

Lecture 9

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1 Last time

Last time we talked about random walks on graphs, and in particular random walks on *undirected* graphs. We saw that performing a random walk on an undirected graph would solve the undirected $s-t$ connectivity problem in randomized logspace. We also saw that the time it takes for a random walk on a graph (that's undirected, d -regular and ergodic) to approach the stationary distribution (which is the uniform distribution over the nodes) within ϵ in ℓ_2 norm takes time about $\log \epsilon / \log \lambda_2$.

2 Today

Today we will see how random walks can be used to save randomness in randomized algorithms while attaining very low error. Suppose we had a polynomial time randomized algorithm \mathcal{A} for a language $L \in \text{RP}$ with the following property:

- $x \in L$: $\Pr_w[\mathcal{A}(x, w) \text{ doesn't accept}] \leq 1/100$, and
- $x \notin L$: $\Pr_w[\mathcal{A}(x, w) \text{ accepts}] = 0$, and
- \mathcal{A} uses $r(n)$ bits of randomness.

Previously, we saw two ways in which the error of this algorithm could be reduced to 2^{-k} :

1. Run \mathcal{A} k times, using new randomness for each repetition. This uses $r \cdot k$ random bits, and has running time $k \cdot T(n)$ (where $T(n)$ is the running time of \mathcal{A} ; or
2. Run \mathcal{A} on pairwise independent randomness. This uses $O(r+k)$ random bits, and has running time $O(2^k \cdot T(n))$.

We will describe a new method that combines the time efficiency of Method 1 with the randomness efficiency of Method 2, and this will be done via random walks on a special type of graph, called an *expander graph*.

3 A Special Kind of Graph

We will postulate that we can construct a family of undirected graphs $G_n = (V_n, E_n)$ with the following properties:

1. G_n has constant degree (i.e. there is a constant d such that the maximum degree of G_n is bounded by d);
2. G_n is regular;
3. $\lambda_2 \leq 1/10$ (where λ_2 is the second largest eigenvalue of the transition matrix corresponding to G_n);
4. G_n is ergodic (this is redundant, because of the λ_2 property above);

5. $|V_n| = 2^{r(n)}$;
6. Given a node $v \in V_n$, the neighbors of v can be computed efficiently (in n).

It turns out that it is possible to construct a graph family with all these magical properties: such constructions are called *expander graph* constructions. Here, we will take the construction of these graphs for granted. Importantly, *any* expander graph construction that has the above properties will suffice for the error reduction procedure we describe next.

4 Error reduction for \mathcal{A}

We describe the algorithm to reduce the error of \mathcal{A} :

- **Algorithm $\mathcal{A}'(x)$:**
 - Pick a random node w in G_n .
 - Repeat k times:
 - * Run $\mathcal{A}(x, w)$. If \mathcal{A} accepts, then halt and accept.
 - * Otherwise, pick a random neighbor of w and set w equal to that neighbor.
 - If the algorithm has not accepted by this time, reject.

Clearly, the running time of \mathcal{A}' is $O(k(T(n) + Q(n)))$, where $T(n)$ is the running time of \mathcal{A} and $Q(n)$ is the time it takes to pick a random neighbor in G_n . We assume here that $Q(n)$ is negligible compared to $T(n)$.

The randomness efficiency is $r(n) + O(k)$: $r(n)$ bits of randomness are required to pick the initial start node w . Then, for each iteration of the loop, a random neighbor has to be picked. Since each node has d neighbors and d is constant, each step of the random walk requires $\log d = O(1)$ random bits.

We now argue that the error of \mathcal{A}' is very low:

Claim 1 *The error probability of \mathcal{A}' is at most $\frac{1}{5^k}$.*

Before diving into the proof, let us spend a few moments to reflect on this algorithm. We have essentially applied a graph structure to the space of all possible random strings that could be used by the original algorithm \mathcal{A} . For a fixed input x , the set of *bad* random strings B_x (random strings that cause \mathcal{A} to err on x) contains at most 1/100 of all the possible nodes. Because of the one-sided error property of \mathcal{A} , we only need to encounter just one *good* random string in order to succeed.

Our algorithm \mathcal{A}' will take a random walk on this graph of random strings, and because of the λ_2 condition stipulated above, this graph has very good *mixing properties*: it is very unlikely for a random walk to be trapped within B_x . The following proof formalizes this intuition.

Proof Suppose input $x \notin L$. Then because of the soundness property of \mathcal{A} , \mathcal{A} will never accept x , no matter the randomness. Thus \mathcal{A}' will never accept on x , and hence the error is 0.

Now suppose $x \in L$. Define N to be the diagonal matrix where $N_{ww} = 1$ if and only if $w \in B_x$. Otherwise, $N_{ww} = 0$ (and 0 everywhere off the diagonal). Let q be some distribution on nodes of G_n . Then $\|q \cdot N\|_1$ is precisely the probability of drawing a bad string from q .

Let P be the transition matrix corresponding to G_n . Then observe that $\|qNP\|_1$ is the probability of drawing a bad string from q , and that taking a random walk with starting distribution q will also yield a bad string. More generally, $\|q(NP)^k\|_1$ is the probability that the first $k + 1$ steps of the random walk yield a bad string.

