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Quantum storage in quantum ferromagnets

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We must protect inherently fragile quantum data to unlock the potential of quantum technologies. A pertinent concern in schemes for quantum storage is their potential for near-term implementation. Since Heisenberg ferromagnets are readily available, we investigate their potential for robust quantum storage. We propose to use permutation-invariant quantum codes to store quantum data in Heisenberg ferromagnets, because the ground space of any Heisenberg ferromagnet must be symmetric under any permutation of the underlying qubits. By exploiting an area law on the expected energy of Pauli errors, we show that increasing the effective dimension of Heisenberg ferromagnets can improve the storage lifetime. When the effective dimension of Heisenberg ferromagnets is maximal, we also obtain an upper bound for the storage error. This result relies on perturbation theory, where we use Davis' divided difference representation for Fréchet derivatives along with the recursive structure of these divided differences. Our numerical bounds allow us to better understand how quantum memory lifetimes can be enhanced in Heisenberg ferromagnets.

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Introduction. Decoherence quickly renders unprotected quantum data unreliable. To combat this, it becomes necessary to encode quantum data into quantum error correction codes. The challenge in designing robust quantum memories arises from the difficulty of simultaneously (i) utilizing an easily accessible physical system, (ii) having a quantum code that lies within the ground space of the system's Hamiltonian H , and (iii) having an increased storage lifetime τ with an increasing number of qubits N in the physical system. Self-correcting quantum memories [1,2] should satisfy (ii) and (iii), but are challenging to implement in a multitude of desirable settings [3–14]. Indeed, constraint (i) does easily not hold, which frustrates the design of reliable quantum storage. For instance, quantum memories based on stabilizer codes which correct at least one error and also satisfy (ii) unfortunately reside in unphysical systems with many-body interactions, and can only be approximately constructed [4,15–17]. Of these constraints, it is most pertinent to satisfy (i), because physically unrealistic quantum memories will be difficult to engineer.

There are two reasons to store quantum data within the ground space and thereby satisfy constraint (ii). First, a growing energy gap can suppress excitations from the ground space [18]. Second, storing quantum data within the ground space avoids unnecessary errors that can occur even in the complete absence of noise. Any state within the ground space is an eigenstate of the physical system, and for such states, they are left unchanged by a unitary operation U_τ that the system's natural dynamics induces, after a storage time of τ elapses. By avoiding the need to uncompute U_τ , we would not suffer

from an imperfect reversal of U_τ caused by our imprecise knowledge of τ .

Storage within the ground space, while satisfying constraint (ii), is not enough to result in self-correcting quantum memories and thereby satisfy constraint (iii). Moreover, many physically realistic systems satisfying constraint (i) comprising of two-local terms are surprisingly incompatible with constraint (iii) [12]. However, this no-go result does not preclude physical systems comprising of noncommuting two-body interactions from satisfying constraint (iii). Consequently, determining whether such physical systems can satisfy constraint (iii) is especially pertinent. In this paper we study Heisenberg ferromagnets as a media for quantum storage because they comprise of noncommuting two-body interactions and therefore sidestep the no-go result of [12]. We also study to what extent Heisenberg ferromagnets satisfy constraint (iii).

The Heisenberg ferromagnet (HF) is a model of quantum magnetism, and is prevalent in many naturally occurring physical systems, and thereby satisfies constraint (i). For instance, the HF is found in various cuprates [19,20], in solid helium-3 [21], and more generally in systems with interacting electrons [22]. Even in many physical systems that cannot be naturally interpreted as ferromagnets, effective HFs can nonetheless be engineered, for instance by symmetrizing systems dominated by dipole interactions using dynamic pulse sequences [23]. Effective HFs have also been engineered in ultracold atomic gases [24] and quantum dots [25]. Specifically, we study spin-half HFs in the absence of an external magnetic field, with Hamiltonian of the form

$$H = - \sum_{\langle i,j \rangle \in E} J(\mathbf{1} - \pi_{i,j}). \quad (1)$$

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Here $\mathbf{1}$ denotes an N -qubit identity operator, $\pi_{i,j}$ denotes a swap operator on the i th and j th qubits, J denotes the exchange constants, and E denotes the set of interactions. Such HFs have $J > 0$ and ground state energy set to zero.

