CS 761: Randomized Algorithms U	University of Waterloo: Winter 2025	
Lecture 1		
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Topics of this Lecture

1	Ran	ndomized Algorithms	1
2	The	e $(\Delta + 1)$ Vertex Coloring Problem	2
3	A F	Taster $(\Delta + 1)$ Coloring Algorithm	2
	3.1	The Algorithm	3
	3.2	Runtime Analysis	4

1 Randomized Algorithms

Welcome to the "Randomized Algorithms" course!

Let me start right away by saying that this is not a good title for this course; a better choice would have been "A Biased Tour of Probabilistic Ideas in Theoretical Computer Science" but that would have been too much a mouthful. So, we are going to stick with the standard name.

So, what are some probabilistic ideas in theoretical computer science? It turns out you can use randomization to design faster and/or simpler randomized algorithms (of course!), but also:

- Design and analyze deterministic algorithms (e.g., *derandomization*),
- Preserve privacy of people affected by the algorithms (e.g., *differential privacy*),
- Ensure truthfulness of people providing inputs to the algorithms (e.g. *mechanism design*),
- Prove entirely non-probabilistic mathematical facts (e.g., *probabilistic method*),
- Allow symmetry breaking and coordination between different processes (e.g., distributed algorithms),
- Analyze enormous amount of data using extremely limited space (e.g., streaming algorithms),

and many many more ways. Think of this course as providing you with basic tools for getting into these exciting areas.

We start our course by studying a recent and surprisingly fast algorithm—using randomization—for one of the oldest problems in graph theory: $(\Delta + 1)$ coloring of a graph with maximum degree Δ .

2 The $(\Delta + 1)$ Vertex Coloring Problem

Given an undirected graph G = (V, E), we use $\Delta := \Delta(G)$ to denote the maximum degree of G. For any integer $c \ge 1$, a **proper c-coloring** of G is an assignment of colors from $\{1, \ldots, c\}$ to the vertices of the graph, such that for every edge $(u, v) \in E$, the color assigned to u and v is different. For instance, think of a group of people with some enmity between some pairs of them; we would like to sit them all at a limited number of tables such that no two enemies sit at the same table. Then, we can think of the people as vertices of a graph G and have edge between any two of them iff they are enemies; then, if we have c tables in total, finding the assignment of the people to the tables is exactly the same as finding a proper c-coloring of G(which may or may not even exist). Figure 1 gives an illustration of these definitions.

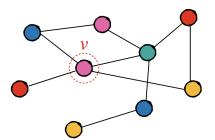


Figure 1: An example of a proper 5-coloring of a graph G with maximum degree $\Delta = 4$.

In this lecture, we are interested in finding a (proper) $(\Delta + 1)$ -coloring of any given graph. A simple observation is that every graph admits such a coloring and in fact there is a very simple greedy algorithm for finding this coloring: color vertices of the graph in some arbitrary order, and when it is vertex v's turn to be colored, check all neighbors of v that are already colored, and find a color missing from all of them. By the pigeonhole principle, since v has at most Δ neighbors but $(\Delta + 1)$ colors to choose from, we can find a missing color for v and move on (see the marked vertex v in Figure 1). Once the algorithm terminates, we will end up with a $(\Delta + 1)$ coloring of G. It is easy to see that this algorithm can be implemented in $O(n \cdot \Delta)$ time (try and prove it yourself).

But, can we find a $(\Delta + 1)$ coloring even faster than $O(n \cdot \Delta)$ time? At first glance, this seems trivially hopeless – after all, an input graph with maximum degree Δ may have $\Omega(n \cdot \Delta)$ edges and thus even reading the entire input once requires us to spend $\Omega(n \cdot \Delta)$ time. It turns out that this intuition is actually not entirely correct and one can indeed obtain even faster algorithms at least for certain ranges of Δ , by using randomization. The first algorithm to achieve this is due to Assadi, Chen, and Khanna [ACK19] in 2019 but this algorithm is quite complicated (and it actually does a lot more than just obtaining a faster $(\Delta + 1)$ coloring algorithm). In this lecture, we see a different and very simple algorithm for this problem inspired by a recent result of Assadi and Yazdanyar [AY25] (although the algorithm of [AY25] is different from this and can achieve some further properties that we will not discuss in this lecture).

3 A Faster $(\Delta + 1)$ Coloring Algorithm

Our goal is to design a *randomized* algorithm for finding a $(\Delta + 1)$ coloring. There are in general two approaches when it comes to randomized algorithms:

• Either the algorithm always outputs a correct answer and thus spend its randomization only to reduce the runtime; in this case, we are interested in the *expected runtime* of the algorithm¹. These algorithms are often called **Las Vegas** (randomized) algorithms.

