HOPPING FORCING NUMBER IN RANDOM *d*-REGULAR GRAPHS

PAWEŁ PRAŁAT AND HARJAS SINGH

ABSTRACT. Hopping forcing is a single player combinatorial game in which the player is presented a graph on n vertices, some of which are initially blue with the remaining vertices being white. In each round t, a blue vertex v with all neighbours blue may hop and colour a white vertex blue in the second neighbourhood, provided that v has not performed a hop in the previous t-1 rounds. The objective of the game is to eventually colour every vertex blue by repeatedly applying the hopping forcing rule. Subsequently, for a given graph G, the hopping forcing number is the minimum number of initial blue vertices that are required to achieve the objective.

In this paper, we study the hopping forcing number for random *d*-regular graphs. Specifically, we aim to derive asymptotic upper and lower bounds for the hopping forcing number for various values of $d \ge 2$.

1. INTRODUCTION AND MAIN RESULTS

1.1. **Definitions.** Zero forcing was introduced in [16], with the purpose of finding bounds on the maximum nullity of a family of matrices associated with any graph. The game starts with a graph G with some vertices coloured **blue** and the remaining **white**. An interpretation of this is that blue vertices contain information, whereas the white vertices are devoid of it. The goal is to use repeated applications of a **colour change rule** with the objective being to (eventually) turn every vertex blue. It quickly gained popularity as a model to study the spread of information on a given graph, with applications being found in several fields including physics and engineering; see, for example [10, 11, 23], survey on the subject [14], and a recent article published in the Notices of the AMS [17].

The standard colour change rule, denoted by \mathcal{Z} , allows a blue vertex b to force a white vertex w to become blue provided that w is the unique white neighbour of b. The wide interest in zero forcing has generated a large volume of work in the last few years, including analyzing it for random graphs [4], analyzing the probabilistic counterpart for deterministic graphs [6] as well as random ones [13]. On the other hand, the **hopping forcing rule** \mathcal{H} , first added to the set of existing colour change rules in [5], allows a blue vertex b to force a white vertex w in the second neighbourhood of b to become blue provided that b has not performed a force yet and each vertex in the first neighbourhood of b is already blue. The new rule was added as a tool in investigations about relations between various zero forcing concepts. The relationships of this graph parameter and other ones, such as vertex connectivity and independence number, as well as other zero forcing parameters were investigated in [12]. For a given graph G, the **hopping forcing number** H(G) denotes the cardinality of the set with the minimum number of initial blue vertices, such that all the vertices of G can eventually be coloured blue.

We borrow the notation used in [12] and call a blue vertex b at **step** t **dormant** if it has neither performed a force yet nor is able to (the reason being that either some neighbour of b is still white or all vertices in the second neighbourhood of b are already blue), **active** if it has not yet performed a force but is able to, or **extinct** if it has already performed a force.

Furthermore, it is important to note that although several vertices can hop simultaneously at step t, for our purpose we only consider **sequential** hopping, that is, at any given step there will be exactly one vertex performing a force, and if no vertex can hop but there are still some vertices that are white, then the process terminates unsuccessfully.

Now, we set up our set notation for this process and carefully lay out the steps. Let B_1 be the set of vertices that are initially blue. At the beginning (that is, before the player makes her move) of time step $t \geq 1$, let B_t denote the set of all blue vertices (either active, dormant or extinct), $R_t \subseteq B_t$ denote the set of vertices that are active (blue vertices that have not yet performed a force whose first neighbourhood is entirely blue and that have at least one white vertex in the second neighbourhood), and W_t denote the set of white vertices. Then, at each sequential hop, the player selects a vertex $x_t \in R_t$ and performs a single force by hopping from x_t to $y_t \in N_2(x_t) \cap W_t$, where $N_2(x_t)$ denotes the second neighbourhood of x_t , that is, the set of vertices that are at distance exactly two from x_t . Subsequently, the sets B_t , R_t , and W_t are updated accordingly. In particular,

$$B_{t+1} = B_t \cup \{y_t\}$$
$$W_{t+1} = W_t \setminus \{y_t\}.$$

Clearly, vertex x_t needs to be removed from R_t since it became extinct but more updates might be necessary as potentially more vertices change their status from active to dormant or vice versa.

Note that a successful **strategy** S can be defined as an initial set B_1 of blue vertices and a sequence of ℓ hops, $x_1 \to y_1, x_2 \to y_2, \ldots, x_\ell \to y_\ell$ where, at the end of the process, every vertex is blue. Such initial sets B_1 will be called **feasible**. Clearly, $B_1 = V(G)$ is feasible which shows that the hopping forcing number is well defined and that $H(G) \leq |V(G)|$. On the other hand, $H(G) \geq 1$ since $B_1 = \emptyset$ is infeasible (unless G is the null graph, that is, $V(G) = \emptyset$ in which case, trivially, H(G) = 0). Furthermore, out of several feasible sets B_1^1, B_1^2, \ldots , the set(s) with least cardinality can be called **optimal** and this cardinality is the hopping number H(G). Observe that given a graph G and B_1 , both feasibility and optimality can be ascertained at the beginning of the process, that is, this is a 'one person game' with perfect information and no randomness. Finally, since the optimal sequential hopping starts with |V(G)| - H(G) white vertices and at each step one of them changes colour to blue, the length of the process is equal to |V(G)| - H(G).

More importantly, the set of initial blue vertices, B_1 , can be updated in an online fashion, that is, we may try to perform the desired sequence of hops and append some vertices to the initial set dynamically instead of knowing the set B_1 a priori. Indeed, we may start with $B_1 = \emptyset$. Then, at time t, if x_t is not extinct and y_t is white and in $N_2(x_t)$, then we may add x_t (if needed) and all white neighbours of x_t (if there are any) to B_1 so that the desired force can be performed.

1.2. Main Results. In this paper, we establish various asymptotic upper and lower bounds for the hopping number of the random *d*-regular graph $\mathcal{G}_{n,d}$ (see Subsection 2.2 for the definition and more details on this model). We say that a random graph has property *P* asymptotically almost surely (or a.a.s.) if the probability that it has property *P* tends to 1 as *n* goes to infinity (see Subsection 2.1 for more on asymptotic notation used in this paper).

Understanding random 2-regular graphs is easy. In Section 3, we prove the following result.

Theorem 1.1. A.a.s. $H(\mathcal{G}_{n,2}) \sim (3/2) \log n$.

Unfortunately, the d = 2 case is the only value of $d \ge 2$ for which we determine an asymptotic behaviour of the hopping number. For the remaining ones, we only have some upper and lower bounds—see Table 1 and Figure 1.1.

Upper bounds are studied in Section 4. We use an (on-line) algorithm to create the initial set of blue vertices, yielding the following upper bounds:

Theorem 1.2. For any integer $d \ge 3$, a.a.s.

$$H(\mathcal{G}_{n,d}) \le (1+o(1)) \frac{(d-1)!(d-2)^{d-1}}{\prod_{i=1}^{d-1} (i(d-2)+1)} n.$$

	Lower]	Upper Bound	
Degree	Expander Mixing Lemma	Configuration Model	Contiguous Model
d = 3	0.0149	0.0699	0.3333
d = 4	0.0372	0.1451	0.4571
d = 5	0.0588	0.2114	0.5341
d = 6	0.0787	0.2678	0.5884
d = 7	0.0968	0.3158	0.6294
d = 8	0.1134	0.3569	0.6618
d = 9	0.1287	0.3924	0.6882
d = 10	0.1429	0.4235	0.7101
d = 20	0.2445	0.6054	0.8231
d = 40	0.3755	0.7437	0.8946
d = 80	0.5556	0.8409	0.9386
d = 160	0.6848	0.9048	0.9649
d = 320	0.7767	0.9446	0.9803
d = 640	0.8420	0.9684	0.9890
d = 1280	0.8882	0.9823	0.9940

TABLE 1. Comparison of upper and lower bounds for the hopping number for small and large values of d.



FIGURE 1.1. Comparison of upper and lower bounds for the hopping number for small and large values of d.

