



Preserving Randomness for Adaptive Algorithms

William M. Hoza*
 Department of Computer Science,
 University of Texas at Austin
 whoza@utexas.edu

Adam R. Klivans
 Department of Computer Science,
 University of Texas at Austin
 klivans@cs.utexas.edu

October 30, 2017

Abstract

Suppose Est is a randomized estimation algorithm that uses n random bits and outputs values in \mathbb{R}^d . We show how to execute Est on k adaptively chosen inputs using only $n + O(k \log(d+1))$ random bits instead of the trivial nk (at the cost of mild increases in the error and failure probability). Our algorithm combines a variant of the INW pseudorandom generator [INW94] with a new scheme for shifting and rounding the outputs of Est . We prove that modifying the outputs of Est is necessary in this setting, and furthermore, our algorithm's randomness complexity is near-optimal in the case $d \leq O(1)$. As an application, we give a randomness-efficient version of the Goldreich-Levin algorithm; our algorithm finds all Fourier coefficients with absolute value at least θ of a function $F : \{0, 1\}^n \rightarrow \{-1, 1\}$ using $O(n \log n) \cdot \text{poly}(1/\theta)$ queries to F and $O(n)$ random bits (independent of θ), improving previous work by Bshouty et al. [BJT04].

1 Introduction

Let Est be a randomized algorithm that estimates some quantity $\mu(C) \in \mathbb{R}^d$ when given input C . The canonical example is the case when C is a Boolean circuit, $d = 1$, $\mu(C) \stackrel{\text{def}}{=} \Pr_x[C(x) = 1]$, and Est estimates $\mu(C)$ by evaluating C at several randomly chosen points. Suppose that Est uses n random bits, and $\Pr[\|\text{Est}(C) - \mu(C)\|_\infty > \varepsilon] \leq \delta$.

Furthermore, suppose we want to use Est as a *subroutine*, executing it on inputs C_1, C_2, \dots, C_k , where each C_i is chosen adaptively based on the previous outputs of Est . The naïve implementation uses nk random bits and fails with probability at most $k\delta$.

In this work, we show how to generically improve the randomness complexity of any algorithm with this structure, without increasing the number of executions of Est , at the expense of mild increases in the error and failure probability. Our algorithm efficiently finds $Y_1, \dots, Y_k \in \mathbb{R}^d$ with $\|Y_i - \mu(C_i)\|_\infty \leq O(\varepsilon d)$ for every i , our algorithm has failure probability $k\delta + \gamma$ for any $\gamma > 0$, and our algorithm uses a total of $n + O(k \log(d+1) + (\log k) \log(1/\gamma))$ random bits.

1.1 The randomness steward model

We model the situation described above by imagining two interacting agents: the *owner* (who plays the role of the outer algorithm) chooses the inputs C_1, \dots, C_k , while the *steward* (who replaces the direct execution of Est) provides the output vectors $Y_1, \dots, Y_k \in \mathbb{R}^d$. The reader might find it useful

*Supported by the National Science Foundation Graduate Research Fellowship Program under Grant No. DGE-1610403.

to think of the steward as a trusted “cloud computing” service. To justify the names, imagine that the owner gives the steward “stewardship” over her random bits. The steward’s job is to “spend” as little randomness as possible without sacrificing too much accuracy.

To describe the model more rigorously, say that a function $f : \{0, 1\}^n \rightarrow \mathbb{R}^d$ is (ε, δ) -concentrated at $\mu \in \mathbb{R}^d$ if $\Pr_{X \in \{0, 1\}^n} [\|f(X) - \mu\|_\infty > \varepsilon] \leq \delta$. In each round i , the chosen input C_i defines a concentrated function $f_i(X) \stackrel{\text{def}}{=} \text{Est}(C_i, X)$, so it is equivalent to imagine that the owner picks an arbitrary concentrated function. In the following definition, ε' is the error of the steward, and δ' is its failure probability.

Definition 1. An (ε', δ') -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ is a randomized algorithm S that interacts with an *owner* O according to the following protocol.

1. For $i = 1$ to k :
 - (a) O chooses $f_i : \{0, 1\}^n \rightarrow \mathbb{R}^d$ that is (ε, δ) -concentrated at some point $\mu_i \in \mathbb{R}^d$ and gives it to S .
 - (b) S chooses $Y_i \in \mathbb{R}^d$ and gives it to O .

Write $O \leftrightarrow S$ (“the interaction of O with S ”) to denote the above interaction. The requirement on S is that for all O ,

$$\Pr[\max_i \|Y_i - \mu_i\|_\infty > \varepsilon' \text{ in } O \leftrightarrow S] \leq \delta'.$$

The probability is taken over the internal randomness of S and O .

From an information-theoretic perspective, stewards as defined above are not particularly interesting, because S could exhaustively examine all outputs of f_i to deterministically compute a point Y_i where f_i is concentrated. But we would like to avoid executing Est more than k times in total, so we will restrict attention to *one-query stewards*:

Definition 2. A *one-query steward* is a steward that only accesses each f_i by querying it at a single point $X_i \in \{0, 1\}^n$. (The point X_i is not seen by the owner.)

1.2 Our results

1.2.1 Main result: A one-query steward with good parameters

Our main result is the explicit construction of a one-query steward that simultaneously achieves low error, low failure probability, and low randomness complexity:

Theorem 1. For any $n, k, d \in \mathbb{N}$ and any $\varepsilon, \delta, \gamma > 0$, there exists a one-query $(O(\varepsilon d), k\delta + \gamma)$ -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ with randomness complexity

$$n + O(k \log(d + 1) + (\log k) \log(1/\gamma)).$$

The total running time of the steward is $\text{poly}(n, k, d, \log(1/\varepsilon), \log(1/\gamma))$.

We also give several variant stewards that achieve tradeoffs in parameters. (See Figure 1.)

ε'	δ'	Randomness complexity	Reference
ε	$k\delta$	nk	Naïve
$O(\varepsilon d)$	$k\delta + \gamma$	$n + O(k \log(d+1) + (\log k) \log(1/\gamma))$	Theorem 1 (main)
$O(\varepsilon)$	$k\delta + \gamma$	$n + O(kd + (\log k) \log(1/\gamma))$	Theorem 4, $d_0 = 1$
$O(\varepsilon d)$	$k\delta + \gamma$	$n + k \log(d+2) + 2 \log(1/\gamma) + O(1)$	Theorem 3*
$O(\varepsilon d)$	$2^{O(k \log(d+1))} \cdot \delta$	n	Theorem 5, $d_0 = d$
$O(\varepsilon)$	$2^{O(kd)} \cdot \delta$	n	Theorem 5, $d_0 = 1$
$O(\varepsilon kd/\gamma)$	$k\delta + \gamma$	$n + O(k \log k + k \log d + k \log(1/\gamma))$	Prop. 1 (based on [SZ99])
$O(\varepsilon)$	$k\delta + k \cdot 2^{-n^{\Omega(1)}}$	$O(n^6 + kd)$	Prop. 2 (based on [IZ89])
Any	Any ≤ 0.2	$n + \Omega(k) - \log(\delta'/\delta)$	Theorem 10 (lower bound)

*Computationally inefficient.

Figure 1: Upper and lower bounds for one-query stewards. Recall that ε, δ are the concentration parameters of f_1, \dots, f_k (i.e. the error and failure probability of the estimation algorithm Est); ε', δ' are the error and failure probability of the steward S; n is the number of input bits to each f_i (i.e. the number of random coins used by Est); k is the number of rounds of adaptivity; d is the dimension of the output of each f_i (i.e. the dimension of the output of Est). Everywhere it appears, γ denotes an arbitrary positive number.

1.2.2 Application: Acceptance probabilities of Boolean circuits

Our first concrete application of Theorem 1 is a time- and randomness-efficient algorithm for estimating the acceptance probabilities of many adaptively chosen Boolean circuits.

Corollary 1. *There exists a randomized algorithm with the following properties. Initially, the algorithm is given parameters $n, k \in \mathbb{N}$ and $\varepsilon, \delta > 0$. Then, in round i ($1 \leq i \leq k$), the algorithm is given a Boolean circuit C_i on n input bits and outputs a number $Y_i \in [0, 1]$. Here, C_i may be chosen adversarially based on Y_1, \dots, Y_{i-1} . With probability $1 - \delta$, every Y_i is $\mu(C_i) \pm \varepsilon$, where $\mu(C_i) \stackrel{\text{def}}{=} \Pr_x[C_i(x) = 1]$. The total running time of the algorithm is*

$$O\left(\frac{\log k + \log(1/\delta)}{\varepsilon^2} \cdot \sum_{i=1}^k \text{size}(C_i)\right) + \text{poly}(n, k, 1/\varepsilon, \log(1/\delta)),$$

and the total number of random bits used by the algorithm is $n + O(k + (\log k) \cdot \log(1/\delta))$.

Corollary 1 should be compared to the case when C_1, \dots, C_k are chosen nonadaptively, for which the randomness complexity can be improved to $n + O(\log k + \log(1/\delta))$ by applying the Goldreich-Wigderson randomness-efficient sampler for Boolean functions [GW97] and reusing randomness. The proof of Corollary 1 works by combining the GW sampler with our steward.

1.2.3 Application: Simulating an oracle for promise-BPP or APP

Recall that **promise-BPP** is the class of promise problems that can be decided in probabilistic polynomial time with bounded failure probability. When an algorithm is given oracle access to a promise problem, it is allowed to make queries that violate the promise, and several models have been considered for dealing with such queries. Following Moser [Mos01], we will stipulate that the oracle may respond in any arbitrary way to such queries. (See, e.g., [BF99] for two other

models.) From these definitions, it is easy to show, for example, that $\mathbf{BPP}^{\text{promise-BPP}} = \mathbf{BPP}$. Using our steward, we give a time- and randomness-efficient simulation of any algorithm with an oracle for **promise-BPP**. (As we will discuss in Section 1.4, the corresponding result for **BPP**-oracle algorithms is trivial.) The algorithm and analysis are almost identical to those used to prove Corollary 1. We also give a similar result for algorithms with an oracle for **APP**, the class introduced by Kabanets et al. [KRC00] of functions $\varphi : \{0, 1\}^n \rightarrow [0, 1]$ that can be approximated to within $\pm\varepsilon$ in probabilistic $\text{poly}(n, 1/\varepsilon)$ time with bounded failure probability.

1.2.4 Application: The Goldreich-Levin algorithm

As a final application, we give a randomness-efficient version of the Goldreich-Levin algorithm [GL89] (otherwise known as the Kushilevitz-Mansour algorithm [KM93]) for finding noticeably large Fourier coefficients. Given oracle access to $F : \{0, 1\}^n \rightarrow \{-1, 1\}$, for any $\theta > 0$, we show how to efficiently find a list containing all U with $|\widehat{F}(U)| \geq \theta$. (Alternatively, thinking of F as an exponentially long bitstring $F \in \{-1, 1\}^{2^n}$, our algorithm finds all Hadamard codewords that agree with F in a $(\frac{1}{2} + \theta)$ -fraction of positions.) Our algorithm makes $O(n \log(n/\delta)) \cdot \text{poly}(1/\theta)$ queries to F , uses $O(n + (\log n) \log(1/\delta))$ random bits, and has failure probability δ . Notice that the number of random bits *does not depend on* θ . To achieve such a low randomness complexity, we first improve the randomness efficiency of each estimate in the standard Goldreich-Levin algorithm using the GW sampler. Then, we reduce the number of rounds of adaptivity by a factor of $\log(1/\theta)$ by making many estimates within each round. Interestingly, we apply our steward with $d = \text{poly}(1/\theta)$, unlike our other applications where we choose $d = 1$. (Recall that d is the number of real values estimated in each round.)

1.2.5 Straightforward application of pseudorandom generators fails

One might hope to design a steward that simply queries each f_i at a pseudorandomly chosen point X_i and returns $Y_i = f_i(X_i)$. We show that a steward of this form (a “pseudorandom generation steward”) must use at least $\Omega(nk)$ random bits, assuming $\delta' \leq 1/2$ and $\delta \geq 2^{-n/2+1}$. So the randomness complexity of any pseudorandom generation steward is not better than that of the naïve steward by more than a constant factor.

1.2.6 Lower bound for one-query stewards

We also prove a randomness complexity lower bound of $n + \Omega(k) - \log(\delta'/\delta)$ for *any* one-query steward. In the case $d \leq O(1)$, this comes close to matching our upper bounds. For example, to achieve $\delta' \leq O(k\delta)$, this lower bound says that $n + \Omega(k)$ random bits are needed; our main steward (Theorem 1) achieves $\varepsilon' \leq O(\varepsilon)$, $\delta' \leq O(k\delta)$ using $n + O(k + (\log k) \log(1/\delta))$ random bits. At the other extreme, if we want a one-query steward that uses only n random bits, this lower bound says that the failure probability will be $\delta' \geq \exp(\Omega(k)) \cdot \delta$; one of our variant stewards (Theorem 5) uses n random bits to achieve $\varepsilon' \leq O(\varepsilon)$ and $\delta' \leq \exp(O(k)) \cdot \delta$.

1.3 Techniques

1.3.1 Block decision trees

A key component in the proof of our main result (Theorem 1) is a pseudorandom generator (PRG) for a new model that we call the *block decision tree* model. Informally, a block decision tree is a decision tree that reads its input from left to right, n bits at a time:

Definition 3. For a finite alphabet Σ , a (k, n, Σ) *block decision tree* is a rooted tree $T = (V, E)$ of height k in which every node v at depth $< k$ has exactly $|\Sigma|$ children (labeled with the symbols in Σ) and has an associated function $v : \{0, 1\}^n \rightarrow \Sigma$. We identify T with a function $T : (\{0, 1\}^n)^{\leq k} \rightarrow V$ defined recursively: $T(\text{the empty string}) = \text{the root node}$, and if $T(X_1, \dots, X_{i-1}) = v$, then $T(X_1, \dots, X_i)$ is the child of v labeled $v(X_i)$.

The standard nonconstructive argument (Appendix D) shows that there exists a γ -PRG for block decision trees with seed length $n + k \log |\Sigma| + 2 \log(1/\gamma) + O(1)$. (See Section 3.1 for the definition of a PRG in this setting.) In Section 3, we explicitly construct a γ -PRG for block decision trees with seed length $n + O(k \log |\Sigma| + (\log k) \log(1/\gamma))$. The generator is constructed by modifying the INW generator for space-bounded computation [INW94].

1.3.2 Shifting and rounding

For a steward S , let $S(X)$ denote S using randomness X . Our main steward is of the form $S(X) \stackrel{\text{def}}{=} S_0(\text{Gen}(X))$. Here, Gen is our PRG for block decision trees, and S_0 is a randomness-inefficient one-query steward. In each round, S_0 queries f_i at a fresh random point $X_i \in \{0, 1\}^n$, but S_0 computes the return value Y_i by carefully *shifting and rounding* each coordinate of $f_i(X_i)$. This deterministic shifting and rounding procedure and its analysis are our main technical contributions. The motivation for shifting and rounding is that by deliberately introducing a small amount of error, we reduce the amount of information about X_i that is leaked by Y_i . This way, Gen can recycle some of the randomness of X_i for future rounds.

To be more quantitative, observe that when any steward and owner interact, it is natural to model the owner’s behavior by a decision tree that branches at each node based on the value Y_i provided by the steward. The *branching factor* of this decision tree is a simple measure of the amount of information leaked. If S_0 simply returned $f_i(X_i)$ without any shifting or rounding, the branching factor for $O \leftrightarrow S_0$ would be 2^n . Three ideas dramatically reduce this branching factor.

