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Tensor-entanglement-filtering renormalization approach and symmetry-protected topological order

Zheng-Cheng Gu and Xiao-Gang Wen

Department of Physics, Massachusetts Institute of Technology, Cambridge, Massachusetts 02139, USA

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We study the renormalization group flow of the Lagrangian for statistical and quantum systems by representing their path integral in terms of a tensor network. Using a tensor-entanglement-filtering renormalization approach that removes local entanglement and produces a coarse-grained lattice, we show that the resulting renormalization flow of the tensors in the tensor network has a nice fixed-point structure. The isolated fixedpoint tensors T_{inv} plus the symmetry group G_{sym} of the tensors (i.e., the symmetry group of the Lagrangian) characterize various phases of the system. Such a characterization can describe both the symmetry breaking phases and topological phases, as illustrated by two-dimensional (2D) statistical Ising model, 2D statistical loop-gas model, and 1+1D quantum spin-1/2 and spin-1 models. In particular, using such a (G_{sym}, T_{inv}) characterization, we show that the Haldane phase for a spin-1 chain is a phase protected by the time-reversal, parity, and translation symmetries. Thus the Haldane phase is a symmetry-protected topological phase. The (G_{sym}, T_{inv}) characterization is more general than the characterizations based on the boundary spins and string order parameters. The tensor renormalization approach also allows us to study continuous phase transitions between symmetry breaking phases and/or topological phases. The scaling dimensions and the central charges for the critical points that describe those continuous phase transitions can be calculated from the fixed-point tensors at those critical points.

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I. INTRODUCTION

For a long time, people believe that all phases and all continuous phase transitions are described by symmetry breaking.¹ Using the associated concepts, such as order parameter, long-range correlation, and group theory, a comprehensive theory for symmetry breaking phases and their transitions is developed. However, nature's richness never stops to surprise us. In a study of chiral spin liquid in high-temperature superconductor,^{2,3} we realized that there exist a large class of phases beyond symmetry breaking.⁴ Since the new phases can be characterized by topology-dependent topologically stable ground-state degeneracies, the new type of orders in those phases are named topological orders.⁵ Later, it was realized that fractional quantum Hall states⁶ are actually topologically ordered states.⁷

Recently, it was shown that topological order is actually the pattern of long-range entanglement in the quantum ground state.^{8–10} The possible patterns of long-range entanglement are extremely rich,^{11–13} which indicate the richness of new states of matter beyond the symmetry breaking paradigm. Such rich new states of matter deserve a comprehensive theory.^{14,15}

But how to develop a comprehensive theory of topological order? First we need a quantitative and concrete description of topological order. The topological ground-state degeneracy^{4,7} is not enough since it does not completely characterize topological order. However, the non-Abelian Berry's phases from the degenerate ground states and the resulting representation of modular group provide a more complete characterization of topological order.⁵ It may even completely characterize topological order. There are other ways to (partially) characterize topological order, such as through the fractional statistics of the quasiparticles^{16,17} or through the string operators that satisfy the "zero-law."¹⁸ Those zero-law string operators appear to be related to another characterization of topological order through a gauge-like symmetry defined on a lower dimensional subspace of the space.^{19,20}

But so far, we have not been able to use those characterizations to develop a theory of topological order. So the key to develop a comprehensive theory of topological order is to find a quantitative and easy-to-use description of topological order. For fractional quantum Hall states, the pattern of zeros appears to provide a quantitative and easy-to-use description of topological order.^{21–23} For more general systems, we may use the renormalization group (RG) approach to find a quantitative and easy-to-use description. We will see that the new description based on the RG approach naturally leads to a theory of topological order and phase transitions between topologically ordered states.

RG approach^{24,25} is a powerful method to study phases and phase transitions for both quantum and statistical systems. For example, for a quantum system, one starts with its Lagrangian and path integral that describe the system at a certain cutoff scale. Then one integrates out the short distance fluctuations in the path integral to obtain a renormalized effective Lagrangian at a longer effective cutoff scale. Repeating such a procedure, the renormalized effective Lagrangian will flow to an isolated fixed-point Lagrangian that represents a phase of the system. As we adjust coupling constants in the original Lagrangian, we may cause the renormalization flow to switch from one fixed point to another fixed point, which will represent a phase transition. Since the fixed-point Lagrangian is unique for each phase and the phase transition is associated with a change in fixed-point Lagrangian, it is natural to use the fixed-point Lagrangian to describe various phases.

But can we use the fixed-point Lagrangian to quantitatively describe topological order? At first sight, the answer appears to be no. When a quantum spin system (or a qubit system) has a topologically ordered ground state, its low energy effective Lagrangian (i.e., the fixed-point Lagrangian) can be a Z_2 gauge theory,^{26–30} a U(1) gauge theory,^{29,31–33} a SU(2) gauge theory,²⁹ a QED theory [with U(1) gauge field and massless fermion field],^{29,34,35} a QCD theory (with non-Abelian gauge fields and massless fermion fields),^{29,34,36,37} a Chern-Simons theory,^{3,38} etc. In the standard implication of RG approach (within the framework of field theory), a bosonic Lagrangian that describes a quantum spin system cannot flow to all those very different possible fixed-point Lagrangian with very different fields (such as gauge fields and anticommuting fermionic fields).

Since lattice spin systems do have those very different topologically ordered phases, we see that the standard implementation of the RG approach is incomplete and needed to be generalized to cover those existing (and exciting) renormalization flows that end at very different fixed-point Lagrangian. In this paper, we will describe one such generalization—tensor-network renormalization approach that should allow us to study any phases (both symmetry breaking and topological phases) of any systems (with a finite cutoff). The resulting fixed-point Lagrangian from the tensor-network renormalization approach give us a quantitative description of topological orders (and the symmetry breaking orders).

The tensor-network renormalization approach is based on an observation that the space-time path integral of a quantum spin system or the partition function of a statistical system on lattice can be represented by a tensor trace over a tensor network after we discretize the space-time and the fields,

$$\int D\varphi(\mathbf{x},t)\exp\left(-\int \mathcal{L}\right) = \sum_{\{\phi_i\}} \prod T_{\phi_i\phi_j\phi_k\phi_l} \equiv t \operatorname{Tr} \otimes_i T,$$
(1)

where *i* label the discretized space-time points and ϕ_i is the discretized field [for detail, see Eq. (2)]. The different choices of tensor $T_{\phi_i\phi_j\phi_k\phi_i}$ in the tensor network correspond to different Lagrangians and/or partition functions. We like to point out that in addition to use it to describe path integral or partition function, tensor network can also be used to describe many-body wave functions.³⁹ Based on such a connection, many variational approaches were developed.³⁹⁻⁴⁵

Since every physical system is described by path integral, so if we can calculate the tensor trace, we can calculate anything. Unfortunately, calculating the tensor trace (tTr) is an exponential hard problem in 1+1D (one dimensional) and higher dimensions.⁴⁶ In Ref. 47, Levin and Nave introduced a tensor renormalization group procedure where a coarse graining of the tensor network is performed to produce a new tensor on the coarse-grained network. This leads to an efficient and accurate calculation of path integral (or partition function) and other measurable quantities. Such a method opens a new direction in numerical calculation of many-body systems.

The above tensor renormalization group procedure also gives rise to a renormalization flow of the tensor. However, the fixed-point tensor obtained from such a flow is not isolated: even starting with different Lagrangians (i.e., the tensors) that describe the same phase, the renormalization flow of those Lagrangians will end up at different fixed-point tensors. This means that the renormalization procedure does not filter out (or integrate out) all the short distance fluctuations so that the fixed-point tensors still contain short distance nonuniversal information. In this paper, we will introduce an improved tensor renormalization group procedure by adding a filtering operation that removes the local entanglements. We show that this additional entanglement-filtering procedure fixes the above problem at least for the cases studied in this paper. We like to mention that real-space renormalization transformations for many-body wave functions have been proposed in Refs. 41, 48, and 49 using the tensor-network representation of many-body quantum wave function. Those real-space renormalization transformations remove local entanglements in the many-body wave function. In some sense, our tensor renormalization group procedure also removes "local entanglements" in the tensor network that represents the path integral.

To test the new tensor-network renormalization method, we study the fixed-point tensors that correspond to the two phases of two-dimensional (2D) statistical Ising model which has a Z_2 spin-flip symmetry. We find that the fixed-point tensor for the disordered phase is a trivial tensor T^{TRI} where the dimensions of all the tensor indices are equal to 1 (we will refer such a trivial tensor as a dimension-one tensor which is represented by a number). While the fixed-point tensor for the symmetry breaking (spin ordered) phase is a dimension-two tensor T^{Z_2} which is a direct sum of two dimension-one tensors: $T^{Z_2}=T^{\text{TRI}}\oplus T^{\text{TRI}}$. We believe that this is a general feature for symmetry breaking phase: a symmetry breaking phase with *n* degenerate states is represented by fixed-point tensors, which is a direct sum of *n* dimension-one trivial tensors.

Since the Z_2 spin-flip symmetry is crucial for the existence and the distinction of the disordered phase and the spin ordered phase, we should use a pair, the symmetry group $G_{\text{sym}}=Z_2$ and the fixed-point tensors $T_{\text{inv}}=T^{\text{TRI}}$, T^{Z_2} , to characterize the two phases. So the disordered phase is characterized by (Z_2, T^{TRI}) and the spin ordered phase is characterized by (Z_2, T^{TRI}) . It turns out that the characterization based on the symmetry group and the fixed-point tensor, $(G_{\text{sym}}, T_{\text{inv}})$, is very general. It can describe both symmetry breaking phases and topological phases. This is one of the main results of this paper.

Even at the critical point between the symmetry breaking phases and/or topological phases, the tensor renormalization approach produces fixed-point tensors (which now have an infinite dimension). From those fixed-point tensors or, more precisely, from a finite-dimension approximation of those fixed-point tensors, we can calculate the critical properties of the critical points with an accuracy of 0.1%. Using such a method, we confirm that the phase transition between the disordered and ordered phases of the Ising model is a central charge c=1/2 critical point.

We also study a 2D statistical loop-gas model. We find that for the small-loop phase, the fixed-point tensor is the dimension-one tensor T^{TRI} , while for the large loop phase, the fixed-point tensor T^{LL} is a tensor with dimension 2 [see Eq. (24)]. However, unlike the symmetry breaking phase, the fixed-point tensor for the large loop phase is *not* a direct sum of two dimension-one tensors: $T^{\text{LL}} \neq T^{\text{TRI}} \oplus T^{\text{TRI}}$. This is consistent with the fact that, as a *statistical* system, the small-loop phase and the large loop phase have the same symmetry and the phase transition is not a symmetry breaking transition. The example demonstrates that fixed-point tensors can be used to study phases and phase transitions beyond symmetry breaking. The phase transition between the two phases is again a central charge c=1/2 critical point.

Next, we study a spin-1/2 model $H=\sum_i(-\sigma_i^x \sigma_{i+1}^x)$, with $Z_2^x \times Z_2^z$ symmetry generated by $i\sigma^x$ and $i\sigma^z$. The spin-1/2 model has two different Z_2 symmetry breaking phases described by, in terms of the new notation, $(Z_2^x \times Z_2^z, T^{Z_2})$ and $(Z_2^x \times Z_2^z, T^{LL})$. The $(Z_2^x \times Z_2^z, T^{Z_2})$ phase breaks the Z_2^x symmetry, while the $(Z_2^x \times Z_2^z, T^{LL})$ phase breaks the Z_2^z symmetry. The phase transition between the two Z_2 symmetry breaking phases is a continuous phase transition that is beyond the Landau symmetry breaking paradigm. We find that the tensor-network renormalization approach can describe such a continuous phase transition. The phase transition between the two phases is a central charge c=1 critical point.

Last we study the T=0 phases of quantum spin-1 model in 1+1D, $H=-\sum_i [S_i \cdot S_{i+1}+U(S_i^z)^2+BS_i^x]$, and found that the model has four different phases for different (U,B): a trivial phase, two Z_2 symmetry breaking phases, and the Haldane phase.^{50–52} We find that the Haldane phase is described by a dimension-four fixed-point tensor T^{Haldane} [see Eq. (31)]. T^{Haldane} is not a direct sum of dimension-one tensors. From the fixed-point tensor T^{Haldane} , we confirm that the Haldane phase has a finite energy gap and the ground state is not degenerate. Thus, despite the fixed-point tensor to be a dimension-four tensor, the Haldane phase, like the trivial phase, does not break any symmetry (as expected).

The tensor-entanglement-filtering renormalization (TEFR) approach allows us to study the stability of the Haldane phase (i.e., the distinction between the Haldane phase and the trivial phase) under a very general setting. Our calculation shows that the fixed-point tensor T^{Haldane} and, hence, the Haldane phase are robust against any perturbations that have time-reversal, parity, and translation symmetries.⁵³ So the Haldane phase is a symmetry-protected topological phase characterized by the pair ($G_{TPT}, T^{\text{Haldane}}$) where G_{TPT} is the symmetry group generated by time-reversal, parity, and translation symmetries.

People have been using the degenerate spin-1/2 degrees of freedom at the ends of spin-1 chain^{54–56} and string order parameter^{57,58} to characterize Haldane phase. Both degenerate boundary spin-1/2 degrees of freedom and the string order parameter can be destroyed by perturbations that break spin rotation symmetry but have time-reversal, parity, and translation symmetries. So if we used the degenerate spin-1/2 degrees of freedom and/or string order parameter to characterize the Haldane phase, we would incorrectly conclude that the Haldane phase is unstable against those perturbations. Since the Haldane phase is indeed stable against those perturbations, we conclude that the (G_{TPT} , $T^{Haldane}$) characteriza-



FIG. 1. A graphic representation of a tensor-network on a square lattice. A vertex represent a tensor T_{ijfe} and the legs of a vertex carries the indices *i*, *j*, etc. Each link carries the same index. The indices on the internal links are summed over, which defines the tensor trace.

tion is more general. We also find that an arbitrary small perturbation that breaks parity symmetry will destabilize the Haldane phase. This confirms a previously known result.⁵⁹

In the first few sections of this paper, we will review and introduce several tensor-network renormalization approaches. Then we will discuss physical picture about phases and phase transitions emerged from the tensor-network renormalization approach through several simple examples.