The algorithm used in the argument is universal but the analysis is slightly different in the degenerate case d = 3. Instead of investigating two cases in one proof, we prepare the reader for

more complicated argument and first prove that a.a.s. $H(\mathcal{G}_{n,3}) \leq (1+o(1))n/3$ (see Subsection 4.1) before analyzing the algorithm for $d \geq 4$ (see Subsection 4.2).

Moreover, note that the constant in the upper bound can be estimated as follows:

$$\frac{(d-1)!(d-2)^{d-1}}{\prod_{i=1}^{d-1}(i(d-2)+1)} = \prod_{i=1}^{d-1} \frac{i(d-2)}{i(d-2)+1} = \prod_{i=1}^{d-1} \left(1 - \frac{1}{i(d-2)+1}\right)$$
$$= \exp\left(-\Theta\left(\sum_{i=1}^{d-1} \frac{1}{id}\right)\right) = \exp\left(-\Theta\left(\frac{\log d}{d}\right)\right)$$
$$= 1 - \Theta\left(\frac{\log d}{d}\right).$$

In particular, it shows that it tends to one as $d \to \infty$.

To get a lower bound for $H(\mathcal{G}_{n,d})$ that explicitly tends to one as $d \to \infty$, we use the expansion properties of random *d*-regular graphs to get a lower bound of $1 - \Theta(1/\sqrt{d})$. Indeed, in Section 5, we prove the following result. (See Subsection 2.4 for the definition of $\lambda(G)$.)

Theorem 1.3. Let G = (V, E) be a d-regular graph with n vertices and set $\lambda = \lambda(G)$. Then,

$$H(G) \ge \max\left(1 - \frac{2\lambda}{d}, \frac{d - \lambda}{d + 3\lambda}\right) n = \left(1 - \min\left(\frac{2\lambda}{d}, \frac{4\lambda}{d + 3\lambda}\right)\right) n.$$

As a result, for any $d \ge 3$ and $\varepsilon > 0$, a.a.s.

$$H(\mathcal{G}_{n,d}) \ge \left(1 - \min\left(\frac{4\sqrt{d-1}}{d}, \frac{8\sqrt{d-1}}{d+6\sqrt{d-1}}\right) - \varepsilon\right)n.$$

In Section 6, the above lower bound is strengthened by applying the configuration model to get the following, stronger but implicit and numerical, lower bound.

Theorem 1.4. For a given integer $d \ge 3$, let

$$g_d(x,z) = \left(\frac{d}{2} - 1 - dz\right) x \log(x) + (d-1)(1-x) \log\left(\frac{1-x}{2}\right) - 2dxz \log(z) - \frac{(1-2z)dx}{2} \log(1-2z) - d\left(\frac{1-x}{2} - zx\right) \log\left(\frac{1-x}{2} - zx\right).$$

For a fixed $x \in (0,1)$, function $g_d(x,z)$ is maximized at

$$z_0(x) := \frac{1 - \sqrt{1 - 2(1 - x)x}}{2x}.$$

Fix $\varepsilon > 0$. Let x_d be the unique $x \in (0, 1)$ for which $h_d(x) = g_d(x, z_0(x)) = 0$. Then, a.a.s. $H(\mathcal{G}_{n,d}) > (x_d - \varepsilon)n$.

Next, we report our attempt to get better upper bounds for the hopping number by introducing a degree-greedy algorithm [28] to create the initial set of blue vertices and then use the differential equation method to analyze it. Unfortunately, the bounds we obtained using this method turned out to be weaker than the ones we established above. Nevertheless, in Section 7 we briefly report our attempt for the case of random 3-regular graphs with the hope that one can modify our algorithm, and use similar techniques to analyze it, to get better bounds than the ones we managed to prove. We finish the paper with a few natural suggestions for future directions (see Section 8).

2. Preliminaries

2.1. Notation. The results presented in this paper are asymptotic by nature. We say that a random graph has property P asymptotically almost surely (or a.a.s.) if the probability that it has property P tends to 1 as n goes to infinity. Given two functions f = f(n) and g = g(n), we will write $f(n) = \mathcal{O}(g(n))$ if there exists an absolute constant $c \in \mathcal{R}_+$ such that $|f(n)| \leq c|g(n)|$ for all n, $f(n) = \Omega(g(n))$ if $g(n) = \mathcal{O}(f(n))$, $f(n) = \Theta(g(n))$ if $f(n) = \mathcal{O}(g(n))$ and $f(n) = \Omega(g(n))$, and we write f(n) = o(g(n)) or $f(n) \ll g(n)$ if $\lim_{n \to \infty} f(n)/g(n) = 0$. In addition, we write $f(n) \gg g(n)$ if g(n) = o(f(n)) and we write $f(n) \sim g(n)$ if f(n) = (1+o(1))g(n), that is, $\lim_{n\to\infty} f(n)/g(n) = 1$.

We will use $\log n$ to denote the natural logarithm of n. For a given $n \in \mathbb{N} := \{1, 2, \ldots\}$, we will use [n] to denote the set consisting of the first n natural numbers, that is, $[n] := \{1, 2, \ldots, n\}$. Finally, as typical in the field of random graphs, for expressions that clearly have to be an integer, we round up or down but do not specify which: the choice of which does not affect the argument.

2.2. Random *d*-regular Graphs. Our main results refer to the probability space of random *d*-regular graphs with uniform probability distribution. This space is denoted $\mathcal{G}_{n,d}$, and asymptotics are for $n \to \infty$ with $d \ge 2$ fixed, and *n* even if *d* is odd.

Instead of working directly in the uniform probability space of random regular graphs on n vertices $\mathcal{G}_{n,d}$, we use the *configuration model* of random regular graphs, first introduced by Bollobás [9], which is described next. Suppose that dn is even, as in the case of random regular graphs, and consider dn points partitioned into n labeled buckets v_1, v_2, \ldots, v_n of d points each. A *pairing* of these points is a perfect matching into dn/2 pairs. Given a pairing P, we may construct a multigraph G(P), with loops and parallel edges allowed, as follows: the vertices are the buckets v_1, v_2, \ldots, v_n , and a pair $\{x, y\}$ in P corresponds to an edge $v_i v_j$ in G(P) if x and y are contained in the buckets v_i and v_j , respectively.

It is an easy fact that the probability of a random pairing corresponding to a given simple graph G is independent of the graph, hence the restriction of the probability space of random pairings to simple graphs is precisely $\mathcal{G}_{n,d}$. Moreover, it is well known that a random pairing generates a simple graph with probability asymptotic to $e^{(1-d^2)/4}$ depending on d, so that any event holding a.a.s. over the probability space of random pairings also holds a.a.s. over the corresponding space $\mathcal{G}_{n,d}$. For this reason, asymptotic results over random pairings suffice for our purposes. One of the advantages of using this model is that the pairs may be chosen sequentially so that the next pair is chosen uniformly at random over the remaining (unchosen) points. For more information on this model, see the survey [29] or any of the books on random graphs [8, 19, 20].

2.3. Contiguous Model. The notion of the union of two random regular graphs on the same vertex set is very useful for proving asymptotic properties of $\mathcal{G}_{n,d}$ with $d \geq 3$. In particular, it is known that, for the purpose of proving statements a.a.s., such a random graph can be viewed as the multigraph formed from the union of a Hamilton cycle and random (d-2)-regular graph on the same vertex set; see [29, Theorem 4.15] for a stronger and more general result. (The probability of multiple edges being created is bounded away from 1, and the resulting graph, conditional upon no multiple edges, is contiguous to a random d-regular graph.)

2.4. Expansion Properties of Random *d*-regular Graphs. We will use the expansion properties of random *d*-regular graphs that follow from their eigenvalues. The adjacency matrix A = A(G)of a given a *d*-regular graph *G* with *n* vertices, is an $n \times n$ real and symmetric matrix. Thus, the matrix *A* has *n* real eigenvalues which we denote by $\lambda_1 \geq \lambda_2 \geq \cdots \geq \lambda_n$. It is known that certain properties of a *d*-regular graph are reflected in its spectrum but, since we focus on expansion properties, we are particularly interested in the following quantity: $\lambda = \lambda(G) = \max(|\lambda_2|, |\lambda_n|)$. In words, λ is the largest absolute value of an eigenvalue other than $\lambda_1 = d$. For more details, see the general survey [18] about expanders, or [3, Chapter 9]. The value of λ for random *d*-regular graphs has been studied extensively. A major result due to Friedman [15] is the following:

Lemma 2.1 ([15]). For every fixed $\varepsilon > 0$ and for $G \in \mathcal{G}_{n,d}$, a.a.s. $\lambda(G) < 2\sqrt{d-1} + \varepsilon$.

