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Noise amplification at spin-glass bottlenecks of quantum annealing: a solvable model

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To gain better insight into the complexity theory of quantum annealing, we propose and solve a class of spin systems which contain bottlenecks of the kind expected to dominate the runtime of quantum annealing as it tries to solve difficult optimization problems. We uncover a noise amplification effect at these bottlenecks, whereby tunneling rates caused by flux-qubit noise scale in proportion to the number of qubits \( N \) in the limit that \( N \to \infty \). By solving the incoherent annealing dynamics exactly, we find a wide range of regimes where the probability that a quantum annealer remains in the ground-state upon exiting the bottleneck is close to one-half. We corroborate our analysis with detailed simulations of the performance of the D-Wave 2X quantum annealer on our class of computational problems.

I. INTRODUCTION

Quantum Annealing (QA), a quantum heuristic for approximately solving NP-hard binary optimization problems, is already in commercial use [1–7] in machine learning and artificial intelligence applications. The algorithm works by mapping Quadratic Unconstrained Binary Optimization (QUBO) problems to the problem of solving for the ground state of a spin glass Hamiltonian. The time-complexity of QA, however, that is, how the required resources for running the algorithm scale with problem size, is still under investigation. The scaling behavior has been computed only for several optimization problems [8–11].

A key benchmarking problem in QA is the question of how an adiabatic quantum computer performs on spin glass bottlenecks. Those are time intervals during the annealing schedule where the gap shrinks exponentially with problem size, see Fig. 1. A cascade of hard bottlenecks was found [11] in the ordered phase of the annealing process. Over a decade later, another work [8] provided an exactly solvable spin glass system, where the scaling of the gap at these bottlenecks was obtained via analytical arguments, as opposed to the usual, numerical treatment. As a result, it is now understood that the time-complexity of QA for large problems is dominated by spin glass bottlenecks [1, 12–16]. There are two general features of these bottlenecks, (i) an exponentially small gap as the system size \( N \) grows, and (ii) a quantum tunneling event that flips \( O(N) \) spins.

These features can be embedded in a simple model, which can then be solved exactly (some properties are accurately accessible only in the asymptotic limit of \( N \to \infty \)). To the best of our knowledge, there have been no analytical studies of the effects of realistic (i.e. longitudinal) qubit noise at bottlenecks of QA in the presence of frustration, in the limit of large \( N \) [17, 18]. In this study, we ultimately find that, at an annealing bottleneck, the effects of frustration on multiqubit tunneling are washed-out in the large-\( N \) limit, leaving behind a large-\( N \) noise amplification effect. In particular, we find that tunneling rates in a wide class of frustrated spin chains diverge as \( N^2 \), where \( M \) is a suitably-defined bulk spontaneous magnetization. This gives analytical confirmation of prior work, which has found that the effective multiqubit noise spectral density at an annealing bottleneck grows as the multiqubit Hamming distance between the crossing states [19, 20].
The rest of the paper is organized as follows. We introduce and discuss basic properties of our exactly solvable model of bottlenecks in QA in Section II. Section III concentrates on the effects of ambient noise at the QA bottlenecks. We solve the Redfield equation to predict the behavior of the annealing processor at the bottleneck at finite temperature in Section IV. We conclude with the main results in Section V.

II. A MINIMAL MODEL OF A SPIN GLASS BOTTLENECK OF QA

We now search for the simplest class of Ising spin systems containing a spin glass bottleneck, where the minimum gap decreases exponentially as the number of qubits $N \to \infty$. Note that spin glass bottlenecks with frustration are impossible to realize with mean-field-like problem Hamiltonians (i.e. problem Hamiltonians with all-to-all interactions), thus motivating an investigation of the one-dimensional case. Defining Pauli matrices $\sigma^\alpha_j$, $\alpha = x, z$ acting on sites $j$ of a lattice, the one-dimensional transverse-field Ising spin glass has Hamiltonian

$$\hat{H}_0(t) \equiv \hat{H}_P + B(t) \sum_j \hat{\sigma}_j^x.$$  

(1)

The parameter $B(t)$ here represents a uniform transverse magnetic field, and the problem Hamiltonian $\hat{H}_P$ for the one-dimensional Ising spin glass is

$$\hat{H}_P \equiv -\sum_{j=1}^{N} J_j \sigma_j^z \sigma_{j+1}^z,$$  

(2)

where here, $J_j$ denotes a coupler connecting qubits $j$ and $j + 1$, and it is understood that the qubit indices $j$ are to be interpreted modulo $N$. The one-dimensional transverse-field Ising model can be mapped to free fermions [21], and is thus a good place to gain some useful physical intuition about these bottlenecks. The simplest one-dimensional annealing problem containing a spin glass bottleneck is the Frustrated Ring, the one-dimensional spin system depicted in Figure 2. The couplings $J_j$ for the Frustrated Ring are given as follows:

$$J_j = \begin{cases} J_L & j = n, n + 1 \\ -J_R & j = 2n + 1 \\ J & \text{otherwise} \end{cases},$$  

(3)

here, $0 < J_R < J_L < J$, and we make the total number of qubits $N \equiv 2n + 1$ odd, to make the problem more symmetric (and thus more amenable to an exact calculation). The Frustrated Ring is a minimal model of a spin glass bottleneck, because, in one dimension, one must modify at least three couplers in an otherwise uniform graph to achieve a spin glass bottleneck; the Frustrated Ring saturates this lower bound.

The Frustrated Ring is frustrated, and therefore has a forced excitation in its groundstate. At zero transverse field $B(t) \equiv 0$, there are two generic positions where this excitation likes to reside: (i) at the antiferromagnetic coupler $J_R$ (forming the frustrated groundstate $|\Psi_R\rangle$), and (ii) at either of the two weak ferromagnetic couplers $J_L$ (forming degenerate first-excited states $|\Psi_L\rangle$), see Fig. 4. At a special value of the transverse-field $B_b$ within the ordered phase of the anneal, the $|\Psi_R\rangle$ states and a pair of $|\Psi_L\rangle$ states form an avoided crossing with a gap that scales as

$$\Delta_{\min} \propto \left( \frac{J(J_L^2 - J_R^2)}{J_R(J^2 - J_L^2)} \right)^N,$$  

(4)

which is exponentially small in $N$, as desired (see Fig. 5). Representative performance of the D-Wave 2X quantum annealer in the bottleneck regime, $JJ_R > J_L^2$, is depicted in Fig. 3. Since $|\Psi_L\rangle$ and $|\Psi_R\rangle$ differ by flipping half
FIG. 4. Lowest energy states of the Frustrated Ring, at the end of the annealing process. The green loop denotes a bond that gives a positive contribution to the energy of the Hamiltonian in Eq. (2). In (a), $|Ψ_R⟩$ is formed by violating the anti-ferromagnetic bond $J_R$. $|Ψ_L⟩$ is obtained by violating either of the two weak ferromagnetic bonds $J_L$, as shown in (b).

of the spins, the avoided crossing of these states has two key features in common with spin glass bottlenecks: (i) an exponentially small gap, and (ii) a quantum tunneling event that flips $O(N)$ spins.

A. Annealing schedule

We now analyze the annealing schedule of the Frustrated Ring, and demonstrate that it runs in exponential time (assuming completely coherent processor dynamics). The first step in the analysis consists of reinterpreting the system of interacting spins as a system of non-interacting fermions. In this new description, the global spin-flip symmetry becomes a symmetry $(-1)^{N_F}$, which counts the number of fermionic excitations modulo 2:

$$(-1)^{N_F} \equiv \prod_{j=1}^{N} \hat{\sigma}^z_j, \quad \left[ \hat{H}_0(t), (-1)^{N_F} \right] = 0 . \quad (5)$$

The appropriate fermionic operators $\hat{\gamma}_1, \hat{\gamma}_2, \ldots, \hat{\gamma}_N$ are Majorana fermions, i.e. they generate the Clifford algebra

$$\hat{\gamma}_i \hat{\gamma}_j + \hat{\gamma}_j \hat{\gamma}_i \equiv 2 \delta_{ij}. \quad (6)$$

These Majorana fermions are written in terms of the original spin operators via the Jordan-Wigner transformation

$$\hat{\sigma}^+_j = -i \hat{\gamma}_{2j-1} \hat{\gamma}_{2j},$$
$$\hat{\sigma}^-_j = (-i)^{j-1} \hat{\gamma}_1 \cdots \hat{\gamma}_{2j-1}. \quad (7)$$

In terms of the fermionic operators, the theory decouples into two free theories:

$$\hat{H}_0^\pm(t) = -i \sum_j J_j^\pm \hat{\gamma}_{2j} \hat{\gamma}_{2j+1} + i B(t) \sum_j \hat{\gamma}_{2j-1} \hat{\gamma}_{2j}, \quad (8)$$

where $J_j^\pm \equiv J_j$ for $j \neq N$; $J_N^\pm \equiv \pm J_N$, and $\hat{H}_0^\pm(t)$ denotes the annealing Hamiltonian $\hat{H}_0(t)$ restricted to the sector with an even and odd number of Majorana fermions, respectively.

FIG. 5. (a) The low-energy spectrum of an $N=7$ Frustrated Ring computed for representative values of couplings $J_j$ in Eq. (1) and Eq. (3). Note the avoided crossing. (b) Due to $Z_2$ Ising symmetry, the annealing bottleneck consists of two crossings, one occurring in each symmetry sector. In the Frustrated Ring, the odd-parity crossing is gapless. However, only the even-parity crossing is seen by the coherent annealing dynamics (see [22] for details), and thus the even-parity gap (see Eq. (28)) determines the QA time-complexity of this problem.

