

Parafermionic phases with symmetry breaking and topological order

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Parafermions are the simplest generalizations of Majorana fermions that realize topological order. We propose a less restrictive notion of topological order in one-dimensional open chains, which generalizes the seminal work by Fendley [J. Stat. Mech. (2012) P11020]. The first essential property is that the ground states are mutually indistinguishable by local, symmetric probes, and the second is a generalized notion of zero edge modes which cyclically permute the ground states. These two properties are shown to be topologically robust, and applicable to a wider family of topologically ordered Hamiltonians than has been previously considered. As an application of these edge modes, we formulate a notion of twisted boundary conditions on a closed chain, which guarantees that the closed-chain ground state is topological, i.e., it originates from the topological manifold of the open chain. Finally, we generalize these ideas to describe symmetry-breaking phases with a *parafermionic* order parameter. These exotic phases are condensates of parafermion multiplets, which generalize Cooper pairing in superconductors. The stability of these condensates is investigated on both open and closed chains.

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I. INTRODUCTION

Parafermions are the simplest generalizations of Majorana fermions [1] that realize topological order [2]. One-dimensional (1D) parafermionic chains have been proposed as emergent degrees of freedom on the edges of two-dimensional (2D), strongly correlated phases, as exemplified by the fractional quantum Hall phases [3]. With open boundaries, a parafermionic chain manifests non-Abelian anyons [4–7] which have envisioned applications in topological quantum computation [8–10]. For practical computation, it is desirable that the anyons robustly survive any generic local perturbation. Following a seminal work by Fendley [2], there is strong evidence that topological order is indeed stable over a family of \mathbb{Z}_N -symmetric Hamiltonians with broken parity and time-reversal symmetries. On an open chain, these so-called chiral Hamiltonians [11–16] satisfy an additional symmetry that renders their *entire* spectra N -fold degenerate [2]. We propose that this symmetry is sufficient but not necessary for a topologically ordered phase since topological order describes the low-energy subspace rather than the entire range of excitations. By relaxing this symmetry constraint, we are led to consider a wider family of topologically ordered Hamiltonians than has been considered in Ref. [2]; these include the nonchiral Hamiltonian which is dual to the widely studied ferromagnetic clock model [17,18]. Despite the loss of degeneracy in the high-energy subspace, the ground-state space remains N -fold degenerate to superpolynomial accuracy in the system size. We are able to demonstrate a stronger statement that the ground states are mutually indistinguishable by *all* local, symmetry-preserving operators, of which the Hamiltonian is but one case in point.

The nontrivial ground-state degeneracy on an open chain is intimately connected to the presence of edge operators which permute the various ground states. The best-known examples

include the dangling Majorana fermion on the edge of the Kitaev chain [7,19], as well as various other proposals to realize Majorana bound states [20–27]. A related signature is the nontrivial degeneracy in the entanglement spectrum, where a real-space bipartition introduces a virtual edge [28,29]. To precisely formulate this “dangling parafermion,” we introduce the notion of a pair of operators $(\mathcal{O}_l, \mathcal{O}_r)$, which localize on left and right edges of an open chain. Each of the pair permutes the ground states cyclically, but does not necessarily commute with the full Hamiltonian, thus differing from past proposals [2,30]. Our generalized edge modes are found to generate a noncommutative algebra within the ground-state space, which for $N = 2$ is the well-known Clifford algebra.

The topological degeneracy that we attribute to “dangling parafermions” is generically lost on a closed chain without edges [29,31]. One might therefore ask the following: Is the nondegenerate, closed-chain ground state continuously connected to the open-chain topological manifold? In this paper, we set out to answer a harder question: What is the closed-chain Hamiltonian (\mathcal{H}) that is *exactly* minimized by a state in the open-chain manifold? We find that \mathcal{H} is obtained by utilizing $\mathcal{Q} \sim \mathcal{O}_l^\dagger \mathcal{O}_r$ as an interedge coupling to close the chain. Since \mathcal{Q} breaks translational invariance, our proposed Hamiltonian differs from conventional methods [32,33] to close the chain. \mathcal{Q} may be interpreted as a topological order parameter, in analogy with the local order parameter (\hat{O}) of traditional broken-symmetry phases. Traditionally, by applying a local, symmetry-breaking “field” $\lambda \hat{O} + \text{H.c.}$, one picks out a broken-symmetry ground state that depends on the phase factor λ . Analogously, by applying $\lambda \mathcal{Q} + \text{H.c.}$ to a topological manifold on an open chain, one selects a closed-chain ground state that originates from this manifold. By exploring the space of λ , we are able to select a different state from the topological manifold, with a different \mathbb{Z}_N charge; this variation of λ is analogous to twisting the boundary conditions in integer

quantum Hall systems [34]. The fundamental difference from \hat{O} is that Q is symmetry preserving but breaks locality.

While topologically ordered parafermions naturally generalize the \mathbb{Z}_2 Kitaev chain [7], some parafermionic phases have no Majorana analog. Indeed, \mathbb{Z}_N parafermions with $N \geq 4$ can also exhibit symmetry breaking, sometimes even in conjunction with topological order [29,31]. Symmetry breaking due to a parafermionic order parameter (\mathcal{Y}) is nonlocal, and thus fundamentally different from traditional symmetry breaking by a local order parameter. Expanding on an interpretation in Ref. [29], we identify these exotic phases as condensates of ‘‘Cooper multiplets,’’ which generalize Cooper pairing in superconductors. A major advance of Ref. [31] is the construction of a class of frustration-free Hamiltonians $H_{(N,n)}$, which realize all possible distinct phases. For these fine-tuned Hamiltonians, the exact form of \mathcal{Y} is now known (when there is symmetry breaking), as is the exact form of the zero edge modes (when there is topological order). On the other hand, the stability of these phases beyond the frustration-free limit $H_{(N,n)}$ has been posed as a conjecture [31]. This lack of quantitative understanding has led to controversy, e.g., with regard to the correct ground-state degeneracy on a closed chain in the presence of symmetry breaking [29,31]. One goal of this paper is to settle this controversy and prove this conjecture.

The outline of our paper: We review the simplest models that realize topological order in Sec. II, and describe the strict notion of a zero edge mode on an open chain, as has been proposed in Ref. [2]. After arguing that this strict notion is sufficient but not necessary for topological order, we then formulate two essential properties of topological phases on an open chain. As described in Sec. III, the first is that the ground-state space is mutually indistinguishable to symmetric, local probes. The second property is a generalized notion of zero edge modes, which we show to have several interesting applications in Sec. IV. One application is the construction of a topological order parameter (Q), which lends us a notion of twisted boundary conditions on a closed chain. This notion is formulated in Sec. V; a critical comparison is made with the conventional method of twisting. In Sec. VI, we generalize these concepts to exotic phases with parafermionic order parameters. The stability of these phases is investigated on both open and closed chains. In the last Sec. VII, we discuss the possible generalizations of our work.

II. REVIEW OF PARAFERMIONIC PHASES WITH TOPOLOGICAL ORDER

Parafermions generalize the Majorana algebra through

$$\gamma_j^N = 1, \quad \gamma_j^\dagger = \gamma_j^{N-1} \quad \text{and} \quad \gamma_j \gamma_l = \omega^{\text{sgn}[l-j]} \gamma_l \gamma_j, \quad (1)$$

for $l \neq j$, and $\omega = e^{i2\pi/N}$. Here, we have denoted γ_j as an operator acting on subcell $j \in \{1, 2, \dots, 2L\}$ on a chain of L sites, in the Hilbert space $(\mathbb{C}^N)^{\otimes L}$. Our convention is to pair subcells $2j-1$ and $2j$ into a single site j , as illustrated in Fig. 1. The simplest, fixed-point Hamiltonians having topological order are dimerized as

$$H_{(N,1)} = -J \sum_{j=1}^{L-1} \sum_{\beta=0}^{N-1} (\omega^{(N-1)/2} \gamma_{2j+1}^\dagger \gamma_{2j})^\beta, \quad (2)$$

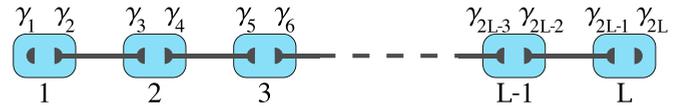


FIG. 1. Parafermion chain of L sites. Each site (blue circle) is split into two subcells.

where each ‘‘even’’ parafermion (on an even subcell) is coupled only to one ‘‘odd’’ parafermion on the adjacent site; these couplings are schematically drawn as black lines in Fig. 1. On an open chain, there is then a dangling parafermion on each end which is uncoupled from the bulk of the chain. For $N = 2$, Eq. (2) is the well-known Kitaev chain with a dangling Majorana mode on each end [7]. $H_{(N,1)}$ has a \mathbb{Z}_N symmetry which is generated by the string operator

$$Q = \prod_{j=1}^L (\omega^{(1-N)/2} \gamma_{2j-1}^\dagger \gamma_{2j}), \quad (3)$$

i.e., $[H_{(N,1)}, Q] = 0$. Q generalizes the fermion parity of Majorana systems through $Q^N = I$. We refer to an eigenstate of Q with eigenvalue ω^α as belonging to the charge sector $\alpha \in \{0, 1, 2, \dots, N-1\} \equiv \mathbb{Z}_N$. Each dangling parafermion $\mathcal{O} \in \{\gamma_1, \gamma_{2L}\}$ is a localized unitary operator satisfying

$$[H_{(N,1)}, \mathcal{O}] = 0 \quad \text{and} \quad \mathcal{O}Q = \omega Q\mathcal{O}; \quad (4)$$

these operators are referred to as ‘‘strong edge modes’’ in a recent review by Alicea and Fendley [3]. The symmetry relations in Eq. (4) imply that the entire spectrum is N -fold degenerate, where each N -multiplet comprises a state in each charge sector. Fendley defines any operator satisfying Eq. (4) as a zero edge mode [2]. We refer to $H_{(N,1)}$ as purely topological, in the sense that it is a fixed point under renormalization group [35], and there is no symmetry breaking by a parafermionic order parameter (cf. Sec. VI). We remark that $H_{(N,1)}$ belongs to a special class of fine-tuned Hamiltonians that is frustration free. To clarify, let us decompose $H_{(N,1)}$ into a sum of two-site operators: $h_{j+1,j} = \sum_{\beta=0}^{N-1} (\omega^{(N-1)/2} \gamma_{2j+1}^\dagger \gamma_{2j})^\beta$. By frustration free, we mean that each of $\{h_{j+1,j}\}$ is a mutually commuting projection, and is individually minimized by each ground state. $h_{j+1,j}$ has an intuitive interpretation in the clock representation of parafermion operators, which we now describe. The clock representation is obtained by the generalized Jordan-Wigner transformation [36]

$$\gamma_{2j-1} = \sigma_j \prod_{k=1}^{j-1} \tau_k, \quad \gamma_{2j} = \omega^{(N-1)/2} \sigma_j \prod_{k=1}^j \tau_k, \quad (5)$$

where σ_j and τ_j act on site j . These clock operators satisfy the algebra

$$\begin{aligned} \sigma_j^N &= \tau_j^N = 1, & \sigma_j^\dagger &= \sigma_j^{N-1}, & \tau_j^\dagger &= \tau_j^{N-1}, \\ \tau_j \sigma_j \tau_j^{-1} &= \omega^* \sigma_j. \end{aligned} \quad (6)$$

By this nonlocal transformation, we find that Eq. (2) is dual to the ferromagnetic clock model [17]:

$$H_{(N,1)} = -J \sum_{j=1}^{L-1} \sum_{\beta=0}^{N-1} (\sigma_j \sigma_{j+1}^\dagger)^\beta, \quad (7)$$

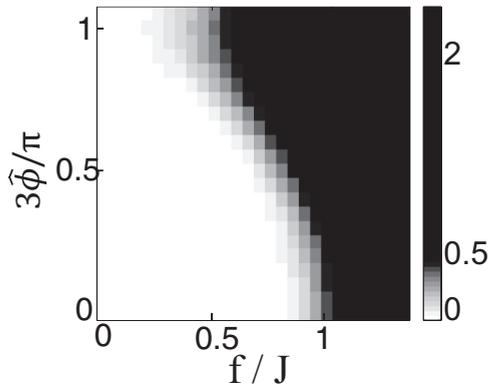


FIG. 2. Numerical simulation of \mathbb{Z}_3 parafermions on a 14-site chain. We plot the absolute spread in energy between three states $\{|\Psi_\alpha\rangle\}$, where $|\Psi_\alpha\rangle$ is the ground state of the open-chain Hamiltonian (9) in the charge sector $\alpha \in \mathbb{Z}_3$.

which has a \mathbb{Z}_N symmetry generated by $Q = \prod_{j=1}^L \tau_j$. To be transparent, let us define an N -dimensional basis on each site j , satisfying

$$\sigma_j|\alpha\rangle_j = \omega^\alpha|\alpha\rangle_j, \quad \tau_j|\alpha\rangle_j = |\alpha + 1 \bmod N\rangle_j, \quad (8)$$

for $\alpha \in \mathbb{Z}_N$. Each of $h_{j+1,j}$ projects to a ferromagnetic alignment of the clock variables on adjacent sites: $|\alpha\rangle_j \otimes |\alpha\rangle_{j+1}$. Thus, $P_{(N,1)}$ exhibits long-range order with respect to the order parameter σ_j , i.e., $\langle \sigma_i \sigma_j^\dagger \rangle$ is nonzero for large $|i - j|$. This type of symmetry breaking by a clock order parameter has a long history [37]. In this paper, we also describe a different type of symmetry breaking by a parafermionic order parameter, which is the subject of Sec. VI.

When deviating from the frustration free $H_{(N,1)}$, it is not obvious if such localized operators (\mathcal{O}) satisfying Eq. (4) still exist. For illustration, we consider a well-studied deformation with \mathbb{Z}_3 symmetry:

$$\hat{H} = -J e^{i\hat{\phi}} \sum_{j=1}^{L-1} \omega \gamma_{2j+1}^\dagger \gamma_{2j} - f \sum_{j=1}^L \omega^* \gamma_{2j-1}^\dagger \gamma_{2j} + \text{H.c.}, \quad (9)$$

which reduces to $H_{(3,1)}$ (up to a constant) if we set $f = \hat{\phi} = 0$. The f term is frustrating, in the sense that it cannot be minimized simultaneously with the J term. This frustration tends to destabilize the ordered phase, but Fendley has nontrivially demonstrated that localized operators satisfying Eq. (4) persist if $\hat{\phi}$ is deformed away from 0 [2]; this is also understood through domain-wall dynamics in the dual clock representation [38]. Since both time-reversal and spatial-inversion symmetries are broken when $\hat{\phi} \neq 0 \bmod \pi/3$, the resulting Hamiltonians are called chiral [11–16]; it has been suggested that the breaking of these symmetries is necessary for topological order [2]. This is peculiar because the nonchiral Hamiltonian (2) exhibits the largest spectral gap as a function of $\hat{\phi}$, thus, one expects that its topological properties are most robust at $\hat{\phi} = 0$. To substantiate this claim, we numerically evaluate the ground-state degeneracy in Fig. 2; this degeneracy is shown to be most persistent at $\hat{\phi} = 0$, where it withstands the largest deformation due to the f term. Our point of view is that topological order is a property of the ground-state space,

and any constraint on the excited states is unnecessary, Eq. (4) being a case in point. This motivates us to formulate two essential characteristics of the ground-state space, which applies to a wider family of Hamiltonians than has been previously considered: (i) a persistent ground-state degeneracy, despite the loss of degeneracy in the excited spectrum (elaborated in Sec. III), and (ii) the existence of generalized zero edge modes (“weak edge modes,” in the language of Ref. [3]) which do not satisfy Eq. (4), but nonetheless generate a noncommutative algebra in the ground-state space, as we will show in Sec. IV.

III. LOCAL INDISTINGUISHABILITY IN PURELY TOPOLOGICAL PHASES

Consider the \mathbb{Z}_3 , nonchiral Hamiltonian in Eq. (9) with $\hat{\phi} = 0$. With $f = 0$ as well, the entire spectrum is threefold degenerate. With nonzero f , it is known that the ground-state splitting is exponentially small in the system size L , while the splittings of the excited triplets are proportional to f/L^2 , for $f/J \ll 1$ [38]. This suggests a general property of topologically ordered systems that a nontrivial ground-state degeneracy can persist without degeneracy in the excited states. In this section, we confirm this general property, and also formulate a stronger statement: that the ground states are mutually indistinguishable by *any* quasilocal, symmetric operator (\mathcal{O}). By quasilocal, we mean that each term in \mathcal{O} has a support that is finite ranged or decaying faster than any power, i.e., superpolynomially (see Appendix A for a precise definition of quasilocality).

For open-chain, quasilocal Hamiltonians that are topologically ordered (without symmetry breaking), we formulate a notion of local indistinguishability in their ground-state spaces. We define $P^{(\alpha)}$ as the nondegenerate ground-state projection in the charge sector labeled by $\alpha \in \mathbb{Z}_N$, i.e., the ground state has eigenvalue ω^α under operation by Q . Let $P_{(N,1)} = \sum_{\alpha=0}^{N-1} P^{(\alpha)}$, and let \mathcal{O} denote *any* quasilocal, symmetry-preserving operator, i.e., $[\mathcal{O}, Q] = 0$. In the thermodynamic limit,

$$P_{(N,1)} \mathcal{O} P_{(N,1)} = c(\mathcal{O}) P_{(N,1)} \quad (10)$$

for some complex number c , as we prove below. Since the open-chain Hamiltonian H_{op} satisfies the same conditions as \mathcal{O} , Eq. (10) also implies that the ground state is N -fold degenerate with energy $c(H_{op})$.

