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1996 J. Phys. A: Math. Gen. 29 7013

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Low-density series expansions for directed percolation on square and triangular lattices

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Received 23 May 1996

Abstract. Greatly extended series have been derived for moments of the pair-connectedness for bond and site percolation on the directed square and triangular lattices. The length of the various series has been at least doubled to more than 110 (100) terms for the square-lattice bond (site) problem and more than 55 terms for the bond and site problems on the triangular lattice. Analysis of the series leads to very accurate estimates for the critical parameters and generally seems to rule out simple rational values for the critical exponents. The values of the critical exponents for the average cluster size, parallel and perpendicular connectedness lengths are estimated by $\gamma = 2.277\,69(4)$, $\nu_{\parallel} = 1.733\,825(25)$ and $\nu_{\perp} = 1.096\,844(14)$, respectively. An improved estimate for the percolation probability exponent is obtained from the scaling relation $\beta = (\nu_{\parallel} + \nu_{\perp} - \gamma)/2 = 0.276\,49(4)$. In all cases the leading correction to scaling term is analytic.

1. Introduction

Models exhibiting critical behaviour similar to directed percolation (DP) are encountered in a wide variety of problems such as fluid flow in porous media, Reggeon field theory, chemical reactions, population dynamics, catalysis, epidemics, forest fires, and even galactic evolution. Directed percolation is thus a model of relevance to a very diverse set of physical problems and it is therefore no wonder that it continues to attract a great deal of attention. Furthermore, two-dimensional directed percolation is one of the simplest models which is not translationally invariant and therefore cannot be treated in the framework of conformal field theory [1]. This leaves open a number of fundamental questions about this model. What should one expect an exact solution to look like and more concretely are the critical exponents rational?

In the absence of an exact solution the most powerful method for studying lattice-statistics models is probably that of series expansions. The method of exact series expansions consists of calculating the first few coefficients in the Taylor expansion of various thermodynamic functions, or, in more abstract terms, various moments of some appropriate generating function. Given such a series, highly accurate estimates can be obtained for the critical parameters using differential approximants [2]. In the most favourable cases one can even find an exact expression for the generating function from the first-series coefficients.

Low-density series in the variable p , which is the probability that bonds or sites are present, were first derived by Blease [3], who used a transfer-matrix method to calculate series for the cluster size and other moments of the pair-connectedness of bond percolation

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on directed square and triangular lattices. These series were greatly extended by Essam *et al* [4], who also studied site percolation. They devised a non-nodal graph expansion, which enabled them to calculate twice as many terms correctly from the basic transfer-matrix calculation, and derived the series to order 49 (48) for the square bond (site) problem and to order 25 (26) for the triangular bond (site) problem. These long series resulted in accurate exponent estimates and led to the conjectured critical exponents $\gamma = 41/18$, $\nu_{\perp} = 79/72$, $\nu_{\parallel} = 26/15$, and $\beta = 199/720$ [4].

High-density series for the percolation probability were derived by Blease [3]. The square bond series was greatly extended by Baxter and Guttmann [5] using a superior transfer-matrix method and an extrapolation procedure based on predicting correction terms from successive calculations on finite lattices of increasing size. The analysis of the resulting series conformed to the conjectured fraction for β . This series and the one for the square site problem were recently extended by Jensen and Guttmann [6] who also studied the triangular bond and site problems [7]. The analysis of these extended series yielded more precise exponent estimates. From these estimates they concluded that there are no simple rational fractions whose decimal expansion agrees with the highly accurate estimates of β obtained from the square bond and triangular site series. In particular, the rational fraction suggested by Essam *et al* [4] is incompatible with the estimates.

In this paper I combine an efficient transfer-matrix calculation with the non-nodal graph expansion and the above-mentioned extrapolation method and have been able to more than double the number of series terms for moments of the pair-connectedness. Most of the series have been extended to order 112 for the square bond problem, 106 for the square site problem, 57 for the triangular bond problem and 56 for the triangular site problem. The series were analysed using differential approximants which can accommodate a wide variety of functional features and certainly should be appropriate in this case. The major result of the analysis is that the exact exponent values conjectured by Essam *et al* [4] generally seems to be incompatible with the numerical estimates from the differential approximant analysis.

The remainder of the article is organized as follows. In section 2 I will give further details of the models studied in this paper. Section 3 contains a description of the series-expansion technique with special emphasis on the transfer-matrix calculation (section 3.1) and the extrapolation procedure for the square bond case (section 3.3). Details of the extrapolation procedure for the remaining problems are given in the appendix. Details of the series analysis are given in section 4 and the results are discussed and summarized in section 5.

2. Specification of the models

Domany and Kinzel [8] demonstrated that site and bond percolation on the directed square lattice are special cases of a one-dimensional stochastic cellular automaton in which the preferred direction t is time. DP is thus a model for a simple branching process in which a site x occupied at time t may give rise to zero or one offspring on each of the sites $x \pm 1$ at time $t + 1$. Whether a site (x, t) is occupied or not depends only on the state of its nearest neighbours in the row above. The evolution of the model on the square lattice is therefore governed by the conditional probabilities $P(\sigma_x | \sigma_l, \sigma_r)$, with $\sigma_i = 1$ if site i is occupied and 0 otherwise. These transition probabilities are the probabilities of finding the site (x, t) in state σ_x given that the sites $(x - 1, t - 1)$ and $(x + 1, t - 1)$ were in states σ_l and σ_r , respectively. One has a very free hand in choosing the transition probabilities as long as one respects conservation of probability, $P(1 | \sigma_l, \sigma_r) = 1 - P(0 | \sigma_l, \sigma_r)$. In addition studies have generally been limited to cases in which the transition probabilities are independent

of both x and t . In this paper I restrict my study to the following two cases corresponding to bond and site percolation:

$$P(0|\sigma_l, \sigma_r) = \begin{cases} (1-p)^{\sigma_l+\sigma_r} & \text{bond} \\ (1-p)^{1-(1-\sigma_l)(1-\sigma_r)} & \text{site.} \end{cases} \quad (2.1)$$

On the triangular lattice the model is described by the probabilities $P(\sigma_x|\sigma_l, \sigma_t, \sigma_r)$ of finding the site (x, t) in state σ_x given that the sites $(x-1, t-1)$, $(x, t-2)$, and $(x+1, t-1)$ were in states σ_l , σ_t and σ_r , respectively, and I study the two cases

$$P(0|\sigma_l, \sigma_t, \sigma_r) = \begin{cases} (1-p)^{\sigma_l+\sigma_t+\sigma_r} & \text{bond} \\ (1-p)^{1-(1-\sigma_l)(1-\sigma_t)(1-\sigma_r)} & \text{site.} \end{cases} \quad (2.2)$$

The behaviour of the model is controlled by the branching probability p . When p is smaller than a critical value p_c the branching process eventually dies out and all space-time clusters remain finite. For $p > p_c$ there is a non-zero probability $P(p)$ that the branching process will survive indefinitely. This percolation probability is the order parameter of the process, and close to p_c it vanishes as a power-law:

$$P(p) \propto (p - p_c)^\beta \quad p \rightarrow p_c^+. \quad (2.3)$$

In the low-density phase ($p < p_c$) many quantities of interest can be derived from the pair-connectedness $C_{x,t}(p)$, which is the probability that the site x is occupied at time t given that the origin was occupied at $t = 0$. The moments of the pair-connectedness may be written as

$$\mu_{n,m}(p) = \sum_{t=0}^{\infty} \sum_x x^n t^m C_{x,t}(p). \quad (2.4)$$

Due to symmetry, moments involving odd powers of x vanish. The remaining moments diverge as p approaches the critical point from below:

$$\mu_{n,m}(p) \propto (p_c - p)^{-(\gamma + n\nu_{\perp} + m\nu_{\parallel})} \quad p \rightarrow p_c^-. \quad (2.5)$$

One generally only studies the lower-order moments such as the mean cluster size $S(p) = \mu_{0,0}(p)$, the first parallel moment $\mu_{0,1}(p)$, the second perpendicular moment $\mu_{2,0}(p)$, and the second parallel moment $\mu_{0,2}(p)$.

3. Series expansions

From (2.4) it follows that the first and second moments can be derived from the quantities

$$S(t) = \sum_x C_{x,t}(p) \quad \text{and} \quad X(t) = \sum_x x^2 C_{x,t}(p) \quad (3.1)$$

as

$$S = \sum_{t=0}^{\infty} S(t) \quad \mu_{0,1} = \sum_{t=1}^{\infty} t S(t) \quad \mu_{0,2} = \sum_{t=1}^{\infty} t^2 S(t) \quad \mu_{2,0} = \sum_{t=0}^{\infty} X(t). \quad (3.2)$$

$S(t)$ and $X(t)$ are polynomials in p obtained by summing the pair-connectedness over all lattice sites whose parallel distance from the origin is t . As shown by Essam [9] the pair-connectedness can be expressed as a sum over all graphs formed by taking unions of directed paths connecting the origin to the site (x, t) ,

$$C_{x,t}(p) = \sum_g d(g) p^e \quad (3.3)$$

where e is the number of random elements (bonds or sites) in the graph g . Any directed path to a site whose parallel distance from the origin is t contains at least $m(t)$ steps with $m(t) = t$ for the square lattice and $m(t) = \lfloor (t+1)/2 \rfloor$ (integer division) for the triangular lattice. From this it follows that if $S(t)$ and $X(t)$ have been calculated for $t \leq t_{\max}$ then one can determine the moments to order $m(t_{\max} + 1) - 1$. One can, however, do much better, as demonstrated by Essam *et al* [4]. They used a non-nodal graph expansion, based on work by Bhatti and Essam [10], to extend the series to order $n(t_{\max})$ approximately equal to $2m(t_{\max})$ (the actual order varies a little from problem to problem). Details of this expansion will be given below, but here it will suffice to note that it works by calculating the contributions $S^N(t)$ and $X^N(t)$ (correct to order $n(t)$) of non-nodal graphs to $S(t)$ and $X(t)$ and using the non-nodal expansions to calculate the final series for $S(p)$ and the various moments. Further extensions of the series can be obtained by using a procedure similar to that of Baxter and Guttmann [5]. One looks at correction terms to the series and tries to identify extrapolation formulae for the first n_r correction terms allowing one to derive a further n_r series terms correctly.

The series expansions for moments of the pair-connectedness is thus obtained as follows:

(i) Calculate the polynomials $S(t)$ and $X(t)$ for $t \leq t_{\max}$ using the transfer-matrix technique to an order greater than $n(t_{\max}) + n_r$.

(ii) For each t use the non-nodal graph expansion to calculate $S_t^N = \sum_{t' \leq t} S^N(t')$ and $X_t^N = \sum_{t' \leq t} X^N(t')$ correct to order $n(t)$.

(iii) From the sequences obtained from $S_t^N - S_{t+1}^N = -S^N(t+1)$ and $X_t^N - X_{t+1}^N = -X^N(t+1)$ for $t < t_{\max}$ identify the first n_r correction terms.

(iv) Use these correction terms to extend the series for S^N and X^N to order $n(t_{\max}) + n_r$.

(v) Finally calculate the series for S , $\mu_{0,1}$, $\mu_{0,2}$ and $\mu_{2,0}$ correct to order $n(t_{\max}) + n_r$.

Details of the transfer-matrix technique, non-nodal graph expansion and extrapolation procedure are given in the following sections.