1. Zero transverse field

The non-interacting theories may be diagonalized via the Bogoliubov-de-Gennes formalism for all $B(t)$. To diagonalize the theories at the end of the anneal (when the transverse field is completely turned off), we form the complex fermions

$$c_{j,\pm} := \gamma_{2j} + i \gamma_{2j+1}, \quad (9)$$
$$\tilde{c}_{j,\pm} := \gamma_{2j} - i \gamma_{2j+1}. \quad (10)$$

In terms of the complex fermions, both free theories take the following diagonalized form at $B = 0$:

$$\hat{H}_0^\pm(t) \underset{B \to 0}{=} \sum_j J_j^\pm c_{j,\pm}^\dagger c_{j,\pm}. \quad (11)$$
\(J^+_j > 0\) for all \(j\), and so, in the odd-fermion sector, all fermionic quasiparticles thus have positive energy at \(B = 0\). The groundstate in this sector is

\[
|\Psi_R^\pm\rangle := c_{N, \pm}^\dagger |\Omega_\pm\rangle, \tag{12}
\]

Similarly, the two degenerate first-excited states in this sector are produced by exciting the bonds \(J_L\) at the left-most portion of the graph:

\[
|\Psi_L^-\rangle := c_{N, -}^\dagger |\Omega_-\rangle, \tag{13}
\]

\[
|\Psi_L^+\rangle := c_{N+1, +}^\dagger |\Omega_+\rangle, \tag{14}
\]

where we have factored \(N = 2N + 1\). The energies of the above states are equal to the odd-parity vacuum energy, plus the energy \(J_{R,L}\) of the bond violated in each case, as expected from examination of the diagonal form of the Hamiltonian in Eq. (11). In the even-fermion sector, the analysis is a bit more complicated, as the presence of an antiferromagnetic bond means that when that bond is violated, the spin chain loses energy. In the fermionic picture, this is reflected by the fact that the fermionic quasiparticle \(c_{N, +}^\dagger\) has negative energy. Accordingly, the groundstate in the even-fermion sector is just vacuum:

\[
|\Psi_R^\pm\rangle := |\Omega_\pm\rangle. \tag{15}
\]

It is impossible to have a single fermionic excitation in this sector. The next-lowest energy eigenstates thus have two excitations each:

\[
|\Psi_L^\pm\rangle := c_{N, \pm}^\dagger c_{N+1, \mp}^\dagger |\Omega_\pm\rangle, \tag{16}
\]

\[
|\Psi_L^\pm\rangle := c_{N, \mp}^\dagger c_{N+1, \pm}^\dagger |\Omega_\pm\rangle, \tag{17}
\]

where we have again factored \(N = 2N + 1\). The energies of the above states are equal to the even-parity vacuum energy, plus the difference \(J_L - J_R\). Again, here, the quasiparticle \(c_{N, +}^\dagger\) has negative energy, as it produces a domain wall at the antiferromagnetic bond \(J_L^+ < 0\).

2. Nonzero transverse field

We now make general comments about the above low-lying spin chain eigenstates in the presence of a nonzero transverse field \(B(t) > 0\). In general, the diagonal form Eq. (11) is now dependent on the transverse-field value \(B(t)\). The odd-parity Hamiltonian \(\hat{H}_L^\pm(t)\) is simplest to analyze, as all of its coupling constants are positive in the limit \(B \to 0\), and thus all fermionic excitations (for sufficiently small transverse field) still have positive energy. For \(B > 0\), the energies \(\epsilon > 0\) of these excitations will in general have some non-trivial dependence on the transverse field, leading to an energy dispersion curve \(\epsilon(B)\).

For instance, for the right-most localized bound state,

\[
\epsilon_{N, \pm}(B) \sim \mp J_R, \quad \text{etc.} \tag{18}
\]

When any one of these dispersion curves crosses the horizontal line \(\epsilon(B) \equiv 0\), the model becomes gapless. We call the smallest transverse field at which this happens \(B_c\), i.e. the critical value of the transverse-field.

We then define \(B(t) < B_c\) to be the ordered phase of the anneal, and the remaining regime \(B(t) > B_c\) to be the paramagnetic phase of the anneal. In the ordered phase, the Bogoliubov de-Gennes (BdG) equations centered on the bond \(J_N \equiv J_R\) may be solved exactly, producing a fermion bound state \(\tilde{c}_R\) exponentially localized at the right-end of the graph, creating the odd-parity ground state.

\[
|\Psi_R^\pm\rangle := c_{R, \mp}^\dagger |\Omega_\pm\rangle. \tag{19}
\]

Similarly, in the limit \(N \to \infty\), the BdG equations centered on the bond \(J_L\) may also be solved exactly (see Appendix A), yielding two low-energy bound states \(\tilde{c}_L, \tilde{c}_L^\prime\) exponentially localized at the left-end of the graph, so that the next-lowest energy states with odd-parity are

\[
|\Psi_L^\pm\rangle := c_{L, \pm}^\dagger |\Omega_\pm\rangle, \tag{20}
\]

\[
|\Psi_L^\pm\rangle := c_{L, \mp}^\dagger |\Omega_\pm\rangle. \tag{21}
\]

The corresponding energies of the states Eqs. (19-21) are \(\tilde{\epsilon}_{vac} + \tilde{\epsilon}_R\), and \(\tilde{\epsilon}_{vac} + \tilde{\epsilon}_L\), and \(\tilde{\epsilon}_{vac} + \tilde{\epsilon}_L^\prime\) respectively, where \(\tilde{\epsilon}_{vac}\) is the energy of the odd-parity vacuum state.
\(|\Omega^-\rangle\). From the definition, it is then clear that, whenever the energy dispersion curves \(\epsilon_R(B)\) and \(\epsilon_L(B)\) cross, the corresponding spin-chain eigenstates Eqs. (19-20) cross in energy.

In the even-parity sector, a similar picture emerges, although the energy considerations are complicated by the fact that some defects have negative energy in the ordered phase, as the even-parity Hamiltonian \(\hat{H}_R^+\) contains a coupling constant \(J_N^+ = -J_R\) which is negative. However, after some book-keeping, one can deduce the three-lowest energy states in the even-fermion sector, which we list below:

\[
|\Psi_R^+\rangle := |\Omega_+\rangle, \quad (22)
\]

\[
|\Psi_L^+\rangle := c_{R,+}^\dagger c_{L,+}^\dagger |\Omega_+\rangle, \quad (23)
\]

\[
|\Psi_L^-\rangle := c_{R,+}^\dagger c_{L,-}^\dagger |\Omega_+\rangle. \quad (24)
\]

As in the odd-parity case, the corresponding energies of the states Eqs. (22-24) are \(\epsilon_R^+\), \(\epsilon_{\text{vac}}^+ + \epsilon_R^+\), \(\epsilon_L^+\), and \(\epsilon_{\text{vac}}^+ + \epsilon_R^+ + \epsilon_L^+\) respectively, where \(\epsilon_{\text{vac}}^+\) is the energy of the even-parity vacuum state \(|\Omega^+\rangle\). Again, this means that when the energy dispersion curves \(|\epsilon_R^+(B)\rangle\) and \(\epsilon_L^+(B)\) cross, the corresponding spin-chain eigenstates Eqs. (19-20) cross in energy. Here, we have taken the absolute value of the energy of the right-localized defect \(c_{R,+}^\dagger\), as this defect has negative energy throughout the ordered phase \(B(t) < B_c\) of the anneal.

The origin of the annealing bottleneck becomes clear once one realizes that the level repulsion between \(\epsilon_L, \epsilon_L'\) deflects the energy dispersion curve \(\epsilon_L(B)\) of the \(c_L\)-defect downwards (see Figure 6), so that, for \(JJ_R > J_L^2\), it crosses with the energy of the \(c_R\) defect at a tunable \(B = B_b\) value of the transverse-field within the spin-glass phase of the anneal:

\[
B_b \equiv \frac{1}{J_R} \left( J^2 - J_L^2 \right) \frac{(J_L^2 - J_R^2)}{J^2 + J_L^2 - 2J_L^2} < B_c. \quad (25)
\]

A rigorous derivation of the identity in Eq. (25) is carried out in Appendix A. The level crossing produces the situation in the odd-fermion sector depicted in Figure 7, where a forced excitation (forced by parity constraints) must quantum-mechanically tunnel from left to right in the graph. A similar situation occurs in the even-fermion sector, leading to a second level crossing at \(B = B_b\). So, in total, at \(B = B_b\), two pairs of energy levels cross.

3. Scaling of the gap

The previous analysis was only exact in the limit \(\mathcal{N} \to \infty\). At finite \(\mathcal{N}\), in each parity sector, there will be hybridization between boundstates localized at opposite ends of the graph (due to finite-size effects), and so, directly at the minimum gap region, the true fermionic eigenmodes will sweep rapidly through a mixture of left- and right-modes, producing an avoided (Landau-Zener) crossing. For example, in the even-fermion sector, one can define a Landau-Zener approach angle

\[
\tan 2\theta_{LZ}(B) \equiv \frac{\Delta(B)}{\epsilon_R^+(B) - \epsilon_L^+(B)},
\]

where \(\Delta(B)\) is defined as the (exponentially-small) overlap between the left- and right-localized even-parity boundstates. In terms of \(\theta_{LZ}\), the two lowest-energy excitations become:

\[
\epsilon_{0,+}^+ = \epsilon_{R,+}^+ \sin \theta_{LZ} + \epsilon_{L,+}^+ \cos \theta_{LZ}, \quad (26)
\]

\[
\epsilon_{1,+}^+ = \epsilon_{L,+}^+ \sin \theta_{LZ} - \epsilon_{R,+}^+ \cos \theta_{LZ}. \quad (27)
\]

In contrast, the odd-parity bound states do not hybridize, as they have different parities under spatial reflection \(j \to \mathcal{N} - (j - 1)\) (see Appendix A), and the crossing there is exact, for all \(\mathcal{N} \leq \infty\). Therefore, at \(B = B_b\), in the large-\(\mathcal{N}\) limit, there is a pair of crossings, with each crossing occurring in a distinct eigenspace of \((-1)^{2\mathcal{N}}\), i.e. each crossing is labelled by a distinct \(\mathbb{Z}_2\) quantum number, see bottom portion of Fig. 5. The scaling of the hybridization of the even-parity boundstates at the bottleneck location gives the inverse QA runtime

\[
\Delta(B_b) \sim N^\infty \mathcal{O}\left( \frac{(J_L^2 - J_R^2)}{J_R^2 + J_L^2 - 2J_L^2} \right)^N, \quad (28)
\]

which is exponentially small in \(\mathcal{N}\), as desired. Note that the spectral data in the odd-parity sector is irrelevant in analyzing the time-complexity of the closed-system annealing dynamics, as the odd-parity sector is never visited during the coherent evolution. This is because the initial state of the QA protocol always has even spin-flip parity (see, e.g. [22]), and spin-flip parity is conserved throughout the annealing schedule.