For finite ranged \mathcal{O} , Eq. (10) is satisfied by the frustration free $H_{(N,1)}$ exactly, without finite-size corrections. This is most conveniently proven in the clock representation (7) of $H_{(N,1)}$. Each of N ground states is a classical product state that is fully polarized in the quantum numbers of σ_j : $|\psi_{\alpha,1}\rangle = \bigotimes_{j=1}^L |\alpha\rangle_j$ for $\alpha \in \mathbb{Z}_N$. The property (10) has a simple explanation in the clock representation: local operators have difficulty transforming one polarized state $|\psi\rangle$ into another. To substantiate this intuition, we first define an alternative ground-state basis which diagonalizes Q . For $\alpha \in \mathbb{Z}_N$, let

$$|\phi_{\alpha,1}\rangle = \frac{1}{\sqrt{N}} \sum_{\beta=0}^{N-1} \omega^{-\alpha\beta} |\psi_{\beta,1}\rangle, \quad (11)$$

such that $Q|\phi_{\alpha,1}\rangle = \omega^\alpha|\phi_{\alpha,1}\rangle$; we refer to α as the \mathbb{Z}_N charge of $|\phi_{\alpha,1}\rangle$. Equation (10) amounts to demonstrating that the matrix $\mathcal{M}_{\alpha\beta}^{(1)} = \langle\phi_{\alpha,1}|\mathcal{O}|\phi_{\beta,1}\rangle$ is proportional to the identity. $[\mathcal{O}, Q] = 0$ implies that the off-diagonal elements vanish. The difference in diagonal elements, $\mathcal{M}_{\alpha\alpha}^{(1)} - \mathcal{M}_{\beta\beta}^{(1)}$, reduces to a sum of terms like $\langle\psi_{\mu,1}|\mathcal{O}|\psi_{\nu,1}\rangle$ for $\mu \neq \nu$. If \mathcal{O} has finite range, this quantity vanishes because \mathcal{O} cannot change the polarization on every site. A variant of this proof was first presented in Ref. [39] to explain the robust ground-state degeneracy of the Ising model. If \mathcal{O} decays superpolynomially instead, $\mathcal{M}_{\alpha\alpha}^{(1)} - \mathcal{M}_{\beta\beta}^{(1)}$ has a finite-size correction that we bound in Appendix B.

While deviating from $H_{(N,1)}$ introduces quantum fluctuations to the classical states $\{|\psi_{\alpha,1}\rangle\}$, these fluctuations have a length scale that is suppressed by a spectral gap, thus preserving the property (10) to superpolynomial accuracy in the system size. We are assuming a gapped, quasilocal interpolation H_s between $H_0 \equiv H_{(N,1)}$ and a deformed, but topologically equivalent, Hamiltonian H_1 ; the \mathbb{Z}_N symmetry is preserved throughout $s \in [0, 1]$. By gapped, we mean that the spectral gap above the lowest N states remains finite throughout the interpolation; this implies that the projection P_s to the lowest N states is uniquely defined at each s . At this point, we do not assume that these N states remain degenerate. These conditions allow us to define a locality- and symmetry-preserving unitary transformation \mathcal{V}_s , that maps the low-energy subspaces as

$$P_s = \mathcal{V}_s P_0 \mathcal{V}_s^\dagger, \quad \text{with } P_0 \equiv P_{(N,1)}, [\mathcal{V}_s, Q] = 0. \quad (12)$$

By locality preserving, we mean that if \mathcal{O} is quasilocal, so is its transformed version $\mathcal{V}_s^\dagger \mathcal{O} \mathcal{V}_s$. \mathcal{V}_s is known as an exact quasiadiabatic continuation [39–41]; its properties are elaborated in Appendix A, and its explicit form is shown in the next section. Under these conditions, we find that if P_0 is locally indistinguishable, so is P_s . To prove this, let \mathcal{O} be quasilocal and symmetry preserving, then $P_s \mathcal{O} P_s = \mathcal{V}_s P_0 \mathcal{V}_s^\dagger \mathcal{O} \mathcal{V}_s P_0 \mathcal{V}_s^\dagger$. Since $\mathcal{V}_s^\dagger \mathcal{O} \mathcal{V}_s$ is also quasilocal and symmetry preserving, we may apply Eq. (10) for P_0 . It follows that $P_s \mathcal{O} P_s = c P_s$ for some constant c , up to superpolynomially small finite-size corrections that we bound in Appendix B.

Finally, we point out that local indistinguishability is a unifying property of many other topologically ordered ground states, as we elaborate in Sec. VII.

IV. GENERALIZED ZERO EDGE MODES FOR PURELY TOPOLOGICAL PHASES

The N -fold degeneracy on an open chain arises solely from degrees of freedom on the edges. To crystallize this notion, we would like to construct localized edge operators which permute the ground states cyclically. We already know such a set of operators for the frustration free $H_0 = H_{(N,1)}$: from (5) and (11), it is simple to deduce that $\langle\phi_{\alpha,1}|\gamma_1|\phi_{\alpha+1 \bmod N,1}\rangle = 1$ and $\langle\phi_{\alpha,1}|\gamma_{2L}|\phi_{\alpha+1 \bmod N,1}\rangle = \omega^{(N+1)/2+\alpha}$. Given this information, we can construct similar operators for any Hamiltonian H_s that is connected to H_0 by a gapped, symmetry-preserving interpolation. This will be accomplished by quasiadiabatic continuation, which is known to preserve the matrix elements of operators within the groundstate space [39]. In

more detail, we decompose the ground-state space of H_s as $P_s = \sum_{\alpha=0}^{N-1} |\phi_{\alpha,1}^s\rangle\langle\phi_{\alpha,1}^s|$, where α denotes the \mathbb{Z}_N charge of $|\phi_{\alpha,1}^s\rangle$. Since $[\mathcal{V}_s, Q] = 0$ for the quasiadiabatic continuation of Eq. (12), the \mathbb{Z}_N charge of each state is preserved. It follows that

$$\begin{aligned} \langle\phi_{\alpha,1}^s|\mathcal{O}_l|\phi_{\alpha+1 \bmod N,1}^s\rangle &= 1, \\ \langle\phi_{\alpha,1}^s|\mathcal{O}_r|\phi_{\alpha+1 \bmod N,1}^s\rangle &= \omega^{(N+1)/2+\alpha}, \end{aligned} \quad (13)$$

with the dressed operators

$$\mathcal{O}_l = \mathcal{V}_s \gamma_1 \mathcal{V}_s^\dagger \quad \text{and} \quad \mathcal{O}_r = \mathcal{V}_s \gamma_{2L} \mathcal{V}_s^\dagger. \quad (14)$$

That \mathcal{O}_l is localized around site 1 follows from \mathcal{V}_s being locality preserving. Since both γ_1 and \mathcal{V}_s are unitary, (13) also implies $\mathcal{O}_l|\phi_{\alpha+1,1}^s\rangle = |\phi_{\alpha,1}^s\rangle$.

To gain some intuition about these edge modes, we construct \mathcal{O}_l for the simplest nontrivial example: a frustrated \mathbb{Z}_2 Kitaev chain, as modeled by

$$H_s = i \sum_{j=1}^{L-1} \gamma_{2j} \gamma_{2j+1} + is \sum_{j=1}^L \gamma_{2j-1} \gamma_{2j}. \quad (15)$$

Decomposing $H_s = H_0 + sV$, we point out that H_0 is the frustration-free Kitaev model. The V term tends to destabilize the topological phase, resulting in a monotonic decrease of the spectral gap Γ^s above the two degenerate ground states; it is known that $\Gamma^0 = 2$ in the frustration-free limit, and $\Gamma^1 = 0$, signaling a phase transition [2]. The quasiadiabatic continuation can thus be carried out for s lying in the real interval $[0, \bar{s}]$, with $\bar{s} < 1$ and $\Gamma^{\bar{s}} > 0$. The quasiadiabatic continuation operator has the form of a path-ordered evolution [39]

$$\mathcal{V}_s = T \exp \left[i \int_0^s \mathcal{D}_{s'} ds' \right] \quad (16)$$

in the ‘‘time’’ variable s , generated by the Hamiltonian

$$\mathcal{D}_s = -i \int_{-\infty}^{\infty} dt F(\Gamma^s t) e^{iH_s t} (\partial_s H_s) e^{-iH_s t}. \quad (17)$$

We refer to $F(t)$ as a filter function, whose purpose is to cut off the time evolution of $\partial_s H_s$ for large $|t|$. It is desirable that $F(t)$ has the fastest decay for large $|t|$, such that \mathcal{D}_s is maximally quasilocal [40,41]; this is elaborated in Appendix A. In addition, $F(t)$ is imaginary, so that \mathcal{D}_s is Hermitian. Let us denote the eigenbasis of H_s by $H_s|j;s\rangle = E_j^s|j;s\rangle$. By choosing the Fourier transform of $F(t)$ as

$$\tilde{F}(\Omega) = \int_{-\infty}^{\infty} dt e^{i\Omega t} F(t) = -\frac{1}{\Omega} \quad \text{for } |\Omega| \geq 1, \quad (18)$$

we obtain the matrix elements of \mathcal{D}_s as

$$i\langle k;s|\mathcal{D}_s|j;s\rangle = -\frac{\langle k;s|\partial_s H_s|j;s\rangle}{E_k^s - E_j^s} = \langle k;s|\partial_s|j;s\rangle \quad (19)$$

between states separated by the spectral gap Γ_s , i.e., for $|E_k^s - E_j^s| \geq \Gamma_s$. This leads to $\partial_s P_s = i[\mathcal{D}_s, P_s]$ and Eq. (12). To first order in s , the edge mode has the form

$$\mathcal{O}_l = \mathcal{V}_s \gamma_1 \mathcal{V}_s^\dagger = \gamma_1 + is[\mathcal{D}_0, \gamma_1] + \dots \quad (20)$$

Employing $[H_0, \gamma_1] = 0$ and $[V, \gamma_{2j-1}] = -2i\gamma_j$, we are thus led to evaluate

$$i[\mathcal{D}_0, \gamma_1] = -2i \int_{-\infty}^{\infty} dt F(2t) e^{iH_0 t} \gamma_2 e^{-iH_0 t}, \quad (21)$$

where we have identified $\Gamma^0 = 2$ as the spectral gap in the frustration-free Kitaev chain. Now, apply that $\gamma_2 = (v_+ + v_-)/2$, where $v_{\pm} = \gamma_2 \pm i\gamma_3$ are constants under evolution by the dimerized H_0 : $[H_0, v_{\pm}] = \mp 2v_{\pm}$. We then find

$$i[\mathcal{D}_0, \gamma_1] = -i \int_{-\infty}^{\infty} dt F(2t) (v_+ e^{-i2t} + v_- e^{i2t}) = \gamma_3.$$

In the last step, we employed $\tilde{F}(\pm 1) = \mp 1$, as follows from Eq. (18). We thus derive that the edge mode spreads as $\mathcal{O}_l = \gamma_1 + s\gamma_3 + \dots$, and a simple computation shows $[H_s, \gamma_1 + s\gamma_3] = O(s^2)$. This result is suggestive of a twofold degeneracy in the entire spectrum, as has been derived alternatively in Ref. [2] for higher orders in s . However, this property is not generic, as evidenced by our next illustration. In Appendix C, we work out the edge mode for the \mathbb{Z}_3 nonchiral parafermion, as modeled by \hat{H} in Eq. (9) with $\hat{\phi} = 0$. We identify the J term as H_0 , and f term as the deformation V . The result is

$$\begin{aligned} \mathcal{O}_l = & \gamma_1 + \frac{f}{3J} (\gamma_3 - \omega \gamma_2^\dagger \gamma_3^\dagger - \omega \gamma_1^\dagger \gamma_3^\dagger) \\ & + \frac{f}{3J} \omega^* \gamma_1^\dagger \gamma_2 \gamma_3 + O\left(\frac{f^2}{J^2}\right), \end{aligned} \quad (22)$$

which demonstrably does not commute with H_s . Instead, \mathcal{O}_l commutes with the projected Hamiltonian $P_s H_s P_s$ to superpolynomial accuracy, as follows from (i) $[\gamma_1, P_0] = 0$ leading to $[\mathcal{O}_l, P_s] = 0$, and (ii) the ground-state degeneracy shown in Sec. III. In Appendix C, we have also derived the edge mode \mathcal{O}_l for the \mathbb{Z}_3 chiral parafermion ($0 < \hat{\phi} < \pi/3$). In this parameter range, we find to first order in s that \mathcal{O}_l may be deformed to commute with the full Hamiltonian H_s , at the cost of having a weaker decay away from the edge. As described in Appendix D, this deformation becomes increasingly singular as $\hat{\phi} \rightarrow 0$, i.e., the deformed edge mode delocalizes in the nonchiral limit. The ‘‘zero edge mode’’ (Ψ_{left}) that is alternatively derived in Eq. (32) of Ref. [2] is constructed so that $[\Psi_{\text{left}}, H_s] \approx 0$. For $\hat{\phi} = 0$, Fendley’s construction is not usable, while ours is able to produce Eq. (22). For $0 < \hat{\phi} < \pi/3$, the edge mode \mathcal{O}_l may be deformed to coincide with Ψ_{left} to first order; we do not know if this coincidence persists to higher orders.

Let us present another application of these generalized edge modes. In the ground-state space, they generate a representation of the \mathbb{Z}_N generator:

$$\mathcal{Q} = \omega^{(1-N)/2} \mathcal{O}_l^\dagger \mathcal{O}_r, \quad (23)$$

which satisfies $\mathcal{Q}^N = I$, and $\mathcal{Q}|\phi_{\alpha,1}^s\rangle = \omega^\alpha |\phi_{\alpha,1}^s\rangle$. This may be compared with the representation (3) in the full Hilbert space, which operates on each site. The difference is that \mathcal{Q} is a fractionalized representation, with support only near the chain edges [28,29,31]. Together, \mathcal{O}_l , \mathcal{O}_r , and \mathcal{Q} generate a

noncommutative algebra in the ground-state space:

$$\begin{aligned} \mathcal{O}_l \mathcal{O}_r &= \omega \mathcal{O}_r \mathcal{O}_l, \quad \mathcal{O}_l \mathcal{Q} = \omega \mathcal{Q} \mathcal{O}_l, \\ \text{and } \mathcal{O}_r \mathcal{Q} &= \omega \mathcal{Q} \mathcal{O}_r. \end{aligned} \quad (24)$$

For $N = 2$, this is the familiar Clifford algebra, with \mathcal{O}_l acting as the Pauli matrix ζ^x in the ground-state space, \mathcal{O}_r as ζ^y , and \mathcal{Q} as ζ^z . Clearly, the charge of \mathcal{O}_l is a topological invariant [28,29,31], i.e., it does not change under quasiadiabatic continuation. We say an operator \mathcal{O} has a charge α if $\mathcal{Q}^\dagger \mathcal{O} \mathcal{Q} = \omega^\alpha \mathcal{O}$. In the present context, $[\mathcal{Q}, \mathcal{V}_s] = 0$ implies that \mathcal{O}_l has unit charge, independent of the deformation parameter s . A case in point is the \mathbb{Z}_3 edge mode (22), where each term in the expansion of \mathcal{O}_l has the same charge; the preceding discussion shows that this is true to all powers of f/J .

Finally, the existence of a fractionalized \mathcal{Q} implies that the ground-state degeneracy is unstable to interedge coupling when we close the chain. It is worth refining our notion of locality on a closed chain. A case in point is \mathcal{Q} in the frustration-free limit:

$$\mathcal{Q} = \omega^{(1-N)/2} \gamma_1^\dagger \gamma_{2L} = \sigma_1^\dagger \sigma_L \prod_{j=1}^L \tau_j. \quad (25)$$

Since sites 1 and L are adjacent on a closed chain, \mathcal{Q} is local in the parafermion representation, but nonlocal in the clock representation. In short, we call \mathcal{Q} parafermion local, but clock nonlocal. Fractionalization implies the existence of a parafermion-local term that distinguishes the ground states. That is, we may legitimately add to the open-chain Hamiltonian a term of the form $\lambda \mathcal{Q} + \text{H.c.}$, which closes the chain and singles out a unique ground state. Which ground state is singled out depends on the phase factor λ , as we elaborate in Sec. V. We remark that this refined notion of locality is not necessary for an open-chain Hamiltonian, where charge neutrality imposes that all parafermion-local terms are also clock local; this is proven in Appendix A. Finally, we point out that the local-indistinguishability condition (10) on an open chain applies only for symmetric, *clock*-local probes, as may be verified in the proof of Sec. III; this allows for a clock-nonlocal but parafermion-local term, e.g., \mathcal{Q} , to close the chain and break the ground-state degeneracy.

V. TOPOLOGICAL ORDER ON A CLOSED CHAIN

While parafermionic chains and their dual ferromagnetic clock models have identically enhanced degeneracy on an open chain, their difference is manifest when the chain is closed. Specifically, the topological degeneracy that we attribute to edge modes is generically lost on a closed chain [29,31], while the degeneracy of symmetry-broken clock models persists. One might ask the following for the parafermionic chain: What interedge coupling guarantees that the nondegenerate, closed-chain ground state ($|\Phi\rangle$) is topological, i.e., that $|\Phi\rangle$ is continuously connected to the topological ground-state manifold of the open chain? A harder question would be as follows: Does a coupling exist such that $|\Phi\rangle$ is *exactly* one of the open-chain ground states? The goal of Sec. V A is to prove the existence of this interedge coupling for Majorana chains, and to describe how the coupling is

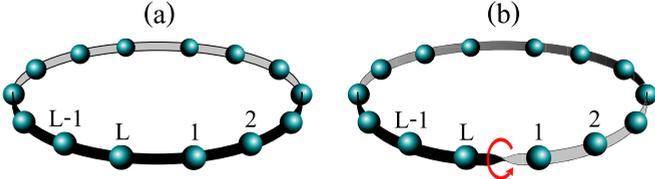


FIG. 3. We formulate a notion of twisted boundary conditions on the closed chain. A topological ground state is especially sensitive to a change in boundary conditions [schematically as (a) to (b)]. Such a twist results in a different ground state with a distinct \mathbb{Z}_N charge.

constructed. In Sec. VB, we perform a critical comparison with the conventional boundary condition that is commonly employed. We then extend our discussion to \mathbb{Z}_N parafermions in Sec. VC.