3.1. Transfer-matrix technique

Figure 1 shows the part of the square and triangular lattices which can be reached from the origin O using no more than five steps. Note that, in keeping with the prescription used by Essam *et al* [4], vertical steps on the triangular lattice correspond to incrementing t by two. The calculation of the pair-connectedness is readily turned into an efficient computer algorithm by use of the transfer-matrix technique. From (2.1) and (2.2) one sees that the evaluation of the pair-connectedness involves only local 'interactions' since the

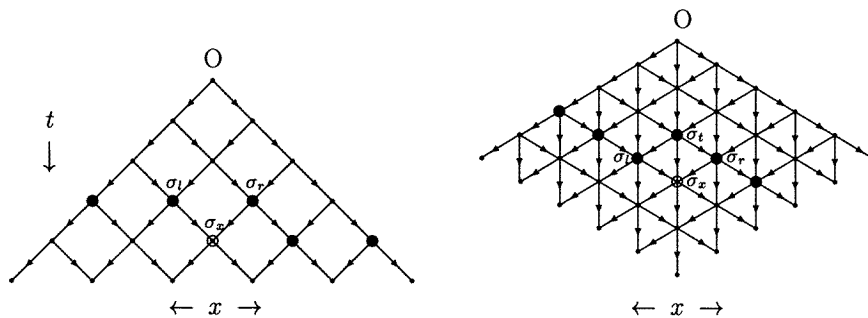


Figure 1. Directed square and triangular lattices with orientation given by the arrows.

transition probabilities depend on neighbouring sites only. The probability of finding a given configuration can therefore be calculated by moving a boundary through the lattice one site at a time. At any given stage this line cuts through a number of, say k , lattice sites thus leading to a total of 2^k possible configurations along this line. Configurations along the boundary line are trivially represented as binary numbers, and the probability of each configuration is represented by a truncated polynomial in p .

Figure 1 shows how the boundary (marked by large filled circles) is moved in order to pick up the weight associated with a given 'face' of the lattice at a position x along the boundary line. On the square lattice the boundary site at σ_r is moved to σ_x and the weight $P(\sigma_x|\sigma_l, \sigma_r)$ is picked up. Similarly on the triangular lattice the boundary site at σ_t is moved to σ_x while picking up the weight $P(\sigma_x|\sigma_l, \sigma_t, \sigma_r)$. In more detail, let $S0 = (\sigma_1, \dots, \sigma_{x-1}, 0, \sigma_{x+1}, \dots, \sigma_k)$ be the configuration of sites along the boundary with 0 at position x and similarly $S1 = (\sigma_1, \dots, \sigma_{x-1}, 1, \sigma_{x+1}, \dots, \sigma_k)$ the configuration with 1 at position x . Then in moving the x 'th site as just described the boundary line polynomials are updated as follows on the square lattice

$$\begin{aligned} P(S0) &= W(0|0, \sigma_l)P(S0) + W(0|1, \sigma_l)P(S1) \\ P(S1) &= W(1|0, \sigma_l)P(S0) + W(1|1, \sigma_l)P(S1) \end{aligned}$$

and as follows on the triangular lattice

$$\begin{aligned} P(S0) &= W(0|\sigma_r, 0, \sigma_l)P(S0) + W(0|\sigma_r, 1, \sigma_l)P(S1) \\ P(S1) &= W(1|\sigma_r, 0, \sigma_l)P(S0) + W(1|\sigma_r, 1, \sigma_l)P(S1). \end{aligned}$$

The pair-connectedness is calculated from the boundary polynomials before the boundary leaves the site by summing over all configurations with a 1 at that site. In practise the data was collected when the boundary reached a horizontal position on the square lattice and a position parallel to the right edge of the triangular lattice. The pair-connectedness is obviously symmetrical in x , $C_{x,t}(p) = C_{-x,t}(p)$, so it suffices to calculate the pair-connectedness for $x \geq 0$. More importantly, due to the directedness of the lattices, if one looks at sites (x, t) with $x \geq 0$ they can never be reached by paths extending onto points (x', t') in the part of the lattice for which $t' > \lfloor t/2 \rfloor$, $x' < -\lfloor t/2 \rfloor$. This effectively means that the pair-connectedness at points with parallel distance t from the origin can be calculated using a boundary which cuts through at most $\lfloor t/2 \rfloor + 1$ sites. Thus the memory (and time) required to derive $S(t)$ and $X(t)$ grows like $2^{\lfloor t/2 \rfloor + 1}$.

For the bond and site problems on the square lattice I was able to calculate the pair-connectedness up to $t_{\max} = 47$ and for the triangular lattice up to $t_{\max} = 45$. Since the integer coefficients occurring in the series expansion become very large the calculation was performed using modular arithmetic [11]. Each run for t_{\max} , using a different prime number, took approximately 12 hours using 64 nodes on an Intel Paragon, and up to eight primes were needed to represent the coefficients correctly. The major limitation of the present calculation was available computer memory rather than time.

3.2. Non-nodal graph expansion

The non-nodal graph expansion has been described in detail in [4] and here I will only summarise the main points and introduce some notation. A graph g is nodal if there is a point (other than the terminal point) through which all paths pass. It is clear that each such nodal point effectively works as a new origin for the cluster growth. This is the essential idea behind the non-nodal graph expansion. $S^N(t)$ is the contribution to $S(t)$ obtained by restricting the sum in (3.3) to non-nodal graphs. The non-nodal expansions are

obtained recursively from the polynomials $S(t)$ and $X(t)$. First one sets $S^N(1) = S(1)$ and $X^N(1) = X(1)$ and then for $2 \leq t \leq t_{\max}$ one calculates $S^N(t)$ and $X^N(t)$ from

$$S^N(t) = S(t) - \sum_{t'=1}^{t-1} S^N(t')S(t-t') \tag{3.4}$$

and

$$X^N(t) = X(t) - \sum_{t'=1}^{t-1} [S^N(t')X(t-t') + X^N(t')S(t-t')]. \tag{3.5}$$

Next form the sums of (3.2) using the truncated non-nodal polynomials $S^N(t)$ and $X^N(t)$ instead of $S(t)$ and $X(t)$. The final series are then obtained from the formulae

$$S = 1/(1 - S^N) \tag{3.6}$$

$$\mu_{0,1} = \mu_{0,1}^N S^2 \tag{3.7}$$

$$\mu_{0,2} = [\mu_{0,2}^N + 2(\mu_{0,1}^N)^2 S] S^2 \tag{3.8}$$

$$\mu_{2,0} = \mu_{2,0}^N S^2. \tag{3.9}$$

3.3. Extrapolation procedure

When forming the sums (3.2) one could have stopped the summation at any t prior to reaching t_{\max} and used the formulae above to derive the series correct to order $n(t)$. Let S_t^N and X_t^N denote the non-nodal expansions obtained in this fashion. As observed by Baxter and Guttmann [5] one can often extend the series considerably by looking at correction terms to such series. The polynomials $S(t)$ and $X(t)$, and thus likewise the non-nodal expansions, will obviously contain terms of much higher order than that to which the final series is correct. One can therefore look at the difference between successive expansions, e.g.

$$S_t^N - S_{t+1}^N = -S^N(t+1) = p^{n(t+1)} \sum_{r \geq 0} s_{t,r} p^r \tag{3.10}$$

which yields sequences of numbers $s_{t,r}$ with $t < t_{\max}$. As observed in [5] the first sequence of numbers $s_{t,0}$ is often quite simple and can readily be conjectured so that a closed form expression or a simple recurrence relation can be found. In the following I will give the details of how this is done in the square bond case. The treatment of the other problems are detailed in the appendix. Note, that if one can find the first n_r correction terms one can use $S_{t_{\max}}^N = \sum_{m \geq 0} a_{N,m} p^m$ to extend the series $S^N = \sum_{m \geq 0} a_m p^m$ to order $n(t_{\max}) + n_r$, via

$$a_{n(t_{\max})+1+k} = a_{N,n(t_{\max})+1+k} - \sum_{m=0}^{\lfloor k/2 \rfloor} s_{t_{\max}+m,k-2m}. \tag{3.11}$$

So in order to find the correct series term $a_{n(t_{\max})+1+k}$ from the ‘partial’ term $a_{N,n(t_{\max})+1+k}$ one first subtracts $s_{t_{\max},k}$ which yields correctly the term $a_{N+1,n(t_{\max}+1)-1+k}$. One continues this process until arriving at $a_{N+\lfloor k/2 \rfloor+1,n(t_{\max}+\lfloor k/2 \rfloor+1)-q}$, where $q = 1(0)$ if k is even (odd), which is the correct term in the series for S^N .

In the square bond case the first sequence of correction terms start out as

$$s_{t,0} = 1, 2, 5, 14, 42, 132, 429, \dots$$

which is immediately recognizable as the Catalan numbers $C_t = (2t)!/(t!(t+1)!)$. These also occurred as the first correction term for the percolation probability series [5]. There is a very simple combinatorial proof for the first correction term. The first correction term arises

from the simplest (containing the minimum number of random elements) non-nodal graphs terminating at level $t + 1$. These graphs are also the ones giving the first term of $S^N(t + 1)$. It is obvious that these graphs are composed of two paths of length $t + 1$ each, which meet at level $t + 1$ but does not cross earlier. These graphs are in one-to-one correspondence with *staircase polyominoes* (or *polygons*) and it is well known that the latter are enumerated by the Catalan numbers [12, 13].

As was the case for the percolation probability series the higher-order correction terms can be expressed as rational functions of $s_{t,0}$. For S^N these extrapolation formulae are

$$s_{t,r} = \frac{2^r}{16^{\lfloor r/2 \rfloor}!} \sum_{k=1}^{\lfloor r/2 \rfloor} b_{r,k} (2t)^k C_{t-r+2} + \sum_{j=1}^{2r} a_{r,j} C_{t-r+j} \quad t \geq r \quad (3.12)$$

which are very similar to the formulae found in the percolation probability case [5]. The extrapolation formulae for $\mu_{0,1}^N$ and $\mu_{0,2}^N$ are simply $(t + 1)s_{t,r}$ and $(t + 1)^2 s_{t,r}$, respectively.

The factor in front of the first sum has been chosen so as to make the leading coefficients particularly simple. I was able to find formulae for all correction terms up to $r = 16$. The coefficients in the extrapolation formulae are listed in table 1.

From (3.12) it is clear that the $t_{\max} - r$ terms available in the sequences for the correction terms are not sufficient to determine all the $2r + \lfloor r/2 \rfloor$ unknown coefficients of the extrapolation formulae for large r . However, from table 1 one immediately sees that the leading coefficients $a_{r,2r}$ and $b_{r,\lfloor r/2 \rfloor}$ in the extrapolation formulae are very simple. In particular one has, $(-1)^r a_{r,2r} = 2$, and

$$b_{r,\lfloor r/2 \rfloor} = \begin{cases} (-1)^{\lfloor r/2 \rfloor} (r - 9) & r \text{ odd} \\ (-1)^{\lfloor r/2 \rfloor} & r \text{ even.} \end{cases}$$

Likewise, $a_{r,1}$ is zero for $r > 2$. In general I find that the leading coefficients $a_{r,2r-m}$ are expressible as polynomials in r of order m :

$$(-1)^r a_{r,2r-m} = \begin{cases} -4r & r > 0, m = 1 \\ 4r^2 - 10 & r > 2, m = 2 \\ -8r^3/3 + 80r/3 - 40 & r > 4, m = 3 \\ 4r^4/3 - 100r^2/3 + 86r - 48 & r > 6, m = 4 \\ -8r^5/15 + 80r^3/3 - 92r^2 - 62r/15 + 350 & r > 8, m = 5. \end{cases}$$

So when calculating the coefficients listed in table 1 I first used the sequences for the correction terms to predict as many of the extrapolation formulae (3.12) as possible. Then I predicted as many of the leading coefficients as possible. This in turn allowed me to find more extrapolation formulae, which I used to find more of the formulae for the leading coefficients $a_{r,2r-m}$. I repeated this until the process stopped with the extrapolation formulae listed in table 1.

