

## Lecture 2: August 30

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## 2.1 Testing Polynomial Identities

Randomized algorithms can be dramatically more efficient than their best known deterministic counterparts. An example of a real-world problem for which there exists a simple and efficient randomized algorithm, but for which there is no known polynomial time deterministic algorithm is that of testing polynomial identities.

The problem takes as input two polynomials  $Q$  and  $R$  over  $n$  variables, with coefficients in some field, and decides whether  $Q \equiv R$ . For example, if we had  $Q(x_1, x_2) = (1+x_1)(1+x_2)$  and  $R(x_1, x_2) = 1+x_1+x_2+x_1x_2$ , the algorithm should output "Yes." This problem arises in many contexts, and is a key primitive in computer algebra packages such as Maple and Mathematica.

An obvious way to attack this problem would be to expand both  $Q$  and  $R$  as sums of monomials and compare coefficients. (Thus, for example,  $Q(x_1, x_2)$  above would be expanded to  $1 + x_1 + x_2 + x_1x_2$ , while  $R(x_1, x_2)$  is already in the correct form.) However, this method in general requires time exponential in the size of the representation of the input polynomials: consider, e.g., the polynomial  $\prod_{i=1}^n (x_i + x_{i+1})$ , which has length  $O(n)$  but expands into  $O(2^n)$  monomials. On the other hand, we may assume that the given polynomials can be *evaluated* efficiently at any given point  $(x_1, \dots, x_n)$ . We will make use of this fact to design an efficient randomized algorithm for the problem.

### 2.1.1 The Schwartz-Zippel Algorithm

This algorithm, due independently to Schwartz [S79] and Zippel [Z79], is a Monte Carlo algorithm with a bounded probability of false positives and no false negatives. For simplicity, we think of the algorithm as receiving a single input polynomial  $P(x_1, \dots, x_n)$  and testing whether  $P \equiv 0$ . (Writing  $P = Q - R$ , this of course, is equivalent to testing  $Q \equiv R$ .) The idea of the algorithm is very simple: assign values  $r_1, \dots, r_n$  chosen independently and uniformly at random from a finite set  $S$  to  $x_1, \dots, x_n$ . Then, using an efficient algorithm for evaluating  $P$ , test if  $P(r_1, \dots, r_n) = 0$ , outputting "Yes" if so and "No" otherwise.

Obviously, if the algorithm outputs "No" then certainly  $P \not\equiv 0$ . However, if the algorithm outputs "Yes" then it is possible that in fact  $P \not\equiv 0$  but  $(r_1, \dots, r_n)$  happens to be a zero of  $P$ . We now bound the probability of such an error occurring.

**Claim 2.1** *If  $P \neq 0$ , then  $\Pr[P(r_1, \dots, r_n) = 0] \leq d/|S|$  where  $d$  is the degree of  $P$ .*

**Proof:** The proof is by induction on  $n$ . For the base case,  $n = 1$ ,  $P$  is a polynomial in one variable and thus has at most  $d$  roots. Hence  $\Pr[P(r_1) = 0] \leq d/|S|$ .

For the inductive step, we let  $k$  be the largest degree of  $x_1$  in  $P$  and write

$$P(x_1, \dots, x_n) = M(x_2, \dots, x_n)x_1^k + N(x_1, \dots, x_n),$$

where the degree of  $M(x_2, \dots, x_n)$  is at most  $d - k$  and the degree of  $x_1$  in  $N$  is strictly less than  $k$ . For the purposes of analysis, we can assume that  $r_2, \dots, r_n$  are chosen first. Then, we let  $\mathcal{E}$  be the event that  $M(r_2, \dots, r_n) = 0$ . There are two cases:

- *Case 1:  $\mathcal{E}$  happens.* From the induction hypothesis applied to  $M$  (a polynomial in  $n - 1$  variables), we know that  $\Pr[\mathcal{E}] \leq (d - k)/|S|$ .
- *Case 2:  $\mathcal{E}$  does not happen.* In this case, we let  $P'$  be the polynomial in one variable  $x_1$  that remains after  $x_2 = r_2, \dots, x_n = r_n$  are substituted in  $P(x_1, \dots, x_n)$ . Since  $M(r_2, \dots, r_n) \neq 0$ , the coefficient of  $x_1^k$  is non-zero, so  $P'$  is a non-zero polynomial of degree  $k$  in one variable. It thus has at most  $k$  roots, so  $\Pr[P'(r_1) = P(r_1, \dots, r_n) = 0 | \neg \mathcal{E}] \leq k/|S|$ .