A. Topological boundary conditions

The simplest illustration lies in the frustration-free Majorana model, whose Hamiltonian on an open chain is

$$H_0 = i \sum_{j=1}^{L-1} \gamma_{2j} \gamma_{2j+1}. \quad (26)$$

As noted in the previous section, the ground-state manifold is topologically ordered, and comprises a state $|\phi_{\alpha,1}\rangle$ in each charge sector $\alpha \in \mathbb{Z}_2$. We would like to construct a closed-chain Hamiltonian ($\mathcal{H}_0^{(\alpha)}$) that is exactly minimized by the open-chain ground state $|\phi_{\alpha,1}\rangle$. This is accomplished by identifying the low-energy degree of freedom on the open chain, which in this example is the complex fermion $\tilde{c} = (\gamma_1 - i\gamma_{2L})/2$. It is simple to verify that $|\phi_{0,1}\rangle$ and $|\phi_{1,1}\rangle$ differ in their occupation of \tilde{c} . Therefore, we are led to $\mathcal{H}_0^{(\alpha)} = H_0 + \delta H_0^{(\alpha)}$, with the interedge coupling

$$\delta H_0^{(\alpha)} = i(-1)^\alpha \gamma_1 \gamma_{2L} = (-1)^\alpha (1 - 2\tilde{c}^\dagger \tilde{c}). \quad (27)$$

The phase factor $(-1)^\alpha$ may be interpreted as flux insertion, and determines the occupation of the fermion (\tilde{c}) in the nondegenerate ground state. Alternatively stated, we obtain

two Hamiltonians by imposing periodic and antiperiodic boundary conditions for the complex fermion $c_j = (\gamma_{2j-1} - i\gamma_{2j})/2$, as schematically illustrated in Fig. 3.

Let us generalize our discussion to frustrated models that remain topologically ordered. For illustration, we may deform the open chain H_0 as

$$H_{XYZ} = \sum_{j=1}^{L-1} (J_x \gamma_{2j-1} \gamma_{2j} \gamma_{2j+1} \gamma_{2j+2} - i J_y \gamma_{2j-1} \gamma_{2j+2} + i \gamma_{2j} \gamma_{2j+1}). \quad (28)$$

We refer to this as the XYZ model, as it can be expressed as a spin-half model through the Jordan-Wigner transformation (5) (cf. Appendix G). As shown in Fig. 4(a) for the range $|J_x|, |J_y| < 1$ (square labeled C), the spectral gap persists above the two lowest-lying states, which we denote by $|\phi_{\alpha,1}^{J_x, J_y}\rangle$ in the charge sector α . The low-energy degree of freedom on the open chain is directly generalized as $\tilde{c} = (\mathcal{O}_l - i\mathcal{O}_r)/2$, where we have dressed the frustration-free edge modes as $\gamma_1 \rightarrow \mathcal{O}_l(J_x, J_y)$ and $\gamma_{2L} \rightarrow \mathcal{O}_r(J_x, J_y)$, following Eq. (14). Such a dressing exists for any deformation within the square C, where a gapped interpolation exists to the frustration-free limit. In the presence of interactions, \tilde{c} represents a many-body excitation. We then close the chain with the interedge coupling

$$\delta H_{XYZ}^{(\alpha)} = (-1)^{\alpha+1} \mathcal{Q} = (-1)^\alpha (1 - 2\tilde{c}^\dagger \tilde{c}), \quad (29)$$

where $\mathcal{Q} = -i\mathcal{O}_l\mathcal{O}_r$ is derived for the XYZ model to be

$$\mathcal{Q} = -i\gamma_1\gamma_{2L} + iJ_y(\gamma_5\gamma_{2L} + \gamma_1\gamma_{2L-4}) - J_x(\gamma_2\gamma_3\gamma_5\gamma_{2L} + \gamma_1\gamma_{2L-4}\gamma_{2L-2}\gamma_{2L-1}) + \dots, \quad (30)$$

to first order in the deformation parameters (cf. Appendix F). Apparently, \mathcal{Q} cannot be derived by translating the bulk terms of Eq. (28) to the edge. Since \mathcal{Q} represents the fermion parity in the open-chain ground-state space (cf. Sec. IV),

$$\delta H_{XYZ}^{(\alpha)} |\phi_{\beta,1}^{J_x, J_y}\rangle = (-1)^{\alpha+\beta+1} |\phi_{\beta,1}^{J_x, J_y}\rangle \quad (31)$$

for $|J_x|, |J_y| < 1$. This implies that $|\phi_{\alpha,1}^{J_x, J_y}\rangle$ separately minimizes the open-chain Hamiltonian (H_{XYZ}) and the interedge

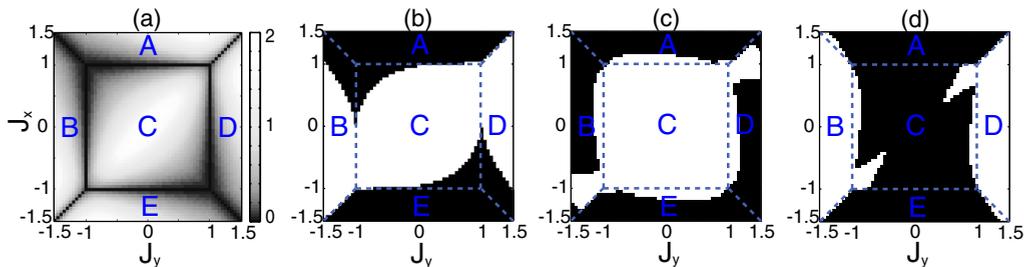


FIG. 4. Characterizing the XYZ model on a chain of $L = 22$ sites, with varying parameters J_x and J_y . (a) The spectral gap above the two lowest states, for the open-chain Hamiltonian (28). As delineated by gapless lines, the regions B, C, and D correspond to topological phases, as explained in the main text. (b)–(d) Evaluate various types of closed boundary conditions by the parity-switch criterion; parameter regions are colored white (resp. black) region where the ground-state parity switches (is invariant) upon twisting the boundary conditions. The blue dashed lines are phase transition lines on the open chain. (b) Translational-invariant boundary conditions, as described in Eq. (35). (c) First-order topological boundary condition (TBC), i.e., Eq. (33) with $\tilde{Q}^{(1)}$ approximating Eq. (30). This boundary condition is constructed for the topological region C only. (d) First-order TBC with $\tilde{Q}^{(1)}$ defined in Appendix F2. This boundary condition is constructed for the topological regions B and D only.

coupling ($\delta H_{XYZ}^{(\alpha)}$), i.e., $|\phi_{\alpha,1}^{J_x, J_y}\rangle$ must be the nondegenerate ground state of the closed-chain Hamiltonian ($H_{XYZ} + \delta H_{XYZ}^{(\alpha)}$). A corollary is that the closed-chain ground state switches parity on twisting the boundary conditions, while the ground-state energy is preserved. In short, we call this permutation by twisting.

While we have focused on one model for specificity, permutation by twisting applies more generally to any closed-chain Hamiltonian that may be decomposed as

$$\tilde{\mathcal{H}}^{(\alpha)} = H_{\text{open}} + (-1)^{\alpha+1} \mathcal{Q}. \quad (32)$$

Here, H_{open} is any open-chain Hamiltonian which is quasiadiabatically connected to H_0 , and $\mathcal{Q} = -i\mathcal{O}_l\mathcal{O}_r$ is the fractionalized representation of the fermion parity in the ground-state space of H_{open} . We have thus established a bulk-edge correspondence: between the existence of topological edge modes on an open chain, to the property of permutation by twisting on a closed chain. We refer to \mathcal{Q} as a topological order parameter, in analogy with the clock-local order parameter (\hat{O}) of traditional broken-symmetry systems. There, the perturbation $\lambda\hat{O} + \text{H.c.}$ picks out a broken-symmetry ground state which depends on λ . The difference is that \hat{O} breaks the symmetry but preserves clock locality, while \mathcal{Q} preserves the symmetry and parafermion locality, but breaks clock locality.

In practice, the exact form of \mathcal{Q} is not easily found. However, the first-order truncation ($\tilde{\mathcal{Q}}^{(1)}$) of \mathcal{Q} is analytically tractable, e.g., Eq. (30) without the dots is an approximation to \mathcal{Q} to first order in the deformation parameters (J_x, J_y). We thus describe the closed-chain Hamiltonian

$$\mathcal{H}_{XYZ}^{(\alpha)} = H_{XYZ} - (-1)^\alpha \tilde{\mathcal{Q}}^{(1)} \quad (33)$$

as having a first-order topological boundary condition (TBC), which we proceed to evaluate. Here and in future sections, we use the parity switch as the simplest diagnostic to evaluate how ‘‘topological’’ a closed-chain Hamiltonian is. Ideally, the nondegenerate ground-state parity switches upon twisting the boundary condition, for all parameters where the open-chain Hamiltonian is topologically ordered. Indeed, we show in Fig. 4(c) that the parity switch occurs (whitened regions) in nearly the entire square C. We remark that the boundary condition (33) is specifically constructed for the topological region C. As we will shortly clarify, regions B and D in Fig. 4 are also topological, but they require a different set of TBC’s.

To critically evaluate the effect of the first-order dressing, we also consider a boundary condition that is zeroth order in J_x and J_y , i.e., we apply the interedge coupling (32) with \mathcal{Q} replaced by $\tilde{\mathcal{Q}}^{(0)} = -i\gamma_1\gamma_{2L}$. For this zeroth-order TBC, we compute the corresponding parity-switch diagram in Fig. 5(a). We find that the parity switch remains robust along $-1 < J_x = J_y < 1$, where the open-chain ground states are completely classical; the ground states of the ferromagnetic XXZ model are well known [42] to be fully polarized. In this classical limit, one may verify that $\tilde{\mathcal{Q}}^{(0)}$ and \mathcal{Q} have identical matrix elements within the ground-state space. This coincidence is lost away from the XXZ line, where quantum fluctuations play an important role. Where fluctuations are strongest (in the orthogonal direction: $-1 < J_x = -J_y < 1$), the parity switch does not always occur: the first-order TBC significantly outperforms its zeroth-order counterpart.

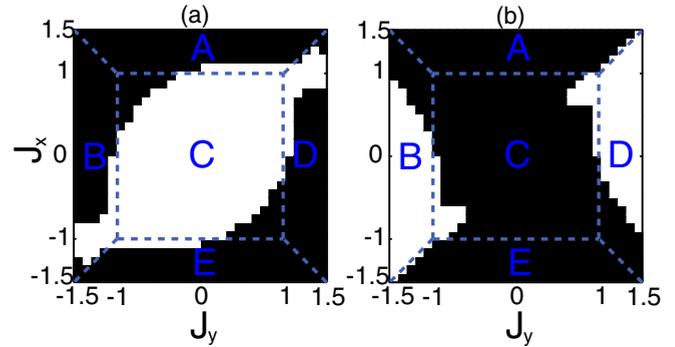


FIG. 5. Evaluation of the zeroth-order topological boundary conditions, for the XYZ model on a closed chain of $L = 22$ sites. (a) We employ the interedge coupling $(-1)^\alpha i\gamma_1\gamma_{2L}$. (b) Here, we employ $(-1)^\alpha J_y i\gamma_2\gamma_{2L-1}$. White (resp. black) region: the ground-state parity switches (is invariant) upon twisting the boundary conditions.

To conclude our discussion of the XYZ model, we remark that regions B and D (see Fig. 4) *also* correspond to a topological phase, by their quasiadiabatic connection to a *different* dimer Hamiltonian:

$$H'_0 = -iJ_y \sum_{j=1}^{L-1} \gamma_{2j-1}\gamma_{2j+2}, \quad (34)$$

which has edge modes γ_2 and γ_{2L-1} . Consequently, the TBC in regions B and D differs from Eq. (30), as we show in Appendix F2. By the parity-switch criterion, the first-order TBC [Fig. 4(d)] is shown to be more effective than the zeroth-order TBC’s [Fig. 5(b)].

B. Comparison with the translational-invariant boundary condition

Having introduced a different type of boundary condition, we would like to make a critical comparison with the conventional method to close chains. Assuming that the open-chain Hamiltonian (H_{open}) is translational invariant up to the edge, it is common practice [32,33] to translate the bulk terms of H_{open} to the edge, and then add phase factors for different flux insertions. For the XYZ model, this procedure produces the closed-chain Hamiltonian:

$$\mathcal{K}_{XYZ}^{(\alpha)} = H_{XYZ} + J_x\gamma_1\gamma_2\gamma_{2L-1}\gamma_{2L} - i(-1)^\alpha J_y\gamma_2\gamma_{2L-1} + i(-1)^\alpha \gamma_1\gamma_{2L}, \quad (35)$$

as we systematically derive in Appendix E; there, we also generalize this procedure for \mathbb{Z}_N parafermions. By construction, $\mathcal{K}^{(\alpha)}$ is translational invariant modulo a ± 1 phase factor. A stronger statement is that $\mathcal{K}^{(0)}$ (resp. $\mathcal{K}^{(1)}$) is completely translational invariant in the even (resp. odd) charge sector, as we demonstrate in Appendix G. Therefore, we describe $\mathcal{K}^{(\alpha)}$ as having a translational-invariant boundary condition (TIBC). In comparison, the TBC generically breaks translational invariance, as exemplified in Eq. (30).

It is commonly believed that if H_{open} is topologically ordered, then the closed chain $\mathcal{K}^{(\alpha)}$ must also be topological, implying that the ground state of $\mathcal{K}^{(\alpha)}$ is especially sensitive to its boundary conditions [32,33]. We numerically test this

expectation for the XYZ model on a finite-size chain [see Fig. 4(b)]. The ground-state parity is invariant where quantum fluctuations are dominant, as we have also seen qualitatively for the zeroth-order TBC [Fig. 5(a)]; both these boundary conditions are outperformed by the first-order TBC [see Fig. 4(c)]. A more thorough finite-size analysis in Appendix G suggests these conclusions are robust in the thermodynamic limit.

C. Closing the chain for general $\mathbb{Z}_{N>2}$ parafermions

For $\mathbb{Z}_{N>2}$ parafermions on an open chain, the low-energy degree of freedom is no longer a complex fermion \tilde{c} . Nevertheless, the N eigenvalues of \mathcal{Q} correspond to distinct states in the ground-state manifold, and the generalized edge modes ($\mathcal{O}_l, \mathcal{O}_r$) induce many-body excitations within this manifold, as per the algebra in Eq. (24). To select a topological ground state on a closed chain, we propose the interedge coupling

$$\delta H^{(\alpha)} = -\frac{1}{N} \sum_{\beta=0}^{N-1} (\omega^\alpha \mathcal{Q})^\beta, \quad (36)$$

which projects to the nondegenerate ground state with charge $N - \alpha$, for $\alpha \in \mathbb{Z}_N$. Twisting the Hamiltonian ($\delta H^{(\alpha)} \rightarrow \delta H^{(\alpha+1)}$) cyclically permutes the \mathbb{Z}_N charge of the nondegenerate ground state, while preserving the ground-state energy.

To see this, let us label the open-chain ground states as $|\phi_{\alpha,1}^s\rangle$ in the charge sector α ; recall that s parametrizes the deformation of the open-chain Hamiltonian: $H_s = H_0 + sV$. By construction, $\delta H^{(\alpha)}$ singles out $|\phi_{N-\alpha,1}^s\rangle$ as having the lowest eigenvalue: $\delta H^{(\alpha)}|\phi_{N-\alpha,1}^s\rangle = -|\phi_{N-\alpha,1}^s\rangle$. $|\phi_{N-\alpha,1}^s\rangle$ has to be the nondegenerate ground state of $H_{op} + \delta H^{(\alpha)}$. This follows from $\mathcal{O}_l, \mathcal{O}_r$, and \mathcal{Q} being unitary, which implies the operator norm $\|\delta H^{(\alpha)}\|_{op} \leq 1$ by the triangle inequality. Thus, $|\phi_{N-\alpha,1}^s\rangle$ separately minimizes H_{op} and $\delta H^{(\alpha)}$.

For illustration, we return to the \mathbb{Z}_3 , open-chain model (9), and now add an interedge coupling in the form of Eq. (36). In the frustration-free limit ($f = 0$), the edge modes are γ_1 and γ_{2L} , and therefore the topological order parameter equals $\omega^* \gamma_1^\dagger \gamma_{2L}$. In our numerical simulation, we ignored the dressing of \mathcal{Q} for nonzero f , and used its frustration-free form throughout; once again, we refer to this as zeroth-order TBC. Despite this approximation, the zeroth-order TBC correctly identifies the topological region in the range of Hamiltonians that we explored; see Fig. 6(a), where the topological region (colored white) indicates a \mathbb{Z}_3 permutation of the nondegenerate ground-state charge, as we twist the boundary conditions. To support our claim that this region is topological, we note its close overlap with the white region of Fig. 2, where the ground state is threefold degenerate on an open chain.

