Asynchronous Variance-reduced Block Schemes for Composite Nonconvex Stochastic Optimization: Block-specific Steplenghts and Adapted Batch-sizes

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Abstract

We consider the minimization of a sum of an expectation-valued coordinate-wise $L_i$-smooth nonconvex function and a nonsmooth block-separable convex regularizer. Prior schemes are characterized by the following shortcomings: (a) Steplengths require global knowledge of Lipschitz constants; (b) Batch-sizes of gradients are centrally updated and require knowledge of the global clock; (c) a.s. convergence guarantees are unavailable; (d) Rates are inferior compared to deterministic counterparts. Specifically, (a) and (b) require coordination across blocks and necessitate global information, leading to potentially larger constants in the rate and the oracle complexity bounds, impeding decentralized implementations, and resulting in relatively poor empirical behavior. We address these shortcomings by proposing an asynchronous variance-reduced algorithm, where in each iteration, a single block is randomly chosen to update its estimates by a proximal variable sample-size stochastic gradient scheme, while the remaining blocks are kept invariant. Notably, each block employs a steplength that is in accordance with its block-specific Lipschitz constant while block-specific batch-sizes are random variables updated at a rate that grows either at a geometric or polynomial rate with the (random) number of times that block is selected. We show that every limit point for almost every sample path is a stationary point and establish the ergodic non-asymptotic rate $O(1/K)$. Iteration and oracle complexity to obtain an $\epsilon$-stationary point are shown to be $O(1/\epsilon)$ and $O(1/\epsilon^2)$, respectively. Furthermore, under a $\mu$-proximal Polyak-Łojasiewicz (PL) condition with the batch size increasing at a geometric rate, we prove that the suboptimality diminishes at a geometric rate, the optimal deterministic rate while iteration and oracle complexity to obtain an $\epsilon$-optimal solution are proven to be $O((L_{\max}/\mu)\ln(1/\epsilon))$ and $O((L_{\text{ave}}/\mu)(1/\epsilon)^{1+c})$ with $c \geq 0$, respectively. In the single block setting, we obtain the optimal oracle complexity bound $O(1/\epsilon)$. In pursuit of less aggressive sampling rates, when the batch sizes increase at a polynomial rate of degree $v \geq 1$, suboptimality decays at a corresponding polynomial rate while the iteration and oracle complexity to obtain an $\epsilon$-optimal solution are provably $O(v(1/\epsilon)^{1/v})$ and $O(e^{v-2v+1}(1/\epsilon)^{1+1/v})$, respectively. Finally, preliminary numerics support our theoretical findings, displaying significant improvements over schemes where steplenghts are based on global Lipschitz constants.

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1 Introduction

In this paper, we consider the following composite stochastic programming:

\[
\min_{x \in \mathbb{R}^d} F(x) \triangleq \bar{f}(x) + \sum_{i=1}^{n} r_i(x_i),
\]

(1)

where the variable \( x_i \in \mathbb{R}^{d_i} \), \( x \) is partitioned into \( n \) blocks as \( x = (x_1, \ldots, x_n) \) with \( d \triangleq \sum_{i=1}^{n} d_i \), \( r_i(x_i) \) is a convex nonsmooth function, \( \bar{f}(x) \triangleq \mathbb{E}_{\xi}[f(x_1, \ldots, x_n, \xi)] \) is an expectation-valued smooth possibly nonconvex function with coordinate-wise \( L_i \)-Lipschitz continuous gradients, the random vector \( \xi : \Omega \rightarrow \mathbb{R}^m \) is defined on the probability space \( (\Omega, \mathcal{F}, \mathbb{P}) \), and \( f : \mathbb{R}^d \times \mathbb{R}^m \rightarrow \mathbb{R} \) is a scalar-valued function. Suppose that problem (1) has at least one solution. Let \( X^* \) and \( F^* \) denote the optimal solution set and the optimal function value, respectively. Nonsmoothness might be addressed through the proximal operator \([24, 33]\), defined as

\[
\text{prox}_{\alpha r}(x) \triangleq \arg\min_y \left( r(y) + \frac{1}{2\alpha} \|y - x\|^2 \right),
\]

(2)

where \( r(\cdot) \) is a closed and convex function, \( \alpha > 0 \), and the argmin is uniquely defined.

Prior research. In this paper, we propose an asynchronous proximal variance-reduced block scheme. We first review the literature on proximal, variance-reduced, and block methods.

(i) Proximal-gradient methods. Proximal-gradient methods and their accelerated variants are among the most important methods for solving composite convex problem \( f(x) + r(x) \) (also see forward-backward splitting (FBS) methods \([6,20,21]\)). While accelerated (or unaccelerated) schemes \([2]\) display non-asymptotic convergence rates in function value of \( O(1/k^2) \) (or \( O(1/k) \)), FBS methods \([4,37]\) display linear convergence when \( \nabla f(x) \) is strongly monotone. Nonconvex extensions have been studied in \([1,12,17]\), where the convergence to a stationary point is shown in \([1]\) while rate statements are provided under both the Kurdyka-Łojasiewicz (KL) property \([12]\) and the Polyak-Łojasiewicz (PL) condition \([17]\) (where a linear rate is proven).

(ii) Variance reduction schemes. A stochastic proximal gradient method was presented in \([34]\) for solving composite convex stochastic optimization, where the a.s. convergence and a mean-squared convergence rate \( O(1/k) \) were developed in strongly convex regimes, in sharp contrast with the linear rate of convergence in deterministic settings. Variance reduction schemes have gained increasing relevance in first-order methods for stochastic convex optimization \([13,16,35]\); in one particular class of schemes, the true gradient is replaced by the average of an increasing batch of sampled gradients, progressively reducing the variance of the sample-average. In strongly convex regimes, linear rates were shown for stochastic gradient methods \([16,35]\) and extragradient methods \([15]\), while for merely convex optimization problems, accelerated rates of \( O(1/k^2) \) and \( O(1/k) \) were proven for smooth \([13,16]\) and nonsmooth \([15]\) regimes, respectively. Mini-batch stochastic approximation methods were developed by \([14]\) for nonconvex stochastic composite optimization. Alternative schemes like SAGA and SVRG, applied to finite-sum machine learning problems \([28,30,39]\), rely on periodic use of the exact gradient, leading to recovery of deterministic convergence rates. For example, geometric rates were provided by \([30]\) for a proximal minibatch-SAGA and a proximal minibatch-SVRG algorithm in nonconvex regimes under the proximal PL inequality.

(iii) Block coordinate descent (BCD) schemes. BCD methods \([10]\) are widely used in machine learning and optimization, where variables are partitioned into manageable blocks and in each iteration, a single block is chosen to update while the remaining blocks remain fixed. Recently, in \([25]\), coordinate-friendly operators
were investigated that perform low-cost coordinate updates and it is shown that a variety of problems in machine learning can be efficiently resolved by such an update. The convergence properties of cyclic BCD methods has been extensively analyzed in [27,38,40]. Nesterov considered a randomized BCD method [22] and proved sublinear and linear convergence in terms of expected objective value for general convex and strongly convex cases, respectively. In [31], proximal (but unaccelerated) extensions were developed to contend with composite problems (also see [9,31,40,42,43]), while in [11], an accelerated, parallel, and proximal RBCD scheme was presented with a rate of $O(1/k^2)$. More recently, in [7], diverse block selection rules are considered and linear statements are provided for deterministic nonconvex problems under the PL condition.

**Motivation.** We consider a class of techniques that combine variance reduction and block-based schemes for solving the nonconvex nonsmooth stochastic programs, drawing inspiration from two seminal papers. Of these, the first by Xu and Yin [41] proposes a block stochastic gradient (BSG) method that cyclically updates blocks of variables. The second paper, by Dang and Lan [8], presents a stochastic block mirror-descent scheme reliant on randomly choosing and updating a single block by a mirror-descent stochastic approximation method. In [41] and [8], rates are provided in the convex setting while in nonconvex regimes, Dang and Lan [8] present non-asymptotic rates. Yet, there are several shortcomings that motivate the present research: (1) **Centralized batch sizes.** The schemes in [8,41] require a centrally specified batch-size across all blocks requiring global knowledge of the global clock, i.e. iteration $k$; (2) **Shorter steps.** Block-invariant steplengths utilize either $L$ or $L_{max}$ leading to significantly shorter steps and poorer performance (see Fig. 1); (3) **Asymptotic guarantees.** No a.s. convergence is available for BCD schemes for general nonconvex problems; (4) **Sub-optimal rate statements.** Optimal deterministic rates via variance-reduced schemes are unavailable but have been alluded to in convex regimes [41, Rem. 7]; Refinements via the PL condition remain open questions.

We address these gaps through a novel Algorithm 1 in Section 2 that combines a randomized BCD method with a proximal VSSG method, reliant on block-specific steplengths based on locally available $L_i$ without knowledge of the central Lipschitz constant $L$ or $L_{max}$, leading to larger steplengths and improved behavior (See Fig. 1 and Fig. 3), and on random block-specific batch-sizes (adapted to its block-selection history) without knowledge of the global clock, leading to lower informational coordination requirements. We make the following contributions supported by numerics in Section 5. Table 1 formalizes the distinctions in our scheme, while Table 2 compares our results with deterministic rates for nonconvex regimes.

(1) In Section 3, we prove that every limit point for almost every sample path is a stationary point under appropriately chosen batch sizes and show that the ergodic mean-squared error of the gradient mapping diminishes at a rate $O(1/K)$. We then establish that for any given $\epsilon > 0$, the iteration complexity (no. of proximal evaluations) and oracle complexity (no. of sampled gradients) to obtain an $\epsilon$-stationary point (measured by the mean-squared gap function) are $O(nL_{max}/\epsilon)$ and $O(n^2\nu^2L_{max}^2L_{min}^{-1}\epsilon^{-2})$ with uniform block selection, where $L_{max} \triangleq \max_i L_i$ and $L_{min} \triangleq \min_i L_i$. When the blocks are chosen as per a non-uniform distribution with probabilities $L_i \left(\sum_{j=1}^n L_j\right)^{-1}$ for any $i = 1, \cdots, n$, the iteration and oracle complexity are improved to $O(nL_{ave}/\epsilon)$ and $O(n^2\nu^2L_{ave}^2/\epsilon^2)$ with $L_{max}/L_{ave} = 2.5$. 

![Figure 1: Oracle Complexity of Algorithm 1 with identical and block-specific steps for (LASSO) with $L_{max}/L_{ave} = 2.5$.](image)
Furthermore, when the batch size increases at a polynomial rate of degree $v \geq 1$, we obtain a convergence rate $\mathbb{E}[F(x(k))] - F^* = O(k^{-v})$ and establish that the iteration and oracle complexity to obtain an $\epsilon$–optimal solution are $O(v(1/\epsilon)^{1/v})$ and $O\left(e^{v^{2v+1}(1/\epsilon)^{1+1/v}}\right)$, respectively.

### Table 1: List of literature on block-structured stochastic optimization

| App. | Metric | Asym/Rate/complexity | Comments |
|------|--------|----------------------|----------|
| C    | $\mathbb{E}[F(x(k)) - F^*]$ | $O(1/\sqrt{E})$ | (i) Centralized batch-size $N_k$; (ii) Central. step depend. on global $L$; (iii) No rate for nonconvex; (iv) sub-linear rates for SC |
| SC   | $\mathbb{E}[\|x(k) - x^*\|^2] = O(1/k)$ | | (i) Centralized batch-size $N_k$; (ii) Central stepsize depend on $L_{\max}$; (iii) Inferior constants dependent on $L_{\max}$; |
| C    | $\mathbb{E}[F(x_k) - F^*]$ | $O(1/\sqrt{K}), O(1/K)$ (SC) | (i) a.s. convergence (unavailable in [8, 41]) (ii) Block-specific stepsize depend on $L_i$; (iii) Optimal rates depend on $L_{ave}$ (Rather than $L, L_{\max}$ [8, 41]) (iv) Block-specific random. batch-sizes $\Rightarrow$ no central coordination of clocks (v) Optimal geometric rate under p-PL |

$L_{ave} \triangleq \sum_{i=1}^{n} L_i / n$. This represents a constant factor improvement in the rate from $L_{\max}$ (in [8]) to $L_{ave}$. Preliminary numerics reveal that block-specific steplengths lead to significant improvements in empirical behavior (see Table 5).

