A formula for crossing probabilities of critical systems inside polygons

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Abstract
In this article, we use our results from Flores and Kleban (2015 Commun. Math. Phys. 333 389–434, 2015 Commun. Math. Phys. 333 435–81, 2015 Commun. Math. Phys. 333 597–667, 2015 Commun. Math. Phys. 333 669–715) to generalize known formulas for crossing probabilities. Prior crossing results date back to Cardy’s prediction of a formula for the probability that a percolation cluster in two dimensions connects the left and right sides of a rectangle at the percolation critical point in the continuum limit (Cardy 1992 J. Phys. A: Math. Gen. 25 L201–6). Here, we predict a new formula for crossing probabilities of a continuum limit loop-gas model on a planar lattice inside a 2N-sided polygon. In this model, boundary loops exit and then re-enter the polygon through its vertices, with exactly one loop passing once through each vertex, and these loops join the vertices pairwise in some specified connectivity through the polygon’s exterior. The boundary loops also connect the vertices through the interior, which we regard as a crossing event. For particular values of the loop fugacity, this formula specializes to FK cluster (resp. spin cluster) crossing probabilities of a critical $Q$-state random cluster (resp. Potts) model on a lattice inside the polygon in the continuum limit. This includes critical percolation as the $Q = 1$ random cluster model. These latter crossing probabilities are conditioned on a particular side-
alternating free/fixed (resp. fluctuating/fixed) boundary condition on the polygon’s perimeter, related to how the boundary loops join the polygon’s vertices pairwise through the polygon’s exterior in the associated loop-gas model. For \(Q \in \{2, 3, 4\}\), we compare our predictions of these random cluster (resp. Potts) model crossing probabilities in a rectangle \((N = 2)\) and in a hexagon \((N = 3)\) with high-precision computer simulation measurements. We find that the measurements agree with our predictions very well for \(Q \in \{2, 3\}\) and reasonably well if \(Q = 4\).

Keywords: crossing probability, conformal field theory, Schramm–Loewner evolution, percolation, random cluster model, Potts model, loop-gas model

(Some figures may appear in colour only in the online journal)

1. Introduction

For a statistical lattice or random-walk/loop model [1–19] in a planar domain, a crossing event is an event in which a ‘crossing’ path along (for example) nearest-neighbor sites in a common state connects nonadjacent segments of the domain’s boundary. The probability of such an event, called a crossing probability, may characterize critical behavior of the system. Indeed, in the continuum limit, obtained by sending the lattice spacing to zero and the number of lattice sites to infinity such that the lattice fills the domain, a crossing probability often equals zero below some particular point (e.g. temperature) and equals one above it (or vice-versa), indicating a phase transition [3, 20, 21]. In particular, at a critical point, the crossing probability is typically neither zero nor one, and to either predict or rigorously derive its precise value is usually non-trivial [22–30]. This class of observables has attracted considerable interest from researchers in recent decades [3, 20–32]. In this article, we use conformal field theory (CFT) [33–35] and results from our previous research [36–40] to predict new formulas for crossing probabilities of various statistical models on a lattice inside a 2N-sided polygon. Our results apply in the continuum limit and at the critical point.

The continuum limit of Bernoulli percolation (henceforth called ‘percolation’) on a lattice inside a rectangle gives one of the simplest examples of a crossing probability at a critical point. We define site (resp. bond) percolation to be the assignment of either one (the ‘activated’ or ‘occupied’ state) or zero (the ‘deactivated’ or ‘vacant’ state) to every vertex (resp. edge) of a given graph. All state assignments are i.i.d. Bernoulli random variables, with common probability \(p \in [0, 1]\) for the activated state. Throughout this article, the vertices of the graph are the sites of a two-dimensional lattice, and the edges of the graph are the bonds between pairs of nearest-neighbor sites. For bond percolation, we define a (bond) cluster to be a connected collection of activated bonds, and in site percolation, we define a (site) cluster to be a collection of activated sites such that the bonds between them, when activated, form a bond cluster.

For a statistical lattice model such as percolation on a lattice inside a rectangle, we define a rectangle crossing event to be the event that a cluster touches the top and bottom sides of a rectangle (figure 1). This event signals critical behavior in the continuum limit of percolation, as follows. First, it is well known that, in the continuum limit, if \(p < p_c\) for some (lattice-dependent) critical value \(p_c\), then the probability \(P(p)\) of the rectangle crossing event equals zero, and if \(p > p_c\), then this probability equals one (figure 2) [20, 41, 42]:
The jump (1) of $P(p)$ signals a phase transition in percolation. At the critical point or critical value $p = p_c$, the crossing probability $P(p_c)$ is strictly between zero and one. Using CFT, Cardy predicted a formula for this probability as a function of the rectangle’s aspect ratio $R$ (bottom to left side). (A non-trivial dependence on the geometry of the domain is typical for observables measured at critical points.) This formula, called Cardy’s formula, is

$$P(p) = \begin{cases} 0, & 0 \leq p < p_c, \\ 1, & p_c \leq p \leq 1, \end{cases}$$

where $p_c \in (0, 1)$.

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$$P(p_c, R) = \frac{3\Gamma(2/3)}{\Gamma(1/3)^2} m^{1/3} \text{Ei} \left( \frac{1}{3}, \frac{2}{3}, \frac{4}{3} \right) m \right),$$

where $R = \frac{K(m)}{K(1 - m)}$.

with $K(m)$ the elliptic function of the first kind [43, 44] (so $R \in (0, \infty)$ corresponds one-to-one with $m \in (0, 1)$, see (141) below). Numerical simulations of $P(p, R)$ [21, 32] verified his prediction (2) with very high accuracy for both site and bond percolation on a variety of...
lattices. Less than a decade after Cardy’s prediction \((2)\), Smirnov discovered a rigorous proof of Cardy’s formula for site percolation on the triangular lattice. His proof elucidated a key property of critical percolation in the continuum limit: the conformal invariance of certain observables, that is, invariance under conformal transformation of the system domain (while preserving the underlying lattice) \([45]\).

In this article, we study two generalizations of these results. First, we consider a range of other lattice models with conformally invariant critical points, and second, we consider polygons \(P\) with an arbitrary even number of sides \(2N\). After enumerating the sides of \(P\) in counterclockwise order and ascending from one to \(2N\), we allow clusters to join the odd sides of \(P\) in any one of the \(C_N\) available topologically distinct connectivities \([31]\), called crossing patterns (figure 3). Here, \(C_N\) is the \(N\)th Catalan number:

\[
C_N = \frac{(2N)!}{N!(N+1)!}.
\]

For example, a cluster may join the two horizontal sides of a rectangle \((N = 2)\) or not, for \(C_2 = 2\) crossing patterns. We regard a ‘crossing event’ as the event that a particular crossing pattern occurs in \(P\), and we regard a ‘crossing probability’ as the probability of a crossing event.

Using CFT \([33–35]\) and results from our previous work \([36–40]\), we predict a formula ((24) of section 2) for crossing probabilities in the continuum limit of the above mentioned models inside even-sided polygons. We always distinguish two adjacent sides of a polygon \(P\) by imposing different boundary conditions (BC) on those sides. Specifically, the models, crossing events, and BCs that we consider are as follows (figure 4):

(i) **The \(Q\)-color random cluster model** \([8–10]\): \(Q \in \{1, 2, 3, 4\}\). This model is the same as bond percolation, except that we independently and uniformly distribute any one of
The $Q$-state Potts model [5, 7]: $Q \in \{2, 3, 4\}$. Here, each lattice site is assigned one of $Q$ values, called spins, and spin configurations are Boltzmann-weighted by the total energy of nearest-neighbor spin coupling. We always condition on the Potts model fluctuating/fixed side-alternating boundary condition (PMB) event (section 3.4). In a PMB event, all sites adjacent to the odd sides of the polygon $\mathcal{P}$ are assigned the first spin state (so each side exhibits the ‘fixed’, or ‘wired’ BC, but fixed to the first spin state) and the sites adjacent to the even sides uniformly sample any but the first state (so each side exhibits the fluctuating BC) [46]. We define a Potts model spin cluster to be a site cluster whose sites share a common spin value, and we define a Potts model crossing event (PMX) to be the event that spin clusters join the odd sides of $\mathcal{P}$ in a particular crossing pattern.

The loop-gas model (or $O(n)$ model) [11–14]: In this model, each configuration consists of a set of nonintersecting loops that follow the bonds of the lattice. Every such loop configuration is weighted by a temperature parameter $x > 0$ raised to the power of the total length of the loops in the configuration, multiplied by a fugacity parameter $n \in (0, 2]$ raised to the power of the number of loops in the configuration. This model includes both items i and ii in the sense that for certain $Q$-dependent values of $n$, its continuum limit is expected to match the continuum limit of the $Q$-state random cluster model or Potts model (section 4). We condition on a loop-gas exterior arc boundary condition (LGB) event, where ‘boundary loops’ (defined in section 4.1 below) join the vertices pairwise in some connectivity via the polygon’s exterior. We define a loop-gas crossing event (LGX) to be the event that boundary loops join the vertices pairwise in some particular connectivity through the polygon’s interior. This definition is natural because these connectivities bijectively correspond to the $C_N$ crossing patterns (figure 16).

We discuss the above mentioned BCs. In order for the side-alternating BCs of items i–iii to be well-defined, the polygon must have an even number of sides. Also, in items i and ii, only one cluster, called a boundary cluster, touches a given fixed side of the polygon, as that entire side is part of the cluster. In section 4, we observe that items i and ii are special cases of item iii in the continuum limit, with $Q$ and $n$ related by

$$n = \sqrt{Q} \text{ in item i, } \quad n = -2 \cos \left( \frac{\pi^2}{\arccos(-\sqrt{Q}/2)} \right) \text{ in item ii.}$$

This observation is an important step in our argument because in section 4, we connect the random cluster model and the Potts model to CFT via the loop-gas model. (In principle, $Q$ could be any positive number in the definition of the random cluster model. In this article, we restrict $Q$ to the range $(0,4]$, corresponding to $n \in (0, 2]$ for the related loop-gas model (4), where the random cluster model is expected to have a conformally invariant critical point [45], and we restrict $Q$ to positive-integer values in order to interpret it as a number of available colors. Thus, $Q \in \{1, 2, 3, 4\}$.)

Finally, items i–iii are the only BC scenarios that we consider in this article, but they are not the only possibilities. Others include:
• Switching the BCs (but not the models) between items i and ii. We are not aware of any formula for the probability of a crossing scenario with these BCs.

• Assigning the free BC (i.e. assign no BC) to all sides of the polygon. Here, some useful results are known, and we summarize those of which we are aware at the end of section 1.1 below.

We do not consider either of these options because they are beyond the scope of the CFT techniques that we use in this article.

Next, we survey recent research on crossing probabilities.

1.1. Crossing probabilities: a survey of known results

In this section, we survey known results about crossing events in models with conformally invariant critical behavior. As mentioned, percolation gives a simple example of conformally invariant critical phenomena: on an infinite square lattice, if \( p > p_c \) with \( p_c \) as in (1), then a sample almost surely has a unique infinite cluster. If \( p < p_c \), then clusters are almost surely finite and typically with diameter less than some \( \xi(p) < \infty \). Finally, if \( p = p_c \), then \( \xi(p_c) = \infty \), meaning that there is no upper bound on the cluster size (figure 5) and also no infinite cluster \([1, 3, 4]\). (See \([47]\) for an equivalent characterization of this phenomenon.) Thus, as we approach the continuum limit, the maximum cluster size does not decrease if \( p = p_c \), one requirement for conformal invariance at criticality. (This pertains to Bernoulli percolation, but see \([48–50]\) for other percolation models with critical behavior.)

Percolation models inside bounded planar domains can also exhibit critical behavior in the continuum limit. Researchers have predicted that two special properties emerge in this limit at criticality. The first property, universality \([2, 21]\), is that all observables, such as crossing probabilities, do not depend on microscopic details such as the type of percolation (e.g. site versus bond) or the lattice (although the value of \( p_c \) does depend on these details \([41, 52–54]\), see table 1). The second property, conformal invariance \([20, 45, 51]\), is that certain observables are invariant under conformal transformation of the system domain (leaving the underlying lattice unchanged) \([45]\). Extensive numerical simulations strongly support these predictions \([20, 21, 32]\).
Crossing probabilities are one tool for probing critical behavior, universality, and conformal invariance in percolation. For example, the phase change \( (\frac{1}{2}) \) of the rectangle crossing probability \( P(p) \) was rigorously established some time ago on certain lattices [41], (see also [3] and references therein), but the value of \( P(p_c) \) was not known in any percolation model until some years later. Later work measured \( P(p_c) \) for site and bond percolation on several lattices and in many different conformal rectangles [20, 21]. These measurements strongly suggest that \( P(p_c) \in (0, 1) \), with the same value for all regular lattices (universality) in any two conformally equivalent conformal rectangles (conformal invariance).

As this deeper understanding of critical percolation emerged, simultaneous advancements in the application of CFT [33–35] to the study of related critical lattice models [1, 3–5, 7, 8, 11] were underway. Similar to percolation, it was conjectured that many of these models have a conformally invariant phase transition, which could be exploited to predict exact formulas for their correlation functions [55–57] via CFT. One of these models, the \( Q \)-state random cluster model [8], admits percolation as the special case \( Q = 1 \). This observation was used by Cardy to predict the formula (2) for the critical crossing probability \( P(p) \) using CFT methods [22]. Cardy’s result evinced CFT as a reliable, yet perhaps mysterious, tool for predicting this and other critical percolation observables, including Schramm’s observable [58, 59], cluster densities [60–63], pinch-point densities [64], and other examples of rectangle crossing events [23, 65].

Less than a decade after Cardy’s prediction (2), mathematicians discovered rigorous proofs for it, as well as many others in critical percolation. Much of this progress stemmed from Schramm’s discovery of a one-parameter \( (\kappa > 0) \) family of conformally invariant planar stochastic processes, now called Schramm–Loewner evolution (SLE) \( \kappa \) [66–68], together with the conjecture that the law of boundary percolation cluster interfaces converges to SLE\( \kappa \) in the continuum limit [69]. Not long afterward, Smirnov proved this conjecture for critical site percolation on the triangular lattice, which implied conformal invariance of the law of the percolation interfaces [24]. In the same article, Smirnov also proved Cardy’s formula for site percolation on the triangular lattice. (Exploiting conformal invariance, Carleson had previously noted that Cardy’s formula (2) simplifies considerably to a linear function if we conformally map it to the equilateral triangle (figure 6). Smirnov proved Cardy’s formula in this setting.)

Subsequent to these discoveries about percolation crossing probabilities, some natural variations were examined. We consider two such generalizations in this article. First, we consider other models with conformally invariant critical points. Here, there has been much progress. For example, using the Ising model fermionic observable [70], Chelkak and Smirnov rigorously derived this \( Q = 2 \) FK-cluster crossing probability formula [71].

### Table 1

Critical points (thresholds) \( p_c \) for bond and site percolation on various lattices. Some are known exactly, and others are found via computer simulation. For site percolation on the triangular lattice, \( p_c \) follows from an easy symmetry argument.

| Lattice    | Site percolation threshold | Bond percolation threshold |
|------------|---------------------------|----------------------------|
| Square     | 0.592 746 … [53]          | 1/2 [41]                   |
| Triangular | 1/2                       | 2 \sin(\pi/18) [52]        |
| Hexagonal  | 0.697 040 … [53]          | 1–2 \sin(\pi/18) [52]     |
| Kagome     | 1–2 \sin(\pi/18) [52]     | 0.524 405 … [54]          |
\[ P(R) = \mathbb{P}\left\{ \text{one FK cluster touches both the top and bottom of the rectangle,} \right. \\
\left. \text{conditioned on both of these sides being fixed to the same color} \right\} \]

\[ = \frac{\sqrt{2} (1 - p(m))}{p(m) + \sqrt{2} (1 - p(m))}, \quad \text{where} \]

\[ p(m) := \frac{\sqrt{1 - m}}{\sqrt{1 - m} + \sqrt{1 - \sqrt{1 - m}}} \quad \text{and} \quad R = \frac{K(m)}{K(1 - m)}, \quad (5) \]

(figure 7), and another formula (133) for a similar observable. (Recently, bounds for similar observables in conformal rectangles were derived in [72, 73].) Another example, using CFT, Arguin and Saint-Aubin [27] predicted this spin-cluster crossing probability formula for the \( Q \)-state Potts model [7] with \( Q = 2 \) (i.e. Ising model):

\[ P(R) = \mathbb{P}\left\{ \text{one spin cluster touches both the top and bottom of the rectangle} \right. \\
\left. \text{conditioned on these two sides being fixed to the first spin value} \right. \\
\left. \text{and the other two sides fixed to the second} \right\} \]

\[ = \frac{2 \Gamma(4/3)}{1 - m + m^2} F_1\left( \frac{1}{3}; \frac{4}{3}; \frac{8}{3}; m \right), \quad \text{where} \quad R = \frac{K(m)}{K(1 - m)}, \quad (6) \]

(figure 7). (This formula (6) is simplified here. See beneath (139).) Subsequently, Kozdron rigorously proved (6) [28]. Interestingly, this formula (6) was pursued by several research groups almost simultaneously some years ago, and it appears as an original result in many articles [26–30], [29] gives a chronological account of this work.

As a second generalization, we consider polygons with an even number of sides greater than four. As mentioned, a \( 2N \)-sided polygon admits \( C_N \) different crossing patterns [31] (figure 3), where \( C_N \) is the \( N \)th Catalan number (3). Recent results of this kind include Dubédat’s discovery [30] that (figure 7)

\[ \mathbb{P}\left\{ \text{a percolation cluster connects} \right. \\
\left. \text{two nonadjacent odd sides} \right. \\
\left. \text{of a regular hexagon} \right\} = \frac{2}{3} \left( \frac{\sqrt{3}}{2^{2/3} \pi} \right)^5 \Gamma(2/3)^5 \mathcal{F}_2\left( \frac{1}{6}, \frac{5}{6}; \frac{5}{2}, \frac{3}{2}; 1 \right) \quad (7) \]

and Simmons’s determination of all \( C_5 = 5 \) hexagon percolation crossing probabilities with free/fixed side-alternating BCs (and free BCs) [25]. Also, Izyurov has rigorously derived some
Formulas for $Q = 2$ FK-cluster crossing probabilities in (conformal) polygons with more general BCs [74].

In the examples of crossing probabilities (2), (5)–(7), the crossing event pairs with its companion BC as in item i or ii above. As such, each of these is a special case of our predicted formula (24) for crossing probability. However, one may consider other scenarios. For example, it is natural to assign the free BC (i.e. no BC) to all sides of the polygon, and for this, there are many interesting results: Watts and Simmons have used CFT to predict formulas for percolation crossing probability in a rectangle and a hexagon respectively, with the free BC [23, 25]. Later, Dubédat and Schramm independently discovered proofs of Watts’s formula [75, 76]. Langlands et al found numerical evidence for the universality of the $Q = 2$ Potts (i.e. Ising) model rectangle crossing probability (6) [77], and Benoist et al rigorously confirmed this fact [78]. Also, Lapalme and Saint-Aubin used CFT and numerical methods to predict a formula for this probability [79]. To determine such crossing probabilities with free BCs is beyond the scope of this article. Other variations that are also beyond the scope of this article include the average number of crossing clusters [80] and crossing cluster densities [65] for critical percolation.

1.2. Objectives and organization

The main purpose of this article is to predict the new formula (24), introduced in section 2, for the probabilities of all $C_N$ distinct crossing events (figure 3) in the critical models of items i–iii inside a $2N$-sided polygon $P$ with the specified BCs in the continuum limit. Our prediction is exact.

Formula (24), our main result, is given in terms of four quantities:

- the Coulomb gas (contour integral) function $\mathcal{F}_\vartheta$ (item iii of section 2), whose subscript $\vartheta \in \{1, 2, \ldots, C_N\}$ indexes the BC,
- the ‘crossing weight’ (i.e. ‘connectivity weight’ [39, 40] or ‘pure SLE$_\kappa$ partition function’ [26, 81]) $\Pi_\varsigma$, (item v of section 2), whose subscript $\varsigma \in \{1, 2, \ldots, C_N\}$ indexes the crossing pattern,
- the number $l_{\varsigma, \vartheta}$ of closed loops in the $(\varsigma, \vartheta)$th ‘product diagram’ (item iv of section 2 and figures 11 and 19),
- and the loop fugacity $n = n(\kappa)$ of the loop-gas model (item i of section 2).
For the random cluster model and Potts model, the loop fugacity $n$ depends on $Q$ via (4), and throughout this article, we parameterize it (27) in terms of the SLE$_\kappa$ [66–68] parameter $\kappa \in [3/8, 8)$. In the random cluster model, the Potts model, and the loop-gas model respectively, the BC (resp. crossing) event is an $RCB$ event, a $PMB$ event, and an $LGB$ event (resp. an $RCX$ event, a $PMX$ event, and an $LGX$ event), as defined in items i–iii above. We first encounter the functions $\Pi$, and $\mathcal{F}_5$ (resp. the quantity $\varsigma$) in section 5 (resp. section 3).

Our strategy is to first relate the lattice models of items i and ii above to the loop-gas model of item iii above. We then consider the continuum limit of the latter, to which CFT applies. Making use of our previous research, we then derive the crossing probability formula (114), which is equivalent to (24). Because CFT should apply directly to the original lattice models of items i and ii, one might wonder why we take this somewhat circuitous route. The application of CFT to all of the cases we consider (various models, BCs, and number of polygon sides) is not completely straightforward. Our route, although a bit indirect, encompasses all these cases in a single scheme.

This article is organized as follows. To begin, section 2, which is self-contained, gives a detailed explanation of the crossing probability formula (24) (but does not derive it), and section 6 applies it to rectangles, reproducing the rectangle crossing probability for critical percolation (2), [22, 24], the critical $Q = 2$ random cluster (i.e. FK-Ising) model (5) [71], and the critical $Q = 2$ Potts (i.e. Ising) model (6) [26–28].

The purpose of sections 3–5 is to derive the predicted crossing probability formula (24). To begin, in section 3, we give formal expressions for various partition functions of a random cluster model or a Potts model on a lattice inside $\mathcal{P}$, and we give formal expressions (48), (56) for crossing probabilities in terms of these partition functions. All of these partition functions either sum exclusively over the BC events described in items i and ii above, or they sum exclusively over these events intersected with the crossing events, also described in items i and ii above. For the random cluster model, in section 3.3, we let $X_{\varsigma, \vartheta}^P$ (46) denote the latter partition function, where $\varsigma, \vartheta \in \{1, 2, \ldots, C_N\}$ index the $RCX$ (crossing) event and $RCB$ (BC) event respectively, and we let $X_{\varsigma, \vartheta}^P = X_{1, \vartheta}^P + X_{2, \vartheta}^P + \cdots + X_{C_N, \vartheta}^P$ (47), a sum over all crossing events, denote the former. For the Potts model, in section 3.4, we let $Y_{\varsigma, 1}^P$ denote the latter partition function, where $\varsigma \in \{1, 2, \ldots, C_N\}$ indexes the $PMX$ (crossing) event (here, the $PMB$ event only allows the same single state on all fixed sides, which results in $\vartheta = 1$), and we let $Y_{\varsigma, 1}^P = Y_{1, 1}^P + Y_{2, 1}^P + \cdots + Y_{C_N, 1}^P$, a sum over all crossing events, denote the former. For the random cluster model, we only use certain $RCB$ events called ‘mutual wiring events’, defined at the beginning of section 3.2. Furthermore, we do not use all mutual wiring events but consider only a subset of them called ‘basic events’, defined in section 3.2. (In appendix B, we show that if $Q \in \{1, 2, 3\}$ or if $Q = 4$ and $N \in \{1, 2, 3\}$, then in the continuum limit, any partition function summing exclusively over a mutual wiring event equals a linear combination of partition functions that all sum exclusively over a basic event. Hence, restricting our attention to basic events does not result in a loss of generality for these $(Q, N) \in \{1, 2, 3, 4\} \times \mathbb{Z}^+$. ) There are $C_N$ basic events, and $\vartheta \in \{1, 2, \ldots, C_N\}$ indexes them. The partition function $X_{\varsigma, \vartheta}^P$ sums over the $\vartheta$th basic event, and we readily obtain a formal expression (50) for this partition function up to a certain combinatorial quantity $C_{\varsigma, \vartheta}$ that counts the number of ways to distribute color to boundary clusters in this event. In section 3.3, we use basic graph theory to find a formula (51) for $C_{\varsigma, \vartheta}$ in terms of the number of loops $l_{\varsigma, \vartheta}$ (figure 19).

Next, in section 4, we give formal expressions for various partition functions of a loop-gas model on a lattice inside $\mathcal{P}$, and we give a formal expression (81) for the crossing probability in terms of these functions. All of these partition functions either sum exclusively over the BC
events described in item iii above, or they sum exclusively over these events intersected with the crossing events, also described in item iii above. For the loop-gas model, in section 4.1, we let $Z_{\zeta,\vartheta}^{i}$ (62) denote the latter partition function, where $\zeta, \vartheta \in \{1, 2, \ldots, C_N\}$ index the LGX (crossing) event and LGB (BC) event respectively, and we let $Z_{\vartheta}^{i} = Z_{1,\vartheta}^{i} + Z_{2,\vartheta}^{i} + \cdots + Z_{C_N,\vartheta}^{i}$, a sum over all crossing events, denote the former. Then in sections 4.2 and 4.3 respectively, we argue that, at the critical point, the random cluster model partition functions and Potts model partition functions, $X_{\zeta,\vartheta}^{i}$ and $Y_{\zeta,1}^{i}$ respectively, are asymptotic to $Z_{\zeta,\vartheta}^{i}$ and $Z_{\zeta,1}^{i}$ respectively in the continuum limit (77). As such, to determine crossing probabilities for the critical random cluster model and the critical Potts model, it suffices to determine them for the corresponding loop-gas model. In section 4.4, we give formal expressions for the latter.

Finally, in section 5, we use physical and mathematical arguments to determine explicit formulas for the critical random cluster model, the critical Potts model, and the loop-gas model crossing probabilities (48), (56), (81) in the continuum limit. In order to do this, in section 5.1, we first relate the loop-gas model partition functions $Z_{\vartheta}^{i}$ in the continuum limit, to boundary CFT correlation functions $\Upsilon_{\vartheta}^{i}$ known to satisfy a particular system of $2N + 3$ partial differential equations (PDEs) (87), (88). The series of articles [36–39] determines the dimension of, as well as explicit formulas for all elements of, the system’s solution space $S_{N}$. (As part of the definition for $S_{N}$, we also require that the solutions are dominated by a product of certain power functions. See (90) of section 5 for the precise bound.) The solutions corresponding to different BCs are distinguished by the choice of contour integrals that define the coulomb gas function $\mathcal{F}_{\vartheta}$ (item iii in section 2), or alternatively, their asymptotic behavior as neighboring corners of the polygon approach each other. (See the proof of lemma 6 in [38].) Then in section 5.2, we argue that $Z_{\zeta,\vartheta}^{i}$ is asymptotic to a special element $\Pi_{\zeta}$ of $S_{N}$ called a crossing weight (or equivalently, ‘connectivity weight’ [39, 40] or ‘pure SLE$_{C_{N}}$ partition function’ [26, 81], see the remarks just below (110) and above (115)). The crossing weights are also solutions of the PDEs just mentioned, i.e. elements of $S_{N}$, linearly related to the $\Upsilon_{\vartheta}^{i}$ by the meander matrix $M_{N}$ [82–85], defined in item v of section 2. Finally, a simple argument gives $\Upsilon_{\vartheta}^{i} = \mathcal{F}_{\vartheta}$, which leads to our main result, the explicit formula (24) (or (114)) for the probability $P_{\zeta,\vartheta}^{i}$ (81) of the $\zeta$th LGX (crossing) event conditioned on the $\vartheta$th LGB (BC) event for the loop-gas model in the continuum limit (item iii above). As special cases, this includes the probability $P_{\zeta,\vartheta}^{i}$ (48) of the $\zeta$th RCX (crossing) event conditioned on the $\vartheta$th basic RCB (BC) event for the critical random cluster model (item i above), and the probability $P_{\zeta,1}^{i}$ (56) of the $\zeta$th PMX (crossing) event conditioned on the PMB (BC) event for the critical Potts model (item ii above), both in the continuum limit. The discussion around (115) also relates (114) to a similar formula for the probability of eventual curve connectivity in multiple SLE$_{C_{N}}$ [86–89].

To support our predictions, we present high-resolution computer results in section 7. The simulations measure crossing probabilities of the $Q \in \{2, 3, 4\}$ random cluster (resp. Potts) model on a large square lattice inside a rectangle (section 7.1) and on a large triangular lattice inside a hexagon (section 7.2), with the side-alternating BCs of item i (resp. item ii). We compare these measurements with our prediction (24), finding very good agreement for $Q \in \{2, 3\}$ and reasonable agreement for $Q = 4$. (For the latter case, much larger lattices and many more samples are required for better agreement [90, 91].) (Similar results for the $Q = 1$ critical random cluster model, i.e. critical percolation, already appear in [22, 92].)

Three appendices accompany this article. In appendix A, we use CFT ‘corner operators’ to regularize the conformal transformation of the boundary CFT correlation functions.
in section 5 from the upper half-plane onto the polygon. The need for regularization arises because this transformation is not conformal at the polygon’s vertices. In appendix B, we study random cluster model partition functions $W^P_c$ that sum exclusively over ‘color schemes’ $c$. These are mutual wiring events in which the color of each fixed side of $P$ is specified. Any partition function summing over an RCB event is a sum of color scheme partition functions $W^P_c$. We show that, in the continuum limit, the span of the partition functions $W^P_c$ (as functions of the shape of $P$) is the same as that of the $X^P_0$ for $Q \in \{1, 2, 3\}$ and any $N \in \mathbb{Z}^+$ or $Q = 4$ and $N \in \{1, 2, 3\}$. Thus, for these $(Q, N)$, working with $X^P_0$ over $W^P_c$, as we do in this article, does not result in loss of generality. In appendix C, we give some simpler formulas for the Coulomb gas function $F_0$ (11) appearing in the denominator of the crossing probability formula (24). They include an alternative formula (C.2) used in [38, 39] for all $\kappa > 0$ and some explicit algebraic formulas (C.7), (C.10), (C.14) for $F_0(\kappa)$ that hold at certain rational values of $\kappa$. Some of these algebraic formulas already appear in the literature ($Q = 2$ [81, 120]), although not in direct connection with $F_0(\kappa)$, and others are, to our knowledge, completely new ($Q \in \{1, 2, 3\}$ for $Q \in \{1, 2\}$, we obtain our algebraic formulas not by explicit evaluation of the multiple contour integrals that appear in $F_0(\kappa)$, but by using a new if indirect method making use of our previous results in [36–39] and other results already in the literature [81, 120]. Thus we rigorously derive algebraic formulas for some rather complicated multiple contour integrals based on a full analysis of the solutions of the system of PDEs (87), (88) that $F_0(\kappa)$ satisfies.

2. Formula for crossing probability

This section contains a step-by-step but unmotivated presentation and explanation of our central result, the formula (24) for the particular crossing probabilities that we consider in various critical statistical mechanics models inside polygons. It is completely self-contained (although reading the introduction section 1 of this article first may be useful) and is illustrated for the rectangle case in section 6. We emphasize that this section simply states and explains (24); its derivation only appears much later at the end of section 5.

The formula (24) for crossing probability has complex contour integrals that use Pochhammer contours, so before presenting this formula, we define a Pochhammer contour and explain how to integrate around it. Before we give this definition, we consider a complex contour integral of the following form:

$$\oint_{\mathcal{C}_0} (u - x_i)^{\beta}(x_j - u)^{\beta} f(u) \, du, \quad \text{where } f(u) \text{ is analytic in a domain containing } \Gamma. \quad (8)$$

The power functions $(u - x_i)^{\beta}$ and $(x_j - u)^{\beta}$ each insert a branch cut into the domain of the integrand in (8), with one endpoint at $x_i$ and $x_j$ respectively. We view $\Gamma$ as a contour on the ‘branched’ domain of the integrand (i.e. different Riemann sheets of the integrand are disjoint subsets of the integrand’s domain), so the integrand varies continuously as $\Gamma$ crosses one of the branch cuts. Hence, tracing a cycle that surrounds $x_i$ (resp. $x_j$) sends the integrand of (8) from some initial value to that same value multiplied by $e^{\pm \pi i \beta}$ (resp. $e^{\pm \pi i \beta}$), where the sign $\pm$ matches the orientation of the cycle. Tracing the same cycle in the opposite direction produces the reciprocal phase factor.

The Pochhammer contour $\mathcal{C}_0(x_i, x_j)$ is the integration contour shown on the left side of figure 8. It comprises four cycles traced in this order: the first cycle circles around $x_j$ clockwise; the second cycle circles around $x_i$ counterclockwise; the third cycle circles around $x_j$
and the fourth cycle circles around \( x_i \) clockwise. According to the previous paragraph, the phases accumulated by these cycles are \( \pi \beta_i e^{2i\pi j} \), \( \pi \beta_j e^{2i\pi i} \), \( \pi \beta_j e^{2i\pi i} \), and \( \pi \beta_i e^{2i\pi j} \) respectively. Because the product of these phases equals one, if we approach the ‘start’ point from the right or the ‘end’ point from the left on the Pochhammer contour in figure 8, then in either limit, the integrand of (8) yields the same value. Viewed as a contour on the larger ‘branched’ domain of the integrand, the Pochhammer contour is thus closed.