VI. GENERALIZATION TO PHASES WITH SYMMETRY BREAKING

A new type of phase can occur for \mathbb{Z}_4 parafermions, where the ground state $|\theta\rangle$ has a reduced \mathbb{Z}_2 symmetry. This means that $|\theta\rangle$ no longer has a conserved charge under \mathcal{Q} , but it does so with \mathcal{Q}^2 , i.e., it is a condensate [29] of parafermion pairs. A general classification of \mathbb{Z}_N parafermions allows for these

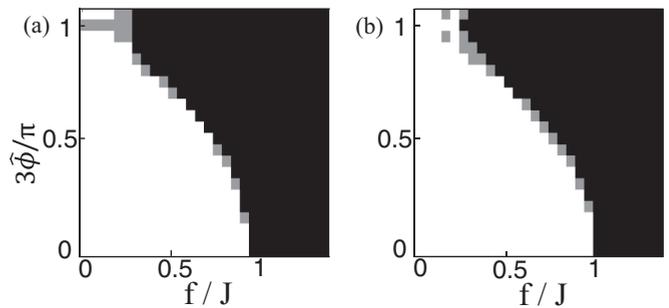


FIG. 6. Simulation of the \mathbb{Z}_3 model (9) on a 14-site chain. (a) The chain is closed with an interedge coupling $\delta H^{(\alpha)} = -\omega^{\alpha-1} \gamma_1^\dagger \gamma_{2L}/3 + \text{H.c.}$, which we call zeroth-order TBC. The figure represents the nondegenerate-ground-state charge $Q_0^{(\alpha)}$, as we twist the closed-ring Hamiltonian through all \mathbb{Z}_3 values of the twist parameter α . White region: Q_0 is maximally permuted, i.e., $\{Q_0^{(\alpha)}\}$ takes on three values, indicating a topological phase. Gray: two. Black: one. (b) For comparison, we have also simulated the translational-invariant boundary condition, which involves the coupling $\delta H^{(\alpha)} = -J\omega^{\alpha-1} e^{i\phi} \gamma_1^\dagger \gamma_{2L} + \text{H.c.}$

coherent states, sometimes in conjunction with topological order. All distinct phases are uniquely labeled by n , which is a divisor of N [29,31]. The $n = 1$ phase is purely topological, and has been characterized in the previous sections; we now turn our attention to $1 < n < N$. It is useful to define $m = N/n$ and the greatest common divisor of m and n : $g = \text{gcd}(m, n)$. If $g > 1$, there is symmetry breaking with a parafermionic order parameter; if $m > g$, the phase is topological ordered; if $m > g > 1$, symmetry breaking coexists with topological order. A major advance of Ref. [31] is the construction of frustration-free models which realizes each of these phases on an open chain:

$$H_{(N,n)} = -\frac{1}{m} \sum_{j=1}^{L-1} \sum_{\beta=0}^{m-1} (\omega^{(N-1)/2} \gamma_{2j+1}^\dagger \gamma_{2j})^{\beta n} - \frac{1}{n} \sum_{j=1}^L \sum_{\alpha=0}^{n-1} (\omega^{(1-N)/2} \gamma_{2j-1}^\dagger \gamma_{2j})^{\alpha m}. \quad (37)$$

For example, $H_{(4,2)}$ realizes the above-mentioned condensate of parafermion pairs. $H_{(N,n)}$ is a sum of mutually commuting projections, and generalizes the dimer model (2). Many ground-state properties are now known in the frustration-free limit [31], including the form of the parafermionic order parameters and the topological edge modes; these properties are reviewed in Sec. VIA. The aim of Sec. VIB is to demonstrate that the phases remain stable even if frustrated on an open chain. Section VIC extends the stability analysis to a closed chain.

A. Review of general parafermionic phases

We begin by describing the ground-state properties of $H_{(N,n)}$, of which many are known from Ref. [31], then extend them to more general Hamiltonians by quasiadiabatic continuation. It is instructive to interpret $H_{(N,n)}$ in the clock representation, where *all* nontrivial phases arise from traditional symmetry breaking with a clock-local order

parameter. For illustration, we consider the clock analog of the parafermion-pair condensate. One possible ground state of $H_{(4,2)}$ is fully polarized as $\bigotimes_{j=1}^L (|0\rangle_j + |2\rangle_j)$. This state breaks the Z_4 symmetry under $|\alpha\rangle_j \rightarrow |\alpha + 1 \bmod 4\rangle_j$, but retains a reduced symmetry under $|\alpha\rangle_j \rightarrow |\alpha + 2 \bmod 4\rangle_j$. More generally,

$$H_{(N,n)} = -\frac{1}{m} \sum_{j=1}^{L-1} \sum_{\beta=0}^{m-1} (\sigma_j \sigma_{j+1}^\dagger)^{\beta n} - \frac{1}{n} \sum_{j=1}^L \sum_{\alpha=0}^{n-1} \tau_j^{\alpha m} \quad (38)$$

has m classical ground states $|\psi_{\alpha,n}\rangle$, which are fully polarized in the quantum numbers of σ_j^n :

$$|\psi_{\alpha,n}\rangle = \bigotimes_{j=1}^L \frac{1}{\sqrt{n}} \sum_{\beta=0}^{n-1} |\alpha + m\beta\rangle_j, \quad (39)$$

where $\alpha \in \mathbb{Z}_m$, and $\sigma_j |\alpha\rangle_j = \omega^\alpha |\alpha\rangle_j$. The \mathbb{Z}_N generator (3) acts in this basis as $Q |\psi_{\alpha,n}\rangle = |\psi_{\alpha+1 \bmod m,n}\rangle$, as follows from (8). In the charge eigenbasis,

$$|\phi_{\alpha,n}\rangle = \frac{1}{\sqrt{m}} \sum_{\beta=0}^{m-1} \omega^{-n\alpha\beta} |\psi_{\beta,n}\rangle \quad (40)$$

with $\alpha \in \mathbb{Z}_m$ and $Q |\phi_{\alpha,n}\rangle = \omega^{n\alpha} |\phi_{\alpha,n}\rangle$. Since $|\psi_{\alpha,n}\rangle$ are fully polarized in the clock representation, clock-local operators have difficulty transforming one $|\psi\rangle$ into another; this transformation would involve turning the clocks on all sites. This observation leads to the local indistinguishability of the ground-state space $P_{(N,n)} = \sum_{\alpha=0}^{m-1} |\phi_{\alpha,n}\rangle \langle \phi_{\alpha,n}|$, i.e., $P_{(N,n)}$ satisfies the same condition as $P_{(N,1)}$ in Eq. (10), when probed by any symmetric, clock-local operator (\mathcal{O}). To show this, we first suppose the probe \mathcal{O} has a finite range. We would like to show that the matrix $\mathcal{M}_{\alpha\beta}^{(n)} = \langle \phi_{\alpha,n} | \mathcal{O} | \phi_{\beta,n} \rangle$ is proportional to the identity. The off-diagonal elements vanish because $[\mathcal{O}, Q] = 0$. The difference in diagonal elements, $\mathcal{M}_{\alpha\alpha}^{(n)} - \mathcal{M}_{\beta\beta}^{(n)}$, equals a sum of terms proportional to $\langle \psi_{\mu,n} | \mathcal{O} | \psi_{\nu,n} \rangle$, with $\mu \neq \nu$. This quantity vanishes due to the above-mentioned observation. If \mathcal{O} is not finite ranged but decays superpolynomially, $\mathcal{M}_{\alpha\alpha}^{(n)} - \mathcal{M}_{\beta\beta}^{(n)}$ vanishes up to finite-size corrections that we bound in Appendix B. An instructive basis for $P_{(N,n)}$ is

$$|\theta_{\beta,\delta}\rangle = \frac{1}{\sqrt{g}} \sum_{\alpha=0}^{g-1} e^{-i2\pi\alpha\beta/g} |\phi_{\alpha m/g + \delta,n}\rangle, \quad (41)$$

where $\alpha m/g + \delta \in \mathbb{Z}_m$, $\beta \in \mathbb{Z}_g$ may be interpreted as a broken-symmetry index, and $\delta \in \mathbb{Z}_{m/g}$ as a topological index; recall $g = \gcd(m,n)$. On application of a symmetry-reducing onsite ‘‘field’’: $\lambda \gamma_j^{N/g} + \text{H.c.}$, the m multiplet splits into g number of (m/g) multiplets, as determined by

$$\begin{aligned} \gamma_{2j-1}^{N/g} |\theta_{\beta,\delta}\rangle &= e^{-i2\pi\beta/g} |\theta_{\beta,\delta}\rangle \quad \text{and} \\ \gamma_{2j}^{N/g} |\theta_{\beta,\delta}\rangle &= \omega^{N^2(g-1)/2g^2} e^{-i2\pi\beta/g} |\theta_{\beta,\delta}\rangle. \end{aligned} \quad (42)$$

For any subcell j , $\gamma_j^{N/g}$ are parafermionic order parameters which reduce the \mathbb{Z}_N symmetry of $H_{(N,n)}$ to $\mathbb{Z}_{N/g}$, i.e., each (m/g) multiplet has a residual symmetry generated by Q^g . If $m > g$, there exists a remnant (m/g) -fold degeneracy which originates from topological edge modes. These localized

operators (γ_1^n and γ_{2L}^n) commute with $H_{(N,n)}$ and permute the ground states as $|\theta_{\beta,\delta}\rangle \rightarrow |\theta_{\beta,\delta-1 \bmod m/g}\rangle$.

The exact form of $|\theta\rangle$ allows us to expand upon their interpretation [29] as coherent states. Equation (42) implies that $|\theta\rangle$ is a condensate of the operator $\varphi_j = \gamma_j^{N/g}$. This condensate is not of Bose-Einstein type since

$$[\gamma_i^{pN/g}, \gamma_j^{qN/g}] = 0 \quad \text{for } p, q \in \mathbb{Z}, \forall i, j, \quad (43)$$

implying that $[\varphi_i, \varphi_j^\dagger] = 0 \neq \delta_{ij}$. Nevertheless, off-diagonal long-range order manifests as $\langle \theta_{\beta,\delta} | \varphi_{2i-1}^\dagger \varphi_{2j-1} | \theta_{\beta,\delta} \rangle = 1$, independent of $|i - j|$. Note that $[\varphi_i, \varphi_j] = [\varphi_i^\dagger, \varphi_j^\dagger] = 0$, as would a bosonic operator. The present situation is reminiscent of long-range order in the BCS wave function, where the order parameter is almost bosonic. There are essential differences: (a) the ground-state expectation $\langle \theta_{\beta,\delta} | \varphi_j | \theta_{\beta,\delta} \rangle$ takes on g discrete values [cf. Eq. (42)], in contradiction with conventional BCS wave functions that have a U(1) degree of freedom. (b) The order parameter φ manifests an attraction between N/g parafermions ($g > 1$), thus generalizing Cooper pairs to ‘‘Cooper multiplets’’ if $N/g > 2$.

B. Robustness of general parafermionic phases on an open chain

Let us address the stability of these phases as we symmetrically deform $H_{(N,n)}$ to a new Hamiltonian H_s . As long as the deformation preserves the gap above the lowest m states, there exists a quasiadiabatic continuation (\mathcal{V}_s) which maps their respective ground-state spaces as $P_{(N,n)}$ to P_s . Thus if $P_{(N,n)}$ is indistinguishable to clock-local probes (as shown in Sec. VI A), then this property robustly carries forward to P_s , following a simple generalization of Sec. III. Moreover, in close analogy with Sec. IV, we find for $g > 1$ generalized order parameters $\mathcal{Y}_j = \mathcal{V}_s \varphi_j \mathcal{V}_s^\dagger$, which are dressed versions of $\varphi_j = \gamma_j^{N/g}$. Dressing \mathcal{Y}_j may be interpreted as spreading the Cooper-multiplet wave function, in contrast with the tightly bound φ_j . By construction, \mathcal{Y}_j permutes the ground states of H_s , in the same manner that φ_j would for $H_{(N,n)}$. While $[\varphi_j, H_{(N,n)}] = 0$, \mathcal{Y}_j commutes only with the ground-state projected H_s .

If the deformed Hamiltonian H_s is topologically ordered, there exist generalized edge modes $\{\mathcal{O}_{l,n}, \mathcal{O}_{r,n}\}$ which permute the ground states, and are related by quasiadiabatic continuation to $\{\gamma_1^n, \gamma_{2L}^n\}$; here, the subscript l (resp. r) indicates that the operator is localized on the left (resp. right) edge. These edge modes generate a fractionalized representation of the $\mathbb{Z}_{N/g}$ generator:

$$\mathcal{Q}_g = \omega^{xn(xn-N)/2} (\mathcal{O}_{l,n})^{-x} (\mathcal{O}_{r,n})^x, \quad (44)$$

with integer x uniquely satisfying $xn = g \bmod m$. This operator acts like Q^g in the ground-state space of $H_{(N,n)}$, i.e., in the basis (41),

$$\mathcal{Q}_g |\theta_{\beta,\delta}\rangle = \omega^{g n \delta} |\theta_{\beta,\delta}\rangle. \quad (45)$$

Decomposing \mathcal{Q}_g into two operators with support on opposite ends of the chain, the charge $(-x)$ of the left-localized operator $[(\mathcal{O}_{l,n})^{-x}]$ is a topological invariant. Within each (m/g) multiplet labeled by β , $\{\mathcal{O}_{l,n}, \mathcal{O}_{r,n}, \mathcal{Q}_g\}$ generate a

noncommutative algebra:

$$\begin{aligned} \mathcal{O}_{l,n}\mathcal{O}_{r,n} &= \omega^{n^2}\mathcal{O}_{r,n}\mathcal{O}_{l,n}, & \mathcal{O}_{l,n}\mathcal{Q}_g &= \omega^{ng}\mathcal{Q}_g\mathcal{O}_{l,n}, \\ \text{and } \mathcal{O}_{r,n}\mathcal{Q}_g &= \omega^{ng}\mathcal{Q}_g\mathcal{O}_{r,n}. \end{aligned} \quad (46)$$

By our assumption of topological order ($m > g$), the phase factors ω^{n^2} and ω^{ng} are never trivially unity.

C. Robustness of general parafermionic phases on a ring

On a ring, the degeneracies due to topological edge modes are generically lost; for $g > 1$, there remains a g -fold degeneracy on the ring due to broken symmetry. The instability of the topological degeneracy may be understood in this way: the edge modes themselves imply the existence of a \mathbb{Z}_N singlet \mathcal{Q}_g which splits the topological degeneracy. Suppose we coupled the edges as

$$\delta\mathcal{H}^{(\alpha)} = -\omega^{g\alpha}\mathcal{Q}_g + \text{H.c.} \quad \text{with } \alpha \in \mathbb{Z}_{m/g}. \quad (47)$$

This coupling splits the open-chain m multiplet into (m/g) number of g multiplets, as indexed by $\delta \in \mathbb{Z}_{m/g}$ in Eq. (45). By twisting the interedge coupling $\delta\mathcal{H}^{(\alpha)}$ as $\alpha \rightarrow \alpha + 1$, we permute the ground-state multiplet as $\delta = m/g - \alpha \rightarrow m/g - \alpha - 1$.

It is interesting to determine if this g -fold degeneracy persists as we frustrate the closed-chain Hamiltonian. Reference [29] claims that the degeneracy is broken, while Ref. [31] suggests that the degeneracy is robust, though without proof. The intuition for a general proof can be obtained from the simplest symmetry-broken model ($H_{(4,2)}$) without topological order. Here, $m = n = g = 2$. The \mathbb{Z}_4 symmetry is reduced to \mathbb{Z}_2 due to the parafermion order parameter γ_j^2 , which manifests in a twofold degenerate ground state $|\psi_{\alpha \in \mathbb{Z}_2, 2}\rangle$ on both open and closed chains [cf. Eq. (39)]. We would like to know if this degeneracy is stable under deformations of $H_{(4,2)}$. The allowed perturbations are parafermion local and charge neutral; all such perturbations on an open chain are also clock local, as proven in Appendix A. We have shown in the preceding paragraph that transforming $|\psi_{1,2}\rangle$ into $|\psi_{2,2}\rangle$ is impossible with clock-local operators, thus, the degeneracy is stable on an open chain. However, a closed chain allows for interedge perturbations which are parafermion local but clock nonlocal, a case in point being $\gamma_1^\dagger\gamma_{2L} = \omega^{3/2}\sigma_1^\dagger\sigma_L Q$. What remains is to demonstrate the stability of the closed-chain degeneracy under clock-nonlocal perturbations. Returning to the example of $\gamma_1^\dagger\gamma_{2L}$, while Q transforms $|\psi_{1,2}\rangle$ into $|\psi_{2,2}\rangle$, $\sigma_1^\dagger\sigma_L|\psi_{\alpha,2}\rangle$ is orthogonal to the ground-state space. More generally, any *odd* power of σ_j (for any j) takes us out of the ground-state space. This follows from $\langle\psi_{\alpha,2}|\sigma_j^a|\psi_{\beta,2}\rangle = \delta_{\alpha\beta}\delta_{a,0 \bmod 2}$, which in turn is deducible from Eq. (39). This motivates us to try an interedge coupling with an even power of σ_j : $\tilde{\mathcal{O}} = \gamma_1^2\gamma_{2L}^2 \sim \sigma_1^2\sigma_L^2 Q^2$. However, any *even* power of Q acts trivially as $\langle\psi_{\alpha,n}|Q^2|\psi_{\beta,n}\rangle = \delta_{\alpha\beta}$, thus, $\tilde{\mathcal{O}}$ cannot distinguish between $|\psi_{1,2}\rangle$ and $|\psi_{2,2}\rangle$. Finding a parafermion-local operator that breaks the ground-state degeneracy turns out to be impossible, as we now show. Any interedge coupling may be decomposed into the form $B_l B_r Q^{[lB_r]}$, where B_l (resp. B_r) is a clock-local operator with definite charge, and with support near site 1 (resp. L). The Jordan-Wigner transformation (5) implies that there are as many powers of the string Q as the charge of B_r .

To transform $|\psi_{1,2}\rangle$ into $|\psi_{2,2}\rangle$, one needs an odd power of the string operator Q ; however, an odd power of Q always accompanies an odd power of σ_j (in B_r), which takes us out of the ground-state space. This intuitive observation can be supplemented with quasiadiabatic continuation techniques to prove that the twofold degeneracy is indeed robust. See Appendix H, where we also prove a stronger statement: that the two ground states are mutually indistinguishable by all symmetric, parafermion-local operators, of which the closed-chain Hamiltonian is but one example.