The number of edges |E(S,T)| between sets S and T is expected to be close to the expected number of edges between S and T in a random graph of edge density d/n, namely, d|S||T|/n. A small λ (or large spectral gap) implies that this deviation is small. The following useful bound is essentially proved in [1] (see also [3]):

Lemma 2.2 (Expander Mixing Lemma). Let G = (V, E) be a d-regular graph with n vertices and set $\lambda = \lambda(G)$. Then for all $S, T \subseteq V$,

$$\left| |E(S,T)| - \frac{d|S||T|}{n} \right| \le \lambda \sqrt{|S||T|} \,.$$

(Note that $S \cap T$ does not have to be empty; in general, |E(S,T)| is defined to be the number of edges between $S \setminus T$ and T plus twice the number of edges that contain only vertices of $S \cap T$.)

At some point it will be better to apply stronger estimates for $|E(S, V \setminus S)|$ that can be easily derived from a slightly stronger version of the above lemma for |E(S, S)| (see [1]), namely,

$$\left| |E(S,S)| - \frac{d|S|^2}{n} \right| \le \frac{\lambda |S| |V \setminus S|}{n}$$

for all $S \subseteq V$. Since $|E(S, V \setminus S)| = d|S| - |E(S, S)|$, we immediately get that

$$\left| |E(S, V \setminus S)| - \frac{d|S||V \setminus S|}{n} \right| \le \frac{\lambda|S||V \setminus S|}{n}$$
(1)

for all $S \subseteq V$.

2.5. The Differential Equation Method. In this paper, we will use the differential equation method (see [7] for a gentle introduction) to establish dynamic concentration of our random variables. The origin of the differential equation method stems from work done at least as early as 1970 (see Kurtz [21]), and which was developed into a very general tool by Wormald [26, 27] in the 1990's. Indeed, Wormald proved a "black box" theorem, which gives dynamic concentration so long as some relatively simple conditions hold. Warnke [24] recently gave a short proof of a somewhat stronger black box theorem.

In fact, we will use another general result of Wormald [28] on the average-case performance of certain greedy algorithms. Its proof uses the differential equation method and the result applies to algorithms in which the possible operations performed at each step are prioritised, including heuristic algorithms for finding large subgraphs with special properties in random regular graphs, such as maximum independent sets and minimum dominating sets. This general approach eliminates some of the complications caused by prioritisation. Since stating this general purpose algorithm would take some space and effort, instead we will simply comment in Section 7 on how the result is applied to our scenario when the algorithm and associated random variables that we aim to estimate are introduced.

3. 2-regular Graphs

Let Y = Y(n) be the total number of cycles in a random 2-regular graph on n vertices. Since exactly three vertices need to be initially blue in each cycle (that is, $H(C_i) = 3$ for any $i \ge 3$), $H(\mathcal{G}_{n,2}) = 3Y(n)$.

We know that the random 2-regular graph is a.a.s. disconnected; by simple calculations one can show that the probability of having a Hamiltonian cycle is asymptotic to $\frac{1}{2}e^{3/4}\sqrt{\pi/n} = o(1)$ (see, for

example, [29]). We also know that the total number of cycles Y(n) is sharply concentrated around $(1/2) \log n$. Indeed, it is not difficult to see this by generating the random graph sequentially using the pairing model. The probability of forming a cycle in step *i* is exactly 1/(2n - 2i + 1), so the expected number of cycles is

$$\sum_{i=1}^{n} \frac{1}{2n-2i+1} = \sum_{i=1}^{2n} \frac{1}{i} - \frac{1}{2} \sum_{i=1}^{n} \frac{1}{i} = \log(2n) - \frac{1}{2} \log n + O(1) = \frac{1}{2} \log n + O(1).$$

The variance can be calculated in a similar way. So we get the following result (Theorem 1.1):

A.a.s. $H(\mathcal{G}_{n,2}) \sim (3/2) \log n$.

4. Upper Bounds from the Contiguous Model: *d*-regular Graphs, $d \ge 3$

To provide an upper bound for $H(\mathcal{G}_{n,d})$ for 3-regular graphs, and subsequently for d-regular graphs with $d \ge 4$, we will use the contiguous model introduced in Subsection 2.3. For a given $d \ge 3$, a d-regular graph (generated by the pairing model) can be viewed as the union of a Hamilton cycle (v_1, v_2, \ldots, v_n) and random (d-2)-regular graph on the same vertex set, namely, $\{v_1, v_2, \ldots, v_n\}$. We will call the two neighbours of v_k that are on the Hamilton cycle HC-neighbours; the remaining d-2 neighbours of v_k will be called RG-neighbours.

As explained at the end of Subsection 1.1, it will be easier to update B_1 , the set of initial blue vertices in an online fashion. The strategy will be the same for all $d \ge 3$ but formulas for d = 3, the degenerate case, will be slightly different. Therefore, once we explain the general strategy we will independently deal with the d = 3 case (Subsection 4.1) before moving to the $d \ge 4$ case (Subsection 4.2).

Start with $B_1 = \emptyset$. We attempt to colour vertices blue as we hop along the Hamilton cycle. We start by turning v_1 and all of its neighbours blue; these vertices are added to B_1 . We will try to make v_1 to hop to some neighbour of v_2 . (Of course, v_1 can try to hop through some other neighbour, not necessarily through v_2 , but insisting on this choice will make the analysis of the strategy tractable. But this strategy is certainly suboptimal.) If v_3 is white, then v_1 can hop there and force v_3 to become blue (note that v_3 could be a neighbour of v_1 and so could be blue). Similarly, if any of the (d-2) RG-neighbours of v_2 are white, then v_1 can hop and force one of them to become blue. Note that if v_1 could not hop through v_2 , then all neighbours of v_2 are already blue. Otherwise, we turn the remaining white neighbours of v_2 (if there are any) blue; these vertices are added to B_1 . After that we will try to make v_2 to hop through v_3 to some neighbour of v_3 , and continue hoping along the Hamilton cycle. Once we investigate v_{n-3} , the strategy is finished and we can check how many vertices were added to B_1 during this process. (Note that v_n is a HC-neighbour of v_1 and so when we reach v_{n-2} , all vertices are certainly blue.) By design, the set B_1 that is constructed during this process is feasible and so its size yields the desired upper bound for $H(\mathcal{G}_{n,d})$.

Before we describe the situation when v_t tries to hop through v_{t+1} , let us make a simple observation that will simplify the analysis of the above process. Let X be the random variable counting how many vertices could not hop during the process; clearly, Y = n - X is the number of vertices that hopped. Since each time a vertex hops exactly one white vertex turns blue, $|B_1| + Y = n$ which is equivalent to

$$|B_1| = n - Y = X.$$

Hence, if we prove that a.a.s. $X \sim f(n)$ for some deterministic function f(n), then we may conclude that a.a.s. $H(\mathcal{G}_{n,d}) \leq (1+o(1))f(n)$.

For each $t \in [n-3]$, let X_t be the indicator random variable for the event that v_t cannot hop through v_{t+1} . Clearly, $X = \sum_{t=1}^{n-3} X_t$. When v_t tries to hop through v_{t+1} , vertices v_1, v_2, \ldots, v_t and all of their neighbours are blue (in particular, v_{t+1} is blue); the remaining vertices are white. Vertex v_t cannot hop (that is, $X_t = 1$) if and only if the following two properties hold:

- (P1) at least one RG-neighbour of v_{t+2} is in $\{v_1, v_2, \ldots, v_t\}$ (that is, v_{t+2} is blue so v_t cannot hop there),
- (P2) for some $i \in \{0, 1, \dots, d-2\}$,
 - (P2') i RG-neighbours of v_{t+1} are in $\{v_n, v_1, v_2, \ldots, v_t, v_{t+2}\}$, and
 - (P2") (d-2-i) RG-neighbours of v_{t+1} are in $\{v_{t+3}, v_{t+4}, \dots, v_{n-1}\}$ but all of these (d-2-i)RG-neighbours have at least one RG-neighbour in $\{v_1, v_2, \dots, v_t\}$ (that is all of the (d-2) RC neighbours of v_{t+1} are blue so v_t cannot hen to any of these

(that is, all of the (d-2) RG-neighbours of v_{t+1} are blue so v_t cannot hop to any of these vertices).

