




Discrete Abelian lattice gauge theories on a ladder and their dualities with quantum clock modelsS. Pradhan ^{1,2,*}, A. Maroncelli ^{3,4,*} and E. Ercolessi ^{1,2}¹*Dipartimento di Fisica e Astronomia, Università di Bologna, 40127 Bologna, Italy*²*INFN, Sezione di Bologna, I-40127 Bologna, Italy*³*Dipartimento di Fisica e Astronomia, Università di Firenze, 50019 Firenze, Italy*⁴*INFN, Sezione di Firenze, I-50019 Firenze, Italy*

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We study a duality transformation from the gauge-invariant subspace of a \mathbb{Z}_N lattice gauge theory on a two-leg ladder geometry to an N -clock model on a single chain. The main feature of this mapping is the emergence of a longitudinal field in the clock model, whose value depends on the superselection sector of the gauge model, implying that the different sectors of the gauge theory can show quite different phase diagrams. To investigate this and see if confined phases might emerge, we perform a numerical analysis for $N = 2, 3, 4$, using both exact diagonalization and the density matrix renormalization group algorithm.

DOI: [10.1103/PhysRevB.109.064410](https://doi.org/10.1103/PhysRevB.109.064410)**I. INTRODUCTION**

Gauge theories constitute the baseline in our microscopical description of physical fundamental laws, and they are a cornerstone of contemporary scientific research. Calculations beyond perturbative regimes, as needed to understand, for example, the quark confinement mechanism in quantum chromodynamics, represent a notorious challenge both analytically and numerically. Standard classical computational methods adopt the Wilson's framework of lattice gauge theories (LGTs) [1–3], in which the continuous space-time is replaced by a discrete set of points, and the calculations are performed in the Euclidean path-integral approach. More recently, inspired by Feynman's idea of quantum simulations [4,5], many authors have adopted a Hamiltonian approach in which only spatial coordinates are discretized, and which might be implemented via a quantum platform once the group degrees of freedom are also discretized, by considering a finite group or by suitable approximations (see [6–10] and references therein). Still, in all approaches, enforcing the gauge constraints to restrict the (analytical, numerical, or experimental) evaluation of observables to the gauge-invariant Hilbert subspace is a challenging task, which is dealt with using different strategies.

In this paper, we consider Abelian LGTs, which are known to exhibit confined/deconfined phases [11–19]. More specifically, in Sec. II we introduce a pure \mathbb{Z}_N gauge model on a (quasi-2D) ladder geometry, which includes both electric and magnetic degrees of freedom and admits superselection sectors, similarly to what happens in the Toric code. To tackle the problem of gauge invariance, in Sec. III we make use of a bond algebraic approach [20,21] to introduce a duality transformation that allows for an exact mapping from the LGT on the ladder restricted to the gauge-invariant Hilbert space

to a 1D N -clock model [22–26] with a transversal field and a longitudinal field. When periodic boundary conditions are enforced, the value of the latter turns out to depend on the superselection sector of the ladder LGT, resulting in possible different phase diagrams for the different sectors, differently from what has been found previously with open boundary conditions [19].

In Sec. IV, we resort to numerical analysis to study the phase diagram in the different sectors (labeled by $n = 0, \dots, N - 1$) for the $N = 2, 3$, and 4 cases. We first make use of exact diagonalization to determine (i) the presence of a deconfined-confined phase transition by calculating the value of the Wilson loops; and (ii) the ground-state structure in the different phases. We find that for all considered N , the model in the $n = 0$ sector is always in the confined phase except for the deconfined point $\lambda = 0$, where only the magnetic degrees of freedom are present in the Hamiltonian. Instead, for N even and $n = N/2$ we see a clear phase transition at about $\lambda = 1$, while for $N = 3$ and $n = 1, 2$ as well as for $N = 4$ and $n = 1, 3$ we observe a crossover region for $\lambda \lesssim 0.8$ followed by a phase transition to a doubly degenerate ground state. A careful study of these cases requires longer chains and therefore is carried out via the density matrix renormalization group (DMRG) algorithm. Finally, in Sec. V we review our results and draw some conclusions.

II. THE LATTICE GAUGE MODEL

Following the Hamiltonian approach of Kogut and Susskind [2], we consider a class of pure Abelian lattice gauge theories, with \mathbb{Z}_N gauge group, on a *ladder geometry*, which consists of a lattice \mathbb{L} made of two parallel chains, the *legs*, coupled to each other by *rungs* to form square plaquettes. On the ladder, each rung is identified by a coordinate $i = 1, \dots, L$, where L is the length of the ladder, and the two vertices on the rung are denoted with i^\uparrow and i^\downarrow in the upper and lower leg, respectively. Links are denoted by ℓ . On the

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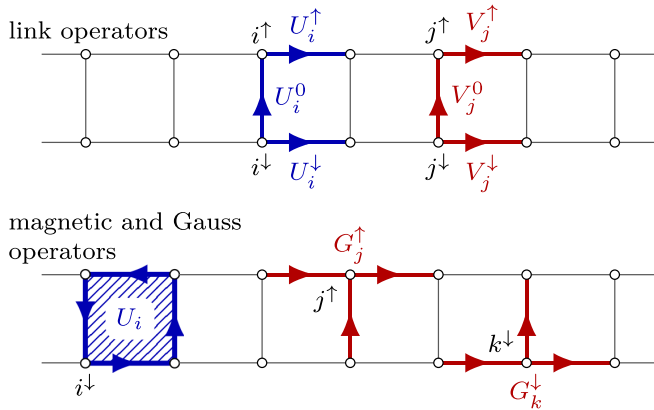


FIG. 1. Operators of the ladder \mathbb{Z}_N LGT. Top panel: Operators U_i^ρ and V_j^ρ ($\rho = \uparrow, \downarrow, 0$) for each link; bottom panel: Gauge-invariant magnetic operator U_i and the Gauss operators G_j^\uparrow and G_k^\downarrow .

legs they are labeled as ℓ_i^\uparrow (upper leg) or ℓ_i^\downarrow (lower leg), while those on the rungs are labeled ℓ_i^0 .

The gauge group degrees of freedom are defined on the links. For a finite group like \mathbb{Z}_N , the notion of infinitesimal generators loses any meaning, and we are led to directly consider, for each link $\ell \in \mathbb{L}$, a pair of conjugate operators U_ℓ and V_ℓ which are unitary and defined by the algebraic relations [27–29]

$$V_\ell U_\ell = \omega U_\ell V_\ell, \quad U_\ell^N = V_\ell^N = \mathbb{1}_N \quad (1)$$

with $\omega = e^{i(\frac{2\pi}{N} + \phi)}$, where the angle ϕ is arbitrary and corresponds to the physical situation in which on each link there is a background electric field [29,30]. In this manuscript, we do not consider this situation and will set $\phi = 0$. Also, these operators commute on different links. This algebra admits a faithful finite-dimensional representation of dimension N [28,31]. To each link ℓ we associate an N -dimensional Hilbert space \mathcal{H}_ℓ generated by an orthonormal basis $\{|v_{k,\ell}\rangle\}$, with $k = 0, \dots, N-1$, the *electric basis*, that diagonalizes V_ℓ :

$$V_\ell |v_{k,\ell}\rangle = \omega^k |v_{k,\ell}\rangle. \quad (2)$$

On this basis, U_ℓ acts as a shift operator,

$$U_\ell |v_{k,\ell}\rangle = |v_{k+1,\ell}\rangle, \quad (3)$$

where $k+1$ is taken mod N .

As shown in the top panel of Fig. 1, we use the symbols V_i^ρ , U_i^ρ for the operators defined on the rung i , and V_i^ρ , U_i^ρ with $\rho = \uparrow, \downarrow$ for the operators on the horizontal links of the upper and lower leg to the right of the rung. The links on the legs are oriented from left to right, while those on the rungs are oriented from bottom to top. To construct a LGT, in addition to the *electric field operators* V 's defined above, we need the following:

(i) The *magnetic operators*, which are defined on each plaquette to the right of the rung i via the formula

$$U_i = U_i^\downarrow U_{i+1}^0 (U_i^\uparrow)^\dagger (U_i^0)^\dagger. \quad (4)$$

(ii) The *Gauss operators*, which are defined on each vertex i^\uparrow , i^\downarrow of the lattice as

$$G_i^\uparrow = V_i^\uparrow (V_{i-1}^\uparrow)^\dagger (V_i^0)^\dagger, \quad G_i^\downarrow = V_i^\downarrow V_i^0 (V_{i-1}^\downarrow)^\dagger \quad (5)$$

and implement local gauge transformations, by imposing that physical states should satisfy $G_i^\rho |\Psi_{\text{phys}}\rangle = |\Psi_{\text{phys}}\rangle$ for $\rho = \uparrow, \downarrow$ and $\forall i$.