II. REVIEW OF A TENSOR NETWORK RENORMALIZATION METHOD

In the following review, we will concentrate on 1+1D quantum systems. However, our discussion can be generalized to 2D statistical systems and the systems in higher dimensions.

To describe the tensor renormalization approach introduced in Ref. 47, we first note that, after we discretize the space time, the partition function represented by a space-time path integral Tr $e^{-\beta H} = \int \exp(-\int \mathcal{L})$ (in imaginary time) can be written as a tensor trace over a tensor network (see Fig. 1 for an example on 2D square lattice),

$$\operatorname{Tr} e^{-\beta H} = \sum_{ijkl...} T_{jfei} T_{hgjk} T_{qklr} T_{lits} \cdots = t \operatorname{Tr} \otimes_i T, \qquad (2)$$

where the indices of the tensor run from 1 to D (in other words the tensor T is a rank-four dimension-D tensor). Choosing different tensors T corresponds to choosing different Lagrangian \mathcal{L} . We see that calculating the tensor trace allows us to obtain properties of any physical systems.

Unfortunately, calculating the tTr is an exponential hard problem in 1+1D and higher dimensions.⁴⁶ To solve this problem, Levin and Nave⁴⁷ introduced an approximate real-space renormalization group approach which accelerates the calculation exponentially. The basic idea is quite simple and is illustrated in Fig. 2. After finding the reduced tensor T'', we can express $tTr[T \otimes T \cdots] \approx tTr[T' \otimes T'' \cdots]$ where the second tensor trace only contains a quarter of the tensors in the first tensor trace. We may repeat the procedure until there are only a few tensors in the tensor trace. This allows us to reduce the exponential long calculation to a polynomial long calculation.

The actual implementation of the renormalization is a little more complicated.^{45,47} For a uniform tensor network on



FIG. 2. The tensor *T* in the tensor network (a) has a dimension *D*. After we combine the two legs on each side into a single leg, the four linked tensors in (a) can be viewed as a single tensor *T'* in (b) with dimension D^2 . *T'* can be approximately reduced to a "smaller" tensor *T''* in (c) with dimension *D* and satisfies $\text{tTr}[T' \otimes T' \cdots] \approx \text{tTr}[T' \otimes T' \cdots]$.

2D square lattice (see Fig. 1), the renormalization group procedure can be realized in two steps. The first step is decomposing the rank-four tensor into two rank-three tensors. We do it in two different ways on the sublattices purple and green [see Fig. 3(a)]. On purple sublattice, we have T_{ruld} $=\sum_{s=1}^{D^2} S_{1uls}S_{3drs}$ and on green sublattice, we have T_{ruld} $=\sum_{s=1}^{D^2} S_{2lds}S_{4rus}$. Note that r, l, u, d run over D values, while srun over D^2 values. For such a range of s, the decomposition can always be exact.

Next we try to reduce the range of *s* through an approximation.⁴⁷ Say, on purple sublattice, we view T_{ruld} as a matrix $M_{lu;rd}^{p} = T_{ruld}$ and do singular value decomposition (SVD),

$$M^{\rm p} = U\Lambda V^{\dagger}, \quad \Lambda = \begin{pmatrix} \lambda_1 & \\ & \lambda_2 \\ & & \ddots \end{pmatrix}, \tag{3}$$

where $\lambda_1 > \lambda_2 > \lambda_3$ We then keep only the largest D_{cut} singular values λ_s and define $S_{3drs} = \sqrt{\lambda_s} V_{s,rd}^{\dagger}$, $S_{1uls} = \sqrt{\lambda_s} U_{lu,s}$. Thus, we can approximately express T_{ruld} by two rank-three tensors S_1 , S_3 ,

$$T_{ruld} \simeq \sum_{s=1}^{D_{cut}} S_{1uls} S_{3drs}.$$
 (4)

Similarly, on green sublattice we may also define T_{ruld} as a matrix $M_{ld;ru}^{g}$, do SVDs, keep the largest D_{cut} singular values, and approximately express T_{ruld} by two rank-three tensors S_2, S_4 ,

$$\mathbb{T}_{\alpha\beta\mu\nu} \simeq \sum_{\gamma=1}^{D_{\text{cut}}} S_{2\nu\alpha\gamma} S_{4\beta\mu\gamma}.$$
 (5)



FIG. 3. (Color online) RG transformation of tensor network produces a coarse-grained tensor network.



FIG. 4. (Color online) (a) We represent the original rank-four tensor by two rank-three tensors, which is an *approximate* decomposition. (b) Summing over the indices around the square produces a single tensor T'. This step is *exact*.

After such decompositions, the square lattice is deformed into the form in Fig. 3(b). The second step is to simply contract the square and get a new tensor T' on the coarsegrained lattice [see Fig. 4(b)]. The dimension for the reduced tensor T' is only D_{cut} which can be chosen to be D or some other values. This defines the (approximated) renormalization flow of the tensor network based on the SVD. Let us call such an approach SVD based tensor renormalization group (SVDTRG) approach.

III. PHYSICAL MEANING OF FIXED-POINT TENSORS

Why the SVDTRG flow fails to produce an isolated fixedpoint tensors? Levin pointed out that one large class of tensors—corner double-line (CDL) tensors—are fixed points of the SVDTRG flow.^{47,60} A CDL tensor has a form [see Fig. 5(a)]

$$T_{ruld}(M^1, M^2, M^3, M^4) = T_{r_1 r_2, u_2 u_1, l_2 l_1, d_1 d_2}$$
$$= M^1_{r_2 u_1} M^2_{u_2 l_1} M^3_{l_2 d_1} M^4_{d_2 r_1}, \qquad (6)$$

where the pair (r_1, r_2) represents the index r, etc. Such type of tensor (assuming \sqrt{D} =integer, r_1 , etc., run from 1 to \sqrt{D}) has exact decomposition,

$$T_{ruld} = \sum_{s=1}^{D_{cut}} S_{3drs} S_{1lus} = \sum_{s=1}^{D_{cut}} S_{2lds} S_{4rus},$$

with $D_{\text{cut}}=D$. This makes the CDL tensor to be the fixedpoint tensor under the SVDTRG flow at least when M^1 $=M^2=M^3=M^4$. If the initial tensor is changed slightly, the tensor may flow to a slight different CDL tensor, and hence the fixed point is not isolated.

A tensor network of CDL tensor is plotted in Fig. 5(b). If the system described by a tensor network does flow to a



FIG. 5. (a) A CDL tensor. (b) A network of CDL tensor.

tensor network of CDL tensor, what does this means? From Fig. 5(b), we see that, at the fixed point described by the CDL tensor, the degrees of freedom form clusters that correspond to the disconnected squares. Only the degrees of freedom within the same square interact and there are no interactions between different clusters. Such a fixed-point tensor (i.e., Lagrangian) describes a system with no interaction between neighbors and has a finite energy gap. So the system is in a trivial phase without symmetry breaking and without topological order (in the sense of long-range entanglement). The $h \gg J$ phase of the transverse-field Ising model,

$$H_{\text{tIsing}} = -\sum_{i} \left(J \sigma_i^x \sigma_{i+1}^x + h \sigma_i^z \right), \tag{7}$$

is such a phase. We can check that the tensor network that describes the above transverse-field Ising model indeed flows to a network of CDL tensor when $h \ge J$.

On the other hand, the low energy physics of such phases is really trivial. J=0 and $h=\infty$ limit, the partition function of the transverse-field Ising model is described by a dimensionone tensor,

$$T_{ruld}^{\text{TRI}} = \delta_{r,1} \delta_{u,1} \delta_{l,1} \delta_{d,1}.$$
 (8)

Thus we expect that the fixed-point tensor for the $h \ge J$ phase should be the above trivial dimension-one tensor. Unfortunately, the renormalization flow generated by SVDTRG fails to produce such simple fixed-point tensor.

The $h \ll J$ phase of the transverse-field Ising model is a symmetry breaking phase with two degenerate ground states. Under the SVDTRG flow, the tensor network of the transverse-field Ising model will flow to a tensor network described by a tensor of the form

$$T = T_{\text{CDL}} \oplus T_{\text{CDL}},\tag{9}$$

where T_{CDL} is a CDL tensor. On the other hand, in the coarse-grained limit the symmetry breaking phase should be described by a simpler tensor [Eq. (9)] where T_{CDL} is the dimension-one tensor T^{TRI} given by Eq. (8).

However, the SVDTRG flow does a remarkable job of squeezing all the complicated entanglement into a single plaquette. Note that the initial tensor network may contain entanglement between lattice sites separated by a distance less than the correlation length. A CDL tensor contains only entanglement between nearest-neighbor lattice sites, while the ideal fixed-point tensor [Eq. (8)] has no entanglement at all (not even with next neighbors).

What should be the fixed-point tensor for topological phases under SVDTRG flow? The essence of topological phases is pattern of long-range entanglement.^{8–10} Since the CDL tensor has no long-range entanglement, it cannot be used to describe topologically ordered phases. We believe that topological phases are described by fixed-point tensors of the following form (see Fig. 6):

$$T = T_{\text{CDL}} \otimes T_{top}, \tag{10}$$

where T_{CDL} is CDL tensor and T_{top} is a tensor that cannot be decomposed further as $T_{top}=T'_{\text{CDL}}\otimes T'_{top}$. Here T_{top} will be the fixed-point tensor that characterizes the topological order.



FIG. 6. The fixed-point tensor and its tensor network for topological phases.

IV. FILTERING OUT THE LOCAL ENTANGLEMENTS

We have seen that the SVDTRG flow can produce a network of CDL tensor which squeezes the nontopological short-range entanglement into single plaquette. To obtain ideal fixed-point tensor [such as the one in Eq. (8)] we need to go one step further to remove the corner entanglement represented by the CDL tensor, i.e., we need to reduce corner matrix M^i in the CDL tensor [see Eq. (8)] to a simpler form,

$$M_{ab}^{i} \to \delta_{a,1} \delta_{b,1}.$$
 (11)

In the following, we will describe one way to do so.

In this approach, we modify the SVDTRG transformation described by Fig. 3 to a new transformation described by Fig. 7. We will call the new approach TEFR approach. In the TEFR transformation, we first deform the square lattice [Fig. 7(a)] into the lattice [Fig. 7(b)] by expressing the rank-four tensor T in terms of two rank-three tensors S_1, S_3 or S_2, S_4 [see Fig. 4(a)]. Then we replace the four rank-three tensors S_1, S_2, S_3, S_4 by a new set of simpler rank-three tensors S'_1, S'_2, S'_3, S'_4 so that the tensor trace of S'_1, S'_2, S'_3, S'_4 (approximately) reproduces that of S_1, S_2, S_3, S_4 , as shown in Fig. 8. In this step, we try to reduce the dimensions of the rank-three tensors from $D \times D \times D$ of S_i to $D \times D \times D'$ of S'_i with D'< D. If T is a CDL tensor, S_1, S_2, S_3, S_4 will also be CDL tensors [see Fig. 10(a)]. In this case, the above procedure will result in a new set of tensors S'_1, S'_2, S'_3, S'_4 with a lower dimension, as shown in Fig. 10(b). (For details, see Appendix, Sec. 1.) We then use the SVD (see Fig. 9) to deform the lattice in Fig. 7(c) to that in Fig. 7(d). Last we trace out the indices on the small squares in Fig. 7(d) to produce new tensor network on a coarse-grained square lattice [see Fig. 7(e)] with a new tensor T'. The tensor T' will have a lower local entanglement than T. In fact, when T is a CDL tensor, the combined procedures described in Figs. 3 and 7 (see also Figs. 10 and 11) can reduce T to a dimension-one tensor. Appendix, Sec. 1 contains a more detailed description of TEFR method.

V. EXAMPLES

In this section, we will apply the TEFR method to study several simple well known statistical and quantum systems to test the TEFR approach. We will calculate the free energy (for statistical systems) or the ground-state energy (for quantum systems). The singularities in free energy and/or groundstate energy in the large system limit will indicate the pres-



FIG. 7. (Color online) The TEFR procedure. We first deform the tensor network on square lattice (a) to the tensor network on lattice (b) using the transformations in Fig. 4(a). We then replace the four rank-three tensors S_1, S_2, S_3, S_4 in (b) by a new set of simpler rank-three tensors S'_1, S'_2, S'_3, S'_4 in (c) while keeping the tensor trace of the tensor network approximately the same (see Fig. 8). The index on a solid link has a dimension D. The index on a dashed link has a dimension D'. Thus the tensor S_i has a dimension $D \times D \times D$, while the tensor S'_i has a dimension $D \times D \times D'$. The SVD method is used to go from (a) to (b) [see Fig. 4(a)] and from (c) to (d) (see Fig. 9). The four S''_i 's on the small squares in (d) are traced over to obtain T' in (e).

ence of phase transitions. We can also calculate the fixedpoint tensors under the TEFR flow for each phase. The fixedpoint tensors will allow us to identify those phases as symmetry breaking and/or topological phases. If we find a continuous phase transition in the above calculation, we can also use the fixed-point tensor at the critical point to calculate the spectrum of the scaling dimensions h_i and the central charge c for the critical point (see Appendix, Sec. 2).



FIG. 8. (Color online) A tensor trace of a product of four rankthree tensors S_1, S_2, S_3, S_4 gives rise to one rank-four tensor. Such a rank-four tensor can also be produced (approximately) by a different set of rank-three tensors S'_1, S'_2, S'_3, S'_4 . We may choose S'_i trying to minimize the dimension of the index on the dashed links. Or we may choose S'_i trying to optimize some other quantities.



FIG. 9. (Color online) The SVD method is used to implement the above two transformations.

A. 2D statistical Ising model

Our first example is the 2D statistical Ising model on the square lattice described by

$$E = -\sum_{\langle i,j \rangle} \sigma_i \sigma_j, \quad \sigma_i = \pm 1,$$
(12)

where $\Sigma_{\langle ij \rangle}$ sums over nearest neighbors. The partition function is given by $Z=\text{Tr}(e^{-\beta H})$. Such a partition function can be expressed as a tensor trace over a tensor network on a (different) square lattice where the tensor *T* is given by

$$T_{1,2,1,2}^{\text{Ising}} = e^{-4\beta}, \quad T_{2,1,2,1}^{\text{Ising}} = e^{-4\beta},$$
$$T_{1,1,1,1}^{\text{Ising}} = e^{4\beta}, \quad T_{2,2,2,2}^{\text{Ising}} = e^{4\beta},$$
others = 1. (13)

Note that the spins are located on the links of the tensor network [i.e., i in Eq. (12) labels the links of the tensor network].