4.1. d = 3 case. The case d = 3 is the degenerate case (slightly different and easier to analyze), and we will deal with it independently. We will prove the following (Theorem 1.2 for d = 3):

A.a.s.
$$H(\mathcal{G}_{n,3}) \le (1+o(1))n/3$$
.

Proof of Theorem 1.2 for d = 3. Fix any $t \in [n-3]$. To compute the probability that $X_t = 1$, we first expose the unique RG-neighbour of v_{t+2} . Property (P1) holds with probability t/(n-1). Conditioning on this event, we expose the RG-neighbour of v_{t+1} to determine whether property (P2) holds or not. An important observation is that if the unique RG-neighbour of v_{t+1} is in $\{v_{t+3}, v_{t+4}, \ldots, v_{n-1}\}$, then the property (P2") cannot hold—this RG-neighbour has only one RG-neighbour, namely, v_{t+1} which is not in $\{v_1, v_2, \ldots, v_t\}$. In other words, the only chance that property (P2) holds is when i = 1. This makes the case d = 3 degenerate and distinguishes it from the case $d \ge 4$. The conditional probability that (P2) holds is then equal to t/(n-3). We get that

$$\mathbb{P}(X_t = 1) = \frac{t}{n-1} \cdot \frac{t}{n-3}$$

and so

$$\mathbb{E}[X] = \sum_{t=1}^{n-3} \mathbb{P}(X_t = 1) = \sum_{t=1}^{n-3} \frac{t^2}{(n-1)(n-3)}$$
$$= \frac{1}{(n-1)(n-3)} \cdot \frac{(n-3)(n-4)(2n-5)}{6} = \frac{n}{3} + \mathcal{O}(1).$$
(2)

It remains to show that X is well-concentrated around its expectation. We demonstrate this by estimating the variance. First, note that

$$\operatorname{Var}[X] = 2 \sum_{1 \le k < \ell \le n-3} \left(\mathbb{P}(X_k = X_\ell = 1) - \mathbb{P}(X_k = 1) \mathbb{P}(X_\ell = 1) \right) + \sum_{k=1}^{n-3} \left(\mathbb{P}(X_k = 1) - \mathbb{P}(X_k = 1)^2 \right).$$

Clearly, the second term is $\mathcal{O}(n)$. Moreover, the first term can be split further depending on whether $\ell = k + 1$ or $\ell \geq k + 2$. If $\ell = k + 1$, then each term is trivially at most one, thus the corresponding sum is again $\mathcal{O}(n)$. It follows that

$$\begin{aligned} \mathbb{V}\mathrm{ar}[X] &= \mathcal{O}(n) + 2\sum_{\substack{1 \le k < \ell \le n-3\\ \ell \ge k+2}} \left(\mathbb{P}(X_k = X_\ell = 1) - \mathbb{P}(X_k = 1)\mathbb{P}(X_\ell = 1) \right) \\ &= \mathcal{O}(n) + 2\sum_{\substack{1 \le k < \ell \le n-3\\ \ell \ge k+2}} \left(\frac{k^2}{(n-1)(n-3)} \cdot \frac{(\ell + \mathcal{O}(1))^2}{(n-5)(n-7)} - \frac{k^2}{(n-1)(n-3)} \cdot \frac{\ell^2}{(n-1)(n-3)} \right) \\ &= \mathcal{O}(n) + \sum_{k=1}^{n-5} \mathcal{O}\left(\frac{k^2}{n^2}\right) \sum_{\ell=k+2}^{n-3} \mathcal{O}\left(\frac{\ell}{n^2}\right) \end{aligned}$$

$$= \mathcal{O}(n) + \sum_{k=1}^{n-5} \mathcal{O}\left(\frac{k^2}{n^2}\right) = \mathcal{O}(n) = o(n^2)$$

Since $\mathbb{V}ar[X] = o((\mathbb{E}[X])^2)$, X is well-concentrated around its expectation by the second moment method. The proof of the theorem is finished.

4.2. $d \ge 4$ case. The bound from Theorem 1.2 is replicated here for easier reference: for any integer $d \ge 3$, a.a.s.

$$H(\mathcal{G}_{n,d}) \le (1+o(1)) \,\frac{(d-1)!(d-2)^{d-1}}{\prod_{i=1}^{d-1} (i(d-2)+1)} \,n. \tag{3}$$

The exact values for $3 \le d \le 10$ are presented in Table 2. Approximated values can be found in Table 1.

Degree	Upper Bound	Degree	Upper Bound
d = 3	1/3	d = 7	78125/124124
d = 4	16/35	d = 8	40310784/60911435
d = 5	243/455	d = 9	40353607/58640175
d = 6	8192/13923	d = 10	17179869184/24192643475

TABLE 2. Explicit upper bounds for the hopping number for small values of d.

The case of $d \ge 4$ follows a similar structure to the case where d = 3. The main distinction here is that, since there are always at least two RG-neighbours in addition to the two HC-neighbours on the cycle for $d \ge 4$, a vertex could be a RG-neighbour of more than one vertex in $\{v_1, v_2, \ldots, v_t\}$ (see Property (P1)). Similarly, v_{t+1} could have (d - 2 - i) RG-neighbours in $\{v_{t+3}, v_{t+4}, \ldots, v_{n-1}\}$ and all of these RG-neighbours can have at least one RG-neighbour in $\{v_1, v_2, \ldots, v_t\}$ for all values of $i \in \{0, 1, \ldots, d-2\}$. In other words, the Property (P2) holds with a non-zero probability even when i = 0.

Proof of Theorem 1.2 for $d \ge 4$. Fix any $t \in [n-3]$. To compute the probability that $X_t = 1$, we first expose the (d-2) RG-neighbours of v_{t+2} , one by one. Property (P1) holds with probability

$$1 - \prod_{i=1}^{d-2} \left(1 - \frac{(d-2)t}{(d-2)n - 2i + 1} \right) + \mathcal{O}(n^{-1}) = 1 - \left(1 - \frac{t}{n} \right)^{d-2} + \mathcal{O}(n^{-1}).$$
(4)

(The first $\mathcal{O}(n^{-1})$ term corresponds to the event that there is a loop at v_{t+2} .) Conditioning on Property (P1), we expose the RG-neighbours of v_{t+1} to determine whether property (P2) holds or not. The conditional probability that (P2) holds is equal to

$$\mathcal{O}(n^{-1}) + \sum_{i=0}^{d-2} \binom{d-2}{i} \left(\frac{(d-2)t + \mathcal{O}(1)}{(d-2)n + \mathcal{O}(1)} \right)^{i} \left(1 - \frac{(d-2)t + \mathcal{O}(1)}{(d-2)n + \mathcal{O}(1)} \right)^{d-2-i} \\ \cdot \left(1 - \left(1 - \frac{(d-2)t + \mathcal{O}(1)}{(d-2)n + \mathcal{O}(1)} \right)^{d-3} \right)^{d-2-i} \\ = \mathcal{O}(n^{-1}) + \sum_{i=0}^{d-2} \binom{d-2}{i} \left(\frac{t}{n} \right)^{i} \left(1 - \frac{t}{n} \right)^{d-2-i} \left(1 - \left(1 - \frac{t}{n} \right)^{d-3} \right)^{d-2-i}.$$
(5)

