On the Complexity of Distributed Graph Coloring with Local Minimality Constraints

Cyril Gavoille and Ralf Klasing
LaBRI - Université Bordeaux 1 - CNRS, 351 cours de la Libération, 33405 Talence cedex, France

Adrian Kosowski
Department of Algorithms and System Modeling, Gdańsk University of Technology, Narutowicza 11/12, 80952 Gdańsk, Poland

Łukasz Kuszner
Department of Algorithms and System Modeling, Gdańsk University of Technology, Narutowicza 11/12, 80952 Gdańsk, Poland

Alfredo Navarra
Dipartimento di Matematica e Informatica, Università degli Studi di Perugia, Via Vanvitelli 1, 06123 Perugia, Italy

Distributed greedy coloring is an interesting and intuitive variation of the standard coloring problem. Given an order among the colors, a coloring is said to be greedy if there does not exist a vertex for which its associated color can be replaced by a color of lower position in the fixed order without violating the property that neighboring vertices must receive different colors. We consider the problems of Greedy Coloring and Largest First Coloring (a variant of greedy coloring with strengthened constraints) in the Linial model of distributed computation, providing lower and upper bounds and a comparison to the $(\Delta + 1)$-Coloring and Maximal Independent Set problems, with $\Delta$ being the maximum vertex degree in $G$. © 2009 Wiley Periodicals, Inc. NETWORKS, Vol. 54(1), 12–19 2009

1. INTRODUCTION

To color the vertices of a graph $G = (V, E)$ means to give each vertex a positive integer color value in such a way that no two adjacent vertices get the same color. In many practical considerations, it is desirable to minimize the number of colors used.

Many applications ranging over code assignment in wireless networks [3], scheduling problems [10,23], track assignment in railway optimization [5–7], and so forth, can be solved by means of minimum coloring.

If at most $k$ colors are used, the result is called a $k$-coloring. The smallest possible positive integer $k$ for which there exists a $k$-coloring of $G$ is called the chromatic number $\chi(G)$. It has to be remembered that even in a centralized setting, approximating $\chi(G)$ within a factor of $|V|^{1-\varepsilon}$ is an NP-hard problem, for any $\varepsilon > 0$ [30].

In the context of distributed computing, the performance of a graph coloring algorithm on a system graph $G$ is characterized by at least two main parameters: the number of colors used by the algorithm to color the graph $G$ and the number of rounds required to obtain a coloring. Designing a fast distributed algorithm (i.e., an algorithm running within $o(D)$ rounds where $D$ is the diameter of the graph $G$) which always uses a number of colors in some way bounded from above with respect to $\chi(G)$ appears to be difficult; e.g., in [22] it is proved that coloring trees of diameter $D$ and maximum degree $\Delta$ requires $\Omega(D^2)$ rounds even if we allow the algorithm to use $\sqrt{\Delta}$ colors. For this reason, it is natural to pose the problem of constructing distributed algorithms having properties analogous to certain centralized graph coloring algorithms.
heuristics, which are known to work well in practice. The
class of greedy algorithms considered herein is of special
significance due to the elegance of their formulation, and
the number of graph classes for which they always produce
optimal or near-optimal results.

1.1. Preliminaries: Greedy Coloring in a
Centralized Setting

For a given graph \( G \) and a sequence \( K \) of all its vertices,
\( K = (v_1, v_2, \ldots, v_n) \), we will use the term greedy coloring
to describe the following procedure of locally minimal color
assignment:

**Algorithm** Greedy-Color\( (G, K) \):

\[
\text{for } v := v_1 \text{ to } v_n \text{ do} \\
\text{give vertex } v \text{ the smallest possible color not used by any} \\
of the already colored neighbors of } v;
\]

Different graph coloring algorithms are obtained by choosing
the sequence \( K \) in a specific way and then applying the
Greedy-Color procedure. Below we briefly recall some of
the types of sequences most often applied in practice (cf.
[17, 18]).

- **General Greedy** (G) sequence: \( K \) is an arbitrarily chosen
sequence of vertices.

- **Largest-First** (LF) sequence: \( K \) is formed by arranging the
vertices of graph \( G \) in non-ascending order of degrees.

- **Smallest-Last** (SL) sequence: \( K \) is formed by iteratively
removing a vertex of minimal degree from the graph and
placing it at the end of \( K \).

All the considered algorithms obviously lead to correct color-
ings of the graph, but the converse does not hold — not every
coloring of a graph can be obtained by applying a specific
algorithm. We will call a coloring of a graph an A-coloring
if it can be obtained by means of greedy coloring with a
sequence of some type A.