III. ANALYSIS OF FLUX QUBIT NOISE AT THE ANNEALING BOTTLENECK

In this section we turn our attention to the effects of ambient flux qubit noise at a frustrated spin glass bottleneck. In the D-Wave 2X quantum annealer, fluctuations in onsite qubit flux bias form the dominant source of noise [19, 20]. This flux bias noise is accurately modelled by the following system-bath Hamiltonian:

\[
\hat{H}_{\text{D-Wave}} = \hat{H}_0 + \sum_{j=1}^{\mathcal{N}} \hat{Q}_j \hat{\sigma}^z_j + \hat{H}_B, \quad (29)
\]

\[
\hat{Q}_j = \sum_u \lambda_u (\hat{b}_{j,u} + \hat{b}_{j,u}^\dagger), \quad (30)
\]

\[
\hat{H}_B = \sum_{j,u} \hbar \omega_u (\hat{b}_{j,u}^\dagger \hat{b}_{j,u} + 1/2), \quad (31)
\]
where the bath operators \( \{ \hat{b}_{j,u} \} \) for each flux qubit satisfy standard bosonic commutation relations \( \{ \hat{b}_{j,u}, \hat{b}_{j,u}^\dagger \} = \delta_{u,u'} \), and \( \omega_u \) is the frequency of mode \( u \) and \( \lambda_u \) sets the interaction strength between that mode and its corresponding qubit. For weak noise strength, the effects of the flux bias fluctuations on the system dynamics are uniquely characterized by the noise spectral density of these fluctuations:

\[
S(\omega) = \int_0^\infty dt e^{i\omega t} \langle e^{iH_B t/\hbar} \hat{Q}_j e^{-iH_B t/\hbar} \hat{Q}_j \rangle \quad (32)
\]

which we assume to be identical for each qubit. For simplicity, in this study we assume the spectral density of the noise to be Ohmic:

\[
S(\omega) = \frac{\hbar^2}{1 - \exp(-\hbar\omega/k_B T)} \quad (33)
\]

at temperature \( T \). Here, \( \eta \) characterizes the strength of the fluctuations in the flux bias, and is typically measured in macroscopic resonant tunneling experiments on the individual qubits in the processing unit. For the D-Wave 2X annealer at NASA QuAIL, \( \eta \) was measured to be \( \sim 0.24 \) in the regime \( B \equiv 0 \) [19]. We use this value of \( \eta \) for our numerical simulations of the D-Wave 2X at Los Alamos National Laboratory.

We treat the open-system dynamics specified by Eqs. (29-31) in the Bloch-Redfield approximation, which, at the qualitative level, closely predicts the probability \( P \) that the D-Wave 2X machine returns a global minimum of the Frustrated Ring MAXCUT problem:

\[
P \equiv \langle 00\cdots 0 | \hat{\rho}_f | 00\cdots 0 \rangle + \langle 11\cdots 1 | \hat{\rho}_f | 11\cdots 1 \rangle \quad (34)
\]

where here, \( \hat{\rho}_f \) is the collective density matrix of the qubits in the annealer at the end of the anneal, and \( \{ 000\cdots 0, 111\cdots 1 \} \) is the complete set of bit strings which solve the Frustrated Ring MAXCUT problem. The actual success probability observed in the D-Wave 2X quantum annealer, compared with the simulated success probability (according to the Bloch-Redfield simulation of the quantum processor) is plotted in Fig. 8.

Assuming the annealing processor is completely incoherent, i.e. the off-diagonal matrix elements of \( \hat{\rho} \) vanish in the energy eigenbasis, then the Redfield master equation degenerates into a kinetic equation involving tunneling rates between instantaneous eigenstates. These are given by Fermi’s golden rule as (c.f. [20]):

\[
\Gamma_{i \rightarrow f} = \frac{1}{2\hbar^2} S(\omega_{i \rightarrow f}) \cdot O_{i \rightarrow f} \quad (35)
\]

\[
O_{i \rightarrow f} = \sum_{j=1}^N \langle \langle \Psi_i | \hat{\sigma}_j^z | \Psi_f \rangle \rangle^2 \quad (36)
\]

where here, \( \omega_{i \rightarrow f} \) is the gap frequency \( (E_i - E_f)/\hbar \) between the initial \( |\Psi_i\rangle \) and final eigenstate \( |\Psi_f\rangle \). Directly at the bottleneck or in our frustrated spin system, we will find this incoherent evolution to be analytically solvable in the large-\( N \) limit. Symmetry simplifies the problem: since \( \hat{\sigma}_z \) flips the \( Z \)-parity number corresponding to spin-flip parity, the relevant form-factors that need to be calculated are (assuming the temperature is sufficiently low so that we can assume that the four-lowest levels are populated during the evolution):

\[
O_{0+ \rightarrow L^-} = \sum_j \langle |\Psi_0^+ | \hat{\sigma}_j^z | \Psi_L^- \rangle^2 \quad (37)
\]

\[
O_{0+ \rightarrow R^-} = \sum_j \langle |\Psi_0^+ | \hat{\sigma}_j^z | \Psi_R^- \rangle^2 \quad (38)
\]

\[
O_{1+ \rightarrow L^-} = \sum_j \langle |\Psi_1^+ | \hat{\sigma}_j^z | \Psi_L^- \rangle^2 \quad (39)
\]

\[
O_{1+ \rightarrow R^-} = \sum_j \langle |\Psi_1^+ | \hat{\sigma}_j^z | \Psi_R^- \rangle^2 \quad (40)
\]

All other matrix elements vanish by symmetry, because \( \hat{\sigma}_z \) mixes fermionic \( Z \)-parity symmetry. To calculate the transition rates, we expand them so that they are written completely in terms of the basis \( \{ |\Psi_L^-\rangle, |\Psi_R^-\rangle, |\Psi_L^+\rangle, |\Psi_R^+\rangle \} \) of crossing states:

\[
O_{0+ \rightarrow L,R^-} = \sum_j \sin \theta_{LZ} \langle |\Psi_0^+ | \hat{\sigma}_j^z | \Psi_{L,R}^- \rangle^2 + \cos \theta_{LZ} \langle |\Psi_0^+ | \hat{\sigma}_j^z | \Psi_{L,R}^- \rangle^2 \quad (41)
\]
\[ O_{1 \rightarrow L,R} = \sum_j \left[ \cos \theta_{LZ} \langle \Psi_R^\dagger \sigma_j^z | \Psi_{L,R} \rangle - \sin \theta_{LZ} \langle \Psi_R^\dagger \sigma_j^x | \Psi_{L,R} \rangle \right]^2. \] (42)

Some comments are in order. In general, tunneling due to non-fermionizable (i.e. \( \hat{\sigma}^z \)) noise in a one-dimensional quantum spin glass is analytically intractable due to the fact that the interacting portion of the Hamiltonian in Eq. (29) maps to a Jordan-Wigner string

\[ \sum_j \sigma_j^z \hat{Q}_j = \sum_j (-i)^j \gamma_1 \cdots \gamma_{2j-1} \sum_u \lambda_u (b_{j,u} + b_{j,u}^\dagger). \] (43)

Therefore, solving the full D-Wave dynamics even at the perturbative, Markovian level is widely considered to be analytically intractable [17, 18]. For example, one of the tunneling matrix elements above involves an inner-product of the form

\[ \langle \Psi_R^\dagger | \hat{\sigma}_j^x | \Psi_R \rangle = \Omega_+ |(-i)^j \gamma_1 \cdots \gamma_{2j-1} c_{R,u}^\dagger \Omega_- \rangle. \] (44)

The Jordan-Wigner string in the above equation means that the matrix element evaluates to a determinant of the matrix element 

\[ \langle \Psi_R^\dagger | \hat{\sigma}_j^x | \Psi_R \rangle = \langle \Omega_+ |(-i)^j \gamma_1 \cdots \gamma_{2j-1} c_{R,u}^\dagger \Omega_- \rangle. \]

The non-perturbative level by performing a field-theoretic calculation (see Appendix B), leading to an analytical understanding of flux qubit noise at a spin glass annealing bottleneck.

A. Non-perturbative large-\( \mathcal{N} \) calculation of tunneling rates

We begin by computing the off-diagonal, i.e. \( \langle \Psi_L \rangle \rightarrow \langle \Psi_R \rangle \), matrix elements. These vanish in the large-\( \mathcal{N} \)-limit as a consequence of Lieb-Robinson bounds [23]. For example, we can factor

\[ \langle \Psi_L^\dagger | \hat{\sigma}_j^z | \Psi_R \rangle = \langle \Psi_L^\dagger(0) | \hat{\sigma}_j^z(B) | \Psi_R(0) \rangle, \] (45)

where \( \hat{\sigma}_j^z(B) = \hat{U}^\dagger(B) \hat{\sigma}_j^z(B) \hat{U}(B) \), and the unitary \( \hat{U}(B) \) has the general form

\[ \hat{U}(B) = T e^{-i \int_{B_0}^B dB' \tilde{H}(B')} \] (46)

with \( \tilde{H} \) defined in [23]. Now, suppose \( B < B_0 \) is fixed. Since \( \tilde{H} \) has a spectral gap which is at least \( \mathcal{O}(1) \) for all \( B' < B \), \( \hat{U}(B) \) is a constant-depth unitary circuit [24]. That is, \( \hat{\sigma}_j^z(B) \), up to exponentially small corrections constant in \( \mathcal{N} \), is supported on a region of constant size. However, \( \langle \Psi_L^\dagger \rangle \) and \( \langle \Psi_R^\dagger \rangle \) at \( B = 0 \) are separated by \( \sim \mathcal{N}/2 \) spin flips. Therefore, for any fixed \( B < B_0 \), the matrix elements mixing \( \langle \Psi_L \rangle \) with \( \langle \Psi_R \rangle \) are exponentially small in \( \mathcal{N} \), i.e.

\[ \langle \Psi_L^\dagger | \hat{\sigma}_j^z(B) | \Psi_R \rangle \sim O(e^{-c \mathcal{N}/2}), \] (47)

with \( c \) a constant. Within a sufficiently small neighborhood of the crossing, we can assume the diabatic crossing states \( \langle \Psi_L^\dagger \rangle \) and \( \langle \Psi_R^\dagger \rangle \) to be approximately independent of \( B \) (with the \( B \)-dependence of the true eigenstates due to mixing within the subspace spanned by these crossing states). Therefore, for \( B > B_0 \), sufficiently near the crossing, the off-diagonal matrix elements are also exponentially small. Via the exact same reasoning, we also have the asymptotic behavior

\[ \langle \Psi_R^\dagger | \hat{\sigma}_j^z(B) | \Psi_L \rangle \sim O(e^{-c' \mathcal{N}/2}), \] (48)

for the other off-diagonal matrix element with \( c' \) another constant. Again, this follows from the fact that at \( B = 0 \) the state \( \langle \Psi_R^\dagger \rangle \) is separated from the state \( \langle \Psi_L^\dagger \rangle \) by \( \sim \mathcal{N}/2 \) spin flips.

