The Computational Complexity Column

BY

Michal Koucký

Computer Science Institute, Charles University
Malostranské nám. 25, 118 00 Praha 1, Czech Republic
koucky@iuuk.mff.cuni.cz
https://iuuk.mff.cuni.cz/~koucky/

Recent Progress on Derandomizing Space-Bounded Computation

William M. Hoza*

Abstract

Is randomness ever necessary for space-efficient computation? It is commonly conjectured that $\mathsf{L} = \mathsf{BPL}$, meaning that halting decision algorithms can always be derandomized without increasing their space complexity by more than a constant factor. In the past few years (say, from 2017 to 2022), there has been some exciting progress toward proving this conjecture. Thanks to recent work, we have new pseudorandom generators (PRGs), new black-box derandomization algorithms (generalizations of PRGs), and new non-black-box derandomization algorithms. This article is a survey of these recent developments. We organize the underlying techniques into four overlapping themes:

1. The *iterated pseudorandom restrictions* framework for designing PRGs, especially PRGs for functions computable by arbitrary-order read-once branching programs.
2. The *inverse Laplacian* perspective on derandomizing BPL and the related concept of local consistency.
3. *Error reduction* procedures, including methods of designing low-error weighted pseudorandom generators (WPRGs).
4. The continued use of *spectral expander graphs* in this domain via the derandomized square operation and the Impagliazzo-Nisan-Wigderson PRG (STOC 1994).

We give an overview of these ideas and their applications, and we discuss the challenges ahead.

*Simons Institute for the Theory of Computing at the University of California, Berkeley. Email: williamhoza@berkeley.edu

1 Introduction

In an effort to solve problems as efficiently as possible, algorithm designers often introduce randomness into their algorithms. This paradigm is undoubtedly ingenious and beautiful. However, random bits can themselves be considered a computational "resource" that might be costly or unavailable. At best, randomization trades one type of inefficiency for another. We therefore want to distinguish between cases in which randomization gives an intrinsic advantage and cases in which algorithms can be derandomized with little to no penalty. In this article, we focus on the question of how randomization affects *space complexity*.

1.1 Randomized Space-Bounded Computation

Informally, $\mathsf{BPSPACE}(S)$ is everything that can be decided using randomness and $O(S)$ bits of space. More precisely, for a function $S : \mathbb{N} \to \mathbb{N}$, a language L is in $\mathsf{BPSPACE}(S)$ if there exists a Turing machine A with the following features.

1. The machine A has three tapes: a read-only input tape, a read-write work tape, and a read-once "random tape" that is initially filled with uniform random bits.

2. For every $N \in \mathbb{N}$,[1] every input $\sigma \in \{0,1\}^N$, and every assignment to the random tape $x \in \{0,1\}^\infty$, the machine touches at most $O(S(N))$ cells of the work tape and eventually halts, outputting a Boolean value $A(\sigma, x) \in \{0,1\}$.

3. For every input $\sigma \in \{0,1\}^*$, we have

$$\sigma \in L \implies \Pr_x[A(\sigma, x) = 1] \geq 2/3$$
$$\sigma \notin L \implies \Pr_x[A(\sigma, x) = 1] \leq 1/3.$$

Let us assume that $S \geq \log N$, so the machine has enough space to store a pointer to an arbitrary location in its input. Note that we assume that the algorithm halts for *every* assignment to the random tape (not merely with high probability). Using this assumption, one can show that the algorithm halts within $2^{O(S)}$ steps.[2] We use

[1] In this article, we use uppercase N to denote the length of the input to a space-bounded algorithm. We use lowercase n to denote the number of random bits that the algorithm uses.

[2] Historically, there was more early interest in the alternative "non-halting" model in which we merely require the algorithm to halt with high probability [33, 72, 73, 45, 13, 52, 69]. Indeed, in the older literature, notation along the lines of "$\mathsf{BPSPACE}(S)$" typically refers to the non-halting model, whereas the halting model is discussed using augmented notation such as "$\mathsf{BP_HSPACE}(S)$." Today, the halting model is standard.

BPL to denote BPSPACE($\log N$). The classes RSPACE(S) and RL are defined the same way, except that we only allow one-sided error.

These models were first studied by Aleliunas, Karp, Lipton, Lovász, and Rackoff [4] more than four decades ago. They presented a randomized algorithm showing that the undirected connectivity problem is in RL, and they asked whether L = RL. Today, the specific problem of undirected connectivity is indeed known to be in L, thanks to Reingold's famous algorithm [65] (the climax of a long sequence of papers studying the space complexity of undirected connectivity [4, 14, 10, 57, 59, 9, 77, 65, 68]). It is commonly believed that more generally L = RL = BPL. By a padding argument, if L = BPL, then DSPACE(S) = BPSPACE(S) for every space-constructible $S \geq \log N$.

Superficially, this sounds like the same frustrating story that pervades complexity theory. "We have been studying these important complexity classes for many decades, and at this point we think we know the relationship between them, but we don't know how to prove it." The same can be said regarding P vs. NP, or P vs. BPP, or L vs. P, or countless other fundamental problems.

However, there is a widespread feeling that *the* L *vs.* BPL *problem is different.* Compared to (say) the problem of proving P = BPP, there is a great deal of *optimism* about the possibility of unconditionally proving L = BPL. This optimism is sensible because the BPL model has a crucial weakness: the read-once random tape.

1.2 The Read-Once Assumption

In the definition of BPL, the machine A is only permitted to read each cell of the random tape a single time; the tape head can move right but not left. The motivation for this assumption is that we are modeling a problem-solving agent who has access to a single fair coin. The agent can see the outcome of only the most recent coin flip. If they want to know the outcome of a previous coin flip, they ought to have written it down at the time that it occurred (and paid for it in terms of space complexity).

As a consequence of the read-once assumption, the action of A on its random bits can be modeled by a polynomial-width standard-order *read-once branching program* (ROBP), defined below.

Definition 1 (Standard-order ROBPs)**.** A *width-w length-n standard-order ROBP* f is defined by a start state $v_0 \in [w]$, a sequence of n transition functions $f_1, \ldots, f_n \colon [w] \times \{0,1\} \to [w]$, and a set of accepting states $V_{\text{acc}} \subseteq [w]$. An input $x \in \{0,1\}^n$ determines a sequence of states $v_0, v_1, \ldots, v_n \in [w]$ by the rule

$v_i = f_i(v_{i-1}, x_i)$ for $i > 0$. The output of the program is given by

$$f(x) = \begin{cases} 1 & \text{if } v_n \in V_{\text{acc}} \\ 0 & \text{if } v_n \notin V_{\text{acc}}. \end{cases} \tag{1}$$

Equivalently, we can think of f as a directed graph with vertices arranged in $n + 1$ layers, $V_0, \ldots, V_n$, where $|V_i| = w$. For $i < n$, each vertex $u \in V_i$ has two outgoing edges leading to V_{i+1}, one labeled 0 and the other labeled 1. There is a designated start vertex $v_0 \in V_0$, and there is a set of designated accepting vertices $V_{\text{acc}} \subseteq V_n$. An input $x \in \{0,1\}^n$ is interpreted as a sequence of edge labels, identifying a path $(v_0, v_1, \ldots, v_n) \in V_0 \times V_1 \times \cdots \times V_n$. The output of the program is once again given by Equation (1).

The term "standard-order ROBP" is not standard. In typical papers on derandomizing space-bounded computation, standard-order ROBPs are simply called "ROBPs."[3] In this article, we include the modifier "standard-order" to emphasize that the program reads the input bits from left to right: first x_1, then x_2, then x_3, etc.

If A is a randomized, halting log-space algorithm and σ is an input of length N, then the function $f(x) \stackrel{\text{def}}{=} A(\sigma, x)$ can be computed by a width-n length-n standard-order ROBP for a suitable value $n = \text{poly}(N)$; each state of the program encodes a configuration of the machine A. An appealing approach to derandomizing A is to design a *pseudorandom generator* (PRG) that fools standard-order ROBPs.

