ON EDGE EXCHANGEABLE RANDOM GRAPHS

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ABSTRACT. We study a recent model for edge exchangeable random graphs introduced by Crane and Dempsey; in particular we study asymptotic properties of the random simple graph obtained by merging multiple edges. We study a number of examples, and show that the model can produce dense, sparse and extremely sparse random graphs. One example yields a power-law degree distribution. We give some examples where the random graph is dense and converges a.s. in the sense of graph limit theory, but also an example where a.s. every graph limit is the limit of some subsequence. Another example is sparse and yields convergence to a non-integrable generalized graphon defined on $(0,\infty)$.

1. Introduction

A model for edge exchangeable random graphs and hypergraphs was recently introduced by Crane and Dempsey [11, 12], who also gave a representation theorem showing that every infinite edge exchangeable random hypergraph can be constructed by this model. An equivalent model, using somewhat different formulations, was given by Broderick and Cai [7] and Campbell, Cai and Broderick [8], see Remark 4.7.

The idea of the model is that random i.i.d. edges, with an arbitrary distribution, are added to a fixed vertex set; see Section 4 for a detailed definition (slightly modified but equivalent to the original definition).

The general model defines a random hypergraph. In the present paper, we concentrate on the graph case, although we state the definitions in Section 4 more generally for hypergraphs.

Since edges can be repeated, the model defines a random multigraph, but this can as always be reduced to a random simple graph by identifying parallel edges and deleting loops. Typically, many of the edges will be repeated many times, see e.g. Remark 6.7, and thus the multigraph and the simple graph versions can be expected to be quite different. Both versions have interest and potential, possibly different, applications, and we consider both versions. Previous papers concentrate on the multigraph version; in contrast and as a complement, in the present paper we study mainly the simple graph version.

The model is, as said above, based on an arbitrary distribution of edges. Different choices of this distribution can give a wide range of different types of random graphs, and the main purpose of the paper is to investigate the types of random graphs that may be created by this model; for this purpose we give some general results on the numbers of vertices and edges, and a

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number of examples ranging from dense to very sparse graphs. The examples show that the model can produce very different graphs. In some dense examples we show that the random graphs converge in the sense of graph limit theory. However, that is not always the case, and we even give a chameleon example (Theorem 8.7) that has every graph limit as the limit of some subsequence. We give also a sparse example (Example 9.1) with a power-law degree distribution and convergence to a generalized graphon in the sense of Veitch and Roy [40].

An important tool in our investigations is a Poisson version of the construction by Crane and Dempsey [12], see Section 4.2, which seems interesting also in its own right.

After some preliminaries in Sections 2–3, we give the definitions of the random hypergraphs in detail in Section 4. The graph case is discussed further in Section 5. Section 6 studies the numbers of vertices and edges in the graphs. Section 7 considers an important special case of the model, called *rank 1*; we study two multigraph examples previously considered by Crane and Dempsey [11] and Pittel [36] and show that they are of this type.

The remaining sections consider various examples of the simple graph version, with dense examples in Section 8, and sparse examples in Sections 9 and 10. Finally, we give some tentative conclusions in Section 11.

2. Some notation

In general, we allow hypergraphs to have multiple edges; we sometimes (but usually not) say multihypergraph for emphasis. Moreover, the edges in a hypergraph may have repeated vertices, i.e., the edges are in general multisets of vertices, see Remark 4.3. An edge with repeated vertices is called a loop. A simple hypergraph is a hypergraph without multiple edges and loops. (Warning: different authors give different meanings to "simple hypergraph".)

The vertex and edge sets of a multigraph G are denoted by V(G) and E(G), and the numbers of vertices and edges by v(G) := |V(G)| and e(G) := |E(G)|.

 $f(x) \sim g(x)$ means $f(x)/g(x) \to 1$ (as x tends to some limit, e.g. $x \to \infty$). We also use $v \sim w$ for adjacency of two vertices v and w in a given graph, and $X \sim \mathcal{L}$ meaning that the random variable X has distribution \mathcal{L} ; there should not be any risk of confusion between these (all standard) uses of \sim .

 $f(x) \approx g(x)$ for two non-negative functions or sequences f(x) and g(x) (defined on some common set S) means that f/g and g/f both are bounded; equivalently, there exist constants c, C > 0 such that $cg(x) \leqslant f(x) \leqslant Cg(x)$ for every $x \in S$. $f(x) \approx g(x)$ as $x \to \infty$ means that $f(x) \approx g(x)$ for x in some interval $[x_0, \infty)$.

We use 'increasing' (for a function or a sequence) in its weak sense i.e., $x \leq y \implies f(x) \leq f(y)$, and similarly with 'decreasing'.

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x \wedge y is min\{x, y\} and x \vee y is max\{x, y\}.
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 $\mathbb{N} := \{1, 2, \dots\} \text{ and } \mathbb{N}_0 := \{0, 1, 2, \dots\}. [n] := \{1, \dots, n\}.$

If μ is a measure on a set \mathcal{S} , then $\|\mu\| := \mu(\mathcal{S}) \leqslant \infty$.

 $\operatorname{Exp}(\lambda)$ denotes the exponential distribution with $\operatorname{rate} \lambda$, i.e., the first point in a Poisson process with rate λ ; this is thus the exponential distribution with mean $1/\lambda$. For convenience we extend this to $\lambda=0$: $X\sim\operatorname{Exp}(0)$ means $X=+\infty$ a.s.

We say that a sequence G_n of simple graphs with $v(G_n) \to \infty$ is dense if $e(G_n) \approx v(G_n)^2$, sparse if $e(G_n) = o(v(G_n)^2)$, and extremely sparse if $e(G_n) \approx v(G_n)$ as $n \to \infty$, and similarly for a family G_t of graphs with a continuous parameter.

We let C, c, C_1, c_1, \ldots denote various unspecified positive constants.

3. Some preliminaries on graph limits, graphons and cut metric

We recall some basic facts on graph limits and graphons. For further details, see e.g. [5; 6], [14] and the comprehensive book [32].

A (standard) graphon is a symmetric measurable function $W: \Omega \times \Omega \to [0,1]$, where $\Omega = (\Omega, \mathcal{F}, \mu)$ is a probability space. (Ω may without loss of generality be taken as [0,1] with Lebesgue measure, but it is sometimes convenient to use other probability spaces too.)

If $\varphi : \Omega_1 \to \Omega_2$ is a measure-preserving map between two probability spaces Ω_1 and Ω_2 , and W is a graphon on Ω_2 , then $W^{\varphi}(x,y) := W(\varphi(x), \varphi(y))$ is a graphon on Ω_1 called the *pull-back* of W.

If W is an integrable function on Ω^2 , then its cut norm is

$$||W||_{\square} := \sup \left| \int_{T \times U} W(x, y) \, \mathrm{d}\mu(x) \, \mathrm{d}\mu(y) \right|, \tag{3.1}$$

taking the supremum over all measurable sets $T, U \subseteq \Omega$.

For two graphons W_1 and W_2 , defined on probability spaces Ω_1 and Ω_2 , their *cut distance* is defined as

$$\delta_{\square}(W_1, W_2) = \inf_{\varphi_1, \varphi_2} \|W_1^{\varphi_1} - W_2^{\varphi_2}\|_{\square}, \tag{3.2}$$

taking the infimum over all pairs (φ_1, φ_2) of measure-preserving maps φ_j : $\Omega \to \Omega_j$ defined on some common probability space Ω .

Two graphons W_1 and W_2 are equivalent if $\delta_{\square}(W_1, W_2) = 0$. Note that a graphon W and any pullback W^{φ} of it are equivalent. For characterizations of equivalent graphons, see [4] and [22, Section 8]. The cut distance δ_{\square} can be regarded as a metric on the set W of equivalence classes of graphons, and makes W into a compact metric space.

A graph limit can be identified with an equivalence class of graphons, so we can regard W as the space of graph limits. Thus, every graphon defines a graph limit, and every graph limit is represented by some graphon, but this graphon is unique only up to equivalence.

For every finite graph G, there is a corresponding graphon W_G that can be defined by taking $\Omega = V(G)$ with the uniform probability measure $\mu\{i\} = 1/v(G)$ for every $i \in V(G)$ and letting $W_G(i,j) := \mathbf{1}_{\{i \sim j\}}$; thus W_G equals the adjacency matrix of G, regarded as a function $V(G)^2 \to \{0,1\}$. (W_G is often defined as an equivalent graphon on [0,1]; for us this makes no

difference.) We identify G and W_G when convenient, and write for example $\delta_{\square}(G, W) = \delta_{\square}(W_G, W)$ for a graph G and a graphon W.¹

Remark 3.1. Let G be a finite graph. A blow-up G^* of G is the graph obtained by taking, for some integer $m \ge 1$, the vertex set $V(G^*) = V(G) \times [m]$ with $(v,i) \sim (w,j)$ in G^* if and only if $v \sim w$ in G. Then, W_{G^*} is a pull-back of W_G (for $\varphi: V(G^*) \to V(G)$ the natural projection), and thus $\delta_{\square}(G^*,G) = \delta_{\square}(W_G,W_{G^*}) = 0$. Hence the graphs G and G^* , which are different (if m > 1) are equivalent when regarded as graphons.

There are several, quite different but nevertheless equivalent, ways to define convergence of a sequence of graphs, see e.g. [5; 6; 14; 32]. For our purposes it suffices to know that a sequence G_n with $v(G_n) \to \infty$ is convergent if and only if there exists a graphon W such that $\delta_{\square}(G_n, W) \to 0$ as $n \to \infty$. We then say that G_n converges to W, or to the corresponding graph limit.

Remark 3.2. The standard graphons defined above are appropriate for dense graphs. For sparse graphs, other, more general, graphons have been constructed by several authors. We will in Section 5.1 compare the edge exchangeable graphs studied in the present paper with random graphs defined by graphons that are defined on \mathbb{R}_+ or another infinite (σ -finite) measure space instead of a probability space, see [39; 3]. Furthermore, in Section 9 we consider an example of edge exchangeable graphs that yields sparse graphs, where we show that the graphs converge in a suitable sense (see [40]) to such a graphon defined on \mathbb{R}_+ . We postpone the definitions to these sections.

4. Constructions of random hypergraphs

In this section, we define the random hypergraphs. We give several versions; we define both multihypergraphs and simple hypergraphs, and we give both the original version with a fixed number of edges and a Poisson version. In later sections we consider only the graph case, but we give the definitions here in greater generality.

Note that the edge exchangeable random hypergraphs constructed here are quite different from the vertex exchangeable graphs in e.g. [5; 6; 32; 9; 39; 3], see Section 5.1.

We begin with some preliminaries.

Let (S, \mathcal{F}) be a measurable space, for convenience usually denoted simply by S. To avoid uninteresting technical complications, we assume that S is a Borel space, i.e., isomorphic to a Borel subset of a complete separable metric space with its Borel σ -field.

Let S^* be the set of all finite non-empty multisets of points in S. We can regard a multiset with n elements as an equivalence class of sequences $(x_1, \ldots, x_n) \in S^n$, where two such sequences are equivalent if one is a permutation of the other. Denoting this equivalence relation by \cong and the set of multisets of n elements in S by $S^{\vee n}$, we thus have $S^{\vee n} = S^n/\cong$ and

¹Usually I keep a distiction between graphs and graphons (and graph limits); this is easiest done by identifying a graph G with the pair $(W_G, 1/v(G)) \in \mathcal{W} \times [0, 1]$ (and a graphon W with (W, 0)), see also [14]. In the present paper, this point of view is not needed.

 $\mathcal{S}^* = \bigcup_{n=1}^{\infty} \mathcal{S}^{\vee n}$. Note that $\mathcal{S}^{\vee n}$ and \mathcal{S}^* are Borel spaces. (One way to see this is to recall that every Borel space is isomorphic to a Borel subset of [0,1]. We may thus assume that $\mathcal{S} \subseteq [0,1]$, and then we can redefine $\mathcal{S}^{\vee n}$ as $\{(x_1,\ldots,x_n)\in\mathcal{S}^n:x_1\leqslant\ldots\leqslant x_n\}$, which is a Borel subset of $[0,1]^n$.)

- Remark 4.1. Definitions 4.2 and 4.8 below use a probability measure μ to define the random (hyper)graphs. In general, this measure may be a random measure, and then the constructions should be interpreted by conditioning on μ , i.e., by first sampling μ , and then using the obtained measure throughout the construction. In other words, the distribution of the random hypergraphs constructed by a random measure μ is a mixture of the distributions given by deterministic μ . For convenience, and because most examples will be with deterministic μ , we usually tacitly assume that μ is deterministic; results in the general case with random μ then follow by conditioning on μ . (See Remark 4.11 for a typical example, where this for once is stated explicitly.)
- 4.1. Random hypergraphs with a given number of edges. We give a minor modification of the original definition by Crane and Dempsey [11, 12]; we will see at the end of this subsection that our definition is equivalent to the original one.

Definition 4.2. Given a Borel space S and a probability measure μ on S^* , define a sequence of finite random (multi)hypergraphs $(G_m^*)_{m=1}^{\infty}$ as follows.

(i) Draw $Y_1, Y_2, \ldots \sim_{\text{iid}} \mu$.

(ii) Let
$$V(G_m^*) := \bigcup_{i \leq m} Y_i$$
 and $E(G_m^*) := \{Y_1, \dots, Y_m\}$ (multiset). (4.1)

Note that $V(G_m^*)$ is the vertex set spanned by the edges; thus there are no isolated vertices in G_m^* . (The same holds for the related definitions in (4.2), (4.6), (4.7) below.)

We also similarly define the infinite (multi)hypergraph G_{∞}^* having edges $(Y_i)_{i=1}^{\infty}$.

The edges in G_m^* may be repeated, so G_m^* is in general a random multi-hypergraph. We define G_m as the simple hypergraph obtained by merging each set of parallel edges in G_m^* to a single edge and deleting loops; thus the simple hypergraphs $(G_m)_1^{\infty}$ are defined by:

(i) Draw $Y_1, Y_2, \ldots \sim_{\text{iid}} \mu$.

(ii) Let
$$E(G_m) = \{Y_i : i \leq m, Y_i \text{ not loop}\}$$
 and
$$V(G_m) = \bigcup_{Y_i \in E(G_m)} Y_i.$$
 (4.2)

Thus $V(G_m) \subseteq V(G_m^*)$, and strict inequality is possible if there are loops.

Note that $G_1^* \subset G_2^* \subset \ldots$, and thus $G_1 \subseteq G_2 \subseteq \ldots$, i.e., $(G_m^*)_m$ and $(G_m)_m$ are increasing sequences of random hypergraphs.

Remark 4.3. We follow [12] and allow for increased generality Y_i to be a multiset (see e.g. the examples in Section 7); thus the edges in G_m^* and G_m are multisets and may contain repeated vertices. If we choose μ with support in the set $\mathcal{S}^{**} := \bigcup_{n=1}^{\infty} \mathcal{S}^{\wedge n} \subset \mathcal{S}^*$ of finite subsets of \mathcal{S} , where $\mathcal{S}^{\wedge n} \subset \mathcal{S}^{\vee n}$ is

the set of subsets of S with n distinct elements, then the edges in G_m^* and G_m are ordinary sets of vertices (i.e., without repeated vertices). (This is commonly assumed in the definition of hypergraphs.)

In particular, if μ has support in $\mathcal{S}^{\wedge 2} = \{\{x,y\} : x,y \in \mathcal{S}, x \neq y\}$, then G_m^* is a multigraph without loops, and G_m is a simple graph with $V(G_m) = V(G_m^*)$.

The construction above yields hypergraphs with vertices labelled by elements of S. We (usually) ignore these labels and regard G_m^* and G_m as unlabelled hypergraphs.

Remark 4.4. We usually also ignore the labels on the edges. If we keep the labels i on the edges Y_i , then the distribution of G_m^* is obviously edge exchangeable, i.e., invariant under permutations of these edge labels, because $(Y_i)_i$ is an i.i.d. sequence. Conversely, as shown by Crane and Dempsey [12, Theorem 3.4], every infinite edge exchangeable hypergraph is a mixture of random hypergraphs G_∞^* , i.e., it can be constructed as above using a random measure μ . In the present formulation, the proof in [12] simplifies somewhat: Give the vertices in the edge exchangeable hypergraph random labels that are i.i.d. and U(0,1) (uniformly distributed on [0,1]), and independent of the edges. Then the edges become multisets in $[0,1]^*$, and their distribution is clearly exchangeable, so by de Finetti's theorem, the edges are given by the construction above for some random probability measure μ on \mathcal{S}^* , taking $\mathcal{S} = [0,1]$.

It is obvious from the definition that if $\psi: \mathcal{S} \to \mathcal{S}_1$ is an injective measurable map of \mathcal{S} into another measurable (Borel) space \mathcal{S}_1 , then μ is mapped to a probability measure μ_1 on \mathcal{S}_1^* , which defines the same random hypergraphs G_m^* and G_m as μ . Hence, the choice of Borel space \mathcal{S} is not important, and we can always use e.g. $\mathcal{S} = [0, 1]$. Moreover, we can simplify further.

Define the *intensity* of μ as the measure on (S, \mathcal{F})

$$\bar{\mu}(A) := \mathbb{E}|A \cap Y|, \qquad A \in \mathcal{F},$$
 (4.3)

where Y has distribution μ . Note that for a singleton set $\{x\}$, $|\{x\} \cap Y| = \mathbf{1}_{\{x \in Y\}}$, and thus (4.3) yields

$$\bar{\mu}(\{x\}) = \mathbb{P}(x \in Y). \tag{4.4}$$

We have $\bar{\mu}(A) = \sum_{n=1}^{\infty} \bar{\mu}_n(A)$, where $\bar{\mu}_n(A) := \mathbb{E}(|A \cap Y| \cdot \mathbf{1}_{\{|Y|=n\}})$, and since each $\bar{\mu}_n$ is a finite measure, it follows that the set of atoms

$$\mathcal{A} := \{ x \in \mathcal{S} : \bar{\mu}(\{x\}) > 0 \} \tag{4.5}$$

is a countable (finite or infinite) subset of S. By (4.4) and (4.5), if $x \notin A$, then $\mathbb{P}(x \in Y) = 0$. Hence, in the construction of G_m^* , if an edge Y_i has a vertex $x \notin A$, then a.s. $x \notin Y_j$ for every $j \neq i$. Consequently, a vertex $x \notin A$ of G_{∞}^* a.s. appears in only one edge. (Such a vertex is called a *blip* in [12].) On the other hand, if $x \in A$, so $\mathbb{P}(x \in Y) = \bar{\mu}(\{x\}) > 0$, then by the law of large numbers, a.s. x belongs to infinitely many edges Y_i of G_{∞}^* .

