Lecture 14
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## 1 Introduction

In this lecture, we discuss linearity testing, self-correction and introduce the Boolean cube.

## 2 Problem setup and definition

We begin by defining terms associated with linearity testing:

Definition 1 (Linear function) Given a function $f$ that maps from a finite set $G$ to another finite set $H$, we say $f$ is linear if $\forall x, y f(x)+_{H} f(y)=f\left(x+_{G} y\right)$.

Here, we define $+_{H}$ to be addition as defined within the set $H$ and $+_{G}$ to be addition as defined within the set $G$. Examples of such linear functions is the family of functions parameterised by $a$ $f_{a}(x)=a x \bmod m, \forall x \in Z_{m}{ }^{\dagger}$. In this example $G=H$. This is linear because $f_{a}\left(x_{1}\right)+f_{a}\left(x_{2}\right)=$ $a x_{1}+a x_{2}=a\left(x_{1}+x_{2}\right)=f_{a}\left(x_{1}+x_{2}\right)$

We also define a closely related concept called $\epsilon$-linear:

Definition 2 ( $\epsilon$-Linear function) Given a function $f$ that maps from a finite set $G$ to another finite set $H$, we say $f$ is $\epsilon$-linear if there exists a linear function $g$ such that $\operatorname{Pr}_{x \in G}[f(x)=g(x)] \geq 1-\epsilon$.

Given these established definitions, we want to understand how we could test if a function is linear. Since both $G$ and $H$ are finite, we could simply run the check on all possible combinations of $x$ and $y$. However, $G$ and $H$ could turn out to be very large. We seek a more tractable way to do this.

## 3 Testing for linearity

As seen in previous lectures, one way to deal with intractability is to introduce randomness. To this end, we present the following randomized algorithm:

```
Do k times:
    Pick random x, y in G
    Test f(x+y)=f(x)+f(y)
```

Figure 1: Call this algorithm $A_{1}$
An important question here, is what should $k$ be? That is, how many times should we run this algorithm so that we can effectively test if $f$ is linear? This is important because it is possible that for some specific $k$ we choose, the function $f$ could pass the linearity test at all those points but could be far from linear. In that case, we want to be able to express how useful $A_{1}$ is. To answer these questions, we make some useful observations and discuss the concept of self-correction.

[^0]
### 3.1 Useful Observations

Suppose we choose some $a, y$ and we want to know $P r_{x \in_{u} G}[y=a+x]$. That is, we want to know the probability that for some $a$ and $y, y=a+x$ for any $x$ chosen uniformly at random from $G$. But with some rearranging, we realise that we can rewrite this probability as follows:

$$
\operatorname{Pr}_{x \in_{u} G}[y=a+x]=\operatorname{Pr}_{x \in_{u} G}[x=y-a]
$$

But since $x$ chosen uniformly at random from $G$, the probability $x$ equals any value is simply $\frac{1}{|G|}$. Therefore, it follows that if $x$ is picked uniformly at random from $G$, then $y=a+x$ is also uniformly distributed in $G$.
This is an important observation and comes up in many places. Consider another example where $G=Z_{2}^{n \dagger}$ and addition under this set is bitwise i.e. $\left(a_{1}, \ldots, a_{n}\right)+\left(b_{1}, \ldots, b_{n}\right)=\left(a_{1}+b_{1}, a_{2}+b_{2}, \ldots, a_{n-1}+b_{n-1}, a_{n}+b_{n}\right)$. Then if we have $a=(0,1,1,0)$ and $x=\left(b_{1}, b_{2}, b_{3}, b_{4}\right)$, under the assumption that $x$ was picked uniformly at random from $G$, then the result $\left(0+b_{1}, 1+b_{2}, 1+b_{3}, 0+b_{4}\right)$ is also uniformly distributed.

### 3.2 Self correction aka "random self-reducibility"

Suppose you are given a function $f: G \rightarrow G^{*}$ such that there exists a linear function $g: G \rightarrow G$ and $\operatorname{Pr}_{x \in G}[f(x)=g(x)] \geq \frac{7}{8}$. Some important notes on this statement:

- This definition is closely related to the definition of $\epsilon$-linear. $\left(\epsilon=\frac{1}{8}\right.$ here)
- We do not know what $f$ looks like; it is handed to us as a blackbox whose functional form is unknown. However, by making calls to $f$, we can approximate what $g$ is.
- There are likely to be multiple such $g$ but we focusing on show that we can approximate one such $g$ with only $O\left(\log \frac{1}{\beta}\right)$ calls to $f$ but with probability of error at most $\beta$.