- The first idea is to round. Suppose that S_0 rounded each coordinate of $f_i(X_i)$ to the nearest multiple of 2ε (with no shifting). Then the branching factor would be reduced to $2^d + \delta 2^n$. The $\delta 2^n$ term corresponds to the outputs of f_i that are far from its concentration point μ_i . The 2^d term corresponds to outputs $f_i(X_i)$ that are close to μ_i ; if each coordinate of μ_i is approximately equidistant from two multiples of 2ε , then the corresponding coordinate of $f_i(X_i)$ could be rounded to either of those two values.
- The second idea is to *shift* each coordinate of $f_i(X_i)$ before rounding. In particular, S_0 finds a single value Δ_i such that after adding $\Delta_i \cdot 2\varepsilon$ to each coordinate of $f_i(X_i)$, every coordinate is ε -far from every rounding boundary. Then, S_0 rounds the shifted coordinates to obtain Y_i . This procedure reduces the branching factor down to $d + 1 + \delta 2^n$. To understand why, think of Δ_i as a *compressed* representation of Y_i . Assuming $f_i(X_i)$ is close to μ_i , given unlimited computation time, O could recover Y_i from Δ_i by computing the *true* vector μ_i , shifting *it* according to Δ_i , and rounding. Hence, each node of the tree just needs to have one child for each possible Δ_i value (along with the $\delta 2^n$ children for the case that $f_i(X_i)$ is far from μ_i).
- The third idea is to *relax* the requirement that the tree perfectly computes $O \leftrightarrow S_0$. In particular, for every owner O , we construct a block decision tree T_O that merely *certifies correctness* of $O \leftrightarrow S_0$. That is, for any X_1, \dots, X_k , if the node $T_O(X_1, \dots, X_k)$ indicates “success”, then the error $\max_i \|Y_i - \mu_i\|_\infty$ in $O \leftrightarrow S_0(X_1, \dots, X_k)$ is small. On the other hand, if $T_O(X_1, \dots, X_k)$ does not indicate success, then “all bets are off”: the error $\max_i \|Y_i - \mu_i\|_\infty$ in

$O \leftrightarrow S_0(X_1, \dots, X_k)$ may be small or large. Our certification tree has the additional property that

$$\Pr_{X_1, \dots, X_k} [T_O(X_1, \dots, X_k) \text{ indicates success}] \geq 1 - k\delta.$$

This relaxation allows us to reduce the branching factor down to just $d + 2$, because for each node, the $\delta 2^n$ children corresponding to outputs of f_i that are far from μ_i can all be merged into a single “failure” node.

Putting everything together, to save random bits, we don’t need to try to fool $O \leftrightarrow S_0$. Instead, it suffices for Gen to fool the certification tree T_O . The small branching factor of T_O allows Gen to have a correspondingly small seed length.

1.3.3 Lower bounds

Our lower bound for pseudorandom generation stewards uses elementary tools from information theory. For a steward that uses $o(nk)$ random bits, the entropy in the query points (X_1, \dots, X_k) is sufficiently low that some X_j can be predicted reasonably well given X_1, \dots, X_{j-1} . So in rounds $1 \leq i \leq j - 1$, our owner chooses f_i so that X_i is *encoded* in $f_i(X_i)$. In round j , the owner predicts X_j and chooses f_j so that $f_j(X_j)$ is far from the concentration point of f_j .

Our lower bound for one-query stewards follows a similar intuition as our upper bounds: we show that in each round, by carefully choosing f_i , the owner can learn $\Omega(1)$ bits of information about the steward’s randomness. To conclude, we argue that if the steward has fewer than n bits of randomness remaining from the owner’s perspective, then the owner can choose a function that causes the steward’s failure probability to be large.

1.4 Why can’t we just reuse the random bits?

Notwithstanding our lower bounds, the reader might be tempted to think that randomness stewards are trivial: why not just pick $X \in \{0, 1\}^n$ uniformly at random *once* and reuse it in every round? For the purpose of discussion, let us generalize, and suppose we are trying to execute an n -coin algorithm A (not necessarily an estimation algorithm) on k inputs C_1, \dots, C_k . If C_1, \dots, C_k are chosen *non-adaptively* (i.e. all in advance), then we really can use the same X for each execution. By the union bound, the probability that $A(C_i, X)$ fails for any i is at most $k\delta$.

That argument breaks down in the adaptive case, because C_2 is chosen based on $A(C_1, X)$, and hence C_2 may be *stochastically dependent* on X , so $A(C_2, X)$ is not guaranteed to have a low failure probability. For example, if X is encoded in the output $A(C_1, X)$, then an adversarially chosen C_2 could *guarantee* that $A(C_2, X)$ fails.

Even if C_1, \dots, C_k are chosen adaptively, randomness can be safely reused in an important special case: Suppose A is a **BPP** algorithm. Then we can let $\widehat{C}_1, \widehat{C}_2, \dots, \widehat{C}_k$ be the inputs that would be chosen if A never failed. Then each \widehat{C}_i really is independent of X , so by the union bound, with probability $1 - k\delta$, $A(\widehat{C}_i, X)$ does not fail for any i . But if $A(\widehat{C}_i, X)$ does not fail for any i , then by induction, $C_i = \widehat{C}_i$ for every i . So the overall failure probability is once again at most $k\delta$.

More generally, randomness can be safely reused if A is *pseudodeterministic*, i.e. for each input, there is a unique correct output that A gives with probability $1 - \delta$.¹ (Pseudodeterministic algorithms were introduced by Gat and Goldwasser [GG11].) A **BPP** algorithm is trivially pseudodeterministic.

¹These two conditions (inputs are chosen nonadaptively, A is pseudodeterministic) are both special cases of the following condition under which the randomness X may be safely reused: for every $1 \leq i \leq k$, C_i is a pseudodeterministic function of $(C_0, C_1, \dots, C_{i-1})$, where C_0 is a random variable that is independent of X .

Observe, however, that a **promise-BPP** algorithm is only guaranteed to be pseudodeterministic on inputs that satisfy the promise. This is why the result we mentioned in Section 1.2.3 is in terms of an oracle for **promise-BPP**. Similarly, estimation algorithms (including **APP** algorithms) are typically not pseudodeterministic.

In the standard Goldreich-Levin algorithm, randomness is used to estimate $\sum_{U \in \mathcal{U}} \hat{F}(U)^2$ for certain collections of subsets \mathcal{U} . The algorithm’s behavior depends on how the estimate compares to $\theta^2/2$. This process is not pseudodeterministic, because if the true value $\sum_{U \in \mathcal{U}} \hat{F}(U)^2$ is very close to $\theta^2/2$, the estimate falls on each side of $\theta^2/2$ with noticeable probability.

1.5 Related work

1.5.1 Adaptive data analysis

The notion of a randomness steward is inspired by the closely related *adaptive data analysis* problem [HU14, SU15, DFH⁺15c, DFH⁺15a, DFH⁺15b, BH15, BNS⁺16, CLN⁺16], introduced by Dwork et al. [DFH⁺15c]. In the simplest version of this problem, there is an unknown distribution \mathcal{D} over $\{0, 1\}^n$ and a *data analyst* who wishes to estimate the mean values (with respect to \mathcal{D}) of k adaptively chosen functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow [0, 1]$ using as few samples from \mathcal{D} as possible. In this setting, these samples are held by a *mechanism* and are not directly accessible by the data analyst. In round i , the data analyst gives f_i to the mechanism, and the mechanism responds with an estimate of $\mathbb{E}_{x \sim \mathcal{D}}[f_i(x)]$. The mechanism constructs the estimate so as to leak as little information as possible about the sample, so that the same sample points can be safely reused for future estimates.

The data analyst and mechanism in the adaptive data analysis setting are analogous to the owner \mathbf{O} and steward \mathbf{S} in our setting, respectively. In each case, the idea is that the mechanism or steward can intentionally introduce a small amount of error into each estimate to hide information and thereby facilitate future estimates. Note, however, that in the adaptive data analysis problem, there is just one unknown distribution \mathcal{D} and we are concerned with sample complexity, whereas in the randomness stewardship problem, we can think of each concentrated function f_i as defining a new distribution over \mathbb{R}^d and we are concerned with randomness complexity.

1.5.2 The Saks-Zhou algorithm

Another highly relevant construction is the algorithm of Saks and Zhou [SZ99] for simulating randomized logspace algorithms in deterministic space $O(\log^{3/2} n)$. The key component in this algorithm can be reinterpreted as a one-query randomness steward. Saks and Zhou also constructed a randomized algorithm **Est** that approximates a large power of a given substochastic matrix using Nisan’s pseudorandom generator [Nis92]. (In fact, Nisan’s generator can be replaced with any pseudorandom generator for small space [Arm98, HU17].) The Saks-Zhou algorithm works by applying **Est** repeatedly to approximate a much larger power of a given substochastic matrix. The “Saks-Zhou steward” reduces the randomness complexity of this process.

The Saks-Zhou steward works by *randomly perturbing and rounding* the output of each f_i , and then reusing the same random query point X in each round. The perturbation and rounding are somewhat similar to our construction, but note that we shift the outputs of each f_i deterministically, whereas the Saks-Zhou steward uses random perturbations. The analysis of the Saks-Zhou steward is substantially different than the analysis of our steward. Instead, the analysis of the Saks-Zhou steward is similar to the proof that randomness can be safely reused for a pseudodeterministic subroutine; one can show that random perturbation and rounding effectively breaks the dependence between X and Y_i . (See Appendix B for the description and analysis of the Saks-Zhou steward.)

Our steward achieves better parameters than the Saks-Zhou steward (see Figure 1). In particular, to achieve failure probability $k\delta + \gamma$, the error ε' of the Saks-Zhou steward is $O(\varepsilon kd/\gamma)$ – the error grows linearly with k , the number of rounds of adaptivity, as well as with $1/\gamma$. This implies, for example, that if we tried to use the Saks-Zhou steward to estimate the acceptance probabilities of k adaptively chosen Boolean circuits to within $\pm\varepsilon$ with failure probability δ in a randomness-efficient way, we would need to evaluate each circuit on $\Theta(k^2\delta^{-2}\varepsilon^{-2}\log(k/\delta))$ inputs. In contrast, because of our steward’s low error, the algorithm of Corollary 1 evaluates each circuit on just $O(\varepsilon^{-2}\log(k/\delta))$ inputs – an exponential improvement in both k and $1/\delta$. Furthermore, our steward has better randomness complexity than the Saks-Zhou steward.

1.5.3 Pseudorandom generators for adaptive algorithms

Impagliazzo and Zuckerman [IZ89, Imp92] were the first to consider the problem of saving random bits when executing a randomized algorithm A on many adaptively chosen inputs. Instead of assuming that A is an estimation algorithm, Impagliazzo and Zuckerman’s result assumes a known bound on the Shannon entropy of the output distribution of A (e.g., the number of bits output by A). They constructed a pseudorandom generator for this setting; for $k \gg n^6$, the seed length is approximately the sum of the entropy bounds for all the executions of A .

In contrast, we make no assumptions about the entropy of $\text{Est}(C)$. Since $\text{Est}(C)$ is a vector of arbitrary-precision real numbers, the entropy could be as large as n , the number of random bits used by Est . And indeed, our lower bound in Section 7.1 implies that the approach of Impagliazzo and Zuckerman fails in our setting.

One might protest that the entropy of $\text{Est}(C)$ can be reduced by simple rounding. In Appendix C, we construct and analyze a steward that straightforwardly rounds each output and then uses the Impagliazzo-Zuckerman generator in a black-box way. Our main steward achieves much better randomness complexity and failure probability than this “Impagliazzo-Zuckerman steward” (see Figure 1). Our main steward admittedly has larger error than the Impagliazzo-Zuckerman steward ($O(\varepsilon d)$ vs. $O(\varepsilon)$), but one of our variant stewards beats or matches the Impagliazzo-Zuckerman steward in every parameter. (See Appendix C.2 for details.) The improvements come from our more powerful PRG and the fact that we shift before rounding.

1.5.4 Decision trees and branching programs

In the most common decision tree model, the branching factor $|\Sigma|$ is just 2, and each node reads an arbitrary bit of the input. In the more general *parity decision tree* model, each node computes the parity of some subset of the input bits. Kushilevitz and Mansour showed [KM93] that the Fourier ℓ_1 norm of any Boolean function computed by a parity decision tree is at most 2^k , the number of leaves in the tree. It follows immediately that a γ -biased generator is a $(2^k\gamma)$ -PRG for parity decision trees. Using, e.g., the small-bias generator of Naor and Naor [NN93], this gives an efficient PRG for parity decision trees with asymptotically optimal seed length.

Decision trees in which each node computes a more complicated function have also been studied previously. Bellare [Bel92] introduced the *universal decision tree* model, in which each node computes an arbitrary Boolean function of the input bits. He gave a bound on the ℓ_1 norm of any Boolean function computed by a universal decision tree in terms of the ℓ_1 norms of the functions at each node. Unfortunately, for block decision trees, his bound is so large that it does not immediately imply any nontrivial pseudorandom generators for block decision trees.

A block decision tree can be thought of as a kind of space-bounded computation. Indeed, a block decision tree is a specific kind of *ordered branching program* of width $|\Sigma|^k$ and length k that reads n

bits at a time. Hence, we could directly apply a pseudorandom generator for ordered branching programs, such as the INW generator [INW94]. For these parameters, the INW generator has seed length of $n + O(k \log k \log |\Sigma| + \log k \log(1/\gamma))$. This seed length can be slightly improved by instead using Armoni’s generator [Arm98] (a generalization of the Nisan-Zuckerman generator [NZ96]), but even that slightly improved seed length is larger than the seed length of the generator we construct.

1.5.5 Finding noticeably large Fourier coefficients

Our randomness-efficient version of the Goldreich-Levin algorithm should be compared to the results of Bshouty et al. [BJT04], who gave several algorithms for finding noticeably large Fourier coefficients, all closely related to one another and based on an algorithm of Levin [Lev93].

- Bshouty et al. gave one algorithm [BJT04, Figure 4] that makes $O(\frac{n}{\theta^2} \log(\frac{n}{\delta\theta}))$ queries and uses $O(n \log(\frac{n}{\theta}) \log(\frac{1}{\delta\theta}))$ random bits. Our algorithm has better randomness complexity, but worse query complexity.
- Bshouty et al. gave another algorithm [BJT04, Figure 5] that makes only $O(n/\theta^2)$ queries and uses just $O(\log(n/\theta) \cdot \log(1/\theta))$ random bits, but it merely outputs a list such that with probability $1/2$, some U in the list satisfies $|\widehat{F}(U)| \geq \theta$, assuming such a U exists.

We also remark that there is a *deterministic* version of the Goldreich-Levin algorithm for functions with bounded ℓ_1 norm; this follows easily from the work of Kushilevitz and Mansour [KM93] (see also [O’D14, Section 6.4]). In contrast, our algorithm works for all functions $F : \{0, 1\}^n \rightarrow \{-1, 1\}$.

1.6 Outline of this paper

In Section 2, we describe the shifting and rounding steward S_0 and prove that it admits certification trees with a small branching factor. Then, in Section 3, we construct and analyze our pseudorandom generator for block decision trees, mimicking the construction and analysis of the INW generator. In Section 4, we put these pieces together to prove our main result (Theorem 1). In Section 5, we show how to construct our variant stewards. In Section 6, we explain our applications of our main steward. Finally, in Section 7, we prove our randomness complexity lower bounds for stewards.

2 The shifting and rounding steward S_0

As a building block for our main steward constructions, we first construct our randomness-inefficient one-query steward S_0 . Recall that any one-query steward makes two choices in each round: the input X_i to f_i and the estimate $Y_i \in \mathbb{R}^d$. The steward S_0 focuses on the second choice: each X_i is chosen uniformly at random, but S_0 carefully shifts and rounds the output $f_i(X_i)$. (See Figure 2.)