It is simple to verify that the U_i -operators commute with all Gauss operators, making them gauge-invariant. The operators defined so far are shown in the bottom panel of Fig. 1.

The gauge-invariant Hamiltonian we use to build a \mathbb{Z}_N LGT on the ladder can be written as

$$H_{\mathbb{Z}_N} = - \sum_i [U_i + \lambda(V_i^\uparrow + V_i^\downarrow + V_i^0) + \text{H.c.}], \quad (6)$$

with $\lambda > 0$, which is the relative strength between the electric and the magnetic fields. One can choose to work with two separate couplings λ_e and λ_m for the electric and magnetic fields, respectively. Nonetheless, we decided to use the ratio $\lambda = \lambda_e/\lambda_m$ for convenience and fix $\lambda_m = 1$ in order to work with just one parameter. We use periodic boundary conditions on legs, which turns out to be an essential step for the duality map provided in the next section.

Similarly to what happens in the two-dimensional Toric code [13,32] (see in Appendix A), the Hilbert space of physical states $\mathcal{H}_{\text{phys}}$ can be decomposed as a direct sum of superselection sectors $\mathcal{H}_{\text{phys}}^{(n)}$, where $n = 0, \dots, N-1$, that can be distinguished by means of the operators

$$\bar{S} = V_{i_0}^\uparrow V_{i_0}^\downarrow, \quad \bar{W} = \prod_{i \in \mathcal{C}_0} U_i^\downarrow, \quad (7)$$

where i_0 label the position of an arbitrary rung in the lattice, while \mathcal{C}_0 is any noncontractible loop around the ladder. They satisfy the relations $\bar{W} \bar{S} = \omega \bar{S} \bar{W}$. Each physical state in a sector $\mathcal{H}_{\text{phys}}^{(n)}$ is an eigenstate of \bar{S} with eigenvalue ω^n , while \bar{W} maps $\mathcal{H}_{\text{phys}}^{(n)}$ into $\mathcal{H}_{\text{phys}}^{(n+1)}$.

III. DUALITY BETWEEN LADDER LGT AND CLOCK MODELS

Clock models [22–24] are a class of models that can be thought of as a generalization of the quantum Ising model. A p -state clock model on a chain has a local p -dimensional Hilbert space for each site $i = 1, \dots, L$ and employs $p \times p$ unitary matrices X_i and Z_i that commute on different sites, while on the same site

$$X_i Z_i = \omega Z_i X_i, \quad (X_i)^p = (Z_i)^p = \mathbb{1}_p, \quad (8)$$

with $\omega = e^{i2\pi/p}$. For example, one can choose a basis where the Z_i 's are diagonal, i.e., $(Z_i)_{mn} = \delta_{m,n} \omega^m$ and $(X)_{mn} = \delta_{m,n+1} \pmod{p}$, with $m, n = 0, \dots, p-1$. The p -clock Hamiltonian is given by

$$H_p(h) = - \sum_i (Z_{i-1}^\dagger Z_i + h X_i + \text{H.c.}), \quad (9)$$

where periodic boundary conditions are assumed, and h is the coupling of the transverse field.

We use the bond-algebraic approach to dualities [20] to introduce a *gauge-reducing duality* mapping between the \mathbb{Z}_N gauge model (with redundant degrees of freedom) on the ladder and an N -clock model on a single chain. Similarly to what can be done in 2D [3,20,33], we associate with each plaquette of the LGT a site of the chain of the clock model,

in such a way that the gauge-invariant magnetic operator U_i is mapped into the single-body operator X_i . The duality of the two-dimensional gauge theories cannot be straightforwardly applied because here the links ℓ^0 have a different role when compared with the links ℓ^\uparrow and ℓ^\downarrow , only the former being *domain walls* between two plaquettes. Also, the electric operators V^\downarrow/V^\uparrow on the top/bottom links $\ell^\uparrow/\ell^\downarrow$ have to be treated separately because they have different commutation relations with the plaquette operators $U_i \equiv U_i^\downarrow U_{i+1}^0 (U_i^\uparrow)^\dagger (U_i^0)^\dagger$:

$$U_i V_i^\downarrow = \omega V_i^\downarrow U_i, \quad U_i V_i^\uparrow = \omega^{-1} V_i^\uparrow U_i. \quad (10)$$

The duality transformation is defined through the following steps. First, the electric field on a vertical link ℓ^0 is mapped to $Z_{i-1}^\dagger Z_i$, as it is the result of the difference of the magnetic states of the two adjacent plaquettes. This can be verified, since from the definition of the plaquette operator U_i we get

$$V_i^0 U_i = \omega^{-1} U_i V_i^0, \quad V_i^0 U_{i-1} = \omega U_{i-1} V_i^0,$$

therefore the maps

$$U_i \mapsto X_i, \quad V_i^0 \mapsto Z_i^\dagger Z_{i-1} \quad (11)$$

conserve the commutation relations of U_i and V_i^0 . Notice that, since from (5) we have

$$\prod_i G_i^\downarrow |\psi_{\text{phys}}\rangle = \prod_i V_i^0 |\psi_{\text{phys}}\rangle = |\psi_{\text{phys}}\rangle, \quad (12)$$

we expect that, after the duality, the product of all V_i^0 is mapped to the identity, as it is from (11). This works as a check of consistency for the duality map.

Second, we consider V^\uparrow and V^\downarrow , which commute with V^0 while satisfying relations (10) with respect to U_i . This allows us to assume

$$V_i^\downarrow \mapsto c_i^\downarrow Z_i, \quad V_i^\uparrow \mapsto c_i^\uparrow Z_i^\dagger, \quad (13)$$

where c_i^\downarrow and c_i^\uparrow are complex numbers, with $|c_i^\downarrow| = |c_i^\uparrow| = 1$ to guarantee unitarity. To further constrain the value of these coefficients, we have to impose that the Gauss constraints (5) become the identity: $G_i^\uparrow \mapsto \mathbb{1}$ and $G_i^\downarrow \mapsto \mathbb{1}$ for all i . Since

$$\begin{aligned} G_i^\uparrow &\mapsto (c_i^\uparrow Z_i^\dagger)(c_{i-1}^\uparrow Z_{i-1}^\dagger)(Z_i^\dagger Z_{i-1})^\dagger = c_i^\uparrow (c_{i-1}^\uparrow)^*, \\ G_i^\downarrow &\mapsto (c_i^\downarrow Z_i)(Z_i^\dagger Z_{i-1})(c_{i-1}^\downarrow Z_{i-1}^\dagger) = c_i^\downarrow (c_{i-1}^\downarrow)^*, \end{aligned} \quad (14)$$

we find that [34]

$$c_i^\downarrow = c_i^\downarrow, \quad c_i^\uparrow = c_i^\uparrow, \quad \forall i \quad (15)$$

for the following reason. Given that all $|c_i| = 1$, the condition $c_i (c_{i-1})^* = 1$ from (14) is equivalent to $c_i = c_{i-1}$, which has to be true for all i .

Finally, since the superselection sectors are identified by the eigenvalue of the operator \bar{S} in (7), whose eigenvalues are simply ω^k , for $k = 0, \dots, N-1$, we get

$$\bar{S} \mapsto (c_i^\uparrow Z_i^\dagger)(c_i^\downarrow Z_i) = c_i^\uparrow c_i^\downarrow = \omega^k. \quad (16)$$

This allows us to fix these coefficients as follows:

$$c_i^\uparrow = 1, \quad c_i^\downarrow = \omega^k. \quad (17)$$

We stress that this freedom of choice for the coefficients c_i^\uparrow and c_i^\downarrow is due to the global \mathbb{Z}_N of the system, not an effect of the already resolved gauge symmetry.