For X^N the sequence determining the first correction formula starts out as

$$x_{t,0} = 0, 2, 8, 30, 112, 420, 1584, 6006, 22\,880, \dots$$

from which one sees that $x_{t,0} = 2(t - 1)C_{t-1}$. The proof of this formula is a little more involved. First one needs the number of configurations, $w(t, x)$, of two non-crossing paths terminating at (x, t) . Essam and Guttmann [14] gives a formula for the number of non-crossing watermelon configurations with p chains which join s steps and at height q from the origin

$$w_s(0) = 1 \quad w_s(s - q) = w_s(q)$$

and

$$w_s(q) = \prod_{i=1}^q \frac{(p+i)_{s-2i+1}}{(i)_{s-2i+1}} \quad 1 \leq q \leq \lfloor s/2 \rfloor \quad (3.13)$$

where $(a)_k = a(a+1)(a+2)\cdots(a+k-1)$, is Pochhammer's symbol. A watermelon configuration with two chains is in one-to-one correspondence with the configuration obtained from the two non-crossing paths by deleting the two bonds connected to the origin and the two bonds connected to the terminal point, so that $w(t, x) = w_{t-2}(x)$. In the case $p = 2$ (3.13) reduces to a simple product of binomial coefficients,

$$\begin{aligned} w_s(q) &= \prod_{i=1}^q \frac{(s-i+2)(s-i+1)}{i(i+1)} = \frac{s!(s+1)!}{(s-q)!q!(s+1-q)!(q+1)!} \\ &= \frac{1}{s+2} \binom{s}{q} \binom{s+2}{q+1}. \end{aligned} \quad (3.14)$$

The correction term $s_{t,0}$ can easily be derived from (3.14) as (remembering that $s_{t,0}$ arises from paths terminating at level $t+1$)

$$\begin{aligned} s_{t,0} &= \sum_{q=0}^{t-1} w_{t-1}(q) = \frac{1}{t+1} \sum_{q=0}^{t-1} \binom{t-1}{q} \binom{t+1}{q+1} \\ &= \frac{1}{t+1} \sum_{q=0}^t \binom{t-1}{q} \binom{t+1}{t-q} = \frac{1}{t+1} \binom{2t}{t} = C_t. \end{aligned}$$

In this derivation I have used only standard properties of binomial coefficients, the main one being the formula

$$\sum_{q=0}^p \binom{m}{q} \binom{n}{p-q} = \binom{m+n}{p}. \quad (3.15)$$

After this little diversion I return to the calculation of $x_{t,0}$. From (3.1) and the measurement of x with respect to the centre line it is clear that

$$x_{t,0} = \sum_{q=0}^s (s-2q)^2 w_s(q) \quad (3.16)$$

where $s = t - 1$. By simple expansion of the square and insertion of $w_s(q)$ one finds

$$\begin{aligned} x_{t,0} &= \frac{1}{s+2} \left[s^2 \sum_{q=0}^{s+1} \binom{s}{q} \binom{s+2}{q+1} - 4s \sum_{q=0}^{s+1} q \binom{s}{q} \binom{s+2}{q+1} \right. \\ &\quad \left. + 4 \sum_{q=0}^{s+1} q(q+1) \binom{s}{q} \binom{s+2}{q+1} - 4 \sum_{q=0}^{s+1} q \binom{s}{q} \binom{s+2}{q+1} \right] \\ &= \frac{1}{s+2} \left[s^2 \binom{2s+2}{s+1} - 4s^2 \binom{2s+1}{s+1} - 4s(s+2) \binom{2s}{s} - 4s \binom{2s+1}{s+1} \right] \\ &= \frac{1}{s+2} \left[-\frac{2s^2(2s+1)}{s+1} \binom{2s}{s} + 4s(s+2) \binom{2s}{s} - \frac{4s(2s+1)}{s+1} \binom{2s}{s} \right] \\ &= \frac{1}{(s+2)(s+1)} \binom{2s}{s} [2s^2 + 4s] = \frac{2s}{(s+1)} \binom{2s}{s} \\ &= 2sC_s = 2(t-1)C_{t-1}. \end{aligned}$$

The major step was the use of (3.15) to get rid of the sum over q . For the rest of the calculations I only used the definition and well known properties of the binomial coefficients.

In this case I find that the general extrapolation formulae can be written as

$$x_{t,r} = \frac{2^r}{16[r/2]!} \sum_{k=1}^{\lfloor r/2 \rfloor + 1} b_{r,k} (2t)^k C_{t-r+2} + \sum_{j=0}^{2r} a_{r,j} C_{t-r+j} \quad t \geq r. \quad (3.17)$$

The coefficients are not reproduced here due to the excessive length of this material, but are available from the author (please see end of article for details). Again I found that the leading coefficients are very simple, so a procedure similar to that used to find more extrapolation formulae for S^N was applied for X^N also. Though in this case it is slightly more complicated because different polynomials are found for $a_{r,2r-m}$ depending on whether r is odd or even. I was able to find the extrapolation formulae for $r \leq 15$.

From the polynomials for $S^N(t_{\max})$ and $X^N(t_{\max})$, using the extrapolation formulae given above, I extended the series for $S(p)$, $\mu_{0,1}(p)$ and $\mu_{0,2}(p)$ to order 112 and the series for $\mu_{2,0}(p)$ to order 111. The new series terms are listed in table 2, while the terms for $n \leq 49$ can be found in [4]. The full series are available from the author via e-mail or can be retrieved from the authors homepage on the world wide web (see later for details).

For the square site problem I have identified the first 12 extrapolation formulae for S^N and the first nine for X^N . This allowed me to derive the series correctly to order 106 and 103, respectively. For the triangular bond and site cases the first 10–12 extrapolation formulae were found and the series calculated to orders 55–57 depending on the particular problem. Details of the extrapolation formulae and lists of the new series coefficients can be found in the appendix. The full series and tables of the coefficients in the extrapolation formulae can be obtained from the author.

4. Analysis of the series

In the vicinity of the critical point one expects the moments of the pair-connectedness to have the functional form

$$f(p) \propto A(p_c - p)^\lambda [1 + a_1(p_c - p)^{\Delta_1} + b_1(p_c - p) \dots] \quad (4.1)$$

where λ is the critical exponent, Δ_1 the leading confluent exponent and the \dots represents higher-order correction terms. By universality we expect λ to be the same for all the percolation problems. In addition to the physical singularity, the series may have non-physical singularities for other values (real or complex) of p .

The series for moments of the pair-connectedness were analysed using inhomogeneous first- and second-order differential approximants. A comprehensive review of these and other techniques for series analysis may be found in [2]. Here it suffices to say that a K th-order differential approximant to a function f is formed by matching the earliest series coefficients to an inhomogeneous differential equation of the form (see [2] for details)

$$\sum_{i=0}^K Q_i(x) \left(x \frac{d}{dx} \right)^i f(x) = P(x) \quad (4.2)$$

where Q_i and P are polynomials of order N_i and L , respectively. First- and second-order approximants are denoted by $[L/N_0; N_1]$ and $[L/N_0; N_1; N_2]$, respectively.

4.1. The square bond series

In this section I will give a detailed account of the analysis of the square bond series which leads to the most accurate estimates. The analysis of the series for the other problems are described summarily in the following sections. In addition to the moment series I have also analysed the series $\mu_{0,2}(p)/\mu_{0,1}(p) \sim (p_c - p)^{-\nu_{\parallel}}$ and the series $\mu_{2,0}(p)\mu_{0,2}(p)/(\mu_{0,1}(p))^2 \sim (p_c - p)^{-2\nu_{\perp}}$.

In order to locate the singularities of the series in a systematic fashion I used the following procedure: I calculate all $[L/N; M]$ and $[L/N; M; M]$ first- and second-order inhomogeneous differential approximants with $|N - M| \leq 1$ and $L \leq 35$, which use more than 95 or 90 terms, respectively. Each approximant yields M possible singularities and associated exponents from the M zeroes of Q_1 or Q_2 , respectively (many of these are of course not actual singularities of the series but merely spurious zeros.) Next these zeroes are sorted into equivalence classes by the criterion that they lie at most a distance 2^{-k} apart. An equivalence class is accepted as a singularity if it contains more than N_c approximants, and an estimate for the singularity and exponent is obtained by averaging over the approximants (the spread among the approximants is also calculated). I used $N_c = 20$ (15) for first-order (second-order) approximants, which means that at least two-thirds to three-quarters of all approximants had to be included before an equivalence class was accepted. The calculation was then repeated for $k - 1, k - 2, \dots$ until a minimal value of 8 or so was reached. To avoid outputting well-converged singularities at every level, once an equivalence class has been accepted, the approximants which are members of it are removed, and the subsequent analysis is carried out on the remaining data only. One advantage of this method is that spurious outliers, a few of which will almost always be present when so many approximants are generated, are discarded systematically and automatically.

In table 3 I have listed the estimates for the physical critical point p_c and the associated exponents obtained from the six series that I studied. The errors listed in the parentheses are calculated from the spread among the approximants and equals one standard deviation. Note that these error estimates should *not* be seen as accurately representing the true errors. N_a is the number of approximants included in the estimates.

Generally the estimates for various orders L of the inhomogeneous polynomial are exceptionally well converged and excellent agreement is observed both between the various estimates for each series as well as between the p_c -estimates from the different series. Apart from the first-order approximants for small L to $\mu_{2,0}(p)\mu_{0,2}(p)/(\mu_{0,1}(p))^2$ all estimates for p_c are consistent with the highly accurate value $p_c = 0.644\,700\,15(15)$. This slight discrepancy is not important since one generally would expect large L first-order approximants and second-order approximants to yield more reliable estimates. These approximants are better at dealing with analytic background terms or other features which might possibly slow down the convergence of the estimates to the true critical values. Further note that N_a generally is well above the cut-off N_c showing that in most cases only a few approximants are discarded. The uncertainty in the last digits of the p_c -estimate, given in parentheses, is probably on the conservative side, and is mostly due to the tendency of $\mu_{0,1}$ and $\mu_{0,2}$ to favour a somewhat lower estimate for the critical point.

Before proceeding I will consider possible sources of systematic errors. First and foremost the possibility that the estimates might display a systematic drift as the number of terms used is increased and secondly the possibility of numerical errors. The latter possibility is quickly dismissed. The calculations were performed using 128-bit real numbers (REAL*16 on an IBM RISC work station). The estimates from a few approximants were compared to values obtained using MAPLE with up to 100 digits accuracy and this clearly

Table 3. Estimates of p_c and critical exponents for the square bond problem.