2.1 The shifting and rounding algorithm

We now describe the algorithm by which S_0 computes $Y_i \in \mathbb{R}^d$ from $W_i \stackrel{\text{def}}{=} f_i(X_i)$. Fix $n, k, d \in \mathbb{N}$ and $\varepsilon, \delta > 0$. Let $[d]$ denote the set $\{1, 2, \dots, d\}$. Partition \mathbb{R} into half-open intervals of length $(d+1) \cdot 2\varepsilon$. Let \mathcal{I} denote the set of these intervals. For $w \in \mathbb{R}$, let $\text{Round}(w)$ denote the midpoint of the interval in \mathcal{I} containing w . Given $W_i \in \mathbb{R}^d$:

1. Find $\Delta_i \in [d+1]$ such that for every $j \in [d]$, there is some $I \in \mathcal{I}$ such that

$$[W_{ij} + (2\Delta_i - 1)\varepsilon, W_{ij} + (2\Delta_i + 1)\varepsilon] \subseteq I.$$

(We will show that such a Δ_i exists.)

1. For $i = 1$ to k :
 - (a) \mathcal{O} chooses $f_i : \{0, 1\}^n \rightarrow \mathbb{R}^d$ and gives it to \mathcal{S}_0 .
 - (b) \mathcal{S}_0 picks *fresh randomness* $X_i \in \{0, 1\}^n$ and queries to obtain $W_i \stackrel{\text{def}}{=} f_i(X_i)$.
 - (c) \mathcal{S}_0 computes Y_i by shifting and rounding W_i according to the algorithm in Section 2.1.
 - (d) \mathcal{S}_0 gives Y_i to \mathcal{O} .

Figure 2: Outline of $\mathcal{O} \leftrightarrow \mathcal{S}_0$.

2. For every $j \in [d]$, set $Y_{ij} = \text{Round}(W_{ij} + 2\Delta_i \varepsilon)$.

We must show that this algorithm is well-defined:

Lemma 1. *For any $W \in \mathbb{R}^d$, there exists $\Delta \in [d + 1]$ such that for every $j \in [d]$, there is a single interval in \mathcal{I} that entirely contains $[W_j + (2\Delta - 1)\varepsilon, W_j + (2\Delta + 1)\varepsilon]$.*

Proof. Consider picking $\Delta \in [d + 1]$ uniformly at random. For each j , the probability that two distinct intervals in \mathcal{I} intersect $[W_j + (2\Delta - 1)\varepsilon, W_j + (2\Delta + 1)\varepsilon]$ is precisely $1/(d + 1)$ by our choice of the length of the intervals. The union bound over d different j values completes the proof. \square

2.2 Analysis: Certification trees

As outlined in Section 1.3.2, the key lemma says that for any owner \mathcal{O} , there exists a block decision tree $T_{\mathcal{O}}$ with a small branching factor that certifies correctness of $\mathcal{O} \leftrightarrow \mathcal{S}_0$:

Lemma 2. *Assume $\delta < 1/2$. Let $\Sigma = [d + 1] \cup \{\perp\}$. For any deterministic owner \mathcal{O} , there exists a (k, n, Σ) block decision tree $T_{\mathcal{O}}$ with the following properties.*

1. For any internal node v , $\Pr_{X \in \{0, 1\}^n} [v(X) = \perp] \leq \delta$.
2. Fix $X_1, \dots, X_k \in \{0, 1\}^n$, and suppose that the path from the root to $T_{\mathcal{O}}(X_1, \dots, X_k)$ does not include any \perp nodes. Then $\max_i \|Y_i - \mu_i\|_{\infty} \leq O(\varepsilon d)$ in $\mathcal{O} \leftrightarrow \mathcal{S}_0(X_1, \dots, X_k)$.

Notice that Lemma 2 does not assert that $T_{\mathcal{O}}$ computes the transcript of $\mathcal{O} \leftrightarrow \mathcal{S}_0$. In fact, for the analysis, we will define *another* steward \mathcal{S}'_0 , and $T_{\mathcal{O}}$ will compute a sequence of values that arise in $\mathcal{O} \leftrightarrow \mathcal{S}'_0$. This new steward \mathcal{S}'_0 will be computationally inefficient; it will *compress and decompress* the output of \mathcal{S}_0 (with some chance of failure) before giving it to \mathcal{O} , as we suggested in Section 1.3.2.

Proof of Lemma 2. For a function $f : \{0, 1\}^n \rightarrow \mathbb{R}^d$ that is (ε, δ) -concentrated at some point $\mu \in \mathbb{R}^d$, define $\mu(f)$ to be the *smallest* vector in \mathbb{R}^d (under, say, the lexicographical order) at which f is (ε, δ) -concentrated. (This exists, because $\{0, 1\}^n$ is finite, so the set of points where f is concentrated is a compact subset of \mathbb{R}^d .) For a vector $Y \in \mathbb{R}^d$, say that a value $\Delta \in [d + 1]$ is *f-compatible* with Y if $Y_j = \text{Round}(\mu(f)_j + 2\Delta\varepsilon)$ for every $j \in [d]$. Just for the analysis, let \mathcal{S}'_0 be the following (many-query) steward:

1. For $i = 1$ to k :
 - (a) Give f_i to \mathcal{S}_0 , allowing it to make its one query and choose its output vector $Y_i \in \mathbb{R}^d$.
 - (b) Query f_i at *every* point in its domain, thereby learning the entire function.

(c) Compute

$$\widehat{\Delta}_i = \begin{cases} \text{the smallest } \Delta \in [d+1] \text{ } f_i\text{-compatible with } Y_i & \text{if any such } \Delta \text{ exists} \\ \perp & \text{otherwise.} \end{cases}$$

(d) Output $\widehat{Y}_i = (\widehat{Y}_{i1}, \dots, \widehat{Y}_{id})$, where for each $j \in [d]$,

$$\widehat{Y}_{ij} = \begin{cases} \text{Round}(\mu(f)_j + 2\widehat{\Delta}_i\varepsilon) & \text{if } \widehat{\Delta}_i \neq \perp \\ 0 & \text{otherwise.} \end{cases}$$

We are now ready to formally define $T_{\mathcal{O}}$ as a function. Because $S'_0(X_1, \dots, X_k)$ looks at X_i only in round i , we can sensibly speak of the first i rounds of $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_i)$ even for $i < k$. This allows us to define $T_{\mathcal{O}}(X_1, \dots, X_i)$ to be the node v in $T_{\mathcal{O}}$ such that the path from the root to v is described by the values $\widehat{\Delta}_1, \dots, \widehat{\Delta}_i$ that arise in $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_i)$.

Now, we must show that this function $T_{\mathcal{O}}$ can be realized as a block decision tree, i.e. that each internal node v can be assigned a transition function $v : \{0, 1\}^n \rightarrow \Sigma$ that is compatible with the definition of $T_{\mathcal{O}}$ as a function. Indeed, observe that $\widehat{\Delta}_1, \dots, \widehat{\Delta}_{i-1}$ fully determine the state of \mathcal{O} after the first $i-1$ rounds of $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_i)$ and hence determine the function f_i . Furthermore, S_0 is “memoryless”, i.e. Y_i is fully determined by f_i and X_i . Thus, $\widehat{\Delta}_i$ is fully determined by $\widehat{\Delta}_1, \dots, \widehat{\Delta}_{i-1}$ and X_i . So there is a function $\varphi : (\widehat{\Delta}_1, \dots, \widehat{\Delta}_{i-1}, X_i) \mapsto \widehat{\Delta}_i$, and if the path from the root to v is described by $\widehat{\Delta}_1, \dots, \widehat{\Delta}_{i-1}$, we can set $v(X_i) \stackrel{\text{def}}{=} \varphi(\widehat{\Delta}_1, \dots, \widehat{\Delta}_{i-1}, X_i)$.

Analysis of $T_{\mathcal{O}}$ By the definition of $T_{\mathcal{O}}$ as a function, to prove Condition 1 in the lemma statement, we must show that in each round of $\mathcal{O} \leftrightarrow S'_0$, $\Pr[\widehat{\Delta}_i = \perp] \leq \delta$. Indeed, by concentration, with probability $1 - \delta$, for every j , $|W_{ij} - \mu(f_i)_j| \leq \varepsilon$. In this case, by the construction of S_0 , $W_{ij} + 2\Delta_i\varepsilon$ and $\mu(f_i)_j + 2\Delta_i\varepsilon$ are in the same interval in \mathcal{I} for every $j \in [d]$. Therefore, in this case, there is at least one Δ value that is f_i -compatible with Y_i , namely the value Δ_i used by S_0 .

Finally, to prove Condition 2 in the lemma statement, suppose the path from the root node to $T_{\mathcal{O}}(X_1, \dots, X_k)$ does not include any \perp nodes. Then in $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_k)$, for every i , $\widehat{\Delta}_i \neq \perp$. This implies that every Y_{ij} is of the form $\text{Round}(\mu(f_i)_j + 2\widehat{\Delta}_i\varepsilon)$ for some $\widehat{\Delta}_i \in [d+1]$. Therefore, $|Y_{ij} - \mu(f_i)_j| \leq 3(d+1)\varepsilon$, since $2\widehat{\Delta}_i\varepsilon \leq 2(d+1)\varepsilon$ and rounding introduces at most $(d+1)\varepsilon$ additional error.

Of course, so far the analysis has treated S'_0 , not S_0 . But the crucial point is, for every i , since $\widehat{\Delta}_i \neq \perp$, we can be sure that $Y_i = \widehat{Y}_i$. Therefore, the values $f_1, \dots, f_k, Y_1, \dots, Y_k$ in $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_k)$ are *exactly the same* as they are in $\mathcal{O} \leftrightarrow S_0(X_1, \dots, X_k)$! Therefore, in $\mathcal{O} \leftrightarrow S_0(X_1, \dots, X_k)$, for every i , $\|Y_i - \mu(f_i)\|_{\infty} \leq (3d+3)\varepsilon$. Finally, since $\delta < 1/2$, if μ_i is *any* point where f_i is (ε, δ) -concentrated, $\|\mu(f_i) - \mu_i\|_{\infty} \leq 2\varepsilon$. Therefore, for every i , $\|Y_i - \mu_i\|_{\infty} \leq 3(d+1)\varepsilon + 2\varepsilon = (3d+5)\varepsilon$. \square

Notice that in $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_k)$, if $\widehat{\Delta}_i = \perp$ for some i , then the interaction might diverge from $\mathcal{O} \leftrightarrow S_0(X_1, \dots, X_k)$, in which case $T_{\mathcal{O}}(X_1, \dots, X_k)$ does not encode the transcript of $\mathcal{O} \leftrightarrow S_0(X_1, \dots, X_k)$ in any way.

3 Pseudorandomness for block decision trees

Recall that our goal is to modify the internal parameters of the INW generator, thereby constructing a γ -PRG for (k, n, Σ) block decision trees with seed length $n + O(k \log |\Sigma| + (\log k) \log(1/\gamma))$. The construction and analysis mimic the standard treatment of the INW generator, and the reader

who is familiar with the INW generator is encouraged to skip to Section 3.4 to just see the new parameters. In words, the only new feature is that we use extractors for a geometrically growing entropy deficit at each level of the recursion to match the geometrically growing width of the block decision tree.

3.1 Formal definitions and theorem statement

Let U_n denote the uniform distribution on $\{0, 1\}^n$. For two probability distributions μ, μ' on the same measurable space, write $\mu \sim_\gamma \mu'$ to indicate that μ and μ' have total variation distance at most γ .

Definition 4. We say that $\text{Gen} : \{0, 1\}^s \rightarrow \{0, 1\}^{nk}$ is a γ -PRG for (k, n, Σ) block decision trees if for every such tree T , $T(\text{Gen}(U_s)) \sim_\gamma T(U_{nk})$.

Theorem 2. For every $n, k \in \mathbb{N}$, every finite alphabet Σ , and every $\gamma > 0$, there exists a γ -PRG $\text{Gen} : \{0, 1\}^s \rightarrow \{0, 1\}^{nk}$ for (k, n, Σ) block decision trees with seed length

$$s \leq n + O(k \log |\Sigma| + (\log k) \log(1/\gamma)).$$

The PRG can be computed in $\text{poly}(n, k, \log |\Sigma|, \log(1/\gamma))$ time.

3.2 Concatenating PRGs for block decision trees

Toward proving Theorem 2, for a (k, n, Σ) block decision tree $T = (V, E)$ and a node $v \in V$, let T_v denote the subtree rooted at v , and observe that we can think of T_v as a (k', n, Σ) block decision tree, where $k' = k - \text{depth}(v)$. This simple observation – after a block decision tree has been computing for a while, the remaining computation is just another block decision tree – implies that pseudorandom generators for block decision trees can be *concatenated* with mild error accumulation. This fact and its easy proof are perfectly analogous to the situation with ordered branching programs. We record the details below.

Lemma 3. Suppose $\text{Gen}_1 : \{0, 1\}^{s_1} \rightarrow \{0, 1\}^{nk_1}$ is a γ_1 -PRG for (k_1, n, Σ) block decision trees and $\text{Gen}_2 : \{0, 1\}^{s_2} \rightarrow \{0, 1\}^{nk_2}$ is a γ_2 -PRG for (k_2, n, Σ) block decision trees. Let $\text{Gen}(x, y) = (\text{Gen}_1(x), \text{Gen}_2(y))$. Then Gen is a $(\gamma_1 + \gamma_2)$ -PRG for $(k_1 + k_2, n, \Sigma)$ block decision trees.

Proof. Fix a $(k_1 + k_2, n, \Sigma)$ block decision tree T . For a node u at depth k_1 and a leaf node v , define

$$\begin{aligned} p(u) &= \Pr[T(U_{nk_1}) = u] & p(v | u) &= \Pr[T_u(U_{nk_2}) = v] \\ \tilde{p}(u) &= \Pr[T(\text{Gen}_1(U_{s_1})) = u] & \tilde{p}(v | u) &= \Pr[T_u(\text{Gen}_2(U_{s_2})) = v]. \end{aligned}$$

To prove correctness of Gen , recall that ℓ_1 distance is twice total variation distance. The ℓ_1 distance between $T(\text{Gen}(U_{s_1+s_2}))$ and $T(U_{n(k_1+k_2)})$ is precisely $\sum_{u,v} |p(u)p(v | u) - \tilde{p}(u)\tilde{p}(v | u)|$. By the triangle inequality, this is bounded by

$$\begin{aligned} & \sum_{u,v} |p(u)p(v | u) - p(u)\tilde{p}(v | u)| + \sum_{u,v} |p(u)\tilde{p}(v | u) - \tilde{p}(u)\tilde{p}(v | u)| \\ &= \sum_{u,v} p(u) \cdot |p(v | u) - \tilde{p}(v | u)| + \sum_{u,v} |p(u) - \tilde{p}(u)| \cdot \tilde{p}(v | u) \\ &= \sum_u p(u) \sum_v |p(v | u) - \tilde{p}(v | u)| + \sum_u |p(u) - \tilde{p}(u)|. \end{aligned}$$

By the correctness of Gen_1 and Gen_2 , this is bounded by $(\sum_u p(u) \cdot 2\gamma_2) + 2\gamma_1 = 2(\gamma_1 + \gamma_2)$. \square

3.3 Recycling randomness

We find it most enlightening to think of the INW generator in terms of extractors, as suggested by Raz and Reingold [RR99] and in the spirit of the Nisan-Zuckerman generator [NZ96]. The analysis is particularly clean if we work with *average-case extractors*, a concept introduced by Dodis et al. [DORS08].

Definition 5. For discrete random variables X, V , the *average-case conditional min-entropy* of X given V is

$$\tilde{H}_\infty(X | V) = -\log_2 \left(\mathbb{E}_{v \sim V} \left[2^{-H_\infty(X|V=v)} \right] \right),$$

where H_∞ is (standard) min-entropy.

