

## Expander Graphs &amp; The Monte Carlo Method

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## 1 Review

**Definition 1.** An  $(n, d, c)$ -expander is a  $d$ -regular bipartite multigraph  $G(X, Y, E)$  with  $|X| = |Y| = n/2$  such that for any  $S \subseteq X$ ,  $|\Delta(S)| \geq (1 + c(1 - 2|S|/n))|S|$

**Definition 2.** For some  $(n, d, c)$ -expander,  $G$ . Let  $P = \text{Adj}/d$  be the probability matrix, and  $P_{i,j} = \text{Adj}_{i,j}/d$ . Let  $Q = (I + P)/2$ .

**Observation 3.**  $Q$  is doubly stochastic and has the uniform distribution as its stationary distribution.

We saw in the previous lecture that the state probability vector will approach the uniform distribution with error decreasing at an exponential rate on the number of steps taken in a random walk.

We will return to random walks later, at which point  $(n, d, c)$ -expander graphs will be used for probability amplification.

## 2 Probability amplification

The goal is to amplify the probability to be exponentially small in the number of trials, while using a small number of random bits, which can be very expensive to generate.

### 2.1 Introduction

#### Naive approach

Modified (and equivalent) definition of **BBP** that has a lower probability of failure:

**Definition 4.** **BBP** is the class of languages  $L$  that have a poly time algorithm  $A$  that for any string  $x$ , and a long random string  $r$ ,

- $x \in L \Rightarrow \Pr[A(x, r) \text{ rejects}] \leq 1/100$
- $x \notin L \Rightarrow \Pr[A(x, r) \text{ accepts}] \leq 1/100$

**Observation 5.** By repeating this algorithm  $k$  times with random strings  $r_1 \dots r_k$ , the probability that the majority of the outcomes is incorrect is  $1/2^{\Omega(k)}$  (proved this with Chernoff bound on a problem set). This uses  $k * r$  random bits.

The goal of this probability amplification exercise is to get a similar exponentially small error, while using many fewer bits. Our target is to use  $O(k) + n$  random bits.

## Walks on expander graphs

The general idea is to construct an expander graph with  $N$  vertices, where  $N = 2^n$ . Each vertex is associated with a unique bit-string from  $\{0, 1\}^n$ . After some number of random steps, the probability vector is almost uniform, and we will be able to generate length  $n$  random bit strings by taking a small number of random steps and returning the label of the current vertex.

**Definition 6.** A  $(d, c)$ -family of expanders is a sequence of  $(n, d, c)$ -expander with  $n = 1 \dots \infty$ .

## Gabber-Galil expanders

Gabber and Galil gave an explicit construction for creating a  $(n, d, c)$ -expander, where  $n = 2m^2$  for some  $m \in \mathbb{Z}$ .

In this construction,  $|X| = |Y| = m^2$ . Assign each vertiex in  $X$  a unique label  $(x, y)$ , such that  $x, y \in \mathbb{Z}_m$ . Do the same for  $Y$ . Since this is an expander family with degree  $d = 5$ , for each vertex  $v = (x, y) \in X$  add an edge from  $v$  to

1.  $(x, y) \in Y$
2.  $(x, x + y) \in Y$
3.  $(x, x + y + 1) \in Y$
4.  $(x + y, y) \in Y$
5.  $(x + y + 1, y) \in Y$

Note that all of the additions above are mod  $n$ . This graph is 5-regular because each of elements in the above list define a permutation on  $X$ , which is a perfect matching from  $X$  to  $Y$ .

The expansion of this graph can be proven to be  $c = (2 - \sqrt{3})/4$ .

## 2.2 Walks on expander graphs

**Definition 7.** Let  $G$  be an  $(N, d, c)$ -expander, where  $N = 2^n$  and  $G$  is created with the Gabber-Galil construction. Let  $A$  be the adjacency matrix of this graph. Its eigenvalues are:  $5 = d = \lambda_1 \geq \lambda_2 \geq \dots \geq \lambda_N = -d = -5$ . The eigenvectors are  $e_1, e_2, \dots, e_N$ , and form an orthonormal basis. Note that there is a constant gap between  $\lambda_1$  and  $\lambda_2$  (shown in the previous lecture).

**Definition 8.**  $P = Adj/d$ . Self loops are added to make the Markov process non-periodic, so  $Q = (I + P)/2$  is the new probability matrix. The  $i^{\text{th}}$  eigenvalue of  $Q$  is  $\lambda'_i = (1 + \lambda_i/d)/2$ . Therefore  $1 = \lambda'_1 \geq \lambda'_2 \geq \dots \geq \lambda'_N = 0$ . The eigenvectors  $e'_1, e'_2, \dots, e'_N$  form an orthonormal basis.