By storing quantum data using permutation-invariant (PI) codes in HFs, we automatically satisfy constraint (ii). This is because symmetric states lie within the ground space of any HF, and quantum data in PI codes, by being invariant under any permutation of their underlying qubits, are symmetric states. Such codes are well studied both in the context of perfect quantum error correction [26–29], and approximate quantum error correction [28,30,31]. PI codes have not only been studied in the qubit setting, but have also recently been considered in bosonic settings [32]. While prior research on PI codes shows that quantum error correction is possible within the ground space of HFs, this is only suggestive that constraint (iii) can be compatible with PI codes. This is because the coding parameters of PI codes alone, being independent of the parameters in HFs, are not enough to determine what happens when physical noise applies to PI codes stored in HFs. To better understand the extent in which HFs with PI codes can satisfy constraint (iii), we study bounds on the storage error of PI codes under the action of two different noise models, where both bounds depend on properties of the underlying HF.

Our first noise model applies to HFs of any geometry, and introduces Pauli errors. These Pauli errors occur with probabilities that are thermodynamically related to their expected energies on the codespace of a specific family of PI codes [28]. To derive an upper bound on the storage error, we find an area law on the expected energy of a Pauli error, which demonstrates that a quantum memory based in a HF can have a macroscopic energy barrier for Pauli errors. This allows us to show that the storage error decreases with increasing dimensionality of the HF.

Our second noise model introduces unitary errors probabilistically, where each unitary arises from a local perturbation of the underlying Hamiltonian. Such a noise model can describe the effects of unwanted physical interactions, such as spurious local fields afflicting each particle independently. We use perturbation theory to bound the storage error by using Davis' divided difference representation of these taking Fréchet derivatives. Because we require complete knowledge of the Hamiltonian's spectrum, we restrict our analysis to exactly solvable mean-field HFs. In such HFs every pair of spins interacts with equal strength. Since such HFs have an infinite effective dimension, analyzing them is indicative of the ultimate limits of robust quantum storage in HFs.

With respect to both noise models, we provide upper bounds for the storage error of quantum memories in HFs that are numerically tractable. In both cases we find that quantum memories in HFs are partially self-correcting in the sense that is an optimal system size for fixed noise parameters that minimizes our upper bounds on the storage error.

Energy of Pauli errors and their geometry. We use GNU codes [28] to elucidate the dependence of a HF's dimension with respect to the storage error of quantum data. GNU codes depend on three parameters g , n , and u , and encode a single qubit into $N = gnu$ qubits. Here g and n quantify the distance of the GNU code with respect to bit-flip and phase-flip errors,

respectively, while u is a scaling parameter where $u \geq 1$. When $g = n = 2t + 1$, the GNU code corrects t errors. A GNU code has logical codewords

$$|r_L\rangle = \sum_{\substack{0 \leq j \leq n \\ \text{mod}(j,2)=r}} \sqrt{\frac{\binom{n}{j}}{2^{n-1}}} |D_{gj}^{gnu}\rangle, \quad (2)$$

where $r = 0, 1$, and $|D_w^N\rangle$ are Dicke states of weight w [33,34].¹

We quantify a HF's dimension using properties of its underlying graph of interactions [35]. This graph G has vertices labeled from 1 to N , and edges E that correspond to the interactions in the HF's Hamiltonian H . Given a subset S of $\{1, \dots, N\}$, let $\partial_E S$ denote its edge boundary with respect to the edge set E , which is the set of edges in E with exactly one vertex in S . When every subset S satisfies the inequality $|\partial_E S| \geq c \min(|S|, N - |S|)^{1-1/\delta}$, the graph and HF have dimension δ with isoperimetric constant c .

Given a set \mathcal{P} of N -qubit Pauli errors that afflict at most $N/2$ qubits, let

$$\langle P \rangle = \min_{|\psi\rangle \in \mathcal{C}} \langle \psi | PHP | \psi \rangle \quad (3)$$

denote the minimum expected energy of $P \in \mathcal{P}$ on the code \mathcal{C} . When \mathcal{C} is a GNU code, we derive a lower bound on $\langle P \rangle$ in terms of the edge boundary $V(P)$, where $V(P)$ denotes the set of vertices on which P acts nontrivially. In particular, Theorem 1 below gives an area law on the minimum size of $\langle P \rangle$, which we prove in the Supplemental Material [36].