Intuitively, $\tilde{H}_\infty(X | V)$ measures the amount of randomness in X from the perspective of someone who knows V . The output of an *average-case extractor* is required to look uniform even from the perspective of someone who knows V , as long as its first input is sampled from a distribution that has high min-entropy conditioned on V :

Definition 6. We say that $\text{Ext} : \{0, 1\}^s \times \{0, 1\}^d \rightarrow \{0, 1\}^m$ is an *average-case $(s - t, \beta)$ -extractor* if for every X distributed on $\{0, 1\}^s$ and every discrete random variable V such that $\tilde{H}_\infty(X | V) \geq s - t$, if we let $Y \sim U_d$ be independent of (X, V) and let $Z \sim U_m$ be independent of V , then $(V, \text{Ext}(X, Y)) \sim_\beta (V, Z)$.

Average-case extractors are the perfect tools for *recycling randomness* in space-bounded computation. We record the details for block decision trees below.

Lemma 4 (Randomness recycling lemma for block decision trees). *Suppose $\text{Gen} : \{0, 1\}^s \rightarrow \{0, 1\}^{nk}$ is a γ -PRG for (k, n, Σ) block decision trees and $\text{Ext} : \{0, 1\}^s \times \{0, 1\}^d \rightarrow \{0, 1\}^s$ is an average-case $(s - k \log |\Sigma|, \beta)$ -extractor. Define*

$$\text{Gen}'(x, y) = (\text{Gen}(x), \text{Gen}(\text{Ext}(x, y))).$$

Then Gen' is a $(2\gamma + \beta)$ -PRG for $(2k, n, \Sigma)$ block decision trees.

Proof. Let T be a $(2k, n, \Sigma)$ block decision tree. Let $X \sim U_s$ and let $V = T(\text{Gen}(X))$. By [DORS08, Lemma 2.2b], the fact that V can be described using $k \log |\Sigma|$ bits implies that $\tilde{H}_\infty(X | V) \geq s - k \log |\Sigma|$. Therefore, by the average-case extractor condition, if we let $Y \sim U_d$ be independent of X and $Z \sim U_d$ be independent of V , then

$$(V, \text{Ext}(X, Y)) \sim_\beta (V, Z).$$

Applying a (deterministic) function can only make the distributions closer. Apply the function $(v, z) \mapsto T_v(\text{Gen}(z))$:

$$T(\text{Gen}'(X, Y)) \sim_\beta T(\text{Gen}(X), \text{Gen}(Z)).$$

By Lemma 3, the right-hand side is (2γ) -close to $T(U_{2nk})$. The triangle inequality completes the proof. \square

To actually construct a generator, we will need to instantiate this randomness recycling lemma with an explicit average-case extractor:

Lemma 5. *For every $s, t \in \mathbb{N}$ and every $\beta > 0$, there exists an average-case $(s - t, \beta)$ -extractor $\text{Ext} : \{0, 1\}^s \times \{0, 1\}^d \rightarrow \{0, 1\}^s$ with seed length $d \leq O(t + \log(1/\beta))$ computable in time $\text{poly}(s, \log(1/\beta))$.*

Proof sketch. It is standard (and can be proven using expanders, see, e.g., [Vad12]) that there exists an *ordinary* $(s - t - \log(2/\beta), \beta/2)$ -extractor $\text{Ext} : \{0, 1\}^s \times \{0, 1\}^d \rightarrow \{0, 1\}^s$ with seed length $d \leq O(t + \log(1/\beta))$ computable in time $\text{poly}(s, \log(1/\beta))$. By the same argument as that used to prove [DORS08, Lemma 2.3], Ext is automatically an average-case $(s - t, \beta)$ -extractor. \square

3.4 The recursive construction

Proof of Theorem 2. Define $\beta = \gamma/2^{\lceil \log k \rceil}$. For $i \geq 0$, define $s_i \in \mathbb{N}$, $d_i \in \mathbb{N}$, $G_i : \{0, 1\}^{s_i} \rightarrow \{0, 1\}^{n \cdot 2^i}$, and $\text{Ext}_i : \{0, 1\}^{s_i} \times \{0, 1\}^{d_i} \rightarrow \{0, 1\}^{s_i}$ through mutual recursion as follows. Start with $s_0 = n$ and $G_0(x) = x$. Having already defined s_i and G_i , let Ext_i be the average-case $(s_i - 2^i \log |\Sigma|, \beta)$ -extractor of Lemma 5, and let d_i be its seed length. Then let $s_{i+1} = s_i + d_i$, and let

$$G_{i+1}(x, y) = (G_i(x), G_i(\text{Ext}_i(x, y))).$$

We show by induction on i that G_i is a $(\beta \cdot (2^i - 1))$ -PRG for $(2^i, n, \Sigma)$ block decision trees. In the base case $i = 0$, this is trivial. For the inductive step, apply Lemma 4, and note that $2\beta(2^i - 1) + \beta = \beta(2^{i+1} - 1)$. This completes the induction. Therefore, we can let $\text{Gen} = G_{\lceil \log k \rceil}$, since $\beta \cdot (2^{\lceil \log k \rceil} - 1) < \gamma$. The seed length $s_{\lceil \log k \rceil}$ of Gen is

$$\begin{aligned} n + \sum_{i=0}^{\lceil \log k \rceil} d_i &\leq n + O\left(\sum_{i=0}^{\lceil \log k \rceil} (2^i \log |\Sigma| + \log k + \log(1/\gamma))\right) \\ &\leq n + O(k \log |\Sigma| + (\log k) \log(1/\gamma)). \end{aligned}$$

The time needed to compute $\text{Gen}(x)$ is just the time needed for $O(k)$ applications of Ext_i for various $i \leq O(\log k)$, which is $\text{poly}(n, k, \log |\Sigma|, \log(1/\gamma))$. \square

4 Proof of main result (Theorem 1)

Without loss of generality, assume $\delta < 1/2$. (If $\delta \geq 1/2$, then either $k = 1$ or $k\delta \geq 1$; in either case, the result is trivial.) Let \mathbf{S}_0 be the steward of Section 2, let Σ be the alphabet of Lemma 2, and let Gen be the γ -PRG for (k, n, Σ) block decision trees of Theorem 2. The steward is $\mathbf{S}(X) \stackrel{\text{def}}{=} \mathbf{S}_0(\text{Gen}(X))$.

Consider any owner \mathbf{O} . We may assume without loss of generality that \mathbf{O} is deterministic, because a randomized owner is just a distribution over deterministic owners. By Condition 1 of Lemma 2 and the union bound,

$$\Pr[\text{some node in the path from the root to } T_{\mathbf{O}}(U_{nk}) \text{ is labeled } \perp] \leq k\delta.$$

Therefore, when $T_{\mathbf{O}}$ reads $\text{Gen}(U_s)$ instead of U_{nk} , the probability is at most $k\delta + \gamma$. By Condition 2 of Lemma 2, this proves the correctness of \mathbf{S} . The randomness complexity of \mathbf{S} is just the seed length of Gen , which is indeed $n + O(k \log |\Sigma| + (\log k) \log(1/\gamma)) = n + O(k \log(d + 1) + (\log k) \log(1/\gamma))$. The total runtime of \mathbf{S} is clearly $\text{poly}(n, k, d, \log(1/\varepsilon), \log(1/\gamma))$.² \square

²We assume here that our computational model allows the necessary arithmetic and rounding of Section 2.1 to be performed efficiently, even if the owner chooses an f_i that outputs vectors whose coordinates are very large numbers.

5 Variant stewards

Theorem 3. *For any $n, k, d \in \mathbb{N}$, for any $\varepsilon, \delta, \gamma > 0$, there exists a (computationally inefficient) one-query $(O(\varepsilon d), k\delta + \gamma)$ -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ with randomness complexity*

$$n + k \log(d + 2) + 2 \log(1/\gamma) + O(1).$$

Proof sketch. Mimic the proof of Theorem 1, but use a PRG obtained by the standard nonconstructive argument (Appendix D). \square

The shifting and rounding steward S_0 can be generalized to achieve a tradeoff between low error ε' and low branching factor $|\Sigma|$ of the certification tree T_O . In particular, for any factorization $d = d_0 d_1$, one can reduce the error from $O(\varepsilon d)$ down to $O(\varepsilon d_0)$ at the cost of increasing the branching factor of T_O from $d + 2$ up to $(d_0 + 1)^{d_1} + 1$. This is achieved by simply partitioning the d coordinates into d_1 groups of d_0 coordinates and shifting each group individually; the details are in Appendix A. This immediately implies the following generalization of Theorem 1, which achieves a tradeoff between error and randomness complexity:

Theorem 4. *For any $n, k, d, d_0 \in \mathbb{N}$ with $d_0 \leq d$, for any $\varepsilon, \delta, \gamma > 0$, there exists a one-query $(O(\varepsilon d_0), k\delta + \gamma)$ -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ with randomness complexity*

$$n + O\left(\frac{kd \log(d_0 + 1)}{d_0} + (\log k) \log(1/\gamma)\right).$$

The total running time of the steward is $\text{poly}(n, k, d, \log(1/\varepsilon), \log(1/\gamma))$.

Recall from the introduction that if f_1, \dots, f_k are chosen nonadaptively, then we can reuse randomness and just union bound over the k functions. We now show that we can reuse the randomness in S_0 , as long as we union bound over *all the nodes* in the certification tree. (This is similar to the analysis of the Saks-Zhou steward, except that in the Saks-Zhou case, the branching factor of the tree is just 1. It is also similar to the analysis in [BH15].) This gives a steward with very low randomness complexity but large failure probability:

Theorem 5. *For any $n, k, d, d_0 \in \mathbb{N}$ with $d_0 \leq d$, for any $\varepsilon, \delta > 0$, there exists a one-query $(O(\varepsilon d_0), \delta')$ -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ with randomness complexity n , where*

$$\delta' \leq \exp\left(O\left(\frac{kd \log(d_0 + 1)}{d_0}\right)\right) \cdot \delta.$$

The total running time of the steward is $\text{poly}(n, k, d, \log(1/\varepsilon))$.

Proof. Assume without loss of generality that d is a multiple of d_0 and that $\delta < 1/2$. The steward is $S(X) \stackrel{\text{def}}{=} S_0(X, X, X, \dots, X)$, where S_0 is the steward of Section 2 generalized as in Appendix A. To prove correctness, fix any deterministic owner O . Let T_O be the block decision tree of Lemma 12. By Condition 1 of Lemma 12, from any internal node, if T_O reads X , the probability that it moves to the \perp child is at most δ . Therefore, by the union bound over all nodes, the probability that there is some node from which T_O would move to the \perp child upon reading X is at most the value δ' in the lemma statement. By Condition 2 of Lemma 12, if no node in T_O takes a \perp transition upon reading X , then $\max_i \|\mu_i - Y_i\|_\infty \leq O(\varepsilon d_0)$ in $O \leftrightarrow S(X)$. \square

6 Applications

6.1 Acceptance probabilities of Boolean circuits

A (ε, δ) -sampler for Boolean functions on n bits is a randomized oracle algorithm **Samp** such that for any Boolean function $C : \{0, 1\}^n \rightarrow \{0, 1\}$, if we let $\mu(C) \stackrel{\text{def}}{=} 2^{-n} \sum_x C(x)$, then

$$\Pr[|\mathbf{Samp}^C - \mu(C)| > \varepsilon] \leq \delta.$$

We will use a near-optimal sampler constructed by Goldreich and Wigderson [GW97]:

Lemma 6 ([GW97, Theorem 6.5]). *For every $n \in \mathbb{N}$ and every $\varepsilon, \delta > 0$, there is an (ε, δ) -sampler for Boolean functions on n bits that makes $O(\log(1/\delta)/\varepsilon^2)$ queries, uses $n + O(\log(1/\delta))$ random bits, and runs in time $\text{poly}(n, 1/\varepsilon, \log(1/\delta))$.*

Proof of Corollary 1. Let c be the constant under the $O(\cdot)$ of the error ε' in the steward of Theorem 1. When given parameters $n, k, \varepsilon, \delta$, let **Samp** be the Boolean $(\varepsilon/c, \delta/(2k))$ -sampler of Lemma 6, and say it uses m coins. Let **S** be the (ε, δ) -steward of Theorem 1 for k adaptively chosen $(\varepsilon/c, \delta/(2k))$ -concentrated functions $f_1, \dots, f_k : \{0, 1\}^m \rightarrow \mathbb{R}$. (So $\gamma = \delta/2$.) When given circuit C_i , define $f_i(X) = \mathbf{Samp}^{C_i}(X)$, i.e. the output \mathbf{Samp}^{C_i} with randomness X . Give f_i to **S**, and output the value Y_i that it returns.

Proof of correctness: The definition of a sampler implies that each f_i is $(\varepsilon/c, \delta/(2k))$ -concentrated at $\mu(C_i)$. Furthermore, each f_i is defined purely in terms of C_i , which is chosen based only on Y_1, \dots, Y_{i-1} . Therefore, the steward guarantee implies that with probability $1 - \delta$, every Y_i is within $\pm\varepsilon$ of $\mu(C_i)$.

Randomness complexity analysis: The number of bits m used by the sampler is $n + O(\log(k/\delta))$. Therefore, the number of bits used by the steward is

$$n + O(\log(k/\delta)) + O(k + (\log k) \log(1/\delta)) = n + O(k + (\log k) \log(1/\delta)).$$

Runtime analysis: The runtime of the steward is $\text{poly}(m, k, \log(1/\gamma)) = \text{poly}(n, k, \log(1/\delta))$. The runtime of the sampler is $\text{poly}(n, 1/\varepsilon, \log k, \log(1/\delta))$. The time required to evaluate each query of the sampler in round i is $O(\text{size}(C_i))$ (assuming we work with a suitable computational model and a suitable encoding of Boolean circuits.) The number of queries that the sampler makes in each round is $O(\log(k/\delta)/\varepsilon^2)$. Therefore, the total runtime of this algorithm is

$$O\left(\frac{\log k + \log(1/\delta)}{\varepsilon^2} \cdot \sum_{i=1}^k \text{size}(C_i)\right) + \text{poly}(n, k, 1/\varepsilon, \log(1/\delta)). \quad \square$$

6.2 Simulating a promise-BPP oracle

Theorem 6. *Suppose a search problem Π can be solved by a deterministic **promise-BPP**-oracle algorithm that runs in time T and makes k queries, and suppose that (regardless of previous oracle responses) each query of this algorithm can be decided by a randomized algorithm that runs in time T' , uses n coins, and has failure probability $1/3$. Then for any δ , Π can be solved by a randomized (non-oracle) algorithm that runs in time*

$$T + O(T' \cdot k \log(k/\delta)) + \text{poly}(n, k, \log(1/\delta)),$$

has randomness complexity

$$n + O(k + (\log k) \log(1/\delta)),$$

and has failure probability δ .

(Recall that search problems generalize decision problems and function problems. In reality, the theorem generalizes to just about any kind of “problem”, but we restrict ourselves to search problems for concreteness.) The theorem can easily be extended to randomized oracle algorithms by considering the problem of executing the randomized oracle algorithm using a given randomness string.

As a reminder, as discussed in Section 1.4, Theorem 6 would be trivial if it involved a **BPP** oracle instead of a **promise-BPP** oracle. Indeed, in the **BPP** case, the randomness can be reduced to just $n + O(\log k + \log(1/\delta))$. This is because a **BPP** algorithm is pseudodeterministic, so the randomness can be safely reused from one query to the next. A **promise-BPP** algorithm is not pseudodeterministic in general – it is only guaranteed to be pseudodeterministic on inputs that satisfy the promise.