It follows that when constructing the hypergraphs G_m^* , if the edge $Y_i = \{y_{i1}, \ldots, y_{in_i}\}$, we do not have to keep track of the vertex labels y_{ij} unless they belong to \mathcal{A} ; any $y_{ij} \notin \mathcal{A}$ will be a blip not contained in any other edge and the actual value of y_{ij} may be forgotten. (Except that if we allow

repeated vertices in the edges, see Remark 4.3, then we still have to know whether two vertex labels y_{ij} and y_{ik} on the same edge are the same or not.)

Now, enumerate \mathcal{A} as $\{a_i\}_{i=1}^N$, where $N \leq \infty$, and replace, for every multiset $Y = (y_1, \ldots, y_\ell) \in \mathcal{S}^*$, every vertex label $y_j = a_k$ for some $a_k \in \mathcal{A}$ by the new label $y_j' = k$, and the vertex labels $y_j \notin \mathcal{A}$ on Y by $0, -1, \ldots$ (For definiteness, we may assume that $\mathcal{S} \subseteq [0,1]$ so \mathcal{S} is ordered, and take the labels in order in case Y has more than one vertex label not in \mathcal{A} .) This maps μ to a probability measure μ' on the set \mathbb{Z}^* of finite multisets of integers, and it follows from the discussion above that we can recover the random hypergraphs G_m^* from μ' by the construction in Definition 4.2, if we first replace each vertex label $y_j' \in \{0, -1, \ldots\}$ by a random label with a continuous distribution in some set, for example U(0,1), making independent choices for each Y_i . Equivalently, and more directly, we obtain G_m^* from the probability measure μ' on \mathbb{Z}^* by the following construction, which is the original definition by Crane and Dempsey [11, 12].

Definition 4.5 (Crane and Dempsey [11, 12]). Given a probability measure μ on \mathbb{Z}^* , we define a sequence of finite random (multi)hypergraphs $(G_m^*)_{m=1}^{\infty}$ as in Definition 4.2 with the modification that in every edge $Y_i = \{y_{i1}, \ldots, y_{i\ell_i}\}$ we replace every vertex label $y_{ij} \leq 0$ (if any) with a new vertex that is not used for any other edge.

Since we ignore the vertex labels in G_m^* , it does not matter what labels we use as replacements for $0, -1, \ldots$ in Definition 4.5. Crane and Dempsey [11, 12] use the same set $0, -1, \ldots$ of integers, taking the first label not already used. An alternative is to take random labels, e.g. i.i.d. U(0,1) as above.

Remark 4.6. To be precise, Definition 4.5 is the definition in [12]. The definition in Crane and Dempsey [11] treats only the binary case $|Y_n| = 2$ in detail; and differs in that only labels $y_i \ge 0$ are used, and that an edge $\{0,0\}$ is replaced by an edge $\{z_1, z_2\}$ with two new vertex labels z_1 and z_2 .

This version is essentially equivalent; apart from a minor notational difference, the only difference is that this version does not allow for "loop dust", where a positive fraction of the edges are isolated loops. Cf. Remark 5.2.

We have shown that Definition 4.2 is essentially equivalent to the original definitions by Crane and Dempsey [11, 12]. One advantage of Definition 4.2 is that no special treatment of vertex labels ≤ 0 is needed; the blips (if there are any) come automatically from the continuous part of the label distribution; a disadvantage is that this continuous part is arbitrary and thus does not contain any information. Another advantage with Definition 4.2 is that it allows for arbitrary Borel spaces \mathcal{S} ; even if it usually is convenient to use $\mathcal{S} = \mathbb{N}$ to label the vertices, it may in some examples be natural to use another set \mathcal{S} .

Remark 4.7. The construction in Campbell, Cai and Broderick [8] is stated differently, but is equivalent. It uses a generalization of Kingman's paintbox construction of exchangeable partitions; in the version in Campbell, Cai and Broderick [8], the paintbox consists of families $(C_{kj})_{k,j\geq 1}$ and $(C'_{jl})_{j,l\geq 1}$ of subsets of [0, 1]; it is assumed that every $x \in [0, 1]$ is an element of only

finitely many of these sets, and that for each j and $k \neq l$, $C_{jk} \cap C_{jl} = \emptyset$ and $C'_{jk} \cap C'_{jl} = \emptyset$. (In general these sets may be random, but similarly as above, in the construction we condition on these sets so we may assume that they are deterministic.) Furthermore, we generate i.i.d. U(0,1) random labels ϕ_k and ϕ_{Njl} for $k, N, j, l \geqslant 1$. For each $N \geqslant 1$ we construct a edge Y_N by taking a uniformly random point $V_N \in [0,1]$, independent of everything else; then, for each (j,k) such that $V_N \in C_{jk}$, Y_N contains k vertices labelled ϕ_j , and for each (j,k) such that $V_N \in C'_{jk}$ and every $l \leqslant k$, Y_N contains j vertices labelled ϕ_{Njl} . (The latter vertices are thus blips.)

Note that this gives the vertices random labels as in Remark 4.4; however, we then ignore the vertex labels. (Actually, in [8], each vertex is represented by a multiset of edge labels (called a *trait*), which contains the label of each edge that contains the vertex, repeated as many times as the vertex occurs in the edge. This is obviously an equivalent way to describe the hypergraph.)

It is obvious that, conditioned on the labels ϕ_k and ϕ_{Njl} , this construction gives a random multiset with some distribution μ ; conversely, every distribution μ of a random (finite) multiset can easily be obtained in this way by suitable choices of C_{jk} and C'_{jk} . Hence, the construction is equivalent to the one above. (In our opinion, it is more natural to focus on the distribution of the edges, since the sets C_{jk} and C'_{jk} in the paintbox construction have no intrinsic meaning; they are just used to describe the edge distribution.)

4.2. The Poisson version. The multihypergraph G_m^* has exactly m edges (not necessarily distinct). It is often convenient to instead consider a Poisson number. (This was done by Broderick and Cai in [7, Example 2.7].) It is then natural to consider a continuous-parameter family of hypergraphs, which we define as follows. We may think of the second coordinate t as time.

Definition 4.8. Given a probability measure μ on \mathcal{S}^* , we define a family of random (multi)hypergraphs $(\tilde{G}_t^*)_{t\geq 0}$ as follows. Recall that a Poisson point process on an infinite, σ -finite measure space is a random countably infinite set of points that can be enumerated as in (i), in our case, for some random $Y_i \in \mathcal{S}^*$ and $\tau_i \in [0, \infty)$.

(i) Let $\Xi = \{(Y_i, \tau_i) : i \ge 1\}$ be a Poisson point process on $\mathcal{S}^* \times [0, \infty)$ with intensity $\mu \times dt$;

(ii) Let
$$E(\tilde{G}_t^*) := \{Y_i : \tau_i \leqslant t\}$$
 (multiset), and $V(\tilde{G}_t^*) := \bigcup_{Y \in E(\tilde{G}_t^*)} Y$.
$$(4.6)$$

Define \tilde{G}_t as the simple hypergraph obtained by merging each set of parallel edges in \tilde{G}_t^* to a single edge, and deleting loops (together with their incident vertices, unless these also belong to some non-loop). Hence, with (i) as in (4.6),

(ii) Let
$$E(\tilde{G}_t) := \{Y_i \text{ not loop} : \tau_i \leqslant t\}$$
, and $V(\tilde{G}_t) := \bigcup_{Y \in E(\tilde{G}_t)} Y$.

(4.7)

Note that the random hypergraphs \tilde{G}_t^* and \tilde{G}_t are a.s. finite for every $t < \infty$.

The projection $\Xi'' := \{\tau_i\}_{i=1}^{\infty}$ of the Poisson process Ξ to the second coordinate is a Poisson point process on $[0, \infty)$ with intensity 1, and we may and will assume that the points of Ξ are enumerated with τ_i in increasing order; thus a.s. $0 < \tau_1 < \tau_2 < \ldots$ Let N(t) be the number of points of Ξ in $\mathcal{S}^* \times [0, t]$, i.e.

$$N(t) := \left| \Xi \cap (\mathcal{S}^* \times [0, t]) \right| = \max\{i : \tau_i \leqslant t\}; \tag{4.8}$$

this is a Poisson counting process on $[0, \infty)$ and $N(t) \sim \text{Po}(t)$. Conversely, τ_m is the time the process N(t) reaches m, so the increments $\tau_m - \tau_{m-1}$ (with $\tau_0 := 0$) are i.i.d. and Exp(1), and τ_m has the Gamma distribution $\Gamma(m)$. Moreover, the random multisets Y_i are i.i.d. with distribution μ and independent of $\{\tau_i\}$, so they can be taken as the Y_i in Definition 4.2, which leads to the following simple relation between the two definitions.

Proposition 4.9. If μ is a probability measure on S^* , then we may couple the random hypergraphs constructed in Definitions 4.2 and 4.8 such that $G_m^* = \tilde{G}_{\tau_m}^*$ and thus $G_m = \tilde{G}_{\tau_m}$ for all $m \ge 1$, and conversely $\tilde{G}_t^* = G_{N(t)}^*$ and $\tilde{G}_t = G_{N(t)}$ for all $t \ge 0$.

Although we usually tacitly consider $t < \infty$, we may here also take $t = \infty$: $G_{\infty}^* = \tilde{G}_{\infty}^*$ and $G_{\infty} = \tilde{G}_{\infty}$.

Note that the relations in Proposition 4.9 hold not just for a single m or t, but also for the entire processes. Hence, asymptotic results, and in particular a.s. limit results, are (typically) easily transferred from one setting to the other.

Remark 4.10. Instead of stopping at the random time τ_m , we can also obtain G_m^* and G_m from \tilde{G}_t^* and \tilde{G}_t by conditioning on N(t) = m, for any fixed t > 0.

Remark 4.11. One reason that the Poisson version is convenient is that different edges appear independently of each other. If we for convenience assume that there are no blips, we may as explained above assume that $\mathcal{S} = \mathbb{N}$, so $V(\tilde{G}_t^*) \subseteq \mathbb{N}$. In this case, the number of copies of an edge $I \in \mathcal{S}^*$ in \tilde{G}_t^* has the Poisson distribution $\operatorname{Po}(t\mu(\{I\}))$, and these numbers are independent for different $I \in \mathcal{S}^*$. Hence, different edges $I \in \mathcal{S}^*$ appear independently in \tilde{G}_t . (In the case μ is random, this holds conditionally on μ , but not unconditionally.)

Note that this independence does not hold for G_m ; the stopping in Proposition 4.9 or the conditioning in Remark 4.10 destroys the independence of different edges.

4.3. Unnormalized measures. We have so far assumed that μ is a probability measure. This is very natural, but we can make a trivial extension to arbitrary finite measures. This will not produce any new random hypergraphs but it is convenient; for example, it means that we do not have to normalize the measure in the examples in later sections.

When necessary, we denote the measure used in the construction of our random hypergraphs by a subscript; we may thus write e.g. $G_{m,\mu}$.

Definition 4.12. Let μ be a finite measure on \mathcal{S}^* , not identically zero. Let μ_0 be the probability measure $\mu_0 := \|\mu\|^{-1}\mu$, and define $G^*_{m,\mu} := G^*_{m,\mu_0}$.

Furthermore, define $\tilde{G}_{t,\mu}^*$ as in Definition 4.8. Let, as usual, $G_{m,\mu}$ and $\tilde{G}_{t,\mu}$ be the corresponding simple graphs.

Thus, $\mu = c\mu_0$, where $c := \|\mu\| = \mu(\mathcal{S}^*)$. It is obvious that, using obvious notation, the Poisson process Ξ_{μ} can be obtained from Ξ_{μ_0} by rescaling the time: If $\Xi_{\mu_0} = \{(Y_i, \tau_i^0)\}$, we can take $\Xi_{\mu} = \{(Y_i, c^{-1}\tau_i^0)\}$, and thus $G_{t,\mu}^* = G_{ct,\mu_0}^*$. Hence, the random hypergraph process defined by μ is the same as for μ_0 , except for a simple deterministic change of time. This implies the following result.

Proposition 4.13. Proposition 4.9 extends to arbitrary finite measures μ (not identically zero), with stopping times τ_m that are the partial sums $\sum_{i=1}^m T_i$ of i.i.d. random variables $T_i \sim \text{Exp}(\|\mu\|)$.

In particular, the law of large numbers yields, as $m \to \infty$,

$$\tau_m/m \xrightarrow{\text{a.s.}} \|\mu\|^{-1}.$$
 (4.9)

Remark 4.14. Definition 4.8 can be employed also when μ is an infinite and, say, σ -finite measure. In this case, \tilde{G}_t^* has a.s. an infinite number of edges for every t > 0. We will not consider this case further.

5. Random Graphs

From now on, we consider the graph case, where μ is a finite measure on $\mathcal{S}^{\vee 2} = \{\{x,y\} : x,y \in \mathcal{S}\}$. This allows for the presence of loops; often we consider μ supported on $\mathcal{S}^{\wedge 2} = \{\{x,y\} : x \neq y\}$, and then there are no loops.

As explained in Section 4, in particular Definition 4.5, if there are no blips (i.e., if the intensity $\bar{\mu}$ is discrete), we may without loss of generality assume that $S = \mathbb{N}$, and if there are blips, we may assume that $S = \mathbb{N} \cup \{0, -1\}$ with the special convention that 0 and -1 are interpreted as blips. Unless stated otherwise, we use this version, and we then write μ_{ij} for $\mu(\{i,j\})$; we say that μ_{ij} is the *intensity* of edges ij. Thus, (μ_{ij}) is an infinite symmetric matrix of non-negative numbers, with indices in $\mathbb{N} \cup \{0, -1\}$ (or in \mathbb{N} if there are no blips); note that, because we consider undirected edges, the total mass of μ is

$$\|\mu\| = \frac{1}{2} \sum_{i,j: i \neq j} \mu_{ij} + \sum_{i} \mu_{ii}.$$
 (5.1)

We assume that $0 < \|\mu\| < \infty$, or equivalently that $\sum_{i,j} \mu_{ij}$ is finite (and non-zero), but we do not insist on μ being a probability measure. As described in Section 4.3, we can always normalize μ to the probability measure $\|\mu\|^{-1}\mu$ when desired.

We also define (for $i \ge 1$)

$$\mu_i := \sum_j \mu_{ij},\tag{5.2}$$

this is the total intensity of edges adjacent to vertex i.

Remark 5.1. The diagonal terms μ_{ii} correspond to loops. Loops appear naturally in some examples, see e.g. Example 7.1 below, but we are often interested in examples without loops, and then take $\mu_{ii} = 0$. Moreover, in

the construction of the simple graphs G_m and \tilde{G}_t we delete loops, so it is convenient to take $\mu_{ii}=0$ and avoid loops completely. Note that, since different edges appear independently in \tilde{G}_t^* , see Remark 4.11, deleting all loops from \tilde{G}_t^* is equivalent to conditioning \tilde{G}_t^* on containing no loops; this is also equivalent to changing every μ_{ii} to 0. (For G_m this is not quite true, since the number of non-loop edges may change; however, the difference is minor.)

Note also the in the construction leading to Definition 4.5, in the graph case, vertex label -1 is used only for the edge $\{0, -1\}$, so we may (and will) assume that $\mu_{i,-1} = 0$ unless i = 0.

Suppose now that we are given such a matrix $(\mu_{ij})_{i,j\geqslant -1}$. We can decompose the matrix into the three parts $(\mu_{ij})_{i,j\geqslant 1}$, $(\mu_{i0})_{i\geqslant 1}$, $(\mu_{i0})_{i\in\{0,-1\}}$, which by the construction and properties of Poisson processes correspond to a decomposition of the Poissonian multigraph \tilde{G}_t^* as a union of three parts, which are independent random graphs:

Central part: The edges $ij \in \tilde{G}_t^*$ with $i, j \in \mathbb{N}$.

Attached stars: For each $i \ge 1$ a star with $Po(t\mu_{i0})$ edges centred at i.

Dust: Po $(t\mu_{00})$ isolated loops and Po $(t\mu_{0,-1})$ isolated edges.

Moreover, the Poisson random variables above, for different i and for the two types of dust, are independent. The vertex set is by definition the set of endpoints of the edges, so there are no isolated vertices. The edges and loops in the dust are always isolated, i.e. with endpoints that are blips (have no other edges). Similarly, the peripheral vertices in the attached stars are blips without other edges, while the central vertex i may, or may not, also belong to the central part.

Note that multiple edges only occur in the central part.

Remark 5.2. We have here discussed the model in full generality, but it is obvious that the main interest is in the central part, and all our examples will be with μ supported on $\mathbb{N} \times \mathbb{N}$, i.e., without dust and attached stars. (Of course, there may be other stars or isolated edges, created in the central part.)

In particular, the dust part is quite trivial, and the dust loops are even less interesting than the dust edges. In a case with dust but no loops in the dust, it is convenient to relabel $\mu_{0,-1}$ as μ_{00} , so μ is a symmetric matrix with index set \mathbb{N}_0 ; this corresponds to using the version of the definition in [11], see Remark 4.6.

5.1. A comparison with vertex exchangeable graphs. Consider the case without dust, attached stars and loops, so μ is supported on $\mathbb{N} \times \mathbb{N}$, with $\mu_{ii} = 0$. Then \tilde{G}_t^* has $\text{Po}(t\mu_{ij})$ edges ij, for every pair of distinct integers $i, j \in \mathbb{N}$.

We may compare this to the vertex exchangeable random graphs studied by e.g. Borgs, Chayes, Lovász, Sós and Vesztergombi [5, 6], Diaconis and Janson [14], Lovász [32] and their generalizations by Caron and Fox [9], Veitch and Roy [39] and Borgs, Chayes, Cohn and Holden [3], see also Orbanz and Roy [34], Herlau, Schmidt and Mørup [17], and Janson [23].

