

## Lecture 15: Expanders

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In the last lecture we introduced randomized computation in terms of machines that have access to a source of random bits and that return correct answers somewhat more than half of the time. We showed that BPP has polynomial-size circuits and the conjecture in the community is that  $\text{BPP} = \text{P}$ .

Today we'll introduce expanders, a type of graph that is useful in improving (amplifying) randomized algorithms with little or no additional random bit overhead. First we'll prove a theorem relating BPP to the polynomial time hierarchy.

## 1 BPP and the Polynomial Time Hierarchy

Although we don't know if  $\text{BPP} = \text{P}$ , or even if  $\text{BPP} \subseteq \text{NP}$ , we do know that  $\text{BPP} \subseteq \text{PH}$ :

**Theorem 1.**  $\text{BPP} \subseteq \Sigma_2^p \cap \Pi_2^p$

*Proof.* Fix  $M$  a randomized polytime machine that accepts  $L \in \text{BPP}$ , and let  $r$  be the number of random bits  $M$  uses when running on an input  $x$  of size  $n = |x|$ . We will now show that  $\text{BPP} \subseteq \Sigma_2^p$ . Since  $\text{BPP} = \text{coBPP}$  it follows that  $\text{BPP} \subseteq \text{co}\Sigma_2^p = \Pi_2^p$ , completing the proof.

We need to remove randomness from the computation. For a given input  $x$ , the space  $\{0, 1\}^r$  of  $r$ -bit strings gets partitioned into two pieces:  $\text{Acc}(x)$ , the set of random strings on which  $M$  accepts  $x$ , and  $\text{Rej}(x)$ , the set of strings on which  $M$  rejects  $x$ . If the error rate  $\varepsilon$  of  $M$  is small, then  $\text{Acc}(x)$  will be much larger than  $\text{Rej}(x)$  when  $x \in L$ , and  $\text{Acc}(x)$  will be much smaller than  $\text{Rej}(x)$  when  $x \notin L$ . See Figure 1. If our random computation has a small error rate then it will be correct on most random bit strings. We'll turn "most" into "all." The idea is that when  $x \in L$  a few "shifts" of  $\text{Acc}(x)$  will cover the whole space  $\{0, 1\}^r$  of  $r$ -bit random sequences, while the same few shifts of  $\text{Acc}(x)$  will fail to cover the whole space when  $x \notin L$ .

If  $S \subseteq \{0, 1\}^r$  and  $\sigma \in \{0, 1\}^r$ , then  $S \oplus \sigma = \{s \oplus \sigma \mid s \in S\}$ , the *shift* of  $S$  by  $\sigma$ . Since shifting is invertible we see that  $|S \oplus \sigma| = |S|$ .

We will use shifts to give a  $\Sigma_2$ -predicate using the intuition discussed above. Namely, consider

$$x \in L \iff \exists \sigma_1, \dots, \sigma_t \forall \rho \in \{0, 1\}^r [\rho \in \bigcup_{i=1}^t \text{Acc}(x) \oplus \sigma_i]. \quad (1)$$

Now,  $r$  is poly in  $n$ , and so if we can pick  $t$  poly in  $n$  as well, then the above will be a  $\Sigma_2^p$ -predicate, provided that we can verify the membership of  $\rho$  in  $\bigcup_{i=1}^t \text{Acc}(x) \oplus \sigma_i$  in time poly in  $n$ . The membership check is no problem since we can equivalently check that  $\rho \oplus \sigma_i \in \text{Acc}(x)$  for some  $i$ , which is polytime since it corresponds to running  $M$  on  $x$  with the random bit string  $\rho \oplus \sigma_i$ , for poly-many  $\sigma_i$ .