Proof sketch of Theorem 6. Let B be the algorithm of Corollary 1 with $\varepsilon = 1/10$ and the desired failure probability δ . When the oracle algorithm makes query i , define $f_i(X)$ to be the value outputted by the **promise-BPP** algorithm on that query string using randomness X . Give B the “circuit” f_i . (The algorithm B treats the circuits as black boxes, so we don’t need to bother implementing f_i as a literal Boolean circuit; the important thing is that $f_i(X)$ can be evaluated in time T' .) When B outputs a value Y_i , give the oracle algorithm the response 0 if $Y_i < 1/2$ and 1 if $Y_i \geq 1/2$. \square

6.3 Simulating an APP oracle

Following Moser [Mos01], we model oracle access to $\varphi \in \mathbf{APP}$ by requiring the oracle algorithm to provide $w \in \{0, 1\}^n$ and a unary representation of $1/\varepsilon \in \mathbb{N}$; the oracle is guaranteed to respond with a value that is within $\pm\varepsilon$ of $\varphi(w)$. From these definitions, it is easy to show, for example, that $\mathbf{BPP}^{\mathbf{APP}} = \mathbf{BPP}$. Just like we did with **promise-BPP**, we now use our steward to construct a time- and randomness-efficient simulation of any algorithm with an oracle for **APP**.

Theorem 7. *Suppose $\varphi \in \mathbf{APP}$ and a search problem Π can be solved by a deterministic φ -oracle algorithm that runs in time T and makes k queries $(w_1, \varepsilon), \dots, (w_k, \varepsilon)$ (where w_i depends on previous oracle responses, but ε is the same for every query.) Let c be the constant under the $O(\cdot)$ in the error ε' in Theorem 1. Suppose that (regardless of the oracle responses) $\varphi(w_i)$ can be approximated to within $\pm\varepsilon/c$ by a randomized algorithm that runs in time T' , uses n coins, and has failure probability $1/3$. Then for any δ , Π can be solved by a randomized (non-oracle) algorithm that runs in time*

$$T + O(T' \cdot k \log(k/\delta)) + \text{poly}(n, k, \log(1/\delta)),$$

has randomness complexity

$$n + O(k + (\log k) \log(1/\delta)),$$

and has failure probability δ .

The proof of Theorem 7 is similar to the proofs of Corollary 1 and Theorem 6. The difference is that a sampler as defined previously is no longer quite the right tool for deterministic amplification; to amplify an **APP** algorithm, we are not trying to estimate the *acceptance probability* of a Boolean function, but rather the point where a $[0, 1]$ -valued function is *concentrated*. For this, we use an *averaging sampler*.

An *averaging (ε, δ) -sampler* for Boolean functions on n bits is an algorithm $\text{Samp} : \{0, 1\}^m \rightarrow (\{0, 1\}^n)^t$ such that for any Boolean function $C : \{0, 1\}^n \rightarrow \{0, 1\}$, if we let $\mu(C) \stackrel{\text{def}}{=} 2^{-n} \sum_x C(x)$

be the acceptance probability of C , then

$$\Pr_{X \in \{0,1\}^m} \left[\left| \mu(C) - \frac{1}{t} \sum_{i=1}^t C(\text{Samp}(X)_i) \right| > \varepsilon \right] \leq \delta.$$

(Note that an averaging sampler induces a sampler of a very specific form: query the oracle at several points and output the empirical mean.) We now show that an averaging sampler can be used to decrease the failure probability of a concentrated function by taking a *median*. This observation (in a different form) is due to Bellare, Goldreich, and Goldwasser [BGG93].

Lemma 7. *Suppose $f : \{0,1\}^n \rightarrow \mathbb{R}$ is (ε, δ_0) -concentrated at $\mu \in \mathbb{R}$ and $\text{Samp} : \{0,1\}^m \rightarrow (\{0,1\}^n)^t$ is an averaging (ε', δ) -sampler for Boolean functions on n bits, where $\varepsilon' + \delta_0 < 1/2$. Define $g : \{0,1\}^m \rightarrow \mathbb{R}$ by*

$$g(x) = \text{median}_{i \in [t]} f(\text{Samp}(x)_i).$$

Then g is (ε, δ) -concentrated at μ .

Proof. Let $C : \{0,1\}^n \rightarrow \{0,1\}$ be the indicator function for $\{x : |f(x) - \mu| \leq \varepsilon\}$. Then by the concentration of f , $2^{-n} \sum_x C(x) \geq 1 - \delta_0$. Therefore, by the averaging sampler condition, with probability $1 - \delta$ over x , $\frac{1}{t} \sum_i C(\text{Samp}(X)_i) \geq 1 - \delta_0 - \varepsilon' > 1/2$. If this is the case, then more than half of the values $f(\text{Samp}(x)_1), \dots, f(\text{Samp}(x)_t)$ are within $\pm \varepsilon$ of μ , which implies that their median is within $\pm \varepsilon$ of μ . \square

The following lemma gives the parameters achieved by the famous “random walk on expanders” averaging sampler; see, e.g., [Vad12, Corollary 4.41].

Lemma 8. *For every $n \in \mathbb{N}$ and every $\varepsilon, \delta > 0$, there is an averaging (ε, δ) -sampler for Boolean functions on n bits with $m \leq n + O(\log(1/\delta)/\varepsilon^2)$ and $t \leq O(\log(1/\delta)/\varepsilon^2)$, computable in time $\text{poly}(n, 1/\varepsilon, \log(1/\delta))$.*

Corollary 2 (Deterministic amplification for **APP**). *Suppose $\varphi \in \mathbf{APP}$ via an algorithm that on input (w, ε) uses n coins and t time steps to compute $\varphi(w) \pm \varepsilon$ with failure probability $1/3$. Then for any δ , is possible to compute $\varphi(w) \pm \varepsilon$ with failure probability δ using $O(t \log(1/\delta)) + \text{poly}(n, \log(1/\delta))$ time steps and $n + O(\log(1/\delta))$ coins.*

Proof. On input (w, ε) :

1. Let $\text{Samp} : \{0,1\}^m \rightarrow (\{0,1\}^n)^t$ be the averaging $(1/10, \delta)$ -sampler for Boolean functions on n bits of Lemma 8.
2. Define $f : \{0,1\}^n \rightarrow [0,1]$ by letting $f(X)$ be the output of the $1/3$ -error-probability algorithm for computing $\varphi(w) \pm \varepsilon$ on randomness X .
3. Pick $X \in \{0,1\}^m$ uniformly at random and return $\text{median}_{i \in [t]} f(\text{Samp}(X)_i)$.

Correctness follows immediately from Lemma 7, since f is $(\varepsilon, 1/3)$ -concentrated at $\varphi(w)$. Efficiency follows immediately from Lemma 8. \square

Proof of Theorem 7. By Corollary 2, there is an algorithm Φ for computing $\varphi(w_i) \pm \varepsilon/c$ with failure probability $\delta/(2k)$ that runs in time $O(T' \cdot \log(k/\delta)) + \text{poly}(n, \log k, \log(1/\delta))$ and uses $m \leq n + O(\log(k/\delta))$ coins. Let \mathbf{S} be the (ε, δ) -steward of Theorem 1 for k adaptively chosen $(\varepsilon/c, \delta/(2k))$ -concentrated functions $f_1, \dots, f_k : \{0,1\}^m \rightarrow \mathbb{R}$. (So $\gamma = \delta/2$.) When the oracle

algorithm makes query i about string w_i , let $f_i(X) = \Phi(w_i, \varepsilon/c, X)$ and give f_i to S . When S outputs a value Y_i , give it to the oracle algorithm.

Proof of correctness: Each f_i is $(\varepsilon/c, \delta/(2k))$ -concentrated at $\varphi(w_i)$. Furthermore, each f_i depends only on the previous oracle responses, i.e. Y_1, \dots, Y_{i-1} . Therefore, the steward guarantee implies that with probability $1 - \delta$, every Y_i is within $\pm\varepsilon$ of $\varphi(w_i)$. If this occurs, then the oracle algorithm is guaranteed to give a correct output.

Randomness complexity analysis: The number of bits used by the steward is

$$m + O(k + (\log k) \log(1/\delta)) = n + O(k + (\log k) \log(1/\delta)).$$

Runtime analysis: The runtime of the steward is $\text{poly}(m, k, \log(1/\gamma)) = \text{poly}(n, k, \log(1/\delta))$. Therefore, the total runtime is bounded by

$$T + k \cdot (O(T' \cdot \log(k/\delta)) + \text{poly}(n, \log k, \log(1/\delta))) + \text{poly}(n, k, \log(1/\delta)),$$

which is bounded by the expression in the theorem statement. \square

6.4 The Goldreich-Levin algorithm

Theorem 8 (Randomness-efficient Goldreich-Levin algorithm). *There is a randomized algorithm that, given oracle access to $F : \{0, 1\}^n \rightarrow \{-1, 1\}$ and given input parameters $\delta, \theta > 0$, outputs a list L of subsets of $[n]$ such that with probability $1 - \delta$,*

1. every U satisfying $|\widehat{F}(U)| \geq \theta$ is in L , and
2. every $U \in L$ satisfies $|\widehat{F}(U)| \geq \theta/2$.

The number of queries made by the algorithm is

$$O\left(\frac{n}{\theta^{11} \log(1/\theta)} \log\left(\frac{n}{\delta\theta}\right)\right),$$

the number of random bits used by the algorithm is

$$O(n + (\log n) \log(1/\delta)),$$

and the runtime of the algorithm is $\text{poly}(n, 1/\theta, \log(1/\delta))$.

For comparison, using standard techniques (the GW sampler, reusing randomness within each round of adaptivity), the Goldreich-Levin algorithm can be implemented in a straightforward way to use $O(\frac{n}{\theta^6} \log(\frac{n}{\delta\theta}))$ queries and $O(n^2 + n \log(\frac{n}{\delta\theta}))$ random bits. So our algorithm significantly improves the randomness complexity at the expense of substantially increasing the exponent of $1/\theta$ in the query complexity.

Toward proving Theorem 8, for a string $x \in \{0, 1\}^{\leq n}$, define

$$\mathcal{U}(x) = \{U \subseteq [n] : \forall j \leq |x|, j \in U \iff x_j = 1\}.$$

(That is, we think of $x \in \{0, 1\}^\ell$ as specifying $U \cap [\ell]$ in the natural way.) Define $W_x[F] = \sum_{U \in \mathcal{U}(x)} \widehat{F}(U)^2$. One of the key facts used in the standard Goldreich-Levin algorithm is that $W_x[F]$ can be estimated using few queries to F ; here, we use the GW sampler to improve the randomness efficiency of that estimation.

Lemma 9. *There is a randomized algorithm that, given oracle access to F and inputs $x \in \{0,1\}^{\leq n}$, $\varepsilon, \delta > 0$, estimates $W_x[F]$ to within $\pm\varepsilon$ with failure probability δ . The number of queries is $O(\log(1/\delta)/\varepsilon^2)$, the number of random bits is $O(n + \log(1/\delta))$, and the runtime is $\text{poly}(n, 1/\varepsilon, \log(1/\delta))$.*

Proof. Let $\ell = |x|$. As shown in the proof of [O'D14, Proposition 3.40],

$$W_x[F] = \mathbb{E}_{\substack{y, y' \in \{0,1\}^\ell \\ z \in \{0,1\}^{n-\ell}}} [F(y, z) \cdot F(y', z) \cdot \chi_x(y) \cdot \chi_x(y')],$$

where $\chi_x(y) \stackrel{\text{def}}{=} \prod_{j: x_j=1} (-1)^{y_j}$. Let $C : \{0,1\}^{n+\ell} \rightarrow \{0,1\}$ be the function

$$C(y, y', z) = \frac{1}{2} + \frac{1}{2} \cdot F(y, z) \cdot F(y', z) \cdot \chi_x(y) \cdot \chi_x(y'),$$

so that $W_x[F] = 2 \mathbb{E}_{y, y', z} [C(y, y', z)] - 1$. We can estimate the expectation of C to within $\pm\varepsilon/2$ with failure probability δ using the GW sampler of Lemma 6, which implies an estimate of $W_x[F]$ to within $\pm\varepsilon$. The number of queries made by the GW sampler is $O(\log(1/\delta)/\varepsilon^2)$, and each query to C can be evaluated by making 2 queries to F . The randomness complexity of the GW sampler is $n + \ell + O(\log(1/\delta))$, which is $O(n + \log(1/\delta))$. \square

The standard Goldreich-Levin algorithm proceeds by finding, for $\ell = 1$ to n , the set of all x with $|x| = \ell$ such that $W_x[F] \gtrsim \theta^2$. In each round, the algorithm estimates $W_x[F]$ for all strings x formed by appending a single bit to a string x' that was previously found to satisfy $W_{x'}[F] \gtrsim \theta^2$. This adaptive structure is exactly suited for saving random bits using a steward. To further drive down the randomness complexity, we reduce the number of rounds of adaptivity by appending $\log(1/\theta)$ bits at a time instead of 1 bit.

Proof of Theorem 8. Algorithm:

1. Let $u = \lfloor \log(1/\theta) \rfloor$, let $k = \lceil n/u \rceil$, and let $d = \lfloor 2^u \cdot 4/\theta^2 \rfloor$.
2. Let \mathbf{S} be a $(\theta^2/4, \delta)$ -steward for k adaptively chosen $(\varepsilon, \delta/(2n))$ -concentrated functions $f_1, \dots, f_k : \{0,1\}^m \rightarrow \mathbb{R}^d$, where $\varepsilon \geq \Omega(\theta^2/d)$ and m will become clear later.
3. Set $L_0 := \{\text{empty string}\}$.
4. For $i = 1$ to k :
 - (a) If $|L_{i-1}| > d/2^u$, abort and output “fail”.
 - (b) Observe that every string in L_{i-1} has length $\ell = u(i-1) < n$. Let x_1, \dots, x_t be the set of all strings obtained from strings in L_{i-1} by appending $\min\{u, n - \ell\}$ bits, so $t \leq 2^u |L_{i-1}| \leq d$.
 - (c) Define $f_i : \{0,1\}^m \rightarrow \mathbb{R}^t$ by letting $f_i(X)_j$ be the estimate of $W_{x_j}[F]$ to within $\pm\varepsilon$ provided by the algorithm of Lemma 9 with failure probability $\delta/(2dn)$ using randomness X . Observe that by the union bound, f_i is $(\varepsilon, \delta/(2n))$ -concentrated at $(W_{x_1}[F], \dots, W_{x_t}[F])$.
 - (d) By giving f_i to \mathbf{S} , obtain estimates μ_1, \dots, μ_t for $W_{x_1}[F], \dots, W_{x_t}[F]$.
 - (e) Set $L_i := \{x_j : \mu_j \geq \theta^2/2\}$.
5. Output $L \stackrel{\text{def}}{=} \bigcup_{x \in L_k} \mathcal{U}(x)$.

As hopefully became clear, m is the number of random bits used by the algorithm of Lemma 9. With probability $1 - \delta$, all of the responses of \mathbf{S} are accurate, i.e. every μ_j value is within $\pm\theta^2/4$ of the corresponding $W_{x_j}[F]$ value. Assume from now on that this has happened.

By the definition of L_i , every x in every L_i satisfies $W_x[F] \geq \theta^2/4$. By Parseval's theorem (see, e.g., [O'D14, Section 1.4]), this implies that $|L_i| \leq 4/\theta^2 \leq d/2^u$ for every i . Therefore, the algorithm does not abort. Let ℓ_i be the length of all the strings in L_i , so $\ell_i = ui$ for $i < k$ and $\ell_k = n$. Suppose $\widehat{F}(U)^2 \geq \theta^2$. By induction on i , the unique string $x \in \{0, 1\}^{\ell_i}$ with $U \in \mathcal{U}(x)$ is placed in L_i , because the estimate of $W_x[F]$ is at least $3\theta^2/4 > \theta^2/2$. This shows that $U \in L$. Conversely, if U ends up in L , then the estimate of $\widehat{F}(U)^2$ in iteration $i = n$ was at least $\theta^2/2$, so $\widehat{F}(U)^2 \geq \theta^2/4$. This completes the proof of correctness of the algorithm.