In the classical case [5; 6; 32], with a standard graphon W defined on a probability space (Ω, ν) as in Section 3, the vertex exchangeable random

graph G(n, W) has a given number n of vertices and is constructed as follows.

- (i) Draw $x_1, x_2, \dots x_n \sim_{\text{iid}} \nu$. (Think of x_i as the "type" of i.)
- (ii) Let $V(G(n, W)) := \{1, \dots, n\}$. For each pair (i, j), i < j, (5.3) let $ij \in E(G(n, W))$ with probability $W(x_i, x_j)$.

(Different edges are added independently, given the types x_i .)

The generalization to graphons on \mathbb{R}_+ or another σ -finite measure space (Ω, ν) [9; 39; 3] is similar. This type of graphon is still a symmetric measurable function $W: \Omega^2 \to [0,1]$ (satisfying some conditions to make the graphs G(t,W) defined below a.s. finite). Ω can be regarded as a space of types, and the random graph, here denoted $\bar{G}(t,W)$, is defined as follows. The number of vertices $N \sim \text{Po}(t\nu(\Omega))$ is a random variable, with $N = \infty$ if ν is an infinite measure.

- (i) Let $\Xi = \{x_i\}_{i=1}^N$ be a Poisson point process on Ω with intensity $t\nu$.
- (ii) Let $V(\bar{G}(t,W)) := \{i \in \mathbb{N} : i \leq N\}$. For each pair (i,j), i < j, (5.4) let $ij \in E(\bar{G}(t,W))$ with probability $W(x_i,x_j)$.

Finally, we may delete all isolated vertices, giving a graph G(t, W) without isolated vertices (as in the construction in Section 4 above):

(iii) Let
$$V(G(t, W)) := \{i : ij \in E(\bar{G}(t, W)) \text{ for some } j\},\ E(G(t, W)) := E(\bar{G}(t, W)).$$
 (5.5)

In both cases (5.3) and (5.4)–(5.5), a natural multigraph version is to modify (ii) by instead taking a Poisson number $Po(W(x_i, x_j))$ of copies of the edge ij. (Cf. e.g. [2, Remark 2.4]. One might also take $Po(-\log(1 - W(x_i, x_j)))$ copies, keeping the probability of no edge the same as for the Bernoulli model. Note that if the $W(x_i, x_j)$ are small, then the standard (Bernoulli) and the two Poisson versions are almost the same.)

The Poisson versions of the edge exchangeable and vertex exchangeable random graphs thus add edges in the same way, if we condition on the types of the vertices in the latter and let $\mu_{ij} = t^{-1}W(x_i, x_j)$. However, the vertices are constructed in very different ways. To see the similarities and differences clearly, consider the case where the type space $\Omega = \mathbb{N}$, with some (finite or infinite) measure ν , and consider the Poisson multigraph version of the vertex exchangeable graphs, which we denote by $\bar{G}^*(t, W)$ and $G^*(t, W)$. Then the vertex exchangeable $G^*(t, W)$ has a Poisson number $Po(t\nu\{i\})$ of vertices of type i, for each $i \in \mathbb{N}$, while the edge exchangeable \tilde{G}_t^* has at most one vertex i for each $i \in \mathbb{N}$. (We can reformulate the construction of G_t^* and say that we start with exactly one vertex of type i for every $i \in \mathbb{N}$, and then remove all isolated vertices after having added edges.) Moreover, although for a fixed t, each pair of distinct vertices of types i and j has Po(W(i,j))edges between them in $\bar{G}^*(t, W)$ or $G^*(t, W)$ and $Po(t\mu_{ij})$ edges in \tilde{G}_t^* , which coincide if $W(i,j) = t\mu_{ij}$, we see that if we keep W and μ fixed and increase t, the two families $\bar{G}^*(t,W)$ and \tilde{G}_t^* behave differently: In \tilde{G}_t^* the number of edges between each pair of vertices increases linearly as t increases, the number of vertices increases more slowly (by Corollary 6.6 below; recall that we only keep vertices with at least one edge), and there is at most one vertex of each type. In $\bar{G}^*(t, W)$ and $G^*(t, W)$, the number of vertices of each type increases linearly, while the number of edges between each pair of vertices remains the same.

6. Numbers of vertices and edges

By construction, the number of edges is m in the multigraph G_m^* and random $Po(t||\mu||)$ in the multigraph \tilde{G}_t^* . The numbers of vertices in the graphs and the numbers of edges in the simple graphs G_m and \tilde{G}_t are somewhat less immediate, and are studied in this section.

We use the notation of Section 5, and assume that we are given a (deterministic) matrix $\mu = (\mu_{ij})$ of intensities. Moreover, for simplicity we assume that μ is concentrated on $\mathbb{N} \times \mathbb{N}$, so there are no attached stars and no dust, and that $\mu_{ii} = 0$ for every i, so there are no loops. We consider briefly the case with dust or attached stars in Section 6.1.

Note that G_m is a simple graph without isolated vertices, and thus

$$\frac{1}{2}v(G_m) \leqslant e(G_m) \leqslant \binom{v(G_m)}{2} \leqslant \frac{1}{2}v(G_m)^2. \tag{6.1}$$

Recall that G_m is dense if $e(G_m) \simeq v(G_m)^2$, sparse if $e(G_m) = o(v(G_m)^2)$, and extremely sparse if $e(G_m) \simeq v(G_m)$ as $m \to \infty$. By Propositions 4.9 and 4.13, these are equivalent to the corresponding conditions for \tilde{G}_t .

The number of edges in G_m is the number of different values taken by the i.i.d. sequence Y_1, \ldots, Y_m . Equivalently, it is the number of occupied bins if m balls are thrown independently into an infinite number of boxes, with the probability μ_{ij} (normalized if necessary) for box $\{i, j\}$. Such numbers have been studied in, for example, [13; 30; 31; 15; 19], where central limit theorems have been proved under various assumptions, see Theorem 6.8 below. These results are often proved using Poissonization, which in our setting is equivalent to considering \tilde{G}_t instead of G_m . We too find it convenient to first study the Poisson version.

The Poisson model is convenient because, as said before, edges ij arrive according to a Poisson process with intensity μ_{ij} and these Poisson processes are independent for different pairs $\{i, j\}$. Let $N_{ij}(t)$ be the number of copies of the edge ij in \tilde{G}_t^* , and let $N_i(t)$ be the degree of vertex i in \tilde{G}_t^* . Then

$$N_{ij}(t) \sim \text{Po}(t\mu_{ij}),$$
 (6.2)

and, recalling (5.2),

$$N_i(t) = \sum_{i \neq i} N_{ij}(t) \sim \text{Po}(t\mu_i). \tag{6.3}$$

Moreover, let $T_i \sim \text{Exp}(\mu_i)$ and $T_{ij} \sim \text{Exp}(\mu_{ij})$ be the random times that the first edge at i and the first edge ij appear, respectively. Thus, $N_i(t) \ge 1 \iff T_i \le t$ and $N_{ij}(t) \ge 1 \iff T_{ij} \le t$.

By the construction of \tilde{G}_t ,

$$v(\tilde{G}_t) = v(\tilde{G}_t^*) = \sum_{i} \mathbf{1}_{\{N_i(t) \ge 1\}} = \sum_{i} \mathbf{1}_{\{T_i \le t\}},$$
(6.4)

$$e(\tilde{G}_t) = \sum_{i < j} \mathbf{1}_{\{N_{ij}(t) \ge 1\}} = \sum_{i < j} \mathbf{1}_{\{T_{ij} \le t\}}.$$
 (6.5)

Recall that for every fixed t, the numbers $N_{ij}(t)$ are independent random variables, and thus the indicators in the sums (6.5) are independent. However, the numbers $N_i(t)$ and the indicators in the sums in (6.4) are dependent, which is a complication. (For example, $v(\tilde{G}_t) = 1$ is impossible, since there are no isolated vertices and no loops.)

We give first a simple lemma for the type of sums in (6.5), where the terms are independent.

Lemma 6.1. Let $Z_i \sim \operatorname{Exp}(\lambda_i)$, $i = 1, 2, \ldots$, be independent exponential random variables with $\lambda_i \geqslant 0$ and $0 < \sum_{i=1}^{\infty} \lambda_i < \infty$, and let $W(t) := \sum_{i=1}^{\infty} \mathbf{1}_{\{Z_i \leqslant t\}}$. Then, the following hold.

(i) For every $t \ge 0$,

$$\mathbb{E}W(t) = \sum_{i=1}^{\infty} \mathbb{P}(Z_i \leqslant t) = \sum_{i=1}^{\infty} (1 - e^{-\lambda_i t}) < \infty$$
 (6.6)

and thus a.s. $W(t) < \infty$ for every $t \ge 0$. Furthermore, $\mathbb{E}W(t)$ is a strictly increasing and concave continuous function of $t \ge 0$ with $\mathbb{E}W(0) = 0$ and $\mathbb{E}W(t)/t \to 0$ as $t \to \infty$.

(ii) For t > 0,

$$\mathbb{E} W(t) \approx \sum_{i=1}^{\infty} (1 \wedge (\lambda_i t)). \tag{6.7}$$

(iii) For every $t \geqslant 0$,

$$\operatorname{Var}(W(t)) = \sum_{i=1}^{\infty} e^{-\lambda_i t} (1 - e^{-\lambda_i t}) \leqslant \mathbb{E} W(t). \tag{6.8}$$

(iv) Let $L := |\{i : \lambda_i > 0\}| \leq \infty$. Then as $t \to \infty$, $\mathbb{E}W(t) \to L$, $W(t) \xrightarrow{\text{a.s.}} L$ and

$$\frac{W(t)}{\mathbb{E}W(t)} \xrightarrow{\text{a.s.}} 1. \tag{6.9}$$

(v) If (t_n) and (t'_n) are two sequences of positive numbers with $t'_n/t_n \to 1$, then $\mathbb{E} W(t'_n)/\mathbb{E} W(t_n) \to 1$.

Proof. This is presumably all known, but it seems easier to give a proof than to find references. Note that W(t) is increasing as a function of t.

(i): The calculation (6.6) of the expectation is immediate, and the sum is finite because $1 - e^{-\lambda_i t} \leq \lambda_i t$. Hence W(t) is a.s. finite for, say, each integer t, and thus for all $t \geq 0$.

It follows by (6.6) that $\mathbb{E}W(t)$ is strictly increasing and concave. Moreover, the sum converges uniformly on every finite interval [0,T], and thus $\mathbb{E}W(t)$ is continuous. Finally, $\mathbb{E}W(t)/t = \sum_{i=1}^{\infty} (1 - e^{-\lambda_i t})/t$, where each summand tends to 0 as $t \to \infty$, and is bounded by λ_i . Hence $\mathbb{E}W(t)/t \to 0$ as $t \to \infty$ by dominated convergence of the sum.

(ii): An immediate consequence of (6.6) and $1 - e^{-x} \approx 1 \wedge x$.

(iii): Since the summands in W are independent,

$$\operatorname{Var}(W(t)) = \sum_{i=1}^{\infty} \mathbb{P}(Z_i \leqslant t) (1 - \mathbb{P}(Z_i \leqslant t)) = \sum_{i=1}^{\infty} e^{-\lambda_i t} (1 - e^{-\lambda_i t})$$

$$\leqslant \sum_{i=1}^{\infty} \mathbb{P}(Z_i \leqslant t) = \mathbb{E} W(t). \tag{6.10}$$

(iv): First, by (6.6) and monotone convergence, as $t \to \infty$,

$$\mathbb{E} W(t) \to \sum_{i=1}^{\infty} \mathbb{P}(Z_i < \infty) = \sum_{i=1}^{\infty} \mathbf{1}_{\{\lambda_i > 0\}} = L.$$
 (6.11)

Furthermore, if $L < \infty$, then a.s. W(t) = L for all large t, and thus (6.9) holds.

Suppose now that $L = \infty$. Then $\mathbb{E} W(t) \to \infty$ by (6.11). Let $\delta \in (0,1)$, let $a := 1 + \delta$ and choose, for $n \ge 1$, $t_n > 0$ such that $\mathbb{E} W(t_n) = a^n$. (This is possible by (i).)

By (6.8) and Chebyshev's inequality, for any t > 0,

$$\mathbb{P}\left(\left|\frac{W(t)}{\mathbb{E}W(t)} - 1\right| > \delta\right) \leqslant \frac{\operatorname{Var}(W(t))}{(\delta \mathbb{E}W(t))^2} \leqslant \frac{1}{\delta^2 \mathbb{E}W(t)}.$$
 (6.12)

Hence, by our choice of t_n and the Borel–Cantelli lemma, a.s. there exists a (random) n_0 such that $1 - \delta \leq W(t_n) / \mathbb{E} W(t_n) \leq 1 + \delta$ for $n \geq n_0$. This, and the fact that W(t) is increasing, implies that if $t \geq t_{n_0}$, and we choose $n \geq n_0$ such that $t_n \leq t < t_{n+1}$, then

$$W(t) \leqslant W(t_{n+1}) \leqslant (1+\delta)a^{n+1} = (1+\delta)^2 \mathbb{E} W(t_n) \leqslant (1+\delta)^2 \mathbb{E} W(t),$$
 (6.13) and similarly

$$W(t) \geqslant W(t_n) \geqslant (1 - \delta)a^n \geqslant (1 - \delta)^2 \mathbb{E} W(t_{n+1}) \geqslant (1 - \delta)^2 \mathbb{E} W(t).$$
 (6.14)

Consequently, a.s.

$$(1 - \delta)^2 \leqslant \liminf_{t \to \infty} \frac{W(t)}{\mathbb{E}W(t)} \leqslant \limsup_{t \to \infty} \frac{W(t)}{\mathbb{E}W(t)} \leqslant (1 + \delta)^2.$$
 (6.15)

Since δ is arbitrarily small, (6.9) follows.

(v): By (i), $\mathbb{E}W(t)$ is increasing, and furthermore it is concave with $\mathbb{E}W(0) = 0$, and thus $\mathbb{E}W(t)/t$ is decreasing on $(0, \infty)$. Hence,

$$\min\{1, t_n'/t_n\} \leqslant \mathbb{E} W(t_n') / \mathbb{E} W(t_n) \leqslant \max\{1, t_n'/t_n\}$$

$$(6.16)$$

and the result follows.

In order to extend this to the dependent sum (6.4), we use a lemma.

Lemma 6.2. Let $(I_{ij})_{i,j=1}^N$ be a finite or infinite symmetric array of random indicator variables, with $\{I_{ij}\}_{i\leqslant j}$ independent. Let $I_i := \max_j I_{ij}$, and $W := \sum_i I_i$. Then

$$Var W \le 2 \mathbb{E} W. \tag{6.17}$$

Proof. Assume first that $N < \infty$. Let $\bar{I}_{ij} := 1 - I_{ij}$ and $\bar{I}_i := 1 - I_i = \prod_j \bar{I}_{ij}$. Let $q_{ij} := \mathbb{E} \bar{I}_{ij} = 1 - \mathbb{E} I_{ij}$.

Fix i and j with $i \neq j$, and let $\bar{I}'_i := \prod_{k \neq j} \bar{I}_{ik}$ and $\bar{I}'_j := \prod_{k \neq i} \bar{I}_{jk}$. Then $\bar{I}_i = \bar{I}_{ij}\bar{I}'_i$ and $\bar{I}_j = \bar{I}_{ij}\bar{I}'_j$, with \bar{I}_{ij} , \bar{I}'_i and \bar{I}'_j independent, and thus

$$Cov(I_i, I_j) = Cov(\bar{I}_i, \bar{I}_j) = \mathbb{E}(\bar{I}_{ij}\bar{I}_i'\bar{I}_j') - \mathbb{E}(\bar{I}_{ij}\bar{I}_i') \mathbb{E}(\bar{I}_{ij}\bar{I}_j')$$

$$= q_{ij} \mathbb{E}\bar{I}_i'\bar{I}_j' - (q_{ij} \mathbb{E}\bar{I}_i')(q_{ij}\bar{I}_j') = q_{ij}(1 - q_{ij}) \mathbb{E}\bar{I}_i'\bar{I}_j'.$$
(6.18)

In particular,

$$Cov(I_i, I_j) \leqslant (1 - q_{ij}) \mathbb{E} \bar{I}'_i = \mathbb{P}(I_{ij} = 1, I_{ik} = 0 \text{ for } k \neq j).$$

$$(6.19)$$

Summing over j, we obtain for every i, since the events $\mathcal{E}_j := \{I_{ij} = 1, I_{ik} = 0 \text{ for } k \neq j\}$ in (6.19) are disjoint and with union $\{\sum_i I_{ij} = 1\} = \{I_i = 1\}$,

$$\sum_{j\neq i} \operatorname{Cov}(I_i, I_j) \leqslant \sum_{j\neq i} (1 - q_{ij}) \mathbb{E} \, \bar{I}_i' = \mathbb{P} \Big(\sum_j I_{ij} = 1 \Big) \leqslant \mathbb{P}(I_i = 1) = \mathbb{E} \, I_i.$$

Furthermore, $Cov(I_i, I_i) = \mathbb{E} I_i - (\mathbb{E} I_i)^2 \leq \mathbb{E} I_i$. Consequently, for every i,

$$\sum_{j=1}^{n} \operatorname{Cov}(I_i, I_j) \leqslant 2 \mathbb{E} I_i, \tag{6.20}$$

and (6.17) follows by summing over i.

Lemma 6.3. Let $(Z_{ij})_{ij}$ be a symmetric array of exponential random variables with $\{Z_{ij}\}_{i\leqslant j}$ independent and $Z_{ij} \sim \operatorname{Exp}(\lambda_{ij})$, where $\lambda_{ij} \geqslant 0$ and $0 < \sum_{ij} \lambda_{ij} < \infty$. Let $Z_i := \inf_j Z_{ij}$ and $W(t) := \sum_i \mathbf{1}_{\{Z_i \leqslant t\}}$. Then $Z_i \sim \operatorname{Exp}(\lambda_i)$ with $\lambda_i := \sum_j \lambda_{ij}$. Moreover, all results of Lemma 6.1 hold except (iii), which is replaced by

$$\operatorname{Var}(W(t)) \leq 2 \mathbb{E} W(t).$$
 (6.21)

Proof. It is well-known and elementary that $Z_i \sim \text{Exp}(\lambda_i)$, since $(Z_{ij})_j$ are independent for every i. Parts (i), (ii) and (v) of Lemma 6.1 deal only with the expectation, and their proofs do not need Z_i to be independent.

