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Preferential Attachment with Location-Based Choice: Degree Distribution in the Noncondensation Phase

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Abstract

We consider the preferential attachment model with location-based choice introduced by Haslegrave et al. (Random Struct Algorithms 56(3):775–795, 2020) as a model in which condensation phenomena can occur. In this model, each vertex carries an independent and uniformly distributed location. Starting from an initial tree, the model evolves in discrete time. At every time step, a new vertex is added to the tree by selecting *r* candidate vertices from the graph with replacement according to a sampling probability proportional to these vertices' degrees. The new vertex then connects to one of the candidates according to a given probability associated to the ranking of their locations. In this paper, we introduce a function that describes the phase transition when condensation can occur. Considering the noncondensation phase, we use stochastic approximation methods to investigate bounds for the (asymptotic) proportion of vertices inside a given interval of a given maximum degree. We use these bounds to observe a power law for the asymptotic degree distribution described by the aforementioned function. Hence, this function fully characterises the properties we are interested in. The power law exponent takes the critical value one at the phase transition between the condensation–noncondensation phase.

Keywords Choice · Phase transition · Power law · Simulation

Mathematics Subject Classification 05C80

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1 Introduction

The study of complex networks is a prevalent area of interest for researchers as many seemingly dissimilar structures observable in the real world can be modelled using a common set of techniques. This is due to many large networks sharing similar topological properties. For instance, it has been observed that the empirical degree distribution of many large-scale real world networks follows an approximate *power law* over a large finite range of degrees. Hence, we seek families of models that imitate this behaviour.

A probabilistic approach is to build networks as a growing sequence of graphs in which the degree distribution follows a power-law when the number of vertices is going to infinity. That is, the tail of the asymptotic proportion of vertices of degree at least k behaves like $k^{-\tau}$ for some power-law exponent τ . We call such a network *scale-free*.

In 1999 Barabási and Albert popularised *preferential attachment* [1] as a method of growth which utilises the famous rich get richer concept. As a new vertex joins the network, it forms an edge to already existing vertices with probability proportional to the degrees of current vertices. This mechanism was generalised by Dorogovtsev et al. [8] by biasing the selection mechanism to enhance or suppress the influence of the degrees. It was shown by various authors that this building mechanism indeed leads to scale-free networks [3,8,12]. Although preferential attachment is often an accurate method of modelling scale-free networks, it fails to consider a new vertex's potential to attract new edges. In order to tackle this issue, Bianconi and Barabási [2] suggested the addition of *vertex fitness* as an additional parameter. Here, each vertex joins the network with its own randomly chosen fitness, allowing for a new level of competition between vertices, separate from their current edge-based popularity. Many models have been devised which include this 'attractiveness' coefficient, most notably by Borgs et al. [4] and Dereich and Ortgiese [6]. Another way of incorporating a vertex's inherent potential for growth is by introducing the notion of choice. In [15-17] preferential attachment is used to sample a set of vertices from the network as candidates for connection. Afterwards, a preassigned attachment rule based on the degrees of the sampled vertices is used to decide where new edges are formed.

Furthermore, a feature of interest is the *condensation* phenomenon. Condensation occurs if the total degree of an o(n) subset of vertices grows linearly in time n. Loosely speaking, at any time there exists some vertex whose degree dramatically dominates the others. Whereas in classical preferential attachment condensation cannot occur, it was shown that both preferential attachment with choice and models with fitness can exhibit condensation [4–7,10].

In this paper, we consider the preferential attachment with location-based choice model introduced by Haslegrave, Jordan and Yarrow in [11] which can be seen as a generalised variant of [9]. This model combines the ideas of both fitness and choice in a natural way. Starting from an initial tree graph, at each time step a new vertex joins the graph and is assigned its own location which is uniformly chosen from (0, 1). When this vertex joins the network, a subset of *r* neighbour candidates is sampled with probability proportional to their degree plus some constant α . The sampled vertices are ranked according to their locations. Following this, a single vertex from the sample is chosen for connection to the new vertex according to some probability measure Ξ . Here, Ξ can be used to make different regions of (0, 1) more or less appealing and thus incorporates more flexibility than in previous models. As in [11], we refer to location as opposed to fitness in order not to give the false impression of preferring the 'fittest' vertex. We could choose any continuous distribution on the real line but we do not expect any changes in the results as the connection mechanism only depends

on the ordering of the vertices' locations and not their actual value. Hence, there is no loss in generality by restricting the locations distribution to the uniform one on (0, 1). This has the added benefit that it matches previous work our results build on. From [11], it can be derived that there exists a critical value α_c such that condensation can occur if $\alpha < \alpha_c$.

In this article, we give a new description of α_c and study the degree-distribution of this model in the noncondensation regime. We show that in the noncondensation case, $\alpha \ge \alpha_c$, the model is asymptotically scale-free with a heavy tailed degree distribution with power-law exponent $\tau = \frac{2+\alpha}{2+\alpha_c}$. Hence, the critical value α_c for the condensation phase transition matches the one for which the power-law exponent is large enough for the degree distribution's first moment to exist. This behaviour coincides with our understanding of condensation. In the condensation phase, with positive probability a proportionally small number of vertices dominate the others. The noncondensation phase is 'regular' in the sense that a typical vertex has finite expected degree. As the behaviour of the degree distribution dramatically changes between the two phases, we lose the finite moments at that phase transition, even though for $\alpha = \alpha_c$ the network is still scale-free with $\tau = 1$. The same behaviour can be observed in similar models with choice [15]. Although a power-law distribution is what one would hope for in the considered regime, it is notable that this is not the case in the original preferential attachment model with choice of Malyshkin and Paquette for more than two options [16].

To derive the degree distribution, we introduce a function f on the location space (0, 1) depending only on Ξ that plays a key role in understanding the influence of location on the degree of a vertex. Given a vertex with location x, the expected probability of choosing that vertex with respect to Ξ , out of a sample containing this vertex and r - 1 uniformly located vertices is given by f(x)/r. We show that the condensation phase transition as well as the power-law exponent can be derived from the maximum value of f. To get this, we determine the concrete degree distribution of a vertex at a given location whose tail behaviour follows a power-law distribution dependent on f from which we derive the final result. The function f in a specific region, the more likely we are to find high degree vertices there. The question of the degree distribution in the condensation phase is also of some interest but cannot be achieved with our methods since we rely on some continuity properties in our proof that are not fulfilled in the condensation regime.