Now, observe that the total number of queries to F is at most kd times the $O(\log(nd/\delta)/\varepsilon^2)$ queries that the algorithm of Lemma 9 makes, i.e. the total number of queries to F is

$$O\left(\frac{kd^3 \log(nd/\delta)}{\theta^2}\right) = O\left(\frac{n}{\theta^{11} \log(1/\theta)} \log\left(\frac{n}{\delta\theta}\right)\right).$$

The randomness complexity of the algorithm is just the randomness complexity of \mathbf{S} . We will use the steward of Theorem 1 with $\gamma = \delta/2$, so the randomness complexity is $m + O(k \log(d+1) + (\log k) \log(1/\delta))$. Since $m \leq O(n + \log(n/(\delta\theta)))$, the total randomness complexity is

$$O\left(n + \frac{n}{\log(1/\theta)} \log(1/\theta) + (\log n) \log(1/\delta) + \log(1/\theta)\right) = O(n + (\log n) \log(1/\delta) + \log(1/\theta)).$$

To get rid of the $\log(1/\theta)$ term as claimed in the theorem statement, just notice that we can assume without loss of generality that $\theta \geq 2^{-n+1}$, because any nonzero Fourier coefficient of a $\{-1, 1\}$ -valued function has absolute value at least 2^{-n+1} . The total runtime of the algorithm is clearly $\text{poly}(n, 1/\theta, \log(1/\delta))$. \square

7 Randomness complexity lower bounds

7.1 Lower bound for pseudorandom generation stewards

Definition 7. A *pseudorandom generation steward* is a one-query steward with the additional property that in each round i , the value Y_i that the steward gives to \mathbf{O} is simply the query response $f_i(X_i)$ that it receives.

Notice that this definition is more general than the usual notion of a pseudorandom generator in two respects. First, the query point X_i may depend on the previous responses $f_1(X_1), \dots, f_{i-1}(X_{i-1})$. Second, rather than making a statistical indistinguishability requirement, we merely impose the standard steward correctness requirement, i.e. with high probability, every Y_i has low ℓ_∞ error.

We begin with an elementary lemma. Let $H(\cdot)$ denote Shannon entropy. If a random variable X takes values in a set of size t , then $H(X) \leq \log_2 t$. The following lemma is a partial converse: if $H(X)$ is much smaller than $\log_2 t$, then there is a set of size t that X is likely to land in.

Lemma 10. *Suppose X is a discrete random variable. Let W be the set consisting of the t most likely values of X . Then*

$$\Pr[X \in W] \geq 1 - \frac{H(X)}{\log_2 t}.$$

Proof. Let p be the probability mass function of X , so that $H(X) = \mathbb{E}[-\log_2(p(X))]$. By Markov's inequality,

$$\Pr[-\log_2(p(X)) > \log_2 t] \leq \frac{H(X)}{\log_2 t}.$$

Therefore, if we let $W' = \{x : p(x) \geq 1/t\}$,

$$\Pr[X \in W'] \geq 1 - \frac{H(X)}{\log_2 t}.$$

Finally, $W' \subseteq W$, because there can be at most t values x such that $p(x) \geq 1/t$. \square

Theorem 9. *Suppose S is an m -coin pseudorandom generation (ε', δ') -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$. Then*

$$m \geq (1 - \delta') \cdot (n - \log_2(2/\delta)) \cdot k.$$

Proof. Without loss of generality, assume $d = 1$. Fix an injective $(\varepsilon, 0)$ -concentrated function $f : \{0, 1\}^n \rightarrow \mathbb{R}$. (E.g., we could set the binary expansion of $f(x)$ to be $\log(1/\varepsilon)$ zeroes concatenated with x .) Let O_0 be the owner who always chooses $f_i = f$ for every i . Let $X_i \in \{0, 1\}^n$ be the query point that S chooses in round i of $\mathsf{O}_0 \leftrightarrow \mathsf{S}$, so that X_1, \dots, X_k are functions of the internal randomness of S .

Since S uses only m coins, $H(X_1, \dots, X_k) \leq m$. By the chain rule, this implies that there is some $j \in [k]$ such that

$$H(X_j \mid X_1, \dots, X_{j-1}) \leq \frac{m}{k}.$$

Let O be the following owner:

1. In each round $i < j$, pick $f_i = f$, and obtain the value Y_i . Compute $x_i = f_i^{-1}(Y_i)$.
2. Let $p(x)$ be the distribution $p(x) = \Pr[X_j = x \mid X_1 = x_1, \dots, X_{j-1} = x_{j-1}]$.
3. In round j , let W be the set of the $\lfloor \delta 2^n \rfloor$ points in $\{0, 1\}^n$ to which p assigns the most mass. Let $f_j : \{0, 1\}^n \rightarrow \mathbb{R}$ be the function

$$f_j(x) = \begin{cases} 0 & \text{if } x \notin W \\ 2\varepsilon' & \text{if } x \in W. \end{cases}$$

Observe that f_j is $(0, \delta)$ -concentrated at zero. Give f_j to S .

4. In rounds $i > j$, choose f_i to be (say) the constant zero function.

The probability that S fails in $\mathsf{O} \leftrightarrow \mathsf{S}$ is precisely $\Pr[X_j \in W]$, which we now lower bound. We can write

$$\Pr[X_j \in W] = \mathbb{E}_{(x_1, \dots, x_{j-1}) \sim (X_1, \dots, X_{j-1})} [\Pr[X_j \in W \mid X_1 = x_1, \dots, X_{j-1} = x_{j-1}]].$$

By Lemma 10 and the definition of W , for any particular values x_1, \dots, x_{j-1} ,

$$\Pr[X_j \in W \mid X_1 = x_1, \dots, X_{j-1} = x_{j-1}] \geq 1 - \frac{H(X_j \mid X_1 = x_1, \dots, X_{j-1} = x_{j-1})}{\log_2 \lfloor \delta 2^n \rfloor}.$$

Therefore, by linearity of expectation,

$$\begin{aligned}\Pr[X_j \in W] &\geq 1 - \frac{H(X_j \mid X_1, \dots, X_{j-1})}{\log_2 \lceil \delta 2^n \rceil} \\ &\geq 1 - \frac{m}{k \log_2 \lceil \delta 2^n \rceil}.\end{aligned}$$

If $\delta < 2^{-n}$, the theorem statement is trivial, so assume that $\delta \geq 2^{-n}$. In this case, $\log_2 \lceil \delta 2^n \rceil \geq n - \log_2(2/\delta)$, so

$$\Pr[X_j \in W] \geq 1 - \frac{m}{(n - \log_2(2/\delta)) \cdot k}.$$

By the correctness of the steward, this value must be at most δ' . Rearranging completes the proof. \square

7.2 Lower bound for one-query stewards

To understand the following lemma, imagine the perspective of \mathcal{O} after $i - 1$ rounds of $\mathcal{O} \leftrightarrow \mathcal{S}(Z)$, where Z was chosen uniformly at random from $\{0, 1\}^m$. Let R be the set of z such that the hypothesis that $Z = z$ is compatible with everything that \mathcal{O} has seen so far. Then at this point, \mathcal{O} 's posterior distribution for Z is uniform over R . The following lemma says that with respect to this posterior distribution, \mathcal{O} can choose f_i such that either \mathcal{O} will learn $\Omega(1)$ bits of information about Z based on Y_i , or else \mathcal{S} will have a failure probability of $\Omega(1)$ in round i .

Lemma 11. *Suppose \mathcal{S} is a one-query m -coin (ε', δ') -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}$ and \mathcal{O} is a deterministic owner. Fix $i \in [k]$. For a function $g : \{0, 1\}^n \rightarrow \mathbb{R}$, let $\mathcal{O}[g]$ be the owner that simulates \mathcal{O} for rounds $1, 2, \dots, i - 1$, but chooses g in round i regardless of what \mathcal{O} would have chosen. Let $R \subseteq \{0, 1\}^m$ be a nonempty set such that the transcript of the first $i - 1$ rounds of $\mathcal{O} \leftrightarrow \mathcal{S}(Z)$ is the same for every $Z \in R$. Assume $\delta \geq 2^{-n}$. Then there exists g that is (ε, δ) -concentrated at μ such that either*

1. $\max_{y \in \mathbb{R}} \Pr_{Z \in R}[Y_i = y \text{ in } \mathcal{O}[g] \leftrightarrow \mathcal{S}(Z)] \leq 0.8$, or
2. $\Pr_{Z \in R}[|Y_i - \mu_i| > \varepsilon' \text{ in } \mathcal{O}[g] \leftrightarrow \mathcal{S}(Z)] \geq 0.2$.

Proof. For each $j \in \mathbb{Z}$, let $g_j : \{0, 1\}^n \rightarrow \mathbb{R}$ be constant at εj . If some g_j satisfies Condition 1, we're done. So assume that for each g_j , there is some $y_j \in \mathbb{R}$ such that $\Pr_{Z \in R}[Y_j = y_j \text{ in } \mathcal{O}[g] \leftrightarrow \mathcal{S}(Z)] > 0.8$. If y_j does not depend on j , then since $0.2 < 0.8$, there is some g_j that satisfies Condition 2, so we are again done. Therefore, assume there is some j such that $y_j \neq y_{j+1}$.

Define $q : R \rightarrow \{0, 1\}^n$ by letting $q(Z) =$ the value X_i chosen by \mathcal{S} in $\mathcal{O} \leftrightarrow \mathcal{S}(Z)$. First, assume there is some x^* such that $\Pr_{Z \in R}[q(Z) = x^*] \geq 0.4$. For $s \in \{\pm 1\}$, define $g^s : \{0, 1\}^n \rightarrow \mathbb{R}$ by

$$g^s(x) = \begin{cases} 0 & \text{if } x = x^* \\ s \cdot 2\varepsilon' & \text{otherwise.} \end{cases}$$

Then $g^s(x)$ is $(0, 2^{-n})$ -concentrated at $s \cdot 2\varepsilon'$. Let \mathcal{O}' be the randomized owner that tosses a coin to decide whether to simulate $\mathcal{O}[g^{+1}]$ or $\mathcal{O}[g^{-1}]$. Then when $Z \in R$ is chosen uniformly at random, in $\mathcal{O}' \leftrightarrow \mathcal{S}(Z)$, there is a 0.4 chance that $f_i(X_i) = 0$, in which case $\mathcal{S}(Z)$ has only a 50% chance of correctly guessing s . This shows that $\Pr_{Z \in R}[|Y_i - \mu_i| > \varepsilon' \text{ in } \mathcal{O}' \leftrightarrow \mathcal{S}(Z)] \geq 0.2$, and hence either g^{+1} or g^{-1} satisfies Condition 2, so we are again done. Therefore, assume that for every x^* , $\Pr_{Z \in R}[q(Z) = x^*] < 0.4$.

For $t \in \{j, j + 1\}$, let

$$A_t = \{Z \in R : Y_i = y_t \text{ in } \mathbf{O}[g_t] \leftrightarrow \mathbf{S}(Z)\},$$

so that $|A_t| > 0.8|R|$. We define g by the following greedy algorithm. Two players, which we identify with A_j and A_{j+1} , alternate taking turns. When it is A_t 's turn, she finds the string $x \in \{0, 1\}^n$ such that $g(x)$ is not yet defined that maximizes $q^{-1}(x) \cap A_t$, and defines $g(x) = \varepsilon t$. This continues for 2^n turns until g is defined everywhere.

Clearly, g thus defined is $(\varepsilon, 0)$ -concentrated. We will show that g satisfies Condition 1. Proof: Say $z \in \{0, 1\}^m$ is *good for* A_t if $z \in A_t$ and $g(q(z)) = \varepsilon t$. In these terms, on A_t 's turn, she defines g on one more point in order to maximize the number of z that become good for A_t . Say that $z \in \{0, 1\}^m$ is *bad for* A_t if $z \in A_t$ and $g(q(z)) \neq \varepsilon t$. When it is not A_t 's turn, some z may become bad for A_t , but the crucial point is that the number of z that become bad for A_t is *at most* the number of z that became *good* for A_t in the *previous* turn (simply because of the greedy choice that A_t made in the previous turn.) This would show that half of A_t is good for A_t , except for one annoyance: the first turn, where some z become bad for the second player with no corresponding previous turn, when z became good. But we already showed that for every x , $|q^{-1}(x)| \leq 0.4|R|$, so the first turn does not matter too much: at the end of the construction, for each t , the number of z that are good for A_t is at least

$$\frac{1}{2}(|A_t| - 0.4|R|) \geq \frac{1}{2}(0.8|R| - 0.4|R|) = 0.2|R|.$$

By construction, if z is good for A_t , then $Y_i = y_t$ in $\mathbf{O}[g] \leftrightarrow \mathbf{S}(z)$. Therefore, for each t , $\Pr_{Z \in R}[Y_i = y_t \text{ in } \mathbf{O}[g] \leftrightarrow \mathbf{S}(Z)] \geq 0.2$, which implies Condition 1 since $y_j \neq y_{j+1}$. \square

Having proved Lemma 11, we are ready to prove our randomness complexity lower bound. The idea is that \mathbf{O} will spend the first $k - 1$ rounds learning as much information as possible about \mathbf{S} 's randomness string using Lemma 11 (unless she gets lucky and is able to cause \mathbf{S} to have an $\Omega(1)$ failure probability in one of these rounds, in which case she will take the opportunity.) Then, in round k , \mathbf{O} uses everything she's learned about \mathbf{S} 's randomness string to choose f_k so as to maximize \mathbf{S} 's failure probability in that round.

Theorem 10. *Suppose \mathbf{S} is a one-query m -coin (ε', δ') -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$. Assume $\delta' < 0.2$ and $\delta \geq 2^{-n}$. Then $m \geq n + \Omega(k) - \log_2(\delta'/\delta)$.*

Proof. Without loss of generality, assume $d = 1$. Let \mathbf{O} be the following owner:

1. For $i = 1$ to k :
 - (a) Let y_1, y_2, \dots, y_{i-1} be the responses received so far.
 - (b) Let $R \subseteq \{0, 1\}^m$ be the set of z such that in $\mathbf{O} \leftrightarrow \mathbf{S}(z)$, $Y_j = y_j$ for every $j < i$. (By induction, we have already defined the behavior of \mathbf{O} in rounds $1, 2, \dots, i - 1$, so R is well-defined. In other words, R is the set of z that are compatible with what \mathbf{O} has seen so far.)
 - (c) If $i < k$:
 - i. Choose $f_i = g$, where g is the function guaranteed by Lemma 11. (Again, \mathbf{O} is already defined and deterministic for rounds $1, 2, \dots, i - 1$, so we can sensibly apply the lemma.)

(d) Otherwise, if $i = k$:

- i. Pick $S \subseteq R$, $|S| = \min\{\lfloor \delta 2^n \rfloor, |R|\}$ uniformly at random, pick $s \in \{\pm 1\}$ independently and uniformly at random, and choose

$$f_k(x) = \begin{cases} 0 & \text{if } x \in q(S) \\ s \cdot 2\varepsilon' & \text{otherwise.} \end{cases}$$

(Note that f_k is $(0, \delta)$ -concentrated at $s \cdot 2\varepsilon'$, because $|q(S)| \leq |S| \leq \delta 2^n$.)

