Entanglement entropy of topological phases with multipole symmetry

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(Received 23 February 2023; revised 21 May 2024; accepted 10 July 2024; published 25 July 2024)

We study entanglement entropy of unusual \mathbb{Z}_N topological stabilizer codes which admit fractional excitations with restricted mobility constraint in a manner akin to fracton topological phases. It is widely known that the subleading term of the entanglement entropy of a disk geometry in conventional topologically ordered phases is related to the total number of the quantum dimension of the fractional excitations. We show that, in our model, such a relation does not hold, i.e., the total number of the quantum dimension varies depending on the system size, whereas the subleading term of the entanglement entropy takes a constant number irrespective to the system size. We give a physical interpretation of this result in the simplest case of the model. More thorough analysis on the entanglement entropy of the model on generic lattices is also presented.

DOI: 10.1103/PhysRevB.110.045146

I. INTRODUCTION

Since their discovery, topologically ordered phases have had a tremendous impact on the condensed matter physics community [1-4]. These phases have drawn a lot of attention mainly because of the exotic fractionalized excitations they carry, called anyons [2,5]. For practical purposes, they may find application in quantum computers [6-8].

While a lot of efforts have been made to understand these phases from both a theoretical and experimental point of view, recently new types of topologically ordered phases have been introduced, referred to as the fracton topological phases [9-11]. These phases cannot be described by the preexisting frameworks of the topologically ordered phases, due to the distinctive feature that the phases are sensitive to the local geometry of the system. Key insight to intuitively understand these phases is that a mobility constraint is imposed on the fractional excitations, giving rise to the subextensive ground-state degeneracy (GSD). Establishing a complete theoretical framework of these phases is an active research topic.

One of the attempts to handle this problem is to introduce new types of symmetries, *multipole symmetries*. The multipole symmetry, in particular, the U(1) multipole symmetry, is the generalization of the global U(1) symmetry; a theory is invariant under the global phase rotation depending on the spatial coordinate in polynomial form. As an example, in the case of a scalar theory which respects the global and dipole U(1) symmetries, the Lagrangian is invariant under $\Phi \rightarrow e^{ia+ibx}\Phi$, where *a* and *b* represent constants, and *x* does the spatial coordinate [12]. Investigating topological phases in this viewpoint of the new symmetry has been recently started [12–18].

Motivated by this new interest in the interplay between topological phases and multipole symmetries, in [19], a simple stabilizer model of unconventional topological phases with dipole symmetry (which is one of the multipole symmetries) was studied. As we review in the next section, a unique feature of the model is that each term of the stabilizer model involves next-nearest-neighboring qubits, which is in contrast with the stabilizer models of conventional topological phases where each term acts on nearest-neighboring qubits. As a consequence, a pair of *dipole* of fractional charges are created [see (12)] as an excited state. Also, the model has a unit cell consisting of N lattice site spacing, corresponding to the dipole symmetry [see (13)]. Due to this unusual fusion rule, the model admits dipole of the Wilson loops (viz., Wilson loops whose charge intensity changes linearly as function of x coordinate) of the fractional excitations, giving rise to the unusual ground-state degeneracy (GSD) dependence on the system size, as opposed to the conventional topological phase which admits homogeneous Wilson loops (i.e., Wilson loops which does not depend on position.). In this work, we dub topological phases with multipole symmetries higherrank topological phases. Note that the higher-rank topological phases are distinct from previous fracton models in that these models stem from another type of symmetry, subsystem symmetry, a symmetry acting on a submanifold.

While the GSD has been thoroughly investigated in the model in [19] as well as other higher-rank topological phases, other physical properties of the models have yet to be explored. To address this problem, in this paper, we investigate the interplay between the higher-rank topological phases and the *entanglement entropy*, which is an important physical quantity to diagnose the quantum entanglement.

A plethora of progress has been made in understanding quantum entanglement of conventional topologically ordered phases. It is well known that the entanglement entropy of the two-dimensional (2D) topologically ordered phases in the disk bipartite geometry A, S_A scales as [20,21]

$$S_A \sim \alpha l - \gamma + \cdots$$
 (1)

The first term, which is the leading term, is the so-called area-law term, proportional to the perimeter of the disk geometry l with α being nonuniversal constant, whereas the second constant term, which is the subleading term, is of

TABLE I. Summary of the entanglement entropy S_A of various subsystem geometries with respect to the ground state $|\xi_{00,00}\rangle$ and the one with the generic ground state (45). Here, $\Gamma = N \times \text{gcd}(N, n_x)$ and (Area) represents the number of clock states with type 1 surrounding the disk geometry.

| | $ \xi_{00,00} angle$ | Generic ground state $ \zeta\rangle$ [Eq. (45)] |
|---------------------------|-----------------------------|---|
| Single row I [Fig. 4(a)] | $n_x \log N - \log \Gamma$ | $n_x \log N - \log \Gamma + \mu_1$ [Eq. (47)] |
| Single row II [Fig. 4(b)] | $n_x \log N$ | $n_x \log N + \mu_2$ [Eqs. (49) and (50)] |
| Single column [Fig. 4(c)] | $2n_{\rm y}\log N - \log N$ | $2n_{v} \log N - \log N + \mu_{3}$ [Eqs. (56) and (57)] |
| Disk geometry [Fig. 5(a)] | $(Area)\log N - 2\log N$ | $(Area)\log N - 2 \log N$ |

particular importance as it is universal number, referred to as the *topological entanglement entropy*. It is rewritten as

$$\gamma = \log\left(\sqrt{\sum_{a} d_{a}^{2}}\right),\tag{2}$$

where d_a represents the quantum dimension of the anyon labeled by a. The ellipsis in (1) represents terms that vanish in the limit of $l \rightarrow \infty$.]

We study entanglement entropy of various geometries of the subsystems in the unconventional topological phase with the dipole symmetry. By making use of the formalism of the entanglement entropy for stabilizer codes [22], jointly with the one in the graph theory, we show that the entanglement entropy is described by the same form as (1), yet Eq. (2), which is an important relation between entanglement entropy and the total quantum dimension of fractional excitations, does not hold. Indeed, while the total quantum dimension varies depending on the geometry of the lattice, the topological entanglement entropy γ takes the constant number, irrespective to the lattice. We give an intuitive interpretation of the result by focusing on the simplest case by setting N = 2. We further study the entanglement entropy of the higher-rank topological phase on the generic lattices. We show that the entanglement entropy of the subgraph consists of the two types of the terms, the leading-order area-law term and the subleading constant term, both of which depend on the the number of vertices surrounding the subgraph and the invariant factors of the Laplacian, which is the matrix describing how the vertices are connected in the graph. The result is summarized in Table I and Theorem 1.

There has been an intimate relation between quantum entanglement and graph theory. For instance, multiparty entangled states are described by a graph and its properties are studied based on graph theory (see, e.g., [23,24]). This work addresses the interplay between graph theory and quantum entanglement in the context of (new type) topological stabilizer models by studying the Laplacian. Our work would comply with recent interest in unconventional topological phases, in particular, higher-rank topological phases with multipole symmetries, from the viewpoint of the quantum entanglement and graph theory.

The rest of the paper is organized as follows. In Sec. II, we review the model and its properties studied in [19]. After reviewing the model, in Sec. III, we calculate entanglement entropy of various geometries of subsystems. In Sec. IV, we give an intuitive understanding of our result on the entanglement entropy. In Sec. V, we further investigate the entanglement entropy of a subgraph in the case where the

model is defined on a generic lattice. We make a brief comment on other cases of topological phases in Sec. VI. Finally, we conclude our work in Sec. VII. Technical details are relegated to Appendixes.

II. STABILIZER MODEL

In this section, we review the model constructed by stabilizers given in [19]. Also, we go over physical properties of the model. As we mentioned in the previous section, the model exhibits unusual behavior of the excitations compared with the conventional topologically ordered phases, yielding the unusual GSD dependence on the lattice.

A. Hamiltonian

To start, we introduce 2D square lattice where we place two types of *N*-qubit state (\mathbb{Z}_N clock states) on each vertex and vertical link. The first clock states are located at vertices of the lattice whereas the second ones are at vertical links. We label the coordinate of the first clock states by (x, y) and the ones of the second clock states are denoted by $(x, y + \frac{1}{2})$, where the second element corresponds to the links between vertices located at (x, y) and (x, y + 1). We term these two types of the clock states clock state with type 1 and clock state with type 2.

We represent basis of the two types of the clock states as $|\omega\rangle_1$ and $|\omega\rangle_2$ with ω being *N*th root of unity, i.e., $\omega = e^{2\pi i/N}$ [red square and blue dot in Fig. 1(a)], and \mathbb{Z}_N shift and clock operators (they become Pauli operators when N = 2) of the first and second clock states as $\{Z_i, X_i\}$ (i = 1, 2). Here we have introduced the subscript i = 1, 2 to distinguish operators



FIG. 1. (a) Two types of \mathbb{Z}_N clock states, defined on each vertex (red square) and vertical link (blue dot). (b), (c) Two types of terms introduced in (4).



FIG. 2. (a) Two terms defined in (7) constituting the Hamiltonian in the case of N = 2. (b) Configurations of stabilizers corresponding to (8) in the case of n_x even. In this case, the model is decomposed into two \mathbb{Z}_2 toric codes. The periodic boundary condition is imposed in the *x* direction so that left and right edges are identified. The left (right) geometry corresponds to configuration of the vertex and plaquette terms which belong to (i) [(ii)] defined in (8). Such decomposition is not possible in the case of n_x being odd. We set the coordinate of the vertex at the bottom left to be (1,1).

that act on the clock clock states with type 1 and the ones with type 2. These operators satisfy the following relation (I_i denotes the identity operator):

$$X_i^N = Z_i^N = I_i, \ Z_i |\omega\rangle_i = \omega |\omega\rangle_i, \ X_i Z_j = \omega Z_j X_i \delta_{i,j}.$$
(3)

With these notations, we define the following two types of operators at each vertex and link:

$$V_{(x,y)} := X_{2,(x,y+1/2)} X_{2,(x,y-1/2)}^{\dagger} (X_{1,(x,y)}^{\dagger})^{2} \\ \times X_{1,(x+1,y)} X_{1,(x-1,y)},$$

$$P_{(x,y+1/2)} := Z_{1,(x,y+1)} Z_{1,(x,y)}^{\dagger} (Z_{2,(x,y+1/2)}^{\dagger})^{2} \\ \times Z_{2,(x-1,y+1/2)} Z_{2,(x+1,y+1/2)}.$$
(4)

These terms are portrayed in Figs. 1(b) and 1(c). We interchangeably call these two terms $V_{(x,y)}$ and $P_{(x,y+1/2)}$ vertex and *plaquette operators*, respectively. It is straightforward to check that each term in (4) commute, forming the stabilizer group. The Hamiltonian is defined by

$$H = -\sum_{(x,y)} V_{(x,y)} - \sum_{(x,y+1/2)} P_{(x,y+1/2)} + \text{H.c.}$$
(5)

This model shares the same feature as the \mathbb{Z}_N toric code [7]. The ground state is the stabilized state: the ground state $|\Omega\rangle$ satisfies

$$V_{(x,y)}|\Omega\rangle = P_{(x,y+1/2)}|\Omega\rangle = |\Omega\rangle \ \forall \ V_{(x,y)}, P_{(x,y+1/2)}.$$
 (6)

Also, our model admits two types of fractional excitations when the condition $V_{(x,y)} = 1$ or $P_{(x,y+1/2)} = 1$ is violated at a vertex or link. The crucial difference between our model and the toric code is that each terms in the Hamiltonian involves not only nearest-neighboring sites but also nextnearest-neighboring ones. As a consequence, the fractional excitations are subject to mobility constraint, giving rise to unusual GSD dependence on the system size. We dub the topological phases with this feature *higher-rank topological phases*. Throughout this paper, we consider the model placed on the torus geometry with length in the x (y) direction being n_x (n_y).

B. Simplest case: N = 2

Before discussing the model for generic case of N, we consider the simplest case by setting N = 2 to make a more

intuitive understanding of our model (5). The argument presented here will be useful to interpret our result on the entanglement entropy (Sec. IV). In this case, the terms defined in (4) become

$$V_{(x,y)} = X_{2,(x,y+1/2)} X_{2,(x,y-1/2)} X_{1,(x-1,y)} X_{1,(x+1,y)},$$

$$P_{(x,y+1/2)} = Z_{1,(x,y+1)} Z_{1,(x,y)} Z_{2,(x-1,y+1/2)} Z_{2,(x+1,y+1/2)}.$$
 (7)

One can regard each term as the plaquette term in a rhombus shape, as portrayed in Fig. 2(a). The Hamiltonian (5) with (7) resembles the \mathbb{Z}_2 toric code. However, there is a crucial difference between the two: each term (7) involves *next*-nearest-neighboring Pauli operators in the *x* direction, as opposed to the regular toric code where each term consists of nearest-neighboring Pauli operators. Due to this property, one can separate the mutually commuting terms (7) into two groups without considering the boundary:

(i)
$$\{V_{(2m,y)}, P_{(2m'-1,y'+1/2)}\},\$$

(ii) $\{V_{(2m-1,y)}, P_{(2m',y'+1/2)}\}\$ $(m,m'\in\mathbb{Z}).$ (8)

Hence, without thinking the boundary, the model that we describe with N = 2 amounts to be two decoupled \mathbb{Z}_2 toric codes. Now we impose the boundary condition on the lattice. We impose the periodic boundary condition in both of x and y directions, assuming the length of the lattice in the x direction is even. In this case, the Hamiltonian (5) consists of two decoupled \mathbb{Z}_2 toric codes on the torus. Such decomposition is portrayed in Fig. 2(b). In each \mathbb{Z}_2 toric code, there are two types of \mathbb{Z}_2 excitations: electric and magnetic charge. The nonlocal string of these charges yields logical operators, giving rise to fourfold GSD on the torus [7,25]. Therefore, the GSD of our model is found to be $4^2 = 16$. On the contrary, the situation is drastically different when the length of the lattice in the x direction is odd, instead of even. In this case, one cannot separate the terms (7) into two groups, thus the decomposition (8) is no longer true. Indeed, the terms belonging to (i) in (8) are "connected" with the ones belonging to (ii). For instance, the terms $V_{(n_x-1,y)}$ which belong to (i) in (8) and $V_{n_x+1,y} = V_{1,y}$ belonging to (ii) (equality follows from the periodic boundary condition) are located adjacent to each other. Therefore, the model is equivalent to one \mathbb{Z}_2 toric code on torus, giving GSD = 4.