We now extend our discussion to more general (N,n) phases with a parafermionic order parameter, possibly in conjunction with topological order. Let us decompose the ground-state space $P[H_{op}]$ of an open-chain Hamiltonian into eigenstates of the \mathbb{Z}_N generator: each of $|\phi_{\alpha \in \mathbb{Z}_m, n}^s\rangle$ has a conserved charge $n\alpha$. For $g > 1$, the parafermionic order parameter $\gamma_j^{N/g}$ (or its dressed version \mathcal{Y}_j) permutes the ground states as $|\phi_{\alpha, n}^s\rangle \rightarrow |\phi_{\alpha - m/g, n}^s\rangle$, as evidenced from $\gamma_j^{N/g} Q = \omega^{-N/g} Q \gamma_j^{N/g}$. Thus, $P[H_{op}]$ divides into m/g number of symmetry-broken subspaces, which we label by $\delta \in \mathbb{Z}_{m/g}$ and denote as

$$\bar{P}_\delta = \sum_{\alpha=0}^{g-1} |\phi_{\alpha m/g + \delta, n}^s\rangle \langle \phi_{\alpha m/g + \delta, n}^s|. \quad (48)$$

We propose that each of \bar{P}_δ is indistinguishable under any symmetric, *parafermion*-local probe $\tilde{\mathcal{O}}$, i.e., in the thermodynamic limit,

$$\bar{P}_\delta \tilde{\mathcal{O}} \bar{P}_\delta = \bar{c}(\tilde{\mathcal{O}}) \bar{P}_\delta \quad (49)$$

for a complex number \bar{c} . The proof of Eq. (49) generalizes our intuition from the $H_{(4,2)}$ model, and may be found in Appendix H. We note that Eq. (49) is a stronger statement than the local-indistinguishability condition (10), which applies *only* to symmetric, *clock*-local probes of $P[H_{op}]$. By “stronger” indistinguishability we mean that a certain ground-state space (\bar{P}_δ) cannot be distinguished by both clock-local and -nonlocal operators (which are symmetric and parafermion local). Suppose we now close the chain symmetrically; any legal term in the closed-chain parafermion Hamiltonian satisfies the same conditions on $\tilde{\mathcal{O}}$. Since each of \bar{P}_δ satisfies the strong-indistinguishability condition, but their direct sum $P[H_{op}] = \sum_{\delta=0}^{m/g-1} \bar{P}_\delta$ does not, the open-chain ground state splits into multiplets. A multiplet indexed by δ is robustly degenerate, so long as the gap persists within each of the g charge sectors of \bar{P}_δ .

VII. DISCUSSION

Topological order is now known to characterize the low-energy subspace for a variety of condensed-matter systems [28,29,43–46]. In this paper, we have precisely characterized the ground state of 1D parafermionic chains, and identified the properties which are robustly associated with topological order and/or symmetry breaking. Our work generalizes a previous notion of topological order that imposed constraints on the high-energy states [2]. These constraints cannot be realized in nature since parafermions only emerge as low-energy quasiparticles [47–49].

A unifying property of topologically ordered ground states is their indistinguishability by local probes. From this fundamental property, one may derive [50] (i) a nontrivial ground-state degeneracy depending on the topology of the manifold, and (ii) well-known signatures in the entanglement entropy [43,44]. Aside from parafermions, a close variant [41,51] of Eq. (10) applies to other models with topological order, including the toric code [8] and Levin-Wen string-net models [52]; a stronger version of Eq. (10) is known to stabilize the spectral gap of frustration-free Hamiltonians under generic local perturbations, and produces an area law for the entanglement entropy [53].

A few works [32,54] have demonstrated that Majorana edge modes may exist with weaker localization properties in particle-number-conserving superconductors, where the $U(1)$ symmetry (associated with the electron charge) is not spontaneously broken to \mathbb{Z}_2 . For one particular soluble model, it has been shown that the ground state of the topological phase switches parity when the boundary condition is changed [54]; it is interesting to determine if this property is more generally true away from the soluble limit, perhaps with similar techniques that are presented in this paper.

In the final stages of this work, the phase diagram for our \mathbb{Z}_3 model has been alternatively derived from an entanglement perspective in Ref. [55]. The comparison is favorable.

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APPENDIX A: LOCAL OPERATORS AND QUASIADIABATIC CONTINUATION

To define a local operator (\mathcal{O}), we first decompose it as

$$\mathcal{O} = \sum_{r \geq 1} \sum_{A \in \mathcal{S}(r)} V_{r,A}, \quad (\text{A1})$$

where $V_{r,A}$ acts nontrivially on the chain A of length r , and $\mathcal{S}(r)$ is the set of all chains with length r . For a chain of length L , \mathcal{O} acts in the Hilbert space $\mathcal{Z} = (\mathbb{C}^N)^{\otimes L}$. The operator norm $\|\mathcal{L}\|_{\text{op}}$ of an operator \mathcal{L} is defined by $\sup\{\|\mathcal{L}v\| : v \in \mathcal{Z} \text{ with } \|v\| = 1\}$. We call \mathcal{O} quasilocal if it can be characterized by strength K and decay $f(r)$, such that the operator norm

$\|V_{r,A}\|_{\text{op}} \leq Kf(r)$ for all $r \geq 1$ and for any chain (A) of length r ; $f : \mathbb{Z}_+ \mapsto [0,1]$ and decays superpolynomially in r , i.e., faster than any power. Included are finite-ranged and exponentially decaying interactions. For example, the strength of the finite-ranged Hamiltonian (2) is J . In this example and the rest of the paper, we consider only quasilocal Hamiltonians, which are characterized by a finite Lieb-Robinson velocity [56], i.e., information effectively propagates at finite speed, in analogy with the speed of light. This implies that any finite-time evolution by a quasilocal Hamiltonian preserves the locality of an operator [40]. We now elaborate on an important example of such an evolution: a quasiadiabatic continuation. For illustration, let us symmetrically deform $H_{(N,1)}$ of Eq. (2), while preserving the spectral gap above the lowest N states (which we do not assume to be degenerate). We assume there exists a family of symmetry-preserving Hamiltonians H_s , which are differentiable in s for $s \in [0,1]$, and $H_0 \equiv H_{(N,1)}$. The spectral gap allows us to uniquely define P_s as the projection into the lowest N states of H_s . We define an exact quasiadiabatic continuation \mathcal{V}_s that maps $P_s = \mathcal{V}_s P_0 \mathcal{V}_s^\dagger$ [39]; \mathcal{V}_s is the unitary transformation (16) with the Hermitian Hamiltonian (17). The infinite-time evolution in Eq. (17) is filtered by a function $F(t)$ that decays faster than any power, for large $|t|$; an explicit construction of $F(t)$ is provided in the appendixes of Refs. [40,41]. We assume for simplicity that the interactions in H_s decay exponentially, or faster. It follows from Lemma 2 of Ref. [41] that \mathcal{D}_s is a quasilocal Hamiltonian. By interpreting $s \in [0,1]$ as a time variable, Lemma 1 of Ref. [41] informs us that the finite-time evolution \mathcal{V}_s generated by \mathcal{D}_s preserves locality, i.e., $\mathcal{V}_s^\dagger \mathcal{O} \mathcal{V}_s$ is a quasilocal if \mathcal{O} is quasilocal. Since $[H_s, Q] = 0$, it follows that the quasiadiabatic continuation is symmetry preserving, i.e., $[\mathcal{D}_s, Q] = [\mathcal{V}_s, Q] = 0$.

It is useful to distinguish between two notions of locality on a closed chain. We say that an operator is clock local (parafermion local) if it has strength J and superpolynomial decay $f(r)$ in the clock (parafermion) representation. We describe parafermions (clocks) only with a parafermion-local (clock-local) Hamiltonian. We also insist that all Hamiltonian terms are charge neutral, by which we mean they commute with the \mathbb{Z}_N generator. On an open chain, we now demonstrate that all parafermion-local terms are also clock local, due to the charge-neutral constraint. Decomposing a generic parafermion-local term as in Eq. (A1), we consider each $V_{r,A}$ separately:

$$V_{r,A} \propto \gamma_i^{n_i} \gamma_{i+1}^{n_{i+1}} \cdots \gamma_{i+r-1}^{n_{i+r-1}}, \quad (\text{A2})$$

where $n_j \in \mathbb{Z}_N$, and by assumption of an open chain, $i \geq 1$ and $i+r-1 \leq L$. Let us express this term in the clock representation through Eq. (5). Loosely speaking, this transformation introduces strings of τ_j for $j < i$, assuming $i > 1$. However, if $V_{r,A}$ is charge neutral, the strings all cancel out. More precisely, we operate on each site labeled by $j < i$ as $\tau_j^{\bar{n}}$ with $\bar{n} = \sum_{k=i}^{i+r-1} n_k$. However,

$$Q^\dagger V_{r,A} Q = \omega^{\bar{n}} V_{r,A}, \quad (\text{A3})$$

and charge neutrality imposes that $\bar{n} = 0 \pmod N$.

The discussion thus far applies to an open chain. On a closed chain, charge-neutral interedge couplings can be either

(i) parafermion local but clock nonlocal [e.g., Eq. (25)], or (ii) clock local but parafermion nonlocal (e.g., $\sigma_1^\dagger \sigma_L$). Refining our notion of locality thus distinguishes between parafermions on a ring and clocks on a ring.

APPENDIX B: INDISTINGUISHABILITY BY GENERAL QUASILocal PROBES

In Sec. III, we claimed that (a) $P_{(N,1)}$ satisfies the local-indistinguishability condition (10), for any \mathbb{Z}_N -symmetry-preserving, clock-local operator \mathcal{O} . More precisely, \mathcal{O} has strength K and superpolynomial decay $f(r)$ in the clock representation, as we describe in Appendix A. Later in Sec. VIA, we also claimed that (b) $P_{(N,n)}$ is locally indistinguishable. The proofs of both statements can be combined. In what follows, n is a divisor of N , and $m = N/n$. It suffices to show that the matrix $\mathcal{M}_{\alpha\beta}^{(n)} = \langle \phi_{\alpha,n} | \mathcal{O} | \phi_{\beta,n} \rangle$ is proportional to the identity, where $|\phi_{\alpha,n}\rangle \in P[H_{(N,n)}]$ is defined in Eq. (40). The off-diagonal elements vanish because $[\mathcal{O}, Q] = 0$ by assumption. The difference in diagonal elements is

$$\begin{aligned} \mathcal{M}_{\alpha\alpha}^{(n)} - \mathcal{M}_{\beta\beta}^{(n)} &= \sum_{\mu, \nu \in \mathbb{Z}_m; \mu \neq \nu} \frac{\omega^{\alpha n(\mu-\nu)} - \omega^{\beta n(\mu-\nu)}}{m} \langle \psi_{\mu,n} | \mathcal{O} | \psi_{\nu,n} \rangle. \end{aligned} \quad (\text{B1})$$

We claim that

$$\mathcal{M}_{\alpha\alpha}^{(n)} - \mathcal{M}_{\beta\beta}^{(n)} \leq 2(m-1)Kf(L) \quad (\text{B2})$$

for a chain of length L .

Proof. For \mathcal{O} to have nonvanishing matrix elements between $\langle \psi_{\mu,n} | \mathcal{O} | \psi_{\nu,n} \rangle$ for $\mu \neq \nu$, we require that \mathcal{O} flips the clock variable on all L sites. Thus, if we decompose \mathcal{O} as in Appendix A, we find $\langle \psi_{\mu,n} | \mathcal{O} | \psi_{\nu,n} \rangle = \langle \psi_{\mu,n} | V_{L,A} | \psi_{\nu,n} \rangle$, where $V_{L,A}$ acts nontrivially on the entire chain. By the Cauchy-Schwartz inequality,

$$\begin{aligned} |\langle \psi_{\mu,n} | V_{L,A} | \psi_{\nu,n} \rangle| &\leq \| |\psi_{\mu,n}\rangle \| \cdot \| V_{L,A} | \psi_{\nu,n} \rangle \| \\ &= \| V_{L,A} | \psi_{\nu,n} \rangle \|. \end{aligned} \quad (\text{B3})$$

Now applying the definitions of the operator norm and quasilocality (cf. Appendix A),

$$\| V_{L,A} | \psi_{\nu,n} \rangle \| \leq \| V_{L,A} \|_{\text{op}} \leq K f(L). \quad (\text{B4})$$

It follows from this and Eq. (B1) that

$$\begin{aligned} &|\langle \phi_{\alpha,n} | \mathcal{O} | \phi_{\alpha,n} \rangle - \langle \phi_{\beta,n} | \mathcal{O} | \phi_{\beta,n} \rangle| \\ &\leq \frac{2}{m} \sum_{\mu, \nu \in \mathbb{Z}_m; \mu \neq \nu} |\langle \psi_{\mu,n} | \mathcal{O} | \psi_{\nu,n} \rangle| \leq 2(m-1)K f(L). \end{aligned} \quad (\text{B5})$$

APPENDIX C: GENERALIZED EDGE MODE FOR THE \mathbb{Z}_3 PARAFERMION MODEL

Let us evaluate the zero edge mode \mathcal{O}_l for the \mathbb{Z}_3 model (9), which we rewrite as $H_s = H_0 + sV$, with

$$\begin{aligned} H_0 &= -\hat{\lambda} \sum_{j=1}^{L-1} \omega \gamma_{2j+1}^\dagger \gamma_{2j} + \text{H.c.}, \quad \hat{\lambda} = e^{i\hat{\phi}}, \\ \text{and } V &= -\sum_{j=1}^L \omega^* \gamma_{2j-1}^\dagger \gamma_{2j} + \text{H.c.} \end{aligned} \quad (\text{C1})$$

We need certain relations introduced in Sec. IV, including the definitions of the quasiadiabatic continuation operator \mathcal{V}_s in Eq. (16), which is generated by the Hamiltonian \mathcal{D}_s in Eq. (17). The form of \mathcal{O}_l in Eq. (20) motivates us to evaluate the operator $i[\mathcal{D}_0, \gamma_1]$. Employing $[H_0, \gamma_1] = 0$ and $[V, \gamma_1] = (\omega^* - \omega)(\gamma_2 - \omega \gamma_1^\dagger \gamma_2^\dagger)$, we are led to

$$i[\mathcal{D}_0, \gamma_1] = (\omega^* - \omega)(X - \omega \gamma_1^\dagger Y) \quad (\text{C2})$$

$$\text{with } X = \int dt F(3t) e^{iH_0 t} \gamma_2 e^{-iH_0 t}$$

$$\text{and } Y = \int dt F(3t) e^{iH_0 t} \gamma_2^\dagger e^{-iH_0 t}. \quad (\text{C3})$$

Here, we have applied that the spectral gap Γ^0 (above the lowest three eigenstates of H_0) equals 3. The problem is reduced by noting $Y = -X^\dagger$, as follows from F being imaginary. Further progress is made by realizing that γ_2, γ_3 , and $\omega \gamma_2^\dagger \gamma_3^\dagger$ form a closed linear algebra under commutation by H_0 . That is, the coefficients defined by

$$[H_0, A_0 \gamma_2 + B_0 \gamma_3 + C_0 \omega \gamma_2^\dagger \gamma_3^\dagger] = A_1 \gamma_2 + B_1 \gamma_3 + C_1 \omega \gamma_2^\dagger \gamma_3^\dagger \quad (\text{C4})$$

are linearly related as

$$\begin{pmatrix} A_1 \\ B_1 \\ C_1 \end{pmatrix} = 3\mathcal{M} \begin{pmatrix} A_0 \\ B_0 \\ C_0 \end{pmatrix} \text{ with Hermitian matrix } \mathcal{M} = \frac{1}{i\sqrt{3}} \begin{pmatrix} 0 & -\hat{\lambda} & \hat{\lambda}^* \\ \hat{\lambda}^* & 0 & -\hat{\lambda} \\ -\hat{\lambda} & \hat{\lambda}^* & 0 \end{pmatrix}. \quad (\text{C5})$$

Diagonalizing \mathcal{M} , we find eigenoperators

$$v_j = \frac{1}{\sqrt{3}} (\gamma_2 + \omega^j \gamma_3 + \omega^{2j+1} \gamma_2^\dagger \gamma_3^\dagger), \text{ with corresponding eigenvalues } \zeta_j(\hat{\phi}) = -\frac{2}{\sqrt{3}} \sin\left(\hat{\phi} + \frac{2\pi}{3}j\right), \quad (\text{C6})$$

such that $[H_0, v_j] = 3\zeta_j v_j$. Expanding γ_2 in this eigenbasis as $\gamma_2 = (v_0 + v_1 + v_2)/\sqrt{3}$,

$$X = \frac{1}{\sqrt{3}} \int dt F(3t) \sum_{j=0}^2 v_j e^{i3\zeta_j t} = \frac{1}{3\sqrt{3}} \sum_{j=1}^3 \tilde{F}(\zeta_j) v_j. \quad (\text{C7})$$

To proceed, we thus need the Fourier components \tilde{F} for the frequencies $\pm\zeta_j$. How are they determined? The two essential properties of a quasiadiabatic continuation \mathcal{V}_s is that (a) it evolves the ground-state space P_s as a function of the deformation parameter s , and (b) that it is locality preserving. The first property (a) determines the Fourier components $\tilde{F}(\Omega) = -1/\Omega$ for $|\Omega| \geq 1$ (as shown in Sec. IV), while the second (b) determines the components for $|\Omega| < 1$, up to a degree of arbitrariness. As will shortly be clarified, this arbitrariness lies in the operators which have zero matrix elements within the ground-state space. Now, we address separately the chiral and nonchiral cases: the nonchiral limit $\hat{\phi} = 0$ is discussed in Appendix C 1, and $0 < \hat{\phi} < \pi/3$ is elaborated in Appendix C 2.