Observe that all G-colorings have the much desired local
minimality property described by Grundy [13], namely, no
single vertex may have its color value decreased without
affecting the color of some other (neighboring) vertex. As
a direct consequence, all G-colorings use no more than \( \Delta + 1 \)
colors, with \( \Delta \) being the maximum vertex degree in \( G \). LF-
colorings and SL-colorings are special types of G-colorings,
thus they inherit this property.

For graph classes found in practice greedy colorings
obtained according to specific sequences may admit even
stronger bounds [18, 25, 29]. For example, any SL-coloring
of a planar graph uses no more than six colors, whereas
any LF-coloring is optimal or within a fixed number of col-
ors of the optimum for numerous graph classes, including
complete \( k \)-partite graphs, caterpillars, crowns, and bipartite
wheels [18].

1.2. Model of Distributed Computation

We consider restricted variants of the vertex coloring prob-
lem in a distributed network, assuming the so-called Linial
model of computation, which is widely used in previous
research on the subject [4, 21, 22, 27]. Such a distributed
network consists of a set \( V \) of processors and a set \( E \) of bi-
directional communication links between pairs of processors.
It can be modeled by an undirected graph \( G = (V, E) \). We
denote \( n = |V| \).

Vertices are identified by means of unique labels. Each
vertex has its local state, described by a certain number of
integer variables. The system functions in so-called synchro-
nized rounds, consisting of three steps: a vertex first reads
its own local state variables and the local state variables of
all its neighbors, then performs an arbitrary amount of local
computations, and finally updates its local state variables
accordingly.

Although local computations are unbounded, the case of
randomized algorithms, i.e., algorithms that can make use of
a so-called coin-flip function, is usually separately studied.
Section 4 is devoted to such a variant.

In all further considerations we assume that two global
parameters are known to all vertices: some constant-factor
upper bounds on the number of vertices \( n \) and on the max-
imum vertex degree \( \Delta \). This assumption is used only to
provide a simple mechanism for executing subroutines of
known complexity with respect to \( n \) and \( \Delta \) and waiting a
known number of rounds for their completion.

For convenience of notation, we will assume that the spe-
cial local state variable \( c \) always stores the outcome of the
algorithm; for example, in a coloring process, \( c(v) \) is the color
value assigned to vertex \( v \). The algorithm is considered to be
complete when the values of all variables \( c \) are correctly set
and the algorithm will not modify them in any subsequent
round.

For each vertex \( v \) define its neighborhood \( N(v) = \{u : \}
\{u, v\} \in E \text{ and vertex degree } \deg_G(v) = |N(v)| \). The
maximum degree among all vertices in \( G \) is denoted by
\( \Delta = \max_{v \in V} \deg_G(v) \). To distinguish among neighbors
of higher degree, we will use the symbol \( N_{\geq}(v) = \{u \in N(v) : \)
\deg(u) \geq \deg(v) \} \) and similarly \( N_{<}(v) = \{u \in N(v) : \)
\deg(u) > \deg(v) \} \). The distance-d neighborhood \( N_d(v) \) of
vertex \( v \) is defined as the set of all vertices at a distance
between 0 and \( d \) from \( v \). For any subset of vertices \( S \subseteq V \),
we denote the subgraph of \( G \) induced by vertex set \( S \) as \( G[S] \).
The length of the shortest cycle in graph \( G \) is known as its
girth and denoted by \( g(G) \).

1.3. Problem Definitions

In this article, we consider two variants of the graph col-
oring problem in a distributed setting, those of obtaining a
G-coloring and of obtaining an LF-coloring of the system
graph, denoted G-COL and LF-COL, respectively. As men-
tioned in Section 1.1, G-COL and LF-COL are refinements
of the problem of finding a \((\Delta + 1)\)-coloring of the system

\[ \text{NETWORKS—2009—DOI 10.1002/net 13} \]
graph $G$ (simply denoted COL), and thus are not easier in the sense of computation time. In fact, G-COL and LF-COL are also not easier than the problem of finding a maximal independent set in $G$ (denoted MIS), because for any G-coloring of $G$ the set of all vertices having color 1 is clearly a maximal independent set.

Nevertheless, it is possible to put forward a definition of G-COL and LF-COL which demonstrates the local nature of the imposed constraints; for completeness we provide local formulations of COL and MIS as well. For a set of vertices $S \subseteq V$, the symbol $c(S)$ is used to denote the set of values $\{c(v) : v \in S\}$.