In summary, the off-diagonal contributions to multi-qubit tunneling asymptotically vanish near the crossing in the large-\( \mathcal{N} \) limit. Therefore, provided that we are sufficiently close to the crossing point so that neither \( \theta_{LZ} \) nor cot \( \theta_{LZ} \) are exponentially small in \( \mathcal{N} \), we can ignore these off-diagonal contributions. In this limit, the large-
$\mathcal{N}$ tunneling form factors have the asymptotics
\begin{align}
O_{0^+ \to R^-} & \sim B_{0 \to B_0} \sum_j \sin^2 \theta_{LZ} |\langle \Psi_R^+ | \sigma_j^z | \Psi_R^- \rangle|^2 , \quad (49) \\
O_{0^+ \to L^-} & \sim B_{0 \to B_0} \sum_j \cos^2 \theta_{LZ} |\langle \Psi_L^+ | \sigma_j^z | \Psi_L^- \rangle|^2 . \quad (50)
\end{align}

with analogous expressions for $O_{1^+ \to R^-, L^-}$. To calculate the simplified form factors given by Eqs. (49-50), we begin with the following generic observation: for each time-dependent annealing Hamiltonian $H_0(t)$ specified in Eq. (1), let $H_0(t)$ denote the annealing Hamiltonian obtained by flipping the sign of $J_{\mathcal{N}}$. Upon taking the Jordan-Wigner transformation (c.f. Eq. (8)), we then have the following relations:

\[ H_0^\dagger(t) = \bar{H}_0^\tau(t) . \quad (51) \]

These relations are completely general and hold for any quantum spin glass on a 2-regular graph. Crucially, if $H_0$ is frustrated, $\bar{H}_0$ lacks frustration, and thus perhaps easier to characterize. For the Frustrated Ring, we can use the above relation to obtain a complete solution to the incoherent tunneling rates (c.f. Eqs. (37-40)) in the large-$\mathcal{N}$ limit.

Indeed, in Appendix B, we derive the following crucial identities relating low-lying eigenstates of the Frustrated Ring benchmark with those of its ferromagnetic counterpart $\bar{H}$:

\[ \langle \Psi_{R,L}^+ | \sigma_j^z | \Psi_{R,L}^- \rangle = -\langle \bar{\Psi}_{R,L}^+ | \sigma_j^z | \bar{\Psi}_{R,L}^- \rangle + O(e^{-\kappa_R |j-j_R|}) , \]

(52)

here, $j_R$ is the position of the antiferromagnetic coupler $J_R$, and $\kappa_R$ is the wavenumber of the boundstate $\psi_{R,L}^\dagger$. Note that these errors are localized at the position of the $J_R$-coupler, and thus do not grow if we sum over all qubits in the graph.

To summarize, by the replacements $|\Psi_{L,R}^\pm\rangle \to |\bar{\Psi}_{L,R}^\pm\rangle$, we can relate our frustrated tunneling form factors to those in an unfrustrated spin system $\bar{H}$ (c.f. Eq. 51), at the cost of inducing an error which does not grow with the total number of qubits $\mathcal{N}$. We thus have

\begin{align}
O_{0^+ \to R^-} & \sim \sum_j \sin^2 \theta_{LZ} |\langle \Psi_R^+ | \sigma_j^z | \Psi_R^- \rangle|^2 , \quad (53) \\
O_{0^+ \to L^-} & \sim \sum_j \cos^2 \theta_{LZ} |\langle \Psi_L^+ | \sigma_j^z | \Psi_L^- \rangle|^2 . \quad (54)
\end{align}

with analogous expressions for $O_{1^+ \to R^-, L^-}$. Note that the above asymptotics are not valid unless the expressions in Eqs. (53-54) are asymptotically greater than $O(1)$; we will find that this is the case (c.f. Eqs. (58-59)), so that our calculation is self-consistent. The frustration-free version ($\bar{H}_0$) of our problem is much easier to solve: in particular, any frustration-free spin system is gauge-equivalent to a ferromagnet via a local $\mathbb{Z}_2$ gauge-transformation of the form

\[ \hat{U}(g) \equiv \prod_j (\sigma_j^x)^{g_j} , \]

(55)

with $g_j \in \{0,1\}$. In fact, we find that the matrix elements in Eqs. (53-54) are related to the spontaneous magnetization of the ferromagnetic spin chain $\bar{H}$. Indeed, at $B = 0$, we have

\[ \langle \bar{\Psi}_R^+ | \sigma_j^z | \bar{\Psi}_R^- \rangle = 1 , \quad (56) \\
\langle \bar{\Psi}_L^+ | \sigma_j^z | \bar{\Psi}_L^- \rangle \sim \text{sgn}(j) . \quad (57)
\]

Therefore, at zero transverse field, the above matrix elements have the physical meaning of being the local magnetization of each state (in Eqs. (56-57), we have implicitly re-indexed the qubits from $j = -n, \ldots, 0, \ldots, n$, where $\mathcal{N} = 2n + 1$). Exact analytical expressions for the matrix elements in Eqs. (56-57) in the more general case $B > 0$ can be obtained by performing a field-theoretic calculation in the corresponding two-dimensional classical Ising model (see Appendix B). There, we find that the absolute values of the magnetizations in Eqs. (56-57) approach the bulk value $M \equiv (1 - (B/J)^2)^{1/8}$. Note that $M$ is equal to the bulk spontaneous magnetization of a quantum Ising chain with uniform ferromagnetic coupling $J$ [25]. A transfer matrix argument in Appendix B is used to show that this convergence is exponentially fast in the distance from the $J_L, J_R$ defects. We confirm this prediction with exact diagonalization for $\mathcal{N}$ up to 23 sites, see Fig. 9.

As a result, the multiqubit tunneling rates at the spin glass bottleneck have the following large-$\mathcal{N}$ asymptotic
form:

\[ \Gamma_{0+ \rightarrow L^-} \sim N \to \infty \frac{N M^2}{2 \hbar^2} \cos^2 \theta_{LZ}, \quad (58) \]

\[ \Gamma_{0+ \rightarrow R^-} \sim N \to \infty \frac{N M^2}{2 \hbar^2} \sin^2 \theta_{LZ}, \quad (59) \]

which is \( \mathcal{O}(N) \), with similar expressions for \( \Gamma_{1+ \rightarrow L^-, R^-} \). Again, \( M \equiv (1 - (B_0/J)^2)^{1/8} \), the bulk value of the spontaneous magnetization of the one-dimensional transverse-field Ising model, sets the coefficient of scaling. The above expressions are in excellent agreement with exact-diagonalization for \( N \) up to 23 sites, as shown in Fig. 10.

The physical implications of Eqs. (58-59) above are clear: in the Landau-Zener formalism, at \( t = -\infty \), \((\sin \theta, \cos \theta) = (0, 1)\), and then, at \( t = \infty \), we have that \((\sin \theta, \cos \theta) = (1, 0)\). Therefore, at \( t = \pm \infty \), the transition rate is predominantly the exponentially-small \(|\Psi_L \rangle \rightarrow |\Psi_R \rangle \) cross-terms (which were neglected in deriving (58-59)), and the approximation breaks down. However, at the bottleneck \( t \sim 0 \), the transition rate quickly reaches a peak which is asymptotic to \( N M^2 / 2 \), times the noise-spectral density evaluated at the minimum gap frequency.

**IV. LARGE-N LIMIT OF QUANTUM ANNEALING THROUGH THE BOTTLENECK AT FINITE-TEMPERATURE**

The simple analytical formulae (58-59) demonstrate the existence of a linear \( \mathcal{O}(N) \) tunneling peak near a quantum annealing bottleneck, and establish the relation of the scaling coefficient to a suitably defined bulk spontaneous magnetization, \( M \), in our benchmark spin system. Using these asymptotics, we can rigorously analyze the effects of flux-bias noise on a quantum annealing chip at a spin glass bottleneck, in the limit that the number of qubits tends to infinity.

Assuming that the annealing processor is fully incoherent, the density matrix \( \rho \) of the system is diagonal in the eigenbasis \( \{|\Psi_0^+, |\Psi_1^+, |\Psi_R^-, |\Psi_L^-\} \) of the coherent portion \( \hat{H}_0 \) of the quantum annealing Hamiltonian (29):

\[
\dot{\rho} = P_0^+ |\Psi_0^+\rangle \langle \Psi_0^+ | + P_1^+ |\Psi_1^+\rangle \langle \Psi_1^+ | + P_R^- |\Psi_R^-\rangle \langle \Psi_R^- | + P_L^- |\Psi_L^-\rangle \langle \Psi_L^- |. \quad (60)
\]

Under those assumptions, the density matrix satisfies incoherent time-evolution in terms of rates of the form \( \Gamma_{i \rightarrow j} \)

\[
\partial_t P_{0}^{+} = \Gamma_{L^{-}\rightarrow 0^{+}, L^{-}} + \Gamma_{R^{-}\rightarrow 0^{+}, R^{-}} - \sum_{L^{-}, R^{-}} \Gamma_{0^{+}, L^{-}\rightarrow R^{-}, L^{-}, 0^{+}} P_{0}^{+}, \quad (61)
\]

constituting a kinetic equation of Pauli type. Via Eqs. (58-59), the multiqubit tunneling rates in the above equation can then be computed exactly in the large-\( N \) limit, within a sufficient radius of the crossing point such that the approximation (49-50) is valid. It is thus convenient to define a tunneling region \( t_i < t < t_f \) within which this assumption holds. In this case, the instantaneous gap is much smaller than the temperature, and the Ohmic noise spectral density \( S(\omega) \) defined in (33) saturates at its low-frequency value \( S(\omega) \approx \frac{\hbar \omega}{k_B T} \to 0 \). Therefore, in the tunneling region \( t_i < t < t_f \), the incoherent master
equation Eq. (61) takes the rather symmetric form
\[
\begin{align*}
\frac{\partial P^+_{0}}{\partial t} & \sim \frac{S(0)N^2M^2}{2\hbar^2} (P^+_L \cos^2 \theta_{LZ} + P^+_R \sin^2 \theta_{LZ} - P^+_0), \\
\frac{\partial P^+_1}{\partial t} & \sim \frac{S(0)N^2M^2}{2\hbar^2} (P^+_L \sin^2 \theta_{LZ} + P^+_R \cos^2 \theta_{LZ} - P^+_1), \\
\frac{\partial P^-_R}{\partial t} & \sim \frac{S(0)N^2M^2}{2\hbar^2} (P^-_0 \sin^2 \theta_{LZ} + P^-_1 \cos^2 \theta_{LZ} - P^-_R), \\
\frac{\partial P^-_L}{\partial t} & \sim \frac{S(0)N^2M^2}{2\hbar^2} (P^-_0 \cos^2 \theta_{LZ} + P^-_1 \sin^2 \theta_{LZ} - P^-_L).
\end{align*}
\] (62)