Select a constant  $\beta$  such that  $\lambda_2\beta < 1/100$ .  $\beta$  is a number of fresh bits we will need.

## Probability amplification walk on $(n, d, c)$ -expander

**Definition 9.** *Algorithm:*

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Select a vertex  $p$  uniformly at random from the set of  $N$  vertices (using  $n$  random bits);
for  $i \leftarrow 1$  to  $k$  do
  walk  $\beta$  steps;
   $v_i =$  current vertex;
   $results[i] = A(x, labelOf(v_i))$ ;
  if  $count1's(results) > k/2$  then
    return 1; //assume that  $k$  is odd, so there are no ties
  end
return 0;
end

```

**Observation 10.** *This algorithm runs in polynomial time. The graph has exponential size, but there is no need to explicitly construct the graph, because given a vertex, its neighbors can be computed in polynomial time.*

*Each step of the random walk selects from 6 choices, only a constant number of random bits are needed per step. Therefore, the total number of bits used by this algorithm is  $n + O(\beta k) = n + O(k)$ .*

Work with BBP as a class with witnesses.

**Definition 11.** *Let  $W = \{r \in \{0, 1\}^n | A(x, r) = 1\}$  and  $\bar{W} = \{r \in \{0, 1\}^n | A(x, r) = 0\}$  From the definition of BPP above,  $|W| = 99N/100$  and  $|\bar{W}| = N/100$ .*

**Definition 12.** *Let  $p^{(i)}$  be the distribution for the  $i^{th}$  step (which is also  $r_i$ ).*

*Let  $M$  be a diagonal matrix, where  $M_{i,i} = 1$  if  $i \in W$  and  $M_{i,i} = 0$  otherwise.*

*Let  $\bar{M}$  be a diagonal matrix where  $\bar{M}_{i,i} = 1$  if  $i \in \bar{W}$  and  $\bar{M}_{i,i} = 0$  otherwise.  $\bar{M} = I - M$ .*

Since this is a Markov process,  $p^{(i)} = p^{(0)}B^i$ .

$\|p^{(i)}M\|_1 = \Pr[r_i \text{ is a witness}]$ .

$\|p^{(i)}\bar{M}\|_1 = \Pr[r_i \text{ is not a witness}]$ .

Let  $B = Q^\beta$ .

$\|p^{(0)}BM\|_1 = \Pr[\text{first sample is a witness}]$ . because  $B$  encapsulates  $\beta$  steps.

Select  $S = (S_1, \dots, S_k)$  where  $S_i \in \{M, \bar{M}\}$ , such that  $s_i = m$  iff  $r_i \in W$  and  $s_i = \bar{M}$  iff  $r_i \notin W$ .

**Lemma 13.**  $\Pr[S \text{ occurs}] = \|p^{(0)}(BS_1)(BS_2)\dots(BS_{k-1})(BS_k)\|_1$ .

**Lemma 14.** *for any distribution  $p$ :*

1.  $\|pBM\| \leq \|p\|$

2.  $\|pB\bar{M}\| \leq (1/5)\|p\|$

**Claim 15.** *The probability of error for this algorithm is  $\approx 1/2^k$ .*

*Proof.* The algorithm fails when more than half of the samples fail. Let  $\kappa$  be the number of  $S_i$  terms in  $S$ , not equal to  $\bar{M}$ . The algorithm fails when  $\kappa > k/2$ .

$$\Pr[S \text{ occurs}] = \|p^{(0)}(BS_1)(BS_2)\dots(BS_{k-1})(BS_k)\|_1$$

$$\begin{aligned}
&\leq \sqrt{N} \|p^{(0)}(BS_1)(BS_2)\dots(BS_{k-1})(BS_k)\| \\
&\leq \sqrt{N}(1/5)^k \|p^{(0)}\| - \text{this is by repeated application of Lemma 14.} \\
&\leq \sqrt{N}(1/5)^{k/2} \|p^{(0)}\|.
\end{aligned}$$

Now we estimate how many events  $S$  there are, such that the algorithm fails. An overestimate on this number is  $2^k$ . Thus,  $\Pr[\text{majority of samples fail}] \leq 2^k \sqrt{N}(1/5)^{k/2} \|p^{(0)}\|$ . Since  $p^{(0)}$  is selected from the uniform distribution, its norm is  $1/\sqrt{N}$ .

Therefore  $\Pr[\text{algorithm fails}] \leq 2^k/5^{k/2}$ . □

## Proof of Lemma 14

1.

Express  $p$  as a linear combination of eigenvectors,  $p = \sum_{i=1}^N c_i e_i$ .

$pB = \sum_{i=1}^N c_i e_i B = \sum_{i=1}^N c_i \lambda_i^\beta e_i$ , where  $\lambda_i^\beta$  is in the interval  $[0, 1]$ .