The statistical Ising model has a Z_2 spin-flip symmetry. Such a symmetry implies that the tensor is invariant under the following Z_2 transformation:

$$T_{ruld}^{\text{Ising}} = W_{rr'} W_{uu'} W_{ll'} W_{dd'} T_{r'u'l'd'}^{\text{Ising}},$$
(14)

where

$$W = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}.$$

We performed 14 steps of TEFR iteration described in Fig. 7 with $D_{\text{cut}}=32$. The calculated free energy (per site) $F=-\beta^{-1} \ln(Z)/N=-T \ln(Z)/N$ is presented in Fig. 12. For temperature T>2.38, the fixed-point tensor is a dimension-one tensor,

$$T_{1111}^{\text{TRI}} = 1$$
, others = 0. (15)

For temperature T < 2.19, the fixed-point tensor is a dimension-two tensor,

$$T_{1111}^{Z_2} = 1, \quad T_{2222}^{Z_2} = 1, \quad \text{others} = 0.$$
 (16)

For temperature 2.22 < T < 2.32, the fixed-point tensor is complicated with many nonzero elements. (We like to point

$$\frac{\Box}{T} = \frac{\Box}{S_3} = \frac{S_1}{S_4} = \frac{S_1}$$

FIG. 10. TEFR transformation of a CDL tensor. Here the index on a dashed line has a dimension 1.



FIG. 11. (Color online) A tensor network of CDL tensor can be reduced to a tensor network of dimension-one tensor. (a) \rightarrow (b) uses the transformation in Fig. 10(a). (b) \rightarrow (c) uses the transformation in Fig. 10(b). (c) \rightarrow (d) uses the transformation in Fig. 9. (d) \rightarrow (e) uses the transformation in (b) \rightarrow (c) in Fig. 4. Here the index on a dashed line has a dimension 1.

out that the above fixed-point tensors are actually invariant fixed-point tensors. For details, see Appendix, Sec. 2.)

To visualize the structure of fixed-point tensors more quantitatively in different regions, we introduce the following two quantities (see Fig. 13),

$$X_1 = \frac{\left(\sum_{ru} T_{ruru}\right)^2}{\sum_{ruld} T_{rulu} T_{ldrd}}, \quad X_2 = \frac{\left(\sum_{ru} T_{ruru}\right)^2}{\sum_{ruld} T_{ruld} T_{ldru}},$$
(17)

for the fixed-point tensors. Note that X_1 and X_2 is independent of the scale of the tensor $T \rightarrow \Gamma T$. Also X_1 and X_2 is invariant under the field-redefinition transformation [Eq. (19)]. When *T* is the direct sum of *n* dimension-one trivial tensors, we find that $X_1 = X_2 = n$. So a jump in X_1 and/or X_2 indicates a phase transition.

We plotted X_1 , X_2 and the central charge obtained from Eq. (A18) for the fixed-point tensors T_{inv} at different temperatures in Fig. 14. (Note that $\tau_0 = \tau_1 = i$ here.) We see that the different values of X_1 and X_2 mark the different phases and the steps in X_1 and X_2 , as well as the peak in *c* marks the point of continuous transition.

From the structure of the fixed-point tensor, we see that the low-temperature phase T < 2.19 is a Z_2 symmetry break-



FIG. 12. (Color online) The temperature dependence of free energy (per site) F(T) and $F'' = \partial^2 F / \partial T^2$ for the 2D statistical Ising model. A phase transition can be seen around the temperature $T \approx 2.3$. The vertical line marks the exact $T_c = 2.269$ 19.



FIG. 13. The graphic representations of (a) $\Sigma_{ru} T_{ruru}$, (b) $\Sigma_{ruld} T_{rulu} T_{ldrd}$, and (c) $\Sigma_{ruld} T_{ruld} T_{ldru}$.

ing phase (where the fixed-point tensor has a form $T^{Z_2} = T^{\text{TRI}} \oplus T^{\text{TRI}}$) and the high-temperature phase T > 2.38 is a trivial disordered phase. Since $\partial F / \partial T$ appears to be continuous and $\partial^2 F / \partial T^2$ appears to have an algebraic divergence at the transition point, all those suggest that the phase transition is a continuous transition, which is a totally expected result. From the Onsager's solution⁶¹ sinh $(2/T_c) = 1$, one obtains the exact critical temperature to be $T_c = \frac{2}{\ln(\sqrt{2}+1)} = 2.269$ 19, which is consistent with our numerical result.

There is another interpretation of the fixed-point tensors. The fixed point in the low-temperature phase $T^{Z_2}=T^{\text{TRI}} \oplus T^{\text{TRI}}$ happens to be the initial tensor [Eq. (13)] in the zero-temperature limit $\beta = \infty$. Thus the flow of the tensor in the low-temperature phase can be viewed as a flow of the temperature $\beta \rightarrow \infty$. Using the same reasoning, we expect that the flow of the tensor in the high-temperature phase can also be viewed as a flow of temperature $\beta \rightarrow 0$ in the opposite direction. This would suggest that the fixed-point tensor should have a form

$$\widetilde{T}_{ruld}^{\text{TRI}} = \frac{1}{4}, \quad r, u, l, d = 1, 2,$$
 (18)

which is very different from the trivial dimension-one tensor T^{TRI} in Eq. (15). We will explain below that the two tensors \tilde{T}^{TRI} and T^{TRI} are equivalent.

To see the equivalence, we would like to point out that two tensors T and T' related by

$$T'_{r'u'l'd'} = (A^{-1})_{l'l} (B^{-1})_{u'u} T_{ruld} A_{rr'} B_{dd'}$$
(19)

give rise to same tensor trace for any square-lattice tensor network. In particular, if the singular values in the SVD decomposition in Fig. 4(a) have degeneracies, the coarsegrained tensor will have the ambiguity described by the above transformation with A and B being orthogonal matrices. Physically, the transformation [Eq. (19)] corresponds to



FIG. 14. (Color online) The temperature dependence of X_1 , X_2 , and the central charge *c* for the 2D statistical Ising model. The vertical line marks the exact T_c =2.269 19.

a field redefinition. One can check explicitly that

$$\widetilde{T}_{ruld}^{\mathrm{TRI}} = R_{rr'}^{-1} R_{uu'}^{-1} T_{r'u'l'd'}^{\mathrm{TRI}} R_{l'l} R_{d'd},$$

where

$$R = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1\\ 1 & -1 \end{pmatrix}.$$
 (20)

Thus, \tilde{T}^{TRI} and T^{TRI} are equivalent under a field redefinition.

We have seen that the coarse-grained tensors after the TEFR transformations may contain a field-redefinition ambiguity [Eq. (19)]. As a result, the Z_2 transformation on the coarse-grained lattice may be different from that on the original lattice. The Z_2 transformation on the coarse-grained lattice is related to the original Z_2 transformation by a fieldredefinition transformation.

However, one can choose a basis on the coarse-grained lattice by doing a proper field-redefinition transformation such that the Z_2 spin-flip transformation has the same form on the coarse-grained lattice as that on the original lattice. In such a basis, the disordered and the Z_2 symmetry breaking phases are described by $(G_{\text{sym}}, T_{\text{inv}}) = (Z_2, \tilde{T}^{\text{TRI}})$ and $(G_{\text{sym}}, T_{\text{inv}}) = (Z_2, T^{Z_2})$, respectively, where the Z_2 transformation is given by Eq. (14). This example shows how to use the pair $(G_{\text{sym}}, T_{\text{inv}})$ to describe the two phases of the statistical Ising model.

Next we like to use the fixed-point tensor T_{inv} at the critical point to calculate the critical properties of the above continuous transition. We can use Eq. (A14) to obtain the invariant fixed-point tensor $T_{inv}^{(i)}$ even away from the critical point. From $T_{inv}^{(i)}$ we calculate the matrices M^{ud} , M^{lr} , M^{ldru} , and M^{lurd} from Eq. (A15) and obtain the eigenvalues λ_n of those matrices. Using those eigenvalues, we can calculate the central charge and the scaling dimensions from Eq. (A18), where $\tau_0 = \tau_1 = i$ for the Ising model discussed here. By choosing $D_{cut}=64$ and doing nine iterations of Fig. 7 (that correspond to a system with 1024 spins) at the critical temperature $T_c = \frac{2}{\ln(\sqrt{2}+1)}$, we find that

| с | h_1 | h_2 | h_3 | h_4 |
|---------|---------|---------|---------|----------|
| 0.49942 | 0.12504 | 0.99996 | 1.12256 | 1.12403, |
| 1/2 | 1/8 | 1 | 9/8 | 9/8 |

where we also listed the exact values of the central charge and the scaling dimensions. Such a calculation takes about 10 h on a desktop computer with 2.4 GHz Intel Q6600 CPU and Intel Fortran compiler.

We would like to mention that the TEFR approach is not supposed to work well at the critical point where the truncation error of keeping only a finite D_{cut} singular values is large. It turns out that for $D_{\text{cut}}=64$, the estimated relative truncation error is less than 10^{-5} through nine iterations. This allows us to obtain quite accurate central charge and scaling dimensions. The truncation error will grow for more iterations. As a result, the agreement with the exact result will worsen.



FIG. 15. (Color online) The temperature dependence of X_1 , X_2 , and the central charge *c* for the 2D statistical loop-gas model. The vertical line marks the exact T_c =2.269 19.

From this example, we see that the TEFR approach is an effective and efficient way to study phases and phase transitions. The TEFR approach allows us to identify spontaneous symmetry breaking from the structure of fixed-point tensors. The critical properties for the continuous phase transition can also be obtained from the fixed-point tensor at the critical point.

B. 2D statistical loop-gas model

Next we consider a 2D statistical loop-gas model on square lattice. The loop-gas model can be viewed as an Ising model with spins on the links of the square lattice. However, the allowed spin configurations must satisfies the following hard constraint: the number of down spins next to each vertex must be even. Here an up spin is viewed as the absence of a loop segment on the link and a down spin is viewed as the presence of a loop segment on the link. The energy of spin configuration is given by

$$E = -\sum_{\langle i \rangle} \sigma_i.$$
 (21)

The partition function $Z=Tr(e^{-\beta H})$ can be expressed as a tensor trace over a tensor network on the square lattice where the tensor *T* is given by

$$T_{ruld}^{\rm LG} = e^{\beta(2d-3)/2} e^{\beta(2u-3)/2} e^{\beta(2r-3)/2} e^{\beta(2l-3)/2} \times \delta_{d+u+r+l=\rm even},$$
(22)

where the indices take the values r, u, l, d=1, 2 and

$$\delta_{d+u+r+l=\text{even}} = 1$$
 if $d+u+r+l=\text{even}$, others = 0.

To use the TEFR approach to calculate the partition function, it is very important to implement the TEFR in such a way that the closed-loop condition

$$T_{ruld} = 0$$
 when $r + u + l + d = \text{odd}$,
 $S_{rul} = 0$ when $r + u + l = \text{odd}$ (23)

are satisfied at every step of iteration.

We performed 14 steps of TEFR iteration described in Fig. 7 with $D_{\text{cut}}=32$. Figure 15 shows the resulting X_1 , X_2 and the central charge $c = \frac{6}{\pi} \ln(\lambda_0)$ obtained from Eq. (A18)

for the invariant fixed-point tensors $(T_{inv})_{ruld}$ at different temperatures. We find that the low-temperature phase is described by the trivial dimension-one fixed-point tensor T^{TRI} [Eq. (8)] as indicated by $X_1=X_2=1$. The high-temperature phase has a nontrivial fixed-point tensor since $X_1=X_2=2$. We find that the invariant fixed-point tensor have a form

$$T_{ruld}^{LL} = \frac{1}{2}$$
 if $l + r + u + d = \text{even}$, $r, u, l, d = 1, 2$,
others = 0. (24)

The two fixed-point tensors T^{TRI} and T^{LL} correspond to the zero-temperature and infinite-temperature limits of the initial tensor T^{LG} in Eq. (22). Both fixed-point tensors satisfy the closed-loop condition [Eq. (23)].

It is well known that the high-temperature phase of the statistical loop-gas model is dual to the low-temperature phase of the statistical Ising model and the low-temperature phase of the statistical loop-gas model is dual to the hightemperature phase of the statistical Ising model. The fixedpoint tensors of the two models are related directly,

$$T_{ruld}^{\text{LL}} = R_{rr'}^{-1} R_{uu'}^{-1} T_{r'u'l'd'}^{2} R_{l'l} R_{d'd},$$
$$T_{ruld}^{\text{TRI}} = R_{rr'}^{-1} R_{uu'}^{-1} \widetilde{T}_{r'u'l'd'}^{\text{TRI}} R_{l'l} R_{d'd},$$

where R is given in Eq. (20). The critical properties at the continuous transition are also identical.

One may wonder if T^{LL} and T^{Z_2} are related by a field redefinition, then why the TEFR flow produces T^{LL} as the fixed-point tensor of the high-temperature phase of the loop gas rather than T^{Z_2} . Indeed, if we implement the TEFR flow in the space of generic tensors, then both T^{LL} and T^{Z_2} , together with many other tensors related by some field redefinitions, will appear as the fixed-point tensors of the hightemperature phase of the loop gas. However, if we implement the TEFR flow in the space of tensors that describe closed loops [i.e., satisfying Eq. (23)], then only T^{LL} will appear as the fixed-point tensor.

C. Generalized 2D statistical loop-gas model

Third, let us discuss a more general 2D statistical loop-gas model described by a tensor network with tensor

$$T_{ruld}^{\text{GLG}} = e^{(r+l+u+d-6)\beta + V\delta_{rlud}}\delta_{d+u+r+l=\text{even}},$$
(25)

where
$$r, u, l, d=1, 2$$
 and δ_{rlud} is given by

$$\delta_{1111} = \delta_{2222} = 1$$
, others = 0.

We note that when V=0, the generalized loop-gas model becomes the one discussed above: $T^{LG}=T^{GLG}|_{V=0}$.