We'll show we can pick  $t$  a suitable poly to make (1) true by showing that we can choose the  $\sigma_i$ s randomly with a high rate of success. There are two cases to consider:

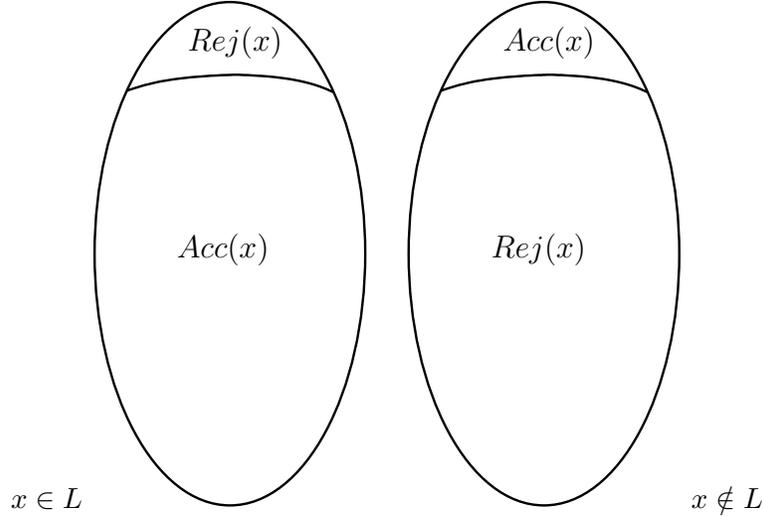


Figure 1: If the error rate  $\varepsilon$  of  $M$  is small, then  $Acc(x)$  will be much larger than  $Rej(x)$  when  $x \in L$ , and  $Acc(x)$  will be much smaller than  $Rej(x)$  when  $x \notin L$ .

1.  $x \in L$ : For  $A \subseteq \{0, 1\}^r$ , define

$$\mu(A) = \frac{|A|}{2^r}.$$

For  $\rho$  fixed,

$$\Pr_{\sigma_1, \dots, \sigma_t} [\rho \notin \bigcup_{i=1}^t Acc(x) \oplus \sigma_i] \leq (\mu(Rej(x)))^t \leq \varepsilon^t,$$

where  $\mu(Rej(x))$  is the probability that  $M$  rejects when it should accept, and is hence not larger than  $\varepsilon$ . So, by an union bound,

$$\Pr[\{0, 1\}^r \not\subseteq \bigcup_{i=1}^t Acc(x) \oplus \sigma_i] \leq |\{0, 1\}^r| \varepsilon^t = 2^r \varepsilon^t,$$

and hence if  $2^r \varepsilon^t < 1$  then some choice of the  $\sigma_i$ s must work.

2.  $x \notin L$ : We need to be sure that for any choice of the  $\sigma_i$ s that none of the shifts of  $\rho$  make  $M$  accept  $x$ . But

$$\mu\left(\bigcup_{i=1}^t Acc(x) \oplus \sigma_i\right) \leq \sum_{i=1}^t \mu(Acc(x) \oplus \sigma_i) = t\varepsilon,$$

since  $\mu(Acc(x)) \leq \varepsilon$  when  $x \notin L$ , and hence we need  $t\varepsilon < 1$ .

So, we need to choose  $t$  so that both  $t\varepsilon$  and  $2^r \varepsilon^t$  are less than 1. So  $t = r$  works, provided that  $\varepsilon < \frac{1}{r}$ . From the definition of BPP we only know that  $\varepsilon < \frac{1}{3}$ , but, using the “majority vote” trick introduced during the last lecture, we can in  $k$  runs reduce the error to  $e^{-2k\delta^2}$ , where  $\delta = \frac{1}{2} - \varepsilon > 0$ . Since  $k$  runs will need  $kr$  random bits, we see that it suffices to choose  $k$  poly and large enough that  $e^{-2k\delta^2} < \frac{1}{rk}$ . In fact,  $k = O(\log r)$  works, and so an examination of the above shows that the time complexity is  $O(T^2 \log T)$  if the original run-time was  $T$ .

□

## 2 Expanders

The proof of Theorem 1 used the “majority vote” trick to get an exponential in  $k$  increase in accuracy using a linear in  $k$  increase in random bit usage. Expander graphs, which are introduced here, lead to an “amplification” technique that gives an exponential in  $k$  accuracy improvement using only a constant increase in random bit usage ( $rk$  versus  $r + k$ ). Expanders have many other uses, some of which we mention in this lecture, and others we may see in later lectures.