Definition 2 (PRGs)**.** Let $\mathcal{F}$ be a class of functions $f\colon \{0,1\}^n \to \{0,1\}$, let X be a distribution over $\{0,1\}^n$, and let $\varepsilon > 0$. We say that X *fools* $\mathcal{F}$ *with error* ε if for every $f \in \mathcal{F}$,

$$|\Pr[f(X) = 1] - \Pr[f(U_n) = 1]| \leq \varepsilon,$$

where U_n denotes the uniform distribution over $\{0,1\}^n$. An *ε-PRG* for $\mathcal{F}$ is a function $G\colon \{0,1\}^s \to \{0,1\}^n$ such that $G(U_s)$ fools $\mathcal{F}$ with error ε. The value s is called the *seed length* of G.

If we could construct a PRG G that 0.1-fools width-n length-n standard-order ROBPs with seed length $O(\log n)$ and space complexity $O(\log n)$, then we could conclude that $\mathsf{L} = \mathsf{BPL}$, because we could deterministically estimate the acceptance probability of an algorithm A on an input σ to within ± 0.1 by computing $A(\sigma, G(x))$ for every seed x.

[3] Standard-order ROBPs are also sometimes referred to as "ordered branching programs," "layered branching programs," or "sequential-access ROBPs."

1.2.1 PRGs and Lower Bounds

Some readers might have an intuition that says that designing unconditional PRGs is hopelessly difficult. This intuition is indeed sensible in many contexts. For example, consider the problem of designing a PRG that fools *general* size-n branching programs, i.e., programs that may read their input bits any number of times and in any order. Such a PRG $G\colon \{0,1\}^s \to \{0,1\}^n$ would induce a corresponding "hard function" $h\colon \{0,1\}^{s+1} \to \{0,1\}$ that cannot be computed by size-n branching programs.[4] If G is computable in space $O(s)$, then so is h. Therefore, the problem of designing explicit PRGs for general branching programs is even harder than the problem of proving *branching program lower bounds* for explicit functions. Perhaps someday our grandchildren will manage to prove optimal lower bounds for branching programs, but until that day, we should probably consider optimal PRGs for general branching programs to be out of reach.[5]

The good news is that the *read-once* assumption is an absolute game-changer. In the read-once setting, optimal lower bounds are already known. For example, using standard communication complexity arguments, one can show that every standard-order ROBP computing the function

$$h(x_1, \ldots, x_{2n}) = x_1 \cdot x_{n+1} \oplus x_2 \cdot x_{n+2} \oplus \cdots \oplus x_n \cdot x_{2n}$$

has width $2^{\Omega(n)}$. To design optimal PRGs for standard-order ROBPs, we "merely" need to bridge the gap between lower bounds and PRGs.[6] There is no clear "barrier" preventing us from designing optimal PRGs for standard-order ROBPs. This is one of the reasons that a proof that $\mathsf{L} = \mathsf{BPL}$ seems vastly more attainable than, say, a proof that $\mathsf{P} = \mathsf{BPP}$.

1.2.2 Nisan's PRG and Beyond

So far, we do have several explicit PRGs that unconditionally fool standard-order ROBPs, but they do not achieve the optimal seed length. Most famously, Nisan designed an explicit PRG that ε-fools width-w length-n standard-order ROBPs with seed length $O(\log(wn/\varepsilon) \cdot \log n)$ [58]. The optimal seed length would be $\Theta(\log(wn/\varepsilon))$.

Admittedly, at this point it has been over three decades since Nisan's work [58], and we still do not have explicit PRGs for polynomial-width standard-order ROBPs with seed length better than Nisan's $O(\log^2 n)$ bound. However, an extensive body

[4] Specifically, h is the indicator function of the set $\{G(x)_{1\ldots s+1} : x \in \{0,1\}^s\}$.

[5] Currently, the best lower bound known is Nečiporuk's near-quadratic lower bound [56]. Explicit PRGs for size-n branching programs are known with a near-matching seed length of $\tilde{O}(\sqrt{n})$ [42, 38].

[6] Note that the Nisan-Wigderson reduction [60] does not work here, because it does not preserve the read-once property.

of research on the L vs. BPL problem has shown how to "go beyond" Nisan's work [58] in one sense or another. This rich and sophisticated literature is full of valuable insights that profoundly clarify the role of randomness in computing, even though the central questions remain open.

In the remainder of this article, we survey exciting progress that has been made on the L vs. BPL problem in just the past few years. (See Saks' survey [69] for an overview of older work.) We structure our discussion around four recurring technical themes: the *iterated pseudorandom restrictions* framework (Section 2), the *inverse Laplacian* perspective (Section 3), *error reduction* procedures (Section 4), and *expander graphs* (Section 5).

2 Iterated Pseudorandom Restrictions

2.1 Arbitrary-Order ROBPs

Nisan's classic PRG [58] suffers from a strange weakness. It turns out that permuting the output bits does not, in general, preserve the pseudorandomness property [78]. In other words, Nisan's PRG does not fool "arbitrary-order ROBPs." An *arbitrary-order ROBP* is defined just like a standard-order ROBP (Definition 1), except that instead of reading the input bits in the standard order $x_1, x_2, \ldots, x_n$, it reads the input bits in the order $x_{\pi(1)}, x_{\pi(2)}, \ldots, x_{\pi(n)}$ for some permutation $\pi\colon [n] \to [n]$.[7]

An interesting line of work has shown how to construct alternative PRGs for ROBPs that work even in the arbitrary-order setting [12, 76, 20, 50, 31]. We highlight a breakthrough paper by Forbes and Kelley [31]. Building on several earlier papers [66, 37, 20], Forbes and Kelley constructed two explicit PRGs for arbitrary-order ROBPs.

Theorem 1 (PRGs for arbitrary-order ROBPs [31])**.** *For every $w, n \in \mathbb{N}$ and $\varepsilon > 0$, there exist explicit ε-PRGs for width-w length-n arbitrary-order ROBPs with seed lengths*

$$O(\log(wn/\varepsilon) \cdot \log^2 n) \tag{2}$$

and

$$\tilde{O}(w \cdot \log(n/\varepsilon) \cdot \log n). \tag{3}$$

These seed lengths are only a little worse than Nisan's seed length [58], yet the PRGs fool a more powerful model.

For our main application (derandomizing BPL), it is no loss of generality to assume that the random bits are read in the standard order $x_1, x_2, x_3, \ldots$, so why

[7] To be clear, these programs are still "oblivious," meaning that vertices in the same layer read the same input bit. Arbitrary-order ROBPs are also called "unordered ROBPs" or "unknown-order ROBPs."

study arbitrary-order ROBPs? One reason is that they capture other interesting models of computation such as read-once formulas [12, 32, 22, 28, 29, 30]. Another reason is that studying arbitrary-order ROBPs forces us to develop *new techniques* for fooling ROBPs. Indeed, the ideas underlying Forbes and Kelley's PRGs [31] are completely different than those underlying Nisan's PRG [58]. Forbes and Kelley's PRGs [31] are based on the framework of *iterated pseudorandom restrictions* – our first "theme."

2.2 Forbes-Kelley Restrictions

Ajtai and Wigderson introduced the iterated restrictions framework in the context of pseudorandomness for AC^0 circuits [3]. Much later, Gopalan, Meka, Reingold, Trevisan, and Vadhan brought the framework to the world of L vs. BPL [36]. The idea is as follows. Our goal is to sample a string $X \in \{0,1\}^n$ that fools some function of interest $f\colon \{0,1\}^n \to \{0,1\}$. Our first step is to design a pseudorandom *restriction* $X \in \{0,1,\star\}^n$, i.e., we pseudorandomly assign values to a *pseudorandom subset* of the variables. We ensure that X "preserves the expectation" of f, meaning that $X \circ U$ fools f, where $X \circ U$ denotes the string obtained by sampling X and then replacing each $\star$ with a fresh truly random bit. Intuitively, designing such an X is easier than designing a full PRG, because in the analysis, some helpful truly random bits (U) are sprinkled in among the pseudorandom bits.