The above master equation represents the open-system dynamics of a fully incoherent quantum annealing processor at the spin glass bottleneck, in the limit $\mathcal{N} \to \infty$. We can vectorize the density matrix populations by defining $P \equiv [P^+_0, P^+_1, P^-_R, P^-_L]^T$. The kinetic equation (62) then takes the following matrix form:
\[
\frac{\partial P}{\partial t} = \frac{S(0)N^2M^2}{2\hbar^2} (\xi^T_1 \sin^2 \theta_{LZ} + \xi^T_2 \cos^2 \theta_{LZ}) - 4_{\times 4} \mathcal{P}.
\] (63)

where here, we have introduced a pseudospin
\[
\begin{align*}
\xi^T_1 & \equiv \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \otimes 1_{2 \times 2}, \\
\xi^T_2 & \equiv 1_{2 \times 2} \otimes \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}.
\end{align*}
\] (64)

The matrix representation Eq. (63) manifestly diagonalizes the Liouvillian for the effective classical master equation (61), allowing us to solve for the density matrix at all times. We begin by defining (anti-)symmetric combinations of the populations via
\[
P_{\tau_1 \tau_2} \equiv (P^+_0 + \tau_2 P^+_1) + \tau_1 (P^-_R + \tau_2 P^-_L).
\] (66)

Equivalently, $P_{\tau_1 \tau_2}$ represents the projection of the vectorized density matrix $P$ onto an arbitrary eigenspace of the time-dependent Liouvillian (63), with $\tau_1, \tau_2 \in \{+1, -1\}$ denoting the eigenvalues under application of $\tau^T_1, \tau^T_2$, respectively. In this eigenbasis, the ground state population upon exiting the crossing is given by
\[
P_{GS}(t_f) \equiv (P^-_R + P^+_0) \bigg|_{t=t_f} = \frac{1}{2} (1 + P_{+-}(t))
\] (67)

The ground state population (67) upon exiting the crossing can be computed analytically with arbitrary initial conditions. In particular, defining an effective rate $\Gamma_{\text{eff}}(t) \equiv S(0)M^2 \cos^2 \theta_{LZ} (B(t))/2\hbar^2$ which sets the timescale of the multiqubit dynamics, we then have
\[
P_{GS}(t_f) \sim \frac{1}{2} \left[ 1 + e^{-2N f^T_1 \Gamma_{\text{eff}}(t) dt} P_{+-}(t_i) \right].
\] (68)

Since this prediction only gives the ground-state population upon exiting the bottleneck region, the above analytical formula is relevant provided that $P_{GS}(t = t_f) = P_{GS}(t = t_{QA}) \equiv P$ (c.f. Eq. (34)). This is the case, e.g. if the density matrix stops evolving non-trivially after the bottleneck (so-called freeze-out, see e.g. [26–28]). Crucially, we witness this stoppage of evolution in numerical simulations of the Redfield master equation in the range $1.55 - 15.5$ mK, as shown Fig. 11. As we can see, up to $\mathcal{N} = 10$ (the performance limit of our simulations), population transfer is nontrivial and limits the performance of the quantum annealer.

We now focus on a specific set of initial conditions, supposing that the four populations in the density matrix are pairwise thermalized. Before the crossing point, we have $E^-_L - E^+_0 < 1$ and $E^-_R - E_1 < 1$. Also, the rates in the annealing process are larger earlier in the annealing (due to larger transverse-field $B$). It is thus reasonable to assume that, entering the crossing (i.e. at $t = t_i$), the density matrix populations satisfy
\[
P^-_R - P^+_0 < 1, \quad P^-_R - P^+_1 < 1.
\] (69)

Since the levels in each pair correspond to states with the fermionic excitation localized at the same location, this is a reasonable assumption to make. In this limit, $P_{+-}(t_i) \rightarrow 0$, and so the success probability (68) is asymptotic to one half.

V. DISCUSSION

In this paper we proposed and analytically solved a model that exhibits the effects of frustration on bottleneck dynamics of quantum annealing. By investigating a simple class of one-dimensional annealing bottlenecks, we are able to rigorously extract the scaling of tunneling rates caused by longitudinal qubit noise, in the large-$\mathcal{N}$ (i.e. complexity-theoretic) limit. In conclusion, in our model, we have found that the effective noise spectral density at an annealing bottleneck is $O(\mathcal{N})$, which is of the order of the Hamming distance between the crossing states. This exact result provides analytical confirmation of and is in agreement with results obtained in [19, 20]. Furthermore, in spite of the non-integrability of the flux bias noise (c.f. (44)), by treating the transverse-field $B$ at the non-perturbative level, we were able to extract the scaling coefficient accurate to all orders in the transverse-field and elucidate its relation to spontaneous magnetization. Finally, we have confirmed using Redfield-type simulations that the ground-state occupation $P_{GS}$ upon exiting a spin glass bottleneck is especially pertinent for the performance of a quantum annealer, and so we have identified a range of temperatures where our analytical scaling formula should directly predict the annealing performance, in the limit of completely incoherent system dynamics. In the future, we will investigate the effects of including coherences, i.e. off-diagonal elements of the density matrix, on the open-system dynamics. This will allow us to gain more nuanced insight into the scaling
performance of combinatorial optimization and sampling problems on near-term quantum annealers.

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where the one-dimensional transverse-field Ising spin glass:

Our starting point is the free-fermion representation of

\[ \begin{align*}
T. Schultz, D. Mattis, and E. Lieb, & Two-Dimensional &
\text{Planar Ising Model as a Soluble Problem of Many Fermions}, &
\text{Reviews of Modern Physics} 85, 808–816 (1952).
\end{align*} \]

This maps a free-fermion model to its corresponding Bogoliubov de-Gennes Hamiltonian. The Bogoliubov
de-Gennes Hamiltonian is a much more lucid representation of a free-fermion model, because the physical data of interest, namely, the quasiparticle operators expressed in the Majorana basis, are given by the eigenvectors of this matrix, and the quasiparticle dispersion corresponds to the eigenvalues. The BdG representation of the one-dimensional transverse-field Ising spin glass is

\[ H_{\text{BdG}}^\pm = \begin{pmatrix} 0 & B(t) & \pm J_N \\
-B(t) & 0 & J_1 \\
-J_1 & \ddots & \ddots \\
\pm J_N & -J_{N-1} & 0 & B(t) \end{pmatrix} \]

(A4)

In the particular case of the Frustrated Ring, after a change-of-basis

\[ \tilde{\Gamma}_j \equiv \Gamma_j + \tilde{\Gamma}_{2N-(j-1)} \]

(A5)

the matrices specified by (A4) admit a \( \mathbb{Z}_2 \) symmetry. It reverses the order of the basis elements, corresponding to the reflection symmetry of the Frustrated Ring \( J_j \rightarrow J_{N-(j-1)} \). As is standard, the symmetry splits our eigenvalue problem into two subspaces, indexed by the eigenvalue under reversion (which we denote by \( \mu \)). The symmetric subspace (i.e. \( \mu = 1 \)) is \( N \)-dimensional, with basis

\[ \tilde{\Gamma}_j \equiv \tilde{\Gamma}_j - \tilde{\Gamma}_{2N-(j-1)} \]

(A6)

whereas the antisymmetric subspace (i.e. \( \mu = -1 \)) is also \( N \)-dimensional, with basis

\[ \hat{\Gamma}_j \equiv \hat{\Gamma}_j - \hat{\Gamma}_{2N-(j-1)} \]

(A7)

Under this splitting, the pair of BdG Hamiltonians (A4) take the block-diagonal form

\[ H_{\text{BdG}}^\sigma \sim \begin{pmatrix} H_{\text{BdG}}^\sigma & 0 \\
0 & H_{\text{BdG}}^\sigma \end{pmatrix} \]

(A8)

where

\[ H_{\text{BdG}}^\sigma,\mu = \begin{pmatrix} -\sigma \mu J_R & B(t) & \pm J_N \\
B(t) & 0 & J_1 \\
-J_1 & \ddots & \ddots \\
\pm J_N & -J_{N-1} & 0 & B(t) \end{pmatrix} \]

(A9)

Where, here, \( \sigma = \pm 1 \) denotes fermion parity. As outlined in the main body of the text, we now find exact low-energy solutions to the BdG equations in the limit \( N \rightarrow \infty \). In this limit, we can treat the BdG boundstate problem as a pair of semi-infinite bound state problems,

\[ H_{\text{BdG}}^\sigma,\mu \psi_{\sigma,\mu} = \epsilon_{\sigma,\mu} \psi_{\sigma,\mu} \]

(A10)
The first such problem looks for a boundstate localized at the right end of the graph:

\[
\psi^R_{\sigma \mu} = \begin{pmatrix}
\alpha_{\sigma \mu} \\
\beta_{\sigma \mu} \\
e^{-\kappa \sigma \mu} \alpha_{\sigma \mu} \\
e^{-\kappa \sigma \mu} \beta_{\sigma \mu} \\
e^{-2\kappa \sigma \mu} \alpha_{\sigma \mu} \\
e^{-2\kappa \sigma \mu} \beta_{\sigma \mu} \\
\vdots
\end{pmatrix}, \quad (A11)
\]

where here, \(\alpha_{\sigma \mu}, \beta_{\sigma \mu}\) are written so that it is clear that they only depend on \(\sigma\) and \(\mu\) through their product. This is because, in the limit \(N \to \infty\), the decay of the mode (A11) means that the matrix element \(\mu B(t)\) in the bottom-right corner of (A9) does not enter the eigenvalue problem (A10). Similarly, we can solve the left-localized eigenvalue problem:

\[
\psi^L_{\mu} = \begin{pmatrix}
\alpha_{\mu} \\
\beta_{\mu} \\
e^{+\kappa \mu} \alpha_{\mu} \\
e^{+\kappa \mu} \beta_{\mu} \\
e^{+2\kappa \mu} \alpha_{\mu} \\
e^{+2\kappa \mu} \beta_{\mu} \\
\vdots
\end{pmatrix}, \quad (A12)
\]

Again, here, \(\alpha_{\mu}, \beta_{\mu}\) are written so that it is clear that they are independent of \(\sigma\). This is because, in the limit \(N \to \infty\), the decay of the mode (A12) means that the only matrix element containing \(\sigma\) (the matrix element in the upper-left corner of (A9)) does not enter the eigenvalue problem (A10).