1. Nonchiral \mathbb{Z}_3 parafermions: $\hat{\phi} = 0$

As defined in Eq. (C6), the eigenvalues of \mathcal{M} are $\zeta_0 = 0$, $\zeta_1 = -1$, and $\zeta_2 = 1$. For the first eigenvalue, the oddness of \tilde{F} implies that $\tilde{F}(\zeta_0) = 0$; for $j \in \{1, 2\}$, $\tilde{F}(\zeta_j) = -1/\zeta_j$ since $|\zeta_j| \geq 1$. Equation (C7) leads to

$$X = \frac{1}{3\sqrt{3}}(v_1 - v_2) = \frac{1}{3(\omega^* - \omega)}(\gamma_3 - \omega\gamma_2^\dagger\gamma_3^\dagger). \quad (\text{C8})$$

Inserting the expressions of X and $Y = -X^\dagger$ into Eq. (C2), we thus derive from Eq. (20) the form of \mathcal{O}_l , to first order in the deformation parameter s . This result is found in Eq. (22).

2. Chiral \mathbb{Z}_3 parafermions: $0 < \hat{\phi} < \pi/3$

In this range of $\hat{\phi}$, both $|\zeta_0|$ and $|\zeta_1| < 1$, while $|\zeta_2| > 1$. To determine the edge mode, we thus need the Fourier component $\tilde{F}(\Omega)$ for $0 < |\Omega| < 1$, which enters the expression (C7). In the conventional quasiadiabatic continuation, \tilde{F} is chosen so that \mathcal{V}_s maximally preserves locality, which amounts to finding $F(t)$ with the fastest possible decay, which is proven to be of subexponential type [40,41]. We thus refer to subexponentially decaying edge modes (\mathcal{O}_l) as maximally localized. On the other hand, one may choose an edge mode $\tilde{\mathcal{O}}_l$ with worse localization properties, but with the desirable property that it commutes with the Hamiltonian. There is thus a tradeoff between localizability and commutativity. Indeed, choosing $\tilde{F}(\Omega) = -1/\Omega$ for $|\Omega| = |\zeta_i| < 1$ worsens the decay of $F(t)$ (and thus of the resultant edge mode), but leads to an edge mode which commutes with the Hamiltonian, as we elaborate in Appendix D. One may verify that the difference $\Delta = \tilde{\mathcal{O}}_l - \mathcal{O}_l$ has zero matrix elements in the ground-state space. To illustrate this for the \mathbb{Z}_3 chiral example, we have shown that the first-order correction to the edge mode has a term proportional to $\sum_{j=1}^3 \tilde{F}(\zeta_j)v_j$; this follows from Eqs. (C2) and (C7). Since $|\zeta_0|, |\zeta_1| < 1$, the coefficients of v_0, v_1 [as defined in Eq. (C6)] are indeterminate in the manner that we have been discussing. This lack of determinacy has no physical consequence in the ground-state space, as $v_j|\psi_{\alpha,1}\rangle = 0$ for $j \in \{0, 1\}$ and $\alpha \in \mathbb{Z}_3$; here, recall ψ are the ground states of H_0 , as defined in Sec. III.

APPENDIX D: ZERO EDGE MODES THAT COMMUTE WITH THE HAMILTONIAN

Let us define the first-order truncation of \mathcal{O}_l as $\mathcal{O}_l^{(1)} = \gamma_1 + is[\mathcal{D}_0, \gamma_1]$, where s is the deformation parameter. This operator depends on our choice of the filter function F , which enters the expression for \mathcal{D}_s in Eq. (17). For a specific choice of F (with corresponding edge mode $\tilde{\mathcal{O}}_l^{(1)}$), we find that $\tilde{\mathcal{O}}_l^{(1)}$ and H_s commute to first order in s . The following discussion clarifies this choice of F . Let v_j be operators which satisfy two properties: (i) they belong in the charge sector α , i.e., $Q^\dagger v_j Q = \omega^\alpha v_j$, and (ii) they are constants under time evolution by H_0 , i.e., $[H_0, v_j] = 2\zeta_j v_j$ with real eigenvalues $2\zeta_j$. These operators v_j span the space (\mathcal{X}_α) of operators in the same charge sector, i.e., for each element $x \in \mathcal{X}_\alpha$, $Q^\dagger x Q = \omega^\alpha x$. Labeling each element by a subscript (e.g., x_i), commutation by H_0 may be interpreted as an eigenvalue problem: $[H_0, x_i] = \mathcal{B}_{ij}x_j$. The matrix \mathcal{B} is known to be Hermitian in the frustration-free limit [2] where $H_0 = H_{N,1}$, as described in Eq. (2). Applying the completeness property, we expand $[V, \gamma_1] = \sum_n r_n v_n$, with c numbers r_n . This leads to

$$\begin{aligned} i[\mathcal{D}_0, \gamma_1] &= \sum_n r_n \int dt F(\Gamma_0 t) e^{iH_0 t} v_n e^{-iH_0 t} \\ &= \sum_n \frac{r_n v_n}{\Gamma_0} \tilde{F}(2\zeta_n / \Gamma_0). \end{aligned} \quad (\text{D1})$$

If the Fourier transform $\tilde{F}(\Omega)$ is chosen as $-1/\Omega$ for all relevant frequencies $\Omega \in \{2\zeta_n / \Gamma_0\}$ in the above sum, we denote the resultant edge mode by $\tilde{\mathcal{O}}_l$. This choice of \tilde{F} , which we call Fendley's choice, is only possible when none of the relevant frequencies are vanishing, i.e., $\zeta_n \neq 0$ for any nonzero r_n . The proof of commutativity to first order is elementary:

$$\begin{aligned} [H_s, \tilde{\mathcal{O}}_l^{(1)}] &= s[V, \gamma_1] + s[H_0, i[\mathcal{D}_0, \gamma_1]] + O(s^2) \\ &= s[V, \gamma_1] + s \sum_n \frac{r_n [H_0, v_n]}{\Gamma_0} \tilde{F}(2\zeta_n / \Gamma_0) + O(s^2) \\ &= s[V, \gamma_1] + s \sum_n \frac{2r_n \zeta_n v_n}{\Gamma_0} \left(-\frac{\Gamma_0}{2\zeta_n} \right) + O(s^2) = O(s^2). \end{aligned} \quad (\text{D2})$$

In cases where all the relevant frequencies are large, by which we mean they satisfy $|2\zeta_n / \Gamma_0| \geq 1$, Fendley's choice is not so much a choice but a necessity. Indeed, Fendley's choice then coincides with Eq. (18) in Sec. IV; we have shown therein that Eq. (18) leads to the correct evolution of the ground-state space: $\partial_s P_s = i[\mathcal{D}_s, P_s]$. The set of cases where all the relevant frequencies are large include all quadratic Majorana models, as we show in Appendix D 1. On the other hand, if some of the relevant frequencies satisfy $|2\zeta_n / \Gamma_0| < 1$, Fendley's choice produces an edge mode that is deformed from the maximally localized edge mode. In Appendix D 2, we carry out this deformation for the \mathbb{Z}_3 chiral parafermion.

1. Zero edge mode for quadratic Majorana models

Suppose we deform the topological dimer model described in Eq. (15). If we limit our deformations to terms which are quadratic in the Majorana operators, we can derive a general expression for the zero edge mode in the first-order approximation. To begin, let us express the dimer model (H_0) and the deformation (V) as

$$H_0 = \frac{i}{2} \gamma_i A_{ij} \gamma_j, \quad V = \frac{i}{2} \gamma_i B_{ij} \gamma_j, \quad (\text{D3})$$

with A and B real, skew-symmetric, and even-dimensional matrices. It follows that the eigenvalues of A (and B) are completely imaginary, and come in complex-conjugate pairs. The eigenproblem $[H_0, \gamma_j] = -2i \sum_m A_{jm} \gamma_m$ thus has real eigenvalues ζ_n : $[H_0, v_n] = 2\zeta_n v_n$. The eigenvectors v_n form a complete basis for all linear Majorana operators, thus, we may define the expansion $\gamma_j = \sum_l z_{jl} v_l$; alternatively, z_{jl} may be obtained from diagonalizing A . The eigenvalues $\{\zeta_j\}$ of $-iA$ are interpreted as half the single-particle energies of H_0 , i.e., H_0 has single-particle excitations with energy $2\zeta_j$. Since H_0 is the topological dimer model described in Eq. (15), we know that A has two zero modes (denoted $\zeta_1 = \zeta_2 = 0$), corresponding to the edge operators $v_1 = \gamma_1$ and $v_2 = \gamma_{2L}$. All other single-particle excitations are gapped, i.e., $|2\zeta_j| \geq \Gamma_0$ for $j > 2$, where Γ_0 is the many-body spectral gap of H_0 . The edge mode is derived as

$$\begin{aligned} i[\mathcal{D}_0, \gamma_1] &= \int dt F(\Gamma_0 t) e^{iH_0 t} [V, \gamma_1] e^{-iH_0 t} \\ &= -2i \sum_{m \neq 1, 2L} B_{1m} \int dt F(\Gamma_0 t) e^{iH_0 t} \gamma_m e^{-iH_0 t} \end{aligned}$$

$$\bar{X} = -\frac{1}{3\sqrt{3}} \sum_{j=0}^2 \frac{v_j}{\zeta_j} = -\frac{1}{9} \left[\left(\frac{1}{\zeta_0} + \frac{1}{\zeta_1} + \frac{1}{\zeta_2} \right) \gamma_2 + \left(\frac{1}{\zeta_0} + \frac{\omega}{\zeta_1} + \frac{\omega^*}{\zeta_2} \right) \gamma_3 + \left(\frac{1}{\zeta_0} + \frac{\omega^*}{\zeta_1} + \frac{\omega}{\zeta_2} \right) \omega \gamma_2^\dagger \gamma_3^\dagger \right]. \quad (\text{D6})$$

Here, we have introduced \bar{X} (to distinguish Fendley's choice) in place of X (which we reserve for the conventional quasiadiabatic continuation). A useful relation is $\zeta_1 \zeta_2 \zeta_3 = 2 \sin(3\hat{\phi})/(3\sqrt{3})$, which is derived from taking the determinant of \mathcal{M} as defined in Eq. (C5). This relation, in combination with elementary trigonometric identities, leads to

$$(\omega^* - \omega) \bar{X} = -\frac{i}{2 \sin(3\phi)} (\gamma_2 + \hat{\lambda}^2 \gamma_3 + \hat{\lambda}^{-2} \omega \gamma_2^\dagger \gamma_3^\dagger). \quad (\text{D7})$$

The expressions for \bar{X} and $\bar{Y} = -\bar{X}^\dagger$, in combination with Eqs. (C2) and (20), lead to the final result

$$\begin{aligned} \bar{\mathcal{O}}_l &= \gamma_1 - \frac{is}{2 \sin(3\phi)} [\gamma_2 + \hat{\lambda}^2 \gamma_3 + \hat{\lambda}^{-2} \omega \gamma_2^\dagger \gamma_3^\dagger \\ &\quad + \omega \gamma_1^\dagger (\gamma_2^\dagger + \hat{\lambda}^{-2} \gamma_3^\dagger + \hat{\lambda}^2 \omega \gamma_2 \gamma_3)] + \dots \end{aligned} \quad (\text{D8})$$

This deformed edge mode $\bar{\mathcal{O}}_l$ is identical to the expression Ψ_{left} alternatively derived in Ref. [2], up to minor typographic errors in that reference. At least to first order in s , Fendley's edge mode can thus be understood as a deformed quasiadiabatic continuation of γ_1 ; the resultant operator is not optimally localized. This deformation cannot always be justified: as $\hat{\phi}$ (and ζ_0)

$$\begin{aligned} &= -2i \sum_{m \neq 1, 2L} \sum_{n \neq 1, 2} B_{1m} z_{mn} v_n \int dt F(\Gamma_0 t) e^{i2\zeta_n t} \\ &= -2i \Gamma_0^{-1} \sum_{m \neq 1, 2L} \sum_{n \neq 1, 2} B_{1m} z_{mn} v_n \tilde{F}(2\zeta_n / \Gamma_0). \end{aligned} \quad (\text{D4})$$

Since V is skew symmetric, the sum over m excludes 1; this sum also excludes $2L$ since γ_{2L} is a zero mode and F an odd function (recall Sec. IV). Since H_0 does not contain a term proportional to either γ_1 or γ_{2L} , it follows that the expansion $\gamma_{m \neq 1, 2L} = \sum_{n \neq 1, 2} z_{mn} v_n$ excludes the zero modes $v_1 = \gamma_1$ and $v_2 = \gamma_{2L}$. Alternatively stated, all the relevant frequencies are gapped, i.e., $|2\zeta_n / \Gamma_0| \geq 1$. Now applying that $\tilde{F}(\Omega) = -1/\Omega$ for these relevant frequencies, we find

$$i[\mathcal{D}_0, \gamma_1] = i \sum_{m \neq 1, 2L} \sum_{n \neq 1, 2} B_{1m} z_{mn} \frac{v_n}{\zeta_n}. \quad (\text{D5})$$

Applying this to the quadratic Majorana model (15), we obtain $i[\mathcal{D}_0, \gamma_1] = \gamma_3$, as we have previously derived in less-general fashion (cf. Sec. IV).

2. Deformed zero edge mode for the \mathbb{Z}_3 chiral parafermion

For illustration, we consider the \mathbb{Z}_3 chiral parafermion with $0 < \hat{\phi} < \pi/3$. Recall that the filter function has not been specified in Eq. (C7). If we choose $\tilde{F}(\zeta_j) = -1/\zeta_j$, we alternately derive Fendley's edge mode. Substituting this choice into Eq. (C7) and applying Eq. (C6), we obtain

tends to zero, the localization of $\bar{\mathcal{O}}_l$ worsens dramatically, as evidenced by the diverging first-order coefficient in Eq. (D8), and the increasingly singular \tilde{F} . In contrast, the conventional quasiadiabatic continuation chooses a filter function that is infinitely differentiable [40]; the resultant edge mode is well defined for $\hat{\phi} = 0$, as shown in Appendix C 1.

APPENDIX E: CLOSING THE CHAIN BY IMPOSING TRANSLATIONAL INVARIANCE

Our goal is to map a parafermion Hamiltonian (H_{op}) on an open chain of L sites to a parafermion Hamiltonian ($\mathcal{K}^{(\omega)}$) on a closed ring. For simplicity, we assume that the open-chain Hamiltonian is translational invariant up to edge corrections, i.e., we can decompose $H_{op} = \sum_{r \geq 1} \sum_{j=1}^{L-r+1} V_{j,r}$, where r is the range of the operator $V_{j,r}$, i.e., $V_{j,r}$ acts nontrivially on sites $\{j, j+1, \dots, j+r-1\}$. Furthermore, we assume $V_{j,r}$ is related to $V_{j+1,r}$ by translation. For a \mathbb{Z}_3 example, consider $H_{op} = -\sum_{j=1}^{L-1} \omega \gamma_{2j+1}^\dagger \gamma_{2j} + \text{H.c.}$, where $V_{j,2} = -\omega \gamma_{2j+1}^\dagger \gamma_{2j} + \text{H.c.}$ It is surprising that for $N > 2$, a \mathbb{Z}_N Hamiltonian on a closed chain cannot be uniquely defined by the identification $\gamma_{2L+j} = \gamma_j$, due to the nontrivial

commutation (1) [31]. In our example, we might have tried to extend H_{op} by translational symmetry: $H_{op} \rightarrow H_{op} + V_{L,2}$, then identified $\gamma_{2L+1} \equiv \gamma_1$ directly. The result would be a closed-chain Hamiltonian

$$\tilde{\mathcal{K}}^{(\alpha)} = -J \sum_{j=1}^{L-1} \omega \gamma_{2j+1}^\dagger \gamma_{2j} - J \omega^{\alpha-1} \gamma_1^\dagger \gamma_{2L} + \text{H.c.} \quad (\text{E1})$$

with $\alpha = 2$. On the other hand, first interchanging $\gamma_{2L+1}^\dagger \gamma_{2L} = \omega \gamma_{2L}^\dagger \gamma_{2L+1}$, and then making the same identification, we arrive at a different Hamiltonian (E1) with $\alpha = 1$. If there exists more than one interedge coupling, the phase of each is ambiguous. These ambiguities motivate us to identify $\sigma_{L+j} = \sigma_j$ instead; since clock operators on different sites commute, the clock Hamiltonian is uniquely defined on a closed chain. Then, by attachment of Jordan-Wigner strings we convert this clock-local Hamiltonian to a parafermion-local Hamiltonian. Let us illustrate this procedure with the family of open-chain Hamiltonians $H_{(N,1)}$ [from Eq. (2)], which we extend and then identify $\sigma_{L+1} = \sigma_1$, to obtain clocks on a ring:

$$\mathcal{G}_{(N,1)} = H_{(N,1)} - J \sum_{\beta=0}^{N-1} (\sigma_L \sigma_1^\dagger)^\beta. \quad (\text{E2})$$

In the next step, we decompose the interedge coupling into a form $B_{cl} = B_l B_r$, such that B_l (resp. B_r) is a clock-local operator with a definite charge, and with support near site 1 (resp. L). To convert B_{cl} to an interedge coupling (B_{pf}) of

parafermions, we must attach as many powers of the string Q as the charge of B_r . The resultant interedge coupling

$$B_{pf}^{(\alpha)} = B_l B_r (\omega^\alpha Q)^{[B_r]} \quad (\text{E3})$$

can always be expressed locally in terms of parafermions, as follows from the Jordan-Wigner transformation (5). Here, in Eq. (E3), $[[B_r]]$ refers to the charge of B_r , i.e., $Q^\dagger B_r Q = \omega^{[B_r]} B_r$, and we have also introduced the twist parameter $\alpha \in \mathbb{Z}_N$. In our example (E2), $[[B_r]] = [[\sigma_L]] = 1$, so we attach $\sigma_L \rightarrow \sigma_L Q \sim \gamma_{2L}$. Then, from $\sigma_L \sigma_1^\dagger Q = \omega^{(1-N)/2} \gamma_1^\dagger \gamma_{2L}$, we obtain parafermions on a ring:

$$\mathcal{K}_{N,1}^{(\alpha)} = H_{(N,1)} - J \sum_{\beta=0}^{N-1} (\omega^{\alpha+(1-N)/2} \gamma_1^\dagger \gamma_{2L})^\beta. \quad (\text{E4})$$

APPENDIX F: DERIVING THE TOPOLOGICAL ORDER PARAMETER IN THE XYZ MODEL

Our aim is to derive the topological order parameter for the XYZ Majorana model (28), which we parametrize by J_x and J_y . There are two parameter regions with distinct order parameters: (i) $|J_x|, |J_y| < 1$, as described in Appendix F 1, and (ii) $|J_y| > 1, |J_x| > |J_y|$, described in Appendix F 2.