We get that

$$\mathbb{P}(X_t = 1) = \mathcal{O}(n^{-1}) + \left(1 - \left(1 - \frac{t}{n}\right)^{d-2}\right)$$
$$\cdot \sum_{i=0}^{d-2} {\binom{d-2}{i}} \left(\frac{t}{n}\right)^i \left(1 - \frac{t}{n}\right)^{d-2-i} \left(1 - \left(1 - \frac{t}{n}\right)^{d-3}\right)^{d-2-i}$$
$$= \mathcal{O}(n^{-1}) + \left(1 - \left(1 - \frac{t}{n}\right)^{d-2}\right) \cdot \left(\frac{t}{n} + \left(1 - \frac{t}{n}\right) - \left(1 - \frac{t}{n}\right)^{d-2}\right)^{d-2}$$
$$= \mathcal{O}(n^{-1}) + \left(1 - \left(1 - \frac{t}{n}\right)^{d-2}\right)^{d-1}$$

and so

$$\mathbb{E}[X] = \sum_{t=1}^{n-3} \mathbb{P}(X_t = 1) = \mathcal{O}(1) + \sum_{t=1}^{n-3} \left(1 - \left(1 - \frac{t}{n}\right)^{d-2} \right)^{d-1}$$
$$= \mathcal{O}(1) + n \int_0^1 (1 - (1 - x)^{d-2})^{d-1} dx$$
$$= \mathcal{O}(1) + n \frac{(d-1)!(d-2)^{d-1}}{\prod_{i=1}^{d-1} (i(d-2) + 1)},$$
(6)

where the last integral follows via a recursive reduction approach. As with the case where d = 3, we would like to show that X is well-concentrated around its expectation by estimating the variance. First, note that

$$\mathbb{V}ar[X] = 2 \sum_{1 \le k < \ell \le n-3} \left(\mathbb{P}(X_k = X_\ell = 1) - \mathbb{P}(X_k = 1) \mathbb{P}(X_\ell = 1) \right) + \sum_{k=1}^{n-3} \left(\mathbb{P}(X_k = 1) - \mathbb{P}(X_k = 1)^2 \right).$$

Clearly, the second term is $\mathcal{O}(n)$. As in the proof for the d = 3 case, the first term can be split further depending on whether $\ell = k + 1$ or $\ell \ge k + 2$. If $\ell = k + 1$, then each term is trivially at most one, thus the sum is again $\mathcal{O}(n)$. To estimate $\mathbb{P}(X_k = X_\ell = 1)$ we first expose the (d - 2)RG-neighbours of v_{k+2} and compute the probability of Property (P1) holding for X_k as in (4). Conditioning on that, Property (P1) holds for X_ℓ with probability

$$1 - \prod_{i=1}^{d-2} \left(1 - \frac{(d-2)\ell + \mathcal{O}(1)}{(d-2)n - 2i - 2(d-2) + 1} \right) + \mathcal{O}(n^{-1}) = 1 - \left(1 - \frac{\ell}{n} \right)^{d-2} + \mathcal{O}(n^{-1}).$$

Conditioning on that, the probability that Property (P2) holds for both X_k and X_ℓ can be computed the same way as in (5). Clearly, the fact that $X_k = 1$ affects the probability that $X_\ell = 1$ but the difference is hidden in the $\mathcal{O}(1)$ terms. We get that

$$\mathbb{P}(X_k = X_\ell = 1) = \mathcal{O}(n^{-1}) + \left(1 - \left(1 - \frac{k}{n}\right)^{d-2}\right)^{d-1} \cdot \left(1 - \left(1 - \frac{\ell}{n}\right)^{d-2}\right)^{d-1}$$

It follows that $\mathbb{P}(X_k = X_\ell = 1) - \mathbb{P}(X_k = 1)\mathbb{P}(X_\ell = 1) = \mathcal{O}(n^{-1})$ and so $\mathbb{V}ar[X] = \mathcal{O}(n)$. Since $\mathbb{V}ar[X] = o((\mathbb{E}[X])^2)$, X is well-concentrated around $\mathbb{E}[X]$ by the second moment method. The proof of the theorem is finished.

5. Lower Bounds From the Expansion Properties: d-regular Graphs, $d \geq 3$

In this section, we provide rudimentary lower bounds by invoking the Expander Mixing Lemma. However, let us first present an observation that will be used not only here but also for stronger numerical bounds presented in the next section.

Lemma 5.1. Suppose that $H(G) \leq k$ for some graph G = (V, E). Then, V can be partitioned into sets S, T, and U such that |S| = |T| = (n - k)/2, |U| = k, and there is no edge between S and T.

Proof. Consider any graph G = (V, E) with $H(G) \leq k$. By the definition of the hopping number, there exists a set $B_1 \subseteq V$ of cardinality k such that one can initially colour vertices of B_1 blue, and then turn everything blue after a sequence of hops.

Recall that at time step $t \ge 1$, B_t denotes the set of blue vertices (either active, dormant or extinct) and W_t denotes the set of white vertices. Let $B'_t \subseteq B_t$ be the set of extinct blue vertices at time step t, that is, the set of blue vertices that already performed a force. Note that initially $B'_1 = \emptyset$ and $W_1 = V \setminus B_1$. More importantly, each hop increases the cardinality of B'_t by one (one active blue vertex becomes extinct) and decreases the cardinality of W_t by one (one white vertex becomes blue, either active or dormant but certainly not extinct). In particular, at time t = (n - k)/2, $|B'_{(n-k)/2}| = |W_{(n-k)/2}| = (n - k)/2$. On the other hand, $B_t \setminus B'_t$ keeps changing during the process but its cardinality is equal to k for all t.

Another important property is that for any t, there is no edge between vertices in B'_t and vertices in W_t . Indeed, when an active vertex $v \in A_t \subseteq B_t$ performs a force at time t, all of its neighbours have to be blue (that is, none of them is in W_t). This vertex is moved to B'_{t+1} and, since W_t is only shrinking (that is, $W_1 \supset W_2 \supset \ldots$), the desired property will be preserved to the end of the process.

The lemma follows by taking the following partition of the vertex set $V: S = B'_{(n-k)/2}, T = W_{(n-k)/2}$, and $U = B_{(n-k)/2} \setminus B'_{(n-k)/2} = V \setminus (S \cup T)$.

The above lemma informally tells us that if the hopping number of some graph G is small, then there are two large sets S and T without any edges in between. Such situation does not happen in good expanders, in particular, this property is not satisfied a.a.s. for dense random d-regular graphs. It provides a lower bound for the hopping number of $\mathcal{G}_{n,d}$ that holds a.a.s. We provide two arguments yielding two corresponding lower bounds. The first bound is stronger for $d \geq 35$.

For convenience, we repeat the statement of Theorem 1.3. Let G = (V, E) be a *d*-regular graph with *n* vertices and set $\lambda = \lambda(G)$. Then,

$$H(G) \ge \max\left(1 - \frac{2\lambda}{d}, \frac{d - \lambda}{d + 3\lambda}\right) n = \left(1 - \min\left(\frac{2\lambda}{d}, \frac{4\lambda}{d + 3\lambda}\right)\right) n.$$

As a result, for any $d \geq 3$ and $\varepsilon > 0$, a.a.s.

$$H(\mathcal{G}_{n,d}) \ge \left(1 - \min\left(\frac{4\sqrt{d-1}}{d}, \frac{8\sqrt{d-1}}{d+6\sqrt{d-1}}\right) - \varepsilon\right)n.$$

Proof of Theorem 1.3. Consider any d-regular graph G on n vertices and set $\lambda = \lambda(G)$. It follows immediately from the Expander Mixing Lemma (Lemma 2.2) that for any two disjoint sets S and T of cardinality (n - k)/2,

$$|E(S,T)| \ge \frac{d|S||T|}{n} - \lambda \sqrt{|S||T|} = \frac{d(n-k)^2}{4n} - \lambda \frac{n-k}{2} = \frac{d(n-k)}{4} \left(\frac{n-k}{n} - \frac{2\lambda}{d}\right) > 0,$$

as long as $k < (1 - 2\lambda/d)n$. Hence, by Lemma 5.1, we get that

$$H(G) \ge \left(1 - \frac{2\lambda}{d}\right)n.$$

To get the second lower bound, let us again consider any two disjoint sets S and T of cardinality (n-k)/2. It follows from (1) that

$$|E(S, V \setminus S)| \ge (d - \lambda) \frac{|S||V \setminus S|}{n} = (d - \lambda) \frac{(n - k)(n + k)}{4n},$$

and the same lower bound holds for $|E(T, V \setminus T)|$. Similarly, using (1) one more time, after setting $U = V \setminus (S \cup T)$, we get that

$$|E(U, V \setminus U)| \le (d+\lambda)\frac{|U||V \setminus U|}{n} = (d+\lambda)\frac{k(n-k)}{n},$$

since |U| = n - (n - k)/2 - (n - k)/2 = k. We get that

$$\begin{aligned} |E(S,T)| &= \frac{1}{2} \Big(|E(S,V\setminus S)| + |E(T,V\setminus T)| - |E(U,V\setminus U)| \Big) \\ &\geq \frac{1}{2} \Big((d-\lambda) \frac{(n-k)(n+k)}{4n} + (d-\lambda) \frac{(n-k)(n+k)}{4n} - (d+\lambda) \frac{k(n-k)}{n} \Big) \\ &= \frac{n-k}{4n} \Big((d-\lambda)(n+k) - 2(d+\lambda)k \Big) \\ &= \frac{n-k}{4n} \Big((d-\lambda)n - (d+3\lambda)k \Big) > 0, \end{aligned}$$

provided that $k < \frac{d-\lambda}{d+3\lambda}n$, and so by Lemma 5.1 we get that

$$H(G) \ge \frac{d-\lambda}{d+3\lambda} \, n.$$

The conclusion for $\mathcal{G}_{n,d}$ follows immediately from Lemma 2.1 which finishes the proof of the theorem.