Using the Cauchy integral theorem, we may deform the Pochhammer contour \( \mathcal{P}(x_i, x_j) \), so its cycles are very small circles surrounding the branch points \( x_i \) and \( x_j \) (which we call ‘endpoints’ of \( \mathcal{P}(x_i, x_j) \)), and the rest of the contour is simple paths along the real axis and joining two cycles. We find four cycles joined by four simple paths, and the integration goes from path to cycle to path to cycle, etc. The right part of figure 8 shows the decomposition of \( \mathcal{P}(x_i, x_j) \) into these eight constituents. If \( \beta_i, \beta_j > -1 \), then the integration around the circles vanishes as their radii vanish, leaving only the integration along the simple paths. In this situation, the bottom line of figure 8 leads to the following simplification, which holds for any function \( f(u) \) is analytic in a domain containing \( \mathcal{P}(x_i, x_j) \) (figure 9):

\[
4e^{i(\beta_i-\beta_j)} \sin(\pi \beta_i) \sin(\pi \beta_j) \text{ if } \beta_i, \beta_j > -1
\]
\[
\int_{\mathbb{R}(x_j,x_k)} (u - x_j)^{\beta} (x_j - u)^{\beta} f(u) \, du = 4e^{\pi i (\beta - \beta')} \sin(\pi \beta) \sin(\pi \beta') \\
\quad \times \int_{\text{int}} (u - x_j)^{\beta} (x_j - u)^{\beta} f(u) \, du \quad \text{if } \beta, \beta' > -1, \quad (9)
\]

In steps i–vi below, we identify the interior of a 2N-sided polygon with the upper half-plane, and we identify the polygon’s vertices with 2N marked points \(x_1 < x_2 < \ldots < x_{2N}\) on the real axis. In step vii, we conformally map the upper half-plane onto a conformal 2N-sided polygon \(\mathcal{D}\) in the complex plane.

(i) We let \(x_1 < x_2 < \ldots < x_{2N}\) be 2N marked points on the real axis and let \(\kappa \in (0, 8)\) (the SLE\(_\kappa\) parameter) \([66–68]\). We also define the loop-gas model fugacity formula \([51, 93]\) to be

\[n(\kappa) = -2 \cos(4\pi/\kappa).\] \(10\)

(ii) We enumerate 1 through \(C_N\) (with \(C_N = (2N)!/N!(N + 1)!\) the \(N\)th Catalan number) the distinct arc ‘connectivities’ or ‘crossing patterns’ (figure 10). Each is a collection of \(N\) curves (arcs) in the upper half-plane

- (a) that do not intersect themselves or each other,
- (b) with all of their endpoints among the \(x_j\),
- (c) with two distinct endpoints per arc, and
- (d) with no two arcs sharing any endpoints.

In particular, we let the first (resp. \(C_N\)th) connectivity be such that the \(j\)th arc has its endpoints at \(x_{2j-1}\) and \(x_{2j}\) for \(1 \leq j < N\) and the \(N\)th arc has its endpoints at \(x_1\) and \(x_{2N}\) (resp. \(j\)th arc has its endpoints at \(x_{2j-1}\) and \(x_{2j}\)).

(iii) For each \(\vartheta \in \{1, 2, \ldots, C_N\}\), we let \(\mathcal{F}_\vartheta\) be the \(\vartheta\)th Coulomb gas function, given by the following formula (with explanation of the integration contours and the symbol \(\mathcal{N}\) beneath it):

\[
\mathcal{F}_{\vartheta}(\kappa \mid x_1, x_2, \ldots, x_{2N}) := \left[\frac{n(\kappa)\Gamma(2 - 8/\kappa)}{4 \sin^2(4\pi/\kappa)\Gamma(1 - 4\kappa)}\right]^N \left(\prod_{j < k} (x_k - x_j)^2/\kappa\right) \\
\quad \times \oint_{\gamma_1} \cdots \oint_{\gamma_2} \oint_{\gamma_3} \int_{\mathcal{N}} \left[\prod_{j=1}^{2N} (x_k - x_{2j})/\kappa\right] \\
\quad \times \left[\prod_{p=1}^{N} (u_p - u_{2j})/\kappa\right] \right]^{(11)}
\]

Figure 10. An enumeration of the \(C_3 = 5\) arc connectivities involving six points.
(a) **Contours:** \( \Gamma_1, \Gamma_2, \ldots, \Gamma_N \) are nonintersecting Pochhammer contours with endpoints among \( x_1, x_2, \ldots, x_N \) (figure 8) and bent to lie completely in the upper half-plane (except for where they wrap around their endpoints).

(b) **Endpoints:** The two ‘endpoints’ of \( \Gamma_m \) (i.e. the two points surrounded by a cycle of \( \Gamma_m \)) are the endpoints of the \( m \)th arc in the \( \vartheta \)th connectivity (as we have enumerated the arc connectivities in item ii above).

(c) **Branching:** We choose the branch of the logarithm for the power functions in (11) so \(-\pi < \arg z \leq \pi \) for all \( z \) in the complex plane. Thus, the branch cut of \( f(z) = z^{\beta} \) for \( \beta \notin \mathbb{Z} \) is \(( -\infty, 0)\).

(d) **Simplification:** If \( \kappa > 4 \), then we use (9) to simplify (11): we replace each contour by a simple curve (that bends into the upper half-plane) with the same endpoints, and we drop all factors of \( (\pi\kappa/4 \sin 4\beta/\kappa) \) from (11).

(e) **Integrand:** The integrand of (11), presently represented as a certain Coulomb-gas-type function enclosed between the square brackets of \( \ldots \) \( \cdot \), is given by the following. (Throughout, if \( \iota \) \( \in \{\ldots, \} \) \( \cdot \), then we write \( \iota = \eta \) \( \cdot \), and as usual, \( u_m \) denotes the integration variable that traces \( \Gamma_m \).

1. For each \( m \in \{1, 2, \ldots, N\} \), we use the following factors in the integrand of (11):

\[
(u_m - x_{(2m-1)})^{-4\kappa/\kappa} (x_{(2m)} - u_m)^{-4\kappa/\kappa}.
\]

The branch cuts (drawn as zig-zags) of (12), seen as a function of \( u_m \), are thus arranged as follows: (Here, we show \( \Gamma_m \) as a simple curve as in item (iii)d. If \( \kappa \leq 4 \), then we replace \( \Gamma_m \) with the Pochhammer contour (figure 8) of item (iii)a.)

\[
\begin{array}{c}
\bullet \quad u_m \\
\bullet \quad x_{(2m-1)} \\
\bullet \quad x_{(2m)}
\end{array}
\]

2. For each pair \( p, q \in \{1, 2, \ldots, N\} \) with \( p \neq q \), if \( \Gamma_q \) does not nest \( \Gamma_p \) and vice versa, then taking \( \Gamma_p \) to be left of \( \Gamma_q \) by pressing these contours so they are sufficiently close to the real line, we ensure that

\[
\arg(u_p - u_q) \in (\pi/4, \pi/4).
\]

In this situation, we use the following factors in the integrand of (11):

\[
\begin{align*}
(u_p - x_{(2q-1)})^{-4\kappa/\kappa} & (u_p - u_q)^{-4\kappa/\kappa} \\
(x_{(2p)} - u_q)^{-4\kappa/\kappa} & (x_{(2p)} - u_q)^{-4\kappa/\kappa} (u_p - u_q)^{8\kappa/\kappa}.
\end{align*}
\]

The branch cuts (drawn as zig-zags) of (15), seen as a function of \( u_p \) and \( u_q \) separately, are thus arranged as follows: (here, we show \( \Gamma_m \) as a simple curve as in item (iii)d. If \( \kappa \leq 4 \), then we replace \( \Gamma_m \) with the Pochhammer contour (figure 8) of item (iii)a.)

\[
\begin{array}{c}
\bullet \quad u_p \\
\bullet \quad x_{(2p-1)} \\
\bullet \quad x_{(2p)}
\end{array}
\]

viewing (15) as a function of \( u_p \),

\[
\begin{array}{c}
\bullet \quad u_q \\
\bullet \quad x_{(2q-1)} \\
\bullet \quad x_{(2q)}
\end{array}
\]

viewing (15) as a function of \( u_q \),
3. For each pair \( p, q \in \{1, 2, \ldots, N\} \) with \( p \neq q \), if \( \Gamma_q \) nests \( \Gamma_p \) then by pressing these contours so they are sufficiently close to the real line, we ensure that

\[
\arg(u_p - u_q) \in (-\pi, \pi/4) \cup (3\pi/4, \pi].
\]  

By item (iii)c, the function \( f(z) := z^{8/\kappa} \) has a branch cut along the negative real axis (for generic \( \kappa \)). The following function, identical to \( f(z) \) except for a difference in phase, has its branch cut along the positive imaginary axis:

\[
\Delta(z) := z^{8/\kappa} \begin{cases} 
\frac{e^{-2i\pi(8/\kappa)}}{\kappa}, & \pi/2 < \arg(z) \leq \pi, \\
1, & -\pi < \arg(z) \leq \pi/2.
\end{cases}
\]  

In light of (18), we avoid the branch cut along the negative real axis by replacing \( \Delta(u_p - u_q) \) with \( \Delta - \kappa u_p \). Altogether, if \( \Gamma_q \) nests \( \Gamma_p \), then we use the following factors in the integrand of (11):

\[
(u_p - x_{(2q-1)})^{-4i/\kappa}(x_{(2q)} - u_p)^{-4i/\kappa} \\
(x_{(2p-1)} - u_q)^{-4i/\kappa}(x_{(2p)} - u_q)^{-4i/\kappa} \Delta(u_p - u_q).
\]  

Because \( u_p - u_q \) is bounded away from the positive imaginary axis (18), using \( \Delta \) instead of \( f \) ensures that (20) is a continuous function of \( (u_p, u_q) \). The branch cuts (drawn as zig-zags) of (20) are arranged as follows:

viewing (20) as a function of \( u_p \).

viewing (20) as a function of \( u_q \).

(Here, we show \( \Gamma_m \) as a simple curve as in item (iii)d. If \( \kappa \leq 4 \), then we replace \( \Gamma_m \) with the Pochhammer contour (figure 8) of item (iii)a.)

Then the integrand of (11) equals the product of all factors in (12) with \( m \in \{1, 2, \ldots, N\} \), multiplied by the product of all factors in (15) with \( p, q \in \{1, 2, \ldots, N\} \) such that \( p \neq q \) and \( \Gamma_q \) is to the left of \( \Gamma_p \), multiplied by the product of all factors in (20) with \( p, q \in \{1, 2, \ldots, N\} \) such that \( p \neq q \) and \( \Gamma_q \) nests \( \Gamma_p \). A straightforward adaptation of the work in appendix B of [38] shows that \( F_0 \) (11) is real-valued. Alternatively, we may use the formula (C.2) for \( F_0 \) given in appendix C. Although the latter appears more complicated, it is really shorter and has one less integration contour. In appendix C, we show that these formulas are equivalent.

(iv) For each \( \varsigma, \vartheta \in \{1, 2, \ldots, C_N\} \), we draw the \( \varsigma \)th connectivity in the upper half-plane, draw the \( \vartheta \)th connectivity in the lower half-plane, delete the real axis, and let \( I_{\varsigma, \vartheta} \) be the number of loops left (figure 11 shows the case \( N = 2 \)).

(v) We let \( M_0(\nu) \) be the \( C_N \times C_N \) meander matrix [82–85] whose \( (\varsigma, \vartheta) \)th element is \( \nu^{I_{\varsigma, \vartheta}} \). If \( \nu \neq 4qlq' \) for any coprime pair of positive integers \( q \) and \( q' \) such that \( q' > 1 \) and \( q \leq N + 1 \), then \( \det(M_0 \circ n)(\kappa) \neq 0 \), and we solve
for the $C_N$ unknown functions $\Pi_i$, called ‘crossing weights’. (The notation $M_N \circ n$ means composition of the real-valued function $n$ with the matrix-valued function $M_N$, i.e. $(M_N \circ n)(\kappa) = M_N(n(\kappa))$.)

If $\kappa = 4q_lq'$, then the system is not invertible. Instead, we replace $\kappa$ with $\kappa \approx \kappa_0$, solve the system, and send $\kappa_0 \to \kappa$ to find the crossing weights for this $\kappa$, or we use $[40]$.)

(vi) **Formula for the crossing probability**: As we will derive in section 5, the following function gives the probability of the $\varsigma$th crossing event conditioned on the $\vartheta$th BC event for models described below:

$$P_{\varsigma|\vartheta}(\kappa | x_1, x_2, \ldots, x_{2N}) = n(\kappa)^{\Pi_{x_1, x_2, \ldots, x_{2N}}}$$

(24)

(a) *Q*-state random cluster model: With $Q \in \{1, 2, 3, 4\}$, we suppose that the continuum limit of a critical random cluster model on a lattice fills the upper half-plane and that the boundary segments $(x_j, x_{j+1})$ with $j$ odd exhibit the fixed BC (but with the state of each fixed-BC segment unspecified). With

$$Q = 4 \cos^2(4\pi/\kappa), \quad \kappa \in [4, 8], \quad Q \in \{1, 2, 3, 4\}$$

$$\Rightarrow \begin{cases} Q = 1 & \Rightarrow \kappa = 6, \\ Q = 2 & \Rightarrow \kappa = 16/3, \\ Q = 3 & \Rightarrow \kappa = 24/5, \\ Q = 4 & \Rightarrow \kappa = 4, \end{cases}$$

(25)

the formula (24) gives the probability that the FK clusters connect the fixed-BC segments together in such a way that the cluster boundaries join the points $x_1, x_2, \ldots, x_{2N}$ pairwise in the $\varsigma$th connectivity, conditioned on the event that any two fixed-BC segments joined together by an arc in the $\vartheta$th connectivity exhibit the same state.

(b) *Q*-state Potts model: With $Q \in \{2, 3, 4\}$, we suppose that the continuum limit of a critical Potts model on a lattice fills the upper half-plane and that the boundary segments $(x_j, x_{j+1})$ with $j$ odd exhibit the first spin state while sites in all other boundary segments sample all but the first spin state uniformly. With

$$Q = 4 \cos^2(\pi \kappa/4), \quad \kappa \in (0, 4], \quad Q \in \{2, 3, 4\}$$

$$\Rightarrow \begin{cases} Q = 2 & \Rightarrow \kappa = 3, \\ Q = 3 & \Rightarrow \kappa = 10/3, \\ Q = 4 & \Rightarrow \kappa = 4, \end{cases}$$

(26)

the formula (24) with $d = 1$ gives the probability that the spin clusters connect the fixed-BC segments together in such a way that the cluster boundaries join the points $x_1, x_2, \ldots, x_{2N}$ pairwise in the $\varsigma$th connectivity, conditioned on the BC event described. (Having all fixed boundary segments exhibit the same state corresponds to $d = 1$.)

(c) Loop-gas model: With $n \in (0, 2]$, we suppose that the continuum limit of a dense-phase or dilute-phase loop-gas model on a lattice fills the upper half-plane. In addition,
boundary loops’ exit and then re-enter the upper half-plane through $x_1, x_2, \ldots, x_{2N}$, with exactly one loop passing once through each point. With

$$n(\kappa) = -2 \cos(4\pi/\kappa), \quad \begin{cases} \kappa \in (8/3, 4], & \text{dilute phase,} \\ \kappa \in (4, 8), & \text{dense phase,} \end{cases} \quad (27)$$

formula (24) gives the probability that the parts of the boundary loops in the upper half-plane (‘interior arcs’) connect the points $x_1, x_2, \ldots, x_{2N}$ pairwise in the $\varsigma$th connectivity, conditioned on the event that the parts of the boundary loops inside the lower-half plane (‘exterior arcs’) join these points pairwise in the $\varsigma$th connectivity. The formula (24) for crossing probability is Möbius invariant in the sense that if $f$ is a conformal bijection of the half-plane onto itself such that $x'_j := f(x_j)$, then

$$P_{\varsigma | \theta}(\kappa | x'_1, x'_2, \ldots, x'_{2N}) = P_{\varsigma | \theta}(\kappa | x_1, x_2, \ldots, x_{2N}). \quad (28)$$

(If $f$ is not conformal, then this invariance is generally violated.) Hence, (24) depends only on the $2N - 3$ independent cross-ratios that we may form from $x_1, x_2, \ldots, x_{2N}$. For example, we may choose these cross-ratios to be

$$\lambda_j := \frac{(x_{j+1} - x_j)(x_{2N} - x_{2N-1})}{(x_{2N-1} - x_j)(x_{2N} - x_{j+1})}, \quad j \in \{1, 2, \ldots, 2N - 3\}. \quad (29)$$

(vii) If $f$ is a conformal map from the upper half-plane onto a Jordan domain $\mathcal{D}$ (such as a polygon) with $w_j := f(x_j)$, and if the same statistical mechanics model lives on the same lattice in $\mathcal{D}$ as in item vi above, then similar to (28),

$$P_{\varsigma | \theta}^\mathcal{D}(\kappa | w_1, w_2, \ldots, w_{2N}) = P_{\varsigma | \theta}(\kappa | x_1, x_2, \ldots, x_{2N}) \quad (30)$$
gives the crossing probability for the above model in $\mathcal{D}$, with the $\theta$th BC or boundary-loop condition involving $x_j$ sent to the same condition involving $w_j$. In this sense, the crossing probability is conformally invariant.

3. Random cluster model and Potts model with boundary conditions

Now we begin to study specific models. This, combined with the continuum limit, will eventually lead to our main result (24). In this section, we give formal expressions for various partition functions of the random cluster (resp. Potts) model on a lattice inside a $2N$-sided polygon $\mathcal{P}$, summing exclusively over the BC events and/or crossing events in item i (resp. item ii) of the introduction section 1. We assign BCs exclusively to the sites and/or bonds adjacent to the sides of $\mathcal{P}$, as illustrated in figure 12. (We note that lattice sites and bonds adjacent to a side
of $P$ approach that side in the continuum limit.) Anticipating the continuum limit, we regard the lattice spacing $a$ as very small compared to the lengths of these sides. As we discuss in section 1.1, universality obviates the need to specify the lattice in the continuum limit, but for clarity, our figures use either the square lattice or the triangular lattice.

### 3.1. Random cluster model and Potts model with free boundary conditions

Among the random cluster models, (Bernoulli) bond percolation on a lattice is the simplest. Here, the bonds between nearest-neighbor sites (i.e. two sites separated by exactly one lattice space) are activated, each with probability $p$ and independently of the other bonds (figure 13).

The bond percolation partition function is

$$X_P(p, Q = 1) = \sum_B p^{N_b} (1 - p)^{N_b - N_j} = 1,$$

where $B$ is the entire sample space of the model (i.e. the collection of all possible nearest-neighbor bond configurations on the lattice inside $P$ (figure 13)), and

$$N_b = \{\text{total number of bonds between nearest-neighbor sites in the entire system}\},$$

$$N_j = \{\text{total number of activated bonds in sample } \{j\} \in B\}.$$ (We include the $Q = 1$ in (31) to conform with notation for the $Q$-state random cluster model partition function.)
The \textit{Q-state random cluster model} generalizes bond percolation. In this model, all bond clusters in every percolation sample are randomly assigned, uniformly and independently, one of \(Q\) possible colors (figure 13). As mentioned in the introduction section 1, we call these colored bond clusters ‘FK clusters’. The partition function for this model is

\[
X^P(p, Q) = \sum_{\mathcal{B}} p^{N_1}(1 - p)^{N_2} Q^{N_3}, \quad \text{where}
\]

\[
N_k = \begin{cases} 
\text{number of activated bond clusters in sample } \{\beta\} \\
\text{in } \mathcal{B} \text{ (including size-zero clusters, i.e. isolated lattice sites)}
\end{cases}
\]

The new factor of \(Q^{N_3}\) that is absent from (31) uniformly distributes the \(Q\) colors to the bond clusters in every sample of \(\mathcal{B}\), including to size-zero clusters (or isolated lattice sites). We recover bond percolation in the special case \(Q = 1\).

The random cluster model is closely related to the Potts model \cite{7} on the same lattice. In the \(Q\)-state Potts model, we randomly assign any one of \(Q\) available values, called ‘spins’, to each lattice site (figure 13). In a given spin configuration (i.e. sample) \(\{\sigma\}\), each pair of lattice sites \(i\) and \(j\), with respective spins \(\sigma_i\) and \(\sigma_j\), contribute energy

\[
E_{ij} = \begin{cases} 
-J \delta_{i,j}, & \text{sites } i \text{ and } j \text{ are nearest-neighbors}, \\
0, & \text{sites } i \text{ and } j \text{ are not nearest-neighbors},
\end{cases}
\]

where \(\sigma_i, \sigma_j \in \{e^{2\pi i \theta}/Q \mid \theta \in [1, 2, \ldots, Q]\}\),

to the configuration, where \(J \in \mathbb{R}\) is the spin-coupling parameter.

The \textit{Q-state Potts model} weights each spin configuration by the Boltzmann distribution, with the total energy for its Hamiltonian. Then the Potts model partition function is

\[
Y^P(K, Q) = \sum_S e^{-\beta \sum_{\langle ij \rangle} E_{ij}}, \quad \text{where } \beta > 0.
\]

\[
= \sum_S \prod_{\langle ij \rangle} (1 + (e^K - 1) \delta_{i,j}), \quad \text{where } K := \beta J, \text{ and}
\]

where \(S\) is the collection of all spin configurations, \(\beta > 0\) is the (inverse) temperature, and the sum \(\sum_{\langle ij \rangle}\) and product \(\prod_{\langle ij \rangle}\) are over all pairs \(\langle ij \rangle\) of nearest-neighbor sites \(i\) and \(j\). After expanding the product in (36), we find \cite{7}

\[
Y^P(K, Q) = (1 - p)^{N_1} X^P(p, Q), \quad \text{where } e^K = 1/(1 - p).
\]

This identifies the ferromagnetic \((J > 0, \text{ so } p \in (0, 1))\) \(Q\)-state Potts model with the \(Q\)-state random cluster model.

For every \(Q \in \{1, 2, 3, 4\}\), the ferromagnetic \(Q\)-state Potts model undergoes a second-order phase transition at a critical temperature \(K_c\). This critical behavior manifests itself in many ways. For example, on an infinite lattice, the typical spin cluster size (i.e. collection of sites in a common spin state and such that the bonds between them form a connected cluster) is finite for \(K < K_c\) but becomes infinite at \(K = K_c\) \cite{10}. Alternatively, for the \(Q = 2\) Potts (i.e. Ising) model inside a Jordan domain \(\mathcal{D}\) with all boundary spins fixed to the \(\sigma = +1\) state, the magnetization (i.e. mean spin) vanishes for \(K \leq K_c\), but is positive for \(K > K_c\) \cite{6, 94, 95}. Explicit formulas for the critical temperature as a function of \(Q\) are known for the square lattice \cite{96} and the triangular lattice \cite{97}:
From (37), we infer that the FK-cluster size in the $Q$-state random cluster model also undergoes a phase transition at critical probability $p_c$ related to $K_c$ via (37). For the square lattice [96] and the triangular lattice [97], these are

$$e^{K_{e}^{\text{spr}}} = 1/(1 - p_{e}^{\text{spr}}) \implies p_{e}^{\text{spr}} = \frac{\sqrt{Q}}{1 + \sqrt{Q}},$$

$$e^{K_{e}^{\text{tri}}} = 1/(1 - p_{e}^{\text{tri}}) \implies p_{e}^{\text{tri}} = \left[1 + \frac{2}{\sqrt{Q}} \cos \left(\frac{1}{3} \arctan \left(\frac{4}{Q} - 1\right)\right)\right]^{-1}. \quad (40)$$

In the partition functions (31), (33), (35) encountered so far, the sums are not restricted to events with a BC imposed on lattice sites adjacent to a boundary. Because we call this lack of a BC constraint the ‘free BC’ in item i of the introduction section 1, we naturally call the partition functions of this section free BC partition functions.

### 3.2. Random cluster model with free/fixed boundary conditions

In this section, we give formal expressions for RCB partition functions, or random cluster model partition functions that sum exclusively over an RCB event (as defined in item i in the introduction section 1) inside a $2N$-sided polygon $\mathcal{P}$. A natural RCB event is a color scheme $c = (c_1, c_2, \ldots, c_N)$. This event comprises all samples where the $i$th fixed side has color $c_i \in \{1, 2, \ldots, Q\}$ for each $i \in \{1, 2, \ldots, N\}$ (figure 14). There are $Q^N$ different color schemes, and color schemes are disjoint. We define a color scheme partition function $W_p^{\text{spr}}(p, Q)$ to be an RCB partition function that sums exclusively over a given color scheme $c$. Any random cluster model RCB event is a union of color schemes, so any RCB partition function is a sum of color scheme partition functions. Thus, it is ostensibly fruitful to study the latter. However, in this article, we instead work with a certain subset of RCB partition functions that we call ‘basic’ partition functions. In appendix B, we show that for $Q \in \{1, 2, 3\}$, any color scheme partition function equals a linear combination of basic partition functions. Furthermore, in section 4.2, we show that basic partition functions are more tractable to use in the continuum limit. For this reason, we work with basic partition functions.
The basic partition functions are examples of mutual wiring partition functions, or RCB partition functions that sum exclusively over a mutual wiring event. A mutual wiring event is an RCB event where every fixed side of $P$ belongs to a collection of fixed sides that are mutually wired, i.e. constrained to have the same color (figure 15). In such an event, the actual color of a collection of mutually wired sides is not specified; it can be anything. (If one such collection has only one fixed side of $P$, then that side is not constrained to exhibit the color of any other fixed side, and we say that it is independently wired.) A mutual wiring event $E$ samples (i.e. is a union of) all color schemes $c$ consistent with it. By 'consistent', we mean $c \leftrightarrow E$, where by definition,

$$c = (c_1, c_2, …, c_N) \leftrightarrow E \quad \Rightarrow \quad \text{if the } i^{th} \text{ and } j^{th} \text{ fixed sides of } P \text{ are mutually wired together in } E, \text{ then } c_i = c_j.$$  

(41)

Thus, the mutual wiring partition function $X^E_P(p, Q)$ summing exclusively over $E$ equals a sum of all color scheme partition functions consistent with $E$:  

$$X^E_P(p, Q) = \sum_{c \leftrightarrow E} W^E_{c}(p, Q).$$  

(42)

Writing $N_c(E)$ for the total number of independent clusters in a given sample, that is the number of clusters after identifying all boundary clusters that touch a common collection of mutually wired sides as one, (33) becomes  

$$X^p_E(p, Q) = \sum_{B} p^{N_B} (1-p)^{N_B-N_c(E)} Q^{N_c(E)}.$$  

(43)

In this article, we consider only a special subset of mutual wiring events $RCB_1$, $RCB_2$, …, $RCB_{C_B}$, called basic events. To define them, we consider all of the distinct ways (up to homotopy) to join the vertices of $P$ pairwise with $N$ nonintersecting planar curves (exterior...
arcs), drawn outside $\mathcal{P}$. We call each possibility an \textit{exterior arc connectivity}, and figure 16 shows that they bijectively correspond to the $C_N(3)$ crossing patterns in $\mathcal{P}$. (We interpret exterior arc connectivities as crossing events in the exterior of $\mathcal{P}$. This analogy is not obvious, but the work that we do below justifies it.) After enumerating these connectivities one through $C_N$, we let the basic event $\vartheta_{RCB}$ be the mutual wiring event where any two fixed sides joined by an exterior arc in the $\vartheta$th connectivity are mutually wired (figure 15). Then the basic partition functions are

$$X_\vartheta^P := X_{RCB}^P, \quad \text{for all } \vartheta \in \{1, 2, \ldots, C_N\}. \quad (44)$$

We note that $\{RCB_1, RCB_2, \ldots, RCB_{C_N}\}$ comprises all mutual wiring events only if $N < 3$. Indeed, for $N = 4$, no exterior arc connectivity generates the mutual wiring event in which the octagon’s top and bottom sides are mutually wired together while the octagon’s left and right sides are also mutually wired together, independent of the former. Therefore, we do not include this mutual wiring event. For larger $N$, even more possibilities are not included.

\section*{3.3. Random cluster model crossing events with free/fixed boundary conditions}

In this section, we give formal expressions for the probabilities of random cluster model crossing events conditioned on the mutual wiring events $RCB_1, RCB_2, \ldots, RCB_{C_N}$ defined in the previous section 3.2. In order to do this, we restrict the partition function sums of the previous

\[\text{Figure 16. Crossing patterns (top) bijectively correspond to interior arc connectivities (middle), which bijectively correspond to exterior arc connectivities (bottom). The bijection respects our enumeration, one through $C_N$, of these three entities.}\]
We enumerate the $CN$ crossing events (and corresponding arc connectivities (figure 16)) such that a single (resp. no) boundary cluster connects all (resp. any two) fixed sides of $P$ in the first (resp. $CN$th) event (figure 17).

We let $\varsigma_{RCX}$ denote the $\varsigma$th FK-cluster crossing event, and we represent it by the $\varsigma$th interior arc connectivity (i.e. the $\varsigma$th exterior arc connectivity, but with interior arcs drawn inside $P$) (figures 16 and 18).

We let $\vartheta_{RCB}$ denote the $\vartheta$th random cluster model mutual wiring event (fixed sides joined by an arc in the $\vartheta$th connectivity are mutually wired, figure 15) and represent it by the $\vartheta$th exterior arc connectivity (figures 16 and 18).

We let $(\varsigma\vartheta)p_{XpQ}$ denote the partition function for the $Q$-state random cluster model on a lattice in $P$, summing exclusively over the intersection $\varsigma_{RCX}\cap \vartheta_{RCB}$ of the $\varsigma$th crossing event with the $\vartheta$th RCB event.

Although section 3.2 already has item iii, we reiterate it here for easy reference. We infer from item i with figure 17 that $RCB_1$ (resp. $RCX_1$) in which all fixed sides are mutually wired together (resp. a single cluster connects all fixed sides). The $C_N$th exterior (resp. interior) arc connectivity corresponds to the event $RCB_{C_N}$ (resp. $RCX_{C_N}$) in which all fixed sides are independently wired (resp. no cluster connects any two fixed sides).

Turning to crossing probabilities of the critical random cluster model, we write a formal expression for the probability of the $\varsigma$th FK-cluster crossing event $\varsigma_{RCX}$ conditioned on the $\vartheta$th RCB event $\vartheta_{RCB}$. First defining

$$ B_\varsigma = \{ \varsigma$th percolation (i.e. $Q = 1$ FK cluster) crossing event $\} \subseteq B, \quad (45) $$

and then replacing $B \mapsto B_\varsigma$ and $\mathcal{E} \mapsto RCB_\varsigma$ in (33), we obtain these formal expressions for $X_{\varsigma,\vartheta}^P$ and $X_{\vartheta}^P$: 

$$ X_{\varsigma,\vartheta}^P(p, Q) = \sum_{B_\varsigma} p\gamma_{\varsigma}(1 - p)\gamma_{\vartheta}Q^{N_{\varsigma}(RCB_{\vartheta})}, \quad (46) $$

Figure 17. The first exterior (resp. interior) arc connectivity corresponds to the event $RCB_1$ (resp. $RCX_1$) in which all fixed sides are mutually wired together (resp. a single cluster connects all fixed sides). The $C_N$th exterior (resp. interior) arc connectivity corresponds to the event $RCB_{C_N}$ (resp. $RCX_{C_N}$) in which all fixed sides are independently wired (resp. no cluster connects any two fixed sides).
Then the probability of the $\varsigma$th FK-cluster crossing event $\text{RC}_\varsigma$ conditioned on the $\vartheta$th RCB event $\text{RCB}_\vartheta$ is given by the ratio of these partition functions (abbreviating ‘random cluster’ with ‘RC’):

$$P_{\varsigma\vartheta}(p, Q) := \mathbb{P}_{\text{RC}}(\text{RC}_\varsigma \mid \text{RCB}_\vartheta) = \frac{X^\text{P}_{\vartheta}(p, Q)}{X^\text{P}_\varsigma(p, Q)}.$$  \hfill (47)

Then the probability of the $\varsigma$th FK-cluster crossing event $\text{RC}_\varsigma$ conditioned on the $\vartheta$th RCB event $\text{RCB}_\vartheta$ is given by the ratio of these partition functions (abbreviating ‘random cluster’ with ‘RC’):

$$P_{\varsigma\vartheta}(p, Q) := \mathbb{P}_{\text{RC}}(\text{RC}_\varsigma \mid \text{RCB}_\vartheta) = \frac{X^\text{P}_{\vartheta}(p, Q)}{X^\text{P}_\varsigma(p, Q)}. \hfill (48)$$

Our primary goal in this article is to find a formula for the continuum limit of (48) at the critical point $p = p_c$. This poses two tasks: the first is to determine the numerator and denominator of (48), which by (46), (47), amounts to determining $N(\text{RCB}_\vartheta)$. The second task, taken up in section 5, is to determine the limit of (48) as the lattice spacing $a$ vanishes.

Taking up the first task, we let $N^\text{free}_c$ denote the number of free clusters, or clusters of activated bonds in configuration $\{\beta\} \in \mathcal{B}_c$ that are not boundary clusters. As such, no bond in a free cluster touches a fixed side of $P$. (These include bulk clusters, which do not touch any side of $P$.) Furthermore, we let $\varsigma C_{\vartheta}$ denote the number of boundary clusters in $\{\beta\}$, with all boundary clusters attached to a common collection of mutually wired sides identified as one. Then

$$N(\text{RCB}_\vartheta) = C_{\varsigma\vartheta} + N^\text{free}_c \quad \text{for each } \{\beta\} \in \mathcal{B}_c. \hfill (49)$$

Because $C_{\varsigma\vartheta}$ is the same for all bond configurations in the $\varsigma$th crossing event, (46) factors into

$$X^\text{P}_{\varsigma\vartheta}(p, Q) = Q^{C_{\varsigma\vartheta}} \sum_{\mathcal{B}_c} p^{N(\beta)} (1 - p)^{N(\beta) - N(\text{RCB}_\vartheta)}.$$  \hfill (50)

Although the prefactor $Q^{C_{\varsigma\vartheta}}$ is not physically relevant to this partition function (50), we need it in order to compute the sum (47), which appears in the denominator of (48).