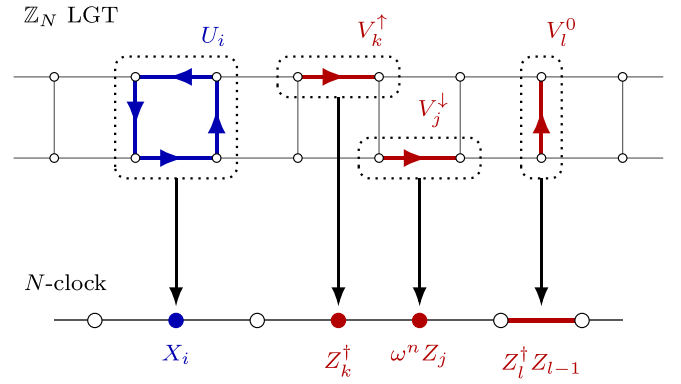


FIG. 2. Visual representation of duality transformation from the \mathbb{Z}_N ladder LGT to the N -clock model.

In summary, the duality mapping for the superselection sector ω^k of the \mathbb{Z}_N LGT on a ladder reads

$$\begin{aligned} U_i &\mapsto X_i, & V_i^0 &\mapsto Z_{i-1}^\dagger Z_i, \\ V_i^\uparrow &\mapsto Z_i^\dagger, & V_i^\downarrow &\mapsto \omega^k Z_i. \end{aligned} \quad (18)$$

A sketch of this duality is given in Fig. 2. The transformed Hamiltonian is

$$H_{\text{lad}}(\lambda) \mapsto \lambda H_N(\lambda^{-1}), \quad (19)$$

where

$$H_N(\lambda^{-1}) = - \sum_i \left(Z_{i-1}^\dagger Z_i + \frac{1}{\lambda} X_i + (1 + \omega^k) Z_i + \text{H.c.} \right). \quad (20)$$

The novelty of (20) is the appearance of a *longitudinal field* term, with a coupling $(1 + \omega^n)$ that depends explicitly on the superselection sector n . Notice that when N is even, the longitudinal field is zero for $n = N/2$. This simple fact makes it reasonable to think that different superselection sectors of the same ladder model can have drastically different phase diagrams.

Let us remark that the complex coupling $(1 + \omega^n)$ does not make the Hamiltonian (20) necessarily chiral [35,36]. In fact, one can get the real Hamiltonian

$$H_N = H_p(1/\lambda) - 2 \cos\left(\frac{\pi n}{N}\right) \sum_i (Z_i + Z_i^\dagger) \quad (21)$$

by absorbing the complex phase in the Z_i -operators, with the transformation $Z_i \mapsto \omega^{-n/2} Z_i$. This transformation globally rotates the eigenvalues of the Z_i -operators while preserving the algebra relations. For n even, this is just a permutation of the eigenvalues, meaning that it does not affect the Hamiltonian spectrum. Instead, for n odd, up to a reorder, the eigenvalues are shifted by an angle π/N , i.e., half the phase of ω . The energy contribution of the extra term in (21) depends on the real part of these eigenvalues, and for n odd we obtain that the lowest energy state is no longer unique. In fact it is doubly degenerate. This means that for $\lambda \rightarrow \infty$, where the extra term becomes dominant, we expect an ordered phase with a doubly degenerate ground state. Finally, one can prove that the sectors n and $N-n$ are equivalent [37].

IV. NUMERICAL INVESTIGATIONS

We wish to investigate the presence of a *deconfined-confining phase transition* (DCPT) for a given \mathbb{Z}_N ladder LGT. In a pure gauge theory, these phases can be detected with the perimeter/area law for Wilson loops [1], which can be expressed as the products of magnetic operators over a given region. Unfortunately, in a ladder geometry there is not much difference between the area and the perimeter of a loop, since they both grow linearly in the size system L .

Nonetheless, we expect a phase transition by varying λ [13,15,16] that can still be captured by an operator like $W_{\mathcal{R}} = \prod_{i \in \mathcal{R}} U_i$, the product of magnetic operators U 's over a (connected) region \mathcal{R} . Indeed, when $\lambda = 0$, the Hamiltonian (6) is analogous to a Toric code [32] which is known to be in a deconfined phase, where the (topologically distinct) ground states are obtained as uniform superpositions of the gauge-invariant states, i.e., closed electric loops. On these ground states $\langle W_{\mathcal{R}} \rangle = 1$, hence a value $\langle W_{\mathcal{R}} \rangle \approx 1$ signals a deconfined phase. On the other hand, when $\lambda \rightarrow \infty$, the electric loops are suppressed, hence $\langle W_{\mathcal{R}} \rangle \approx 0$, signaling a confined phase.

In the dual clock model picture, the Wilson loop translates to a disorder operator [33], which means that a deconfined phase can be thought of as a paramagnetic (or disordered) phase, while the confined phase is like a ferromagnetic (or ordered) phase. Moreover, the longitudinal field breaks the N -fold symmetry of the ferromagnetic phase into a onefold or twofold degeneracy, depending on the superselection sector.

We first start by studying the \mathbb{Z}_N LGT on a ladder numerically through *exact diagonalization* by evaluating the half-ladder Wilson loop, i.e.,

$$W = U_1 U_2 \cdots U_{L/2}, \quad (22)$$

on the ground state while working in the restricted physical Hilbert space $\mathcal{H}_{\text{phys}}^{(n)}$ ($n = 0, \dots, N-1$), which has dimension N^L , much smaller than N^{3L} (the dimension of the total Hilbert space). Then, we proceed to analyze some region of interest by means of the DMRG [38], but on the corresponding dual clock model.

In the following, we will present the results for $N = 2, 3$, and 4, but before doing so we will discuss the implementation of the physical Hilbert space for the exact diagonalization.

A. Implementation of the Gauss law

When considering a LGT, one would like to work within the physical subspace, which is obtained by imposing the Gauss law at every site. A straightforward but inefficient method, in which one generates all the possible states and then filters out all the states that violate Gauss' law, is not efficient, even for moderately small lattices. To better exemplify this, consider a \mathbb{Z}_2 theory on an $L \times L$ periodic lattice, which has L^2 sites and $2L^2$ links, and only 2^{L^2} physical states. There are therefore 2^{2L^2} possible states, and for each one up to L^2 checks (one per site) have to be performed. As a result, the construction of the physical Hilbert space involves $O(L^2 2^{2L^2})$ operations in a search space of 2^{2L^2} objects for finding only 2^{L^2} elements. Here, we exploit the gauge-reducing duality map described in Sec. III for the ladder one to devise an algorithmic procedure that generates physical configurations

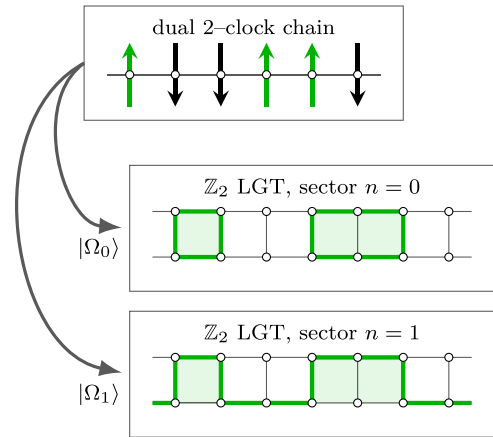


FIG. 3. Duality between the states of a 2-chain and the states of a \mathbb{Z}_2 ladder LGT in the different sectors $n = 0$ (no noncontractible electric loop) and $n = 1$ (one noncontractible loop around the ladder). In the sector $n = 0$ it is evident that all the physical states contain closed electric loops. On the other hand, in the sector $n = 1$ the physical states are all the possible deformations of the electric string that goes around the ladder.

starting from the states of the dual clock model. This is not a search or pattern-matching algorithm, and it gives a *major speedup* with respect to the direct method just described. A similar approach has been used in [39] for a $U(1)$ gauge theory, where, however, an overcomplete basis is found.