6. Lower Bounds From the Configuration Model: d-regular Graphs, $d \ge 3$

We will continue exploiting Lemma 5.1 to get stronger (but numerical, not explicit) lower bounds for the hopping number. This time we will use the configuration model (see Subsection 2.2) to show that there are no two large disjoint sets in $\mathcal{G}_{n,d}$ with no edge between them.

For a given integer $d \geq 3$, let

$$g_{d}(x,z) = \left(\frac{d}{2} - 1 - dz\right) x \log(x) + (d-1)(1-x) \log\left(\frac{1-x}{2}\right) - 2dxz \log(z) - \frac{(1-2z)dx}{2} \log(1-2z) - d\left(\frac{1-x}{2} - zx\right) \log\left(\frac{1-x}{2} - zx\right).$$
(7)

For a fixed $x \in (0, 1)$, function $g_d(x, z)$ is maximized at

$$z_0(x) := \frac{1 - \sqrt{1 - 2(1 - x)x}}{2x} \,. \tag{8}$$

(Note that $z_0(x)$ does not depend on d; see the proof of Theorem 6.1 for more details.) Let $h_d(x) = g_d(x, z_0(x))$ be the corresponding maximum value. (For an illustration, we plot functions $h_3(x)$ and $h_{10}(x)$ in Figure 6.1.)

Function $h_d(x)$ has the following properties: $\lim_{x\to 0} h_d(x) = -(d-2)\log(2)/2 < 0$, it is increasing on the interval $(0, \hat{x}]$ (for some $\hat{x} \in (0, 1)$), decreasing on the interval $[\hat{x}, 1)$, and $\lim_{x\to 1} h_d(x) = 0$. As a result, there is a unique value $x_d \in (0, 1)$ for which $h_d(x_d) = 0$. This value can be easily



FIGURE 6.1. Function $h_3(x)$ and $h_{10}(x)$.

approximated numerically and will play a crucial role in our next result. In light of the above definitions, the following theorem is equivalent to Theorem 1.4.

Theorem 6.1. Fix any integer $d \ge 3$ and $\varepsilon > 0$. Let x_d be the unique $x \in (0,1)$ for which $h_d(x) = g_d(x, z_0(x)) = 0$, where $g_d(x, z)$ and $z_0(x)$ are defined in (7) and in (8), respectively. Then, a.a.s.

$$H(\mathcal{G}_{n,d}) > (x_d - \varepsilon)n.$$

Proof. Fix $d \geq 3$ and consider the configuration model generating $\mathcal{G}_{n,d}$. Suppose that for some carefully chosen function x = x(n) (0 < x < 1), the expected number S(x) of partitions of n vertices of $\mathcal{G}_{n,d}$ into sets S, T, and U such that |S| = |T| = (1-x)n/2, |U| = xn, with the property that there is no edge between S and T is o(1). Then, it follows from the first moment method that a.a.s. there is no such partition in $\mathcal{G}_{n,d}$ and we immediately get from Lemma 5.1 that a.a.s. $H(\mathcal{G}_{n,d}) > xn$.

Let us fix some auxiliary functions y = y(n) and z = z(n) such that y > 0, z > 0, y + z < 1, yx < (1-x)/2, and zx < (1-x)/2. Let S(x, y, z) be the expected number of partitions into sets S, T, and U such that |S| = |T| = (1 - x)n/2, |U| = xn, with the properties that |E(U, S)| = ydxn, |E(U, T)| = zdxn, and |E(S, T)| = 0. It is clear that

$$S(x, y, z) = \binom{n}{xn} \binom{(1-x)n}{(1-x)n/2} \binom{dxn}{ydxn} \binom{d(1-x)n/2}{ydxn} (ydxn)!$$
$$\cdot \binom{(1-y)dxn}{zdxn} \binom{d(1-x)n/2}{zdxn} (zdxn)!$$
$$\cdot M((1-y-z)dxn)M(d((1-x)/2-yx)n)M(d((1-x)/2-zx)n)/M(dn),$$

where M(i) is the number of perfect matchings on *i* points, that is,

$$M(i) = \frac{i!}{(i/2)!2^{i/2}}.$$

Indeed, we first need to select vertices to form set $U(\operatorname{term} \binom{n}{xn})$ and partition the remaining vertices into S and $T(\operatorname{term} \binom{(1-x)n}{(1-x)n/2})$. After that we need to select points in $U(\operatorname{term} \binom{dxn}{ydxn})$ and points in $S(\operatorname{term} \binom{d(1-x)n/2}{ydxn})$ that will be matched in the configuration model with edges between U and S; there are (ydxn)! ways to do that. Then, we need to select points from the remaining points in U(term $\binom{(1-y)dxn}{zdxn}$) and points in T (term $\binom{d(1-x)n/2}{zdxn}$) and match them (term (zdxn)!) to form edges between U and T. Finally, we independently and arbitrarily pair the remaining points in U (term M((1-y-z)dxn)), in S (term M(d((1-x)/2-yx)n)), and in T (term M(d((1-x)/2-zx)n))). Since each of these configurations occurs with the same probability, namely, with probability 1/M(dn), we get the expected value by dividing the product of all above terms by M(dn).

After simplification we get

$$\begin{split} S(x,y,z) &= n!(dxn)!(d(1-x)n/2)!^2 2^{dn/2}(dn/2)! \\ &\cdot (xn)!^{-1}((1-x)n/2)!^{-2}(ydxn)!^{-1}(zdxn)!^{-1} \\ &\cdot 2^{-(1-y-z)dxn/2}((1-y-z)dxn/2)!^{-1} \\ &\cdot 2^{-d((1-x)/2-yx)n/2}(d((1-x)/2-yx)n/2)!^{-1} \\ &\cdot 2^{-d((1-x)/2-zx)n/2}(d((1-x)/2-zx)n/2)!^{-1}(dn)!^{-1} \end{split}$$

Using Stirling's formula $(i! \sim \sqrt{2\pi i}(i/e)^i)$ we obtain

$$S(x, y, z) = \Theta(n^{-2}) \exp\left(f_d(x, y, z)n\right),$$

where

$$\begin{aligned} f_d(x,y,z) &= \left(d-1-d(y+z) - \frac{(1-y-z)d}{2}\right) x \log(x) \\ &+ (d-1)(1-x) \log\left(\frac{1-x}{2}\right) \\ &- dxy \log(y) - dxz \log(z) \\ &- \frac{(1-y-z)dx}{2} \log(1-y-z) \\ &- \frac{d}{2} \left(\frac{1-x}{2} - yx\right) \log\left(\frac{1-x}{2} - yx\right) - \frac{d}{2} \left(\frac{1-x}{2} - zx\right) \log\left(\frac{1-x}{2} - zx\right). \end{aligned}$$

Not surprisingly, for a fixed $x \in (0, 1)$, the function $f_d(x, y, z)$ is maximized when z = y. To see it, for example, we may compute the directional derivative of $f_d(x, y, z)$ in the direction (0, 1, -1), which is equal to

$$\frac{dx}{2}\left(2\left(\log(z) - \log(y)\right) + \log\left(\frac{1-x}{2} - xy\right) - \log\left(\frac{1-x}{2} - xz\right)\right).$$

Clearly, if y < z, then this derivative is positive which implies that the maximum is obtained when z = y. This motivates the definition (7) of $g_d(x, z)$, since $g_d(x, z) = f_d(x, z, z)$. To maximize $g_d(x, z)$, we compute the derivative with respect to z:

$$\frac{\partial g(x,z)}{\partial z} = dx \log\left(\frac{(1-x-2zx)(1-2z)}{2xz^2}\right).$$

It follows that $\frac{\partial g(x,z)}{\partial z} = 0$ if and only if $(1 - x - 2zx)(1 - 2z) = 2xz^2$. By solving this quadratic equation we get that

$$z = \frac{1/x \pm \sqrt{\Delta}}{2}$$
, where $\Delta = \frac{1}{x^2} - \frac{2}{x} + 2$

The constraint y + z < 1 implies that z < 1/2 and, since the larger root is at least 1/(2x) > 1/2, it is not a feasible solution. This motivates the definition of (8), as $z_0(x) = (1/x - \sqrt{\Delta})/2$.

