

Extracting the Luttinger Parameter from a Single Wave Function

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The low-energy physics of Tomonaga-Luttinger liquids (TLLs) is controlled by the Luttinger parameter. We demonstrate that this parameter can be extracted from a single wave function for one-component TLLs with periodic boundary condition. This method relies on the fact that TLLs are described by conformal field theory in which crosscap states can be constructed. The overlaps between the crosscap states and the ground state as well as some excited states are proved to be universal numbers that directly reveal the Luttinger parameter. In microscopic lattice models, crosscap states are formed by putting each pair of antipodal sites into a maximally entangled state. Analytical and numerical calculations are performed in a few representative models to substantiate the conformal field theory prediction. The extracted Luttinger parameters are generally quite accurate in finite-size systems with moderate lengths, so there is no need to perform data fitting and/or finite-size scaling.

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Introduction—The theory of Tomonaga-Luttinger liquid (TLL) is a great triumph of strongly correlated physics [1]. Based on previous work of Tomonaga [2], Luttinger introduced a model to study interacting fermions in one dimension [3]. It clearly exemplifies the peculiarity of reduced dimensionality because infinitesimal interaction in a one-dimensional system drives it away from the Fermi liquid. In subsequent works [4–8], the bosonization framework was established to provide a unified treatment for many strongly correlated problems in one dimension. Spin chains are usually studied using the Jordan-Wigner transformation that converts spins to fermions. Gapless phases of bosons or fermions can be understood using free boson fields that correspond to density fluctuations. Instabilities of the TLL lead to symmetry-breaking ordered phases. It is not easy to realize one-dimensional systems, but experimental investigations have been carried out in solid state and cold atom platforms [9–18]. Another context in which TLL thrives is the edge of certain two-dimensional topological states [19–22].

The low-energy physics of one-component TLLs is captured by the Hamiltonian

$$H = \frac{v}{8\pi} \int_0^L dx [(\partial_x \hat{\phi})^2 + (\partial_x \hat{\theta})^2], \quad (1)$$

which is a system of compactified bosons in the language of conformal field theory (CFT). Here L is the length of the system, v is the characteristic velocity, $\hat{\phi}$ is a compactified boson field with radius R (i.e., $\hat{\phi} \sim \hat{\phi} + 2\pi R$), and $\hat{\theta}$ is the conjugate boson field of $\hat{\phi}$ with radius $R' = 2/R$. The Luttinger parameter K can be expressed using the radius as $K = R^2/4$. Its value is 1 for free fermions and deviates from 1 when interaction is turned on. For a given TLL, many properties are determined by K so it is routinely computed in numerical studies. If the Hamiltonian has a microscopic U(1) symmetry (e.g., particle number conservation), K can be obtained by comparing the ground-state energy of multiple systems with different U(1) charges [23]. This is feasible because K can be interpreted as the stiffness of a superfluid. The powerlaw decay of certain correlation functions can be used to deduce K when their bosonized expressions are known [1]. It has been found that K also appears in the U(1) charge fluctuations of a subsystem [24,25], the U(1) symmetry-resolved entanglement spectrum [26,27], different types of entanglement entropy [28–32], and finite entanglement scaling analysis [33,34].

In this Letter, we propose that the Luttinger parameter in microscopic models can be extracted from the eigenstate

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crosscap overlap (ESCO). It has three favorable features: (i) a single wave function instead of multiple states is used; (ii) data fitting or finite-size scaling is not needed; (iii) microscopic $U(1)$ symmetry is not required. To the best of our knowledge, no other method enjoys all three advantages simultaneously. Our method is motivated by classical results and recent progresses in CFT and integrable models. When a CFT is defined on the Klein bottle, important information can be inferred from the thermal entropy [35–47]. The Klein bottle can be viewed as a special cylinder with crosscaps at its ends. A simple circle with all pairs of antipodal points identified becomes a crosscap. It is possible to define a variety of crosscap states in the framework of boundary CFT [48]. For a rational CFT, the crosscap states have universal overlaps (“crosscap coefficients”) with its ground state [49–51]. In some integrable models, the crosscap states have been identified as “integrable boundary states” [52–56]. The overlaps between all eigenstates of these integrable models and the crosscap states can be expressed as determinants, but their physical significance is not transparent. We aim to establish a general framework for studying crosscap states in CFT (without and with perturbations) such that physical properties can be deduced from their overlaps with eigenstates.