Given a \mathbb{Z}_N LGT on a lattice of size $L \times L$, we consider the dual N -clock model on a similar lattice with $A = L^2$ sites. A basis for the Hilbert space of the clock-model is the set of states $|\{s_i\}\rangle \equiv |s_0 s_1 \cdots s_{A-1}\rangle$, with $s_i = 0, \dots, N-1$. The corresponding gauge-invariant state in each superselection sector $\mathcal{H}_{\text{phys}}^{(n,m)}$ of the Hilbert space of the dual LGT model is given by

$$|\{s_i\}\rangle \mapsto \prod_{i=0}^{A-1} U_i^{s_i} |\Omega_{(n,m)}\rangle, \quad (23)$$

where U_i is the plaquette operator on the i th plaquette, and $|\Omega_{(n,m)}\rangle$ is the ‘‘Fock vacuum’’ of the $\mathcal{H}_{\text{phys}}^{(n,m)}$ subspace. Moreover, the ‘‘Fock vacuums’’ $|\Omega_{(n,m)}\rangle$ can be obtained as

$$|\Omega_{(n,m)}\rangle = (\overline{W}_1)^n (\overline{W}_2)^m |\Omega_{(0,0)}\rangle, \quad (24)$$

where $|\Omega_{(0,0)}\rangle$ is the vacuum in the $(0,0)$ -sector, i.e., the state $|000 \cdots 0\rangle$. For more details, see Appendix A. In the case of a ladder geometry, where m is always zero, we shorten the notation of the vacuum states to $|\Omega_n\rangle$. Figure 3 show some examples of physical states in different superselection sectors.

Let us quantify the obtained speedup with this method. In the case of a \mathbb{Z}_2 theory on a square lattice $L \times L$, there are 2^{L^2} possible clock configurations. For each configuration, there are at most L^2 magnetic fluxes to apply. This translates into $O(L^2 2^{L^2})$ operations: Notice that the exponent does not contain the factor 2 that is present in the direct but inefficient method, thus reducing the number of operations by an order of $O(L^2)$. The procedure is generalizable for any \mathbb{Z}_N . This algorithm has been developed independently, but similar techniques for different models can be found, for example, in [39].

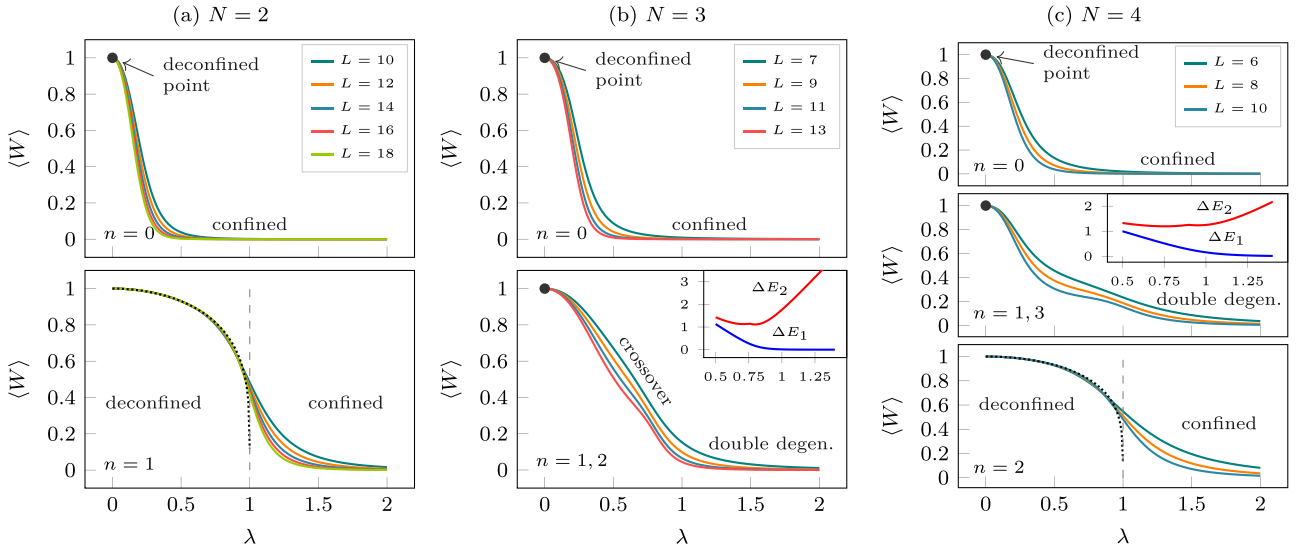


FIG. 4. Expectation values of the half-ladder Wilson loop (22) of the ground state of the \mathbb{Z}_N ladder LGT. They have been computed using exact diagonalization in each superselection sector for different lattice sizes, using 100 values of λ in the range $0 \leq \lambda \leq 2$. (a) Case $N = 2$ for sizes $L = 10, 12, \dots, 18$ and sectors $n = 0$ (top) and $n = 1$ (bottom). Only for $n = 1$ do we have a clear phase transition for $\lambda \simeq 1$, while $n = 0$ is always confined for $\lambda \neq 0$. The limit for large L of the Wilson loop (27) is shown with a dotted line. (b) Case $N = 3$ for sizes $L = 7, 9, 11, 13$ and sectors $n = 0$ (top) and $n = 1, 2$ (bottom, which are equivalent). In the latter, we see the appearance of a crossover region and a doubly degenerate ordered phase. Inset: Energy gaps $\Delta E_i = E_i - E_0$ for $i = 1, 2$ and size $L = 13$, as a function of the coupling λ , in the sectors $n = 1, 2$, showing the emergence of a doubly degenerate ground state for $\lambda > 1$. (c) Case $N = 4$ for sizes $L = 6, 8, 10$ and sectors $n = 0$ (top), $n = 1, 3$ (middle, which are equivalent), and $n = 2$ (bottom). Only the sector $n = 2$ has a clear deconfined-confined phase transition, as expected from the duality with the 4-clock model, while for $n = 1, 3$ the situation is similar to $N = 3$ and $n = 1, 2$. The limit for large L of the Wilson loop for $n = 2$ is the same as (27), shown with a dotted line. This is because a 4-clock model is equivalent to two Ising models [23]. Inset: Energy gaps $\Delta E_i = E_i - E_0$ for $i = 1, 2$ and size $L = 10$. A situation similar to $N = 3$ and $n = 1, 2$ arises.

B. Exact diagonalization for $N = 2$

As a warmup, we consider the \mathbb{Z}_2 ladder LGT, with lengths $L = 10, 12, \dots, 18$. This model is equivalent to a $p = 2$ clock model, which is just the quantum Ising chain, with only two superselection sectors for $n = 0$ and 1. When $n = 1$, the Hamiltonian contains only the transverse field and is integrable [40]. Thus, we expect a critical point for $\lambda \simeq 1$, which will be a DCPT in the gauge model language. This is seen in the behavior of the half-ladder Wilson loop, as shown in the lower panel of Fig. 4(a). For $n = 0$, both the transverse and longitudinal fields are present, the model is no longer integrable [41–43], and we expect to always see a confined phase, except for $\lambda = 0$. This is indeed confirmed by the behavior of the half-ladder Wilson loop shown in the upper panel of Fig. 4(a).

Regarding the Wilson loop in the sector $n = 1$, it is possible to compute its behavior for large L using a Kramers-Wannier duality. Consider $\lambda H_N(\lambda^{-1})$ in (20) for $N = 2$ and $n = 1$. This model is self-dual under the following Kramers-Wannier transformation:

$$X_i \mapsto Z_i Z_{i+1}, \quad Z_{i-1} Z_i \mapsto X_i. \quad (25)$$

Note that $Z = Z^\dagger$ and $X = X^\dagger$ for $N = 2$. With the above map, we obtain

$$\lambda H_{N=2}(\lambda^{-1}) \mapsto H_{N=2}(\lambda),$$

where now the coupling λ acts as a transverse field, hence for $\lambda < 1$ we have a ferromagnetic phase signaled by the magnetization $M = \sum_i \langle Z_i \rangle / L$.

Furthermore, the half-ladder Wilson loop (22) after the gauge-reduction can be written as $W = X_1 \cdots X_{L/2}$, which under the map (25) becomes

$$W = Z_1 Z_{1+L/2}. \quad (26)$$

Such a correlator in the limit of large L reduces to the magnetization squared:

$$\langle W \rangle \sim \langle Z \rangle^2 = M^2 \quad \text{for } L \gg 1.$$

Thanks to the Onsager formula, we have a closed analytical expression for the magnetization, which in the one-dimensional quantum Ising model translates to [44]

$$M \sim (1 - \lambda^2)^{1/8},$$

therefore in the limit of large L we have

$$\langle W \rangle \sim (1 - \lambda^2)^{1/4}. \quad (27)$$

This curve is shown in the lower panel of Fig. 4(a) (dotted line) to compare it with the numerical results. Close to the phase transitions ($\lambda = 1$), the numerical data suffer from strong finite-size corrections, while they are in good agreement with the theoretical result in the region where λ is small and we are deep inside the deconfined phase.