Putting the two cases together, we have

$$\begin{aligned} \Pr[P(r_1, \dots, r_n) = 0] &= \Pr[P(r_1, \dots, r_n) = 0 | \mathcal{E}] \Pr[\mathcal{E}] + \Pr[P(r_1, \dots, r_n) = 0 | \neg \mathcal{E}] \Pr[\neg \mathcal{E}] \\ &\leq (d - k)/|S| + k/|S| = d/|S|. \end{aligned}$$

■

Thus, if we take the set  $S$  to have cardinality at least twice the degree of our polynomial, we can bound the probability of error by  $1/2$ . This can be reduced to any desired small number by repeated trials, as usual.

### Remarks

1. The Schwartz-Zippel algorithm works fine in finite fields, provided only that we can satisfy the condition  $|S| > d$ . Thus in particular the degree  $d$  of the polynomial being tested must be less than the size of the field.
2. A paper by Gonnet [G84] generalizes the above technique to obtain efficient randomized identity checking algorithms for a much wider range of functions, including trigonometric and logarithmic functions.
3. Is there an efficient *deterministic* algorithm for identity testing? Some efforts in this direction can be found in [LV98], and a sequence of papers by Amir Shpilka and co-authors give deterministic algorithms for various restricted classes of polynomials, such as those described by non-commutative arithmetic formulas, or by depth-3 arithmetic circuits [RS05,KS08,SV08]. On the other hand, a rather devastating negative result was proved by Kabanets and Impagliazzo [KI03], who showed that if there exists a deterministic polynomial time algorithm for checking polynomial identities, then either:
  - NEXP does not have polynomial size circuits; or
  - the Permanent cannot be computed by polynomial-size arithmetic circuits.

This means that an efficient derandomization of the above Schwartz-Zippel algorithm (or indeed, any efficient deterministic algorithm for identity testing) would necessarily entail a major breakthrough in complexity theory. (The point here is that, although both of the above conclusions are generally believed to hold, proving either one of them would be a big deal.)

### 2.1.2 Application to Bipartite Matching

We begin with a nonstandard application of the Schwartz-Zippel algorithm to bipartite matching. Although bipartite matching is easy to solve in deterministic polynomial time using flow techniques, it remains an important open question whether there exists an efficient *parallel* deterministic algorithm for this problem.

Bipartite matching is the following problem: Given a bipartite graph  $G = (V_1, V_2, E)$ , where  $|V_1| = |V_2| = n$ , and all edges connect vertices in  $V_1$  to vertices in  $V_2$ , does  $G$  contain a perfect matching, that is, a subset of exactly  $n$  edges such that each vertex is contained in exactly one edge?

To obtain an efficient randomized algorithm for this problem, we begin with a definition:

**Definition 2.2 (Tutte matrix)** *The Tutte matrix  $A_G$  corresponding to the graph  $G$  is the  $n \times n$  matrix  $[a_{ij}]$  such that  $a_{ij}$  is a variable  $x_{ij}$  if  $(i, j) \in E$ , and 0 otherwise.*

**Claim 2.3**  *$G$  contains a perfect matching if and only if  $\det(A_G) \neq 0$ .*

**Proof:** By definition,  $\det(A_G) = \sum_{\sigma} \text{sgn}(\sigma) \prod_{i=1}^n a_{i\sigma(i)}$ , where the sum is over all permutations  $\sigma$  of  $\{1, \dots, n\}$ . Note that each monomial in this sum corresponds to a *possible* perfect matching in  $G$ , and that the monomial will be non-zero if and only if the corresponding matching is present in  $G$ . Moreover, every pair of monomials differs in at least one (actually, at least two) variables, so there can be no cancellations between monomials. This implies that  $\det(A_G) \neq 0$  iff  $G$  contains a perfect matching. ■

The above Claim immediately yields an efficient algorithm for testing whether  $G$  contains a perfect matching: Simply run the Schwartz-Zippel algorithm on  $\det(A_G)$ , which is a polynomial in  $n^2$  variables of degree  $n$ . Note that the determinant can be computed in  $O(n^3)$  time by Gaussian elimination. Moreover, the algorithm can be efficiently parallelized using the standard fact that an  $n \times n$  determinant can be computed in  $O(\log^2 n)$  time on  $O(n^{3.5})$  processors [Ber84].