To summarize the argument presented in this subsection, in the simplest case by setting N = 2, we learn that our model is identified as the \mathbb{Z}_2 toric code where each plaquette term is in a rhombus shape and that its GSD on torus drastically changes depending on the length of the lattice in the *x* direction, i.e.,

$$GSD = \begin{cases} 16 \ (n_x \text{ even}), \\ 4 \ (n_x \text{ odd}). \end{cases}$$
(9)

Such behavior can be understood by the fractional excitations with mobility constraint, analogously to the fracton topological phases. (See [19] for more details.) In the subsequent subsections, we turn to the generic cases of N and discuss its topological properties.

C. Fusion rules

To get a handle on our model (5), we also review the behavior of the fractional excitations via fusion rules [19]. The ground state $|\Omega\rangle$ of the model is the projected state, satisfying (6). When acting an operator $Z_{1,(x,y)}$ on the ground state, it violates the condition $V_{(x,y)} = 1$, namely,

$$V_{(x,y)}(Z_{1,(x,y)}|\Omega\rangle) = \omega^{-2}(Z_{1,(x,y)}|\Omega\rangle),$$

$$V_{(x\pm 1,x)}(Z_{1,(x,y)}|\Omega\rangle) = \omega(Z_{1,(x,y)}|\Omega\rangle)$$
(10)

giving rise to electric charges. This indicates that by acting on an operator $Z_{1,(x,y)}$, an electric charge is induced at the coordinate $(x \pm 1, x)$ and two conjugates of the electric charges are obtained at (x, y). Denoting the \mathbb{Z}_N eclectic charge as $e_{(x,y)}$ (and its conjugate as $\overline{e}_{(x,y)}$), (10) can be described by the following fusion rule:

$$I \rightarrow e_{(x-1,y)} \otimes \overline{e}_{(x,y)}^2 \otimes e_{(x+1,y)},$$
 (11)

which can be rewritten as

$$I \rightarrow (e_{(x-1,y)} \otimes \overline{e}_{(x,y)}) \otimes \overline{(e_{(x,y)} \otimes \overline{e}_{(x+1,y)})}$$
(12)

indicating that a pair of dipoles (i.e., the term in the big parentheses) of electric charges are created, which is distinct from conventional fusion rules where a pair of single charges is generated. Also, from the fusion rule (12), it can be shown that by applying a string of Z_1 operators $\prod_{a=1}^{N-1} (Z_{1,(x+a,y)})^a$ to the ground state, we have

$$I \to e_{(x,y)} \otimes \overline{e}_{(x+N,y)},$$
 (13)

indicating that an electric charge can hop in the *x* direction in the unit of *N* lattice sites rather than a single site. This corresponds to the fact that the U(1) dipole symmetry is reduced to \mathbb{Z}_N , as alluded to in [18,26,27]. Likewise, by acting on a single $Z_{2,(x,y+1/2)}$ on the ground state, we have electric charges subject to the following fusion rule:

$$I \to \overline{e}_{(x,y+1)} \otimes e_{(x,y)},\tag{14}$$

which is identical to the one in the conventional toric code. The fusion rule of the magnetic charges can be similarly discussed.

To recap the argument, while fusion rules of the charges in the y direction are the same as the conventional ones, we have an unusual fusion rule in the x direction, mirroring the fact that each stabilizer involves next-nearest-neighboring sites. By the fusion rules, one can evaluate the form of Wilson loops, which are noncontractible loops of fractional excitations (in the context of the stabilizer model, they are also known as logical operators), contributing to GSD. We defer the discussion on this point to Appendix A.

D. Logical operators and ground-state degeneracy

In this subsection, we discuss our model in the generic cases of *N*. As opposed to the case with N = 2, identifying the GSD and the logical operators is not so immediate in the generic case of *N*. Thus, we employ an alternative approach to accomplish this task. We adopt an algebraic tool of the graph theory, which will play a pivotal role in the subsequent discussion on the entanglement entropy. As we mentioned below (5), each term of the Hamiltonian involves the clock states in the next-nearest-neighbor as well as the nearest-neighbor ones. Such a feature can be succinctly described by the *Laplacian matrix* (*Laplacian*, in short), the graph theoretical analog of the second-order derivatives [28].

At each *y* on the lattice, we define the Laplacian *L* as the $n_x \times n_x$ matrix indexed by vertices (x, y) $(1 \le x \le n_x) \exists y$ running in the *x* direction, which has the following form:

$$L = \begin{pmatrix} 2 & -1 & & -1 \\ -1 & 2 & -1 & & \\ & -1 & 2 & \ddots & \\ & & \ddots & \ddots & -1 \\ -1 & & & -1 & 2 \end{pmatrix}.$$
 (15)

Equivalently,

$$(L)_{x,x'} = \begin{cases} 2 \ (x = x'), \\ -1 \ (x = x' \pm 1), \\ 0 \ (\text{else}), \end{cases}$$
(16)

where we have conventionally defined $x = n_x + 1 = 1$ due to the periodic boundary condition. The term $V_{(x,y)}$ at given y, constituting the Hamiltonian (4), now can be rewritten in terms of the matrix element of the Laplacian (16) via

$$V_{(x,y)} = X_{2,(x,y+1/2)} X_{2,(x,y-1/2)}^{\dagger} X_{1,(x,y)}^{-(L)_{x,x}} X_{1,(x+1,y)}^{-(L)_{x+1,x}} X_{1,(x-1,y)}^{-(L)_{x-1,x}},$$
(17)

where we conventionally set $X_{1,(x,y)}^{-1} = X_{1,(x,y)}^{\dagger}$. Likewise, each term of $P_{(x,y+1/2)}$ can be rewritten in the similar manner as

$$P_{(x,y)} = Z_{1,(x,y+1)} Z_{1,(x,y)}^{\dagger} Z_{2,(x,y+1/2)}^{-(L)_{x,x}} Z_{2,(x-1,y+1/2)}^{-(L)_{x-1,x}} Z_{2,(x+1,y+1/2)}^{-(L)_{x+1,x}}.$$
(18)

With this preparation, we now turn to counting the GSD on the torus geometry. The GSD is obtained by the number of total clock states divided by the number of independent stabilizers. The number of the clock states is found to be $N^{2n_xn_y}$. Also, the total number of the stabilizers (4) is given by $N^{2n_xn_y}$. However, we have overcounted the number of stabilizers; we need to take into account the constraint on the stabilizers as multiplication of some of stabilizers becomes identity. To find such constraint, we resort to the formalism of the Laplacian (15). At given *y*, introducing n_x -dimensional vector $\mathbf{r} := (r_1, \ldots, r_{n_x})^T \in \mathbb{Z}_N^{n_x}$, we consider the following product of the vertex operators in the *x* direction:

$$\prod_{x=1}^{n_x} V_{(x,y)}^{r_x}$$

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FIG. 3. (a) (Left) Configuration of the product of the operators $V_{(x,y)}$ (blue stars) given in (19), $\prod_{x=1}^{n_x} V_{(x,y)}^{r_x}$ with $\mathbf{s} = \mathbf{0}$. Since $\mathbf{s} = \mathbf{0}$, it does not have X_1 operators in the horizontal direction, leaving X_2 's in the vertical links (dots). (Middle) If we multiply the product (19) with the adjacent one in the *y* direction, some of X_2 's are canceled out. (Right) Iterating the similar procedure, one obtains the product of (19) all along in the *y* direction, $\prod_{y=1}^{n_x} (\prod_{x=1}^{n_x} V_{(x,y)}^{r_x})$, which is identity due to the periodic boundary condition. (b) (Top) Configurations of the logical operators constructed by noncontractible loops of X_1 or X_2 operators defined in (29) in the case of N = 3 and $n_x = 6$. (Bottom) Logical operators defined in (31) with N = 3 and $n_x = 6$. The periodic boundary condition is imposed so that the left and right edges are identified.

Referring to (17), we transform it by using the Laplacian L as

$$\prod_{x=1}^{n_x} V_{(x,y)}^{r_x} = \left[\prod_{x=1}^{n_x} \{ X_{2,(x,y+1/2)}^{\dagger} X_{2,(x,y-1/2)} \}^{r_x} \right] \\ \times \prod_{x=1}^{n_x} X_{1,(x,y)}^{s_x} (s_x \in \mathbb{Z}_N)$$
(19)

with

$$\mathbf{s} := \left(s_1, \ldots, s_{n_x}\right)^T = -L\mathbf{r}.$$
 (20)

Suppose $s = 0 \pmod{N}$. Then, multiplying the operators (19) along the *y* direction gives [see also Fig. 3(a)]

$$\prod_{y=1}^{n_y} \left(\prod_{x=1}^{n_x} V_{(x,y)}^{r_x} \right) = \prod_{y=1}^{n_y} \left[\prod_{x=1}^{n_x} \{ X_{2,(x,y+1/2)}^{\dagger} X_{2,(x,y-1/2)} \}^{r_x} \right] = I,$$
(21)

where the last equation holds due to the periodic boundary condition. Therefore, to find constraints of the stabilizers, we need to evaluate the solution of

$$L\mathbf{r} = \mathbf{0} \mod N,\tag{22}$$

i.e., the kernel of the Laplacian.

To proceed, we transform the Laplacian into the diagonal form, known as the *Smith normal form*. Introducing invertible integer matrices P and Q, one can transform the Laplacian (15) as

$$PLQ = D, D = \text{diag}(\underbrace{1, \dots, 1}_{n_x - 2}, n_x, 0).$$
 (23)

With this transformation, one can rewrite (22) as

$$(22) \Leftrightarrow P^{-1}DQ^{-1}\mathbf{r} = \mathbf{0} \mod N$$

 $\Leftrightarrow D\tilde{\mathbf{r}} = \mathbf{0} \mod N. \tag{24}$

When moving from the second to the third equation, we have used the fact that *P* is the integer matrix. Also, we have defined $\tilde{\mathbf{r}} := Q^{-1}\mathbf{r}$. From (23) and (24), it follows that

$$\tilde{r}_i = 0 \ (1 \le i \le n_x - 2), \ n_x \tilde{r}_{n_x - 1} = 0 \mod N.$$
 (25)

The last entry of the vector \tilde{r} , \tilde{r}_{n_x} takes N distinct values since the last diagonal element of the matrix D is zero [Eq. (23)].

Taking this fact as well as (25) into account, we have

$$\widetilde{\mathbf{r}} = (\underbrace{0\cdots,0}_{n_x-2}, N'\alpha_1, \alpha_2)^T, \ N' := \frac{N}{\gcd(N, n_x)},$$
$$(\alpha_1 \in \mathbb{Z}_{\gcd(N, n_x)}, \ \alpha_2 \in \mathbb{Z}_N).$$
(26)

Here, gcd stands for the greatest common divisor. By evaluating the form of the matrix Q [19], one finds

$$\mathbf{r} = Q\tilde{\mathbf{r}} = N'\alpha_1 \begin{pmatrix} n_x - 1\\ n_x - 2\\ \vdots\\ 1\\ 0 \end{pmatrix} + \alpha_2 \begin{pmatrix} 1\\ 1\\ \vdots\\ 1\\ 1 \end{pmatrix} \mod N.$$
(27)

Therefore, there are $gcd(N, n_x) \times N$ constraints on the stabilizers $V_{(x,y)}$.

One can analogously discuss the number of the constraints on the stabilizers $P_{(x,y+1/2)}$, yielding that there are gcd $(N, n_x) \times N$ constraints. Overall, there are [gcd $(N, n_x) \times N$]² constraints on the stabilizers, hence, the GSD is given by

$$GSD = \frac{N^{2n_x n_y}}{N^{2n_x n_y} / [\gcd(N, n_x) \times N]^2} = [\gcd(N, n_x) \times N]^2.$$
(28)

It is emphasized that our model exhibits unusual GSD dependence on the system size which was not seen in fracton topological phases which exhibit the subextensive GSD. Also, it is known that the Wen's \mathbb{Z}_2 plaquette model exhibits the similar gcd dependence of the GSD [29]. However, our model discussed in this paper is qualitatively different from the one in [29] as our model admits the dipole of the fractional charges due to the multipole symmetries.

One can make use of the formalism of the Laplacian to identify the form of the logical operators. By evaluating the kernel and cokernel of the Laplacian (see [19] for derivation and Appendix A), one can introduce logical operators consisting of noncontractible loops of X_1 or X_2 operators as

$$\eta_{1}^{x} = \prod_{x=1}^{n_{x}} X_{2,(x,y+1/2)}^{p_{x}}, \quad \gamma_{1}^{x} = \prod_{x=1}^{n_{x}} X_{2,(x,y+1/2)}^{q_{x}},$$

$$\eta_{2}^{x} = \left(\prod_{y=1}^{n_{y}} X_{1,(n_{x}-1,y+1/2)}\right) \times \left(\prod_{y=1}^{n_{y}} X_{1,(n_{x},y+1/2)}^{\dagger}\right),$$

$$\gamma_{2}^{x} = \prod_{y=1}^{n_{y}} X_{1,(n_{x},y+1/2)}$$
(29)

with

$$\mathbf{p} = (p_1, p_2, \dots, p_{n_x})^T = N' \alpha_1 \begin{pmatrix} n_x - 1 \\ n_x - 2 \\ \vdots \\ 1 \\ 0 \end{pmatrix},$$
$$q = (q_1, q_2, \dots, q_{n_x})^T = \alpha_2 \begin{pmatrix} 1 \\ 1 \\ \vdots \\ 1 \\ 1 \end{pmatrix} \mod N.$$
(30)

We portray them in Fig. 3(b) in the case of N = 3 and $n_x = 6$. These operators are represented as η_m^x , γ_m^x (m = 1, 2), where two types of the loop are labeled by η and γ , whereas the index m denotes the direction of the loops (x or y). In addition to the logical operators, γ_1^x , γ_2^x that one can find in the toric code, there are new forms of the logical operators in our model, η_1^x , η_2^x . In the horizontal direction, η_1^x is formed by the inhomogeneous string of X_2 's whereas in the vertical direction, the model admits the "dipole of the loops," i.e., a pair of the noncontractible loops of X_1 's in the vertical direction with opposite sign located adjacent to each other, giving η_2^x . The logical operators η_1^x and γ_1^x can be deformed upward or downward in the y direction and any logical operator can be generated by the combination of η_1^x and γ_1^x . Also, any logical operator consisting of the product X_2 's can be generated by η_2^x and γ_2^x . In Appendix A, we give more details of these logical operators by investigating the behavior of the fractional charges.