Since  $W$  is a diagonal 0-1 matrix, it can only reduce the norm of a vector that it is applied to. Therefore  $\|pBW\| \leq \|pB\|$ .

Rewrite as  $\|\sum_{i=1}^N c_i \lambda_i^\beta e_i\| = \sqrt{\sum_{i=1}^N c_i^2 \lambda_i^{2\beta}} \leq \sqrt{\sum_{i=1}^N c_i^2}$ . this last step is because the eigenvalue is between 0 and 1.

This last term,  $\sqrt{\sum_{i=1}^N c_i^2} = \|p\|$  □

2.

Decompose  $p$  into  $x$  and  $y$  components:  $p = \sum_{i=1}^N c_i e_i = x + y = c_1 e_1 + \sum_{i=2}^N c_i e_i$ .

By the Pythagorean Inequality,  $\|x\| \leq \|p\|$  and  $\|y\| \leq \|p\|$ .

Focusing on  $x$ ,

since  $\lambda_1 = 1$ ,  $xB = x\lambda_1 c_1 = x$ .

$\|x\bar{W}\| \leq \|x\|/10$ , because  $\bar{W}$  will zero out at least 99/100 of the entries. which will reduce the number of nonzero entries to at most 1/100 of the original amount, which will reduce the norm by at least  $\sqrt{1/100} = 1/10$ .

Focusing on  $y$ ,

$yB = \sum_{i=2}^N c_i e_i B = \sum_{i=2}^N c_i \lambda_i^\beta e_i$ .

$\|yB\bar{W}\| \leq \|yB\|$  because the diagonal matrix can only zero out elements.

Since  $\lambda_2^\beta \leq 1/10$  is the second largest eigenvalue,

$\|yB\bar{W}\| \leq \sqrt{\sum_{i=2}^N c_i^2 \lambda_i^{2\beta}} \leq \lambda_2^\beta \sqrt{\sum_{i=2}^N c_i^2} \leq \|y\|/10$ .

Combining the parts with triangle inequality:  $\|pB\bar{W}\| \leq \|xB\bar{W}\| + \|yB\bar{W}\| \leq (1/10)(\|x\| + \|y\|)$ .

Since  $\|x\| \leq \|p\|$  and  $\|y\| \leq \|p\|$ ,  $(1/10)(\|x\| + \|y\|) < 2\|p\|/10 = \|p\|/5$  □

## 3 Monte Carlo method

Monte carlo method, (and later, the Markov Chain Monte Carlo) estimates values with random sampling.

Start with a simple example for estimating  $\pi$  and move onto to try estimating the number of satisfying assignments of a DNF formula, which will raise some problems with the naive sampling used in the first example.

## Approximating $\pi$

**Example 16.** *Approximating  $\pi$ .*

*Use a unit circle with center at the origin embedded in a square with sides of length 2.*

*Sample  $(x, y)$  pairs,  $n$  times, such that  $x, y$  are in the interval  $[-1, 1]$  and are chosen uniformly at random.*

*Define indicator random variable  $Z_i$  as follows:*

- $Z_i = 1$  if  $x_i^2 + y_i^2 \leq 1$  which is the case that the  $i^{\text{th}}$  sample falls in the circle.
- $Z_i = 0$  in all other cases.

$Pr[Z_i = 1] = \text{area of circle} / \text{area of square} = \pi/4.$

*Let  $Z = \sum_{i=1}^n Z_i.$*

$E[Z] = \sum_{i=1}^n E[Z_i] = n\pi/4.$

*Let  $Z' = 4Z/n.$*

$E[Z'] = \pi.$

*Applying Chernoff bound:*

$Pr(|Z' - \pi| \geq \epsilon\pi)$

$= Pr(|Z - (n\pi)/4| \geq (n\epsilon\pi)/4)$

$= Pr(|Z - E[Z]| \geq \epsilon E[Z])$

$\leq 2e^{-n\pi\epsilon^2/12}$

*If we want to approximate within a certain error probability ( $\delta$ ) simple algebraic manipulation yields:  
 $2e^{-n\pi\epsilon^2/12} \leq \delta \rightarrow n \geq 12\ln(2/\delta)/(\pi\epsilon^2)$*

The example above is quite natural and efficient. Note the ratio of events that cause  $Z_i$  to succeed and fail:  $\pi/4$ .