Theorem 1. Let \mathcal{C} be an N -qubit GNU code with parameters $g = n = 2t + 1$ and $u = 2$, where $N = 2(2t + 1)^2$ and $t \geq 1$. Then with respect to the Hamiltonian H given by (1) with exchange constants J and set of interactions E , for every N -qubit Pauli P in \mathcal{P} , we have

$$\langle P \rangle \geq \chi J |\partial_E [V(P)]|,$$

where $\chi = \min\{2\mu, 1 - 4\mu\}$ and

$$\mu = (1 + 5t + 6t^2)/(4 + 32t + 32t^2).$$

The significance of Theorem 1 lies in the geometric interpretation it imparts to $\langle P \rangle$. Namely, when the graph G has dimension δ and isoperimetric number c , we have the isoperimetric inequality

$$\langle P \rangle \geq J \chi c |P|^{1-1/\delta}, \quad (4)$$

where $|P| = |V(P)|$ denotes the weight of P . For a HF on a one-dimensional (1D) spin chain, $\langle P \rangle \geq J \chi$. For a HF on a square lattice [37], with our result implies that $\langle P \rangle \geq J \chi \sqrt{|P|}$. Whenever $\delta > 1$, the expected energy of P grows with its weight, and we have a macroscopic energy barrier [38]. This suggests that when $\delta > 1$, HFs can be good quantum memories. To see this, consider a noisy channel \mathcal{T} that introduces Pauli errors $P \in \mathcal{P}$ with probability proportional to $e^{-\beta\langle P \rangle}$ and

¹Dicke states are uniform superpositions of computational basis states labeled by binary vectors of Hamming weight w .

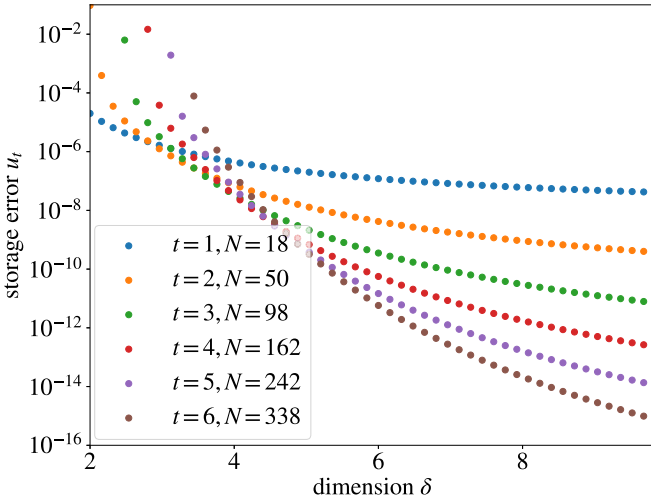


FIG. 1. When a HF stores an encoded qubit within a GNU code on $N = 2(2t + 1)^2$ physical qubits that corrects t errors, we use (7) to obtain upper bounds on the storage error u_t with respect to the HF's dimension δ and t . Here $\beta J = 13$ and $c = 1$.

with effective inverse temperature β . Explicitly,

$$\mathcal{T}(\rho) = \sum_{P \in \mathcal{P}} (e^{-\beta(P)} / \mathcal{Z}) P \rho P, \quad (5)$$

where $\mathcal{Z} = \sum_{P \in \mathcal{P}} e^{-\beta(P)}$. The corresponding probability of obtaining an uncorrectable error, which is the storage error under the assumption of perfect error correction, is then

$$u_t = \frac{1}{\mathcal{Z}} \sum_{\substack{P \in \mathcal{P} \\ |P| \geq t+1}} e^{-\beta(P)}. \quad (6)$$

From the isoperimetric inequality (4) and the bound $|\partial_E V(P)| \leq \Delta |P|$, where Δ is the maximum vertex degree of G , we obtain

$$u_t \leq \left(\sum_{w=t+1}^{N/2} \binom{N}{w} 3^w e^{-\beta J \chi c w^{1-1/\delta}} \right) \left(\sum_{w=0}^{N/2} \binom{N}{w} 3^w e^{-\beta J \Delta w} \right)^{-1}. \quad (7)$$

We illustrate (7) in Fig. 1 with $\Delta = 4$, $c = 1$, $\beta J = 13$, and vary the dimension δ and number of correctible errors t . We see that increasing δ decreases u_t . Moreover, when $2 \leq \delta < 4$, the optimal PI code has $1 \leq t \leq 4$. This shows that for low-dimensional systems with $\delta \geq 2$, our is partially self-correcting, where increasing the system size cannot indefinitely reduce the storage error.