Analogously to (29), logical operators involving Z_1 or Z_2 operators are defined by [Fig. 3(b)]

$$\eta_{1}^{z} = \prod_{x=1}^{n_{x}} Z_{1,(x,y)}^{p_{x}}, \ \gamma_{1}^{z} = \prod_{x=1}^{n_{x}} Z_{1,(x,y)}^{q_{x}},$$
$$\eta_{2}^{z} = \left(\prod_{y=1}^{n_{y}} Z_{2,(n_{x}-1,y)}\right) \left(\prod_{y=1}^{n_{y}} Z_{2,(n_{x},y)}^{\dagger}\right),$$
$$\gamma_{2}^{z} = \prod_{y=1}^{n_{y}} Z_{2,(n_{x},y)},$$
(31)

where the integers p_x and q_x are determined by (30). One can check that

$$\eta_{1}^{x}\eta_{2}^{z} = \omega^{N'}\eta_{2}^{z}\eta_{1}^{x}, \ \gamma_{1}^{x}\gamma_{2}^{z} = \omega\gamma_{2}^{z}\gamma_{1}^{x}, \ \eta_{2}^{x}\eta_{1}^{z} = \omega^{N'}\eta_{1}^{z}\eta_{2}^{x},$$
$$\gamma_{2}^{x}\gamma_{1}^{z} = \omega\gamma_{1}^{z}\gamma_{2}^{x}, \ (\eta_{1}^{x})^{\text{gcd}(N,n_{x})} = (\eta_{2}^{x})^{\text{gcd}(N,n_{x})}$$
$$= (\eta_{1}^{z})^{\text{gcd}(N,n_{x})} = (\eta_{2}^{z})^{\text{gcd}(N,n_{x})} = I,$$
$$(\gamma_{1}^{x})^{N} = (\gamma_{2}^{x})^{N} = (\gamma_{1}^{z})^{N} = (\gamma_{2}^{z})^{N} = I$$
(32)

with commutation relation between other combinations of the logical operators being trivial. Logical operators with the commutation relation (32) yield the GSD $[gcd(N, n_x) \times N]^2$.

Ground states with such GSD can be constructed by the logical operators. Labeling stabilizers involving X_1 or X_2 operators that are obtained by the product of vertex operators $V_{(x,y)}$ as

$$G = \left\{ g | g \in \prod_{x,y} V_{(x,y)}^{a_{x,y}}, a_{x,y} \in \mathbb{Z}_N \right\},$$

we define the following ground state:

$$|\psi\rangle := \frac{1}{\sqrt{|G|}} \sum_{g \in G} g|0\rangle, \tag{33}$$

where $|0\rangle$ is the trivial state in the diagonal basis of the Z_1 and Z_2 operators, satisfying $Z_{1,(x,y)}|0\rangle = Z_{2,(x,y+1/2)}|0\rangle = |0\rangle$ (the generalization of the "spin-up state" to the clock state). Also, |G| denotes the total number of the stabilizers *G*. One can

verify that this state is the stabilized state. The $[gcd(N, n_x) \times N]^2$ ground states are labeled by

$$|\xi_{ab,cd}\rangle := (\eta_1^x)^a (\gamma_1^x)^b (\eta_2^x)^c (\gamma_2^x)^d |\psi\rangle [0 \le a, c \le \gcd(N, n_x)]$$

$$-1, 0 \leqslant b, d \leqslant N - 1]. \tag{34}$$

It is known that the number of distinct fractional excitations in a topological phase on torus geometry is identical to GSD [30]. Moreover, in our model, all of the fractional excitations are Abelian, implying that quantum dimension d_a of any excitation is one. Taking these into account, one can evaluate the total quantum dimension as

$$\sqrt{\sum_{a} d_a^2} = N \times \gcd(N, n_x), \tag{35}$$

where we have used the fact that there are $[N \times \text{gcd}(N, n_x)]^2$ distinct fractional excitations, all of which carry quantum dimension one.

In the next section, we calculate the entanglement entropy of our model with respect to the ground states. In particular, we investigate whether the total quantum dimension (35)enters in the form of the entanglement entropy in the disk geometry.

III. ENTANGLEMENT ENTROPY

Now we come to the main part of this paper. In this section, we study entanglement entropy of our model defined in the previous section. Our calculations rely on the formulation of the entanglement entropy in the stabilizer codes, invented in [22].

A. Review of the bipartite entanglement in lattice spin systems

Before we systematically discuss the entanglement entropy of the higher-rank topological phases, let us review the formalism of the bipartite entanglement in lattice spin systems proposed in [22]. Readers familiar with this formalism may skip this subsection.

1. Reduced density matrix

For a qubit system, the Hilbert space is $\mathscr{H} = \bigotimes_j \mathscr{H}_j$, where $\mathscr{H}_j = \operatorname{span}\{|0\rangle_j, |1\rangle_j, \ldots, |N-1\rangle_j\}$ in the σ_j^z basis, where σ_j^z is generalized Pauli matrix in the *N*-clock model. We define a reference state $|0\rangle := \bigotimes_j |0\rangle_j$, namely, all-spin-up state. We define $E := \bigotimes_j \{1, \sigma_j^x, \ldots, (\sigma_j^x)^{N-1}\}$ to be the Abelian group that rotates spins in an onsite manner. Any state in the Hilbert space can be written as $g|0\rangle$ for some $g \in E$ with $g^N = 1$. A generic state can be written as

$$|\psi\rangle = \sum_{g \in G \subset E} a_g g |0\rangle, \ \sum_g |a_g|^2 = 1.$$
(36)

Here G is a subgroup of E and a_g is the wave function of $|\psi\rangle$ in the computational basis, namely, $a_g = \langle 0|g|\psi\rangle$. Given the state $|\psi\rangle$, we can define the corresponding density matrix as

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$$p_{\psi} := |\psi\rangle\langle\psi| = \sum_{g,g'} a_g \bar{a}_{g'} g|0\rangle\langle0|g'^{\dagger}$$
$$= \sum_{g,g'} a_g \bar{a}_{gg'} g|0\rangle\langle0|(gg')^{\dagger}, \qquad (37)$$

where in the last equality we redefined $g' \rightarrow gg'$. Once a bipartition of the lattice into subsystem *A* and its complement *B*, any element *g* can be unambiguously decomposed into the product form $g = g_A \otimes g_B$, where $g_{A/B}$ only acts nontrivially on subsystems *A* and *B*, respectively. With the decomposition of the reference state $|0\rangle = |0\rangle_A \otimes |0\rangle_B$, we obtain the reduced density matrix of subsystem *A* by tracing over subsystem *B* as

$$\rho_{A} = \sum_{n} {}_{B} \langle n|\rho|n\rangle_{B} = \sum_{n} \sum_{g,g'} a_{g} \bar{a}_{gg'} g_{A}|0\rangle_{AA} \langle 0|(g_{A}g'_{A})^{\dagger}{}_{B}$$

$$\times \langle n|g_{B}|0\rangle_{BB} \langle 0|(g_{B}g'_{B})^{\dagger}|n\rangle_{B}$$

$$= \sum_{g,g'} a_{g} \bar{a}_{gg'} g_{A}|0\rangle_{AA} \langle 0|(g_{A}g'_{A})^{\dagger}{}_{B} \langle 0|g'^{\dagger}_{B}|0\rangle_{B}, \qquad (38)$$

where we used the completeness relation $\sum_{n} |0\rangle_{BB} \langle 0| = I$. It is easy to see that nonzero contributions only come from $g'_{B} = I_{B}$. Define

$$G_A := \{g \in G | g = g_A \otimes I_B\}, \ G_B := \{g \in G | g = I_A \otimes g_B\}.$$
(39)

Then $g'_B = I_B$ implies $g' \in G_A$. The reduced density matrix is rewritten as

$$\rho_A = \sum_{g \in G, g' \in G_A} a_g \bar{a}_{gg'} g_A |0\rangle_{AA} \langle 0| (g_A g'_A)^{\dagger}.$$
(40)

The equivalences between the following statements are proven [22]: (1) ρ_A is diagonal, (2) $G_A = \{1\}$, (3) $\nexists g = g_A g_B$ with both $g_A \in G_A$, $g_B \in G_B$ nontrivial for $g \in G$.

2. Entanglement entropy of a subsystem

We will study the entanglement entropy of subsystem A:

$$S_A = -\mathrm{tr}\,\rho_A \mathrm{log}\,\rho_A. \tag{41}$$

Without loss of generality, we can assume the total density matrix ρ is a pure state. Let *S* be the stabilizer group, which contains mutually commuting operators formed from *G* that are called *stabilizers*. The codeword space is defined as $\mathscr{L} = \{|\psi\rangle \in \mathscr{H} | U | \psi \rangle = |\psi\rangle, \forall U \in S\}$. Let us take $|\psi\rangle = \frac{1}{\sqrt{|G|}} \sum_{g \in G} g |0\rangle$ for the moment. We will study the generic ground states in the Appendix C. Apparently, $|\psi\rangle; \in \mathscr{L}$ as $g | \psi \rangle = |\psi\rangle, \forall g \in G$. From (40), we have

$$\rho_A = \frac{1}{|G|} \sum_{g \in G, g' \in G_A} g_A |0\rangle_{AA} \langle 0| (g_A g'_A)^{\dagger}.$$
(42)

Let G/G_B and $G_{AB} := G/(G_A \cdot G_B)$ be quotients that contain elements that act freely on A and $A \cup B$, respectively. It is shown that for the ground state (42), the entanglement entropy is described by [22]

$$S_A = \log |G_{AB}|. \tag{43}$$

The details are given in Appendix **B**. The entanglement entropy of the subsystem A is given by

$$S_A = \log |G| - \log |G_A| - \log |G_B|,$$
 (44)

where |G| denotes the total number of independent stabilizers containing X_1 and X_2 , whereas $|G_A|$ ($|G_B|$) represents the number of stabilizers containing X_1 and X_2 that have support on A (B).



FIG. 4. (a) The subsystem A is a single row on the horizontal sites (clock states inside the red frame). (b) The subsystem A is a single row on the vertical links (red dots inside the red frame). The blue triangles represent one of the n_x constraints along the horizontal direction. (c) The subsystem A is a single column (red dots and squares inside the frame). The blue triangles represent one of the n_y constraints along the vertical sites. (d) Multiplication of the vertex operators (represented by black stars) that act within B. (e) Such product of the stabilizers is redundant due to the constraint that the product of all of the vertex operators becomes identity (52).

In the following subsections, we study the entanglement entropy of various geometries based on the formula (44).

B. Single-row and -column geometry

After reviewing formulation of the entanglement entropy, we calculate it in various geometries of subsystems in our model. We evaluate the entanglement entropy with respect to the ground state $|\xi_{00,00}\rangle$ given in (34). One could study the entanglement entropy for more generic ground state $|\zeta\rangle$ which is superposition of $|\xi_{ab,cd}\rangle$, defined by

$$\begin{aligned} |\zeta\rangle &= \sum_{a,c=0}^{\gcd(N,n_x)-1} \sum_{b,d=0}^{N-1} \alpha_{ab,cd} |\xi_{ab,cd}\rangle \\ &\times \left(\sum_{abcd} |\alpha_{ab,cd}|^2 = 1, \; \alpha_{ab,cd} \in \mathrm{U}(1) \right). \end{aligned}$$
(45)

In this case, depending on the geometry of the subsystem *A*, there is an additional contribution to (44). While we focus on the entanglement entropy with respect to the ground state $|\xi_{00,00}\rangle$, we make a brief comment on the case of the generic ground state (45), deferring the details to Appendix C.

In this subsection, we consider the single-row or-column subsystem geometries which go around the torus in the x or y direction, as shown in Fig. 4. When calculating the entanglement entropy based on (44), there is a technical caveat: when evaluating $|G_A|$ and $|G_B|$, we need to take into account

the product of vertex operators which have support only on *A* and *B*.

1. Single row I

Let us first look at the subsystem A in Fig. 4(a): we have $|G| = N^{n_x n_y} / \Gamma$, $|G_A| = 1$, $|G_B| = N^{n_x n_y - n_x}$, where the factor Γ corresponds to the two independent constraints from (27), $\Gamma := N \times \text{gcd}(N, n_x)$. Thus, (44) gives us

$$S_A = n_x \log N - \log N - \log[\gcd(N, n_x)]$$

= $n_x \log N - \log \Gamma$. (46)

For the generic ground state (45), there is an additional constant μ_1 in (46):

$$S_A = n_x \log N - \log \Gamma + \mu_1,$$

$$\mu_1 = -\sum_{c,d} \left[\left(\sum_{a,b} |\alpha_{ab,cd}|^2 \right) \log \left(\sum_{a,b} |\alpha_{ab,cd}|^2 \right) \right]. \quad (47)$$

Derivation of (47) is given in Appendix C.

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2. Single row II

For the subsystem A in Fig. 4(b), we have

$$|G_A| = 1, |G_B| = (|G|/N^{2n_x})N^{n_x} = |G|/N^{n_x},$$

When evaluating $|G_B|$, the N^{n_x} in the first equality comes from the n_x -independent constraints, one of which is labeled by the

blue triangles in Fig. 4(b). Following (44), we have

$$S_A = n_x \log N. \tag{48}$$

In the case of the generic ground state (45), we have an additional contribution to (48):

$$S_A = n_x \log N + \mu_2, \quad \mu_2 = -\sum_{p=0}^{\gcd(N,n_x)-1} \sum_{q=0}^{N-1} \lambda_{p,q} \log \lambda_{p,q}$$
(49)

with $(v = e^{2\pi i / \gcd(N, n_x)}, \omega = e^{2\pi i / N})$

$$\lambda_{p,q} = \frac{1}{\Gamma} \sum_{k=0}^{\gcd(N,n_x)-1} \sum_{l=0}^{N-1} \nu^{pk} \omega^{ql} \sigma_{kl},$$

$$\sigma_{kl} = \sum_{\substack{a,a',b,b',c,d \\ a-a'=k \mod \gcd(N,n_x) \\ b-b'=l \mod N}} \alpha_{ab,cd} \bar{\alpha}_{a'b',cd}.$$
(50)

The derivation of this constant is relegated to Appendix C.