After assigning values according to X, our remaining task is to fool the restricted function $f|_X$. Therefore, we repeat the process, i.e., we sample a restriction X' that preserves the expectation of $f|_X$. Iterating in this way, we assign values to more and more variables. Eventually, we have assigned values to all the variables and hence we have a full PRG.

Forbes and Kelley's primary contribution is to show how to accomplish the first step, i.e., how to sample a pseudorandom restriction that assigns values to many variables while preserving the expectation of every bounded-width arbitrary-order ROBP [31]. Indeed, they prove the following.

Theorem 2 (Restrictions for arbitrary-order ROBPs [31])**.** *Let $w, n \in \mathbb{N}$ and $\varepsilon > 0$, and let $k = 4\log(wn/\varepsilon)$. Let D and T be k-wise independent n-bit strings (with uniform marginals), let U be uniform random over $\{0,1\}^n$, and assume that D, T, and U are mutually independent. Then $D + (T \wedge U)$ fools width-w length-n arbitrary-order ROBPs with error ε, where $+$ denotes bitwise* XOR *and $\wedge$ denotes bitwise* AND.

The strings D and T define a restriction X by letting T indicate the $\star$ positions and using D to assign values to the non-$\star$ positions. The statement that X preserves the expectation of f is equivalent to the statement that $D+(T \wedge U)$ fools f. The way

of thinking exemplified by the latter statement can be called the "pseudorandomness plus noise" perspective [37, 49].

Using standard constructions of k-wise independent random variables [44], one can explicitly sample D and T using $O(k \log n) = O(\log(wn/\varepsilon) \cdot \log n)$ truly random bits. In expectation, each restriction assigns values to half of the living variables, so after roughly $\log n$ iterations, we should intuitively expect that all the variables have been assigned values. Indeed, a more careful argument shows how to achieve an overall seed length of $O(\log(wn/\varepsilon) \cdot \log^2 n)$ (see Forbes and Kelley's work for details [31, Section 7]).

The proof of Theorem 2 is a beautiful application of Boolean Fourier analysis. Forbes and Kelley's techniques [31] work particularly well in the constant-width setting. By leveraging "Fourier growth bounds" for ROBPs [66, 76, 20, 48], Forbes and Kelley obtain restrictions for constant-width arbitrary-order ROBPs with better parameters. In the constant-width case, rather than k-wise independent distributions, Forbes and Kelley use "δ-biased distributions," i.e., distributions that fool parity functions with error $\delta/2$ [55].

Theorem 3 (Restrictions for constant-width arbitrary-order ROBPs [31])**.** *Let $w \in \mathbb{N}$ be a constant. For every $n \in \mathbb{N}$ and $\varepsilon > 0$, there exists a value $\delta = \exp\left(-\tilde{O}(\log(n/\varepsilon))\right)$ such that the following holds. Let D and T be δ-biased random variables distributed over $\{0,1\}^n$, let U be uniform random over $\{0,1\}^n$, and assume that D, T, and U are mutually independent. Then $D + (T \wedge U)$ fools width-w length-n arbitrary-order ROBPs with error ε, where $+$ denotes bitwise* XOR *and $\wedge$ denotes bitwise* AND.

Using standard constructions of δ-biased distributions [55, 5], the random variables D and T of Theorem 3 can be sampled explicitly using $O(\log(n/\delta)) = \tilde{O}(\log(n/\varepsilon))$ truly random bits. This leads to a PRG for constant-width arbitrary-order ROBPs with seed length $\tilde{O}(\log(n/\varepsilon) \cdot \log n)$.

2.3 The Early Termination Technique

Forbes and Kelley's PRGs [31] are examples of restrictions-based PRGs with seed length polylog(n), similar to the seed length of Nisan's PRG [58]. In some cases, we can use the iterated restrictions framework to get seed lengths as low as $\tilde{O}(\log n)$ or even $O(\log n)$. The key idea is to show that after applying a few pseudorandom restrictions (say, poly($\log \log n$) many), the function f that we are trying to fool "simplifies" in some sense with high probability. When this occurs, we can *terminate the restriction process early*, and use some other approach to fool the restricted function, taking advantage of its simplicity.

This "early termination" technique was introduced by Gopalan, Meka, Reingold, Trevisan, and Vadhan [36], and it has turned out to be useful for quite a few PRG problems [36, 50, 28, 47, 49, 29, 30]. Let us briefly discuss three examples.

- Gopalan, Meka, Reingold, Trevisan, and Vadhan designed an explicit PRG for *read-once* CNF *formulas* with near-optimal seed length $\tilde{O}(\log(n/\varepsilon))$ [36].
- Doron, Hatami, and Hoza designed an explicit PRG for *read-once* AC^0 *formulas* with near-optimal seed length $\tilde{O}(\log(n/\varepsilon))$ [28].
- Doron, Meka, Reingold, Tal, and Vadhan designed an explicit PRG for constant-width arbitrary-order *monotone ROBPs* (defined next) with near-optimal seed length $\tilde{O}(\log(n/\varepsilon))$ [30].

Definition 3 (Monotone ROBPs)**.** Let f be a width-w length-n arbitrary-order ROBP with transition functions $f_1, \ldots, f_n \colon [w] \times \{0,1\} \to [w]$. We say that f is *monotone* if, for each $i \in [n]$ and each bit $b \in \{0,1\}$, the transition function $f_i(\cdot, b)$ is a monotone function $[w] \to [w]$ [51, 30].

It turns out that constant-width arbitrary-order monotone ROBPs can simulate read-once AC^0 formulas [22, 30]. In turn, obviously read-once AC^0 formulas generalize read-once CNF formulas. Thus, the classes fooled by the three PRGs mentioned above form a hierarchy:

$$\begin{aligned}&\text{read-once CNFs}\\ &\quad\subseteq \text{read-once } \mathsf{AC}^0\\ &\qquad\subseteq \text{constant-width arbitrary-order monotone ROBPs.}\end{aligned}$$

Over time, we are gradually figuring out how to fool *more and more powerful classes* with near-optimal seed length, building our way up toward the class of general (arbitrary-order) ROBPs. This type of progress (steadily improving the class of functions fooled) has turned out to be more feasible than insisting on fooling *all* (standard-order) ROBPs and trying to improve the seed length.

Recall that Forbes and Kelley's work (Theorem 3) shows how to assign values to half the input variables of a constant-width arbitrary-order ROBP at a cost of only $\tilde{O}(\log(n/\varepsilon))$ truly random bits. To get a full PRG in the monotone case, Doron, Meka, Reingold, Tal, and Vadhan show that after a few Forbes-Kelley restrictions, monotone ROBPs are likely to simplify [30]. Roughly speaking, the notion of simplification is that the *width* of the program steadily decreases until the function is trivial.

We remark that Doron, Meka, Reingold, Tal, and Vadhan's work [30] is one example where *techniques designed for the arbitrary-order case have turned out to*

be useful even for the standard-order case.[8] This demonstrates the counterintuitive wisdom of working on problems that are even *more* difficult than the problems that we care about most.

2.4 A Challenge: Parity Gates

The iterated restrictions paradigm is flexible and powerful, especially when it is combined with the early termination technique. All of the recent work using these techniques is certainly exciting and encouraging. Unfortunately, however, we still do not have a clear path toward fooling all constant-width ROBPs (let alone polynomial-width ROBPs) with near-logarithmic seed length. Indeed, it seems that this line of work is perhaps "running out of steam."