Let us consider the phases of the generalized loop-gas model along the β =0 line. In this case, the loop-gas model also has a Z_2 symmetry: the tensor T^{GLG} is invariant under switching the 1 and 2 indices,



FIG. 16. (Color online) The V dependence of $\ln z = \ln(Z)/N$ and $(\ln z)'' = d^2 \ln z/dV^2$ for the generalized 2D loop-gas model. The sharpest peak for $(\ln z)''$ and the $\ln z$ curve are for a tensor network with $N=4^7$ tensors. Other peaks are for tensor networks with $N=4^6$, $N=4^5$, $N=4^4$, and $N=4^3$ tensors.

$$T_{ruld}^{\text{GLG}} = W_{rr'} W_{uu'} W_{ll'} W_{dd'} T_{r'u'l'd'}^{\text{GLG}},$$
$$W_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}.$$
(26)

We will call such a symmetry Z_2^x symmetry. The closed-loop condition [Eq. (23)] can also be represented as a Z_2 symmetry of the tensor,

$$T_{ruld}^{\text{GLG}} = W_{rr'} W_{uu'} W_{ll'} W_{dd'} T_{r'u'l'd'}^{\text{GLG}},$$
$$W_{z} = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}.$$
(27)

The second Z_2 symmetry will be called Z_2^z symmetry. So along the $\beta=0$ line, the model has a $Z_2^x \times Z_2^z$ symmetry.

We performed the TEFR calculation for the generalized loop-gas model along the $\beta=0$ line with $D_{\text{cut}}=32$. Figure 16 shows the V dependence of the resulting logarithmic partition function per tensor, $\ln z=\ln(Z)/N$ and $d^2 \ln z/dV^2$. Figures 17 and 18 show the V dependence of X_1, X_2 and central charge c for the resulting fixed-point tensor. We see that there is a phase transition at $V_c=1.0985$. The fixed-point tensor for the $V < V_c$ phase is found to be T^{LL} and the fixed-point tensor for the $V > V_c$ phase is found to be T^{Z_2} . The central charge is zero for the two phases on the two sides of the transition.



FIG. 17. (Color online) The V dependence of X_1 , X_2 for the generalized 2D loop-gas model.



FIG. 18. (Color online) The V dependence of the central charge c for the generalized 2D loop-gas model. The sharpest peak is for a tensor network with $N=4^7$ tensors. Other peaks are for tensor networks with $N=4^6$, $N=4^5$, $N=4^4$, and $N=4^3$ tensors. The horizontal line marks c=1 (the expected central charge value at the critical point).

Thus both phases have short-range correlations. Using the $(G_{\text{sym}}, T_{\text{inv}})$ notation, the two phases are characterized by $(Z_2^x \times Z_2^z, T^{\text{LL}})$ and $(Z_2^x \times Z_2^z, T^{Z_2})$.

Since $\partial \ln(Z)/\partial V$ has no discontinuous jump at V_c , the transition is a continuous phase transition. Also the central charge is nonzero at the transition point which suggests that the correlation length diverges at V_c . This also implies that the transition is a continuous phase transition. From Fig. 18, we see that the transition is a central charge c=1 critical point.

We note that the two fixed-point tensors T^{LL} and T^{Z_2} give rise to the same X_1 and X_2 . In fact, T^{LL} and T^{Z_2} are related by a field-redefinition transformation [Eq. (19)],

$$T_{ruld}^{\text{LL}} = R_{rr'}^{-1} R_{uu'}^{-1} T_{r'u'l'd'}^{Z_2} R_{l'l} R_{d'd},$$

where *R* is given in Eq. (20). Since X_1 and X_2 are invariant under the above transformation, so they are the same in the two phases.

The generalized loop-gas model provides us an interesting example that the fixed-point tensors T^{LL} and T^{Z_2} that describe the two phases are related by a field-redefinition transformation. In this case, one may wonder should we view the V $< V_c$ and $V > V_c$ phases as the same phase? A simple direct answer to the above question is no. It is incorrect to just use fixed-point tensors to characterize a phase. We should use the pair (G_{sym}, T_{inv}) to characterize a phase. If we use the $(G_{\rm sym}, T_{\rm inv})$ notation, the two phases are more correctly described by $(Z_2^x \times Z_2^z, T^{LL})$ and $(Z_2^x \times Z_2^z, T^{Z_2})$. The above fieldredefinition transformation transforms T^{LL} to T^{Z_2} . It also transform the symmetry transformations $Z_2^x \times Z_2^z$ to a different form $\tilde{Z}_2^x \times \tilde{Z}_2^z$. Thus the pair $(Z_2^x \times Z_2^z, T^{LL})$ is transformed to $(\tilde{Z}_2^x \times \tilde{Z}_2^z, T^{Z_2})$ under the above field-redefinition transformation. Therefore, we cannot transform the pair (Z_2^x) $\times Z_2^z, T^{LL}$ to $(Z_2^x \times Z_2^z, T^{Z^2})$ using any field-redefinition transformation. This implies that the two phases $(Z_2^x \times Z_2^z, T^{LL})$ and $(Z_2^x \times Z_2^z, T^{Z_2})$ are different if we do not break the Z_2^x $\times Z_2^z$ symmetry. If we do break the $Z_2^x \times Z_2^z$ symmetry, then the two phases are characterized by $(1, T^{LL})$ and $(1, T^{Z_2})$, where 1 represents the trivial group with only identity implying that the system has no symmetry. Without any symmetry, the two labels $(1, T^{LL})$ and $(1, T^{Z_2})$ are related under a field-redefinition transformation. This implies that the two phases $(1, T^{LL})$ and $(1, T^{Z_2})$ are the same phase if the system has no symmetry. In Appendix, Sec. 3, we will give a more detailed and general discussion about the definition of phases and its relation to the symmetry of the system.

D. Quantum spin-1/2 chain and phase transition beyond symmetry breaking paradigm

The generalized loop-gas model is closely related to the following 1D quantum spin-1/2 model,

$$H = \sum_{i} \left[-\sigma_{i}^{x} \sigma_{i+1}^{x} - J \sigma_{i}^{z} \sigma_{i+1}^{z} - h \sigma_{i}^{z} \right].$$
(28)

Note that $\Sigma(1-\sigma_i^z)/2=\Sigma_i n_i \mod 2$ is a conserved quantity and the above model can be viewed as a 1D hard-core boson model with next-neighbor interaction. Such a 1D quantum model can be simulated by a 2D statistical loop-gas model described by a tensor network with tensor T^{GLG} [Eq. (25)]. The loops in the loop gas correspond to the space-time trajectories of the hard-core bosons. The $\beta=0$ limit of the loopgas model corresponds to h=0 limit of the 1D quantum model and V corresponds to J.

In the following, we will concentrate on the h=0 case. When h=0, Hamiltonian (28) has a $Z_2^x \times Z_2^z$ symmetry, where Z_2^x is the Z_2 group generated by σ^x and Z_2^z is the Z_2 group generated by σ^z . Such a $Z_2^x \times Z_2^z$ symmetry corresponds to the $Z_2^x \times Z_2^z$ symmetry discussed in Sec. VI.

When $J < J_c = 1$, the system is in a symmetry breaking phase. We will use a pair of groups $(G_{\text{sym}}, G_{\text{grnd}})$ to label such a symmetry breaking phase, where G_{sym} is the symmetry group of the Hamiltonian and G_{grnd} is the symmetry group of a ground state. So the $J < J_c$ phase of our system is labeled by $(G_{\text{sym}}, G_{\text{grnd}}) = (Z_2^x \times Z_2^z, Z_2^x)$, where the ground state breaks the Z_2^z symmetry but has the Z_2^x symmetry. When $J > J_c$, our system is in a different symmetry breaking phase, which is labeled by $(G_{\text{sym}}, G_{\text{grnd}}) = (Z_2^x \times Z_2^z, Z_2^z)$.

Under the correspondence between the loop gas and our quantum spin-1/2 model, the $V < V_c$ and $V > V_c$ phases of the loop-gas model discussed above will correspond to the above two Z_2 symmetry breaking phases for $J < J_c = 1$ and $J > J_c$, respectively. The results from the loop gas imply that the two Z_2 symmetry breaking phases are connected by a *continuous* phase transition. Such a continuous phase transition can be viewed as two Z_2 symmetry breaking transitions happening at the same point.

Strictly speaking, the above continuous transition does not fit within the standard symmetry breaking paradigm for continuous phase transitions. According to Landau symmetry breaking theory, a continuous phase transition can only happen between two phases labeled by $(G_{\text{sym}}, G_{\text{grnd}})$ and $(G_{\text{sym}}, G'_{\text{grnd}})$, where G_{grnd} is a subgroup of G'_{grnd} or G'_{grnd} is a subgroup of G_{grnd} . We see that the continuous transition between the $(Z_2^x \times Z_2^z, Z_2^x)$ phase and $(Z_2^x \times Z_2^z, Z_2^z)$ phase does not fit within the standard Landau symmetry breaking theory. Such a continuous transition is another example of continuous transitions that is beyond the Landau symmetry breaking



FIG. 19. (a) The graphic representation of $(e^{-\delta \tau H_i})_{m_i m_{i+1}, m'_i m'_{i+1}}$. (b) $e^{-\delta \tau H_i}$ can be viewed as a rank-four tensor.

paradigm.^{29,62–67} Even though the above continuous phase transition does not fit within the standard symmetry breaking theory, such a transition can still be studied through the tensor-network renormalization approach. In particular, the tensor-network renormalization approach reveals that the transition is a central charge c=1 critical point.

E. Quantum spin-1 chain

Last, we consider a quantum spin-1 chain at zero temperature. The Hamiltonian is given by

$$H = \sum_{i} \left[S_{i} \cdot S_{i+1} + U(S_{i}^{z})^{2} + BS_{i}^{x} \right].$$
(29)

The model has a translation symmetry, a Z_2 time-reversal symmetry $S^y \rightarrow -S^y$, and a Z_2 parity symmetry. This imaginary-time path integral of the model can be written as

$$\operatorname{Tr} e^{-\tau H} = \lim_{M \to \infty} \operatorname{Tr} (e^{-\delta \tau H_e} e^{-\delta \tau H_o})^M,$$

where $\delta \tau = \tau / M$ and

$$H_{e} = \sum_{i=\text{even}} \left(S_{i} \cdot S_{i+1} + U(S_{i}^{z})^{2} + \frac{B}{2}(S_{i}^{x} + S_{i+1}^{x}) \right),$$
$$H_{o} = \sum_{i=\text{odd}} \left(S_{i} \cdot S_{i+1} + U(S_{i}^{z})^{2} + \frac{B}{2}(S_{i}^{x} + S_{i+1}^{x}) \right).$$

Note that H_e and H_o are sum of nonoverlapping terms and thus

$$e^{-\delta\tau H_e} = \prod_{i=\text{even}} e^{-\delta\tau H_i},$$
$$e^{-\delta\tau H_o} = \prod_{i=\text{odd}} e^{-\delta\tau H_i}.$$

 H_i in the above expression has a form

$$H_i = S_i \cdot S_{i+1} + U(S_i^z)^2 + \frac{B}{2}(S_i^x + S_{i+1}^x).$$

The matrix elements of $e^{-\delta \tau H_i}$ can be represented by a rank-four tensor (see Fig. 19),

$$T_{m_{i+1},m_i,m_i',m_{i+1}'} = (e^{-\delta\tau H_i})_{m_i m_{i+1},m_i' m_{i+1}'},$$
(30)

where m_i, m'_i label the quantum states of S_i . Therefore the imaginary-time path-integral Tr $e^{-\tau H}$ can be expressed as a



FIG. 20. (Color online) A T_a - T_b tensor network of size $L \times M$ is reduced to a T'_a - T'_b tensor network of size $L \times M/2$. From (a) \rightarrow (b), the local deformation Fig. 4(a) is used. From (b) \rightarrow (c), the local deformation Fig. 9 is used. From (c) \rightarrow (d), a local deformation similar to that in Fig. 4(b) is used.

tensor trace over a tensor network of T [see Fig. 20(a)],

Tr
$$e^{-\tau H} = t Tr \otimes_i T$$
.

Now we can use the TEFR approach to evaluate the tensor trace and to obtain the fixed-point tensors. This allows us to find the zero-temperature phase diagram and the quantum phase transitions of the spin-1 model [Eq. (29)].

However, for small $\delta\tau$, the tensor network is very anisotropic with very different correlation lengths (measured by the distance of the tensor network) in the time and the space directions. This causes large truncation errors in the TEFR steps, (a) \rightarrow (b) and (c) \rightarrow (d) in Fig. 7 [or Fig. 4(a)]. We need to perform anisotropic coarse graining to make the system more or less isotropic before performing the TEFR coarse graining. One such anisotropic coarse graining in time direction is described in Fig. 20. Each coarse graining step of Fig. 20 triples the effective $\delta\tau$. After making the effective $\delta\tau$ to be of order 1, we switch to the isotropic coarse graining in Fig. 7.

We performed 11 steps of TEFR iteration in Fig. 2 (which is implemented as 22 steps of TEFR iteration in Fig. 7) with $D_{\text{cut}}=37$ for various values of U and B. After the TEFR iteration, an initial tensor T(U,B) flows to a fixed-point tensor. As we vary U and B, the different initial tensors may flow to different fixed-point tensors which represent different phases. A more detailed description of the renormalization flow of the tensors can be found in Appendix, Sec. 4.

To quantitatively plot the fixed-point tensors, we can use X_1 , X_2 [see Eq. (17)] and λ_{sum} . Here λ_{sum} is given by $\lambda_{sum} = \sum_s |\lambda_s/\lambda_1|$ and λ_i are the singular values of the matrix M^{lurd} obtained from the fixed-point tensor [see Eqs. (A15) and (3)]. In Fig. 21, we plot the X_1 , X_2 , and λ_{sum} of the resulting fixed-point tensors for different values of U and B.



FIG. 21. (Color online) The color-map plot of $(r,g,b) = (X_1/3, X_2/3, 1/2\lambda_{sum})$ shows the phase diagram of our spin-1 system [Eq. (29)] which contains four phases, a trivial $S^z=0$ phase "TRI," two Z_2 symmetry breaking phases Z_2^x and Z_2^y , and a Haldane phase "Haldane," The transition between those phases are all continuous phase transitions.