(II) In Section 4, we consider a class of nonconvex functions satisfying the **proximal PL condition** with parameter $\mu$ (see Assumption [3]) and prove that when the block-specific batch size is random and increases at a suitable geometric rate with the number of times the block is selected, the expectation-valued optimality gap $\mathbb{E}[F(x(k))] - F^*$ diminishes at a geometric rate. In addition, with uniform block selection, the iteration and oracle complexity to obtain an $\epsilon$–optimal solution are $O((nL_{\max}/\mu) \ln(1/\epsilon))$ and $O\left(nL_{ave}/\mu \ln(1/\epsilon) \left(1 + \frac{1}{n_{\min}^{-1}}\right) \frac{L_{\max}}{L_{\min}}\right)$ respectively when $n \geq 2$, where $\kappa_{\min} \triangleq L_{\min}/\mu$. While in the smooth regimes with a non-uniform block selection, the iteration and oracle complexity bounds are improved to $O\left(\frac{nL_{ave}}{\mu} \ln \left(\frac{1}{\epsilon}\right)\right)$ and $O\left(\frac{nL_{ave}}{\mu} (1/\epsilon) \left(1 + \frac{1}{n_{\min}^{-1}}\right) \frac{L_{\max}}{L_{\min}}\right)$, respectively. Specifically, when $n = 1$, the **optimal** oracle complexity $O\left(\frac{L}{\mu \epsilon}\right)$ is obtained. Notably, these rates match the deterministic versions in [7]. Furthermore, when the batch size increases at a polynomial rate of degree $v \geq 1$, we obtain a convergence rate $\mathbb{E}[F(x(k))] - F^* = O(k^{-v})$ and establish that the iteration and oracle complexity to obtain an $\epsilon$–optimal solution are $O(v(1/\epsilon)^{1/v})$ and $O\left(e^{v^{2v+1}(1/\epsilon)^{1+1/v}}\right)$, respectively.
a Poisson process with rate \( nL_{\mu} \epsilon \). Assume that the local Poisson clocks are independent, then the global clock ticks according to \( \varrho \) parameter selection rule that each block \( i \in \mathcal{N} \) at time \( n \) with probability one, there is a single block whose Poisson clock ticks according to a Poisson process with rate \( nL_{\mu} \epsilon \). The block selection rule (S.1.) accommodates the Poisson model employed by [3] as a special case, where Remark 1 In fact, Algorithm [7] might not require a global coordinator to coordinate the block selection. The block selection rule (S.1.) accommodates the Poisson model employed by [3] as a special case, where each block \( i \) is activated according to a local Poisson clock that ticks according to a Poisson process with rate \( q_i \). Suppose that there is a virtual global clock which ticks whenever any of the local Poisson clocks tick. Assume that the local Poisson clocks are independent, then the global clock ticks according to a Poisson process with rate \( \sum_{i=1}^{n} q_i \). Let \( Z_k \) denote the time of the \( k \)-th tick of the global clock. Since the local Poisson clocks are independent, with probability one, there is a single block whose Poisson clock ticks at time \( Z_k \) with probability \( \mathbb{P}(i_k = i) = \frac{q_i}{\sum_{i=1}^{n} q_i} \triangleq p_i \). In addition, note by (S.2.) and the algorithmic parameter selection rule that each block \( i \) maintains a local clock counting the number of its block updates, and that the update (3) does not necessitate knowledge of the global Lipschitz constant and the virtual global clock. As such, Algorithm [7] is an asynchronous scheme with limited coordination across blocks. Thus, in practical applications, Algorithm [7] might be helpful for the decentralized implementations compared to those designed in [8] and [41].

Table 2: Comparison with deterministic rates for nonconvex block methods

(a) Iteration complexity in smooth case \( (r(x) = 0) \)

| block selection rule | PL | general nonconvex |
|----------------------|----|------------------|
| unif. deterministic [7] | \( \mathcal{O}\left( \frac{nL_{\text{max}}}{\mu} \ln \left( \frac{F(x_1)-F^*}{\epsilon} \right) \right) \) | \( \mathcal{O}\left( \frac{nL_{\text{max}}(F(x(0))-F^*)}{\epsilon} \right) \) |
| stoch. (This work) | \( \mathcal{O}\left( \frac{nL_{\text{max}}}{\mu} \ln \left( \frac{F(x_1)-F^*}{\epsilon} + \frac{n\mu^2}{\epsilon} \right) \right) \) | \( \mathcal{O}\left( \frac{nL_{\text{max}}(F(x(0))-F^*)}{\epsilon} \right) \) |
| non-unif. deterministic [7] | \( \mathcal{O}\left( \frac{nL_{\text{max}}}{\mu} \ln \left( \frac{F(x_1)-F^*}{\epsilon} \right) \right) \) | \( \mathcal{O}\left( \frac{nL_{\text{max}}(F(x(0))-F^*)}{\epsilon} \right) \) |
| stoch. (This work) | \( \mathcal{O}\left( \frac{nL_{\text{max}}}{\mu} \ln \left( \frac{F(x_1)-F^*}{\epsilon} + \frac{n\mu^2}{\epsilon} \right) \right) \) | \( \mathcal{O}\left( \frac{nL_{\text{max}}(F(x(0))-F^*)}{\epsilon} \right) \) |

(b) Iteration complexity in nonsmooth case \( (r(x) \neq 0) \)

| block selection rule | PL | general nonconvex |
|----------------------|----|------------------|
| unif. deterministic [7] | \( \mathcal{O}\left( \frac{nL_{\text{max}}}{\mu} \ln \left( \frac{F(x_1)-F^*}{\epsilon} \right) \right) \) | \( \mathcal{O}\left( \frac{nL_{\text{max}}(F(x(0))-F^*)}{\epsilon} \right) \) |
| stoch. (This work) | \( \mathcal{O}\left( \frac{nL_{\text{max}}}{\mu} \ln \left( \frac{F(x_1)-F^*}{\epsilon} + \frac{n\mu^2}{\epsilon} \right) \right) \) | \( \mathcal{O}\left( \frac{nL_{\text{max}}(F(x(0))-F^*)}{\epsilon} \right) \) |

2 Asynchronous Block Proximal Stochastic Gradient Algorithm

We assume access to a proximal oracle (PO) that outputs prox_{\alpha\xi}(x_i) at any \( x_i \in \mathbb{R}^d_i \) for any \( \alpha > 0 \). Since the exact gradient \( \nabla f(x) \) is unavailable in a closed form, we assume there exists a stochastic first-order oracle (SFO) such that for every \( i \in \mathcal{N} \) and for any given \( x, \xi \), a sampled gradient \( \nabla_{x,i} f(x, \xi) \) is returned, which is an unbiased estimator of \( \nabla f(x) \). We aim to develop efficient algorithms for obtaining an \( \epsilon \)-optimal solution, where the efficiency is measured by the iteration complexity (no. of PO calls) and the oracle complexity (no. of SFO calls). Time is slotted at \( k = 0, 1, 2, \ldots \). Block \( i \) at time \( k \) holds a state \( x_i(k) \in \mathbb{R}^d_i \) that is an estimate for the corresponding coordinates of the optimal solution. We propose an asynchronous variance-reduced block stochastic gradient scheme (Algorithm 1) where at time instant \( k \), a block \( i \in \mathcal{N} \) is randomly chosen with probability \( p_i \) to compute the proxima update (3), where \( \alpha_i \) is the constant steplength and \( N_i(k) \) is the number of sampled gradients utilized at block \( i \), respectively. To be specific, the steplength \( \alpha_i \) depends on its block-specific Lipschitz constant \( L_i \) and the batch-size \( N_i(k) \) is a function of the random number of times block \( i \) is selected up to time \( k \). We will specify the selections of \( \alpha_i \) and \( N_i(k) \) upon the performance analysis in Sections 3 and 4.
Algorithm 1: Asynchronous variance-reduced block stochastic gradient algorithm

Let \( k := 0, x_i(0) \in \mathbb{R}^d \) and \( 0 < p_i < 1 \) for \( i = 1, \ldots, n \) such that \( \sum_{i=1}^{n} p_i = 1 \).

(S.1) Pick \( i_k = i \in \mathcal{N} \) with probability \( p_i \).

(S.2) If \( i_k = i \), then block \( i \) updates the state \( x_i(k+1) \) as follows:

\[
x_i(k+1) = \text{prox}_{\alpha_i r_i} \left( x_i(k) - \frac{\sum_{j=1}^{N_i(k)} \nabla x_i f(x(k), \xi_j(k))}{N_i(k)} \right),
\]

where \( \alpha_i > 0 \) is the steplength of block \( i \), \( N_i(k) \) is the number of sampled gradients utilized at block \( i \), and \( \{\xi_j(k)\}_{j=1}^{N_i(k)} \) are randomly generated from the probability space \((\Omega, \mathcal{F}, \mathbb{P})\). Otherwise, \( x_j(k+1) := x_j(k) \) if \( j \neq i_k \).

(S.3) If \( k > K \), stop and return \( \{x(k)\}_{k=0}^{K} \); Else, \( k := k + 1 \) and return to (S.1).

If the observation noise \( w_i(k) \) of the exact gradient is defined as

\[
w_i(k+1) = \frac{\sum_{j=1}^{N_i(k)} \nabla x_i f(x(k), \xi_j(k))}{N_i(k)} - \nabla x_i \bar{f}(x(k)),
\]

then (3) may be rewritten as

\[
x_i(k+1) = \text{prox}_{\alpha_i r_i} \left( x_i(k) - \alpha_i \left( \nabla x_i \bar{f}(x(k)) + w_i(k+1) \right) \right).
\]

By taking \( r_i(x_i) \) as an indicator function of a convex set \( X_i \), i.e., \( r_i(x_i) = 0 \) if \( x_i \in X_i \) and \( r_i(x_i) = +\infty \) otherwise, then the problem [1] reduces to the stochastic programming \( \min_{x_i \in X_i, i \in \mathcal{N}} \mathbb{E} [f(x_1, \ldots, x_n, \xi)] \). In this case, the update (3) reduces to the variable sample-size projected stochastic gradient method:

\[
x_i(k+1) = \mathbf{P}_{X_i} \left( x_i(k) - \alpha_i \left( \nabla x_i \bar{f}(x(k)) + w_i(k+1) \right) \right),
\]

where \( \mathbf{P}_{X_i}(x_i) \) denotes the projection of \( x_i \) onto the set \( X_i \). This can be thought as a generalization of the schemes proposed in [15, 35] for solving constrained stochastic convex program.

We impose the following conditions on the objective functions and observation noises.

**Assumption 1** (i) \( r_i \) is a proper lower semicontinuous and convex function with effective domain \( \mathcal{R}_i \) required to be compact. (ii) There exists a constant \( L_i > 0 \) such that for any \( x_i' \in \mathcal{R}_i \) and any \( x \in \prod_{j=1}^{n} \mathcal{R}_j \),

\[
\| \nabla x_i \bar{f}(x) - \nabla x_i \bar{f}(x_1, \ldots, x_{i-1}, x_i', x_{i+1}, \ldots, x_n) \| \leq L_i \| x_i - x_i' \|.
\]

**Assumption 2** (i) There exists \( \nu > 0 \) such that for any \( i \in \mathcal{N} \) and all \( k \geq 1 \), \( \mathbb{E}[\| w_i(k+1) \|^2 | \mathcal{F}_k ] \leq \frac{\nu^2}{N_i(k)} \) a.s.; (ii) \( i_k \) is independent of \( \mathcal{F}_k \) for all \( k \geq 1 \).

**Remark 2** Assumption [ii] requires the effective domain of each block’s nonsmooth convex regularizer to be a compact set. Suppose \( r_i(x_i) \) is set as an indicator function of a convex set \( X_i \), then it is required that \( X_i \) is a compact set. Such a condition guarantees that the iterates \( \{x(k)\} \) produced by Algorithm 1 are uniformly bounded. Assumption [ii] necessitates the gradient function of the nonconvex smooth term \( \bar{f}(x) \) to be block-wise Lipschitz continuous. Assumption [iii] can be satisfied if the sampled gradients are independently generated and \( \mathbb{E}[\| \nabla x_i f(x, \xi) - \nabla x_i \bar{f}(x) \|^2 | \mathcal{F}_k ] \) (the variance of the stochastic gradient noise) is uniformly bounded.
3 Convergence to Stationary Points

In this section, we will prove the almost sure convergence of iterates to a stationary point and establish the non-asymptotic rate of Algorithm 1.