To understand the limitations of these techniques, observe that for any arbitrary-order ROBP $f\colon \{0,1\}^n \to \{0,1\}$, we can define a more complicated function $g\colon \{0,1\}^{n^2} \to \{0,1\}$ by block-composing with the parity function, i.e.,

$$g(x_{11},\ldots,x_{nn}) = f\left(\bigoplus_{i=1}^{n} x_{1i}, \bigoplus_{i=1}^{n} x_{2i}, \ldots, \bigoplus_{i=1}^{n} x_{ni}\right).$$

If the initial ROBP f has width $w = O(1)$, then g can be computed by an arbitrary-order ROBP of width $2w = O(1)$, but the early termination technique seems to break down when we try to apply it to g. It seems that (pseudo)random restrictions have very little effect on g, because a restriction of the parity function is always either the parity function or its complement. Fooling a typical restriction of g is thus at least as difficult as fooling f.

More concretely, consider the problem of fooling read-once AC^0 formulas *with parity gates* (Figure 1). Doron, Hatami, and Hoza gave an explicit PRG for this class with seed length $\tilde{O}(t + \log(n/\varepsilon))$ where t is the number of parity gates in the formula [28]. For the depth-2 case, we have explicit PRGs with near-optimal seed length [49, 50, 47], and in fact with "partially optimal" seed length $O(\log n) + \tilde{O}(\log(1/\varepsilon))$ [29]. However, when the depth is a large constant and the number of parity gates is unbounded, it seems quite difficult to achieve seed length $\tilde{O}(\log n)$.

3 The Inverse Laplacian Perspective

In light of challenges such as that discussed in Section 2.4, it is worthwhile to take a step back and ask whether we truly need to design better PRGs for ROBPs. After

[8] The monotone ROBP model was first introduced by Meka and Zuckerman [51], who were not concerned with issues of variable ordering. They presented PRGs for the standard-order case [51]; in the constant-width regime, their seed length matches Nisan's [58].

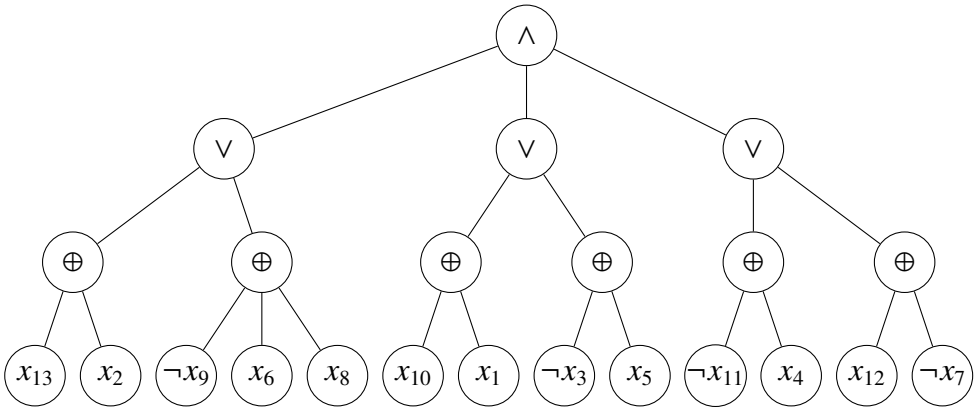

Figure 1: We would like to design explicit PRGs for constant-width (arbitary-order) ROBPs with near-optimal seed length $\tilde{O}(\log(n/\varepsilon))$. The class of read-once AC^0 formulas with parity gates is a challenging special case. Indeed, the case of read-once $\mathsf{AND} \circ \mathsf{OR} \circ \mathsf{PARITY}$ formulas already seems formidable.

all, our primary goal is derandomizing space-bounded computation. In this section, we discuss a non-PRG-based approach to proving $\mathsf{L} = \mathsf{BPL}$.

3.1 The Matrix of Expectations of Subprograms

To derandomize BPL, it suffices to design a deterministic log-space algorithm that is given a width-n length-n standard-order ROBP f and estimates $\mathbb{E}[f]$ to within a small additive error. There is no need to treat f as a black box; it is permissible to inspect the transitions of f and try to thereby gain some advantage. Since we are only concerned with space complexity, if we intend to estimate the expectation of the program, we might as well estimate the expectations of all subprograms, too.

Definition 4 (Subprograms)**.** Suppose f is a width-w length-n standard-order ROBP with layers $V_0, \ldots, V_n$. Let $u \in V_i$ and $v \in V_j$ be vertices with $i \leq j$. We define the *subprogram* $f_{u \to v}$ to be the width-w length-$(j - i)$ standard-order ROBP on layers $V_i, V_{i+1}, \ldots, V_j$ obtained from f by designating u as the start vertex and v as the unique accepting vertex.

Let us collect all the expectations of these subprograms $\mathbb{E}[f_{u \to v}]$ in an $m \times m$ matrix P, where m is the number of vertices in f, namely $m = w \cdot (n + 1)$. That is, for every pair of vertices u, v in f, if $u \in V_i$ and $v \in V_j$, then

$$P_{u,v} = \begin{cases} \mathbb{E}[f_{u \to v}] & \text{if } i \leq j \\ 0 & \text{if } i > j. \end{cases} \tag{4}$$

The following problem is essentially complete for BPL:

- **Input:** A width-n length-n standard-order ROBP f.
- **Output:** A matrix $\hat{P}$ that approximates the matrix of expectations of subprograms (P) to within additive entrywise error 0.1.

(By "essentially complete for BPL," we mean that a *decision version* of the problem is complete for the *promise version* of BPL with respect to deterministic log-space reductions. These technicalities do not seem to be important.)

3.2 The Inverse Laplacian of a Standard-Order ROBP

To try to approximate P, we can start by computing the *random walk matrix* W. By definition, for each pair of vertices u, v in f, the entry $W_{u,v}$ gives the probability of arriving at v when we start at u and take a single random step. Computing W is trivial: $W_{u,v}$ is half the number of edges from u to v.

The expectations of subprograms of f correspond to *powers* of W. Indeed, $(W^t)_{u,v}$ is the probability that a t-step random walk from u arrives at v. Therefore, if $u \in V_i$ and $v \in V_j$, then

$$(W^t)_{u,v} = \begin{cases} \mathbb{E}[f_{u \to v}] & \text{if } j - i = t \\ 0 & \text{otherwise.} \end{cases}$$

Consequently, there is a simple formula for the matrix of expectations of subprograms (P) in terms of the random walk matrix (W):

$$P = W^0 + W^1 + W^2 + \cdots + W^n. \tag{5}$$

Furthermore, $W^{n+1} = 0$, so we can simplify Equation (5) using the geometric series formula:

$$P = (I - W)^{-1}.$$

The matrix $I - W$ is called the (directed) *Laplacian matrix* of the program f and denoted L. Thus, we are looking for an *approximate inverse Laplacian* $\hat{P} \approx L^{-1}$. This way of thinking – the "inverse Laplacian perspective" – was introduced most clearly in work by Ahmadinejad, Kelner, Murtagh, Peebles, Sidford, and Vadhan [1], and it is our second technical "theme."

3.3 Local Consistency

A key benefit of the inverse Laplacian perspective is that it suggests a new way of thinking about error. Suppose that someone gives us a candidate matrix $\hat{P}$. Is $\hat{P}$ a good approximation to P? We cannot directly compare the entries of $\hat{P}$ to those

of P, because we do not know P (remember, approximating P is essentially BPL-complete). However, we *can* compute the error *after multiplying by the Laplacian matrix*. That is, we can compare $\hat{P}L$ to the identity matrix. Define E to be the error matrix $E = I - \hat{P}L$.