Random coherent noise and storage error. A good quantum memory preserves entanglement. Given a quantum code \mathcal{C} with logical codewords $|0_L\rangle, \dots, |(M-1)_L\rangle$, consider the entangled state $|\Psi_{\mathcal{C}}\rangle = \sum_{j=0}^{M-1} |j\rangle \otimes |j_L\rangle / \sqrt{M}$. The storage error of \mathcal{C} with respect to a noisy channel \mathcal{N} is

$$\epsilon(\mathcal{N}, \mathcal{C}) = \min_{\mathcal{R}} \frac{1}{2} \left\| |\Psi_{\mathcal{C}}\rangle \langle \Psi_{\mathcal{C}}| - \overline{\mathcal{R}}(\overline{\mathcal{N}}(|\Psi_{\mathcal{C}}\rangle \langle \Psi_{\mathcal{C}}|)) \right\|_1,$$

where $\overline{\mathcal{R}} = \mathcal{I} \otimes \mathcal{R}$, $\overline{\mathcal{N}} = \mathcal{I} \otimes \mathcal{N}$, \mathcal{I} is an identity channel, the minimization is over all recovery maps \mathcal{R} , and $\|\cdot\|_1$

denotes the trace norm. For simplicity, when the code \mathcal{C} and noise model \mathcal{N} are implicit, we write $\epsilon = \epsilon(\mathcal{N}, \mathcal{C})$.

Let perturbations A_1, \dots, A_α to the Hamiltonian H occur with probabilities p_1, \dots, p_α , respectively. These perturbations model the coupling of qubits to spurious classical fields. Each perturbation A_j is a linear combination of operators that acts nontrivially on a single qubit, and induces a unitary evolution $U_j = g(H + A_j)$, where $g(x) = e^{-ix\tau}$. In what follows, we consider a random coherent noise channel \mathcal{N}_τ , which is parametrized by its noise strength $a = \max_j \|A_j\|/N$, and for any initial state ρ ,

$$\mathcal{N}_\tau(\rho) = \sum_{j=1}^{\alpha} p_j U_j \rho U_j^\dagger. \quad (8)$$

For any perturbation A_j , the Taylor series of the unitary $g(H + A_j)$ gives

$$g(H + A_j) = g(H) + \sum_{k=1}^{\infty} D_g^{[k]}(H, A_j) / k!, \quad (9)$$

where

$$D_g^{[k]}(H, A_j) = \frac{d^k}{d\xi^k} g(H + \xi A_j) |_{\xi=0} \quad (10)$$

are Fréchet derivatives of $g(H)$ in the matrix direction A_j [39,40]. Now the correctible component of $g(H + A_j)$ with respect to a code that corrects t errors comprises of Fréchet derivatives of order at most t , because these Fréchet derivatives are polynomials in A_j of order no more than t . Therefore, we study only the high order Fréchet derivatives. These Fréchet derivatives allow us to bound the storage error.

Lemma 2. Given a quantum code \mathcal{C} that corrects t errors, let $R_j = \sum_{k=t+1}^{\infty} D_g^{[k]}(H, A_j)$ and define $\|R\|_{\mathcal{C}} = \max_j \{\|R_j|\psi\rangle\| : |\psi\rangle \in \mathcal{C}\}$. Then $\epsilon \leq \|R\|_{\mathcal{C}} + \|R\|_{\mathcal{C}}^2$.

We prove Lemma 2 in the Supplemental Material [36]. From the integral representation of R_j [41], we exploit the fact that $g(H + A_j)$ is unitary for Hermitian H and A to get

$$\|R\|_{\mathcal{C}} \leq \max_j \frac{\max \{\|D_g^{[t+1]}(H, A_j)|\psi\rangle\| : |\psi\rangle \in \mathcal{C}\}}{(t+1)!}, \quad (11)$$

which depends only on a single Fréchet derivative instead of infinitely many. Given (11), one can clearly bound $\|R\|_{\mathcal{C}}$ in terms of just $\|H\|$ and aN . However, such a bound increases with increasing $\|H\|$, and exhibits a behavior contrary to the numerical evidence in Fig. 1. Increasing the number of long-range interactions increases both $\|H\|$ and dimensionality, but since increasing dimensionality should decrease the storage error, this suggests that the storage error should instead decrease with increasing $\|H\|$. We solve this conundrum by using Davis' representation [42] of Fréchet derivatives, which reveals the intricate dependence of Fréchet derivatives on the spectral decomposition $H = \sum_{j \geq 0} \lambda_j \Pi_j$. Here λ_j strictly increase with j and Π_j are eigenprojectors. Namely, we can write $D_g^{[k]}(H, A_j) / k!$ as

$$\sum_{n_0, \dots, n_k} g(\lambda_{n_0}, \dots, \lambda_{n_k}) (\Pi_{n_k} A_j) \cdots (\Pi_{n_1} A_j) \Pi_{n_0}. \quad (12)$$