3. Single column

Let us turn to subsystem A in Fig. 4(c). The total number of the independent vertex operators is given by $|G| = N^{n_x n_y} / \Gamma$, where we take the constraint discussed in the previous section into consideration (Sec. II D). Next we evaluate $|G_B|$. Naively, the number of stabilizers that act within B is given by

$$N^{n_x n_y - 3n_y}.$$

However, this counting is incorrect, as one has to take into account the multiplication of the stabilizers that act trivially on *A*. There are two types of such product. The first type is the product of the vertex operators on both sides of the vertical line. One of such configuration is indicated by the blue triangles in Fig. 4(c). There are N^{n_y} such products. Another type is multiplication of the vertex operators which are located inside *A* and the ones surrounding *A*, as shown in black stars in Fig. 4(d). Having identified these products of the stabilizers that contribute to $|G_B|$, one has to check whether these products are "redundant" to the constraint of the stabilizes that we have already considered.

To be more specific to what we have just said, we look at the second type of the product of the stabilizers that are depicted in Fig. 4(d). We schematically write this product as $\prod_{\bigstar} V_{(x,y)}$ in accordance with Fig. 4(d). As we discussed in Sec. II D, the multiplication of all of the vertex operators becomes identity, i.e.,

$$\prod_{\blacklozenge} V_{(x,y)} \times \prod_{\bigstar} V_{(x,y)} = I,$$
(52)

where $\prod_{\mathbf{A}} V_{(x,y)}$ denotes the product of the vertex operators defined on the complement of \mathbf{A} [see Fig. 4(e)]. Hence, the product in question, $\prod_{\mathbf{A}} V_{(x,y)}$, is redundant: it can be generated by the vertex operators belonging to *B*. Likewise, regarding the first type of the product, one of which is depicted in Fig. 4(c), combination of some of them is redundant due to the constraint. Overall, the multiplication of the vertex operators that act on *B* is found to be

$$\frac{N^{n_y} \times N}{[\gcd(N, n_x) \times N]}.$$
(53)

Therefore, $|G_B|$ is obtained by multiplying this number with (51), that is,

$$|G_B| = N^{n_x n_y - 3n_y} \times \frac{N^{n_y} \times N}{[\gcd(N, n_x) \times N]}.$$
 (54)

There is no stabilizer that acts within *A*, giving $|G_A| = 1$. Referring to (44), the entanglement entropy is given by

$$S_A = 2n_v \log N - \log N. \tag{55}$$

In the case of the generic ground state (45),

$$S_A = 2n_y \log N - \log N + \mu_3, \quad \mu_3 = -\sum_{b=0}^{N-1} \sum_{q=0}^{N-1} \lambda_q^{(b)} \log \lambda_q^{(b)},$$
(56)

where

$$\lambda_{q}^{(b)} = \frac{1}{N} \sum_{k=0}^{N-1} \omega^{kq} \sigma_{k}^{(b)}, \quad \sigma_{k}^{(b)} = \sum_{\substack{a,c,d,d'\\d-d'=k \mod N}} \alpha_{ab,cd} \bar{\alpha}_{ab,cd'}.$$
(57)

The derivation is given in Appendix C.

C. Disk geometry

Let us calculate entanglement of the contractible disk geometry, as portrayed in Fig. 5(a). We assume that the width of the disk is more than one. We set the width and height of the disk to be $l_x \ge 2$ and l_y , respectively. Accordingly, the coordinate of \mathbb{Z}_N clock states with type 1 (i.e., clock states located at vertices) inside the disk is denoted as (x, y) ($x_0 \le x \le x_0 + l_x - 1, y_0 \le y \le y_0 + l_y - 1$) and the ones with type 2 (i.e, clock states at vertical links) inside the disk as (x, y + 1/2) ($x_0 \le x \le l_x - 1 + x_0, y_0 \le y \le y_0 + l_y - 2$). See also Fig. 5(a).

In this setting, we have $|G| = N^{n_x n_y} / \Gamma$, $|G_A| = N^{(l_x-2)(l_y-2)}$. As for $|G_B|$, the number of individual stabilizers that have support only on *B* is

$$N^{n_x n_y} / N^{l_y (l_x + 2)}.$$
 (58)

In addition to this number, we need to take the product of the vertex operators that act within *B* into consideration. To find such product, it is useful to resort to the formalism of the Laplacian that we have discussed in Sec. II D. Consider the following product of $V_{(x,y)}$'s in the horizontal direction at given y ($y_0 \le y \le y_0 + l_y - 1$) [purple triangles in Fig. 5(a)]:

$$\prod_{x=x_0}^{x_0+l_x-1} V_{(x,y)}^{t_x} \quad (t_x \in \mathbb{Z}_N),$$
(59)

which can be rewritten as

$$(59) = \left[\prod_{x=x_0}^{x_0+l_x-1} \{X_{2,(x,y+1/2)}^{\dagger}X_{2,(x,y-1/2)}\}^{l_x}\right] \\ \times \prod_{x=1}^{n_x} X_{1,(x,y)}^{u_x} \ (u_x \in \mathbb{Z}_N).$$
(60)



FIG. 5. (a) Disk geometry. The clock states inside the disk are marked by red squares and dots. Purple triangles represent the product given in (59). (b) Decomposition of the Laplacian into submatrices, corresponding to (62) and (66). We need to find multiplication of the vertex operators corresponding to the vector **t** which does not have X_1 operators inside the disk, yielding the condition (65). (c) Configuration of the product given in (64) (purple triangles). With the condition (63), it exclusively acts on *B*. (d) The number of the clock states marked by light-blue squares gives rise to the area-law term (Area), given in (73), which is (Area) = $4l_y + 2(l_x - 2)$ in the present case.

Analogously to the discussion in the previous section (Sec. II D), \mathbb{Z}_N numbers t_x and u_x that appear in (59) and (60) are related via the Laplacian. Defining $(l_x + 2)$ - and n_x -dimensional vectors by $\mathbf{t} := (\underbrace{t_{x_0-1}, \ldots, t_{x_0+l_x}}_{l_x+2})^T$ and $\mathbf{u} :=$

 $(\underbrace{u_1,\ldots,u_{n_x}}_{n_x})^T$, respectively, and referring to (15) and (16), we

have

$$\mathbf{u} = -\tilde{L}\mathbf{t}.\tag{61}$$

Here, \tilde{L} represents a submatrix of the Laplacian L (16). It is obtained by decomposing the Laplacian into three matrices as

$$L = (L' \quad \tilde{L} \quad L''), \tag{62}$$

where L', \tilde{L} , L'' is $n_x \times (x_0 - 2)$, $n_x \times (l_x + 2)$, $n_x \times (n_x - x_0 - l_x)$ submatrix, respectively, and concentrating on the middle matrix \tilde{L} [see also Fig. 5(b)].

Suppose the product we consider, Eq. (60) with (61), has no X_1 operators within the disk A, that is,

$$\mathbf{u} = (u_1, \dots, u_{x_0-1}, \underbrace{0, \dots, 0}_{l_x}, u_{x_0+l_x}, \dots, u_{n_x})^T.$$
(63)

Then, multiplying the product (60) along the y direction gives

$$\prod_{y=y_0}^{y_0+l_y-1} \left[\prod_{x=x_0}^{x_0+l_x-1} V_{(x,y)}^{t_x} \right]$$
(64)

which yields the stabilizers that have support only on *B* [see Fig. 5(c)]. Therefore, to find the product of stabilizers that acts within *B*, we need to evaluate (61) with (63). Such condition can be rewritten as

$$\tilde{L}_{\rm sub}\mathbf{t}=\mathbf{0},\tag{65}$$

where \tilde{L}_{sub} denotes the submatrix of \tilde{L} , which is obtained by decomposing the matrix \tilde{L} into three via

$$\tilde{L} = \begin{pmatrix} \tilde{L}_1 \\ \tilde{L}_{\text{sub}} \\ \tilde{L}_2 \end{pmatrix}, \tag{66}$$

where \tilde{L}_1 , \tilde{L}_{sub} , and \tilde{L}_2 is the $(x_0 - 1) \times (l_x + 2)$, $l_x \times (l_x + 2)$, and $(n_x - l_x - x_0 + 1) \times (l_x + 2)$ matrix, respectively, and focusing on the middle matrix \tilde{L}_{sub} [Fig. 5(b)]. From (15), (62), and (66), the explicit form of \tilde{L}_{sub} is given by

$$\tilde{L}_{\rm sub} = \begin{pmatrix} -1 & 2 & -1 & & & \\ & -1 & 2 & -1 & & & \\ & & -1 & 2 & \ddots & & \\ & & & \ddots & \ddots & -1 & \\ & & & & -1 & 2 & -1 \end{pmatrix}.$$
 (67)

Introducing $l_x \times l_x$ and $(l_x + 2) \times (l_x + 2)$ invertible matrices, \tilde{P} and \tilde{Q} , whose matrix elements are integers, one can transform the matrix \tilde{L}_{sub} into a diagonal form (i.e., the *Smith*

$$\tilde{D} = \begin{pmatrix} 1 & & & & & \\ & 1 & & & & \\ & & 1 & & & \\ & & \ddots & & & \\ & & & & 1 & 0 & 0 \end{pmatrix}.$$
(68)

The $l_x \times (l_x + 2)$ matrix \tilde{D} contains diagonal entries diag $(\underbrace{1, \ldots, 1}_{l_x})$ with other entries being zero. With this form,

it follows that

$$\tilde{L}_{\rm sub}\mathbf{t} = \mathbf{0} \Leftrightarrow \tilde{D}\tilde{\mathbf{t}} = \mathbf{0} \mod N,\tag{69}$$

where $\tilde{\mathbf{t}} := \tilde{Q}^{-1}\mathbf{t}$ and $\mathbf{0}$ represents l_x -dimensional zero vector. We rename the entries of the vector $\tilde{\mathbf{t}}$ as $\tilde{\mathbf{t}} = (\tilde{t}_1, \dots, \tilde{t}_{l_x+2})^T$. From the Smith normal form (68) and (69), the first l_x entries of the vector $\tilde{\mathbf{t}}$ are subject to

$$\tilde{t}_i = 0 \ (1 \leqslant i \leqslant l_x), \tag{70}$$

whereas there is no constraint on the last two entries \tilde{t}_{l_x+1} , \tilde{t}_{l_x+2} , giving

$$\tilde{t}_{l_x+1} = \beta_1, \ \tilde{t}_{l_x+2} = \beta_2 \ (\beta_1, \beta_2 \in \mathbb{Z}_N).$$
 (71)

Hence, there are N^2 solutions of (65).

One has to check that such N^2 product of vertex operators which have support only on *B* is redundant to the constraint of the stabilizers, analogously to the discussion presented around (52). Thus, the number of products of the stabilizers that act within *B* is given by $\frac{N^2}{r}$.

Overall, we obtain $|G_B|$ by multiplying this number with (58):

$$|G_B| = \frac{N^{n_x n_y}}{N^{l_x (l_y + 2)}} \times \frac{N^2}{\Gamma}.$$
 (72)

From (44), we finally arrive at

$$S_A = (\text{Area})\log N - 2\log N. \tag{73}$$

Here, (Area) is the number of vertex operators that have support on both A and its complement B, which can be regarded as the number of clock states with type 1 (i.e., clock states located at vertices) surrounding the disk geometry A [see Fig. 5(d)]. In the present case, it is given by (Area) = $4l_y + 2(l_x - 2)$. One can check that the entanglement entropy of the disk geometry for the generic ground state (45) gives the same answer as (73).

The first term in (73) is nothing but the area-law term and the second term, which is the subleading term of the entanglement entropy, is of particular importance. It is "topological"; one can confirm that the second term remains the constant when we deform the shape of the disk geometry retaining its topology. Therefore, we have found the same scaling behavior of the entanglement entropy as the one obtained in conventional topologically ordered phases (1). However, there is a crucial deference between the two. In the case of the conventional topologically ordered phases, the topological entanglement entropy γ is related to the total quantum dimension via $\gamma = \log \sqrt{\sum_a d_a^2}$, where *a* labels the distinct types of fractional excitations. On the contrary, in our case,



FIG. 6. Example of the tripartition ABC. The clock states belonging to A, B, and C are marked by red, blue, and green color, respectively.

such a relation does not hold as the number of distinct number of fractional excitations varies depending on the system size [Eq. (35)] and the topological entanglement entropy takes a constant number $\gamma = -2 \log N$ irrespective to the system size, implying the incapability to associate γ to the total quantum dimension. In the next section, we give a physical interpretation of this result in the simplest setting, i.e., N = 2. We summarize the results obtained in this section in Table I.

D. Topological entanglement entropy

In this subsection, we resort to the prescription to suppress the area-law term of the entanglement entropy to extract the topological subleading term in our model. It was found that in 2D topologically ordered phases, for a tripartite disk geometry *ABC*, the topological entanglement entropy is given by [20]

$$S_{\rm top} = S_A + S_B + S_C + S_{ABC} - S_{AB} - S_{AC} - S_{BC}.$$
 (74)

The sign in front of each term is designed so that the area-law term cancels out. To carry out this prescription in our model (5), we envisage the disk geometry *ABC* in the square lattice as portrayed in Fig. 6. We set the height of the disk *A* and the one of the *B* and *C* to be l_{1y} and l_{2y} . Here, the height of the disk represents the number of pairs of a clock state on a vertex and the one on a link just above it along each column inside the disk. (In Fig. 6, we portray a geometry with $l_{1y} = 3$, $l_{l_2y} = 4$.) We also set the width of the disks *B* and *C* to be l_{1x} and l_{2x} , respectively (width denotes the number of the clock states along each row inside the disk), with $l_x = l_{1x} + l_{2x}$, where l_x is the width of the disk *A*.