The idea that we can approximate a linear function $g$ from another function $f$ that may or may not be linear gives rise to the idea of a self-corrector. The idea behind a self-corrector is that we would ideally like to compute $g(x)$ but we do not know what it is and do not have blackbox access to it either. However, there exists a function $f(x)$ which usually agrees with $g(x)$ (at least $\frac{7}{8}$ of the time). We can therefore use calls to $f(x)$ to estimate $g(x)$. The algorithm is as follows:

```
For i=1 to c*log(1/\beta):
    Pick y uniformly at random from G
    Record answer = f(y) + f(x-y) for this round
Output the most common answer from all rounds.
```

Figure 2: $A_{2}(x)$ : Algorithm runs on the same input $x$ for each of the $c \log \frac{1}{\beta}$ rounds
We make our first claim for the algorithm:
Claim 1 : $\operatorname{Pr}\left[A_{2}(x)=g(x)\right] \geq 1-\beta$.
Proof: Recall that

$$
\operatorname{Pr}[f(y) \neq g(y)] \leq \frac{1}{8}
$$

[^1]$$
\operatorname{Pr}[f(x-y) \neq g(x-y)] \leq \frac{1}{8}
$$
from our definition above. Using union bound on these two probabilities, we can write that :
$$
\operatorname{Pr}[f(y)=g(y) \& f(x-y)=g(x-y)] \geq 1-\left(\frac{1}{8}+\frac{1}{8}\right)=\frac{3}{4}
$$

Since $\operatorname{Pr}[f(y)+f(x-y)=g(y)+g(x-y)] \geq \operatorname{Pr}[f(y)=g(y) \& f(x-y)=g(x-y)] \geq \frac{3}{4}{ }^{\dagger}$, we can write that:

$$
\operatorname{Pr}[f(y)+f(x-y)=g(y)+g(x-y)] \geq \frac{3}{4}
$$

But notice that $f(y)+f(x-y)$ is the recorded answer for a single round in $A_{2}$ and $g(x)=g(y)+g(x-y)$. Therefore, this tell us the probability that the answer we get on any round is indeed $g(x)$. To analyze this over all the rounds, we use Chernoff bounds where the $X_{i}$ s are indicator variables for round $i$ and is 1 iff in that round the recorded answer is indeed equal to $g(x)$ and 0 otherwise

$$
\operatorname{Pr}[\text { failure }]=\operatorname{Pr}\left[X<\frac{1}{2} * \mu\right]<e^{-\frac{3 k}{32}}, \quad \mu=\frac{3 k}{4}
$$

Since we want an error of at most $\beta$, it follows that:

$$
\beta=e^{-\frac{3 k}{32}}
$$

Rearranging yields:

$$
k=\frac{32}{3} \ln \left(\frac{1}{\beta}\right)
$$

so we indeed require $O\left(\log \left(\frac{1}{\beta}\right)\right.$ rounds to achieve an error of at most $\beta$.

### 3.3 Linearity Testing is non-trivial

The motivation for discussing self-correction and linearity testing is that if a given function is $\epsilon$-linear and we can test for it, we can hopefully correct for it as we did in the previous subsection because we know the $\operatorname{Pr}$ [failure] However, developing tests for linearity is notoriously difficult. To see this, consider the same linearity testing scheme in Figure 1 and suppose we have a function $f: Z_{m} \rightarrow Z_{m}$. We have that $f$ is defined as follows:

$$
f(x)= \begin{cases}1 & x \equiv 1 \bmod 3 \\ 0 & x \equiv 0 \bmod 3 \\ -1 & x \equiv 2 \bmod 3\end{cases}
$$

A graph of the function is found in Figure 3. From the graph, we see that the best linear function $g(x)=0$ that best approximates $f(x)$ will only align with at most $\frac{1}{3}$ of the points in $f(x)$. Therefore, we say that $g(x)$ is $\frac{2}{3}$-far from linear.
Notice that when $g(x)=0$, our algorithm $A_{1}$ only fails ${ }^{\ddagger}$ anytime $x \equiv y \equiv 1 \bmod 3$ and whenever $x \equiv y \equiv 2 \bmod 3$ but it passes for all other $x, y$. Therefore with our algorithm, the function $f$ succeeds on a large number of $x, y$ even though it is clearly far from linear. We call this example the Coppersmith's Example and it is an example of the worst-case.

[^2]

Figure 3: "tough" function $f(x)$ from Coppersmith's Example

### 3.4 Rejection probability

As seen in the previous subsection, there are cases where our algorithm does not perform well for certain choices of $f$ and $x, y$. As a result, our algorithm isn't guaranteed to return the correct answer all the time. To this end, we define $\delta_{f}$ to be the rejection probability of linearity testing which is the probability that for any $x, y f(x)+f(y) \neq f(x+y)$. In the Coppersmith example, $\delta_{f}=\frac{2}{9}$ for a function $f$ that is $\frac{2}{3}$-far from any $g(x)$. Since the Coppersmith Example is a worst case scenario, $\frac{2}{9}$ is in fact a threshold for defining an important theorem:

Theorem 1: If $\delta_{f}<\frac{2}{9}$, then $f$ must be $\delta_{f}$-close to linear $\S$
Given this theorem about the rejection probability, we can set $k$ from $A_{1}$ to $\frac{1}{\delta_{f}}$. The reasoning here is that if a function is far from linear, then there is some probability of rejection and so we exploit that probability to find the number of times we need to repeat the algorithm $A_{1}$ until we encounter a failing $x, y$.