3.1 Preliminary Lemmas

Before presenting the convergence results, we recall a preliminary result from [30, Lemma 2].

**Lemma 1** Suppose \( y \doteq \text{prox}_{\alpha r}(x - \alpha g) \) for some \( g \in \mathbb{R}^d \). Then for any \( z \in \mathbb{R}^d \),

\[
\bar{f}(y) + r(y) \leq \bar{f}(z) + r(z) + (y - z)^T (\nabla \bar{f}(x) - g) + \left( \frac{L}{2} - \frac{1}{2\alpha} \right) \|y - x\|^2 \\
+ \left( \frac{L}{2} + \frac{1}{2\alpha} \right) \|z - x\|^2 - \frac{1}{2\alpha} \|y - z\|^2.
\]

We now give a simple relation on the conditional expectation of the function value in the following lemma, for which the proof can be found in Appendix A. This is an important preliminary result because the convergence results to be presented in this section are essentially obtained through a recursive application of this basic lemma.

**Lemma 2** Let \( \{x(k)\} \) be generated by Algorithm 1 where \( 0 < \alpha_i \leq \frac{1}{L_i} \) for any \( i = 1, \cdots, n \). Suppose that Assumptions 1 and 2(ii) hold. Define

\[
\bar{x}_i(k+1) \doteq \text{prox}_{\alpha_i r_i}(x_i(k) - \alpha_i \nabla_x \bar{f}(x_i(k)))
\]

as the update if the true gradient is used. Then the following holds a.s.:

\[
\mathbb{E}[F(x(k+1)) | \mathcal{F}_k] \leq F(x(k)) - \sum_{i=1}^{n} p_i \left( \frac{1}{2\alpha_i} - L_i \right) \|\bar{x}_i(k+1) - x_i(k)\|^2 + \frac{1}{2} \mathbb{E}[\alpha_{ik} \|w_{ik}(k+1)\|^2 | \mathcal{F}_k], \quad \forall k \geq 1.
\]

Throughout the paper, all inequalities and equalities between random variables are assumed to hold a.s., but we often omit to write “a.s.” for simplicity.

3.2 Asymptotic Convergence

From Assumption 1(i) it is seen that the sequence of estimates \( \{x(k)\} \) produced Algorithm 1 is bounded. We now establish the a.s. convergence by showing that for every limit point of almost every sample path \( \{x(k)\} \) is a stationary point of problem (1).

**Theorem 1** Let \( \{x(k)\} \) be generated by Algorithm 1 Suppose Assumptions 1 and 2 hold, and that for any \( i = 1, \cdots, n \), \( 0 < \alpha_i < \frac{1}{2L_i} \) and \( \sum_{k=0}^{\infty} \frac{1}{N_i(k)} < \infty \) a.s.. Then every cluster point of almost every sample path \( \{x(k)\} \) is a stationary point.
Proof. Note by Assumption 2(i) that
\[
E \left[ \alpha_{ik} \|w_{ik}(k+1)\|^2 | F_k \right] \leq \sum_{i=1}^{n} E \left[ \alpha_i \|w_i(k+1)\|^2 | F_k \right] \leq \sum_{i=1}^{n} \frac{\alpha_i \nu_i^2}{N_i(k)}.
\]

Note that \( \frac{1}{2\alpha_i} - L_i > 0 \) by \( 0 < \alpha_i < \frac{1}{2\alpha_i} \). Then by recalling that \( \sum_{k=0}^{\infty} \frac{1}{N_i(k)} < \infty \text{ a.s.} \), we may then apply [32 Thm. 1] to inequality (6) and conclude that \( F(x(k)) \) converges almost surely and
\[
\sum_{k=1}^{\infty} \sum_{i=1}^{n} p_i \left( \frac{1}{2\alpha_i} - L_i \right) \| \bar{x}_i(k+1) - x_i(k) \|^2 < \infty \text{ a.s.}
\]
This implies that
\[
\sum_{k=0}^{\infty} \| \bar{x}(k+1) - x(k) \|^2 < \infty, \text{ a.s.}
\]

(7)

Then for almost every sample path, we have that \( \| \bar{x}(k+1) - x(k) \| \to 0 \) as \( k \to \infty \). Let \( \hat{x} \) be a cluster point of any such sequence \( \{ x(k) \} \). Then there exists a subsequence \( \{ x(k_t) \} \) such that \( \lim_{t \to \infty} x(k_t) = \hat{x} \) and hence \( \lim_{t \to \infty} \bar{x}(k_t + 1) = \hat{x} \) by (7). For any \( i = 1, \ldots, n \), using the definitions (2) and (5), we obtain that
\[
\bar{x}_i(k + 1) = \arg\min_{y \in \mathbb{R}^{d_i}} \left[ \nabla_{x_i} \tilde{f}(x(k))^T (y - x_i(k)) + \frac{1}{2\alpha_i} \| y - x_i(k) \|^2 + r_i(y) \right].
\]

(8)

Then by using the first-order optimality condition, we obtain for all \( t \):
\[
- \frac{1}{\alpha_i} \left( \bar{x}_i(k_t + 1) - x_i(k_t) \right) \in \nabla_{x_i} \tilde{f}(x(k_t)) + \partial r_i(\bar{x}_i(k_t + 1)).
\]

(9)

By passing to the limit in (9), using \( \| \bar{x}(k_t + 1) - x(k_t) \| \to 0 \text{ a.s.} \), by \( \lim_{t \to \infty} x(k_t) = \lim_{t \to \infty} \bar{x}(k_t + 1) = \hat{x} \), and the continuity of \( \nabla_{x_i} \tilde{f}(\cdot) \) and the closedness of \( \partial r_i \), we obtain that \( 0 \in \nabla_{x_i} \tilde{f}(\hat{x}) + \partial r_i(\hat{x}_i) \) for any \( i = 1, \ldots, n \). Thus, \( 0 \in \nabla \tilde{f}(\hat{x}) + \partial r(\hat{x}) \), implying that \( \hat{x} \) is a stationary point of (1). \( \square \)

Remark 3 (i) For any \( i \in \mathcal{N} \), define \( \Gamma_i(k) \triangleq \sum_{p=0}^{k-1} I_{[y_p=i]} \) for any \( k \geq 1 \), where \( I_{[a=b]} = 1 \) if \( a = b \), and \( I_{[a=b]} = 0 \) otherwise. Thus, \( \Gamma_i(k) \) is adapted to \( F_k \), and \( \sum_{k=0}^{\infty} \frac{1}{N_i(k)} < \infty \text{ a.s.} \) holds by setting \( N_i(k) \triangleq \left\lfloor \Gamma_i(k) + 1 \right\rfloor^{1+\delta} \) for some \( \delta > 0 \). This follows by [18 Lemma 7] that for almost every \( \omega \in \Omega \), there exists a sufficiently large \( \bar{k} \) possibly contingent on the sample path \( \omega \) such that for any \( k \geq \bar{k} \), \( \Gamma_i(k) \geq \frac{kn}{2}, i = 1, \ldots, n \).

(ii) If \( \tilde{f}(x) \) is convex, then Theorem 7 implies that \( F(x(k)) \) converges almost surely to the optimal function value \( F^* \), and every cluster point of almost every sample path \( \{ x(k) \} \) is a global minimum to the problem (1).

3.3 Non-asymptotic Rate

Recall that for convex optimization, a frequently-used metric is the sub-optimality metric \( F(x) - F^* \) or the distance to the optimal solution set \( d(x, X^*) \). However, in nonconvex optimization, the iterates might converge to stationary points which are not necessarily global minima, and as a consequence, the standard metric cannot be applied. Thus, one crucial problem in analyzing Algorithm 1 for nonconvex optimization lies in the selection of the convergence criterion. In smooth regimes, it is typical to use \( \| \nabla \tilde{f}(x) \| \) while
in nonsmooth settings, an appropriate alternative is the proximal gradient mapping \([30]\): \(G_\alpha(x) = \frac{1}{\alpha}(x - \text{prox}_{\alpha f}(x - \alpha \nabla f(x)))\). Then \(x^0 \in \mathbb{R}^d\) satisfying \(G_\alpha(x^0) = 0\) is a stationary point of \((1)\). We now analyze the rate of convergence of Algorithm \([1]\) and establish iteration and oracle complexity bounds to obtain an \(\epsilon\)-stationary point, by using the following metric to measure stationarity.

\[
G_{i,\alpha_i}(x) = \frac{x_i - \text{prox}_{\alpha_i f_i}(x_i - \alpha_i \nabla f_i(x))}{\alpha_i}, \quad i \in \mathcal{N}, \quad G_\alpha(x) \triangleq (G_{i,\alpha_i}(x))_{i=1}^n. \tag{10}
\]

It is seen that any zero of \(G_\alpha(x)\) is a stationary point of \((1)\). Next, we establish a result for Algorithm \([1]\) when the block is chosen according to a uniform distribution.

**Theorem 2** Let \(\{x(k)\}\) be generated by Algorithm \([1]\). Suppose Assumptions \([1]\) and \([2]\) hold, \(p_i = \frac{1}{n}\) and \(\alpha_i = \frac{1}{4L_i}\) for any \(i = 1, \ldots, n\). Let \(x_{\alpha,K}\) be chosen from \(\{x(k)\}_{k=0}^K\) as per a uniform distribution. We have the following bound on the mean-squared error:

\[
\mathbb{E} [\|G_\alpha(x_{\alpha,K})\|^2] \leq \frac{16nL_{\max}(\mathbb{E}[F(x(0))] - F^*)}{K + 1} + \frac{2\nu^2L_{\max}^2}{K + 1} \sum_{i=1}^n \sum_{k=0}^K L_i^{-1}\mathbb{E}[N_i(k)^{-1}]. \tag{11}
\]

Then for any given \(\epsilon > 0\), by setting \(K = \tilde{K}_1(\epsilon) \triangleq \frac{32nL_{\max}(\mathbb{E}[F(x(0))] - F^*)}{\epsilon} \) and \(N_i(k) \equiv \tilde{N}_1(\epsilon) \triangleq \frac{4n\nu^2L_{\max}^2}{\epsilon L_{\min}}\), the iteration and oracle complexity to obtain an \(\epsilon\)-stationary point such that \(\mathbb{E} [\|G_\alpha(x_{\alpha,K})\|^2] \leq \epsilon\) are \(\tilde{K}_1(\epsilon)\) and \(\tilde{K}_1(\epsilon)\tilde{N}_1(\epsilon)\), respectively.