Lemma 6.2 yields (6.21).

Finally, the proof of (iv) holds as before, now using
$$(6.21)$$
.

We return to the random graphs. We define, for a given measure (matrix) μ , using Lemmas 6.1 and 6.3 together with (6.2)–(6.5), the functions

$$v(t) = v(t; \mu) := \mathbb{E} v(\tilde{G}_t) = \sum_{i=1}^{\infty} (1 - e^{-\mu_i t}) \approx \sum_{i=1}^{\infty} (1 \wedge (\mu_i t)),$$
 (6.22)

$$e(t) = e(t; \mu) := \mathbb{E} e(\tilde{G}_t) = \sum_{i \le j} (1 - e^{-\mu_{ij}t}) \simeq \sum_{i \ne j} (1 \wedge (\mu_{ij}t)).$$
 (6.23)

Since \tilde{G}_t has no isolated vertices, $e(\tilde{G}_t) \geqslant \frac{1}{2}v(\tilde{G}_t)$, and thus, cf. (6.1),

$$e(t) \geqslant \frac{1}{2}v(t). \tag{6.24}$$

Theorem 6.4. Assume that $\mu = (\mu_{ij})_{i,j=1}^{\infty}$ is a symmetric non-negative matrix with $\mu_{ii} = 0$ and $0 < \|\mu\| := \sum_{i < j} \mu_{ij} < \infty$.

(i) $As t \to \infty$,

$$v(\tilde{G}_t)/v(t) \xrightarrow{\text{a.s.}} 1,$$
 (6.25)

$$e(\tilde{G}_t)/e(t) \xrightarrow{\text{a.s.}} 1.$$
 (6.26)

Moreover, if $\mu_{ij} > 0$ for infinitely many pairs (i, j), then as $t \to \infty$, $v(t) \to \infty$, $e(t) \to \infty$ and $v(\tilde{G}_t)$, $e(\tilde{G}_t) \xrightarrow{\text{a.s.}} \infty$.

(ii) $As \ m \to \infty$,

$$v(G_m)/v(\|\mu\|^{-1}m) \xrightarrow{\text{a.s.}} 1, \tag{6.27}$$

$$e(G_m)/e(\|\mu\|^{-1}m) \xrightarrow{\text{a.s.}} 1.$$
 (6.28)

In particular, a.s.

$$v(G_m) \approx \sum_{i=1}^{\infty} (1 \wedge (\mu_i m)),$$
 (6.29)

$$e(G_m) \approx \sum_{i,j} (1 \wedge (\mu_{ij}m)).$$
 (6.30)

Consequently, if $\mu_{ij} > 0$ for infinitely many pairs (i, j), then as $m \to \infty$, a.s. $v(G_m), e(G_m) \to \infty$.

Proof. (i): This is an immediate consequence of Lemma 6.1(iv) and Lemma 6.3.

(ii): Part (i) and Propositions 4.9 and 4.13 show that $v(G_m)/v(\tau_m) = v(\tilde{G}_{\tau_m})/v(\tau_m) \xrightarrow{\text{a.s.}} 1$. Furthermore, $\tau_m \sim \|\mu\|^{-1}m$ by (4.9), and thus $v(\tau_m) \sim v(\|\mu\|^{-1}m)$ by Lemmas 6.1(v) and 6.3. Hence (6.27) follows. The proof of (6.28) is the same.

Finally (6.29)–(6.30) follow by (6.22)–(6.23), and the final sentence follows by monotone convergence (or by Lemma 6.1(iv)).

Hence, to find asymptotics of the numbers of vertices and edges in our random graphs, it suffices to study the expectations in (6.22)–(6.23). In particular, we note the following consequences.

Corollary 6.5. Assume that $\mu = (\mu_{ij})_{i,j=1}^{\infty}$ is a symmetric non-negative matrix with $\mu_{ii} = 0$ and $0 < \|\mu\| := \sum_{i < j} \mu_{ij} < \infty$. Then:

- (i) G_m is a.s. dense if and only if $e(t) \approx v(t)^2$ as $t \to \infty$.
- (ii) G_m is a.s. sparse if and only if $e(t) = o(v(t)^2)$ as $t \to \infty$.
- (iii) G_m is a.s. extremely sparse if and only if $e(t) \approx v(t)$ as $t \to \infty$.

Proof. By Theorem
$$6.4(ii)$$
.

Corollary 6.6. Assume that $\mu = (\mu_{ij})_{i,j=1}^{\infty}$ is a symmetric non-negative matrix with $\mu_{ii} = 0$ and $0 < \|\mu\| := \sum_{i < j} \mu_{ij} < \infty$. Then, a.s.,

- (i) $v(G_m) = o(m)$ and $e(G_m) = o(m)$ as $m \to \infty$;
- (ii) $v(\tilde{G}_t) = o(t)$ and $e(\tilde{G}_t) = o(t)$ as $t \to \infty$.

Proof. By Theorem 6.4, since $e(t)/t \to 0$ and $v(t)/t \to 0$ as $t \to \infty$ by Lemma 6.1(i) and Lemma 6.3.

Remark 6.7. If we consider the random multigraph G_m^* we have (at least in the loop-less case, and in general with a minor modification) $v(G_m^*) = v(G_m) = o(m)$ by Corollary 6.6, while by definition there are m edges. Hence, the average degree $2e(G_m^*)/v(G_m^*) \to \infty$ a.s. as $m \to \infty$. Similarly, the average number of copies of each edge $e(G_m^*)/e(G_m) \to \infty$ a.s.

For future use we note also that since v(t) is concave with v(0) = 0, for any $C \ge 1$,

$$v(t) \leqslant v(Ct) \leqslant Cv(t). \tag{6.31}$$

Hence $v(Ct) \approx v(t)$ for any constant C > 0.

We have so far considered only simple first order properties of $v(G_m)$ and $e(G_m)$. For the number of edges, much more follows from the central limit results in the references mentioned above. In particular, the local and global central limit theorems in [19] apply and yield the following. (Although the estimates (6.32) and (6.35) are uniform in all x, the main interest is for x constant, or perhaps tending to infinity very slowly.)

Theorem 6.8. Let μ be as in Theorem 6.4. The following hold with O(1) bounded by an absolute constant C uniformly for all $m \ge 1$, $x \in \mathbb{R}$, and matrices μ .

Let $\sigma_m^2 := \operatorname{Var}(e(G_m))$. Then

$$\mathbb{P}(e(G_m) = \lfloor \mathbb{E} e(G_m) + x\sigma_m \rfloor) = \frac{e^{-x^2/2}}{\sqrt{2\pi}\sigma_m} + \frac{O(1)}{\sigma_m^2}.$$
 (6.32)

Moreover, assuming for simplicity $\|\mu\| = 1$,

$$\mathbb{E} e(G_m) = \mathbb{E} e(\tilde{G}_m) + O(1), \tag{6.33}$$

$$Var(e(G_m)) = Var(e(\tilde{G}_m)) + O(1), \tag{6.34}$$

and, recalling (6.23) and defining $\tilde{\sigma}_t^2 := \operatorname{Var}(e(\tilde{G}_t))$,

$$\mathbb{P}(e(G_m) = \lfloor e(m) + x\tilde{\sigma}_m \rfloor) = \frac{e^{-x^2/2}}{\sqrt{2\pi}\tilde{\sigma}_m} + \frac{O(1)}{\tilde{\sigma}_m^2}.$$
 (6.35)

In particular, if $m \to \infty$ and $\tilde{\sigma}_m^2 \to \infty$, then $(e(G_m) - \mathbb{E} e(G_m))/\sigma_m \stackrel{\mathrm{d}}{\longrightarrow} N(0,1)$ and $(e(G_m) - e(m))/\tilde{\sigma}_m \stackrel{\mathrm{d}}{\longrightarrow} N(0,1)$.

The O(1) in (6.33)–(6.34) can be replaced by o(1) as $m \to \infty$ for a fixed μ .

Proof. By [19, Theorems 2.1, 2.3, 2.4 and Corollary 2.5, together with Section 9]. \Box

Note that $e(m) = \mathbb{E} e(\tilde{G}_m)$ and $\tilde{\sigma}_m^2 = \operatorname{Var} e(\tilde{G}_m)$ are given by (6.23) and (6.8); they are usually simpler and more convenient to handle than $\mathbb{E} e(G_m)$ and $\sigma_m^2 = \operatorname{Var}(e(G_m))$.

We conjecture that similar results holds for $v(G_m)$, the number of vertices. However, we cannot obtain this directly from results on the occupancy problem in the same way as Theorem 6.8, again because the variables $N_i(t)$ are dependent. (The number of vertices corresponds to an occupancy problem where balls are thrown in pairs, with a dependency inside each pair.)

Problem 6.9. Show asymptotic normality for $v(G_m)$ when $Var(v(G_m)) \to \infty$.

6.1. The case with dust or attached stars. We consider briefly the case when the model contains dust (other than loops) or attached stars. In this case, the results are quite different. We may for simplicity assume that there are no loops at all, since loops are deleted in any case. Thus $\mu_{ii} = 0$ for $i \ge 0$ and $\mu_{0i} > 0$ for some $i \in \mathbb{N} \cup \{-1\}$.

The number of edges in the dust and attached stars of \tilde{G}_t is $\operatorname{Po}(ct)$ with $c:=\sum_{i=-1}^\infty \mu_{0i}>0$, and thus this number is a.s. $\sim ct \asymp t$ as $t\to\infty$, by the law of large numbers for the Poisson process. (Recall that all edges in the dust and attached stars of \tilde{G}_t^* are simple, so the number of them is the same in \tilde{G}_t and in \tilde{G}_t^* .) It follows by Proposition 4.9 that the number of edges in the dust and attached stars of G_m a.s. is $\asymp m$. Moreover, since each edge in the dust or an attached star has at least one endpoint that is not shared by any other edge, the same estimates hold for the number of vertices in the dust and attached stars. This leads to the following theorem, which shows that if there is any dust or attached star all, then those parts will dominate the random graphs.

Theorem 6.10. Assume that $\mu_{0i} > 0$ for some $i \in \mathbb{N} \cup \{-1\}$. Then, a.s.,

- (i) $v(G_m) \simeq m$ and $e(G_m) \simeq m$ as $m \to \infty$;
- (ii) $v(\tilde{G}_t) \simeq t$ and $e(\tilde{G}_t) \simeq t$ as $t \to \infty$.

Moreover, a.s., all but a fraction o(1) of the edges and vertices are in the dust or attached stars.

Consequently, the random graphs G_m are a.s. extremely sparse, but in a rather trivial way.

Proof. The argument before the theorem shows (i) and (ii).

Moreover, Corollary 6.6 applies to the central part of G_t and shows that the number of edges and vertices there a.s. are o(t), and thus only a fraction o(1) of all edges and vertices. By Proposition 4.9, the same holds for G_m . \square

7. Rank 1 multigraphs

We turn to considering specific examples of the construction. One interesting class of examples are constructed as follows.

Example 7.1 (Rank 1). Let $(q_i)_1^{\infty}$ be a probability distribution on \mathbb{N} , and construct a sequence of i.i.d. edges e_1, e_2, \ldots , each obtained by selecting the two endpoints as independent random vertices with the distribution $(q_i)_i$. (Thus loops are possible.) Define the random multigraph G_m^* by taking the m edges e_1, \ldots, e_m , letting the vertex set be the set of their endpoints. (Equivalently: start with the vertex set \mathbb{N} and then remove all isolated vertices.)

In other words, let V_1, V_2, \ldots be an i.i.d. sequence of vertices with the distribution $(q_i)_i$, and let the edges of G_m^* be $V_1V_2, V_3V_4, \ldots, V_{2m-1}V_{2m}$.

This is clearly a random multigraph of the type constructed in Section 5, with

$$\mu_{ij} = \begin{cases} 2q_i q_j, & i \neq j, \\ q_i^2, & i = j. \end{cases}$$
 (7.1)

We thus have, by (5.2),

$$\mu_i = \sum_{j \neq i} 2q_i q_j + q_i^2 = 2q_i - q_i^2.$$
 (7.2)

In particular, $\mu_i \simeq q_i$.

The corresponding Poisson model \tilde{G}_t^* is by Proposition 4.9 obtained by taking a Poisson number of edges $e_1, \ldots, e_{N(t)}$, with $N(t) \sim \text{Po}(t)$.

As usual, we obtain the corresponding simple graphs by omitting all repeated edges and deleting all loops.

We call a random multigraph constructed as in Example 7.1, or equivalently by (7.1), for some (possibly random) probability distribution $(q_i)_1^{\infty}$, a rank 1 edge exchangeable multigraph, for the reason that the matrix (7.1) is a rank 1 matrix except for the diagonal entries.

Remark 7.2. The diagonal entries, creating loops, are less important to us. In the multigraph examples below, it is natural, and simplifies the results, to allow loops. However, when we consider the simple graphs \tilde{G}_t and G_m , we ignore loops and, see Remark 5.1, it is then simpler to modify (7.1) by taking $\mu_{ii} = 0$; we still say that the resulting random graphs are rank 1.

Remark 7.3. Note that the rank 1 random graphs in [2] are different; they are simple graphs, and they are vertex exchangeable or modifications of vertex exchangeable random graphs, cf. Section 5.1. Nevertheless, both types of "rank 1" random graphs can be seen as based on the same idea: each vertex is given an "activity" (q_i in our case), and the probability of an edge between two vertices is proportional to the product of their activities. (See the references in [2] for various versions of this idea.)

Recall that the *configuration model* is an important model for constructing random multigraphs with a given degree sequence, which is defined as follows, see e.g. Bollobás [1].

Definition 7.4 (Configuration model). Given a sequence $(d_i)_{i=1}^n$ of nonnegative integers with $\sum_i d_i$ even, the random multigraph $\tilde{G}^*(n, (d_i)_{i=1}^n)$ is defined by considering a set of $\sum_i d_i$ half-edges (or stubs), of which d_i are labelled i for each $i \in [n]$, and taking a uniformly random matching of the half-edges; each pair of half-edges is interpreted as an edge between the corresponding vertices.

By construction, the multigraph $\tilde{G}^*(n,(d_i)_{i=1}^n)$ has degree sequence $(d_i)_{i=1}^n$. (With a loop counted as 2 edges at its only endpoint.) Note that the distribution of $\tilde{G}^*(n,(d_i)_{i=1}^n)$ is not uniform over all multigraphs with this degree sequence. (As is well-known, and easy to see, the probability distribution has a factor (weight) 1/2 for each loop and $1/\ell!$ for each edge of multiplicity $\ell > 1$; in particular, conditioned on being a simple graph, $\tilde{G}^*(n,(d_i)_{i=1}^n)$ has a uniform distribution.) Nevertheless, $\tilde{G}^*(n,(d_i)_{i=1}^n)$ has the right distribution for our purposes.

Theorem 7.5. The random multigraph G_m^* constructed in Example 7.1 has, conditioned on its degree sequence $(d_i)_{i=1}^n$, the same distribution as the

random multigraph $\tilde{G}^*(n,(d_i)_{i=1}^n)$ constructed by the configuration model for that degree sequence.

The same holds for \tilde{G}_t^* .

Proof. In the construction of G_m^* above, the sequence V_1, \ldots, V_{2m} is i.i.d., and thus exchangeable; hence its distribution is unchanged if we replace each V_i by $V_{\pi(i)}$ for a uniformly random permutation π of [2m], independent of everything else. Consequently, the distribution of G_m^* is the same if we modify the definition above and let the edges be $V_{\pi(1)}V_{\pi(2)}, \ldots, V_{\pi(2m-1)}V_{\pi(2m)}$; but this is the same as saying that the edges are obtained by taking a random matching of the multiset $\{V_1, \ldots, V_{2m}\}$, which is precisely what the configuration model does. (Note that the vertex degree d_i is the number of times i appears in V_1, \ldots, V_{2m} .)

The result for \tilde{G}_t^* follows, since the degree sequence tells how many edges there are, so conditioning on the degree sequence implies conditioning on $e(\tilde{G}_t^*) = N(t)$, which reduces to the case of G_m^* just proved, see Remark 4.10.

Remark 7.6. In statistical language, the theorem implies that the degree distribution is a sufficient statistic for the family of distributions of multigraphs G_m^* (or \tilde{G}_t^*) given by Example 7.1 with different distributions $(q_i)_1^{\infty}$.

Example 7.7. A trivial example of the construction in Example 7.1 is obtained by fixing $n \ge 1$ and letting $q_i = 1/n$, $1 \le i \le n$, i.e., the uniform distribution on [n]. This means that we consider a sequence of i.i.d. edges, each obtained by taking the two endpoints uniformly at random, and independently, from [n]. In other words, the endpoints of the edges are obtained by drawing with replacement from [n]. This gives the random multigraph process studied in e.g. [25], which is a natural multigraph version of the (simple) random graph process studied by Erdős and Rényi [16].

The rank 1 random multigraphs in Example 7.1 appear also hidden in some other examples.

Example 7.8 (The Hollywood model). The Hollywood model of a random hypergraph was defined in Crane and Dempsey [11] using the language of actors participating in the same movie, see [11] for details. We repeat their definition in somewhat different words.

The model can be defined by starting with the two-parameter version of the Chinese restaurant process, see e.g. [35, Section 3.2] and [10], which starts with a single table with one customer. New customers arrive, one by one; if a new customer arrives when there are n customers seated at k tables, with $n_i \ge 1$ customers at table i, then the new customer is placed:

$$\begin{cases} \text{ at table } i \ (1 \leq i \leq k) \text{ with probability } (n_i - \alpha)/(n + \theta), \\ \text{at a new table } k + 1 \text{ with probability } (\theta + k\alpha)/(n + \theta). \end{cases}$$
 (7.3)

Here α and θ are parameters, and either

- (i) $0 \le \alpha \le 1$ and $\theta > -\alpha$, or
- (ii) $\alpha < 0$ and $\theta = N|\alpha| > 0$ for some $N \in \mathbb{N}$.