The paper is structured as follows: In Sect. 2 we formally introduce the model and formulate the main theorem. Afterwards we recall the phase transition conditions determined in [11]. We introduce formally the function f and rewrite these conditions. In Sect. 3, we use stochastic approximation methods to deduce bounds of the growth of the empirical degree distribution. We use these bounds to deduce the asymptotic degree distribution, proving the main theorem. In the last section, we show numerical results and simulations for some interesting and important choices of Ξ underlining our understanding and results.

2 Model Description and Main Result

Let $r \ge 2$ be an initial integer model parameter and let Ξ be a probability measure on $\{1, \ldots, r\}$. In the following, we treat Ξ as a probability vector (Ξ_1, \ldots, Ξ_r) . Furthermore, let G_0 be an initial tree graph on $n_0 \ge 2$ vertices $\{v_{1-n_0}, \ldots, v_0\}$. Additionally, let each vertex v_i in G_0 have its own location x_i that is drawn independently and uniformly at random from (0, 1) and is therefore almost surely unique.

At time n+1, a new vertex v_{n+1} assigned its own location x_{n+1} , again drawn independently and uniformly at random from (0, 1) is added to the graph. Given G_n and the locations of all its vertices, we form the graph G_{n+1} by connecting the new vertex v_{n+1} by a single edge to a vertex in G_n . Note that this maintains the tree structure of the graph. The connection mechanism is as follows: First, we sample r candidate vertices with replacement from G_n according to preferential attachment, i.e. proportional to the vertices' degrees plus a fixed constant α . Second, v_{n+1} chooses one vertex for connection out of the sample according to Ξ applied to the ranks of the locations. More precisely, fix $\alpha \in (-1, \infty)$ and denote deg $_{G_n}(v_j)$ as the degree of vertex v_j in G_n . We first select a sample of r candidate vertices from G_n with replacement so that independently for each of the r candidates

$$\mathbb{P}(v_i \text{ is sampled } | G_n) = \frac{\deg_{G_n}(v_i) + \alpha}{(n + n_0 - 1)(2 + \alpha) + \alpha}.$$
(1)

Here, due to the tree structure, the denominator equals the total degree weight of G_n , that is the sum over each vertices' degree plus α . We next order the *r* sampled vertices according to their location. That is, we obtain a sample of vertices $(v_1^{(n+1)}, \ldots, v_r^{(n+1)})$ and associated locations $(x_1^{(n+1)}, \ldots, x_r^{(n+1)})$ such that the locations satisfy $x_1^{(n+1)} \leq \cdots \leq x_r^{(n+1)}$. An important observation is that equality for the locations happens almost surely only if a vertex has been sampled multiple times. Thus, the ordered sample is uniquely determined. Finally, according to Ξ one vertex out of $(v_1^{(n+1)}, \ldots, v_r^{(n+1)})$ is chosen for connection. That is, the probability that vertex $v_i^{(n+1)}$ is chosen for connection is given by Ξ_i .

2.1 Main Result

As mentioned in the introduction, it is known that there exists a threshold α_c such that condensation can only occur if and only if $\alpha < \alpha_c$, see [11] and Proposition 2.3 below. Let μ_k be the asymptotic proportion of vertices of degree at least *k*.

Theorem 2.1 If $\alpha \ge \alpha_c$, then μ_k exists and satisfies

$$\mu_k = k^{-\frac{2+\alpha}{2+\alpha_c} + o(1)},$$

as $k \to \infty$.

In order to prove this result, one has to understand the influence of the location on a vertex's degree. To this end, define $\Psi_n(x)$ as the conditional probability, given the graph G_n and the locations of all the vertices of G_n , that the new vertex v_{n+1} selects under preferential attachment according to equation (1) a vertex which has location at most x. Denote by $V(G_n)$ the vertex set of G_n . Then it holds that

$$\Psi_n(x) = \frac{1}{(n+n_0-1)(2+\alpha)+\alpha} \left(\sum_{v_i \in V(G_n): x_i \le x} (\deg_{G_n}(v_i)+\alpha) \right).$$
(2)

The random measures induced by $\Psi_n(x)$ converge weakly almost surely to a probability measure on [0, 1], whose continuous distribution function we call $\Psi(x)$ [11, Theorem 2.2]. Here, it is important to note that in general $\Psi(x)$ may be random. However, it is not random whenever $\alpha \ge \alpha_c$. Finally, define the function $f : [0, 1] \to \mathbb{R}_+$ by

$$f(x) = \sum_{s=1}^{r} s \Xi_s {\binom{r}{s}} x^{s-1} (1-x)^{r-s} , \qquad (3)$$

(for more details about Ψ_n , Ψ and f, we refer the reader to Sect. 2.2.) Conditioned on the event that there is a vertex at a given location x, we denote by $\nu(k, x)$ the probability that the vertex at location x has asymptotically at least k neighbours.

Theorem 2.2 If $\alpha \ge \alpha_c$ and $x \in (0, 1)$, then v(k, x) is well-defined and satisfies

$$v(k, x) = k^{-\frac{2+\alpha}{f(\Psi(x))} + o(1)}$$

as $k \to \infty$.

2.2 Condensation Phase Transition

According to (2), $\Psi_n(x)$ is almost surely monotonically increasing with $\Psi_n(0) = 0$ and $\Psi_n(1) = 1$. Hence, we can think of $\Psi_n(x)$ as a random distribution function on the location space. The measures induced by $\Psi_n(x)$ converge weakly almost surely to a (possibly random) probability measure on [0, 1]. We call the distribution function of this limit $\Psi(x)$.

We define condensation as a *discontinuity* in Ψ since a jumping point of Ψ implies that Ψ_n increases by O(1) on an interval of length o(1), as $n \to \infty$, matching the condensation description given in the introduction. Here, condensation may arise due to the existence of a persistent hub [11, Theorem 2.3] as well as without a hub where the currently leading vertex is replaced over time [11, Theorem 2.4]. The following proposition summarises arguments from [11] showing that the discontinuity of Ψ can only occur if α is smaller than the threshold α_c . Additionally, it gives a new description of this threshold. We call $(-1, \alpha_c)$ the *condensation* and $[\alpha_c, \infty)$ the *noncondensation phase* of the model (Fig. 1).