We see that our spin-1 system has four different phases: TRI, Z_2^z , Z_2^y , and Haldane. The transition between the Haldane phase and the Z_2 phase happens at $B_c=0.405$ along the U=0 line, which agrees with the Haldane gap Δ =0.4097= B_c obtained in Ref. 68. The transition between the Haldane phase and the TRI phase happens at $U_c=1.0$ along the B=0 line, which agrees with the result $U_c=0.99$ obtained in Ref. 69.

The fixed-point tensor in the TRI phase is the trivial dimension-one tensor T^{TRI} with $(X_1, X_2, \lambda_{\text{sum}}) = (1, 1, 1)$. So the TRI phase has no symmetry breaking. The fixed-point tensor in the Z_2^v and Z_2^z phase, up to a field-redefinition transformation, is the dimension-two tensor T^{Z_2} which is a direct sum of two dimension-one tensors T^{TRI} with $(X_1, X_2, \lambda_{\text{sum}}) = (2, 2, 2)$. So the Z_2^v and the Z_2^z phases are Z_2 symmetry breaking phases. The fixed-point tensor in the Haldane phase is a dimension-four tensor. However, many different dimension-four tensors can appear as the fixed-point tensors. All those dimension-four fixed-point tensors have $(X_1, X_2, \lambda_{\text{sum}}) = (1, 1, 4)$. In fact they are all equivalent to the following CDL tensor,

$$T^{\text{Haldane}} = T(\sigma^{y}, \sigma^{y}, \sigma^{y}, \sigma^{y}), \qquad (31)$$

through the field-redefinition transformation [Eq. (19)], where $T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$ is given in Eq. (6). In fact, one can show that the following four CDL tensors, $T(\sigma^x, \sigma^x, \sigma^x, \sigma^x)$, $T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$, $T(\sigma^z, \sigma^z, \sigma^z, \sigma^z)$, and $T(\sigma^0, \sigma^0, \sigma^0, \sigma^0)$, are all equivalent through the field-redefinition transformation. (Here σ^0 is the 2×2 identity matrix.) So they can all be viewed as the fixed-point tensors in the Haldane phase. Since the fixed-point tensor in the Haldane phase is a CDL tensor, the Haldane phase, like the TRI phase, also does not break any symmetry.

From the fixed-point tensors, we can also obtain the low energy spectrum of our spin-1 system (see Appendix, Sec. 2). Using the fixed-point tensors of the four phases, we find that all four phases have a finite energy gap. We also find that the ground states of the TRI and Haldane phases are nondegenerate, while the ground states of the Z_2^{ν} and Z_2^{τ} phases have a twofold degeneracy. This is consistent with our result that the



FIG. 22. (Color online) The color-map plot of the central charge as a function of U, B. The color is given by $(r, g, b) = (c_8, c_9, c_{10})$, where c_i is the central charge obtained at the *i*th iteration of Fig. 2. The lines with nonzero central charge are the critical lines with gapless excitations. The model [Eq. (29)] has a U(1) spin rotation symmetry along the U=0 line. The U=0 line between the two Z_2 phases is a gapless state.

TRI phase and the Haldane phase do not break any symmetry, while the Z_2^{y} and Z_2^{z} phases break a Z_2 symmetry.

The central charge calculated from the resulting fixedpoint tensors is plotted in Fig. 22 [see Eq. (A18)]. The lines of nonzero central charge in Fig. 22 mark the regions of diverging correlation length, which correspond to line of continuous phase transition. The energy gap closes along those lines. Comparing Fig. 22 with Fig. 21, we see that all the phase transitions in the phase diagram (Fig. 21) correspond to those critical lines. Those phase transitions are all continuous phase transitions. The transitions between Haldane and Z_2^z , Haldane and Z_2^y , TRI and Z_2^y are described by central charge c=1/2 critical point. The transitions between Z_2^y and Z_2^z , Haldane and TRI, are described by central charge c=1 critical point.

The continuous phase transitions between the TRI and Z_2 phases and between the Haldane and Z_2 phases are Z_2 symmetry breaking transitions. The phase transition between the Z_2^y and the Z_2^z phases is of the same type as that discussed in Sec. V D along the h=0 line. Along the U=0 line, our spin-1 model has a U(1) spin rotation symmetry around the S^x axis. The gapless phase at $B > B_c$ and U=0 is close to a U(1) symmetry breaking phase where spins are mainly in the y-z plane with a finite and uniform S_x component. When $U \neq 0$, the U(1) spin rotation symmetry is broken. When U > 0, the spin-1 model is in the Z_2^y phase with $S_i^y \propto (-)^i$. When U < 0, the spin-1 model is in the Z_2^z phase with $S_i^z \propto (-)^i$.

Both the TRI phase and the Haldane phase have the same symmetry, but yet they are separated by phase transitions. This suggests that the TRI phase and the Haldane phase are distinct phases. But we cannot use symmetry breaking to distinguish the two phases. So let us discuss Haldane phase in more detail.

F. Haldane phase—a symmetry-protected topological phase

The first question is that whether the TRI phase and the Haldane phase are really different? (For a detailed discussion on the definitions of phases and phase transitions, see Appendix, Sec. 3.) Can we find a way to deform the two phases into each without any phase transition? At first sight, we note that the fixed-point tensors for the Haldane phase are CDL tensors. According to the discussion in Sec. III, this seems to suggest that the Haldane phase is the trivial TRI phase described by T^{TRI} .

In fact, the answer in the above question depends on the symmetry of the Hamiltonian. If we allow to deform the spin-1 Hamiltonian arbitrarily, then the TRI phase and the Haldane phase will belong to the same phase since the two phases can be deformed into each other without phase transition. If we require the Hamiltonian to have certain symmetries, then the TRI phase and the Haldane phase are different phases (in the sense defined in Appendix, Sec. 3).

To demonstrate the above result more concretely, let us consider a spin-1 system with time-reversal, parity (spatial reflection), and translation symmetries. The time-reversal symmetry requires the Hamiltonian H and the corresponding tensor T [see Eq. (30)] to be real. Since the real Hamiltonian must be symmetric, time-reversal symmetric tensor must also satisfy

$$T_{ruld} = T_{dlur}.$$
 (32)

[Note that the time-reversal transformation exchanges $(m_i, m_{i+1}) \leftrightarrow (m'_i, m'_{i+1})$ in Eq. (30).] The parity symmetry requires the tensor *T* to satisfy

$$T_{ruld} = T_{urdl}.$$
 (33)

[Note that the parity exchanges $(m_i, m'_i) \leftrightarrow (m_{i+1}, m'_{i+1})$ in Eq. (30).] The tensor network considered here is always uniform which ensures the translation symmetry. Our spin-1 Hamiltonian (29) has time-reversal, parity, and translation symmetries.

Using our numerical TEFR calculation, we have checked that the Haldane phase is stable against any perturbations that respect the time-reversal, parity, and translation symmetries. In other words, we start with a tensor T with respect to those symmetries and flows to the fixed-point tensor $T^{Haldane}$. We then add an arbitrary perturbation $T \rightarrow T + \delta T$ with respect to the time-reversal, parity, and translation symmetries. We find that $T + \delta T$ also flows to the fixed-point tensor T^{Haldane} as long as δT is not too large. This result suggests that the time-reversal, parity, and translation symmetries protect the stability of the Haldane phase. This implies that in the presence of those symmetries, the TRI phase and the Haldane phase are always different. We cannot smoothly deform the Haldane phase into the TRI phase through the Hamiltonian that has the time-reversal, parity, and translation symmetries.

If we add a perturbation δT that has time-reversal symmetry but not parity symmetry, our numerical result shows that the perturbed tensor $T + \delta T$ will fail to flow to T^{Haldane} . Instead, it will flow to the trivial fixed-point tensor T^{TRI} . This implies that, without the parity symmetry, the Haldane phase is unstable and is the same as the trivial phase TRI, which agrees the result obtained previously in Ref. 59.

To summarize, for Hamiltonian with time-reversal and parity symmetries, the Haldane phase and the TRI phase are different phases, though both phases do not break any symmetries. So the difference between the Haldane phase and the TRI phase cannot be describe by Landau symmetry breaking theory. In this sense the Haldane phase behaves like the topologically ordered phases.

On the other hand, for Hamiltonian without parity symmetry, the Haldane phase and the TRI phase are the same phase. We see that the Haldane phase exists as a distinct phase only when the Hamiltonian have time-reversal, parity, and translation symmetries. This behavior is very different from the topological phases which are stable against any perturbations, including those that break the parity symmetry. Therefore, we should refer Haldane phase as a symmetryprotected topological phase. We like to stress that the Haldane phase is only one of many possible symmetryprotected topological phases. Many other symmetryprotected topological phases are studied using projected symmetry group in Refs. 17 and 29. For example, without symmetry, there is only one kind of Z_2 spin liquid. In the presence of spin rotation symmetry and the symmetries of square lattice, there are hundreds different Z_2 spin liquids.²⁹ All those Z_2 spin liquids can be viewed as symmetryprotected topological phases.

Since the time-reversal, parity, and translation symmetries play such an important role in the very existence of the Haldane phase, it is not proper to characterize the Haldane phase just using the fixed-point tensor T^{Haldane} . We should use a pair $(G_{TPT}, T^{\text{Haldane}})$ to characterize the Haldane phase. Here G_{TPT} is the symmetry group generated the timereversal, parity, and translation transformations. Using the more complete notation, we can say that $(G_{TPT}, T^{\text{Haldane}})$ and $(G_{TPT}, T^{\text{TRI}})$ correspond to different phases, while $(G_{TT}, T^{\text{Haldane}})$ and (G_{TT}, T^{TRI}) correspond to the same phase. Here G_{TT} is the symmetry group generated only by the timereversal and translation transformations. In terms of the new notation, the four phases, TRI, Haldane, and the two Z_2 , are labeled by $(G_{TPT}, T^{\text{TRI}})$, $(G_{TPT}, T^{\text{Haldane}})$, $(G_{TPT}, T^{\mathbb{Z}_2^{\mathcal{Y}}})$. and $(G_{TPT}, T^{Z_2^{\tilde{z}}})$. For a more detailed discussion on how the symmetry transformations act on the fixed-point tensor, see Appendix, Sec. 5

In this paper, we propose to use the pair $(G_{TPT}, T^{\text{Haldane}})$ to characterize the Haldane phase-a symmetry-protected topological phase. Such a notation allows us to show that the Haldane phase is stable against any perturbations that respect time-reversal, parity, and translation symmetry. We feel that the new characterization captures the essence of Haldane phase. People have proposed several other ways to characterize the Haldane phase. One of them is to use the emergent spin-1/2 boundary spins at the two ends of a segment of the isotropic spin-1 chain in the Haldane phase. However, such a characterization fails to capture the essence of Haldane phase. The Haldane phase still exists in the presence of a uniform magnetic field, $B\sum_i S_i^z$, while the fourfold degeneracy from the emergent spin-1/2 boundary spins is lifted by such a uniform magnetic field. In the presence of a uniform magnetic field, we can no longer use the degeneracy of the boundary spin-1/2 spins to detect the Haldane phase.

People also proposed to use the string order parameter



FIG. 23. (Color online) The color-map plot of $(r,g,b) = (X_1, X_2, 1/\lambda_{sum})$ shows the phase diagram of the spin-1 system [Eq. (34)] which contains four phases, a trivial phase TRI, two Z_2 symmetry breaking phases Z_2^y and Z_2^z , and a Haldane phase Haldane. The transition between those phases are all continuous phase transitions.

$$\lim_{|i-j|\to\infty} \left\langle S_i^z \exp\left(i\pi\sum_{k=i+1}^{j-1} S_k^z\right) S_j^z\right\rangle \neq 0$$

to characterize the Haldane phase. But the string order parameter vanishes in the presence of perturbations that contain odd number of spin operators, such as

$$\begin{split} \delta H &= \sum_{i} A_{1} \left[(S_{i}^{z})^{2} S_{i+1}^{z} + (S_{i+1}^{z})^{2} S_{i}^{z} \right] \\ &+ \sum_{i} A_{2} \left[(S_{i}^{x})^{2} S_{i+1}^{x} + (S_{i+1}^{x})^{2} S_{i}^{x} \right] \end{split}$$

This is because the total number of S^z =+1 and S^z =-1 sites mod 2 is not conserved in the presence of the above perturbation. However, the vanishing string order parameter does not imply the instability of Haldane phase. The above perturbation respects the time-reversal and parity symmetries. The Haldane phase is stable against such a perturbation. Therefore, the string order parameter also fails to capture the essence of Haldane phase.



FIG. 24. (Color online) The central charge as a function of U,B. The lines with nonzero central charge are the critical lines with gapless excitations. The model [Eq. (34)] has a U(1) spin rotation symmetry along the U=0 line. The U=0 line between the two Z_2 phases is a gapless state.



FIG. 25. (Color online) The color-map plot of $(r,g,b) = (X_1/3, X_2/3, 1/2\lambda_{sum})$ shows the phase diagram of the spin-1 system [Eq. (35)] which contains four phases, a trivial phase TRI, two Z_2 symmetry breaking phases Z_2^x and Z_2^y , and a Haldane phase Haldane. The transitions between those phases are all continuous phase transitions.

To illustrate the stability of Haldane phase, we also calculated the phase diagrams of following two spin-1 models. The first model is given by

$$H = \sum_{i} \left[S_{i} \cdot S_{i+1} + U(S_{i}^{z})^{2} + B(S_{i}^{x} + S_{i}^{z}) \right].$$
(34)

The phase diagram (the color-map plot of X_1 , X_2 , and λ_{sum}) and the central charge are plotted in Figs. 23 and 24. The second model is given by

$$H = \sum_{i} \left[S_{i} \cdot S_{i+1} + U(S_{i}^{z})^{2} \right]$$

+ $\frac{B}{2} \sum_{i} \left[S_{i}^{x} (S_{i+1}^{z})^{2} + S_{i+1}^{x} (S_{i}^{z})^{2} + 2S_{i}^{z} \right].$ (35)

The phase diagram and the central charge are plotted in Figs. 25 and 26.

We see that in both models, the Haldane phase and the TRI phases are separated by phase transitions, indicating the stability of the Haldane phase. (Note that the stability of the Haldane phase is defined by the impossibility to deform the



FIG. 26. (Color online) The color-map plot of the central charge as a function of U,B. The color is given by $(r,g,b)=(c_8,c_9,c_{10})$, where c_i is the central charge obtained at the *i*th iteration of Fig. 2. The lines with nonzero central charge are the critical lines with gapless excitations.