Here $g(\lambda_{n_0}, \dots, \lambda_{n_k})$ are divided differences that arise naturally from the theory of Lagrange interpolation. To unravel (12), we leverage on the remarkable properties of divided

differences. First, divided differences are invariant under any permutation of their arguments. Hence, we can always arrange the arguments of a divided difference in nondecreasing order. Second, divided differences generalize scalar derivatives, because the divided difference of a vector with k identical arguments is proportional to the $(k-1)$ th derivative of the underlying function. For the exponential function, $|g(y_1, \dots, y_k)| = \tau^{k-1}/(k-1)!$ when $y_1 = \dots = y_k$. For instance, $|g(2, 2)| \leq \tau$. Third, a divided difference when not evaluated on identical arguments can be recursively defined; whenever y_i and y_j are distinct, $g(\mathbf{y}) = \{g(\mathbf{y}[\text{not } i]) - g(\mathbf{y}[\text{not } j])\}/(y_i - y_j)$, where $\mathbf{y}[\text{not } i]$ denotes a vector obtained from \mathbf{y} by deleting its i th component. From (11) and (12) we find that

$$\|R\|_{\mathcal{C}} \leq (aN)^{t+1} \left(h_{t+1} + \frac{\tau^{t+1}}{(t+1)!} \right). \quad (13)$$

Here $\tau^{t+1}/(t+1)!$ arises from the divided difference with all arguments equal to zero. The term h_{t+1} is the sum of all $|g(0, \lambda_{n_1}, \dots, \lambda_{n_{t+1}})|$ where $n_1 + \dots + n_{t+1} > 0$.

Evaluating a bound on h_{t+1} requires knowing the eigenvalues of H . Since finding the eigenvalues of H for arbitrary E is difficult [35], we study an exactly solvable HF where every pair of spins interacts equally with $J_{i,j} = J$. We call such a HF a mean-field HF, and its ground state energy is $\lambda_0 = 0$ and its higher energy eigenvalues are $\lambda_1 = JN$, $\lambda_2 = 2J(N-1)$, $\lambda_3 = 3J(N-2)$. In general, $\lambda_j = Jj(N+1-j)$ [35]. Now denote δ_j as the minimum energy needed to transition away from λ_j . For instance, $\delta_0 = \lambda_1 - \lambda_0 = JN$, $\delta_1 = \lambda_2 - \lambda_1 = J(N-2)$, and $\delta_2 = \lambda_3 - \lambda_2 = J(N-4)$. In general, $\delta_{\lfloor N/2 \rfloor} = 2 + (N - 2\lfloor N/2 \rfloor)$ and $\delta_j = J(N-2j)$ for all $j = 0, \dots, \lfloor N/2 \rfloor - 1$. Importantly, δ_j is nonincreasing in j and is maximal when $j = 0$. Exploiting the recursive structure of divided differences, one gets

$$|g(0, \lambda_{n_1}, \dots, \lambda_{n_r})| \leq 2^{t+1} \delta_0^{-1} (\delta_{n_1} \dots \delta_{n_{t+1}})^{-1}. \quad (14)$$

We provide the full details of obtaining this upper bound in the Supplemental Material [36]. When a divided difference has repeated arguments, we overestimate its contribution to h_{t+1} by severe overcounting. For this we first use (14) for divided differences even when there are r repeated arguments. Second, for divided arguments with r repeated entries, we count the contributions from leaves that terminate with all possible divided differences with repeated identical arguments.

If the contribution to the divided differences is dominated by leaves with no repeating indices, the total contribution of such leaves to h_{t+1} is at most $S^t 2^{t+1}/\delta_0$, where $S = \delta_0^{-1} + \dots + \delta_{t+1}^{-1}$. The contribution to h_{t+1} by leaves that terminate with r repeated arguments is $[\tau^{r-1}/(r-1)!] S^{t+1-r}/\delta_0$. From this we get $h_{t+1} \leq \theta$ where

$$\theta = \frac{S^{t+1}}{\delta_0} + \frac{n/2 + 1}{\delta_0} \sum_{r=2}^{t+1} \frac{\binom{t+1}{r-1} \tau^{r-1}}{(r-1)!} S^{t+2-r}. \quad (15)$$

From Lemma 2, (13) and (15), we get the following result.