The function f (see (3)) only depends on the model parameter Ξ and plays an important role in characterising the condensation phase transition. It can be observed that f is a probability density on [0, 1]. Rewriting the binomial coefficient, one can interpret f(x)/r as the expected probability of connecting with respect to Ξ to a given vertex of location x where the remaining r - 1 vertices are chosen uniformly.

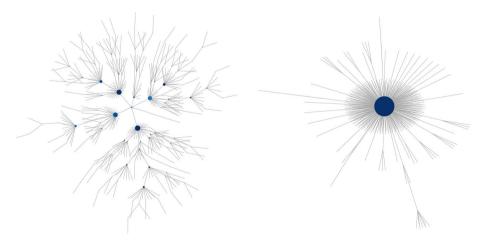


Fig. 1 Plots of a simulated tree for $\Xi = (0, 1, 0)$ after 500 vertices have been added. On the left, a realization for $\alpha > \alpha_c$ and on the right, a realization with $\alpha < \alpha_c$. In both cases, the start configuration consists of a root vertex and a single child, both with uniform drawn location. In the plot, the size of a vertex corresponds to its degree. We use colour saturation to indicate how close to the maximum value of *f* a vertex's location is

Proposition 2.3 There exists $\alpha_c \geq -1$ such that condensation can occur with a positive probability if $\alpha < \alpha_c$ but cannot occur if $\alpha \geq \alpha_c$. Moreover.

$$\alpha_c = \max\{f(y) : y \in [0, 1]\} - 2.$$
(4)

Proof As a function of $y \in [0, 1]$, we define

$$F_1(y; x, \Xi) = x(\alpha + 1) - (2 + \alpha)y + \sum_{s=1}^r \Xi_s \sum_{i=s}^r \binom{r}{i} y^i (1 - y)^{r-i}$$

for $x \in [0, 1]$. By [11, Theorem 2.2], $\Psi_n(x)$ converges almost surely to a zero of the function $F_1(y; x, \Xi)$ and by [11, Theorems 2.3, 2.4] condensation occurs with positive probability, whenever there exists $x \in (0, 1)$ such that $F_1(y; x, \Xi)$ has a touchpoint. Here, we call $p \in (0, 1)$ a touchpoint if $F_1(p; x, \Xi) = 0$ and there exists $\varepsilon > 0$ such that either $F_1(y; x, \Xi) < 0$ for all $y \in (p - \varepsilon, p + \varepsilon) \setminus \{p\}$ or $F_1(y; x, \Xi) > 0$ for all $y \in (p - \varepsilon, p + \varepsilon) \setminus \{p\}$. Note that if $F_1(y; x, \Xi)$ is increasing in y somewhere on [0, 1], one can vary x in such a way that $F_1(y; x, \Xi)$ has a touchpoint. Hence, condensation can occur with positive probability for $\alpha < \alpha_c$, where

$$\alpha_c = \inf\{\alpha > -1 : F_1'(y; x, \Xi) \le 0 \text{ for all } y \in (0, 1)\},\$$

see also [11, p. 792]. Conversely, if $\alpha \ge \alpha_c$, then, for all $x \in [0, 1]$, $F_1(y; x, \Xi)$ has only one zero to which $\Psi_n(x)$ converges almost surely. Since F_1 is continuous and strictly decreasing in the neighbourhood of the root, the zero Ψ is continuous and almost surely no condensation can occur, proving the first part of the proposition. To prove (4), we calculate

$$F'_{1}(y; x, \Xi) = -(2+\alpha) + \sum_{s=1}^{r} \Xi_{s} \sum_{i=s}^{r} i\binom{r}{i} y^{i-1} (1-y)^{r-i} \left(1 - \frac{r-i}{i} y(1-y)^{-1}\right)$$
$$= -(2+\alpha) + \sum_{s=1}^{r} s \Xi_{s} \binom{r}{s} y^{s-1} (1-y)^{r-s} = -(2+\alpha) + f(y).$$

Hence, $F'_1(y; x, \Xi) \le 0$ holds for all $y \in (0, 1)$ if and only if $2 + \alpha \ge \max\{f(y) : y \in [0, 1]\}$.

Proposition 2.3 shows that both preferential attachment and the location-based choice aspect are necessary for condensation to occur. Sampling according to preferential attachment but then choosing one vertex independently of the locations coincides with the choice of $\Xi_s =$ 1/r. Then, $f(x) \equiv 1$ and thus $\alpha_c = -1 < \alpha$ for all $\alpha \in (-1, \infty)$. By Theorem 2.1, the network is then scale-free with power-law exponent $\tau = 2 + \alpha \in (1, \infty)$, matching the results of [8]. On the other hand, sampling without preferential attachment in this model coincides with the case $\alpha \to \infty$. Therefore, it holds $\alpha > \alpha_c$ for all choices of Ξ . Summarizing the above, whenever $\alpha \ge \alpha_c$, no condensation can occur and the limiting distribution Ψ is continuous and non-random. This is shown to be important in following sections.

3 Noncondensation Phase Degree Distribution

We utilize a number of stochastic approximation techniques constructed by Robbins and Monro [19] outlined in Pemantle [18, Section 2]. For a stochastic process $(X_n)_{n \in \mathbb{N}} \subset \mathbb{R}^n$ adapted to a filtration $(\mathcal{F}_n)_{n \in \mathbb{N}}$, the idea of stochastic approximation is to find a representation of the increments $X_{n+1} - X_n$ which fulfills certain properties. This then allows for results on the asymptotic behaviour of the process to be derived. Classically, we call equations of the form

$$X_{n+1} - X_n = \frac{1}{n} \left(F(X_n) + \xi_{n+1} + r_n \right)$$

stochastic approximation equations. Here, F is an \mathbb{R}^n -vector field, ξ_{n+1} is a noise term satisfying $\mathbb{E}[\xi_{n+1} | \mathcal{F}_n] = 0$ and the remainder term r_n is \mathcal{F}_n -measurable, converging to zero and satisfying $\sum_n n^{-1} |r_n| < \infty$ almost surely. Depending on the properties of F and possible further assumptions on the noise ξ_{n+1} , different results for the asymptotic behaviour of the process are known [18, Section 2]. Many results can be further extended to hold, when F is random, see e.g. [11,13].

In our setting, we will need a statement for the asymptotic behaviour when only bounds on the increments are given. To this end, we adapt Lemma 5.4. of [14] by Jordan and Wade.