We can further characterize the phases of the two sectors by looking at the structure of the ground state, for $\lambda < 1$ and $\lambda > 1$, which is possible thanks to the exact diagonalization (see Appendix B). In particular, in the deconfined phase of the sector $n = 1$, the ground state is a superposition of the deformations of the noncontractible electric string that makes the $n = 1$ vacuum $|\Omega_1\rangle$. For this reason, this phase can be

thought of as a *kink condensate* [33] (which is equivalent to a paramagnetic phase), where each kink corresponds to a deformation of the string. Instead, for $\lambda > 1$, where we have confinement (as in the $n = 0$ sector), the ground state is essentially a product state, akin to a ferromagnetic state.

C. Exact diagonalization for $N = 3$

The \mathbb{Z}_3 LGT is studied for lengths $L = 7, 9, 11$, and 13 . This model can be mapped to a three-clock model, which is equivalent to a three-state quantum Potts model, with a longitudinal field present in all sectors, as one can see from (21). This field is expected to disrupt any ordered state in the X -basis, hence any deconfined phase in the gauge model. Thus it is not possible to observe a phase transition, as confirmed by the behavior of the half-ladder Wilson loops in Fig. 4(b). As expected, all the sectors present a deconfined point at $\lambda = 0$.

In the case $n = 0$, for $\lambda > 0$ we recognize a quick transition to a confined phase, similar to what happens in [19], while for $n = 1$ and 2 (which are equivalent) the model exhibits a smoother *crossover* to an ordered phase characterized by a doubly degenerate ground state for $\lambda > 1$. In the crossover region, the Wilson loops decrease much slower with respect to the $n = 0$ sector. This could point to a new phenomenology that appears in the sectors $n = 1, 2$. A more detailed analysis of this crossover region can be found in Sec. IV E. Notice that, as discussed above, the presence of the “skew” longitudinal field breaks the threefold degeneracy, expected in an ordered phase of the three-clock model, into a twofold degeneracy only.

D. Exact diagonalization for $N = 4$

The \mathbb{Z}_4 ladder LGTs have four superselection sectors. The behavior of half-ladder Wilson loops as a function of λ is shown in Fig. 4(c). As in the previous models, for $n = 0$ we see a deconfined point at $\lambda = 0$, followed by a sharp transition to a confined phase. The sector $n = 2$, which has no longitudinal field, is the only one to present a clear DCPT for $\lambda \approx 1$, as it is expected from the fact that the four-clock model is equivalent to two decoupled Ising chains [23].

In the two equivalent sectors $n = 1$ and 3 , where the longitudinal field coupling is complex, the Wilson loop shows a peculiar behavior, at least for the largest size ($L = 10$) of the chain: it decreases fast as soon as $\lambda > 0$ to stabilize to a finite value in the region $0.5 \lesssim \lambda \lesssim 1$ before decreasing to zero. As for the $N = 4$ case, we present a deeper analysis of this situation in Sec. IV E. For $\lambda \gtrsim 1$, the system enters a confined phase with a doubly degenerate ground state, as for the \mathbb{Z}_3 model.

E. DMRG analysis of the crossover region

In this section, we further analyze the crossover regions and the possible transition point that appears in the $N = 3$ and 4 cases. In particular, we focus on the following cases: (i) $N = 3$ sector $n = 1$; (ii) $N = 4$ sector $n = 1$. We chose to do so by directly studying the dual clock model with DMRG techniques, which allow us to access much larger lattice sizes. The results obtained with exact diagonalization on the ladder gauge theory all confirm the duality discussed in Sec. III, henceforth we no longer feel the need to study the gauge

model instead. For the DMRG simulations we used open boundary conditions, a maximum bond dimension $\chi = 800$, and a cutoff of 10^{-10} . Additionally, in order to avoid finite-size effects, all the expectation values have been computed on the ground state in the bulk of the system, i.e., in the region from $L/4$ to $3L/4$.

First of all, we confirm that for large λ we have a doubly degenerate ordered phase by looking at Figs. 5(a) and 6(a). They show, for $N = 3$ and 4 , respectively, the energy gaps of the first and second excited levels with respect to the ground-state energy E_0 . In both cases, $E_1 - E_0$ goes to zero for large λ .

Furthermore, in the crossover region of Figs. 4(b) and 4(c) we notice a small bump, even though the lattice size is quite small. These bumps become more evident for larger L , as can be seen in Figs. 5(b) and 6(b). But some rough finite-size scaling shows that the maxima of the bumps do not reach a positive nonzero value in the limit $L \rightarrow \infty$. To reach a more definite answer, simulations of much larger scales are necessary, as the results presented here are just qualitative. We speculate that the origin of these bumps is due to the transition to the doubly degenerate ordered phase.

Since we have an ordered phase for large λ , we expect a nonzero value of the magnetization. Therefore, we computed the expectation value of $(Z + Z^\dagger)/2$ as a function of λ , and the result is shown in Fig. 5(c) for $N = 3$, and in Fig. 6(c) for $N = 4$. The remarkable feature of this magnetization is that a bifurcation arises in both cases. For $N = 3$ it appears around $0.87 \lesssim \lambda \lesssim 0.92$, while for $N = 4$ it is around $1.0 \lesssim \lambda \lesssim 1.1$. The origin of this bifurcation is explained by the doubly degenerate ground states, each with a different value of the magnetization. Due to the lack of conserved quantum numbers, the DMRG algorithm can end up in either of the two cases randomly for each run.

The two degenerate ground states can be easily described. They have to be ferromagnetic states, due to the minus sign in front of $Z^\dagger Z$ in (20), which means all the sites have to be aligned along the same eigenvector in the Z -basis. Moreover, the two possible alignments are given by the highest weight eigenvectors of the longitudinal field $(1 + \omega^n)Z + (1 + \omega^{-n})Z^\dagger$. Then, the corresponding magnetization is just given by their eigenvalues with respect to $(Z + Z^\dagger)/2$. Let $|s\rangle$ be a ket in the Z -basis with eigenvalue ω^s . With some simple algebra, we find the following:

- (i) For $N = 3$ and $n = 1$ the highest weight eigenvectors are $|0\rangle$ and $|2\rangle$, with magnetization 1 and $-1/2$, respectively.
- (ii) For $N = 4$ and $n = 1$ instead we have $|0\rangle$ and $|3\rangle$, with magnetization 1 and 0 , respectively.

These values are confirmed by Figs. 5(c) and 6(c) in the limit of large λ .

The ground states in the large- λ limit have a clear interpretation in the gauge model. Consider the case $N = 3$ and sector $n = 1$. Following the duality in Sec. III and the procedure in Sec. IV A, one can see that the clock state $|0 \cdots 0\rangle$ corresponds to the situation in which all the links in the lower leg are in the $|1\rangle$ state, while the rest are in $|0\rangle$. Conversely, the clock state $|2 \cdots 2\rangle$ corresponds to the configuration where the upper leg is in the $|2\rangle$ state. Similar pictures can be drawn for $N = 4$ and $n = 1$, where it is either the lower leg in the $|1\rangle$ state or the upper leg in the $|3\rangle$ state.