L	First-order DA			Second-order DA		
	p_c	γ	N_a	p_c	γ	N_a
0	0.644 700 51(60)	2.278 32(77)	25	0.644 700 181(37)	2.277 716(30)	22
5	0.644 700 18(72)	2.278 07(71)	25	0.644 700 169(26)	2.277 708(23)	18
10	0.644 700 04(13)	2.277 602(93)	26	0.644 700 158(41)	2.277 703(34)	23
15	0.644 700 136(29)	2.277 665(56)	23	0.644 700 146(29)	2.776 90(23)	20
20	0.644 700 102(21)	2.277 649(21)	24	0.644 700 146(17)	2.277 689(14)	18
25	0.644 700 097(49)	2.277 646(42)	23	0.644 700 149(20)	2.277 693(15)	21
30	0.644 700 108(29)	2.277 659(24)	26	0.644 700 162(12)	2.277 704(11)	16
35	0.644 700 129(21)	2.277 678(15)	21	0.644 700 29(22)	2.277 92(42)	22
L	p_c	ν_{\parallel}	N_a	p_c	ν_{\parallel}	N_a
0	0.644 700 153(12)	1.733 818 4(50)	22	0.644 700 169(97)	1.733 845(45)	19
5	0.644 700 154(31)	1.733 818(12)	27	0.644 700 178(50)	1.733 846(28)	16
10	0.644 700 115(11)	1.733 807 1(35)	22	0.644 700 171 8(88)	1.733 836 2(42)	20
15	0.644 700 142(33)	1.733 819(21)	22	0.644 700 136(50)	1.733 813(34)	18
20	0.644 700 162(14)	1.733 831 9(78)	25	0.644 700 154(23)	1.733 827(11)	19
25	0.644 700 149(24)	1.733 824(11)	25	0.644 700 142(13)	1.733 821 3(67)	18
30	0.644 700 155 7(63)	1.733 827 9(31)	23	0.644 700 122(34)	1.733 806(25)	21
35	0.644 700 150 3(61)	1.733 825 4(32)	22	0.644 700 164(20)	1.733 831 2(92)	20
L	p_c	$2\nu_{\perp}$	N_a	p_c	$2\nu_{\perp}$	N_a
0	0.644 700 40(13)	2.193 828(55)	22	0.644 700 196(17)	2.193 711(11)	17
5	0.644 700 438(94)	2.193 843(36)	22	0.644 700 192(18)	2.193 708(10)	18
10	0.644 700 41(17)	2.193 826(95)	22	0.644 700 174(47)	2.193 703(29)	17
15	0.644 700 147(17)	2.193 685 2(79)	22	0.644 700 163(23)	2.193 693(12)	18
20	0.644 700 201(17)	2.193 712 6(82)	23	0.644 700 217(40)	2.193 722(22)	16
25	0.644 700 200(10)	2.193 713 2(54)	23	0.644 700 192(28)	2.193 708(13)	16
30	0.644 700 196(10)	2.193 710 7(51)	23	0.644 700 183(12)	2.193 703 9(64)	17
35	0.644 700 195(14)	2.193 711 0(69)	23	0.644 700 182(15)	2.193 703 1(84)	18
L	p_c	$\gamma + \nu_{\parallel}$	N_a	p_c	$\gamma + \nu_{\parallel}$	N_a
0	0.644 700 091(76)	4.011 423(76)	24	0.644 700 091(32)	4.011 434(35)	18
5	0.644 700 042(74)	4.011 375(65)	25	0.644 700 095(20)	4.011 440(23)	18
10	0.644 700 023(97)	4.011 361(79)	25	0.644 700 079(37)	4.011 413(44)	20
15	0.644 700 071(72)	4.011 403(73)	24	0.644 700 105(47)	4.011 455(50)	20
20	0.644 700 015(66)	4.011 350(57)	26	0.644 700 096(32)	4.011 443(34)	18
25	0.644 700 04(15)	4.011 39(15)	21	0.644 700 096(63)	4.011 440(73)	19
30	0.644 700 037(68)	4.011 370(59)	24	0.644 700 101(21)	4.011 448(22)	19
35	0.644 700 038(54)	4.011 369(49)	23	0.644 700 090(20)	4.011 438(22)	18
L	p_c	$\gamma + 2\nu_{\parallel}$	N_a	p_c	$\gamma + 2\nu_{\parallel}$	N_a
0	0.644 700 043(87)	5.745 15(10)	24	0.644 700 079(19)	5.745 208(29)	18
5	0.644 700 079(96)	5.745 20(13)	24	0.644 700 084(25)	5.745 224(35)	16
10	0.644 700 05(11)	5.745 17(13)	21	0.644 700 075(29)	5.745 208(37)	17
15	0.644 700 11(10)	5.745 25(17)	22	0.644 700 075(17)	5.745 213(25)	22
20	0.644 700 051(27)	5.745 156(34)	24	0.644 700 087(38)	5.745 232(51)	17
25	0.644 700 13(17)	5.745 31(32)	25	0.644 700 082(22)	5.745 225(32)	18
30	0.644 700 068(45)	5.745 180(57)	21	0.644 700 082(25)	5.745 231(50)	18
35	0.644 699 99(10)	5.745 10(11)	25	0.644 700 091(45)	5.745 231(75)	19

Table 3. (Continued)

L	First-order DA			Second-order DA		
	p_c	$\gamma + 2\nu_{\perp}$	N_a	p_c	$\gamma + 2\nu_{\perp}$	N_a
0	0.644 700 081 9(37)	4.471 298 8(18)	22	0.644 700 119(52)	4.471 341(57)	20
5	0.644 700 080 6(26)	4.471 298 1(13)	23	0.644 700 117(21)	4.471 329(20)	17
10	0.644 700 085 7(78)	4.471 301 7(62)	24	0.644 700 115(46)	4.471 332(46)	16
15	0.644 700 138(69)	4.471 36(10)	21	0.644 700 094(68)	4.471 319(50)	16
20	0.644 700 101(24)	4.471 315(21)	23	0.644 700 132(40)	4.471 351(42)	16
25	0.644 700 101(29)	4.471 316(25)	25	0.644 700 101(16)	4.471 314(14)	16
30	0.644 700 112(21)	4.471 324(19)	21	0.644 700 121(42)	4.471 340(46)	19
35	0.644 700 119(17)	4.471 330(16)	21	0.644 700 114(41)	4.471 334(44)	18

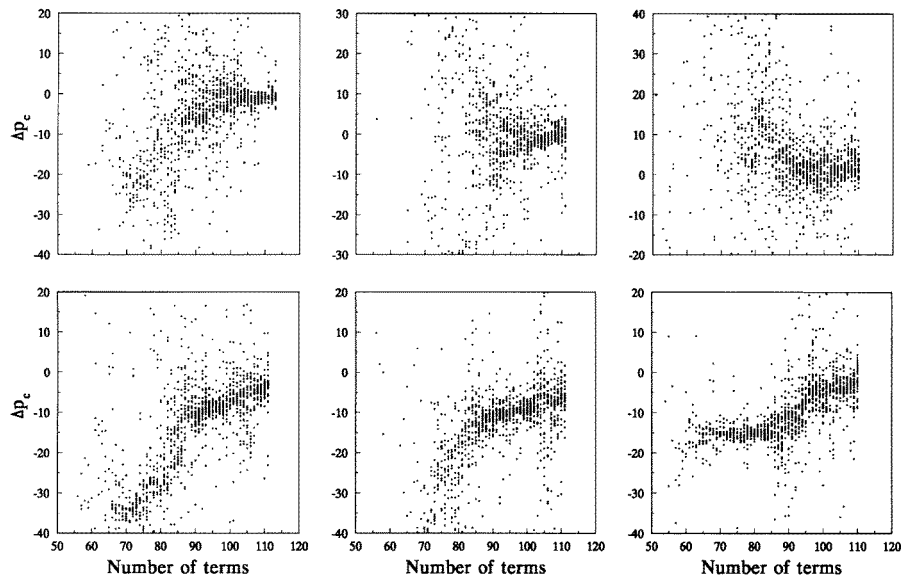


Figure 2. The deviation in the last two digits, $10^8 \Delta p_c$, from the central estimate of the critical point $p_c = 0.644\,700\,15$, of the estimates for the critical point by second-order differential approximants. Shown is (from left to right and top to bottom) estimates from the series $S(p)$, $\mu_{0,2}(p)/\mu_{0,1}(p)$, $\mu_{2,0}(p)\mu_{0,2}(p)/(\mu_{0,1}(p))^2$, $\mu_{0,1}(p)$, $\mu_{0,2}(p)$, and $\mu_{2,0}(p)$.

showed that the program was numerically stable and rounding errors were negligible. In order to address the possibility of systematic drift and lack of convergence to the true critical values I refer to figure 2. In this figure I have plotted the deviation in the last two digits, $10^8 \Delta p_c$, from the critical point $p_c = 0.644\,700\,15$. Included in the figure are estimates from inhomogeneous second-order differential approximants with $L \leq 35$ to the six series that I have studied. From this figure it is evident that the series estimates displayed on the top row are well converged once the number of terms exceeds 90 or so, while the series on the bottom row still show evidence of a systematic drift and the estimates have not yet converged to their asymptotic value. This is particularly manifest for the series $\mu_{0,1}$ and $\mu_{0,2}$ shown in the bottom left and central panels. Since these series were the ones responsible for most of the error on the estimate for p_c , and given the very good convergence of the estimates from the series shown in the top row, it does not seem overly optimistic to adopt

the tighter estimate $p_c = 0.647\,700\,15(5)$. Clearly the large majority of estimates for the first three series lie well within this error-bound as the number of terms increase and likewise the estimates from the remaining series clearly seem to converge towards this value.

Next I turn my attention to the estimates for the critical exponents. Very precise estimates for γ , ν_{\parallel} , and $2\nu_{\perp}$ can be obtained by examining table 3. I have used a slightly more systematic and enlightening procedure. Close to the critical point there is an apparent linear dependence of the estimates for critical exponents on the estimates for p_c . One can use this to obtain improved estimates for the exponents by performing a linear fit of the exponent estimates as a function of Δp_c (the distance from the critical point). The result of such linear fits is listed below. In these fits I used the same set of approximants as those on which the estimates in the tables above were based. But I discarded any approximant for which $|\Delta p_c| = |p_c - 0.644\,700\,15| > 0.000\,000\,15$. The error on the ‘pure’ exponent part of the estimates mainly reflects the slight difference between the first- and second-order approximants (the errors as listed are approximately twice this difference). In the estimates for γ and $\gamma + 2\nu_{\perp}$ I used only the first-order approximants with $L \geq 15$.

$$\begin{aligned}
 \gamma &= 2.277\,690(10) \pm 750\Delta p_c \\
 \nu_{\parallel} &= 1.733\,824(3) \pm 500\Delta p_c \\
 2\nu_{\perp} &= 2.193\,687(2) \pm 500\Delta p_c \\
 \gamma + \nu_{\parallel} &= 4.011\,495(15) \pm 1150\Delta p_c \\
 \gamma + 2\nu_{\parallel} &= 5.745\,308(15) \pm 1400\Delta p_c \\
 \gamma + 2\nu_{\perp} &= 4.471\,368(3) \pm 1000\Delta p_c.
 \end{aligned}
 \tag{4.3}$$

As can be seen the exponent estimates are very precise. Even with the very small error in the p_c -estimate, this is still the major source of error (by an order of magnitude) in the exponent estimates. As previously noted [6], there is no simple rational fraction whose decimal expansion agrees with the estimate of β obtained from the percolation-probability series. The same is true for the estimates of ν_{\parallel} and $2\nu_{\perp}$ listed above. In particular note that the rational fraction suggested by Essam *et al* [4], $\nu_{\parallel} = 26/15 = 1.733\,333\dots$, and $2\nu_{\perp} = 79/36 = 2.19444\dots$, is incompatible with the estimates. The rational fraction suggested for $\gamma = 41/18 = 2.277\,777\dots$ lies within the error bounds for the exponent estimate if the error on p_c exceeds 10^{-7} . So the more conservative error estimate listed earlier would just include the suggested value of γ . However, most of the estimates in table 3 clearly exclude the exact fraction as does the more narrow error estimate on p_c . Finally I note that the better converged estimates for $\gamma + 2\nu_{\perp}$ and $2\nu_{\perp}$ yields the estimate $\gamma = 2.277\,681(5)$, which, within the error, agrees with the direct estimate but points to a possibly slightly lower value of γ .

The estimate for p_c advocated above lies within the error-bounds of that obtained from the percolation probability series [6] $p_c = 0.644\,700\,6(10)$, though a lower central value is favoured by the series analysed in this paper. From the scaling relation $\beta = (\nu_{\parallel} + \nu_{\perp} - \gamma)/2$ I obtain the estimate $\beta = 0.276\,489(7) \pm 750\Delta p_c$, which is consistent with the direct estimate $\beta = 0.276\,43(10)$. It is quite likely that the minor discrepancies between the central values would disappear if the percolation probability series could be extended from the 55 terms in [6] to an order comparable to the series analysed here. Evidence to this effect is provided by the biased estimate $\beta = 0.276\,483(14)$ calculated at $p_c = 0.644\,700\,15$ using Dlog Padé approximants utilizing at least 45 terms of the percolation-probability series.

I also analysed the series in order to estimate the leading confluent exponents Δ_1 . As was the case for the percolation-probability series both the Baker–Hunter transformation and the method of Adler, Moshe and Privman (see [6] and references therein for details

regarding these methods) yielded estimates consistent with $\Delta_1 = 1$. So there are no signs of non-analytic corrections to scaling.