General construction of crosscap states—While the crosscap is easy to visualize, the problem of what kind of states qualify as conformal crosscap states in CFTs has not been completely settled [49–51]. We take the viewpoint that a conformal crosscap state should satisfy two sets of constraints. First, the field operators on antipodal spatial coordinates are identified with each other in some way. If a CFT has Lagrangian description, this can be made precise using the path integral formalism [57–59]. These requisites are physical manifestations of the geometric structure of crosscap and referred to as “sewing conditions.” Second, the partition functions on a cylinder with crosscap boundaries (tree channel) and on a Klein bottle (loop channel) should be equal. This “loop-tree channel correspondence” allows one to determine certain coefficients in crosscap states. As we turn to lattice models, the problem becomes even more complicated because the lattice counterpart of a field operator is generally difficult to find. It is thus unlikely that a conformal crosscap state can be converted rigorously to a lattice one. Based on the intuitive picture of identifying antipodal spatial points, recent works have studied lattice crosscap states in which antipodal sites form Bell states [52–56], but their relation to the conformal crosscap states has not been elucidated.

It is particularly instructive to study a specific lattice realization of the Hamiltonian in Eq. (1), namely, the modified Villain model [60] as illustrated in Fig. 1(a):

$$H_{\text{mV}} = \sum_{j=1}^N [(2\pi R \hat{p}_j)^2 + R'^2 (\hat{X}_{j+1} - \hat{X}_j + 2\pi \hat{n}_{j,j+1})^2]. \quad (2)$$

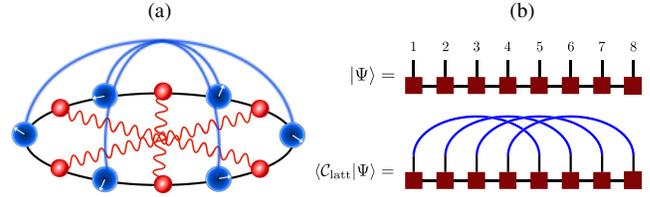


FIG. 1. (a) Schematics of the modified Villain model and its crosscap states $|C_{\pm}\rangle$. (b) Graphic representations of an MPS with eight sites and its crosscap overlap.

There are N sites in the chain (labeled by j), the lattice spacing is a , and periodic boundary condition is imposed. \hat{X}_j ($\hat{E}_{j,j+1}$) is the site (link) variable at site j [link $(j, j+1)$] and \hat{p}_j ($\hat{n}_{j,j+1}$) is the associated canonical momentum, and they satisfy the standard commutation relations $[\hat{X}_j, \hat{p}_l] = [\hat{E}_{j,j+1}, \hat{n}_{l,l+1}] = i\delta_{jl}$. The eigenbasis of these variables is denoted as $|X_j\rangle$ ($X_j \in \mathbb{R}$) and $|E_{j,j+1}\rangle$ [$E_{j,j+1} \in [0, 2\pi)$], which span the local Hilbert space at site j and link $(j, j+1)$, respectively. The tensor product of these local ones is the full Hilbert space. However, there are additional local operators $\hat{G}_j = e^{i(\hat{E}_{j,j+1} - \hat{E}_{j-1,j} - 2\pi \hat{p}_j)}$ that commute with the Hamiltonian. We define the physical subspace \mathcal{H}_{phy} by imposing the Gauss law constraints $\hat{G}_j = 1 \forall j$. The projection operator onto the subspace \mathcal{H}_{phy} is denoted as \mathcal{P}_G . Using new variables $\hat{\theta}_j = R'(\hat{X}_j + 2\pi \sum_{k=1}^j \hat{n}_{k-1,k})$ and $\hat{\phi}_{j,j+1} = R\hat{E}_{j,j+1}$ [60], Eq. (2) can be rewritten as

$$H_{\text{mV}} = \sum_{j=1}^N [(\hat{\phi}_{j+1,j+2} - \hat{\phi}_{j,j+1})^2 + (\hat{\theta}_{j+1} - \hat{\theta}_j)^2]. \quad (3)$$

