Stochastic Bias-Reduced Gradient Methods

Abstract

We develop a new primitive for stochastic optimization: a low-bias, low-cost estimator of the minimizer $x_\star$ of any Lipschitz strongly-convex function f. In particular, we use a multilevel Monte-Carlo approach due to Blanchet and Glynn [9] to turn any optimal stochastic gradient method into an estimator of $x_\star$ with bias δ, variance $O(\log(1/\delta))$, and an expected sampling cost of $O(\log(1/\delta))$ stochastic gradient evaluations. As an immediate consequence, we obtain cheap and nearly unbiased gradient estimators for the Moreau envelope of any Lipschitz convex function. We demonstrate the potential of our estimator through four applications. First, we develop a method for minimizing the maximum of N functions, improving on recent results and matching a lower bound up logarithmic factors. Second and third, we recover state-of-the-art rates for projection-efficient and gradient-efficient optimization using simple algorithms with a transparent analysis. Finally, we show that an improved version of our estimator would yield a nearly linear-time, optimal-utility, differentially-private non-smooth stochastic optimization method.

1 Introduction

Consider the fundamental problem of minimizing a μ-strongly convex function $F : \mathcal{X} \to \mathbb{R}$ given access to a stochastic (sub-)gradient estimator $\hat{\nabla} F$ satisfying $\mathbb{E}\, \hat{\nabla} F(x) \in \partial F(x)$ and $\mathbb{E}\|\hat{\nabla} F(x)\|^2 \le G^2$ for every $x \in \mathcal{X}$. Is it possible to transform the unbiased estimator $\hat{\nabla} F$ into a (nearly) unbiased estimator of the minimizer $x_\star := \operatorname{argmin}_{x \in \mathcal{X}} F(x)$? In particular, can we improve upon the $O(G/(\mu\sqrt{T}))$ bias achieved by T iterations of stochastic gradient descent (SGD)?

In this paper, we answer this question in the affirmative, proposing an *optimum estimator* $\hat{x}_\star$, which (for any fixed $\delta > 0$) has

$$\text{bias } \|\mathbb{E}\hat{x}_\star - x_\star\| = O(\delta) \text{ and variance } \mathbb{E}\|\hat{x}_\star - \mathbb{E}\hat{x}_\star\|^2 = O\!\left(\frac{G^2}{\mu^2}\log\!\left(\frac{G}{\mu\delta}\right)\right),$$

and, *in expectation*, costs $O(\log(\frac{G}{\mu\delta}))$ evaluations of $\hat{\nabla} F$.[1] Setting $\delta = G/(\mu\sqrt{T})$, we obtain the same bias bound as T iterations of SGD, but with expected cost of only $O(\log T)$ stochastic gradient evaluations (the worst-case cost is T).

Our estimator is an instance of the multilevel Monte Carlo technique for de-biasing estimator sequences [23] and more specifically the method of Blanchet and Glynn [9]. Our key observation is that this method is readily applicable to strongly-convex variants of SGD, or indeed any stochastic optimization method with the same (optimal) rate of convergence.

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[1]When $\mathcal{X} = \mathbb{B}_R(x_0) \subset \mathbb{R}^d$, $F(x) = \frac{1}{n}\sum_{i \in [n]} \hat{F}(x; i)$, and $\hat{\nabla} F$ is the subgradient of a uniformly random $\hat{F}(x; i)$ we can also get an estimator with bias 0 and expected cost $O(\log(nd))$. See Appendix A for details.

Objective	Expensive operation $\mathcal{O}$	$\mathcal{N}_{\mathcal{O}}$	$\mathbb{E}\mathcal{N}_{\hat{\nabla} f}$
$\max_{i \in [N]} f_{(i)}(x)$ (Sec. 4)	$f_{(1)}(x), \ldots, f_{(N)}(x)$	$\widetilde{O}(\epsilon^{-2/3})$	
$f(x)$ in domain $\mathcal{X}$ (Sec. 3)	$\mathsf{Proj}_{\mathcal{X}}(x)$	$O(\epsilon^{-1})$	$\widetilde{O}(\epsilon^{-2})$
$\Lambda(x) + f(x)$ for L-smooth Λ (Sec. 5)	$\nabla\Lambda(x)$	$O(\sqrt{L/\epsilon})$	

Table 1. Summary of our applications of accelerated bias-reduced stochastic gradient methods. We use $\mathcal{N}_{\mathcal{O}}$ and $\mathcal{N}_{\hat{\nabla} f}$ to denote the number of expensive operations and subgradient estimations, respectively. The $\widetilde{O}$ notation hides polylogarithmic factors. See Section 1.2 for additional description.

1.1 Estimating proximal points and Moreau-Yoshida envelope gradients

Given a convex function f and regularization level λ, the proximal point of y is $\mathsf{P}_{f,\lambda}(y) :=$ $\operatorname{argmin}_{x \in \mathbb{R}^d}\{f(x) + \frac{\lambda}{2}\|x - y\|^2\}$. Since computing $\mathsf{P}_{f,\lambda}$ amounts to solving a λ-strongly-convex problem, our technique provides low-bias and cheap proximal point estimators. Proximal points are ubiquitous in optimization [34, 18, 40, 32] and estimating them efficiently with low bias opens up new algorithmic possibilities. One of these possibilities is estimating the gradient of the Moreau-Yoshida envelope $f_\lambda(y) := \min_{x \in \mathbb{R}^d}\{f(x) + \frac{\lambda}{2}\|x - y\|^2\}$, which is a λ-smooth, $G^2/(2\lambda)$-accurate approximation of any G-Lipschitz f (see, e.g., [34, 25] and Appendix C). Since $\nabla f_\lambda(y) = \lambda(y - \mathsf{P}_{f,\lambda}(y))$, our optimum estimator provides a low-bias estimator for $\nabla f_\lambda(y)$ with second moment and expected cost greater than those of $\hat{\nabla} f$ by only a logarithmic factor. Thus, for any non-smooth f we can turn $\hat{\nabla} f$ into a gradient estimator for the smooth surrogate f_λ, whose smoothness is independent of the problem dimension, allowing us to perform *dimension-free* randomized smoothing [19].

1.2 Applications via accelerated bias-reduced methods

Our optimum estimator is a new primitive in stochastic convex optimization and we expect it to find multiple applications. We now describe three such applications: the first improves on previously known complexity bounds while the latter two recover existing bounds more easily. For simplicity of presentation we assume (in the introduction only) $\mathbb{E}\|\hat{\nabla} f\|^2 \leq 1$ and unit domain size.

In each application, we wish to minimize an objective function given access to a cheap subgradient estimator $\hat{\nabla} f$ as well as an expensive application-specific operation $\mathcal{O}$ (e.g., a projection to a complicated set). Direct use of the standard stochastic gradient method finds an ϵ-accurate solution using $O(\epsilon^{-2})$ computations of both $\hat{\nabla} f$ and $\mathcal{O}$, and our goal is to improve the $\mathcal{O}$ complexity without hurting the $\hat{\nabla} f$ complexity.