### 2.1 Definitions

Here we introduce two definitions of expander graphs, namely the combinatorial definition and the algebraic definition.

**Definition 1** (Combinatorial definition). *A graph  $G = (V, E)$  is a  $(k, c)$ -expander if  $\forall S \subseteq V$  with  $|S| \leq k$  satisfies  $|\Gamma(S)| \geq c|S|$ , where  $\Gamma(S) = \{v \in V \mid \exists s \in S, (s, v) \in E\}$  is the neighborhood of  $S$  in  $G$ .*

Notice that it is easy to construct a graph which is  $(k, c)$ -expanding for all  $k$  with  $c \leq 1$ . Since  $\Gamma(V) = V$  one can see it is important not to let  $|S|$  be too large. In other words, for  $k \geq |V|$  and  $c > 1$ , no graph is  $(k, c)$ -expanding. A typical setting is  $c$  is a constant  $> 1$  and  $k = \frac{N}{2}$  where  $N = |V|$ .

**Definition 2** (Algebraic definition). *Given a regular  $G = (V, E)$  with degree  $d$  its  $N \times N$  normalized adjacency matrix  $A$  is defined by*

$$A_{ij} = \begin{cases} \frac{1}{d}, & (i, j) \in E \\ 0, & \text{otherwise} \end{cases}$$

The normalized adjacency matrix describes the Markov chain of a random walk on  $G$ :

$$A_{ij} = \Pr[\text{go to state } i \text{ from state } j],$$

and if  $p$  is a column vector with  $p_i$  the probability of being at vertex  $i$ , then  $(Ap)_i$  gives the probability of being at vertex  $i$  after one step of the random walk on  $G$ .

$$(Ap)_i = \sum_{j=1}^N A_{ij} p_j.$$

### 2.2 Properties of the normalized adjacency matrix of a regular graph

The normalized adjacency matrix has a number of properties that can be proved using basic linear algebra. We will use the following properties, but we do not prove them here.

**Proposition 1.** *Let  $A$  be a normalized adjacency matrix of a regular graph  $G$  on  $N$  vertices.*

1.  *$A$  is real and symmetric: all eigenvalues of  $A$  are real, and there exists a full orthonormal basis of eigenvectors.*
2. *All eigenvalues  $\lambda$  of  $A$  satisfy  $|\lambda| \leq 1$ .*

3. 1 is an eigenvalue of  $A$ , with corresponding eigenvector  $u = (\frac{1}{N}, \frac{1}{N}, \dots, \frac{1}{N})$ . The multiplicity of 1 as an eigenvalue equals the number of components of  $G$ .
4. -1 is an eigenvalue of  $A$  iff  $G$  is bipartite.

As the uniform distribution is always an eigenvector with eigenvalue 1, we look only at the remaining eigenvector/eigenvalues. It turns out that the second largest eigenvalue in absolute value is often useful to work with.

**Definition 3.**  $\lambda(G) = \max\{|\lambda| \mid \lambda \text{ is an eigenvalue of } G \text{ with eigenvector } e_\lambda \text{ orthogonal to } u\}$

From Proposition 1 we have  $\lambda(G) \leq 1$ , and

$$\lambda(G) < 1 \iff G \text{ is connected and not bipartite.}$$

So, for expanders, the uniform distribution  $u$  is the only fixed point. The following theorem tells us that the larger the difference between 1 and  $\lambda(G)$  the faster an arbitrary probability distribution converges to the uniform distribution via a random walk on  $G$ .

**Theorem 2.** For any probability distribution vector  $p$ ,

$$\|A^t p - u\|_1 \leq \sqrt{N} \lambda^t,$$

where  $\|v\|_q = [\sum_i |v_i|^q]^{1/q}$  is the  $q$ -norm of  $v$  and  $\lambda = \lambda(G)$ .

One way to understand this theorem is: given an initial distribution for a random walk  $p$ , after  $t$  steps of the random walk  $A^t p$ , we will be exponentially closer to the uniform distribution  $u$ , provided  $\lambda < 1$ , that is  $G$  is connected and not bipartite.