Finally I looked for non-physical singularities of the series. The series have a singularity on the negative axis closer to the origin than p_c . This singularity is quite weak and consequently the estimates for its location and the associated exponents are quite inaccurate. The singularity is located at $p_- = -0.5168(5)$ and the associated exponents are $\gamma = 0.065(15)$, $\nu_{\parallel} = 0.97(3)$ and $2\nu_{\perp} = 0.90(15)$. It is quite possible that the divergence of the cluster length series at p_- is logarithmic and the estimates are certainly consistent with $\gamma = 0$, $\nu_{\parallel} = 1$ and $\nu_{\perp} = \frac{1}{2}$. Finally there is some weak evidence of a pair of singularities in the complex p -plane at $p_{\pm} = -0.2255(15) \pm 0.440(1)i$. Note that this singularity pair also lies within the physical disc. The exponent estimates at p_{\pm} are not very accurate. The cluster size series seems to *converge* with exponent $\gamma \simeq -3$, while $\nu_{\parallel} \simeq 1$ and $\nu_{\perp} \simeq \frac{1}{2}$, but the error on these estimates are as large as 25–50%.

4.2. The square site series

In table 4 I have listed some of the estimates for p_c and critical exponents obtained from an analysis of the square site series. The estimates are based on approximants using at least 85–90 terms with $N_c = 15$. Though the length of the series is comparable to the bond case the estimates are generally less accurate. In particular it should be noted that the p_c -estimates obtained from different series are only marginally consistent leading to the rather poor estimate, $p_c = 0.7054850(15)$, which is at least an order of magnitude less accurate than in the bond case. Some exponent estimates differ significantly from those of the bond case. Particularly γ and $\gamma + 2\nu_{\parallel}$ are generally quite a bit smaller than the bond estimates. However, due to the discrepancy between the various site series, the importance of this deviation is questionable. If the error-bar on p_c is accepted, the resulting exponent estimates from the site series will agree with the bond estimates.

If one accepts the exponent estimates from the bond series one can use the linear dependence between p_c and exponent estimates to obtain improved estimates for p_c . (This is just the reverse of the method used in the previous section to obtain the exponent estimates.) By performing a linear fit of the p_c -estimates as a function of the deviation of the exponent estimate from the central values listed in the previous section I obtain the estimate $p_c = 0.7054853(5)$. In these fits I used the approximants whose exponent estimates differ by less than 0.001 from the central values. This estimate agrees with that obtained from the percolation-probability series [6] $p_c = 0.705485(5)$.

The square site series have a singularity on the negative axis closer to the origin than p_c . In this case the singularity appears to be stronger than in the bond case, i.e. the various estimates are better converged. The singularity is located at $p_- = -0.4519522(3)$ and the associated exponents are quite possibly consistent with $\gamma = -\frac{1}{2}$ (i.e. the cluster-size series *converges*), $\nu_{\parallel} = 1$ and $\nu_{\perp} = \frac{1}{2}$. There is firm evidence of a pair of singularities in the complex p -plane at $p_{\pm} = -0.2263(1) \pm 0.3847(1)i$, which is within the physical disc. The exponent estimates at this pair of singularities are quite accurate. The cluster-size series seems to *converge*, with $\gamma \simeq -3$, while $\nu_{\parallel} \simeq 1$ and $\nu_{\perp} \simeq \frac{1}{2}$, where errors on the estimates are only a few per cent.

4.3. The triangular bond series

Table 5 lists a selection of estimates for p_c and critical exponents obtained from the analysis of the triangular bond series. The estimates are based on approximants using at least 45 or

Table 4. Estimates of p_c and critical exponents for the square site problem.

L	First-order DA			Second-order DA		
	p_c	γ	N_a	p_c	γ	N_a
0	0.705 483 90(20)	2.276 850(66)	19	0.705 485 00(26)	2.277 51(15)	17
5	0.705 484 09(20)	2.276 924(88)	23	0.705 485 16(28)	2.277 60(18)	18
10	0.705 484 41(35)	2.277 21(30)	24	0.705 484 72(19)	2.277 334(95)	17
15	0.705 484 594(68)	2.277 232(33)	23	0.705 484 71(14)	2.277 314(74)	19
20	0.705 484 805(72)	2.277 364(39)	24	0.705 484 86(36)	2.277 42(25)	20
25	0.705 484 723(82)	2.277 319(46)	20	0.705 484 671(58)	2.277 295(35)	16
30	0.705 484 811(34)	2.277 367(18)	21	0.705 484 689(29)	2.277 306(16)	16
35	0.705 484 850(62)	2.277 389(31)	21	0.705 484 713(83)	2.277 313(39)	17
L	p_c	ν_{\parallel}	N_a	p_c	ν_{\parallel}	N_a
0	0.705 484 49(93)	1.733 47(25)	19	0.705 484 96(30)	1.733 70(10)	16
5	0.705 484 27(28)	1.733 416(72)	23	0.705 484 91(23)	1.733 686(84)	16
10	0.705 484 85(36)	1.733 66(14)	20	0.705 485 020(95)	1.733 729(25)	16
15	0.705 485 13(26)	1.733 763(88)	23	0.705 484 91(34)	1.733 69(12)	18
20	0.705 485 65(53)	1.733 97(20)	22	0.705 484 80(17)	1.733 650(66)	19
25	0.705 485 75(33)	1.734 03(12)	23	0.705 484 70(21)	1.733 608(93)	17
30	0.705 485 60(63)	1.733 96(28)	19	0.705 484 43(26)	1.733 50(11)	16
35	0.705 485 45(43)	1.733 88(17)	24	0.705 484 52(21)	1.733 548(84)	16
L	p_c	$2\nu_{\perp}$	N_a	p_c	$2\nu_{\perp}$	N_a
0	0.705 486 9(13)	2.194 45(46)	19	0.705 486 50(23)	2.194 33(21)	19
5	0.705 486 87(57)	2.194 47(16)	19	0.705 486 47(23)	2.194 34(13)	16
10	0.705 485 1(15)	2.193 97(33)	21	0.705 486 49(12)	2.194 254(51)	16
15	0.705 485 7(10)	2.194 00(39)	19	0.705 485 77(24)	2.194 033(76)	20
20	0.705 486 6(16)	2.194 34(53)	19	0.705 485 89(42)	2.194 06(13)	21
25	0.705 486 0(10)	2.194 12(42)	19	0.705 485 85(24)	2.194 048(81)	17
30	0.705 486 0(12)	2.194 10(45)	20	0.705 485 60(65)	2.193 91(28)	18
35	0.705 486 2(13)	2.194 08(53)	20	0.705 485 15(78)	2.193 76(31)	17
L	p_c	$\gamma + \nu_{\parallel}$	N_a	p_c	$\gamma + \nu_{\parallel}$	N_a
0	0.705 483 65(38)	4.009 89(23)	19	0.705 484 03(70)	4.010 23(58)	18
5	0.705 483 81(17)	4.010 00(12)	23	0.705 484 38(33)	4.010 47(39)	16
10	0.705 483 85(42)	4.010 05(29)	25	0.705 484 41(34)	4.010 55(30)	16
15	0.705 483 62(55)	4.009 94(38)	24	0.705 484 30(51)	4.010 46(44)	21
20	0.705 483 49(30)	4.009 79(20)	19	0.705 484 24(34)	4.010 41(28)	18
25	0.705 483 80(43)	4.010 06(30)	22	0.705 484 50(65)	4.010 67(65)	21
30	0.705 483 80(21)	4.009 99(14)	21	0.705 484 28(21)	4.010 43(18)	16
35	0.705 483 78(61)	4.010 02(43)	23	0.705 484 47(33)	4.010 61(32)	19
L	p_c	$\gamma + 2\nu_{\parallel}$	N_a	p_c	$\gamma + 2\nu_{\parallel}$	N_a
0	0.705 483 58(35)	5.743 11(21)	19	0.705 484 60(45)	5.744 20(51)	19
5	0.705 483 55(20)	5.743 07(14)	19	0.705 484 43(18)	5.744 00(20)	17
10	0.705 484 04(60)	5.743 58(65)	23	0.705 484 34(18)	5.743 92(21)	17
15	0.705 483 82(10)	5.743 299(94)	19	0.705 484 31(52)	5.743 90(62)	20
20	0.705 483 79(15)	5.743 27(14)	22	0.705 484 15(22)	5.743 69(24)	18
25	0.705 483 75(16)	5.743 21(13)	22	0.705 484 00(10)	5.743 52(10)	16
30	0.705 483 68(16)	5.743 17(14)	19	0.705 484 22(25)	5.743 77(30)	16
35	0.705 483 87(24)	5.743 34(22)	25	0.705 484 74(65)	5.744 49(85)	19

Table 4. (Continued)

L	First-order DA			Second-order DA		
	p_c	$\gamma + 2\nu_{\perp}$	N_a	p_c	$\gamma + 2\nu_{\perp}$	N_a
0	0.705 483 8(33)	4.472 9(94)	19	0.705 484 57(13)	4.470 71(10)	20
5	0.705 484 58(16)	4.470 69(11)	19	0.705 484 60(10)	4.470 740(93)	16
10	0.705 484 63(16)	4.470 72(10)	20	0.705 484 57(11)	4.470 695(93)	19
15	0.705 484 77(19)	4.470 84(15)	19	0.705 484 73(27)	4.470 84(25)	21
20	0.705 484 43(43)	4.470 61(26)	20	0.705 484 72(17)	4.470 81(15)	17
25	0.705 484 49(47)	4.470 66(30)	20	0.705 484 80(49)	4.470 89(45)	19
30	0.705 484 75(42)	4.470 87(37)	19	0.705 484 2(13)	4.470 4(11)	17
35	0.705 484 69(22)	4.470 78(18)	19	0.705 485 1(13)	4.471 3(12)	20

40 terms with $N_c = 15$ or 10 for first and second order, respectively. As one would expect, due to the shorter series, the estimates are generally encumbered with larger errors than was the case for the square bond series. The estimates for ν_{\parallel} and $2\nu_{\perp}$ are generally consistent with those from the square bond series, while the remaining exponent estimates exceeds those from the square bond case. The linear fit of p_c to the deviation of the exponent estimates from the values favoured by the square bond series yields $p_c = 0.478\,025(1)$, which is in excellent agreement with the estimate $p_c = 0.478\,02(1)$ from the percolation-probability series [7]. The triangular bond series does not appear to have any non-physical singularities.

4.4. The triangular site series

In table 6 I have listed some estimates for p_c and critical exponents obtained from an analysis of the triangular site series similar to that for the bond problem. In this case all exponent estimates are consistent with the square bond case. The biased estimate for p_c based on the usual fitting procedure is $p_c = 0.595\,646\,8(5)$ in excellent agreement with the estimate $p_c = 0.595\,647\,2(10)$ from the percolation probability series [7]. Again there is no compelling evidence for non-physical singularities.

5. Summary and discussion

From the analysis presented in the previous section it was clear that the square bond series yield by far the most accurate p_c -estimates which in turn enables one to obtain very precise estimates for the critical exponents. The remaining cases yielded less accurate estimates. Though the square site and triangular bond cases tended to yield exponent estimates only marginally consistent with the square bond estimates, the p_c estimates showed less consistency among the various series. In the square site case this could possibly be caused by the presence of rather strong non-physical singularities closer to the origin than p_c . The triangular site estimates, though marred by larger error-bars, were fully consistent with the square bond estimates. I have therefore chosen to base my final exponent estimates mainly on the square bond series.

From figure 2 it would appear that the estimate $p_c = 0.644\,700\,15(5)$ is fully consistent with the data and not overly optimistic. With this highly accurate p_c value one can obtain very accurate exponent estimates using the values listed in (4.3). The values of the critical exponents for the average cluster size, parallel and perpendicular connectedness lengths are

Table 5. Estimates of p_c and critical exponents for the triangular bond problem.