FIG. 27. (Color online) The fixed-point wave function of the Haldane phase can be obtained from the imaginary-time evolution of the fixed-point tensor network of the Haldane phase $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$, which is formed by the corner matrix σ^y (i.e., the boundary of the tensor network gives raise to the fixed-point wave function). The dimension 2 of the corner matrix implies that each index of the corner matrix is associated with a pseudo-spin-1/2. Thus each coarse-grained site has two spin-1/2 pseudospins. The corner matrix σ^y ties the two pseudospins on the neighboring sites into a pseudospin singlets. Thus the fixed-point wave function of the Haldane phase can be viewed as a dimer state formed by nearest-neighbor pseudospins.

Haldane phase to the TRI phase without a phase transition within a symmetry class of Hamiltonian.) Both the models have time-reversal and parity symmetries. In particular, the second model has no other symmetries beside translation, time-reversal and parity symmetries. This suggests that the stability of Haldane phase requires only translation, timereversal, and parity symmetries for the Hamiltonian.

Before ending this section, we like to point out that the form of the fixed-point tensor of the Haldane phase $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$ implies the following fixed-point wave function for the Haldane phase. Since the fixed-point tensor $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$ has a dimension 4, the effective degrees of freedom in the fixed-point wave function is four states per site (i.e., per coarse-grained site). The four states can be viewed as two spin-1/2 pseudo spins. The fixed-point wave function is then formed by the pseudospin singlet dimers between nearest-neighbor bonds (see Fig. 27).

VI. SUMMARY

In this paper, we introduce a TEFR approach for statistical and quantum systems. We can use the TEFR approach to calculate phase diagram with both symmetry breaking and topological phases. We can also use the TEFR approach to calculate critical properties of continuous transitions between symmetry breaking and/or topological phases. The TEFR approach suggests a very general characterization of phases based on the symmetry group G_{sym} and the fixed-point tensor T_{inv} . The ($G_{\text{sym}}, T_{\text{inv}}$) characterization can describe both symmetry breaking and topological phases. Using such a characterization and through the stability of the fixed-point tensor, we show that the Haldane phase for spin-1 chain is stable against any perturbations with respect to time-reversal, parity, and translation symmetry. This suggests that the Haldane phase is a symmetry-protected topological phase.



FIG. 28. (Color online) The paradigm of topological order.

The TEFR approach introduced here can be applied to any strongly correlated systems in any dimensions. It has a potential to be used as a foundation to study topological phases and symmetry-protected topological phases in any dimensions. In this paper, we only demonstrate some basic applications of the TEFR approach through some simple examples. We hope those results will constitute the foundation for using the TEFR approach to study more complicated strongly correlated systems in 1+2 and in 1+3 dimensions.

To gain a better understanding of the tensor-network renormalization approach, let us compare the tensor-network renormalization approach with the block-spin renormalization approach.²⁴ Let us consider a spin system whose partition function can be represented by a tensor network of a tensor T(g), where g is a coupling constant of the spin system. After performing a coarse graining of the lattice, the system is described by a coarse-grained tensor network of tensor T'. The transformation $T(g) \rightarrow T'$ represents tensornetwork renormalization flow. In the block-spin renormalization approach, we insist on the coarse-grained systems that have the same form as described by a new coupling constant g'. So we try to find a g' such that T(g') best approximates T'. This way, we obtain a block-spin renormalization flow $g \rightarrow g'$. We see that tensor-network renormalization flow is a transformation in the generic tensor space, while the blockspin renormalization flow is a transformation in the subspace T(g) parametrized by the coupling constant g. We believe that it is this generic flow in the whole tensor space that allows the tensor-network renormalization approach to describe topological phases.

In this paper, we stress that the fixed-point tensor of the tensor-network renormalization flow can be used as a quantitative description of topological order. This is just one aspect of topological order. In fact, theory of topological order has a very rich structure and covers a wide area (see Fig. 28). Turaev-Viro invariant⁷⁰ on three-dimensional manifold can be viewed as a tensor trace of the fixed-point tensors for string-net models.¹² Therefore, the tensor-network renormalization approach should apply to all time-reversal and parity symmetric topological orders in 1+2 dimensions which are described by the string-net condensations.^{11,12} This type of topological order include emergent gauge theory, Cherntheory, Simons and quantum gravity in 1 + 2dimensions.^{12,71,72} The mathematical foundation that classifies this type of fixed-point tensors is the tensor category theory. It is natural to expect that the topological order in 1 +3 dimensions and the associated emergent gauge bosons, emergent fermions, and emergent gravitons^{34,73} can also be studied through the tensor-network renormalization approach. The tensor-network renormalization approach also provides a way to calculate lattice gauge theories. We see that physics is well interconnected. It appears that the tensor-network renormalization is a glue that connects all different parts of physics.

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APPENDIX

1. Detailed discussion of the TEFR algorithm

The key step in the TEFR calculation is the entanglement-filtering operation from Fig. 7(b) to Fig. 7(c). In this section we will discuss several ways to implement such an entanglement-filtering calculation.

a. Linear algorithm

To find those new S'_i described in Fig. 8, we may minimize the following cost function,

$$C_{L} = \sum_{\alpha\beta\gamma\lambda} (\Delta T_{\alpha\beta\gamma\lambda})^{2},$$

$$\Delta T_{\alpha\beta\gamma\lambda} = \sum_{ijkl} S_{1li\alpha} S_{2ij\beta} S_{3jk\gamma} S_{4kl\delta}$$

$$- \sum_{i'j'k'l'} S'_{1l'i'\alpha} S'_{2i'j'\beta} S'_{3j'k'\gamma} S'_{4k'l'\delta}, \quad (A1)$$

where $\alpha, \beta, \delta, \gamma, i, j, k, l=1, ...D$ and i', j', k', l'=1, ...D'(see Fig. 30). This cost function can be minimized by solving a sets of least square problems iteratively. For example, we may first fixed S'_2, S'_3, S'_4 and solve a least square problem for S'_1 , then we fixed S'_1, S'_3, S'_4 to solve a least square problem for S'_2 , etc. In practice, to reduce the risk of being trapped by local minima, we may start with D'=D-1 and gradually decrease D'.

b. Nonlinear algorithm

Although the linear algorithm is an efficient way to do the entanglement-filtering calculation in Fig. 8, however, this algorithm could not avoid being trapped by local minima in generic case. To solve this problem, here we introduce a nonlinear algorithm which can avoid the trapping by local minima.

In the first implementation of TEFR discussed above, we choose S'_i to minimize the dimension of the one of the indices of the rank-three tensors S'_i . The second implementation of TEFR is still described by Fig. 7. But now we choose S'_i



FIG. 29. A rank-four tensor expressed as tensor network.

to minimize the entanglement entropy on the diagonal links in Fig. 7(d).

Let us define the entanglement entropy on a link. We may view the tensor network in Fig. 29 as matrix $M_{i,j}$ where i = (a,b) and j = (c,d). We then perform an SVD of M which produces the singular values λ_i that are ordered as $\lambda_1 \ge \lambda_2$ $\ge ... \ge 0$. We call λ_i the singular values on the horizontal link in Fig. 29. The entanglement entropy on the horizontal link in Fig. 29 is defined as $S = -\sum_i \tilde{\lambda}_i \ln(\tilde{\lambda}_i)$, where $\tilde{\lambda}_i$ $= \lambda_i / \sum \lambda_i$. In the nonlinear algorithm, we choose S'_1, S'_2, S'_3 , and S'_4 to minimize the following cost function:

$$C_{NL} = \lambda S_{\text{sum}} + \sum_{\alpha\beta\gamma\lambda} (\Delta T_{\alpha\beta\gamma\lambda})^{2},$$

$$\Delta T_{\alpha\beta\gamma\lambda} = \sum_{ijkl} S_{1li\alpha} S_{2ij\beta} S_{3jk\gamma} S_{4kl\delta}$$

$$- \sum_{i'j'k'l'} S'_{1l'i'\alpha} S'_{2i'j'\beta} S'_{3j'k'\gamma} S'_{4k'l'\delta}, \quad (A2)$$

where S_{sum} is the sum of entanglement entropy on the four diagonal links in Fig. 7(d). Here λ should be as small as possible.

2

We note that when we reduce the entanglement entropy on the diagonal links in Fig. 7(d), the singular values on those diagonal links decay faster. This allows us to drop indices associated with the very small singular values and hence reduce the dimension on those diagonal links. When we use Fig. 4(b) to transform Fig. 7(d) to Fig. 7(e), the resulting tensor T' will have a small dimension. This approach can again reduce a CDL tensor T to a trivial dimension-one tensor.

The nonlinear algorithm is a very general approach, but the calculation is very expensive. In the paper, we will not use this algorithm.

c. Scaling-SVD algorithm for CDL tensors

In this section, we introduce the third algorithm, which is very simple and efficient. We will call such algorithm scaling-SVD algorithm, which can be used to find S'_1 , S'_2 , S'_3 , and S'_4 in Fig. 8 with smaller dimensions.

In the scaling-SVD algorithm, we introduce four diagonal weighting matrices,

$$\begin{split} & w_{ii'}^{12} = w_i^{12} \delta_{ii'}, \quad w_{jj'}^{23} = w_j^{23} \delta_{jj'}, \\ & w_{kk'}^{34} = w_k^{34} \delta_{kk'}, \quad w_{ll'}^{41} = w_l^{41} \delta_{ll'}, \end{split}$$

on those links i, j, k, l in Fig. 30 and initialize them as 1. We first perform the scaling-SVD calculation on the pair S_1 and S_2 by introducing a matrix



FIG. 30. The entanglement-filtering procedure.

$$M_{\alpha l;j\beta}^{12} = \sum_{i} \sqrt{w_l^{41}} S_{1li\alpha} S_{2ij\beta} \sqrt{w_j^{23}}$$

and do the SVD decomposition $M^{12} = U\Lambda V^{\dagger}$,

$$M_{\alpha l;j\beta}^{12} = \sum_{i} U_{\alpha l;i} \Lambda_{i}^{12} V_{j\beta;i}.$$

We then update S_1, S_2 as

$$S'_{1li\alpha} = \sqrt{\Lambda_i^{12}} U_{\alpha l;i} / \sqrt{w_l^{41}}, \quad S'_{2ij\beta} = \sqrt{\Lambda_i^{12}} V_{j\beta;i} / \sqrt{w_j^{23}}$$

We can update S_2, S_3 in the same way, then we can further update S_3, S_4 and S_4, S_1 . After one loop update for those S_i 's, we update all the weighting factors w^{12} , etc., as $w^{12} = \Lambda^{12}$, etc., on corresponding links.

After several loop iterations, we truncate the smallest singular values of Λ_i^{12} , Λ_j^{23} , Λ_k^{34} , Λ_l^{41} and obtain new rank-three tensors S'_1 , S'_2 , S'_3 , S'_4 with a smaller dimension D' < D on the inner links. If S_1 , S_2 , S_3 , and S_4 are CDL tensor, we find

$$S'_{1li\alpha}S'_{2ij\beta}S'_{3jk\gamma}S'_{4kl\delta} \propto S_{1li\alpha}S_{2ij\beta}S_{3jk\gamma}S_{4kl\delta}$$

after truncation. By proper normalizing those S'_i , we can find the solution of

$$\sum_{ijkl} S_{1li\alpha} S_{2ij\beta} S_{3jk\gamma} S_{4kl\delta} - \sum_{ijkl} S'_{1li\alpha} S'_{2ij\beta} S'_{3jk\gamma} S'_{4kl\delta} \approx 0.$$
(A3)

If S_1 , S_2 , S_3 , and S_4 are not CDL tensor, the truncated S'_1 , S'_2 , S'_3 , and S'_4 cannot satisfy Eq. (A3) even after the rescaling of S'_1 , S'_2 , S'_3 , and S'_4 . In this case, we reject the truncated S'_1 , S'_2 , S'_3 , and S'_4 and the scaling-SVD fails to produce simplified tensors. To summarize, the scaling-SVD algorithm can simplify CDL tensors but not other tensors.

To understand why the above scaling-SVD algorithm can work for CDL tensors, we can consider the simplest CDL tensor $T_{ruld}(M^1, M^2, M^3, M^4)$ with corner matrices,

$$M^{1} = M^{2} = M^{3} = M^{4} = \begin{pmatrix} 1 & 0 \\ 0 & x \end{pmatrix},$$

where 0 < x < 1.

After one loop iteration, the 4×4 diagonal weighting matrix $w = w_{in} \otimes w_{out}$ on each link is updated as

$$w_{in} = \begin{pmatrix} 1 & 0 \\ 0 & x \end{pmatrix}, \quad w_{out} = \begin{pmatrix} 1 & 0 \\ 0 & x \end{pmatrix},$$

where *in* represents the inner line and *out* represent the outer line of the square as shown in Fig. 31. However, after n loop iteration, we find



FIG. 31. For the simplest CDL tensor, the outer line singular values are not scaled after one loop iteration, however, the inner line singular values are scaled to a power. This nontrivial scaling behavior is the key why the scaling-SVD algorithm can filter out those local entanglements for a CDL tensor.

$$w_{in} = \begin{pmatrix} 1 & 0 \\ 0 & x^n \end{pmatrix}, \quad w_{out} = \begin{pmatrix} 1 & 0 \\ 0 & x \end{pmatrix}.$$

The striking thing is that the weighting matrices on the outer lines are never scaled but the weighting matrices (as well as the singular values on the corresponding link) on the inner lines are scaled to its *n* power. After enough times of loop iterations, x^n can be smaller than the machine error. The inner line can be removed and we successfully filter out those local entanglements in the square. In this paper, we do all our TEFR calculations using this scaling-SVD algorithm.

It is also clear that if the corner matrices in a CDL tensor have degenerate eigenvalues, the scaling-SVD cannot simplify such a CDL tensor. This is a desired result since the degeneracy in the singular values may indicate the presence of topological order (see the discussion at the end of Appendix, Sec. 4).