In case (ii), there are never more than N tables; in case (i), the number of tables grows a.s. to ∞ .

In the construction of the Hollywood model hypergraph, the vertices are the tables in the Chinese restaurant process. We furthermore draw the sizes of the edges as i.i.d. random variables X_j with some distribution ν on the non-negative integers \mathbb{N} . The first edge is then defined by (the set of tables of) the first X_1 customers, the second edge by the next X_2 customers, and so on. The random hypergraph \tilde{G}_m with m edges is thus described by the first $X_1 + \cdots + X_m$ customers.

A standard calculation shows that the sequence of table numbers is exchangeable, except that the numbers occur for the first time in the natural order; to be precise, the probability of any finite sequence of table numbers, such that the first 1 appears before the first 2, and so on, depends only on the number of occurences of each number. Consequently, as noted in [11], since we ignore vertex labels, and the sequence X_1, X_2, \ldots is i.i.d. and independent of the Chinese restaurant process, the random hypergraph \tilde{G}_{∞}^* is exchangeable, and by the representation theorem by Crane and Dempsey [11, 12], see Remark 4.4, the Hollywood model can be constructed as in Definition 4.2 for some random measure μ on \mathbb{N} .

We can see this more concretely by replacing the table labels $i \in \mathbb{N}$ by i.i.d. random labels $U_i \sim U(0,1)$; then the sequence of table labels of the customers is exchangeable. Hence, by de Finetti's theorem, there exists a random probability measure \hat{P} on [0, 1] such that conditioned on \hat{P} , the sequence of (new) table labels is an i.i.d. sequence with distribution \hat{P} . Clearly, the random measure $\hat{P} = \sum_{i} \widetilde{P}_{i} \delta_{U_{i}}$ for some random sequence \widetilde{P}_{i} of numbers with $\sum_{i} \widetilde{P}_{i} = 1$. Furthermore, by the law of large numbers, for every $i \in \mathbb{N}$, P_i equals a.s. the asymptotic frequency of customers sitting at the table originally labelled i in the Chinese restaurant process. Hence, the random probability measure $\widetilde{P} = (\widetilde{P}_i)_1^{\infty}$ on \mathbb{N} has the distribution $GEM(\alpha, \theta)$, see [35, Theorem 3.2 and Definition 3.3]. (An alternative version of this argument uses Kingman's paintbox representation for exchangeable random partitions [35, Theorem 2.2] instead of the random lables U_i above; we leave the details to the interested reader.) Consequently, the Hollywood model hypergraph can be constructed as follows: Let the random probability measure P on N have the distribution $GEM(\alpha, \theta)$; conditionally given P take an infinite i.i.d. sequence of vertices with distribution \widetilde{P} ; construct the edges by taking the first X_1 vertices, the next X_2 vertices, ...; finally, ignore the vertex labels.

We specialize to the graph case and assume from now on that $X_j = 2$ (deterministically). Thus edges are constructed by taking the customers pairwise as they arrive. We then see by comparing the constructions above and in Example 7.1 that the Hollywood model yields the same result as the rank 1 model in Example 7.1, based on a random probability distribution with distribution $GEM(\alpha, \theta)$.

Since the order of the probabilities q_i does not matter in Example 7.1, we obtain the same result if we reorder the probabilities \widetilde{P}_i in decreasing order; this gives the Poisson–Dirichlet distribution $PD(\alpha, \theta)$ [35, Definition 3.3], and thus the Hollywood model is also given by the rank 1 model based on $PD(\alpha, \theta)$.

Theorem 7.5 shows that yet another way to define the Hollywood model multigraph G_m^* is to take the configuration model where the degree sequence $(d_i)_1^m$ is the (random) sequence of numbers of customers at each table in the Chinese restaurant process when there are 2m customers.

Example 7.9. Pittel [36] considers the random multigraph process with a fixed vertex set [N], where edges are added one by one (starting with no edges) such that the probability that a new edge joins two distinct vertices i and j is proportional to $2(d_i + \alpha)(d_j + \alpha)$, and the probability that the new edge is a loop at i is proportional to $(d_i + \alpha)(d_i + 1 + \alpha)$; here d_i is the current degree of vertex i and $\alpha > 0$ is a fixed parameter. ([36] considers also the corresponding process for simple graphs; we do not consider that process here.)

It is easily seen that this multigraph process can be obtained as above, with a minor modification of the Chinese restaurant process. Consider now a restaurant with a fixed number N of tables, initially empty, and seat each new customer at table i with probability

$$(n_i + \alpha)/(n + N\alpha), \tag{7.4}$$

where $n_i \ge 0$ is the number of customers at table i and n is their total number. Then construct edges by taking the customers pairwise, as above; this yields the multigraph process just described.

Furthermore, although this construction uses a modification of the Chinese restaurant process, we can relabel the tables in the random order that they are occupied. It is then easily seen that we obtain the Chinese restaurant process (7.3) with parameters $(-\alpha, N\alpha)$. Since the vertex labels are ignored, this means that Pittel's multigraph process is the same as the Hollywood model with parameters $(-\alpha, N\alpha)$. Consequently, it can be defined by the rank 1 model in Example 7.1 with the random probability distribution $\text{GEM}(-\alpha, N\alpha)$ on $[N] \subset \mathbb{N}$, or, equivalently, the random probability distribution $\text{PD}(-\alpha, N\alpha)$.

Moreover, the restaurant process (7.4) can be seen as a Pólya urn process, with balls of N different colours and initially α balls in each colour, where n_i is the number of additional balls of color i in the urn; balls are drawn uniformly at random from the urn, and each drawn ball is replaced together with a new ball of the same colour. Note that then n_i is the number of times colour i has been drawn. (It does not matter whether α is an integer or not; the extension to non-integer α causes no mathematical problem, see e.g. [20, Remark 4.2], [21] or [28].) The sequence of vertex labels is thus given by the sequence of colours of the balls drawn from this urn. It is well-known, by an explicit calculation, see e.g. [33] (where N=2), that this sequence is exchangeable. By de Finetti's theorem it can thus can be seen as an i.i.d. sequence of colours with a random distribution \hat{P} , which equals the asymptotic colour distribution. Moreover, it is well-known [29] (see also [33] and [37] for N=2) that this asymptotic distribution is a symmetric Dirichlet distribution $Dir(\alpha/N, \ldots, \alpha/N)$, with the density function $c \prod x_i^{\alpha/N-1}$ on the (N-1)-dimensional simplex $\{(x_1,\ldots,x_N) \in$ $\mathbb{R}^N_+: \sum_i x_i = 1$. Consequently, the multigraph process \tilde{G}_N^* can be obtained by the rank 1 model in Example 7.1 with the random probability distribution $Dir(\alpha/N, \ldots, \alpha/N)$.

Alternatively, by Theorem 7.5, G_m^* may be obtained by the configuration model, with vertex degrees given by the first 2m draws in the Pólya urn process described above.

See further [27].

7.1. Rank 1 simple graphs. We will in the following sections study several examples of the simple random graphs G_m in the rank 1 case. We note here a few general formulas. We ignore the trivial case when the probability distribution $\{q_i\}$ is supported on one point. (Then \tilde{G}_t and G_m have only a single vertex and no edges. In fact, the interesting case is when the support of $\{q_i\}$ is infinite.) We thus assume $\max q_i < 1$.

Since we ignore loops when constructing the simple graphs \tilde{G}_t and G_m , we modify (7.1) by taking $\mu_{ii} = 0$, see Remark 7.2; this changes (7.2) to $\mu_i = 2q_i - 2q_i^2$, but we still have $\mu_i \approx q_i$. Thus (6.22) and (6.23) yield

$$v(t) \approx \sum_{i} (1 \wedge (tq_i)),$$
 (7.5)

$$e(t) \approx \sum_{i \neq j} (1 \wedge (tq_i q_j)).$$
 (7.6)

Moreover, adding the diagonal terms to the sum in (7.6) does not affect this estimate, since if we assume as we may that $q_1, q_2 > 0$, then $q_i^2 = O(q_1q_i)$ and $q_1^2 = O(q_1q_2)$, and thus $\sum_i \left(1 \wedge (tq_i^2)\right) = O\left(\sum_{i>1} \left(1 \wedge (tq_1q_i)\right)\right) = O\left(e(t)\right)$. Hence also

$$e(t) \simeq \sum_{i,j} (1 \wedge (tq_i q_j)) \simeq \sum_i v(tq_i).$$
 (7.7)

Note that although we are interested in large t, the argument tq_i in (7.7) is small for large i, so (7.7) requires that we consider v(t) for both large and small t.

Similarly, the expected degree of vertex i in \tilde{G}_t is

$$\mathbb{E} D_i = \sum_{j \neq i} \left(1 - e^{-2tq_i q_j} \right) \approx \sum_{j \neq i} \left(1 \wedge (tq_i q_j) \right) \approx \sum_j \left(1 \wedge (tq_i q_j) \right) \approx v(tq_i).$$
(7.8)

8. Dense examples

We may obtain examples where G_m and \tilde{G}_t are dense by letting μ_{ij} decrease very rapidly.

We begin with an extreme case, which gives complete graphs.

Example 8.1 (Complete graphs). Let $\mu = (\mu_{ij})$ be such that for every $k \ge 2$

$$0 < \sup_{\ell > 1} \mu_{k+1,\ell} \leqslant k^{-4} \min_{i < k} \mu_{k,i}. \tag{8.1}$$

For example, we may take $\mu_{ij} = ((i \vee j)!)^{-4}$, or the rank 1 example $\mu_{ij} = q_i q_j$ with $q_i = \exp(-3^i)$.

We will show that a.s., for all large n, $G_{\binom{n}{2}}$ is the complete graphs K_n .

Define $a_i := \sup_j \mu_{ij}$. Then (8.1) implies, for every $k \ge 2$,

$$a_{k+1} \leqslant k^{-4} \mu_{k1} \leqslant k^{-4} a_k.$$
 (8.2)

In particular, for $k \ge 2$, $a_{k+1} \le \frac{1}{2}a_k$. Moreover, $(k-1)^4 a_k \ge a_k \ge k^4 a_{k+1}$; hence the sequence $k^4 a_{k+1}$ is decreasing for $k \ge 1$.

Define $t_n := (n^3 a_{n+1})^{-1}$. Let Y_n be the number of edges in $\tilde{G}_{t_n}^*$ with at least one endpoint outside [n]. Then, since (8.2) implies that $(k+1)a_{k+1} \le \frac{1}{2}ka_k$ when $k \ge 2$,

$$\mathbb{E} Y_n = \sum_{k \geqslant n+1} \sum_{i < k} t_n \mu_{ki} < t_n \sum_{k \geqslant n+1} k a_k \leqslant 2t_n (n+1) a_{n+1} = \frac{2(n+1)}{n^3}.$$
 (8.3)

Consequently, by Markov's inequality and the Borel–Cantelli lemma, a.s. $Y_n = 0$ for all large n.

On the other hand, if Z_n is the number of pairs (i, j) with $i < j \le n$ such that ij is not an edge of $\tilde{G}_{t_n}^*$, i.e., $N_{ij}(t_n) = 0$, then

$$\mathbb{E} Z_n = \sum_{i < j \le n} \mathbb{P}(N_{ij}(t_n) = 0) = \sum_{i < j \le n} e^{-t_n \mu_{ij}}.$$
 (8.4)

Moreover, if $i < j \le n$, then by (8.1) and (8.2), $\mu_{ij} \ge j^4 a_{j+1} \ge n^4 a_{n+1}$ and thus $t_n \mu_{ij} \ge t_n n^4 a_{n+1} = n$. Hence, (8.4) yields $\mathbb{E} Z_n \le \binom{n}{2} e^{-n}$, and we see, by the Borel–Cantelli lemma again, that a.s. also $Z_n = 0$ for all large n.

We have shown that a.s. for all large n, $\tilde{G}_{t_n}^*$ contains at least one edge ij whenever $i < j \le n$, but no other edges; in other words, the simple graph \tilde{G}_{t_n} is the complete graph K_n . Since K_n has $\binom{n}{2}$ edges, this also means that $G_{\binom{n}{2}} = K_n$, as asserted above.

We have shown that a.s., for all large m, G_m is the complete graph K_n if $m = \binom{n}{2}$; since G_n is an increasing sequence of graphs, it follows that for intermediate values $m = \binom{n}{2} + \ell$, $1 \leq \ell < n$, G_m consist of K_n plus an additional vertex joined to ℓ of the other vertices. We thus have a complete description of the process (G_m) for large m. (And thus also of the process \tilde{G}_t .)

In particular, for all large m, G_m differs from the complete graph K_n with $n = v(G_m)$ by less than n edges, and thus, see Section 3, $\delta_{\square}(G_m, K_n) \leq \|W_{G_m} - W_{K_n}\|_{L^1} \leq 2/n = o(1)$. It follows that in the sense of graph limit theory, $G_m \to \Gamma_1$ a.s., where Γ_1 is the graph limit defined as the limit of the complete graphs, which is the graph limit defined by the constant graphon $W_1(x, y) = 1$ (on any probability space Ω).

The assumption (8.1) in Example 8.1 is is not best possible, and may easily be improved somewhat, but we only wanted to give a class of examples.

Problem 8.2. Find necessary and sufficient conditions on μ for G_m to be complete for all large m of the form $m = \binom{n}{2}$.

Here is another example, where the limit is less trivial.

Example 8.3. Consider a rank 1 example $\mu_{ij} = q_i q_j$, $i \neq j$, where q_i has a geometric decay $q_i \approx b^{-i}$ for some b > 1.

Let $n \ge 1$ and suppose $b^n \le t \le b^{n+1}$. Then the expected number of edges ij in \tilde{G}_t with i+j>n is at most, with $C:=\sup_i b^i q_i < \infty$ and letting $\ell=i+j$,

$$\sum_{i+j>n} tq_i q_j \leqslant t \sum_{i+j>n} C^2 b^{-i} b^{-j} \leqslant b^{n+1} \sum_{\ell \geqslant n+1} \ell C^2 b^{-l} = O(n).$$
 (8.5)

Similarly, the expected number of edges ij with $i + j \leq n$ not in \tilde{G}_t is at most, for $c := \inf_i b^i q_i > 0$,

$$\sum_{i+j\leqslant n} \exp(-2tq_i q_j) \leqslant \sum_{2\leqslant \ell\leqslant n} \sum_{i=1}^{\ell-1} \exp(-tc^2 b^{-\ell}) \leqslant n \sum_{2\leqslant \ell\leqslant n} \exp(-c^2 b^{n-\ell})$$

$$= O(n). \tag{8.6}$$

Moreover, the same argument shows that the expected number of edges ij in \tilde{G}_t with $i+j>n+n^{0.1}$ and the number of non-edges ij with $i+j< n-n^{0.1}$ both are $O(nb^{-n^{0.1}})$; hence the Borel–Cantelli lemma shows that a.s. for every large n and every $t\in [b^n,b^{n+1}]$, \tilde{G}_t contains every edge with $i+j< n-n^{0.1}$ and no edge with $i+j>n+n^{0.1}$; a consequence, we also have $[n-n^{0.1}-1]\subseteq V(\tilde{G}_t)\subseteq [n+n^{0.1}]$. It follows that if H_n is the graph with vertex set $\{1,\ldots,n\}$ and edge set $\{ij:i+j\leqslant n\}$, then a.s. the cutdistance $\delta_{\square}(\tilde{G}_t,H_n)=o(1)$, when $b^n\leqslant t\leqslant b^{n+1}$. As $n\to\infty$, $H_n\to\Gamma_{\mathsf{half}}$, the graph limit defined by the graphon $W(x,y)=\mathbf{1}_{\{x+y\leqslant 1\}}$ on [0,1] (known as the "half-graphon"). Consequently, $\tilde{G}_t\to\Gamma_{\mathsf{half}}$ a.s. as $t\to\infty$. By Proposition 4.9, $G_m\to\Gamma_{\mathsf{half}}$ a.s. as $m\to\infty$.

Example 8.4. Example 8.3 can be generalized without difficulty. Consider, for example, a rank 1 case $\mu_{ij} = q_i q_j$ with

$$q_i = \exp(-ci + O(i^{1-\varepsilon})) \tag{8.7}$$

for some constants c>0 and $\varepsilon>0$. Arguing as in Lemma 8.3 we see that a.s., for every large n and all $t\in [e^{cn},e^{c(n+1)}]$, \tilde{G}_t contains all edges ij with $i+j< n-n^{1-\varepsilon/2}$ and no edges ij with $i+j> n+n^{1-\varepsilon/2}$. Consequently, a.s., $\delta_{\square}(\tilde{G}_t,H_n)=o(1)$ and thus $\tilde{G}_t\to \Gamma_{\mathsf{half}}$ as $t\to\infty$ and $G_m\to \Gamma_{\mathsf{half}}$ as $m\to\infty$.

Example 8.5. Consider the simple graphs \tilde{G}_t and G_m given by the Hollywood model in Example 7.8 in the case $\alpha = 0$. As shown there, the resulting random graphs are the same as the ones given by the rank 1 model with a random probability distribution $(q_i)_1^{\infty}$ having the distribution $GEM(0, \theta)$, where $\theta \in (0, \infty)$ is a parameter.

By a well-known characterization of the GEM distribution, see [35, Theorem 3.2], this means that

$$(q_1, q_2, \dots) = ((1 - X_1), X_1(1 - X_2), X_1X_2(1 - X_3), \dots),$$
 (8.8)

where $X_i \sim \text{Beta}(\theta, 1)$ are i.i.d. In other words, $q_i = (1 - X_i) \prod_{j=1}^{i-1} X_j$, and thus

$$\log(q_i) = \log(1 - X_i) + \sum_{j=1}^{i-1} \log(X_j). \tag{8.9}$$

Hence, by the law of iterated logarithm, a.s.

$$\log(q_i) = -ci + O(\sqrt{i \log \log i}) = -ci + O(i^{0.6}), \tag{8.10}$$

where $c := -\mathbb{E} \log(X_1) = 1/\theta$. Hence, by conditioning on $(q_i)_1^{\infty}$, Example 8.4 applies. Consequently, $G_m \to \Gamma_{\mathsf{half}}$ a.s. as $m \to \infty$ for the Hollywood model with $\alpha = 0$ and any $\theta > 0$.