Next, we determine a formula for $C_{\varsigma\vartheta}$ in terms of the number $l_{\varsigma\vartheta}$ of loops in the $(\varsigma, \vartheta)$th product diagram, created by superimposing the $\varsigma$th interior arc connectivity and $\vartheta$th exterior arc connectivity on the same polygon, joining the interior arcs to the exterior arcs at the vertices of the polygon to form loops, and deleting the polygon, leaving only the loops (figure 19). The formula that we find is

$$C_{\varsigma\vartheta} = (l_{\varsigma\vartheta} - l_{\vartheta} + l_{\varsigma\vartheta} + 1)/2. \hfill (51)$$
To derive (51), we identify the $(s, \vartheta)$th product diagram with a graph $G_{s,\vartheta}$, in which the quantity $C_{s,\vartheta}$ naturally arises. This identification goes as follows:

(i) We enumerate the fixed sides of the $2N$-sided polygon $P$ one through $N$ in counterclockwise order. (By convention, we take the lowest side of $P$ to be the first fixed side.)
(ii) We identify each fixed side of $P$ with a vertex of the graph $G_{s,\vartheta}$ under construction, and we enumerate these vertices so the $i$th vertex corresponds with the $i$th fixed side.
(iii) If $\{s_1, s_2, \ldots, s_k\} \subset \mathbb{Z}^+$ is the set of labels for all fixed sides joined by a common boundary cluster in $RCX_s$ (with $1 < k \leq N$), then we join the $s_i$th vertex and the $s_{i+1}$th vertex of $G_{s,\vartheta}$ with an edge for each $i \in \{1, 2, \ldots, k-1\}$.
(iv) If $\{s_1, s_2, \ldots, s_k\} \subset \mathbb{Z}^+$ is the set of labels for all fixed sides of a collection of mutually wired sides in $RCB_s$ (with $1 < k \leq N$), then we join the $s_i$th vertex and the $s_{i+1}$th vertex of $G_{s,\vartheta}$ with an edge for each $i \in \{1, 2, \ldots, k-1\}$.

(We note that an edge joins the first and $N$th vertices of $G_{s,\vartheta}$ via item iii (resp. item iv) if and only if a boundary cluster in $RCX_s$ joins only these two fixed sides (resp. these sides are mutually wired to only each other in $RCB_s$).) Figure 20 gives three examples of product diagrams in its top row and their associated graphs beneath. Now we let
It is evident from steps iii and iv that the number \( C_{c,\theta} \) of components in \( \varpi \) equals the number of independent boundary clusters in the event \( \cap \varpi \), also denoted by \( C_{c,\theta} \). Euler’s formula relates all of these quantities together as

\[
C_{c,\theta} - \mathcal{I}_{c,\theta} + \mathcal{E}_{c,\theta} - \mathcal{V}_{c,\theta} = 0, \quad \text{where} \quad \mathcal{E}_{c,\theta} := \mathcal{E}_{c} + \mathcal{E}_{\varpi}.
\]  

(52)

Figure 20 shows that the loops in the \((\varsigma, \theta)\)th product diagram trace all of the internal faces and surround all of the connected components of \( \varpi \) (seen as embedded on the Riemann sphere), with one loop per face and one loop per component. Also, it is obvious that the number \( \mathcal{V}_{c,\theta} \) of vertices in \( \varpi \) is \( N \). Thus,

\[
I_{c,\theta} = C_{c,\theta} + \mathcal{I}_{c,\theta}, \quad \mathcal{V}_{c,\theta} = N.
\]  

(53)

To derive formula (51) for \( C_{c,\theta} \) from (52), (53) we consider two extreme cases. In the first case, we take \( \varsigma = 1 \). According to item i in section 3.3, a single boundary cluster connects all fixed sides of \( \mathcal{P} \) in the first crossing event \( \text{RCX}_i \). The left column of figure 21 depicts this event, with the \((1, \theta)\)th product diagram on top of it and its associated graph \( \varpi_1 \) beneath. This graph has one component with \( N - 1 \) internal edges joining the vertices. Hence,
\( G_{i, \vartheta} = 1, \quad \mathcal{E}_1^i = N - 1. \)  

(54)

In the second case, we take \( \vartheta = C_N \). According to items i and iii in section 3.3, all fixed sides of \( \mathcal{P} \) are independently wired in the event \( RCB_{C_N} \). The right column of figure 21 depicts this event, with the (\( \varsigma, C_N \))th product diagram on top of it and its associated graph \( \mathcal{G}_{\varsigma, C_N} \) beneath. This graph has no external edges. Furthermore, it has no internal faces because any face is formed by at least one external edge meeting one internal edge. Hence,

\[ T_{\varsigma, C_N} = 0, \quad \mathcal{E}_{C_N}^\varsigma = 0. \]  

(55)

Combining (52)–(55) gives the expression (51) for \( C_{i, \vartheta} \), relating the effective number of independent boundary clusters in the event \( \cap_{\varsigma} RCB_{\varsigma} \) with the number \( l_{\varsigma, \vartheta} \) of loops in the (\( \varsigma, \vartheta \))th product diagram (figure 19). In section 4, we use this connection to interpret the partition functions (47), (50) as instances of loop-gas model partition functions.

3.4. Potts model crossing events with fluctuating/fixed boundary conditions

Although it seems natural to consider free/fixed side-alternating BC events in the Potts model, determining crossing probabilities conditioned on these events is beyond the scope of the methods used in this article. Instead, our methods (section 5) are suited to PMB events for the Potts model, as defined in item ii in the introduction section 1. These are side-alternating BC events in which every fixed side exhibits the first spin state and the remaining sides exhibit the “fluctuating BC”, where sites uniformly sample all but the first spin state \([46]\). (See the left illustration of figure 29.) In parallel with items i–v at the beginning of section 3.3, we define the following for the Potts model:

(i) For the Potts model, crossing events are now events in which spin clusters join the fixed sides of \( \mathcal{P} \) in some particular crossing pattern. We enumerate these events as in item i in section 3.3.

(ii) We let \( PMX_{\varsigma} \) denote the \( \varsigma \)th spin-cluster crossing event, and we represent it by the \( \varsigma \)th interior arc connectivity (figures 16).

(iii) We let \( PMB \) denote the Potts model fluctuating/fixed side-alternating BC event (we note that, as defined, there is only one), and we represent it by the first exterior arc connectivity (figure 17).

(iv) We let \( Y_{i}^{P}(K,Q) \) denote the partition function for the \( Q \)-state Potts model on a lattice inside \( \mathcal{P} \), summing exclusively over the intersection \( \cap_{\varsigma} PMX_{\varsigma} \cap PMB \) of the \( \varsigma \)th crossing event with the PMB event.

(v) We let \( Y_{i}^{P}(K,Q) \) denote the partition function for the \( Q \)-state Potts model on a lattice inside \( \mathcal{P} \), summing exclusively over the PMB event. Hence, \( Y_{i}^{P} = Y_{1,1}^{P} + Y_{2,1}^{P} + \cdots + Y_{C_N,1}^{P} \).

(The subscript 1 in \( Y_{i}^{P}(K,Q) \) and \( Y_{i}^{P}(K,Q) \) of items iv and v is natural to use, as will become evident in section 4.) In terms of these quantities, the probability of the \( \varsigma \)th spin-cluster crossing event conditioned on the PMB event is

\[ P_{i,1}^{P}(K,Q) := \mathbb{P}_{\text{Potts}}(PMX_{\varsigma} \mid PMB) = \frac{Y_{i,1}^{P}(K,Q)}{Y_{i}^{P}(K,Q)}. \]  

(56)

In section 5, we determine an explicit formula for the continuum limit of the crossing probabilities (48), (56) encountered so far. For the former, we condition the random cluster model on an RCB event, and for the latter, we condition the Potts model on a PMB event. As mentioned,
the Potts model with RCB conditioning is outside the scope of our methods. The same is true for the random cluster model with PMB conditioning.

4. Convergence to loop-gas models

In this section, we show that partition functions of the random cluster model and the Potts model that we encountered in sections 3.2, 3.3 and 3.4 are asymptotic to partition functions of certain loop-gas models [11, 93, 98] in the continuum limit. In section 5, these identifications lead us, finally, to the formula (24) for the crossing probabilities (48), (56) in the continuum limit. (Section 2 explains how to make this formula completely explicit for the random cluster model and the Potts model. Sections 6 and 7 give these explicit formulas if the polygon is a rectangle or a hexagon respectively.)

4.1. Definition of loop-gas models

We begin with a definition of a loop-gas model. The sample space of a loop-gas model on a given graph inside is the collection of all sets of non-crossing and non-self-crossing loops that trace the graph’s edges [12, 99]. In this section, we assume that the graph is a square or hexagonal lattice, with edges being bonds between nearest-neighbor lattice sites. Figure 22 shows the five scenarios in which loops may interact with a site of the square lattice:

(i) No loop passes through the lattice site.
(ii) A loop passes through the site without turning.
(iii) A loop touches the site, turns, and never returns to the site.
(iv) A loop touches the site, turns, and eventually returns to the site.
(v) Two distinct loops touch a common site and turn.

In a typical loop-gas model, we endow the kth scenario with some weight , and we endow a sample collection of loops with weight

where

(57)

(One may consider loop-gas models with , but we do not use them here.) Because scenarios iv and v appear identical, (i.e. the information contained within the dashed box of either scenario in figure 22 is the same), we typically set . If we also set and , then (57) simplifies to

(58)

Figure 22 also shows the two different ways in which the loops may interact with a single site on the hexagonal lattice. Because only scenarios i and iii occur, we set and in the expression (57) for the weight of the loop sample . Setting and , this expression again reduces to (58). The partition functions for these models are
\[ Z_f(x, n) = \sum_{\omega}^L \omega(\gamma), \quad \text{where } L := \{ \text{the sample space of the loop-gas model} \}. \quad (59) \]

The subscript \( f \) (for ‘free’) means that we sum over the entire sample space, without imposing any BCs such as those described below. (We mention that this loop-gas model (58), (59) is closely related to the O(\( n \)) model \([11]\) of statistical physics.) In what follows, we identify the loop-gas model partition function with the random cluster model and Potts model partition functions previously encountered in sections 3.2 and 3.4.

The loop-gas model (58) exhibits critical behavior in its continuum limit. This critical behavior is well understood on the hexagonal lattice \([13, 51, 93, 100]\). Under the renormalization group flow induced by shrinking the lattice spacing \( a \), the temperature \( x \) of (58) flows to either zero if it is less than some critical temperatures \( x_c \) and \( n \in [-2, 2] \), or to another value \( x_\sim > x_c \) if it is greater than \( x_c \) and \( n \in [0, 2] \). On the hexagonal lattice, we have [51]

\[ x_c(n) = (2 + \sqrt{2 - n})^{-1/2}, \quad x_\sim(n) = (2 - \sqrt{2 - n})^{-1/2}. \quad (60) \]

If \( x < x_c \) and \( n \in [0, 2] \), then as \( a \downarrow 0 \), the model strongly favors small-diameter loops over large-diameter loops, and its continuum limit contains only infinitesimal loops. However, if \( x > x_c \), then the continuum limit contains loops of all sizes. Moreover, the fraction of all lattice sites that a typical loop visits does not vanish as \( a \downarrow 0 \). For this reason, we call the interval \( x > x_c \), the dense phase. In the continuum limit, the loops are conjectured to have the conformally invariant law of \( \text{CLE}_\kappa \) \([101–103]\), with \( \kappa \in (4, 8) \) and \( (10) \)

\[ n = n(\kappa) := -2 \cos(4\pi/\kappa). \quad (61) \]

Finally, if \( x = x_c \) and \( n \in [-2, 2] \), then \( x \) is invariant under the renormalization group flow. As \( a \downarrow 0 \), the model still strongly favors large-diameter loops over small-diameter loops, but the fraction of all lattice sites that a typical loop visits vanishes. For this reason, we call \( x = x_c \), the dilute phase. In the continuum limit, the loops are conjectured to have the conformally invariant law of \( \text{CLE}_\kappa \) \([101–103]\), for \( \kappa \in (8/3, 4) \) and \( n \in (0, 2) \) related by (61).

A loop-gas model on a lattice inside an even-sided polygon \( P \) admits certain events that are analogous to the random cluster model RCX and RCB events in \( P \). To realize this analogy, we restrict the loop-gas sample space to all configurations such that each odd side of \( P \) (see above (3) for the enumeration of the sides) is traced from vertex to vertex by one loop, called a boundary loop (figure 23). These boundary loops necessarily join the vertices of \( P \).
pairwise through its interior in one of the $C_N$ available connectivities. If it is the $\varsigma$th connectivity, then we regard this event as the analog of the $\varsigma$th crossing event $\varsigma$RCX in the random cluster model. Furthermore, we may bisect the part of the boundary loops that follow the odd sides of $P$ and rejoin the $2N$ dangling ends with nonintersecting exterior arcs outside $P$ (figure 23). If the exterior arcs join the vertices of $P$ in the $\vartheta$th connectivity, then we regard this event as the analog of the $\vartheta$th mutual wiring event $\vartheta$RCB in the random cluster model. Thus, we define the following:

(i) For the loop-gas model, the $\varsigma$th crossing ($\varsigma$LGX) event (resp. $\vartheta$th exterior arc BC ($\vartheta$LGB) event) is the event that the boundary loops join the vertices of $P$ via the $\varsigma$th interior (resp. $\vartheta$th exterior) arc connectivity.

(ii) We let $\varsigma$LGX denote the $\varsigma$th loop-gas model crossing ($\varsigma$LGX) event, and we represent it by the $\varsigma$th interior arc connectivity (figure 16).

(iii) We let $\vartheta$LGB denote the $\vartheta$th exterior arc BC ($\vartheta$LGB) event, and we represent it by the $\vartheta$th exterior arc connectivity (figure 16).

(iv) We let $Z^{\mu}_{c,\vartheta}(x,n)$ denote the partition function (59) for the loop-gas model on a lattice inside $P$, summing exclusively over the intersection $\varsigma$LGX $\cap$ $\vartheta$LGB of the $\varsigma$th LGX event with the $\vartheta$th LGB event.

(v) We let $Z^{\mu}_{c}(x,n)$ denote the partition function (59) for the loop-gas model on a lattice inside $P$, summing exclusively over the $\vartheta$th LGB event $\vartheta$LGB. Hence, $Z^{\mu}_{c} = Z^{\mu}_{c,\vartheta} + Z^{\mu}_{c,\vartheta} + \cdots + Z^{\mu}_{c,\vartheta}$. 

Figure 24 illustrates the event $\varsigma$LGX $\cap$ $\vartheta$LGB. The number of boundary loops in this event always equals $l_{c,\vartheta}$ (figures 19 and 24). Hence, writing the total number of loops as the sum of boundary loops and the remaining free loops gives

\[ Z^{\mu}_{c} = \sum_{\varsigma=L_1,\ldots,L_N} \sum_{\vartheta=L_1,\ldots,L_N} Z^{\mu}_{c,\vartheta} + \sum_{\varsigma=L_1,\ldots,L_N} \sum_{\vartheta=L_1,\ldots,L_N} \sum_{\varsigma=L_1,\ldots,L_N} Z^{\mu}_{c,\vartheta} + \cdots + \sum_{\varsigma=L_1,\ldots,L_N} \sum_{\vartheta=L_1,\ldots,L_N} Z^{\mu}_{c,\vartheta}. \]
The rightmost expression in (62) only depends on our choice of LGB event (i.e. on $\vartheta$) via the factor $n^{\rho,\vartheta}$. The sum that multiplies this factor depends on the crossing event $\text{LG}_X$, and we interpret it as a partition function with the fugacities of all boundary loops equaling one.

4.2. Critical random cluster models in relation to loop-gas models

Turning to the critical random cluster model, our next goal is to identify each basic partition function $X_{c,\vartheta}(p, Q)$ (46) with the corresponding partition function $Z_{c,\vartheta}(x, n)$ of a loop-gas model on the square lattice. This identification is well known for the free partition function (33), (59) [104]; we adapt this argument to our purpose here. To begin, we map each bond configuration in $B_c$ (45) onto a unique loop configuration in which the perimeter of each bond cluster and dual-bond cluster is traced by one loop particular to that cluster (figure 25). (This includes size-zero clusters, which are isolated lattice sites and isolated dual lattice sites.) [104] describes this map in detail. As in section 3.2, Euler’s formula for planar graphs gives

$$2N_c + N_b = N_l + N_n,$$

where

$$N_c = \text{number of activated bond clusters},$$
$$N_b = \text{number of activated bonds},$$
$$N_l = \text{number of loops surrounding clusters},$$
$$N_n = \text{number of lattice sites}.$$

In every bond configuration, we distinguish boundary clusters, which touch the fixed sides of $P$, from the remaining free clusters, which do not, and we distinguish boundary loops, which surround these boundary clusters, from free loops, which surround the remaining free clusters (figure 25). Denoting the number of each of these as $N_c^{\text{bdry}}$, $N_c^{\text{free}}$, $N_l^{\text{bdry}}$, and $N_l^{\text{free}}$ respectively, we have

$$N_c = N_c^{\text{bdry}} + N_c^{\text{free}},$$
$$N_l = N_l^{\text{bdry}} + N_l^{\text{free}},$$
$$N_c^{\text{bdry}} = N_l^{\text{bdry}} = I_c C_b^{\text{bdry}}$$

for $\{\beta\} \in B_c$.

The last equality of (64) follows from the facts that, in the $c$th connectivity, a unique boundary loop surrounds every boundary cluster and that these loops are topologically identical to those in the $(c, C_b)$th product diagram (figure 19), where interior arcs close into boundary loops by wrapping outside and around the fixed (i.e. odd-numbered) sides of $P$. After inserting (39), (49) into the partition function formula (46), we find

$$X_{c,\vartheta}(p_c, Q) = (1 + \sqrt{Q})^{N_c} Q^{C_c} \sum_{B_c} Q^{N_{\text{free}} + N_l^2/2}.$$

The factor of $(1 + \sqrt{Q})^{N_c}$ is physically irrelevant, so we drop it. Then after inserting (63), (64) into (65) and using (51) to simplify the result, we find

$$X_{c,\vartheta}(p_c, Q) = Q^{(I_c + C_b + 1)/2} \sum_{B_c} Q^{(N_c + N_{\text{free}} + N_l^2)/2}.$$
Finally, after dropping irrelevant factors and straightening loops into zig-zag paths along bonds on the lattice medial to the square lattice in figure 25 and (66) becomes a partition function for a loop-gas model on the medial lattice:

\[ \sum \propto = \sum_{\rho}^{\nu} \rho^\nu \]

random cluster model on the square lattice,

\[ = n^L \sum_{\text{loop-gas on the medial lattice}} n^{N_{\text{loop}}} \]

The condition that \( x_1 = x_2 = x_3 = 0 \) for all bulk (i.e. interior) sites implies that only loop configurations with each bulk site visited twice by loops contributes to the sum (68) (figure 22). Because the other loop configurations do not contribute, we cannot equate (68) with \( Z_{x=1}^{\text{loop-gas on the square lattice}}(x,n) \) for \( x = 1 \). However, the temperature \( x = 1 \) is greater than the critical temperature \( x_c(n) \) of the loop-gas model on the square lattice [105], so it flows under renormalization to the dense-phase fixed point \( \tilde{x}_c(n) \) as we take the continuum limit. Because scenarios iv and v statistically dominate scenarios i–iii in the dense phase, (68) is at least asymptotic to (a multiple of) \( Z_{x=1}^{\text{loop-gas on the medial lattice}}(1,n) \) in the continuum limit.

4.3. Critical Potts model in relation to loop-gas models

Turning next to the critical Potts model on some given lattice in \( \mathcal{P} \), our next goal is to show that the PMB partition function \( Y_{\text{PMB}}^\mathcal{P}(K, Q) \) is asymptotic to (a multiple of) the partition function
for the loop-gas model on the dual lattice. This is easy to do rigorously at any temperature if \( Q = 2 \) (i.e. the Ising model), but if \( Q \in \{3, 4\} \), the arguments are not rigorous and are somewhat imprecise. The simulation results of section 7, however, strongly suggest that these arguments are in fact correct, especially if \( Q = 3 \).

We begin with the Ising model. Here, the arguments are already well known. If the Ising model lives on the triangular lattice, then the loops live on the (dual) hexagonal lattice, and they surround clusters of spin + lattice sites. Hence, spin configurations on the triangular lattice bijectively correspond to loop-gas configurations on the hexagonal lattice (figure 26).

With this bijection, it is easy to see that

\[
Y^P(K, Q = 2) \propto \sum_{\mathcal{L}} x^\rho, \quad x := \exp(-K), \rho \text{ as in (58), Ising model on the triangular lattice,}
\]

\[
= \sum_{\mathcal{L}} \omega(\gamma), \quad x_1 = x_2 = x_4 = x_8 = 1, x_3 = x, n = 1, \text{ loop-gas model on the hexagonal lattice.} \quad (69)
\]

Because (69) equates the Ising model on the triangular lattice with the loop-gas model on the hexagonal lattice, the critical points of these two models must match. And indeed, they do: setting \( Q = 2 \) in (38) and \( n = 1 \) in (60) gives

\[
K^\text{ini}_c = \log \sqrt{3} \quad \Rightarrow \quad x(K^\text{ini}_c) = \exp(-K^\text{ini}_c) = 1/\sqrt{3} = x_c(n = 1). \quad (70)
\]

Similarly, every spin configuration on a square lattice corresponds to a unique loop-gas configuration on its dual, also the square lattice, and vice versa, if we exclude scenario iv (figures 22 and 26). Thus, for the square lattice,

\[
Y^P(K, Q = 2) \propto \sum_{\mathcal{L}'} x^\rho, \quad x := \exp(-K), \rho \text{ as in (58), Ising model on the square lattice,}
\]

\[
= \sum_{\mathcal{L}} \omega(\gamma), \quad x_1 = 1, x_3 = x, x_4 = 0, x_8 = x^2, n = 1, \text{ loop-gas model on the dual square lattice,} \quad (71)
\]

where \( \mathcal{L}' \) is the subset of \( \mathcal{L} \) that excludes from \( \mathcal{L} \) all samples with a scenario iv vertex.

By contrast, equating the partition function (35) for the critical \( Q = 3 \) Potts model with that (59) for a loop-gas model in its dilute phase appears to be impossible for finite systems. To understand why, we consider a sample \( \{\sigma_i\} \) of the \( Q = 3 \) Potts model on a triangular lattice in which some given site \( P \) and its six neighbors exhibit the first spin state. From this sample, we generate two more samples \( \{\sigma_2\} \) and \( \{\sigma_3\} \) by changing the spin at \( P \) to the second and third state respectively. Then the ratio of the total weight of the latter two samples to the first sample’s weight is
\[
\frac{\prod \text{factors from interactions between all sites except } P \cdots (Q - 1)}{\prod \text{factors from interactions between all sites except } P \cdots \exp(K_{Q}^{\text{tri}})^{n}} = 0.0453874 \ldots, \text{ where } Q = 3. \tag{72}
\]

Next, we translate this scenario into the loop-gas picture. Numerical results [46] indicate that, at least in the continuum limit, the loops trace the boundaries of clusters with all sites in the first spin state. Therefore, \{σ\} corresponds to a loop configuration \{γ\} with no loop separating P from any of its neighbors, and the event that either \{σ2\} or \{σ3\} occurs corresponds to the loop configuration \{γ\} \cup \{γ0\}, where \{γ\} is the loop configuration for \{σ\} and \{γ0\} is a loop that surrounds P, separating it from its neighbors. By (58), (60), (38), the ratio of the weights of these configurations is

\[
\frac{\omega(\{γ\} \cup \{γ0\})}{\omega(\{γ\})} = nx_c(n)^{\delta} = \frac{4(1 + \sqrt{3})}{(4 + \sqrt{6} - 2\sqrt{3})^{3}} = 0.0901699 \ldots, \text{ where } n = (1 + \sqrt{3})/2. \tag{73}
\]

(\text{figure 27). (See } (75) \text{ for the value of } n \text{ in } (73).) \text{ If the identification of the critical } Q = 3 \text{ Potts model on the triangular lattice with the dilute-phase loop-gas model on the hexagonal lattice exists, then the ratios } (72) \text{ and } (73) \text{ should be equal. Because they are not, such an identification does not exist, at least in the discrete setting. (However, it is true that } (72), (73) \text{ are equal if } Q = 2 \text{ and } n = 1, \text{ so } x_c(n) = 1/\sqrt{3}, \text{ and } K_{Q}^{\text{tri}} = \log \sqrt{3} \text{ (70), as must be the case by (69).})

Because a dense-phase loop-gas representation of the critical random cluster model emerges in the continuum limit for \(Q \in \{1, 2, 3, 4\}\), the association (37) between the random cluster model and the Potts model suggests that a dilute-phase loop-gas representation of the latter, although apparently nonexistent in the discrete setting for \(Q > 2\), emerges in the continuum limit too. As evidence for this claim, we consider a bulk FK cluster on the square lattice (\text{figure 28}). Typically, this cluster has fjords entering it, and a loop on the medial lattice that surrounds this cluster visits the mouth of each fjord twice. Now, as we approach the continuum limit, the loop that surrounds the FK cluster is conjectured to have the law of CLE\(_{\kappa}\) [101–103] for some \(\kappa > 0\). Also, the fjords into the FK cluster do not vanish in this limit, so this loop must retain the self-intersection points at the fjord openings. Now, a loop in a CLE\(_{\kappa}\)
sample has self-intersection points only if $\kappa \in (4, 8)$ (and has points where the loop very nearly self-intersects at $\kappa = 4$). Indeed, this property follows from the fact that the fractal geometry of a $\text{CLE}_\kappa$ loop is locally that of an $\text{SLE}_\kappa$ curve [101–103], for which this property holds in the $\text{SLE}_\kappa$ dense phase $\kappa \in (4, 8)$ [66–68]. These facts together with (61), (67) give

\[
\sqrt{Q} = n(\kappa) = -2 \cos(4\pi/\kappa) \quad \text{with} \quad \kappa \in [4, 8)
\]

\begin{align*}
Q = 1 & \quad \Rightarrow \quad n = 1, \quad \kappa = 6, \\
Q = 2 & \quad \Rightarrow \quad n = \sqrt{2}, \quad \kappa = 16/3, \\
Q = 3 & \quad \Rightarrow \quad n = \sqrt{3}, \quad \kappa = 24/5, \\
Q = 4 & \quad \Rightarrow \quad n = 2, \quad \kappa = 4.
\end{align*}

(74)

Returning to the discrete setting, if we activate the bond at the mouth of each fjord into the FK cluster, then the loop surrounding this cluster becomes simple, and we interpret it as the boundary of the smallest Potts model spin cluster that may contain the given FK cluster. In the continuum limit, the simple loop thus created goes to the previously encountered $\text{CLE}_\kappa$ loop, but with its fjords removed (figure 28). By $\text{SLE}_\kappa$ duality [66, 67, 106–108], this latter simple loop must have the law of $\hat{\text{CLE}}_{\hat{\kappa}}$ with parameter $\hat{\kappa} = 16/\kappa \in (8/3, 4]$, where $\kappa$ is the parameter associated with the $\text{CLE}_\kappa$ loop having fjords (74) (figure 28). For the above Potts models, these parameters are

\[
\hat{\kappa} = 16/\kappa \quad \Rightarrow \quad \begin{cases} 
Q = 2 & \quad \Rightarrow \quad \hat{\kappa} = 3, \quad n = n(\hat{\kappa}) = 1, \\
Q = 3 & \quad \Rightarrow \quad \hat{\kappa} = 10/3, \quad n = n(\hat{\kappa}) = (1 + \sqrt{5})/2, \\
Q = 4 & \quad \Rightarrow \quad \hat{\kappa} = 4, \quad n = n(\hat{\kappa}) = 2.
\end{cases}
\]

(75)

In general, we see from (74) that the relation between $Q$ and $\hat{\kappa} = 16/\kappa$ in (75) is

\[
\sqrt{Q} = -2 \cos(\hat{\kappa}\pi/4) \quad \text{with} \quad \hat{\kappa} \in (8/3, 4]\]

(The $\hat{\kappa} = 8/3$ cutoff arises because $\text{CLE}_\kappa$ is defined only for $\kappa \in (8/3, 8]$ [101, 103].) Combining this with (61) gives

\[
n := -2 \cos\left(\frac{\pi^2}{\arccos(-\sqrt{Q}/2)}\right), \quad \text{where} \quad Q \in \{2, 3, 4\},
\]

(76)

which relates the loop fugacity $n$ of the dilute-phase loop-gas model to the number of states $Q$ in the corresponding critical Potts model in the continuum limit. Thus, we expect that a dilute-phase loop-gas representation of the critical $Q \in \{2, 3, 4\}$ Potts model exists in the continuum limit as $\text{CLE}_\kappa$, with $\hat{\kappa}$ given by (75).
Our identification of the critical Potts model with the dilute-phase loop-gas model in the continuum limit does not involve BCs, so next, we modify this identification to include the event PMB (item iii of section 3.4) if the Potts model is inside an even-sided polygon $P$. Naively, because all fixed sides of $P$ exhibit the same state in PMB, the LGB event corresponding to PMB, should one exist, must be LGB$_1$. Indeed, the latter corresponds to the random cluster model mutual wiring event RCB$_1$ (figure 17). However, verifying this idea is difficult because the identification of the Potts model with the loop-gas model holds only in the continuum limit. (The $Q=2$ exception does not help because, here with $n=1$ (75), the loop-gas model partition functions (62) are identical for all LGB events $\vartheta$.)

In spite of this difficulty, there is a simple argument that identifies PMB with LGB$_1$. To make it, we extend the lattice inside $P$ so it fills a very large planar region $R$ containing $P$, with the boundary of $R$ far from the boundary of $P$. Next, we restrict the state of the lattice sites inside $P$ to exhibit only the state common to all fixed sides of $P$. This addition alters the Potts model partition function $P^{\vartheta}$, summing over PMB, by only a physically irrelevant constant. Furthermore, the boundary loops, surrounding clusters anchored to the fixed sides of $P$, close through exterior arcs that pass just behind the fluctuating sides of $P$ (figure 29). Because LGB$_1$ is the only LGB event with its boundary loops arranged outside $P$ in this way, we identify PMB with LGB$_1$.

4.4. Crossing probability for loop-gas models

Above, we identify RCB random cluster model partition functions (items iv and v in section 3.3) and PMB Potts model partition functions (items iv and v in section 3.4) with LGB loop-gas model partition functions (items iv and v of this section). To summarize, for the dense and dilute phases, we respectively have

\[
\begin{align*}
\left\{ X^P_{c,0}(p_r, Q) \sim \mathcal{Z}^P_{c,0}(\xi_r, n), \right. \\
\left. X^P_{\partial,0}(p_r, Q) \sim \mathcal{Z}^P_{\partial,0}(\xi_{\partial}, n), \right. \\
\left\{ Y^P_{c,0}(K_r, Q) \sim \mathcal{Z}^P_{c,0}(\xi_r, n), \right. \\
\left. Y^P_{\partial,0}(K_{\partial}, Q) \sim \mathcal{Z}^P_{\partial,0}(\xi_{\partial}, n), \right. \\
\end{align*}
\]

dense, $n = \sqrt{Q}$ and $Q \in \{1, 2, 3, 4\}$,

\[
\begin{align*}
\left\{ Y^P_{c,1}(K_r, Q) \sim \mathcal{Z}^P_{c,1}(\xi_r, n), \right. \\
\left. Y^P_{\partial,1}(K_{\partial}, Q) \sim \mathcal{Z}^P_{\partial,1}(\xi_{\partial}, n), \right. \\
\end{align*}
\]

dilute, $n$ given by (76) and $Q \in \{2, 3, 4\}$.

(77)

Figure 29. $Q=3$ Potts model in a rectangle with the PMB (fixed BC on top/bottom, fluctuating BC on the sides). Surrounding the sample with lattice sites whose states match the fixed BC reveals how boundary loops (red) close outside the rectangle.
as the lattice spacing $a$ vanishes in the continuum limit, where we define the symbol $\sim$ as follows:

$$f(x) \sim g(x) \quad \implies \quad \lim_{x \to c} \frac{f(x)}{g(x)} = \lambda \quad \text{for some } \lambda \in \mathbb{R} \setminus \{0\}. \quad (78)$$

In (77), the constant $\lambda$ is not universal, i.e. it depends on microscopic details of the model such as the lattice. The loop fugacity of the dilute phase $\{ n \in [-2, 2], x = x_i(n) \}$ overlaps that of the dense phase $\{ n \in [0, 2], x = \tilde{x}_i(n) \}$, but conveniently, we may simultaneously represent both phases with a single parameter, the SLE/CLE parameter $\kappa$, by replacing $n$ with (61). After doing this, restricting to $n \in (0, 2]$ as usual, and using (74), (75), we write (77) as

$$
\begin{cases}
X^0_{\kappa}(\kappa) \sim Z^0_{\kappa}(\kappa), & \text{dense, } Q = 4 \cos^2(4\pi/\kappa), \kappa \in (4, 8) \text{ and } Q \in \{1, 2, 3, 4\}, \\
X^0_{\kappa}(\kappa) \sim Z^0_{\kappa}(\kappa), & \text{dilute, } Q = 4 \cos^2(\pi\kappa/4), \kappa \in (8/3, 4] \text{ and } Q \in \{2, 3, 4\},
\end{cases} \quad (79)
$$

where we have combined the dense phase and dilute phase loop-gas model partition functions together into a single function, defined by

$$Z^0_{\kappa}(\kappa) := \begin{cases}
Z^0_{\kappa, a}(x_i(n), n), & \text{dense, } n = -2 \cos(4\pi/\kappa) \text{ and } \kappa \in (4, 8), \\
Z^0_{\kappa, a}(\tilde{x}_i(n), n), & \text{dilute, } n = -2 \cos(4\pi/\kappa) \text{ and } \kappa \in (8/3, 4],
\end{cases} \quad (80)$$

As such, the two crossing probabilities (48), (56) unify under the following ratio of loop-gas model partition functions, which we may propose as a ‘crossing probability’ for a loop-gas model in $\mathcal{P}$ conditioned on an LGB event:

$$P_{\downarrow_{\kappa}}^\mathcal{P} := \mathbb{P}_{\text{loop-gas}(\mathcal{L}G\mathcal{X}_1 \mid \text{LGB}_a)} = \frac{Z^P_{\downarrow_{\kappa}}(\kappa)}{Z^0_{\kappa}(\kappa)} \quad (81)$$

In the next section, we use methods of CFT and results of [36–39] to predict a formula for the continuum limit of this ratio (81) as a function of $\kappa \in (8/3, 8]$.