1. Topological order parameter for $|J_x|, |J_y| < 1$

Let us parametrize the XYZ model (28) as

$$H_s = H_0 + sV, \quad \text{where } V = \sum_{j=1}^{L-1} (\gamma_{2j-1} \gamma_{2j} \gamma_{2j+1} \gamma_{2j+2} - it \gamma_{2j-1} \gamma_{2j+2}), \quad s = J_x \quad t = \frac{J_y}{J_x}, \quad (\text{F1})$$

and the frustration free H_0 is defined in Eq. (26). H_0 commutes with the zero edge modes γ_1 and γ_{2L} , thus its ground-state space is described by a topological order parameter $\mathcal{Q} = -i \gamma_1 \gamma_{2L}$. If we frustrate this model through V , the spectral gap above the lowest two states remains finite in the parameter region: $|J_x|, |J_y| < 1$. Our task is to derive the dressed order parameter in this region. We will employ the identities

$$\begin{aligned} [V, \gamma_1] &= -2\gamma_2 \gamma_3 \gamma_4 + 2it \gamma_4 \\ &= (-\gamma_2 \gamma_3 + it)(w_+ + w_-), \\ \text{and } [V, \gamma_{2L}] &= (z_+ + z_-)(-it + \gamma_{2L-2} \gamma_{2L-1}), \end{aligned} \quad (\text{F2})$$

where $w_\pm = \gamma_4 \pm i \gamma_5$ and $z_\pm = \gamma_{2L-3} \pm i \gamma_{2L-4}$ are eigenoperators under commutation by H_0 , i.e., $[H_0, w_\pm] = \mp 2w_\pm$ and $[H_0, z_\pm] = \pm 2z_\pm$. It follows that

$$\begin{aligned} i[\mathcal{D}_0, \gamma_1] &= \int dt F(2t) e^{iH_0 t} [V, \gamma_1] e^{-iH_0 t} \\ &= (it - \gamma_2 \gamma_3)(i \gamma_5), \\ i[\mathcal{D}_0, \gamma_{2L}] &= \int dt F(2t) e^{iH_0 t} [V, \gamma_{2L}] e^{-iH_0 t} \\ &= -i \gamma_{2L-4} (-it + \gamma_{2L-2} \gamma_{2L-1}), \end{aligned} \quad (\text{F3})$$

which determine the edge modes \mathcal{O}_l and \mathcal{O}_r to first order in s [see Eq. (20)]. Applying that $\mathcal{Q} = -i \mathcal{O}_l \mathcal{O}_r$, we are immediately led to Eq. (30). In this derivation, we have assumed that $|J_x| < 1, J_x \neq 0$; if instead $J_x = 0$ and $|J_y| < 1, J_y \neq 0$, the final answer (30) is identical.

2. Deriving the topological order parameter for $|J_y| > 1, |J_x| > |J_y|$

Up to a proportionality constant, we parametrize the XYZ model (28) as $H'_s = H'_0 + sV'$, where $s = 1/J_y$,

$$\begin{aligned} H'_0 &= -i \sum_{j=1}^{L-1} \gamma_{2j-1} \gamma_{2j+2} \\ \text{and } V' &= J_x \sum_{j=1}^{L-1} \gamma_{2j-1} \gamma_{2j} \gamma_{2j+1} \gamma_{2j+2} + i \sum_{j=1}^{L-1} \gamma_{2j} \gamma_{2j+1}. \end{aligned} \quad (\text{F4})$$

The frustration free H'_0 commutes with the zero edge modes γ_2 and γ_{2L-1} , thus its ground-state space is described by a topological order parameter $\mathcal{Q}' = -i \gamma_2 \gamma_{2L-1}$, which must be

distinguished from the order parameter ($\mathcal{Q} = -i\gamma_1\gamma_{2L}$) of H_0 . If we frustrate this model through V' , the spectral gap above the lowest two states remain finite in the parameter region: $|J_y| > 1, |J_y| > |J_x|$. In this region, the order parameter is dressed as

$$\begin{aligned} \mathcal{Q}^y = & -i\gamma_2\gamma_{2L-1} + \frac{J_x}{J_y}(\gamma_1\gamma_4\gamma_6\gamma_{2L-1} + \gamma_2\gamma_{2L-5}\gamma_{2L-3}\gamma_{2L}) \\ & + \frac{i}{J_y}(\gamma_6\gamma_{2L-1} + \gamma_2\gamma_{2L-5}) + \dots \end{aligned} \quad (\text{F5})$$

Its derivation straightforwardly generalizes Appendix F 1. Defining $\tilde{\mathcal{Q}}^y$ as the first-order truncation of \mathcal{Q}^y (i.e., dropping the dots), we define a closed-chain Hamiltonian by $H_{XYZ} - (-1)^\alpha J_y \tilde{\mathcal{Q}}^y$, where H_{XYZ} is defined in Eq. (28). We numerically evaluate the ground-state parities of this closed-chain Hamiltonian for both $\alpha \in \mathbb{Z}_2$, on a chain of 22 sites. The result is shown in Fig. 4(d); we have whitened the parameter regions where the ground-state parity switches upon flux insertion. Note that $|J_y| > 1, |J_y| > |J_x|$ corresponds to regions B and D in this figure.

APPENDIX G: FINITE-SIZE ANALYSIS OF THE TRANSLATIONAL-INVARIANT MAJORANA CHAIN

In this appendix, we present a finite-size analysis of the XYZ Majorana model (28) for the parameter region $|J_x|, |J_y| < 1$, which we label as the square C in Fig. 4. Suppose we close the chain with the translational-invariant boundary conditions (TIBC); a detailed description of this procedure is available in Appendix E. We then obtain two Hamiltonians, which we denote by $\mathcal{K}_{XYZ}^{(0)}$ (resp. $\mathcal{K}_{XYZ}^{(1)}$) in the case of periodic (resp. antiperiodic) boundary condition. A signature of topological order is that the groundstate of \mathcal{K}_{XYZ} is sensitive to its boundary condition; in particular, we expect that the ground-state parity switches between $\mathcal{K}_{XYZ}^{(0)}$ and $\mathcal{K}_{XYZ}^{(1)}$ for all parameters within the square C. A finite-size study shows that the parity is invariant for a subregion (colored black) of C where quantum fluctuations are dominant [see Fig. 4(b)]. The aim of this appendix is to explain the effect of finite size on the ground-state sensitivity. Our explanation relies on a Majorana-spin duality which we will first introduce; one motivation for the spin representation is that numerical simulations are often easier with a clock-local basis. (i) To describe this duality, we first relate the open-chain Majorana Hamiltonian (H_{XYZ}) to a closed-chain spin Hamiltonian (\mathcal{G}_{XYZ}) in Appendix G 1. We argue that H_{XYZ} being topologically ordered implies clock-local order in \mathcal{G}_{XYZ} . (ii) Then, in Appendix G 2, we show that the closed-chain Majorana Hamiltonian ($\mathcal{K}_{XYZ}^{(\alpha)}$) is dual to the closed-chain spin Hamiltonian (\mathcal{G}_{XYZ}). Applying this duality, we explain what clock-local order in \mathcal{G}_{XYZ} implies for $\mathcal{K}_{XYZ}^{(\alpha)}$. We then analyze the ground states of $\mathcal{K}_{XYZ}^{(\alpha)}$ and their sensitivity to twisting. (iii) Finally, in Appendix G 3, we describe a finite-size comparative study of the three types of boundary conditions discussed in this paper: (i) TIBC, (ii) zeroth-order topological boundary condition (TBC), and (iii) first-order TBC.

1. Relating topological order in H_{XYZ} to clock-local order in the spin Hamiltonian \mathcal{G}_{XYZ}

As defined in Eq. (28), the open-chain Majorana Hamiltonian can be rewritten as the Heisenberg spin model

$$H_{XYZ} = - \sum_{j=1}^{L-1} \sum_{\mu \in \{x,y,z\}} J_\mu \zeta_j^\mu \zeta_{j+1}^\mu, \quad \text{with } J_z = 1. \quad (\text{G1})$$

Here, ζ_j^μ are Pauli matrices acting on site j . To derive the above spin representation, we identify $\zeta^z = \sigma$, $\zeta^x = \tau$, and apply the Jordan-Wigner transformation (5): for bulk couplings ($1 \leq j < L$),

$$\begin{aligned} \zeta_j^x \zeta_{j+1}^x &= -\gamma_{2j-1} \gamma_{2j} \gamma_{2j+1} \gamma_{2j+2}, \\ \zeta_j^y \zeta_{j+1}^y &= i\gamma_{2j-1} \gamma_{2j+2}, \quad \text{and} \quad \zeta_j^z \zeta_{j+1}^z = -i\gamma_{2j} \gamma_{2j+1}. \end{aligned} \quad (\text{G2})$$

We define \mathcal{G}_{XYZ} as the translational-invariant extension of H_{XYZ} to a ring:

$$\mathcal{G}_{XYZ} = - \sum_{j=1}^L \sum_{\mu \in \{x,y,z\}} J_\mu \zeta_j^\mu \zeta_{j+1}^\mu. \quad (\text{G3})$$

Here, periodic spin boundary conditions are imposed as $\zeta_{L+1}^\mu = \zeta_1^\mu$. The phase diagram of \mathcal{G}_{XYZ} is well understood analytically [57,58]. \mathcal{G}_{XYZ} is gapless where $|J_\mu| = |J_\nu| \geq |J_\kappa|$, i.e., for any two parameters having equal magnitude that is greater than or equal to the third magnitude; otherwise, the phase is gapped with a clock-local order parameter ζ^μ , corresponding to J_μ of the largest magnitude. We illustrate this symmetry breaking in Fig. 7(a), for the parameters $(J_x, J_y, J_z) = (0.1, -0.2, 1)$.

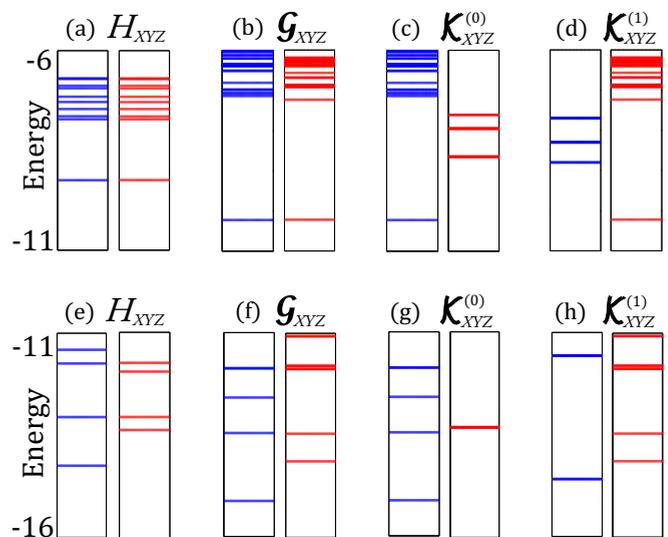


FIG. 7. Lowest-lying spectra of the XYZ model on a chain of 10 sites. Two parametrizations are considered: (a)–(d) correspond to the same parameters $(J_x, J_y, J_z) = (0.1, -0.2, 1)$, while (e)–(h) correspond to $(J_x, J_y, J_z) = (0.75, -0.75, 1)$. The spectrum of each Hamiltonian is split into even (blue) and odd (red) charge sectors. (a), (e) Spectrum of the open chain $H_{XYZ}(J_x, J_y, J_z)$, (b), (f) \mathcal{G}_{XYZ} , (c), (g) $\mathcal{K}_{XYZ}^{(0)}$, (d), (h) $\mathcal{K}_{XYZ}^{(1)}$.

We argue that H_{XYZ} being topologically ordered (in the fermionic language) indicates that \mathcal{G}_{XYZ} is symmetry broken (in the spin language). This implies that H_{XYZ} and \mathcal{G}_{XYZ} simultaneously manifest a degenerate ground-state manifold, as illustrated in Figs. 7(a) and 7(b) for the parameters $(J_x, J_y, J_z) = (0.1, -0.2, 1)$. This follows because \mathcal{G}_{XYZ} is the translational-invariant extension of H_{XYZ} , thus, both Hamiltonians are approximately minimized by the same wave functions. Alternatively stated, we expect that the spectral gaps of H_{XYZ} and \mathcal{G}_{XYZ} (above their two lowest-lying states) are related multiplicatively: a quantum phase transition in H_{XYZ} indicates a simultaneous transition in \mathcal{G}_{XYZ} . To support our hypothesis, we numerically evaluate the spectral gap of H_{XYZ} in Fig. 4(a); the gapless lines (colored black) separate the plot into five regions labeled by A to E, with B, C, and D corresponding to topological phases (cf. Sec. V). Our hypothesis is supported by the exact coincidence of gapless lines in both Hamiltonians (H_{XYZ} and \mathcal{G}_{XYZ}). Here, we are comparing our numerical simulation of H_{XYZ} with a known analytical result about \mathcal{G}_{XYZ} ; cf. the previous paragraph.

2. Majorana-spin duality on a closed chain

We extend the open chain H_{XYZ} onto a ring through the prescription (E3) to obtain the closed-chain Hamiltonian $\mathcal{K}_{XYZ}^{(\alpha)} = H_{XYZ} + \Delta H_{XYZ}^{(\alpha)}$, as shown in Eq. (35). The interedge

$$|\phi_{0,1}^{J_x, J_y}\rangle \text{ minimizes } [\mathcal{K}_{XYZ}^{(0)}]_0 \quad \text{and} \quad |\phi_{1,1}^{J_x, J_y}\rangle \text{ minimizes } [\mathcal{K}_{XYZ}^{(1)}]_1. \quad (\text{G7})$$

For $(J_x, J_y, J_z) = (0.1, -0.2, 1)$,

$$|\phi_{0,1}^{J_x, J_y}\rangle \text{ also minimizes } \mathcal{K}_{XYZ}^{(0)} \quad \text{and} \quad |\phi_{1,1}^{J_x, J_y}\rangle \text{ also minimizes } \mathcal{K}_{XYZ}^{(1)}. \quad (\text{G8})$$

This is illustrated in the spectra of Figs. 7(b)–7(d), and explains why the ground-state parity switches as we twist $\mathcal{K}_{XYZ}^{(\alpha)}$.

Due to the duality (G6), the first statement (G7) is exactly satisfied for any finite size; the second statement (G8) is not necessarily satisfied on finite-size chains, i.e., the actual ground state of $\mathcal{K}_{XYZ}^{(\alpha)}$ need not originate from the symmetry-broken manifold of \mathcal{G}_{XYZ} . For example, we consider the spectra at $(J_x, J_y, J_z) = (0.75, -0.75, 1)$ for a 10-site chain, as plotted in Figs. 7(e)–7(h). Figure 7(f) demonstrates the finite-size splitting of the low-energy subspace of \mathcal{G}_{XYZ} ; in particular, $|\phi_{0,1}^{3/4, -3/4}\rangle$ is energetically favored in comparison with $|\phi_{1,1}^{3/4, -3/4}\rangle$. While $|\phi_{0,1}^{3/4, -3/4}\rangle$ also minimizes $\mathcal{K}_{XYZ}^{(0)}$ [see Fig. 7(g)], $|\phi_{1,1}^{3/4, -3/4}\rangle$ does not minimize $\mathcal{K}_{XYZ}^{(1)}$ [see Fig. 7(h)]. A corollary is that the parity does not switch, and we have verified that this behavior persists for a chain of 30 sites.

As a side remark, we note that the duality described in Eq. (G6) is easily generalized to a parafermion-clock duality: if $\mathcal{K}^{(\alpha)}$ is the Hamiltonian for parafermions on a ring, with a particular twist $\alpha \in \mathbb{Z}_N$, and \mathcal{G} is the Hamiltonian for clocks on a ring, then

$$[\mathcal{K}^{(N-\alpha)}]_\alpha = [\mathcal{G}]_\alpha \quad \text{with } \alpha \in \mathbb{Z}_N. \quad (\text{G9})$$

Let us exemplify this duality for the frustration-free models $H_{(N,1)}$ defined in Eq. (2). The closed-chain parafermion

coupling can be rewritten as

$$\Delta H_{XYZ}^{(\alpha)} = -J_x S_L^x S_1^x - J_y S_L^y S_1^y (-1)^\alpha Q - J_z S_L^z S_1^z (-1)^\alpha Q. \quad (\text{G4})$$

Let us define for any Hamiltonian H a projected Hamiltonian $[H]_\beta$ in the charge sector β , i.e., $[H]_\beta$ acts in the space of functions with eigenvalue ω^β under conjugation by Q . Clearly,

$$[\Delta H_{XYZ}^{(\alpha)}]_\beta = -J_x S_L^x S_1^x - J_y S_L^y S_1^y (-1)^{\alpha+\beta} - J_z S_L^z S_1^z (-1)^{\alpha+\beta}. \quad (\text{G5})$$

By inspecting the matrix elements of H_{XYZ} , \mathcal{G}_{XYZ} , and $\mathcal{K}_{XYZ}^{(\alpha)}$ in each charge sector, we derive a duality between Majorana fermions and spins on a closed chain:

$$[\mathcal{K}_{XYZ}^{(\alpha)}]_\alpha = [\mathcal{G}_{XYZ}]_\alpha \quad \text{for } \alpha \in \mathbb{Z}_2. \quad (\text{G6})$$

This duality on a closed chain generalizes an open-chain duality [59,60] between the transverse-field Ising model and the Kitaev wire. \mathcal{G}_{XYZ} being translational invariant then also implies the same for $\mathcal{K}_{XYZ}^{(0)}$ (resp. $\mathcal{K}_{XYZ}^{(1)}$) in the even (resp. odd) charge sector, as we alluded to in Sec. VB.

