

Randomized Algorithms

Lecture 6

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Max-Cut

Recall: for an undirected graph $G = (V, E)$, a *cut* is a partition of V into two non-empty sets, $(S, V \setminus S)$. The *capacity* of the cut $(S, V \setminus S)$ is the total number $|C_S|$ of edges that cross the cut, where

$$C_S = \{\{u, v\} \in E \mid u \in S, \& v \in V \setminus S\}$$

A *maximum cut* in G is a cut with maximum capacity.

The **Max-Cut problem**: given G , find a maximum cut.

This is a classic NP-hard problem.

Indeed, it is **NP-complete** to decide, given G and k , whether the size of the max-cut is $\geq k$. ([Karp,1972])

So we believe there is no polynomial-time algorithm to compute this exactly.

Question: How about **approximating** a max-cut? Can we do so efficiently? For example, can we get within factor $\frac{1}{2}$ of a max cut in polynomial time?

We will next show, using a trivial randomized algorithm, that every graph $G = (V, E)$ has a cut of size *at least* $|E|/2$, which can be found efficiently.

Max-Cut: random partitioning

Consider the following trivial **random partition** algorithm:

Algorithm RANDOMCUT($G = (V, E)$)

1. $S \leftarrow \emptyset$
2. **for** every $v \in V$ **do**
3. Flip an (independent) fair coin, X_v .
4. **if** ($X_v = \text{“Heads”}$) **then**
5. $S \leftarrow S \cup \{v\}$
6. **return** $S, V \setminus S$

Question: What is the *expected size* of C_S ?

Max-Cut: analyzing random partitioning

Theorem (6.3)

Let $(S, V \setminus S)$ be the output of *RANDOMCUT*. Then $E[|C_S|] = \frac{|E|}{2}$.

Proof.

Consider the random S created by *RANDOMCUT*.

For each edge $e = \{u, v\} \in E$, let I_e be the indicator variable of whether e is in C_S or not. There are 4 possibilities for each $e = \{u, v\} \in E$:

- (a) $u, v \in S$; (b) $u \in S, v \notin S$; (c) $u \notin S, v \in S$; (d) $u, v \notin S$.

Note: in exactly 2 out of the 4 cases (namely, (b) and (c)) we have $e \in C_S$. Thus, $E[I_e] = \Pr[e \in C_S] = \frac{1}{2}$.

Note that $|C_S| = \sum_{e \in E} I_e$. Hence, summing over all $e \in E$, and by linearity of expectation,

$$E[|C_S|] = \sum_{e \in E} E[I_e] = \frac{|E|}{2}.$$



Max-Cut: analyzing the random partition algorithm

Corollary

For any graph $G = (V, E)$, there exists some cut $(S, V \setminus S)$ such that $|C_S| \geq |E|/2$.

Proof.

Basic but useful observation: If the expected size of C_S is $|E|/2$, then there certainly must exist at least one cut of at least that size. □

The probabilistic method

- ▶ The proof that every graph has a cut of cardinality $\geq |E|/2$ is a very very simple example of **the probabilistic method**.
- ▶ With the probabilistic method, we use randomness and the laws of probability/expectation to prove that a certain combinatorial object must exist.

The probabilistic method

The (basic) probabilistic method:

- ▶ Draw a random object from a set of candidate objects Ω ;
- ▶ Prove that the probability that the random object satisfies a certain property is **strictly positive**;
- ▶ Therefore, an object satisfying that property **must exist!**

This is a non-constructive method of proving the existence of combinatorial objects, pioneered by [Paul Erdős](#).

Although this approach uses probability, the result (that some object with the property exists) is a definite fact, not a probabilistic statement.

Although it only tells us that some object satisfying some desired property exist, in many cases we can also find / construct the object efficiently.

More on the Probabilistic Method later in the course.

De-randomization

- ▶ We did not analyse the probability that RANDOMCUT gives a good (high cardinality) cut, and are not going to do that.
- ▶ Instead, we will in fact *de-randomize* the algorithm using conditional expectation, to obtain a deterministic algorithm that always produces a cut with capacity at least $\frac{|E|}{2}$.

De-randomization

We derandomize via “conditional expectation”.

We are interested in the value of $|C_S|$, and the (conditional) expected value of this quantity will change throughout the algorithm, as vertices get added to S or \bar{S} .