Lemma 3.1 Let $(\mathcal{F}_n)_{n \in \mathbb{N}_0}$ be a filtration. Furthermore, let $X = (X_n)_{n \in \mathbb{N}_0}$, $A_1 = (A_1^{(n)})_{n \in \mathbb{N}_0}$, $A_2 = (A_2^{(n)})_{n \in \mathbb{N}}$, $K_1 = (K_1^{(n)})_{n \in \mathbb{N}_0}$, $K_2 = (K_2^{(n)})_{n \in \mathbb{N}_0}$, $\xi = (\xi_n)_{n \in \mathbb{N}_0}$, $r_1 = (r_1^{(n)})_{n \in \mathbb{N}_0}$ and $r_2 = (r_2^{(n)})_{n \in \mathbb{N}_0}$ be real-valued stochastic processes adapted to $(\mathcal{F}_n)_{n \in \mathbb{N}_0}$ where X, A_1 , A_2 , K_1 and K_2 are non-negative and bounded. Let $(\gamma_n)_{n \in \mathbb{N}_0}$ be a sequence of non-negative constants and suppose that

$$\gamma_n (A_1^{(n)} - K_1^{(n)} X_n + \xi_{n+1} + r_1^{(n)}) \le X_{n+1} - X_n \le \gamma_n (A_2^{(n)} - K_2^{(n)} X_n + \xi_{n+1} + r_2^{(n)}).$$
(5)

Assume further that

- (i) $\mathbb{E}[\xi_{n+1} | \mathcal{F}_n] = 0$ and $\mathbb{E}[\xi_{n+1}^2 | \mathcal{F}_n] \leq C$ for a finite constant C,
- (ii) $\sum_{n=0}^{\infty} \gamma_n = \infty$, $\sum_{n=0}^{\infty} \gamma_n^2 < \infty$ and $\sum_{n=0}^{\infty} |r_i^{(n)}| \gamma_n < \infty$ almost surely (i = 1, 2),
- (iii) $0 < \ell_i \le K_i^{(n)} \le u_i$ for some finite constants ℓ_i , u_i (i = 1, 2),
- (iv) $\liminf_{n \to \infty} \frac{A_1^{(n)}}{K_1^{(n)}} \ge L_1 \text{ and } \limsup_{n \to \infty} \frac{A_2^{(n)}}{K_2^{(n)}} \le L_2 \text{ almost surely.}$

Then, almost surely,

$$L_1 \leq \liminf_{n \to \infty} X_n \leq \limsup_{n \to \infty} X_n \leq L_2.$$

Proof We only prove the lower bound for the lim inf as the upper bound for the lim sup works with analogous argumentation. For $\varepsilon > 0$ there exists an almost surely finite N_1 such that $L_1 \le A_1^{(n)}/K_1^{(n)} + \varepsilon/2$ for $n \ge N_1$ by (iv). For each $x < L_1 - \varepsilon$ and $n \ge N_1$, we have

$$A_1^{(n)} - K_1^{(n)} x \ge A_1^{(n)} - K_1^{(n)} (L_1 - \varepsilon) \ge A_1^{(n)} - K_1^{(n)} \left(\frac{A_1^{(n)}}{K_1^{(n)}} - \frac{\varepsilon}{2}\right) \ge \frac{\ell_1 \varepsilon}{2} > 0,$$

using (iii). Now, summing (5), we get $X_n - X_0 \ge M_n + O_n$, where

$$M_n = \sum_{k=0}^{n-1} \gamma_k \xi_{k+1}$$
 and $O_n = \sum_{k=0}^{n-1} \gamma_k (A_1^{(k)} - K_1^{(k)} X_k + r_1^{(k)}).$

Here, O_n is \mathcal{F}_{n-1} -measurable and M_n is a martingale satisfying

$$\mathbb{E}[M_{n+1}^2 - M_n^2 \mid \mathcal{F}_n] = \mathbb{E}[(M_{n+1} - M_n)^2 \mid \mathcal{F}_n] = \mathbb{E}[\xi_{n+1}^2 \gamma_n^2 \mid \mathcal{F}_n] \le C\gamma_n^2$$

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by (i). Since γ_n^2 is summable by (ii), M_n is L^2 -bounded and hence there exists a finite M_∞ such that $M_n \to M_\infty$ almost surely as $n \to \infty$. Moreover, by (ii), we have

$$R_n := \sum_{k=0}^{n-1} \gamma_k r_1^{(k)} \to R_\infty < \infty,$$

almost surely as $n \to \infty$. Hence, for the ε above, there exists N_2 such that

$$\sup_{n \ge N_2} \sup_{m \ge 0} |M_{n+m} - M_n| \le \frac{\varepsilon}{4} \text{ and } \sup_{n \ge N_2} \sup_{m \ge 0} |R_{n+m} - R_n| \le \frac{\varepsilon}{4}.$$

Now, fix some $n_0 \ge N := N_1 \lor N_2$ for which $X_{n_0} < L_1 - \varepsilon$. Let $\kappa_{n_0} := \min\{t > n_0 : X_t \ge L_1 - \varepsilon\}$ be the first time after n_0 for which X returns to $[L_1 - \varepsilon, \infty)$. Then, for $m \ge 0$, using (5), we have

$$\begin{aligned} X_{(n_0+m)\wedge\kappa_{n_0}} - X_{n_0} &\geq \left(M_{(n_0+m)\wedge\kappa_{n_0}} - M_{n_0} \right) + \left(R_{(n_0+m)\wedge\kappa_{n_0}} - R_{n_0} \right) \\ &+ \sum_{k=n_0}^{((n_0+m)\wedge\kappa_{n_0})-1} \gamma_k (A_1^{(k)} - K_1^{(k)} X_k) \\ &\geq -\frac{\varepsilon}{2} + \frac{\ell_1 \varepsilon}{2} \sum_{k=n_0}^{((n_0+m)\wedge\kappa_{n_0})-1} \gamma_k. \end{aligned}$$