**Exercise:** Generalize the above to a non-bipartite graph  $G$ . For this, you will need the skew-symmetric matrix  $A_G = [a_{ij}]$  defined as:

$$a_{ij} = \begin{cases} x_{ij} & \text{if } (i, j) \in E \text{ and } i < j; \\ -x_{ij} & \text{if } (i, j) \in E \text{ and } i > j; \\ 0 & \text{otherwise.} \end{cases}$$

**Hint:** You should show that the above Claim still holds, with this modified definition of  $A_G$ . This requires quite a lot more care than in the bipartite case, because the monomials in  $\det(A_G)$  do not necessarily correspond to perfect matchings, but to *cycle covers* in  $G$ . In this case some cancellations will occur and have to be carefully accounted for.

### 2.1.3 Fast Parallel Algorithms for Finding a Perfect Matching

We now turn to the problem of actually finding a perfect matching. One way to do this would be to reuse the previous algorithm in the following way (assuming we have previously tested  $G$  and discovered that it almost certainly has a perfect matching):

```

Input:  $G = (V, E)$  and  $M = \emptyset$ 
while  $M$  is not a perfect matching
    pick an arbitrary edge  $(i, j)$  of  $G$ 
    test if  $G' = G \setminus \{(i, j)\}$  has a perfect matching using the Schwartz-Zippel algorithm
    if Schwartz-Zippel outputs "Yes"
        then let  $M = M \cup \{(i, j)\}$  and  $G = G'$ 
        else let  $G = G \setminus \{(i, j)\}$ 
    end if
end while
output  $M$ 

```

Since the size of  $E$  decreases by 1 at each iteration of the while loop, after at most  $|E|$  stages we will have constructed a matching. (A detail here is that, since the Schwartz-Zippel test is being invoked several times, we need to reduce its error probability on each invocation to less than  $1/|E|$  using repeated trials, so that the overall error probability remains bounded.) Unfortunately, this approach is inherently sequential and does not yield an efficient parallel algorithm. Instead, we present a randomized algorithm due to Mulmuley, Vazirani and Vazirani [MVV87] that is parallelizable.

We first prove a key lemma; this version is slightly stronger than the original and is due to Noam Ta-Shma.

**Lemma 2.4 [Isolation Lemma]** *Let  $S_1, \dots, S_k$  be subsets of a set  $S$  of cardinality  $m$ . Let each element  $x \in S$  have a weight  $w_x$ , chosen independently and uniformly at random from the set  $\{1, \dots, \ell\}$ . Then*

$$\Pr[\exists \text{ a unique set } S_i \text{ of minimum weight}] \geq \left(1 - \frac{1}{\ell}\right)^m \geq 1 - \frac{m}{\ell}.$$

(Here the weight of set  $S_i$  is defined as  $\sum_{x \in S_i} w_x$ .)

**Proof:** First note that we may assume w.l.o.g. that no set  $S_i$  is a subset of any other set  $S_j$ ; for if so then  $S_j$  cannot have minimum weight and so may be safely ignored.

Now let  $\mathcal{W}$  denote the set of all weight functions  $w = \{w_x\}$ , and  $\mathcal{W}^+$  the subset of  $\mathcal{W}$  satisfying  $w_x > 1$  for all  $x \in S$ . We claim that to each function  $w \in \mathcal{W}^+$ , we can associate a *distinct* function  $w' \in \mathcal{W}$  such that  $w'$  has a unique minimum weight set. This will complete the proof since then

$$\Pr[\exists \text{ a unique set } S_i \text{ of minimum weight}] \geq \frac{|\mathcal{W}^+|}{|\mathcal{W}|} = \left(1 - \frac{1}{\ell}\right)^m \geq 1 - \frac{m}{\ell}.$$

To verify the above claim, given a function  $w \in \mathcal{W}^+$ , pick an arbitrary subset  $S_i$  of minimum weight under  $w$  and define

$$w'_x = \begin{cases} w_x - 1 & \text{if } x \in S_i; \\ w_x & \text{otherwise.} \end{cases}$$

Then clearly  $S_i$  is the unique minimum weight set under  $w'$ . And moreover, we can recover  $w$  from  $w'$  simply by increasing the weights of all elements in  $S_i$  by 1, so the  $w$ 's are distinct. ■

Note that if we take  $\ell = 2m$  (as we will in our application below) then we get a unique minimum weight set with probability at least  $1/2$ .