**Definition 1.** The considered distributed problems are defined by the following constraints on the local variable $c$ at any vertex $v$:

- $(\Delta + 1)$-Coloring (COL): $c(v) \in [1, \ldots, \Delta + 1] \setminus c(N(v))$.
- Maximal Independent Set (MIS): $c(v) \in [0, 1] \land c(v) = 1 \iff c(N(v)) = \emptyset$.
- Greedy Coloring (G-COL): $c(v) = \min \{1, \ldots, \Delta + 1\} \setminus c(N(v))$.
- Largest-First Coloring (LF-COL): $c(v) = \min \{1, \ldots, \Delta + 1\} \setminus c(N(v))$.

The proof of the equivalence of the above definitions with the prior characterizations is straightforward. For the G-COL problem, consider any G-coloring $c$ of graph $G$, and an arbitrarily chosen vertex $v$. Because at the time of coloring some of the vertices from $N(v)$ are already colored and $v$ obtains the minimum possible color, we have $c(v) \leq \min[1, \ldots, \Delta + 1] \setminus c(N(v))$, but also $c(v) \notin c(N(v))$ by the legality of the coloring, thus $c(v) = \min[1, \ldots, \Delta + 1] \setminus c(N(v))$. Conversely, given a set of color values fulfilling $c(v) = \min[1, \ldots, \Delta + 1] \setminus c(N(v))$ we can always construct a corresponding sequence $K$ for the Greedy-Color procedure simply by ordering vertices according to nondecreasing values of $c(v)$. For the definition of LF-COL the argument is analogous; an appropriate sequence $K$ can be constructed by ordering vertices according to nonincreasing values of $\deg(v)$, breaking ties according to nondecreasing values of $c(v)$.

As a side note, it is interesting to observe that the problem of finding an SL-coloring of graph $G$ does not admit any local definition. Indeed, when $G$ is a ring of even length, any such coloring uses exactly two colors. Taking into account that the 2-coloring of a ring requires $\Omega(n)$ rounds (cf. e.g. [22]), this means that finding a distributed SL-coloring of a graph may also require $\Omega(n)$ rounds; such a problem is of little interest in a distributed setting. The same also holds for well-known algorithms with dynamic sequences, such as DSATUR [17].

### 1.5. Our Contribution

We provide lower and upper bounds on the time complexity of Greedy Coloring (G-COL) and Largest First Coloring (LF-COL) with respect to Coloring (COL) and Maximal Independent Set (MIS).

A summary of the results is contained in Table 1. The obtained lower and upper bounds for the G-COL and LF-COL problems are expressed in terms of parameters $n$ and $\Delta$. In particular, we prove a lower bound of $\Omega\left(\frac{\log n}{\log \log n}\right)$ for G-COL and $\Omega(\sqrt{n})$ for LF-COL, an improvement upon the current MIS-derived lower bounds [19]. We note that our lower bounds also apply for randomized algorithms (cf. Section 4).

The rest of the article is organized as follows. The next section provides lower bounds for the considered problems. Upper bound results for deterministic algorithms can be found in Section 3, whereas a discussion concerning
TABLE 1. The time complexity of Greedy Coloring and LF-Coloring with respect to other well-known problems, \((\Delta + 1)-\)Coloring and Maximal Independent Set, in the distributed setting.

<table>
<thead>
<tr>
<th></th>
<th>COL</th>
<th>MIS</th>
<th>G-COL</th>
<th>LF-COL</th>
</tr>
</thead>
<tbody>
<tr>
<td>Lower bounds (deterministic and randomized)</td>
<td>(\Omega(\log^* n)) [21]</td>
<td>(\Omega\left(\frac{\log n}{\log \log n}\right)) [19]</td>
<td>(\Omega\left(\frac{\log n}{\log \log n}\right)) [Thm 5]</td>
<td>(\Omega(\sqrt{n})) [Thm 6]</td>
</tr>
<tr>
<td>Upper bounds (deterministic algorithms)</td>
<td>(2^{O(\sqrt{\log n})}) [1, 27]</td>
<td>(O(\Delta n)) [1]</td>
<td>(O(\Delta + T_{\text{COL}})) [Prop 9]</td>
<td>(O(\Delta n)) [Thm 6]</td>
</tr>
<tr>
<td></td>
<td>(O(\Delta \log \Delta + \log^* n)) [20]</td>
<td>(O(\Delta + T_{\text{COL}})) [Prop 9]</td>
<td>(O(\Delta + T_{\text{COL}})) [Prop 9]</td>
<td>(O(\Delta \log T_{\text{COL}})) [Thm 11]</td>
</tr>
<tr>
<td>Upper bounds (randomized algorithms)</td>
<td>(O(Nd)) [15, 24]</td>
<td>(O(\log n)) [24]</td>
<td>(O(D + T_{\text{COL}})) [Prop 13]</td>
<td>(O(D + T_{\text{COL}})) [Prop 13]</td>
</tr>
<tr>
<td></td>
<td>(O(\Delta \log \Delta + \log^* n)) [20]</td>
<td>(O((\Delta + T_{\text{COL}})^2)) [Prop 13]</td>
<td>(O(\Delta \log \Delta + \log^* n)) [Prop 14]</td>
<td>(O(\Delta \log \Delta n)) [Prop 14]</td>
</tr>
</tbody>
</table>