To that end, we design stochastic accelerated methods consisting of T iterations, each one involving only a constant number of $\mathcal{O}$ and proximal point computations, which we approximate by averaging copies of our optimum estimator.[2] The low bias of our estimator allows us to bound $T \ll \epsilon^{-2}$ as though our proximal points were exact, while maintaining an $\widetilde{O}(\epsilon^{-2})$ bound on the total expected number of $\hat{\nabla} f$ calls.[3] Thus, we save expensive operations without substantially increasing the gradient estimation cost. Table 1 summarizes each application, and we briefly describe them below.

Minimizing the maximal loss. Given N convex, 1-Lipschitz functions $f_{(1)}, \ldots, f_{(N)}$ we would like to find an ϵ-approximate minimizer of their maximum $f_{\max}(x) = \max_{i \in [N]} f_{(i)}(x)$. This problem naturally arises when optimizing worst-case behavior, as in maximum margin classification and robust optimization [41, 16, 36, 7]. We measure complexity by the number of individual function and subgradient evaluations, so that the expensive operation of evaluating $f_{(1)}, \ldots, f_{(N)}$ at a single point has complexity $O(N)$. Carmon et al. [14] develop an algorithm for minimizing $f_{\max}$ with complexity $\widetilde{O}(N\epsilon^{-2/3} + \epsilon^{-8/3})$, improving on the subgradient method for sufficiently large N. Using our bias-reduced Moreau gradient envelope estimator in a Monteiro-Svaiter-type accelerated proximal point method [13, 12, 32], we obtain improved complexity $\widetilde{O}(N\epsilon^{-2/3} + \epsilon^{-2})$. This matches (up to logarithmic factors) a lower bound shown in [14], settling the complexity of minimizing the maximum of N non-smooth functions. Our result reveals a surprising fact: for $N \ll (GR/\epsilon)^{-4/3}$, minimizing the maximum of N functions is no harder than minimizing their average.

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[2]While averaging is parallelizable, our optimum estimator itself is sequential. Consequently, our approach does not yield improve parallelism; see Appendix B.1 for further discussion

[3]It is easy to turn expected complexity bounds into deterministic ones; see Appendix B.2.

Projection-efficient stochastic optimization via dimension-free randomized smoothing. Consider the problem of minimizing a convex function f using an unbiased gradient estimator $\tilde{\nabla} f$ over convex set $\mathcal{X}$ for which Euclidean projections are expensive to compute (for example, the cone of PSD matrices). When f is L-smooth, a stochastic version of Nesterov's accelerated gradient descent (AGD) [17] performs only $O(\sqrt{L/\epsilon})$ projections. For non-smooth f we instead instead apply AGD to the Moreau envelope smoothing of f (with appropriate $\lambda = O(\epsilon^{-1})$) using our nearly-unbiased stochastic estimator for ∇f_λ. This yields a solution in $O(\epsilon^{-1})$ projections and $\widetilde{O}(\epsilon^{-2})$ evaluations of $\tilde{\nabla} f$. Our algorithm provides a simple alternative to the recent work of Thekumparampil et al. [39] whose performance guarantees are identical up to a logarithmic factor.

Gradient-efficient composite optimization. We would like to minimize $\Psi(x) = \Lambda(x) + f(x)$, where Λ is convex and L-smooth but we can access it only via computing (expensive) exact gradients, while f is a non-smooth convex functions for which we have a (cheap) unbiased subgradient estimator $\hat{\nabla} f$. Problems of this type include inverse problems with sparsity constraint and regularized loss minimization in machine learning. To save $\nabla \Lambda$ computations, it is possible to use a composite version of AGD which solves $O(\sqrt{L/\epsilon})$ subproblems of the form $\text{minimize}_x \{ \langle \nabla \Lambda(y), x \rangle + f(x) + \frac{\beta}{2} \|x - x'\|^2 \}$. Lan [29] designed a careful subgradient method called "gradient sliding" for which the total cost of solving these subproblems is $O(\epsilon^{-2})$ evaluations of $\hat{\nabla} f$. We show that a simple alternative—estimating the subproblem solutions via our low-bias optimum estimator—recovers the same guarantee up to logarithmic factors.

1.3 Non-smooth differentially private stochastic convex optimization

We now discuss a potential application of our technique that is conditional on the existence of an improved optimum estimator. In it, we minimize the population objective function $f(x) = \mathbb{E}_{S \sim P} \hat{f}(x; S)$ under the well-known constraint of differential privacy [21]. Given n i.i.d. samples $S_i \sim P$ and assuming that each $\hat{f}$ is 1-Lipschitz, convex and sufficiently smooth, Feldman et al. [22] develop algorithms that obtain the optimal error and compute $O(n)$ subgradients of $\hat{f}$. The non-smooth case is more challenging and the best existing bound is $O(n^{11/8})$ for the high-dimensional setting $d = n$ [27, 3]. In Section 6 we show that our optimum estimator, combined with recent localization techniques [22], reduces the problem to private mean estimation. Unfortunately, our estimator has unbounded outputs, causing large sensitivity and leading to insufficient utility. Nevertheless, assuming a version of our estimator that has bounded outputs, we give an algorithm that queries $\widetilde{O}(n)$ subgradients for non-smooth functions, solving a longstanding open problem in private optimization [15, 4]. This motivates the study of improved versions of our estimators that have constant sensitivity.

1.4 Related work

Multilevel Monte-Carlo (MLMC) techniques originate from the literature on parametric integration for solving integral and differential equations [23]. We use an MLMC variant put forth by Blanchet and Glynn [9] for estimating functionals of expectations. Among several applications, Blanchet and Glynn [9, Section 5.2] propose an estimator for $\text{argmin}_x \mathbb{E}_{S \sim P} \hat{f}(x; S)$ where $\hat{f}(\cdot; s)$ is convex for all s and assuming access to minimizers of empirical objectives of the form $\sum_{i \in [N]} \hat{f}(x; s_i)$. They provide a preliminary analysis of the estimator's variance (later elaborated in [10]) using an asymptotic Taylor expansion around the population minimizer. In comparison, we study the more general setting of stochastic gradient estimators and provide a complete algorithm based on SGD, along with a non-asymptotic analysis and concrete settings where our estimator is beneficial.

A number of works have used the Blanchet-Glynn estimator in the context of optimization and machine learning. These applications include estimating the ratio of expectations for semisupervised learning [8], estimating gradients of distributionally robust optimization objectives [30], and estimating gradients in deep latent variable models [38]. Our estimator is similar to that of Levy et al. [30] in that we also have to pick a "critical" doubling probability for the (random) computational budget, which makes the expected cost and variance of our estimators depend logarithmically on the bias.