Our randomized algorithm considered the vertices in fixed order. Let X_1, \dots, X_n be the indicator random variables ($X_i = 1$ means that v_i is added to S , whereas $X_i = 0$ means v_i is added to \bar{S}).

Our deterministic derandomized algorithm will construct a specific cut (defined inductively by assigning values x_1, \dots, x_n to X_1, \dots, X_n) of size $\geq \frac{|E|}{2}$, by making decisions for the vertices sequentially one-by-one. At each step we will ensure we choose x_{k+1} so that

$$\mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_{k+1} = x_{k+1}] \geq \mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_k = x_k].$$

Derandomization cont'd.

Suppose $V = \{v_1, \dots, v_n\}$, and suppose we have considered vertices v_1, \dots, v_k sequentially so far, and we have taken decisions x_1, \dots, x_k for these vertices.

Suppose (the induction hypothesis) we know that

$$\mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_k = x_k] \geq \mathbb{E}[|C_S|].$$

Think about the (random) process of adding v_{k+1} . There are two choices for x_{k+1} , of equal probability. Hence,

$$\begin{aligned} \mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_k = x_k] &= \frac{\mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_k = x_k, X_{k+1} = 1]}{2} \\ &+ \frac{\mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_k = x_k, X_{k+1} = 0]}{2}. \end{aligned}$$

Hence, one of the two conditional expectations on the right hand side must be $\geq \mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_k = x_k]$, which (by induction) is $\geq \mathbb{E}[|C_S|] = \frac{|E|}{2}$.

Derandomization cont'd.

In our de-randomized algorithm, how do we decide the value of X_{k+1} ?

For $i \in \{0, 1\}$, we want to compute the conditional expectations

$$Z_i := \mathbb{E}[|C_S| \mid X_1 = x_1, \dots, X_{k+1} = i].$$

Recall the linearity of conditional expectation, and $|C_S| = \sum_{e \in E} I_e$. Hence, we just need to compute

$$Z_{i,e} := \mathbb{E}[I_e \mid X_1 = x_1, \dots, X_{k+1} = i]$$

for each $e \in E$, and then note that $Z_i = \sum_{e \in E} Z_{i,e}$.

Derandomization cont'd.

$$Z_{i,e} = \mathbb{E}[I_e \mid X_1 = x_1, \dots, X_{k+1} = i]$$

There are three possibilities of the two endpoints of e :

- ▶ Both have been determined – the conditional expectation is 0 or 1;
- ▶ One of them is determined – the conditional expectation is $1/2$;
- ▶ None of them is determined – the conditional expectation is $1/2$.

Moreover, these values are easy to compute.

Thus we can compute $Z_{i,e}$ for each e , and sum them up to compute the desired Z_j .

Then we compare Z_0 and Z_1 , and choose the larger of the two.

Derandomization cont'd.

To decide the value of X_{k+1} , all we care actually is whether or not $Z_1 - Z_0 \geq 0$. In particular, we will let $x_{k+1} := 1$ (respectively $x_{k+1} := 0$) precisely when $Z_1 - Z_0 \geq 0$ (respectively, $Z_1 - Z_0 < 0$). Note that

$$Z_1 - Z_0 = \sum_{e \in E} Z_{1,e} - Z_{0,e}.$$

Back to the possibilities for e :

- ▶ If neither endpoints of e is v_{k+1} , $Z_{1,e} = Z_{0,e}$;
- ▶ If one endpoint of e is v_{k+1} and the other end is not determined, then $Z_{1,e} = Z_{0,e} = \frac{1}{2}$.
- ▶ If one endpoint of e is v_{k+1} and the other end is determined, then $Z_{1,e} \neq Z_{0,e}$.

Thus we only need to care about the last case.

Derandomization cont'd.

For $i = 0, 1$, let $A_i := \{v_j \mid j \in [k], X_j = i, \{v_j, v_{k+1}\} \in E\}$.

Namely A_1 (respectively A_0) is the set of neighbours of v_{k+1} from among the already determined nodes $\{v_1, \dots, v_k\}$ that are in S (respectively, in \bar{S}).

Each vertex in A_1 contributes 1 to Z_0 ,
and each vertex in A_0 contributes 1 to Z_1 .

Thus, $Z_1 - Z_0 = |A_0| - |A_1|$.

So, what is the de-randomized algorithm?