In this configuration, recalling that the entanglement entropy of a disk geometry is given by (73), where the (Area) denotes the number of stabilizers $V_{(x,y)}$ that act on both the clock states inside the disk and the ones in its complement



FIG. 7. (a) In the case of n_x being even. One can decompose the lattice into two. The periodic boundary condition is imposed so that the left and right edges are identified. (b) In the case of n_x being odd. We double the lattice by going around the torus twice, and omit vertices and links as explained in the main text. The red frame corresponds to the subsystem A with disk shape considered in the previous section.

around the border, we have

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$$S_{A} = (2l_{x} + 4l_{1y} - 2) \log N - 2 \log N,$$

$$S_{B} = (2l_{1x} + 4l_{2y} - 2) \log N - 2 \log N,$$

$$S_{C} = (2l_{2x} + 4l_{1y} - 2) \log N - 2 \log N,$$

$$S_{AB} = (2l_{x} + 4l_{1y} + 4l_{2y} - 3) \log N - 2 \log N,$$

$$S_{AC} = (2l_{x} + 4l_{1y} + 4l_{2y} - 3) \log N - 2 \log N,$$

$$S_{BC} = (2l_{x} + 4l_{2y} - 2) \log N - 2 \log N,$$

$$S_{ABC} = (2l_{x} + 4l_{1y} + 4l_{2y} - 2) \log N - 2 \log N,$$

$$S_{ABC} = (2l_{x} + 4l_{1y} + 4l_{2y} - 2) \log N - 2 \log N.$$
(75)

Therefore, the topological entanglement entropy is given by

$$S_{\rm top} = -2 \, \log N. \tag{76}$$

IV. ALTERNATIVE INTERPRETATION OF THE RESULT IN THE SIMPLEST CASE

In this section, we give an intuitive understanding of the topological entanglement entropy found in the second term of (73) by focusing on the simplest case of N, i.e., N = 2. To this end, discussion presented in Sec. II B is of usefulness.

Recalling the argument in Sec. II B, in the case of N = 2, depending on whether the length of the lattice in the *x* direction is even or odd, the GSD, accordingly, the total quantum dimension changes whereas the topological entanglement entropy gives $-2 \log 2$. Even though we have the topological entanglement entropy $-2 \log 2$ irrespective to the length of the lattice in the *x* direction being even or odd, its topological origin to have such a number is rather different in these two cases.

As seen from (18), in the case of N = 2 and n_x being even, the mutually commuting terms of Hamiltonian get simplified, allowing us to decompose the Hamiltonian into two sectors, each of which describes the \mathbb{Z}_2 toric code. Such decomposition is shown in Fig. 7(a) where we separate vertices with even number *x* coordinate (white squares) and the ones with odd number (black squares) as well as vertical links with odd number *x* coordinate (white dots) and the ones with even number (black dots). Correspondingly, the entanglement entropy S_A is decomposed into two, each of which is the one of the disk geometry in the \mathbb{Z}_2 toric code. Since the topological entanglement entropy of the disk geometry in the \mathbb{Z}_2 toric code is known to be $-\log 2$ [20,21], in total, the topological entanglement entropy gives $2 \times (-\log 2) = -2 \log 2$. On the contrary, in the case of n_x being odd, such decomposition (8) is not valid, thus the Hamiltonian describes one \mathbb{Z}_2 toric code. To see this more clearly, recalling the fact that each commuting term of the Hamiltonian involves next-nearest-neighboring Pauli operators in the *x* direction, we think of doubling the lattice, obtained by going around the torus in the *x* direction twice and rearranging the vertices and links.

To be more specific to what we have just mentioned, consider a geometry portrayed in Fig. 7(b). We double the lattice obtained by traveling around the torus in the *x* direction twice. In the first half of the doubled lattice, we omit the vertices with even number *x* coordinate and vertical links with odd number *x* coordinate whereas in the latter half, we discard the vertices with with odd number *x* coordinate. Recalling (7), one is convinced that the Hamiltonian describes one \mathbb{Z}_2 toric code on the torus. By this rearrangement, the entanglement entropy S_A amounts to be the one of two spatially separated disks in the \mathbb{Z}_2 toric code. The topological entanglement entropy is found to be $-2 \log 2$, the same value as the case of n_x even.

V. GENERIC LATTICES

As we have seen in the previous argument, our model contains the second-order spatial derivative, involving the next-nearest-neighboring sites. Such a property can be straightforwardly described by the Laplacian, the graph the-oretical analog of the second-order spatial derivatives with which one can evaluate the entanglement entropy. In this section, we further explore the entanglement entropy of the model on *generic lattices*, rather than standard square one to see the interplay between quantum entanglement and geometry of the system.

A. Laplacian and Hamiltonian

To this end, let us start with reviewing a few terminologies in the graph theory. A graph G = (V, E) which is a pair consisting of a set of vertices V and a set of edges E comprised of pairs of vertices $\{v_i, v_i\}$. In the rest of the work, we assume that the graph is *connected*, meaning there is a path from a vertex to any other vertex, and that the graph does not have an edge that emanates from and terminates at the same vertex. We also introduce two quantities, $deg(v_i)$ and l_{ii} . The former one, $deg(v_i)$, denotes the *degree* of the vertex v_i , i.e., the number of edges emanating from the vertex v_i and the latter one l_{ii} represents the number of edges between two vertices v_i and v_i . (We have $l_{ij} = 0$ when there is no edge between two vertices v_i and v_i .). Using these two quantities, the Laplacian *matrix* of the graph, which is the analog of the second-order derivative operator ∂_r^2 on a graph, is defined. For a given graph G = (V, E), the Laplacian matrix L (which we abbreviate as Laplacian in the rest of this work) is the matrix with rows and columns indexed by the elements of vertices $\{v_i\} \in V$, with

$$L_{ij} = \begin{cases} \deg(v_i) \ (i = j), \\ -l_{ij} \ (i \neq j). \end{cases}$$
(77)

The Laplacian is singular due to the connectivity of the graph. (Summing over all rows or columns gives zero.) As an example, the Laplacian of the cycle graph C_4 (i.e., a square) consisting of four vertices and four edges, where there is a single edge between a pair of vertices, is given by

$$L = \begin{pmatrix} 2 & -1 & & -1 \\ -1 & 2 & -1 & \\ & -1 & 2 & -1 \\ -1 & & -1 & 2 \end{pmatrix}.$$

It is known that by introducing invertible matrices over integer *P*, *Q* corresponding to linear operations on rows and columns of the Laplacian, respectively, the Laplacian of any connected graph can be transformed into a diagonal form $(Smith normal form)^1$

$$PLQ = diag(u_1, u_2, \dots, u_{n-1}, 0).$$
 (78)

The diagonal entries u_i are referred to as the *invariant* factors.

With these terminologies, we introduce generic 2D lattice on which the higher-rank topological phase is placed. We consider 2D lattices constructed by the product of the graph G(V, E) and 1D line. An example of such a lattice is shown in Fig. 8, left. The lattice is constructed in such a way that copies of the graphs are stacked along the vertical direction.²

We introduce two types of the generalized *N*-qubit states $(\mathbb{Z}_N \text{ clock states})$ on this 2D lattice. The first type of the clock states is located on the vertices of the graph in the horizontal direction, whereas the ones with the second type are on the vertical links. We denote the coordinate of the first clock states by (v_i, y) where v_i represents a vertex of the graph and y does the height taking integer values in the unit of lattice spacing. Analogously, the coordinates of the second clock states are denoted by $(v_i, y + \frac{1}{2})$, where the second element corresponds to the edge between vertices located at (v_i, y) and $(v_i, y + 1)$.

Having defined the 2D lattice, we introduce the Hamiltonian which is generalization of the one studied in the previous sections to an arbitrary graph. We represent the local Hilbert space of the two types of the clock states as $|a\rangle_{(v_i,y)}$ and $|b\rangle_{(v_i,y+1/2)}$ $(a, b \in \mathbb{Z}_N)$. Further, we define operators $X_{1,(v_i,y)}, Z_{1,(v_i,y)}$, and $X_{2,(v_i,y+1/2)}, Z_{2,(v_i,y+1/2)}$ that act on these states as

$$Z_{1,(v_{i},y)}|a\rangle_{(v_{i},y)} = \omega^{a}|a\rangle_{(v_{i},y)},$$

$$Z_{2,(v_{i},y+1/2)}|b\rangle_{(v_{i},y+1/2)} = \omega^{b}|b\rangle_{(v_{i},y+1/2)},$$

$$X_{1,(v_{i},y)}|a\rangle_{(v_{i},y)} = |a+1\rangle_{(v_{i},y)},$$

$$X_{2,(v_{i},y+1/2)}|b\rangle_{(v_{i},y+1/2)} = |b+1\rangle_{(v_{i},y+1/2)}$$
(79)

¹The last entry is zero, mirroring the fact that the graph is connected.

²Note that the lattice obtained this way is not embeddable in flat 2D space. Rather, it is defined on an abstract 2D cell complex. Indeed, the graph consists of vertices and edges, corresponding to 0- and 1-simplices. Due to this reason, we regard our lattice as 2D.



FIG. 8. Left: One example of the 2D lattice consisting of a connected graph G_x and the 1D line. Right: One of the terms $V_{(v_i,y)}$ defined in (80) is shown in the same example of the lattice.

with $\omega = e^{2\pi i/N}$. With these notations, we define the following two types of operators that we dub vertex and plaquette operators at each vertex and vertical link:

$$V_{(v_{i},y)} := X_{2,(v_{i},y+1/2)} X_{2,(v_{i},y-1/2)}^{\dagger} (X_{1,(v_{i},y)}^{\dagger})^{\deg(v_{i})} \\ \times \prod_{j} (X_{1,(v_{j},y)})^{l_{ij}}, \\ P_{(v_{i},y+1/2)} := Z_{1,(v_{i},y+1)}^{\dagger} Z_{1,(v_{i},y)} Z_{2,(v_{i},y+1/2)}^{\deg(v_{i})} \prod_{j} (Z_{2,(v_{j},y+1/2)}^{\dagger})^{l_{ij}}.$$

$$(80)$$

Here, $\deg(v_i)$ is the number of edges in the *x* direction which emanate from the vertex v_i whereas l_{ij} gives the number of edges in the *x* direction between two vertices with coordinates (v_i, y) and (v_j, y) , in accordance with the matrix element of the Laplacian (77). It is straightforward to check terms given in (80) commute with each other. We demonstrate one example of the term in Fig. 8, right. The Hamiltonian is defined by

$$H = -\sum_{(v_i, y)} V_{(v_i, y)} - \sum_{(v_i, y+1/2)} P_{(v_i, y+1/2)} + \text{H.c.}$$
(81)

When we impose the periodic boundary condition in the y direction, by the same reasoning presented in Sec. II D, one finds that the GSD is given by [19]

$$GSD = [N \times \gcd(N, u_1) \times \cdots \times \gcd(N, u_{n-1})]^2.$$
(82)

Here, positive integers u_i denote the invariant factors of the Laplacian (78).

B. Entanglement entropy

Now we are in a good place to study the entanglement entropy of our model on a graph. We set the subsystem Aas the "cylinder geometry" consisting of the subgraph of Gand the vertical line with length l_y , i.e., the cylinder encloses l_y copies of the subgraphs. The clock states belonging to the subsystem A are the ones located on the vertices of the subgraph within the cylinder and the ones on the vertical links inside the cylinder and vertical links that cross the top and bottom faces of the cylinder. We also denote the complement of the subsystem A as B. For simplicity, we set the subgraph to be connected. An example is shown in Fig. 9(a). We set the y coordinate of vertices inside the subsystem A to be $(y_0 \le y \le y_0 - 1 + l_y)$.



FIG. 9. (a) Example of the subsystem *A* in a shape of a cylinder geometry consisting of the subgraph and the 1D line. The clock states belonging to the subsystem *A* are marked by red colors. (b) Classification of the vertices of the graph at $(y_0 \le y \le y_0 - 1 + l_y)$. Inside the subsystem *A*, the vertices which are (are not) connected with the ones outside *A* are represented by $A_I(A_0)$, marked by pink (red) color. Also, outside *A*, the vertices which are (are not) connected with the ones in *A* are denoted by $B_I(B_0)$, marked by light blue (blue).