2. Partition function and critical properties from the tensornetwork renormalization flow of the tensors

In this section, we will give a more detailed discussion about the renormalization flow of the tensor. The more detailed discussion allows us to see the relation between the fixed-point tensor and the energy spectrum for a quantum system. It also allows us to see the relation between the fixed-point tensor at the critical point and the scaling dimensions and central charge in the critical point. To be concrete, we will consider the renormalization transformation described by Fig. 2, which represents a single iteration step. Such an iteration step roughly corresponds to two iteration steps described by Fig. 4.

We start with a tensor network formed by $N_0=N$ tensors $T^{(0)}=T$. After *i* iterations, the tensor network is deformed into a tensor network of $N_i T^{(i)}$ tensors. We formally denote the tensor-network renormalization flow as

$$T^{(i)} \rightarrow T^{(i+1)}$$
.

Those tensors satisfy the following defining property:

$$\mathbf{t}\mathrm{Tr}(\otimes T^{(i)})^{N_i} = \mathbf{t}\mathrm{Tr}(\otimes T^{(i+1)})^{N_{i+1}}.$$

The scales of the tensors $T^{(i)}$ changes under each iteration. It is more convenient to describe the tensor-network renormalization flow in terms of the normalized tensor. The normalized tensor is defined as

$$\mathcal{T}^{(i)} = T^{(i)} / \Gamma_i, \tag{A4}$$

which satisfies



FIG. 32. A square-lattice tensor network in a 2D space (or 1 + 1D space time).

$$\sum_{l,u} \mathcal{T}_{l,l,u,u}^{(i)} = 1$$

In terms of the normalized tensor, after a tensor-network renormalization transformation, a tensor network of N_i normalized $\mathcal{T}^{(i)}$ tensors is transformed into a tensor network of N_{i+1} normalized $\mathcal{T}^{(i+1)}$ tensors,

$$\mathcal{T}^{(i)} \to f_{i+1} \mathcal{T}^{(i+1)}.$$

The N_i normalized $\mathcal{T}^{(i)}$ tensors and the $N_{i+1} f_i \mathcal{T}^{(i+1)}$ tensors have the same tensor trace,

$$\operatorname{tTr}(\otimes \mathcal{T}^{(i)})^{N_i} = f_{i+1}^{N_{i+1}} \operatorname{tTr}(\otimes \mathcal{T}^{(i+1)})^{N_{i+1}}.$$
 (A5)

 f_i is a quantity that can be directly obtained from the numerical tensor-network renormalization calculation.

From Eq. (A4) we see that the partition function at *i*th iteration can be expressed as a tensor trace of the normalized tensor $\mathcal{T}^{(i)}$ up to an over all factor,

$$Z_i = \Gamma_i^{N_i} \operatorname{tTr}(\otimes \mathcal{I}^{(i)})^{N_i}.$$
 (A6)

Note that by definition, the partition function is invariant under the tensor-network renormalization transformation Z_i = Z_{i+1} . From Eq. (A5), we find that Γ_i satisfies

$$\Gamma_i^{N_i} f_{i+1}^{N_{i+1}} = \Gamma_{i+1}^{N_{i+1}}, \tag{A7}$$

which allows us to calculate Γ_i from f_i . We can keep performing the tensor-network renormalization until N_i is of order 1. This way, we can calculate the partition function $Z = Z_i$ from Eqs. (A6) and (A7).

To calculate the density of the logarithmic partition function $\ln(Z)$, we need to describe how tensor positions in the tensor network are mapped into the physical coordinates $\mathbf{r} = (x_1, x_2)$. Let X_1 be the vector in physical space connecting the two neighboring tensors in the horizontal direction and X_2 be the vector connecting the two neighboring tensors in the vertical direction (see Fig. 32). Let us denote the coordinates of X_i as

$$X_{\alpha} = \begin{pmatrix} W_{1\alpha} \\ W_{2\alpha} \end{pmatrix}, \quad \alpha = 1, 2.$$
 (A8)

The 2×2 matrix W describes the shape of the tensor network.

We note that the tensor network is changed under each iteration. So $W_{\alpha\beta}$ is also changed under each iteration. After the *i*th iteration, the shape of the resulting tensor network is

described by $W_{\alpha\beta}^{(i)}$, where $W_{\alpha\beta}^{(0)} = W_{\alpha\beta}$ is for the initial tensor network. Under the renormalization transformation described in Fig. 2, $W^{(i)}$ transforms in a simple way $W^{(i+1)} = 2W^{(i)}$. Under the renormalization transformation described in Fig. 3, $W^{(i)}$ transforms as $W^{(i+1)} = W^{(i)} \begin{pmatrix} 1 & 1 \\ -1 & 1 \end{pmatrix}$. The above recursion relations allow us to calculate $W^{(i)}$ for each step of iteration. Such information will be useful later.

The area occupied by each tensor $T^{(i)}$ after the *i*th iteration is

$$A_i = \det(W^{(i)}),$$

where $W^{(i)}$ is the 2×2 matrix formed by $W^{(i)}_{\alpha\beta}$. It is also clear that

$$A_i N_i = A_{i+1} N_{i+1} = A_{tot}$$

is the total area of the system.

Let us introduce

$$z_i = \Gamma_i^{1/A_i}.\tag{A9}$$

From Eq. (A7), we see that z_i satisfies the following recursion relation:

$$z_{i+1} = z_i f_{i+1}^{1/A_{i+1}} \tag{A10}$$

or

$$\ln(z_i) = \ln(z_0) + \sum_{j=1}^{i} \frac{\ln(f_j)}{A_j},$$
 (A11)

which will allows us to calculate the density of the logarithmic partition function,

$$A_{\text{tot}}^{-1}\ln(Z_i) = \ln(z_i) + A_{\text{tot}}^{-1}\ln(\text{tTr}(\otimes \mathcal{T}^{(i)})^{N_i}).$$

Note that, after a number of iterations, we can reduce N_i to be of order 1. In this case $\ln(z_i)$ itself is the density of the logarithmic partition function in the large A_{tot} limit.

In the large iteration limit, $\mathcal{I}^{(i)}$ and f_i become independent of *i*. This allows us to introduce scale invariant tensor,

$$T_{\rm inv}^{(i)} = \gamma^{-1} \mathcal{T}^{(i)},$$

which satisfy

$$T_{\rm inv}^{(i)} \rightarrow T_{\rm inv}^{(i+1)}$$

in large *i* limit. We note that the invariant tensor has the following defining property:

$$\mathrm{tTr}(\otimes T_{\mathrm{inv}}^{(i)})^{N_i} = \mathrm{tTr}(\otimes T_{\mathrm{inv}}^{(i+1)})^{N_{i+1}}.$$

From Eq. (A5), we see that γ must satisfy $\gamma^{N_i} = \gamma^{N_{i+1}} f_{i+1}^{N_{i+1}}$ or

$$\gamma = f_i^{1/3}, \qquad (A12)$$

where we have used

$$N_i = 4N_{i+1}.$$
 (A13)

Thus the invariant tensor is simply given by

$$T_{\rm inv}^{(i)} = f_i^{-1/3} T^{(i)}.$$
 (A14)

In the large *i* limit, f_i approaches a constant and the recursion relation [Eq. (A10)] implies that

$$\ln\{z_{i+1}f_{i+1}^{(1/3)[1/(A_{i+1})]}\} - \ln[z_if_i^{(1/3)(1/A_i)}]$$
$$= \frac{1}{3A_i}[\ln(f_{i+1}) - \ln(f_i)] = o(1/A_i).$$

This result allows us to calculate the finite-size correction to the partition function. For a tensor network of *N* tensors *T*, from the relation $Z=tTr(\otimes T)^N=\Gamma_i^{N_i} tTr(\otimes T^{(i)})^{N_i}$, we find that the partition function is given by

$$Z = z_i^{A_{\text{tot}}} \operatorname{tTr}(\otimes \mathcal{T}^{(i)})^{N_i} = (z_i f_i^{(1/3)(1/A_i)})^{A_{\text{tot}}} \operatorname{tTr}(\otimes \mathcal{T}^{(i)}_{\text{inv}})^{N_i}.$$

When A_{tot} and A_i are of the same order (i.e., when N_i is of order 1), we find that

$$\ln(Z) = A_{\text{tot}} \ln(z_{\infty}) + \ln[\text{tTr}(\otimes T_{\text{inv}}^{(i)})^{N_i}] + o(1),$$

where $z_{\infty} = \lim_{i \to \infty} z_i f_i^{1/3A_i}$. Also note that the invariant tensor $T_{\text{inv}}^{(i)}$ is independent of *i* in the large *i* limit (because $T_{\text{inv}}^{(i)}$ approaches to a fixed-point tensor). We see that $\ln(Z)$ contains a term $A_{\text{tot}} \ln(z_{\infty})$ that is proportional to the system area A_{tot} . $\ln(Z)$ also contains a A_{tot} independent term $\ln[\text{tTr}(\otimes T_{\text{inv}}^{(i)})^{N_i}]$. The remaining o(1) term approaches to zero as $A_{\text{tot}} \to \infty$.

It is interesting to note that although the $\ln[tTr(\otimes T_{inv})^{N_i}]$ term in the partition function does not depend on the total area A_{tot} of the tensor network, it can depend on the shape of the network if the system is at the critical point. We can use this shape-dependent partition function to obtain the critical properties of the critical point. In the following, we will study the shape dependence of the partition function at a critical point. We will assume that in the 1+1D space time, the velocity is equal to 1 (or the critical system is isotropic in the 2D space).

One way to obtain shape-dependent partition function (i.e., the A_{tot}^0 part of the partition function) is to introduce four matrices associated with the invariant fixed-point tensor $T_{inv}^{(i)}$,

$$(M^{ud})_{u,d} = \sum_{r} (T^{(i)}_{inv})_{rurd}, \quad (M^{lr})_{l,r} = \sum_{u} (T^{(i)}_{inv})_{rulu},$$
$$(M^{ldru})_{ld,ru} = (T^{(i)}_{inv})_{ruld}, \quad (M^{lurd})_{lu,rd} = (T^{(i)}_{inv})_{ruld}.$$
(A15)

The eigenvalues of those matrices encode the information about that shape-dependent partition function.

To understand the physical meaning of those eigenvalues, let us consider a $T_{inv}^{(i)}$ -tensor network that has a rectangular shape $N_i = L \times M$. The shape-dependent part of the partition function is given by $tTr(\otimes T_{inv}^{(i)})^{N_i}$. To describe the shape quantitatively, we note that the $T_{inv}^{(i)}$ -tensor network corresponds to a parallelogram spanned by two vectors $l = l_{\alpha} x_{\alpha}$ $= LW_{1\alpha}^{(i)} x_{\alpha}$ and $m = m_{\alpha} x_{\alpha} = MW_{2\alpha}^{(i)} x_{\alpha}$, where x_{α} , $\alpha = 1, 2$ form an orthonormal basis of the space. The shape, by definition, is described by a complex number

$$\tau^{(i)} = \frac{m_1 + \mathrm{i}m_2}{l_1 + \mathrm{i}l_2} = \frac{W_{21}^{(i)} + \mathrm{i}W_{22}^{(i)}}{W_{11}^{(i)} + \mathrm{i}W_{12}^{(i)}} \frac{M}{L} \,.$$

Many critical properties of the critical point can be calculated from the τ dependence of $tTr(\otimes T_{inv}^{(i)})^{N_i}$. In particular (according to conformal field theory)⁷⁴

$$[tTr(\otimes T_{inv}^{(i)})^{N_i}] = \sum_{n=0} e^{-2\pi [(h_n^R + h_n^L - c/12)Im \tau^{(i)} + i(h_n^R - h_n^L)Re \tau^{(i)}]},$$
(A16)

where *c* is the central charge and $\{h_n^R, h_n^L\}$ is the spectrum of right and left scaling dimensions of the critical point.

Since $T_{inv}^{(i)}$ is already a fixed-point tensor, we may evaluate Eq. (A16) by choosing L=1,

$$t \operatorname{Tr}(\otimes T_{\text{inv}}^{(i)})^{M} = \operatorname{Tr}(M^{ud})^{M}$$
$$= \sum_{n=0} e^{-2\pi M [(h_{n}^{R} + h_{n}^{L} - c/12) \operatorname{Im} \tau_{0}^{(i)} + i(h_{n}^{R} - h_{n}^{L}) \operatorname{Re} \tau_{0}^{(i)}]},$$

where

$$\tau_0^{(i)} = \frac{W_{21}^{(i)} + iW_{22}^{(i)}}{W_{11}^{(i)} + iW_{12}^{(i)}}.$$
(A17)

Using the above result and some other results obtained with similar methods, we find that

$$\begin{pmatrix} h_n^R + h_n^L - \frac{c}{12} \end{pmatrix} \operatorname{Im}(\tau_0^{(i)}) + i(h_n^R - h_n^L) \operatorname{Re}(\tau_0^{(i)}) \\ = -\frac{\ln(\lambda_n^{ud})}{2\pi} + \operatorname{mod} i, \\ \begin{pmatrix} h_n^R + h_n^L - \frac{c}{12} \end{pmatrix} \operatorname{Im}\left(\frac{-1}{\tau_0^{(i)}}\right) + i(h_n^R - h_n^L) \operatorname{Re}\left(\frac{-1}{\tau_0^{(i)}}\right) \\ = -\frac{\ln(\lambda_n^{lr})}{2\pi} + \operatorname{mod} i, \\ \begin{pmatrix} h_n^R + h_n^L - \frac{c}{12} \end{pmatrix} \operatorname{Im}(\tau_1^{(i)}) + i(h_n^R - h_n^L) \operatorname{Re}(\tau_1^{(i)}) \\ = -\frac{\ln(\lambda_n^{lurd})}{\pi} + \operatorname{mod} 2i, \\ \begin{pmatrix} h_n^R + h_n^L - \frac{c}{12} \end{pmatrix} \operatorname{Im}\left(\frac{-1}{\tau_1^{(i)}}\right) + i(h_n^R - h_n^L) \operatorname{Re}\left(\frac{-1}{\tau_1^{(i)}}\right) \\ = -\frac{\ln(\lambda_n^{lurd})}{\pi} + \operatorname{mod} 2i, \end{cases}$$

$$(A18)$$

where n=0,1,2,... and λ_n^{ud} , λ_n^{lr} , λ_n^{lurd} , and λ_n^{ldru} are the eigenvalues of M^{ud} , M^{lr} , M^{lurd} , and M^{ldru} , respectively. Also

$$\tau_1^{(i)} = \frac{W_{21}^{(i)} + iW_{22}^{(i)} - W_{11}^{(i)} - iW_{12}^{(i)}}{W_{11}^{(i)} + iW_{12}^{(i)} + W_{21}^{(i)} + iW_{22}^{(i)}} = \frac{\tau_0^{(i)} - 1}{\tau_0^{(i)} + 1}.$$
 (A19)

Since $h_0=0$, we see that the central charge *c* and the scaling dimensions h_1^L, h_1^R, \ldots can be calculated from the invariant fixed-point tensor $T_{inv}^{(i)}$.