Its continuum limit is taken by sending $N \rightarrow \infty$ and $a \rightarrow 0$ with $L = Na$ kept unchanged. After the substitutions $\hat{\phi}_{j+1,j+2} - \hat{\phi}_{j,j+1} \rightarrow a\partial_x \hat{\phi}(x)$ and $\hat{\theta}_{j+1} - \hat{\theta}_j \rightarrow a\partial_x \hat{\theta}(x)$, the TLL Hamiltonian emerges from Eq. (3). The operators $e^{i\hat{\theta}_j/R'}$ and $e^{i\hat{\phi}_{j,j+1}/R}$ act within \mathcal{H}_{phy} thanks to the commutation relations $[e^{i\hat{\theta}_j/R'}, \hat{G}_j] = [e^{i\hat{\phi}_{j,j+1}/R}, \hat{G}_j] = 0 \forall j, j'$. The compactification radii of $\hat{\theta}_j$ and $\hat{\phi}_{j,j+1}$ are R' and R , respectively.

Motivated by the intuitive picture mentioned above and previous works [49–56], we construct the *Ansätze*

$$|C_{\pm}\rangle = \mathcal{P}_G \prod_{j=1}^{N/2} |B_{\pm}^{\text{V}}\rangle_{j,j+N/2} \otimes |B_{\mp}^{\text{L}}\rangle_{j,j+N/2} \quad (4)$$

for crosscap states by putting antipodal links and sites in the generalized Bell pairs

$$\begin{aligned} |B_{\mp}^{\text{L}}\rangle_{j,j+N/2} &= \int_0^{2\pi} dE |E\rangle_{j,j+1} | \mp E \rangle_{j+N/2, j+N/2+1}, \\ |B_{\pm}^{\text{V}}\rangle_{j,j+N/2} &= \int_{-\infty}^{\infty} dX |X\rangle_j | \pm X \rangle_{j+N/2}, \end{aligned} \quad (5)$$

respectively. It can be verified that the relations

$$\begin{aligned} e^{i(\hat{\theta}_j \mp \hat{\theta}_{j+N/2})/R'} |\mathcal{C}_\pm\rangle &= |\mathcal{C}_\pm\rangle, \\ e^{i(\hat{\varphi}_{j,j+1} \pm \hat{\varphi}_{j+N/2, j+N/2+1})/R} |\mathcal{C}_\pm\rangle &= |\mathcal{C}_\pm\rangle \end{aligned} \quad (6)$$

are fulfilled for all j , which in the continuum limit becomes

$$\begin{aligned} e^{i(\hat{\theta}(x) \mp \hat{\theta}(x+L/2))/R'} |\mathcal{C}_\pm\rangle &= |\mathcal{C}_\pm\rangle, \\ e^{i(\hat{\varphi}(x) \pm \hat{\varphi}(x+L/2))/R} |\mathcal{C}_\pm\rangle &= |\mathcal{C}_\pm\rangle. \end{aligned} \quad (7)$$

These identities are sewing conditions in the sense that field operators on antipodal spatial coordinates are combined to form certain operators under which the crosscap states are invariant. It is amusing that they can be derived rigorously using our formalism.