We conclude that for any $0 \le y, z \le 1$ satisfying our constraints,

$$f_d(x, y, z) \le h_d(x) = g_d(x, z_0(x)) = f_d(x, z_0(x), z_0(x))$$

Finally, recall that $x_d \in (0, 1)$ is the unique value of x for which $h_d(x) = 0$. Moreover, $h_d(x_d - \varepsilon) < 0$. As a result, the expected number $S(x_d - \varepsilon)$ of partitions of the vertex set into sets S, T, and U such that $|S| = |T| = (1 - x_d + \varepsilon)n/2$, $|U| = (x_d - \varepsilon)n$, with the property that there is no edge between S and T can be estimated as follows

$$S(x_d - \varepsilon) = \sum_{y,z} S(x_d - \varepsilon, y, z) = \sum_{y,z} \mathcal{O}(n^{-2}) \exp\left(f_d(x_d - \varepsilon, y, z)n\right)$$
$$= \sum_{y,z} \mathcal{O}(n^{-2}) \exp\left(h_d(x_d - \varepsilon)n\right) = \mathcal{O}(1) \exp\left(h_d(x_d - \varepsilon)n\right)$$
$$= \mathcal{O}(1) \exp\left(-\Omega(n)\right) = o(1).$$

Hence, a.a.s. there is no partition with this property and Lemma 5.1 implies that a.a.s. $H(\mathcal{G}_{n,d}) > (x_d - \varepsilon)n$. This finishes the proof of the theorem.

7. Upper Bound from the Degree-Greedy Algorithm: 3-regular graphs

In this section, we assume that d = 3 is fixed with dn even. In order to get an asymptotically almost sure upper bound on the hopping number, we study an algorithm that selects random vertices of minimum degree and tries to hop from them. This algorithm is called *degree-greedy* because the vertex is chosen from those with the lowest degree. Similar technique was successfully used to estimate the brush number of random *d*-regular graphs [2]. We are not as successful in our present problem but we do hope that one can modify our algorithm, and use similar techniques to analyze it, to get better bounds than the ones we managed to prove. We illustrate our ideas in the simplest case, namely, for d = 3.

We start with a random 3-regular graph G = (V, E) on n vertices, and we will work with the configuration model. During the process, we will keep track of a set D_t of vertices that have at least one point that is still unmatched. Vertices in D_t will be considered as potential candidates to hop from. The process will ensure that these vertices have not perform a force yet. Moreover, white vertices are easy to identify, namely, they have 3 unmatched points.

Initially, $D_0 = V$. In every step t of the process, we select a random vertex α_t , chosen uniformly at random from those vertices with the lowest degree in the induced subgraph $G[D_{t-1}]$ on unexposed points. We expose the neighbours associated with the remaining points of α_t , make them blue (if they are still white), and try to hop through one of them. To be able to do it, one of these neighbours has to have at least one remaining unchosen point which is associated with a white vertex. Regardless, whether we succeed to hop or not, α_t has to be removed from D_t together with other vertices that got all points exposed.

In the first step, a vertex of degree 3 is selected to hop from. Three of its neighbours become blue and we hop through one of them, making another vertex blue. (A.a.s. there is no triangle around the initial vertex.) The induced subgraph $G[D_1]$ now has 1 vertex of degree 1, 3 vertices of degree 2, and n-5 vertices of degree 3. Note that α_1 is a.a.s. the only vertex whose degree in $G[D_t]$ is 3 at the time of selecting a vertex to hop from. Indeed, if α_t ($t \ge 2$) has degree 3 in $G[D_{t-1}]$, then $G[D_{t-1}]$ consists of some connected components of G and thus G is disconnected. It was proven in [25] that for constant d, G is disconnected with probability o(1) (this also holds when d is growing with n, as shown in [22]).

In the second step, we try hop from the vertex of degree 1. A.a.s. its neighbour is white and after hopping through it, this vertex becomes of degree 1. However, when vertices of degree 2 become plentiful, we will be hopping through them often (making them of degree 0) and we might run out of vertices of degree 1. We will start hopping from vertices of degree 2 but eventually come back to hopping from vertices of degree 1. Our goal is to control such "hiccups". The details of the application of the differential equations method to such degree-greedy algorithms have been omitted, but can be found in [28]. For $0 \le i \le 3$, let $Y_i = Y_i(t)$ denote the number of vertices with *i* unmatched points at time *t*. (Note that $Y_0(t) = n - \sum_{i=1}^3 Y_i(t)$ so $Y_0(t)$ does not need to be calculated, but it is useful in the discussion.) Let $S(t) = \sum_{i=1}^3 iY_i(t)$ be the total number of points that are unmatched at time *t*. It is tedious but not so difficult to consider all the cases to see that for $r \in \{1, 2\}$ and $i \in \{0, 1, 2, 3\}$,

$$\mathbb{E}\Big(Y_i(t+1) - Y_i(t) \mid G[D_t] \land \deg_{G[D_t]}(\alpha_{t+1}) = r\Big) \\
= f_{i,r}(t/n, Y_1(t)/n, Y_2(t)/n, Y_3(t)/n) \\
= \sum_j p_r^{(j)} \Delta Y_i^{(j)},$$
(9)

where $p_r^{(j)}$ and $\Delta Y_i^{(j)}$ are presented in Tables 3 and 4 (the sum is over all possible cases: $j \in [5]$ for r = 1, and $j \in [12]$ for r = 2). Indeed, let us explain, for example, the very first case, r = 1 and j = 1. We try to hop from a vertex of degree 1. Its neighbour had degree 3 with probability $3Y_3(t)/(S(t)-1) \sim 3Y_3(t)/S(t)$ but will have degree 1 once we hop through it. Its neighbour is white with probability $3(Y_3(t) - 1)/(S(t) - 3) \sim 3Y_3(t)/S(t)$. Hence, this case happens with probability asymptotic to $p_1^{(1)} = (3Y_3(t)/S(t))^2$. We can hop there, making this effort count. Three vertices changed their degrees. Initially we have one vertex of degree 1 and 2 of degree 3. After the hop, one of them became of degree 0, one of degree 1, and one of degree 2, explaining the $\Delta Y_i^{(1)}$'s. Note that sometimes we are not able to make a hop, as none of the vertices in the second neighbourhood were initially white. We indicate whether the hop occurred or not in the last column in the two corresponding tables.

j	$p_1^{(j)}$	$\Delta Y_0^{(j)}$	$\Delta Y_1^{(j)}$	$\Delta Y_2^{(j)}$	$\Delta Y_3^{(j)}$	succesful hop?
1	$\left(\frac{3Y_3(t)}{S(t)}\right)^2$	1	0	1	-2	Yes
2	$\left(\frac{3Y_3(t)}{S(t)}\right)^2 \left(\frac{2Y_2(t)}{S(t)}\right)$	2	0	0	-2	Yes
3	$\left(\frac{3Y_3(t)}{S(t)}\right) \left(\frac{2Y_2(t)}{S(t)}\right)^2$	2	1	-2	-1	No
4	$\left(\frac{3Y_3(t)}{S(t)}\right)\left(\frac{2Y_2(t)}{S(t)}\right)$	2	-1	0	-1	Yes
5	$\left(\frac{2Y_2(t)}{S(t)}\right)^2$	2	0	-2	0	No

TABLE 3. Vertex of degree r = 1 tries to hop (5 cases)

Suppose that at some step t of the process, hopping from a vertex of degree 2 creates, in expectation, β vertices of degree 1 and hopping from a vertex of degree 1 decreases, in expectation, the number of vertices of degree 1 by τ . After hopping from a vertex of degree 2, we expect to then hop (on average) from β/τ vertices of degree 1. Thus, the proportion of steps which hop from vertices of degree 2 is $1/(1 + \beta/\tau) = \tau/(\beta + \tau)$. If τ falls below zero, vertices of degree 1 begin to build up and do not decrease under repeated hoppings of this type and we move to the next phase.