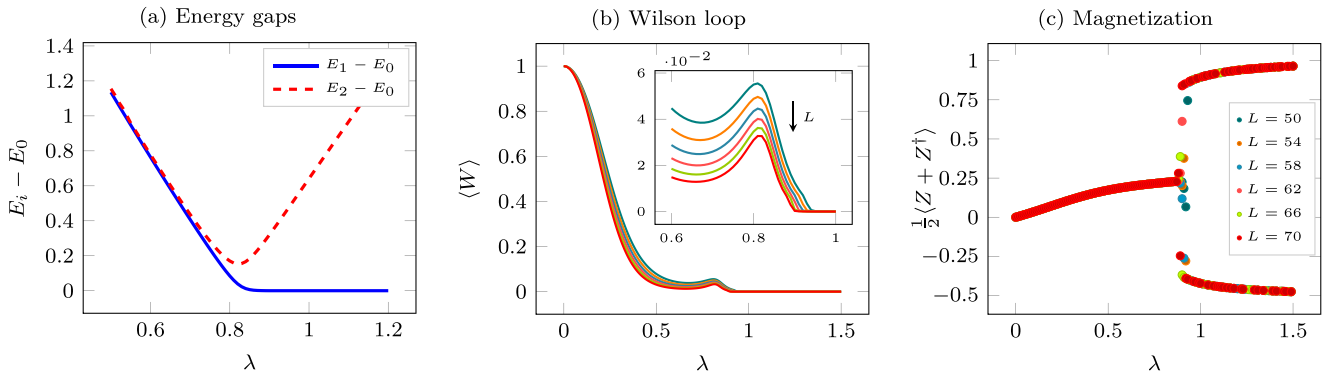


FIG. 5. Numerical analysis of the dual clock model for $N = 3$ and $n = 1$ with DMRG. The expectation values have been computed on the ground state in the bulk of the chain, by which we mean the region from $L/4$ to $3L/4$, in order to avoid finite-size effects. The shown quantities have been calculated for 150 values of λ in the region $0 \leq \lambda \leq 1.5$. (a) Energy gaps of the first two excited states at size $L = 70$; we can see that the large λ region is doubly degenerate. (b) Equivalent of the half-ladder Wilson loop, which is the expectation value of the disorder operator $W = X_{L/4} \cdots X_{3L/4}$. The lattice sizes are $L = 50, 54, \dots, 70$, following the arrow direction. In the inset, a focus on the region $0.6 \leq \lambda \leq 1$ is shown, where bumps are present. In the region $\lambda \lesssim 0.6$ the Wilson loop quickly decreases to zero, while for $\lambda > 1$ it is vanishing. Notice, however, that the heights of the bumps are still small when compared to the deconfined phase. (c) Magnetization, by which we mean the expectation value of $\frac{1}{2}(Z + Z^\dagger)$ averaged over the lattice sites in the bulk. Notice that we have a bifurcation of the order parameter for $0.87 \lesssim \lambda \lesssim 0.9$.

A surprising feature of the crossover region is the behavior of the magnetization. Consider the case $N = 3$. If the phase $\lambda \lesssim 0.87$ is indeed paramagnetic, one would expect a vanishing magnetization. Instead, we find that it follows a profile where it slowly grows, until it reaches a maximum close to 0.25 around $\lambda \approx 0.87$, just before the bifurcation. We exclude the possibility of finite-size effects, because the magnetization is computed in the bulk and its behavior is independent of the chain size, as can be seen in Fig. 5(c). This maximum is close to the average of the two magnetizations in the large- λ limit, which suggest that ground state may be close to a superposition $(|0 \cdots 0\rangle + |2 \cdots 2\rangle)/\sqrt{2}$. In the region close to the transition point ($\lambda \sim 1$), we observe that the numerical data for the magnetization are very scattered, while Wilson loops show a nonzero bump in the expectation value. We

interpret this as a signal that, in this region, the numerics is strongly affected by the presence of low-energy disordered states, which are instead suppressed deeper in the confined region.

The same argument can be repeated for $N = 4$, with the only noticeable difference being a clear plateau of a vanishing Wilson loop before the bifurcation. This means that the ground state is much closer to the superposition $(|0 \cdots 0\rangle + |3 \cdots 3\rangle)/\sqrt{2}$ in this case.

V. CONCLUSIONS AND OUTLOOKS

In this work, we proposed an exact gauge reducing duality transformation that maps the \mathbb{Z}_N lattice gauge theory on a ladder onto a 1D N -clock model in a transversal field, coupled

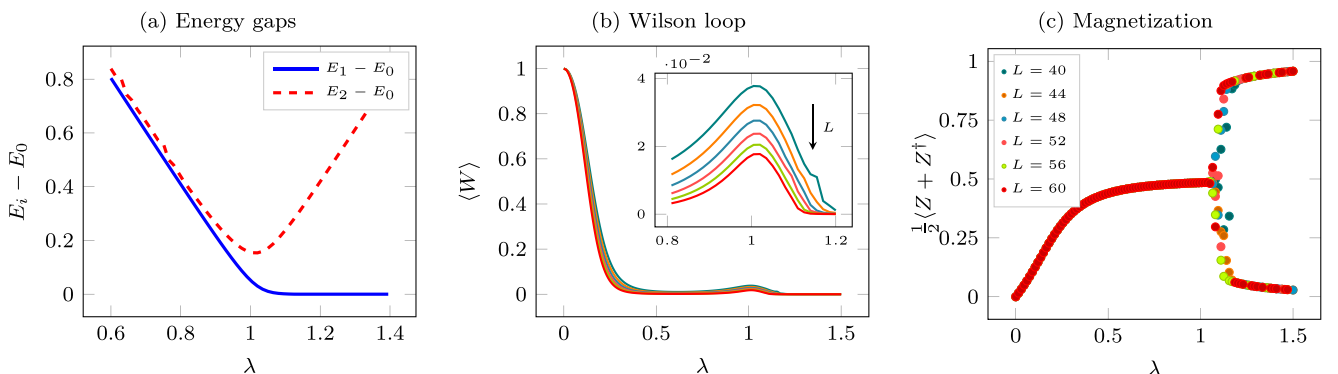


FIG. 6. Numerical analysis of the dual clock model for $N = 4$ and $n = 1$ with DMRG. The expectation values have been computed on the ground state in the bulk of the chain, by which we mean the region from $L/4$ to $3L/4$, in order to avoid finite-size effects. The shown quantities have been computed for 100 values of λ in the region $0 \leq \lambda \leq 1.5$. (a) Energy gaps of the first two excited states at size $L = 60$; we can see that the large λ region is doubly degenerate. (b) Equivalent of the half-ladder Wilson loop, which is the expectation value of the disorder operator $W = X_{L/4} \cdots X_{3L/4}$. The lattice sizes are $L = 50, 54, \dots, 70$, following the arrow direction. In the inset, a focus on the region $0.8 \leq \lambda \leq 1.2$ is shown, where bumps are present. The only clear difference with the $N = 3$ case is the presence of a more consistent plateau, for $\lambda \lesssim 0.8$, where the Wilson loop vanishes. (c) Magnetization, by which we mean the expectation value of $\frac{1}{2}(Z + Z^\dagger)$ averaged over the lattice sites in the bulk. Here the bifurcation point has moved to $\lambda \approx 1$.

to a possibly complex longitudinal field that depends on the superselection sector.

This map allowed us to perform numerical simulations with an exact diagonalization algorithm with sizes up to $L = 18, 13, 10$ for $N = 2, 3, 4$, respectively. To study the phases of the model and a possible DCPT point, we calculated the Wilson loops in the different superselection sectors. For N even and $n = N/2$, we obtain a DCPT point; for $n = 0$ and any N , we are always in a confined phase when $\lambda \neq 0$; while we find an unusual behavior in the other cases ($N = 3$ with $n = 1, 2$ and $N = 4$ with $n = 1, 3$). In particular, the case of $n = 1$ for $N = 3, 4$ has been further analyzed using DMRG techniques, and the results suggest the emergence of a phase that cannot be properly called paramagnetic.

We have shown that the phase diagram of these gauge models depends heavily on the superselection sectors. Such sectors exist only with periodic boundary conditions, but one can obtain a similar setup with open boundary conditions instead. It is sufficient to fix the electric flux at the ends of the ladder, which will be effectively equivalent to looking at the bulk of the periodic lattice. Therefore, boundary conditions play a key role in the phenomenology of these models.

The results presented here regarding the odd sectors are just qualitative and they deserve a proper analysis, which we plan to do in the future. Another possible direction for future work is the inclusion of matter (static or dynamical) in these gauge models, and how they affect the correspondence with quantum clock models.

ACKNOWLEDGMENTS

The numerical analyses have been performed with the QuSpin library [45,46] for the exact diagonalization, and the ITensor library [47,48] for the DMRG simulations. We thank M. Burrello and O. Pomponio for useful discussions. This research is partially supported by Istituto Nazionale di Fisica Nucleare (INFN) through the project ‘‘QUANTUM,’’ the project ‘‘SFT,’’ and the project ‘‘QuantHEP’’ of the QuantERA ERA-NET Cofund in Quantum Technologies (GA No. 731473).

APPENDIX A: REVIEW OF TWO-DIMENSIONAL LGTS AND THE TORIC CODE

In this Appendix, we review some aspects of LGTs in two dimensions. For a more general review, we suggest [13,32].