L	First-order DA			Second-order DA		
	p_c	γ	N_a	p_c	γ	N_a
0	0.478 026 8(13)	2.278 50(35)	21	0.478 025 48(13)	2.277 976(80)	15
4	0.478 025 96(10)	2.278 170(47)	16	0.478 025 78(42)	2.278 09(21)	14
8	0.478 026 14(10)	2.278 242(64)	16	0.478 025 60(16)	2.278 054(48)	11
12	0.478 026 02(42)	2.278 19(14)	20	0.478 025 79(27)	2.278 093(91)	14
16	0.478 025 99(29)	2.278 19(10)	18	0.478 026 05(50)	2.278 20(19)	17
L	p_c	ν_{\parallel}	N_a	p_c	ν_{\parallel}	N_a
0	0.478 027 2(19)	1.734 35(30)	17	0.478 026 24(79)	1.734 13(18)	17
4	0.478 025 5(10)	1.734 04(33)	17	0.478 025 85(59)	1.734 04(17)	12
8	0.478 025 51(57)	1.733 98(16)	16	0.478 026 4(10)	1.734 17(30)	15
12	0.478 025 6(18)	1.734 03(53)	19	0.478 025 36(79)	1.733 92(22)	11
16	0.478 024 4(25)	1.733 65(65)	18	0.478 027 3(19)	1.734 41(52)	15
L	p_c	$2\nu_{\perp}$	N_a	p_c	$2\nu_{\perp}$	N_a
0	0.478 027 16(70)	2.194 29(16)	18	0.478 026 0(10)	2.193 89(23)	17
4	0.478 026 83(80)	2.194 20(15)	17	0.478 026 1(17)	2.193 95(54)	14
8	0.478 024 74(53)	2.193 55(15)	16	0.478 024 6(12)	2.193 55(33)	14
12	0.478 025 1(28)	2.193 67(71)	18	0.478 024 4(12)	2.193 49(36)	14
16	0.478 024 7(11)	2.193 54(35)	17	0.478 025 22(40)	2.193 69(11)	11
L	p_c	$\gamma + \nu_{\parallel}$	N_a	p_c	$\gamma + \nu_{\parallel}$	N_a
0	0.478 026 76(52)	4.012 59(28)	18	0.478 026 65(24)	4.012 624(79)	13
4	0.478 026 70(47)	4.012 61(14)	20	0.478 026 86(12)	4.012 693(33)	13
8	0.478 026 45(51)	4.012 51(22)	19	0.478 026 66(17)	4.012 649(45)	11
12	0.478 026 12(59)	4.012 36(30)	17	0.478 026 53(68)	4.012 44(54)	16
16	0.478 026 22(45)	4.012 43(21)	16	0.478 026 82(16)	4.012 688(36)	11
L	p_c	$\gamma + 2\nu_{\parallel}$	N_a	p_c	$\gamma + 2\nu_{\parallel}$	N_a
0	0.478 025 4(17)	5.7456(17)	17	0.478 026 4(16)	5.7464(14)	13
4	0.478 025 1(10)	5.745 66(95)	19	0.478 026 6(24)	5.7460(20)	13
8	0.478 025 2(11)	5.7457(11)	17	0.478 026 4(19)	5.7461(16)	17
12	0.478 025 66(33)	5.746 23(26)	16	0.478 025 4(10)	5.7457(11)	16
16	0.478 025 88(78)	5.746 33(52)	18	0.478 026 3(18)	5.7463(12)	17
L	p_c	$\gamma + 2\nu_{\perp}$	N_a	p_c	$\gamma + 2\nu_{\perp}$	N_a
0	0.478 026 16(38)	4.472 28(18)	16	0.478 025 85(24)	4.472 04(14)	13
4	0.478 026 32(82)	4.472 34(41)	17	0.478 025 70(52)	4.471 91(33)	14
8	0.478 025 89(47)	4.472 14(23)	17	0.478 026 37(54)	4.472 35(31)	11
12	0.478 025 66(48)	4.471 96(31)	18	0.478 026 24(50)	4.472 28(31)	13
16	0.478 026 18(31)	4.472 28(15)	17	0.478 026 10(42)	4.472 18(23)	12

estimated by $\gamma = 2.277 69(4)$, $\nu_{\parallel} = 1.733 825(25)$ and $\nu_{\perp} = 1.096 844(14)$, respectively. An improved estimate for the percolation probability exponent is obtained from the scaling relation $\beta = (\nu_{\parallel} + \nu_{\perp} - \gamma)/2 = 0.276 49(4)$. As already noted these estimates are generally incompatible with the exact fractions conjectured by Essam *et al* [4]. Only γ is marginally consistent with the suggested fraction, $\gamma = 41/18 = 2.77 777 \dots$, if a larger error-bar were adopted for p_c .

Table 6. Estimates of p_c and critical exponents for the triangular site problem.

L	First-order DA			Second-order DA		
	p_c	γ	N_a	p_c	γ	N_a
0	0.595 647 31(31)	2.277 848(67)	16	0.595 645 98(71)	2.277 49(16)	18
4	0.595 646 41(30)	2.277 597(79)	18	0.595 646 5(13)	2.277 55(64)	16
8	0.595 646 64(41)	2.277 67(12)	18	0.595 646 81(10)	2.277 708(28)	12
12	0.595 646 53(27)	2.277 628(81)	16	0.595 646 67(20)	2.277 672(64)	13
16	0.595 646 84(78)	2.277 72(22)	18	0.595 646 59(32)	2.277 662(84)	12
L	p_c	ν_{\parallel}	N_a	p_c	ν_{\parallel}	N_a
0	0.595 646 56(15)	1.733 766(15)	16	0.595 646 75(45)	1.733 796(53)	15
4	0.595 645 4(11)	1.733 58(18)	16	0.595 646 62(60)	1.733 78(11)	11
8	0.595 645 9(88)	1.7336(17)	16	0.595 644 8(32)	1.733 44(74)	11
12	0.595 647 6(31)	1.734 07(68)	16	0.595 645 7(13)	1.733 61(29)	11
16	0.595 650 7(29)	1.734 77(65)	16	0.595 643 2(58)	1.7328(15)	15
L	p_c	$2\nu_{\perp}$	N_a	p_c	$2\nu_{\perp}$	N_a
0	0.595 650(12)	2.1943(37)	16	0.595 647 0(38)	2.1938(12)	14
4	0.595 655 5(49)	2.1958(11)	16	0.595 647 7(10)	2.193 97(25)	11
8	0.595 648 9(14)	2.194 25(30)	17	0.595 647 53(88)	2.193 97(24)	11
12	0.595 646 9(73)	2.1938(15)	16	0.595 645 7(22)	2.193 57(42)	12
16	0.595 647 3(10)	2.193 87(22)	16	0.595 648 5(18)	2.194 11(37)	16
L	p_c	$\gamma + \nu_{\parallel}$	N_a	p_c	$\gamma + \nu_{\parallel}$	N_a
0	0.595 643 5(26)	4.010 06(80)	18	0.595 645 3(22)	4.0108(10)	15
4	0.595 644 6(16)	4.010 36(54)	16	0.595 647 6(46)	4.0122(24)	17
8	0.595 645 42(67)	4.010 64(27)	17	0.595 647 29(73)	4.011 68(46)	11
12	0.595 644 89(48)	4.010 41(20)	16	0.595 647 19(88)	4.011 68(49)	11
16	0.595 644 95(28)	4.010 47(10)	17	0.595 645 0(12)	4.010 57(55)	11
L	p_c	$\gamma + 2\nu_{\parallel}$	N_a	p_c	$\gamma + 2\nu_{\parallel}$	N_a
0	0.595 648 4(66)	5.7469(60)	17	0.595 644 4(17)	5.743 86(91)	11
4	0.595 644 0(29)	5.7437(10)	16	0.595 644 2(28)	5.7438(16)	12
8	0.595 649 2(45)	5.7468(31)	18	0.595 643 2(32)	5.7433(12)	13
12	0.595 646 3(37)	5.7448(24)	17	0.595 646 2(20)	5.7448(13)	12
16	0.595 645 7(15)	5.744 40(85)	17	0.595 646 5(13)	5.745 02(80)	12
L	p_c	$\gamma + 2\nu_{\perp}$	N_a	p_c	$\gamma + 2\nu_{\perp}$	N_a
0	0.595 647 7(11)	4.471 67(39)	16	0.595 647 15(31)	4.471 61(13)	12
4	0.595 647 48(19)	4.471 776(73)	17	0.595 647 06(43)	4.471 56(17)	14
8	0.595 647 49(26)	4.471 770(98)	17	0.595 647 24(29)	4.471 64(12)	12
12	0.595 647 56(33)	4.471 79(12)	16	0.595 647 44(81)	4.471 70(29)	14
16	0.595 647 58(42)	4.471 80(15)	17	0.595 647 29(15)	4.471 670(61)	12

Below I have listed improved estimates for a number of critical exponents obtained using various scaling relations.

$$\Delta = \beta + \gamma = 2.554 18(8)$$

$$\tau = \nu_{\parallel} - \beta = 1.457 34(7)$$

$$z = \nu_{\parallel} / \nu_{\perp} = 1.580 74(4)$$

$$\gamma' = \gamma - \nu_{\parallel} = 0.543\,86(7)$$

$$\delta = \beta/\nu_{\parallel} = 0.159\,47(3)$$

$$\eta = \gamma/\nu_{\parallel} - 1 = 0.313\,68(4).$$

Here Δ is the exponent characterizing the scale of the cluster size distribution, τ is the cluster length exponent, z is the dynamical critical exponent, γ' the exponent characterizing the steady-state fluctuations of the order parameter, while δ and η characterize the behaviour at p_c as $t \rightarrow \infty$ of the survival probability and average number of particles, respectively.

Assuming that the exponent estimates from the square bond case are correct, improved p_c -estimates were obtained for the three other problems studied in this paper. These are:

$$\begin{aligned} p_c &= 0.705\,485\,3(5) && \text{square site} \\ p_c &= 0.478\,025(1) && \text{triangular bond} \\ p_c &= 0.595\,646\,8(5) && \text{triangular site.} \end{aligned}$$

Finally I note, that the analysis of the various series, in order to determine the value of the confluent exponent, yielded estimates consistent with $\Delta_1 \simeq 1$. Thus there is no evidence of non-analytic confluent correction terms. This provides a hint that the models might be exactly solvable.

E-mail or WWW retrieval of series

The series and the coefficients in the extrapolation formulae for the directed percolation problems on the various lattices can be obtained via e-mail by sending a request to iwan@maths.mu.oz.au or via the world wide web on the URL <http://www.maths.mu.oz.au/~iwan/> by following the relevant links.

Acknowledgments

I would like to thank G X Viennot for providing the reference to the article by Shapiro. The series enumerations were performed on the Intel Paragon parallel computer at the University of Melbourne. The work was supported by a grant from the Australian Research Council.

Appendix. The extrapolation formulae and series for the square site, triangular bond and triangular site problems

A.1. The square site problem

The sequence determining the first correction term for S^N starts out as

$$s_{t,0} = 1, 0, 1, 2, 6, 18, 57, 186, 622, 2120, 7338, \dots$$

from which one sees that $2s_{t,0} + s_{t-1,0} = C_{t-1}$. Shapiro [15] has given an interpretation of this sequence by adding diagonals in a certain Catalan triangle.

At first glance one might find it strange that the correction term differs from the bond case, since clearly all the non-nodal bond graphs that give rise to the first correction term have their counterparts as site graphs. In the following I shall always be talking only of non-nodal graphs consisting of two equal-length paths. The reason for the difference is quite simply that for some graphs the d -weight in (3.3) is 0 for the *site* graph but non-zero for the bond graph. A schematic representation of such a graph is shown in figure A1. A proof of this was given by Arrowsmith and Essam [16], who showed that $d(g)$ is non-zero

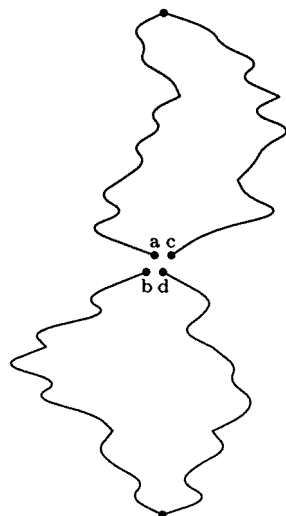


Figure A1. Schematic pictorial representation of a non-nodal graph which contributes to S^N in the bond problem but *not* in the site problem.

if and only if g is coverable by a set of directed paths *and has no circuit (or loop)*. From figure A1 we see that in the bond case the graph obtained by putting in the bonds a–b and c–d has no loops. However, in the site case there is a loop from the origin to point d and this graph does, therefore, not contribute in the site case. On the other hand it is clear that for any contributing site graph there is a corresponding contributing bond graph. So the contributing site graphs form a subset of the bond graphs.