### 5. The continuum limit and conformal field theory

In this section, we use boundary CFT to predict a formula for the continuum limit of the crossing probability (81). For this, we conformally map the interior of the $2N$-sided polygon $\mathcal{P}$ onto the upper half-plane via a Schwarz–Christoffel transformation. This map extends to a continuous bijection that sends the sides and vertices $w_1, w_2, \ldots, w_{2N}$ of $\mathcal{P}$ to the real axis and the respective points $x_1 < x_2 < \ldots < x_N$. It is widely believed [20, 45, 51], and in some cases rigorously proven [24, 70, 71], that in the continuum limit, the statistics of each of the lattice models considered in this article (Potts model, random cluster model, and loop-gas model) are invariant under a conformal transformation of the model’s domain (but not its lattice). Hence, this transformation does not affect our determination of crossing probabilities, and it has the advantage of sending all polygon interiors to a common domain, the upper half-plane.
Smeared partition functions and conformal field theory

If a Potts model or random cluster model occupies a lattice inside $\mathcal{P}$, then under a conformal transformation sending the interior of $\mathcal{P}$ onto the upper half-plane, the BC associated with each side carries over to its image segment on the real axis, and the boundary condition changes (BCCs) occurring at the vertices carry over to their image points. In the loop-gas representation of these models, the BCCs mark points along the real axis where loops pass into or out of the upper half-plane. We let $Z_{\vartheta}(\text{LG}_X)$, resp. $Z(\text{LG}_X)$ denote the partition function (80) for the loop-gas model in the upper half-plane, summing exclusively over the event $\cap_{\vartheta}(\text{LG}_X \cap \text{LG}_X)$ also sent to the upper half-plane (items ii and iii beneath (62)), and we let $Z_f$ denote the free partition function (59) for the critical loop-gas model in the upper half-plane. As usual, we have

$$Z_{\vartheta}=Z_{\vartheta}+Z_{\vartheta}+\ldots+Z_{\vartheta}.$$  

Before continuing, we address a subtlety regarding the BCC locations on the real axis. In the discrete setting, the BCCs occur exactly at the lattice sites nearest to the image points $x_j$. However, in the continuum limit where the lattice spacing vanishes, the BCCs occur exactly at $x_j$, an event whose probability is zero. To avoid summing over events with vanishing probabilities, we replace our partition function $Z_{\vartheta}(\text{LG}_X)$, resp. $Z_{\vartheta}(\text{LG}_X)$ that sums over all $\text{LG}_X$ events of the type $\text{LG}_X$ that have the $j$th BCC occurring within some small distance $\epsilon_j \sim |x_j - x_{j+1}|$ of $x_j$ (figure 30).

A standard prediction of physics literature [34, 109] is that, after shrinking the lattice spacing $a$ to zero to arrive with the continuum limit, the loop-gas model partition function behaves as

$$\lim_{a \to 0} Z_{\vartheta}(\{x_j\}; \{\epsilon_j\})/Z_{\vartheta} \sim \epsilon_{j_1}^{\theta_1} \epsilon_{j_2}^{\theta_2} \ldots \epsilon_{j_{2N}}^{\theta_{2N}} \Upsilon_{\vartheta}(\{x_j\}),$$  

(82)

where $\Upsilon_{\vartheta}$ is some universal function, and where $\theta_1$ is given in (86) below. Throughout this article, we abbreviate (82) by writing

$$Z_{\vartheta}(\epsilon) \approx \Upsilon_{\vartheta} \text{ as } \epsilon \downarrow 0,$$  

(83)

where the precise meaning of the symbol $\approx$ follows from rewriting (82) in the form (83).

Another well-known prediction of the physics literature [22, 26, 34, 57, 109–112] is that $\Upsilon_{\vartheta}$ equals the $2N$-point boundary CFT correlation function

$$\Upsilon_{\vartheta}(x_1, x_2, \ldots, x_{2N}) = \langle \psi(x_1)\psi(x_2)\cdots\psi(x_{2N}) \rangle_{\vartheta},$$  

(84)
The one-leg boundary operator $\psi_{(j)}$, so-called because it induces the event that a boundary loop passes into or out of the upper half-plane through (or near) the point $x_{j}$, is a primary operator in a CFT with central charge $c$.

$$c(\kappa) = (6 - \kappa)(3 - 8\kappa)2\kappa,$$

where $\kappa$ is the SLE$_{\kappa}$ parameter. The conformal weight $\theta_1$ of $\psi_1$ equals the $(1, 2)$ (resp. $(2, 1)$) Kac weight in the dense (resp. dilute) phase of SLE$_{\kappa}$.

$$\theta_1 = \frac{6 - \kappa}{2\kappa} = \begin{cases} h_{1.2}(c), & \kappa > 4, \\ h_{2.1}(c), & \kappa \leq 4. \end{cases}$$

The $(1, 2)$ (resp. $(2, 1)$) CFT null-state condition implies that (84) satisfies the following elliptic system of null-state PDEs for all $\{j\} \in \ldots$:

$$\sum_{k} \frac{\partial_j \psi_j}{\partial x_k} - \frac{(6 - \kappa)2\kappa}{(x_k - x_j)^2} \psi_j(x_1, x_2, \ldots, x_{2N}) = 0.$$  

(87)

Also, $\psi_\theta$ satisfies three conformal Ward identities that require it to be invariant under uniform translation of the points $x_1, x_2, \ldots, x_{2N}$ and covariant, with weight $\theta_1$, under dilations and special conformal transformations:

$$\sum_{k} [x_k^2 \partial_k + 2\theta_{1}\kappa] \psi_j(x_1, x_2, \ldots, x_{2N}) = 0.$$  

(88)

Altogether, (88) implies that if $f$ is a conformal bijection of the upper half-plane onto itself (it is easy to show that this is a Möbius transformation with real coefficients) such that $x'_i < x'_j$ if $i < j$, where $x'_j := f(x_j)$, then $\psi_\theta$ satisfies

$$\psi_\theta(x'_1, x'_2, \ldots, x'_{2N}) = \partial f(x'_1)^{-\theta_1} \partial f(x'_2)^{-\theta_1} \cdots \partial f(x'_{2N})^{-\theta_1} \psi_\theta(x_1, x_2, \ldots, x_{2N}),$$

(89)

where $\partial f$ is the derivative of $f$. (We note that $\partial f(x) > 0$ for all $x \in \mathbb{R}$.) We call this property (89) Möbius covariance. In light of (82), (84), this transformation rule (89) is natural. Indeed, under $f$, the radius of the $\epsilon_j$-ball centered on $x_j$ dilates to $\partial f(x_j)\epsilon_j$. Combining this fact with (89) gives the conformal invariance of the continuum limit of the ratio $Z_\theta(\{\epsilon_j\})/Z_\theta$ on the left side of (82). This is consistent with our expectations of conformal invariance in these models.

A typical indication of a critical point in a statistical mechanics system such as the Potts model or random cluster model is the existence of certain correlation functions with power-law decay or growth (exclusively) at that critical point. Because we are working only at the critical point of these models, where conformal invariance and non-trivial crossing observables are expected (see the introduction section 1), it is reasonable to assume that $\psi_\theta$ satisfies the bound

$$|\psi_\theta(x_1, x_2, \ldots, x_{2N})| \leq C \prod_{i < j} \max[(x_j - x_i)^p, (x_j - x_i)^{-p}]$$

(90)

for all $x_1 < x_2 < \ldots < x_{2N}$ and some positive constants $p$ and $C$ that may depend on $\psi_\theta$. As in [36–39], we let $S_N$ denote the vector space (over the real numbers) of solutions to the system.
(87), (88) that satisfy the power-law bound (90) for some constants \( p \) and \( C \) (that may depend on the solution). Thus, we have \( \Upsilon_{\vartheta} \in S_{N} \).

In [36–39], two authors of this article completely determine the solution space \( S_{N} \) for all \( \kappa \in (0, 8) \). In [36, 37], they rigorously prove that \( \dim S_{N} \leq C_{\kappa} \) with \( C_{\kappa} \) the \( \kappa \)th Catalan number (3), and in [38], they construct a linearly independent collection of \( C_{\kappa} \) explicit solutions via the CFT Coulomb gas (contour integral) formalism [113, 114], proving that \( \dim S_{N} = C_{\kappa} \) and that such solutions span \( S_{N} \). With the solution space \( S_{N} \) completely understood, we seek a property of \( \Upsilon_{\vartheta} \) that identifies it with a particular explicit element of \( S_{N} \).

### 5.2. Crossing weights and crossing probability

An explicit formula for \( \Upsilon_{\vartheta} \) is not sufficient to determine the continuum limit of the crossing probability (81); we also need a formula for the asymptotic behavior of \( Z_{\vartheta}(\{ \epsilon_{j} \}) \) as \( \epsilon_{j} \to 0 \). Similar to (82), we expect

\[
Z_{\vartheta}(\epsilon) \simeq T_{\vartheta, \vartheta}, \quad \text{as } \epsilon \downarrow 0
\]

for some universal function \( T_{\vartheta, \vartheta} \) (where the meaning of the symbol \( \simeq \) is given in and explained beneath (83)). Then the continuum limit of the crossing probability (81), now in the upper half-plane setting, is given by

\[
P_{\vartheta, \vartheta} = \lim_{\epsilon \to 0} \left( \frac{Z_{\vartheta}(\{ \epsilon_{j} \})}{Z_{\vartheta}(\{ \epsilon_{j} \})} \right) = \frac{\Upsilon_{\vartheta, \vartheta}}{T_{\vartheta, \vartheta}}.
\]

Furthermore, we only need to determine the numerator \( T_{\vartheta, \vartheta} \) of (92), as summing over the crossing index \( \vartheta \) determines the denominator:

\[
\Upsilon_{\vartheta} = Z_{1, \vartheta} + Z_{2, \vartheta} + \cdots + Z_{C_{\kappa}, \vartheta}
\]

\[
\Rightarrow \quad \Upsilon_{\vartheta} = T_{1, \vartheta} + T_{2, \vartheta} + \cdots + T_{C_{\kappa}, \vartheta} \quad \text{for all } \vartheta \in \{1, 2, \ldots, C_{\kappa}\}.
\]

To determine the universal functions \( T_{\vartheta, \vartheta} \), it is useful to study the quantity \( 3_{\vartheta} := n^{-\vartheta} Z_{\vartheta, \vartheta} \), which we identify with the second sum in (62) in the continuum limit. As mentioned beneath (62), we interpret \( 3_{\vartheta} \) as the usual loop-gas model partition function \( Z_{\vartheta, \vartheta} \) except that the boundary loops of each sample have fugacity \( n = 1 \). As such, \( 3_{\vartheta} \) does not depend on the LGB index \( \vartheta \). If we define a smeared version \( 3_{\vartheta}(\{ \epsilon_{j} \}) \) relative to \( 3_{\vartheta} \) in the usual way, then

\[
3_{\vartheta}(\epsilon) \simeq \Pi_{\vartheta}, \quad \text{as } \epsilon \downarrow 0, \quad \text{where } \Pi_{\vartheta} := n^{-\vartheta} T_{\vartheta, \vartheta}.
\]

We call the new universal function \( \Pi_{\vartheta} \) the \( \vartheta \)th crossing weight. Altogether, the partition functions \( Z_{\vartheta, \vartheta} \) and \( 3_{\vartheta} \) and the universal part of their respective asymptotic behaviors \( T_{\vartheta, \vartheta} \) and \( \Pi_{\vartheta} \) are related via the diagram

\[
\begin{align*}
Z_{\vartheta, \vartheta}(\{ \epsilon_{j} \} & \to 0) \\
\Upsilon_{\vartheta, \vartheta} & \downarrow \Pi_{\vartheta}(\epsilon_{j} \to 0) \to 0
\end{align*}
\]

The crossing weights \( \Pi_{\vartheta} \) are useful because we may express all of the universal functions \( T_{\vartheta, \vartheta} \) and \( \Upsilon_{\vartheta} \) and the crossing probability (92) in terms of them. Indeed, from (94), we have for all \( \vartheta \in \{1, 2, \ldots, C_{\kappa}\} \)
\[ \Upsilon_{\beta, \gamma} = n^{k_1} \Pi_1 \implies \Upsilon_{\beta, \gamma} = n^{k_1} \Pi_1 + n^{k_2} \Pi_2 + \cdots + n^{k_{C_N}} \Pi_{C_N}, \quad (96) \]

\[ P_{\beta, \gamma} = n^{k_1} \Pi_1 \implies P_{\beta, \gamma} = n^{k_1} \Pi_1 + n^{k_2} \Pi_2 + \cdots + n^{k_{C_N}} \Pi_{C_N}. \quad (97) \]

(This formula (97) immediately gives the necessary property \( P_{\beta, \gamma} = P_{\beta, \gamma} + P_{\beta, \gamma} + \cdots + P_{\beta, \gamma} = 1 \) for all \( \beta \in \{1, 2, \ldots, C_N\} \).)

Our goal now is to find explicit formulas for all crossing weights \( \Pi_i \). We first show that these functions are elements of \( \mathcal{S}_N \) by solving the system (96) for all of the \( \Pi_i \) and expressing each as a linear combination of \( \Upsilon_1, \Upsilon_2, \ldots, \Upsilon_{C_N} \in \mathcal{S}_N \). This approach is possible if and only if the \( C_N \times C_N \) meander matrix \((M_{\beta, \gamma} \circ n)(\kappa)\) (23) [82–85], whose \((c, \varphi)\)th entry is by definition \( n(\kappa)^{c, \varphi} \) (figure 19), is invertible. Francesco et al [82] calculates the determinant of this matrix, and from this result, we find [37] that \((M_{\beta, \gamma} \circ n)(\kappa)\) is invertible if and only if \( n(\kappa) \) (10) is not among \( n_{q, q''} = -2 \cos(\pi q''/q), \quad \{ \text{with } q, q'' \in \mathbb{Z}^+, 1 \leq q'' < q < N + 1, \text{ and } q, q'' \text{ coprime.} \} \quad (98) \)

Hence, if \( \kappa \in (0, 8) \) does not satisfy \( n(\kappa) = n_{q, q''} \) for any \( q \leq N + 1 \), then the crossing weights \( \Pi_i \) are elements of \( \mathcal{S}_N \). The solutions to this equation are the \textit{exceptional speeds} [38], given by

\[ \kappa_{q, q'} := 4qq', \quad \text{with } q, q' \in \mathbb{Z}^+, q, q' \text{ coprime, and } q > 1. \quad (99) \]

These speeds inconveniently include the random cluster model with \( Q = 1 \) (\( \kappa = \kappa_{3, 2} = 1 \)), \( Q = 2 \) (\( \kappa = \kappa_{4, 3} = 16/3 \)), and \( Q = 3 \) (\( \kappa = \kappa_{6, 5} = 24/5 \) [74]), and other conformally invariant critical models [38]. We assume for now that \( \kappa \) is not an exceptional speed. Later, in the paragraph below (110), we extend the results obtained under this assumption to the case that \( \kappa \) is an exceptional speed.

Asymptotic arguments identify each crossing weight with a unique element of \( \mathcal{S}_N \). We begin with the case \( N = 1 \). With only two BCCs on the real axis, at \( x_1 \) and \( x_2 \) respectively, there is an arc connectivity, in which an arc joins \( x_1 \) with \( x_2 \) and the corresponding \((1, 1)\) th product diagram (figure 19) has \( l_{1, 1} \) boundary loop that passes into the upper half-plane through \( x_1 \) and out of the upper half-plane through \( x_2 \). Thus, there is only one \( N = 1 \) crossing weight \( \Pi_1^{(1)} \), related to the loop-gas partition functions via (93), (95):

\[ Z_{1, 1}^{(1)} = Z_{1, 1}^{(1)} = nZ_{1, 1}^{(1)} \iff \Upsilon_{1, 1}^{(1)} = \Upsilon_{1, 1}^{(1)} = n\Pi_1^{(1)}. \quad (100) \]

In (100) and below, the superscript 1 emphasizes that \( N = 1 \). According to previous arguments, \( \Pi_1^{(1)} \) solves the system (87), (88) with \( N = 1 \), the solution space \( \mathcal{S}_1 \) of which we easily see to be all multiples of \((x_2 - x_1)^1 - b/\kappa \) [36]. So far, \( Z_1 \) and \( \Pi_i \) are defined only up to normalization, so we choose

\[ \Pi_i^{(1)}(x_1, x_2) = (x_2 - x_1)^1 - b/\kappa. \quad (101) \]

For \( N > 1 \) there are \( C_N > 1 \) crossing events to consider, and thus \( C_N \) crossing weights to determine. We determine all of the latter from the \( N = 1 \) case and an asymptotic argument. To begin, we note that for any \( \varsigma \in \{1, 2, \ldots, C_N\} \), the \( \varsigma \)th interior arc connectivity has at least one interior arc with endpoints \( x_i < x_{i+1} \) for some \( i < 2N \). This interior arc corresponds to part of a boundary loop \( \gamma \) joining \( x_i \) to \( x_{i+1} \) in the underlying loop-gas model. As such, if \( x_i \) and \( x_{i+1} \) are very close, then \( \gamma \) only explores a region very close to its two endpoints with high probability. Thus, as \( x_{i+1} \rightarrow x_i \), the system fragments into two statistically independent subsystems: one microscopic with an infinitesimal-sized curve joining \( x_i \) with \( x_{i+1} \), and the other macroscopic.
with the boundary loops joining the remaining $2N - 2$ points as before (figure 31). Because these two subsystems are expected to be statistically independent, the partition function $Z_\varsigma$ for the entire system factors into

$$
Z_\varsigma \sim \prod_{\varsigma' \in \mathcal{S}_{N-1}} Z_\varsigma^{(N-1)}(x_{i}, x_{i+1}; \{\epsilon_j\}_{j=0}^{i-1}),
$$

where $\varsigma' \in \{1, 2, \ldots, C_{N-1}\}$.

(102)

The superscript $N$ indicates the number of points $2N$ and where the $\varsigma'$th connectivity follows from removing the interior arc joining $x_i$ with $x_{i+1}$ from the $\varsigma$th connectivity. Similarly, if $\Pi_\varsigma^{(N-1)} \in \mathcal{S}_{N-1}$ (resp. $\Pi_\varsigma^{(N)} \in \mathcal{S}_{N}$) gives the universal part of the asymptotic behavior of $Z_\varsigma^{(\{\epsilon_j\})}$ (resp. $Z_\varsigma^{(\{\epsilon_j\})}$) as $\epsilon_j \to 0$, then by (94), (101), (102), we have

$$
\Pi_\varsigma^{(N)}(x_1, x_2, \ldots, x_{2N}) \sim \lim_{x_{i+1} \to x_i} \Pi_\varsigma^{(1)}(x_i, x_{i+1}) \times \Pi_\varsigma^{(N-1)}(x_i, x_2, \ldots, x_{j-1}, x_{i+2}, \ldots, x_{2N}).
$$

(103)

Because $\Pi_\varsigma^{(N-1)}$ is a crossing weight, we may apply this same asymptotic argument to it and repeat. Altogether, we find that if we enumerate the arcs in the $\varsigma$th connectivity such that the $j$th arc nests the $k$th arc only if $j > k$ and let $x_{i-1}$ and $x_i$ be the endpoints of the $j$th arc, then for any such enumeration,

$$
[\varsigma^{(N)}] : \lim_{x_{2N} \to x_{2N-1}, x_{2N-2} \to x_{2N-3}} \ldots \lim_{x_i \to x_{i+1}} \lim_{x_{j-1} \to x_j} \ldots \lim_{x_{k-1} \to x_k} \Pi_\varsigma^{(N)}(x_i, x_{i+1}, \ldots, x_{2N}) = 1,
$$

(104)

where $[\varsigma^{(N)}]$ denotes the sequence of limits (actually, an equivalence class of them [38]) shown above. In [38], theorem 8 states that the collection of all $C_N$ of these functionals is a basis for the dual space $\mathcal{S}_N^*$, and we denote this basis as

$$
\mathcal{B}^*_N := \{[\mathcal{A}_1], [\mathcal{A}_2], \ldots, [\mathcal{A}_{C_N}]\} \subset \mathcal{S}_N^*.
$$

(105)

By lemma 12 of [36], reordering the limits in (104) does not change the value of the limit, assuming that the reordering does not cause points to collide (as happens if, e.g., $x_{i-1} < x_{i-2} < x_i < x_{i+1}$ and the $k$th limit precedes the $j$th).

With the quantity $[\varsigma^{(N)}]$, known (103), we next determine $[\varsigma^{(N)}]\Pi_\varsigma$ for all $\vartheta \neq \varsigma$. Combined with (104), this information is sufficient to determine $\Pi_\varsigma$ because $\mathcal{B}^*_N$ is a basis for $\mathcal{S}_N^*$. To this
end, we note that for all $\vartheta \in \{1, 2, \ldots, C_N\}$, $P_{\vartheta, \zeta}$ (92) goes to the probability of the $\zeta$th crossing event conditioned on some LGB event for a $(2N - 2)$-sided polygon as $x_{i+1} \rightarrow x_i$, where $x_i$ and $x_{i+1}$ are still endpoints of a common arc in the $\zeta$th connectivity. Then (96), (97), (103) imply

$$T_\vartheta(x_1, x_2, \ldots, x_{2N}) \sim (x_{i+1} - x_i)^{-6/\kappa} \Psi_\vartheta(x_1, x_2, \ldots, x_{i-1}, x_{i+1}, \ldots, x_{2N}), \quad (106)$$

for all $\vartheta \in \{1, 2, \ldots, C_N\}$ for some function $\Psi_\vartheta \neq 0$ because $P_{\vartheta, \zeta}$ does not vanish as $x_{i+1} \rightarrow x_i$. (Recalling the identification of $T_\vartheta$ with the CFT correlation function in (84), the power-law behavior in (106), with its power $-2\theta_1 + \theta_2 = 1 - 6/\kappa$, implies the Kac fusion rule $\psi_1 \psi_1 = \psi_2$, where $\psi_0$ is the identity operator with conformal weight $\theta_0 = 0$ [33–35].) The index $i$ depends on our choice of crossing weight $\Pi_i$, but because this choice is arbitrary, it is evident that (106) holds for any $i \in \{1, 2, \ldots, 2N - 1\}$ (although the function $\Psi_\vartheta$, while not zero, may depend on $i$).

Meanwhile, $P_{\vartheta, \zeta}$ (92) should vanish as we bring together two endpoints of two different arcs in the $\zeta$th connectivity, such as by sending $x_i \rightarrow x_{i-1}$ (assuming that $i > 1$). Indeed, if $x_{i-1}$ and $x_i$ are close, then it is very probable that a boundary loop joins them together through the upper half-plane in the underlying loop-gas model, making the $\zeta$th crossing event very unlikely. In order for $P_{\vartheta, \zeta}$ (97) to vanish as $x_i \rightarrow x_{i-1}$, the asymptotic behavior (106) of its denominator (96) must dominate that of its numerator $\Pi_i$. According to theorem 2 of [39], $\Pi_i$ then behaves as

$$\Pi_i(x_1, x_2, \ldots, x_{2N}) \sim (x_i - x_{i-1})^{2/\kappa} \Lambda_i(x_1, x_2, \ldots, x_{i-2}, x_{i+1}, \ldots, x_{2N}), \quad (107)$$

for some function $\Lambda_i \neq 0$. (Here, the power-law behavior in (107), with its power $-2\theta_1 + \theta_2 = 2/\kappa$, implies the Kac fusion rule $\psi_1 \psi_1 = \psi_2$, where $\psi_2$ is the two-leg boundary operator with conformal weight $\theta_2 = 8/\kappa - 1$ [37].)

The asymptotic behaviors (103) and (107) of $\Pi_i$ let us determine $[\vartheta, \zeta] \Pi_i$ for all $\vartheta, \zeta \in \{1, 2, \ldots, C_N\}$. To begin, by (103), (104), $[\vartheta, \zeta] \Pi_i = 1$ if $\vartheta = \zeta$. And if $\vartheta \neq \zeta$, then at least one limit of $[\vartheta, \zeta] \Pi_i$ brings together two adjacent endpoints of different interior arcs in the $\zeta$th connectivity. If the $j$th limit of $[\vartheta, \zeta] \Pi_i$ is the first to do this, then (107) gives

$$[\vartheta, \zeta] \Pi_i = \left( \frac{j+1}{\ldots} \frac{\text{last limit in } [\vartheta, \zeta]}{\ldots} \right) \times \lim_{x_{2j} \rightarrow x_{2j-1}} \left( x_{2j} - x_{2j-1} \right)^{6/\kappa - 1} \times \frac{\text{crossing weight}}{\text{with } 2N-j+1 \text{ points}} = 0 \quad (108)$$

for $\vartheta \neq \zeta$ and $\kappa \in (0, 8)$. In summary, we see that the complete collection of crossing weights is dual to the collection (105) of (equivalence classes of) sequences of limits in $B_N$ in the sense that

$$[\vartheta, \zeta] \Pi_i = \delta_{\vartheta, \zeta}. \quad (109)$$

Because $B_N$ (105) is a basis for $S_N$ [38], the duality relation (109) is sufficient to identify each crossing weight with a unique element of $S_N$ and proves that the collection $B_N$ of all $C_N$ crossing weights is a basis for $S_N$:

$$B_N := \{ \Pi_1, \Pi_2, \ldots, \Pi_{C_N} \} \subset S_N. \quad (110)$$

Flores and Kleban [39] calls $\Pi_i \in S_N$ satisfying the duality condition (109) the $\zeta$th connectivity weight. Thus, the terms ‘crossing weight’ and ‘connectivity weight’ are synonymous (as are ‘crossing’ and ‘connectivity’ by figure 16).
So far, our results presume that the crossing weights are elements of \( S_N \), a fact that we proved only for \( \kappa \in (0, 8) \) not an exceptional speed (99) with \( q \leq N + 1 \). To extend our results to these speeds, we reasonably assume that each crossing weight is a continuous function of \( \kappa \). Flores and Kleban [39] shows that each connectivity weight is continuous, so by continuous extension, crossing weights are identified with connectivity weights for exceptional speeds also.

To find explicit formulas for all crossing weights from (109), we use the collection of \( C_N \) elements \( \mathcal{F}_1, \mathcal{F}_2, \ldots, \mathcal{F}_{C_N} \) of \( S_N \) given by item iii of section 2. Their formulas (11) follow from the CFT Coulomb gas (contour integral) formalism [34, 35, 113, 114]. (Dubédat constructed such solutions in [30].) [38] proves that the set

\[
\mathcal{B}_N := \{ \mathcal{F}_1, \mathcal{F}_2, \ldots, \mathcal{F}_{C_N} \} \subset S_N
\]

is a basis for \( S_N \) if and only if \( \kappa \in (0, 8) \) is not an exceptional speed (99) with \( q \leq N + 1 \) and that \( [\mathcal{F}_\varrho, \mathcal{F}_0] = n(\kappa)^{\varrho_0} \) for \( \kappa \in (0, 8) \) (figure 19). As such, we have (see also (23) in item v of section 7)

\[
[\mathcal{F}_\varrho, \mathcal{F}_0] = n^{\varrho_0} \Rightarrow \quad \mathcal{F}_\varrho = n^{\varrho_0} \Pi_1 + n^{\varrho_0} \Pi_2 + \cdots \\
\cdots + n^{\varrho_\kappa} \Pi_{C_N}, \quad \text{for all } \varrho, \varrho \in \{1, 2, \ldots, C_N\}.
\]

As stated above (98), this system (112) is invertible if and only if \( \kappa \in (0, 8) \) is not an exceptional speed (99) with \( q \leq N + 1 \) [38], so for all such \( \kappa \), we obtain formulas for all crossing weights \( \Pi_\varrho \), and the crossing probability (97).

If \( \kappa \) is an exceptional speed with \( q \leq N + 1 \), then the system (112) is not invertible. In this case, we may either isolate the crossing weights for some SLE\( \kappa \) parameter \( \kappa \) near \( \kappa \) and then send \( \kappa \to \kappa \), or we may verify that a certain choice of integration contours in (11) gives a solution satisfying (109) for all \( \varrho \in \{1, 2, \ldots, C_N\} \) and some \( \xi \) in that range. Both approaches have drawbacks. The first is tedious, as the finiteness of the limit of \( \Pi_\varrho(\kappa) \) as \( \kappa \to \kappa \) relies on the cancelation of many infinite quantities. The second is used in [40] for \( N \in \{1, 2, 3, 4\} \), but it is impractical for \( N \geq 5 \).

Equation (112) offers a useful physical interpretation of the Coulomb gas functions in (111). Comparing (112) to (96) gives an explicit formula for the universal function \( \Upsilon_\varrho \) (item iii of section 2):

\[
\Upsilon_\varrho = \mathcal{F}_\varrho, \quad \text{for all } \varrho \in \{1, 2, \ldots, C_N\}.
\]

After inserting (113) into (82), we find a physical interpretation of the Coulomb gas function \( \mathcal{F}_\varrho \). Moreover, using (112), (113), we can sum the denominator of (97) to find

\[
P_{\varrho, \varrho} = \frac{n^{\varrho_0} \Pi_\varrho}{\mathcal{F}_\varrho}
\]

This is a very important step because it justifies the crossing probability formula (24) previously stated in section 2. This summation is a significant simplification because the formulas for the crossing weights are often considerably more complicated than those of \( \mathcal{F}_\varrho \).

Flores and Kleban [39] posits that each crossing weight \( \Pi_\varrho(\kappa) \) is positive-valued for all \( \kappa \in (0, 8) \). Assuming that this is true, the crossing probability formula (97), (114) satisfies \( 0 < P_{\varrho, \varrho} < 1 \) if \( n(\kappa) > 0 \), or \( \kappa \in (8/3, 8) \). This restricts the fugacity of the loop-gas model to \( n \in (0, 2] \), or \( \kappa \in (8/3, 4] \), cutting off the \( n \in [-2, 0] \) range of the dilute phase. To our knowledge,
has no application if $n(\kappa) \leq 0$. Indeed, we cannot interpret it as a probability if it is negative.

It is interesting to interpret (114) in terms of multiple-SLE$_\kappa$ [26], the random process that is expected to give the law of boundary cluster interfaces in the continuum limit of the critical random cluster model and the critical Potts model (figure 32). [39, 112] predict that in a multiple-SLE$_\kappa$ process with ‘SLE$_\kappa$ partition function’ $F \in S_N$, the $2N$ curves exploring the upper half-plane join together pairwise in the $j$th connectivity with probability

$$ P_j = \frac{a_j \Pi_j}{a_1 \Pi_1 + a_2 \Pi_2 + \cdots + a_{2N} \Pi_{2N}} = \frac{a_j \Pi_j}{F}, \quad a_j := [\ell_j]F. $$

(115)

If $F = F_0$, then (115) matches (114). This implies the following: if we choose $F = F_0$ for the SLE$_\kappa$ partition function with $\kappa$ among (74), then the generated multiple-SLE$_\kappa$ curves match the FK boundary cluster interfaces in the corresponding (continuum limit) critical random cluster model with the $\nu$th mutual wiring event $\text{RCB}_\nu$. Or if we choose $F = F_1$ for the SLE$_\kappa$ partition function with $\kappa$ among (75), then the generated multiple-SLE$_\kappa$ curves match the spin boundary cluster interfaces in the corresponding (continuum limit) critical Potts model with the $\text{PMB}$ event.

To finish, we note that the half-plane crossing probability formulas (97), (114) are manifestly Möbius invariant. Indeed, the crossing weights, as elements of $S_N$, are Möbius covariant, satisfying the functional equation (89) for any conformal bijection $f$ of the upper half-plane onto itself that preserves coordinate order. Therefore, for any such $f$, (97), (114) give

$$ P_{\rho}(x_1, x_2, \ldots, x_{2N}) = P_{\rho}(x_1', x_2', \ldots, x_{2N}'), \quad x_i' := f(x_i). $$

(116)

An example of this is the following Möbius transformation $f$ that sends the points $x_1 < x_2 < x_3 < \ldots < x_{2N-2} < x_{2N-1} < x_{2N}$ in that order to the points $0 < \lambda_1 < \lambda_2 < \ldots < \lambda_{2N-3} < 1 < \infty$ respectively. Here, (116) gives

$$ f(x_j) = \frac{(x_j - x_1)(x_{2N} - x_{2N-1})}{(x_{2N-1} - x_1)(x_{2N} - x_j)}, \quad \lambda_j := f(x_{j+1}) $$

$$\implies\quad P_{\rho}(x_1, x_2, x_3, \ldots, x_{2N-2}, x_{2N-1}, x_{2N}) = \lim_{w \to \infty} P_{\rho}(0, \lambda_1, \lambda_2, \ldots, \lambda_{2N-3}, 1, w). $$

(117)

This shows that the crossing probability $P_{\rho}(x)$ depends only on the $2N - 3$ cross-ratios $\lambda_j$ (117) that can be formed from the points $x_1, x_2, \ldots, x_{2N}$. These cross-ratios are Möbius invariants.