To analyze O , in $\mathsf{O} \leftrightarrow \mathsf{S}$, say that O *tries to win* in round i if either $i = k$ or else $i < k$ and the function f_i chosen satisfies Condition 2 in Lemma 11. For a string $z \in \{0, 1\}^m$, let $w(z) \in [k]$ be the index of the first round in which O tries to win in $\mathsf{O} \leftrightarrow \mathsf{S}(z)$, and let $\tau(z)$ be the transcript of rounds $1, 2, \dots, w(z) - 1$ in $\mathsf{O} \leftrightarrow \mathsf{S}(z)$. Note that since O is deterministic in rounds $1, 2, \dots, k - 1$, $w(z)$ and $\tau(z)$ are not random variables. Define an equivalence relation on $\{0, 1\}^m$ by saying that $z \sim z'$ if and only if $\tau(z) = \tau(z')$. Say O uses v random bits. We first show that for each equivalence class \bar{z} ,

$$\Pr_{Z \in \bar{z}, V \in \{0, 1\}^v} [\|Y_{w(\bar{z})} - \mu_{w(\bar{z})}\|_\infty > \varepsilon' \text{ in } \mathsf{O}(V) \leftrightarrow \mathsf{S}(Z)] \geq \min\{0.2, \delta \cdot (1/0.8)^{k-1} \cdot 2^{n-m-2}\}. \quad (1)$$

Proof: Observe that in round $w(\bar{z})$, O 's set R is precisely \bar{z} . If $w(\bar{z}) < k$, then Condition 2 of Lemma 11 immediately implies that the failure probability in Equation 1 is at least 0.2. Suppose instead that $w(\bar{z}) = k$. Then in every previous round, O did not try to win, i.e. O chose a function satisfying Condition 1 of Lemma 11. This implies that in every previous round, O 's set R decreased in size by a factor of 0.8. So at the beginning of round k , $|R| \leq 0.8^{k-1} \cdot 2^m$. The probability (over $Z \in \bar{z}$) that S chooses X_k such that $f_k(X_k) = 0$ is

$$\begin{aligned} \frac{|S|}{|R|} &= \frac{\min\{\lfloor \delta 2^n \rfloor, |R|\}}{|R|} \\ &\geq \min\left\{1, \delta 2^{n-m-1} (1/0.8)^{k-1}\right\}. \end{aligned}$$

Conditioned on $f_k(X_k) = 0$, the probability of the event in Equation 1 is at least 0.5, because conditioned on $f_k(X_k) = 0$, s is independent of everything S has seen. Therefore, the probability of the event in Equation 1 is at least $\min\{0.5, \delta 2^{n-m-2} (1/0.8)^{k-1}\}$, completing the proof of Equation 1.

Now, to prove the theorem, observe that

$$\begin{aligned} \delta' &\geq \Pr_{Z \in \{0, 1\}^m, V \in \{0, 1\}^v} [\max_i \|\mu_i - Y_i\|_\infty > \varepsilon' \text{ in } \mathsf{O}(V) \leftrightarrow \mathsf{S}(Z)] \\ &\geq \Pr_{Z \in \{0, 1\}^m, V \in \{0, 1\}^v} [\|\mu_{w(Z)} - Y_{w(Z)}\|_\infty > \varepsilon' \text{ in } \mathsf{O}(V) \leftrightarrow \mathsf{S}(Z)] \\ &= \sum_{\bar{z}} \Pr_{Z \in \{0, 1\}^m} [Z \in \bar{z}] \cdot \Pr_{Z' \in \bar{z}, V \in \{0, 1\}^v} [\|\mu_{w(\bar{z})} - Y_{w(\bar{z})}\|_\infty > \varepsilon' \text{ in } \mathsf{O}(V) \leftrightarrow \mathsf{S}(Z')] \\ &\geq \sum_{\bar{z}} \Pr_{Z \in \{0, 1\}^m} [Z \in \bar{z}] \cdot \min\{0.2, \delta \cdot (1/0.8)^{k-1} 2^{n-m-2}\} \\ &= \min\{0.2, \delta \cdot (1/0.8)^{k-1} 2^{n-m-2}\}. \end{aligned}$$

We assumed that $\delta' < 0.2$, so we can conclude that $\delta' \geq \delta \cdot (1/0.8)^{k-1} 2^{n-m-2}$. Rearranging proves that

$$\begin{aligned} m &\geq (n - 2) + (k - 1) \log_2(1/0.8) - \log_2(\delta'/\delta) \\ &\geq n + \Omega(k) - \log_2(\delta'/\delta), \end{aligned}$$

completing the proof. \square

8 Directions for further research

The problem of randomness stewardship is fundamental, and the main open problem left by this work is to construct optimal stewards. The following are examples of concrete questions along these lines.

- Does every one-query steward with failure probability $\delta' \leq O(k\delta)$ have randomness complexity $n + \Omega(k \log(d + 1))$? (Is the randomness complexity of our main steward near-optimal?)
- Does there exist a one-query $(O(\varepsilon), k\delta + 0.1)$ -steward with randomness complexity $n + O(k \log(d + 1))$? (Can the error of our main steward be improved?)

We explained in this work how the steward model captures some older derandomization constructions, and we gave new applications of stewards. We hope that future researchers find more connections and applications.

9 Acknowledgments

We thank David Zuckerman for observations about block decision trees. We thank an anonymous reviewer for pointing out Impagliazzo and Zuckerman’s previous work on this subject [IZ89, Imp92].

References

- [Arm98] R. Armoni. On the derandomization of space-bounded computations. In *Randomization and Approximation Techniques in Computer Science*, pages 47–59. Springer, 1998.
- [Bel92] M. Bellare. A technique for upper bounding the spectral norm with applications to learning. In *Proceedings of the 5th Annual Workshop on Computational Learning Theory*, pages 62–70. ACM, 1992.
- [BF99] H. Buhrman and L. Fortnow. One-sided versus two-sided error in probabilistic computation. In *Annual Symposium on Theoretical Aspects of Computer Science*, pages 100–109. Springer, 1999.
- [BGG93] M. Bellare, O. Goldreich, and S. Goldwasser. Randomness in interactive proofs. *Computational Complexity*, 3(4):319–354, 1993.
- [BH15] A. Blum and M. Hardt. The ladder: A reliable leaderboard for machine learning competitions. In *International Conference on Machine Learning, ICML ’15*, pages 1006–1014, 2015.
- [BJT04] N. H. Bshouty, J. C. Jackson, and C. Tamon. More efficient PAC-learning of DNF with membership queries under the uniform distribution. *Journal of Computer and System Sciences*, 68(1):205–234, 2004.
- [BNS⁺16] R. Bassily, K. Nissim, A. Smith, T. Steinke, U. Stemmer, and J. Ullman. Algorithmic stability for adaptive data analysis. In *Proceedings of the 48th Annual Symposium on Theory of Computing, STOC ’16*, pages 1046–1059. ACM, 2016.
- [CLN⁺16] R. Cummings, K. Ligett, K. Nissim, A. Roth, and Z. S. Wu. Adaptive learning with robust generalization guarantees. In *Proceedings of the 29th Conference on Learning Theory, COLT ’16*, pages 23–26, 2016.

- [DFH⁺15a] C. Dwork, V. Feldman, M. Hardt, T. Pitassi, O. Reingold, and A. Roth. Generalization in adaptive data analysis and holdout reuse. In *Advances in Neural Information Processing Systems*, pages 2350–2358, 2015.
- [DFH⁺15b] C. Dwork, V. Feldman, M. Hardt, T. Pitassi, O. Reingold, and A. Roth. The reusable holdout: Preserving validity in adaptive data analysis. *Science*, 349(6248):636–638, 2015.
- [DFH⁺15c] C. Dwork, V. Feldman, M. Hardt, T. Pitassi, O. Reingold, and A. L. Roth. Preserving statistical validity in adaptive data analysis. In *Proceedings of the 47th Annual Symposium on Theory of Computing*, STOC '15, pages 117–126. ACM, 2015.
- [DORS08] Y. Dodis, R. Ostrovsky, L. Reyzin, and A. Smith. Fuzzy extractors: How to generate strong keys from biometrics and other noisy data. *SIAM journal on computing*, 38(1):97–139, 2008.
- [GG11] E. Gat and S. Goldwasser. Probabilistic search algorithms with unique answers and their cryptographic applications. In *Electronic Colloquium on Computational Complexity (ECCC)*, volume 18, page 136, 2011.
- [GL89] O. Goldreich and L. A. Levin. A hard-core predicate for all one-way functions. In *Proceedings of the 21st Annual ACM Symposium on Theory of Computing*, STOC '89, pages 25–32. ACM, 1989.
- [GW97] O. Goldreich and A. Wigderson. Tiny families of functions with random properties: A quality-size trade-off for hashing. *Random Structures & Algorithms*, 11(4):315–343, December 1997.
- [HU14] M. Hardt and J. Ullman. Preventing false discovery in interactive data analysis is hard. In *Proceedings of the 55th Annual Symposium on Foundations of Computer Science*, FOCS '14, pages 454–463. IEEE, 2014.
- [HU17] W. M. Hoza and C. Umans. Targeted pseudorandom generators, simulation advice generators, and derandomizing logspace. In *Proceedings of the 49th Annual Symposium on Theory of Computing*, STOC '17, pages 629–640. ACM, 2017.
- [Imp92] R. Impagliazzo. *Pseudo-random generators for cryptography and for randomized algorithms*. PhD thesis, University of California, Berkeley, 1992.
- [INW94] R. Impagliazzo, N. Nisan, and A. Wigderson. Pseudorandomness for network algorithms. In *Proceedings of the 26th Annual Symposium on Theory of Computing*, STOC '94, pages 356–364. ACM, 1994.
- [IZ89] R. Impagliazzo and D. Zuckerman. How to recycle random bits. In *Proceedings of the 30th Annual Symposium on Foundations of Computer Science*, FOCS '89, pages 248–253. IEEE, 1989.
- [KM93] E. Kushilevitz and Y. Mansour. Learning decision trees using the Fourier spectrum. *SIAM Journal on Computing*, 22(6):1331–1348, 1993.
- [KRC00] V. Kabanets, C. Rackoff, and S. Cook. Efficiently approximable real-valued functions. In *Electronic Colloquium on Computational Complexity (ECCC) Report TR00*, volume 34, 2000.

- [Lev93] L. A. Levin. Randomness and nondeterminism. *Journal of Symbolic Logic*, 58(3):1102–1103, 1993.
- [Mos01] P. Moser. Relative to **P promise-BPP** equals **APP**. In *Electronic Colloquium on Computational Complexity (ECCC) Report TR01*, volume 68, 2001.
- [Nis92] N. Nisan. Pseudorandom generators for space-bounded computation. *Combinatorica*, 12(4):449–461, 1992.
- [NN93] J. Naor and M. Naor. Small-bias probability spaces: Efficient constructions and applications. *SIAM journal on computing*, 22(4):838–856, 1993.
- [NZ96] N. Nisan and D. Zuckerman. Randomness is linear in space. *Journal of Computer and System Sciences*, 52(1):43–52, 1996.
- [O’D14] R. O’Donnell. *Analysis of boolean functions*. Cambridge University Press, 2014.
- [RR99] R. Raz and O. Reingold. On recycling the randomness of states in space bounded computation. In *Proceedings of the 31st Annual ACM Symposium on Theory of Computing*, STOC ’99, pages 159–168. ACM, 1999.
- [SU15] T. Steinke and J. Ullman. Interactive fingerprinting codes and the hardness of preventing false discovery. In *Proceedings of the 28th Conference on Learning Theory*, COLT ’15, pages 1588–1628, 2015.
- [SZ99] M. Saks and S. Zhou. $\mathbf{BP}_H\mathbf{SPACE}(S) \subseteq \mathbf{DSPACE}(S^{3/2})$. *Journal of Computer and System Sciences*, 58(2):376–403, 1999.
- [Vad12] S. P. Vadhan. *Pseudorandomness*, volume 56. Now, 2012.

A Generalized shifting and rounding algorithm

In this section, we show how to generalize the steward S_0 to achieve a tradeoff between its error and the branching factor of the certification tree T_O . Fix a factorization $d = d_0 d_1$. Partition $[d]$ as $[d] = J_1 \cup J_2 \cup \dots \cup J_{d_1}$, where $|J_t| = d_0$ for each t . Instead of partitioning \mathbb{R} into intervals of length $2(d+1)\varepsilon$, partition \mathbb{R} into intervals of length $2(d_0+1)\varepsilon$. Let \mathcal{I} denote the set of these intervals. The following algorithm for computing Y_i from W_i generalizes that of Section 2.1:

1. For each $t \in [d_1]$:
 - (a) Find $\Delta_{it} \in [d_0+1]$ such that for every $j \in J_t$, there is a single interval in \mathcal{I} that entirely contains $[W_{ij} + (2\Delta_{it} - 1)\varepsilon, W_{ij} + (2\Delta_{it} + 1)\varepsilon]$. (Such a Δ_{it} exists by Lemma 1.)
 - (b) For every $j \in J_t$, set $Y_{ij} = \text{Round}(W_{ij} + 2\Delta_{it}\varepsilon)$.

The following lemma is the appropriate generalization of Lemma 2:

Lemma 12. *Assume $\delta < 1/2$. Let $\Sigma = [d_0+1]^{d_1} \cup \{\perp\}$. For any deterministic owner O , there exists a (k, n, Σ) block decision tree T_O with the following properties.*

1. For any internal node v , $\Pr_{X \in \{0,1\}^n}[v(X) = \perp] \leq \delta$.
2. Fix $X_1, \dots, X_k \in \{0,1\}^n$, and suppose that the path from the root to $T_O(X_1, \dots, X_k)$ does not include any \perp nodes. Then in $O \leftrightarrow S_0(X_1, \dots, X_k)$, $\max_i \|Y_i - \mu_i\|_\infty \leq O(d_0\varepsilon)$.

The proof of Lemma 12 is essentially the same as the proof of Lemma 2; we record the details below.

Proof of Lemma 12. For a vector $Y \in \mathbb{R}^d$ and function $f : \{0, 1\}^n \rightarrow \mathbb{R}^d$ that is (ε, δ) -concentrated at some point, say that a vector $(\Delta_1, \dots, \Delta_{d_1})$ is f -compatible with Y if $Y_j = \text{Round}(\mu(f)_j + 2\Delta_t \varepsilon)$ for every $t \in [d_1]$ and every $j \in J_t$. Just for the analysis, let S'_0 be the following (many-query) steward:

1. For $i = 1$ to k :

- (a) Give f_i to S_0 , allowing it to make its one query and choose its output vector $Y_i \in \mathbb{R}^d$.
- (b) Query f_i at *every* point in its domain, thereby learning the entire function.
- (c) Compute

$$\widehat{\Delta}_i = \begin{cases} \text{the first } (\Delta_1, \dots, \Delta_{d_1}) \in [d_0 + 1]^{d_1} \text{ } f_i\text{-compatible with } Y_i & \text{if any exist} \\ \perp & \text{otherwise.} \end{cases}$$

(d) Output $\widehat{Y}_i = (\widehat{Y}_{i1}, \dots, \widehat{Y}_{id})$, where for each $t \in [d_1]$ and each $j \in J_t$,

$$\widehat{Y}_{ij} = \begin{cases} \text{Round}(\mu(f_i)_j + 2\widehat{\Delta}_{it}\varepsilon) & \text{if } \widehat{\Delta}_i \neq \perp \\ 0 & \text{otherwise.} \end{cases}$$

The definition of $T_{\mathcal{O}}$ is exactly the same as in the proof of Lemma 2, except that S'_0 now refers to the above steward. To prove Condition 1 in the lemma statement, we must show that in each round of $\mathcal{O} \leftrightarrow S'_0$, $\Pr[\widehat{\Delta}_i = \perp] \leq \delta$. Indeed, by concentration, with probability $1 - \delta$, for every j , $|W_{ij} - \mu(f_i)_j| \leq \varepsilon$. In this case, by the construction of S_0 , $W_{ij} + 2\Delta_{it}\varepsilon$ and $\mu(f_i)_j + 2\Delta_{it}\varepsilon$ are in the same interval in \mathcal{I} for every $t \in [d_1]$ and every $j \in J_t$. Therefore, in this case, there is at least one vector $(\Delta_1, \dots, \Delta_{d_1})$ that is f_i -compatible with Y_i , namely the vector of Δ_{it} values used by S_0 . To prove Condition 2 in the lemma statement, suppose the path from the root node to $T_{\mathcal{O}}(X_1, \dots, X_k)$ does not include any \perp nodes. Then in $\mathcal{O} \leftrightarrow S'_0(X_1, \dots, X_k)$, for every i , $\widehat{\Delta}_i \neq \perp$. This implies that every Y_{ij} is of the form $\text{Round}(\mu(f_i)_j + 2\widehat{\Delta}_{it}\varepsilon)$ for some $\widehat{\Delta}_{it} \in [d_0 + 1]$. Therefore, $|Y_{ij} - \mu(f_i)_j| \leq 3(d_0 + 1)\varepsilon$, since $2\widehat{\Delta}_{it}\varepsilon \leq 2(d_0 + 1)\varepsilon$ and rounding introduces at most $(d_0 + 1)\varepsilon$ additional error. Just as in the proof of Lemma 2, the same bound holds in $\mathcal{O} \leftrightarrow S_0$. Finally, since $\delta < 1/2$, $\|\mu(f_i) - \mu_i\|_{\infty} \leq 2\varepsilon$, so by the triangle inequality, for every i , $\|Y_i - \mu_i\|_{\infty} \leq 3(d_0 + 1)\varepsilon + 2\varepsilon = (3d_0 + 5)\varepsilon$. \square

B The Saks-Zhou steward

In this section, for completeness, we give the description and analysis of the Saks-Zhou steward. This algorithm and analysis are the same in spirit as what appears in [SZ99], but the presentation has been changed to match our framework. None of our results use this steward, but it is interesting to see how the stewards compare.