We like to stress that the invariant fixed-point tensor T_{inv} can be calculated from Eq. (A14) even away from the critical points. This allows us to calculate the central charge c and the scaling dimensions h_1^L, h_1^R, \ldots from T_{inv} away from the critical point. For phases with short-range correlations, the resulting central charge c will be zero. However, the central charge will be nonzero when correlation length diverges (or energy gap vanishes). Thus plotting central charge as a function of temperature and other parameters is a good way to discover second-order phase transitions. Also for phases with short-range correlations, a finite number of h_n^R, h_n^L are zero (which represent the degenerate ground states) and rest of them are infinite.

We can also use the fixed-point tensors to calculate the low energy spectrum of a quantum system, such as our spin-1 system [Eq. (29)]. From the tensor-network representation of 1D quantum Hamiltonian [see Fig. 20(a)], we see that the shape of the original tensor network is described by a *W* matrix $W = \begin{pmatrix} 1 & -1 \\ \delta \tau & \delta \tau \end{pmatrix}$. After *i* steps of iteration, the *W* matrix becomes $W^{(i)} = 2^{i}W$. From $W^{(i)}$, we find that $\tau_1^{(i)}$ is purely imaginary. When $\tau_1^{(i)} \propto i$, the eigenvalues of the matrix $-\ln(M^{lurd})$ will be proportional to the energy eigenvalues of our spin-1 system, as one can see from the expression of the partition function [Eq. (A16)]. From this result and using the fixed-point tensor, we can calculate the energy spectrum of quantum system.

3. Phases and phase transitions from the fixed-point tensors

In this section, we like to discuss how to define phases and phase transition more carefully and in a very general setting. For a system that depends on some parameters, for example, β and V for our generalized loop-gas model, we can calculate the partition function $Z(\beta, V)$ as a function of those parameters. The phase and the phase transitions are determined from the partition function $Z(\beta, V)$ in the following way: the singularity in the partition function $Z(\beta, V)$ marks the position of a phase transition. (This defines the phase transition.) The regions separated by the phase transition represent different phases. In other words, if a system (β_1, V_1) and another system (β_2, V_2) can be deformed into each other without phase transition, then the two systems are in the same phase. If there is no way to deform system (β_1, V_1) to system (β_2, V_2) without encountering at least one phase transition, then the two systems are in different phases. We like to point out that to avoid the complication of taking thermodynamic limit, here we only consider systems with some translation symmetry. We also like to point out that the above definition is very general. It includes symmetry breaking phases, topological phases, as well as the phases with quantum order.29

However, the above definition is still incomplete. To define phases and phase transitions, it is also very important to specify the conditions that the Hamiltonian or Lagrangian must satisfy. It is well known that without specifying the symmetry of Hamiltonian, the symmetry breaking phases and symmetry breaking phase transitions cannot be defined. After we specify the conditions on the Hamiltonian, the deformations used in defining phases and phase transitions must satisfy those conditions.

Since the conditions on the Hamiltonian or Lagrangian are part of the definition of phases and phase transitions, it is not surprising that the meaning of phases and phase transitions varies as we vary the conditions on the systems. Let us illustrate this point using the generalized loop-gas model along the β =0 line. In fact, weather the fixed-point tensors T^{LL} for $V < V_c$ and T^{Z_2} for $V > V_v$ in the generalized loop-gas model describe two different phases or not depend on the conditions on the system.

Since we are using tensor network to describe our systems, the conditions on the systems (or the Hamiltonian and/or Lagrangian) will be conditions on the tensors. If the only condition on the tensors is that the tensors T_{rudl} must be real, then T^{LL} and T^{Z_2} describe the same phase. To show this, let us consider a family of tensors labeled by θ ,

$$T^{\theta}_{ruld} = R^{-1}_{rr'}(\theta) R^{-1}_{uu'}(\theta) T^{Z_2}_{r'u'l'd'} R_{l'l}(\theta) R_{d'd}(\theta),$$

where $R(\theta) = \begin{pmatrix} \cos \theta \sin \theta \\ \sin \theta - \cos \theta \end{pmatrix}$. We see that the tensors T^{θ} connect $T^{Z_2} = T^{\theta=0}$ and $T^{\text{LL}} = T^{\theta=\pi/4}$ as we change θ from 0 to $\pi/4$. Since $R(\theta)$ is a field-redefinition transformation, the partition function for T^{θ} does not depend on θ and is hence a smooth function of θ . Therefore, T^{LL} and T^{Z_2} describe the same phase.

On the other hand, if we specify that the tensors must satisfy the closed-loop condition [Eq. (23) or (27)] and the Z_2 symmetry condition [Eq. (26)], then T^{LL} and T^{Z_2} describe two different phases. The continuous deformation T^{θ} can no longer be used. This is because T^{θ} does not satisfy Eq. (23) if $0 < \theta < \pi/4$.

Before ending this section, we like to point out that due to field-redefinition ambiguity in the TEFR calculation, some time, a single phase may correspond to group fixed-point tensors that are related by field redefinitions. To understand this point more concretely, let us consider in more detail the $V > V_c$ phase described by the fixed-point tensor T^{Z_2} . In the decomposition of rank-four tensor T into two rank-three tensors S_2 and S_4 ,

$$T_{ruld} = \sum_{a} (S_2)_{ldk} (S_4)_{ruk},$$

we may have a ambiguity that T can also be decomposed into two different rank-three tensors S'_2 and S'_4 ,

$$T_{ruld} = \sum_{a} (S_2')_{ldk} (S_4')_{ruk},$$

where S_2 and S'_2 , S_4 and S'_4 are related by an invertible transformations,

$$(S_2)_{ldk} = A_{kk'}(S'_2)_{ldk'} \quad (S_4)_{ldk} = (S'_4)_{ldk'}(A^{-1})_{k'k'}$$

Here A represents the field-redefinition ambiguity in the TEFR renormalization approach.

If the tensors T and S must satisfy certain conditions, such as the closed-loop condition [Eq. (23)], those conditions will limit the ambiguity A. For the $V > V_c$ phase, the fixed-point tensor is T^{Z_2} . The closed-loop condition [Eq. (23)] limits the ambiguity of the decomposition of T^{Z_2} to a diagonal A matrix. The field redefinition [Eq. (19)] does not change T^{Z_2} when A and B in Eq. (19) are diagonal matrices. This way, we show that the fixed-point tensor in the $V > V_c$ phase has no field-redefinition ambiguity. Similarly, for the $V < V_c$ phase, the fixed-point tensor T^{LL} also has no fieldredefinition ambiguity if we implement the closed-loop condition [Eq. (23)]. Those results are confirmed by our numerical calculation.

4. Detailed discussion of the renormalization flow of tensors

Let us describe the TEFR flow in our spin-1 model in more detail which allows us to see how an initial timereversal and parity symmetric tensor T flows to the various fixed-point tensors. Let us first consider only the SVDTRG transformation described in Fig. 3.

It is most convenient to describe the tensor in terms of its singular values λ_s under the SVD decomposition [see Eq. (3)]. At first a few steps of iterations, the singular values λ_s are in general not degenerate. If the initial tensor *T* is in the Haldane phase, then after several iterations, the four largest singular values become more and more degenerate. If the initial tensor *T* is in the TRI phase, then after several iterations, the largest singular values remain nondegenerate. We see that the Haldane phase and the TRI phase can be distinguished from their different tensor flowing behaviors under SVDTRG transformation.

Now let us consider the tensor flow under the TEFR transformation in Fig. 7. Again at first a few steps of iterations, the singular values λ_s are in general not degenerate. Also the tensors in first a few steps are quite different from CDL tensor. As a result, the scaling-SVD operation in (b) \rightarrow (c) in Fig. 7 is a null operation which does not simplify the tensors. After several iterations, the tensors become closer and closer to CDL tensors. So after a certain step of iterations, the scaling-SVD operation becomes effective which further simplify the tensor.

If the initial tensor *T* is in the TRI phase, even after the tensor *T* becomes CDL tensor, the singular values λ_s are still not degenerate. This implies that the corner matrices M_i in the CDL tensors [see Eq. (6)] have nondegenerate eigenvalues. The entanglement-filtering operation in (b) \rightarrow (c) in Fig. 7 actually simplify the tensors by removing the smaller eigenvalues in the corner matrices M_i . So the scaling-SVD operation reduces the corner matrices M_i to a trivial one, $(M_i)_{ii'} = \delta_{1,i} \delta_{1,i'}$. This way the scaling-SVD operation reduces the initial tensor to the trivial dimension-one tensor T^{TRI} .

If the initial tensor *T* is in the Haldane phase, after the tensor *T* becomes CDL tensor after a few steps of TEFR iteration, the largest four singular values λ_s also become degenerate. This implies that the largest two eigenvalues of the corner matrices M_i in the CDL tensors are also degenerate. In this case, the scaling-SVD operation cannot reduce such a degenerate corner matrices to a trivial one. But it can still

simplify the corner matrices to a certain degree by removing other smaller eigenvalues. This way, the combinations of the SVD coarse graining in (d) \rightarrow (e) in Fig. 7 and the entanglement filtering in (b) \rightarrow (c) in Fig. 7 reduce the initial tensor *T* to a simple fixed-point tensor *T*^{Haldane}. Therefore the Haldane phase and the TRI phase can be distinguished by their different fixed-point tensors under the TEFR renormalization transformation.

We like to point out that the emergence of the fourfold degenerate singular values requires the initial tensors to satisfy the time-reversal and parity symmetries. If we start with an initial tensor T that has time-reversal symmetry but not parity symmetry, our numerical results show that the largest singular values do not become degenerate after many iterations. In this case the entanglement-filtering operation through the scaling-SVD always reduces the initial tensor T to T^{TRI} . This implies that, without the parity symmetry, the Haldane phase is the same as the trivial phase TRI.

We see that the emergence of fourfold degeneracy in the spectrum of singular values is closely related to the Haldane phase—a topologically ordered phase. In Ref. 75, an entanglement spectrum is introduced to identify topological order. One may wonder if there are any relations between the spectrum of singular values and the entanglement spectrum.

We also like to point out that, only away from critical points, the tensors flow to CDL tensor. Near critical points, tensors are always quite different from CDL tensors. As a result, the entanglement-filtering operation in $(b) \rightarrow (c)$ in Fig. 7 is a null operation which does not simplify the tensors. So near a critical point, the SVDTRG in Fig. 3 and the TEFR in Fig. 7 produce the same result.

5. Projective realization of symmetries on fixed-point tensors

We like to point out that in the original tensor network of T for the spin-1 system [Eq. (30)], the time-reversal and parity symmetries are given in Eqs. (32) and (33). However, as discussed near the end of Appendix, Sec. 3, the TEFR transformation has a field-redefinition ambiguity. As a result, the time-reversal and parity transformations contain a field-redefinition ambiguity when act on the coarse-grained tensor network. Similarly, the fixed-point tensor also contains such field-redefinition ambiguity. If we choose a basis that the fixed-point tensor in the Haldane phase is given by $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$, the time-reversal and parity transformations that act on the fixed-point tensor are given by

parity:
$$T \rightarrow T_p$$
, where $(T_p)_{ruld}$
= $\sum_{r'u'l'd'} W_{rr'}W_{uu'}W_{ll'}W_{dd'}T_{u'r'd'l'}$, (A20)

time reversal:
$$T \rightarrow T_t$$
, where $(T_t)_{ruld} = T_{dlur}$.
(A21)

Here W is a field-redefinition transformation,

$$W = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix}.$$
 (A22)

The time-reversal and parity transformations on the coarsegrained tensor network and those on the initial tensor network [see Eqs. (32) and (33)] differ by a field-redefinition transformation. The fixed-point tensor T^{Haldane} is invariant under the above projective symmetry group (PSG) transformations.²⁹ We also numerically confirm that the fixed-point tensor T^{Haldane} is stable if it is perturbed by a tensor δT that is invariant under the same PSG defined through Eqs. (A20) and (A21). We note that the PSG is used to define quantum order (a generalization of topological order) in Ref. 29. The tensor-network renormalization naturally allows the generalization from topological order to quantum order.

We like to remark that the field-redefinition transformation should be derived by keeping track of the parity and time-reversal symmetry at each TEFR flow step. As a result, the field-redefinition transformation W should satisfy Eq. (A20) for the fixed-point tensor $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$. However, not all the W that satisfy Eq. (A20) for the fixedpoint tensor $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$ correspond to the parity symmetry for the original tensor [Eq. (33)]. For example,

$$W' = \begin{pmatrix} 0 & 0 & 0 & 1 \\ 0 & -1 & 0 & 0 \\ 0 & 0 & -1 & 0 \\ 1 & 0 & 0 & 0 \end{pmatrix}$$
(A23)

also satisfies Eq. (A20), but this is not the correct fieldredefinition transformation for parity symmetry of $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$. In particular, perturbations with the wrong parity symmetry,

$$\delta T_{ruld} = \sum_{r'u'l'd'} W'_{rr'} W'_{uu'} W'_{ll'} W'_{dd'} \delta T_{u'r'd'l'}, \quad (A24)$$

and time-reversal symmetry [Eq. (A21)] can destabilize the fixed-point tensor $T^{\text{Haldane}} = T(\sigma^y, \sigma^y, \sigma^y, \sigma^y)$.

For the fixed-point tensor of Haldane phase, the field-redefinition transformations that correspond to parity and time-reversal symmetries are also base dependent. If we choose $T^{\text{Haldane}} = T(\sigma^0, \sigma^0, \sigma^0, \sigma^0)$, then the correct field-redefinition transformation corresponding to the parity symmetry should be W' not W. We can also choose other bases that T^{Haldane} is invariant in a trivial way under parity transformation but invariant up to a field redefinition under time-reversal transformation.

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