Numerical evidence [20, 21] and analytic results [24, 71] promote Möbius invariance (116) of crossing probabilities to conformal invariance. (Alternatively, we may posit, as we do in appendix A, that the partition functions appearing in (92) are conformally invariant, which...
immediately implies conformal invariance of crossing probabilities.) Indeed, if $f$ is a conformal map from the upper half-plane onto a Jordan domain $D$, then
\[ P^D_{\gamma}(\kappa \mid w_1, w_2, \ldots, w_{2N}) = P^D_{\gamma}(\kappa \mid x_1, x_2, \ldots, x_{2N}), \quad \text{where } w_j := f(x_j), \quad (118) \]
gives the crossing probability for the same model (on the same lattice) in $D$, with the BC or boundary-loop condition at $x_j$ sent to the same condition at $w_j \in \partial D$. Unlike (116), (118) is not a functional equation $\kappa$ because the left side of (118) is a function completely different from that on the right side. Indeed, $P^D_{\gamma}$ and $P^D_{\gamma}$ have different domains. Thanks to (117), we see that $P^D_{\gamma}$ depends on only $2N - 3$ degrees of freedom that determine the shape, but not the position, orientation, or scale, of $D$ in the complex plane.

Appendix A shows that conformal covariance of boundary CFT correlation functions is consistent with the conformal invariance law (118). This justification encounters a difficulty in our application: if $f$ is a conformal map that sends the upper half-plane onto the interior of a polygon, then its continuous extension to the real axis is not conformal at the vertex pre-images $x_j$. In appendix A, we circumvent these problems formally with the introduction of CFT corner operators [57, 115].

6. Formula for rectangle crossing probability

Section 2 gives a step-by-step construction of the formula (24) for crossing probability. As an example, we calculate all crossing probabilities (24) for the rectangle ($N = 2$), generalizing Cardy’s formula for percolation (2). At times, we introduce some helpful but nonessential simplifications that are typically not available for $N > 2$. The steps of this section mirror those of section 2, and except for this reliance, this section is completely self contained. To keep it so, we avoid any notations and terminology not introduced in section 2.

(i) We let $x_1 < x_2 < x_3 < x_4$ be four marked points on the real axis, let $\kappa \in (0, 8)$ (the SLE$_\kappa$ parameter) [66–68], and let $n(\kappa) = -2 \cos(4\pi/\kappa)$ (the ‘loop-gas model fugacity formula’) [51, 93].

(ii) We let the first arc connectivity comprise an arc joining $x_2$ and $x_3$ and another arc joining $x_1$ with $x_4$. We let the second arc connectivity comprise an arc joining $x_1$ with $x_2$ and another arc $x_3$ with $x_4$.

(iii) We use the formulas for $\mathcal{F}_\theta$ given in section III of [38] (or (C.2) of appendix C) instead of those in item iii of section 2 because the former are a little simpler. According to our enumeration of connectivities in step ii, we have

\[
\mathcal{F}_\theta(\kappa \mid x_1, x_2, x_3, x_4) = \frac{n(\kappa)^3(2 - 8/\kappa)}{4\sin^2(4\pi/\kappa)(1 - 4/\kappa)^2}[x_2 x_3 x_2]^2[x_4 x_2 x_4]^2[1 - 6/\kappa]^{-3/\kappa} \times \int_\Gamma (u - x_1)^{-4/\kappa} \left\{ \begin{array}{ll} (u - x_2)^{-4/\kappa}, & \theta = 1 \\ (x_3 - u)^{-4/\kappa}, & \theta = 2 \end{array} \right\} (x_3 - x_4)^{-4/\kappa}(x_4 - u)^{-1/\kappa - 2} du,
\]

where $\Gamma = \begin{cases} \mathcal{P}(x_2, x_3), & \theta = 1, \\ \mathcal{P}(x_1, x_2), & \theta = 2, \end{cases}$ (119)
\( x_i := x_j - x_i \), and \( \mathcal{P}(x_i, x_j) \) is the Pochhammer contour of figure 8. It is useful to express these functions (119) in terms of hypergeometric functions \([43, 44]\). To begin, the quantities

\[
I_0(\kappa | x_1, x_2, x_3, x_4) := \frac{\mathcal{F}_1(\kappa | x_1, x_2, x_3, x_4)}{[x_3 x_4]^2^{1 - 6/\kappa}},
\]
\[
I_0(\kappa | x_1, x_2, x_3, x_4) := \frac{\mathcal{F}_2(\kappa | x_1, x_2, x_3, x_4)}{[x_3 x_4]^2^{1 - 6/\kappa}}
\]

are Möbius invariants because their numerators and denominators transform as in (89) (using (86)) for a conformal bijection \( f \) of the upper half-plane onto itself preserving coordinate order. So with \( \lambda = x_1 x_3 x_4 / \lambda = x_2 x_3 x_4 \),

\[
I_0(\kappa | x_1, x_2, x_3, x_4) = I_0(\kappa | 0, \lambda, 1, \infty) = \frac{n(\kappa)^2 (2 - 8/\kappa)}{4 \sin^2(4\pi / \kappa) \Gamma(1 - 4/\kappa) \Gamma(1 - 4/\kappa)}
\]

\[
\times \int_0^1 u^{-4/\kappa} \left( (u - \lambda)^{-4/\kappa}, \vartheta = 1 \right) (1 - u)^{-4/\kappa} \mathrm{d}u, \quad \Gamma = \left\{ \mathcal{P}(\lambda, 1), \vartheta = 1, \right\}
\]
\[
\times \left\{ \mathcal{P}(0, \lambda), \vartheta = 2, \right\}
\]

\[
= n(\kappa)^2 \eta^{2/\kappa} (1 - \eta)^{1 - 6/\kappa} \left( 4 \kappa, 1 - 4 \kappa, 8 \kappa, 1 - \eta \right) \eta = \left\{ \lambda, \vartheta = 1, \right\}
\]
\[
\times \left\{ 1 - \lambda, \vartheta = 2, \right\}
\]

which, when inserted into (120), gives a formula for \( \mathcal{F}_0 \) in terms of hypergeometric functions.

(iv) We have \( h_1 = b_2 = 2 \) and \( h_2 = b_1 = 1 \) (figure 11).

(v) Using our results from items iii and iv (figure 11), we may calculate the crossing weights \( \Pi \) by inverting the \( 2 \times 2 \) meander matrix \( M_2(n) \), as follows:

\[
\begin{align*}
\left( \begin{array}{c}
\mathcal{F}_1 \\
\mathcal{F}_2
\end{array} \right) &= \left( \begin{array}{cc}
n^2 & n \\
n & n^2
\end{array} \right) \left( \begin{array}{c}
\Pi_1 \\
\Pi_2
\end{array} \right) \\
\implies \Pi_1 &= \frac{n \mathcal{F}_1 - \mathcal{F}_2}{n(n^2 - 1)}, \quad \Pi_2 = \frac{n \mathcal{F}_2 - \mathcal{F}_1}{n(n^2 - 1)}.
\end{align*}
\]

By using a standard linear transformation formula of the hypergeometric function \([44]\), we may express these crossing weights in terms of another hypergeometric function:

\[
\Pi_1(\kappa | x_1, x_2, x_3, x_4) := \frac{\Gamma(12/\kappa - 1)}{\Gamma(8/\kappa) \Gamma(8/\kappa - 1)} \left( [x_3 x_4]^2 \right)^{1 - 6/\kappa}
\]

\[
\times \eta^{2/\kappa} (1 - \eta)^{1 - 6/\kappa} \left( 4 \kappa, 1 - 4 \kappa, 8 \kappa, 1 - \eta \right) \eta = \left\{ \lambda, \vartheta = 1, \right\}
\]
\[
\times \left\{ 1 - \lambda, \vartheta = 2, \right\}
\]

\[
(123)
\]

The meander matrix \( (M_2 \circ n)(\kappa) \) is not invertible if \( n(\kappa) = 1 \), that is, if \( \kappa = 6 \) or \( \kappa = 6(3m \pm 1) < 4 \) for some \( m \in \mathbb{Z}^+ \). If \( \kappa \) is any of these values, then the limit of \( \Pi_1(\kappa) \) as \( \kappa \to \kappa \) is subtle to calculate from the formula given in (122) because \( n(\kappa) = 1 \), but trivial to find from the formula in (123).

(vi) We assemble these results as in (24) to find formulas for crossing probabilities of the models described in item vi of section 2. Because these formulas depend only on \( \lambda \), we express all probabilities as functions of only this parameter:
we interpret these probabilities as follows. our descriptions here are equivalent to those in items (vi)a–(vi)c of section 2 above, but we state them a little differently.

(a) Q-state random cluster model: With Q ∈ {1, 2, 3, 4}, we suppose that the continuum limit of a critical random cluster model on a lattice fills the upper half-plane and that the segments (x1, x2) and (x3, x4) are fixed to some state. With (25) relating Q and κ ∈ [4, 8), we have (figure 33)

\[ P_{1|2}(κ | λ) = P_{1|2}(κ | 1 − λ) = \frac{\Gamma(12/κ − 1)\Gamma(4/κ)}{\Gamma(8/κ)\Gamma(8/κ − 1)} \frac{\text{2F1}(\frac{2}{κ}; 1 − \frac{4}{κ}; \frac{8}{κ} | λ)}{\text{2F1}(\frac{2}{κ}; 1 − \frac{4}{κ}; 2 − \frac{8}{κ} | 1 − λ)}, \]  

(124)

\[ P_{1|2}(κ | λ) = P_{2|1}(κ | 1 − λ) = \frac{\Gamma(12/κ − 1)\Gamma(4/κ)}{2 \cos(4πκ/8)\Gamma(8/κ)\Gamma(8/κ − 1)} \times \frac{\text{2F1}(\frac{2}{κ}; 1 − \frac{4}{κ}; \frac{8}{κ} | λ)}{(1 − λ)^{8/κ − 1} \text{2F1}(\frac{2}{κ}; 1 − \frac{4}{κ}; 2 − \frac{8}{κ} | λ)}. \]  

(125)

We interpret these probabilities as follows. Our descriptions here are equivalent to those in items (vi)a–(vi)c of section 2 above, but we state them a little differently.

(a) Q-state random cluster model: With Q ∈ {1, 2, 3, 4}, we suppose that the continuum limit of a critical random cluster model on a lattice fills the upper half-plane and that the segments (x1, x2) and (x3, x4) are fixed to some state. With (25) relating Q and κ ∈ [4, 8), we have (figure 33)

\[ P_{1|1}(κ | λ) = \frac{\text{the probability that one FK cluster touches both (x1, x2) and (x3, x4), conditioned on both of these segments being fixed to the same state}}{\text{the probability that both (x1, x2) and (x3, x4), conditioned on both of these segments being fixed to the same state}}, \]  

(127)

\[ P_{1|2}(κ | λ) = \frac{\text{the probability that one FK cluster touches both (x1, x2) and (x3, x4), conditioned on both of these segments being independently fixed}}{\text{the probability that these segments being independently fixed}}, \]  

(128)

\[ P_{2|1}(κ | λ) = \frac{\text{the probability that no FK cluster touches both (x1, x2) and (x3, x4), conditioned on both of these segments being fixed to the same state}}{\text{the probability that these segments being fixed to the same state}}, \]  

(129)
\[ P_{2|2}(\kappa \mid \lambda) = \left\{ \text{the probability that no FK cluster touches both } (u_1, x_2) \text{ and } (x_3, x_4), \text{ conditioned on both of these segments being independently fixed} \right\}. \quad (130) \]

It is interesting to examine some special cases. First, the case \( Q = 1 \) corresponds to critical percolation. Setting \( \kappa = 6 \) in (124), (125), we recover Cardy’s formula (2):

\[
P_{1|1}(6 \mid \lambda) = P_{2|2}(6 \mid \lambda) = P_{1|1}(6 \mid 1 - \lambda) = P_{2|2}(6 \mid 1 - \lambda) = \frac{3\Gamma(2/3)}{(1/3)^2} \begin{pmatrix} \frac{2}{3} & \frac{4}{3} \\ \lambda & \lambda \end{pmatrix}. \quad (131)
\]

(See (151) below for a formula for the derivative of \( P_{2|2}(6 \mid \lambda) \) in terms of the Dedekind eta function.) These formulas are independent of the second index \( \vartheta \) because, with one color distributed to all bond clusters, there is no difference between the BC for \( \vartheta = 1 \) and the BC for \( \vartheta = 2 \). Next, the case \( Q = 2 \) corresponds to the \( Q = 2 \) random cluster (i.e. FK-Ising) model. Setting \( \kappa = 16/3 \) in (124), (125) gives

\[
P_{1|1}(16/3 \mid \lambda) = P_{2|2}(16/3 \mid 1 - \lambda) = \frac{\sqrt{2}\lambda}{\sqrt{1 + \sqrt{1 - \lambda}} \sqrt{1 + \sqrt{\lambda}}},
\]

\[
P_{1|2}(16/3 \mid \lambda) = P_{2|1}(16/3 \mid 1 - \lambda) = \frac{\sqrt{\lambda}}{1 + \sqrt{1 - \lambda}}. \quad (132)
\]

(After transforming to the rectangular geometry in item vii below, we give alternative formulas (149), (150) for these probabilities in terms of the Dedekind eta function.) In [71], Chelkak and Smirnov rigorously calculated the above Ising-FK crossing probabilities, and in our notation, their results read

\[
P_{1|1}(16/3 \mid \lambda) = P_{2|2}(16/3 \mid 1 - \lambda) = \frac{\sqrt{2} \, q(\lambda)}{p(\lambda) + \sqrt{2} \, q(\lambda)},
\]

\[
P_{1|2}(16/3 \mid \lambda) = P_{2|1}(16/3 \mid 1 - \lambda) = \frac{q(\lambda)}{\sqrt{2} \, p(\lambda) + q(\lambda)},
\]

where \( 1 - q(\lambda) = p(\lambda) := \frac{\sqrt{1 - \sqrt{\lambda}}}{\sqrt{1 - \sqrt{1 - \lambda}} + \sqrt{1 - \sqrt{1 - \lambda}}} \). \quad (133)

In order to verify that these formulas (133) agree with ours (132), we consider the difference of these two formulas for \( P_{1|1}(16/3 \mid \lambda) \):

\[
\begin{align*}
\frac{(131)}{p(\lambda) + \sqrt{2} \, q(\lambda)} - \frac{(132)}{\sqrt{2} \, q(\lambda)} & = \frac{(131)}{\sqrt{1 + \sqrt{1 - \lambda}} \sqrt{1 + \sqrt{\lambda}}} \\
& = \frac{(133)}{\sqrt{2} \, q(\lambda)} \left( \frac{g(\lambda)}{\sqrt{2} \, q(\lambda)} - \frac{1 - g(\lambda)}{\sqrt{2} \, q(\lambda)} \right) = \frac{(133)}{\sqrt{2} \, q(\lambda)} \left( \frac{g(\lambda)}{\sqrt{2} \, q(\lambda)} - \frac{1 - g(\lambda)}{\sqrt{2} \, q(\lambda)} \right) \\
& = h(\lambda) = \frac{(134)}{f(\lambda) = g(\lambda) + h(\lambda) \text{ satisfies}}.
\end{align*}
\]

To show that this difference vanishes, we may either note that for all \( \lambda \in [0, 1] \), \( g(\lambda)^2 = h(\lambda)^2 \) and \( g(1) = h(1) \), from which \( g(\lambda) = h(\lambda) \) follows, or we may observe that \( f = g + h \) satisfies
\( \lambda(1 - \lambda)\partial^2 f(\lambda) + (1/2 - \lambda)\partial f(\lambda) + f(\lambda)/16 = 0 \) \hspace{1cm} (135)

and the initial conditions \( f(1/2) = \partial f(1/2) = 0 \), which leads directly to \( f = 0 \). Equivalence of the two formulas for \( P_{2|1}(16/3 | \lambda) \) then follows from the fact that the first (132) satisfies

\[
P_{1|1}(16/3 | \lambda) + P_{2|1}(16/3 | \lambda) = \frac{\sqrt{2\lambda}}{\sqrt{1 + \sqrt{1 - \lambda}} \sqrt{1 + \sqrt{\lambda}}} + \frac{\sqrt{1 - \lambda}}{1 + \sqrt{\lambda}} = 1. \hspace{1cm} (136)
\]

We may also verify this algebraic identity (136) by moving the second term to the right side, cross-multiplying the denominators, and squaring both sides of the resulting equation.

(b) \textit{Q-state Potts model:} With \( Q \in \{2, 3, 4\} \), we suppose that the continuum limit of a critical Potts model on a lattice fills the upper half-plane and that the segments \((x_1, x_2)\) and \((x_3, x_4)\) are fixed to the first state while boundary sites outside these segments freely sample any but the first state. With \((26)\) relating \( Q \) and \( \kappa \in (0, 4) \),

\[
P_{1|1}(\kappa \mid \lambda) = \left\{ \begin{array}{l}
\text{the probability that one spin cluster touches both } (x_1, x_2) \\
\text{and } (x_3, x_4), \text{conditioned on both of these segments being}
\text{fixed to the first spin state and the rest of the boundary}
\text{freely sampling any but the first spin state}
\end{array} \right. \hspace{1cm} (137)
\]

\[
P_{2|1}(\kappa \mid \lambda) = \left\{ \begin{array}{l}
\text{the probability that no spin cluster touches both } (x_1, x_2)
\text{and } (x_3, x_4), \text{conditioned on both of these segments being}
\text{fixed to the first spin state and the rest of the boundary}
\text{freely sampling any but the first spin state}
\end{array} \right. \hspace{1cm} (138)
\]

It is interesting to look at the special case \( Q = 2 \) corresponding to the Ising model. Setting \( \kappa = 3 \) in (124), (125), we find the following formula for crossing probability of Ising spin clusters:

\[
P_{1|1}(3 \mid \lambda) = P_{2|1}(3 \mid 1 - \lambda) = \frac{2\Gamma(4/3)}{\Gamma(5/3)\Gamma(8/3)} \left( \frac{\lambda^{5/3}}{1 - \lambda + \lambda^2} \right) \left[ \begin{array}{l}
\text{2F}_1 \left( \begin{array}{l}
\frac{1}{3}, \frac{4}{3}, \frac{8}{3} \\
\lambda
\end{array} \right)
\end{array} \right]. \hspace{1cm} (139)
\]

The literature [26–30] contains at least three different formulas for this probability (139), summarized in section 7 of [28] and discussed in [29]. In particular, the second formula in [28] simplifies to (139) via linear transformation formulas [44] of the hypergeometric function.

(c) \textit{Loop-gas model:} With \( \kappa \in (0, 2] \), we suppose that the continuum limit of a dense-phase or dilute-phase loop-gas model on a lattice fills the upper half-plane. In addition, ‘boundary’ loops exit and then re-enter the upper half-plane through \( x_1, x_2, x_3, \) and \( x_4 \), with exactly one loop passing once through each point. With

\[
n(\kappa) = -2\cos(4\pi/\kappa), \hspace{1cm} \begin{array}{c}
\kappa \in (8/3, 4], \\
\kappa \in (4, 8),
\end{array} \hspace{1cm} \text{dilute phase,}
\hspace{1cm} \text{dense phase,}
\hspace{1cm} (140)
\]

formulas (124), (125) for \( P_{i|0} \) give the probability that the parts of the boundary loops in the upper-half plane connect the points \( x_1, x_2, x_3, \) and \( x_4 \) pairwise in the \( \kappa \)th connectivity, conditioned on the event that the parts of the boundary loops inside the lower-half plane join these points pairwise in the \( \kappa \)th connectivity. Figure 34 illustrates some of these interpretations for the random cluster model, the Potts model, and the loop-gas model. (We note that our formulas apply only in the continuum limit, not shown in the figure.)
We map the upper half-plane onto the rectangle $\mathcal{R} = \{ x + iy \mid 0 < x < \mathcal{R}, 0 < y < 1 \}$ via the conformal bijection that sends $x_1 < x_2 < x_3 < x_4$ to the corners $w_1 = 0$, $w_2 = \mathcal{R}$, $w_3 = \mathcal{R} + i$, and $w_4 = i$ respectively. The fixed segments $(x_1, x_2)$ and $(x_3, x_4)$ go to the bottom and top sides of $\mathcal{R}$ respectively, and the BC of the half-plane system goes to the equivalent BC for the image system too.

Our goal is to write the crossing probability formulas (124), (125) as explicit functions of the aspect ratio $\mathcal{R}$. The following functions are useful for this purpose: (Although these functions are well-known in the literature [43, 44], conventions for their formulas vary. Therefore, we explicitly write out the formulas that we use here.)

- The complete elliptic function of the first kind [43, 44]:
  \[ K(m) := \int_0^{\pi/2} \frac{d\theta}{\sqrt{1 - m \sin^2 \theta}}. \]  
  \[ (141) \]

- The second, third, and fourth Jacobi theta functions [43, 44]:
  \[ \vartheta(z, q) := 2q^{1/4} \sum_{n=0}^{\infty} q^{(n+1)/2} \cos((2n + 1)z) = \sum_{n=-\infty}^{\infty} q^{(n+1)/2} \omega^{2n+1}, \]  
  \[ (142) \]
  \[ \vartheta_3(z, q) := \sum_{n=-\infty}^{\infty} q^n \cos(2nz) = \sum_{n=-\infty}^{\infty} q^n \omega^{2nz}, \]  
  \[ (143) \]
  \[ \vartheta_4(z, q) := \sum_{n=-\infty}^{\infty} (-1)^n q^n \cos(2nz) = \sum_{n=-\infty}^{\infty} (-1)^n q^n \omega^{2nz}. \]  
  \[ (144) \]

- The Dedekind eta function [116]:
  \[ \eta(\tau) := e^{i\pi\tau/12} \prod_{n=1}^{\infty} (1 - e^{2\pi in\tau}) = \sum_{n=-\infty}^{\infty} (-1)^n q^{(6n+1)^2/24} \text{ where } q := e^{2\pi i \tau}. \]  
  \[ (145) \]
The following identities relate these functions [115, 116]:

\[ R = \frac{K(\lambda)}{K(1 - \lambda)} \iff \lambda = \left( \frac{\partial_3(0, e^{-\tau R})}{\partial_3(0, e^{-\tau R})} \right)^4 = 1 - \left( \frac{\partial_2(0, e^{-\tau R})}{\partial_3(0, e^{-\tau R})} \right)^4, \]

\[ \partial_2(0, e^{\tau R}) = \frac{2\eta(2\tau)^2}{\eta(\tau)}, \quad \partial_3(0, e^{\tau R}) = \frac{\eta(\tau)^5}{\eta(\tau/2)^2\eta(2\tau)^2}, \quad \partial_4(0, e^{\tau R}) = \frac{(\tau/2)^4}{\eta(\tau)}. \]

(Also, the transformation \( R \mapsto 1/R \) sends \( \lambda \mapsto 1 - \lambda \) in (146).) Then because the crossing probability (24) is (predicted to be) conformally invariant, we have the following rectangle crossing probability (with \( P_{\alpha\beta} \) given by (124), (125)):

\[ P_{1|1}^R(\kappa \mid R) = P_{1|1}(\kappa \mid 1/R), \quad P_{2|2}^R(\kappa \mid R) = P_{2|2}^R(\kappa \mid 1/R). \]

After replacing \( (x_1, x_2) \) with ‘the bottom side of \( \mathcal{R} \)’ and \( (x_3, x_4) \) with ‘the top side of \( \mathcal{R} \)’, (127)–(130) and (137), (138) interpret (147), (148) for the random cluster model and the Potts model in \( \mathcal{R} \) respectively. Also, after replacing ‘\( x_j \)’ with ‘\( w_j \)’ and ‘upper half-plane’ with ‘\( \mathcal{R} \)’, item (vi) c interprets (147), (148) for the loop gas model in \( \mathcal{R} \).

Using (146), we may also write the rectangle crossing probability formulas (147), (148) in terms of the Dedekind eta function (145). Furthermore, in the case of the \( Q = 2 \) random cluster model (\( \kappa = 16/3 \)), we may use (132), (146) to write these formulas as algebraic functions of the Dedekind eta function. For example, we find

\[ P_{1|2}^R(16/3 \mid R) = \frac{\eta(\tau/2)^8\eta(2\tau)^4}{\eta(\tau)^12 + 4\eta(2\tau)^8\eta(\tau/2)^2}, \quad \text{where } \tau := iR. \]

Using (149) and the relations \( P_{1|2}^R + P_{2|1}^R = 1 \) and \( P_{1|1}^R + P_{2|2}^R = 1 \), we find similar formulas for \( P_{1|1}^R \) and \( P_{2|2}^R \) in terms of Dedekind eta functions. Working numerically in Mathematica, we have also found
The relations $P_{1|2}^R + P_{2|1}^R = 1$ and $P_{1|1}^R + P_{2|1}^R = 1$ applied to (149), (150) leads to some identities for the Dedekind eta function which seem to be unknown.

Finally, we recall another well-known appearance of the Dedekind eta function in crossing probability. The derivative of Cardy’s formula (131) has an explicit expression in terms of the Dedekind eta function, namely

$$\frac{\partial}{\partial R} \cdot \frac{\partial}{\partial R} = \frac{2}{\sqrt{3} \Gamma(1/3)^3} \eta(iR)^2. \quad (151)$$

7. Simulation results for rectangle and hexagon crossing probability

In this section, we present computer simulation results for FK cluster (resp. spin cluster) crossing probabilities for the critical $\{Q \in \mathbb{Q} \in [2, 3, 4] \}$ random cluster (resp. Potts) model on a square lattice inside a rectangle $R$ and inside a hexagon $H$. Item i (resp. item ii) of the introduction section 1 explains the corresponding side-alternating BCs. Here, we assign the top and bottom sides of $R$ the fixed BC, and we assign the upper left/right and bottom sides of $H$ the fixed BC. (Flores et al [92] presents simulation results for critical percolation ($Q = 1$) crossing probabilities for the rectangle.)

Our simulations sampled thirty-three rectangles, each with two successive side lengths, $\ell$ for the length of (i.e. number of lattice sites in) the top and bottom sides, and $\ell'$ for the length of the left and right sides. For the $i^{th}$ rectangle, we chose

$$\ell_i \approx 2048 - \ell_i \approx \ell_j \approx 1633, \quad \text{for } i \in \{1, 2, \ldots, 33\}. \quad (152)$$

These side-lengths (152) are almost uniformly distributed between $\ell_1$ and $\ell_{33}$, and with $R_i := \ell_i \ell'_i$ the aspect ratio of the $i^{th}$ rectangle, we have $R_1 \approx 1/16, R_{17} = 1$, and $R_{33} \approx 16$.

The formula (24) for crossing probability presumes that we condition the system inside the rectangle on a side-alternating BC event, as described in items i or ii of the introduction.
For the random cluster model, these are RCB events, and the bottom row of figure 35 illustrates them with exterior arc rectangle diagrams (item iii of section 3.2), where any two fixed sides joined by an exterior arc are mutually wired together (i.e. constrained to exhibit the same state). So with the top and bottom sides of \( R \) fixed, these two sides are mutually wired together in the first RCB event, \( RCB_1 \), and independently wired (i.e. not constrained to exhibit the same state) in the second, \( RCB_2 \). For the Potts model (section 3.4), we condition on the PMB event with the top and bottom sides of \( R \) mutually wired together and the left and right sides having the fluctuating BC. Conditioned as usual on either of the two RCB events described above, there are \( C_2 = 2 \) distinct FK-cluster crossing events of the rectangle, the event \( RCX_1 \) that a vertical crossing cluster joins the two fixed sides of \( R \), and the event \( RCX_2 \) that there is no such cluster (items i and ii of section 3.2). The top row of figure 35 illustrates these crossing events with interior arc rectangle diagrams (item ii of section 3.2). There are also \( C_2 = 2 \) distinct spin-cluster crossing events of the rectangle, \( PMX_1 \) and \( PMX_2 \). These events are similarly defined, but for spin clusters rather than FK clusters. Because the probabilities of the first and second crossing events sum to one, our simulations measure the probability of only the first, the vertical crossing event \( RCX_1 \) or \( PMX_1 \).

We used the Swendsen–Wang algorithm [118], with bond activation probability \( p = p_{c_{sq}}^{MF} \) (39), to generate \( 2^{22} = 4194304 \) samples of the critical random cluster model and the critical Potts model in each rectangle. The time required to generate all samples for one rectangle, using a single 2GHz processor, was about 5 d. For each model, we seeded the FK clusters in a way that respects the corresponding side-alternating BC:

(i) **The \( Q \)-color random cluster model:** \( Q \in \{2, 3, 4\} \). In each sample, we seeded the FK boundary clusters that touch a fixed side of \( R \) before we seeded any other FK cluster. The RCB event requires that we activate all FK bonds within the top and bottom sides of \( R \).

If these sides were mutually wired together, then we assigned them the same color, and if they were wired independently of each other, then we did not impose this constraint.

(ii) **The \( Q \)-state Potts model:** \( Q \in \{2, 3, 4\} \). In each sample, we seeded the spin boundary clusters that touch a fixed side of \( R \) first, giving them the first spin state. Then we seeded the spin clusters that touched a fluctuating side (item ii of the introduction section 1) of \( R \).
next, giving each of them any but the first spin state with uniform probability. Finally, we seeded all other spin clusters, giving each any spin state with uniform probability.

The wiring within the fixed sides of $\mathcal{R}$ ensured that one FK (resp. spin) boundary cluster touched the bottom side of $\mathcal{R}$ in case i (resp. case ii). If this cluster touched the top side of $\mathcal{R}$ too, then the sample exemplified a vertical crossing event ($\varsigma = 1$), and if otherwise, then it did not. We tallied the number of samples with a vertical crossing and divided it by the total number of samples to measure the vertical crossing probability.

Formula (24) with $x_1 = 0, x_2 = m, x_3 = 1,$ and $x_4 = \infty$ gives the probability $P_{i|\vartheta}$ of the $\varsigma$th FK (resp. spin) cluster crossing event $RCX_\varsigma$ (resp. $PMX_\varsigma$) conditioned on the $\vartheta$th RCB event $RCB_\vartheta$ (resp. the one PMB event, with $\vartheta = 1$) for the rectangle $\mathcal{R}$ with aspect ratio $R$ (bottom to left side) (146). (There, we have $\lambda = m$. Also, section 6 casts this formula in the simpler form (124), (125).) Then according to figure 35, $P_{i|1}$ (resp. $P_{i|2}$) is the probability of the vertical crossing event in $\mathcal{R}$, conditioned on the mutual wiring (resp. independent wiring) BC event.

Figure 36 compares our simulation measurements of these rectangle crossing probabilities to our predictions (124), (125), (147), (148). We plot the probabilities as functions of $x = (R - 1)/(R + 1) \in (0,1)$, where $R$ is the rectangle’s aspect ratio (bottom to left side). Our plots show excellent agreement for $Q \in \{2,3\}$ and more modest agreement for $Q = 4$. (Actually, we did not need to test the $Q = 2$ cases because their crossing probability formulas (133), (139) [28, 71] have been proven rigorously. But for completeness, we include them here.) Table 2 gives the average relative error of our measurements and the standard deviation of our relative errors from the average relative error. (See the last two paragraphs of section 7.2 for an explanation of these results.)

### 7.2. Simulation results for hexagon crossing probability

In this section, we present computer simulation measurements of FK cluster (resp. spin cluster) crossing probabilities for the critical $Q \in \{2,3,4\}$ random cluster (resp. Potts) model on a triangular lattice inside an equiangular hexagon $\mathcal{H}$. (We presented simulation measurements of critical percolation ($Q = 1$) crossing probabilities for the hexagon in [92].) Our simulations sampled thirty-three hexagons, each with two successive side lengths: $\ell$ for the length of (i.e. number of lattice sites in) the bottom and upper left/right sides, and $\ell'$ for the length of the remaining sides. For the $i$th hexagon, we chose

$$\ell_i = [60.2353 + 28.2353(i - 1)], \quad \ell'_i \approx 1024 - \ell_i \approx \ell_{34-i}, \quad \text{for } i \in \{1, 2, \ldots, 33\}. \quad (153)$$

These side-lengths (153) are almost uniformly distributed between $\ell_1$ and $\ell_{33}$, and with $R_i := \ell_i/\ell'_i$ the side-length ratio of the $i$th hexagon, we have $R_i \approx 1/16, R_{17} = 1$, and $R_{33} \approx 16$.

In our simulations, the triangular lattice is represented by a square lattice with six nearest neighbors to each lattice site, as figure 37 shows. Also, the hexagon is not equiangular, but its interior angles measure $3\pi/4, \pi/2, 3\pi/4, 3\pi/4, \pi/2,$ and $3\pi/4$ radians respectively, in counterclockwise order starting from the bottom-left vertex. These angles ensure that the hexagon’s sides follow along nearest-neighbor bonds. The transformation of figure 37 takes us from this setup to the usual triangular lattice inside the equiangular hexagon $\mathcal{H}$.

