CS 6815: Pseudorandomness and Combinatorial Constructions

Fall 2022

Lecture 20: Nov 1, 2022

Lecturer: Jyun-Jie Liao

Scribe: Surendra Ghentiyala

## 1 Introduction

Expander graphs are graphs that are both sparse and well connected. By sparse we mean that they have O(N) edges (where N is the number of edges in the graph). There are several different (yet connected) definitions of well connected which we will see thoroughout the lecture.

## 2 Vertex Expansion

For the duration of the lecture, we will be considering directed *D*-regular graphs.

**Definition 2.1.** For a graph G = (V, E), we define the neighbor set of a vertex  $u \in V$  as  $N(u) = \{v : (u, v) \in E\}$ . We define the neighbor set of a set of vertices  $S \subseteq V$  as  $N(S) = \bigcup_{u \in S} N(u)$ .

**Definition 2.2.** G is a (K, A) vertex expander if for all  $S \subseteq V$  such that  $|S| \leq K$ ,  $|N(S)| \geq A|S|$ .

Notice that the property that  $|S| \leq K$  is necessary because otherwise A = 1 (because then we can have the case where S = V).

Intuitively, a G is a vertex expander if when we look at a subset of vertices S, we can always reach more vertices if we take a step from inside the set S.

**Remark 2.3.** Edge expansion is very similar, it just requires you to think about the neighbors as the edges you can reach quickly form S rather than the vertices you can reach quickly from S.

Ideally, we would like D = O(1),  $A \approx D - 1$  and  $K = \Omega(N)$ .

We can show the existence of expanders by the probabilistic method.

However, we will cheat slightly and show the existence of bipartite expanders. A bipartite expander only requires that the expansion property only hold for subsets in the left side of the graph. We will also only require that the graph is left *D*-regular.

**Theorem 2.4.** for all D = O(1), there exists an  $\alpha = O(1)$  such that a random left D-regular bipartite digraph is an  $(\alpha N, D - 1.1)$  vertex expander with high probability.

Let us consider the probability that for a fixed set S with |S| = k that a random bipartite graph violates the expansion property. In other words that  $|N(S)| \leq (D - 1.1)K$ .

Notice that for  $|N(S)| \leq DK - 1.1K$ , there must be at least 1.1K repetitions in the edges that repeat a node, go to a node which has already been reached by a previous edge.

The probability that a given edge goes to a node that has been covered by a previous edge is bounded above by  $\frac{KD}{N}$ . Thus, the probability that a given set of edges cover 1.1K or more already covered nodes is bounded above by  $(\frac{KD}{N})^{1.1K}$ . Furthermore, there are  $\binom{N}{1.1K}$  choices for the edges which will be repetitions. Thus, for a given S, the probability that there will be 1.1K or more repetitions is upper bounded by  $\binom{N}{1.1K}(\frac{KD}{N})^{1.1K}$ .

Now we will look at the probability that there exists any S with |S| = k, which violates the expansion property.

$$\mathbb{P}[\exists S, |S| = K, |N(S)| \le (D - 1.1)K]$$

By the union bound this is

$$\leq \binom{N}{K} \binom{N}{1.1K} \left(\frac{KD}{N}\right)^{1.1K}$$

By the approximation for the binomial we can see that the above is less than or equal to

$$\leq \left(\frac{eN}{K}\right)^{K} \left(\frac{eKD}{1.1K}\right)^{1.1K} \left(\frac{KD}{N}\right)^{1.1K} \\ = \left(\frac{e^{2.1}D^{2.2}}{1.1^{1.1}} \frac{K^{0.1}}{N^{0.1}}\right)^{K}$$

By making  $\alpha$  very small, we can make K arbitrarily small, thus we can make the following less than or equal to

$$\leq 11^{-K}$$

Finally, we can arrive at the probability that the expansion property is not violated for any set of size less than K as less than or equal to

$$\sum_{i=1}^{\alpha N} 11^{-1} \le 0.1$$

Thus with high probability (at least 90%) our graph is an expander. Thus, it exists.

## 3 Spectral Expansion

**Definition 3.1.** *M* is the random walk matrix of a graph G if  $M_{i,j} = \frac{\text{number of edges from i to } j}{D}$ .

 $\mathbf{M}_{i,j}$  can be thought of as the probability you go from to node j if you are at node i and choose to travel along one of the edges with equal probability.

let  $\boldsymbol{\pi} = [p_1, p_2, \dots, p_n]$ , where  $p_i$  is the probability that you are at node *i*.

 $(\boldsymbol{\pi}\mathbf{M})_i$  is then the probability that you are at node *i* after taking a random step in the graph having been in node *x* with probability  $p_x$  previously.

let **u** be the vector that represents the uniform distribution on vertices,  $\mathbf{u} = \begin{bmatrix} \frac{1}{N}, \frac{1}{N}, \dots, \frac{1}{N} \end{bmatrix}$ Then we can define the expansion of the graph as

$$\lambda(G) = \max_{\mathbf{x} \perp \mathbf{u}} \frac{\|\mathbf{x}\mathbf{M}\|}{\|\mathbf{x}\|}$$

**Remark 3.2.** Usually  $xM^n$  (the distribution after taking n steps in the graph) converges to a stationary distribution.