**Proof.** Note by \(\mathbb{P}(i_k = i) = p_i\) and Assumption \([2]\), we obtain that

\[
\mathbb{E} [\alpha_i \|w_{i}(k + 1)\|^2] = \sum_{i=1}^n p_i \alpha_i \mathbb{E} [\|w_i(k + 1)\|^2] \leq \sum_{i=1}^n \alpha_i p_i \nu^2 \mathbb{E} [N_i(k)^{-1}].
\]

By taking unconditional expectations of \((6)\) and rearranging the terms, there holds

\[
\mathbb{E} \left[ \sum_{i=1}^n p_i \left( \frac{1}{2\alpha_i} - L_i \right) \|\bar{x}_i(k + 1) - x_i(k)\|^2 \right] \leq \mathbb{E} [F(x(k))] - \mathbb{E} [F(x(k + 1))] + \frac{1}{2} \sum_{i=1}^n \alpha_i p_i \nu^2 \mathbb{E} [N_i(k)^{-1}]. \tag{12}
\]

Thus, by summing up \((12)\) from \(k = 0\) to \(K\), we have that

\[
\sum_{k=0}^K \mathbb{E} \left[ \sum_{i=1}^n p_i \left( \frac{1}{2\alpha_i} - L_i \right) \|\bar{x}_i(k + 1) - x_i(k)\|^2 \right] \leq \mathbb{E} [F(x(0))] - \mathbb{E} [F(x(K + 1))] + \frac{\nu^2}{2} \sum_{k=0}^K \sum_{i=1}^n \alpha_i p_i \mathbb{E} [N_i(k)^{-1}]. \tag{13}
\]

By the definitions \((5)\) and \((10)\), \(\alpha_i = \frac{1}{4L_i}, p_i = \frac{1}{n}\), we have that

\[
p_i \left( \frac{1}{2\alpha_i} - L_i \right) \|\bar{x}_i(k + 1) - x_i(k)\|^2 = p_i \alpha_i \left( \frac{1}{2} - \alpha_i L_i \right) \left( \|\bar{x}_i(k + 1) - x_i(k)\| / \alpha_i \right)^2 \geq \frac{1}{16nL_i} ||G_{i,\alpha_i}(x(k))||^2.
\]
This combined with (13), \( F(x_{K+1}) \geq F^* \), and \( p_i = \frac{1}{n} \) implies that
\[
\frac{1}{16nL_{\max}} \sum_{k=0}^{K} \mathbb{E} \left[ \|G_{\alpha}(x(k))\|^2 \right] \leq \mathbb{E}[F(x(0))] - F^* + \frac{\nu^2}{8n} \sum_{i=1}^{n} \frac{1}{L_i} \sum_{k=0}^{K} \mathbb{E} \left[ N_i(k)^{-1} \right].
\]

Therefore, by multiplying both sides of the above equation by \( \frac{16nL_{\max}}{K+1} \), by using
\[
\mathbb{E} \left[ \|G_{\alpha}(x,\alpha,K)\|^2 \right] = \frac{1}{K+1} \sum_{k=0}^{K} \mathbb{E} \left[ \|G_{\alpha}(x(k))\|^2 \right]
\]
we obtain (11). Since \( N_i(k) = \bar{N}_i(\epsilon) \), we have \( \frac{1}{n+1} \sum_{k=0}^{K} \mathbb{E} \left[ N_i(k)^{-1} \right] = \frac{1}{\bar{N}_i(\epsilon)} \). This combined with (11) produces
\[
\mathbb{E} \left[ \|G_{\alpha}(x,\alpha,K)\|^2 \right] \leq \frac{16nL_{\max} (\mathbb{E}[F(x(0))] - F^*)}{K_1(\epsilon) + 1} + \frac{2\nu^2 L_{\max}}{N_1(\epsilon)L_{\min}} \leq \epsilon.
\]

Since a single block is chosen to update in each iteration, the total number of samples used to the update of \( x(K) \) is \( \sum_{k=0}^{K-1} \sum_{i=1}^{n} N_i(k)I_{i=1} = \bar{K}_1(\epsilon)\bar{N}_1(\epsilon) \). Then the number of PO and SFO calls required to ensure that \( \mathbb{E} \left[ \|G_{\alpha}(x,\alpha,K)\|^2 \right] \leq \epsilon \) are \( \bar{K}_1(\epsilon) \) and \( \bar{K}_1(\epsilon)\bar{N}_1(\epsilon) \), respectively. \( \square \)

We now analyze the rate of convergence of Algorithm 1 with the active block chosen via a non-uniform distribution constructed using block-specific Lipschitz constants.

**Theorem 3** Let \( \{x(k)\} \) be generated by Algorithm 1 where \( p_i = \frac{L_i}{\sum_{i=1}^{n} L_i} \) and \( \alpha_i = \frac{1}{4L_i} \) for \( i = 1, \ldots, n \). Suppose Assumptions 1 and 2 hold. Then
\[
\mathbb{E} \left[ \|G_{\alpha}(x,\alpha,K)\|^2 \right] \leq \frac{16nL_{\text{ave}} (\mathbb{E}[F(x(0))] - F^*)}{K+1} + \frac{2\nu^2}{K+1} \sum_{i=1}^{n} \sum_{k=0}^{K} \mathbb{E} \left[ N_i(k)^{-1} \right].
\]

Thus, for any given \( \epsilon > 0 \), by setting \( K = K_2(\epsilon) = \frac{32nL_{\text{ave}} (\mathbb{E}[F(x(0))] - F^*)}{\epsilon} \) and \( N_i(k) \equiv \bar{N}_2(\epsilon) = \frac{4\nu^2}{\epsilon} \), the iteration and oracle complexity to obtain an \( \epsilon \)-stationary point such that \( \mathbb{E}[\|G_{\alpha}(x,\alpha,K)\|^2] \leq \epsilon \) are \( \bar{K}_2(\epsilon) \) and \( \bar{K}_2(\epsilon)\bar{N}_2(\epsilon) \), respectively.

**Proof.** By definitions (5), \( \alpha_i = \frac{1}{4L_i} \), \( p_i = \frac{L_i}{nL_{\text{ave}}} \), we have that
\[
p_i \left( \frac{1}{2\alpha_i} - L_i \right) \|\bar{x}(k+1) - x_i(k)\|^2 = p_i\alpha_i \left( \frac{1}{2} - \alpha_i L_i \right) \left( \frac{\|\bar{x}(k+1) - x_i(k)\|}{\alpha_i} \right)^2 = \frac{1}{16nL_{\text{ave}}} \|G_{i,\alpha_i}(x(k))\|^2.
\]

Then using (13), the definition of \( G_{\alpha}(x) \) and \( \alpha_i p_i = \frac{1}{4L_{\text{ave}}} \), we have the following:
\[
\sum_{k=0}^{K} \frac{1}{16nL_{\text{ave}}} \mathbb{E} \left[ \|G_{\alpha}(x(k))\|^2 \right] \leq \mathbb{E}[F(x(0)) - F(x(K+1))] + \frac{\nu^2}{8nL_{\text{ave}}} \sum_{i=1}^{n} \sum_{k=0}^{K} \mathbb{E} \left[ N_i(k)^{-1} \right].
\]

Then by using \( F(x_{K+1}) \geq F^* \) and multiplying both sides of the above equation with \( \frac{16nL_{\text{ave}}}{K+1} \), we obtain (15). The rest of the proof is the same as that of Theorem 2. \( \square \)
Remark 4 We observe the following regarding Theorems 2–3:

(i) Note that \( G_\alpha(x) = \nabla f(x) \) when \( r(x) \equiv 0 \). Suppose \( \sum_{k=0}^K \mathbb{E} [N_i(k)^{-1}] \) is bounded for any \( i = 1, \ldots, n \). Then we attain the non-asymptotic rate \( \mathbb{E} [\| \nabla f(x_{\alpha,K}) \|^2] = O(1/K) \), the best known rate for first-order methods for nonconvex programs [13].

(ii) Note that iteration and oracle complexity bounds of Algorithm 1 with uniform block selection are respectively \( O(nL_{\max}/\epsilon) \) and \( O(n^2\nu^2L_{\max}^2/(L_{\min}\epsilon^2)) \), while if the blocks are selected with a likelihood proportional to the block-specific Lipschitz constant, the bounds are respectively reduced to \( O(nL_{\text{ave}}/\epsilon) \) and \( O(n^2\nu^2L_{\text{ave}}/(\epsilon^2)) \).

(iii) The iteration complexity (no. partial proximal evaluations) is \( O(n/\epsilon) \). Since the variable is partitioned into \( n \) blocks, the iteration complexity (no. full proximal evaluations) is \( O(1/\epsilon) \), which is optimal for deterministic gradient descent methods.

4 Global Linear Convergence under PL-Inequality

In this section, we will prove the global linear convergence of iterates and derive complexity bounds when the proposed scheme is applied to a class of nonsmooth nonconvex composite functions satisfying the proximal PL inequality. The PL inequality \( \| \nabla f(x) \|^2 \geq 2\mu(f(x) - \min_x f(x)) \) requires the gradient norm to grow faster than a quadratic function when moving away from the optimal value. It was first proposed in [26] that the global linear convergence of the gradient descent method can be obtained under the PL condition. Its generalization, called the proximal PL inequality, was proposed in [17] for the composite function. It has been shown in [17] that several important classes of functions satisfy this proximal PL condition, e.g., (i) \( f \) is strongly convex; (ii) \( f \) has the form \( f(x) = h(Ax) \) for a strongly convex function \( h \) and a matrix \( A \) while \( r \) is an indicator function for a polyhedral set; and (iii) \( F \) is convex and satisfies the quadratic growth property. We impose the following condition on the problem [1].

Assumption 3 (\( \mu \)-PL) There is a \( \mu > 0 \) satisfying \( \frac{1}{2}D_r(x, L_{\max}) \geq \mu(F(x) - F^*) \), where

\[
D_r(x, L) \triangleq -2L \min_y \left[ \nabla \tilde{f}(x)^T (y - x) + \frac{L}{2} \| y - x \|^2 + r(y) - r(x) \right].
\]

4.1 Rate Analysis

We first present a preliminary lemma, based on which we show in Theorem 4 and Proposition 1 that \( F(x(k)) \) converges in mean to the optimal value \( F^* \) at a geometric rate and a polynomial rate when the number of the sampled gradients increases at a geometric rate and a polynomial rate, respectively. The proof of the Lemma can be found in Appendix B.

Lemma 3 Let \( \{x(k)\} \) be generated by Algorithm 1. Suppose Assumptions 2 and 3 hold. Let \( \beta \in (\frac{1}{2}, 1) \) and \( 0 < \alpha_i \leq \frac{2\beta-1}{L_i(1+\beta)} \). Define \( \alpha_{\min} = \min_{i \in \mathcal{N}} \alpha_i \). Then for all \( k \geq 1 \),

\[
\mathbb{E} [F(x(k+1)) - F^*] \leq (1 - \alpha_{\min}(1 - \beta)\mu_{N})\mathbb{E} [F(x(k)) - F^*] + \frac{\mu^2}{2} \sum_{i=1}^n \alpha_i p_i \mathbb{E} [N_i(k)^{-1}] . \tag{16}
\]

We now discuss the optimal selection of parameters \( \alpha_i \) and \( \beta \). Define \( \rho(\alpha, \beta) \triangleq 1 - \alpha(1 - \beta)\mu_{\min} \). Then by \( 0 < \alpha_i \leq \frac{2\beta-1}{L_i(1+\beta)} \) and \( \beta \in (\frac{1}{2}, 1) \), we have that \( 0 < \alpha_i(1 - \beta) \leq \frac{(2\beta-1)(1 - \beta)}{L_i(1+\beta)} \). We set \( \beta \) to be
the maximizer of \( \frac{(2\beta-1)(1-\beta)}{1+\beta} \), given by \( \beta^* = \sqrt{3} - 1 \). Then by setting \( \alpha_i^* = \frac{2\beta^*-1}{L_i(1+\beta^*)} = \frac{2-\sqrt{3}}{L_i} \), we get
\[ 0 < \rho(\alpha_{\min}, \beta^*) = 1 - \frac{(2-\sqrt{3})^2\mu\rho_{\min}}{L_{\max}} < 1. \]
By setting \( \alpha_i = \alpha_i^* \) and \( \beta = \beta^* \) in Algorithm 1, we obtain the geometric rate under the proximal PL condition with the geometrically increasing sample-sizes.

**Theorem 4 (Geometric rate of convergence)** Let \( \{x(k)\} \) be generated by Alg. 1 where \( \alpha_i = \frac{2-\sqrt{3}}{L_i} \) and \( N_i(k) = [(1- q_i)^{-1} - \Gamma_i(k)] \) for some \( q_i \in (0, 1) \) with \( \Gamma_i(k) := \sum_{p=0}^{k-1} I[p=q_i] \). Suppose \( p_i = \frac{1}{n} \forall i \in N \), Assumptions 2 and 3 hold. Let \( q_{\min} := \min_{i \in N} q_i \) and \( \rho^* := 1 - \frac{(2-\sqrt{3})^2\mu}{nL_{\max}} \).