**Claim 17.** *In general, there must be a dense space to run a polynomial time Monte Carlo simulation, or else interesting objects will most likely not be sampled. The ratio should be  $1/p(n)$  for some polynomial  $p(n)$ .*

## #DNF

**Definition 18.** **#P** is the class of counting problems for languages whose decision problems are in **NP**. Given an instance of a problem from **NP** the analogous **#P** solution is the number of solutions that satisfy the NP decision problem. It is easy to see that **#P** is at least as hard to solve as the analogous NP problem, because if there is a known **#P** algorithm, the NP algorithm could run the **#P** algorithm and return true if the result was not 0.

**Definition 19.** **#DNF**: Count the number of satisfying assignments for a given boolean formula which is represented in DNF.

DNF is the disjunction (or) of conjunctive (and) clauses. for example:  $(x_1 \wedge x_2 \wedge x_3) \vee (x_1 \wedge x_2 \wedge x_4) \vee (x_1)$  is in DNF.

**Observation 20.** Note that the decision problem is easy. Check if any of the clauses are true. The counting problem is harder, and  $\#P$  complete. The formula can be negated and by De Morgan's law, the resulting formula will be in CNF. The relationship between these two counting problems is  $a = 2^n - b$ , where  $a$  is answer to  $\#DNF$ ,  $b$  is answer to  $\#CNF$ , and  $n$  is the number of variables in the formula. Thus, if there an algorithm for the  $\#DNF$  problem, is can be used to compute the answer to  $\#CNF$ . Since  $\#CNF$  is  $\#P$  complete, so is this problem. Therefore, a polynomial algorithm can only exist if  $P = NP$ .

Since it is unlikely that  $P = NP$ , let us try to find an approximation algorithm that runs on polynomial time.

## Approximating $\#DNF$

Try the naive approach, which worked for estimating  $\pi$ .

**Example 21.** Let  $F$  be a formula in DNF.

Randomly assign true/false to each of the  $n$  variables in  $F$ .

Define random variable  $X_i$  as follows:

- $X_i = 1$  if the  $i^{\text{th}}$  assignment is satisfying.
- $X_i = 0$  in all other cases.

Let  $c(F)$  be the number of satisfying assignments.

$$Pr[X_i = 1] = c(F)/2^n.$$

$$E[X_1 + \dots + X_m] = mc(F)/2^n.$$

Applying a Chernoff bound and algebra we derive that the number of trials for desired error prob is:

$$m \geq (3 * 2^n \ln(2/\delta)) / (\epsilon^2 c(F)).$$

Therefore, if  $c(F) \approx 2^n$ , then only a polynomial number of steps is needed. If  $c(F)$  is much smaller than  $2^n$ , then  $m$  is exponential.

**Observation 22.** When the space we are sampling is dense with "interesting" events, Monte Carlo method takes polynomial samples, else it takes exponential samples.

**Definition 23.** The solution to sparse samples space are *Densifiers*. The goal is to define a set of assignments  $S$  such that:

1.  $S$  contains all of the satisfying assignments.
2.  $S$  is not too big compared to  $c(F)$ , this should be bounded by an inverse polynomial.
3. Can uniformly randomly sample  $S$ .

## Definitions for next lecture

The lecture closed with some definitions that will be helpful for the next week:

$C_1 \dots C_t$  are clauses.

For each  $i$ , let  $SC_i$  be the set of assignments that satisfy  $C_i$ .

If  $C_i$  has  $l$  literals, then there are  $2^{n-l}$  satisfying assignments, because the clause only has one satisfying assignment with its variables, and all other variables are free.

$$c(F) = \cup_{i=1}^t SC_i$$

Let  $S = (i, a) : 1 \leq i \leq t$  and  $SC_i$

$$|S| = \sum_{i=1}^t |SC_i|.$$

Let  $i(a)$  be the smallest  $i$ , such that  $a \in SC_i$ .

$$S^c = \{(i(a), a) : \forall \text{ satisfying assignment}\}.$$

$$|S^c| = c(F).$$

Next lecture will show how to sample this set in polynomial time for an accurate estimate.