In order to prove the formula for $s_{t,0}$ it is convenient to give another interpretation of the loop-free non-nodal graphs. Let us first characterize the graphs by the distance k between the paths. Since the graphs start and end with $k = 0$, and the distance zero appears nowhere else along the graph, these two ‘steps’ can be deleted. It is clear that in each step (increase of t by one) k changes by 0 or ± 1 . When k is unchanged there are two configurations corresponding to both paths moving either south-east or south-west, while for changes of ± 1 there is just one configuration. The non-nodal graphs are thus in bijection with paths of length $t - 1$ starting and ending at the ground level, which can take north-east, east and south-east steps, and where east steps come in two varieties or colours (such paths are known as *two-colour Motzkin paths*). It is one of the fundamental results of combinatorics that the number of two-colour Motzkin paths of length $n - 1$ is C_n . It is easy to see that loop-free non-nodal graphs form the subset where the distance between paths is never 1 twice in a row, i.e. if $k_n = 1$ then $k_{n+1} = 2$. These graphs are in bijection with two-colour Motzkin paths with no east steps on the ground level.



Figure A2. Typical two-colour Motzkin path with no east steps on the ground level.

Figure A2 shows an example of a two-colour Motzkin path with no east steps on the ground level. It is clear that all paths formed by taking the parts of the original path lying one level above the ground level (those above the dotted line), are ordinary unrestricted two-

colour Motzkin paths, and these paths are therefore enumerated by the Catalan numbers. The number of no-loop non-nodal graphs can therefore be expressed in terms of Catalan numbers, by summing over the number of times m the associated restricted two-colour Motzkin path meets the ground level prior to the terminal point. Let D_n denote the number of two-colour Motzkin paths of length n with no east steps on the ground level. The number of such two-colour Motzkin paths, $D_{n,0}$, which does not hit the ground level prior to n is simply C_{n-1} because the path obtained by deleting the first and last step is an ordinary two-colour Motzkin path of length $n - 2$. The number of restricted two-colour Motzkin paths $D_{n,1}$ which hit the ground level once is,

$$D_{n,1} = \sum_{k=0}^{n-4} C_{k+1} C_{n-4-k+1} = \sum_{i+j=n-2} C_i C_j \quad i, j \geq 1.$$

This formula is simply obtained by noting that the path to the left of the point where the restricted path meets the ground level for the first time can have a length k ranging from 0 to $n - 4$ (the four steps connecting the ground level to the level above are discarded) while the length of the second path is $n - 4 - k$. Obviously the number of left and right paths are just C_{k+1} and $C_{n-4-k+1}$, independently, which leads to the formula above once we sum over the length of the left path. The generalization to $D_{n,m}$ is obvious

$$D_{n,m} = \sum_{i_1+i_2+\dots+i_m=n-m-1} C_{i_1} C_{i_2} \dots C_{i_m} \quad i_1, \dots, i_m \geq 1, m \leq \lfloor n/2 \rfloor - 1.$$

The sum $D_n = \sum_{m=0}^{\lfloor n/2 \rfloor - 1} D_{n,m}$ is exactly the same as that obtained by Shapiro [15] by adding diagonals in the Catalan triangle.

The higher-order correction terms are quite complicated though still expressible as linear functions of $s_{t,0}$,

$$2^r (r + 1)! s_{t,r} = \sum_{k=1}^{n_a} a_{r,k} s_{t-r+k-1,0} + \sum_{k=1}^r \binom{t-r}{k} [b_{r,k} (s_{t-r-1,0} + 2s_{t-r,0}) + c_{r,k} s_{t-r,0}] \tag{A.1}$$

where $n_a = r - 1 + \max(\lfloor r/2 \rfloor, 2)$. This representation leads to particularly simple coefficients $c_{r,k}$, since $c_{r,r-m} 2^4 / (r + 1)!$ are expressible as polynomials in r of order m for $r > m$.

The sequence determining the first correction term for X^N starts out as

$$x_{t,0} = 0, 0, 0, 2, 8, 34, 136, 538, 2112, 8264, \dots$$

In this case $x_{t,0} = u(t + 1)$ is determined by the following recurrence relation

$$\begin{aligned} u(0) = 0 \quad u(1) = 0 \quad u(2) = 0 \quad u(3) = 2 \quad u(4) = 8 \\ u(t + 5) = [(2 + 4t)u(t) + (10 + 13t)u(t + 1) + (63/2 + 25/2t)u(t + 2) \\ + (4 + 2t)u(t + 3) + (-53/2 - 11/2t)u(t + 4)] / (t + 6). \end{aligned}$$

The formulae for the higher-order correction terms are complicated though still expressible as functions of $x_{t,0}$,

$$\begin{aligned} 6^{r+1} (r + 1)! x_{t,r} = \sum_{k=0}^{2r} a_{r,k} x_{t-r+k-3,0} + \sum_{k=1}^r \binom{t-r}{k} [b_{r,k} x_{t-r-4,0} + c_{r,k} x_{t-r-3,0}] \\ + (t-r)([d_{r,1} + (t-r-1)d_{r,2}/2] x_{t-r-2,0} \\ + [d_{r,3} + (t-r-1)d_{r,4}/2] x_{t-r-1,0}). \end{aligned} \tag{A.2}$$

Table A1. New series terms for the directed square lattice site problem.

n	$S(n)$	$\mu_{0,1}(p)$	$\mu_{0,2}(p)$	$\mu_{2,0}(p)$
49	-3989867121261	-89398788718610	-1997686213754238	-333247646465974
50	8580315717912	196529687645088	4051882104618304	74612004326120
51	-18450741974659	-430495003001046	-10037202119113882	-162365675913426
52	39714445919946	945547880790844	2256702296650172	359589023440940
53	-85497506155974	-207341425248340	-50294843206169288	-786814538064912
54	184179126806512	4552510986519760	11272362269382724	1735941752380332
55	-396886210803357	-99883878134103694	-2157121426296384	-380866741860840
56	835734249183509	21952260509297964	562892405947427952	838867741860840
57	-1845825000749297	-444989742917392	-1256868058330309144	-184288358054805456
58	3983384787346219	10572675215995148	2809559318045191606	40560117237064492
59	-8599553568965444	-232161203280983216	-6273640218845134156	-89163845350458120
60	18572320618806137	50989256031942796	14014639196725078868	19617995919781548
61	-4012452964388180	-1119197082819437502	-312887055698610990	-431410135430855468
62	86720156055109481	2460128423304262496	69862837357022993636	949102253844386388
63	-187496448569247473	-5404525046384365902	-15594007760170801964	-2087589817858270048
64	405534728909684450	11874838693213744120	348067895959380971188	4592790116256019372
65	-877426805166151991	-2609530626241552830	-77672638516503678130	-10103619566512091258
66	1899030133479082710	5734175392024263200	1733139806171397955248	2229577279488169258
67	-893689949790473757	-276984891538277258264	-836650128702863924260	-48907979301629134116
68	41296381922966439485	138844148831888382992	8624938634269764601752	107612749966418186776
69	-187426805166151991	-2942388348045572798924	-95660151087626485248682	-236787215239641195332
70	9059801149611631009	646984538401094502448	21328844875719496518288	48907979301629134116
71	-19643465771734195605	-14226181769514806962484	-475483808306730269156	-107612749966418186776
72	4260203192481411080	3128338155480838424564	1059887209300551355726068	1222271233383676378204
73	-20053567830421930042	-68797304028652689357280	-23623074775292445297394	-26902131263080690712678
74	92417248164200231190	15130848894126286057844	520465866276025029764160	59214386284087515891824
75	-4352442757004987018437	-332803965349353164833862	-1173171053838350921986156	-1303450921796401770801016
76	9449685901219086599817	13130848894126286057844	26140275338901762860484152	286925229989719714385800
77	-20518626196069405747527	-161037946973601504828142	-382392797578892149588184	-631637066943482168917668
78	44566294136459275950057	55427346389406871182184	1297417904728603074722100	13903430175504621295804
79	-9681818801218797898273	-1794285613503270838084276	-2890058837749845236747636	-30613948059752325054690
80	210381380688137675218788	1714913383106237923255720	64370935671258070365747636	6740206924126201296159132
81	-45724537516014414583903	-3773436434875621675302994	-146368977286469020295660	-148402323958751245043080
82	993998306718319199641	83034631348600236802398900	319260405331942765760444	326766045409443338862812
83	-216128909232339668165416	-18272485519529461160385126	-109392867626269148317469716	-148402323958751245043080
84	472003141403811031841113	10281149989876160385126	15329392487491413168166444	326766045409443338862812
85	-22247628069387596364783	-882642754529706751251248	-35243862148923843353861729	-484453331607331531822372
86	48244880069078246295706	1980622754529706751251248	78461674603860667924	58130939103038683030
87	-10531453634386083245453	-42838217180721733338482	-17466507241925214397360404	-892327091012941214
88	7931600864883250562	207351172104602760319200	3881923080511302343466	1092327091012941214
89	-469354322488309731598443	-4571681209713397983774228	-1025075109858607204334684	-18079495388276010384778
90	1087158909200163145760110	72170276476261351549340488	42361077448068108903549086	3771010905830307547520
91	-3267710909700530145760110	-72170276476261351549340488	-9137017158046036803356768	-8771010905830307547520
92	5157620497801231748845302303	107601053745140120112880988	-212217701740108069073003250	1092327091012941214
93	-113365604973511633094605416	-23669880120680053708219197562	-105070718844833469761805026	-435657049576803345388688
94	241842772053440925872069789	53122032311757013195608191058	33375871638133464934956203524	4977980397401630313004813880
95	-5335917711909810925872069789	-23669880120680053708219197562	-570020930546833508057109532542	20658062587097660145209370436
96	11633048811909810925872069789	53122032311757013195608191058	11568101389043820100553763168	45512001050464810884157750210
97	-25351064645433334857454772	-252700926082097840376086116580	-247519138306315830732243352154	1002713007529312666003007897
98	552702631363759277069727202340	5567139104414627702110090307636	11568101389043820100553763168	220025768896141879850402046812
99	-120519189490701280671238485	-2706479972852824952050183785	-572345448502435400276024258044360	4867684393234585770213494872
100	3628908431680539278383640619768	2706479972852824952050183785	12762030951814209187110090307636	1002713007529312666003007897
101	-572430669326900169571754951803	-59409788210010411406482763781852	-1272978976108650012894512185508	20658062587097660145209370436
102	150921800072863751154085013	59409788210010411406482763781852	2831193097045200748014663	4467684393234585770213494872
103	-2720255306071400069382730103133	-310437925264638135260564966662	-601619898963702063671866626049704	107254816832237290531015089289476
104	5955440926167295408476354898380	288668826410070209140674320354752	140014554490116449591789	656686308352

Table A2. New series terms for the directed triangular lattice bond problem.