Notation. We let $\mathbb{B}_R(x) = \{ y \in \mathbb{R}^d : \|y - x\| \leq R \}$ denote the ball of radius R around x, where $\|\cdot\|$ is the Euclidean norm throughout. We write $\mathsf{Proj}_\mathcal{S}$ for the Euclidean projection to $\mathcal{S}$. We use $\mathsf{P}_{f,\lambda}$ to denote the proximal operator (2) and f_λ to denote the Moreau envelope (3). Finally, we use $\mathcal{N}_f$ and $\mathcal{N}_{\hat{\nabla} f}$ to denote function and subgradient estimator evaluation complexity, respectively.

Algorithm 1: OPTEST($\hat{\nabla}f, \psi, \mu, \delta, \sigma^2, \mathcal{X}$)	**Algorithm 2:** MORGRADEST($\hat{\nabla}f, y, \lambda, \delta, \sigma^2, \mathcal{X}$)
$\triangleright$ δ, σ^2 are required bias and square error	$\triangleright$ δ, σ^2 are required bias and square error
$\triangleright$ c is the ODC algorithm constant	$\triangleright$ λ is the regularization level
$T_{\max} = \left\lceil \frac{4cG^2}{\mu^2 \min\{\delta^2, \frac{1}{2}\sigma^2\}} \right\rceil$	$\triangleright$ y is the point at which to estimate $\nabla f_\lambda(y)$
$N = \left\lceil \frac{32cG^2 \log(T_{\max})}{\mu^2 \sigma^2} \right\rceil$	$\psi_\lambda(x) = \frac{\lambda}{2}\|x - y\|^2$
for $i = 1, \ldots, N$ **do**	$\hat{x}_\star = \text{OPTEST}\left(\hat{\nabla}f, \psi_\lambda, \lambda, \frac{\delta}{\lambda}, \frac{\sigma^2}{\lambda^2}, \mathcal{X}\right)$
$\quad \hat{x}_\star^{(i)}$ = a draw of the estimator (1)	**return** $\lambda(y - \hat{x}_\star)$
return $\frac{1}{N} \sum_{i=1}^{N} \hat{x}_\star^{(i)}$	

2 A multilevel Monte-Carlo optimum estimator

In this section, we construct a low-bias estimator for the minimizer of any strongly convex function $F : \mathcal{X} \to \mathbb{R}$. This estimator is the key component of our algorithms in the subsequent sections, which use it to approximate proximal points and Moreau envelope gradients. We assume that F is of the form $F = f + \psi$, where the function ψ is "simple"[4] and that f satisfies the following.

Assumption 1. *The function $f : \mathcal{X} \to \mathbb{R}$ is convex and is accessible only via an unbiased subgradient estimator $\hat{\nabla}f$ which satisfies $\mathbb{E}\|\hat{\nabla}f(x)\|^2 \leq G^2$ for all $x \in \mathcal{X}$.*

2.1 ODC algorithms

Our estimator can use, in a black-box fashion, any method for minimizing F with sufficiently fast convergence to $x_\star = \operatorname{argmin}_{x \in \mathcal{X}} F(x)$. We highlight the required convergence property as follows.

Definition 2.1. *An* optimal-distance-convergence *algorithm* ODC *takes as input $\hat{\nabla}f$ satisfying Assumption 1, a simple function ψ and a budget $T \geq 1$. If $F = f + \psi$ is μ-strongly convex with minimizer $x_\star$, the algorithm's output $x = \text{ODC}(\hat{\nabla}f, \psi, T)$ requires at most T evaluations of $\hat{\nabla}f$ to compute and satisfies $\mathbb{E}\|x - x_\star\|^2 \leq c\frac{G^2}{\mu^2 T}$ for some constant $c > 0$.*

Standard lower bound constructions imply that the $O(\frac{G^2}{\mu^2 T})$ squared distance convergence rate is indeed optimal; see Appendix B.3 for additional discussion. Conversely, ODC algorithms are readily available in the literature [35, 26] since any point x satisfying $\mathbb{E}F(x) - F(x_\star) = O(\frac{G^2}{\mu T})$ (the optimal rate of convergence in strongly convex, Lipschitz optimization) also satisfies $\mathbb{E}\|x - x_\star\|^2 \leq O(\frac{G^2}{\mu^2 T})$ by virtue of the strong convexity of F. We provide a concrete ODC algorithm consisting of a generalization of epoch SGD [26].

Lemma 2.1. EPOCHSGD *(Algorithm 8 in Appendix D.1) is an ODC algorithm with constant $c = 32$.*

2.2 Constructing an optimum estimator

To turn any ODC algorithm into a low-bias, low-cost and near-constant variance optimum estimator, we use the multilevel Monte Carlo (MLMC) technique of Blanchet and Glynn [9]. Given a problem instance $\hat{\nabla}f, \psi$, an algorithm ODC and a cutoff parameter $T_{\max} \in \mathbb{N}$, our estimator $\hat{x}_\star$ is:

Draw $J \sim \text{Geom}\left(\frac{1}{2}\right)$ and, writing $x_j := \text{ODC}(\hat{\nabla}f, \psi, 2^j)$, compute

$$\hat{x}_\star = x_0 + \begin{cases} 2^J(x_J - x_{J-1}) & 2^J \leq T_{\max} \\ 0 & \text{otherwise.} \end{cases} \quad (1)$$

We note that for certain ODC algorithms it is possible to extract x_0, x_{J-1} from the intermediate steps of computing x_J, so that we only need to invoke ODC once. This is particularly simple to do for EPOCHSGD, as we explain in Appendix D.1. The key properties of our estimator are as follows.

Proposition 1. *Let f and $\hat{\nabla}f$ satisfy Assumption 1, $F = f + \psi$ be μ-strongly convex with minimizer $x_\star$ and $T_{\max} \in \mathbb{N}$. For any ODC algorithm with constant c, the estimator (1) has bias $\|\mathbb{E}\hat{x}_\star - x_\star\| \leq$*

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[4]Our applications only use ψ of the form $\psi(x) = \frac{\lambda}{2}\|x - x'\|^2$ but our estimator applies more broadly to cases where $\operatorname{argmin}_x\{\langle v, x\rangle + \psi(x) + \frac{1}{2\eta}\|x - y\|^2\}$ is easy to compute for all v and y.

152 $\sqrt{2c}\frac{G}{\mu\sqrt{T_{\max}}}$ *and variance* $\mathbb{E}\|\hat{x}_\star - \mathbb{E}\hat{x}_\star\|^2 \le 16c\frac{G^2}{\mu^2}\log_2(T_{\max})$. *Moreover, the expected number of*
153 $\hat{\nabla}f$ *evaluations required to compute* $\hat{x}_\star$ *is* $O(\log T_{\max})$.