For later convenience, at each y, $(y_0 \le y \le y_0 - 1 + l_y)$, we label vertices of the graph classified by the following four groups:

 $A_0 = \{$ vertices inside A which are not connected with the ones outside $A\},$

- $A_I = \{$ vertices inside A which are connected with the ones outside A $\},$
- $B_I = \{$ vertices outside A which are connected with the ones inside $A\},\$
- $B_0 = \{ \text{vertices outside } A \text{ which are not connected with the ones inside } A \}.$ (83)

An example is demonstrated in Fig. 9(b). We denote the number of such vertices by n_{A_0} , n_{A_I} , n_{B_I} , and n_{B_0} . In accordance with the vertices of the four classes [Eq. (83)], we rearrange the order of the vertices so that vertices of A_0 come first, followed by the ones of A_I , B_I , and B_0 . In this order of the vertices, we write the Laplacian of the graph in the form

$$L = \begin{pmatrix} L_{A_0A_0} & L_{A_0A_I} & & \\ L_{A_IA_0} & L_{A_IA_I} & L_{A_IB_I} & \\ & & L_{B_IA_I} & L_{B_IB_I} & L_{B_IB_0} \\ & & & & L_{B_0B_I} & L_{B_0B_0} \end{pmatrix},$$
(84)

where L_{**} describes the submatrix of the Laplacian. For instance, the matrix $L_{A_0A_0}$ is the $n_{A_0} \times n_{A_0}$ matrix indexed by vertices which belong to A_0 . After preparing terminologies, we are going to show the following theorem:

Theorem 1. Entanglement entropy of a subgraph. Consider the stabilizer model on the 2D lattice (81) obtained by the product of the graph and one-dimensional (1D) line with periodic boundary condition in the vertical direction. The entanglement entropy of the subsystem A with cylinder geometry, constructed by a subgraph and the 1D line with length l_y with respect to the ground state $|\psi\rangle$, is given by

$$S_{A} = \left[(\operatorname{Area}) \log N - 2l_{y} \log \left(\prod_{i} \gcd(N, r_{i}) \right) - l_{y} |n_{A_{I}} - n_{B_{I}}| \log N \right] - \log \left(\frac{N^{\min(n_{A_{I}}, n_{B_{I}})}}{\prod_{i} \gcd(N, r_{i})} \right) - 2 \log \left(\prod_{j} \gcd(N, s_{j}) \right).$$
(85)

Here, (Area) represents the number of the vertex operators that have support on both of *A* and *B*, which is associated with the number of the clock states located around the border of *A* and *B*. Also, the numbers r_i and s_j represent invariant factors of the submatrix of the Laplacian via (SNF represents the Smith normal form)

$$L_{A_{I}B_{I}} \xrightarrow{\mathrm{SNF}} \begin{pmatrix} r_{1} & & \\ & \ddots & \\ & & r_{p} \\ & & \end{pmatrix}, \quad \begin{pmatrix} L_{A_{0}A_{0}} & L_{A_{0}A_{I}} \\ L_{A_{I}A_{0}} & L_{A_{I}A_{I}} & L_{A_{I}B_{I}} \end{pmatrix} \xrightarrow{\mathrm{SNF}} \begin{pmatrix} s_{1} & & \\ & \ddots & \\ & & s_{n_{A_{0}}+n_{A_{I}}} \end{pmatrix}.$$
(86)

Proof. In order to find the entanglement entropy, we resort to the formula (44). Generically, it is cumbersome to evaluate $|G_A|$ and $|G_B|$ due to the constraint on the stabilizers. As we have seen in Sec. III, the complication stems from the fact that there is nontrivial GSD when we impose the periodic boundary condition on the lattice, giving rise to constraints on the multiplication of the stabilizers on the entire lattice. While it is still possible to implement a similar approach as the one discussed in Sec. III to identify the entanglement entropy, here we employ an alternative simpler way to obtain the entanglement entropy. We accommodate an appropriate boundary condition on the system in the vertical direction, instead of the periodic boundary condition. Consider the 2D lattice that we introduced in (81) with finite length in the vertical direction so the bottom boundary terminates with a graph G and the top boundary ends with vertical links (see Fig. 10). With this boundary condition, one can check that the GSD is trivial and there is no constraint on the multiplication

of stabilizers on the entire lattice, simplifying the problem. The boundary that we consider here is intriguing in its own right. It is reminiscent of smooth and rough boundary of the toric code, where anyons with magnetic and electric charge are condensed [31].

After accommodating the boundary, now we turn to calculation of the entanglement entropy based on (44) assuming $n_{B_I} \leq n_{A_I}$. (The proof in the case of $n_{B_I} > n_{A_I}$ is similarly discussed.). From (44), one needs to evaluate the number of products of vertex operators that act exclusively on *A* or *B*. We denote the number of *individual* vertex terms that act within *A* and *B* as $|\tilde{G}_A|$ and $|\tilde{G}_B|$. To find $|G_A|$ and $|G_B|$, we also need to think of the multiplication of the vertex terms which act exclusively on *A* and *B*. Representing these numbers as Γ_A and Γ_B , $|G_A|$ and $|G_B|$ are written as

$$|G_A| = |\tilde{G}_A| \times \Gamma_A, \quad |G_B| = |\tilde{G}_B| \times \Gamma_B. \tag{87}$$

Below we identify Γ_A and Γ_B .



FIG. 10. We accommodate the boundary condition on the lattice so that in the bottom the lattice terminates the graph whereas on the top, the lattice terminates with the vertical links. One can verify that the there is no constraint on the stabilizer on the entire lattice, simplifying the problem.

At given y ($y_0 \le y \le y_0 - 1 + l_y$), we consider the following product of the vertex terms:

$$\prod_{v_i \in A_I} V_{(v_i, y)}^{\alpha_i} \ (\alpha_i \in \mathbb{Z}_N), \tag{88}$$

which can be rewritten as

$$\prod_{v_i \in A_I} V_{(v_i, y)}^{\alpha_i} = \left[\prod_{i \in A_I} X_{2, (v_i, y+1/2)}^{-\alpha_i} X_{2, (v_i, y-1/2)}^{\alpha_i} \right] \\ \times \prod_{j \in A_0 + A_I + B_I} X_{1, (v_j, y)}^{\beta_j} \ (\beta_j \in \mathbb{Z}_N).$$
(89)

Defining a n_{A_I} - and $n_{A_0+A_I+B_I}$ -dimensional vectors $\boldsymbol{\alpha}$ and $\boldsymbol{\beta}$ whose element is given by α_i and β_j , respectively, and referring to (84), it follows that the vectors $\boldsymbol{\alpha}$ and $\boldsymbol{\beta}$ are related via submatrix of the Laplacian as

$$\boldsymbol{\beta} = \begin{pmatrix} L_{A_0 A_I} \\ L_{A_I A_I} \\ L_{B_I A_I} \end{pmatrix} \boldsymbol{\alpha}.$$
 (90)

Suppose the last n_{B_I} entries of β are zero. Then the product (89) is what we want: multiplication of the vertex operators that act exclusively on A_I , contributing to Γ_A . Thus, the number of such product amounts to be the number of solutions of

$$L_{B_i A_i} \boldsymbol{\alpha} = \boldsymbol{0} \mod N. \tag{91}$$

Introducing $n_{B_I} \times n_{B_I}$ and $n_{A_I} \times n_{A_I}$ invertible integer matrices U and V, one can transform the $n_{B_I} \times n_{A_I}$ matrix $L_{B_IA_I}$ into the Smith normal form as

$$UL_{B_{I}A_{I}}V = \begin{pmatrix} r_{1} & & \\ & \ddots & \\ & & r_{n_{A_{I}}} \\ & & & \end{pmatrix}.$$
 (92)

From this form (92), we have

$$L_{B_{I}A_{I}}\boldsymbol{\alpha} = \boldsymbol{0} \Leftrightarrow \begin{pmatrix} r_{1} & & \\ & \ddots & \\ & & r_{n_{A_{I}}} \end{pmatrix} \tilde{\boldsymbol{\alpha}} = \boldsymbol{0} \mod N, \quad (93)$$

where $\tilde{\boldsymbol{\alpha}} := V^{-1} \boldsymbol{\alpha}$. The *i*th element of $\tilde{\boldsymbol{\alpha}}$ is subject to

$$r_i \tilde{\alpha}_i = 0 \mod N, \tag{94}$$

from which it follows that r_i takes $gcd(N, r_i)$ distinct values, where gcd stands for the greatest common divisor. Therefore, the number of the solution of (91) is given by

$$\prod_{i=1}^{n_{A_I}} \gcd(N, r_i).$$

Since we have considered the product of the vertex operators defined on A_I that act trivially on B_I at given y ($y_0 \le y \le y_0 - 1 + l_y$), there are in total

$$\left[\prod_{i=1}^{n_{A_l}} \gcd(N, r_i)\right]^{l_y} (:= \Gamma_1)$$
(95)

of the product of the vertex operators defined on A_I which acts exclusively on A_I , contributing to Γ_A .

We can analogously discuss the number of product of the vertex operators located on B_I which act trivially on A_I , contributing to Γ_B . Suppose the following product at given $y (y_0 \le y \le y_0 - 1 + l_y)$

$$\prod_{j \in B_I} V_{(v_j, y)}^{\eta_j} (\eta_j \in \mathbb{Z}_N)$$
(96)

acts trivially on A_I . By the similar augment presented around (91), the number of such product amounts to be the number of solution of

$$L_{A_I B_I} \boldsymbol{\eta} = \boldsymbol{0} \mod N, \tag{97}$$

where η denotes n_{B_I} -dimensional vector whose entry is given by η_j . Similarly to (92), we transform the matrix $L_{A_IB_I}$ into the Smith normal form. Introducing $n_{A_I} \times n_{A_I}$ and $n_{B_I} \times n_{B_I}$ invertible integer matrices U' and V', the Smith normal form reads as

$$U'L_{A_{I}B_{I}}V' = \begin{pmatrix} r_{1} & & & & 0 & \\ & \ddots & & & \\ & & r_{n_{A_{I}}} & & & 0 \end{pmatrix}.$$
 (98)

Since $L_{B_IA_I}^T = L_{A_IB_I}$, the smith normal form of the $n_{A_I} \times n_{B_I}$ matrix $L_{A_IB_I}$ contains the same diagonal entries as (92). Equation (97) is rewritten as

$$\begin{pmatrix} r_1 & & & \\ & \ddots & & \\ & & r_{n_{A_I}} & & \ddots & \\ & & & r_{0} \end{pmatrix} \tilde{\boldsymbol{\eta}} = \boldsymbol{0} \mod N \quad (99)$$

with $\tilde{\eta} := V'^{-1}\eta$. The first n_{A_I} entries of $\tilde{\eta}$ are subject to $\tilde{\eta}_i = r_i \ (1 \le i \le n_{A_I})$ whereas there is no constraint on the last $n_{B_I} - n_{A_I}$ entries of $\tilde{\eta}$. Hence, the number of solution satisfying (97) is given by

$$\prod_{i=1}^{n_{A_I}} \gcd(N, r_i) \times N^{(n_{B_I} - n_{A_I})}.$$
 (100)

Taking into account the y direction, we have

$$\prod_{i=1}^{n_{A_{I}}} \gcd(N, r_{i}) \times N^{(N_{B_{I}} - n_{A_{I}})} \bigg]^{l_{y}} (:= \Gamma_{2})$$
(101)

product of the vertex operators defined on B_I which act trivially on A_I , contributing to Γ_B .

In addition to (95) and (101), there are products of the vertex operators along both horizontal and vertical directions, which contributes to Γ_A and Γ_B . At given *y*, let us consider the following product:

$$\prod_{i \in A_0 + A_I} V_{(v_i, y)}^{\delta_i} \ (\delta_i \in \mathbb{Z}_N).$$
(102)

Suppose (102) does not have X_1 operators in the horizontal direction. Recalling the previous argument presented around (91) and (97), such a condition is described by

$$\begin{pmatrix} L_{A_0A_0} & L_{A_0A_I} \\ L_{A_IA_0} & L_{A_IA_I} \\ & L_{B_0A_I} \end{pmatrix} \boldsymbol{\delta} = \boldsymbol{0} \mod N.$$
(103)

By transforming the matrix on the left-hand side of (103) into the Smith normal form as

$$\begin{pmatrix} L_{A_0A_0} & L_{A_0A_I} \\ L_{A_IA_0} & L_{A_IA_I} \\ & L_{B_0A_I} \end{pmatrix} \rightarrow \begin{pmatrix} s_1 & & \\ & \ddots & \\ & & s_{n_{A_I}+A_0} \\ & & & \end{pmatrix}, \quad (104)$$

there are

$$\Gamma_3 := \prod_{i=1}^{n_{A_I+A_I}} \gcd(N, s_i) \tag{105}$$

solutions. Multiplying (102) along the vertical direction gives rise to stabilizers, which contribute to Γ_A and Γ_B . Indeed,

$$\prod_{y=y_0-1}^{y_0+l_y} \left[\prod_{i \in A_0+A_I} V_{(v_i,y)}^{\delta_i} \right]$$
(106)

has support only on *B*, which contributes to Γ_B (Fig. 11, left). Likewise, the product

$$\prod_{y=1}^{y_0} \left[\prod_{i \in A_0 + A_I} V_{(v_i, y)}^{\delta_i} \right]$$
(107)

acts exclusively on A, contributing to Γ_A (Fig. 11, right).

We also consider the product of the vertex operators defined on $A_0 + A_I + B_I$:

$$\prod_{i\in A_0+A_I+B_I} V_{(v_i,y)}^{\sigma_i},\tag{108}$$

assuming it does not have X_1 operators on $A_0 + A_I$. Such a condition can be described by

$$\begin{pmatrix} L_{A_0A_0} & L_{A_0A_I} \\ L_{A_IA_0} & L_{A_IA_I} & L_{A_IB_I} \end{pmatrix} \boldsymbol{\sigma} = \boldsymbol{0} \mod N.$$
(109)

Multiplying (108) along *y*,

$$\prod_{y=y_0}^{y_0-1+l_y} \prod_{i \in A_0+A_I+B_I} V_{(v_i,y)}^{\sigma_i},$$
(110)

yields the stabilizer that acts within *B*, contributing to Γ_B . Thus, we need to evaluate the number of solutions satisfying (109). However, some of solutions are redundant to what we have already identified previously [Eqs. (100) and (105)]. Therefore, the number of products in question which contribute to Γ_B [Eq. (110)] is equivalent to the number of solutions of (109) divided by the numbers given in (100) and (105):

$$\frac{\text{The number of solutions of (109)}}{\prod_{i=1}^{n_{A_I}} \gcd(N, r_i) \times N^{(N_{B_I} - N_{A_I})} \times \prod_{j=1}^{n_{A_I + n_{A_0}}} \gcd(N, s_j)}.$$
 (111)



FIG. 11. Left: Multiplication of the stabilizer $V_{v_i,y}$ with blue stars gives rise to the operators X_1 and X_2 that act exclusively on *B*. Right: Multiplication of the stabilizer $V_{v_i,y}$ with blue stars gives rise to the operators X_1 and X_2 that act exclusively on *A*. The clock states with red color belong to the subsystem A.

We transform the matrix in (105) into the Smith normal form as

$$\begin{pmatrix} L_{A_0A_0} & L_{A_0A_I} \\ L_{A_IA_0} & L_{A_IA_I} & L_{A_IB_I} \end{pmatrix}$$

$$\xrightarrow{\text{SNF}} \begin{pmatrix} s_1 & & & & \\ & \ddots & & & \\ & & s_{n_{A_0}+n_{A_I}} & & & 0 \end{pmatrix}, \quad (112)$$

from which it follows that the denominator of (111) is given by $\prod_{j=1}^{n_{A_j+A_l}} \operatorname{gcd}(N, s_j) \times N^{n_{B_l}}$. Substituting it into (111) yields

$$(111) = \frac{N^{n_{A_I}}}{\prod_{i=1}^{n_{A_I}} \gcd(N, r_i)} (:= \Gamma_4).$$
(113)

Overall,

$$\Gamma_A = \Gamma_1 \times \Gamma_3, \ \Gamma_B = \Gamma_2 \times \Gamma_3 \times \Gamma_4.$$
 (114)

Referring to (44) and (87), we finally arrive at

$$S_{A} = \left[\log \left(\frac{|G|}{|\tilde{G}_{A}||\tilde{G}_{B}|} \right) - 2l_{y} \log \left(\prod_{i=1}^{n_{A_{I}}} \gcd(N, r_{i}) \right) - l_{y}(n_{B_{I}} - n_{A_{I}}) \log N \right] - \log \left(\frac{N^{n_{A_{I}}}}{\prod_{i=1}^{n_{A_{I}}} \gcd(N, r_{i})} \right) - 2\log \left(\frac{n_{A_{I}} + n_{A_{0}}}{\prod_{j=1}^{n_{A_{I}}} \gcd(N, s_{j})} \right) (n_{A_{I}} \leq n_{B_{I}}). \quad (115)$$

Proof of the theorem (85) is completed by rewriting $\frac{|G|}{|\tilde{G}_A||\tilde{G}_B|} = N^{(\text{Area})}$, where (Area) denotes the number of vertex operators that have support both on *A* and *B*, which is interpreted as the number of vertices surrounding the subsystem *A*. Analogous argument leads to (85) in the case of $n_{B_I} < n_{A_I}$.