n	$S(p)$	$\mu_{0,1}(p)$	$\mu_{0,2}(p)$	$\mu_{2,0}(p)$
26	5 337 497 418	209 994 728 977	28 038 948 604 228	357 799 862 456
27	11 678 931 098	487 411 142 729	68 883 587 787 794	841 629 097 226
28	25 513 719 388	1 127 362 924 089	168 327 542 017 154	1 972 059 110 234
29	55 663 119 018	2 599 086 582 635	409 289 987 873 146	4 604 235 247 626
30	121 272 163 372	5 973 768 053 766	990 554 419 328 610	10 713 215 525 118
31	263 864 408 629	13 690 809 855 903	2 386 824 242 808 628	24 848 543 707 616
32	573 556 848 773	31 292 824 198 260	5 727 568 988 920 190	57 462 309 456 098
33	1 245 063 650 267	71 342 703 947 141	13 690 818 307 565 964	132 505 664 249 544
34	2 700 144 659 216	162 261 360 324 560	32 605 625 326 065 898	304 737 782 904 598
35	5 851 221 147 909	368 214 693 911 431	77 383 096 278 813 208	699 075 297 747 540
36	12 660 942 847 609	833 758 529 144 166	183 049 343 643 929 384	1 599 836 631 974 088
37	27 392 697 005 550	1 884 144 109 110 908	431 652 603 971 595 032	3 652 954 620 022 208
38	59 166 631 983 818	4 249 400 422 872 492	1 014 868 412 269 977 442	8 322 867 118 585 614
39	127 777 294 036 668	9 566 581 135 474 702	2 379 355 385 563 105 336	18 923 690 215 681 768
40	275 696 162 276 153	21 499 276 492 272 919	5 563 403 530 205 036 262	42 943 367 206 142 286
41	594 048 482 357 433	48 233 388 196 399 900	12 974 964 525 963 569 978	97 265 602 603 253 438
42	1 281 000 979 206 493	108 047 966 744 458 962	30 186 354 559 080 349 712	219 921 104 676 935 224
43	2 755 074 940 142 431	241 645 525 989 717 809	70 064 113 568 387 529 280	496 383 864 923 234 468
44	5 932 229 201 093 542	539 692 019 601 879 166	162 259 519 144 323 831 762	1 118 569 140 266 192 598
45	12 754 620 464 996 577	1 203 634 572 376 367 923	374 966 937 946 540 768 796	2 516 752 401 957 810 240
46	27 393 502 356 280 237	2 680 685 119 486 373 279	864 732 112 976 429 729 296	5 653 852 976 905 997 716
47	58 904 482 286 533 364	5 963 270 787 963 481 223	1 990 292 162 650 597 920 198	12 683 846 242 039 392 030
48	126 300 979 513 067 199	13 247 560 344 786 965 319	4 572 211 932 174 265 999 574	28 413 833 808 390 157 526
49	271 153 388 225 432 487	29 397 708 611 765 878 122	10 484 509 048 736 986 795 242	63 570 493 940 799 673 654
50	581 799 707 017 985 602	65 162 373 599 194 694 838	23 999 926 816 621 820 432 406	142 041 285 657 057 320 738
51	1 245 200 040 883 711 881	144 265 291 339 186 480 170	54 845 072 436 992 120 826 262	316 981 854 770 124 968 722
52	2 672 296 117 689 586 731	319 107 834 898 349 284 317	125 131 020 334 445 948 974 496	706 573 223 473 121 970 044
53	5 721 610 946 798 161 890	705 067 186 518 337 735 671	285 043 213 836 022 414 418 910	1 573 161 190 417 955 836 862
54	12 219 537 226 294 787 605	1 556 202 374 122 366 410 976	648 336 112 166 418 027 074 000	3 498 618 026 159 745 044 592
55	26 278 769 763 797 603 705	3 432 580 531 634 699 049 051	1 472 529 893 791 471 135 605 612	7 773 224 302 066 420 178 488
56	55 868 130 245 151 778 098	7 561 145 873 732 448 408 790	3 339 705 956 678 263 537 822 184	17 250 739 435 533 913 221 856
57	120 005 563 753 505 676 014	16 647 643 650 693 934 045 389	7 564 345 024 108 961 163 420 714	

Table A3. New series terms for the directed triangular lattice site problem.

n	$S(p)$	$\mu_{0,1}(p)$	$\mu_{0,2}(p)$	$\mu_{2,0}(p)$
27	31 086 416	2 537 201 920	180 162 619 784	3 493 604 968
28	54 484 239	4 696 226 432	351 465 799 212	6 578 499 844
29	95 220 744	8 662 963 994	682 372 429 474	12 255 365 130
30	166 451 010	15 938 662 652	1 319 072 709 540	22 945 871 212
31	290 209 573	29 236 920 460	2 539 112 346 126	42 418 505 522
32	506 071 134	53 506 963 048	4 868 795 865 052	79 065 895 100
33	880 465 145	97 662 175 022	9 301 026 350 316	145 071 334 272
34	1 532 283 109	177 894 354 832	17 707 215 868 596	269 543 696 068
35	2 660 274 891	323 249 218 548	33 597 579 475 250	490 798 690 662
36	4 621 898 737	586 336 769 144	63 552 411 513 904	910 306 336 312
37	8 009 846 706	1 061 171 804 692	119 850 074 633 534	1 644 056 437 386
38	13 891 471 400	1 917 510 976 440	225 393 528 150 372	3 049 141 333 676
39	24 041 215 812	3 457 940 539 676	422 719 590 219 566	5 456 382 479 138
40	41 625 532 064	6 226 878 220 792	790 809 981 499 104	10 141 493 117 240
41	71 931 529 791	11 192 318 698 210	1 475 724 176 635 586	17 948 875 370 594
42	124 411 612 350	20 092 269 205 896	2 747 568 614 463 200	33 532 113 165 512
43	214 621 391 390	36 004 956 808 838	5 103 796 857 539 224	58 529 997 237 324
44	370 839 553 549	64 452 114 092 524	9 460 996 104 306 040	110 351 718 228 800
45	639 024 696 294	115 182 948 294 020	17 501 002 169 903 066	189 161 996 834 038
46	1 102 419 174 084	205 638 719 322 044	32 311 701 334 358 584	361 978 973 535 312
47	1 898 477 439 658	366 587 483 305 266	59 540 588 349 689 460	605 431 024 385 712
48	3 271 434 676 999	652 904 591 166 608	109 522 752 581 367 792	1 185 609 582 832 608
49	5 624 820 363 027	1 161 134 164 194 872	201 098 347 347 198 582	1 916 175 057 214 282
50	9 693 710 116 271	2 063 632 450 148 240	368 654 569 738 994 916	3 885 789 400 216 356
51	16 634 472 160 666	3 661 795 173 290 544	674 667 552 855 892 942	5 981 962 784 372 730
52	28 649 053 574 116	6 494 555 752 892 524	1 232 887 441 544 215 856	12 779 152 925 915 688
53	49 158 925 607 599	11 502 147 999 885 690	2 249 412 773 359 085 386	18 336 104 911 125 754
54	84 477 695 445 892	20 358 932 047 872 636	4 098 441 587 758 882 072	42 326 707 460 800 448
55	144 947 819 272 120	35 990 408 059 294 200	7 456 350 674 610 337 790	54 742 323 913 847 946
56	249 148 051 950 911	63 598 870 606 450 408	13 548 513 117 372 733 000	

From the polynomials for $S^N(t_{\max})$ and $X^N(t_{\max})$ with $t_{\max} = 47$, and using the extrapolation formulae, I extended the series for $S(p)$, $\mu_{0,1}(p)$ and $\mu_{0,2}(p)$ to order 106 and the series for $\mu_{2,0}(p)$ to order 103. The new series terms are listed in table A1.

A.2. The triangular bond problem

The correction terms for the triangular bond problem are very simple. The first correction term for S^N is just a constant $s_{t,0} = 2$, while the first correction term for X^N alternates between 0 and 2. The non-nodal graphs responsible for these correction terms are almost trivial. It is clear (see figure 1) that the non-nodal graphs terminating at level $t + 1$ having the smallest possible number of bonds are those composed of two paths meeting on the centre line (t odd) or on the site next to the centre-line (t even), with each path having as few south-east and south-west steps as possible. These sites can be reached by a non-nodal graph with $t + 1$ bonds. For t odd the only two such graphs are those consisting of a path with $\lfloor t/2 \rfloor + 1$ south steps and a path starting with a south-east (south-west) step followed by $\lfloor t/2 \rfloor$ south steps, while ending with a south-west (south-east) step. For t even, the two graphs are those consisting of a path with $\lfloor t/2 \rfloor$ south steps terminating with a south-east (south-west) step and a path starting with a south-east (south-west) step followed by $\lfloor t/2 \rfloor$

south steps. It is easy to check that any other non-nodal graph contains more bonds. So $s_{t,0} = 2$ while $x_{t,0}$ alternate between 0 and 2 since for t odd the non-nodal graphs terminate on the centre-line and therefore do not contribute to X^N .

The sequence determining the second correction terms for S^N is

$$1, 2, 5, 10, 17, 26, 37, 50, 65, \dots$$

from which it is clear that $s_{t,1}$ grows as a polynomial in t , $s_{t,1} = t^2 - 2t + 2$. In general the correction terms can be represented as a polynomial in t of order $2r$. The alternation between odd and even values of t seen in $x_{t,0}$ eventually also manifests itself in the correction terms for S^N . The general formulae for the correction term is,

$$s_{t,r} = \frac{1}{r!(r+1)!} \sum_{j=0}^{2r} a_{r,j}(t-1)^j + \frac{t \bmod 2}{r!(r+1)!} \sum_{j=0}^{\lfloor (r-3)/2 \rfloor} b_{r,j}(t-1)^j \quad t \geq 2r - 2. \quad (\text{A.3})$$

The prefactors and the expression of the polynomials in terms of $t - 1$ has been chosen to make the leading coefficients particularly simple. Once again it should be noted that the leading coefficients $a_{r,2r-m}$ are polynomials in r of order $m + \lfloor m/2 \rfloor$ (this is valid for $m \leq 5$), which again was used to obtain a few additional correction formulae.

The extrapolation formulae for X^N are very similar to the ones above,

$$x_{t,r} = \frac{1}{r!(r+1)!} \sum_{j=0}^{2r} a_{r,j}(t-1)^j + \frac{t \bmod 2}{r!(r+1)!} \sum_{j=0}^r b_{r,j}(t-1)^j \quad t \geq 2r - 2. \quad (\text{A.4})$$

In this case the leading coefficients of both $a_{r,2r-m}$ and $b_{r,r-m}$ can be predicted. For $r > m$ I find that $a_{r,2r-m}$ can be expressed as a polynomial in r of order $\leq m + 2$. Likewise $(-1)^r b_{r,r-m}/(r+1)!$ is a polynomial in r of order $2m$ for $r > 2m$.

As stated earlier, the non-nodal contribution to the series for the triangular bond case were calculated up to $t_{\max} = 45$. With the extrapolation formulae I derived the series for $S(p)$, $\mu_{0,1}(p)$ and $\mu_{0,2}(p)$ to order 57 and the series for $\mu_{2,0}(p)$ to order 56. The resulting new series terms are listed in table A2.

A.3. The triangular site problem

In this case the first correction term for S^N alternates between 0 and 1 while the first correction term for X^N is 0. The graphs giving rise to these correction terms are very simple. First note that the graphs giving rise to the bond correction terms all have loops when viewed as site graphs. The non-nodal site graphs with fewest elements for t odd consist of the two paths starting with a south-east (south-west) step followed by $\lfloor t/2 \rfloor$ south steps and ending with a south-west (south-east) step. These graphs have $t + 2$ random elements (remember that the origin is not a random element). For t even one can easily see that there are no loop-free non-nodal graphs with $t + 2$ or fewer elements. This fully accounts for the first correction terms.

The other extrapolation formulae for the triangular site problem are very similar to those for the bond case. The only difference is that the order of the polynomials correcting the odd- t values is somewhat higher. Once again the leading coefficients are low-order polynomials in r . With the help of the extrapolation formulae I extended the series for $S(p)$, $\mu_{0,1}(p)$ and $\mu_{0,2}(p)$ to order 56 and the series for $\mu_{2,0}(p)$ to order 55. The new series terms are listed in table A3.

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