154 *Proof.* Let $j_{\max} = \max\{j \in \mathbb{N} \mid 2^j \le T_{\max}\} = \lfloor \log_2 T_{\max} \rfloor$. The expectation of $\hat{x}_\star$ is

$$\mathbb{E}\hat{x}_\star = \mathbb{E}x_0 + \sum_{j=1}^{j_{\max}} \mathbb{P}(J = j)2^j(\mathbb{E}x_j - \mathbb{E}x_{j-1}) = \mathbb{E}x_{j_{\max}},$$

155 where the second equality follows from $\mathbb{P}(J = j) = 2^{-j}$ and the sum telescoping. Noting that
156 $x_{j_{\max}} = \text{ODC}(\hat{\nabla}f, \psi, T)$ for $T = 2^{j_{\max}} \ge T_{\max}/2$, we have that

$$\|\mathbb{E}x_{j_{\max}} - x_\star\| \le \sqrt{\mathbb{E}\|x_{j_{\max}} - x_\star\|^2} \le \sqrt{c}\frac{G}{\mu\sqrt{T_{\max}/2}}$$

157 by Definition 2.1. To bound the variance we use $\|a + b\|^2 \le 2\|a\|^2 + 2\|b\|^2$ and note that

$$\mathbb{E}\|\hat{x}_\star - \mathbb{E}\hat{x}_\star\|^2 \le \mathbb{E}\|\hat{x}_\star - x_\star\|^2 \le 2\mathbb{E}\|\hat{x}_\star - x_0\|^2 + 2\mathbb{E}\|x_0 - x_\star\|^2.$$

158 The ODC property implies that $\mathbb{E}\|x_0 - x_\star\|^2 \le cG^2/\mu^2$. For the term $\mathbb{E}\|\hat{x}_\star - x_0\|^2$ we have

$$\mathbb{E}\|\hat{x}_\star - x_0\|^2 = \sum_{j=1}^{j_{\max}} \mathbb{P}(J = j)2^{2j}\mathbb{E}\|x_j - x_{j-1}\|^2 = \sum_{j=1}^{j_{\max}} 2^j\mathbb{E}\|x_j - x_{j-1}\|^2, \text{ and}$$

$$\mathbb{E}\|x_j - x_{j-1}\|^2 \le 2\mathbb{E}\|x_j - x_\star\|^2 + 2\mathbb{E}\|x_{j-1} - x_\star\|^2 \le 6c\frac{G^2}{\mu^2}2^{-j}.$$

159 Substituting, we get $\mathbb{E}\|\hat{x}_\star - x_0\|^2 \le 6c\frac{G^2}{\mu^2}j_{\max}$ and $\mathbb{E}\|\hat{x}_\star - \mathbb{E}\hat{x}_\star\|^2 \le 16c\frac{G^2}{\mu^2}\log_2(T_{\max})$. Finally,
160 the expected number of $\hat{\nabla}f$ evaluations is $1 + \sum_{j=1}^{j_{\max}} \mathbb{P}(J = j)(2^j + 2^{j-1}) = O(j_{\max})$. $\qquad\square$

161 The function OPTEST in Algorithm 1 computes an estimate of $x_\star$ with and desired bias δ and square
162 error σ^2 by averaging independent draws of the MLMC estimator (1). The following guarantees are
163 immediate from Proposition 1; see Appendix D.2 for a short proof.

164 **Theorem 1.** *Let f and $\hat{\nabla}f$ satisfy Assumption 1, $F = f + \psi$ be μ-strongly convex with minimizer*
165 *$x_\star \in \mathcal{X}$, and $\delta, \sigma > 0$. The function* OPTEST$(\hat{\nabla}f, \psi, \mu, \delta, \sigma^2, \mathcal{X})$ *outputs $\hat{x}_\star$ satisfying $\|\mathbb{E}\hat{x}_\star - x_\star\| \le$*
166 *δ and $\mathbb{E}\|\hat{x}_\star - x_\star\|^2 \le \sigma^2$ using $\mathcal{N}_{\hat{\nabla}f}$ stochastic gradient computations, where*

$$\mathbb{E}\mathcal{N}_{\hat{\nabla}f} = O\left(\frac{G^2}{\mu^2\sigma^2}\log^2\left(\frac{G}{\mu\min\{\delta,\sigma\}}\right) + \log\frac{G}{\mu\min\{\delta,\sigma\}}\right).$$

2.3 Estimating proximal points and Moreau envelope gradients

168 Recall that the proximal point of function f with regularization level λ at point y is

$$\mathsf{P}_{f,\lambda}(y) := \operatorname*{argmin}_{x\in\mathcal{X}}\left\{f(x) + \tfrac{\lambda}{2}\|x - y\|^2\right\}. \tag{2}$$

169 When f satisfies Assumption 1, we may use OPTEST (with $\psi(x) = \frac{\lambda}{2}\|x - y\|^2$ and $\mu = \lambda$) to obtain
170 a reduced-bias proximal point estimator. The proximal point $\mathsf{P}_{f,\lambda}(y)$ is closely related to the Moreau
171 envelope

$$f_\lambda(y) := \min_{x\in\mathcal{X}}\left\{f(x) + \tfrac{\lambda}{2}\|x - y\|^2\right\} \tag{3}$$

172 via the relationship $\nabla f_\lambda(y) = \lambda(y - \mathsf{P}_{f,\lambda}(y))$ (see Appendix C). Therefore, we can use our optimum
173 estimator to turn $\widetilde{O}(1)$ calls to $\hat{\nabla}f$ into a nearly unbiased estimator for ∇f_λ. We formulate this as:

174 **Corollary 2.1.** *Let f and $\hat{\nabla}f$ satisfy Assumption 1, let $y \in \mathcal{X}$ and let $\lambda, \sigma, \delta > 0$. The func-*
175 *tion* MORGRADEST$(\hat{\nabla}f, \lambda, y, \delta, \sigma^2, \mathcal{X})$ *outputs $\hat{\nabla}f_\lambda(y)$ satisfying $\|\mathbb{E}\hat{\nabla}f_\lambda(y) - \nabla f_\lambda(y)\| \le \delta$ and*
176 *$\mathbb{E}\|E\hat{\nabla}f_\lambda(y) - \nabla f_\lambda(y)\|^2 \le \sigma^2$ and has complexity $\mathbb{E}\mathcal{N}_{\hat{\nabla}f} = O\left(\log\left(\frac{G}{\lambda\delta}\right) + \frac{G^2}{\sigma^2}\log^2\left(\frac{G}{\lambda\delta}\right)\right)$.*

3 Projection-efficient convex optimization

In this section, we combine the bias-reduced Moreau envelope gradient estimator with a standard accelerated gradient method to recover the result of Thekumparampil et al. [39]. We consider the problem of minimizing a function f satisfying Assumption 1 over the domain $\mathbb{B}_R(0)$ subject to the constraint $x \in \mathcal{X}$, where $\mathcal{X} \subset \mathbb{B}_R(0)$ is a complicated convex set that we can only access via (expensive) projections of the form $\mathsf{Proj}_{\mathcal{X}}(x) := \operatorname{argmin}_{y \in \mathcal{X}} \|y - x\|$. We further assume that an initial point $x_0 \in \mathcal{X}$ satisfies $\|x_0 - x_\star\| \le D$.