The crossing probability formula (24) presumes that we condition the system inside the hexagon on a side-alternating BC event, as described in items i or ii of the introduction section 1. In our present situation, these BCs are as follows:
Table 2. Summary of errors in crossing probability measurements for the rectangle with mutual wiring (θ = 1) and independent wiring (θ = 2), and for the hexagon with mutual wiring (θ = 1), mixed wiring (θ = 3), and independent wiring (θ = 5). We averaged each relative error over all thirty-three rectangles and hexagons, except that we dropped the points for rectangles and hexagons with crossing probability less than 0.01 because our simulations may have generated too few samples of such rare crossing events, resulting in artificially high relative errors. We averaged the deviation of each relative error from the average relative error over all thirty-three rectangles and hexagons, but again, we dropped the points for the same rectangles and hexagons as before for the same reason as before. Figure 35 (resp. figure 38) illustrates the $\varsigma$th crossing event for the rectangle (resp. hexagon) respectively, with the top and bottom sides (resp. bottom and upper left/right sides) fixed.

| Polygon | BC | $\varsigma$ | $\theta$ | Random cluster model (FK clusters) | Potts model (spin clusters) |
|---------|----|------------|---------|---------------------------------|-----------------------------|
|         |     | $Q = 2$ rel. err. | $Q = 2$ std. dev. | $Q = 3$ rel. err. | $Q = 3$ std. dev. | $Q = 4$ rel. err. | $Q = 4$ std. dev. | $Q = 2$ rel. err. | $Q = 2$ std. dev. | $Q = 3$ rel. err. | $Q = 3$ std. dev. | $Q = 4$ rel. err. | $Q = 4$ std. dev. |
| Rectangle | mut. | 2 | 1 | -0.002 065 | 0.003 407 | -0.002 349 | 0.006 693 | -0.071 972 | 0.161 850 | 0.001 901 | 0.003 128 | 0.022 325 | 0.039 424 | 0.133 829 | 0.241 286 |
|         | indep. | 2 | 2 | -0.001 054 | 0.001 852 | -0.004 215 | 0.006 788 | -0.049 402 | 0.128 972 |                   |                 |                   |                 |                   |                 |
| Hexagon | mut. | 3 | 1 | -0.002 468 | 0.004 573 | -0.007 783 | 0.013 127 | -0.227 979 | 0.208 866 | -0.000 516 | 0.008 806 | 0.029 516 | 0.027 194 | 0.365 309 | 0.27 194 |
|         | indep. | 5 | 3 | -0.001 861 | 0.001 454 | 0.001 402 | 0.007 845 | -0.048 882 | 0.156 690 |                   |                 |                   |                 |                   |                 |
|         |      | 1 | 5 | -0.002 535 | 0.003 067 | -0.003 441 | 0.006 427 | -0.008 541 | 0.116 411 |                   |                 |                   |                 |                   |                 |
|         |     | 3 | 5 | -0.001 637 | 0.006 212 | -0.004 639 | 0.007 255 | -0.218 566 | 0.156 977 |                   |                 |                   |                 |                   |                 |

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Figure 36. Comparison of computer simulation results to our prediction (24), (124), (125) of vertical cluster crossing probability in a rectangle with aspect ratio (bottom side to left side) $R$. Item i (resp. item ii) of the introduction section 1 defines the BCs and crossing events for random cluster model FK (resp. Potts model spin) clusters. The rectangle to the right of each graph shows the vertical crossing event whose probability $P_{i/0}$ appears in the graph. In it, the crossing boundary cluster is colored gray, sides touching a boundary cluster are fixed, and any two fixed sides (resp. one fixed side) with endpoints joined by a dashed exterior arc are mutually wired together (resp. wired independently of the other fixed side).
(i) The $Q$-color random cluster model: Item i of the introduction section 1 and items i–v of section 3.2 describe the available side-alternating BC events. They are RCB events, and there are $C_3 = 5$ of them, $\text{RCB}_1, \text{RCB}_2, \ldots, \text{RCB}_5$, enumerated as follows: in $\text{RCB}_1$, the three fixed sides of $\mathcal{H}$ are mutually wired together (i.e. constrained to exhibit the same state). In $\text{RCB}_5$, the three sides are independently wired (i.e. not constrained to exhibit the same state). Finally, in $\text{RCB}_2, \text{RCB}_3,$ and $\text{RCB}_4$, two fixed sides are mutually wired together while the remaining fixed side is wired independently of the others. Because the hexagon is invariant under rotations of $2\pi/3$ (thanks to its alternating side-lengths), we may identify these last three events as one. All in all, our simulations sampled the $\text{RCB}$ events $\text{RCB}_1, \text{RCB}_3,$ and $\text{RCB}_5$, which we respectively call ‘mutual wiring’, ‘independent wiring’, and ‘mixed wiring’.

(ii) The $Q$-state Potts model: Item ii of the introduction section 1 and section 3.4 describe the one available side-alternating BC event. It is the event PMB with all fixed sides of $\mathcal{H}$ mutually wired together and the remaining three sides of $\mathcal{H}$ exhibiting the fluctuating BC.

In our simulations, we imposed the fixed BC on the bottom and upper-left/right sides of $\mathcal{H}$. As such, the remaining three sides of $\mathcal{H}$ exhibited either the free BC or the fluctuating BC. The bottom row of figure 38 illustrates these side-alternating BCs with exterior arc hexagon.
diagrams (item iii of section 3.2). There, any two fixed sides that are joined by an exterior arc are mutually wired together.

There are \( C_3 = 5 \) distinct FK-cluster crossing events of the hexagon, \( \text{RCX}_1, \text{RCX}_2, \ldots, \text{RCX}_5 \) (items i and ii of section 3.2), and we enumerate them as follows: in \( \text{RCX}_1 \), a single cluster connects all fixed sides of \( \mathcal{H} \). In \( \text{RCX}_5 \), no cluster connects any two fixed sides of \( \mathcal{H} \). Finally, in \( \text{RCX}_2, \text{RCX}_3, \) and \( \text{RCX}_4 \), a single cluster connects two fixed sides of \( \mathcal{H} \) but does not touch the remaining fixed side. The top row of figure 38 illustrates these crossing events with interior arc hexagon diagrams (item ii of section 3.2).

Thanks to some symmetries of the hexagon, we only needed to measure probabilities of some of these crossing events if we conditioned on an RCB event (item i). For example, because the mutual (resp. independent) wiring RCB event \( \text{RCB}_1 \) (resp. \( \text{RCB}_5 \)) is invariant under rotations of \( 2\pi/3 \), the probabilities of the FK-cluster crossing events \( \text{RCX}_2, \text{RCX}_3, \) and \( \text{RCX}_4 \) conditioned on this RCB event are equal. As such, we only measured the conditioned probability of the event \( \text{RCX}_3 \) in which a single FK cluster connects the bottom and upper-right sides of \( \mathcal{H} \). Although the mixed wiring RCB event \( \text{RCB}_3 \) does not have this rotational symmetry, it is evident that the probabilities of \( \text{RCX}_2 \) and \( \text{RCX}_4 \) conditioned on \( \text{RCB}_3 \) are equal, so we only measured the conditioned probability of the former event.

There are also \( C_3 = 5 \) distinct spin-cluster crossing events of the hexagon, \( \text{PMX}_1, \text{PMX}_3, \ldots, \text{PMX}_5 \) (items i–ii of section 3.4). In the event \( \varsigma_{\text{PMX}} \), spin clusters join the fixed sides of \( \mathcal{H} \) as do FK clusters in the event \( \varsigma_{\text{RCX}} \). Again, thanks to symmetries of the hexagon, we only needed to measure probabilities of events \( \text{PMX}_1, \text{PMX}_3, \) and \( \text{PMX}_5 \) when we condition on the one available PMB event (item ii).

We used the Swendsen–Wang algorithm [118], with bond activation probability \( p = p^{ni.} \) (40), to generate \( 2^{22} = 4194304 \) samples of the critical random cluster model and the critical Potts model in every hexagon. The time required to generate all samples for one hexagon, using a 2 GHz processor, was about five days. We seeded the FK (resp. spin) clusters of case i (resp. case ii) for the critical random cluster (resp. Potts) model inside the hexagon, as we did in case i (resp. case ii) for the same model inside the rectangle. As we generated the FK (resp. spin) boundary clusters of a sample, we noted which fixed sides of \( \mathcal{H} \) these clusters touched to determine the crossing pattern of the sample. We tallied the number of samples with the \( \varsigma \) th crossing pattern and divided this total by the total number of samples to measure the \( \varsigma \) th FK (resp. spin) cluster crossing probability.

The formula (24) for the crossing probability is given in terms of six coordinates \( x_j \), corresponding to the images of the six vertices \( w_j \) of the hexagon \( \mathcal{H} \) under a conformal bijection from the interior of \( \mathcal{H} \) onto the upper half-plane \( \mathbb{H} \). We chose the inverse of this bijection to be the Schwarz–Christoffel transformation \( f(A.6) \) with parameters (A.14) having \( C = 1 \) (appendix A). With these choices,

\[
f(z) = \frac{2}{3} \int_0^z \zeta^{-1/3}(m_1 - \zeta)^{-1/3}(m_2 - \zeta)^{-1/3}(m_3 - \zeta)^{-1/3}(1 - \zeta)^{-1/3} d\zeta,
\]

This map (154) sends the boundary points (figure 39)

\[
x_1 = 0 < x_2 = m_1 < x_3 = m_2 < x_4 = m_3 < x_5 = 1 < x_6 = \infty
\]

(155)
counterclockwise to the hexagon’s vertices $w_1, w_2, \ldots, w_6$ respectively. With $w_1 = f(0) = 0$, $w_2 = f(m_1) > 0$, and $w_3 = f(m_2)$ in $\mathbb{H}$, the hexagon resides in $\mathbb{H}$ with its base flush against the positive-real axis.

According to (117), the crossing probability actually depends on just three parameters, $\lambda_j = m_j$ (117) for $j \in \{1, 2, 3\}$, that determine the shape of $\mathcal{H}$. Restricting to $m = -\frac{m}{m}$, $m = -\frac{m}{m}$, $m = -\frac{m}{m}$, and $m = -\frac{m}{m}$, ensures that the side-lengths of $\mathcal{H}$ alternate. Let us briefly explain why: it is easy to see that the following conformal bijection $\phi : \mathbb{H} \mapsto \mathbb{H}$ continuously extended to $\mathbb{R}$ and thus sending $m = 0, 0 \mapsto \infty$, and $\infty \mapsto 1$,

$$\phi(m \mid z) := \frac{z - m}{z}$$

(157)

rotates the hexagon $\mathcal{H}$ from (154) by $-\pi/3$ radians. If the side-lengths of $\mathcal{H}$ alternate, then $\mathcal{H}$ must be invariant under two consecutive such rotations, corresponding to $z \mapsto \phi(\phi(m \mid m_2) \mid \phi(m \mid z))$. This implies the system of equations

\[
\begin{align*}
0 &= \phi(\phi(m_1 \mid m_2) \mid \phi(m_1 \mid m_2)), & m_1 &= \phi(\phi(m_1 \mid m_2) \mid \phi(m_1 \mid m_3)), \\
2 &= \phi(\phi(m_1 \mid m_2) \mid \phi(m_1 \mid 1)), & m_3 &= \phi(\phi(m_1 \mid m_2) \mid \phi(m_1 \mid \infty)), \\
1 &= \phi(\phi(m_1 \mid m_2) \mid \phi(m_1 \mid 0)), & \infty &= \phi(\phi(m_1 \mid m_2) \mid \phi(m_1 \mid m_1)).
\end{align*}
\]

(158)

Solving this system gives (156). This reduces the number of independent parameters that control the shape of $\mathcal{H}$ from three to one. Hence, if $\ell$ (resp. $\ell'$) is the length of the bottom (resp. top) side, then the side-length ratio $R = \ell/\ell'$ is

$$\ell = f(m_1), \quad \ell' = \left| f(m_2) - f(m_1) \right| \quad \Rightarrow \quad R = \frac{f(m_1)}{|f(m_2) - f(m_1)|}.$$  

(159)

Relations (154)–(159) put $R \in (0, \infty)$ in one-to-one correspondence with $m_2 \in (0, 1)$. We can manipulate these relations using the symmetry of the hexagon to find...
\[
R = \int_0^1 \left[ \zeta (1 - \zeta) \left( 1 - \frac{m_2^2}{1 - m_2 + m_2^2} \zeta \right) \left( 1 - m_2 \zeta \left( 1 - \frac{m_2^2}{1 - m_2 + m_2^2} \zeta \right) \right) \right]^{-1/3} \, d\zeta \\
\times \left\{ \int_0^1 \left[ \zeta (1 - \zeta) \left( 1 - \left( 1 - \frac{m_2^2}{1 - m_2 + m_2^2} \zeta \right) \right) \right]^{-1/3} \, d\zeta \right\}.
\]

Formula (24) with (155), (156) gives the probability \( P_{\Phi \Theta} \) of the \( \zeta \)th FK (resp. spin) cluster crossing event \( \text{RCX}_\zeta \) (resp. \( \text{PMX}_\zeta \)) conditioned on the \( \vartheta \)th RCB event \( \text{RCB_\vartheta} \) (resp. the one PMB event with \( \vartheta = 1 \)) for the hexagon \( \mathcal{H} \) with alternating side lengths and side-length ratio \( R \). Items i–v of section 2 determine formulas for the functions \( \mathcal{F}_\varphi \) and \( \Pi \) that appear in this formula. However, simpler formulas are available for the hexagon. Indeed, we may replace the formula (11) for \( \mathcal{F}_\varphi \) by its alternative (C.2) with one less integration contour. With connectivities labeled as in figure 10, we have

\[
\mathcal{F}_1 = \text{(C.2) with } \Pi_1 = \mathcal{P}(x_2, x_3), \ \Pi_2 = \mathcal{P}(x_4, x_5), \ \text{and } c = 6, \tag{161}
\]

\[
\mathcal{F}_3 = \text{(C.2) with } \Pi_1 = \mathcal{P}(x_2, x_3), \ \Pi_2 = \mathcal{P}(x_5, x_6), \ \text{and } c = 4, \tag{162}
\]

\[
\mathcal{F}_5 = \text{(C.2) with } \Pi_1 = \mathcal{P}(x_2, x_3), \ \Pi_2 = \mathcal{P}(x_3, x_4), \ \text{and } c = 6, \tag{163}
\]

where \( \mathcal{P}(x_i, x_j) \) is the Pochhammer contour entwining points \( x_i \) and \( x_j \) (figure 8). Our choices of \( c \) values in (161)–(162) ensure that integration contours entwine adjacent branch points and that no integration contour arcs over or wraps around another contour or branch point. This choice facilitates the numerical integration. In addition, section II.C of [40] gives simpler formulas for the hexagon crossing weights \( \Pi_1 \) and \( \Pi_2 \). These formulas use the functions

\[
I_{\nu}^\lambda (\kappa \mid m_1, m_2, m_3) := n(\nu)^2 \frac{\Gamma(2 - \frac{8}{\kappa})^2}{\Gamma(1 - \frac{4}{\kappa})^2} \prod_{k \neq j}^{N} \left( x_k - x_\pi \right)^{2/\kappa} \prod_{m = 1}^{6} \left| x_c - x_m \right|^{\frac{\lambda - 6/\kappa}{\nu}} \int_{x_i}^{x_{i+1}} \int_{x_j}^{x_{j+1}} \mathcal{N} \left( (u_2 - u_1)\frac{2}{\kappa} \prod_{p = 1}^{2} (x_c - u_p) \frac{2}{\kappa} - \prod_{q = 3}^{6} (x_q - u_p) \right) \, du_1 \, du_2,
\]

with \( \{ x_1 = 0, x_2 = m_1, x_3 = m_2, x_4 = m_3, x_5 = 1, x_6 = \infty, i \neq j, i, j, c \in \{ 1, 2, 3, 4, 5, 6 \} \} \) and \( \int_{x_i}^{x_i} := \int_{-\infty}^{0} \). \tag{164}

where \( \mathcal{N} \[ \ldots \] \) orders the differences inside its brackets so the integrand is real-valued over \( [x_i, x_{i+1}] \times [x_j, x_{j+1}] \). From the formulas for \( \Pi_1 \) and \( \Pi_2 \) we find formulas for the other crossing weights, \( \Pi_\zeta \) with \( \zeta \in \{ 3, 4, 5 \} \), by rotating the hexagon to send either the first or second connectivity to the \( \zeta \)th connectivity and then cyclically permuting the arguments of \( \Pi_1 \) or \( \Pi_2 \) and the value of \( c \) consistently with this rotation.

Next, we adjust the formulas for the crossing weights given in the previous paragraph in order to minimize the number of contour integrals that must be computed numerically. Assuming \( \kappa > 4 \) (so we may replace all Pochhammer contours with simple curves via (9))
Indeed, (165) gives a formula for \( \Pi_1 \) that is apparently even simpler than the one presented in section II.C of [40]. Also, we use (166) for our formula for \( \Pi_2 \), which is slightly different from the formula given in section II.C of [40]; this difference arises from a changed value of \( c \) in (164). Using the techniques presented in [40], it is easy to verify that (165)–(168) are indeed formulas for the crossing weights \( \Pi_1, \Pi_2, \Pi_3, \) and \( \Pi_5 \).

Upon inserting (161)–(163), (165)–(168) into (24) with \( N = 3 \) and (155), we find explicit formulas for crossing probabilities in the hexagon. Altogether, for the hexagon with mutual wiring, the probability of the \( \varsigma \)th FK (resp. spin) cluster crossing event with \( \varsigma \in \{1, 3, 5\} \) is

\[
P^\varsigma \Pi_1'(\kappa \mid m_1, m_2, m_3) = \frac{n^2 I^{(2)}_{13} - n I^{(2)}_{12}}{I^{(6)}_{24}}(\kappa \mid m_1, m_2, m_3),
\]

(169)

\[
P^\varsigma \Pi_3'(\kappa \mid m_1, m_2, m_3) = \frac{n I^{(2)}_{34}}{I^{(6)}_{24}}(\kappa \mid m_1, m_2, m_3),
\]

(170)

\[
P^\varsigma \Pi_5'(\kappa \mid m_1, m_2, m_3) = \frac{I^{(4)}_{24} - n I^{(4)}_{34}}{I^{(6)}_{24}}(\kappa \mid m_1, m_2, m_3),
\]

(171)

where \( \kappa \) is related to \( Q \) via (74) (resp. (75), denoted there as \( \hat{\kappa} \)). For the hexagon with mixed wiring, the probability of the \( \varsigma \)th FK-cluster crossing event with \( \varsigma \in \{1, 2, 3, 5\} \) is given by

\[
P^\varsigma \Pi_1'(\kappa \mid m_1, m_2, m_3) = \frac{n I^{(2)}_{13} - n^2 I^{(2)}_{12}}{I^{(6)}_{25}}(\kappa \mid m_1, m_2, m_3),
\]

(172)

\[
P^\varsigma \Pi_2'(\kappa \mid m_1, m_2, m_3) = \frac{I^{(2)}_{12}}{I^{(6)}_{25}}(\kappa \mid m_1, m_2, m_3),
\]

(173)

\[
P^\varsigma \Pi_3'(\kappa \mid m_1, m_2, m_3) = \frac{n I^{(4)}_{34}}{I^{(6)}_{25}}(\kappa \mid m_1, m_2, m_3),
\]

(174)

\[
P^\varsigma \Pi_5'(\kappa \mid m_1, m_2, m_3) = \frac{n I^{(4)}_{24} - n^2 I^{(4)}_{34}}{I^{(6)}_{25}}(\kappa \mid m_1, m_2, m_3),
\]

(175)

where, again, \( \kappa \) is related to \( Q \) via (74). Finally, for the hexagon with independent wiring, the probability of the \( \varsigma \)th FK-cluster crossing event with \( \varsigma \in \{1, 3, 5\} \) is
\[ P_{15}^\kappa(\kappa \mid m_1, m_2, m_3) = \left[ \frac{r_{13}^{(2)}(k) - nI_{13}^{(2)}}{r_{13}^{(2)}} \right](\kappa \mid m_1, m_2, m_3), \tag{176} \]

\[ P_{35}^\kappa(\kappa \mid m_1, m_2, m_3) = \left[ \frac{nI_{13}^{(2)}}{r_{13}^{(2)}} \right](\kappa \mid m_1, m_2, m_3), \tag{177} \]

\[ P_{55}^\kappa(\kappa \mid m_1, m_2, m_3) = \left[ \frac{n^2 I_{23}^{(4)}(\kappa) - n^2 I_{34}^{(4)}(\kappa)}{I_{13}^{(4)}} \right](\kappa \mid m_1, m_2, m_3), \tag{178} \]

where, again, \( \kappa \) is related to \( Q \) via (74). Restricted to (156), these formulas (169)–(178) are easy to integrate numerically for \( \kappa > 4 \) and values of \( m_2 \in (0, 1) \) sufficiently far from zero and one, using Mathematica. However, numerical integration can give poor results for the first \( (R \approx 1/16, m_2 \approx 0) \) and last \( (R \approx 16, m_2 \approx 1) \) two hexagons in (153).

The improper integrals \( I_{ij}^{(1)}(\kappa) \) (164) diverge for all \( \kappa \in (0, 4) \), which includes the \( Q \in (2, 3, 4) \) critical Potts models (75). To resolve this issue, we replaced all simple integration contours with Pochhammer contours as in (9) and divided each Pochhammer contour integral by the normalization factor on the right side of the equation in that figure. In order to numerically integrate the result, we decomposed each Pochhammer contour as in figure 8. The integrals \( I_{ij}^{(1)}, I_{ij}^{(2)} \), having two simple integration contours that span adjacent intervals, need to be evaluated with nested Pochhammer contours as in figure 40. These nested contours correspond to the ‘rainbow’ configurations discussed in [40]. We evaluated the outer Pochhammer contour of figure 40 as we do in figure 8, but with the left point \( x_i \) of the latter figure replaced by the inner Pochhammer contour of figure 40, so \( \beta_1 = -4/\kappa + 12/\kappa - 2 + 8/\kappa = 16/\kappa - 2 \). Thus,

\[ I_{ij}^{(2)}(\kappa \mid x_1, x_2, \ldots, x_6) = \frac{1}{4e^{16\pi i/\kappa} \sin(-4\pi/\kappa) \sin(12\pi/\kappa)} \int_{\Gamma_2} \int_{\Gamma_1} \prod_{i=1}^{6} (u_i - x_1)^{4/\kappa} (u_i - x_2)^{2/\kappa - 2} \prod_{j=1}^{6} (x_j - u_i)^{4/\kappa} du_i \prod_{j=1}^{6} du_j, \tag{179} \]

where \( \Gamma_2 = \mathcal{D}(x_1, x_2) \) and \( \Gamma_1 = \mathcal{D}(\Gamma_2, x_3) \) is the Pochhammer contour with its left (resp. right) cycle exclusively surrounding \( \Gamma_1 \) (resp. \( \Gamma_2 \)). The first (resp. second) prefactor on the right side of (179) arises from dividing out by the prefactor on the right side of the equation in (9) associated with \( I_1 \) (resp. \( I_2 \)), as mentioned above. Using the techniques of section II.E in [40], we may verify that inserting the right side of (179) gives an alternative formula for \( I_{ij}^{(1)}(\kappa) \) that holds for all \( \kappa \in (0, 8) \). Thus, the right side of (179) is the analytic continuation of \( I_{ij}^{(1)}(\kappa) \) to \( \kappa \in (0, 4] \).

This formula (179) for \( I_{ij}^{(2)}(\kappa) \) has an inconvenience: among the values of \( \kappa \) in (74), (75), it is singular at \( \kappa = 16/3 \) (corresponding to \( Q = 2 \) in the random cluster model). Fortunately, this singularity is removable. To see this, we first observe that because \( \kappa = 16/3 > 4 \), we may replace \( \Gamma_2 \) with the simple contour \([x_1, x_2]\) via figure (9), so (179) simplifies to
\[
I_{12}^{(2)}(\kappa | x_1, x_2, \ldots, x_6) = \frac{1}{4e^{2\pi i/\kappa} \sin(16\pi/\kappa) \sin(-4\pi/\kappa)} \oint_{\Gamma_1} \int_{s_1} \prod_{i=1}^{2} (u_i - x_i)^{-4i/\kappa} (u_1 - x_2)^{2i/\kappa - 2} (u_2 - x_1)^{-4i/\kappa} (u_3 - x_4)^{2i/\kappa - 2} (u_4 - x_3)^{-4i/\kappa} \prod_{j=3}^{6} (x_j - u_j)^{-4i/\kappa} \, du_2 \, du_i,
\]

where \( \Gamma_1 = \mathcal{P}(\{x_1, x_2\}, x_3) \) is the Pochhammer contour with its left (resp. right) cycle exclusively surrounding \([x_1, x_2]\) (resp. \(x_3\)). After decomposing \( \Gamma_1 \) into two loops via figure 8 with \( x_i \mapsto [x_1, x_2] \), \( \beta_1 = 16/\kappa - 2 \), \( x_j = x_3 \), and \( \beta_j = -4/\kappa \) (here, we give the straight contour in figure 8 length zero, so its corresponding term does not contribute anything and the two loops mentioned are tangent to each other at a point on the real axis), we find

\[
I_{12}^{(2)}(\kappa | x_1, x_2, \ldots, x_6)
= \left[ \frac{e^{-2\pi i(-4/\kappa)} - 1}{4e^{2\pi i/\kappa} \sin(16\pi/\kappa) \sin(-4\pi/\kappa)} \oint_{\Gamma} \int_{s_1} \prod_{i=1}^{2} (u_i - x_i)^{-4i/\kappa} (u_1 - x_2)^{2i/\kappa - 2} (u_2 - x_1)^{-4i/\kappa} (u_3 - x_4)^{2i/\kappa - 2} (u_4 - x_3)^{-4i/\kappa} \prod_{j=3}^{6} (x_j - u_j)^{-4i/\kappa} \, du_2 \, du_i \right]
\]

\[
= \prod_{i=1}^{2} \prod_{j=3}^{6} (x_j - u_i)^{-4i/\kappa} \, du_2 \, du_i,
\]

where \( \Gamma \) (resp. \( \Gamma' \)) is a counterclockwise-oriented simple loop exclusively surrounding \([x_1, x_2]\) (resp. \(x_3\)). The limit as \( \kappa \to 16/3 \) of the term with integration around \( \Gamma' \) obviously exists. To see that the limit as \( \kappa \to 16/3 \) of the term with integration around \( \Gamma \) exists too, we note that at \( \kappa = 16/3 \), the integrand does not acquire a phase factor as \( u_1 \) traces \( \Gamma \). Thus, \( \Gamma \) closes (i.e. its start-point and endpoint lie on the same Riemann sheet of the integrand), so the function

\[
F(\kappa | x_1, x_2, \ldots, x_6) = \prod_{i<j}^{6} (x_j - x_i)^{2/\kappa} \prod_{k=2}^{6} \left| x_k - x_2 \right|^{|-6/\kappa|} \oint_{s_1} \int_{s_1} \prod_{i=1}^{2} (u_i - x_i)^{-4i/\kappa} (u_1 - x_2)^{2i/\kappa - 2} (u_2 - x_1)^{-4i/\kappa} (u_3 - x_4)^{2i/\kappa - 2} (u_4 - x_3)^{-4i/\kappa} \prod_{j=3}^{6} (x_j - u_j)^{-4i/\kappa} \, du_2 \, du_i,
\]

is an element of \( \mathcal{S}_N \) if \( \kappa = 16/3 \), according to theorem 2 of [38]. However, \([x_1, x_2], [x_3, x_4], [x_4, x_5], \) and \([x_5, x_6]\) are two-leg intervals (definition 13 of [36]) of \( F(\kappa = 16/3) \), so we have \( \left\lfloor F(\kappa = 16/3) \right\rfloor = 0 \) for all equivalence classes \( \zeta \in \{1, 2, 3, 4, C_3 = 5\} \). (We describe \( \left\lfloor F(\kappa = 16/3) \right\rfloor \) beneath (104). See definitions 9 and 11 of [36] for a formal definition.) As a consequence of
this fact, lemma 15 of [36] implies that $F(\kappa = 16/3) = 0$, so because $F$ is an analytic function of $\kappa$, we must have $F(\kappa) = O(\kappa - 16/3)$ as $\kappa \to 16/3$. Altogether, this implies that the limit as $\kappa \to 16/3$ of the term with integration around $\Gamma$ in (181), and therefore of $I_{12}^{(2)}(\kappa)$, exists, which is what we wanted to show.

The case with $\kappa = 4$, corresponding to the $Q = 4$ critical random cluster model (74) and the $Q = 4$ critical Potts model (75), also requires special treatment. In this case, the improper integrals of (164) diverge, so we must regularize them by replacing all simple integration contours with Pochhammer contours, as in (9). But if $\kappa = 4$, then the branch points $x_j$ of these contour integrals are, in fact, simple poles. Hence, we may use the Cauchy integral formula to explicitly evaluate all of these contour integrals and find algebraic expressions for all crossing probabilities. (See appendix C for an example of such an evaluation.)

The formulas given in (169)–(178) hold whenever $0 < m_1 < m_2 < m_3 < 1$. However, as a result of the $2\pi/3$ rotation invariance of the hexagon that arises from (156), some of the ostensibly different contour integrals in (164) are in fact equal. For example, we have $I_{12}^{(2)}(\kappa | m_4, m_2, m_3) = I_{34}^{(2)}(\kappa | m_4, m_2, m_3)$ whenever this rotational symmetry (156) holds. Using this and similar equalities, we minimized the number of different contour integrals needed to numerically evaluate (169)–(178).

A final alteration exploited the Möbius symmetry of (28) to use the points $x_1 = 1 - 1/(1 - m_2)$, $x_2 = 0$, $x_3 = 1 - m_2$, $x_4 = 1$, $x_6 = \infty$ rather than (155), (156). This is equivalent to the previous choice (155), (156) up to a scaling and translation. This change was useful because the nested Pochhammer contours of figure 40 were difficult to evaluate in the coordinates of (155), (156) whenever $m_2$ approached 0 or 1 because groups of four points approached each other, with integration contours between them.

Figure 41 compares our simulation measurements of the critical Potts model spin-cluster crossing probabilities for the hexagon with the mutual wiring PMB to our predictions (169)–(171). Figure 42 (resp. figure 43, resp. figures 44 and 45) compares our simulation measurements of the critical random cluster model FK-cluster crossing probabilities for the hexagon with the mutual wiring (resp. mixed wiring, resp. independent wiring) RCB to our predictions (169)–(171) (resp. (176)–(178), resp. (172)–(175)). We plot the probabilities as functions of $x = (R - 1)/(R + 1) \in (0, 1)$, where $R$ is the hexagon’s ratio of alternating side-lengths (bottom to top side).

Table 2 gives the relative error of our measurements, averaged over all thirty-three rectangles and hexagons, except that we drop the points for hexagons and rectangles with crossing probability less than 0.01 because our simulations may have generated too few samples of such rare crossing events, resulting in artificially high relative errors. For $Q \in \{2, 3\}$, the relative errors are all small, on the order of $10^{-3}$, except the $Q = 3$ critical Potts model with relative error on the order of $10^{-2}$. These low relative errors support our prediction of the formula.
Figure 41. Comparison of computer simulation measurements to our prediction (24), (169)–(171) of spin-cluster crossing probability in a hexagon with side-length ratio (bottom side to bottom-left side) $R$. Item ii of the introduction section I defines the BCs and crossing events for Potts model spin clusters. The hexagon to the right of each graph shows the $\varsigma$th crossing event whose probability $P_{\varsigma|1}$ appears in the graph. In it, all boundary clusters are colored gray, sides touching a boundary cluster are fixed, and all fixed sides, with their endpoints joined together by dashed exterior arcs, are mutually wired together.

(24) for crossing probability in rectangles and hexagons. For $Q = 4$, the relative errors are significantly greater, on the order of $10^{-1}$. These greater relative errors exemplify the well-known fact [90, 91] that the convergence of the discrete $Q = 4$ critical random cluster model and critical Potts model to their continuum limits is much slower than it is for $Q \in \{2, 3\}$. 
To compensate this effect and improve our numerical results would require sampling the model many more times on a much larger lattice. Table 2 also gives the standard deviation of our relative errors from the average relative error. (We omit the same rectangles and hexagons as in the previous paragraph.) Again, for

\begin{figure}[h]
\centering
\includegraphics[width=\textwidth]{figure42.pdf}
\caption{Comparison of computer simulation measurements to our prediction (24), (169)–(171) of FK-cluster crossing probability in a hexagon with side-length ratio (bottom side to bottom-left side) \( R \). Item i of the introduction section 1 defines the BCs and crossing events for random cluster model FK clusters. The hexagon to the right of each graph shows the \( \varsigma \)th crossing event whose probability \( P_{1,\varsigma} \) appears in the graph. In it, all boundary clusters are colored gray, sides touching a boundary cluster are fixed, and all fixed sides, with their endpoints joined together by dashed exterior arcs, are mutually wired together.}
\end{figure}
$69 \in \mathbb{Q}$, these deviations are all small, typically on the order of $10^{-3}$, and in a few cases, $10^{-2}$. For $Q = 4$, the deviations are significantly higher, on the order of $10^{-1}$. This reflects the fact that for $Q = 4$, the deviations between theory and simulation are often much larger for ‘middle’ data points than than they are for ‘extreme end’ data points (figures 41 and 42). Indeed, the extreme end data points correspond to very irregular hexagons whose shapes

Figure 43. Comparison of computer simulation measurements to our prediction (24), (176)–(178) of FK-cluster crossing probability in a hexagon with side-length ratio (bottom side to bottom-left side) $R$. Item i of the introduction section 1 defines the BCs and crossing events for random cluster model FK clusters. The hexagon to the right of each graph shows the $\kappa$th crossing event whose probability $P_{\kappa}$ appears in the graph. In it, all boundary clusters are colored gray, sides touching a boundary cluster are fixed, and fixed sides, each with endpoints joined by a dashed exterior arc, arc wired independently of each other.
Figure 44. Comparison of computer simulation measurements to our prediction (24), (172)–(175) of FK-cluster crossing probability in a hexagon with side-length ratio (bottom side to bottom-left side) $R$. Item i of the introduction section 1 defines the BCs and crossing events for random cluster model FK clusters. The hexagon to the right of each graph shows the $\varsigma$th crossing event whose probability $P_{3|\varsigma}$ appears in the graph. In it, all boundary clusters are colored gray, sides touching a boundary cluster are fixed, and any two fixed sides (one fixed side) with endpoints joined by a dashed exterior arc are mutually wired together (resp. wired independently of other fixed sides). (Figure 45 shows the graph for the second crossing probability $P_{2|3}$.)
approximate an equilateral triangle, with sides (resp. corners) corresponding to the very long (resp. very short) sides of the hexagons. For these hexagons, the crossing probabilities are always near zero or one, so they are easier to measure. (Indeed, if the long sides of the very irregular hexagons are fixed, then the probability that a single boundary cluster joins all of these fixed sides nearly equals one, and the other crossing probabilities nearly equal zero. This is because inside an equilateral triangle with the fixed BC on all three of its sides, the system strongly favors configurations in which those sides are in the same state, if they are not already required to be so via mutual wiring. Because the three sides of the triangle touch each other, their sites are trivially always part of a boundary cluster that joins all three sides. (One should not confuse this with what happens inside a triangle with free BCs on all sides. For example, for critical site percolation on a triangular lattice, a simple duality argument shows that the probability of a cluster connecting all three free sides of a triangle equals 1/2.) On the other hand, if the short sides of the very irregular hexagons are fixed, then the probability that any cluster joins two or more of these fixed sides nearly equals zero, and the probability that no such cluster joins any pair of these sides nearly equals one. This is because inside a triangle with free BCs on all three of its sides, the probability that a single cluster touches all three corners equals zero in the continuum limit.)