The first three terms in (85) depend on the system size, whereas the subleading-order terms corresponding to the last two terms depend on the number of clock states inside A which are connected with B (i.e., n_{A_I}) and invariant factors of the Laplacian of the subgraph.

While we corroborate that the subleading-order term of the entanglement entropy (73) in the case of the square lattice is topological by an intuitive argument given in Sec. IV, one can understand the topological origin of the subleading term in the case of the square lattice from a different perspective based on the result (85). In the case of the square lattice, for a disk geometry, we have $n_{A_I} = n_{B_I} = 2$ and invariant factors of the subgraph (i.e., disk geometry) become trivial (namely, $r_i = s_j = 1$). Thus, (85) is identical to (73). The subleading term $-n_{A_I} \log N = -2 \log N$ retains the same value, regardless of the width of the disk as the number n_{A_I} is always constant 2.

VI. BRIEF COMMENTS ON OTHER CASES OF THE HIGHER-RANK TOPOLOGICAL PHASES

We have studied entanglement entropy of the higher-rank topological phases defined in (5) by making use of formalism of the stabilizers [22] jointly with the one of the Laplacian. One could study entanglement entropy of other higher-rank topological phases by resorting to a similar approach.

One example is the model studied recently in [32], which is introduced as follows. At each vertex of the square lattice, we introduce two clock states $|a\rangle_{(x,y)}|b\rangle_{(x,y)}$, $a, b \in \mathbb{Z}_N$. Also, we define operators acting on these states as (recall that $\omega = e^{2\pi i/N}$)

$$Z_{1,(x,y)}|a\rangle_{(x,y)}|b\rangle_{(x,y)} = \omega^{a}|a\rangle_{(x,y)}|b\rangle_{(x,y)},$$

$$Z_{2,(x,y)}|a\rangle_{(x,y)}|b\rangle_{(x,y)} = \omega^{b}|a\rangle_{(x,y)}|b\rangle_{(x,y)},$$

$$X_{1,(x,y)}|a\rangle_{(x,y)}|b\rangle_{(x,y)} = |a+1\rangle_{(x,y)}|b\rangle_{(x,y)},$$

$$X_{2,(x,y)}|a\rangle_{(x,y)}|b\rangle_{(x,y)} = |a\rangle_{(x,y)}|b+1\rangle_{(x,y)}.$$
(116)



FIG. 12. Two types of terms defined in (117).

Introducing the following terms by (Fig. 12)

$$V_{(x,y)} := X_{1,(x+1,y)} X_{1,(x-1,y)} (X_{1,(x,y)}^{\dagger})^2 X_{2,(x,y+1)} X_{2,(x,y-1)} \times (X_{2,(x,y)}^{\dagger})^2,$$

$$P_{(x,y)} := Z_{1,(x,y+1)}^{\dagger} Z_{1,(x,y-1)}^{\dagger} Z_{2,(x+1,y)}^2 Z_{2,(x-1,y)} \times (Z_{2,(x,y)}^{\dagger})^2,$$
(117)

the Hamiltonian reads as

$$H_{\mathbb{Z}_N} = -\sum_{x,y} (V_{(x,y)} + P_{(x,y)}) + \text{H.c.}$$
(118)

The GSD of the phase on the torus geometry with length of the lattice in the *x* and *y* directions being n_x and n_y , is found to be [32]

$$GSD = [N \times \gcd(N, n_x) \times \gcd(N, n_y) \times \gcd(N, n_x, n_y)]^2.$$
(119)

By making use of the similar line of thoughts outlined in Sec. III, one obtains the entanglement entropy of the disk geometry as

$$S_A = (\text{Area})\log N - 4\log N. \tag{120}$$

Here, Area is the number of vertices surrounding the disk. Compared with the result (73), the absolute value of the topological entanglement entropy is increased. Also, the total quantum dimension of the model is given by $\sqrt{\sum_a d_a^2} = [N \times \text{gcd}(N, n_x) \times \text{gcd}(N, n_y) \times \text{gcd}(N, n_x, n_y)]$, thus, analogously to (73), the relation between total quantum dimension and the topological entanglement entropy does not hold.

It would be interesting to study other higher-rank topological phases, such as the ones studied in [26,33–36]. Generically, we make a conjecture that one cannot associate the topological entanglement entropy in the higher-rank topological phases with the total quantum dimension of the fractional excitations and also that the absolute value of the topological entanglement entropy becomes larger when the model contains the terms which involve clock states in the longer range. We leave confirmation of this speculation for future studies.

VII. CONCLUSION

In this work, we study entanglement entropy of unusual topological stabilizer models with dipole symmetry, which is one of the multipole symmetries, admitting dipole of the fractional charges. As opposed to the entanglement entropy of the fracton topological phases, where it is not immediate to identify the entanglement entropy due to the subsystem symmetries [37–39], entanglement entropy of our model is easily obtained based on the formulation of the combinatorics. While the GSD and the total quantum dimension of the fractional excitations drastically changes depending on the system size, topological entanglement entropy γ takes a constant, given by $\gamma = -2 \log N$. Due to this result, the wellknown relation between the topological entanglement entropy and the total quantum dimension is not valid in the case of the higher-rank topological phases. Such a result can be understood by decomposition of the model into two or rearrangement of the lattice in the simplest case N = 2. We further study the entanglement entropy of the model on generic lattices and have found that the entanglement entropy depends on the number of clock states surrounding the system and invariant factors of the Laplacian of the subgraph. The result presented in this work complies with the growing interests in topological phases with multipole symmetries in view of quantum entanglement and combinatorics.

It would be interesting to see if our model can be decomposed into layers of the toric codes since such a way of thinking is crucial in understanding physics of the fracton topological phases. Indeed, it is known that the X-cube model, which is the prototype of the fracton topological phase realized on a cubic lattice is decomposed into layers of the toric codes via unitary operations [40,41]. Furthermore, based on this fact, a new type of topological field theories (foliated topological field theories) have been recently introduced, where layers of the (2 + 1)D BF theories are stacked, giving rise to quasiparticle excitations with mobility restriction [42]. In our model, one can see that a single fractional charge can hop in the x direction in the unit of N lattice spacing rather than a single site [see the discussion around (13) and [18,26,27]], tantalizing to decompose the model into N layers of the toric codes. Also, it is proved that for prime p, all translation-invariant \mathbb{Z}_p stabilizer codes in 2D are equivalent to copies of toric codes by a local Clifford circuit of constant depth [43]. On the other hand, based on the recent study of multipole symmetries [44], one can show that the effective field theory description of the model (5) turns out to be a foliated topological field theory constructed by two layers of the BF theories. It would be intriguing and important to see whether these facts reconcile with each other and, in particular, to identify a proper unitary quantum circuit to decompose our model into layers of the toric codes in the case of N > 2.

Regarding the issue in the previous paragraph, it was recently discussed that a model which exhibits the same GSD of the the model (13) is introduced by preparing two copies of the toric codes with arrays of Dijkgraaf-Witten (DW) twist terms (gauged symmetry-protected topological phases) [45,46]. This consideration is consistent with our result (73). Generally, the DW twist terms change the fractional statistics of excitations, yet do not alter the total quantum dimension. (Indeed, it was investigated that topological entanglement entropy of a contractible disk in quantum double with DW twist is the same as the one without DW twist [47].) Hence, regardless of the presence of the DW twist, the topological entanglement entropy amounts to the one of layers of toric codes $-2 \log N$, which is consistent with (73). More detailed discussions can be found in [46].

There are several future research directions regarding this study. It is known that capability of the error correction of the toric code is characterized by the classical Ising universality class [6]. It would be interesting to study how our model (5) can be utilized for quantum error corrections, in particular, to identify what kind of universality class characterizes the capability of the error correction in our model. One naively expects that due to the fact that the model has the multipole symmetries, such a universal class is qualitatively different from the one in the conventional toric code. While we focus on the entanglement entropy in the higher-rank topological phases, one would be curious to elucidate behavior of other observable to quantify the quantum entanglement. One candidate would be entanglement negativity [48,49], an important observable to extract quantum correlation rather than a classical one. It would be intriguing to investigate the entanglement negativity of our model to see how different it is compared with the case of the conventional topologically ordered phases [50].

ACKNOWLEDGMENTS

The author thanks B. Han, K. Shiozaki, T. Nakanishi, M. Honda, and Y. Oreg for helpful discussions. The author also thanks G. Delfino and J. Hoon Han for notification of their papers [34,36]. This work is in part supported by KAKENHI-Project No. 23H01097.

APPENDIX A: DERIVATION OF THE LOGICAL OPERATORS (WILSON LOOPS)

In this Appendix, we present an intuitive understanding of the logical operators of the model (29) and (31) by investigating the behavior of the fractional excitations (i.e., fusion rules) following [19]. Based on the fusion rules, one can study the logical operators which are the noncontractible Wilson loops when the model is placed on a torus geometry. To this end, we consider the model (5) on a 2D lattice with periodic boundary condition and the system size being $n_x \times n_y$. In this lattice, we first focus on the logical operators consisting of a string of Z_1 operators in the *x* direction. We think of a closed loop of the electric charge in the horizontal direction at *y*, described by $\prod_{x=1}^{n_x} Z_{1,(x,y)}^{a_x}$ with $a_x \in \mathbb{Z}_N$. From (12), the electric charges induced by acting this operator on the ground state are described by the fusion rule

$$I \to \prod_{x=1}^{n_x} \otimes e_{(x,y)}^{r_x} (r_x \in \mathbb{Z}_N)$$
(A1)

with

$$\mathbf{r} = -L\mathbf{a}.\tag{A2}$$

Here, $\mathbf{a} := (a_1, \ldots, a_{n_x})^T$, $\mathbf{r} := (r_1, \ldots, r_{n_x})^T$, and $n_x \times n_x$ matrix *L* is the Laplacian (15). Since the closed noncontractible loop $\prod_{x=1}^{n_x} Z_{1,(x,y)}^{a_x}$ has to commute with the Hamiltonian, the fusion rules (A1) and (A2) have to be trivial, i.e., $\mathbf{r} = \mathbf{0} \mod N$. Thus, the closed loops of the electric charges are characterized by the *kernel* of the Laplacian, which is derived in (27):

$$\mathbf{a} = \frac{N}{\gcd(n_x, N)} \alpha_1 \begin{pmatrix} n_x - 1\\ n_x - 2\\ \vdots\\ 1\\ 0 \end{pmatrix} + \alpha_2 \begin{pmatrix} 1\\ 1\\ \vdots\\ 1\\ 1 \end{pmatrix} \mod N \ (\alpha_1 \in \mathbb{Z}_{\gcd(N, n_x)}, \ \alpha_2 \in \mathbb{Z}_N).$$

Based on this result, the logical operators consisting of the closed string of the Z_1 operators are described by

$$\eta_1^z = \prod_{x=1}^{n_x} Z_{1,(x,y)}^{p_x}, \ \gamma_1^z = \prod_{x=1}^{n_x} Z_{1,(x,y)}^{q_x}$$

with

$$\mathbf{p} = (p_1, p_2, \dots, p_{n_x})^T = \frac{N}{\gcd(n_x, N)} \alpha_1 \begin{pmatrix} n_x - 1 \\ n_x - 2 \\ \vdots \\ 1 \\ 0 \end{pmatrix},$$
$$\mathbf{q} = (q_1, q_2, \dots, q_{n_x})^T = \alpha_2 \begin{pmatrix} 1 \\ 1 \\ \vdots \\ 1 \\ 1 \end{pmatrix} \mod N$$

which correspond to the first two terms in (31). One can check that the loops are deformable in the *y* direction, and that any form of the logical operators comprised of the product of the Z_1 operators in the *x* direction is generated by the combination of the two η_1^z , γ_1^z .