 $\lambda(G)$  can be thought of as how fast you converge to the uniform distribution for any  $\pi$ . We will now see why.

 $\pi = \mathbf{u} + \mathbf{x}$ , observe that  $\mathbf{u} \perp \mathbf{x}$  since  $\|\mathbf{x}\| = 0$ 

Therefore  $\pi \mathbf{M} = (\mathbf{u} + \mathbf{x})\mathbf{M} = \mathbf{u} + \mathbf{x}\mathbf{M}$  and we know that  $\|\mathbf{x}\mathbf{M}^t\| \leq \lambda(G)^t \|\mathbf{x}\|$ . So how quickly the distribution converges to the uniform distribution is dependent in  $\mathbf{x}\mathbf{M}^t$  which is bounded by  $\lambda(G)$ . So  $\lambda(G)$  tells us how quickly the distribution converges to uniform.

**Remark 3.3.** If you take an undirected graph and let  $\lambda_1 \ge \lambda_2 \ge \cdots \ge \lambda_n$  be the eigenvalues of M, then  $\lambda(G) = \lambda_2$ .

**Theorem 3.4.** Spectral expansion implies vertex expansion. More precisely, for  $\alpha \in [0, 1]$ , G is a  $(\alpha N, \frac{1}{(1-\alpha)\lambda^2+\alpha})$  vertex expander.

Notice that if  $\lambda < 1$ , you have  $\frac{1}{(1-\alpha)\lambda^2+\alpha} > 1$ , which means we have expansion. Otherwise, we do not.

## 4 Mixing

**Definition 4.1.** A graph G has the mixing property if for 2 sets S and T where  $|S| = \alpha N$  and  $|T| = \beta N$ ,  $\frac{e(S,T)}{ND} \approx \alpha \beta$ . e(S,T) is the number of edges between S and T.

Notice that in a random graph, you would expect the mixing property to hold.

**Theorem 4.2.** Spectral Expansion implies Mixing. More precisely  $\left|\frac{e(S,T)}{ND} - \alpha\beta\right| \le \lambda \sqrt{\alpha\beta(1-\alpha)(1-\beta)}$  or the more useful bound  $\left|\frac{e(S,T)}{ND} - \alpha\beta \le \sqrt{\alpha\beta}\right|$ 

Note that if  $\lambda \approx 0$ , then the density between S and T is very close to  $\alpha\beta$ .

**Remark 4.3.** Vertex expansion with strong parameters implies spectral expansion. In fact, to a certain degree any of the 3 definitions of expansion given in this lecture imply the other 2.

We will now present the proof that mixing implies spectral expansion  $\mathbb{1}_S = [\mathbb{1}_1, \mathbb{1}_2, \dots, \mathbb{1}_n]$ , where  $\mathbb{1}_i = 1$  if and only if  $i \in S$ .

$$e(S,T) = \mathbb{1}_S^T A \mathbb{1}_T$$

Where A is the adjacency matrix of M. A = DM. Therefore

$$e(S,T) = \mathbb{1}_{S}^{T} DM \mathbb{1}_{T}$$

The above is true because the left hand side is equal to the following

$$\sum_{i,j} (\mathbb{1}_S)_i (D\mathbf{M})_{i,j} (\mathbb{1}_T)_j$$

Notice that the expression expression in the sum gives the number of edges between i and j if i is in S and j is in T, and zero otherwise. Thus the sum gives the total number of edges between S and T.

Recall from linear algebra that we can write any vector v as  $kv + v^{\perp}$ , where  $k = \sum v_i$ . Using this fact, we can rewrite out expression as

$$(\alpha Nv + (\mathbb{1}_S^{\perp}))^T D\mathbf{M} (\beta Nv + (\mathbb{1}_T^{\perp}))^T$$

Expanding and combing terms we get

$$\alpha\beta N^2 Dv^T \mathbf{M} v \frac{1}{N} + ((\mathbb{1}_S)^{\perp})^T D \mathbf{M} (\mathbb{1}_T)^{\perp}$$
$$= \alpha\beta N^2 D \frac{1}{N} + ((\mathbb{1}_S)^{\perp})^T D \mathbf{M} (\mathbb{1}_T)^{\perp}$$

The first term now simplifies to  $\alpha\beta ND$ . All that remains is to bound the error term  $((\mathbb{1}_S)^{\perp})^T D\mathbf{M}(\mathbb{1}_T)^{\perp}$ . By Cauchy-Swartz

 $((\mathbb{1}_S)^{\perp})^T D\mathbf{M}(\mathbb{1}_T)^{\perp} \le \|(\mathbb{1}_S)^{\perp}\|\|D\mathbf{M}\mathbb{1}_T^{\perp}\|$ 

 $\leq \lambda D \|\mathbb{1}_S^{\perp}\| \|\mathbb{1}_T^{\perp}\|$  By Pythagoras, we know  $\|\mathbb{1}_S^{\perp}\| = \sqrt{\alpha(1-\alpha)N}$ . Thus we have

$$\leq \lambda D \sqrt{\alpha(1-\alpha)N} \sqrt{\beta(1-\beta)N}$$
$$= \lambda N D \sqrt{\alpha(1-\alpha)\beta(1-\beta)}$$