(i) If \( q_{\min} \neq \frac{(2-\sqrt{3})^2\mu}{L_{\max}} \), then for all \( k \geq 0 \):
\[
\mathbb{E}[F(x(k)) - F^*] \leq \left( 1 - \frac{1}{n} \min \left\{ q_{\min}, \frac{(2-\sqrt{3})^2\mu}{L_{\max}} \right\} \right)^k \times \left( \mathbb{E}[F(x(0)) - F^*] + \frac{\nu^2 \sum_{i=1}^{n} \alpha_i}{2q_{\min} - (2-\sqrt{3})^2\mu/L_{\max}} \right). \tag{17}
\]

(ii) If \( q_{\min} = \frac{(2-\sqrt{3})^2\mu}{L_{\max}} \), then the following holds for any \( \tilde{\rho} \in (\rho^*, 1) \) and all \( k \geq 0 \):
\[
\mathbb{E}[F(x(k)) - F^*] \leq \rho^k \left( \mathbb{E}[F(x(0)) - F^*] + \frac{\nu^2 \sum_{i=1}^{n} \alpha_i}{2n\rho^* \ln((\tilde{\rho}/\rho^*)^\rho)} \right). \tag{18}
\]

**Proof.** Since \( \alpha_i = \alpha_i^* \) and \( p_i = \frac{1}{n} \), by setting \( \beta = \beta^* \), we have that \( \rho(\alpha_{\min}, \beta^*) = 1 - \frac{(2-\sqrt{3})^2\mu\rho_{\min}}{L_{\max}} = \rho^* \). Then by the definition \( \Gamma_i(k) := \sum_{p=0}^{k-1} I[p=q_i] \), we know that
\[
\mathbb{P}(\Gamma_i(k) = m) = \binom{k}{m} p_i^m (1-p_i)^{k-m}. \tag{19}
\]

Then using \( p_i = 1/n \), we have the following for any \( k \geq 1 \) and \( i \in N \):
\[
\mathbb{E}\left[ N_i(k)^{-1} \right] \leq \mathbb{E}\left[ (1 - q_i)\Gamma_i(k) \right] = \sum_{m=0}^{k} (1 - q_i)^m \mathbb{P}(\Gamma_i(k) = m) \tag{20}
\]
\[
= \sum_{m=0}^{k} \binom{k}{m} (p_i(1 - q_i))^m (1-p_i)^{k-m} = (p_i(1-q_i) + 1-p_i)^k = (1-p_iq_i)^k.
\]

By combining (20) with (16) and by recalling that \( q_{\min} = \min_{i \in N} q_i \) we obtain that
\[
v_{k+1} \leq \rho^* v_k + \sum_{i=1}^{n} \alpha_i \nu^2 \left( 1 - \frac{q_{\min}}{n} \right)^k,
\]
where \( v_k := \mathbb{E}[F(x(k)) - F^*] \). Then by defining \( q^* := 1 - q_{\min}/n \), we have
\[
v_{k+1} \leq (\rho^*)^{k+1} v_0 + \sum_{m=0}^{k} (\rho^*)^m (q^*)^{k-m} \sum_{i=1}^{n} \alpha_i \nu^2 \left( 2n \right)^k. \tag{21}
\]
Proof. 
By (ii).
\[ \rho \] and hence (17) holds by definitions of \( \rho^* \).
For any
\[ \text{Lemma 4 (Eq. (17) and Lemma 4 \cite{19})} \]
investigate the rate of convergence of Algorithm 1 with polynomially increasing sample sizes based on a convergence rate of Algorithm 1 when the batch-size is increased at a slower polynomial rate. Next, we will impossible to increase the batch-size too fast at a geometric rate. As such, we also try to explore the
\[ \Pi \]
Consider Algorithm 1, where
\[ (ii) \text{ By (i)} \]
\[ \text{Similarly, for the case where } 0 < q_{\text{min}} < (2-\sqrt{3})^2 \mu \text{, we have that } q^* > \rho^* \text{ and } \sum_{m=0}^{k} (\rho^*)^m (q^*)^{k-m} \leq (q^*)^k \frac{1}{1-q^*/\rho^*} = (\rho^*)^{k+1} \frac{1}{\rho^* - q^*}. \]

Similarly, for the case where \( 0 < q_{\text{min}} < (2-\sqrt{3})^2 \mu \), we have that \( q^* > \rho^* \) and \( \sum_{m=0}^{k} (\rho^*)^m (q^*)^{k-m} \leq (q^*)^k \frac{1}{\rho^* - q^*} \). As a result, by using (21) we obtain that
\[ \mathbb{E}[F(x_{k+1}) - F^*] \leq \max \{q^*, \rho^*\}^{k+1} \left( \mathbb{E}[F(x(0)) - F^*] + \sum_{i=1}^{n} \alpha_i \nu_i^2 \right), \]
and hence (17) holds by definitions of \( \rho^* \) and \( q^* \).

(ii) By \( q_{\text{min}} = (2-\sqrt{3})^2 \mu \), we have that \( q^* = \rho^* \). Then by (21), we have that \( v_k \leq (\rho^*)^k v_0 + (\rho^*)^k \sum_{i=1}^{n} \alpha_i \nu_i^2 \).

By using \( k (\rho^*)^k \leq \hat{\rho}^k / \ln ((\hat{\rho}/\rho^* \nu_i)^) \) (see \cite[Lemma 2]{19}) and \( v_0 \triangleq \mathbb{E}[F(x(0)) - F^*] \), we obtain the result (ii).

In some settings, the evaluation of sampled gradients might be costly. Hence it is unreasonable or impossible to increase the batch-size too fast at a geometric rate. As such, we also try to explore the convergence rate of Algorithm 1 when the batch-size is increased at a slower polynomial rate. Next, we will investigate the rate of convergence of Algorithm 1 with polynomially increasing sample sizes based on a preliminary result from \cite{19}.

**Lemma 4 (Eq. (17) and Lemma 4 \cite{19})** For any \( q \in (0, 1) \) and \( v > 0 \), the following hold:
\[ \sum_{m=1}^{k+1} q^{k+1-m} m^{-v} \leq q^{k+1} e^{2v} \frac{q^{1-v} - 1}{1 - q} + \frac{2(k+1)^{1-v}}{q \ln(1/q)} \quad \forall k \geq 0 \]
\[ (i) \]
\[ q^* \leq c_{q,v} x^{-v} \text{ for all } x > 0 \text{ where } c_{q,v} \triangleq e^{-v} \left( \frac{q}{\ln(1/q)} \right)^{v}. \]
\[ (ii) \]

**Proposition 1 (Polynomial rate of convergence)** Consider Algorithm 1 where \( \alpha_i = \frac{2-\sqrt{3}}{L_{\text{max}}} \) and \( N_i(k) = \Pi_{t=1}^{\nu}(\Gamma_i(k)+t) \) for some positive integer \( v \geq 1 \). Suppose \( p_i = \frac{1}{n} \), Ass. 2 and 3 hold. Define \( C_v \triangleq \frac{(2-\sqrt{3})^v}{n L_{\text{max}}} \), \( \rho^* \leq 1 - \frac{(2-\sqrt{3})^2 \mu}{n L_{\text{max}}} \), and \( C_f \triangleq c_{\rho^*,v} C_v \frac{e^{2v}}{1-\rho^*} + c_{\rho^*,v} \mathbb{E}[F(x(0)) - F^*] \).

Then
\[ \mathbb{E}[F(x(k)) - F^*] \leq C_f k^{-v}, \quad \forall k \geq 1. \]
\[ (22) \]

**Proof.** By \( N_i(k) = \Pi_{t=1}^{\nu}(\Gamma_i(k)+t) \) and (19), we obtain that for all \( i \in N \) :
\[ \mathbb{E} \left[ N_i(k)^{-1} \right] = \sum_{m=0}^{k} \Pi_{i=1}^{\nu}(m+t)^{-1} \mathbb{P}(\Gamma_i(k) = m) = \sum_{m=0}^{k} \frac{k! p_i^m (1-p_i)^{k-m}}{m! (k-m)! \Pi_{i=1}^{\nu}(m+t)} \]
\[ \leq \sum_{m=0}^{k} \frac{k! p_i^m (1-p_i)^{k-m}}{(m+v)! (k-m)!} \leq \Pi_{i=1}^{\nu}(k+t)^{-1} k^{-v} \sum_{m=0}^{k} \frac{(k+v)! p_i^{m+v} (1-p_i)^{k-m}}{(m+v)! (k-m)!} \]
\[ \leq \Pi_{i=1}^{\nu}(k+t)^{-1} k^{-v} \leq (k+1)^{-v} p_i^{-v}. \]
Then by setting $\beta = \sqrt{3} - 1$, $\alpha_i = \frac{2 - \sqrt{3}}{L_i}$, and $p_i = \frac{1}{n}$ in (16), and using the above bound on $\mathbb{E}[N_i(k) \cdot k]$, we obtain the following recursion:

$$
\mathbb{E}[F(x(k+1)) - F^*] \leq \rho^k \mathbb{E}[F(x(k)) - F^*] + (k+1)^{-v} \left( \frac{2 - \sqrt{3}}{\sqrt{3}} v^2 n^{-1} \right) \sum_{i=1}^{n} L_i^{-1}
$$

$$
\leq (\rho^*)^{k+1} \mathbb{E}[F(x(0)) - F^*] + C \rho^{k+1} n^{-1} m^{-v}.
$$

(23)

This together with Lemma (4)(i) produces that

$$
\mathbb{E}[F(x(k)) - F^*] \leq (\rho^*)^k \left( C \rho^{k} + \mathbb{E}[F(x(0)) - F^*] \right) + \frac{2C \rho^{-v}}{\rho^* \ln(1/\rho^*)}.
$$

(24)

Since $\rho^* \in (0, 1)$ and $v \geq 1$, by Lemma (4)(ii), $(\rho^*)^k \leq c_{\rho^*, v} n^{-v} \forall k \geq 1$ with $c_{\rho^*, v} \triangleq e^{-\left( \frac{v}{\ln(1/\rho^*)} \right) v}$. This combined with (24) proves (22).

### 4.2 Iteration and Oracle Complexity

Next, we derive the iteration complexity (measured by the number of proximal oracle) and oracle complexity (measured by the expected number of stochastic first-order oracle) bounds for obtaining an $\epsilon$-optimal solution such that $\mathbb{E}[F(x)] - F^* \leq \epsilon$.

**Theorem 5 (Iteration Complexity)** Let $\{x(k)\}$ be generated by Algorithm 1 where $\alpha_i = \frac{2 - \sqrt{3}}{L_i}$, $p_i = \frac{1}{n}$ and $N_i(k) = \left( 1 - q_i \right)^{-\Gamma_i(k)}$ for some $q_i \in (0, \frac{2 - \sqrt{3}}{L_i \mu})$. Suppose Assumptions 2 and 3 hold. Define $q_{\min} \triangleq \min_{i \in \mathcal{N}} q_i$, $q^* = 1 - \frac{q_{\min}}{\mu}$, and $\eta_i \triangleq \frac{q_i}{n (1-q_i)} + 1$. Then the iteration complexity and oracle complexity required to obtain an $\epsilon$-optimal solution are respectively $\frac{\ln(1/\epsilon)}{\ln(1/\rho^*)}$ and $\frac{1}{n} \sum_{i=1}^{n} \frac{\eta_i}{\ln(1/\rho^*)} \frac{\ln(1/\epsilon)}{\ln(1/\rho^*)}$ with $\bar{\epsilon}$ defined by (25).

**Proof.** By $q_i < \frac{(2 - \sqrt{3})^2 \mu}{L_i}$ and $\forall i$ there holds $q_{\min} \leq \frac{(2 - \sqrt{3})^2 \mu}{L_{\max}}$. Define $K_1(\epsilon) \triangleq \left[ \frac{\ln(1/\epsilon)}{\ln(1/\rho^*)} \right]$, where $\bar{\epsilon} \triangleq \epsilon \left( F(x(0)) - F^* + \frac{(1 - q_{\min})}{(2 - \sqrt{3})^2 \mu} \frac{\sum_{i=1}^{n} L_i}{L_{\max}} \right)^{-1}$.