¹More formally, for each input x and choice of random bits r for the algorithm, if T(x, r) denote the runtime of the algorithm on this particular input x and this choice of random bits, then we are interested in minimizing $\mathbb{E}_r[T(x, r)]$.

• The other option is to sometimes err and thus spend its randomization to obtain the correct answer with high probability². These algorithms are often called **Monte Carlo** (randomized) algorithms.

In many cases, one can also translate algorithms of one type into the other type, and we will see that later in the course.

For this lecture, we will be designing a randomized algorithm that always outputs a $(\Delta + 1)$ coloring of any given graph and its expected runtime is

$$O(\frac{n^2}{\Delta} \cdot \log n).$$

Notice that whenever $\Delta = \omega(\sqrt{n \log n})$, the runtime of this algorithm will be $o(n \cdot \Delta)$ and thus (possibly) even faster than reading the entire input. In general, we can combine this algorithm and the standard greedy algorithm to obtain an algorithm for $(\Delta + 1)$ coloring that runs in expected $O(n \cdot \sqrt{n \log n})$ time³- for any graph with more than these many edges, this algorithm runs in *sublinear* time! The rest of this lecture is dedicated to the presentation and analysis of this algorithm.

3.1 The Algorithm

The new algorithm also follows the approach of the greedy algorithm quite closely. It will iterate over the vertices in <u>some order</u> (to be determined later). For every vertex v, as in the greedy algorithm, it attempts to find a color to assign to v and will not change its decision afterwards. The difference between this algorithm and the standard greedy algorithm is here: instead of going over all neighbors of v to find a missing color for v, the new algorithm picks a color $c \in [\Delta + 1]$ uniformly at random for v and then go over all vertices that are colored c to see if they are neighbor to v; if not, v will be colored c and the algorithm moves on, otherwise, it simply picks another random color (again, from all of $[\Delta + 1]$ without even discarding "bad" colors) and continue like this until it can actually color v.

Algorithm 1.

- 1. Let $C_1, C_2, \ldots, C_{\Delta+1} = \emptyset$ (we maintain the invariant that C_c always contains vertices colored c).
- 2. Iterate over vertices in some order (to be determined later); for each vertex v in this order:
 - (a) Sample a color $c \in [\Delta + 1]$ uniformly at random;
 - (b) For every vertex $u \in C_c$, i.e., colored c so far, check if u is a neighbor of v;
 - (c) If none of C_c is neighbor to v, color v with c, i.e., update $C_c \leftarrow C_c \cup \{v\}$; otherwise, go to the sampling step again.

We can easily observe that if the algorithm does indeed terminate, then the coloring it finds is a proper $(\Delta+1)$ -coloring of the input graph. Any new vertex colored does not create a conflict with previously colored vertices (given the algorithm explicitly checks to not color v with a color c if one of its neighbors is already colored c) and thus at the end, there cannot be any monochromatic edge in the graph. As such, the only actual part of the analysis is to figure out how long this algorithm takes (in expectation). Before getting to this however, we need to take care of an important issue.

Representation of the input. Graphs are typically presented to algorithms as adjacency list or adjacency matrix. In many cases, the choice of representation does not matter much as we can simply switch between them without much extra cost by simply reading the input once (especially going from adjacency list to adjacency matrix is quite straightforward). Yet, in our setting, we are interested in an algorithm that aims

²More formally, for each input x and choice of random bits r, if $O(x, r) \in \{\text{`correct'}, \text{`incorrect'}\}\$ denote whether the output of the algorithm on this particular input and this choice of random bits is correct or not, then, we are in interested in maximizing $\Pr_r(O(x, r) = \text{`correct'})$.

³Basically, run this algorithm when $\Delta = \omega(\sqrt{n \log n})$ and otherwise run the original greedy algorithm.

to solve the problem faster than even reading the input once. As such, the representation of the input actually matters in our case. So, what does our algorithm need? It is easy to see that the only access of the algorithm to the graph is in Line (2b): it needs to check for a given vertex v and another vertex $u \in C_c$, whether or not (u, v) is an edge in the graph. For this purpose, having access to the adjacency *matrix* is more helpful as it immediately allows for implementing such a *query* to the input in O(1) time⁴. Thus, for the rest of this lecture, we assume the input is given via an adjacency matrix.