Substituting the ansätze (A11-A12) into the eigenvalue problem (A10), and letting

\[
e^{-\lambda} = \frac{\alpha}{\beta}
\]

parametrize the ratio of \(\alpha\) to \(\beta\), we get two main types of equations: bulk conditions, and boundary conditions. The equations in the bulk give

\[
\epsilon = e^{-\lambda}(B - Je^{+\kappa}), \quad (A13)
\]

\[
\epsilon = e^{\lambda}(B - Je^{-\kappa}). \quad (A14)
\]

From Eqs. (A13-A14), we get the dispersion relation, as well as (after some hyperbolic trigonometry) some useful bulk identities involving \(\lambda\):

\[
\epsilon^2 = J^2 + B^2 - 2JB \cosh \kappa, \quad (A15)
\]

\[
B^2 = \epsilon^2 + J^2 - 2J \epsilon \cosh (\lambda - \kappa), \quad (A16)
\]

\[
\tanh \lambda = \frac{J \sinh \kappa}{J \cosh \kappa - B}. \quad (A17)
\]

For the boundary conditions, we get different conditions at opposite ends (as would be expected; c.f. (A9)): for the right-localized boundstate ansatz (A11), the boundary conditions are

\[
Be^{-\lambda} = \epsilon + \sigma \mu J_R, \quad (A18)
\]

whereas, for the left-localized boundstate ansatz (A12), the boundary conditions are

\[
B e^{\lambda} = \epsilon - \frac{J^2}{\epsilon - \mu B}. \quad (A19)
\]

The right-boundary condition (A18) generically yields one solution, \(\psi^R_{\sigma \mu}\) for each value of the product \(\sigma \mu\). When \(\sigma \mu = +1\), we denote this solution as \(\psi^R_+\), and when \(\sigma \mu = -1\), we denote this solution as \(\psi^R_-\). For \(B \to 0\), these solutions have the limiting form

\[
\psi^R_{\pm} \bigg|_{B \to 0} \approx \begin{pmatrix} 1 \\ 0 \\ \vdots \end{pmatrix}. \quad (A20)
\]

Similarly, the left-boundary condition (A19) generically yields two solutions for each value of \(\mu\). For \(B \to 0\), the solution whose eigenvalue is least in magnitude limits to

\[
\psi^L_{\mu} \bigg|_{B \to 0} \approx \begin{pmatrix} \vdots \\ 0 \\ 1 \end{pmatrix}. \quad (A21)
\]

In summary, putting together both the left- and right-eigenvalue problems, there are six subgap states in total for the Frustrated Ring benchmark.

To write down the quasiparticle/quasihole excitations corresponding to a given boundstate, we simply contract the boundstate vector with the Gamma matrices. This is succinctly captured by the Feynman slash notation, which is the standard physics notation for such expressions:

\[
\hat{\psi}^j = \sum_{j=1}^{2N} \psi_j \hat{\gamma}_j. \quad (A23)
\]

with \(\psi_j\) the components of the BdG eigenvector in the original basis \(\hat{\gamma}_j\) (c.f. (A4)). Having enumerated the subgap states in equations (A20-A22), we must now physically identify them as localized excitations in our frustrated spin system. To do this, we define

\[
\hat{H}_0^\pm \bigg|_{B \to 0} \approx \sum_{j=1}^{N} J_{j}^\pm \hat{\epsilon}_{j,\pm} \hat{c}_{j,\pm}. \quad (A24)
\]
Cross-matching with Eqs. (A20-A22), we get, in the $\mu = +1$ sector:

$$\hat{\mathcal{E}}_R^R \propto \hat{c}_{2n+1,\pm}, \quad \mu = +1 \quad (A25)$$

$$\hat{\mathcal{E}}_L^R \propto \hat{c}_n, + \hat{c}_{n+1,\pm}, \quad \mu = +1 \quad (A26)$$

$$\hat{\mathcal{E}}_L^L \propto \hat{c}_{n,\pm} - \hat{c}_{n+1,\pm}, \quad \mu = +1 \quad (A27)$$

where here, we have factored $\mathcal{N} \equiv 2n + 1$. Similarly, in the $\mu = -1$ sector, we get:

$$\hat{\mathcal{E}}_R^R \propto \hat{c}_{2n+1,\mp}, \quad \mu = -1 \quad (A28)$$

$$\hat{\mathcal{E}}_L^L \propto \hat{c}_{n,\pm} + \hat{c}_{n+1,\pm}, \quad \mu = -1 \quad (A29)$$

$$\hat{\mathcal{E}}_L^L \propto \hat{c}_{n,\pm} - \hat{c}_{n+1,\pm}, \quad \mu = -1 \quad (A30)$$

This concludes the main exposition of the boundstate problem for the Frustrated Ring. In the following sections, we will apply this knowledge to derive all identities used in the main text.

### a. Definitions of spin-glass excitations used in the main text

Our analysis thus culminates in our first result, which is to give a precise definition of the excitations used in the main text: the identities (A25-A30) derived in the previous subsection physically motivate the following definitions at $B = 0$:

$$\hat{c}_{R,\pm}^\dagger | B=0 \equiv \hat{c}_{2n+1,\pm}, \quad (A31)$$

$$\hat{c}_{L,\pm}^\dagger | B=0 \equiv \hat{c}_{n,\pm} + \hat{c}_{n+1,\pm}, \quad (A32)$$

$$\hat{c}_{L,\pm}^\dagger | B=0 \equiv \hat{c}_{n,\pm} - \hat{c}_{n+1,\pm} \quad (A33)$$

We can now use (A25-A30) to analytically continue the above definitions to non-zero transverse-field $B \neq 0$, by simply tracking the BdG boundstates as they evolve:

$$\hat{c}_{R,\pm}^\dagger (B) := \sum_j (\psi_{\pm}^R_j) \hat{\Gamma}_{j,\pm}^R \quad (A34)$$

$$\hat{c}_{L,\pm}^\dagger (B) := \sum_j (\psi_{\pm}^L_j) \hat{\Gamma}_{j,\pm}^L \quad (A35)$$

$$\hat{c}_{L,\pm}^\dagger (B) := \sum_j (\psi_{\pm}^L_j) \hat{\Gamma}_{j,\pm}^L \quad (A36)$$

Eqs. (A34-A36), along with the definition of the BdG eigenvectors (A11-A12), give a mathematically precise notion of these excitations existing throughout the ordered phase $B < B_c$.

Armed with Eqs. (A34-A36), we can now prove that, in the Frustrated Ring benchmark, the crossing in the odd-parity sector is exact, whereas the even-parity crossing is not. It suffices to look at the positive-energy crossing, as states with energies of opposite sign cannot cross in the spin glass phase $B < B_c$ (by definition). For $\sigma = +1$, the only right-localized boundstate with positive eigenvalue is the quasihole

$$\hat{c}_{R,\pm}^\dagger \quad (A37)$$

By taking the adjoint of (A34), we see that this excitation consists only of $\mu = -1$ Gamma matrices, as (c.f. A5-A7)

$$(\hat{\Gamma}_{j,\pm}^+) \dagger \propto \hat{\Gamma}_{j,-} \quad (A38)$$

In contrast, for $\sigma = -1$, the only right-localized boundstate with positive eigenvalue is the quasiparticle

$$\hat{c}_{R,-}^\dagger \quad (A39)$$

Again, the reversion quantum number for this boundstate satisfies $\mu = +1$, which follows from direct inspection of (A34). On the other hand, on the left-hand-side of the graph, the positive-energy excitation

$$\hat{c}_{L,\pm}^\dagger \quad (A40)$$

has reversion number $\mu = -1$ (c.f. (A35)). Crucially, this quantum number is insensitive to the value of $\sigma$, as $\sigma$ does not show up in the left-boundary condition (c.f. (A19)). In summary, therefore, when $\sigma = -1$, the left- and right-boundstates with positive energy have differing reversion quantum number $\mu$, and thus cannot hybridize. In contrast, when $\sigma = +1$, the left- and right-boundstates with positive energy are both antisymmetric ($\mu = -1$) under reversion, and can thus hybridize. The analysis for the case of the negative energy crossing proceeds in the exact same fashion.

### b. Location of crossings, and scaling of the gap

We now derive the conditions for the existence of a spin glass bottleneck in the Frustrated Ring annealing schedule, as well as compute its location $B = B_h$. Furthermore, we analytically calculate the scaling of the gap at the bottleneck location. To obtain the crossing point of the left- and right-bound state energies, we set them equal to each other, i.e. $\epsilon_{R,\mu}^\dagger = \epsilon_{L,\mu}^\dagger = \epsilon_b$. We will deal with both crossings at the same time: from the discussion in the previous paragraph, we have that the left-localized boundstate involved in the positive-energy crossing has $\mu = -1$ for both the even and odd-crossing, so the left-localized boundstate involved in the positive-energy crossing is always

$$\psi_{L,-} \quad (A41)$$

For the right-localized boundstate involved in the positive-energy crossing, from the discussion in the previous paragraph, the value of $\mu$ depends on the fermion parity $\sigma$. However, we can compute the product $\sigma \mu = -1$, which is the same in both cases. Therefore, the
right-localized boundstate involved in the positive-energy crossing is
\[ \psi^R \]
(A42)

Furthermore, we note that these states have the same energy \( \epsilon = \epsilon_b \), so by the dispersion relation (A15), we have that these crossing states have equal and opposite \( \kappa = \pm \kappa_b \). Therefore, by (A17), the states also have equal and opposite \( \lambda = \pm \lambda_b \). Thus, at the crossing, these states therefore satisfy the boundary conditions
\[ J_b e^{\lambda_b} = \epsilon_b - J_R \]
(A43)
\[ J_b e^{\lambda_b} = \epsilon_b - \frac{J_L^2}{J_R} \]
(A44)
which yield the identity
\[ \epsilon_b + B_b = \frac{J_L^2}{J_R} \]
(A45)

Combining this with one of the bulk equation (A13) yields
\[ e^{\kappa_b - \lambda_b} = \frac{J_R}{J} \]
(A46)

Lastly, we can also consider using the bulk equation (A16). The triplet of equations (A45), (A46) and (A16), considered together, yield the bottleneck location, as well as the crossing energy (e.g., using Solve in Mathematica):
\[ B_b = \frac{1}{J_R} \frac{(J^2 - J_L^2)(J_L^2 - J_R^2)}{J_L^2 + J^2 - 2J_L^2}, \quad \epsilon_b = \frac{1}{J_R} \frac{J_R^2 J^2 - J_L^2}{J_L^2 + J^2 - 2J_L^2} \]
(A47)

Note that \( JJ_R > J_L^2 \) is thus a necessary condition for the crossing point to exist. Also, crucially, the bottleneck location is independent of \( \sigma \), and thus happens in the same location for both the even- and odd-fermion sectors. Using expressions for \( B_b \) and \( \epsilon_b \) to solve for \( \cosh \kappa_b \), we obtain
\[ \kappa_b = \log \frac{J_R(J^2 - J_L^2)}{J(J_L^2 - J_R^2)} \]
(A48)

which determines the scaling of the hybridization of the boundstates, \( \Delta_{\text{min}} \sim \mathcal{O}(e^{-\kappa_bN}) \). The location and the value of the gap can be tuned by adjusting the parameters of the model.