On $\{\kappa_{n_0} = \infty\}$, for $m \to \infty$, the left-hand-side remains finite since X is a bounded process; however the right-hand-side diverges to infinity by (ii). Hence, κ_{n_0} is almost surely finite. Furthermore, since $X_{(n_0+m)\wedge\kappa_{n_0}} \ge X_{n_0} - \varepsilon/2$ using the above calculation and $\gamma_k \ge 0$ for all k, the process X returns to $[L_1 - \varepsilon, \infty)$ without dropping below $X_{n_0} - \varepsilon/2$. Moreover, for large enough $n \ge N$,

$$X_{n+1} - X_n \ge -\frac{\varepsilon}{2} + \gamma_n (A_1^{(n)} - K_1^{(n)} X_n) \ge -\varepsilon,$$

as $A_1 - K_1 X$ is bounded and γ_n tends to zero. Therefore, almost surely, $X_n \ge L_1 - \varepsilon$ infinitely often, and for all except a finite number of *n* any exit from $[L_1 - \varepsilon, \infty)$ cannot drop under $L_1 - 2\varepsilon$; but starting from $[L_1 - 2\varepsilon, L_1 - \varepsilon)$, the process *X* returns to $[L_1 - \varepsilon, \infty)$ before hitting $L_1 - 3\varepsilon$. Hence, $\lim \inf X_n \ge L_1 - 3\varepsilon$, almost surely. Since ε was chosen arbitrarily, this concludes the proof.

3.1 Bounds on the Empirical Degree Distribution

The aim of this section is to find bounds for the proportion of the vertices of degree at least k located inside $[x_1, x_2] \subset (0, 1)$. To this end, we define $P_{x_1, x_2}^{(n)}(k)$ as the proportion of vertices in G_n that have degree at most k and are located inside the interval $[x_1, x_2]$, that is

$$P_{x_1,x_2}^{(n)}(k) = \frac{1}{n+n_0} \sum_{(v,x)\in G_n} \mathbb{1}_{\{\deg_{G_n}(v)\leq k\}} \mathbb{1}_{\{x\in[x_1,x_2]\}}.$$

To get bounds on $P_{x_1,x_2}^{(n)}(k)$, we define the event that the new vertex v_{n+1} , arriving at time n+1, connects to a vertex of degree k in G_n which is located inside $[x_1, x_2]$. We denote this event by E_{n+1} . We cannot give a precise description of the probability of E_{n+1} , however we can bound it from above and below in a natural way. The estimations are made in the part in

which one of the r candidates of the sample is chosen for connection. First, we consider the function

$$f_1(y_1, y_2) := \sum_{s=1}^r \Xi_s \left(\sum_{j=0}^{s-1} \sum_{i=s}^r \binom{r}{i} \binom{i}{j} y_1^j (y_2 - y_1)^{i-j-1} (1 - y_2)^{r-i} \right).$$
(6)

Upon multiplying f_1 by $(y_2 - y_1)$, the term inside the outer brackets states that for r points ranked by 1, ..., r, the first j points are sampled from the interval $(0, y_1)$, the next i - j points (including point s) are sampled from $[y_1, y_2]$ and the remaining i points are from $(y_2, 1)$. Since we do not consider the precise ordering within the intervals, an upper bound for the probability of choosing a vertex for connection from the candidate sample with location inside $[x_1, x_2]$ is given by

$$f_1(\Psi_n(x_1), \Psi_n(x_2))(\Psi_n(x_2) - \Psi_n(x_1)).$$

Secondly, we denote \mathbb{P}_n as the conditional probability measure given by the graph G_n and all locations of vertices contained within. We also denote w_n as the vertex which has been chosen for connection at time n + 1. We have

$$\mathbb{P}_n(w_n \text{ has degree } k \mid w_n \text{ is located inside } [x_1, x_2])$$

$$= \frac{\mathbb{P}_n(w_n \text{ has degree } k \text{ and is located inside } [x_1, x_2])}{\Psi_n(x_2) - \Psi_n(x_1)}$$

$$= \frac{k + \alpha}{(2 + \alpha)(n + n_0) - 2}(n + n_0) \left(P_{x_1, x_2}^{(n)}(k) - P_{x_1, x_2}^{(n)}(k - 1)\right) \frac{1}{\Psi_n(x_2) - \Psi_n(x_1)}.$$

This holds because the first factor is the probability in which a vertex of degree k is sampled according to (1) and $(n + n_0) (P_{x_1,x_2}^{(n)}(k) - P_{x_1,x_2}^{(n)}(k - 1))$ counts the number of degree k vertices with corresponding locations inside the interval $[x_1, x_2]$ in G_n . Therefore, both parts together yield

$$\mathbb{P}_{n}(E_{n+1}) \leq \frac{(k+\alpha)(n+n_{0})}{(2+\alpha)(n+n_{0})-2} \left(P_{x_{1},x_{2}}^{(n)}(k) - P_{x_{1},x_{2}}^{(n)}(k-1) \right) f_{1}(\Psi_{n}(x_{1}),\Psi_{n}(x_{2})).$$
(7)

Similarly, we can achieve a lower bound for $\mathbb{P}_n(E_{n+1})$ if we only consider samples of candidates where exactly one vertex is located inside $[x_1, x_2]$. Thus, we obtain

$$\mathbb{P}_{n}(E_{n+1}) \geq \frac{(k+\alpha)(n+n_{0})}{(2+\alpha)(n+n_{0})-2} \left(P_{x_{1},x_{2}}^{(n)}(k) - P_{x_{1},x_{2}}^{(n)}(k-1) \right) f_{2}(\Psi_{n}(x_{1}),\Psi_{n}(x_{2})), \quad (8)$$

where f_2 is given by

$$f_2(y_1, y_2) = \sum_{s=1}^r s \Xi_s \binom{r}{s} y_1^{s-1} (1 - y_2)^{r-s}.$$
(9)

With these bounds, which will be crucial for the asymptotic degree later, we are ready to proceed to the stochastic approximation.