Algorithm 3: Stochastic accelerated gradient descent on the Moreau envelope

Input: A gradient estimator $\hat{\nabla} f$ satisfying Assumption 1 in $\mathbb{B}_R(x)$, projection oracle $\mathsf{Proj}_{\mathcal{X}}$, and initial point $x_0 = v_0$ with $\|x_0 - x_\star\| \le D$.

Parameters : iteration budget T, Moreau regularization λ, approximation parameters δ_k, σ_k^2

1 **for** $k = 0, 1, \cdots, T-1$ **do**
2 $\quad$ $y_k = \frac{k}{k+2} x_k + \frac{2}{k+2} v_k$
3 $\quad$ $g_k = \mathrm{MorGradEst}(\hat{\nabla} f, y_k, \lambda, \delta_k, \sigma_k^2, \mathbb{B}_R(0))$
4 $\quad$ $x_{k+1} = \mathsf{Proj}_{\mathcal{X}}\left(x_k - \frac{1}{3\lambda} g_k\right)$
5 $\quad$ $v_{k+1} = \mathsf{Proj}_{\mathbb{B}_R(0)}\left(v_k - \frac{k+1}{6\lambda} g_k\right)$
6 **return** x_T

Algorithm 3 applies Nesterov's accelerated gradient descent method on the (λ-smooth) Moreau envelope f_λ defined in eq. (3). Since computing the Moreau envelope does not involve projection to $\mathcal{X}$, for sufficiently accurate approximation of ∇f_λ we require only $T = O(\sqrt{\lambda D^2/\epsilon})$ projections to $\mathcal{X}$ for finding an $O(\epsilon)$-suboptimal point of f_λ constrained to $\mathcal{X}$. For that point to be also ϵ-suboptimal for f itself, we must choose λ of the order of G^2/ϵ, so that the number of projections is $O(GD/\epsilon)$.

As noted in [39] computing ∇f_λ to accuracy $O(\epsilon/R)$ is sufficient for the above guarantee to hold, but doing so using a stochastic gradient method requires $O((GD/\epsilon)^2)$ evaluations of $\hat{\nabla} f$ per iteration, and $O((GD/\epsilon)^3)$ evaluations in total. To improve this, we employ Algorithm 2 to compute nearly-unbiased estimates for ∇f_λ and bound the error incurred by their variance. Our result matches the gradient sliding-based technique of Thekumparampil et al. [39] up to polylogarithmic factors while retaining the conceptual simplicity of directly applying AGD on the Moreau envelope. We formally state the guarantees of our method below, and provide a self-contained proof in Appendix E.

Theorem 2. *Let* $f : \mathbb{B}_R(0) \to \mathbb{R}$ *and* $\hat{\nabla} f$ *satisfy Assumption 1. Let* $\mathcal{X} \subseteq \mathbb{B}_R(0)$ *be a convex set admitting a projection oracle* $\mathsf{Proj}_{\mathcal{X}}$*. Let* $x_0 \in \mathcal{X}$ *be an initial point with* $\|x - x_\star\| \le D$ *for some* $x_\star \in \mathcal{X}$*. With* $\lambda = \frac{2G^2}{\epsilon}$*,* $\delta_k = \frac{\epsilon}{8R}$*,* $\sigma_k^2 = \frac{2\epsilon\lambda}{k+2}$*, and* $T = \frac{7GD}{\epsilon}$ *Algorithm 3 computes* $x \in \mathcal{X}$ *with* $\mathbb{E}\left[f(x)\right] \le f(x_\star) + \epsilon$ *with complexity* $\mathbb{E}\mathcal{N}_{\hat{\nabla} f} = O\left(\frac{G^2 D^2}{\epsilon^2} \log^2\left(\frac{GR}{\epsilon}\right)\right)$ *and* $O\left(\frac{GD}{\epsilon}\right)$ *calls to* $\mathsf{Proj}_{\mathcal{X}}$*.*

4 Stochastic accelerated proximal methods and minimizing the maximal loss

In this section we apply our estimator in an accelerated proximal point method and use it to obtain an optimal rate for minimizing the maximum of N convex functions (up to logarithmic factors).

4.1 Accelerated proximal point method via Moreau envelope gradient estimation

Algorithm 4 uses our reduced-bias Moreau envelope gradient estimator to minimize a function f satisfying Assumption 1. To explain the method, we contrast it with stochastic AGD on the Moreau envelope (Algorithm 3). First and foremost, Algorithm 3 provides a suboptimality bound on the Moreau envelope f_λ (which for small λ is far from f) while Algorithm 4 minimizes f itself. Second, while Algorithm 3 uses a fixed regularization parameter λ, Algorithm 4 handles an arbitrary sequence $\{\lambda_k\}$ given by a black-box function NextLambda. To facilitate our application of the method to minimizing the maximal loss—where gradient estimation is only tractable in small Euclidean balls—we include an optional parameter r such that the proximal point movement bound $\|\mathsf{P}_{f,\lambda_{k+1}}(y_k) - y_k\| \le r$ holds for all k. However, most of our analysis of Algorithm 4 does not require this parameter (i.e., holding for $r = \infty$), making it potentially applicable to other settings that use accelerated proximal point methods [13, 12, 32].

Algorithm 4: Stochastic accelerated proximal point method

Input: Gradient estimator $\hat{\nabla} f$, function NEXTLAMBDA, initialization $x_0 = v_0$ and $A_0 \geq 0$.
Parameters: Approximation parameters $\{\varphi_k, \delta_k, \sigma_k\}$, stopping parameters $A_{\max}$ and $K_{\max}$,
    optional movement bound $r > 0$.