(25)

Then by (17) and $\alpha_i = \frac{2 - \sqrt{3}}{L_i}$, $p_i = \frac{1}{n}$, $q^* = 1 - \frac{q_{\min}}{n}$, the following holds:

$$
\mathbb{E}[F(x_k) - F^*] \leq \left( 1 - \frac{q_{\min}}{n} \right)^k \left( F(x(0)) - F^* + \frac{\sum_{i=1}^{n} \alpha_i \mu^2}{(2 - \sqrt{3})^2 \mu} \right) \leq \epsilon \quad \forall k \geq K_1(\epsilon).
$$

Then the number of PO to obtain an $\epsilon$-optimal solution is $K_1(\epsilon)$. By $p_i = 1/n$ and (19),

$$
\mathbb{E}[N_i(k)] \leq \mathbb{E} \left[ \left( 1 - q_i \right)^{-\Gamma_{i,k}} \right] + 1 = \sum_{m=0}^{k} \left( \frac{k}{m} \right) \left( p_i (1 - q_i)^{-1} \right)^m (1 - p_i)^{k-m} + 1
$$

$$
= (p_i (1 - q_i)^{-1} + 1 - p_i)^k + 1 = \eta_i^k + 1.
$$

(26)
Note that for \( \lambda > 1 \), \( \sum_{k=0}^{K} \lambda^k \leq \int_{0}^{K+1} \lambda^x dx \leq \frac{\lambda^{K+1}}{\ln \lambda} \). Since \( i_k \) is independent of \( N_i(k) \), by using (25) and (26), the expected number of SFO calls required to approximate an \( \epsilon \)–optimal solution is bounded as follows.

\[
\mathbb{E} \left[ K_1(\epsilon) - 1 \right] = \sum_{k=0}^{K} \sum_{i=1}^{n} N_i(k) I_{[i_k = i]} = \sum_{k=0}^{K} \sum_{i=1}^{n} \mathbb{E} [N_i(k)] \mathbb{E} [I_{[i_k = i]}]
\]

Thus, the expected number of SFO to obtain an \( \epsilon \)–optimal solution is bounded by

\[
\frac{1}{n} \sum_{i=1}^{n} \frac{\eta_i}{\ln(\eta_i)} \left( \frac{\ln(\eta_i)}{\ln(1/q^*)} \right) + \left[ \frac{\ln(1/\epsilon)}{\ln(1/q^*)} \right].
\]

Note that for any \( 0 < \epsilon, q < 1 \), we have the following relations:

\[
\eta_i \ln(1/q^*) = \left( e^{\ln(\eta_i)} \right) \ln(1/q^*) = e^{\ln(1/\epsilon)} \ln(1/q^*) = (1/\epsilon) \ln(1/q^*) .
\]

Thus, the expected number of SFO to obtain an \( \epsilon \)–optimal solution is bounded by

\[
\frac{1}{n} \sum_{i=1}^{n} \frac{\eta_i}{\ln(\eta_i)} \left( \frac{\ln(\eta_i)}{\ln(1/q^*)} \right) + \left[ \frac{\ln(1/\epsilon)}{\ln(1/q^*)} \right],
\]

giving us the required oracle complexity.

The following corollary emerges from Theorem 5 with \( q_i \) taking a specific form.

**Corollary 1** Let \( N_i(k) = \left[ (1 - q_i) - \Gamma_i(k) \right] \) with \( q_i = \frac{\theta_i \mu}{L_i} \) for some \( \theta_i \in (0, (2 - \sqrt{3})^2) \) in Algorithm 4 while \( n > 1 \) and the other conditions of Theorem 5 still hold. Define \( \theta_{\min} \triangleq \min_i \theta_i \) and \( \theta_{\max} \triangleq \max_i \theta_i \). Then the iteration and oracle complexity bounds to obtain an \( \epsilon \)–optimal solution are \( \mathcal{O} \left( \frac{L_{\max} \ln(1/\epsilon)}{\mu} \right) \) and \( \mathcal{O} \left( \frac{n L_{\max} \ln(1/\epsilon)}{\mu} \right) \), respectively.

**Proof.** We begin by deriving a bound on \( K_1(\epsilon) \):

\[
\frac{\ln(1/\epsilon)}{\ln(1/q^*)} = \frac{\ln(1/\epsilon)}{-\ln(q^*)} = \frac{\ln(1/\epsilon)}{-\ln(1 - q_{\min}/n)} \leq \frac{\ln(1/\epsilon)}{q_{\min}/n} = \frac{n L_{\max} \theta_{\min}}{\mu} \ln(1/\epsilon),
\]

where \( -\ln(1 - q_{\min}/n) \geq q_{\min}/n \) and \( q_{\min} \geq \frac{\theta_{\min} \mu}{L_{\max}} \), implying that \( K_1(\epsilon) = \mathcal{O} \left( \frac{n L_{\max} \ln(1/\epsilon)}{\mu} \right) \). Next, we analyze the two terms necessary for bounding the oracle complexity.

\[
\frac{\ln(\eta_i)}{\ln(1/q^*)} \leq \frac{n L_{\max} \theta_{\min}}{\mu \theta_{\max}} \ln(\eta_i) = \frac{n L_{\max} \theta_{\min}}{\mu \theta_{\max}} \ln(1 + q_i/(n - q_i)) \leq \frac{n L_{\max} \theta_{\min}}{\mu \theta_{\max}} \frac{q_i}{n - q_i} \leq \frac{n L_{\max} \theta_{\max} \theta_{\min}}{\theta_{\min} \mu L_i (n - \theta_{i}/L_i)} \leq \frac{n L_{\max} \theta_{\max}}{(n - 1/\kappa_{\min}) L_{\min} \theta_{\min}},
\]

where the second inequality holds by \( \ln(1 + x) \leq x, \forall x \in [0, 1] \), the fourth inequality follows from \( \theta_i \leq 1 \), and the last equality follows from \( \kappa_{\min} = L_{\min}/\mu \). In addition, we derive a bound on \( \eta_i/\ln(\eta_i) \) where
\[ \eta_i = 1 + \frac{q_i}{n(1-q_i)}. \]

\[ \frac{\eta_i}{\ln(\eta_i)} = \frac{1 + q_i/(n(1-q_i))}{\ln(1 + q_i/(n(1-q_i)))} = \frac{(1 + x^0)}{\ln(1 + x^0)} \leq \frac{(x^0 + 1)^2}{x^0} = x^0 + 2 + \frac{1}{x^0} = 2 + \frac{q_i}{n(1-q_i)} + \frac{n(1-q_i)}{q_i} \leq 2 + \frac{1}{n} + \frac{nL_i}{\theta_i \mu}, \]

where \( x^0 = q_i/(n(1-q_i)), q_i/(1-q_i) \leq 1 \) (since \( q_i \leq (2 - \sqrt{3})^2 \)), and the first inequality follows from \( \ln(1+x) \geq x/(x+1) \) for any \( x \geq 0 \). We prove the result by deriving a bound on the oracle complexity:

\[ \frac{1}{n} \sum_{i=1}^{n} \frac{\eta_i}{\ln(\eta_i)} \left( \frac{\ln(\eta_i)}{\ln(\eta_i^\mu)} + \left[ \frac{\ln(1/\epsilon)}{\ln(1/\rho^* \mu)} \right] \right) = \mathcal{O} \left( \frac{nL_i}{\mu} \left( \frac{\mathbb{E}[F(x(0))] - F^*}{\epsilon} + \frac{n\mu^2}{\epsilon} \right) \right). \]

\[ \Box \text{It can be seen that if } \theta_{\max} = \theta_{\min} (\text{by choosing } \theta_i = \theta \text{ for all } i) \text{ and } L_{\max} = L_{\min} = L, \text{ the oracle complexity reduces to } \mathcal{O}(n \kappa (1/\epsilon) + n / \mu) \text{ with } \kappa \triangleq \frac{L}{\mu}, \text{ which tends to the optimal oracle complexity of } \mathcal{O} \left( \frac{n}{\epsilon} \right) \text{ for large } n. \text{ From Theorem 5, we may obtain the optimal oracle complexity for } n = 1 \text{ by noting that } \ln(\eta_i) / \ln(1/\rho^*) = \ln(1/(1-q_i))/\ln(1/(1-q_i)) = 1. \]

**Corollary 2** Let \( \{x(k)\} \) be generated by Algorithm 1 where \( n = 1 \) and \( \alpha = \frac{2 - \sqrt{3}}{L} \). Suppose Assumptions 2 and 3 hold. Set \( N(k) = \left( (1-q)^{-k} \right) \) for some \( q \in (0, (2 - \sqrt{3})^2 \mu) \). Then the iteration and oracle complexity bounds for obtaining an \( \epsilon \)-optimal solution are \( \mathcal{O} \left( \frac{L}{\mu} \ln(1/\epsilon) \right) \) and \( \mathcal{O} \left( \frac{L}{\mu} \right) \), respectively.

In Theorems 4 and 5, we establish the rate as well as the iteration and oracle complexity bounds of Algorithm 1 when each block is randomly picked with equal probability. When blocks are chosen by a non-uniform distribution, we state a result in the smooth regime but omit the proof since it is similar to Theorems 4 and 5.

**Corollary 3** Suppose \( r(x) \equiv 0 \) and \( \bar{f} \) satisfies \( \|\nabla \bar{f}(x)\|^2 \geq 2\mu \left( \frac{\bar{f}(x) - F^*}{\mu} \right) \) with \( \mu > 0 \). Let \( \{x(k)\} \) be generated by Algorithm 1 where \( \alpha_i = \frac{2 - \sqrt{3}}{L_i}, p_i = \frac{\theta_i}{\sum_{j=1}^{i} L_j}, \) and \( N_i(k) = \left( (1-q_i)^{-1} \right) \) with \( q_i = \frac{\theta_i \mu}{L_i} \), for some \( \theta_i \in (0, (2 - \sqrt{3})^2) \). Then the iteration and oracle complexity bounds for obtaining an \( \epsilon \)-optimal solution are \( \mathcal{O} \left( \frac{nL_{\max}}{\mu} \ln \left( \frac{1}{\epsilon} \right) \right) \) and \( \mathcal{O} \left( \frac{nL_{\max}}{\mu} \left( \frac{1}{1 + \frac{1}{\mu \min \theta_{\min}}} \right) \right) \), respectively.

Finally, we investigate the iteration and oracle complexity of Algorithm 1 when the number of the sampled gradients increases at a slower polynomial rate.

**Proposition 2** Consider Algorithm 4 where \( \alpha_i = \frac{2 - \sqrt{3}}{L_i} \) and \( N_i(k) = \prod_{i=1}^{k} (\Gamma_i(k) + t) \) for some integer \( v \geq 1 \). Suppose \( p_i = \frac{1}{n} \forall i \in N \), Assumptions 2 and 3 hold. Define \( \rho^* = 1 - \frac{(2 - \sqrt{3})^2 \mu}{nL_{\max}} \) and \( C_v = \frac{(2 - \sqrt{3})^2 v^2}{2} \sum_{i=1}^{n} L_i^{-1} \). Then the iteration and oracle complexity bounds required to obtain an \( \epsilon \)-optimal solution are \( \mathcal{O}(v (1/\epsilon)^{1/v}) \) and \( \mathcal{O} \left( e^{v} v^{2v+1} (1/\epsilon)^{1+1/v} \right) \), respectively.
Proof. From (22) it follows that for any \( k \geq K(\epsilon) \triangleq \left( \frac{C_f}{\epsilon} \right)^{1/v} \), \( \mathbb{E}[F(x(k)) - F^*] \leq \epsilon \). By the definition of \( C_f \) in Prp. 1, it follows that \( C_f = O(e^v v^v) \), hence the iteration complexity is \( O((v/\epsilon)^{1/v}) \). Note that \((m+v)^v = \sum_{t=0}^v \frac{v!}{t!(v-t)!} m^t v^{v-t} \). Then by \( N_i(k) = \Pi_{t=1}^e (\Gamma_i(k) + t) \leq (\Gamma_i(k) + v)^v \) and using (20) we obtain that for any \( i \in \mathcal{N} \):

\[
\mathbb{E} [N_i(k)] \leq \sum_{m=0}^k (m + v)^v \mathbb{P}(\Gamma_i(k) = m)
\]

\[
= \sum_{t=0}^v \frac{v!}{t!(v-t)!} \sum_{m=0}^k m^t \frac{k!}{m!(k-m)!} p_i^m (1 - p_i)^{k-m} = \sum_{t=0}^v \frac{v!}{t!(v-t)!} \mathbb{P}[\Gamma_i(k) = t].
\]

By [23, p.154] we know that the \( t \)-th moment of the binomial distribution \( \Gamma_i(k) \) equals the \( t \)-th derivative of \( M_i(y) \) at \( y = 0 \), where \( M_i(y) = (p_i e^y + 1 - p_i)^k \). Thus, we can show that \( \mathbb{E}[\Gamma_i(k)^t] = O(k^t) \), hence by (28) and \( \left( \begin{array}{c} v \\ t \end{array} \right) v^v = O(v^v) \) we obtain \( \mathbb{E} [N_i(k)] = O(v^v(1/k)^v) \). Note by \( C_f = O(e^v v^v) \) that

\[
\sum_{k=0}^{K(\epsilon)-1} k^v \leq \int_{1}^{K(\epsilon)} t^v dt = \frac{t^{v+1}}{v+1} \bigg|_{1}^{K(\epsilon)} \leq (v+1)^{-1} \left( \frac{C_f}{\epsilon} \right)^{1/v+1} = O(v^v(1/e)^{1+1/v}).
\]

This together with \( \mathbb{E} [N_i(k)] = O(v^v(1/k)^v) \) implies that the expectation of the total number of sampled gradients required to obtain an \( \epsilon \)-optimal solution is bounded by

\[
\mathbb{E} \left[ \sum_{k=0}^{K(\epsilon)-1} \sum_{i=1}^n N_i(k) I_{[i=k]} \right] = \sum_{k=0}^{K(\epsilon)-1} \sum_{i=1}^n \mathbb{P} \mathbb{E} [N_i(k)] = O(v^v(1/e)^{1+1/v}).
\]

\[\square\]

5 Numerical Experiments

In this section, we will examine empirical algorithm performance on the sparse and nonlinear least squares problems to demonstrate the behavior of Algorithm 1. Throughout this section, the empirical mean error is based on averaging across 50 trajectories.