8. Summary

The purpose of this article is to generalize previous results on crossing probabilities in planar geometries. Specifically, we predict, via boundary conformal field theory (CFT) the formula (24) for the probability $\varpi_{\vartheta} \mid P$ of the $\vartheta$th crossing event in the critical random cluster model (where crossing paths are part of boundary FK clusters), the critical Potts model (where crossing paths are part of boundary spin clusters), and the loop-gas model (where the crossing paths follow ‘boundary loops’) on a lattice inside a $2N$-sided polygon $P$, conditioned on a side-alternating boundary condition (BC). Our prediction is exact in the continuum limit. In the formula (24) for $P_{\vartheta} \mid P$, $\vartheta \in \{1, 2, \ldots, C_N\}$ indexes the $C_N$ possible crossing patterns (figure 3), where $C_N$ is the $N$th Catalan number (3), and $\vartheta \in \{1, 2, \ldots, C_N\}$ indexes the side-alternating BCs that we exclusively consider in this article. Items i–iii of the introduction section

Figure 45. Comparison of computer simulation measurements to our prediction (24), (173) of FK-cluster crossing probability $P_{23}$ in a hexagon with side-length ratio (bottom side to bottom-left side) $R$. See the caption of figure 44 for further description.
explain the crossing events and the side-alternating BCs for each model, and section 3 gives a full exposition.

Formula (24) is expressed in terms of these four quantities: the crossing weight $\Pi_\varsigma$ (which, in the context of multiple SLE$_\kappa$, is also called the ‘pure SLE$_\kappa$ partition function’), the Coulomb gas (contour integral) function $\mathcal{F}_\mathrm{g}$, which arises from the solution of the partial differential equations (PDE) describing the affiliated CFT correlation functions (see the discussion of section 5 starting around (82)), the number of closed loops $l_{\varsigma,\vartheta}$ in the $(\varsigma, \vartheta)$th product diagram (see the discussion around (51) for the definition and meaning of $l_{\varsigma,\vartheta}$), and the loop fugacity $n(\kappa)$ of the related loop-gas model (item i of section 2), parameterized by the SLE$_\kappa$ parameter $\kappa \in (8/3, 8)$.

This article begins with an introduction section 1, which surveys the large body of research and recent advances on crossing probabilities for the models of interest. Section 2, which is self-contained, gives a detailed explanation of the formula (24) (but does not derive it), and section 6 contains its application to rectangles, reproducing the rectangle crossing probability for critical percolation (2), [22, 24], the critical $Q = 2$ random cluster (i.e. FK-Ising) model (5) [71], and the critical $Q = 2$ Potts (i.e. Ising) model (6) [26–30]. Section 3 gives formal expressions for the discrete random cluster model and Potts model partition functions. We consider the case of the free BC first, and then sections 3.2 and 3.3 treat, respectively, the random cluster model with side-alternating BCs and crossing events in it; then subsection 3.4 considers the Potts model. In section 3, we give formal expressions (48), (56) for the crossing probabilities $P_{\varsigma,\vartheta|\varsigma}'$ and $P_{\varsigma,\vartheta|\varsigma}'$ of the random cluster model and the Potts model respectively, inside $\mathcal{P}$. The former involves the number $C_{\varsigma,\vartheta}$ of independent boundary clusters (i.e. clusters that touch a fixed side of $\mathcal{P}$) inside $\mathcal{P}$, and using basic graph theory, we find an equation (51) that relates $C_{\varsigma,\vartheta}$ to the number $l_{\varsigma,\vartheta}$ of closed loops formed by the intersection of external and internal arcs in the $(\varsigma, \vartheta)$th product diagram (figure 20). We use this last result in section 4 to express the random cluster model partition function as a certain partition function of a loop-gas model.

Section 4 begins with a definition of the loop-gas model on a lattice inside $\mathcal{P}$ and its partition functions $Z_{\vartheta}^P$ and $Z_{\varsigma,\vartheta}^P$, summing exclusively over events that we interpret as analogous to crossing and side-alternating BC events of the random cluster model. More precisely, for the loop-gas model, the $\varsigma$th crossing event is the event that certain loops, called ‘boundary loops’, join the vertices of $\mathcal{P}$ through its interior in the $\varsigma$th connectivity (corresponding to the $\varsigma$th crossing pattern, see figure 16), and the $\vartheta$th side-alternating BC event is the event that those same loops join the vertices of $\mathcal{P}$ through its exterior in the $\vartheta$th connectivity. Next, we show that for $Q \in \{1, 2, 3, 4\}$, these loop-gas model partition functions are asymptotic, in the continuum limit, to partition functions of random cluster models and Potts models (79) summing exclusively over certain side-alternating BC events. Finally, we give a formal expression (81) for the probability $P_{\varsigma|\varsigma}$ of the $\varsigma$th loop-gas model crossing event conditioned on the $\varsigma$th loop-gas model side-alternating BC event in $\mathcal{P}$. In the continuum limit, this probability equals the corresponding crossing probabilities of the random cluster model and the Potts model, discussed above, when evaluated at appropriate values of $\kappa$.

The main purpose of section 5 is to combine results from previous sections with CFT arguments to derive the crossing probability formula (24), which is the main result of this article. First, we relate the loop-gas model partition functions $Z_{\vartheta}^P$, in the continuum limit, to the boundary CFT correlation function $\Upsilon_{\vartheta}^P$ (84), known to satisfy a particular system of
2N + 3 PDEs. The series of articles [36–39] determines the dimension of, as well as explicit formulas for all elements of, the system’s solution space $S_N$. (As part of the definition for $S_N$, we also require that the solutions are dominated by a certain product (90) of power functions.) Then just above (94), we identify the continuum limit of $Z_{v_0}$ with the ‘crossing weight’ $\Pi_\vartheta$ (or, equivalently, the ‘connectivity weight’, [39, 40] or the ‘pure SLE$\kappa$ partition function’ [26, 81], see the remarks just below (110)). This identification is the main step that leads to our prediction of the explicit formula (24) for polygon crossing probability. The crossing weights are elements of $S_N$, linearly related to the correlation function $\Upsilon\vartheta$ through the meander matrix $M_N$ [38, 39, 82–85]. This fact and the realization that $T_\vartheta = F_\vartheta$ (113) combine to determine explicit formulas for all crossing weights and thus, for all crossing probabilities (24) of every models considered in this article, in any 2N-sided polygon $P$. These probabilities are conditioned on the BC events of items i–iii in the introduction section 1, and they are conformally invariant. Finally, the discussion surrounding (115) interprets this formula as an instance of the conjectured formula for the probability of curve connectivity events in multiple SLE$\kappa$ [86–89].

To support our predictions, we present high-resolution computer simulation results in section 7. The simulations measure crossing probabilities of the $Q \in \{2, 3, 4\}$ random cluster (resp. Potts) model on a large square lattice inside a rectangle and on a large triangular lattice inside a hexagon, using the side-alternating BCs of item i (resp. item ii) in the introduction section 1. We compare these measurements with our prediction (24), finding very good agreement for $Q \in \{2, 3\}$ and reasonable agreement for $Q = 4$. (Similar results for the $Q = 1$ critical random cluster model, i.e. critical percolation, have already appeared in [22, 92].)

Three appendices accompany this article. In appendix A, we use CFT ‘corner operators’ to regularize the conformal transformation of the boundary CFT correlation functions in section 5 from the upper half-plane onto the polygon $P$. The need for regularization arises because this transformation is not conformal at the corners of $P$. In appendix B, we study random cluster model partition functions that sum exclusively over ‘color schemes’, or side-alternating BC events where every other side of $P$ is fixed to a particular color while the other sides are free. Such ‘color scheme partition functions’ are natural because any partition function that sums over a free/fixed side-alternating BC event is necessarily a sum of the former. However, instead of using color scheme partition functions here, we mainly use ‘basic partition functions’, which correspond to simpler loop-gas partition functions in the continuum limit. In appendix B, we show that in the continuum limit and for $Q \in \{1, 2, 3\}$, the span of all color scheme partition functions equals the span of all basic partition functions, as functions of the shape of $P$. In appendix C, we give some simpler formulas for the Coulomb gas function $F_\vartheta(\kappa)$ (11) appearing in the denominator of the crossing probability formula (24). They include an alternative formula (C.2) used in [38, 39] for all $\kappa > 0$ and explicit algebraic (i.e. without contour integrals) formulas for $F_\vartheta(\kappa)$ that hold only at certain rational values of $\kappa$. This latter result leads to a rigorous if indirect evaluation of some rather complicated nested Coulomb gas (contour) integrals.

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Appendix A. Transformation to the polygon via CFT corner operators

The smeared loop-gas partition functions $Z_{\text{g}}(\{\epsilon_j\})$ and $Z_{\text{g},\theta}(\{\epsilon_j\})$ (and the related crossing probabilities $P_{\theta}(92)$) are expected to be invariant with respect to conformal transformations of the system domains. At the same time, the universal part, $\Upsilon_{\text{g}}$ and $\Upsilon_{\text{g},\theta}$ respectively, of their asymptotic behavior (83), (91) as $\epsilon_j \to 0$ is expected to be covariant with respect to these transformations. In this appendix, we explain how the two different transformation laws for these related functions are consistent with each other, and we use the CFT corner operator approach to ‘regularize’ the conformal covariance law for $\Upsilon_{\text{g}}$ and $\Upsilon_{\text{g},\theta}$ at corner points on the boundary of the transformed system domain.

We consider a loop-gas model on a lattice inside a Jordan domain $D \subset \mathbb{C}$, and we let $f$ be a conformal bijection from the upper half-plane $\mathbb{H}$ onto $D$ whose continuous extension to the real axis is conformal at $2N$ marked boundary points $x_1 < x_2 \ldots x_{2N}$. As usual, we require that $N$ boundary loops exit and then re-enter $D$ through its boundary points $w_j := f(x_j)$, with exactly one boundary loop passing once within distance $\delta_j$ from $w_j$. Identical to the case (82) where $D = \mathbb{H}$, we expect the asymptotic behavior

$$
\lim_{\alpha \to 0} Z_{\alpha}^{D}(\{w_j\}; \{\delta_j\})/Z_{\alpha}^{D} = \delta_{\alpha}^{\theta_1} \delta_{\alpha}^{\theta_2} \cdots \delta_{\alpha}^{\theta_{2N}} \Upsilon_{\alpha}^{D}(\{w_j\}),
$$

(A.1)

where $\Upsilon_{\alpha}^{D}$ is some universal function. In (A.1), the superscript $D$ indicates the system domain. (Also, if $D = \mathbb{H}$, then we do not include this superscript, and (A.1) with $w_j \to x_j$ and $\delta_j \to \epsilon_j$ reduces to (82).)

Now we examine how to transform these partition functions and universal functions to their half-plane versions via $f$. The supposition that the partition functions of (A.1) are conformally invariant amounts to

$$
Z_{\alpha}^{D}(\{w_j\}; \{\delta_j\}) = Z_{\alpha}(\{x_j\}; \{\epsilon_j\}),
$$

(A.2)

where $\delta_j = |\partial f(x_j)| \epsilon_j$. (Indeed, under a conformal map, the radius $\epsilon_j$ of the ball centered on $x_j$ dilates by a factor of $|\partial f(x_j)|$. Meanwhile, after we identify the universal function $\Upsilon_{\alpha}^{D}$ with the boundary CFT correlation function

$$
\Upsilon_{\alpha}^{D}(w_1, w_2, \ldots, w_{2N}) = \langle \psi_1(w_1) \psi_1(w_2) \cdots \psi_1(w_{2N}) \rangle_{\alpha}^{D},
$$

(A.3)

this universal function inherits the conformal covariance transformation law of its associated correlation function. By this, we mean that for any conformal bijection $f: \mathbb{H} \to D$ (and with $\theta_1$ (86) the conformal weight of $\psi_1$), we have

$$
\Upsilon_{\alpha}^{D}(\{w_j\}) = |\partial f(x_1)|^{-\theta_1} \cdots |\partial f(x_{2N})|^{-\theta_1} \Upsilon_{\alpha}(\{x_j\}).
$$

(A.4)

These rules (A.2), (A.4) completely determine $Z_{\alpha}^{D}$ and $\Upsilon_{\alpha}^{D}$ in terms of their half-plane versions $Z_{\alpha}$ and $\Upsilon_{\alpha}$ respectively. Now, it is easy to see that the transformation rules (A.2), (A.4) imply one another. Indeed, we have

invariance rule (A.2) for $Z_{\alpha}$ \iff covariance rule (A.4) for $\Upsilon_{\alpha}$.

(A.5)
In this article, the system domain $\mathcal{D}$ is often a 2N-sided polygon $\mathcal{P}$, and here we encounter a subtlety: the continuous extension to the real axis of any conformal bijection $f: \mathbb{H} \to \mathcal{P}$, necessarily a Schwarz–Christoffel transformation

$$ f(z) = A \int_{z_1}^z (x - \zeta)^{\alpha_1/\pi - 1} (x_2 - \zeta)^{\alpha_2/\pi - 1} \cdots (x_{2N} - \zeta)^{\alpha_{2N}/\pi - 1} \, d\zeta + B $$

(A.6)

for some $A, B \in \mathbb{C}, x_1 < x_2 < \ldots < x_{2N}$, and

$$ \alpha_j := \{ \text{interior angle measure at the } j \text{th corner } w_j := f(x_j) \} \in [0, 2\pi), $$

(A.7)

is not conformal at $x_j$ if $\alpha_j \neq \pi$. As such, $\partial f(x_j)$ either vanishes or is infinite, so (A.4) is not applicable. Also, the new radius $\delta_j = |\partial f(x_j)|\epsilon_j$ either vanishes or is infinite, which is physically unacceptable.

To further investigate this singular behavior, we consider the point $\omega_j$ on the $j$th side $[w_j, w_{j+1}]$ of the polygon, offset from the $j$th vertex by a small distance $\epsilon_j > 0$, and its preimage $\xi_j \in \mathbb{R}$:

$$ \omega_j := w_j + \epsilon_j \exp \left( \arg(A) + \sum_{k=1}^j (\pi - \alpha_k) \right) = f(\xi_j). $$

(A.8)

After replacing $w_j \mapsto \omega_j$ and $x_j \mapsto \xi_j$ into (A.4), we send $\epsilon_j \to 0$. Then as the product of derivatives on the right side of (A.4) blows up or vanishes in this limit, so too does $\gamma^{P^D}_\omega(\omega_j)$ on the right side. This is physically unacceptable.

In order to eliminate these singularities from $\gamma^{P^D}_\omega$, which is a physical quantity, we invoke CFT corner operators. To do this, we first note that the asymptotic behavior of the dilation factor $|\partial f(\xi_j)|$ as $\epsilon_j \to 0$ is

$$ |\partial f(\xi_j)| \sim \left( \frac{\alpha_j^{\frac{1}{\pi}}}{\epsilon_j} \right)^{\pi/\alpha_j} \left( \prod_{k=1}^{2N} |x_k - x_j|^{\alpha_k - \pi/\alpha_j} \right). $$

(A.9)

Evidently, (A.9) either blows up or vanishes as $\epsilon_j \to 0$ if $\alpha_j \neq \pi$. To eliminate this singularity, we introduce the following replacements for each $j \in \{1, 2, \ldots, 2N\}$:

$$ \delta_j := |\partial f(x_j)|\epsilon_j \quad \mapsto \quad \delta_j := \epsilon_j \lim_{\epsilon_j \to 0} |\partial f(\xi_j)| \left( \frac{\pi}{\alpha_j} e^{\pi/\alpha_j} \right)^{\delta_j} $$

(A.10)

$$ \gamma^{P^D}_{\delta}(w_1, w_2, \ldots, w_{2N}) \mapsto \gamma^{P^D}_{\delta}(w_1, w_2, \ldots, w_{2N}) = \lim_{\epsilon_1 \to 0} \lim_{\epsilon_2 \to 0} \cdots \lim_{\epsilon_{2N} \to 0} \left( \prod_{k=1}^{2N} \left( \frac{\alpha_k}{\pi} e^{\pi/\alpha_k} \right)^{\delta_k} \right)^{\delta_j} \gamma^{P^D}_{\delta}(\omega_1, \omega_2, \ldots, \omega_{2N}). $$

(A.11)

Without the limit $\epsilon_j \to 0$, the inserted regularization factors of (A.10), (A.11) cancel each other on the right side of (A.1) with $\mathcal{D} = \mathcal{P}$. The advantage is that the new quantities (A.10), (A.11) are finite in the limit $\epsilon_j \to 0$.

We formalize these substitutions (A.10), (A.11) with CFT ‘corner operators’ [80, 115]. If two boundary segments join at an angle with vertex $w \in \mathbb{C}$ and measure $\alpha$ and $\phi_h(w)$ is a primary boundary operator of conformal weight $h$, then

$$ \phi^P_h(w) := \lim_{\epsilon \to 0} \left( \frac{\alpha}{\pi} e^{\pi/\alpha} \right)^{h} \phi_h(w + \epsilon) $$

(A.12)
is the primary corner operator derived from $\phi_x$. (Replacing $w \mapsto \lambda w$ and $\varepsilon \mapsto \lambda \varepsilon$ for any complex $\lambda$ reveals that its conformal weight equals $\pi h/\alpha$.) Using the CFT expression (A.3) for $\Upsilon_\vartheta$ and (A.11), we may write this substitution as

$$\Upsilon_\vartheta^P(w_1, w_2, \ldots, w_{2N}) = (\psi_1(w_1)\psi_2(w_2)\cdots \psi_j(w_{2N}))_\vartheta^P$$

$$\quad \mapsto \tilde{\Upsilon}_\vartheta^P(w_1, w_2, \ldots, w_{2N}) = (\tilde{\psi}_1(w_1)\tilde{\psi}_2(w_2)\cdots \tilde{\psi}_j(w_{2N}))_\vartheta^P. \quad (A.13)$$

To finish, we find an explicit expression for $\tilde{\Upsilon}_\vartheta^P$ if $\mathcal{P}$ is an equiangular $2N$-sided polygon (so all $2N$ interior angles of $\mathcal{P}$ measure $\alpha = (N - 1)\pi/N$). For any $C > 0$, we choose the following parameters for the map (A.6):

$$x_1 = 0, \quad x_j = m_{j-1} \text{ for } j \in \{2, 3, \ldots, 2N - 2\}, \quad x_{2N-1} = 1,$$

$$x_{2N} \to \infty, \quad A = \frac{\alpha}{\pi}C \frac{1}{2N}\sum_{k=1}^{2N-3} e^{-i(\pi - \alpha)} = 0, \quad C > 0, \quad \alpha_j = \frac{(N - 1)\pi}{N} \text{ for all } j \in \{1, 2, \ldots, 2N\}. \quad (A.14)$$

Then the vertices $w_j := f(x_j)$ of $\mathcal{P}$ are such that $w_1 = 0$, $w_2 > 0$, and $w_j \in \mathbb{H}$ for all $j \in \{3, 4, \ldots, 2N\}$, so $\mathcal{P}$ resides in the upper half-plane with its base $[w_1, w_2]$ flush against the positive-real axis. From (A.6) with (A.14), we have

$$\xi_1 \sim (C\varepsilon)^{N(1/N - 1)} \prod_{k=1}^{2N-3} m_k^{1/(N-1)},$$

$$\xi_{2N-1} \sim 1 + (C\varepsilon^{2N-1})^{N(1/N - 1)} \prod_{k=1}^{2N-3} (1 - m_k)^{1/(N-1)},$$

$$\xi_{2N} \sim (C\varepsilon^{2N})^{-N(1/N - 1)}$$

$$\xi_j \sim m_{j-1} + (C\varepsilon^{2N})^{N(1/N - 1)} \prod_{k=1}^{2N-3} (m_k - m_j)^{1/(N-1)} \quad j \in \{2, 3, \ldots, 2N - 2\}, \quad (A.15)$$

as we send $\varepsilon \to 0$. In particular, we find the asymptotic behavior of $\xi_{2N}$ (A.15) after substituting $\xi(t) = \xi_{2N}/t$ into the map (A.6) with the parameters (A.14). This substitution also gives

$$|\partial f(\xi_{2N})| \sim C^{N(1/N - 1)} \frac{e^{2N(1/N - 1)}}{\varepsilon^{2N(1/N - 1)}} \left(\frac{\alpha}{\pi} e^{-\pi/\alpha}\right), \quad \text{where } \alpha = (N - 1)\pi/N. \quad (A.16)$$

After inserting these results (A.15)–(A.16) into (A.11), we find that the regularized universal function $\tilde{\Upsilon}_\vartheta^P$ of the equiangular $2N$-sided polygon $\mathcal{P}$ generated by the conformal bijection $f : \mathbb{H} \to \mathcal{P}$ (A.6) with (A.14) is given by

$$\tilde{\Upsilon}_\vartheta^P(w_1, w_2, \ldots, w_{2N}) \sim \left[\prod_{k=1}^{2N-3} m_k (1 - m_k) \prod_{j<k} (m_j - m_k)\right]^{(1-\delta)/(N-1)}$$

$$\sim \left[C^{N(1-\delta)/(N-1)} \lim_{x \to \infty} x^{\delta/(N-1)} \Upsilon_\vartheta^P(0, m_1, m_2, \ldots, m_{2N-3}, 1, x)\right]. \quad (A.17)$$

(Thanks to its Möbius covariance, we know that this limit (A.17) of $\Upsilon_\vartheta$ exists.) Similarly, we may define a regularized version of the universal function $\Upsilon_{\vartheta, \varphi}$ (91) by replacing the index $\vartheta$ in (A.17) with the double index $\varsigma, \vartheta$. 

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Appendix B. Color scheme partition functions for the random cluster model

In this appendix, we consider the ‘color scheme partition function’ \( W^P_\rho(p, Q) \) defined in the first paragraph of section 3.2. These are natural quantities with which to construct an arbitrary RCB partition function in the critical random cluster model. However, the purpose of this appendix is to show that for most \( Z \in \mathbb{Z}^+ \) and \( Q \in \{1, 2, 3, 4\} \), the color scheme partition functions are linear combinations of the basic partition functions, defined in (44),

\[
X_1^P, X_2^P, \ldots, X_C^P.
\]

(B.1)

Thus, we may work with the latter rather than the former without loss of generality. The latter are more convenient because they converge to natural loop-gas model partition functions in the continuum limit (section 4).

Now obviously, any random cluster model RCB event is a union of (disjoint) color schemes, so any RCB partition function is a sum of color scheme partition functions. Therefore, if we view RCB partition functions as functions of the shape of the polygon \( P \), then

\[
\{ \text{the set of all RCB partition functions} \subseteq \text{span} \{ W^P_\rho(p, Q) | c \text{ a color scheme} \}. \quad (B.2)
\]

The specific purpose of this appendix is to determine for which \( (Q, N) \) we have

\[
\text{span} \{ X_\rho^P(p, Q) | 1 \leq \rho \leq C_N \} = \text{span} \{ W^P_\rho(p, Q) | c \text{ a color scheme} \}
\]

(B.3)

in the continuum limit. If (B.3) is true, then in light of (B.2), restricting our attention to the basic partition functions (B.1) does not result in loss of generality because any RCB partition function is a linear combination of them.

Evidently, (B.2) implies that (B.3) holds if we replace its equality \( = \) with an inclusion \( \subset \).

To show that equality holds for certain \( (Q, N) \), we prove that the dimensions of the spaces in (B.3) agree for these values of \( (Q, N) \) by showing that for these values, we have

\[
\text{rank} \{ X_\rho^P(p, Q) | 1 \leq \rho \leq C_N \} = |\{ W^P_\rho(p, Q) | c \text{ a color scheme} \}|. \quad (B.4)
\]

To find the rank of \( \{ X_\rho^P(p, Q) | 1 \leq \rho \leq C_N \} \) in the continuum limit, we restore the prefactor \( 12^{-\ell+1/2} \) in (66) (omitted from (67), (68), (79) for convenience) to the right side of both left equations in the top line of (77), obtaining

\[
X_\rho^P(p, Q) \sim a_{\rho}^{-\ell+1/2} Z_\rho(Q, \kappa), \quad \text{where } Q = 4 \cos^2(4\pi/\kappa) \text{ and } \kappa \in (4, 8). \quad (B.5)
\]

In order to match previous results in the literature [80] (beneath (B.19)), we insert an extra but physically irrelevant factor of \( Q^{N_C} \) on the right side of (B.5) too. After doing this and recalling (83), (113), we have

\[
X_\rho^P(p, Q) \sim a_{\rho}^{-\ell+1/2} Q^{(N-C_C_1)/2} Z_\rho(Q, \kappa), \quad (B.6)
\]

where \( Z_\rho(Q, \kappa) \sim F_\rho(Q, \kappa) \) as \( \epsilon \downarrow 0 \) for the smeared version of \( Z_\rho(Q, \kappa) \).

From (B.6), we infer that in the continuum limit the rank of \( \{ X_\rho^P(p, Q) | 1 \leq \rho \leq C_N \} \) equals the rank of \( B_N \) (111). Recalling the meander matrix \( M_N \) (23) from section 5, by corollary 7 of [38], we have

\[
\text{rank} \{ X_\rho^P(p, Q) | 1 \leq \rho \leq C_N \} = \text{rank} B_N = \text{rank} M_N. \quad (B.7)
\]
The meander matrix $M_N$ has full rank $C_N$ if $Q = 4$ ($n = 2$, $\kappa = 4$) \cite{82}, so $\text{rank} \ B_N = C_N$ in this case. If $Q \in \{1, 2, 3\}$, then $M_N$ does not have full rank, and we use (53, 54) of \cite{38} and (5.23, 5.24) of \cite{82} to compute its rank. This gives

$$Q = 1 \quad \Rightarrow \quad \kappa = 6, \quad \text{rank} \ B_N = 1 = C_N \quad \text{for} \ N = 1,$$

$$Q = 2 \quad \Rightarrow \quad \kappa = 16/3, \quad \text{rank} \ B_N = 2^{N-1} = C_N \quad \text{for} \ N \leq 2,$$

$$Q = 3 \quad \Rightarrow \quad \kappa = 24/5, \quad \text{rank} \ B_N = (3^{N-1} + 1)/2 = C_N \quad \text{for} \ N \leq 4,$$

$$Q = 4 \quad \Rightarrow \quad \kappa = 4, \quad \text{rank} \ B_N = C_N = C_N \quad \text{for} \ N \in \mathbb{Z}^+.$$

For $Q \in \{1, 2, 3\}$, if $N$ does not satisfy the given condition in (B.8), then $\text{rank} \ B_N < C_N$.

Next, we determine the right side of (B.4). Although there are $Q^N$ color schemes, there are fewer distinct color scheme partition functions, as the latter depend only on differences in color. Indeed, we have (figure B1),

$$W^p_\kappa(p, Q) = W^p_\kappa(p, Q) \{\text{whenever for all } i, j \in \{1, 2, \ldots, N\}, \text{we have } b_i = b_j \text{ if and only if } c_i = c_j.\} \quad (B.9)$$

We say that color schemes $b$ and $c$ are equivalent and write $b \sim c$ if they satisfy the right relation in (B.9), let $[b] = [c]$ denote their common equivalence class, and write $W^p_\kappa(p, Q)$ for $W^p_\kappa(p, Q)$. Then the right side of (B.4) equals

$$\{[W^p_\kappa(p, Q) | c \text{ a color scheme}]\} = s_N(Q) := \left\{\text{number of equivalence classes for color schemes } c = (c_1, c_2, \ldots, c_N) \text{ with } c_i \in \{1, 2, \ldots, Q\}\right\}. \quad (B.10)$$

\begin{figure}[h]
\centering
\includegraphics[width=0.8\textwidth]{figure_B1.png}
\caption{If $Q = 2$ (resp. $Q = 3$), then the four left (resp. five) hexagons show the $s_3(2) = 4$ (resp. $s_3(3) = 5$) \cite{82} ways that we may assign one of $Q$ colors to the odd sides of the hexagon, regarding only differences in color \cite{82} between these sides.}
\end{figure}

\begin{figure}[h]
\centering
\includegraphics[width=0.8\textwidth]{figure_B2.png}
\caption{The upper-left (resp. lower-left) rectangle, with top and bottom sides mutually wired together (resp. wired independently of each other) samples all configurations with these sides having the same color (resp. same color or different colors).}
\end{figure}
Thus, in light of (B.7), in order to determine for what \((Q,N)\) (B.4) holds, we must determine for what \((Q,N)\) we have \(s_N(Q) = \text{rank } B_N\). For this purpose, we let

\[
\#c = \text{the number of distinct colors in color scheme } c, \quad (B.11)
\]

\[
\|c\| = \text{the cardinality of the equivalence class for color scheme } c. \quad (B.12)
\]

By assigning any one of the \(Q\) colors to the first collection of sides with the same color in \(c\), then assigning any one of the \(Q-1\) remaining colors to the second collection of sides with the same color, etc, we find that

\[
\|c\| = Q(Q-1) \cdots (Q-\#c + 1). \quad (B.13)
\]

The total number \(Q^N\) of color schemes obviously equals the sum of the cardinalities (B.13) of all equivalence classes. Because the number of equivalence classes \([c]\) having \(\#c = P\) colors equals \(s_N(P) - s_N(P-1)\), this total is

\[
Q^N = \sum_{P=1}^{Q} (s_N(P) - s_N(P-1))Q(Q-1) \cdots (Q-P + 1), \quad \text{with } s_N(0) = 0. \quad (B.14)
\]

We may use (B.14) to compute \(s_N(Q)\) for all \(Q \in \mathbb{Z}^+\). In this article, we restrict our attention to \(Q \in \{1,2,3,4\}\). For these values, we find

\[
s_N(1) = 1, \quad (B.15)
\]

\[
s_N(2) = 2^{N-1}, \quad (B.16)
\]

\[
s_N(3) = (3^{N-1} + 1)/2, \quad (B.17)
\]

\[
s_N(4) = \left\lfloor (2)(4^{N-2}) + (3)(2^{N-2}) + 1/3. \quad (B.18)
\right\

Upon comparing these values (B.15)–(B.18) with (B.8), we discover that \(s_N(Q) = \text{rank } B_N\) for all \(N \in \mathbb{Z}^+\) if \(Q \in \{1,2,3\}\) and for all \(N \in \{1,2,3\}\) if \(Q = 4\). Thus, (B.3) is true for these values of \((Q,N)\) in \(\{1,2,3,4\} \times \mathbb{Z}^+\):

\[
\span \{X^{P}_{\bar{\vartheta}}(P,Q) | 1 \leq \vartheta \leq C_N = \text{span } \{W^{P}_{c}([P],Q) | c \text{ a color scheme},
\]

\[
\text{if } Q \in \{1,2,3\}, \text{ or if } Q = 4 \text{ and } N \in \{1,2,3\}. \quad (B.19)
\]

Hence, we infer from (B.2), (B.19) that for any \(Q \in \{1,2,3\}\) or for \(Q = 4\) and \(N \in \{1,2,3,4\}\), any RCB partition function is a linear combination of the basic partition functions (B.1), so restricting our attention to the latter, as we do in this article, does not result in loss of generality. On the other hand, if \(Q = 4\) and \(N \geq 4\), then rank \(B_N = C_N < s_N(Q)\) (B.8), (B.15). Hence, not every RCB partition function is necessarily a linear combination of the basic partition functions (B.1), so restricting our attention to the latter, as we do in this article, does result in loss of generality.