3.2 Runtime Analysis

For every vertex $v \in V$, define a random variable X_v , which is equal to the total number of entries of the adjacency matrix queried by Algorithm 1 in Line (2b) when attempting to color the vertex v. Basically, if the algorithm "tries" colors c_1, c_2, \ldots, c_k before it finds a color for v, then $X_v = \sum_{i=1}^k |C_{c_i}|$. Note that it is possible for X_v to be even infinity (although the probability if this event is zero). We have,

Runtime of Algorithm
$$1 = O(1) \cdot \sum_{v \in V} X_v.$$
 (1)

This is simply because the only real time consuming step of the algorithm is this step and all the rest only take O(n) time (deterministically) which is suppressed in the above asymptotic runtime. Thus, our goal is now to upper bound

$$\mathbb{E}\left[\text{Runtime of Algorithm 1}\right] = \mathbb{E}\left[O(1) \cdot \sum_{v \in V} X_v\right] = O(1) \sum_{v \in V} \mathbb{E}\left[X_v\right].$$

Here, the first step is by Eq (1) and the second one is by *linearity of expectation*. Thus, our task is now to simply bound $\mathbb{E}[X_v]$ for any given vertex $v \in V$.

Fix a vertex v and define deg[<](v) as the number of neighbors of v that appear before v in the ordering of Algorithm 1. In other words, deg[<](v) counts the neighbors of v that already received a color before coloring v. We claim the following.

Claim 1. For every $v \in V$

$$\mathbb{E}\left[X_{v}\right] \leqslant \frac{n}{\Delta + 1 - \deg^{<}(v)}.$$

Before proving this (simple) claim, let us start with two warm-up questions that provide more intuition about the algorithm and its analysis.

• Probability a random color is "good" for v: What is the probability that when we sample a color for the vertex v, it can be used to color v, i.e., c is good for v? Well, this color should not be the color of one of the already colored neighbors of v which are $\deg^{<}(v)$ many. Thus, even if they all receive a different color, we have,

 $\Pr(c \text{ is } \underline{\text{not}} \text{ used in the neighborhood of } v) = 1 - \Pr(c \text{ is used in the neighborhood of } v)$

$$\ge 1 - \frac{\deg^{<}(v)}{\Delta + 1} = \frac{\Delta + 1 - \deg^{<}(v)}{\Delta + 1}$$

• Expected number of queries for a single random color c: Another question is that when we sample a single color $c \in [\Delta + 1]$, how many queries the algorithm needs to do (in expectation) for just this single color, regardless of whether or not this color is "good" for v. This can be easily calculated by the definition of expected value as

$$\frac{1}{\Delta+1}\sum_{c=1}^{\Delta+1}|C_c|\leqslant \frac{n}{\Delta+1},$$

⁴Although this is not the only way; for instance, if the input is stored in a way that neighbors of every vertex is stored in a dedicated hash table (or a balanced search trees), we can obtain almost (but not exactly) the same guarantees.

since picking a color c means making $|C_c|$ queries, and the sets $C_1, \ldots, C_{\Delta+1}$ form a partition of (a subset of) vertices of the graph at every point (since every vertex receives a single color).

We now use the same ideas to prove Claim 1 also.

Proof of Claim 1. By the law of conditional expectations, we have,

$$\mathbb{E}\left[X_{v}\right] = \sum_{c=1}^{\Delta+1} \Pr\left(\text{first sampled color for } v \text{ was } c\right) \cdot \mathbb{E}\left[X_{v} \mid \text{first sampled color for } v \text{ was } c\right].$$

Now suppose we know the first sampled color for v was c, what does it tell us about X_v (conditioned on this event)? Well, first of all, we know that X_v involves making $|C_c|$ queries; then, if c is good for v, i.e., can be used to color v, then we are done. In other words, for every c which is good for v, we have,

 $\mathbb{E}\left[X_v \mid \text{first sampled color for } v \text{ was } c\right] = |C_c|.$

But, what if c is not good for v? Then, we spend $|C_c|$ queries and are in exactly the same place as we were when we started (since the algorithm simply does a "goto step" to the beginning of coloring v). Thus, for every c which is not good for v,

 $\mathbb{E}\left[X_{v} \mid \text{first sampled color for } v \text{ was } c\right] = \mathbb{E}\left[|C_{c}| + X_{v}\right] = |C_{c}| + \mathbb{E}\left[X_{v}\right].$

But now, letting B(v) denote the set of colors that are not good for v and noting that $|B(v)| \leq \deg^{<}(v)$ (for the same reason as the one when calculating probability a color is not good for v), we have,

$$\mathbb{E}\left[X_{v}\right] = \sum_{c=1}^{\Delta+1} \Pr\left(\text{first sampled color for } v \text{ was } c\right) \cdot \mathbb{E}\left[X_{v} \mid \text{first sampled color for } v \text{ was } c\right]$$
$$= \sum_{c=1}^{\Delta+1} \frac{1}{\Delta+1} \cdot |C_{c}| + \sum_{c \in B(v)} \frac{1}{\Delta+1} \cdot \mathbb{E}\left[X_{v}\right]$$
$$(\text{since the given probability is } 1/\Delta + 1 \text{ and by the discussions above})$$

$$\leq \frac{n}{\Delta+1} + \frac{\deg^{<}(v)}{\Delta+1} \cdot \mathbb{E}[X_v]$$

Moving $\mathbb{E}[X_v]$ to the LHS in the equation above and simplifying the terms imply the result.