5.1 Sparse Least Squares

We apply Algorithm 1 on the following sparse least squares problem:

\[
\min_{x \in \mathbb{R}^d} \frac{1}{2N} \sum_{i=1}^N (a_i^T x - b_i)^2 + \lambda \|x\|_1 \quad \text{(LASSO)}
\]

with the regularization parameter \( \lambda > 0 \). We first generate a sparse vector \( x^* \) where only 10% components of the vector is nonzero with nonzero ones independently generated from the standard normal distribution. We then generate \( N \) samples \((a_i, b_i)\), where components of \( a_i \in \mathbb{R}^d \) are generated from standard normal distribution while \( b_i = a_i^T x^* + \hat{e} \) with \( \hat{e} \) normally distributed with zero mean and standard deviation 0.01. We partition \( x \in \mathbb{R}^d \) into \( n = 10 \) blocks and set \( \lambda = 0.1 \).

Sensitivity to sample-size policies: We now implement Algorithm 1 with \( \alpha = 0.01 \), \( p_i = \frac{1}{n} \) and the geometric batch-size \( N_i(k) = \left[ q^{-\Gamma_i(k)} \right] \), and investigate how the parameters \( q, N, d \) influence the algorithm.
performance. We ran Algorithm 1 for 50 epochs where each epoch implies the usage of all samples. The results are displayed in Table 3 for the empirical relative error $\frac{1}{N} \sum_{i=1}^{N} (F(x) - F^*)$, the number of proximal evaluations, and CPU times. The results suggest that for given a fixed simulation budget, slower geometric rates of growth of batch-sizes lead to better empirical error while requiring more CPU time since more proximal evaluations are needed. In addition, the running time increases approximately linearly with $N$ and $d$.

### Table 3: Comparison of the different selections of batch-sizes

| $N$  | $q$  | emp.err | prox.eval | CPU(s) |
|------|------|---------|-----------|--------|
| 1000 | 0.85 | 2.46e-02| 86 | 4.22 |
|      | 0.9  | 1.71e-02| 105 | 4.7  |
|      | 0.95 | 5.00e-03| 164 | 6.28 |
| 2000 | 0.85 | 3.71e-02| 90  | 7.11 |
|      | 0.9  | 2.49e-02| 112 | 8.82 |
|      | 0.95 | 6.10e-03| 178 | 11.48|
| 4000 | 0.85 | 1.27e-02| 94  | 16.15|
|      | 0.9  | 7.60e-03| 119 | 18.27|
|      | 0.95 | 1.90e-03| 192 | 24.3 |

### Comparison with BSG [41]:

Let $N = 2000$ and $d = 200$ in the LASSO problem. We compare Algorithm 1 with BSG [41] by running both schemes for 50 epochs. We show the results in Table 4 and plot trajectories in Figure 2, where BSG-$t$ denotes the minibatch BSG algorithm that utilizes $t$ samples at each iteration while in Algorithm 1 we set $N_i(k) = \lceil q - \Gamma_i(k) \rceil$. The empirical rate of convergence in terms of proximal evaluations shown in the upper Figure 2 implicitly supports the iteration complexity statements. We observe the following: (i) at first, minibatch BSG displays a faster decay in objective than Algorithm 1 since the batch-size in our scheme is relatively small at the outset; (ii) Algorithm 1 proceeds to catch up and outperform the minibatch BSG since the variance of the sampled gradient decreases with increasing batch-size; (iii) Both minibatch BSG with larger batch-sizes and Algorithm 1 with faster increasing batch-size display faster empirical rates with fewer proximal evaluations. The empirical algorithm performance in terms of epochs displayed in the lower Figure 2 demonstrates the results of oracle complexity. By comparing the number of samples given the fixed relative error, Algorithm 1 with $N_i(k) = \lceil 0.98^{-T_i(k)} \rceil$ has the best performance, which can also be concluded from Table 4.

### Table 4: Comparison of Algorithm 1 and BSG

|       | Algorithm 1, p=0.95 | Algorithm 1, p=0.98 | BSG-16 | BSG-64 |
|-------|---------------------|---------------------|--------|--------|
| emp.err | 4.30e-3 | 1.73e-4 | 2.60e-3 | 2.75e-4 |
| prox.eval | 178 | 375 | 6251 | 1563 |
| CPU(s)     | 5.94 | 10.25 | 118.65 | 33.36 |

### Influence of block-specific steplengths:

In this experiment, we set $N = 1000, d = 200$, and let the entries of $a_i \in \mathbb{R}^d$ corresponding to different blocks be generated from normal distributions with zero mean but with differing variances. Such data generation implies that the block-wise Lipschitz constants of $\frac{1}{2N} \sum_{i=1}^{N} (a_i^T x - b_i)^2$ can vary. We implement Algorithm 1 with the non-uniform block selection as per a distribution $p_i = \frac{L_i}{\sum_{i=1}^{K} L_i}$ in two settings: (i) the same steplength $\alpha_i \equiv \alpha = \frac{12}{L}$ depending on the
Lipschitz constant of $\nabla f(x)$, and (ii) the block-specific steplength $\alpha_i = \frac{1}{L_i}$ depending on the block-wise Lipschitz constant $L_i$. Such a selection ensures that the steplengths in two settings are approximately the same when the block-wise Lipschitz constants are identical. For a particular set of realizations with the Lipschitz constant satisfying $L_{\text{max}}/L_{\text{ave}} = 2.5$, the empirical iteration and oracle complexity of Algorithm 1 in the two settings are shown in Figures 1 and 3, respectively. These findings reinforce the point that block-specific steplengths, reliant on block-wise Lipschitz constant $L_i$, display better empirical behavior since less proximal evaluations (see Figure 1) and less sampled gradients (see Figure 3) are required for obtaining a solution with similar accuracy. In addition, we generate four sets of data, for which the global Lipschitz constant $L$ of the problem (LASSO) is the same while the ratio $L_{\text{max}}/L_{\text{ave}}$ is different. We then run Algorithm 1 with the identical and block-specific steplengths on the four generated datasets up to 100 epochs and compare the empirical errors. The results are shown in Table 5, where $x^I(K)$ and $x^B(K)$ denote the estimates generated by Algorithm 1 with identical and block-specific steplengths, respectively. Since the ratio $\frac{E[F(x^I(K))]-F^*}{E[F(x^B(K))]-F^*}$ is greater than one, we may conclude that the block-specific steplengths might lead to better algorithm performance compared with identical steplength. We observe that empirical error can be 50 times poorer when $L_{\text{max}}/L_{\text{ave}} = 1.47$.

| $L_{\text{max}}$ | $L_{\text{ave}}$ | $L_{\text{max}}$ | $L_{\text{ave}}$ | $L_{\text{max}}$ | $L_{\text{ave}}$ |
|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|
| 1.15            | 0.85            | 1.27            | 0.85            | 1.34            | 0.85            | 1.47            | 0.85            |
| $E[F(x^I(K))]-F^*$ | $27.5$         | $31.9$          | $52.4$          |
| $E[F(x^B(K))]-F^*$ | $15.3$           | $27.5$          | $31.9$          | $52.4$          |

Table 5: Comparison of Algorithm 1 with identical and block-specific steplengths

5.2 Nonlinear least squares

We consider a binary classification problem on a data set $\{x_i, y_i\}_{i=1}^N$, where $x_i \in \mathbb{R}^d$ and $y_i \in \{0, 1\}$ are the $i$th feature vector and the corresponding label, respectively. We consider the minimization of empirical error:

$$
\min_{w, b} \frac{1}{2N} \sum_{i=1}^N \left( y_i - \phi(w^T x_i + b) \right)^2,
$$

(NLS)
where $\phi(z) = \frac{1}{1+e^{-z}}$ is the sigmoid function. We will apply Algorithm 1 to the gisette from LIBSVM library and how does the batch-size influence the training loss and misclassification rate. We partition the vector $w \in \mathbb{R}^d$ into $n = 10$ blocks. We implement Algorithm 1 with $\alpha = 0.2$, where the batch-sizes are set to be the constant batch-sizes $N_i(k) \equiv 0.02N, 0.05N$, and the increasing batch-sizes $N_i(k) = \max\{0.001N, \Gamma_i(k)\}$, $\max\{0.001N, \Gamma_i(k)^2\}$. From Figure 4 we conclude that smaller batch-sizes would lead to better performance if we run the algorithm with a relatively smaller amount of samples (e.g., $N$); the mini-batch schemes may not perform well if the batch-size is not suitably selected, for instance, $0.05N$. Favorable behavior follows if the batch-size increases at a suitable rate, e.g., linearly.

6 Concluding remarks

Existing block-based techniques for stochastic nonconvex optimization rely on centrally mandated batch-sizes and steplengths bounded by the global Lipschitz constant, leading to larger oracle complexities and poorer performance (because of shorter steps), as well as higher informational coordination requirements. We consider minimizing the sum of an expectation-valued smooth nonconvex function and a nonsmooth separable convex function through a limited coordination asynchronous variance-reduced method, reliant on block-specific steplengths and random decentralized batch-sizes. The almost sure convergence of the generated iterates is established. In addition, the scheme achieves the deterministic rate of $O(1/K)$ with the rate and oracle complexities dependent on $L_{\text{ave}}$ rather than $L_{\text{max}}$. Furthermore, under the proximal PL requirement, the iterates provably converge linearly (polynomially) to the global optimum in a mean sense when batch-sizes grow geometrically (polynomially). Notably, despite using randomized batch-size sequences, we show that the deterministic iteration complexities may be achieved. Specifically, the schemes achieve the optimal oracle complexity when the problem (1) admits a single block. Finally, numerical studies are carried out to support the theoretical findings and reveal that schemes that leverage block-specific Lipschitz constants lead to significantly superior empirical behavior.