We end this appendix by comparing our work for the case \(N = 2\) to similar results in [80]. To begin, we introduce some convenient notation: we write \(c \sim \vartheta \iff c \sim \text{RCB}_\vartheta\) (41). That is,

\[
c = (c_1, c_2, \ldots, c_N) \sim \vartheta \iff \left\{ \begin{array}{l}
\text{if a cluster joins the } \vartheta \text{th and } j \text{th fixed sides} \\
of \text{in } P \text{ in the } \vartheta \text{th connectivity, then } c_i = c_j.
\end{array} \right. \quad (B.20)
\]

\[
\text{Hence, we infer from (B.2), (B.19) that for any } Q \in \{1,2,3\} \text{ or for } Q = 4 \text{ and } N \in \{1,2,3,4\}, \text{any RCB partition function is a linear combination of the basic partition functions (B.1), so restricting our attention to the latter, as we do in this article, does not result in loss of generality. On the other hand, if } Q = 4 \text{ and } N \geq 4, \text{ then rank } B_N = C_N < s_N(Q) \text{ (B.8), (B.15). Hence, not every RCB partition function is necessarily a linear combination of the basic partition functions (B.1), so restricting our attention to the latter, as we do in this article, does result in loss of generality.}
\]

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\text{if a cluster joins the } \vartheta \text{th and } j \text{th fixed sides} \\
of \text{in } P \text{ in the } \vartheta \text{th connectivity, then } c_i = c_j.
\end{array} \right. \quad (B.20)
\]

\[
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\]

\[
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\]

\[
c = (c_1, c_2, \ldots, c_N) \sim \vartheta \iff \left\{ \begin{array}{l}
\text{if a cluster joins the } \vartheta \text{th and } j \text{th fixed sides} \\
of \text{in } P \text{ in the } \vartheta \text{th connectivity, then } c_i = c_j.
\end{array} \right. \quad (B.20)
\]
Furthermore, if \( b, c \in [\ell] \), then \( b \rightsquigarrow \vartheta \) clearly implies \( c \rightsquigarrow \vartheta \), so we simply write \([\ell] \rightsquigarrow \vartheta\) to mean all of these relations. Then using \((\mathbf{B.13})\), we may write \((\mathbf{42})\) as
\[
X^\vartheta_{c}(p_{\vartheta}, Q) = \sum_{c \rightsquigarrow \vartheta} W_{c}^{\vartheta}(p_{\vartheta}, Q)
\]
\[= \sum_{c \rightsquigarrow \vartheta} Q(Q - 1) \cdots (Q - \#c + 1)W_{c}^{\vartheta}(p_{\vartheta}, Q).
\]
If \( N = 2 \), then there are two inequivalent color schemes: the top and bottom sides of a rectangle have the same color \((c = (1,1))\), or they have different colors \((c = (1,2))\). The corresponding basic partition functions are \((\text{figure B2})\)

\[
\text{mutual wiring } (\vartheta = 1, \text{ section } 6): \quad X_{c}^{\vartheta} = QW_{1,1}^{\vartheta}, \quad \text{(B.23)}
\]

\[
\text{independent wiring } (\vartheta = 2, \text{ section } 6): \quad X_{c}^{\vartheta} = QW_{1,1}^{\vartheta} + Q(Q - 1)W_{1,2}^{\vartheta}. \quad \text{(B.24)}
\]

Now, we let \( W_{c}(\varepsilon) \) denote the smeared half-plane version of the color scheme partition function \( W_{c}^{\vartheta} \) \((\text{section } 5)\) for \( c = (1,1) \) and \( c = (1,2) \). With the meaning of \( \approx \) given by \((\mathbf{82})\), \((\mathbf{83})\), we identify \( W_{c}(\varepsilon) \) with a universal function \( \Phi_{c} \):
\[W_{c}(\varepsilon) \simeq \Phi_{c} \quad \text{as } \varepsilon \downarrow 0.
\]

With \( \psi_{1}^{\varepsilon}(x) \) \((\text{resp. } \psi_{2}^{\varepsilon}(x))\) denoting a BCC operator that changes a critical random cluster model \( \text{BC} \) from free \((\text{resp. } \text{fixed with color } c \in [1,2,\ldots,Q])\) for \( y < x \) to fixed with color \( c \in [1,2,\ldots,Q] \) \((\text{resp. free})\) for \( y > x \), one writes
\[
\Phi_{c}(x_{1}, x_{2}, x_{3}, \ldots, x_{2N}) = \left\{ \psi_{1}^{\varepsilon}(x_{1})\psi_{1}^{\varepsilon}(x_{2})\psi_{2}^{\varepsilon}(x_{3})\psi_{1}^{\varepsilon}(x_{4}) \cdots \psi_{1}^{\varepsilon}(x_{2N-1})\psi_{1}^{\varepsilon}(x_{2N}) \right\}.
\]

Like \( \psi_{1}(x) \), these BCC operators are primary CFT operators with conformal weights equaling the \((1,2)\) Kac weight \((\mathbf{86})\) from a CFT with central charge \( c \) given by combining \((\mathbf{74})\), \((\mathbf{85})\). Now, \((\mathbf{6})\), \((\mathbf{22})\), \((\mathbf{25})\) combine to give
\[
Q^{(N-6,\vartheta + 1/2)}_{\vartheta}J_{c}(\kappa) = \sum_{c \rightsquigarrow \vartheta} Q(Q - 1) \cdots (Q - \#c + 1)\Phi_{c}(\kappa).
\]

If \( N = 2 \), then the right side of \((\mathbf{27})\) equals the right side of \((\mathbf{23})\), \((\mathbf{24})\) \((\text{with } \Phi_{c}(\varepsilon) \simeq \Phi_{c})\) for \( \vartheta = 1 \) and \( \vartheta = 2 \) respectively, and the left side of \((\mathbf{27})\) equals \( Q^{1/2}\mathcal{F}_{1}(\kappa) \) and \( Q^{1/2}\mathcal{F}_{2}(\kappa) \) for \( \vartheta = 1 \) and \( \vartheta = 2 \) respectively. Hence,
\[
\begin{cases}
Q^{1/2}\mathcal{F}_{1}(\kappa) = Q\Phi_{1,1}(\kappa), \\
Q^{1/2}\mathcal{F}_{2}(\kappa) = Q\Phi_{1,1}(\kappa) + Q(Q - 1)\Phi_{1,2}(\kappa),
\end{cases}
\]
\[\Rightarrow \begin{cases}
\Phi_{1,1}(\kappa) = Q^{-1/2}\mathcal{F}_{1}(\kappa), \\
\Phi_{1,2}(\kappa) = (Q - 1)^{-1}[\mathcal{F}_{2} - Q^{-1/2}\mathcal{F}_{1}](\kappa) = \Pi_{2}(\kappa).
\end{cases}
\]

After inserting the formula \((\mathbf{119})\) for \( \mathcal{F}_{1} \), now expressed in terms of hypergeometric functions with \((\mathbf{120})\), \((\mathbf{121})\), and the formula \((\mathbf{123})\) for \( \Pi_{2} \) into \((\mathbf{28})\), we find these explicit formulas (with \( x_{y} := x_{i} - x_{j} \) and \( \lambda := x_{21}x_{43}/x_{33}x_{42} \)):
\[ \Phi_{1,1}(\kappa \mid x_1, x_2, x_3, x_4) = n(\kappa)[x_3 x_4]^{1-6/\kappa} \lambda^{2/\kappa}(1 - \lambda)^{1-6/\kappa} \]
\[ \mathcal{F}_1 \left( \frac{4}{\kappa}, 1 - \frac{4}{\kappa}; 2 - \frac{8}{\kappa} \mid 1 - \lambda \right). \]  
(B.29)

\[ \Phi_{1,2}(\kappa \mid x_1, x_2, x_3, x_4) = \frac{\Gamma(12/\kappa - 1)\Gamma(4/\kappa)}{\Gamma(8/\kappa)\Gamma(8/\kappa - 1)}[x_3 x_4]^{1-6/\kappa} \lambda^{1-6/\kappa}(1 - \lambda)^{2/\kappa} \mathcal{F}_1 \left( \frac{4}{\kappa}, 1 - \frac{4}{\kappa}; \frac{8}{\kappa} \mid 1 - \lambda \right). \]  
(B.30)

In [80], Cardy gives formulas for \( \Phi_{1,1} \) and \( \Phi_{1,2} \). They are respectively (50, 51) of that article. In order to cast them in our notation, we must set, in (50, 51) of [80],

\[ Z_{aa} = \Phi_{1,1}, \quad Z_{ab} = \Phi_{1,2}, \]

\[ x = \theta_1 := (6 - \kappa)/2\kappa, \quad \eta = \lambda, \quad \zeta = \frac{1}{1 + 4x}. \]  
(B.31)

After applying the identity \( \mathcal{F}_1(a, b; c \mid z) := (1 - z)^{c-a-b} \mathcal{F}_1(c - a, c - b; c \mid z) \) and the Euler reflection formula for the Gamma function, Cardy’s formula for \( Z_{aa} \) matches our formula (B.29) for \( \Phi_{1,1} \), and (after correcting what seems to be a typo in the factor \( \Gamma(1 + x) \) of \( Z_{ab} \) to \( \Gamma(1 + 4x) \)), Cardy’s formula for \( Z_{ab} \) matches our formula (B.30) for \( \Phi_{1,2} \).

**Appendix C. Alternative formulas for Coulomb gas functions**

In this appendix, we present some simpler formulas for the Coulomb gas function \( F(x) \) (11), and we obtain algebraic formulas for \( F \) in certain cases.

Now, there are two types of simpler formulas for the Coulomb gas function: the alternative formula presented in appendix C.1 agrees with (11) for all \( \kappa > 0 \) (we thank Kytölä and Peltola for sharing this fact with us), and the alternative formulas presented in appendix C.2 hold only at special values of \( \kappa \).

In principle, a direct proof of equality between these different formulas entails performing some of the integrations in (11). However, because this integration appears to be impossible to do in closed form for most values of \( \kappa \), we prove equality indirectly via these two steps:

(i) We show that the functions of either formula are elements of \( S_N \), that is, they satisfy (87), (88), (90).

(ii) We show that the images of these functions under the following vector space isomorphism [38] are equal:

\[ v : S_N \rightarrow \mathbb{R}^{C_N}, \quad v(F) = [\mathcal{L} F], \text{ for all } \zeta \in \{1, 2, \ldots, C_N\}. \]  
(C.1)

(Here, \( [\mathcal{L} F] \) is an equivalence class of a sequence of limits, defined somewhat vaguely in (104). See definitions 9 and 11 of [36] to make this precise.) Thanks to item ii, the desired equality follows from item 4 of theorem 8 in [38]. These results lead to algebraic expressions (C.9), (C.12), (C.15) for the contour integral in (11) at special values of \( \kappa \).

Parts of this appendix assume familiarity with the results of [36–39], rigorously proven only for \( \kappa \in (0, 8) \), and the required main results are the following: the collection \( S_N \) of all functions satisfying (87), (88), (90) has dimension \( C_N \) (3), and the collection \( B_N \subset S_N \) (111) of
C.1. An alternative formula for $F_\vartheta$ (11)

With $x_1 < x_2 < \ldots < x_N$ an ordered collection of points on the real axis and $x := (x_1, x_2, \ldots, x_N)$, [38] defines a Coulomb gas function $F_{c,\vartheta}(\kappa \mid x)$ for each $c \in \{1, 2, \ldots, 2N\}$, $\vartheta \in \{1, 2, \ldots, C_N\}$, and $\kappa > 0$ by the explicit formula

$$
F_{c,\vartheta}(\kappa \mid x) = n(\kappa) \left[ \frac{n(\kappa)\Gamma(2 - 8/\kappa)}{4 \sin^2(4\pi/\kappa)\Gamma(1 - 4/\kappa)^2} \right]^{N-1} \prod_{j<k} (x_j - x_k)^{2\kappa} \\
\left[ \prod_{k=e}^{2N} |x_k - x_e|^{6\kappa} \right] \prod_{j \neq k}^N du_1 \cdots \prod_{j \neq k}^N du_{N-1} \left[ \prod_{j<k} (u_j - u_k)^{8\kappa} \right] \\
\left[ \prod_{i=1}^{2N} \prod_{m=1}^{N-1} (x_i - u_m)^{-4\kappa} \right] \left[ \prod_{m=1}^{N-1} (x_e - u_m)^{2\kappa - 2} \right].
$$

(C.2)

where $n(\kappa)$ is given by (10), $\Gamma_m$ is the Pochhammer contour $\mathcal{P}(x_j, x_j)$ (figure 8) that shares its endpoints $x_j$ and $x_{j'}$ with the $m$th arc of the $i$th connectivity on $x_1 < x_2 < \ldots < x_N$ (item ii of section 2), and no contour shares its endpoints with the $N$th arc of this connectivity, which has an endpoint at $x_e$. Also, the symbol $\mathcal{N}$ selects the branch of the logarithm for each power function in its integrand so $F_{c,\vartheta}$ is real-valued for $\kappa > 0$ [38]. Last, if $x_1 < x_j < x_k < x_e$ then the contour $\mathcal{P}(x_j, x_k)$ arcs over $x_j$ and $x_k$ (and $\mathcal{P}(x_j, x_1)$ if present) in the upper half-plane, and we reverse the orientation of $\mathcal{P}(x_j, x_k)$ if this contour arcs over $x_e$. (See definition 4 and figure 5 of [38] for further elaboration.)

The Coulomb gas function $F_{c,\vartheta}$ is an example of a solution to the system (87), (88) that we produce via the Coulomb gas formalism of CFT [34, 35, 113, 114]. This method invokes a ‘charge neutrality condition’, which produces the distinguished point $x_1$ in the formula (C.2). However, corollary 9 of [38] implies that the distinguished point is, in fact, not so special. Indeed, the corollary states that $F_{c,\vartheta} = F_{c',\vartheta}$ for all $c, c' \in \{1, 2, \ldots, 2N\}$, $\vartheta \in \{1, 2, \ldots, C_N\}$, and $\kappa \in (0, 8)$. In light of this fact, we drop the subscript $c$ and denote this Coulomb gas function by $F_{\vartheta}$.

This independence of (C.2) on our choice of $c$ suggests that an alternative formula for $F_{\vartheta}$, one in which all points $x_i$ appear with equal footing, may exist. Indeed, it does, and it is given by (11). While this formula has greater aesthetic appeal and appears more compact, it is actually more complicated, with a longer integrand and one more integration contour than (C.2).

For this reason, we use (C.2) over (11) as our formula for $F_{\vartheta}$ in most practical situations.

The CFT Coulomb gas formalism also produces the more complicated, yet more symmetric-appearing, formula (11) if we replace the charge neutrality condition with the condition that the total charge of the associated CFT Coulomb gas correlation function equals zero instead. Forsaking the charge neutrality condition puts us at risk of generating functions that, while satisfying the null-state PDEs (87) and translation invariance, violate the last two conformal Ward identities of (88). Somewhat surprisingly, this violation does not occur if we
require a zero total charge; the functions thus generated (11) still satisfy the entire system (87), (88).

Now we move to the main purpose of this section: to rigorously prove that (C.2), (11) are different formulas for the same function $F_\theta$ for all $\kappa > 0$. The proof invokes the two steps i, ii described in the introduction of this appendix:

(i) First, we show that the functions given by (C.2), (11) are elements of $S_N$. In section V of [30], Dubédat proves this claim for the first formula (C.2), and as mentioned above, [119] proves this claim for the second formula (11).

Before moving to the next step, we offer what is perhaps a simpler proof, using the results of [38], that the function $F_\theta$ defined by the second formula (11) is an element of $S_N$ if $\kappa \in (0, 8)$. The idea is to promote $F_\theta$ to another Coulomb gas function $E_\theta' \in S_{N+1}$ by inserting two new points $(x_{2N+1}, x_{2N+2})$ with $x_{2N} < x_{2N+1} < x_{2N+2}$. Here, the subscript $\theta'$ indicates the $\theta'$th connectivity, obtained from the $\theta$th by joining the two inserted end-points $x_{2N+1} < x_{2N+2}$ with an arc. Then, we define $E_\theta'$ by the first formula (C.2) with $N \rightarrow N + 1$ and with $c = 2N + 2$ (so no integration contour of $E_\theta'$ surrounds $x_{2N+2}$ or its partner $x_{2N+1}$) (figure C1).

Next, we consider this limit:

$$\lim_{x_{2N+2} \rightarrow x_{2N+1}} (x_{2N+2} - x_{2N+1})^{6/\kappa - 1} E_\theta'(\kappa | x_1, x_2, \ldots, x_{2N+2}). \quad (C.3)$$

We may obtain a closed expression for this limit (C.3) from the explicit formula of $E_\theta'$, and this expression appears to depend on $x_{2N+1}$. But according to lemma 4 of [36], because $E_\theta' \in S_{N+1}$, this limit (C.3) is actually independent of $x_{2N+1}$. Therefore, sending $x_{2N+1} \rightarrow \infty$ after taking the limit in (C.3) does not alter that function. After taking this second limit, we find that

$$(C.3) = \lim_{x_{2N+2} \rightarrow \infty} \lim_{x_{2N+1} \rightarrow x_{2N+1}} (x_{2N+2} - x_{2N+1})^{6/\kappa - 1} E_\theta'(\kappa | x_1, x_2, \ldots, x_{2N+2})$$

$$= F_\theta(\kappa | x_1, x_2, \ldots, x_N) \quad (11). \quad (C.4)$$

That is, we recover $F_\theta$ (11) from the limit (C.3) of $E_\theta' \in S_{N+1}$ after sending $x_{2N+1} \rightarrow \infty$ in the latter. Furthermore, lemma 5 of [36] says that the limit (C.3) is an element of $S_N$. In light of this fact, it follows from (C.4) that the function $F_\theta$ defined by the second formula (11) is an element of $S_N$ if $\kappa \in (0, 8)$.

(ii) Next, we prove that (C.2) equals (11) if $\kappa \in (0, 8)$ by showing that their images are identical under the isomorphism (C.1). We compute $\nu(F_\theta)$, with $F_\theta$ given by the first formula (C.2), in the proof of lemma 6 in [38], finding (figure 19)
By exactly repeating the analysis of that proof, we prove that (C.5) is also true if \( F_0 \) is given by the second formula (C.2). (There are a few small changes that do not significantly affect the analysis. Case 3 of the proof does not arise. Also, the power \( 12\kappa - 2 \), is present in the integrand of the first formula (C.2) but absent from the integrand of the second formula (11). And finally, infinity is not a branch point of the integrand of the first formula (C.2) but is of the integrand of the second formula (11).) Hence, because \( \nu \) is an isomorphism, it follows that (C.2) and (11) are different formulas for the same function \( F_0(\kappa) \) if \( \kappa \in (0, 8) \). Finally, analytic continuation extends this equality to all \( \kappa > 0 \).

C.2. Algebraic formulas for \( F_0(\kappa) \) at some special values of \( \kappa \)

If the SLE\(_\kappa\) parameter \( \kappa \) is one of the values related to the \( Q \)-color critical random cluster model with \( Q \in \{1, 2, 4\} \) (74) or the \( Q \)-state Potts model with \( Q \in \{2, 4\} \) (75), then \( F_0(\kappa) \) (11) has an algebraic (i.e. without contour integrals) formula for all \( \vartheta \in \{1, 2, \ldots, C_N\} \). In the \( Q = 4 \) case, we may obtain such a formula by explicit integration, but in the \( Q \in \{1, 2\} \) cases, it is not evident that we can explicitly perform the integration. Instead, to prove that \( F_0(\kappa) \) has an algebraic formula, we observe that both \( F_0(\kappa) \) and the formula satisfy the same system of PDEs (87), (88) and the same boundary conditions. From these facts and theorem 8 of [38], we infer that \( F_0(\kappa) \) equals the algebraic formula. Some of these algebraic formulas already appear in the literature (\( Q = 2 \) [81, 120]), although not in direct connection with \( F_0(\kappa) \), and others are, to our knowledge, completely new (\( Q \in \{1, 4\} \)). The \( Q = 3 \) case is ostensibly absent, and we do not know if algebraic formulas for \( F_0(\kappa) \) (11) exist in this case. At least for \( N = 2 \), the hypergeometric expressions (120), (121) for \( F_0(\kappa) \) and \( F_2(\kappa) \) seem to not have such formulas if \( Q = 3 \) (so \( \kappa = 24/5 \) (74) or \( 10/3 \) (75)). For \( Q \in \{1, 2\} \), our results in this section, while rigorous, are obtained in an indirect way, using previous results in [36–39] and other results in the literature [81, 120]. Thus we obtain algebraic formulas for some rather complicated contour integrals based on a full analysis of the solutions of the system of PDEs (87), (88) that \( F_0(\kappa) \) satisfies.

(i) \( Q = 1 \) critical random cluster model (i.e. critical percolation) \( (\kappa = 6) \): Here, we have \( n(\kappa) = 1 \) (10). As such, (112) gives

\[
[\kappa] F_0(\kappa = 6) = 1 \quad \text{for all } \vartheta \in \{1, 2, \ldots, C_N\}
\]

\[
\Rightarrow \quad F_0(\kappa = 6) = [\Pi_1 + \Pi_2 + \cdots + \Pi_{C_N}](\kappa = 6) \quad \text{for all } \vartheta \in \{1, 2, \ldots, C_N\}.
\]

Moreover, it is trivial to see that \( 1 \in S_N \) (i.e. satisfies (87), (88), (90)) if \( \kappa = 6 \) (so \( \theta_1 = 0 \) (86)) and \( [\kappa] \) is an identity for all \( \vartheta \in \{1, 2, \ldots, C_N\} \), so \( 1 = [\Pi_1 + \Pi_2 + \cdots + \Pi_{C_N}](\kappa = 6) \). Thus, (C.6) implies that

\[
F_0(\kappa = 6) = 1
\]

\[
= [\Pi_1 + \Pi_2 + \cdots + \Pi_{C_N}](\kappa = 6), \quad \text{for all } \vartheta \in \{1, 2, \ldots, C_N\},
\]

which is our algebraic result for the \( Q = 1 \) critical random cluster model.

We may use (C.7) to indirectly evaluate the Coulomb gas integral in (11) with \( \kappa = 6 \). Indeed, after inserting the formula (11) for \( F_0(\kappa = 6) \) into (C.7) (and replacing Pochhammer contours with simple contours as in (9), we encounter this algebraic formula for the definite integral in (11) (item (iii)e of section 2 defines \( \mathcal{N} \) …):
\[
\int_{\Gamma_1} \ldots \int_{\Gamma_N} d\nu_N \ldots d\nu_2 \int_{\Gamma_1} d\nu_1 \mathcal{N} \left[ \left( \prod_{i=1}^{2N} \prod_{m=1}^{N} (x_i - u_m)^{-2/3} \right) \left( \prod_{p < q}^{N} (u_p - u_q)^{1/3} \right) \right] = \frac{\Gamma(1/3)2^{2N}}{\Gamma(2/3)^N} \prod_{j<k}^{N} (x_k - x_j)^{-1/3} .
\]

In (C.9), the integration contours are nonintersecting simple contours in the upper half-plane that join their endpoints, all of them among \(x_1, x_2, \ldots, x_{2N}\), pairwise in the \(\nu\)th connectivity for \(\vartheta \in \{1, 2, \ldots, C_\nu\}\) (item ii of section 2). Thus, using results from [36–39], we have indirectly evaluated the Coulomb gas integral in (C.9).

(ii) \(Q = 2\) critical Potts model (i.e. critical Ising model) (\(\kappa = 3\)): Here, we again have \(\nu(\kappa) = 1\) (10). As such, (112) gives (C.6) once again, but with \(\kappa = 3\) instead. Now, we have

\[
\mathcal{F}_{\vartheta}(\kappa = 3 | x_1, x_2, \ldots, x_{2N}) = \sum_{\lambda} \text{sgn } \lambda \left( \prod_{(i,j) \in \lambda} \frac{1}{x_i - x_j} \right) = \left[ \prod_1 + \prod_2 + \cdots + \prod_{C_\lambda}(\kappa = 3 | x_1, x_2, \ldots, x_{2N}) \right]
\]

for all \(\vartheta \in \{1, 2, \ldots, C_\nu\}\), where the sum in (C.10) is over all mutually exclusive pairings \(\lambda\) of indices in the set \(\{1, 2, \ldots, 2N\}\), and where \(\text{sgn } \lambda\) is the sign of the product \(\prod (i - k)(j - l)(j - k)(j - l)\) over pairs of distinct \(i, j, k, l \in \lambda\). Indeed, proposition 4.6 of [81] states that the right side of (C.10) satisfies (87), (88) if \(\kappa = 3\), so because it obviously satisfies (90) too, it is an element of \(\Delta_N\). Furthermore, the cascade property in proposition 4.6 of [81] implies that the action of \([\subset]\) on the right side of (C.10) equals one for all \(\vartheta \in \{1, 2, \ldots, C_\nu\}\). Thus, the right sides of (C.10), (C.11) are equal, so because (C.6) is true for \(\kappa = 3\), both sides of (C.10) are equal. This is our algebraic result for the \(Q = 2\) critical Potts model.

We may use (C.10) to indirectly evaluate the Coulomb gas integral in (11) with \(\kappa = 3\). Indeed, after inserting the formula (11) for \(\mathcal{F}_{\vartheta}(\kappa = 3)\) into (C.10), we find (item (iii)e of section 2 defines \(\Lambda_{\vartheta} \ldots \):)

\[
\int_{\Gamma_N} \ldots \int_{\Gamma_2} d\nu_2 \int_{\Gamma_1} d\nu_1 \mathcal{N} \left[ \left( \prod_{i=1}^{2N} \prod_{m=1}^{N} (x_i - u_m)^{-4/3} \right) \left( \prod_{p < q}^{N} (u_p - u_q)^{2/3} \right) \right] = \frac{3\Gamma(-1/3)^2}{\Gamma(-2/3)} \prod_{j<k}^{N} (x_k - x_j)^{-2/3} \sum_{\lambda} \text{sgn } \lambda \left( \prod_{(i,j) \in \lambda} \frac{1}{x_i - x_j} \right).
\]

In (C.12), the integration contours are nonintersecting Pochhammer contours in the upper half-plane, with endpoints among \(x_1, x_2, \ldots, x_{2N}\), and joining these points pairwise in the \(\nu\)th connectivity for any \(\vartheta \in \{1, 2, \ldots, C_\nu\}\). Thus, using results from [36–39], we have indirectly found an algebraic formula for the Coulomb gas integral in (C.12).

(iii) \(Q = 2\) critical random cluster model (\(\kappa = 16/3\)): Here, we have \(\nu(\kappa) = \sqrt{2}\) (10). Moreover, we have that for all \(\vartheta \in \{1, 2, \ldots, C_\nu\}\) and all \(i \in \{1, 2, \ldots, 2N - 1\}\) (see the proof of lemma 6 in [38]),
\[ \mathcal{F}_\theta(\kappa = 16/3 \mid x_1, x_2, \ldots, x_{2N}) \sim (x_i - x_j)^{-1/8} \]

\[
\begin{cases}
\sqrt{2} \mathcal{G}_\theta(\kappa = 16/3 \mid \ldots, x_{i-1}, x_{i+2}, \ldots), & x_i, x_{i+1} \text{ joined in } \vartheta \text{th connectivity,} \\
\mathcal{G}_\omega(\kappa = 16/3 \mid \ldots, x_{i-1}, x_{i+2}, \ldots), & x_i, x_{i+1} \text{ otherwise,}
\end{cases}
\]

where \( \mathcal{G}_\theta, \mathcal{G}_\omega \in \mathcal{B}_{N-1} \), and where we form the \( g \)th (resp. \( \omega \)th) connectivity from the \( \vartheta \)th connectivity by dropping from the latter the one arc with its endpoints at \( x_i \) and \( x_{i+1} \) (resp. detaching in the latter the two arcs with their endpoints at \( x_i \) or \( x_{i+1} \) from those points and joining their dangling ends together to form one arc). (Thus, \( \vartheta, \omega \in \{1, 2, \ldots, C_N-1\} \).)

Now, we have

\[
\mathcal{F}_\theta(\kappa = 16/3 \mid x_1, x_2, \ldots, x_{2N}) = \prod_{j=1 \text{ odd}}^{2N-1} (y_j - y_j')^{-1/8}
\]

\[
\begin{cases}
\sum_{\mu_1=\pm 1} \sum_{\mu_2=\pm 1} \ldots \sum_{\mu_{2N-1}=\pm 1} \prod_{j \leq \text{odd}}^{2N-1} (y_k - y_j)(y_{k+1} - y_j') \mu_{jk}\mu_{1k}^{1/4}
\end{cases}
\]

where \( y_j, y_j' \in \{x_1, x_2, \ldots, x_{2N}\} \) with \( y_j < y_j' \) are the endpoints of the \( j \)th arc in the \( \vartheta \)th connectivity. Indeed, [120] says that for \( \vartheta = C_N \) (so \( y_j = x_j \) for all \( j \in \{1, 2, \ldots, 2N\} \)), the right side of (C.14) satisfies (87), (88) with \( \kappa = 16/3 \). By appropriately permuting the points \( x_j \), which does not change the system (87), (88), we generate the right side of (C.14) for all \( \vartheta < C_N \) and see that it satisfies the system (87), (88) too. Obeying (90) as well, each solution (C.14) is an element of \( \mathcal{S}_N \). So to prove the equality in (C.14), it suffices to show that the left and right sides have identical images under the action of [\( \subset \)] for each \( \zeta \in \{1, 2, \ldots, C_N\} \). This immediately follows from the fact that the right side of (C.14) has the asymptotic behavior (C.13), the same as of \( \mathcal{F}_\theta(\kappa = 16/3) \) (11), as \( x_{i+1} \rightarrow x_i \) for all \( i \in \{1, 2, \ldots, 2N-1\} \).

In addition to \( \mathcal{F}_\mathcal{C}_\theta \), that is (C.14) with \( y_j = x_j \) and \( \vartheta = C_N \), (6) of [120] gives \( 2^{N-1} - 1 \) more solutions to the system (87), (88) with \( \kappa = 16/3 \), and their formulas are identical to (C.14) with \( y_j = x_j \), except that a factor, equaling one or minus one and depending on \( \mu_\{j\} \), multiplies the product in (C.14). We let \( \mathcal{D}_N \) denote the total collection of \( 2^{N-1} \) solutions in (6) of [120]. Because the rank of \( \mathcal{B}_N \) (111), (C.14) with \( \kappa = 16/3 \) (B.8) equals the cardinality of \( \mathcal{D}_N \), we anticipate that these two sets of solutions, both containing \( \mathcal{F}_\mathcal{C}_\theta \), span the same subspace of \( \mathcal{S}_N \). Indeed, this is true. To see why, we note that the solutions of \( \mathcal{D}_N \) are conformal blocks [34]. Thus, span \( \mathcal{D}_N \) is closed under the braid group action that permutes the points \( x_j \). Now for any \( \vartheta < C_N \), there exists a braiding that sends \( \mathcal{F}_\mathcal{C}_\theta \in \mathcal{D}_N \) to \( \mathcal{F}_\theta \in \mathcal{B}_N \) [121]. Thus, \( \mathcal{B}_N \subset \text{span } \mathcal{D}_N \). But because rank \( \mathcal{B}_N = |\mathcal{D}_N| \) (B.8), we have span \( \mathcal{B}_N = \text{span } \mathcal{D}_N \). For \( N \gg 3 \), this span is a proper subspace of \( \mathcal{S}_N \), and we do not know if any solutions in \( \mathcal{S}_N \setminus \text{span } \mathcal{B}_N \) have algebraic formulas too. Equation (C.14) is our algebraic result for the \( Q = 2 \) critical random cluster model.

We may use (C.14) to indirectly evaluate the Coulomb gas integral in (11) with \( \kappa = 16/3 \). Indeed, after inserting the formula (11) for \( \mathcal{F}_\theta \) into (C.14) (and replacing Pochhammer contours with simple contours as in (9)), we obtain this algebraic formula for the definite integral in (11) (item (iii) of section 2 defines \( N \) ? ):
\[
\int_{1^N} du_N \cdots \int_{1^2} du_2 \int_{1^1} du_1 \Lambda \left[ \prod_{i=1}^{2N} \prod_{m=1}^{N} (x_i - u_m)^{\nu_1} \left( \prod_{p < q}^{N} (u_p - u_q)^{\nu_2} \right) \right] \\
= \left( \prod_{j < k} (x_j - x_k)^{3/8} \right) \frac{\Gamma(1/4)2^{2N}}{\left( 2^{1/2} \right)^{2N}} \prod_{j=1 \text{ odd}}^{2N-1} (y_j+1 - y_j)^{-1/8} \\
\left\{ \sum_{\mu_1^{-1}} \sum_{\mu_2^{-1}} \cdots \sum_{\mu_{2N-1}^{-1}} \prod_{j < k}^{2N} \left( y_j - y_k \right) \left( y_k+1 - y_j \right) \left( y_j+1 - y_k \right) \right\}^{1/2}.
\]

(C.15)

In (C.15), the integration contours are nonintersecting simple contours in the upper half-plane, with endpoints among \(x_1, x_2, \ldots, x_{2N}\) and joining these points pairwise in the \(i\)th connectivity, and we define \(y_j\) under (C.14). Thus, using results from [36–39], we have indirectly found an algebraic expression for the Coulomb gas integral in (C.15).

(iv) \(Q = 4\) critical random cluster model and critical Potts model (\(\kappa = 4\), but more generally, we consider all \(\kappa\) such that \(4/\kappa \in \mathbb{Z}^+\)): If \(\kappa = 4\) then \(n(\kappa) = 2\) (10), and each \(x_r\) is a simple pole of the integrand of \(\mathcal{F}_\rho(\kappa = 4)\) (11). More generally, each \(x_r\) is an order-\(r\) pole of the integrand of \(\mathcal{F}_\rho(\kappa)\) (11) if and only if \(\kappa = 4/r\) for \(r \in \mathbb{Z}^+\). In this case, we decompose each Pochhammer contour into a collection of loops that surrounds the contour’s endpoints (figure 8, (34) of [38]), and we perform the integration around these loops via the Cauchy integral formula. We find

\[
\mathcal{F}_\rho(\kappa = 4/r \mid x_1, x_2, \ldots, x_{2N}) = \left( -1 \right)^{r-1} \frac{(r-1)!}{(2r-2)!} N \prod_{i<j} (x_j - x_i)^{r/2} \sum_{\mu_1=1 \atop \mu_2=1}^{2N} \cdots \sum_{\mu_{2N-1}=1}^{2N} \left( \prod_{m=1}^{N} |x_m - u_m|^{-r} \prod_{p < q}^{N} (u_p - u_q)^{2r} \right),
\]

(C.16)

where \(\omega(\mu) := \left\{ \begin{array}{ll} (-1)^{\mu_1 + \mu_2 + \cdots + \mu_r}, & r \text{ odd}, \\ 1, & r \text{ even}. \end{array} \right.\) (C.17)

(In [39], we found an equivalent formula (35) by integrating the alternative formula (C.2) for \(\mathcal{F}_\rho\).) Specializing to \(\kappa = 4\) (\(r = 1\)), this becomes

\[
\mathcal{F}_\rho(\kappa = 4 \mid x_1, x_2, \ldots, x_{2N}) = N \prod_{i<j} (x_j - x_i)^{1/2} \sum_{\mu_1=1}^{2N} \cdots \sum_{\mu_{2N-1}=1}^{2N} \left( \prod_{m=1}^{N} |x_m - x_{\mu_m}|^{-1} \prod_{p < q}^{N} (x_{\mu_p} - x_{\mu_q})^{2} \right).
\]

(C.18)

This is our algebraic result for the \(Q = 4\) critical Potts model and the \(Q = 4\) critical random cluster model.

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