For subsequent calculations, we introduce the mode expansions [61]

$$\begin{aligned} \hat{\theta}(x) &= \tilde{\varphi}_0 + \frac{4\pi x}{L} \pi_0 + \sum_{k \neq 0} \frac{i}{k} (e^{i\frac{2\pi k}{L}x} a_k - e^{-i\frac{2\pi k}{L}x} \bar{a}_k), \\ \hat{\varphi}(x) &= \varphi_0 + \frac{4\pi x}{L} \tilde{\pi}_0 + \sum_{k \neq 0} \frac{i}{k} (e^{i\frac{2\pi k}{L}x} a_k + e^{-i\frac{2\pi k}{L}x} \bar{a}_k), \end{aligned} \quad (8)$$

where the zero modes and oscillatory modes satisfy $[\varphi_0, \pi_0] = [\tilde{\varphi}_0, \tilde{\pi}_0] = i$ and $[a_k, a_l] = [\bar{a}_k, \bar{a}_l] = k\delta_{k+l,0}$, respectively. φ_0 ($\tilde{\varphi}_0$) is an angular variable with radius R (R'), and π_0 ($\tilde{\pi}_0$) is the associated canonical momentum. The eigenvectors of π_0 and $\tilde{\pi}_0$ that can be annihilated by all bosonic annihilation operators $a_{k>0}$ and $\bar{a}_{k>0}$ are denoted as $|n, m\rangle$ ($n, m \in \mathbb{Z}$); they are highest weight states (Virasoro primary states) of the CFT and satisfy $\pi_0|n, m\rangle = (n/R)|n, m\rangle$ and $\tilde{\pi}_0|n, m\rangle = (m/R')|n, m\rangle$. The TLL Hamiltonian in Eq. (1) becomes

$$H = \frac{2\pi v}{L} \left[\pi_0^2 + \tilde{\pi}_0^2 + \sum_{k=1}^{\infty} (a_{-k} a_k + \bar{a}_{-k} \bar{a}_k) - \frac{1}{12} \right], \quad (9)$$

whose energy eigenstates are just Virasoro primary states $|n, m\rangle$ and the descendant states obtained by applying the creation operators $a_{k<0}$ and $\bar{a}_{k<0}$ on top of $|n, m\rangle$. The crosscap states can be expressed as

$$\begin{aligned} |\mathcal{C}_+\rangle &= \kappa_+ \exp \left[\sum_{k=1}^{\infty} \frac{(-1)^k}{k} a_{-k} \bar{a}_{-k} \right] \sum_{n \in \mathbb{Z}} |2n, 0\rangle, \\ |\mathcal{C}_-\rangle &= \kappa_- \exp \left[- \sum_{k=1}^{\infty} \frac{(-1)^k}{k} a_{-k} \bar{a}_{-k} \right] \sum_{m \in \mathbb{Z}} |0, 2m\rangle. \end{aligned} \quad (10)$$

The overall factors are not fixed by normalization but should be chosen as $\kappa_+ = \sqrt{R'}$ and $\kappa_- = \sqrt{R}$ to fulfill the loop-tree channel correspondence [62].

The conformal crosscap states in Eq. (10) are scale-invariant (i.e., they contain no scale) and cannot be normalized. Their lattice counterparts may be normalizable [Eq. (14) for spin-1 = 2 chains defined below] or not normalizable [Eq. (4) for the modified Villain model], depending on the local Hilbert space dimension. In fact, for a given lattice model whose low-energy theory is the compactified boson CFT, the crosscap states together with the coefficients κ_\pm in Eq. (10) should be understood as the continuum description of their lattice counterparts $|\mathcal{C}_{\text{latt}}^\pm\rangle$. This means that they are fully determined by the overlaps of lattice crosscap states with low-lying eigenstates in the continuum limit: $\langle \mathcal{C}_\pm | n, m; \{n_k\}, \{\bar{n}_k\} \rangle = \lim_{N \rightarrow \infty} \langle \mathcal{C}_{\text{latt}}^\pm | \psi_{n,m; \{n_k\}, \{\bar{n}_k\}} \rangle$, where $|\psi_{n,m; \{n_k\}, \{\bar{n}_k\}}\rangle$ is the lattice counterpart of the CFT eigenstate $|n, m; \{n_k\}, \{\bar{n}_k\}\rangle$.