Lemma 3.2 Let $k \in \mathbb{N}$ and $0 < x_1 < x_2 < 1$. Define for $j \in \{1, 2\}$ the random variables

$$A_j^{(n)} = A_j^{(n)}(k) := \frac{k+\alpha}{2+\alpha} f_j \left(\Psi_n(x_1), \Psi_n(x_2)\right) P_{x_1, x_2}^{(n)}(k-1) + (x_2 - x_1)$$

and

$$K_j^{(n)} = K_j^{(n)}(k) := 1 + \frac{k+\alpha}{2+\alpha} f_j(\Psi_n(x_1), \Psi_n(x_2)),$$

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for $n \in \mathbb{N}_0$. Let \mathcal{F}_n be the filtration generated by the sequence of graphs $(G_i, x_i; i \leq n)$ and, for $n \in \mathbb{N}_0$,

$$\xi^{(n+1)} = (n+n_0+1) \left(P_{x_1,x_2}^{(n+1)}(k) - \mathbb{E} \left[P_{x_1,x_2}^{(n+1)}(k) \big| \mathcal{F}_n \right] \right).$$

Then, for the growth of the proportion of vertices with degree at most k and location inside $[x_1, x_2]$, it holds

$$\frac{A_1^{(n)} - K_1^{(n)} P_{x_1, x_2}^{(n)}(k) + \xi^{(n+1)} - R^{(n)}}{n + n_0 + 1} \le P_{x_1, x_2}^{(n+1)}(k) - P_{x_1, x_2}^{(n)}(k)$$
$$\le \frac{A_2^{(n)} - K_2^{(n)} P_{x_1, x_2}^{(n)}(k) + \xi^{(n+1)}}{n + n_0 + 1},$$

where $R^{(n)}$ is a non-random error term satisfying $R^{(n)}/(n+n_0+1) = O(n^{-2})$ as $n \to \infty$.

Proof Since

$$P_{x_1,x_2}^{(n+1)}(k) - P_{x_1,x_2}^{(n)}(k) = \mathbb{E}\left(P_{x_1,x_2}^{(n+1)}(k)\Big|\mathcal{F}_n\right) + \frac{\xi^{(n+1)}}{n+n_0+1} - P_{x_1,x_2}^{(n)}(k),$$

it is sufficient to find bounds for the expected increase in the number of vertices with degree at most k and location inside the interval $[x_1, x_2]$ when v_{n+1} joins the graph with location x_{n+1} , given G_n . This can be expressed by

$$\mathbb{E}\left((n+n_0+1)P_{x_1,x_2}^{(n+1)}(k)\Big|\mathcal{F}_n\right) = (n+n_0)P_{x_1,x_2}^{(n)}(k) + \mathbb{P}(x_{n+1} \in [x_1,x_2]) - \mathbb{P}_n(E_{n+1}),$$

The first term here counts the number of degree at most k vertices in G_n with locations in the interval $[x_1, x_2]$. The second term is the probability that the location of the new vertex v_{n+1} falls into the same interval. Both \mathbb{P}_n and E_{n+1} are as defined as above. We have utilized here the fact that $P_{x_1,x_2}^{(n)}(k)$ is \mathcal{F}_n -measurable, the new location x_{n+1} is independent of \mathcal{F}_n , and that the event E_{n+1} only depends on the graph G_n and the corresponding locations of vertices contained within. As the locations are i.i.d. uniform, this probability is equal to $x_2 - x_1$. For the probability of the event E_{n+1} an upper and lower bound is given by (7) and (8). Hence, we have

$$\mathbb{E}\left((n+n_0+1)P_{x_1,x_2}^{(n+1)}(k)\Big|\mathcal{F}_n\right) \leq \frac{k+\alpha}{2+\alpha} f_2(\Psi_n(x_1),\Psi_n(x_2))\left(P_{x_1,x_2}^{(n)}(k-1)-P_{x_1,x_2}^{(n)}(k)\right) + (n+n_0)P_{x_1,x_2}^{(n)}(k) + (x_2-x_1)$$

as well as

$$\mathbb{E}\left((n+n_0+1)P_{x_1,x_2}^{(n+1)}(k)\Big|\mathcal{F}_n\right) \ge \frac{k+\alpha}{2+\alpha} f_1(\Psi_n(x_1),\Psi_n(x_2))\left(P_{x_1,x_2}^{(n)}(k-1)-P_{x_1,x_2}^{(n)}(k)\right) + (n+n_0)P_{x_1,x_2}^{(n)}(k) + (x_2-x_1)-R^{(n)},$$

where $R^{(n)} = O(n^{-1})$ is an error term, occurring as the difference of the given bound in (7) and the first summand on the right-hand side of the equation together with the fact that $(P_{x_1,x_2}^{(n)}(k-1) - P_{x_1,x_2}^{(n)}(k)) \ge -1$ and the boundedness of f_1 .

Since the number of vertices with degree at most *k* and location inside $[x_1, x_2]$ can change by at most one if we add a new vertex v_{n+1} to the graph G_n , the noise $\xi^{(n)}$ defined in Lemma 3.2 is absolutely bounded by one. Additionally, it holds $\mathbb{E}\left[\xi^{(n+1)}|\mathcal{F}_n\right] = 0$ by its definition. Therefore, we can use stochastic approximation techniques to construct bounds for the asymptotic behaviour of the proportion of vertices with degree at most *k* and location inside $[x_1, x_2]$. **Lemma 3.3** For $\alpha \geq \alpha_c$ and all $k \in \mathbb{N}$, the proportion of vertices with degree at most k and location inside $[x_1, x_2] \subset (0, 1)$ satisfies

$$L_1(k) \le \liminf_{n \to \infty} P_{x_1, x_2}^{(n)}(k) \le \limsup_{n \to \infty} P_{x_1, x_2}^{(n)}(k) \le L_2(k)$$

almost surely, where

$$L_j(k) = (x_2 - x_1) \left(1 - \frac{\Gamma\left(\alpha + 1 + \frac{2+\alpha}{f_j(\Psi(x_1), \Psi(x_2))}\right) \Gamma(\alpha + 1 + k)}{\Gamma(\alpha + 1) \Gamma\left(\alpha + 1 + k + \frac{2+\alpha}{f_j(\Psi(x_1), \Psi(x_2))}\right)} \right), \text{ for } j \in \{1, 2\}.$$