1 **for** $k = 0, 1, \ldots$ **do**
2 $\quad \lambda_{k+1} = \text{NEXTLAMBDA}(x_k, v_k, A_k) \; \triangleright \text{ guaranteeing that } \|\mathsf{P}_{f,\lambda_{k+1}}(y_k) - y_k\| \leq r$
3 $\quad a_{k+1} = \frac{1}{2\lambda_{k+1}}\sqrt{1 + 4\lambda_{k+1}A_k}$ and $A_{k+1} = A_k + a_{k+1}$ and $\mathcal{X}_k = \mathcal{X} \cap \mathbb{B}_r(y_k)$
4 $\quad y_k = \frac{A_k}{A_{k+1}}x_k + \frac{a_{k+1}}{A_{k+1}}v_k$ and $x_{k+1} = \tilde{\mathsf{P}}_{f,\lambda_{k+1}}^{\varphi_{k+1}}(y_k) \; \triangleright \text{ defined in eq. (4)}$
5 $\quad g_{k+1} = \text{MORGRADEST}(\hat{\nabla} f, y_k, \lambda_{k+1}, \delta_k, \sigma_k^2, \mathcal{X}_k)$ and $v_{k+1} = \text{Proj}_{\mathcal{X}}\left(v_k - \frac{1}{2}a_{k+1}g_{k+1}\right)$
6 $\quad$ **if** $A_{k+1} \geq A_{\max}$ **or** $k + 1 = K_{\max}$ **then return** x_{k+1}

---

215 The third and final notable difference between Algorithms 3 and 4 is the method of updating the x_k
216 iteration sequence. While a projected stochastic gradient descent step suffices for Algorithm 3, here
217 we require a more direct approximation of function value decrease attained by the exact proximal
218 mapping $\mathsf{P}_{f,\lambda}$ (see eq. (2)). For a given accuracy φ, we define the φ-approximate proximal mapping

$$\tilde{\mathsf{P}}_{f,\lambda}^{\varphi}(y) \coloneqq \text{ any } x \in \mathcal{X} \text{ such that } \mathbb{E}F(x) \leq F(\mathsf{P}_{f,\lambda}(y)) + \varphi \text{ for } F(z) \coloneqq f(z) + \frac{\lambda}{2}\|z - y\|^2. \quad (4)$$

219 Note that $\tilde{\mathsf{P}}_{f,\lambda}^0 = \mathsf{P}_{f,\lambda}$ and that for $\varphi > 0$ we can compute $\tilde{\mathsf{P}}_{f,\lambda}^{\varphi}$ with an appropriate SGD variant
220 (such as EPOCHSGD) using $O(G^2/(\lambda\varphi))$ evaluations of $\hat{\nabla} f$.

221 With the differences between the algorithms explained, we emphasize their key similarity: both
222 algorithms update the v_k sequence using our bias reduction method MORGRADEST (Algorithm 2),
223 which holds the key to their efficiency. The following proposition shows that Algorithm 4 has the
224 same bound on $K_{\max}$ as an exact accelerated proximal point method [13], while requiring at most
225 $\widetilde{O}(G^2 R^2 \epsilon^{-2})$ stochastic gradient evaluations; see proof in Appendix F.1.

226 **Proposition 2.** *Let $f : \mathcal{X} \to \mathbb{R}$ and $\hat{\nabla} f$ satisfy Assumption 1, and let $\mathcal{X} \subseteq \mathbb{B}_R(x_0)$. For a*
227 *target accuracy $\epsilon \leq GR$ let $\varphi_{k+1} = \frac{\epsilon}{60\lambda_{k+1}a_{k+1}}$, $\delta_{k+1} = \frac{\epsilon}{120R}$, $\sigma_{k+1}^2 = \frac{\epsilon}{60a_{k+1}}$, $A_0 = \frac{R}{G}$, and*
228 $A_{\max} = \frac{9R^2}{\epsilon}$. *If $\lambda_k \geq \lambda_{\min} \geq \frac{1}{A_{\max}} = \Omega(\frac{\epsilon}{R^2})$ for all $k \leq K_{\max}$, then lines 4 and 5 of Algorithm 4*
229 *have total complexity $\mathbb{E}\mathcal{N}_{\hat{\nabla} f} = O\left(K_{\max} \log \frac{GR}{\epsilon} + \frac{G^2 R^2}{\epsilon^2} \log^2 \frac{GR}{\epsilon}\right)$. If in addition $\|\mathsf{P}_{f,\lambda_k}(y_{k-1}) -$*
230 $y_{k-1}\| \geq 3r/4$ *whenever $\lambda_k \geq 2\lambda_{\min}$ then for $K_{\max} = O\left(\left(\frac{R}{r}\right)^{2/3} \log\left(\frac{GR}{\epsilon}\right) + \sqrt{\frac{\lambda_{\min}R^2}{\epsilon}}\right)$, the*
231 *algorithm's output x_K satisfies $f(x_K) - f(x_\star) \leq \epsilon$ with probability at least $\frac{2}{3}$.*

4.2 Minimizing the maximal loss

233 We now consider objectives of the form $f_{\max}(x) \coloneqq \max_{i \in [N]} f_{(i)}(x)$ where each function $f_{(i)} :$
234 $\mathcal{X} \to \mathbb{R}$ is convex and G-Lipschitz. Our approach to minimizing $f_{\max}$ largely follows Carmon et al.
235 [14] except we replace their approximate accelerated proximal point method with Algorithm 4. The
236 first step is to replace $f_{\max}$ with the "softmax" function, defined for a given target accuracy ϵ as

$$f_{\text{smax}}(x) \coloneqq \epsilon' \log\left(\sum_{i \leq N} \exp\left(f_{(i)}(x)/\epsilon'\right)\right), \text{ where } \epsilon' \coloneqq \frac{\epsilon}{2 \log N}.$$

237 Since $f_{\text{smax}}(x) - f_{\max}(x) \in [0, \frac{\epsilon}{2}]$, any $\frac{\epsilon}{2}$-accurate solution of f_{smax} is ϵ-accurate for $f_{\max}$.

238 The second step is to develop an efficient gradient estimator for f_{smax}; this is non-trivial because
239 f_{smax} is not a finite sum or expectation. In [14] this is addressed via an "exponentiated softmax" trick,
240 but we develop a different rejection sampling-based approach that fits Algorithm 4 more directly (see
241 Algorithm 9). To produce an unbiased estimate for $\nabla f_{\text{smax}}(x)$ for x in a ball of radius $r_\epsilon = \epsilon'/G$
242 we require a single $\nabla f_{(i)}(x)$ evaluation (for some i), $O(1)$ evaluations of $f_{(i)}(x)$ in expectation, and
243 evaluation of the N functions $f_{(1)}(y), \ldots, f_{(N)}(y)$ for pre-processing. Plugging this estimator into
244 Algorithm 4 with $r = r_\epsilon$, the total pre-processing overhead of lines 4 and 5 is $O(K_{\max}N)$.