For a problem \(P\), the notation \(T_p\) (resp. \(T_p^R\)) describes a known upper bound on the complexity of an algorithm (resp. randomized algorithm) for solving \(P\).

upper bound results for randomized algorithms can be found in Section 4. Finally, Section 5 provides concluding remarks.

2. LOWER BOUNDS

2.1. The G-COL Problem

For a given graph \(G\), let \(C_G(v)\) denote the set of all possible colors which may be assigned to vertex \(v \in V\), taken over all greedy colorings of \(G\). For trees, we have the following property.

Lemma 1. Let \(T\) be a tree and \(v\) some vertex of \(T\) such that the maximum distance from \(v\) to another vertex of \(T\) is \(d\). Then \(\max C_T(v) \leq d + 1\).

Proof. Consider an arbitrary greedy coloring of \(T\). If a vertex \(v_i\) receives some color \(c > 1\), then it must be adjacent to a vertex \(v_{i-1}\) which is assigned color \(c - 1\). By induction we easily prove that \(v_i\) must be the endpoint of a path in \(T\) with \(c\) vertices. Since by assumption the longest path of \(T\) with one endpoint in \(v\) has \(d + 1\) vertices, clearly the color assigned to \(v\) in the coloring cannot exceed \(d + 1\). ■

Lemma 2. Suppose that there exists an \(n\)-vertex graph \(G\) such that for some \(\ell\) we have bounds on its girth \(g(G) \geq \ell\) and chromatic number \(\chi(G) > \ell/2\). Then, for any distributed algorithm for the G-COL problem, there exist graphs of at most \(n\) vertices which require no less than \(\ell/2 - 1\) rounds to color.

Proof. Suppose that there exists a distributed algorithm \(A\) which for all graphs of at most \(n\) vertices produces a greedy coloring \(C_A\) in at most \(d\) rounds. Algorithm \(A\) must be able to assign for any vertex \(v \in V\) a color \(C_A(v) \in C_G(v)\), using information about the graph structure at distance at most \(d\) only. Thus, if for two graphs \(G = (V, E)\) and \(G' = (V', E')\) with distinguished vertices \(v \in V\), \(v' \in V'\) we have \(G[N_G(v)] = G'[N_{G'}(v')]\), then the coloring of graph \(G\) produced by algorithm \(A\) has the property that \(C_A(v) \in C_G(v) \cap C_G(v')\).

Now, let \(G\) be the graph of girth \(g(G) \geq \ell\) described in the assumption of the lemma. Let \(v \in V\) be arbitrarily chosen. We now apply the proven property for graphs \(G\) and \(G'_v \equiv G[N_{G_v}(v)]\) with vertices \(v\) and \(v' \equiv v\), obtaining \(C_A(v) \in C_G(v) \cap C_{G'}(v)\). However, the graph \(G'_v\) is clearly a tree (because it is acyclic and connected) and all vertices of \(G'_v\) are at a distance of at most \(\ell/2 - 1\) from \(v\). Hence by Lemma 1 we obtain \(\max C_{G'_v}(v) \leq \ell/2\), and we conclude that for graph \(G\), \(C_A(v) \leq \ell/2\). Since \(v\) was arbitrarily chosen, we have \(C_A(v) \leq \ell/2\) for all \(v \in V\). We have thus obtained a coloring of \(G\) using at most \(\ell/2\) colors, whereas by assumption the chromatic number \(\chi(G)\) is larger than \(\ell/2\), a contradiction. ■

Lemma 3 (\([8]\)). For all \(n, k \in \mathbb{N}\) \((4 \leq k \leq n)\) there exists an \(n\)-vertex graph \(G\), such that \(\chi(G) \geq k\) and \(g(G) \geq \frac{\log n}{\log \log n}\).