At this point, we have proven that

$$\mathbb{E}\left[\text{Runtime of Algorithm 1}\right] = O(1) \cdot \sum_{v \in V} \frac{n}{\Delta + 1 - \deg^{<}(v)}.$$
(2)

To conclude the runtime calculation, we need to bound this RHS. The challenge however is that $\deg^{<}(v)$ can be quite different for different vertices and in fact, the value of this sum can change dramatically depending on the ordering of vertices in the algorithm. This is where we have to fix the notion of picking vertices in <u>some order</u> stated earlier in Algorithm 1.

Fixing the ordering of vertices in Algorithm 1. The solution here is quite simple: we will also pick the ordering of the vertices uniformly at random. Let π denote an ordering of vertices. Eq (2) shows that no matter the choice of π , the runtime of the algorithm can be bounded by the given expression. although now that ordering of vertices is also random, deg[<](v) is a random variable that depends on π ; we denote this by deg[<]_{π}(v)_{π} for more clarity. Thus, by taking the randomness of π into account, we will have

 \mathbb{E} [Runtime of Algorithm 1] = $\mathbb{E}_{\pi} \mathbb{E}$ [Runtime of Algorithm 1 | π] (by the law of conditional probabilities)

$$= O(1) \cdot \sum_{v \in V} \mathbb{E}\left[\frac{n}{\Delta + 1 - \deg_{\pi}^{<}(v)}\right]. \quad (by Eq (2) and linearity of expectation)$$

The final step is to bound this RHS for every vertex $v \in V$, which is done in the following claim.

Claim 2. For every vertex $v \in V$,

$$\mathbb{E}_{\pi}\left[\frac{n}{\Delta+1-\deg_{\pi}^{<}(v)}\right] \leqslant \frac{n}{\Delta+1} \cdot \left(\ln\left(\Delta+1\right)+1\right).$$

Proof. Let deg(v) denote the degree of v in the entire graph. By picking π randomly, we also obtain that the ordering of the vertices $v \cup N(v)$ (where N(v) are neighbors of v) is chosen uniformly at random. In particular, the relative order of v among $v \cup N(v)$ is chosen uniformly. As such, deg[<]_{π}(v) is chosen uniformly at random from $\{0, 1, \ldots, \deg(v)\}$. Hence,

$$\mathbb{E}_{\pi}\left[\frac{n}{\Delta+1-\deg_{\pi}^{<}(v)}\right] = \sum_{d=0}^{\deg(v)} \frac{1}{\deg(v)+1} \cdot \frac{n}{\Delta+1-d}.$$
 (by the discussion above)

Note that for any choice of deg(v), we have,

$$\sum_{d=0}^{\deg(v)} \frac{1}{\deg(v)+1} \cdot \frac{n}{\Delta+1-d} \leqslant \sum_{d=0}^{\Delta} \frac{1}{\Delta+1} \cdot \frac{n}{\Delta+1-d},$$

because if we let $A := \{n/(\Delta + 1 - i)\}_{i=0}^{\Delta}$, then, in the LHS, we are taking the average of the smallest $\deg(v) + 1$ numbers in A, whereas in the RHS we are taking the average of all of A. Continuing the above then we have,

$$\sum_{d=0}^{\Delta} \frac{1}{\Delta+1} \cdot \frac{n}{\Delta+1-d} = \frac{n}{\Delta+1} \cdot \sum_{i=1}^{\Delta+1} \frac{1}{i}$$
 (by a change of variable)
$$\leqslant \frac{n}{\Delta+1} \cdot \left(\ln\left(\Delta+1\right)+1\right),$$

where we used the standard inequality that $\sum_{i=1}^{K} 1/i \leq \ln(K) + 1$ (the series is called the Harmonic series). Putting everything together, we have,

$$\mathbb{E}_{\pi}\left[\frac{n}{\Delta+1-\deg_{\pi}^{<}(v)}\right] \leqslant \frac{n}{\Delta+1} \cdot \left(\ln\left(\Delta+1\right)+1\right),$$

concluding the proof.

Plugging in all the bounds we proved so far, we have that

$$\mathbb{E}\left[\text{Runtime of Algorithm 1}\right] = \sum_{v \in V} \mathbb{E}\left[X_v\right] \leqslant \sum_{v \in V} \frac{n}{\Delta + 1} \cdot \ln\left(\Delta + 1\right) = O(\frac{n^2}{\Delta} \cdot \log\left(\Delta\right)),$$

proving the desired upper bound on the expected runtime of Algorithm 1. Thus, at this point, we saw an algorithm and its full analysis for finding a $(\Delta + 1)$ coloring in *sublinear* time (for certain ranges of Δ).

References

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