From (9) it follows that

$$\beta = \beta(x, y_1, y_2, y_3) = f_{1,2}(x, y_1, y_2, y_3) = f_{1,2}(x, \mathbf{y}),$$

$$\tau = \tau(x, y_1, y_2, y_3) = -f_{1,1}(x, y_1, y_2, y_3) = -f_{1,1}(x, \mathbf{y}),$$

where x = t/n and $y_i(x) = Y_i(t)/n$ for $i \in [3]$. This suggests the following system of differential equations

$$\frac{dy_i}{dx} = F(x, \mathbf{y}, i)$$

where

$$F(x, \mathbf{y}, i) = \frac{\tau}{\beta + \tau} f_{i,2}(x, \mathbf{y}) + \frac{\beta}{\beta + \tau} f_{i,1}(x, \mathbf{y}).$$

j	$p_2^{(j)}$	$\Delta Y_0^{(j)}$	$\Delta Y_1^{(j)}$	$\Delta Y_2^{(j)}$	$\Delta Y_3^{(j)}$	succesful hop?
1	$\left(\frac{3Y_3(t)}{S(t)}\right)^2 \left(\frac{2Y_2(t)}{S(t)}\right)$	1	1	1	-3	Yes
2	$\left(\frac{3Y_3(t)}{S(t)}\right)^3 \left(\frac{2Y_2(t)}{S(t)}\right)$	2	1	0	-3	Yes
3	$\left(rac{3Y_3(t)}{S(t)} ight)^3 \left(rac{2Y_2(t)}{S(t)} ight)^2$	2	3	-2	-3	Yes
4	$\left(\frac{3Y_3(t)}{S(t)}\right)^3 \left(\frac{2Y_2(t)}{S(t)}\right)^3$	3	3	-3	-3	Yes
5	$\left(\frac{3Y_3(t)}{S(t)}\right)^2 \left(\frac{2Y_2(t)}{S(t)}\right)^4$	3	4	-5	-2	No
6	$2\left(\frac{3Y_3(t)}{S(t)}\right)^2\left(\frac{2Y_2(t)}{S(t)}\right)$	2	0	0	-2	Yes
7	$2\left(\frac{3Y_3(t)}{S(t)}\right)^2\left(\frac{2Y_2(t)}{S(t)}\right)^2$	2	2	-2	-2	Yes
8	$2\left(\frac{3Y_3(t)}{S(t)}\right)^2 \left(\frac{2Y_2(t)}{S(t)}\right)^3$	3	2	-3	-2	Yes
9	$2\left(\frac{3Y_3(t)}{S(t)}\right)\left(\frac{2Y_2(t)}{S(t)}\right)^4$	3	3	-5	-1	No
10	$\left(\frac{3Y_3(t)}{S(t)}\right) \left(\frac{2Y_2(t)}{S(t)}\right)^2$	2	1	-2	-1	Yes
11	$\left(\frac{3Y_3(t)}{S(t)}\right) \left(\frac{2Y_2(t)}{S(t)}\right)^3$	3	1	-3	-1	Yes
12	$\left(\frac{2Y_2(t)}{S(t)}\right)^4$	3	2	-5	0	No

TABLE 4. Vertex of degree r = 2 tries to hop (12 cases)

At this point we may formally define the termination point \hat{x} is defined as the infimum of those $x > \hat{x}$ for which at least one of the following holds: $\tau \le 0$, $\tau + \beta = 0$, or $y_2 \le 0$. The initial conditions are $y_3(0) = 1$ and $y_i(0) = 0$ for $i \in \{0, 1, 2\}$.

The general result [28, Theorem 1] studies a deprioritized version of degree-greedy algorithms, which means that the vertices are chosen to process in a slightly different way, not always the minimum degree, but usually a random mixture of two degrees. Once a vertex is chosen, it is treated the same as in the degree-greedy algorithm. The variables Y are defined in an analogous manner. The hypotheses of the theorem are mainly straightforward to verify but require several inequalities involving derivatives to hold at the termination of phases, for the full rigorous conclusion to be obtained. However, in practice, the equations are simply solved numerically in order to find the points \hat{x} , since a fully rigorous bound is not obtained unless one obtains strict inequalities on the values of the solutions. The conclusion is that, for a certain algorithm using a deprioritized "mixture" of the steps of the degree-greedy algorithm, with variables Y_i defined as above, we have that a.a.s.

$$Y_i(t) = ny_i(t/n) + o(n)$$

for $0 \le i \le 3$. We omit all details, pointing the reader to [28] and Subsection 2.5 for a short discussion on the differential equations method which is the main tool in proving the result. In addition, the theorem gives information on an auxiliary variable such as, of importance to our present application, the number of vertices that actually hopped.

The numerical solution to the relevant differential equations is shown in Figure 7.1. During the process, only vertices of degree 1 and 2 attempt to make a hop. Since we prioritize vertices of degree 1, the function y_1 responsible for "monitoring" such vertices is equal to zero. Note that it means that a.a.s. there are only o(n) vertices of degree 1 at any point of the process, not that we do not see them at all (we clearly do). As expected, the number of vertices of degree 3 (represented by y_3) decreases, and the number of vertices of degree 0 (represented by y_0) increases. The number of vertices of degree 2 (represented by y_2) initially increases but then it decreases. A.a.s., at time $\hat{x}n \approx 0.6614n$ the process stops with all vertices being of degree 0. By investigating the auxiliary random variable, and the associated differential equation, we conclude that a.a.s. approximately 0.5159n



FIGURE 7.1. Evolution of $Y_i(t)$, $0 \le i \le 3$ for d = 3 using the Differential Equation Method.

vertices hop, yielding an upper bound for the hoping number of approximately (1 - 0.5159)n = 0.4841n. Unfortunately, it is a much weaker upper bound than the one we obtained by investigating the contiguous model, namely, (1 + o(1))n/3.

8. FUTURE DIRECTIONS

In this paper, we investigated the hopping number of random d-regular graphs. Unfortunately, for all values of d, except the d = 2 case, we only have upper and lower bounds—see Table 1 and Figure 1.1. It would be nice to narrow the gaps between these bounds.

It feels natural to attempt a greedy algorithm to generate a small initial set of blue vertices. We analyzed one such heuristic algorithm for the d = 3 case in Section 7 using the Differential Equation method. Unfortunately, the bound we got turned out to be weaker than the one based on the contiguous model. Nevertheless, one may want to leverage the same tool to analyze similar algorithms with, hopefully, better success.

There are numerous variants of the classic zero forcing rule. Some other colouring rules might be analyzable using similar techniques as the ones used in this paper. Moreover, we concentrated on minimizing the size of an initial set of blue vertices. Some other papers investigate the propagation time instead, assuming that multiple forces are performed simultaneously in a sequence of time steps, and optimize the trade-off between the size of a forcing set and its propagation time—the so-called throttling number. Analyzing the throttling number for hopping forcing (or any other forcing rule) on random structures might be an interesting task.

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DEPARTMENT OF MATHEMATICS, TORONTO METROPOLITAN UNIVERSITY, TORONTO, CANADA *Email address*: pralat@torontomu.ca

DEPARTMENT OF MATHEMATICS, TORONTO METROPOLITAN UNIVERSITY, TORONTO, CANADA *Email address*: harjas.singh@torontomu.ca