245 The final step is to find a function NEXTLAMBDA such that $\|\mathsf{P}_{f_{\text{smax}},\lambda_{t+1}}(y_k) - y_k\| \leq r_\epsilon$ for all k
246 (enabling gradient estimation), and $\|\mathsf{P}_{f_{\text{smax}},\lambda_{t+1}}(y_k) - y_k\| \geq \frac{3}{4}r_\epsilon$ when $\lambda_{k+1} > 2\lambda_{\min}$ (allowing
247 us to bound $K_{\max}$ with Proposition 2). Here we use the bisection subroutine from [14] as is (see
248 Algorithm 10). By judiciously choosing $\lambda_{\min}$—an improvement over the analysis in [14]—we obtain
249 the following complexity guarantee on $\mathcal{N}_{f_{(i)}}$ and $\mathcal{N}_{\partial f_{(i)}}$, the total numbers of individual function and
250 subgradient evaluations, respectively. (See proof Appendix F.2).

251 **Theorem 3.** *Let* $f_{(1)}, \ldots, f_{(N)} : \mathcal{X} \to \mathbb{R}$ *be convex and G-Lipschitz and let* $\mathcal{X} \subseteq \mathbb{B}_R(x_0)$. *For*
252 *any* $\epsilon < \frac{1}{2}GR/\log N$, *Algorithm 4 (with* $\tilde{\mathsf{P}}^\varphi_{f_{\text{smax}},\lambda}$ *implemented in Algorithm 8,* $\hat{\nabla}f_{\text{smax}}$ *given by*
253 *Algorithm 9 and* NEXTLAMBDA *given by Algorithm 10 with* $\lambda_{\min} = \widetilde{O}(\epsilon/(r_\epsilon^{4/3}R^{2/3}))$ *outputs* $x \in$
254 $\mathcal{X}$ *that with probability at least* $\frac{1}{2}$ *is ϵ-suboptimal for* $f_{\max}(x) = \max_{i \in [N]} f_{(i)}(x)$ *and has complexity*

255
$$\mathbb{E}\mathcal{N}_{f_{(i)}} = O\left(\left[N\left(\frac{GR\log N}{\epsilon}\right)^{2/3} + \left(\frac{GR}{\epsilon}\right)^2\right]\log^2\frac{GR}{\epsilon}\right) \text{ and } E\mathcal{N}_{\partial f_{(i)}} = O\left(\left(\frac{GR}{\epsilon}\right)^2\log^2\frac{GR}{\epsilon}\right).$$

256 The rate given by Theorem 3 matches (up to logarithmic factors) the lower bound of $\Omega(N(GR/\epsilon)^2 +$
257 $(GR/\epsilon)^2)$ shown in [14] and is therefore near-optimal.

5 Gradient-efficient composite optimization

259 Consider the problem of finding a minimizer of the following convex composite optimization problem
260

$$\underset{x \in \mathcal{X}}{\text{minimize}} \ \Psi(x) := \Lambda(x) + f(x) \text{ where } \Lambda \text{ is } L\text{-smooth and } f \text{ satisfies Assumption 1,} \quad (5)$$

261 given x_0 such that $\|x_0 - x_\star\| \leq R$ for some $x_\star \in \arg\min_{x \in \mathcal{X}} \Psi(x)$. Lan [29] developed a
262 method called "gradient sliding" that finds an ϵ-accurate solution to (5) with complexities $\mathcal{N}_{\nabla\Lambda} =$
263 $O(\sqrt{LR^2/\epsilon})$ and $\mathcal{N}_{\hat{\nabla}f} = O((GR/\epsilon)^2)$, which are optimal even for each component separately.[5]

264 In this section, we provide an alternative algorithm that matches the complexity of gradient up to
265 logarithmic factors and conceptually simple. Our approach, Algorithm 5, is essentially composite
266 AGD [33], where at the kth iteration we compute a proximal point (2) with respect to a partial
267 linearization of Ψ around y_k. In particular, letting $\bar{\Lambda}_k(v) := \Lambda(y_k) + \langle\nabla\Lambda(y_k), v - y_k\rangle$ and
268 $\beta_k = \frac{2L}{k+1}$, we approximate $\mathsf{P}_{\bar{\Lambda}+f,\beta_k}(v_{k-1})$. Similar to Algorithm 4, Algorithm 5 computes two types
269 of approximations: one is an ϵ_k-approximate proximal point $\tilde{\mathsf{P}}^{\epsilon_k}_{\bar{\Lambda}+f,\beta_k}(v_{k-1})$ as per its definition (4),
270 while the other is our bias-reduced optimum estimator from Algorithm 1. We note, however, that
271 unlike Algorithm 4 which approximates the x_k update, here we approximate v_k the "mirror descent"
272 update. This appears to be a substantial difference between the two algorithms.

273 Below we state the formal convergence guarantee for Algorithm 5; we defer its proof to Appendix G.

Algorithm 5: Stochastic composite accelerated gradient descent

---

Input: A problem of the form (5) with $\Lambda, f, \nabla\Lambda, \hat{\nabla}f$.
Parameters: Step size parameters $\gamma_k = \frac{2}{k}$ and $\beta_k = \frac{2L}{k+1}$, iteration number N, approximation
parameters $\{\epsilon_k, \delta_k, \sigma_k^2\}$ and initial $x_0 = v_0$ with distance bound $\|x_0 - x_\star\| \leq R$.

1 **for** $k = 1, 2, \cdots, N$ **do**
2     $y_k = (1 - \gamma_k)x_{k-1} + \gamma_k\mathsf{Proj}_{\mathcal{X}}(v_{k-1})$
3     $\bar{v}_k = \tilde{\mathsf{P}}^{\epsilon_k}_{\bar{\Lambda}_k+f,\beta_k}(v_{k-1})$ for $\bar{\Lambda}_k(v) := \Lambda(y_k) + \langle\nabla\Lambda(y_k), v - y_k\rangle$
4     $v_k = \mathsf{OPTEST}(\hat{\nabla}f, \psi_k, \beta_k, \delta_k, \sigma_k^2, \mathbb{B}_R(v_0) \cap \mathcal{X})$ for $\psi_k(z) = \frac{\beta_k}{2}\|z - v_{k-1}\|^2 + \bar{\Lambda}_k(z)$
5     $x_k = (1 - \gamma_k)x_{k-1} + \gamma_k\bar{v}_k$
6 **return** x_N

---

274 **Theorem 4.** *Given problem* (5) *with solution* $x_\star$, *a point* x_0 *such that* $\|x_0 - x_\star\| \leq R$ *and target*
275 *accuracy* $\epsilon > 0$, *Algorithm 5 with* $\epsilon_k = LR/2kN$, $\delta_k = R/16N$, $\sigma_k^2 = R^2/4N$, *and* $N =$

---

[5]The gradient sliding result holds under a relaxed Lipschitz assumption [see 29, eq. (1.2)]. It is straightforward to extend EPOCHSGD, and hence all of our results, to that assumption as well.