Algorithm: starting from the first vertex to the last, assign the current vertex to S or to \bar{S} so as to maximize the current cut value.

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This amounts to just the following simple “**greedy**” algorithm: consider the vertices in some specific order v_1, \dots, v_n . Add v_1 to S (arbitrarily). Then, successively, add the next vertex v_i to the side which has fewer of its neighbors (resulting in the larger addition to the size of the cut), breaking ties arbitrarily.

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The greedy algorithm is guaranteed to yield a cut at least as large and the expected size of a random cut, which is $\frac{|E|}{2}$.

Hence, we can efficiently (and *deterministically*) $\frac{1}{2}$ -approximate a MaxCut.

Question: Can we do better than factor $\frac{1}{2}$?

Max-Cut: the Goemans-Williamson algorithm

In fact, there is a (randomized) polynomial-time algorithm for Max-Cut with a better than $\frac{1}{2}$ -approximation ratio.

In a **breakthrough** result, **Goemans and Williamson** (1995) gave a Max-Cut algorithm with approximation ratio ≈ 0.87856 .

Improving upon this ratio would disprove a major conjecture in computational complexity: either the $\mathbf{P} \neq \mathbf{NP}$ conjecture, or the “**Unique Games Conjecture**” ([Khot’02]) which asserts NP-hardness of a certain problem.

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Max-Cut is equivalent to the following quadratic integer program:

$$\text{MaxCut}(G) = \max : \frac{1}{2} \sum_{\{u,v\} \in E} 1 - x_u x_v$$

$$\text{Subject to: } x_v \in \{+1, -1\}, \quad \forall v \in V.$$

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We cannot solve this efficiently. Instead, we can solve the following **Semidefinite Programming relaxation**:

$$\text{MaxCut}^R(G) = \max : \frac{1}{2} \sum_{\{u,v\} \in E} 1 - \langle \vec{x}_u, \vec{x}_v \rangle$$

$$\text{Subject to: } \vec{x}_v \in \mathbb{R}^n \ \& \ \|\vec{x}_v\|_2 = 1, \quad \forall v \in V.$$

Goemans-Williamson algorithm

For $G = (V, E)$, with $|V| = n$, solving

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Subject to: $\vec{x}_v \in \mathbb{R}^n$ & $\|\vec{x}_v\|_2 = 1, \forall v \in V$.

gives us a collection of n vectors, $\vec{x}_v, v \in V$, on the unit sphere in \mathbb{R}^n .

Question: How do we extract an approximately optimal $\{+1, -1\}$ solution x to $\text{MaxCut}(G)$, from a solution to $\text{MaxCut}^R(G)$?

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Question: How do we extract an approximately optimal $\{+1, -1\}$ solution x to $\text{MaxCut}(G)$, from a solution to $\text{MaxCut}^R(G)$?

Answer: We do a kind of “**randomized rounding**” of the vector solution.

Specifically, choose a “**random vector**”, $\vec{r} \in \mathbb{R}^n$, and set

$x_v := +1$ if $\langle \vec{x}_v, \vec{r} \rangle \geq 0$, and otherwise, set $x_v := -1$.

Equivalently: r describes a random hyperplane, H , through the origin that cuts the unit sphere in half, defining a partition of the vertices (via the partition defined by H of the unit vectors associated with vertices).

It can be shown that the approximation ratio of this algorithm is at least:

$$\frac{2}{\pi} \cdot \min_{0 \leq \theta \leq \pi} \frac{\theta}{(1 - \cos \theta)} \approx 0.87856.$$

References and reading

The $\frac{1}{2}$ -approximation algorithm for MaxCut is covered in Sections 6.2, 6.3 of the book.

The book does not cover the Goemans-Williamson algorithm, and we will not expect you to know the Goemans-Williamson algorithm for the exam. (The G-W algorithm has also been derandomized, using a more involved application of the method of conditional expectations ([Mahajan-Ramesh,1995]).)

If you want to learn more about these and other approximation algorithms for NP-hard problems, two very nice books on the subject are:

V. Vazirani, *Approximation Algorithms*, Springer, 2001.

D. P. Williamson and D. B. Shmoys, *The Design of Approximation Algorithms*, Cambridge University Press, 2011.

We will return to the probabilistic method, and derandomization, later in the lectures.

We will cover Chernoff Bounds next. It's a good idea to start reading the early sections of Chapter 4.