An appealing property of $|\mathcal{C}_\pm\rangle$ is that their overlaps with certain highest weight states $|n, m\rangle$ are universal numbers depending only on the compactification radius (equivalently, the Luttinger parameter):

$$\begin{aligned} |\langle \mathcal{C}_+ | 2n, 0 \rangle|^2 &= R' = \frac{1}{\sqrt{K}}, & n \in \mathbb{Z}, \\ |\langle \mathcal{C}_- | 0, 2m \rangle|^2 &= R = 2\sqrt{K}, & m \in \mathbb{Z}. \end{aligned} \quad (11)$$

It was found in Ref. [39] that the Luttinger parameter appears in the thermal entropy of compactified boson CFTs on the Klein bottle. In view of the spacetime symmetry of CFTs, it is natural to expect that the same variable can also be found using the ground state alone. Our derivation not only puts this expectation on a firm ground but also provides further insights into crosscap states. Some excited states are incorporated in the present formalism, and perturbations to a pure CFT can be investigated [66].

Spin chains with microscopic U(1) symmetry—To apply the general theory in specific lattice models, the crosscap states should be expressed in terms of the microscopic degrees of freedom. It is helpful to begin with spin-1/2 chains. The first one is the XXZ model,

$$H_{\text{XXZ}} = \sum_{j=1}^N (S_j^x S_{j+1}^x + S_j^y S_{j+1}^y + \Delta S_j^z S_{j+1}^z), \quad (12)$$

with $K = \pi/[2(\pi - \cos^{-1} \Delta)]$ for $\Delta \in (-1, 1]$ [1]. The second one is the Haldane-Shastry (HS) model,

$$H_{\text{HS}} = \left(\frac{\pi}{N} \right)^2 \sum_{1 \leq j < l \leq N} \frac{\mathbf{S}_j \cdot \mathbf{S}_l}{\sin^2 \left[\frac{\pi}{N} (j-l) \right]}, \quad (13)$$

with $K = 1/2$ [67,68]. Both models have a U(1) symmetry associated with the conservation of the z component of the total spin.

It is natural to speculate that

$$|\mathcal{C}_{\text{latt}}\rangle = \prod_{j=1}^{N/2} (|\uparrow\rangle_j |\uparrow\rangle_{j+N/2} + |\downarrow\rangle_j |\downarrow\rangle_{j+N/2}) \quad (14)$$

is a crosscap state when N is even [52–56], but there is one subtlety: it corresponds to $|\mathcal{C}_+\rangle$ only if $\text{mod}(N,4)=0$ but does not for the cases with $\text{mod}(N,4)=2$. This property can be traced back to the constraint $(S_j^z - S_{j+N/2}^z)|\mathcal{C}_{\text{latt}}\rangle = 0$. The S_j^z operator in bosonization reads $S_j^z \sim \partial_x \hat{\theta}(x) + (-1)^j \alpha \cos[\hat{\theta}(x)/R']$ for XXZ and HS chains (α is a model-dependent constant). For $\text{mod}(N,4)=0$, the lattice constraint $(S_j^z - S_{j+N/2}^z)|\mathcal{C}_{\text{latt}}\rangle = 0$ is hence consistent with Eq. (7).

The validity of Eq. (14) in the XY model (i.e., $\Delta = 0$ in the XXZ model) and the HS model can be substantiated by analytical calculations. In the hardcore boson basis, the Virasoro primary states $|n, 0\rangle$ have the wave functions $\Psi^\lambda(x_1, \dots, x_M) \propto (-1)^{\sum_i x_i} \prod_{j < k} |\sin[\pi(x_j - x_k)/N]|^\lambda$, where $1 \leq x_1 < \dots < x_M \leq N$ denote the positions of spin- \uparrow sites (occupied by hardcore bosons), and $\lambda = 1$ ($\lambda = 2$) for the XY (HS) chain [69]. The number of bosons M is related to the U(1) charge n via $M = N/2 + n$. It is important to note that the cases with $\text{mod}(N,4)=2$ and $\text{mod}(N,4)=0$ are different. Because $|\mathcal{C}_{\text{latt}}\rangle$ only contains states with even M , $\langle \mathcal{C}_{\text{latt}} | \Psi^\lambda(M) \rangle$ vanishes identically for odd M . We also have $\langle \mathcal{C}_{\text{latt}} | \Psi^\lambda(M = N/2 + 2n) \rangle = 0$ when $\text{mod}(N,4)=2$, so $|\mathcal{C}_{\text{latt}}\rangle$ does not correspond to $|\mathcal{C}_\pm\rangle$ in these cases. In contrast, it can be proven that