This error matrix E has a natural probabilistic interpretation. Expanding the definition, we have $E = I - \hat{P} \cdot (I - W) = I + \hat{P}W - \hat{P}$. Therefore, if $u \in V_i$ and $v \in V_j$ where $i < j$, then

$$E_{u,v} = (\hat{P}W)_{u,v} - \hat{P}_{u,v} = \underbrace{\left(\sum_{s \in V_{j-1}} \hat{P}_{u,s} \cdot W_{s,v} \right)}_{(*)} - \hat{P}_{u,v}.$$

The entry $E_{u,v}$ measures the difference between two different methods of using $\hat{P}$ to estimate $\mathbb{E}[f_{u \to v}]$. The first method is to simply consult the (u, v) entry of $\hat{P}$, since after all $\hat{P}$ is intended to be an approximation to P. The second method is to look at $\hat{P}$'s estimates for the probabilities of arriving at vertices in the layer V_{j-1} that *precedes* v, and then propagate those probabilities forward by a single step, leading to quantity $(*)$.

Thus, E measures the extent to which $\hat{P}$ is locally *consistent with itself*; we refer to E as the matrix of *local consistency errors*. The term "local consistency" was introduced by Cheng and Hoza [23]; the connection between local consistency and the Laplacian matrix was observed by subsequent papers [24, 63, 39]. We will discuss an application of the notion of local consistency next. For additional applications of the inverse Laplacian perspective, see Sections 4 and 5.

3.4 One-Sided vs. Two-Sided Derandomization

Cheng and Hoza used the concept of local consistency to prove a new conditional derandomization of BPL [23]. Recall that a *hitting set generator* (HSG) is a one-sided version of a PRG.

Definition 5 (HSGs)**.** Let $\mathcal{F}$ be a class of functions $f\colon \{0,1\}^n \to \{0,1\}$ and let $\varepsilon > 0$. An *ε-HSG* for $\mathcal{F}$ is a function $G\colon \{0,1\}^s \to \{0,1\}^n$ such that for every $f \in \mathcal{F}$,

$$\text{if } \Pr[f(U_n) = 1] \geq \varepsilon, \text{ then there exists } x \text{ such that } f(G(x)) = 1.$$

If G is an ε-PRG for $\mathcal{F}$, then G is also is an ε'-HSG for $\mathcal{F}$ for every $\varepsilon' > \varepsilon$. HSGs are potentially much easier to construct than PRGs, so it is worthwhile to ask, what would be the applications of optimal explicit HSGs? Working through the definitions, one can easily show that an optimal explicit HSG for standard-order ROBPs would imply L = RL (one-sided derandomization). Cheng and

Hoza showed that it would also imply the stronger statement $\mathsf{L} = \mathsf{BPL}$ (two-sided derandomization) [23].

Theorem 4 (HSGs would derandomize BPL [23])**.** *Assume that for every $n \in \mathbb{N}$, there is a $\frac{1}{2}$-HSG for width-n length-n standard-order ROBPs that has seed length $O(\log n)$ and that is computable in space $O(\log n)$. Then $\mathsf{L} = \mathsf{BPL}$.*

Let us briefly sketch the proof of Theorem 4. Suppose we are given a width-n length-n standard-order ROBP f. Let G be an HSG with output length n^c, where c is a large enough constant. For each seed x, we think of $G(x)$ as a long stream of random bits and use it to compute a matrix $\hat{P}^{(x)}$ that is a candidate approximation to the matrix P of expectations of subprograms of f. Using the hitting property of G, one can show that there is at least one "good seed" x such that $\hat{P}^{(x)} \approx P$. To identify such a seed algorithmically, we find an x such that $\hat{P}^{(x)}$ has good local consistency.

We remark that an analogous theorem for time-bounded derandomization has been known for decades [6, 7, 18, 34]. In fact, Buhrman and Fortnow showed generically that derandomizing the promise version of RP would imply $\mathsf{P} = \mathsf{BPP}$, regardless of whether the derandomization is via an HSG [18]. An interesting open problem is to prove the analogous theorem for the space-bounded setting, generalizing Theorem 4.

4 Error Reduction Procedures

In the previous section, we introduced the inverse Laplacian perspective, and we discussed one application (the derandomization of BPL using a hypothetical HSG). There are several other applications of the inverse Laplacian perspective. These other applications take advantage of the rich literature on *fast, randomized* algorithms for approximately solving Laplacian systems of equations, starting with Spielman and Teng's seminal work [74]. Most especially, these other applications work by importing *error reduction* techniques – our third "theme" – to the space-bounded derandomization setting.

4.1 Non-Black-Box Error Reduction

As our first example, let us discuss a theorem by Ahmadinejad, Kelner, Murtagh, Peebles, Sidford, and Vadhan [1] (strengthening prior work by Hoza and Zuckerman [41]). Their theorem shows how to generically decrease the error of space-bounded derandomization algorithms. In the following, think of ε as negligibly small compared to n, such as perhaps $\varepsilon = 2^{-\mathrm{polylog}(n)}$.

Theorem 5 (Error reduction for non-black-box derandomization [1])**.** *Let $S : \mathbb{N} \to \mathbb{N}$ be a function. Assume that given a width-n length-n standard-order ROBP f, it is possible to deterministically compute $\mathbb{E}[f]$ to within $\pm 1/n^3$ in space $S(n)$. Then given f and $\varepsilon > 0$, it is possible to deterministically compute $\mathbb{E}[f]$ to within $\pm\varepsilon$ in space*

$$O(S(n) + \log n \cdot \log \log_n(1/\varepsilon)).$$

Let us sketch the proof of Theorem 5, which uses the inverse Laplacian perspective. Let P be the matrix of expectations of subprograms of f. Using the given $S(n)$-space algorithm, we can construct a matrix $\hat{P}$ such that $\left\|P - \hat{P}\right\|_\infty \leq O(1/n)$.[9] Let W be the random walk matrix of f, let $L = I - W$ be the Laplacian matrix, and let $E = I - \hat{P}L$ be the error matrix after multiplying by L (aka the matrix of local consistency errors). Then, we define a new approximation matrix $\hat{P}'$ by the formula

$$\hat{P}' = \hat{P} + E\hat{P} + E^2\hat{P} + \cdots + E^m\hat{P}$$

for a suitably chosen parameter m. (Intuitively, we start with $\hat{P}$, and then we add a sequence of finer and finer "correction terms" $E\hat{P}, E^2\hat{P}, \ldots, E^m\hat{P}$.) Let us measure the quality of this new approximation. The key, again, is to measure quality *after* multiplying by the Laplacian matrix, which causes a telescoping sum:

$$\hat{P}'L = (I - E) + E \cdot (I - E) + E^2 \cdot (I - E) + \cdots + E^m \cdot (I - E) = I - E^m.$$

Amazingly, we have managed to replace E with E^m, which intuitively should mean that the errors are getting much smaller. This technique for decreasing the error of an approximate matrix inverse is called *preconditioned Richardson iteration*.

Ultimately, what we care about is entrywise closeness to P. We can bound the entrywise errors using the submultiplicative $\|\cdot\|_\infty$ matrix norm:

$$\begin{aligned}
\left\|\hat{P}' - P\right\|_\infty = \left\|\left(\hat{P}'L - I\right) \cdot P\right\|_\infty = \|E^m \cdot P\|_\infty &\leq \|E\|_\infty^m \cdot \|P\|_\infty \\
&= \left\|\left(P - \hat{P}\right) \cdot L\right\|_\infty^m \cdot \|P\|_\infty \\
&\leq \left(\left\|P - \hat{P}\right\|_\infty \cdot \|L\|_\infty\right)^m \cdot \|P\|_\infty \\
&\leq O(1/n)^m \cdot O(n),
\end{aligned}$$

which is at most ε if we choose a suitable value $m = O(\log(1/\varepsilon)/\log n)$. One can compute $\hat{P}'$ deterministically in space $O(S(n) + \log n \cdot \log m)$, completing the proof of Theorem 5.