Corollary 4. For all values of \(n \geq 2^{16}\) there exists a graph \(G\) of order \(n\), such that \(\chi(G) \geq \frac{\log n}{\log \log n}\) and \(g(G) \geq \frac{\log n}{4 \log \log \log n}\).

Proof. For any given value of \(n \geq 2^{16}\), put \(k = \left\lfloor \frac{\log n}{\log \log n} \right\rfloor\) in Lemma 3. We immediately obtain that there exists a graph \(G\) of order \(n\) such that the following conditions are fulfilled:

\[
\chi(G) \geq k \geq \left\lfloor \frac{\log n}{\log \log n} \right\rfloor \geq \frac{\log n}{\log \log n}.
\]

\[
g(G) \geq \frac{1}{4} \frac{\log n}{k} \geq \frac{1}{4} \frac{\log n}{\log \log n} - \frac{1}{4} \log \log n + 1 \geq \frac{1}{4} \frac{\log n}{\log \log n},
\]

which completes the proof. ■
Combining Corollary 4 with Lemma 2 (putting \( l = \frac{1}{4 \log \log n} \)), we immediately obtain that for all values of \( n \geq 2^{16} \), no distributed algorithm can solve the G-COL problem in less than \( \frac{\log n}{\log \log n} - 1 \) rounds for all graphs of order at most \( n \). Thus, we have the following theorem.

**Theorem 5.** The distributed time complexity of G-COL is \( \Omega \left( \frac{\log n}{\log \log n} \right) \).

### 2.2. The LF-COL Problem

**Theorem 6.** The distributed time complexity of LF-COL is \( \Omega(\sqrt{n}) \) and \( \Omega(\Delta) \).

**Proof.** Consider a family of graphs \( \{G_d\}_{d=4,5,6,...} \) having \( n(G_d) = \frac{5}{2}d^2 - \frac{1}{2}d + 1 \) and \( \Delta(G) = 2d \), whose representative is depicted in Figure 1. Vertices \( v_0, v_1, \ldots, v_d \) induce a path of length \( d \). Some additional components are connected to particular vertices to ensure that vertex \( v_i \) obtains a color \( d - i + 1 \) in each LF-coloring of \( G_d \). Each of these additional components \( K_i \) depicts a complete graph with \( r \) vertices, and a bold line between such a component and \( v_i \) illustrates that each of the vertices of \( K_i \) is connected to \( v_i \). Similarly, when two components \( K_{r_j} \) and \( K_{r_j} \) are connected, each vertex from \( K_{r_j} \) is connected to each vertex from \( K_{r_j} \), thus forming a clique \( K_{r_j} \). We have \( \deg(v_i) = d + i \), hence the vertices of the path appear in the LF sequence in the order \( v_d, v_{d-1}, \ldots, v_1, v_0 \). Moreover, we have \( |N_r(v_i)| = d - i \), and it is easy to obtain inductively that in any LF-coloring of \( G_d \) the set of colors used in \( N_r(v_i) \) is \( \{1, 2, \ldots, d - i\} \), hence the only possible color for \( v_i \) in an LF-coloring is \( d - i + 1 \). However, if vertex \( v_d \) were to be removed from the graph, the colors of all other vertices of the path would decrease by 1. Thus, we have shown that the color of \( v_0 \) depends on the existence of vertex \( v_d \); if \( v_d \) exists, \( v_0 \) must obtain color \( d + 1 \), otherwise it must obtain color \( d \). As vertices \( v_d \) and \( v_0 \) are at a distance of \( d \) from each other, and information in our model can propagate only at the speed of one vertex per round, the coloring cannot be completed in less than \( d \) rounds, and the claim follows.

![FIG. 1. A graph which requires \( \Omega(d) \) time to LF-color in the distributed model.](image)

### 3. UPPER BOUNDS FOR DETERMINISTIC ALGORITHMS

#### 3.1. The G-COL Problem

We recall after [1,27] that a \((j,k)\)-decomposition of a graph \( G = (V,E) \) is a partition \( C = \{C_1, C_2, \ldots\} \) of \( V \) such that the following conditions are fulfilled:

- each subgraph \( G[C_i] \), called a cluster, is connected and of diameter at most \( j \);
- the so called cluster graph, having vertex set \( C \) and edges connecting those clusters \( C_i, C_j \in C \) for which there exist vertices \( v_i \in C_i, v_j \in C_j \) such that \( \{v_i, v_j\} \) is an edge of \( G \), is vertex colored with at most \( k \) colors.