Example 8.6. For another generalization of Example 8.3, consider the rank 1 case with $q_i \approx \exp(-ci^{\gamma})$ for some c > 0 and $\gamma > 0$. It follows by a similar argument that a.s. $G_m \to W$, where W is the graphon $\mathbf{1}_{\{x^{\gamma} + y^{\gamma} \leq 1\}}$ on [0, 1].

In Examples 8.1–8.6, G_m converges a.s. to some graph limit. There are also many examples, see e.g. Sections 9–10, for which G_m are sparse, which is equivalent to $G_m \to \Gamma_0$, the zero graph limit defined by the graphon W(x,y)=0. In fact, any graph limit can occur as a limit of G_m , at least along a subsequence. Moreover, the following result shows that there exists a "chameleon" example where every graph limit occurs as the limit of some subsequence. (Note that this includes that there is a subsequence converging to the zero graph limit Γ_0 , which means that $e(G_m)=o(v(G_m)^2)$ along this subsequences; hence this example is neither dense nor sparse.)

Theorem 8.7. There exists a matrix $\mu = (\mu_{ij})$ such that a.s. the graphs G_m are dense in the space of graph limits, in the sense that for every graph limit Γ , there exists a subsequence G_{m_ℓ} that converges to Γ .

Proof. Let F_k , $k \ge 1$, be an enumeration of all finite (unlabelled) simple graphs without isolated vertices, each repeated an infinite number of times. Let $v_k := v(F_k)$ and let $f_k(i,j)$ be the adjacency matrix of F_k .

Let $N_0 := 1$ and, inductively, $N_k := kv_k N_{k-1}$ for $k \ge 1$. Let also

$$a_k := \prod_{j=1}^k N_j^{-4}. \tag{8.11}$$

Clearly, $N_k \ge k!$, $N_k \ge N_{k-1}$ and $a_k \le a_{k-1}$. Finally let, for $i \ne j$,

$$\mu_{ij} = a_k f_k \left(\left\lceil \frac{i}{k N_{k-1}} \right\rceil, \left\lceil \frac{j}{k N_{k-1}} \right\rceil \right), \qquad N_{k-1} < i \lor j \leqslant N_k.$$
 (8.12)

Let $I_k := [1, N_k]$ and divide I_k into the v_k subintervals $I_{k,\ell} := [(\ell-1)kN_{k-1} + 1, \ell kN_{k-1}], \ \ell = 1, \ldots v_k$. Note that (8.12) says that if $i \in I_{k,p}$ and $j \in I_{k,q}$ and not both $i, j \in I_{k-1}$, then $\mu_{ij} = a_k f_k(p,q)$.

Let $t_k := N_k a_k^{-1}$. If n > k, then the expected number of edges ij in $\tilde{G}_{t_k}^*$ with $i \vee j \in I_n \setminus I_{n-1}$ is at most, using (8.11),

$$t_k \sum_{i \lor j \in I_n \setminus I_{n-1}} \mu_{ij} \leqslant t_k a_n N_n^2 \leqslant t_k \frac{a_{n-1}}{N_n^2} = \frac{a_{n-1}}{a_k} \frac{N_k}{N_n^2} \leqslant \frac{1}{N_n} \leqslant \frac{1}{n!}.$$
 (8.13)

Hence the probability that $\tilde{G}_{t_k}^*$ contains some edge with endpoint not in $I_k \times I_k$ is at most

$$\sum_{n>k} \frac{1}{n!} \leqslant \frac{1}{k!},\tag{8.14}$$

and by the Borel–Cantelli lemma, a.s. this happens for only finitely many k.

Similarly, if $(i, j) \in I_k^2 \setminus I_{k-1}^2$, then $\mu_{ij} \in \{0, a_k\}$, and the probability that there exists some such pair (i, j) with $\mu_{ij} = a_k$ but $N_{ij}(t_k) = 0$ is at most

$$N_k^2 e^{-t_k a_k} = N_k^2 e^{-N_k} = O(N_k^{-1}) = O(k!^{-1}).$$
(8.15)

Consequently, again by the Borel–Cantelli lemma, a.s. for every large k, there exists no such pair (i, j).

We have shown that a.s. for every large k, the simple graph \tilde{G}_{t_k} contains no edge with an endpoint outside I_k , and for $(i,j) \in I_k^2 \backslash I_{k-1}^2$, recalling (8.12), if $i \in I_{k,p}$ and $j \in I_{k,q}$, then there is an edge ij if and only if $f_k(p,q) = 1$. In particular, since F_k has no isolated vertices, every $i \in I_k$ is the endpoint of some edge in \tilde{G}_{t_k} and thus a vertex, but no $i \notin I_k$ is; in other words, a.s. for every large k, $V(\tilde{G}_{t_k}) = I_k$. It follows that if F_k^* is the blow-up of F_k with every vertex repeated kN_{k-1} times, then a.s. for every large k, the graphs \tilde{G}_{t_k} and H_k^* have the same vertex set I_k and their adjacency matrices can differ only for $(i,j) \in I_{k-1}^2$. Consequently, using Remark 3.1,

$$\delta_{\square}(\tilde{G}_{t_k}, F_k) = \delta_{\square}(\tilde{G}_{t_k}, F_k^*) \leqslant \|W_{\tilde{G}_{t_k}} - W_{F_k^*}\|_{\square} \leqslant \|W_{\tilde{G}_{t_k}} - W_{F_k^*}\|_{L^1}$$

$$\leqslant \frac{N_{k-1}^2}{N_k^2} = \frac{1}{(kv_k)^2} \leqslant k^{-2},$$
(8.16)

a.s. for all large k.

Now, let Γ be a graph limit, and let $\ell \geq 1$. By graph limit theory (or definition), there exists a sequence of graphs H_j with $v(H_j) \to \infty$ and $\delta_{\square}(H_j,\Gamma) \to 0$ as $j \to \infty$; hence we may take j so large that $H := H_j$ satisfies $v(H) > \ell$ and $\delta_{\square}(H,\Gamma) < 1/\ell$. H may have isolated vertices, so we define H' by choosing a vertex $v \in H$ and adding an edge from v to any other vertex in H. Then at most v(H) - 1 edges are added, and thus, similarly to (8.16),

$$\delta_{\square}(H', H) \leqslant \|W_{H'} - W_H\|_{\square} \leqslant \|W_{H'} - W_H\|_{L^1} \leqslant \frac{2v(H)}{v(H)^2} < \frac{2}{\ell}. \tag{8.17}$$

Moreover, H' has no isolated vertices, and thus H' occurs infinitely often in the sequence (F_k) above. Consequently, a.s., there exists $k > \ell$ such that (8.16) holds and $F_k = H'$. Then, by (8.16) and (8.17),

$$\delta_{\Box}(\tilde{G}_{t_k}, \Gamma) \leq \delta_{\Box}(\tilde{G}_{t_k}, F_k) + \delta_{\Box}(H', H) + \delta_{\Box}(H, \Gamma) < \frac{1}{k^2} + \frac{2}{\ell} + \frac{1}{\ell} < \frac{4}{\ell}.$$
 (8.18)

By Proposition 4.9, this means that if $m_{\ell} := N(t_k)$, then $\delta_{\square}(G_{m_{\ell}}, \Gamma) < 4/\ell$. Since ℓ is arbitrary, this completes the proof. (We may choose m_{ℓ} inductively, and choose k above so large that $m_{\ell} > m_{\ell-1}$.)

The chameleon example in Theorem 8.7 is theoretically very interesting, but it is hardly useful as a model in applications; since the behaviour of G_m changes so completely with m, it is a model of nothing rather than a model of everything.

If we want convergence of the full sequence G_m and not just subsequence convergence as in Theorem 8.7, we do not know whether every graph limit can occur as a limit.

Problem 8.8. For which graph limits Γ does there exist a matrix (μ_{ij}) such that for the corresponding simple random graphs, $G_m \to \Gamma$?

9. Sparse examples

We gave in the preceding section some dense examples. It seems to be more typical, however, that the graph G_m contains many vertices of small degree (maybe even degree 1), and that the graph is sparse. We give here a few, related, rank 1 examples; see also the following section.

Example 9.1. Consider a rank 1 example with $q_i \approx i^{-\gamma}$ for some $\gamma > 1$. Then (7.5) yields

$$v(t) \approx \sum_{i \geqslant 1} \left(1 \wedge (ti^{-\gamma}) \right) = \sum_{i \leqslant t^{1/\gamma}} 1 + t \sum_{i > t^{1/\gamma}} i^{-\gamma} \approx \begin{cases} t^{1/\gamma}, & t \geqslant 1, \\ t, & t < 1. \end{cases}$$
(9.1)

This yields by (7.7), for $t \ge 2$,

$$e(t) \approx \sum_{i \geqslant 1} v(tq_i) \approx \sum_{i \geqslant 1} v(ti^{-\gamma}) \approx \sum_{i \leqslant t^{1/\gamma}} t^{1/\gamma} i^{-1} + \sum_{i > t^{1/\gamma}} ti^{-\gamma} \approx t^{1/\gamma} \log t.$$

$$(9.2)$$

Hence, using Theorem 6.4, a.s. $v(\tilde{G}_t) \approx t^{1/\gamma}$ and $e(\tilde{G}_t) \approx t^{1/\gamma} \log t$ as $t \to \infty$, and $v(G_m) \approx m^{1/\gamma}$ and $e(G_m) \approx m^{1/\gamma} \log m$ as $m \to \infty$. It follows that the average degree in G_m is $\approx \log m$.

In this example we may also show that the degree distribution has a power-law; we state this as a theorem. There is no standard precise definition of what is meant by a power-law degree distribution; we may say that a random variable X has a power-law distribution with exponent τ if $\mathbb{P}(X>x) \asymp x^{-(\tau-1)}$ as $x\to\infty$, but this does not make sense for the degree distribution of a finite graph, so we must either consider the asymptotic degree distribution, provided one exists, or give uniform estimates for a suitable range of x. (See e.g. [18, Sections 1.4.1 and 1.7] for a discussion of power-laws for degree distributions.) We follow here the second possibility.

For a (finite) graph G, let $v_{\geqslant k}(G)$ be the number of vertices of degree at least k, and let $\pi_{\geqslant k}(G) := v_{\geqslant k}(G)/v(G)$, the probability that a random vertex has degree $\geqslant k$.

Theorem 9.2. In Example 9.1, the random graphs G_m have a power-law distribution with exponent 2 in the following sense. There exist positive constants c and C such that a.s. for every large m,

$$\pi_{\geqslant k}(G_m) \leqslant C/k, \qquad 1 \leqslant k < \infty,$$
 (9.3)

$$\pi_{\geqslant k}(G_m) \geqslant c/k, \qquad 1 \leqslant k \leqslant cv(G_m).$$
 (9.4)

As usual, the same result holds for \tilde{G}_t . Note that the restriction $k \leq cv(G_m)$ in (9.4) is necessary, and best possible (up to the value of the constants); we necessarily have $\pi_{\geq k}(G) = 0$ when $k \geq v(G)$. Note also that we have the same exponent $\tau = 2$ for every $\gamma > 1$.

Proof. As usual, we prove the results for \tilde{G}_t ; the results for G_m follow by Proposition 4.9. We then can write (9.3)–(9.4) as $v_{\geqslant k}(\tilde{G}_t) \leqslant Cv(\tilde{G}_t)/k$, $k \geqslant 1$, and $v_{\geqslant k}(\tilde{G}_t) \geqslant cv(\tilde{G}_t)/k$, $1 \leqslant k \leqslant cv(\tilde{G}_t)$, and by Theorem 6.4 and (9.1), it suffices (and is equivalent) to prove that a.s.

$$v_{\geqslant k}(\tilde{G}_t) \leqslant C_1 t^{1/\gamma}/k, \qquad 1 \leqslant k < \infty,$$
 (9.5)

$$v_{\geqslant k}(\tilde{G}_t) \geqslant c_1 t^{1/\gamma}/k, \qquad 1 \leqslant k \leqslant c_2 t^{1/\gamma},$$

$$(9.6)$$

for every large t.

Let I_{ij} be the indicator of an edge ij in \tilde{G}_t ; thus $I_{ij} \sim \text{Be}(1 - e^{-2tq_iq_j})$ Let $D_i := \sum_{i \neq i} I_{ij}$ be the degree of i in the simple graph \tilde{G}_t . (The degree is defined as 0 if i is not a vertex.)

(i) The upper bound (9.5): We fix $t \ge 1$ and an integer $k \ge 1$; for convenience we often omit them from the notation, but note that many variables below depend on them, while all explicit and implicit constant are independent of t and k.

Let
$$J_i := \mathbf{1}_{\{D_i \geqslant k\}}$$
 and $N := \sum_i J_i = v_{\geqslant k}(\tilde{G}_t)$.

Let A be a large constant, chosen later, and assume that $k \ge A$, let $i_0 := At^{1/\gamma}/k$ and let $N^* := \sum_{i>i_0} J_i$. Thus $N \leq N^* + i_0$. If $i \geq i_0$, then using (7.8), (6.31) and (9.1),

$$\mathbb{E} D_i \simeq v(tq_i) \simeq v(ti^{-\gamma}) \leqslant v(A^{-\gamma}k^{\gamma}) \simeq k/A. \tag{9.7}$$

Thus $\mathbb{E} D_i \leqslant C_2 k/A$ for some $C_2 \geqslant 0$, and choosing $A = \max(14C_2, 4)$, we find that $\mathbb{E} D_i \leq k/14 \leq (k-1)/7$. Since D_i is a sum $\sum_j I_{ij}$ of independent Bernoulli variables, a Chernoff bound (see e.g. [26, (2.11) and Theorem 2.8]) yields

$$\mathbb{E} J_i = \mathbb{P}(D_i \geqslant k) \leqslant e^{-k}, \qquad i \geqslant i_0, \tag{9.8}$$

and also, for later use,

$$\mathbb{P}(D_i \geqslant k - 1) \leqslant e^{1 - k}, \qquad i \geqslant i_0. \tag{9.9}$$

For $i \ge t^{1/\gamma}$ we also have, by (9.7) and (9.1),

$$\mathbb{E} D_i \simeq v(ti^{-\gamma}) \simeq ti^{-\gamma}. \tag{9.10}$$

Let $(x)_r := x(x-1)\cdots(x-r+1)$, the falling factorial. Since D_i is a sum of independent indicators, it is easily seen that for any positive integer r, the factorial moment can be bounded by $\mathbb{E}(D_i)_r \leq (\mathbb{E}\,D_i)^r$. Hence, by (9.10) and Markov's inequality, since we assume $k \ge A \ge 4$,

$$\mathbb{E} J_i = \mathbb{P}(D_i \geqslant k) \leqslant \frac{\mathbb{E}(D_i)_4}{(k)_4} \leqslant \frac{(\mathbb{E} D_i)^4}{(k)_4} \leqslant C_3 \frac{(ti^{-\gamma})^4}{k^4} \leqslant C_3 \frac{ti^{-\gamma}}{k^4}, \qquad i \geqslant t^{1/\gamma}.$$

$$(9.11)$$

(This also follows from [26, (2.10) and Theorem 2.8].) Summing (9.8) and (9.11), we obtain

$$\mathbb{E} N^* = \sum_{i>i_0} \mathbb{E} J_i \leqslant \sum_{i_0 < i \leqslant t^{1/\gamma}} e^{-k} + \sum_{i>t^{1/\gamma}} C_3 t i^{-\gamma} / k^4 \leqslant C_4 t^{1/\gamma} / k^4. \tag{9.12}$$

For the variance of N^* , we note that the indicators J_i are not quite independent, since an edge ij influences both J_i and J_j , but conditioned on I_{ij} , J_i and J_j are independent. Hence, for any distinct i and j,

$$\mathbb{E}(J_{i}J_{j}) = \mathbb{P}(I_{ij} = 1) \,\mathbb{E}(J_{i}J_{j} \mid I_{ij} = 1) + \mathbb{P}(I_{ij} = 0) \,\mathbb{E}(J_{i}J_{j} \mid I_{ij} = 0)$$

$$= \mathbb{P}(I_{ij} = 1) \,\mathbb{E}(J_{i} \mid I_{ij} = 1) \,\mathbb{E}(J_{j} \mid I_{ij} = 1)$$

$$+ \mathbb{P}(I_{ij} = 0) \,\mathbb{E}(J_{i} \mid I_{ij} = 0) \,\mathbb{E}(J_{j} \mid I_{ij} = 0)$$

$$\leq \mathbb{P}(I_{ij} = 1) \,\mathbb{P}(D_{i} \geq k - 1) \,\mathbb{P}(D_{j} \geq k - 1) + \mathbb{P}(I_{ij} = 0) \,\mathbb{E}\,J_{i} \,\mathbb{E}\,J_{j}$$

and thus

$$Cov(J_i, J_j) \leqslant \mathbb{P}(I_{ij} = 1) \, \mathbb{P}(D_i \geqslant k - 1) \, \mathbb{P}(D_j \geqslant k - 1). \tag{9.13}$$

By (9.13) and (9.9), for $i, j \ge i_0$ with $i \ne j$,

$$Cov(J_i, J_j) \le 2tq_iq_je^{2(1-k)} \le C_5ti^{-\gamma}j^{-\gamma}e^{-2k}.$$
 (9.14)

Consequently, using also (9.12),

$$\operatorname{Var} N^* = \sum_{i,j>i_0} \operatorname{Cov}(J_i, J_j) \leqslant \mathbb{E} N^* + C_5 t e^{-2k} \sum_{i,j>i_0} i^{-\gamma} j^{-\gamma}$$

$$\leqslant C_4 t^{1/\gamma} k^{-4} + C_6 t e^{-2k} i_0^{2(1-\gamma)} \leqslant C_7 t^{1/\gamma} k^{-4}. \tag{9.15}$$

Hence, by Chebyshev's inequality,

$$\mathbb{P}(N^* - \mathbb{E} N^* > t^{1/\gamma}/k) \leqslant \frac{\text{Var } N^*}{(t^{1/\gamma}/k)^2} \leqslant C_7 t^{-1/\gamma} k^{-2}. \tag{9.16}$$

We have so far kept t and k fixed. We now sum (9.16) over all $k \ge A$ and $t = 2^{\ell}$ for $\ell \in \mathbb{N}$, and find by the Borel–Cantelli lemma that a.s. for every large t of this form and every $k \ge A$, $N^* - \mathbb{E} N^* \le t^{1/\gamma}/k$, and consequently, using also (9.12),

$$N \leqslant N^* + i_0 \leqslant \mathbb{E} N^* + t^{1/\gamma}/k + i_0 \leqslant C_8 t^{1/\gamma}/k.$$
 (9.17)

This is (9.5) for $k \ge A$ and $t \in \{2^{\ell}\}$; since N increases with t, (9.5) follows in general (with a different constant), a.s. for large t and all $k \ge A$.