We will use the following lemmas, which we will prove later, to arrive at our desired claim:

Lemma 2 For all π such that $\sum \pi_i \leq 1$ and $\pi_i \geq 0$ for all i , $\|\pi PN\|_2 \leq \frac{1}{5} \|\pi\|_2$.

Lemma 3 For all $x \in \mathbb{R}^m$, $\|x\|_1 \leq 2^{m/2} \|x\|_2$.

\mathcal{A}' starts the random walk with the uniform distribution π_0 . The probability that after k iterations, \mathcal{A}' does not accept is precisely $\|\pi_0(NP)^k N\|_1$. This is bounded from above by $\|\pi_0(PN)^k\|_1$, where we have ignored the first N factor. This is because without the first N terms we are now allowing the initial choice of random string to fall outside the bad set, which means there are more ways of arriving at the bad set after k steps. Hence,

$$\begin{aligned} \Pr[\mathcal{A}' \text{ does not accept}] &\leq \|\pi_0(PN)^k\|_1 \\ &\leq 2^{r/2} \|\pi_0(PN)^k\|_2 \\ &\leq 2^{r/2} \frac{1}{5^k} \|\pi_0\| \\ &\leq \frac{1}{5^k}. \end{aligned}$$

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Proof (Of Lemma 2). Since G_n is d -regular, we have that P is symmetric and hence diagonalizable. Let $\{v_i\}$ be the orthonormal basis in which P is diagonal, such that $1 = \lambda_1 > |\lambda_2| \geq \dots \geq |\lambda_{2r}|$. Let $\pi = \sum \alpha_i v_i$. Then,

$$\begin{aligned} \|\pi PN\|_2 &= \left\| \alpha_1 v_1 PN + \sum_{i>2} \alpha_i v_i PN \right\|_2 \\ &\leq \|\alpha_1 \lambda_1 v_1 N\|_2 + \left\| \sum_{i>2} \alpha_i \lambda_i v_i N \right\|_2 \end{aligned}$$

where the second line follows from the triangle inequality. We bound each term separately. Since λ_1 , and $v_1 = \frac{1}{2^{r/2}}(1, 1, \dots, 1)$, we have that $\|\alpha_1 \lambda_1 v_1 N\|_2 = \alpha_1 \|v_1 N\|_2 \leq \alpha_1 \sqrt{1/100} \leq \frac{1}{10} \|\pi\|_2$.

For the second term, we note that $\|\sum \alpha_i \lambda_i v_i N\|_2 \leq |\lambda_2| \|\sum \alpha_i v_i N\|_2 \leq \frac{1}{10} \|\pi\|_2$, because multiplying by N does not grow the length of a vector. Combining the two bounds yields the lemma.

■

Proof (Of Lemma 3). This follows from applying Cauchy-Schwarz: $(\sum x_i)^2 = (\sum 1 \cdot x_i)^2 \leq (\sum 1^2) (\sum x_i^2) = 2^m (\sum x_i^2)$. ■

This (temporarily) concludes the section on random walks on graphs. Next, we turn towards the topics of Fourier Analysis and Linearity Testing. We end this class with a prelude of program checking.

5 Program Checking and Linearity Testing

Suppose you can't prove that a program P correctly computes $f(x)$ on all inputs x , but you don't care – you just want to verify that it's correct on a *specific* input. Is it possible to create a *program checker* C that will do this for you? We'd want a program that would tell us if $P(x)$ is correct or not. Of course, the checker itself might be faulty, but what we want is for the checker to be faulty in a way that's *independent* of the way P is faulty. A program that admits such a program checker is called *checkable*. What kind of programs are checkable?

One class of checkable programs are ones computing functions that are *close to linear*. Let's see some definitions.

Definition 4 (Linear function) A function $f : \{0, 1\}^n \rightarrow \{0, 1\}$ is linear if and only if for all $x, y \in \{0, 1\}^n$, $f(x + y) = f(x) + f(y)$, where addition is performed component-wise modulo 2.

Definition 5 (ϵ -close to linear) A function $g : \{0, 1\}^n \rightarrow \{0, 1\}$ is ϵ -close to linear if and only if there exists a linear function f such that $\Pr_x[g(x) = f(x)] \geq 1 - \epsilon$.

We describe one component of a program checker, called a *self-corrector*, for programs that compute linear functions. Suppose we have a program P computing a function g that is ϵ -close to a linear function f , and we intended P to compute f . We can *correct* P so that it computes f exactly.

- **Program Self-Corrector**(x)
 - For $O(\frac{1}{\epsilon})$ iterations:
 - * Pick $y \in \{0, 1\}^n$ uniformly at random.
 - * Let $\alpha \leftarrow P(x + y) + P(y)$.
 - Output the majority of the α 's.

Claim 6 *With very high probability, Self-Corrector will compute f .*

Proof Note that at each iteration, y is chosen uniformly at random, so thus $x + y$ and y are uniformly distributed. Thus, with probability at least $1 - 2\epsilon$, $P(x + y) = f(x + y)$ and $P(y) = f(y)$. By the linearity of f , $P(x + y) + P(y) = f(x + y) + f(y) = f(x)$. Since we're picking y independently with each iteration, the probability that the majority of the α 's agree with $f(x)$ rapidly approaches 1. Thus, with high probability, **Self-Corrector** will output the correct value of $f(x)$. ■

Generally, program checkers will combine self correctors with program “testers”, which we still study next time.