For a discrete group like \mathbb{Z}_N , the notion of infinitesimal generators loses any meaning and we are led to directly consider, for each link $\ell \in \mathbb{L}$, two unitary operators V_ℓ, U_ℓ , such that [27,49]

$$V_\ell U_\ell V_\ell^\dagger = e^{2\pi i/N} U_\ell, \quad U_\ell^N = \mathbb{1}_N, \quad V_\ell^N = \mathbb{1}_N, \quad (\text{A1})$$

while on different links they commute. Thus, by representing \mathbb{Z}_N with the set of the N roots of unity $e^{i2\pi k/N}$ ($k = 1, \dots, N$), commonly referred to as the discretized circle, we see that V plays the role of a ‘‘position operator’’ on the discretized circle, while U plays the role of a ‘‘momentum operator.’’

These algebraic relations admit a faithful finite-dimensional representation of dimension N [31], for any integer N , which is obtained as follows. To each link ℓ , we

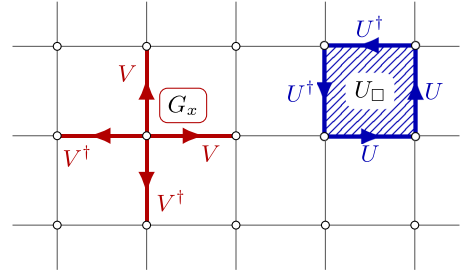


FIG. 7. Pictorial representation of the Gauss operators G_x in (A3) (left) and plaquette operator U_\square in (A7) (right).

can associate an N -dimensional Hilbert space \mathcal{H}_ℓ generated by an orthonormal basis $\{|v_{k,\ell}\rangle\}$ ($k = 1, \dots, N$), called the *electric basis*, that diagonalizes V_ℓ . With this choice, we can promptly write the actions of U_ℓ and V_ℓ :

$$\begin{aligned} U |v_{k,\ell}\rangle &= |v_{k+1,\ell}\rangle, & U |v_{N,\ell}\rangle &= |v_{1,\ell}\rangle, \\ U^\dagger |v_{k,\ell}\rangle &= |v_{k-1,\ell}\rangle, & U^\dagger |v_{1,\ell}\rangle &= |v_{N,\ell}\rangle, \\ V |v_{k,\ell}\rangle &= \omega^k |v_{k,\ell}\rangle, & V^\dagger |v_{k,\ell}\rangle &= \omega^{-k} |v_{k,\ell}\rangle, \end{aligned} \quad (\text{A2})$$

where $\omega = e^{2\pi i/N}$ and $k = 0, \dots, N-1$. We choose to work in this particular basis, and the various k can be interpreted as the quantized values of the electric field on the links.

1. Gauge invariance and physical states

Gauge transformations act on vector potentials while preserving the electric field. In the case of a discrete symmetry, a gauge transformation at a site $x \in \mathbb{L}$ is a product of V 's (and V^\dagger 's) defined on the links that comes out of (and enters) the vertex. More specifically, for a two-dimensional lattice, if the link ℓ at site x is oriented in the positive direction, i.e., either $(x, +\hat{1})$ or $(x, +\hat{2})$, then V is used, otherwise V^\dagger is used. Thus, the single local gauge transformation at the site x is enforced by the operator

$$G_x = V_{(x,\hat{1})} V_{(x,\hat{2})} V_{(x,-\hat{1})}^\dagger V_{(x,-\hat{2})}^\dagger, \quad (\text{A3})$$

as shown in the left part of Fig. 7.

The total Hilbert space \mathcal{H}^{tot} is given by $\otimes_\ell \mathcal{H}_\ell$. A state of the whole lattice $|\Psi_{\text{ph}}\rangle \in \mathcal{H}^{\text{tot}}$ is said to be *physical* if it is a *gauge-invariant state*:

$$G_x |\Psi_{\text{ph}}\rangle = |\Psi_{\text{ph}}\rangle, \quad \forall x \in \mathbb{L}. \quad (\text{A4})$$

This condition can be translated into a constraint on the eigenvalues $v_{(x,\pm\hat{i})} = \omega^{k_{(x,\pm\hat{i})}}$ of the operators V_ℓ on the links $\ell = (x, \pm\hat{i})$ of the vertex x :

$$v_{(x,\hat{1})} v_{(x,\hat{2})} v_{(x,-\hat{1})}^* v_{(x,-\hat{2})}^* = 1, \quad (\text{A5})$$

or, because of (A2),

$$\sum_{i=1,2} (k_{(x,\hat{i})} - k_{(x,-\hat{i})}) = 0 \pmod{N}. \quad (\text{A6})$$

Given the fact that the k 's in (A1) represent the values of the electric field, one can see that (A6) can be interpreted as a discretized version of the Gauss law $\nabla \cdot \vec{E} = 0$ in two dimensions, for a pure gauge theory where there are no electric charges.

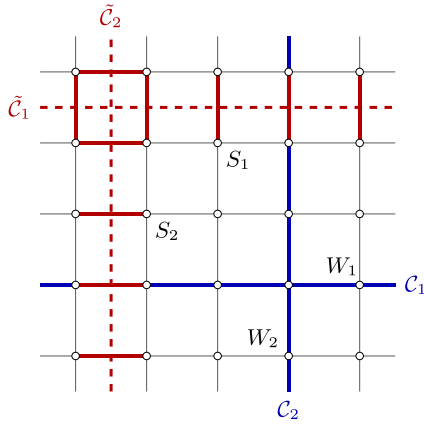


FIG. 8. Graphical representation of the nonlocal order parameters $\bar{W}_{1,2}$ (in blue) and $\bar{S}_{1,2}$ (in red) and their respective paths $\mathcal{C}_{1,2}$ and $\tilde{\mathcal{C}}_{1,2}$.

One can see that the electric operators V_ℓ and the plaquette operators

$$U_\square = U_{(x,\hat{1})}U_{(x+\hat{1},\hat{2})}U_{(x+\hat{1}+\hat{2},-\hat{1})}^\dagger U_{(x+\hat{2},-\hat{2})}^\dagger \quad (\text{A7})$$

are gauge-invariant. Then, the Hamiltonian of the model can be written as

$$H_{\mathbb{Z}_N}(\lambda) = - \sum_{\square} U_\square - \lambda \sum_{\ell} V_\ell + \text{H.c.} \quad (\text{A8})$$

2. Superselection sectors

One of the main features of (A8) is the presence of topologically protected superselection sectors. To illustrate this, we first need to define the noncontractible Wilson loop operators

(pictured in blue in Fig. 8):

$$\bar{W}_i = \prod_{\ell \in \mathcal{C}_i} U_\ell, \quad i = 1, 2, \quad (\text{A9})$$

where \mathcal{C}_i are noncontractible loops around the lattice \mathbb{L} , along the \hat{i} direction. A simple calculation shows that both \bar{W}_1 and \bar{W}_2 commute with all G_x , thus they are gauge-invariant, but one also finds out that none of them can be written as a product of U_\square nor V_ℓ .

Besides \bar{W}_i , another type of nonlocal operators has to be introduced. They are defined on *cuts* of the lattice \mathbb{L} , i.e., paths on the dual lattice $\tilde{\mathbb{L}}$. Consider *noncontractible* cuts $\tilde{\mathcal{C}}_1$ and $\tilde{\mathcal{C}}_2$ along the directions $\hat{1}$ and $\hat{2}$, respectively. On these cuts, the 't Hooft string operators \bar{S}_1 and \bar{S}_2 are constructed as

$$\bar{S}_i = \prod_{\ell \in \tilde{\mathcal{C}}_i} V_\ell, \quad i = 1, 2, \quad (\text{A10})$$

in a similar fashion to (A9) (shown in red in Fig. 8).