For k < A, (9.3) and (9.5) follow trivially from $v_{\geq k}(\tilde{G}_t) \leq v(\tilde{G}_t)$.

(ii) The lower bound (9.6): Fix again $t \ge 1$ and $k \ge 1$, let B be a large constant chosen later, and assume that $k \le t^{1/\gamma}/B$.

Let L be the set of odd integers i with $1 \leq i \leq i_1 := B^{-1}t^{1/\gamma}/k$, and let R be the set of even integers j with $1 \leq j \leq 6k$. By our assumption on $k, i_1 \geq 1$, and thus $|L| = \lfloor (i_1+1)/2 \rfloor \geq i_1/3$. Note that the indicators $\{I_{ij}\}_{i \in L, j \in R}$ are independent. For $i \in L$, let $D_i' := \sum_{j \in R} I_{ij}$ and $J_i' = \mathbf{1}_{\{D_i' \geq k\}}$. Thus the indicators $\{J_i'\}_{i \in L}$ are independent. Also, let $N' := \sum_{i \in L} J_i'$. Since $J_i' \leq J_i$, we have $N' \leq \sum_{i \in L} J_i \leq \sum_{i \geq 1} J_i = N = v_{\geq k}(\tilde{G}_t)$.

If $i \in L$ and $j \in R$, then $ij \leqslant 6ki_1 = 6B^{-1}t^{1/\gamma}$, and thus

$$tq_iq_j \geqslant c_3ti^{-\gamma}j^{-\gamma} \geqslant c_4B^{\gamma}. \tag{9.18}$$

Choose $B := c_4^{1/\gamma}$; then by (9.18), when $i \in L$ and $j \in R$, $tq_iq_j \geqslant 1$ and thus

$$\mathbb{P}(I_{ij} = 0) = e^{-2tq_iq_j} \leqslant e^{-2}.$$
(9.19)

Since |R| = 3k, it follows that if $i \in L$, then $\mathbb{E} D_i' \geqslant 3(1 - e^{-2})k > 2.5k$, and moreover, by a Chernoff bound (e.g. [26, (2.12)]),

$$\mathbb{P}(J_i' = 0) = \mathbb{P}(D_i' < k) \leqslant e^{-k} \leqslant e^{-1}. \tag{9.20}$$

Since the indicators J_i' are independent for $i \in L$, another Chernoff bound shows that

$$\mathbb{P}(N' < |L|/2) \leqslant e^{-c_5|L|} \leqslant e^{-c_6 i_1}. \tag{9.21}$$

Alternatively, (9.20) and a union bound yield

$$\mathbb{P}(N' < |L|/2) \leqslant \mathbb{P}(N' < |L|) \leqslant \sum_{i \in L} \mathbb{P}(J_i' = 0) \leqslant i_1 e^{-k}. \tag{9.22}$$

If $1 \leqslant k \leqslant t^{1/2\gamma}$, then $i_1 \geqslant B^{-1}t^{1/2\gamma}$, and thus (9.21) yields $\mathbb{P}(N' < |L|/2) \leqslant e^{-c_7t^{1/2\gamma}}$. If $t^{1/2\gamma} < k \leqslant t^{1/\gamma}/B$, then (9.22) yields $\mathbb{P}(N' < |L|/2) \leqslant i_1e^{-c_8t^{1/2\gamma}} \leqslant C_9e^{-c_9t^{1/2\gamma}}$. Consequently, for every $k \leqslant t^{1/\gamma}/B$,

$$\mathbb{P}(N < |L|/2) \leqslant \mathbb{P}(N' < |L|/2) \leqslant C_{10}e^{-c_{10}t^{1/2\gamma}}.$$
 (9.23)

We have kept k and t fixed, but we now sum (9.23) over all $k
leq t^{1/\gamma}/B$ and $t = 2^{\ell}$ for some $\ell \in \mathbb{N}_0$. It follows by the Borel-Cantelli lemma that a.s. for every large t of this form and every $k
leq t^{1/\gamma}/B$, $N
ge |L|/2
ge i_1/6
ge c_{11}t^{1/\gamma}/k$. This proves (9.6) for t of the form 2^{ℓ} , and again the general case follows since N is monotone in t.

Furthermore, assuming $q_i \sim ci^{-\gamma}$ in Example 9.1, we can show that \tilde{G}_t and G_m converge a.s. to a graphon of the type defined by Veitch and Roy [39] and mentioned in Section 5.1; these graphons are measurable functions $W: \mathbb{R}^2_+ \to [0,1]$, such that the random graphs G(t,W) defined in (5.5) are a.s. finite. (See [39] for precise conditions; see also [9] and [3] for related versions.) Recall that the standard graphons discussed in Section 3 are useful for dense graphs, but not for sparse graphs as here; the more general graphons in [39] are intended for sparse graphs.

Veitch and Roy [40] defined two notions \to_{GP} and \to_{GS} of convergence for such general graphons on \mathbb{R}_+ (and the even more general graphexes defined in [39]) based on convergence in distribution of the corresponding random graphs G(t,W). We can define $W_n \to_{\mathsf{GP}} W$ as meaning $G(r,W_n) \stackrel{\mathrm{d}}{\longrightarrow} G(r,W)$ for each fixed $r < \infty$, see further [40] and [24].

Furthermore, the random graphs G(r,W) are naturally coupled for different r and form an increasing graph process $(G(r,W))_{r\geqslant 0}$. Let $(G_{\tau_k}(W))_k$ be the sequence of different graphs that occur among G(r,W) for $r\geqslant 0$. Then $W_n\to_{\mathsf{GS}} W$ if $(G_{\tau_k}(W_n))_k\stackrel{\mathrm{d}}{\longrightarrow} (G_{\tau_k}(W))_k$; again see further [40] and [24].

Recall that for a finite graph G, we defined a corresponding graphon W_G in Section 3. In the context of graphons on \mathbb{R}_+ , Veitch and Roy [40] define for every s>0 a modification $W_{G,s}$, called the dilated empirical graphon, as follows. We may assume that G has vertices labelled $1, \ldots, v(G)$; then $W_G(i,j) := \mathbf{1}_{\{i \sim j\}}$ for $i,j \leq v(G)$; we extend this by $W_G(i,j) := 0$ when $i \vee j > v(G)$. Then, for every s>0, let the dilated graphon $W_{G,s}$ be the function $\mathbb{R}^2_+ \to \{0,1\}$ given by $W_{G,s}(x,y) := W_G(\lceil sx \rceil, \lceil sy \rceil)$. Hence, every vertex in G corresponds to an interval of length 1/s in the domain of $W_{G,s}$.

If G_n is a sequence of graphs and W a graphon, then $G_n \to_{\mathsf{GS}} W$ means that $W_{G_n} \to_{\mathsf{GS}} W$; furthermore, the convergence \to_{GS} is insensitive to dilations, so $G_n \to_{\mathsf{GS}} W$ is equivalent to $W_{G_n,s_n} \to_{\mathsf{GS}} W$ for any sequence $s_n > 0$.

Remark 9.3. We have in Section 5.1 given the version of G(r, W) without loops; more generally, one can allow i = j in (5.4) and thus allow loops. The loopless case considered here then is obtained by imposing W(x, x) = 0 for x > 0. Hence, for the version with loops, Theorem 9.4 below still holds, provided we redefine W to be 0 on the diagonal.

Theorem 9.4. In Example 9.1, assume that $q_i \sim ci^{-\gamma}$ as $i \to \infty$, with c > 0. Then the dilated empirical graphon $W_{\tilde{G}_t,t^{1/2\gamma}} \to_{\mathsf{GP}} W$ a.s. as $t \to \infty$, where W is the graphon $W(x,y) = 1 - \exp(-2c^2x^{-\gamma}y^{-\gamma})$ on \mathbb{R}^2_+ .

As a consequence, $\tilde{G}_t \to_{\mathsf{GS}} W$ a.s. as $t \to \infty$.

Note that $W(x,y) \ge 1 - \exp(-2c^2) > 0$ when $xy \le 1$, and thus $\int W = \infty$. We prove first two lemmas.

Lemma 9.5. Let $(Z_{kl})_{k,l}$ be an array of i.i.d. random variables. Furthermore, let $x_1, \ldots, x_n > 0$ be distinct and let X be a random variable, independent of the array $(Z_{kl})_{k,l}$, with $X \sim U(a,b)$ where $0 < a < b < \infty$. Then,

$$\mathcal{L}((Z_{\lceil tx_i \rceil, \lceil tX \rceil})_{i=1}^n \mid (Z_{kl})_{k,l}) \xrightarrow{\text{a.s.}} \mathcal{L}((Z_{i,n+1})_{i=1}^n)$$
(9.24)

as $t \to \infty$.

In other words, conditionally on $(Z_{kl})_{k,l}$ and for a.e. every realization of $(Z_{kl})_{k,l}$, the random vector $(Z_{\lceil tx_i \rceil, \lceil tX \rceil})_{i=1}^n$ converges in distribution to $(Z'_{i,n+1})_{i=1}^n$, where $(Z'_{kl})_{k,l}$ is an independent copy of $(Z_{kl})_{k,l}$.

Proof. It suffices to prove that for every fixed rational z_1, \ldots, z_n ,

$$\mathbb{P}(Z_{\lceil tx_i \rceil, \lceil tX \rceil} \leqslant z_i, 1 \leqslant i \leqslant n \mid (Z_{kl})_{k,l}) \xrightarrow{\text{a.s.}} \pi := \mathbb{P}(Z_{i,n+1} \leqslant z_i, 1 \leqslant i \leqslant n),$$
(9.25)

where

$$\pi = \prod_{i=1}^{n} \mathbb{P}(Z_{i,n+1} \leqslant z_i) = \prod_{i=1}^{n} \mathbb{P}(Z_{11} \leqslant z_i).$$
 (9.26)

Let further $I_{k,l,i} := \mathbf{1}_{\{Z_{kl} \leq z_i\}}$ and $J_l := \prod_{i=1}^n I_{\lceil tx_i \rceil,l,i}$. Then, with the error term coming from edge effects,

$$P_t := \mathbb{P}\Big(Z_{\lceil tx_i \rceil, \lceil tX \rceil} \leqslant z_i, 1 \leqslant i \leqslant n \mid (Z_{kl})_{k,l}\Big) = \mathbb{E}\Big(\prod_{i=1}^n I_{\lceil tx_i \rceil, \lceil tX \rceil, i} \mid (Z_{kl})_{k,l}\Big)$$

$$= \mathbb{E}\left(J_{\lceil tX \rceil} \mid (Z_{kl})_{k,l}\right) = \frac{1}{(b-a)t} \sum_{ta < l \le tb} J_l + o(1). \tag{9.27}$$

If t is sufficiently large, then $\lceil tx_1 \rceil, \ldots, \lceil tx_n \rceil$ are distinct, and then, see (9.26),

$$\mathbb{E} J_l = \prod_{i=1}^n \mathbb{E} I_{\lceil tx_i \rceil, l, i} = \prod_{i=1}^n \mathbb{P} \left(Z_{\lceil tx_i \rceil, l} \leqslant z_i \right) = \pi. \tag{9.28}$$

Furthermore, then the variables $J_l \sim \text{Be}(\pi)$ are i.i.d., so their sum in (9.27) has a binomial distribution, and a Chernoff bound shows that for every $\varepsilon > 0$, there is a $c = c(\varepsilon) > 0$ such that for large t,

$$\mathbb{P}(|P_t - \pi| > \varepsilon) \leqslant e^{-ct}. \tag{9.29}$$

This shows that P_t converges to π in probability as $t \to \infty$. In order to show convergence a.s., we note that if 0 < t < u, and t(b-a) > 1, then (for fixed a and b) $\mathbb{P}(\lceil tX \rceil \neq \lceil uX \rceil) = O(u-t)$, and consequently, for some C > 0,

$$|P_t - P_u| \le \mathbb{P}(\lceil tX \rceil \ne \lceil uX \rceil) \le C(u - t).$$
 (9.30)

Let $\varepsilon > 0$, let $N := \lceil C/\varepsilon \rceil$ and let $t_n := n/N$. By (9.29) and the Borel–Cantelli lemma, a.s. $|P_{t_n} - \pi| \leq \varepsilon$ for all large n. Furthermore, if n is large

and $t_n \leqslant t \leqslant t_{n+1}$, then (9.30) implies $|P_t - P_{t_n}| \leqslant \varepsilon$, and thus $|P_t - \pi| \leqslant 2\varepsilon$. Consequently, a.s., $|P_t - \pi| \leqslant 2\varepsilon$ for every large t. Since ε is arbitrary, this proves (9.25) and thus the lemma.

Lemma 9.6. Let $(Z_{kl})_{k,l}$ be an array of i.i.d. random variables, and let (X_1, \ldots, X_n) be a random vector in \mathbb{R}^n_+ with an absolutely continuous distribution, independent of the array $(Z_{kl})_{k,l}$. Then,

$$\mathcal{L}((Z_{\lceil tX_i \rceil, \lceil tX_j \rceil})_{1 \leqslant i < j \leqslant n} \mid (Z_{kl})_{k,l}) \xrightarrow{\text{a.s.}} \mathcal{L}((Z_{i,j})_{1 \leqslant i < j \leqslant n})$$
(9.31)

as $t \to \infty$.

Proof. Step 1. Assume first that X_1, \ldots, X_n are independent with $X_i \sim U(I_i)$ for some intervals I_1, \ldots, I_n . In this case we prove (9.31) by induction on n, so we may assume that

$$\mathcal{L}((Z_{\lceil tX_i \rceil, \lceil tX_j \rceil})_{1 \leqslant i < j \leqslant n-1} \mid (Z_{kl})_{k,l}) \xrightarrow{\text{a.s.}} \mathcal{L}((Z_{i,j})_{1 \leqslant i < j \leqslant n-1}). \tag{9.32}$$

Furthermore, by Lemma 9.5 and conditioning on X_1, \ldots, X_{n-1} ,

$$\mathcal{L}((Z_{\lceil tX_i \rceil, \lceil tX_n \rceil})_{1 \leqslant i \leqslant n-1} \mid (Z_{kl})_{k,l}, X_1, \dots, X_{n-1}) \xrightarrow{\text{a.s.}} \mathcal{L}((Z_{i,n})_{1 \leqslant i \leqslant n-1}).$$

$$(9.33)$$

The result (9.31) follows by (9.32) and (9.33), which shows the induction step and completes the proof of this step.

Step 2. Suppose that there exists a finite family of disjoint intervals I_k such that the density function $f(x_1, \ldots, x_n)$ of (X_1, \ldots, X_n) is supported on $\left(\bigcup_k I_k\right)^n$ and constant on each $\prod_{i=1}^n I_{k_i}$. Then Step 1 shows that for each sequence k_1, \ldots, k_n of indices, (9.31) holds conditioned on $(X_1, \ldots, X_n) \in \prod_{i=1}^n I_{k_i}$. Hence (9.31) holds unconditioned too.

Step 3. The general case. Let $f(x_1, ..., x_n)$ be the density function of $(X_1, ..., X_n)$, and let $\varepsilon > 0$. Then there exists a density function $f_0(x_1, ..., x_n)$ of the type in Step 2 such that $\int |f - f_0| dx_1 ... dx_n < \varepsilon$. We can interpret f_0 as the density function of a random vector $\mathbf{X}^0 = (X_1^0, ..., X_n^0)$, and we can couple this vector with $\mathbf{X} = (X_1, ..., X_n)$ such that $\mathbb{P}(\mathbf{X} \neq \mathbf{X}^0) < \varepsilon$.

Since Step 2 applies to \mathbf{X}^0 , it follows that

$$\mathbb{P}(\text{the convergence in } (9.31) \text{ holds}) \geqslant \mathbb{P}((X,Y) = (X_0,Y_0)) > 1 - \varepsilon.$$
 (9.34) Since $\varepsilon > 0$ is arbitrary, (9.31) follows.

Proof of Theorem 9.4. Let $w_x := c^{-1}q_{\lceil x \rceil}x^{\gamma} = 1 + o(1)$, as $x \to \infty$.

We can construct G_t for all t > 0 by taking i.i.d. random variables $Z_{kl} \sim \text{Exp}(1)$ and letting there be an edge kl in \tilde{G}_t if $2tq_kq_l \geqslant Z_{kl}$, for every pair (k,l) with k < l.