Proof We prove the result by applying Lemma 3.1 to the observed bounds in Lemma 3.2. Here, we focus on the lower bound for the lim inf; the upper bound for the lim sup follows by replacing f_1 by f_2 in the arguments. For each fixed k, the boundedness conditions on $A_1^{(n)}$ and $K_1^{(n)}$, defined in Lemma 3.2, and $P_{x_1,x_2}^{(n)}(k)$ as well as the assumptions (ii) and (iii) are straight forward to check. We have already shown (i) above. We must now show that $A_1^{(n)}$ and $K_1^{(n)}$ converge such that $\lim \inf_{n\to\infty} (A_1^{(n)}/K_1^{(n)}) \ge L_1(k)$, for every k. First note that, since $\alpha \ge \alpha_c$, $\Psi_n(x)$ converges almost surely to $\Psi(x)$ as defined in Sect. 2.2 for all $x \in [0, 1]$. Hence, $K_1^{(n)}$ converges almost surely to $1 + \frac{k+\alpha}{2+\alpha} f_1(\Psi(x_1), \Psi(x_2))$. Now, the theorem can be derived by induction. Let k = 1, then by definition $A_1^{(n)} = x_2 - x_1$. Hence, $\liminf_{n\to\infty} P_{x_1,x_2}^{(n)}(1) \ge L_1(1)$ almost surely. Assume that for an arbitrary fixed $k \in \mathbb{N}$ the stated lower bound holds. Then, for k + 1, we get

$$\liminf_{n \to \infty} A_1^{(n)} \ge \frac{k+1+\alpha}{2+\alpha} f_1(\Psi(x_1), \Psi(x_2)) L_1(k) + (x_2 - x_1)$$

almost surely and hence

$$\begin{split} \liminf_{n \to \infty} P_{x_1, x_2}^{(n)}(k+1) &\geq \frac{\frac{k+1+\alpha}{2+\alpha} f_1(\Psi(x_1), \Psi(x_2)) L_1(k) + (x_2 - x_1)}{1 + \frac{k+1+\alpha}{2+\alpha} f_1(\Psi(x_1), \Psi(x_2))} \\ &= (x_2 - x_1) \left(1 - \frac{\Gamma\left(\alpha + 1 + \frac{2+\alpha}{f_j(\Psi(x_1), \Psi(x_2))}\right) \Gamma(\alpha + 2 + k)}{\Gamma(\alpha + 1) \Gamma\left(\alpha + 2 + k + \frac{2+\alpha}{f_j(\Psi(x_1), \Psi(x_2))}\right)} \right) \\ & \text{most surely.} \end{split}$$

almost surely.

3.2 Limiting Degree Distribution

In this section, we use the established bounds of Lemma 3.3 to prove the main results stated in Sect. 2.1. To this end, we consider now the proportion of vertices located within some interval that have a given maximum degree. We show, that in the late time regime this proportion converges, by shrinking the interval to a single point, to some probability kernel μ on $\mathcal{P}(\mathbb{N}) \times (0, 1)$. Here, $\mathcal{P}(\mathbb{N})$ denotes the set of all subsets of \mathbb{N} . We show that this probability kernel μ is heavy tailed, proving Theorem 2.2 as $\nu(k, x) = \mu(\{k, k+1, ...\}, x)$

Lemma 3.4 If $\alpha \geq \alpha_c$, there exists a probability kernel $\mu : \mathcal{P}(\mathbb{N}) \times (0, 1) \rightarrow [0, 1]$ such that (i) Almost surely,

$$\lim_{x_1 \downarrow x} \lim_{n \to \infty} \frac{P_{x,x_1}^{(n)}(k)}{x_1 - x} = \mu(\{1, \dots, k\}, x).$$

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Proof Note that the functions f_1 and f_2 , defined in (6) and (9), both converge to the same limit as $y_1 \downarrow y$, namely

$$\lim_{y_1 \downarrow y} f_1(y, y_1) = \lim_{y_1 \downarrow y} f_2(y, y_1) = \sum_{s=1}^r s \Xi_s \binom{r}{s} y^{s-1} (1-y)^{r-s},$$

that is f(y), the function used to describe the condensation phase transition in Sect. 2.2. Sending first $n \to \infty$ and applying then the limit $x_1 \to x$ on the bounds observed in Lemma 3.3, we get by continuity of f and Ψ ,

$$\lim_{x_1 \downarrow x} \lim_{n \to \infty} \frac{P_{x,x_1}^{(n)}(k)}{x_1 - x} = 1 - \frac{\Gamma\left(\alpha + 1 + \frac{2+\alpha}{f(\Psi(x))}\right)}{\Gamma(\alpha + 1)} \frac{\Gamma(\alpha + 1 + k)}{\Gamma\left(\alpha + 1 + k + \frac{2+\alpha}{f(\Psi(x))}\right)}.$$
 (10)

Note that

$$\frac{\Gamma(\alpha+1+k)}{\Gamma\left(\alpha+1+k+\frac{2+\alpha}{f(\Psi(x))}\right)} \sim (\alpha+1+k)^{-\frac{2+\alpha}{f(\Psi(x))}} \sim k^{-\frac{2+\alpha}{f(\Psi(x))}}, \text{ as } k \uparrow \infty$$
(11)

by Stirling's formula. For fixed $x \in (0, 1)$, the right-hand side of (10) converges to one as $k \to \infty$, and hence defines a distribution function. Moreover, for fixed k, the right-hand side of (10) is continuous in x. Therefore, the desired probability kernel μ exists, proving (i). The tail behaviour stated in (ii) is an immediate consequence of (11).

Since the empirical distribution of the vertices' locations converges to the uniform distribution on (0, 1), we can now use the probability kernel to properly describe μ_k , the asymptotic proportion of vertices with degree at least *k* by integrating $\mu(\{k, k + 1, ...\}, x)$ with respect to the location. Namely,

$$\mu_k = \int_0^1 \mu(\{k, k+1, \dots\}, x) dx.$$

We are now ready to prove Theorem 2.1.

Proof of Theorem 2.1 Since we only consider the case when k is (very) large, we want to apply a saddle point method approach. To this end, write

$$g(x) = \frac{\Gamma\left(\alpha + 1 + \frac{2+\alpha}{f(x)}\right)}{\Gamma(\alpha + 1)}$$

and consider

$$\mu_{k} = \int_{0}^{1} \mu(\{k, k+1, \dots\}, x) dx = \int_{0}^{1} g(\Psi(x)) \frac{\Gamma(\alpha+1+k)}{\Gamma\left(\alpha+1+k+\frac{2+\alpha}{f(\Psi(x))}\right)} dx$$

given from the proof of Lemma 3.4. Using (11) this reads

$$\mu_k \sim \int_0^1 g(\Psi(x)) \exp\left(-\frac{2+\alpha}{f(\Psi(x))}\log(k)\right) dx$$

as $k \to \infty$. If $f \equiv c$ is constant (e.g. when $\Xi_s = 1/r$), the exponent does not depend on x and the claim follows immediately. Hence, we assume that f is non constant. Now, since

