

The generation of connected random graphs is non-trivial and important to many applications. In particular, it is not easy to sample a connected random graph from the space  $\mathcal{G}$  of all connected graphs with uniformity; i.e. without a bias for graphs of a special kind. This entry contains two algorithms for sampling random connected graphs that do not overcome a special kind of bias. A third algorithm which overcomes all bias, though at great computational expense, is presented.

## 1 Bottom-up approach

Before proceeding, quick definitions.

- Let  $\mathcal{T}_n$  the set of all trees of  $n$  vertices,  $\mathcal{G}_n$  the set of all graphs with  $n$  vertices. We shall assume the vertices of these graphs are labeled  $1, \dots, n$ .
- For any  $T \in \mathcal{T}_n$ , we define  $\mathcal{U}_T := \{G \in \mathcal{G}_n : T \subseteq G\}$  and refer to it as *the universe* of  $T$ .
- Let  $\Gamma(n) = \{\{x, y\} : x, y \in \{1, \dots, n\}\}$ .

In general, whenever I write of a tree  $T$ , I mean an arbitrary  $T \in \mathcal{T}_n$ ; and whenever I write of a graph  $G$ , I mean an arbitrary  $G \in \mathcal{G}_n$ .

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There are several properties of trees which make them attractive for graph generation. On one hand, any finite  $G$  contains a finite number of spanning trees  $T \subseteq G$ ; on the other, any tree spans a finite number of graphs. It is trivial to observe that  $T \subseteq G$  if and only if there is some  $S \in \Gamma(n)$  s.t.  $E(G) = E(T) \cup S$ . We can make use of this special relationship between  $\Gamma(n)$  and  $\mathcal{U}_T$  to produce random connected graphs.

Each tree is perfectly identified by its Prüfer sequence. Then, for a fixed  $n$ , the language  $\{1, \dots, n\}^{n-2}$  indexes a family of functions  $\mathcal{F}$  defined as:

$$\begin{aligned}\mathcal{F}(w) : \mathcal{U}_{T_w} &\rightarrow \Gamma(n) \\ \mathcal{F}(w)(G) &= E(G) - E(T_w)\end{aligned}$$

where  $T_w$  is the tree corresponding to the Prüfer sequence  $w$ . We shall use  $\mathcal{F}_w$  to abbreviate  $\mathcal{F}(w)$ .

Since there is a perfect correspondence between any given  $G \in \mathcal{U}_T$  and the set of edges which, if aggregated to  $T$ , produce  $G$ ,  $\mathcal{F}_w$  is a bijection. (This is easy to

prove.) This entails that any  $G \in \mathcal{G}_n$  which may be spanned from  $T_w$  corresponds to a unique set in  $\Gamma(n)$ —namely, that which extends  $T$  into  $G$ .

Thus, we have established a series of generational relations. A Prüfer sequence maps to a unique tree  $T$ , the tree maps to a universe of connected graphs  $\mathcal{U}_T$ , and each graph in this universe is perfectly determined by a set in  $\Gamma(n)$ . This inspires the following effective procedure for generating random connected graphs of  $n$  vertices.

- (1) Generate randomly  $p = p_1 \dots p_{n-2} \in \Sigma^{n-2}$ .
- (2) Span the tree  $T = (V, E)$  of the Prüfer sequence  $p$ .
- (3) Let  $k \in_R \left\{0, \dots, \frac{n(n-1)}{2}\right\}$ .
- (4) Let  $\ell_1, \dots, \ell_k \in_R \Gamma(n) - E(T)$ , all distinct.
- (5) Let  $E = E \cup \{\ell_1, \dots, \ell_k\}$

Because all trees of  $n$  vertices correspond to a sequence, all trees can be sampled. And all connected graphs can be derived from the set of all spanning trees. Then this procedure generates all graphs in  $\mathcal{G}_n$ .

The question is whether it is equally likely to generate any two graphs in  $\mathcal{G}_n$ . It is obvious that it is equally likely to generate any tree. And the probability that a given graph is generated depends entirely on the number of spanning trees it contains. Not all graphs have the same number of spanning trees.  $\therefore$  It is more likely to generate a graph with many spanning trees than a graph with few spanning trees.

The second algorithm extends the input from only  $n$ , the number of vertices, to  $m$ , the number of edges. Thus, it specifies the problem further into the question of how to generate denser or sparser connected graphs of  $n$  edges. The effective procedure to do this is:

- (1) Generate randomly  $p = p_1 \dots p_{n-2} \in \Sigma^{n-2}$ .
- (2) Span the tree  $T = (V, E)$  of the Prüfer sequence  $p$ .
- (3) If  $V = \{v_1, \dots, v_n\}$ , list all its non-neighbours; i.e. generate  $\Gamma^c(v_1), \dots, \Gamma^c(v_n)$ .
- (4) Sample a random, non-saturated vertex  $v \in V$ , and sample a random vertex  $w$  from  $\Gamma^c(v)$ .
- (5) Add  $\{v, w\}$  to  $E$ . Remove  $v$  from the list of non-neighbours of  $w$ , and  $w$  from the list of non-neighbours of  $v$ .

(6) If  $|E| = m$ , finish. Otherwise go to 4.

It is clear that the procedure always finishes, and since the tree is connected the generated graph is connected. A pseudo-code algorithm may look as follows.

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Input:  $n, m$ 
 $(V, E) = \text{gen\_random\_tree}(n)$ 
 $S = [ \Gamma^c(v_1), \dots, \Gamma^c(v_n) ]$                                 {Non-neighbours}
 $C = [ |S[1]|, \dots, |S[n]| ]$                                 {Cardinalities}
 $V = [1, \dots, n]$                                 {Non-saturated vertices }
 $n' = n$ 
while  $|E(T)| < m$  do
   $i := \text{random}(1, n')$ 
   $v := V[i]$ 
  if  $(d(v) == n - 1)$  do
    delete_at $(V, i)$ 
     $n' := n' - 1$ 
  else
     $j = \text{random}(1, C[v])$ 
     $w := S[v][j]$ 
     $E(T) := E(T) \cup \{v, w\}$ 
     $C[v] := C[v] - 1$ 
    delete_at $(S[v], j)$ 
    delete_element $(S[w], v)$ 
  fi
od
return  $T$ 

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Generating a tree from a random Prüfer sequence is  $O(n^2)$ . Listing all the non-neighbours is also  $O(n^2)$ . Within the while loop there are simply index manipulations, so the complexity of the loop is  $\varphi \times O(1) = \varphi$ , with  $\varphi$  the complexity of the number of iterations.

All iterations add an edge except those where a saturated vertex is chosen. A saturated vertex may be chosen at most once.  $\therefore$  There are  $O(n)$  iterations where a saturated vertex is chosen. Since the rest of the iterations add an edge, their

number is fixed:  $m - (n - 1)$ , where  $(n - 1)$  is the number of edges in the spanned tree.  $\therefore$  There are exactly  $m - n + 1$  edge-adding iterations.

$$\therefore \varphi = O(n) + O(m - n + 1) = O(n) + O(m) = O(n^2)$$

$$\therefore \text{The complexity of the algorithm is } O(n^2) + O(n^2) = O(n^2).$$