Let $\hat{W}_t := W_{\tilde{G}_t,t^{1/2\gamma}}$ be the dilated empirical graphon in the statement. Fix r > 0, and consider the random graph $G(r, \hat{W}_t)$; this is by (5.4)–(5.5) obtained by taking a Poisson process $\{\eta_i\}_i$ on \mathbb{R}_+ with intensity r (where we assume $\eta_1 < \eta_2 < \ldots$), and then taking an edge ij if and only if $\hat{W}_t(\eta_i, \eta_j) = 1$. By the definition of \hat{W}_t , this is equivalent to \tilde{G}_t having an edge between $\lfloor t^{1/2\gamma} \eta_i \rfloor$ and $\lfloor t^{1/2\gamma} \eta_j \rfloor$, and thus by the construction of \tilde{G}_t to (assuming that t is large so that $\lfloor t^{1/2\gamma} \eta_i \rfloor \neq \lfloor t^{1/2\gamma} \eta_j \rfloor$)

$$2tq_{\lceil t^{1/2\gamma}\eta_i\rceil}q_{\lceil t^{1/2\gamma}\eta_i\rceil}\geqslant Z_{\lceil t^{1/2\gamma}\eta_i\rceil,\lceil t^{1/2\gamma}\eta_i\rceil} \tag{9.35}$$

or, equivalently,

$$2c^{2}\eta_{i}^{-\gamma}\eta_{j}^{-\gamma} \geqslant w_{t^{1/2\gamma}\eta_{i}}^{-1}w_{t^{1/2\gamma}\eta_{j}}^{-1}Z_{\lceil t^{1/2\gamma}\eta_{i}\rceil,\lceil t^{1/2\gamma}\eta_{j}\rceil}.$$
(9.36)

Fix $n < \infty$ and consider the edge indicators $I_{i,j,t}$ in $G(r, \hat{W}_t)$ for $1 \le i < j \le n$. Furthermore, fix a large integer N and condition (η_1, \ldots, η_n) on $\lceil N\eta_1 \rceil, \ldots, \lceil N\eta_n \rceil$. By Lemma 9.6, and recalling $w_x = 1 + o(1)$, the distribution of the right-hand side of (9.36) converges a.s. to independent Exp(1) variables, jointly for $1 \le i < j \le n$. Since $I_{i,j,t}$ equals the indicator of (9.36), it follows by first replacing the left-hand side of (9.36) by upper and lower bounds obtained by rounding each η_i down or up to the nearest multiple of 1/N, applying Lemma 9.6 and then letting $N \to \infty$, that

$$\mathcal{L}((I_{i,j,t})_{1 \leqslant i < j \leqslant n} \mid \tilde{G}_t) \to \mathcal{L}((\mathbf{1}\{2c^2\eta_i^{-\gamma}\eta_j^{-\gamma} \geqslant Z_{ij}\})_{1 \leqslant i < j \leqslant n}). \tag{9.37}$$

Here, conditioned on η_1, \ldots, η_n , the indicators in the right-hand side are independent, and have (conditional) expectations

$$\mathbb{P}(2c^{2}\eta_{i}^{-\gamma}\eta_{j}^{-\gamma} \geqslant Z_{ij} \mid \eta_{i}, \eta_{j}) = 1 - \exp(-2c^{2}\eta_{i}^{-\gamma}\eta_{j}^{-\gamma}) = W(\eta_{i}, \eta_{j}). \quad (9.38)$$

This equals the (conditional) probability of an edge ij in G(r, W). Consequently, if $I_{i,j}$ is the indicator of an edge ij in G(r, W), (9.37) shows that a.s., as $t \to \infty$,

$$\left((I_{i,j,t})_{1 \leqslant i < j \leqslant n} \mid \tilde{G}_t \right) \xrightarrow{\mathrm{d}} (I_{i,j})_{1 \leqslant i < j \leqslant n} \tag{9.39}$$

This shows the desired convergence $G(r, \hat{W}_t) \stackrel{\mathrm{d}}{\longrightarrow} G(r, W)$, provided we restrict the graphs to a fixed finite set of vertices.

To extend this to the infinite number of potential vertices, we need a tightness argument. (Unfortunately, we did not find a really simple argument.) Let a, b > 0, and let $V_{a,b,t}$ denote the number of edges in $G(r, \hat{W}_t)$ with endpoints labelled η_i, η_j with $\eta_i \in (a, 2a]$ and $\eta_j \in (b, 2b]$. Then, cf. (9.35),

$$\mathbb{E}(V_{a,b,t} \mid (Z_{kl})_{k,l}) \leqslant \sum_{k=\lceil at^{1/2\gamma} \rceil}^{\lceil 2at^{1/2\gamma} \rceil} \sum_{l=\lceil bt^{1/2\gamma} \rceil}^{\lceil 2bt^{1/2\gamma} \rceil} r^2 t^{-1/\gamma} \mathbf{1}\{2tq_k q_l \geqslant Z_{k,l}\}.$$
(9.40)

For k, l in the ranges in (9.40), $q_k \leqslant C_1 a^{-\gamma} t^{-1/2}$ and $q_l \leqslant C_1 b^{-\gamma} t^{-1/2}$. Define $J_{kl} := \mathbf{1}\{2C_1^2 a^{-\gamma} b^{-\gamma} \geqslant Z_{k,l}\}, \ S_{m,n} := \sum_{k \leqslant m, l \leqslant n} J_{kl}, \ \bar{S}_{m,n} := S_{m,n}/mn \text{ and } S^* := \sup_{m,n \geqslant 1} \bar{S}_{m,n}.$ Then (9.40) implies, assuming $t \geqslant t_{a,b} := \max\{a^{-2\gamma}, b^{-2\gamma}\},$

$$\mathbb{E}(V_{a,b,t} \mid (Z_{kl})_{k,l}) \leqslant \sum_{k=\lceil at^{1/2\gamma} \rceil}^{\lceil 2at^{1/2\gamma} \rceil} \sum_{l=\lceil bt^{1/2\gamma} \rceil}^{\lceil 2bt^{1/2\gamma} \rceil} r^2 t^{-1/\gamma} J_{kl} \leqslant r^2 t^{-1/\gamma} S_{\lceil 2at^{1/2\gamma} \rceil, \lceil 2bt^{1/2\gamma} \rceil}
\leqslant r^2 t^{-1/\gamma} \lceil 2at^{1/2\gamma} \rceil \lceil 2bt^{1/2\gamma} \rceil S^* \leqslant 9r^2 abS^*.$$
(9.41)

Fix p > 1 with $p < \gamma$. Then by the multi-dimensional version of Doob's L^p inequality, see [38, Lemma 3], (9.41) implies, for fixed r,

$$\mathbb{E} \sup_{t \geqslant t_{a,b}} \mathbb{E} \big(V_{a,b,t} \mid (Z_{kl})_{k,l} \big) \leqslant C_2 ab \, \mathbb{E} \, S^* \leqslant C_2 ab \big(\mathbb{E} (S^*)^p \big)^{1/p} \leqslant C_3 ab (\mathbb{E} \, I_{11})^{1/p}$$

$$\leqslant C_4 a^{1-\gamma/p} b^{1-\gamma/p}. \tag{9.42}$$

Let $\varepsilon > 0$, and use (9.42) with $a = 2^m \varepsilon$ and $b = 2^n \varepsilon$. Then summing over all $(m, n) \in \mathbb{Z}_+$ with $m \vee n \geqslant N$ implies, using Markov's inequality,

 $\mathbb{E}\sup_{t\geqslant t_{\varepsilon,\varepsilon}}\mathbb{P}\big(\hat{W}_t \text{ has an edge labelled } (x,y)\in [\varepsilon,\infty)^2\setminus [\varepsilon,2^N\varepsilon]^2\mid (Z_{kl})_{k,l}\big)$

$$\leqslant C_5 2^{-(\gamma/p-1)N} \varepsilon^{2(1-\gamma/p)}. \tag{9.43}$$

Choosing N large enough, this is less than ε . Furthermore, the probability that \hat{W}_t has a vertex labelled $< \varepsilon$ is at most $\mathbb{P}(\eta_1 < \varepsilon) < \varepsilon$, and we can choose n such that $\mathbb{P}(\eta_n \leq 2^N \varepsilon) < \varepsilon$.

It now follows from (9.39) that for any finite graph H,

$$\left| \mathbb{P} \left(G(r, \hat{W}_t) = H \mid \tilde{G}_t \right) - \mathbb{P} \left(G(r, W) = H \right) \right| \leqslant 3\varepsilon + o(1) \tag{9.44}$$

a.s. as $t \to \infty$. Since $\varepsilon > 0$ is arbitrary, this shows $(G(r, \hat{W}_t) \mid \tilde{G}_t) \xrightarrow{d} G(r, W)$ a.s. as $t \to \infty$, for every fixed $r < \infty$, which is the same as $\hat{W}_t \to_{\mathsf{GP}} W$.

Finally, we note that $\hat{W}_t \to_{\mathsf{GP}} W$ implies $\hat{W}_t \to_{\mathsf{GS}} W$, see [40; 24], and that \to_{GS} is not affected by dilations of the graphons; hence a.s. also $W_{\tilde{G}_t} \to_{\mathsf{GS}} W$, i.e., $\tilde{G}_t \to_{\mathsf{GS}} W$.

Example 9.7. Consider the simple graphs \tilde{G}_t and G_m given by the Hollywood model in Example 7.8 in the case $0 < \alpha < 1$. As shown there, the resulting random graphs are the same as the ones given by the rank 1 model with a random probability distribution $(q_i)_1^{\infty}$ having the distribution $PD(\alpha, \theta)$, where $\theta > -\alpha$ is the second parameter. This implies that a.s. $q_i \sim Zi^{-1/\alpha}$ for some (random) Z > 0, see [35, Theorem 3.13]. Consequently, Example 9.1 applies with $\gamma = 1/\alpha$ (after conditioning on (q_i)). In particular, a.s. $v(G_m) \approx m^{\alpha}$ and $e(G_m) \approx m^{\alpha} \log m$ as $m \to \infty$.

Moreover, G_m has a.s. a power-law degree distribution with exponent $\tau = 2$ in the sense of Theorem 9.2.

Furthermore, Theorem 9.4 shows that $G_m \to_{\mathsf{GS}} W$ a.s. as $m \to \infty$ and that the dilated empirical graphon converges a.s. in the sense $W_{\tilde{G}_t,t^{\alpha/2}} \to_{\mathsf{GP}} W$, where W is the random graphon $W(x,y) = 1 - \exp\left(-2Z^2x^{-1/\alpha}y^{-1/\alpha}\right)$ on \mathbb{R}_+ .

Problem 9.8. In the simple graph Hollywood model with $0 < \alpha < 1$ as in Example 9.7, does the degree distribution of G_m converge (a.s., or at least in probability) as $m \to \infty$? If so, what is the asymptotic distribution? Is it random or deterministic?

10. Extremely sparse examples

We can obtain extremely sparse examples in several ways.

First, Theorem 6.10 shows that any example including dust or attached stars is extremely sparse.

Another way to obtain extremely sparse graphs is to force the degrees to be bounded, as follows.

Example 10.1. Let $\mu = (\mu_{ij})_{i,j=1}^{\infty}$ be a symmetric non-negative matrix with $0 < \|\mu\| < \infty$ and assume that each row contains at most d non-zero

entries, for some $d < \infty$. (For example, let μ be a band matrix, with $\mu_{ij} = 0$ unless $0 < |i - j| \le d/2$.)

Since an edge ij can exist only when $\mu_{ij} > 0$, it follows that every vertex in G_m has degree at most d. Hence the sequence G_m has bounded degree, and in particular G_m is sparse; to be more precise we have

$$v(G_m) \leqslant 2e(G_m) \leqslant dv(G_m). \tag{10.1}$$

Less obviously, it is also possible to obtain extremely sparse graphs in the rank 1 case, with a sequence q_i that decreases very slowly (remember that $\sum_i q_i = 1$ by assumption). We give one such example.

Example 10.2. Consider the rank 1 case (Section 7.1) with $q_i = c/(i \log^2 i)$ for $i \ge 2$, where c is the appropriate normalization constant. (Any (q_i) with $q_i \approx 1/(i \log^2 i)$ would yield the same results below.) Recall that, by comparison with an integral, $\sum_{i \ge k} 1/(i \log^2 i) \sim 1/\log k$ as $k \to \infty$.

We will see that (in a sense made precise below) almost all edges belong to stars, and that, moreover, most edges and vertices belong to a small (finite) number of stars; in particular, most vertices have degree 1.

For large t, let $\ell(t) := \lfloor t/\log^2 t \rfloor$. Then $\ell(t) \log^2 \ell(t) \sim t$, and thus, using (7.5),

$$v(t) \approx \sum_{i=1}^{\infty} ((q_i t) \wedge 1) \approx \sum_{i \leq \ell(t)} 1 + \sum_{i > \ell(t)} \frac{t}{i \log^2 i} \approx \ell(t) + \frac{t}{\log \ell(t)} \approx \frac{t}{\log t}.$$
 (10.2)

The expected number of edges is by (7.6),

$$e(t) \approx \sum_{i \neq j} \left((q_i q_j t) \wedge 1 \right) \approx \sum_{i \neq j} \frac{t}{i \log^2(i+1) j \log^2(j+1)} \wedge 1 \tag{10.3}$$

We split the sum in (10.3) into three (overlapping) parts. The case $j \ge t^{0.4}$ yields at most

$$\sum_{i \geqslant 1} \sum_{j \geqslant t^{0.4}} \frac{t}{i \log^2(i+1) j \log^2(j+1)} \leqslant C_1 \frac{t}{\log t}.$$
 (10.4)

The case $i \ge t^{0.4}$ yields the same, and finally the case $i, j < t^{0.4}$ yields at most

$$\sum_{i < t^{0.4}} \sum_{j < t^{0.4}} 1 < t^{0.8} = o\left(\frac{t}{\log t}\right). \tag{10.5}$$

By (10.3)–(10.5), and the lower bound (6.24), we find

$$e(t) \approx \frac{t}{\log t} \approx v(t).$$
 (10.6)

Thus Theorem 6.4 yields $e(\tilde{G}_t) \simeq v(\tilde{G}_t)$ and $e(G_m) \simeq v(G_m)$ a.s. In other words, G_m is extremely sparse.

We can be more precise. Recall that $N_i(t)$ is the degree of vertex i in \tilde{G}_t^* ; by (6.3) $N_i(t) \sim \text{Po}(\mu_i t)$. Consequently, using also (7.2),

$$\mathbb{E}(N_i(t)\mathbf{1}_{\{N_i(t)>1\}}) \leqslant \mathbb{E}(N_i(t)(N_i(t)-1)) = (\mu_i t)^2 \leqslant 4q_i^2 t^2.$$
 (10.7)

Summing over $i > t/\log^2 t$ we obtain

$$\sum_{i>t/\log^2 t} \mathbb{E}(N_i(t)\mathbf{1}_{\{N_i(t)>1\}}) \leqslant C_2 \sum_{i>t/\log^2 t} \frac{t^2}{i^2 \log^4 i} = O(\frac{t}{\log^2 t}).$$
 (10.8)

Hence the expected number of edges that have one endpoint in $(t/\log^2 t, \infty)$ and that endpoint is not isolated is $O(t/\log^2 t)$. Moreover, the expected number of edges with both endpoints in $[1, t/\log^2 t]$ is at most

$$\sum_{i < j \le t/\log^2 t} 2((tq_i q_j) \wedge 1) \le t^{0.8} + \sum_{t^{0.4} < j \le t/\log^2 t} \sum_{i < j} 2((tq_i q_j) \wedge 1)$$
 (10.9)

where the last sum is at most a constant times, cf. (10.2),

$$\sum_{j=t^{0.4}}^{t/\log^2 t} v(tq_j) \approx \sum_{j=t^{0.4}}^{t/\log^2 t} \frac{tq_j}{\log(tq_j+2)} \approx \sum_{j=t^{0.4}}^{t/\log^2 t} \frac{t/j\log^2 t}{\log(t/j\log^2 t+2)}$$

$$\approx \int_{x=t^{0.4}}^{t/\log^2 t} \frac{t/\log^2 t}{\log(t/x\log^2 t+2)} \frac{dx}{x} = \int_{y=1}^{t^{0.6/\log^2 t}} \frac{t/\log^2 t}{\log(y+2)} \frac{dy}{y}$$

$$= O\left(\frac{t}{\log^2 t} \log \log t\right) = o\left(\frac{t}{\log t}\right). \tag{10.10}$$

It follows by (10.8), (10.9) and (10.10) that, in \tilde{G}_t^* and thus in \tilde{G}_t , all but $o_p(t/\log t)$ edges have one endpoint isolated. If the number of such edges in \tilde{G}_t is $e'(\tilde{G}_t)$, then thus the total number of edges is $e(\tilde{G}_t) = e'(\tilde{G}_t) + O_p(t/\log^2 t)$, and since each edge has at most two endpoints, the number of vertices is at least $e'(\tilde{G}_t)$ and at most $2e'(\tilde{G}_t) + O_p(t/\log^2 t)$. Moreover, it is easily seen that the expected number of edges with both endpoints in $(t/\log^2 t, \infty)$ is $O(t/\log^2 t)$, and it follows that, in fact, $v(\tilde{G}_t) = e'(\tilde{G}_t) + O_p(t/\log^2 t)$; we omit the details. Consequently, using also (6.26) and (10.6), it follows that $e(\tilde{G}_t)/v(\tilde{G}_t) \stackrel{p}{\longrightarrow} 1$ as $t \to \infty$. Moreover, we see also that almost all edges belong to stars. (These are not attached stars in the sense of Section 5, since our example contains no attached stars, but they have a similar effect on the graph.) As a consequence, at least in probability, most vertices have degree 1, so the asymptotic degree distribution is concentrated at 1.

Furthermore, a large fraction of the edges (and thus vertices) belong to a finite number of such stars. To be precise, let $\varepsilon > 0$; then there exists an integer $K = K(\varepsilon) < \infty$ such that summing over i > K only in (10.4) yields $< \varepsilon t/\log t$, which together with (10.5) and (10.8)–(10.10) shows that the expected number of edges that are not in a star with centre at i for some $i \le K$ is $O(\varepsilon t/\log t) = O(\varepsilon e(t))$.

Since a.s. $G_m \subseteq \tilde{G}_{2m}$ for all large m, the same results follow also for G_m . Unfortunately, these properties make the random graphs in this example rather uninteresting for applications.

11. Conclusions

For the multigraph version, the examples in Section 7 seem very interesting, but perhaps a bit special. We do not know whether they are typical for a large class of interesting examples or not.

For the simple graph version, the examples above show a great variety of different behaviour. Nevertheless, the results are somewhat disappoining for applications; the relations between the intensity matrix (μ_{ij}) and properties of the random graphs G_m such as edge density and degree distribution are far from obvious, and it is not clear how one can choose the intensity matrix to obtain desired properties; for example, we do not know any example of a power-law degree distribution with an exponent $\tau \neq 2$.

Consequently, for both versions, it seems desirable to study more examples, as well as to find more general theorems.

The present paper is only a first step (or rather second step, after [7; 8; 11; 12]), of the investigation of these random graphs, and it seems too early to tell whether they will be useful as random graph models for various applications or not.

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