**Lemma 7.** Given a \((j,k)\)-decomposition of a graph, there exists a distributed \( O(j \cdot k) \) time algorithm for solving the G-COL problem.

**Proof.** Once the \((j,k)\)-decomposition is given, we have a situation in which the clustered graph is already colored (perhaps not greedily). The algorithm then proceeds according to the assigned colors: first clusters of color 1 are processed in parallel, then clusters of color 2 are processed in parallel, and so forth. Inside each cluster, a leader is elected in order to collect all the information from the related cluster. This operation clearly requires \( O(j) \) rounds to retrieve all the information about the cluster. The leader then greedily assigns a suitable color to each vertex of its cluster respecting the greedy coloring property, i.e., each vertex gets the smallest color possible with respect to its already colored neighbors from other clusters. Again this requires \( O(j) \) rounds. Note that clusters of color 1 have no constraints with respect to the other clusters. In the case of clusters of color 2, the greedy coloring may begin only after waiting \( O(k) \) steps for all the neighbors from clusters of color 1 to be colored. Consequently, a coloring of the last clusters is complete after \( O(j \cdot k) \) rounds.

The following theorem is obtained directly from Lemma 7, taking into account the results of [27] where a \( \left( 2^{O(\sqrt{\log n})} \right) \)-decomposition of any \( n \)-vertex graph is provided in \( 2^{O(\sqrt{\log n})} \) distributed time.

**Theorem 8.** There exists a distributed algorithm for the G-COL problem running in \( 2^{O(\sqrt{\log n})} \) time.

**Proposition 9.** There exists a distributed \( O(\Delta + T_{COL}) \) algorithm for G-COL, where \( T_{COL} \) denotes a known upper bound on the execution time of an algorithm for the COL problem in \( G \).

**Proof.** Suppose that a \((\Delta + 1)\)-coloring of graph \( G \) is already provided. This induces a partition of the vertex set into independent sets \( V = IS_1 \cup IS_2 \cup \cdots \cup IS_{\Delta+4} \). It is easy to see that in the \( i \)-th round we can simultaneously assign the minimum possible color to all vertices from \( IS_i \), taking into account the constraints imposed by the colors of vertices.
from \( IS_1 \cup \cdots \cup IS_{\Delta - 1} \). Thus, a greedy coloring of the graph can be obtained \( \Delta + 1 \) rounds after the initial \((\Delta + 1)\)-coloring is known to be complete.

3.2. The LF-COL Problem

Before discussing the details of the distributed implementation of the LF algorithm, we present an equivalent characterization of a correct LF-coloring. For a given coloring of \( G \), let \( IS_{(j,k)} \in V \), for any \( 1 \leq j \leq \Delta, 1 \leq k \leq \Delta + 1 \), denote the independent set of vertices of \( G \) of degree \( j \) and colored with color \( k \). We define \( H_{(j,k)} \subseteq G \) as the subgraph induced by the set of all vertices \( v \) such that \( \deg_G(v) = j \) and \( v \notin (\bigcup_{k < j} IS_{(j,k)}) \cup N(\bigcup_{j > k} IS_{(j,k)}) \).

**Lemma 10.** Given an assignment of colors \( c : V(G) \to \mathbb{N} \), if for all \( j,k \) the set \( IS_{(j,k)} \) is a maximal independent set in \( H_{(j,k)} \), then \( c \) is an LF-coloring of \( G \).