*Proof.* Since  $Au = u$  we have  $A^t p - u = A^t(p - u)$ . Now,  $(p - u) \perp u$  because

$$\langle p - u, u \rangle = \langle p, u \rangle - \langle u, u \rangle = \sum_{i=1}^N p_i / N - \sum_{i=1}^N 1 / N^2 = 1/N - 1/N,$$

which follows since  $p$  is a probability distribution.

Write  $p - u = \sum_i a_i e_i$  for some real  $a_i$ , where the  $\{e_i\}$  form an orthogonal basis of eigenvectors for  $A$  with  $Ae_i = \lambda_i e_i$ . Then, since the  $u$ -component of  $p - u$  is 0, we have the following inequality

$$\begin{aligned} \|A^t(p - u)\|_2^2 &= \left\| A^t \sum_i a_i e_i \right\|_2^2 = \left\| \sum_i \lambda_i^t a_i e_i \right\|_2^2 = \sum_i \|\lambda_i^t a_i e_i\|_2^2 = \sum_i |\lambda_i|^{2t} \|a_i e_i\|_2^2 \\ &\leq \sum_i \lambda^{2t} \|a_i e_i\|_2^2 = \lambda^{2t} \sum_i \|a_i e_i\|_2^2 = \lambda^{2t} \|p - u\|_2^2. \end{aligned}$$

Now

$$\|p - u\|_2^2 + \|u\|_2^2 = \|p\|_2^2 \leq \|p\|_1^2 = 1,$$

where  $(p - u) \perp u$  implies the first equality, and  $\|v\|_2^2 \leq \|v\|_1^2$  for all vectors  $v$  implies the inequality. By the Cauchy-Schwartz inequality, we have

$$\begin{aligned} \|A^t(p - u)\|_1 &= |\langle (-1^{\sigma_1}, \dots, -1^{\sigma_N}), A^t(p - u) \rangle| \\ &\leq \|(-1^{\sigma_1}, \dots, -1^{\sigma_N})\|_2 \|A^t(p - u)\|_2 \\ &= \sqrt{N} \|A^t(p - u)\|_2, \end{aligned}$$

where  $\sigma_k = \begin{cases} 0, & (A^t(p - u))_k \geq 0 \\ 1, & \text{otherwise} \end{cases}$ . Combining all of the above we get

$$\begin{aligned} \|A^t(p - u)\|_1 &\leq \sqrt{N} \|A^t(p - u)\|_2 \\ &\leq \sqrt{N} \lambda^t \|p - u\|_2 \\ &\leq \sqrt{N} \lambda^t, \end{aligned}$$

completing the proof. □

This can be used to prove that the random walk algorithm for the undirected path problem needs only polynomially many steps, i.e. that  $\text{PATH} \in \text{BPL}$ . The proof uses Theorem 2 and the following exercise.

**Exercise 1.** *If  $G$  is connected and not bipartite then  $\lambda(G) < 1 - \frac{1}{dN^2}$ .*

### 3 Next Time

In the next lecture we will prove expander mixing lemma.

**Lemma 1** (Expander Mixing Lemma). *For every pair of subsets  $S, T \subseteq V$ ,*

$$\left| \frac{|E(S, T)|}{dN} - \mu(S)\mu(T) \right| \leq \lambda \sqrt{\mu(S)(1 - \mu(S))\mu(T)(1 - \mu(T))}.$$

This lemma can be used in the following applications:

1. *Deterministic error reduction:* Given a randomized algorithm  $R$  that uses  $r$  random bits we reduce the error to be less than some arbitrary  $\varepsilon$ , without using any additional random bits. This requires running  $R$   $\text{poly}(1/\varepsilon)$  times.
2. *Randomness efficient error reduction:* With  $R$  and  $r$  as above we reduce the error to be less than some arbitrary  $\varepsilon$  using  $r + O(\log(1/\varepsilon))$  additional random bits and running  $R$   $O(\log(1/\varepsilon))$  times.

We will discuss these applications and prove their correctness in the next lecture.

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