[9] Indeed, $\|P - \hat{P}\|_\infty \stackrel{\text{def}}{=} \max_u \sum_v |P_{u,v} - \hat{P}_{u,v}| \leq n \cdot (n+1) \cdot \max_{u,v} |P_{u,v} - \hat{P}_{u,v}| \leq n \cdot (n+1) \cdot n^{-3}$.

4.2 Weighted Pseudorandom Generators (WPRGs)

The parameters of Theorem 5 are impressive; we pay very little penalty for error reduction, even when the target error ε is extremely small. The algorithm of Theorem 5 is non-black-box, because we must inspect the graph structure of the given ROBP to compute the matrices W, L, E, etc.

As discussed previously, non-black-box algorithms are sufficient for proving $\mathsf{L} = \mathsf{BPL}$. However, black-box algorithms are stronger, and they tend to be more useful as building blocks inside larger algorithms. What is the best way to compute $\mathbb{E}[f]$ to within a tiny additive error ε if we only have *query* access to a standard-order ROBP f? An ε-PRG clearly suffices for this task, but could there be an easier approach? This motivates the intriguing concept of a *weighted pseudorandom generator* (WPRG), introduced by Braverman, Cohen, and Garg [15].

Definition 6 (WPRGs)**.** Let $\mathcal{F}$ be a class of functions $f\colon \{0,1\}^n \to \{0,1\}$ and let $\varepsilon > 0$. An *ε-WPRG* for $\mathcal{F}$ is a pair (G,ρ), where $G\colon \{0,1\}^s \to \{0,1\}^n$ and $\rho\colon \{0,1\}^s \to \mathbb{R}$, such that for every $f \in \mathcal{F}$,

$$\left| \mathop{\mathbb{E}}_{x \sim U_n}[f(x)] - \mathop{\mathbb{E}}_{x \sim U_s}[f(G(x)) \cdot \rho(x)] \right| \leq \varepsilon. \tag{6}$$

A PRG is the special case $\rho \equiv 1$. Crucially, Definition 6 allows for $\rho(x) < 0$, which opens the door for the possibility of *error cancellation* in Equation (6).[10] One can think of these negative weights as effectively introducing a kind of "negative probability" into the picture; WPRGs are also called *pseudorandom pseudodistribution generators*.[11]

One can show that if (G,ρ) is an ε-WPRG for $\mathcal{F}$, then G is an ε'-HSG for $\mathcal{F}$ for every $\varepsilon' > \varepsilon$. Thus, we have a hierarchy,

$$\text{PRG} \implies \text{WPRG} \implies \text{HSG}.$$

When they introduced the concept of a WPRG, Braverman, Cohen, and Garg presented an explicit construction of an ε-WPRG for polynomial-width standard-order ROBPs [15] with seed length

$$\tilde{O}(\log^2 n + \log(1/\varepsilon)). \tag{7}$$

For comparison, recall that Nisan's PRG ε-fools polynomial-width standard-order ROBPs with seed length $O(\log^2 n + \log n \cdot \log(1/\varepsilon))$ [58]. Thus, Braverman, Cohen,

[10] WPRGs with nonnegative weight functions $\rho\colon \{0,1\}^s \to [0,\infty)$ are essentially equivalent to unweighted PRGs [63, Appendix C of ECCC version].

[11] Braverman, Cohen, and Garg coined the term "pseudorandom pseudodistribution" [15]. The alternative term "weighted pseudorandom generator" was introduced later, by Cohen, Doron, Renard, Sberlo, and Ta-Shma [24].

and Garg's seed length [15] is superior when ε is very small (again, the case $\varepsilon = 2^{-\operatorname{polylog}(n)}$ is good to have in mind). Prior to their work [15], it was not even known how to construct an ε-HSG with the seed length that they achieve.

Braverman, Cohen, and Garg's work [15] is quite complex. This spurred a search for simpler approaches [41, 21, 24, 63, 39]. In addition to achieving improved simplicity, this line of work was also able to remove the lower-order terms hiding under the $\tilde{O}$ in Equation (7).

Theorem 6 (Optimal-error WPRGs [39])**.** *For every $w, n \in \mathbb{N}$ and $\varepsilon > 0$, there is an explicit ε-WPRG for width-w length-n standard-order ROBPs with seed length $O(\log(wn) \cdot \log n + \log(1/\varepsilon))$.*

To prove Theorem 6, we start with Nisan's PRG with error $1/\operatorname{poly}(nw)$ and seed length $O(\log(wn) \cdot \log n)$. Then, we use the *preconditioned Richardson iteration* technique that we discussed in Section 4.1 to decrease the error of the PRG. Implementing this technique is not completely straightforward, because we are in the black-box setting, and hence we can no longer compute the matrices W, L, E, etc. However, two independent papers (one by Cohen, Doron, Renard, Sberlo, and Ta-Shma [24] and the other by Pyne and Vadhan [63]) contributed the insight that one can set up the WPRG *construction* in such a way that preconditioned Richardson iteration happens in the *analysis*. Finally, to achieve the seed length of Theorem 6, we combine these ideas with a suitable sampler trick [39].

In general, starting from an explicit PRG for width-w length-n standard-order ROBPs with error $1/(wn)^c$ and seed length s (for a suitable constant $c > 1$), we get an explicit WPRG for such programs with arbitrarily small error ε and seed length $O(s + \log(1/\varepsilon))$ [39]. There are other, related error reduction procedures that achieve slightly better parameters in some cases [41, 24, 63]. For example, consider standard-order ROBPs of width w and length $\log^c w$ for a constant $c \in \mathbb{N}$. Nisan and Zuckerman showed how to fool these short, wide programs with seed length $O(\log w)$ and a relatively large error such as $2^{-(\log w)^{0.99}}$ [61]. By applying an error-reduction procedure to the Nisan-Zuckerman PRG [61], Hoza and Zuckerman designed an explicit ε-HSG for these programs with asymptotically optimal seed length $O(\log(w/\varepsilon))$, even when ε is small [41]. It remains an interesting open problem to match this seed length with a WPRG.

4.3 Improving the Saks-Zhou Algorithm

Let us now discuss an application of low-error WPRGs. Recall our original derandomization goal: we want to deterministically decide languages in $\mathsf{BPSPACE}(S)$, for $S \geq \log N$, using as little space as possible.

Savitch's theorem [71] implies that $\mathsf{RSPACE}(S) \subseteq \mathsf{DSPACE}(S^2)$. The more general inclusion $\mathsf{BPSPACE}(S) \subseteq \mathsf{DSPACE}(S^2)$ follows from early work on the

non-halting version of $\mathsf{BPSPACE}(S)$ [13, 45]. Later, Saks and Zhou used Nisan's PRG [58] in a sophisticated way to prove $\mathsf{BPSPACE}(S) \subseteq \mathsf{DSPACE}(S^{3/2})$ [70]. Now, decades later, we can finally improve Saks and Zhou's bound.

Theorem 7 (Improved derandomization of $\mathsf{BPSPACE}$ [39])**.** *Let $S : \mathbb{N} \to \mathbb{N}$ be a function satisfying $S(N) \geq \log N$. Then*

$$\mathsf{BPSPACE}(S) \subseteq \mathsf{DSPACE}\left(\frac{S^{3/2}}{\sqrt{\log S}}\right). \tag{8}$$

Admittedly, the bound of Equation (8) is only barely better than Saks and Zhou's $O(S^{3/2})$ bound [70]. Still, Theorem 7 potentially has some "psychological" value, because it demonstrates that Saks and Zhou's result [70] is not the "end of the road." There is no particular reason to think that Theorem 7 is the end of the road either. No compelling barriers to further progress are known; humanity has no real excuse for having not yet proven $\mathsf{L} = \mathsf{BPL}$.