**Proof.** Observe that if a coloring \( c \) of \( G \) fulfills the assumption of the lemma, then it is identical to the coloring \( c' \) obtained using the greedy algorithm with the largest-first sequence of vertices: \( K = (IS_{(\Delta,1)}, IS_{(\Delta,2)}, \ldots, IS_{(\Delta,\Delta + 1)}, IS_{(\Delta - 1,1)}, IS_{(\Delta - 1,2)}, \ldots, IS_{(\Delta - 1,\Delta - 1)}, IS_{(1,1)}, IS_{(1,2)}, \ldots) \), where the elements of each independent set may be enumerated in arbitrary order. Indeed, suppose that for all vertices \( u \) which appear in sequence \( K \) before some fixed vertex \( v \) both colorings are identical, \( c(u) = c'(u) \). Putting \( j = \deg_G(v) \), for all \( j_i < c(v), IS_{(j,v)} \) is a maximal independent set in \( H_{(j,v)} \), \( v \in V(H_{(j,v)}) \), and \( v \notin IS_{(j,v)} \), hence \( IS_{(j,v)} \cup \{v\} \) is not an independent set in \( G \). This means that \( v \) has at least one neighbor \( u \in IS_{(j,v)} \) which appears earlier in \( K \), and \( c'(u) = k_i \), so \( c'(v) \neq k_i \). In this way we obtain \( c'(v) \geq c(v) \). On the other hand, all vertices \( u \) which appear before \( v \) in \( K \) and have color \( c(u) = c(v) \) either belong to the same independent set \( IS_{c'(v)} \) as \( v \), or to some independent set \( IS_{c(v)} \), \( j_i > j \), and \( v \notin N(IS_{c(v)}) \). Consequently, color \( c(v) \) is the smallest legal color for \( v \) in the greedy coloring, so we have \( c'(v) = c(v) \), which completes the proof by an inductive argument.

We now propose a distributed algorithm which defines the maximal independent sets \( IS_{(j,v)} \subseteq H_{(j,v)} \). Let \( r \) be a known upper bound on the number of rounds required to compute an independent set in any subgraph of \( G \). Then the execution of the proposed algorithm is divided into \( \Delta + 1 \) steps, each of which lasts \( r \) rounds. In the \( i \)-th step, \( 1 \leq i \leq \Delta + 1 \), we compute at once all of the independent sets \( IS_{(j,v)} \) such that \( 1 \leq j \leq \Delta, 1 \leq k \leq \Delta + 1 \), and \( j - k = \Delta - i \). Thus, in step 1 we compute only a maximal independent set \( IS_{(\Delta,1)} \) in graph \( H_{(\Delta,1)} \), i.e. in the subgraph of \( G \) induced by all vertices of degree \( \Delta \). In step 2 we compute a maximal independent set \( IS_{(\Delta - 1,1)} \) in the subgraph of \( G \) induced by all vertices of degree \( \Delta - 1 \), and at the same time a maximal independent set \( IS_{(\Delta,2)} \) in the subgraph of \( G \) induced by all vertices of degree \( \Delta \) not belonging to \( IS_{(\Delta,1)} \), and so on. An illustration of the time ordering of the independent set construction is shown in Figure 2. By Lemma 10, when all vertices have been assigned to some independent set, the result can be interpreted as a distributed LF-coloring of the graph. The time complexity of the algorithm is determined by \( \Delta + 1 \) steps of an algorithm for the MIS problem, hence we have the following theorem.

**Theorem 11.** There exists a distributed \( O(\Delta \cdot T_{MIS}) \) algorithm for LF-COL, where \( T_{MIS} \) denotes a known upper bound on the execution time of an algorithm for the MIS problem in graphs of order at most \( n \) and degree at most \( \Delta \).

**Theorem 12.** There exists a distributed \( O(\sqrt{n} \cdot T_{MIS}) \) algorithm for LF-COL, where \( T_{MIS} \) denotes a known upper bound on the execution time of an algorithm for the MIS problem in graphs of order at most \( n \) and degree at most \( \Delta \).

**Proof.** Consider a partition of the vertex set \( V(G) = V_1 \cup V_2 \), where \( V_1 \) contains all vertices of degree at least \( a \), and \( V_2 \) contains all vertices of degree less than \( a \), for some value of parameter \( a \). Let \( H \) be the spanning subgraph of \( G \) with all edges having at least one endpoint of degree at least \( a \), \( H = G \setminus E(G[V_2]) \). Let \( P = (v_1 v_2 \ldots v_d) \subseteq H \) be a shortest path in \( H \) connecting some two vertices \( v_1 \) and \( v_d \). In \( P \) we must have at least \( d/4 \) vertices which belong to \( V_1 \) and whose neighborhoods are pairwise disjoint in \( H \), hence also in \( G \). Because each of these vertices is of degree at least \( a \), we obtain \( a \cdot d/4 \leq n \), which means that the diameter of each of the connected components of \( H \) is at most \( 4n/a \).

The coloring of graph \( G \) is obtained by first coloring all vertices from \( V_1 \), and then all vertices from \( V_2 \). The former stage can be performed by electing a leader in each of the connected components of \( H \), computing a centralized LF-coloring of all vertices from \( V_1 \) within this component, and disseminating the coloring to all vertices. In view of the obtained bound on the diameter, this step takes \( O(n/a) \) time. In the latter stage, the coloring of \( V_2 \) is easily completed using an approach similar to that described by Theorem 11 in \( O(a \cdot T_{MIS}) \) time, because all vertices are of degree less than \( a \) and obtain color at most \( a \). The overall complexity of the algorithm is therefore \( O(n/a + a \cdot T_{MIS}) \), and the proof of the claim is complete when we put \( a = \sqrt{n}/T_{MIS} \).

