher than the computational error threshold. Finally, it is important to develop methods to deal with losses on *S* qubits with lower overhead than that presented here, perhaps by dynamically routing defect regions and *S* qubits around any revealed losses.

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