The starting point for proving Theorem 7 is work by Armoni from more than two decades ago [8]. Armoni designed an explicit ε-PRG for width-w length-n standard-order ROBPs based on a generalization of Nisan and Zuckerman's techniques [61]. Armoni's seed length is slightly better than Nisan's seed length [58] in the regime $n \ll w$ and $\varepsilon \gg 1/w$ [8]. By combining his PRG with recent error reduction techniques [24, 63], we get an explicit WPRG with a seed length that is slightly better than Nisan's seed length [58] in the regime $n \ll w$, *even for low error* such as $\varepsilon = 1/\operatorname{poly}(w)$.

The original Saks-Zhou algorithm [70] uses Nisan's PRG with parameters in this regime ($n \ll w$ and $\varepsilon = 1/\operatorname{poly}(w)$) as a subroutine. Armoni showed how to use a generic PRG in place of Nisan's PRG [8], and Chattopadhyay and Liao showed more generally how to use WPRGs [21], building on an earlier suggestion by Braverman, Cohen, and Garg [15]. Combining these results proves Theorem 7. (See Figure 2.) This argument appears in work by Hoza [39], but to be clear, the ingredients all come from prior work [70, 8, 21, 24, 63]. Hoza's contribution to the proof of Theorem 7 is merely to put the pieces together [39].

Cohen, Doron, and Sberlo recently designed an algorithm that improves on Saks and Zhou's work [70] in a different direction [25]. Consider the following natural computational problem.

- **Input:** A value $n \in \mathbb{N}$ and a stochastic matrix $M \in \mathbb{R}^{w \times w}$, where each entry has bit complexity $O(\log(wn))$.
- **Output:** A matrix that approximates M^n to additive entrywise error 0.1.

When we restrict to the case $n = w$, the problem above is essentially complete for BPL. One can think of the Saks-Zhou algorithm as a method of solving the problem

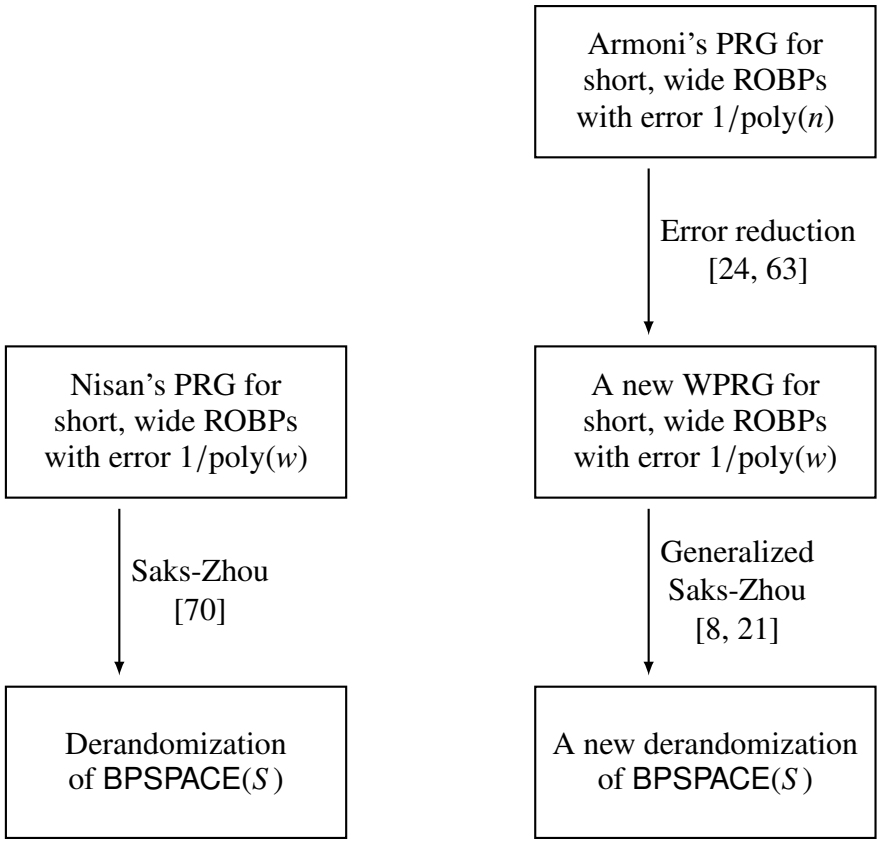

Figure 2: Saks and Zhou's derandomization of BPSPACE [70] (left) vs. the new and improved derandomization of BPSPACE (Theorem 7, right).

in space $O(\log(wn) \cdot \sqrt{\log n})$. Cohen, Doron, and Sberlo show how to solve the problem in space $\tilde{O}(\log w \cdot \sqrt{\log n} + \log n)$ [25], which is a significant improvement in the regime $n \gg w$. Their algorithm combines the Saks-Zhou algorithm with Richardson iteration, but in a different way than the proof of Theorem 7.

5 Spectral Expander Graphs

Let us consider one more natural problem that is essentially complete for BPL.

- **Input:** A directed graph G, two vertices s and t, and two positive integers k and m (represented in unary).
- **Output:** The probability that a k-step random walk starting at s ends at t, to within an additive error of $1/m$.

An appealing special case is when G is undirected. As mentioned previously, Reingold designed a deterministic log-space algorithm to determine whether there *exists* a path from s to t in an undirected graph G [65], which, intuitively, corresponds to the case $k = \infty$. A recent line of work has studied the case that k is finite, and in particular, k might be smaller than the mixing time of G [53, 54, 1]. For any k, Ahmadinejad, Kelner, Murtagh, Peebles, Sidford, and Vadhan gave an algorithm for computing k-step random walk probabilities in undirected graphs that runs in near-logarithmic space [1].

Theorem 8 (Estimating random-walk probabilities in undirected graphs [1])**.** *Given an undirected graph (or, more generally, an Eulerian digraph) G, two vertices s and t, and positive integers k and m represented in unary, it is possible to deterministically compute the probability that a length-k random walk starting at s arrives at t to within additive error $1/m$ in space $\tilde{O}(\log N)$, where N is the bit-length of the input.*

One of the (many) ideas in the proof of Theorem 8 is to use *expander graphs* to take a certain type of pseudorandom walk through G instead of a truly random walk. There is a long history of using expanders as tools for space-bounded derandomization, going back to work by Ajtai, Komlós, and Szemerédi [2]. Modern work on L vs. BPL continues to develop new ways of using and analyzing expanders – our fourth technical "theme."

5.1 The Derandomized Square

In more detail, the proof of Theorem 8 uses expanders via Rozenman and Vadhan's *derandomized square* operation [68]. For simplicity, consider a D-regular

undirected graph G. By definition, a single step in the *square graph* G^2 consists of two steps in the original graph G. Effectively, this means that squaring G places a clique on the D neighbors of each vertex v. The idea of the derandomized square is to instead place an expander graph on the neighbors of v, thereby producing a sparse approximation to G^2.

In Rozenman and Vadhan's original paper, they prove that the *spectral expansion* of the derandomized square is almost as good as that of G^2 [68]. They use this bound to derive an alternative proof that undirected connectivity is in L [68]. Recent work [53, 54, 1] shows that the derandomized square approximates G^2 in much stronger senses. The proof of Theorem 8 combines this analysis with several other techniques, including the inverse Laplacian perspective and error reduction methods.

5.2 The INW Generator

The derandomized square operation also has connections to the PRG approach to L vs. BPL, and in particular to a PRG by Impagliazzo, Nisan, and Wigderson [43] (the "INW generator"). The INW generator samples n pseudorandom bits as follows:

1. Recursively construct a PRG $G\colon \{0,1\}^s \to \{0,1\}^{n/2}$.
2. Sample a uniform random vertex X and a uniform random neighbor Y in a low-degree expander graph on 2^s vertices.
3. Output the concatenation $G(X) \circ G(Y)$.

Several decades after its introduction [43], we are still learning more and more about what the INW generator is capable of. It has been shown to work particularly well for (standard-order) *regular* and *permutation* ROBPs, defined next.