Appendix B: Field-theoretic calculation of the tunneling matrix elements

a. Relation to the corresponding ferromagnetic problem

We now turn to the relations used in the text, which transform a calculation in a transverse-field Ising spin glass to one in a corresponding ferromagnetic model \( \tilde{H} \).

We then take advantage of this transformation to calculate the tunneling rates exactly via the quantum-classical correspondence. Indeed, recall that the transverse-field Ising chain maps to two free fermion models \( H^\pm \) (c.f. (A1)) (Ramond/Neveu-Schwarz), leading to an unphysical doubling of the number of eigenstates. In our calculations up to this point, we have only considered the low-energy states
\[ |\Psi^R_R\rangle = c^\dagger_{R,-} |\Omega^-\rangle, \quad |\Psi^L_L\rangle = c^\dagger_{L,-} |\Omega^-\rangle \]
(B1)
\[ |\Psi^R_R\rangle = |\Omega^-\rangle, \quad |\Psi^L_L\rangle = c^\dagger_{L,+} c^\dagger_{R,+} |\Omega^+\rangle \]
(B2)
and have implicitly discarded the low-energy states
\[ |\tilde{\Psi}^R_R\rangle \equiv |\Omega^-\rangle, \quad |\tilde{\Psi}^L_L\rangle \equiv c^\dagger_{R,-} c^\dagger_{L,-} |\Omega^-\rangle \]
(B3)
\[ |\tilde{\Psi}^R_R\rangle \equiv c^\dagger_{R,+} |\Omega^+\rangle, \quad |\tilde{\Psi}^L_L\rangle \equiv c^\dagger_{L,+} |\Omega^+\rangle \]
(B4)
as they are not genuine eigenstates of the original spin chain Hamiltonian (they have the wrong parity). However, utilizing the identity \( H^\pm = \tilde{H}^\mp (\tilde{H} \text{ is the ferromagnetic modification of the original frustrated Hamiltonian; see Eq. (51) in the main text, c.f. Figure 12 for a depiction of this modification for the case of the Frustrated Ring}) \), we can reinterpret the unphysical states of \( H \) as physical states for \( \tilde{H} \):
\[ |\tilde{\Psi}^R_R\rangle = |\tilde{\Omega}^+\rangle, \quad |\tilde{\Psi}^L_L\rangle = c^\dagger_{R,+} c^\dagger_{L,-} |\tilde{\Omega}^+\rangle \]
(B5)
\[ |\tilde{\Psi}^R_R\rangle = c^\dagger_{R,-} c^\dagger_{L,-} |\tilde{\Omega}^-\rangle, \quad |\tilde{\Psi}^L_L\rangle = c^\dagger_{L,-} |\tilde{\Omega}^-\rangle \]
(B6)
with similar relations for the remaining (i.e. bulk) eigenstates. Therefore, one can interpret the general situation in the following way: when one diagonalizes \( H^+ \) and \( H^- \) in (A1), yielding two full sets of fermionic Fock states, half unphysical, it is actually valid to say that one obtains two full sets of physical spin chain eigenstates, one for the original spin chain, and the remaining set corresponding to the eigenstates of its frustration-free version.

Now, consider the central problem encountered in the main text, namely, that of computing
\[ \langle \Psi^R_R | \sigma_j^z | \Psi^L_L \rangle, \quad \langle \Psi^L_L | \sigma_j^z | \Psi^R_R \rangle \]
(B7)

Naively, we can try to take advantage of the above observation and convert the matrix elements in our glassy problem \( H \) into matrix elements in our ferromagnetic problem \( \tilde{H} \). Indeed, substituting, we get
\[ \langle \Psi^R_R | \sigma_j^z | \Psi^R_R \rangle = \langle \Omega^+ | \sigma_j^z c^\dagger_{R,-} | \Omega^- \rangle \]
\[ = \langle \Omega^- | \sigma_j^z c^\dagger_{R,+} | \Omega^+ \rangle \]
(B8)
\[ \langle \Psi^L_L | \sigma_j^z | \Psi^L_L \rangle = \langle \Omega^- | c^\dagger_{L,+} c^\dagger_{R,+} \sigma_j^z c^\dagger_{R,-} | \Omega^- \rangle \]
\[ = \langle \Omega^- | c^\dagger_{L,-} \sigma_j^z c^\dagger_{R,+} | \Omega^+ \rangle \]
(B9)

Unfortunately, all of these matrix elements are between unphysical states, suggesting that one must pass excitations across the \( \sigma^z \) operator, to amend the situation.
In general, since $\sigma^z$, when written-out in terms of $\hat{\gamma}$-matrices, is an element of the $Pin$ group $Pin(2N)$ (see [29] for a standard reference), passing $\sigma^z$ past a fermionic excitation will rotate that excitation by an orthogonal matrix [29], producing a new fermionic excitation:

$$\psi' = R_j \psi$$

where

$$H' \equiv \sigma^z H \sigma^z$$

with $R_j \in O(2N)$ a rotation matrix. Utilizing this fact, we have

$$\langle \Psi^+_R | \sigma^z_j | \Psi^-_R \rangle = \langle \Omega^+_+ | \sigma^z_j c^\dagger_{R,-} | \Omega^-_+ \rangle$$

$$\langle \Psi^+_L | \sigma^z_j | \Psi^-_L \rangle = \langle \Omega^+_+ | c^\dagger_{L,+} \sigma^z_j | \Omega^-_+ \rangle$$

where in deriving (B13), we have utilized the fact that $R^2_j = 1_{2N \times 2N}$. Computation of the rotation $R_j \in O(2N)$ implemented by $\sigma^z_j$ is straightforward. In particular, by writing everything out in terms of $\hat{\gamma}$ matrices, we find that, for $i < j$, $-1 < 2N - i$,

$$R_j(\hat{\Gamma}_i) = -\hat{\Gamma}_i$$

i.e. $R_j$ flips the reversion-symmetry number quantum. From this, due to the localized nature of the $R$-bound states, we obtain our desired result:

$$R_j(c^\dagger_{R,\pm}) = R_j \left( \sum_i (\psi^R_i \hat{\Gamma}_i^\dagger) \right)$$

$$= -\sum_i (\psi^R_i \hat{\Gamma}_i^\dagger) + O(e^{-\kappa j})$$

which is essential to derive the main results of the paper. Substituting the identity (B15) into Eq.’s (B12-B13) and comparing with Eqs. (B5-B6), we get the relations Eqs. (52) used in the main text:

$$\langle \Psi^+_R | \sigma^z_j | \Psi^-_R \rangle = -\langle \tilde{\Psi}^+_R | \sigma^z_j | \tilde{\Psi}^-_R \rangle + O(e^{-\kappa |j-j'|})$$

$$\langle \Psi^+_L | \sigma^z_j | \Psi^-_L \rangle = -\langle \tilde{\Psi}^+_L | \sigma^z_j | \tilde{\Psi}^-_L \rangle + O(e^{-\kappa |j-j'|})$$

b. Field-theoretic treatment of the ferromagnetic problem

We now give exact expressions for the matrix elements given in Eqs. (B16-B17), involving eigenstates of the ferromagnetic Hamiltonian $H$. We begin by calculating the matrix element corresponding to the ferromagnetic groundstate:

$$\langle \Psi^+_R | \sigma^z_j | \Psi^-_R \rangle$$

Since the frustration-free model is globally gapped in the ordered (i.e. ferromagnetic) phase $B < B_c$, there are no crossings involving the groundstate manifold (See Figure 12), and

$$\langle \Psi^+_R \rangle = | \tilde{\Psi}^+_R \rangle$$

$$\langle \Psi^-_R \rangle = | \tilde{\Psi}^-_R \rangle$$

span the degenerate ground state manifold in the thermodynamic limit of infinite-chain length $N \to \infty$. We begin by showing that the quantity we wish to compute, is really the spontaneous magnetization of our quantum spin chain, in disguise. Since the model is completely ferromagnetic (as we have removed frustration), if we perturb the model with an appropriately-aligned longitudinal field, i.e.

$$\vec{H}_0(\hbar) = \hbar \frac{1}{N} \sum_{j=1}^N \hat{\sigma}_j^z$$

then the groundstate degeneracy is broken. Here, the external field is scaled as $\sim O(1/N)$ in order to ensure
the perturbation is bounded in the thermodynamic limit \( \mathcal{N} \to \infty \). Since the perturbation mixes fermion parity, simple degenerate perturbation theory in the ground-state manifold Eqs. (B19-B20) yields that the perturbed ground-state, in the limit \( h \to 0^+ \), is

\[
|\Psi^{(0)}_{\text{GS}}(0)\rangle = \lim_{N \to \infty} \lim_{h \to 0^+} |\Psi^{(0)}_{\text{GS}}(h)\rangle = \lim_{N \to \infty} \frac{1}{\sqrt{2}} (|\tilde{\Psi}^+_R\rangle + |\tilde{\Psi}^-_R\rangle) \quad (B22)
\]

The reason we consider this perturbation of the frustration-free model is because this allows us to relate our transition matrix element to the \( \text{spontaneous magnetization of this model. Indeed, since} \quad \sigma_j^z \text{ mixes fermion parity,} \quad (B23)

\[
\lim_{N \to \infty} \langle \tilde{\Psi}^+_R | \sigma_j^z | \tilde{\Psi}^-_R \rangle = \lim_{N \to \infty} \frac{1}{2} \langle \tilde{\Psi}^+_R \rangle + \langle \tilde{\Psi}^-_R \rangle \rangle | \sigma_j^z \rangle \rangle = \langle \Psi^{(0)}_{\text{GS}} | \sigma_j^z | \Psi^{(0)}_{\text{GS}} \rangle. \quad (B24)
\]

To compute this spontaneous magnetization, we write

\[
\lim_{N \to \infty} \langle \tilde{\Psi}^+_R | \sigma_j^z | \tilde{\Psi}^-_R \rangle = \lim_{h \to 0^+} \lim_{N \to \infty} \langle \Psi^{(0)}_{\text{GS}}(h) | \sigma_j^z | \Psi^{(0)}_{\text{GS}}(h) \rangle = \lim_{h \to 0^+} \lim_{N \to \infty} \langle \Psi^{(0)}_{\text{GS}}(h) | \sigma_j^z | \Psi^{(0)}_{\text{GS}}(h) \rangle = \langle \Psi^{(0)}_{\text{GS}} | \sigma_j^z | \Psi^{(0)}_{\text{GS}} \rangle. \quad (B25)
\]