Proposition 1. *For any $n, k, d \in \mathbb{N}$ and any $\varepsilon, \delta, \gamma > 0$, there exists a one-query $(O(kd\varepsilon/\gamma), k\delta + \gamma)$ -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ with randomness complexity*

$$n + O(k \log k + k \log d + k \log(1/\gamma)).$$

The total running time of the steward is $\text{poly}(n, k, d, \log(1/\varepsilon), \log(1/\gamma))$.

Proof. Let u be the smallest power of two such that $u \geq 2kd/\gamma$. (The only reason we choose a power of two is so that we can cleanly draw a uniform random element of $[u]$ using $\log u$ random bits.) Partition \mathbb{R} into half-open intervals of length $\ell = u\varepsilon$. For $w \in \mathbb{R}$, let $\text{Round}(w)$ be the midpoint of the interval containing w . Algorithm S:

1. Pick $X \in \{0, 1\}^n$ uniformly at random *once*.
2. For $i = 1$ to k :
 - (a) Obtain $W_i = f_i(X) \in \mathbb{R}^d$.
 - (b) Pick $\Delta_i \in [u]$ uniformly at random.
 - (c) Return $Y_i = (Y_{i1}, \dots, Y_{id})$, where $Y_{ij} = \text{Round}(W_{ij} + \Delta_i\varepsilon)$.

Proof of correctness: Just for the analysis, define a (many-query) steward S' by the following algorithm:

1. For $i = 1$ to k :
 - (a) Query f_i at *every* point in its domain, thereby learning the entire function.
 - (b) Compute a point $\hat{\mu}_i \in \mathbb{R}^d$ where f_i is (ε, δ) -concentrated.
 - (c) Pick $\Delta_i \in [u]$ uniformly at random.
 - (d) Return $Y_i = (Y_{i1}, \dots, Y_{id})$, where $Y_{ij} = \text{Round}(\hat{\mu}_{ij} + \Delta_i\varepsilon)$.

Now fix any deterministic owner O . For a vector $\vec{\Delta} = (\Delta_1, \dots, \Delta_k) \in [u]^k$, let $f_1^{[\vec{\Delta}]}, \dots, f_k^{[\vec{\Delta}]}$ be the functions that O chooses in $O \leftrightarrow S'(\vec{\Delta})$, and let $\hat{\mu}_i^{[\vec{\Delta}]}$ be the point at which $f_i^{[\vec{\Delta}]}$ is concentrated that S' chooses in $O \leftrightarrow S'(\vec{\Delta})$. Observe that

$$\Pr_{\substack{\vec{\Delta} \in [u]^k \\ X \in \{0,1\}^n}} [\text{for some } i, \|f_i^{[\vec{\Delta}]}(X) - \hat{\mu}_i^{[\vec{\Delta}]}\|_\infty > \varepsilon] \leq k\delta. \quad (2)$$

(Imagine picking $\vec{\Delta}$ first, and then apply the union bound over the k different values of i .) Next, observe that

$$\Pr_{\substack{\vec{\Delta} \in [u]^k \\ X \in \{0,1\}^n}} [\text{for some } i, j, [\hat{\mu}_{ij}^{[\vec{\Delta}]} + (\Delta_i - 1)\varepsilon, \hat{\mu}_{ij}^{[\vec{\Delta}]} + (\Delta_i + 1)\varepsilon] \text{ is not entirely contained in one interval}] \leq \gamma. \quad (3)$$

(Indeed, for each i, j , the probability is just $2/u$, so by the union bound, the probability is at most $2kd/u \leq \gamma$.) Now, by the union bound, assume from now on that $\vec{\Delta}, X$ are such that neither the event of Equation 2 nor the event of Equation 3 takes place. Assume without loss of generality that $\delta < 1/2$. We will show that in $O \leftrightarrow S(X, \vec{\Delta})$, for every i , $\|Y_i - \mu_i\|_\infty \leq 1.5\ell + 3\varepsilon$.

We first show by induction on i that in $O \leftrightarrow S(X, \vec{\Delta})$, every f_i is precisely $f_i^{[\vec{\Delta}]}$. In the base case $i = 1$, this is trivial. For the inductive step, since the bad event of Equation 2 did not occur, we know that $f_i(X)$ is $\hat{\mu}_i^{[\vec{\Delta}]} \pm \varepsilon$. Therefore, since the bad event of Equation 3 did not occur, for every j , $\text{Round}(f_{ij}(X) + \Delta_i\varepsilon) = \text{Round}(\hat{\mu}_{ij}^{[\vec{\Delta}]} + \Delta_i\varepsilon)$. Therefore, the value Y_i in $O \leftrightarrow S(X, \vec{\Delta})$ is the same as the value Y_i in $O \leftrightarrow S'(\vec{\Delta})$, and hence O chooses the same f_{i+1} in both cases. This completes the induction.

Again using the fact that the bad event of Equation 2 did not occur, this immediately implies that in $\mathcal{O} \leftrightarrow \mathcal{S}(X, \vec{\Delta})$, every $f_i(X)$ is within ℓ_∞ distance ε of a point where f_i is (ε, δ) -concentrated. Since $\delta < 1/2$, this implies that every $f_i(X)$ is within ℓ_∞ distance 3ε of μ_i . Shifting by $\Delta_i\varepsilon$ and rounding introduce at most 1.5ℓ additional error, showing that $\|Y_i - \mu_i\| \leq 1.5\ell + 3\varepsilon$ as claimed. To complete the proof of correctness, note that $1.5\ell + 3\varepsilon \leq O(kd\varepsilon/\gamma)$.

The randomness complexity of this steward is n bits (for X) plus the randomness needed for $\vec{\Delta}$, for a total randomness complexity of

$$n + k \log u \leq n + O(k \log k + k \log d + k \log(1/\gamma)).$$

The steward clearly runs in $\text{poly}(n, k, d, \log(1/\varepsilon), \log(1/\gamma))$ time. \square

C The Impagliazzo-Zuckerman steward

C.1 Description and analysis of the steward

As discussed in Section 1.5.3, this steward is formed by combining the Impagliazzo-Zuckerman generator [IZ89, Imp92] with straightforward rounding. None of our results use this steward.

Proposition 2. *For any $n, k, d \in \mathbb{N}$ and $\varepsilon, \delta > 0$, there exists a one-query $(O(\varepsilon), k\delta + k \cdot 2^{-n^{\Omega(1)}})$ -steward for k adaptively chosen (ε, δ) -concentrated functions $f_1, \dots, f_k : \{0, 1\}^n \rightarrow \mathbb{R}^d$ with randomness complexity $O(n^6 + kd)$. The total running time of the steward is $\text{poly}(n, k, d, \log(1/\varepsilon))$.*

Proof. Let A be the “bit provider” of [Imp92, Theorem 2.6.8]. Partition \mathbb{R} into intervals of length 2ε . The steward:

1. For $i = 1$ to k :
 - (a) Give A the entropy bound $\delta n + d + 1$. It outputs some $X_i \in \{0, 1\}^n$.
 - (b) By querying, obtain the vector $W_i \stackrel{\text{def}}{=} f_i(X_i) \in \mathbb{R}^d$.
 - (c) Output $Y_i \in \mathbb{R}^d$, where Y_{ij} is the midpoint of the interval containing W_{ij} .

The randomness complexity of this steward is just the randomness complexity of A . From the proof of [Imp92, Theorem 2.6.8], this randomness complexity can be seen to be $O(n^6 + k/n + k(\delta n + d + 1))$. We may assume without loss of generality that $k\delta < 1$, because otherwise the proposition is trivial. Thus, the randomness complexity is $O(n^6 + kd)$ as claimed.

Now, for correctness, suppose $f : \{0, 1\}^n \rightarrow \mathbb{R}^d$ is (ε, δ) -concentrated at $\mu \in \mathbb{R}^d$. Let $g(X)_j$ be the midpoint of the interval containing $f(X)_j$. Let $H(\cdot)$ denote Shannon entropy. To bound $H(g(U_n))$, let $S = \{x : \|f(x) - \mu\|_\infty \leq \varepsilon\}$. Let $X \sim U_n$, and let E indicate whether $X \in S$. Then by the chain rule,

$$\begin{aligned} H(g(X)) &= H(g(X) \mid E) + H(E) \\ &\leq H(g(X) \mid X \in S) \cdot \Pr[X \in S] + H(g(X) \mid X \notin S) \cdot \Pr[X \notin S] + 1. \end{aligned}$$

Observe that $|g(S)| \leq 2^d$, because for each j , $[\mu_{ij} - \varepsilon, \mu_{ij} + \varepsilon]$ only intersects at most 2 intervals. Therefore, $H(g(X) \mid X \in S) \leq d$. Furthermore, $H(X \mid X \notin S) \leq n$ and applying a (deterministic) function can only reduce entropy, so $H(g(X) \mid X \notin S) \leq n$. Therefore,

$$\begin{aligned} H(g(X)) &\leq d \cdot \Pr[X \in S] + n \cdot \delta + 1 \\ &\leq \delta n + d + 1. \end{aligned}$$

By the guarantee of A , it follows that for any owner O , the sequence of responses (Y_1, Y_2, \dots, Y_k) is $(k \cdot 2^{-n^{\Omega(1)}})$ -close (in total variation distance) from how it would be distributed if A chose each X_i independently and uniformly at random. In this “fresh randomness” case, by the union bound, with probability $1 - k\delta$, for every i , $\|W_i - \mu_i\|_\infty \leq \varepsilon$. Rounding introduces at most ε additional error, so by the triangle inequality, in this case, for every i , $\|Y_i - \mu_i\|_\infty \leq 2\varepsilon$. \square

C.2 Comparison to our stewards

As noted in Section 1.5.3, our main steward (Theorem 1) has much better randomness complexity than the Impagliazzo-Zuckerman steward. Furthermore, the Impagliazzo-Zuckerman steward has failure probability $\delta' = k\delta + k \cdot 2^{-n^{\Omega(1)}}$, which becomes trivial when k is $O(2^{n^c})$ for a small constant $c > 0$. In contrast, our main steward has failure probability $\delta' = k\delta + \gamma$ for an arbitrary $\gamma > 0$; our steward is nontrivial for γ as small as $2^{-\Omega(nk/\log k)}$ and k as large as $\Omega(1/\delta)$, which could be as large as $2^{\Omega(n)}$.

We also noted in Section 1.5.3 that our main actually has worse accuracy than the Impagliazzo-Zuckerman steward. But one of our variant stewards, namely that of Theorem 4, is strictly superior to the Impagliazzo-Zuckerman steward, as we now demonstrate:

Proof of Proposition 2 from Theorem 4. Let S be the steward of Theorem 4 with $d_0 = 1$ and $\gamma = k \cdot 2^{-n^{1/2}}$. Then S is a one-query $(O(\varepsilon), k\delta + k \cdot 2^{-n^{1/2}})$ -steward. We may assume without loss of generality that $k \leq 2^{n^{1/2}}$, because otherwise Proposition 2 is trivial (the failure probability bound is more than 1). Therefore, the randomness complexity of S is bounded by

$$n + O(kd + \log k \log(1/\gamma)) \leq O(n + kd). \quad \square$$

D Nonconstructive PRG for block decision trees

For completeness, we record the details of the standard nonconstructive argument that there exists a PRG for block decision trees with a small seed length.

Lemma 13. *Suppose \mathcal{C} is a class of Boolean functions $f : \{0, 1\}^n \rightarrow \{0, 1\}$ such that a function in \mathcal{C} can be specified using t bits, i.e. $|\mathcal{C}| \leq 2^t$. Then for any γ , there exists a γ -PRG $\text{Gen} : \{0, 1\}^s \rightarrow \{0, 1\}^n$ for \mathcal{C} with seed length*

$$s \leq \log t + 2 \log(1/\gamma) + O(1).$$

Proof. Consider picking Gen uniformly at random from the set of all functions $\{0, 1\}^s \rightarrow \{0, 1\}^n$. Fix $C \in \mathcal{C}$, and let $\mu(C) = \Pr_x[C(x) = 1]$. Then for each fixed seed $x \in \{0, 1\}^s$, the probability (over Gen) that $C(\text{Gen}(x)) = 1$ is precisely $\mu(C)$. Therefore, the expected fraction of x such that $C(\text{Gen}(x)) = 1$ is precisely $\mu(C)$, and by Hoeffding’s inequality,

$$\Pr_{\text{Gen}} \left[\left| \frac{\#\{x : C(\text{Gen}(x)) = 1\}}{2^s} - \mu(C) \right| > \gamma \right] \leq 2^{-\Omega(\gamma^2 2^s)}.$$

Therefore, by the union bound, the probability that the above bad event holds for *any* C is at most $2^{t-\Omega(\gamma^2 2^s)}$. If we choose s large enough, this probability will be less than 1, showing that there exists a Gen that works for all C . How large do we need to choose s ? There is some constant c such that it is sufficient to have $c\gamma^2 2^s > t$. Taking logarithms completes the proof. \square

Proposition 3. *For any $k, n \in \mathbb{N}$, any finite alphabet Σ , and any $\gamma > 0$, there exists a γ -PRG $\text{Gen} : \{0, 1\}^s \rightarrow \{0, 1\}^{nk}$ for (k, n, Σ) -block decision trees with seed length*

$$s \leq n + k \log |\Sigma| + 2 \log(1/\gamma) + O(1).$$

Proof. Let \mathcal{C} be the class of all Boolean functions $f : \{0, 1\}^{nk} \rightarrow \{0, 1\}$ of the form $f(x) = g(T(x))$, where T is a (k, n, Σ) block decision tree. To specify a function $f \in \mathcal{C}$, we need to specify (1) a bit for each leaf of T and (2) a function $v : \{0, 1\}^n \rightarrow \Sigma$ for each internal node of T . In total, this number of bits t is

$$\begin{aligned}
 t &= |\Sigma|^k + 2^n \lceil \log |\Sigma| \rceil \cdot \sum_{i=0}^{k-1} |\Sigma|^i \\
 &\leq |\Sigma|^k + 2^{n+1} \log |\Sigma| \cdot \frac{|\Sigma|^k - 1}{|\Sigma| - 1} \\
 &\leq |\Sigma|^k + 2^{n+1} |\Sigma|^k \\
 &\leq 2^{n+2} |\Sigma|^k.
 \end{aligned}$$

By Lemma 13, this implies that there is a γ -PRG $\text{Gen} : \{0, 1\}^s \rightarrow \{0, 1\}^{nk}$ for \mathcal{C} with seed length $n + k \log |\Sigma| + 2 \log(1/\gamma) + O(1)$. The “operational” characterization of total variation distance implies that Gen is also a γ -PRG for (k, n, Σ) block decision trees as defined in Section 3. \square