Definition 7 (Regular and permutation ROBPs)**.** Let f be a width-w length-n standard-order ROBP with transition functions $f_1, \ldots, f_n\colon [w] \times \{0,1\} \to [w]$. We say that f is a *permutation ROBP* if, for every $i \in [n]$ and every $b \in \{0,1\}$, the function $f_i(\cdot, b)$ is a permutation on $[w]$. More generally, we say that f is *regular* if, for every $i \in [n]$ and every $u \in [w]$, we have $|f_i^{-1}(u)| = 2$.

Regular and permutation ROBPs have been studied extensively over the course of roughly the past decade [16, 17, 27, 46, 75, 66, 19, 1, 40, 62, 63, 26, 64, 11, 35, 48]. We now have various types of pseudorandomness results for regular or permutation ROBPs that are superior to the best corresponding results for general ROBPs, including constructions of PRGs [16, 17, 27, 46, 75, 66, 19, 40, 48], WPRGs [63], and HSGs [16, 11]. In many cases, the proofs consist of improved analyses of the classic INW construction [43] (with modified parameters). In other cases, the INW generator is one of multiple ingredients.

5.3 Unbounded-Width ROBPs

The first few papers on regular and permutation ROBPs [16, 17, 27, 46, 75, 66] focused on constant-width programs. Arguably the most important case is that of polynomial-width programs. The trend recently has been to study the intriguing setting of *unbounded-width* programs [40, 62, 63, 64, 11, 35, 48].

Without further constraints, unbounded-width standard-order permutation ROBPs are too powerful to be interesting: they can compute all Boolean functions. Therefore, we assume that the program has a bounded number of *accepting states* in the final layer. Admittedly, width is a more natural complexity measure than the number of accepting states, but it turns out that programs with a bounded number of accepting states have some interesting properties. Even with just *one* accept state, exponential-width standard-order permutation ROBPs can compute doubly-exponentially many distinct functions:

Proposition 1 ([40])**.** *Let $n \in \mathbb{N}$ be a positive even integer, and let $\pi\colon \{0,1\}^{n/2} \to \{0,1\}^{n/2}$ be a permutation. There exists a width-$(2^{n/2})$ length-n standard-order permutation ROBP f computing the following function:*

$$f(x,y) = 1 \iff \pi(x) = y.$$

(Briefly, to prove Proposition 1, we use the state space $\{0,1\}^{n/2}$. The all-zeroes state is the start state and the unique accepting state. We XOR x into our state, then apply π to our state, then XOR y into our state.) On the other hand, one can check that the majority function on three bits cannot be computed by a standard-order regular ROBP with a single accept vertex, no matter how wide the program is. Thus, these strange unbounded-width models have both dramatic strengths and dramatic weaknesses. One of the most striking results in this area is the following theorem by Pyne and Vadhan [63].

Theorem 9 (WPRGs for unbounded-width permutation ROBPs [63])**.** *For every $n \in \mathbb{N}$ and $\varepsilon > 0$, there is an explicit ε-WPRG for unbounded-width standard-order permutation ROBPs with a single accept state with seed length*

$$\tilde{O}\left(\log^{3/2} n + \log n \cdot \sqrt{\log(1/\varepsilon)} + \log(1/\varepsilon)\right).$$

Theorem 9 has implications for the more conventional setting of bounded-width standard-order permutation ROBPs. Every ε-WPRG for programs with one accepting state automatically (εm)-fools programs with m accepting states. Therefore, Theorem 9 implies an explicit WPRG for width-n length-n standard-order permutation ROBPs (with any number of accepting vertices) with error $1/n$ and seed length $\tilde{O}(\log^{3/2} n)$, compared to Nisan's $O(\log^2 n)$ bound.

Theorem 9 also helps to clarify the importance of weights. When $\varepsilon = 1/n$, the seed length in Theorem 9 is $\tilde{O}(\log^{3/2} n)$. In contrast, Hoza, Pyne, and Vadhan proved that every *unweighted* PRG that $(1/n)$-fools unbounded-width standard-order permutation ROBPs with a single accept vertex must have seed length $\Omega(\log^2 n)$ [40]. Therefore, in at least one natural setting, WPRGs are *intrinsically more powerful* than traditional PRGs.

The proof of Theorem 9 uses the INW generator, the inverse Laplacian perspective, and error reduction techniques, among other ideas.

5.4 The Permutation Case and the Monotone Case: Opposite Extremes

Why study regular and permutation ROBPs? The main reason is the hope that studying these special cases will lead to improvements in the general case. Indeed, there is a reduction showing that good PRGs or HSGs for polynomial-width standard-order *regular* ROBPs imply good PRGs or HSGs for *all* polynomial-width standard-order ROBPs [67, 11].[12]

In addition to that reduction [67, 11], there is another approach for constructing PRGs for constant-width standard-order ROBPs (albeit a vague and speculative one). At an intuitive level, one can argue that permutation ROBPs and monotone ROBPs are "opposites" of one another. In a permutation ROBP, edges with the same label never collide, whereas in a monotone ROBP, the only way that a layer can do any nontrivial computation is by introducing collisions. Now, we have one set of techniques that works well for (standard-order) permutation ROBPs: spectral expanders and the INW generator. Meanwhile, we have another set of techniques that works well for (arbitrary-order) monotone ROBPs: iterated restrictions with early termination. Given that these two sets of techniques cover two "extreme" classes of constant-width ROBPs, it is natural to try to combine the two sets of techniques. Could this approach yield an explicit PRG that fools *all* width-w standard-order ROBPs, with seed length $\tilde{O}(\log n)$ when w is a constant? The idea might sound a bit naïve or fantastical, especially considering the difficulty discussed in Section 2.4. Remarkably, however, Meka, Reingold, and Tal proved that the answer is yes for the case $w = 3$ [50].

Theorem 10 (PRGs for width-3 ROBPs [50])**.** *For every $n \in \mathbb{N}$ and $\varepsilon > 0$, there is an explicit ε-PRG for width-3 standard-order ROBPs with seed length $\tilde{O}(\log n \cdot \log(1/\varepsilon))$.*

[12] Note that Theorem 8 implies a *non-black-box* algorithm for estimating the expectation of a given standard-order regular ROBP in near-logarithmic space. Unfortunately, the reduction from the general case to the regular case does not work in the non-black-box setting.

To prove Theorem 10, Meka, Reingold, and Tal first show how to sample pseudorandom restrictions that preserve the expectation of width-3 arbitrary-order ROBPs. For this first step, one can alternatively use Forbes and Kelley's analysis (Theorem 3), which works more generally for width-w arbitrary-order ROBPs where w is small. (The papers of Forbes and Kelley [31] and Meka, Reingold, and Tal [50] are independent.)

Next, Meka, Reingold, and Tal show that width-3 arbitrary-order ROBPs simplify after a few pseudorandom restrictions [50]. And what does "simplify" mean in this context? Roughly speaking, they show that the program becomes *more and more permutation-like* as the restrictions are applied. After $\text{poly}(\log\log(n/\varepsilon))$ many restrictions, they terminate the restriction process and apply the INW generator [43] as the final step. Building on Braverman, Rao, Raz, and Yehudayoff's analysis [16], they show that the INW generator fools these "highly permutation-like" ROBPs with seed length $\tilde{O}(\log n \cdot \log(1/\varepsilon))$ [50] (in the standard-order case).

It remains an open problem to design an explicit PRG (or WPRG or HSG) for width-4 standard-order ROBPs with seed length $o(\log^2 n)$.

6 Conclusions

We continue to make steady, substantial progress toward proving $\mathsf{L} = \mathsf{BPL}$. The past few years alone have yielded many exciting results and developments. The problem remains challenging, but there does not seem to be any firm obstacle preventing further breakthroughs.

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