This allows us to probe this matrix element using the quantum-classical correspondence: we begin by defining a partition function via

\[
Z_{N,h} \equiv \lim_{T \to 0} \text{Tr}[e^{-\tilde{H}_0(h)/T}]. \quad (B26)
\]

As is standard, to obtain the correspondence with a classical model, we apply the Suzuki-Trotter transformation with a time-step \( \tau_c > 0 \) to the partition function, producing a family of effective actions \( \{S[\tau_c]\}_{\tau_c > 0} \) describing (classical) stochastic fluctuations of an Ising spin system on a cylindrical spacetime lattice. After a straightforward manipulation, one gets [30]:

\[
Z_{N,h} = \sum_{\{s_j\} = \pm 1} e^{-S[\tau_c,h]} \quad (B27)
\]

Here, \( \tau_c \ll 1 \) is a UV cutoff defining a non-perturbative renormalization group flow in imaginary time [31]. For small values of the UV cutoff, the action of the statistical field theory simplifies to [30]:

\[
S[\tau_c,h] \sim_{\tau_c \to 0} \sum_{j, \tau \in \tau_c Z} (J_j[\tau_c] s_j s_{j+1,\tau} + J_{\perp}[\tau_c] s_j s_{j,\tau + \tau_c} + H[\tau_c] s_j,\tau) \quad (B28)
\]

where the coupling constants in our theory have the following dependence on the cutoff (for small values of the cutoff):

\[
J_{\perp}[\tau_c] \equiv \ln \tanh(B \tau_c) , \quad H[\tau_c] \equiv \tau_c h/N \quad \text{and} \quad J_j[\tau_c] \equiv \tau_c J_j.
\]

Thus, we have

\[
\lim_{N \to \infty} \langle \tilde{\Psi}^+_R | \sigma_j^z | \tilde{\Psi}^-_R \rangle = \lim_{J \to 0^+} \lim_{N \to \infty} \lim_{N_\perp \to \infty} \langle s_j,\tau \rangle , \quad (B29)
\]

which is exactly Yang’s definition (as reviewed in [32]) of the spontaneous magnetization of the two-dimensional classical Ising model (B28). Following [32], we denote this with the shorthand \( M_j \), so that Eq. (B29) is equivalently stated as

\[
\lim_{N \to \infty} \langle \tilde{\Psi}^+_R | \sigma_j^z | \tilde{\Psi}^-_R \rangle = M_j \quad (B30)
\]

where it is understood that we are applying Yang’s definition of the spontaneous magnetization to the action Eq. (B28), which lacks translational symmetry in the spatial direction.

In the case that the weights in our MAXCUT problem are uniform in absolute value, i.e. when \( |J_j| \equiv J \) for all \( j \), then the two-dimensional Ising model corresponding to the ferromagnet \( \tilde{H} \) is uniform, and we can cite Yang’s result [33] here for the exact matrix element, which is independent of \( j \):

\[
M_j = M, \quad M \equiv (1 - k^{-2})^{1/8}. \quad (B31)
\]

Here, \( k \) is called the \textit{spectral parameter}, and has the following exact form [29]:

\[
k \equiv \sinh J[\tau_c] \sinh J_{\perp}[\tau_c]. \quad (B32)
\]

In this (uniform) case, we can calculate the spectral parameter in the UV limit of our field theory, in which case we get the ratio \( J/B \) coming from the quantum spin chain:

\[
k = \sinh J[\tau_c] \sinh J_{\perp}[\tau_c] \sim_{\tau_c \to 0} J_{\tau_c} e^{-J_{\perp}[\tau_c]} \sim_{\tau_c \to 0} \frac{J}{B} \quad (B33)
\]

In other words, in the uniform case \( |J_j| \equiv \text{const} \), the matrix element in the spin chain comes out to, in the large-\( N \) limit:

\[
\lim_{N \to \infty} \langle \tilde{\Psi}^+_R | \sigma_j^z | \tilde{\Psi}^-_R \rangle = (1 - k^{-2})^{1/8} \quad (B34)
\]

for all \( B(t) < J \) (i.e. the ordered phase for the ferromagnetic problem).
c. Transfer matrix calculation

In the Frustrated Ring, the couplings are non-uniform. However, the broken translation invariance in the spin chain is due to the modification of only 3 couplers, namely the couplers

\[ J_n = J_{n+1} \equiv J_L, \]
\[ J_{2n+1} \equiv J_R. \]  
(B35)

(Note that here, we are using the values of these couplers in the ferromagnetic version \( H \) of our model). Due to the finite correlation length in the classical model, the effect of local changes to the coupling constants in the theory is washed-out in the thermodynamic limit, when we sum the spontaneous magnetization over all sites \( j \). That is, we can expect the behavior

\[ \sum_{j=1}^{N} M_j^2 \sim N M^2. \]  
(B36)

where, here, \( M \equiv (1 - k^{-2})^{1/8} \) is the result for the uniform chain. To demonstrate the asymptotic result Eq. (B36), we must demonstrate a boundary effect in the classical model. To do this, we calculate the local spontaneous magnetization using the row transfer matrix.

Indeed, consider computing the spontaneous magnetization (representing the tunneling matrix element in the frustrated spin system) using Yang’s algorithm (as reviewed in [32]), but now applied to the row transfer matrix, as opposed to the column transfer matrix:

\[ M_j = \lim_{\alpha \to 0^+} \lim_{N \to \infty} \lim_{N_\perp \to \infty} \frac{\langle s_{j, \tau} \rangle_H}{\langle s_{j, \tau} \rangle_H}. \]
(B37)

Here, the row transfer matrix is (see, e.g. [29])

\[ T[J; H] \equiv \left( \prod_{\tau} e^{J_{\tau}[\tau_\tau] \hat{\sigma}_\tau^z \hat{\sigma}_{\tau+1}^z + H \hat{\sigma}_\tau^z} \right) \left( \prod_{\tau} e^{J_{\tau}[\tau_\tau] (1 + e^{-2J_{\tau}[\tau_\tau]} \hat{\sigma}_\tau^z)} \right). \]  
(B38)

The spectral parameter for this transfer matrix when \( H \equiv 0 \) can be calculated, and comes out to

\[ k_l = \sinh J_l[\tau_\tau] \sinh J_\perp[\tau_\tau] \]
\[ \sim J_l[\tau_\tau] \frac{e^{-J_\perp[\tau_\tau]}}{\tau_\tau} \sim J_l \frac{J_\perp}{B}. \]  
(B39)

Note that, since the model is no longer translation-invariant, this spectral parameter is now dependent on the qubit location \( l \in \{ 1, \cdots, N \} \). We can then proceed with the calculation, letting \( d \) denote the distance between the site \( j \) (where we are calculating the local spontaneous magnetization), and the nearest defect (e.g. the \( J_L \) or \( J_R \) coupler):

\[ M_j = \lim_{\alpha \to 0^+} \lim_{N \to \infty} \lim_{N_\perp \to \infty} \frac{\text{Tr}[T[J; H] \hat{\sigma}_z^d T[J; H] N \text{Tr}[J; H] N \text{Tr}[J; H]]}{\text{Tr}[T[J; H] N \text{Tr}[J; H] N \text{Tr}[J; H]]}. \]

with \( N \equiv n - 1 \) equal to the bulk chain length. We can now see the emergence of a boundary effect in the classical model (and thus, by the quantum-classical correspondence, in the quantum ferromagnet \( \tilde{H} \) as well): letting \( |\Psi_+\rangle \) denote the maximal eigenvector of \( T[J; H] \), we have, by analogous arguments to Eqs. (B19-B23), the following limiting behavior:

\[ \lim_{\alpha \to 0^+} \lim_{N \to \infty} \lim_{N_\perp \to \infty} \frac{|\Psi_+\rangle}{\langle \Psi_+ |} = \frac{1}{\sqrt{2}} (|\Psi_0^+\rangle + |\Psi_{k=0}^-\rangle), \]  
(B40)

where here, \( |\Psi_0^+\rangle \) and \( |\Psi_{k=0}^-\rangle \) are the even- and odd-parity maximal-eigenvectors of \( T[J; H \equiv 0] \) (following the notation of [32]). Therefore, we can write

\[ \lim_{\alpha \to 0^+} \lim_{N \to \infty} T[J; H] = |\Psi_+\rangle \langle \Psi_+ | + O(e^{-d\Delta}), \]  
(B41)

where, here, \( \Delta \) is the spectral gap for the unperturbed transfer matrix \( T[J; H \equiv 0] \), which sets the correlation length in the spatial direction of the lattice. Using Eqs. (B40-B41), we have

\[ \frac{\langle \Psi_+ | \hat{\sigma}_z^d | \Psi_+ \rangle}{\langle \Psi_+ | \Psi_+ \rangle} = (1 - k_0^{-2})^{1/8}, \]  
(B43)

where \( k_0 \) is the spectral parameter for the bulk transfer matrix \( T[J; H \equiv 0] \), which is simply \( J/B \). Therefore, in total, at a distance \( d \) away from either the right- or left-end of the chain,

\[ \langle \tilde{\Psi}_R | \hat{\sigma}_j^z | \tilde{\Psi}_R \rangle \sim N \rightarrow M + O(e^{-\Delta d}). \]  
(B44)

This bulk convergence behavior is confirmed by exact diagonalization with up to \( N = 23 \) sites, see Figure 9 in the main text. The other tunneling matrix

\[ \langle \tilde{\Psi}_L^+ | \hat{\sigma}_j^z | \tilde{\Psi}_L \rangle \]  
(B45)
can be computed similarly: one begins by noting that $|\Psi^+_L\rangle$ is related to $|\Psi^+_R\rangle$ by applying two quasiparticle operators (and the same is true for the relationship between $|\Psi^-_L\rangle$ and $|\Psi^-_R\rangle$). Since these operators are localized at opposite locations of the graph, their product gets mapped, under the Jordan-Wigner transformation, to a string of spin flips $\prod_j \sigma_j^x$ between the centers $j_L$ and $j_R$ of the corresponding bound state wave functions. Therefore, because conjugation by a product of spin flips only has the potential to flip the sign of the magnetization, the tunneling form factors corresponding to (B45) has the exact same asymptotics (a tunneling form factor is given by the squared absolute value of a matrix element of the type considered above).