Theorem 3. Let H be a mean-field HF with exchange constant J , and \mathcal{C} be any PI code that corrects t errors. Let \mathcal{N}_τ be the random coherent noise channel (8). Then $\epsilon(\mathcal{N}_\tau, \mathcal{C}) \leq \Theta + \Theta^2$, where $\Theta = (aN)^{t+1}[\theta + \tau^{t+1}/(t+1)!]$, and θ is given in (15).

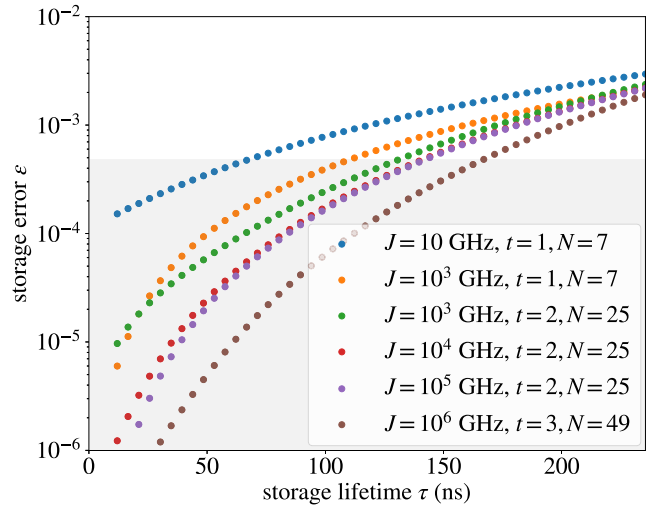


FIG. 2. When a mean-field HF stores an encoded qubit within an N -qubit PI that corrects t errors, we use Theorem 3 to obtain upper bounds for the corresponding storage error ϵ after a target storage lifetime of τ . The baseline lifetime and storage error for an unprotected qubit are 12 ns and 0.00048, respectively. The shaded region indicates where ϵ is smaller than the baseline.

Theorem 3 implies that the quantum memory is partially self-correcting, because the bound on ϵ contains a term $(N\tau)^t/t!$ which diverges for large t , since N is quadratic in t . Hence for fixed noise parameter a and exchange constant J , our scheme for a quantum memory has an optimal system size. Figure 2 illustrates only results for optimal system sizes.

Recently, a superconducting qubit was stored between 12 to 20 ns with a fidelity of 0.9995 [43]. Using our noise model, these experimental parameters can be recast into a baseline storage error of 5×10^{-4} with a memory lifetime of 12 ns and a noise strength of $a = 0.04$ MHz. Given this noise model, we use Theorem 3 to obtain upper bounds on the storage error ϵ of an encoded qubit within a PI code in a mean-field HF, and we depict these numerical results in Fig. 2. Here the number of qubits for $t = 1$ is seven [27], and when $t \geq 2$, $N = (2t + 1)^2$. From Fig. 2, if one uses a seven-qubit PI code with $J = 10^3$ GHz, the qubit's storage lifetime can be improved to over 100 ns. In addition, if one uses a 25 qubit PI code that corrects two errors [28], the qubit's storage lifetime can be enhanced to over 120 ns when $J = 10^4$ GHz. Similarly, if $J = 10^6$ GHz, the qubit's storage lifetime can be enhanced to over 150 ns using a 49 qubit PI code that corrects three errors. From this we can see how increasing J and the number of qubits in HFs can enhance the storage lifetime.

Discussions. Here we study quantum storage in a physically abundant physical system, the HF. Since our scheme is a physical model that is simple to realize, it will be easier to implement than those built upon many-body interactions. Because Pauli errors on PI codes exhibit a macroscopic energy barrier, we see evidence that a quantum memory based in a HF can become increasingly robust with increasing dimensionality of the HF. Moreover, we find that strengthening the coupling strengths can extend the storage lifetime of HF-based quantum memories when used in concert with PI codes. For this we analyze an infinite-dimensional HF, namely the

mean-field HF, and find numerically tractable upper bounds on the storage error. Our derivation of the bounds relies on a novel approach based on the connection between matrix perturbation theory and divided differences.

Since our analysis technique extends to any physical system with a completely understood spectral structure, it applies also to other code-inspired Hamiltonians, and lays the foundations for analyzing quantum memories using our new methodology. While PI codes can be prepared in superconducting charge qubits [44], it remains to integrate the initialization Hamiltonian with HFs. Furthermore, constructing explicit protocols for the decoding of PI codes can bring quantum memories in HFs closer to implementation.

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