$$\begin{aligned} |\langle \mathcal{C}_{\text{latt}} | \Psi^{\lambda=1}(M) \rangle|^2 &= 1, \\ |\langle \mathcal{C}_{\text{latt}} | \Psi^{\lambda=2}(M) \rangle|^2 &= \frac{2^M (M!)^3}{[(M/2)!]^2 (2M)!} = \sqrt{2} + \mathcal{O}(1/M) \end{aligned} \quad (15)$$

if $\text{mod}(N,4)=0$ and $M = N/2 + 2n$ [62]. This is a perfect agreement with the field theory prediction Eq. (11), giving a strong indication that $|\mathcal{C}_{\text{latt}}\rangle$ corresponds to $|\mathcal{C}_+\rangle$ for $\text{mod}(N,4)=0$.

For the XXZ model with $\Delta \neq 0$, the ground state can be computed numerically using the density matrix renormalization group (DMRG) [70]. When the ground state $|\Psi\rangle$ is expressed as a matrix product state (MPS) [71,72], its crosscap overlap $\langle \mathcal{C}_{\text{latt}} | \Psi \rangle$ would be the tensor contraction shown in Fig. 1(b). Numerical results at many different N 's and Δ 's are displayed in Fig. 2. The MPS bond dimension is 2000 for $N = 100$ and 6000 for $N = 400$. The accuracy is very good in a wide range of Δ for a moderate $N = 100$. The relative error of K is of the order 10^{-4} or smaller when $\Delta \in [-0.95, 0.8]$. As Δ approaches 1, clear deviation from exact values is observed [not visible in Fig. 2(a) due to its scale] and the relative error reaches 7.20% at $\Delta = 1$. This

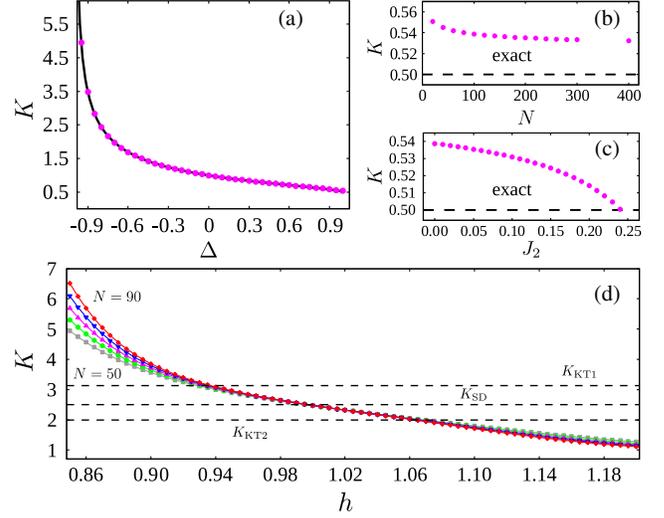


FIG. 2. Numerically extracted Luttinger parameters in three models. Exact values are indicated using solid or dashed black lines for comparison. (a) The XXZ model with $N = 100$ and various Δ . (b) The XXZ model with $\Delta = 1$ and various N . (c) The next-nearest-neighbor model with $N = 100$ and various J_2 . (d) The five-state clock model with various h and N .