4. NOTES ON RANDOMIZED ALGORITHMS

The introduction of randomization consists of allowing local computations of vertices to involve a random coin-flip.
function, which returns a value 0 or 1 (a so-called random bit) with equal probability. We consider randomized algorithms that always stop with a correct result, whereas the time complexity of an algorithm is its expected running time.

Although randomized solutions to some problems may be significantly faster than the best known deterministic approaches (e.g. MIS in Table 1), certain lower time bounds related to the speed of dissemination of information in the network still apply. In particular, the arguments used when proving lower bounds in Section 2 are still valid: if the distance $d$-neighborhoods of a vertex $v$ in some two system graphs $G_1$ and $G_2$ containing $v$ are identical, and the sets of permissible values of color $c(v)$ are disjoint for graphs $G_1$ and $G_2$, then for any coloring algorithm, for at least one of the graphs $G_1$, $G_2$, the probability of obtaining a correct solution within $d$ rounds cannot exceed 1/2. Hence, by Markov’s inequality, the expected execution time of any algorithm is $\Omega(d)$. In this way we obtain that all the lower bounds in Table 1 also hold for expected execution times in the randomized model.

On the other hand, deterministic algorithms can also be applied in the randomized model. Hence, all upper complexity bounds in Table 1 which are expressed in terms of $\Delta$ still hold. In particular, the G-COL problem can still be solved in $2^{O(\sqrt{\log n})}$ time. Upper bounds which include bounds on the execution time of other algorithms need to be approached more carefully; we show the following simple results.

Proposition 13. There exists a distributed randomized $O(\Delta + T_{COL})$ expected time algorithm for G-COL, where $T_{COL}$ denotes the complexity of a randomized algorithm for the COL problem in $G$.

Proof. The $O(\Delta + T_{COL})$-time algorithm for G-COL given by Proposition 9 is easily adapted to allow for any randomized subroutine for the COL problem. Indeed, the initial $(\Delta + 1)$-coloring is used only to define independent sets $\{I_i : 1 \leq i \leq \Delta + 1\}$, and instead of waiting for the COL algorithm to complete, we can start the greedy coloring phase even before the independent sets are fully defined (for example, in round $(\Delta + 1)k + i$, $k \in \mathbb{N}$, we activate all vertices from set $I_i$, none of whose neighbors belong to $I_i$).

Proposition 14. There exist distributed randomized algorithms for LF-COL running in $O(\Delta \log \Delta \log n)$ and $O(\sqrt{n} \log \log n)$ expected time.

Proof. The deterministic subroutine for MIS applied in the algorithms for LF-COL (Theorems 11 and 12) can be replaced by a randomized subroutine, provided that this algorithm for MIS has the property that at every stage of execution the partial solution described by local variables $c$ of those vertices which have already completed the algorithm induces a (not necessarily maximal) independent set. The randomized distributed algorithm for MIS proposed by Luby [24] has this property. The complexity of this algorithm is given as $O(\log n)$, and the time distribution of executions of the Luby algorithm is dominated by the negative binomial distribution (a direct corollary of [24] Theorem 1). When computing an LF-coloring, the subroutine for MIS is called independently $O(\Delta^2)$ times, and thus by a simple calculation we have that the slowest of the iterations of MIS will almost certainly complete in $O(\log n \log \Delta)$ time. We may therefore use the value $O(\log n \log \Delta)$ (or $O(\log^2 n)$ for simplicity) in place of $T_{MIS}$ in Theorems 11 and 12, which completes the proof.

5. CONCLUSION

The number of colors used by most distributed algorithms for $(\Delta + 1)$-coloring, such as Johansson’s algorithm [15], is close to $\Delta$ even if the graph is bipartite. This is not surprising, because such algorithms have no mechanism for economizing on the number of colors. Distributed greedy coloring (G-COL) is a natural approach for optimizing the number of colors. We have shown that such greedy coloring can be obtained within a time bound as good as the best known $(\Delta + 1)$-coloring algorithm [27]. However, and maybe surprisingly, no polylog randomized algorithm for greedy coloring is known, unlike the case of $(\Delta + 1)$-coloring.

REFERENCES