These two classes of nonlocal operators resemble the same operators of the Toric code [32] that distinguish the degenerate ground states. One key difference here is that the operators \bar{W}_i do not commute with the Hamiltonian (A8), which contains an electric field term. Thus, unlike the Toric code, we no longer have degenerate ground states when $\lambda \neq 0$. But we can still use the \bar{S}_i operators to decompose the Hilbert space $\mathcal{H}_{\text{phys}}$, since they still commute with all the *local operators* U_\square and V_ℓ (thus also with $H_{\mathbb{Z}_N}$). Now one can see that the operator \bar{S}_i ($i = 1, 2$) of (A10) has N eigenvalues ω^n , with $n = 1, \dots, N - 1$. Hence, one can decompose $\mathcal{H}_{\text{phys}}$ as a sum of superselection sectors,

$$\mathcal{H}_{\text{phys}} = \bigoplus_{n,m=0}^{N-1} \mathcal{H}_{\text{phys}}^{(n,m)}, \quad (\text{A11})$$

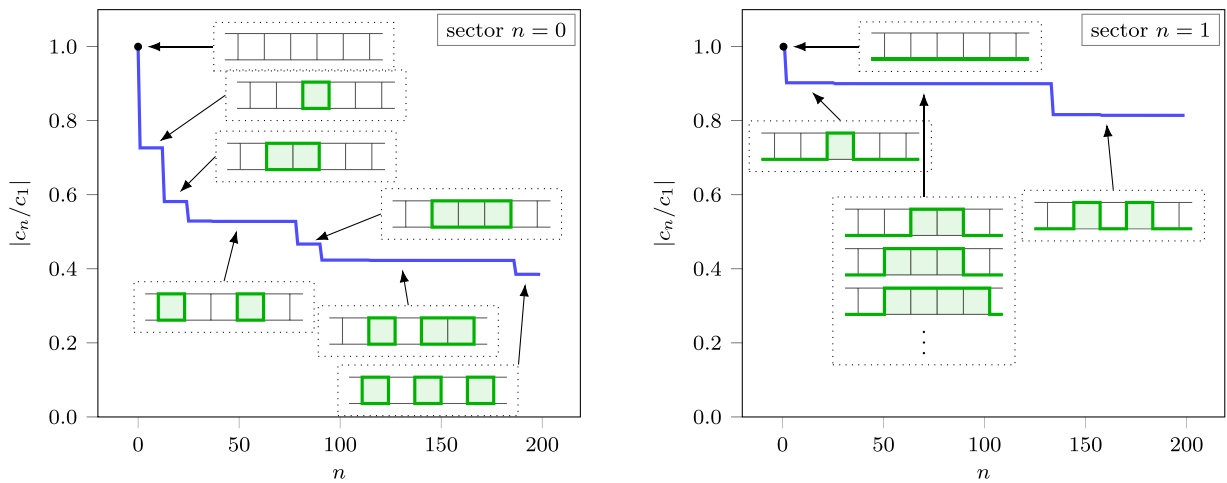


FIG. 9. \mathbb{Z}_2 ground-state amplitude distribution for $\lambda = 0.1$ of the first 200 states and with lattice size 12×2 . Left: Distribution of the ratios $|c_n/c_1|$ for the sector $n = 0$ [see (B1)]. We see that the heaviest states that enter the ground state, apart from the vacuum that sets the scale, are made of small electric loops, typical of a confined phase. Right: The same distribution of ratios for the sector $n = 1$. We see that the heaviest states are made of bigger and bigger deformations of the electric string that goes around the ladder. This happens because the energy contributions depend only on the number of domain walls between two regions with different flux content. The length of each step in the (blue) curves can be found by means of combinatorial methods. For example, the first step in the left panel corresponds to the 12 possible states that contain just one closed loop encircling a single plaquette.

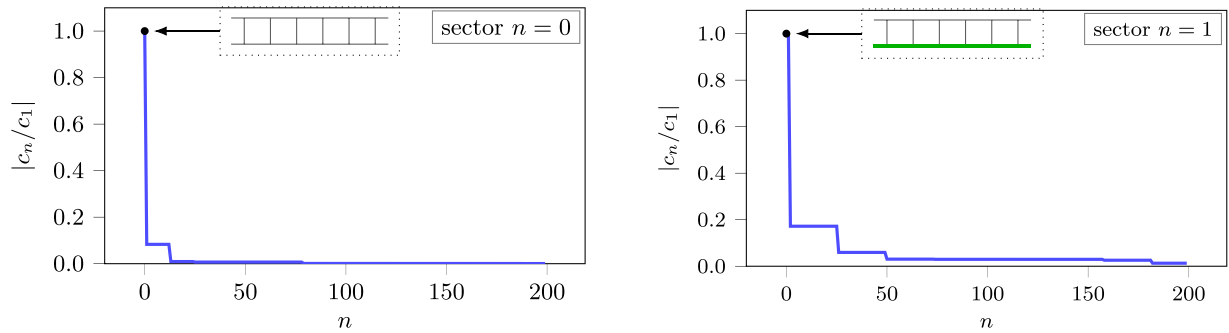


FIG. 10. \mathbb{Z}_2 ground-state amplitude distribution for $\lambda = 1.5$ of the first 200 states and with lattice size 12×2 . For both sectors $n = 0$ (left) and $n = 1$ (right) we are in a confined phase, which corresponds to a ferromagnetic phase in the Ising chain. Here we see a polarized state where the domain walls are suppressed and the ground state is essentially a product state.

where for each $|\phi\rangle \in \mathcal{H}_{\text{phys}}^{(n,m)}$ we have

$$\bar{S}_1 |\phi\rangle = \omega^m |\phi\rangle, \quad \bar{S}_2 |\phi\rangle = \omega^n |\phi\rangle. \quad (\text{A12})$$

Let us consider now the role of the Wilson loops \bar{W}_i . One can see that

$$\bar{W}_2 \bar{S}_1 = \omega \bar{S}_1 \bar{W}_2, \quad \bar{W}_1 \bar{S}_2 = \omega \bar{S}_2 \bar{W}_1. \quad (\text{A13})$$

It follows that $\bar{W}_{1,2}$ acts as a shift operator for the eigenspaces of $\bar{S}_{2,1}$:

$$\bar{W}_1 : \mathcal{H}_{\text{phys}}^{(n,m)} \rightarrow \mathcal{H}_{\text{phys}}^{(n+1,m)}, \quad \bar{W}_2 : \mathcal{H}_{\text{phys}}^{(n,m)} \rightarrow \mathcal{H}_{\text{phys}}^{(n,m+1)}, \quad (\text{A14})$$

where the integers $n+1$ and $m+1$ have to be taken mod N .

From a physical point of view, the Wilson loops operators \bar{W}_1 and \bar{W}_2 create noncontractible electric loops around the lattice, while the 't Hooft strings \bar{S}_2 and \bar{S}_1 detect the presence and the strength of these electric loops. Therefore, it is clear that the Hilbert subspace $\mathcal{H}_{\text{phys}}^{(n,m)}$ is the subspace of all the states that contains an electric loop of strength ω^n and ω^m along the $\hat{1}$ and $\hat{2}$ direction, respectively. Furthermore, the evolution of a state in $\mathcal{H}_{\text{phys}}^{(n,m)}$ with the Hamiltonian in (A8) is confined in $\mathcal{H}_{\text{phys}}^{(n,m)}$.

APPENDIX B: DISTRIBUTION OF THE AMPLITUDES IN THE GROUND STATE

In the $N = 2$ case, we further differentiate the phase diagrams of the two sectors by looking at the ground-state amplitudes distribution for $\lambda < 1$ and $\lambda > 1$. The ground state

can be written as a superposition of the gauge-invariant states of $\mathcal{H}_{\text{phys}}$ in the given sector

$$|\Psi_{\text{g.s.}}\rangle = \sum_n c_n |n\rangle. \quad (\text{B1})$$

The basis $|n\rangle$ and the amplitudes c_n are sorted in decreasing order with respect to their modulus. The first state of the list, with amplitude c_1 , is always the Fock vacua $|\Omega_n\rangle$ of the sector n , hence we consider the distribution of the ratios $|c_n/c_1|$, which are plotted in Figs. 9 and 10 for $\lambda = 0.1$ and 1.5, respectively.

The most interesting one is at $\lambda = 0.1$ in Fig. 9, where the difference between the deconfined phase in the sector $n = 1$ and the confined one in the sector $n = 0$ can be seen. In particular, in the sector $n = 1$ the ground state is a superposition of deformations of the noncontractible electric loop (that makes the Fock vacuum). For this reason, the ground state can be thought of as a *kink condensate* [33] (which is a paramagnetic state), where each kink corresponds to a deformation of the electric loop. This behavior differs from the confined state of the sector $n = 0$, where most of the contributions to the ground state come from states with small electric loops. In other words, the creation of magnetic fluxes is suppressed. The peculiar behavior of the sector $n = 1$ is due to the fact that the energy contributions depend only on the domain walls between two plaquettes with different flux content, while in the sector $n = 0$ we also receive energy contributions from the presence of the fluxes themselves.

Meanwhile, for $\lambda > 1$, where we have confinement in both sectors, the ground state is essentially a product state, akin to a ferromagnetic state, as is shown in Fig. 10.

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