we work in the noncondensation phase, $\Psi(x)$ is the unique zero of $F_1(y; x, \Xi)$, defined in Sect. 2.2. Due to the structure of $F_1(\Psi(x); x, \Xi) = 0$, we can see that the inverse of Ψ exists and that it is a polynomial. Thus, it is differentiable. Together with $\Psi(0) = 0$, almost surely, and $\Psi(1) = 1$ a change of variable leads to

$$\mu_k \sim \int_0^1 (\Psi^{-1})'(y) \cdot g(y) \exp\left(-\frac{2+\alpha}{f(y)}\log(k)\right) dy$$

as $k \to \infty$. For $k \to \infty$, this integral gets dominated by its largest peak that is located at the minimum value of $(2 + \alpha)/f(y)$, occurring at the maximum value of f(y). Since fis a non-negative polynomial, there exists some $x_0 \in [0, 1]$ that maximizes f. In the case that x_0 is not uniquely determined, we can split [0, 1] in finitely many disjoint subintervals such that each subinterval only contains exactly one maximizer. We then integrate these subintervals separately which leads to a sum of integrals all of the same order. Moreover, we know that the second derivative of f exists and that $-((2 + \alpha)/f)''(x_0) > 0$ as well as $(\Psi^{-1})'(x_0) \cdot g(x_0) > 0$. Hence, we get by the saddle point method, for some constant C and with $2 + \alpha_c = f(x_0)$ that

$$\mu_k \sim C_{\sqrt{\frac{2\pi}{\log(k+\alpha)}}} \exp\left(-\frac{2+\alpha}{2+\alpha_c}\log(k+\alpha)\right) \cdot \left(1+O\left(\frac{1}{\log(k+\alpha)}\right)\right),$$

as $k \to \infty$, which yields the desired result.

4 Examples and Simulations

In this section, we discuss a number of examples of the model and use the stated results to calculate the critical value α_c and the power law exponent τ . Simulations of the model back up those results and showcase the different behaviour of the local degree distribution. For this, the different examples are simulated for an initial tree graph of 100 vertices where 1,000,000 new vertices are added to the graph. The code for the simulations can be freely accessed at: http://www.mi.uni-koeln.de/~agrauer/files/code/PA_with_location.R

Throughout this section, we denote by f_{max} the global maximum of f on [0, 1]. The first example is the *middle of three model* introduced in [11]. This model corresponds to the selection vector $\Xi = (0, 1, 0)$, which implies f(y) = 6y(1 - y) due to Eq. (3). This function is maximized at y = 1/2 giving $f_{\text{max}} = 3/2$. As seen in Fig. 2a, y coincides with the maximiser of the local degree distribution $\mu([k, \infty), x)$, for any $k \in \mathbb{N}$. Using the method introduced in Sect. 2.2, the critical value is $\alpha_c = -1/2$, agreeing with the results in [11]. By Theorem 2.1 it can be seen that the degree distribution associated with the middle of three model follows

$$\mu_k = k^{-\frac{2+\alpha}{3/2} + o(1)}$$
, as $k \uparrow \infty$.

Introduced in [11] is the second or sixth of seven model, corresponding to $\Xi = (0, 1/2, 0, 0, 0, 1/2, 0)$. Hence, the associated function is $f(y) = 21y(1-y)((1-y)^4 + y^4)$. This leads to $f_{\text{max}} = \frac{7(5\sqrt{10}-14)}{9}$. Unlike the middle of three model, f has two maximisers which are also peaks of the local degree distribution, see Fig. 2b. The critical value for this example is $\alpha_c = \frac{35\sqrt{10}-116}{9} \approx -0.591$ and it holds

$$u_k = k^{-\frac{9(2+\alpha)}{7(5\sqrt{10}-14)} + o(1)}$$
, as $k \uparrow \infty$.

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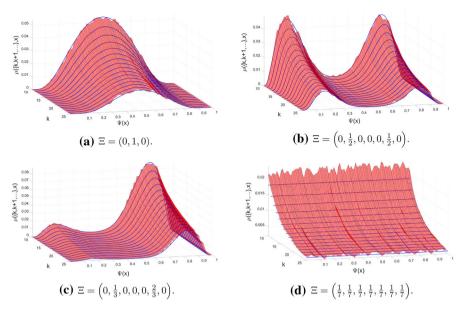


Fig. 2 Simulation of the local degree distribution for the three examples of this section. We have inserted picture (d), which coincides to standard preferential attachment, for comparison. The red surface shows the simulation results while the blue curves depicts the analytical result of Lemma 3.4 for each k. Each plot is generated for $\Psi(x) \in (0, 1)$ and $k \in [10, 25]$ and $\alpha = 0$

The final example is an *asymmetric* version of the second or sixth of seven model, i.e. $\Xi = (0, 1/3, 0, 0, 0, 2/3, 0)$ as selection vector leading to $f(y) = 14y(1 - y)((1 - y)^4 + 2y^4)$. Although this function has two local maximisers, we only care about the global maximum point with $f_{\text{max}} \approx 1.8769$. Figure 2c shows that the mass of the local degree distribution vanishes for large k at the non-global maximiser but concentrates at the global one. The estimation of f_{max} leads to the critical value $\alpha_c \approx -0.1231$ and

$$\mu_k \approx k^{-\frac{2+\alpha}{1.8769}+o(1)}$$
, as $k \uparrow \infty$.

Although the proof of Theorem 2.1 only shows slow convergence to the stated result, our simulations show the stated power law behaviour. For the following figure the simulated degree distribution of the models is fitted to $k^{-\tau}$, considering the logarithmic correction term arising in the proof of Theorem 2.1. For large α , it is necessary to consider simulations of bigger graphs, since the degree is less important for the preferential attachment mechanism, which leads to a small maximum degree of the model. Note that in Fig. 3 the power law exponent of the simulations in each example converges to 1 as $\alpha \rightarrow \alpha_c$.

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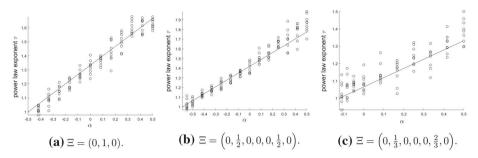


Fig. 3 Simulations of the power law exponent of the degree distribution for each example for α between the corresponding α_c and 1/2. The lines show the analytical result of Theorem 2.1

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