276 $\Theta(\sqrt{LR^2/\epsilon})$ *finds an approximate solution x satisfying* $\mathbb{E}\Psi(x) \leq \Psi(x_\star) + \epsilon$ *and has complexity*
277 $\mathcal{N}_{\nabla\Lambda} = O\left(\sqrt{\frac{LR^2}{\epsilon}}\right)$ *and* $\mathbb{E}\mathcal{N}_{\hat{\nabla}f} = O\left(\left(\frac{GR}{\epsilon}\right)^2 \log^2 \frac{GR}{\epsilon} + \sqrt{\frac{LR^2}{\epsilon}} \log\left(\frac{GR}{\epsilon}\right)\right).$

6 Efficient non-smooth private convex optimization

279 We conclude the paper with a potential application of our optimum estimator for differentially private
280 stochastic convex optimization (DP-SCO), where we have $f(x) = \mathbb{E}_{S\sim P}[\hat{f}(x;S)]$ for a convex
281 function $\hat{f} : \mathcal{X} \times \mathbb{S} \to \mathbb{R}$, $\mathcal{X} \subset \mathbb{R}^d$, and a set of n i.i.d. samples $s_i \sim P$. We wish to find $\hat{x} \in \mathcal{X}$ that
282 minimizes the excess loss $f(\hat{x}) - \min_{x\in\mathcal{X}} f(x)$ while preserving differential privacy.

283 **Definition 6.1** ([21]). *A randomized algorithm $\mathcal{A}$ is (α, β)-differentially private ((α, β)-DP) if, for*
284 *all datasets $\mathcal{S}, \mathcal{S}' \in \mathbb{S}^n$ that differ in a single data element and for every event $\mathcal{O}$ in the output space*
285 *of $\mathcal{A}$, we have $P[\mathcal{A}(\mathcal{S}) \in \mathcal{O}] \leq e^\alpha P[\mathcal{A}(\mathcal{S}') \in \mathcal{O}] + \beta$.*

286 DP-SCO has received increased attention over the past few years. Bassily et al. [5] developed
287 (inefficient) algorithms that attain the optimal excess loss $1/\sqrt{n} + \sqrt{d}/n\alpha$. Feldman et al. [22]
288 gave efficient algorithms for smooth functions that query $O(n)$ gradients. In the non-smooth setting,
289 however, their algorithms require $O(n^2)$ subgradients. Subsequently, [3] and [27] developed more
290 efficient algorithms for non-smooth functions which need $O(\min(n^2/\sqrt{d}, n^{5/4}d^{1/8}, n^{3/2}/d^{1/8}))$
291 subgradients which is $O(n^{11/8})$ for the high-dimensional setting $d = n$. Whether a linear gradient
292 complexity is achievable for DP-SCO in the non-smooth setting is still open. In this section, we
293 develop an efficient algorithm for non-smooth DP-SCO that queries $\widetilde{O}(n)$ subgradients conditional
294 on the existence of an optimum estimator with the following properties.

295 **Conjecture 1.** *Let $F = f + \psi$ be μ-strongly convex with minimizer $x_\star$ and f is G-Lipschitz. For*
296 *$\delta > 0$, there is an oracle $\mathcal{O}_\delta$ that returns $\hat{x}_\star = \mathcal{O}_\delta(F)$ such that $\|\mathbb{E}[\hat{x}_\star - x_\star]\|^2 \leq \delta^2$, $\|\hat{x}_\star - x_\star\|^2 \leq$*
297 *$C_1 G^2 \log(G/\mu\delta)/\mu^2$, and the expected number of gradient queries is $O(\log(G/\mu\delta))$.*

298 Comparing to our MLMC estimator (1) and Proposition 1, we note that the only place our current
299 estimator falls short is the probability 1 bound on $\|\hat{x}_\star - x_\star\|^2$, which for (1) holds only in expectation.
300 Indeed, for our estimator, $\|\hat{x}_\star - x_\star\|$ can be as large as $O(G/(\mu\delta))$, meaning that it is heavy-tailed.

301 Conditional on Conjecture 1, Algorithm 6 solves the DP-SCO problem with a near-linear amount of
302 gradient computations. The algorithm builds on the recent localization-based optimization methods
303 in [22] which iteratively solve regularized minimization problems. We average multiple draws of the
304 (hypothetical) bounded optimum estimator to solve the regularized problems, and apply private mean
305 estimation procedures to preserve privacy. We defer the proof of the following results Appendix H.

306 **Theorem 5.** *If Conjecture 1 holds, then for $\mathcal{X} \in \mathbb{B}_R(x_0)$, convex and G-Lipschitz $\hat{f}(x;s)$, Al-*
307 *gorithm 6 is (α, β)-DP, queries $\widetilde{O}(n)$ subgradients and has (hiding logarithmic factors in n)*
308 $$\mathbb{E}[f(x_k) - \min_{x\in\mathcal{X}} f(x)] \leq GR \cdot \widetilde{O}\left(\frac{1}{\sqrt{n}} + \frac{\sqrt{d\log^3(1/\beta)}}{n\alpha}\right).$$

309 Theorem 5 provides a strong motivation for proving Conjecture 1 by constructing a bounded optimum
310 estimator. In Appendix H.3, we discuss the challenges in making our MLMC estimator bounded, as
311 well as some directions to overcome them.

Algorithm 6: Differentially-private stochastic convex optimization via optimum estimation

Input: $(s_1, \ldots, s_n) \in \mathbb{S}^n$, domain $\mathcal{X} \subset \mathbb{B}_R(x_0)$, optimum estimator $\mathcal{O}$ satisfying Conjecture 1.

1 Set $k = \lceil \log n \rceil$, $B = 20(\log^2(n) + \log \frac{1}{\beta})$, $\bar{n} = \frac{n}{k}$ and $\eta = \frac{R}{G} \min\left\{\frac{1}{\sqrt{n}}, \frac{\alpha}{B\log(n)\sqrt{d\log(1/\beta)}}\right\}$

2 **for** $i = 1, 2, \cdots, k$ **do**

3 Let $\eta_i = 2^{-4i}\eta$, $f_i(x) = \frac{1}{\bar{n}} \sum_{j=1+(k-1)\bar{n}}^{k\bar{n}} \hat{f}(x; s_j)$, $\psi_i(x) = \|x - x_{i-1}\|^2/\eta_i\bar{n}$

4 Let $\tilde{x}_i = \frac{1}{\bar{n}} \sum_{j=1}^{\bar{n}} \mathcal{O}_{\delta_i}(F_i)$ with $F_i = f_i + \psi_i$, $\delta_i^2 = G^2\eta_i^2\bar{n}$

5 Set $x_i = \tilde{x}_i + \zeta_i$ where $\zeta_i \sim \mathsf{N}(0, \sigma_i^2 I_d)$ with $\sigma_i = 4B(\sqrt{C_1\log n} + 2)\eta_i\sqrt{\log(1/\beta)}/\alpha_i$

6 **return** x_k

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