This duality is further illustrated in Figs. 7(b)–7(d) for the parameters $(J_x, J_y, J_z) = (0.1, -0.2, 1)$. If we define $|\phi_{\alpha,1}^{J_x, J_y}\rangle$ as the ground state of \mathcal{G}_{XYZ} in the charge sector $\alpha \in \mathbb{Z}_2$, then clearly

Hamiltonian (E4) is dual to the closed-chain clock Hamiltonian ($\mathcal{G}_{(N,1)}$) in Eq. (E2). That is, $\mathcal{K}_{N,1}^{(\alpha)}$ and $\mathcal{G}_{(N,1)}$ satisfy Eq. (G9), as follows from the identity (25).

3. Comparing the three types of closed-chain boundary conditions

Applying the parity switch as a criterion, we would like to evaluate the various types of closed-chain boundary conditions. We focus on the $J_x = -J_y$ line, for $J_x \in [0, 1)$, and define \bar{J}_x such that the parity switches for $0 \leq J_x < \bar{J}_x$, and is invariant for $\bar{J}_x < J_x < 1$. With the topological boundary condition (TBC) of Eq. (32), the parity switches for all parameters where the open-chain Hamiltonian is topological, i.e., $\bar{J}_x = 1$, as proven in Sec. VA. With the translational-invariant boundary condition (TIBC) of Eq. (E3), it has also been argued [33] that $\bar{J}_x = 1$ for system sizes much larger than the correlation length; in practice, this can only be verified with a much larger system size than can be simulated with Lanczos. In Fig. 8, we numerically evaluate \bar{J}_x for three different types of boundary conditions: (i) TIBC, (ii) zeroth-order TBC (see Sec. VA), and (iii) first-order TBC [see Eq. (33)]. For all the system sizes that we simulate (up to $L = 30$), the ground states with (i) and (ii) are comparably sensitive to twisting, while the ground state with (iii) is most sensitive. A naive

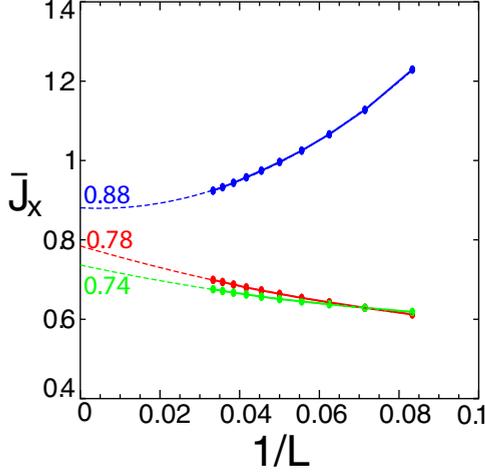


FIG. 8. Finite-size analysis of the closed-chain XYZ model along the $J_x = -J_y$ line, for $J_x \in [0, 1)$. We close the chain using three types of boundary conditions: (i) translational invariant (colored red), (ii) first-order topological (blue), and (iii) zeroth-order topological (green). We define \bar{J}_x such that upon twisting the boundary conditions, the ground-state parity switches for $0 \leq J_x < \bar{J}_x$, but is invariant for $\bar{J}_x < J_x < 1$. We numerically evaluate \bar{J}_x for varying system sizes (up to $L = 30$), as shown by solid dots in the plot. Each dashed line is a fit to the function $a + b/L + c/L^2$, and by extrapolation we obtain an estimate of \bar{J}_x in the thermodynamic limit. Since the fitting function is not theoretically motivated, we emphasize that these extrapolations are more suggestive than rigorous.

extrapolation suggests that these conclusions are robust in the thermodynamic limit.

APPENDIX H: INDISTINGUISHABILITY BY SYMMETRIC, PARAFERMION-LOCAL OPERATORS

Let $|\phi_{\alpha \in \mathbb{Z}_m, n}\rangle$ denote the ground states of the frustration-free Hamiltonian $H_{(N, n)}$. Our basis is chosen as $Q|\phi_{\alpha, n}\rangle = \omega^{n\alpha}|\phi_{\alpha, n}\rangle$. Assuming $g = \text{gcd}(m, n) > 1$, divide $P[H_{(N, n)}]$ into (m/g) subspaces labeled by $\delta \in \mathbb{Z}_m/g$: $\bar{P}_\delta = \sum_{\alpha=0}^{g-1} |\phi_{\alpha m/g + \delta, n}\rangle \langle \phi_{\alpha m/g + \delta, n}|$. Let us use \bar{O} to denote a symmetric, parafermion-local probe. We would like to show that the $g \times g$ matrix $\bar{\mathcal{M}}_{\alpha\beta}^{(\delta)} = \langle \phi_{\alpha m/g + \delta, n} | \bar{O} | \phi_{\beta m/g + \delta, n} \rangle$ is proportional to the identity, up to superpolynomially small finite-size corrections; this is equivalent to the condition (49). The off-diagonal elements vanish since $[\bar{O}, Q] = 0$. The difference between diagonal elements is

$$\begin{aligned} \bar{\mathcal{M}}_{\alpha\alpha}^{(\delta)} - \bar{\mathcal{M}}_{\beta\beta}^{(\delta)} &= \frac{1}{m} \sum_{\mu, \nu=0}^{m-1} \omega^{n\delta(\mu-\nu)} (\omega^{\alpha N(\mu-\nu)/g} - \omega^{\beta N(\mu-\nu)/g}) \\ &\quad \times \langle \psi_{\mu, n} | \bar{O} | \psi_{\nu, n} \rangle, \end{aligned} \quad (\text{H1})$$

where $|\psi_{\mu \in \mathbb{Z}_m, n}\rangle$ are fully polarized states (39) related to $|\phi\rangle$ by the Fourier transform (40). Crucially, the difference in phase

factors vanishes when $\mu = \nu \bmod g$, thus, $\bar{\mathcal{M}}_{\alpha\alpha}^{(\delta)} - \bar{\mathcal{M}}_{\beta\beta}^{(\delta)}$ reduces to a sum of terms proportional to $\mathcal{T}_{\mu, \nu} = \langle \psi_{\mu, n} | \bar{O} | \psi_{\nu, n} \rangle$, with $\mu \neq \nu \bmod g$. A straightforward generalization of Appendix B shows that $\mathcal{T}_{\mu, \nu}$ vanishes superpolynomially in the system size, if \bar{O} is clock local. Now, we address the case when \bar{O} is clock nonlocal but parafermion local, as is possible if \bar{O} is an interedge coupling with support near sites 1 and L . Any interedge coupling may be decomposed into the form $B_l B_r Q^{[[B_r]]}$, where B_l (resp. B_r) is a clock-local operator with a definite charge, and with support near site 1 (resp. L). The nonlocal clock-parafermion transformation (5) implies that there are as many powers of the string Q as the charge of B_r . Once again, our notation for the charge is $Q^\dagger B_r Q = \omega^{[[B_r]]} B_r$. We provide an example of this decomposition for \mathbb{Z}_4 parafermions:

$$\gamma_1 \gamma_2 \gamma_{2L-1} \gamma_{2L} = \omega \sigma_1^2 \tau_1 \sigma_L^2 \tau_L Q^2, \quad (\text{H2})$$

where $B_L = \omega \sigma_1^2 \tau_1$, $B_r = \sigma_L^2 \tau_L$, and $[[B_r]] = 2$.

We then claim that

$$\begin{aligned} \mathcal{T}_{\mu\nu} &= \langle \psi_{\mu, n} | B_l B_r Q^{[[B_r]]} | \psi_{\nu, n} \rangle \\ &= 0 \quad \text{for } \mu, \nu \in \mathbb{Z}_m, \quad \mu \neq \nu \bmod g, \end{aligned} \quad (\text{H3})$$

and for finite-ranged B_r and B_l ; our claim will be proven in Appendix H1. Equation (H3) implies the vanishing of Eq. (H1), and thus a proof of Eq. (49) in the frustration-free limit. If B_l and B_r are superpolynomially decaying into the bulk, Eq. (H1) vanishes to up to superpolynomially small finite-size corrections, which can be bound in a manner analogous to Appendix B. Finally, let us consider deviating symmetrically from the frustration-free limit, while preserving the gap within each of the g charge sectors of \bar{P}_δ . Then, the deformed subspace is related to the old by a quasiadiabatic continuation $V_{\delta, s}$, which is distinct for each g multiplet labeled by δ : $\bar{P}_{\delta, s} = V_{\delta, s} \bar{P}_\delta V_{\delta, s}^\dagger$. Then, applying $[V_{\delta, s}, Q] = 0$,

$$\begin{aligned} \bar{P}_{\delta, s} \bar{O} \bar{P}_{\delta, s} &= V_{\delta, s} \bar{P}_\delta V_{\delta, s}^\dagger \bar{O} V_{\delta, s} \bar{P}_\delta V_{\delta, s}^\dagger \\ &= V_{\delta, s} \bar{P}_\delta (V_{\delta, s}^\dagger B_l V_{\delta, s}) (V_{\delta, s}^\dagger B_r V_{\delta, s}) Q^{[[B_r]]} \bar{P}_\delta V_{\delta, s}^\dagger. \end{aligned} \quad (\text{H4})$$

Since $V_{\delta, s}$ is locality preserving, the dressed operators $V_{\delta, s}^\dagger B_l V_{\delta, s}$ and $V_{\delta, s}^\dagger B_r V_{\delta, s}$ are superpolynomially decaying into the bulk, thus we may apply our frustration-free result $\bar{P}_\delta (V_{\delta, s}^\dagger B_l V_{\delta, s}) (V_{\delta, s}^\dagger B_r V_{\delta, s}) Q^{[[B_r]]} \bar{P}_\delta \approx c \bar{P}_\delta$. The final result is that the indistinguishability condition is robust to superpolynomial accuracy.

1. Vanishing of Eq. (H3) for finite-ranged B_l and B_r

Let us decompose $B_l B_r Q^{[[B_r]]}$ into a sum of terms like $r \prod_{j=1}^L \mathcal{O}_j$, where r is a nonzero complex number, and \mathcal{O}_j is an onsite operator acting on site j , with the general form $\sigma_j^{a_j} \tau_j^{b_j}$ for integers a_j and b_j . In our \mathbb{Z}_4 example (H2),

$$\mathcal{O}_j = \begin{cases} \sigma_j^2 \tau_j^3 & \Rightarrow a_j = 2, \quad b_j = 3 \quad \text{for } j \in \{1, L\}, \\ \tau_j^2 & \Rightarrow a_j = 0, \quad b_j = 2 \quad \text{for } j \in \{2, 3, \dots, L-2, L-1\}. \end{cases} \quad (\text{H5})$$

We deduce two useful properties of $\{a_j, b_j\}$ from the general form of $B_l B_r Q^{[[B_r]]}$:

(a) B_r has support over a set of sites near L , thus, we may express $B_r \propto \prod_{j=L-R+1}^L \sigma_j^{a_j} \tau_j^{b_j}$ for some range $R \ll L$. Since $[[\sigma_j]] = 1$ while $[[\tau_j]] = 0$, $[[B_r]] = \sum_{j=L-R+1}^L a_j$. In our example, $B_r = \sigma_L^2 \tau_L$ and $[[B_r]] = a_L = 2$.

(b) B_l and B_r have no support on any site j far away from any edge, where the only nontrivial operation ($\tau_j^{[[B_r]]}$) arises due to the string operator $Q^{[[B_r]]}$. Thus, for j in the bulk, $b_j = [[B_r]]$ and $a_j = 0$. In example (H5), the bulk sites are labeled by $j \in \{2, 3, \dots, L-2, L-1\}$.

Applying Eq. (39), we now express Eq. (H3) as a sum of products of onsite elements, with the form

$$\mathcal{T}_{\mu\nu} = r \prod_{j=1}^L \langle \tilde{\mu}, n | \sigma_j^{a_j} \tau_j^{b_j} | \tilde{\nu}, n \rangle_j, \quad \text{where } |\tilde{\alpha}, n\rangle_j = \frac{1}{\sqrt{n}} \sum_{\beta=0}^{n-1} |\alpha + m\beta\rangle_j, \quad \sigma_j |\alpha\rangle_j = \omega^\alpha |\alpha\rangle_j. \quad (\text{H6})$$

Each onsite element may be evaluated as

$$\begin{aligned} \langle \tilde{\mu}, n | \sigma_j^{a_j} \tau_j^{b_j} | \tilde{\nu}, n \rangle_j &= \sum_{\alpha, \beta=0}^{n-1} \frac{\omega^{\alpha_j(v+b_j+m\beta)}}{n} \langle \mu + m\alpha | \nu + b_j + m\beta \rangle \\ &= \sum_{\alpha=0}^{n-1} \frac{\omega^{\alpha_j(\mu+m\alpha)}}{n} \delta_{\mu, \nu+b_j \bmod m} = \omega^{\alpha_j \mu} \delta_{\mu, \nu+b_j \bmod m} \delta_{a_j, 0 \bmod n}. \end{aligned} \quad (\text{H7})$$

The second equality follows from μ and $\nu \in \mathbb{Z}_m$. In our \mathbb{Z}_4 example, where $g = m = n = 2$,

$$\langle \tilde{0}, 2 | \mathcal{O}_j | \tilde{1}, 2 \rangle_j = \begin{cases} \frac{1}{2} (\langle 0|_j + \langle 2|_j) \sigma_j^2 \tau_j^3 (|1\rangle_j + |3\rangle_j) = \frac{1}{2} (\langle 0|_j + \langle 2|_j) (|0\rangle_j + |2\rangle_j) = 1 & \text{for } j \in \{1, L\}, \\ \frac{1}{2} (\langle 0|_j + \langle 2|_j) \tau_j^2 (|1\rangle_j + |3\rangle_j) = \frac{1}{2} (\langle 0|_j + \langle 2|_j) (|1\rangle_j + |3\rangle_j) = 0 & \text{for } j \in \{2, 3, \dots, L-1\}. \end{cases} \quad (\text{H8})$$

From the vanishing of onsite elements in the bulk, we deduce that $\mathcal{T}_{01} = 0$. Suppose we try to have the bulk terms be nonvanishing; in the following, we show that one of the edge terms must consequently vanish! From the second line of Eq. (H8), we would need a bulk operator $\tau_j^{b_j}$ with odd b_j . From property (b) above, we would then need that $[[B_r]]$ is also odd. This implies from property (a) above that $\sum_{j=L-R+1}^L a_j$ is odd, for R the range of the Hamiltonian, i.e., for at least one site labeled by $\bar{j} \in \{L-R+1, \dots, L\}$, $a_{\bar{j}}$ is odd. For this particular site, we evaluate

$$\langle \tilde{0}, 2 | \mathcal{O}_{\bar{j}} | \tilde{1}, 2 \rangle_{\bar{j}} = \frac{1}{2} (\langle 0|_{\bar{j}} + \langle 2|_{\bar{j}}) \sigma_{\bar{j}}^{a_{\bar{j}}} \tau_{\bar{j}}^{b_{\bar{j}}} (|1\rangle_{\bar{j}} + |3\rangle_{\bar{j}}). \quad (\text{H9})$$

If $b_{\bar{j}}$ is even, then this term reduces to a sum of $\langle 0|1\rangle$, $\langle 0|3\rangle$, $\langle 2|1\rangle$, and $\langle 2|3\rangle$, which individually vanishes. Suppose we choose $b_{\bar{j}}$ to be odd, then

$$\begin{aligned} \langle \tilde{0}, 2 | \mathcal{O}_{\bar{j}} | \tilde{1}, 2 \rangle_{\bar{j}} &= \frac{1}{2} (\langle 0|_{\bar{j}} + \langle 2|_{\bar{j}}) \sigma_{\bar{j}}^{a_{\bar{j}}} (|0\rangle_{\bar{j}} + |2\rangle_{\bar{j}}) \\ &= \frac{1}{2} (\langle 0|_{\bar{j}} + \langle 2|_{\bar{j}}) (|0\rangle_{\bar{j}} + (-1)^{a_{\bar{j}}} |2\rangle_{\bar{j}}) = 0. \end{aligned} \quad (\text{H10})$$

The last equality follows from $a_{\bar{j}}$ being odd. We have thus proven that one edge term must vanish, and therefore $\mathcal{T}_{01} = 0$ no matter how hard we try.

More generally, we would like to show that $\mathcal{T}_{\mu,\nu} = 0$ for $\mu \neq \nu \bmod g$. If we assume the converse, we require that each onsite element must be nonvanishing, i.e., from Eq. (H7),

$$\begin{aligned} \forall j \in \{1, 2, \dots, L\}, \\ \mu = \nu + b_j \bmod m \quad \text{and} \quad a_j = 0 \bmod n. \end{aligned} \quad (\text{H11})$$

From our assumption, we now prove a contradiction:

(i) For a site j far away from the edge, we derive from Eq. (H11) and property (b) that $[[B_r]] = \mu - \nu \bmod m$. The condition on μ and ν then implies that $[[B_r]]$ is *not* an integer multiple of g .

(ii) $\mathcal{T}_{\mu,\nu} \neq 0$ implies from Eq. (H7) that each of a_j is a multiple of n . With property (a), we derive that $[[B_r]] = \sum_{j=L-R+1}^L a_j = nk$, where $k \in \mathbb{Z}_m$. Since $g = \gcd(m, n)$, this implies $[[B_r]]$ is an integer multiple of g , in contradiction with (i). Thus, we conclude that $\mathcal{T}_{\mu,\nu} = 0$ for $\mu \neq \nu \bmod g$.

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