## 4 Linearity Testing for Boolean functions

In this section, we restrict ourselves to looking at linearity testing for Boolean functions i.e. $f: G \rightarrow$ $\{0,1\}$. to be able to analyze this case, we need a few important definitions and tools like Fourier analysis over Boolean cube. We define these tools and definitions today and will finish up Fourier analysis over Boolean cube next lecture.

[^3]
### 4.1 Boolean cube

The Boolean cube of size $n$ is best described by example:


Figure 4: Boolean cube for $n=3$.
It consists of $i 0 \mathrm{~s}$ and $n-i 1 \mathrm{~s}$ at each level $i \mathbb{I}$. It is funnel shaped at both ends with 111 at the top and 000 at the bottom. The arrows are drawn between levels to connect bits that differ by a single bit flip from 0 to 1 .

### 4.2 Inner product and notation change

We define the inner product on two $n$-bit binary digits $x, y$ as:

$$
x \cdot y=\sum_{i=1}^{n} x_{i} y_{i} \bmod 2
$$

From here, we define a set of linear functions on the domain $\{0,1\}^{n}$ :

$$
L_{a}(x)=a x
$$

for a fixed $a \in\{0,1\}^{n}$. Since these functions are parameterized by $a$ and there are $2^{n}$ choices for $a$, there are also $2^{n}$ of these functions. However, using our inner product notation, we can rewrite this as:

$$
L_{a}(x)=L_{A}(x)=\sum_{i \in A} x_{i} \bmod 2
$$

where $A$ are the set of indices $i$ where $a_{i}=1$
The change of notation we propose is for the function $f$. Before we had $f:\{0,1\}^{n} \rightarrow\{0,1\}$. Now we change $f$ such that $f:\{+1,-1\}^{n} \rightarrow\{+1,-1\}$. Essentially, we are letting $0 \rightarrow+1$ and $1 \rightarrow-1$. This means we have that:

$$
\begin{equation*}
a \rightarrow(-1)^{a} \tag{A}
\end{equation*}
$$

[^4]$$
a+b \rightarrow(-1)^{(a+b)}=(-1)^{a}(-1)^{b} \quad(B)
$$

Essentially by making this change, we are mapping addition to multiplication ${ }^{\|}$. This also changes our linearity condition:

$$
f(a)+f(b)=f(a+b) \rightarrow f(a) f(b)=f(a * b)
$$

As a result, $A_{1}$ changes also. It's condition for passing the test is as follows:

$$
f(x) f(y) f(x * y)= \begin{cases}1 & \text { test accepts } \\ -1 & \text { test rejects }\end{cases}
$$

We can redefine this test for a passed round as follows:

$$
\frac{1-f(x) f(y) f(x * y)}{2}= \begin{cases}0 & \text { test accepts } \\ 1 & \text { otherwise }\end{cases}
$$

We also redefine the rejection probability from $\delta_{f}=P r_{x, y}[f(x)+f(y) \neq f(x+y)]$ to $\delta_{f}=\operatorname{Pr}_{x, y}[f(x) f(y) \neq f(x * y)]$

## 5 Conclusion

In the next lecture, we look at Fourier analysis on the Boolean cube which will help us get stronger bounds than we achieved with Theorem 1.

[^5]
[^0]:    ${ }^{\dagger}$ Recall that $Z_{m}=[0, \ldots, m-1]$

[^1]:    ${ }^{\dagger}$ This notation simply means that $G$ consists of vectors of length $n$ whose entries are either 0 or 1 .
    ${ }^{*}$ Could also do this definition in terms of $G \rightarrow H$

[^2]:    ${ }^{\dagger}$ One way to see this is to consider probability from a counting perspective. The number of ways that $f(y)+f(x-y)=$ $g(y)+g(x-y)$ includes the case where $f(y)=g(y) f(x-y)=g(x-y)$. As a result, the probability of the former occurring is larger than just the latter.
    ${ }^{\ddagger}$ Here fails means that the test on the third line with $f$ fails

[^3]:    ${ }^{\S}$ Or equivalently $\left(1-\delta_{f}\right)$-far from linear.

[^4]:    ICount levels from the top downward as you would a recursion tree

[^5]:    $\|$ We see this clearly in $(B)$