Armed with the Isolation Lemma, we can give a parallel algorithm to find perfect matchings in bipartite graphs. The set  $S$  in the Isolation Lemma will simply be the edge set  $E$  of the graph, and  $S_1, \dots, S_k$  are the perfect matchings in  $G$ . We pick a weight  $w_{ij}$  for each edge  $(i, j)$  uniformly at random from the set  $\{1, \dots, 2|E|\}$ . The weight  $w(M)$  of a matching  $M$  is the sum of the weights of the edges in  $M$ . We then form the matrix  $B$  by replacing  $x_{ij}$  in the Tutte matrix by  $2^{w_{ij}}$ .

The role of the Isolation Lemma is to “break symmetry” between the perfect matchings in  $G$ . The uniqueness of the minimum weight matching means that all processors are able to work towards the same goal in parallel, without explicit coordination. Here is the algorithm:

```

assign a random weight  $w_{ij}$  to each edge  $(i, j) \in E$  as above
define the matrix  $B$  by replacing each  $x_{ij}$  in the Tutte matrix by  $2^{w_{ij}}$ 
calculate  $2^w$ , the largest power of 2 that divides  $\det(B)$ 
for each edge  $(i, j)$  in parallel do
    compute  $t_{ij} = 2^{w_{ij}} \det(B_{ij})$ , where  $B_{ij}$  is the  $(i, j)$  minor of  $B$ 
    place  $(i, j)$  in  $M$  iff the largest power of 2 that divides  $t_{ij}$  is  $2^w$ 
if  $M$  is a perfect matching then output  $M$  else output fail.

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We note that this algorithm can be implemented to run in polylog time on polynomially many processors, since this applies to each determinant computation and the  $t_{ij}$  can be computed in parallel. The next Claim shows that, if there is a unique minimum weight matching  $M$ , then the algorithm always outputs it. Together with the Isolation Lemma, this implies that the algorithm succeeds with probability at least  $1/2$ . Note that the algorithm ends by checking that it has indeed found a perfect matching, because if the minimum weight matching is not unique we have no control over what the algorithm does. This ensures that any non-fail output is always correct.

**Claim 2.5** *If the minimum weight perfect matching  $M$  is unique, the above algorithm outputs it.*

**Proof:** Let  $M_0$  be the unique minimum weight perfect matching. Then the value  $w$  computed in the first line of the algorithm is just the weight of  $M_0$ ; this follows from the fact that

$$\det(B) = \sum_{M \in \mathcal{M}(G)} \pm 2^{w(M)},$$

where  $\mathcal{M}(G)$  is the set of perfect matchings in  $G$ , and  $M_0$  is unique, so  $\det(B)/2^{w(M_0)}$  is odd.

Moreover, for any edge  $(i, j) \in E$ ,  $\det(B_{ij})$  corresponds to the perfect matchings in  $G \setminus \{i, j\}$ , i.e.,

$$\begin{aligned} \det(B_{ij}) &= \sum_{M \in \mathcal{M}(G \setminus \{i, j\})} \pm 2^{w(M)} \\ &= 2^{-w(i, j)} \sum_{M \cup \{i, j\} \in \mathcal{M}(G)} 2^{w(M \cup \{i, j\})}. \end{aligned}$$

This implies that the largest power of 2 dividing the quantity  $t_{ij}$  computed by the algorithm will still be  $2^w$  iff  $(i, j)$  is an edge of  $M_0$ . Hence the algorithm correctly identifies  $M_0$  assuming that it is unique. ■

**Exercise:** Generalize this algorithm to nonbipartite graphs.

The question of whether a perfect matching can be found *deterministically* in polylog time on polynomially many processors remains one of the central open problems in parallel algorithms. Recently, major progress has been made on this question by Fenner, Gurjar and Thierauf [FGT16] and by Svensson and Tarnawski [ST17]. Both papers achieve polylog time but require *quasi*-polynomially many (i.e.,  $n^{\text{polylog } n}$ ) processors. The key ingredient in both papers is a derandomization of the Isolation Lemma in the particular context where the  $S_i$  are perfect matchings in a graph. [FGT16] does this for bipartite graphs, and [ST17] extends it (non-trivially) to general graphs. Even more recently, Anari and Vazirani [AV18] gave a polylog time algorithm with polynomially many processors for the special case of *planar* graphs, using quite different techniques.

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