As for the logical operators comprised of product of the Z_2 operators in the y direction, from the fusion rule (59), which is identical to the one in the toric code, where a pair of excitations are created. From this feature, one can construct a noncontractible closed loop of magnetic charge formed by string of Z_2 operators defined by

$$W_{ey}(x) := \prod_{y=1}^{n_y} Z_{2,(x,y+1/2)}.$$
 (A3)

Although the loop (A3) resembles the one in the toric code, the way it is deformed is crucially different. To see this, we think of the product of the operators $P_{(x,y+1/2)}$ by

 $\prod_{y=1}^{n_y} P_{(x,y+1/2)}$, and acting it on the loop as

$$\prod_{y=1}^{n_y} P_{(x,y+1/2)} W_{ey}(x) = W_{ey}(x-1) [W_{ey}^{\dagger}(x)]^2 W_{ey}(x+1), \quad (A4)$$

which can be understood that the loop $W_{ey}(x)$ is transferred to the adjacent *x* coordinate (i.e., ± 1), compensated by loop with opposite charge at *x*. In order to find the number of distinct logical operators consisting of Z_2 's, we need to evaluate the distinct configuration of the loop $W_{ey}(x)$ up to the deformation (A4). To do this, let us consider composite of the loops in the *y* direction described by

$$\prod_{x=1}^{n_x} W_{ey}^{b_x}(x) \quad (b_x \in \mathbb{Z}_N).$$
(A5)

Also, we introduce product of the operators $P_{(x,y+1/2)}$ by

$$\prod_{x=1}^{n_x} \left(\prod_{y=1}^{n_y} P_{(x,y+1/2)} \right)^{c_x} \quad (c_x \in \mathbb{Z}_{\mathbb{N}}).$$
 (A6)

Acting [Eq. (A6)] on the loops $\prod_{x=1}^{n_x} W_{ey}^{b_x}(x)$ ($b_x \in \mathbb{Z}_N$), jointly with (A4) gives rise to the loops in the form of

$$\prod_{x=1}^{n_x} W_{ey}^{b'_x}(x) \ (b'_x \in \mathbb{Z}_N)$$
(A7)

with

$$\mathbf{b}' = \mathbf{b} - L\mathbf{c}.\tag{A8}$$

Here, the vector $\mathbf{b} := (b_1, \dots, b_{n_x})^T$ with other vectors being similarly defined, and *L* represents the Laplacian (15). The number of the distinct configuration of the loops is equivalent to the configuration of \mathbf{b} under the identification $\mathbf{b} \sim \mathbf{b}'$. Hence, we need to evaluate $\mathbf{f} \in \mathbb{Z}_N/\text{Im}(L)$, which is *cokernel* of the Laplacian. It can be shown that the cokernel is labeled by [19]

$$\mathbf{f} = \beta_1 \begin{pmatrix} 0\\0\\\vdots\\1\\-1 \end{pmatrix} + \beta_2 \begin{pmatrix} 0\\0\\\vdots\\0\\1 \end{pmatrix}$$

$$\times \mod N \ (\beta_1 \in \mathbb{Z}_{\gcd(N,n_r)}, \ \beta_2 \in \mathbb{Z}_N).$$
(A9)

It is interesting to note that depending on N and the system size, the model admits the dipole of the loops (corresponding to the first term), i.e, a pair of the loops with opposite charges located at adjacent sites. Based on this result one finds that any logical operator consisting of the Z_2 's in the y direction is generated by

$$\eta_2^z = \left(\prod_{y=1}^{n_y} Z_{2,(n_x-1,y)}\right) \times \left(\prod_{y=1}^{n_y} Z_{2,(n_x,y)}^{\dagger}\right), \ \gamma_2^z = \prod_{y=1}^{n_y} Z_{2,(n_x,y)}$$

with condition $(\eta_2^z)^{\text{gcd}(n_x,N)} = (\gamma_2^z)^N = 1$. These are nothing but the last two terms in (31). By a similar line of thought, one can derive the logical operators comprised of X_1 's or X_2 [Eq. (29)].

APPENDIX B: PROOF OF EQ. (43)

Given $g, g' \in G$, we have $g'_A = g_A \Leftrightarrow g' = hg, h \in G_B$. Then (42) is reduced to

$$\rho_A = \frac{|G_B|}{|G|} \sum_{g \in G/G_B, g' \in G_A} g_A |0\rangle_{AA} \langle 0| (g_A g'_A)^{\dagger}.$$
(B1)

We have

$$\begin{split} \rho_A^2 &= \left(\frac{|G_B|}{|G|}\right)^2 \sum_{g,\tilde{g}\in G/G_B,\tilde{g}',\tilde{g}'\in G_A} g_A|0\rangle_{AA} \\ &\times \langle 0|(g_A\tilde{g}_A)^{\dagger}g'_A|0\rangle_{AA} \langle 0|(g'_A\tilde{g}'_A)^{\dagger} \\ &= \left(\frac{|G_B|}{|G|}\right)^2 \sum_{g\in G/G_B,\tilde{g},\tilde{g}'\in G_A} g_A|0\rangle_{AA} \langle 0|(g_A\tilde{g}_A\tilde{g}'_A)^{\dagger} \\ &= \left(\frac{|G_B|}{|G|}\right)^2 |G_A| \sum_{g\in G/G_B,\tilde{g}\in G_A} g_A|0\rangle_{AA} \langle 0|(g_A\tilde{g}_A)^{\dagger} \\ &= \frac{|G_B|}{|G|} |G_A|\rho_A, \end{split}$$
(B2)

from which one finds

$$\rho_A^n = \left(\frac{|G_A||G_B|}{|G|}\right)^{(n-1)} \rho_A.$$
 (B3)

The entanglement entropy is obtained as

$$S_A = \lim_{n \to 1} \frac{1}{1 - n} \log \operatorname{tr} \rho_A^n = \log \frac{|G|}{|G_A||G_B|}.$$
 (B4)

APPENDIX C: ENTANGLEMENT ENTROPY WITH RESPECT TO THE GENERIC GROUND STATE

In this Appendix, we give a derivation of the entanglement entropy of three subsystems discussed in Sec. III B in the case of the generic ground state (45). The density matrix ρ_A^{ζ} of a subsystem A in our model with respect to the generic ground state (45) reads as

$$\rho_A^{\zeta} = \mathrm{Tr}_B |\zeta\rangle \langle \zeta|, \tag{C1}$$

where $|\zeta\rangle$ is given by (45). From (34), it is explicitly written as

$$\rho_{A}^{\zeta} = \sum_{\substack{a,b,c,d \\ a',b',c',d'}} \alpha_{ab,cd} \bar{\alpha}_{a'b',c'd'} \operatorname{Tr}_{B} \\
\times \left[\left(\eta_{1}^{x} \right)^{a} \left(\gamma_{1}^{x} \right)^{b} \left(\eta_{2}^{x} \right)^{c} \left(\gamma_{2}^{x} \right)^{d} \rho_{0} \left(\eta_{1}^{x} \right)^{a'} \left(\gamma_{1}^{x} \right)^{b'} \left(\eta_{2}^{x} \right)^{c'} \left(\gamma_{2}^{x} \right)^{d'} \right] \\$$
(C2)

with $\rho_0 = |\zeta_{00}\rangle\langle\zeta_{00}|$. Below we look at the three cases corresponding to Figs. 4(a)-4(c).

1. Single row I

In Fig. 4(a), recalling the form of the logical operators (29) [see also Fig. 3(b)], the logical operators η_1^x and η_1^x act within subsystem *B*, hence the term Tr_{*B*}[...] in (C2) is transformed

as

$$\frac{1}{|G|} \sum_{g,g' \in G} (\eta_{2A}^{x})^{c} (\gamma_{2A}^{x})^{d} g_{A} |0\rangle_{AA} \langle 0| (g_{A}g'_{A})^{\dagger} (\eta_{2A}^{x\dagger})^{c'} (\gamma_{2A}^{x\dagger})^{d'}_{B} \\ \times \langle 0| g'^{\dagger}_{B} (\eta_{1}^{x})^{a-a'} (\gamma_{1}^{x})^{b-b'} (\eta_{2B}^{x})^{c-c'} (\gamma_{2B}^{x})^{d-d'} |0\rangle_{B}, \quad (C3)$$

where we rewrite the logical operator by the product form $\eta_2^x = \eta_{2A}^x \otimes \eta_{2B}^x$, and similarly for γ_2^x . From the inner product of the state $|0\rangle_B$ in (C3), one finds that

$$g'_B = I_B, \ a = a', b = b', c = c', d = d'.$$
 (C4)

Hence, the density matrix ρ_A^{ζ} becomes

$$\rho_{A}^{\zeta} = \sum_{c,d} \sum_{g \in G, g' \in G_{A}} \frac{1}{|G|} \sum_{a,b} |\alpha_{ab,cd}|^{2} (\eta_{2}^{x})^{c} (\gamma_{2}^{x})^{d} g_{A} |0\rangle_{AA}$$
$$\times \langle 0| (\eta_{2}^{x^{\dagger}})^{c} (\gamma_{2}^{x^{\dagger}})^{d} (g_{A}g'_{A})^{\dagger}.$$
(C5)

From this form we obtain

$$\operatorname{Tr}_{A}\left[\left(\rho_{A}^{\zeta}\right)^{n}\right] = \sum_{c,d} \left[\sum_{a,b} |\alpha_{ab,cd}|^{2}\right]^{n} \times \left(\frac{|G_{A}||G_{B}|}{|G|}\right)^{n-1}.$$
(C6)

_ ..

The entanglement entropy is given by

$$S_{A} = \lim_{n \to 1} \frac{1}{1 - n} \log \operatorname{tr}_{A} (\rho_{A}^{\zeta})^{n}$$

= $\log \frac{|G|}{|G_{A}||G_{B}|} - \sum_{c,d} \left[\left(\sum_{a,b} |\alpha_{ab,cd}|^{2} \right) \times \log \left(\sum_{a,b} |\alpha_{ab,cd}|^{2} \right) \right].$ (C7)

Therefore, (47) follows.

2. Single row II

In the case of Fig. 4(b), from (29) [see also Fig. 3(b)], it follows that two logical operators running in the *y* direction, η_2^x and γ_2^x exclusively act on *B*. Regarding the other two, η_1^x and γ_1^x which go around in the *x* direction, one can deform these operators so that they act within *A*. Therefore, the term Tr_{*B*}[...] in (C2) becomes

$$\frac{1}{|G|} \sum_{g,g' \in G} (\eta_1^x)^a (\gamma_1^x)^b g_A |0\rangle_{AA} \langle 0| (g_A g'_A)^{\dagger} (\eta_1^{x^{\dagger}})^{a'} (\gamma_1^{x^{\dagger}})^{b'}{}_B \\
\times \langle 0| g'_B (\eta_2^x)^{c-c'} (\gamma_2^x)^{d-d'} |0\rangle_B,$$
(C8)

from which one finds

$$g'_B = I_B, \ c = c', \ d = d'.$$
 (C9)

The density matrix is then rewritten as

$$\rho_A^{\zeta} = \sum_{a,b,a',b'} \sum_{c,d} \sum_{g \in G, g' \in G_A} \frac{\alpha_{ab,cd} \bar{\alpha}_{a'b',cd}}{|G|} \times (\eta_1^x)^a (\gamma_1^x)^b g_A |0\rangle_{AA} \langle 0| (\eta_1^{x\dagger})^{a'} (\gamma_1^{x\dagger})^{b'} (g_A g'_A)^{\dagger}.$$
(C10)

Introducing orthogonal states by

$$|p,q\rangle_{A} := \frac{1}{\sqrt{N \times \gcd(N,n_{x})}} \sum_{a=0}^{\gcd(n,n_{x})-1} \sum_{b=0}^{N-1} v^{pa} \omega^{qb}$$
$$\times (\eta_{1}^{x})^{a} (\gamma_{1}^{x})^{b} |0\rangle_{A} \ [0 \leq p \leq \gcd(N,n_{x})$$
$$-1, 0 \leq q \leq N-1]$$
(C11)

with $v = e^{2\pi i/\gcd(N,n_x)}$, $\omega = e^{2\pi i/N}$, Eq. (C10) is transformed as

$$\rho_A^{\zeta} = \frac{1}{|G|} \sum_{g \in G, g' \in G_A} \sum_{p=0}^{\gcd(N, n_x) - 1} \sum_{q=0}^{N-1} \left(\frac{1}{\Gamma} \sigma_{kl} \nu^{kp} \omega^{lq} \right) \\ \times g_A |p, q\rangle_{AA} \langle p, q | (g_A g'_A)^{\dagger}, \qquad (C12)$$

where

$$\sigma_{kl} := \sum_{\substack{a,a',b,b',c,d\\a-a'=k \mod \gcd(N,n_x)\\b-b'=l \mod N}} \alpha_{ab,cd} \bar{\alpha}_{a'b',cd}.$$
(C13)

Recall that we have defined $\Gamma = N \times \text{gcd}(N, n_x)$. Defining $\lambda_{p,q} := \frac{1}{\Gamma} \sigma_{kl} v^{kp} \omega^{lq}$, we have

$$\operatorname{Tr}_{A}\left[\left(\rho_{A}^{\zeta}\right)^{n}\right] = \sum_{p,q} \lambda_{p,q}^{n} \times \left(\frac{|G_{A}||G_{B}|}{|G|}\right)^{n-1}.$$
 (C14)

Substituting this into

$$S_A = \lim_{n \to 1} \frac{1}{1 - n} \log \operatorname{tr}_A \left(\rho_A^{\zeta}\right)^n \tag{C15}$$

yields (49).

3. Single column

In the case of Fig. 4(c), one can deform the logical operators (29) so that η_1^x and η_2^x act exclusively on *B* and γ_2^x acts within *A*, with γ_1^x acting on both *A* and *B*. The term $\text{Tr}_B[\ldots]$ in (C2) is described by

$$\frac{1}{|G|} \sum_{g,g'\in G} (\gamma_{1A}^{x})^{b} (\gamma_{2}^{x})^{d} g_{A} |0\rangle_{AA} \langle 0| (g_{A}g'_{A})^{\dagger} (\gamma_{1A}^{x\dagger})^{b'} (\gamma_{2}^{x\dagger})^{d'}{}_{B} \langle 0| g'_{B}^{\dagger} \\ \times (\eta_{1}^{x})^{a-a'} (\gamma_{1B}^{x})^{b-b'} (\eta_{2}^{x})^{c-c'} |0\rangle_{B}.$$
(C16)

× (η_1) (γ_{1B}) (η_2) $|0\rangle_B$. The inner product gives a constraint

$$g'_B = I_B, \ a = a', \ b = b', \ c = c'.$$
 (C17)

We define orthogonal states by

$$|b,q\rangle_A := \frac{1}{\sqrt{N}} \sum_{d=0}^{N-1} \omega^{qd} \left(\gamma_{1A}^x\right)^b \left(\gamma_2^x\right)^d |0\rangle_A, \qquad (C18)$$

and the density matrix ρ_A^{ζ} is rewritten as

$$\rho_A^{\zeta} = \frac{1}{|G|} \sum_{g \in G, g' \in G_A} \sum_{b,q} \lambda_q^{(b)} g_A | b, q \rangle_{AA} \langle b, q | (g_A g'_A)^{\dagger}$$
(C19)

with

$$\lambda_{q}^{(b)} = \frac{1}{N} \sum_{k=0}^{N-1} \omega^{kq} \sigma_{k}^{(b)}, \quad \sigma_{k}^{(b)} = \sum_{\substack{a,c,d,d'\\d-d'=k \mod N}} \alpha_{ab,cd} \bar{\alpha}_{ab,cd'}.$$
(C20)

By calculating (C15), one obtains (56).

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