deviation is tentatively attributed to the marginal terms in the low-energy theory at $\Delta = 1$ [73], which usually cause strong finite-size effects. It does get smaller as N increases but the convergence is very slow [see Fig. 2(b)]. To validate this conjecture, we study the model $H_{\text{next}} = \sum_{j=1}^N (\mathbf{S}_j \cdot \mathbf{S}_{j+1} + J_2 \mathbf{S}_j \cdot \mathbf{S}_{j+2})$ with a next-nearest-neighbor Heisenberg interaction, for which the marginal term disappears when $J_2 \approx 0.24$ [74]. The results for this model with $N = 100$ are presented in Fig. 2(c). As J_2 increases toward 0.24, the relative error of K decreases and eventually becomes 0.11%. Further analysis based on conformal perturbation theory reveals that finite-size results for the XXZ model with $\Delta = 1$ should be fitted as $K = a[\ln(N)]^{-1} + b$ [62,66]. This leads to $b = 0.510$ using the data points with $N \in [100, 400]$ in Fig. 2(b).

Spin chains without microscopic U(1) symmetry—A fundamental improvement of our method is that microscopic U(1) symmetry is not required. To demonstrate this advantage, we consider the quantum q -state clock model [75–80]

$$H_{q\text{SC}} = - \sum_{j=1}^N (\sigma_j^\dagger \sigma_{j+1} + \sigma_j \sigma_{j+1}^\dagger) - h \sum_{j=1}^N (\tau_j + \tau_j^\dagger), \quad (16)$$

where the \mathbb{Z}_q spin operators are defined as $\sigma|\alpha\rangle = |\text{mod}(\alpha - 1 + q, q)\rangle$ and $\tau|\alpha\rangle = e^{i2\pi\alpha/q}|\alpha\rangle$ with $\alpha = 1, \dots, q$. This model is self-dual (SD) at $h = 1$. For $q \geq 5$ and intermediate h (including the SD point), it hosts a Luttinger liquid that is separated by two Kosterlitz-Thouless (KT) transitions at h_{KT1} and h_{KT2} from gapped phases [81,82].

The transition points are related to each other by duality as $h_{\text{KT1}}h_{\text{KT2}} = 1$. There is no analytical expression for the Luttinger parameter, but its value at the SD point and the KT transition points was predicted to be $K_{\text{KT1}} = q^2/8$, $K_{\text{SD}} = q/2$, $K_{\text{KT2}} = 2$ [40,77]. As in previous models, the *Ansatz*

$$|C_{\text{latt}}\rangle = \prod_{j=1}^{N/2} \sum_{\alpha=1}^q |\alpha\rangle_j |\alpha\rangle_{j+N/2} \quad (17)$$

is proposed as a crosscap state. Numerical results for the $q = 5$ case with $N = 50, \dots, 90$ and various h are presented in Fig. 2(d). The MPS bond dimension is 1500 for $N = 50$ and 2500 for $N = 90$. At the SD point, finite-size systems are described by CFT with irrelevant perturbations, and the Luttinger parameter has a relative error 0.79% at $N = 90$. In the Luttinger liquid regime, the crosscap overlap only exhibits weak dependence on the system size, so the data points for different N almost coincide. For the small and large h phases, the overlap strongly depends on N .

Summary and outlook—To summarize, we have unveiled that the Luttinger parameter can be extracted from the overlaps between individual eigenstates and the crosscap states. The advantages of our method are corroborated by analytical and numerical results. This method may be applied to check if a trial wave function indeed describes the physics of TLL or other CFTs (see, e.g., Refs. [83,84]) in the absence of a Hamiltonian. From the theoretical perspective, our results open up an exciting new avenue and many interesting directions are yet to be explored. The crosscap state used here is constructed from a certain type of Bell states. For integrable spin chains, other types of crosscap states have been studied [53]. Do they have clear counterparts in CFTs? An extension of our method to multicomponent Luttinger liquids, such as the fermionic Hubbard model and SU(3) spin chains, is certainly very desirable. Finally, experimental protocols for measuring the crosscap overlap should be pursued. We hope that interesting results on these topics will be reported in the future.

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