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A Proof of Lemma 2

Let \( x_{-i} \) denote the coordinates of the variable \( x \) except those correspond to block \( i \). By applying Lemma 1 to the function \( \bar{f}(\cdot, x_{-i_k}(k)) \) and Eqn. (5) with \( y = \bar{x}_{i_k}(k+1), z = x_{i_k}(k) \), and \( g = \nabla x_{i_k} \bar{f}(x(k)) \), we obtain the following inequality:

\[
\bar{f}(x_{-i_k}(k), \bar{x}_{i_k}(k+1)) + r_{i_k}(\bar{x}_{i_k}(k+1)) \\
\leq \bar{f}(x(k)) + r_{i_k}(x_{i_k}(k)) + \left( \frac{L_{i_k}}{2} - \frac{1}{\alpha_{i_k}} \right) \| \bar{x}_{i_k}(k+1) - x_{i_k}(k) \|^2.
\]

Define \( \bar{x}(k+1) \) as follows:

\[
\bar{x}_{i_k}(k+1) \triangleq \bar{x}_{i_k}(k+1) \quad \text{and} \quad \bar{x}_j(k+1) \triangleq x_j(k) \quad \forall j \neq i_k.
\]

Then \( r_j(\bar{x}_j(k+1)) = r_j(x_j(k)) \quad \forall j \neq i_k \), and hence we obtain the following bound:

\[
F(\bar{x}(k+1)) \leq F(x(k)) + \left( \frac{L_{i_k}}{2} - \frac{1}{\alpha_{i_k}} \right) \| \bar{x}_{i_k}(k+1) - x_{i_k}(k) \|^2.
\]

(30)
By applying Lemma 1 to the function $\tilde{f}(\cdot, x_{-i_k}(k))$ and the update (4) with $y = x_i(k+1)$, $z = x_{ik}(k+1)$, $x = x_{ik}(k)$, and $g = \nabla_{x_{ik}} \tilde{f}(x(k)) + w_{ik}(k+1)$, one obtains

$$
\tilde{f}(x_{-i_k}(k), x_{ik}(k+1)) + r_{ik}(x_{ik}(k+1)) \
- (x_{ik}(k+1) - \bar{x}_{ik}(k+1))^T w_{ik}(k+1) + \left(\frac{L_{ik}}{2} - \frac{1}{2\alpha_{ik}}\right) \|x_{ik}(k+1) - x_{ik}(k)\|^2 \
- \left(\frac{L_{ik}}{2} + \frac{1}{2\alpha_{ik}}\right) \|\bar{x}_{ik}(k+1) - x_{ik}(k)\|^2.
$$

By recalling that $-a^T b \leq -a^T a/2 + \alpha/2 \|b\|^2$, the following holds:

$$
- (x_{ik}(k+1) - \bar{x}_{ik}(k+1))^T w_{ik}(k+1) \
\leq \frac{1}{2\alpha_{ik}} \|x_{ik}(k+1) - \bar{x}_{ik}(k+1)\|^2 + \frac{\alpha_{ik}}{2} \|w_{ik}(k+1)\|^2.
$$

Therefore, by substituting (32) into (31), we obtain the following bound:

$$
F(x(k+1)) \leq F(\bar{x}(k+1)) + \left(\frac{L_{ik}}{2} - \frac{1}{2\alpha_{ik}}\right) \|x_{ik}(k+1) - x_{ik}(k)\|^2 \
+ \left(\frac{L_{ik}}{2} + \frac{1}{2\alpha_{ik}}\right) \|\bar{x}_{ik}(k+1) - x_{ik}(k)\|^2.
$$

By adding inequalities (30) and (33),

$$
F(x(k+1)) \leq F(x(k)) + \left(\frac{L_{ik}}{2} - \frac{1}{2\alpha_{ik}}\right) \|x_{ik}(k+1) - x_{ik}(k)\|^2 \
+ \frac{\alpha_{ik}}{2} \|w_{ik}(k+1)\|^2.
$$

Note that for all $i = 1, \cdots, n$, $L_{ik}/2 - 1/2\alpha_{ik} \leq 0$ by $\alpha_{ik} \leq 1/L_{ik}$. Then the second term on the right-hand side of Eqn. (34) is nonpositive, hence we can take out this term from the upper bound of $F(x(k+1))$. Since $x(k)$ is adapted to $\mathcal{F}_k$, by taking expectations conditioned on $\mathcal{F}_k$ on both sides of (34), we obtain that

$$
E[F(x(k+1))|\mathcal{F}_k] \leq F(x(k)) + \left(\frac{L_{ik}}{2} - \frac{1}{2\alpha_{ik}}\right) \|\bar{x}_{ik}(k+1) - x_{ik}(k)\|^2|\mathcal{F}_k|
$$

+ $\frac{1}{2} E[\alpha_{ik} \|w_{ik}(k+1)\|^2|\mathcal{F}_k].
$$
Note that for any $i \in \mathcal{N}$, $\bar{x}_i(k + 1)$ is adapted to $\mathcal{F}_k$ by the definition (5), and $i_k$ is independent of $\mathcal{F}_k$ by Assumption 3(ii). Therefore, by [5, Corollary 7.1.2] and $\mathbb{P}(i_k = i) = p_i$, the following holds a.s.:

$$
\mathbb{E} \left[ \left( L_{i_k} - \frac{1}{2\alpha_{i_k}} \right) \| x_{i_k}(k + 1) - x_{i_k}(k) \|^2 | \mathcal{F}_k \right] = \sum_{i=1}^{n} p_i \left( L_i - \frac{1}{2\alpha_i} \right) \| x_i(k + 1) - x_i(k) \|^2. 
$$

(36)

Then by substituting (36) into (35), we obtain Eqn. (6).

\[ \square \]

**B Proof of Lemma 3.**

By recalling that the gradient map $\nabla_x \tilde{f}(x)$ is $L_i$-Lipschitz continuous from Assumption 1(ii) and that $\tilde{x}_j(k + 1) = x_j(k) \forall j \neq i_k$ by definition (29), we have the following inequality:

$$
\tilde{f}(\tilde{x}(k + 1)) \leq \tilde{f}(x(k)) + (\tilde{x}_{i_k}(k + 1) - x_{i_k}(k))^T \nabla_{x_{i_k}} \tilde{f}(x(k)) + \frac{L_{i_k}}{2} \| x_{i_k}(k + 1) - x_{i_k}(k) \|^2.
$$

Using the definition of $\tilde{x}_{k+1}$ in (29), we have that $r_j(\tilde{x}_j(k + 1)) = r_j(x_j(k)) \forall j \neq i_k$, $\tilde{x}_{i_k}(k + 1) = \tilde{x}_{i_k}(k + 1)$. Thus, the following relation holds

$$
\tilde{F}(\tilde{x}(k + 1)) \leq F(x(k)) + (\tilde{x}_{i_k}(k + 1) - x_{i_k}(k))^T \nabla_{x_{i_k}} \tilde{f}(x(k)) + \frac{L_{i_k}}{2} \| x_{i_k}(k + 1) - x_{i_k}(k) \|^2 + r_{i_k}(\tilde{x}_{i_k}(k + 1)) - r_{i_k}(x_{i_k}(k)),
$$

(37)

where the last inequality holds by $\alpha_i < 1/L_i \forall i \in \mathcal{N}$. Since for any $i \in \mathcal{N}$, $\tilde{x}_{i}(k + 1)$ is adapted to $\mathcal{F}_k$ by its definition (5), and $i_k$ is independent of $\mathcal{F}_k$. Then, by [5, Corollary 7.1.2] and $\mathbb{P}(i_k = i) = p_i$, we have that

$$
\mathbb{E} \left[ (\tilde{x}_{i_k}(k + 1) - x_{i_k}(k))^T \nabla_{x_{i_k}} \tilde{f}(x(k)) + \| \tilde{x}_{i_k}(k + 1) - x_{i_k}(k) \|^2 / (2\alpha_{i_k}) + r_{i_k}(\tilde{x}_{i_k}(k + 1)) - r_{i_k}(x_{i_k}(k)) | \mathcal{F}_k \right] 
$$

$$
= \sum_{i=1}^{n} p_i \left( (\tilde{x}_i(k + 1) - x_i(k))^T \nabla_{x_i} \tilde{f}(x(k)) + \frac{1}{2\alpha_i} \| \tilde{x}_i(k + 1) - x_i(k) \|^2 + r_i(\tilde{x}_i(k + 1)) - r_i(x_i(k)) \right)
$$

\[ \leq \mathbb{E} \left[ \min_{y_i \in \mathbb{R}^{d_i}} \left( \nabla_{x_i} \tilde{f}(x(k))^T (y_i - x_i(k)) + \frac{1}{2\alpha_i} \| y_i - x_i(k) \|^2 + r_i(y_i) - r_i(x_i(k)) \right) \right],
$$

\[ \leq p_{\min} \sum_{i=1}^{n} \min_{y_i \in \mathbb{R}^{d_i}} \left[ \nabla_{x_i} \tilde{f}(x(k))^T (y_i - x_i(k)) + \frac{1}{2\alpha_i} \| y_i - x_i(k) \|^2 + r_i(y_i) - r_i(x_i(k)) \right].
$$

Let the random vectors $X \in \mathbb{R}^m$ and $Y \in \mathbb{R}^m$ be independent of one another and let $f$ be a Borel function on $\mathbb{R}^{m \times n}$ with $|\mathbb{E}[f(X,Y)]| \leq \infty$. If for any $x \in \mathbb{R}^m$, $g(x) = \mathbb{E}[f(x,Y)]$ if $|\mathbb{E}[f(x,Y)]| \leq \infty$ and $g(x) = 0$ otherwise, then $g$ is a Borel function with $g(X) = \mathbb{E}[f(X,Y)|\sigma(X)]$.
where the last inequality follows by $\min_{y_i \in \mathbb{R}^d} [\nabla_x f(x(k))^T (y_i - x_i(k)) + \frac{1}{2\alpha_i} \|y_i - x_i(k)\|^2 + r_i(y_i) - r_i(x_i(k))] \leq 0$. Then by $\alpha_i^{-1} \leq \alpha_{\min}$ and Assumption 3, the above equation is further bounded by

$$
p_{\min} \sum_{i=1}^{n} \min_{y_i \in \mathbb{R}^d} \left[ \nabla_x f(x(k))^T (y_i - x_i(k)) + \frac{1}{2\alpha_{\min}} \|y_i - x_i(k)\|^2 + r_i(y_i) - r_i(x_i(k)) \right]
= -\frac{p_{\min} \alpha_{\min}}{2} D_r(x(k), \alpha_{\min}^{-1}) \leq -\frac{\alpha_{\min} p_{\min}}{2} D_r(x(k), \ell_{\max})
\leq -\alpha_{\min} \mu_{p_{\min}} (F(x(k)) - F^*)
$$

(38)

where the first inequality follows from [17, Lemma 1] since $D_r(x, \cdot)$ is nonnegative and nondecreasing in $(0, \infty)$ and $\alpha_{\min}^{-1} \geq \ell_{\max}$. Then by taking unconditional expectations on both sides of (37) and using (38), we obtain that

$$\mathbb{E}[F(\bar{x}(k + 1))] \leq \mathbb{E}[F(x(k))] - \alpha_{\min} \mu_{p_{\min}} \mathbb{E}[F(x(k)) - F^*].
$$

(39)

By taking unconditional expectations on both sides of (30) and using $\mathbb{P}(i_k = i) = p_i$, we obtain

$$\mathbb{E}[F(\bar{x}(k + 1))] \leq \mathbb{E}[F(x(k))] + \sum_{i=1}^{n} p_i \left( \frac{L_i}{2} - \frac{1}{\alpha_i} \right) \|x_i(k + 1) - x_i(k)\|^2.
$$

(40)

Adding $(1 - \beta) \times (39)$ to $\beta \times (40)$ with $\beta \in (0.5, 1)$, we obtain the following inequality:

$$\mathbb{E}[F(\bar{x}(k + 1))] \leq \mathbb{E}[F(x(k))] + \beta \sum_{i=1}^{n} p_i \left( \frac{L_i}{2} - \frac{1}{\alpha_i} \right) \|x_i(k + 1) - x_i(k)\|^2
- \alpha_{\min} (1 - \beta) \mu_{p_{\min}} \mathbb{E}[F(x(k)) - F^*].
$$

(41)

Using $\alpha_i < \frac{1}{L_i}$, Assumption 2 and $\mathbb{P}(i_k = i) = p_i$, and by taking unconditional expectations on both sides of (33), the following holds:

$$\mathbb{E}[F(x_{k+1})] \leq \mathbb{E}[F(\bar{x}(k + 1))]
+ \sum_{i=1}^{n} p_i \left( \frac{L_i}{2} + \frac{1}{2\alpha_i} \right) \|x_i(k + 1) - x_i(k)\|^2 + \frac{L_i^2}{2} \sum_{i=1}^{n} \alpha_i p_i \mathbb{E}[N_i(k)^{-1}].
$$

(42)

Therefore, by adding inequality (42) to (41) yields the following bound:

$$\mathbb{E}[F(x(k + 1))] \leq \mathbb{E}[F(x(k))] + \sum_{i=1}^{n} p_i \left( \frac{L_i (1 + \beta)}{2} - \frac{2\beta - 1}{2\alpha_i} \right) \|x_i(k + 1) - x_i(k)\|^2
- \alpha_{\min} (1 - \beta) \mu_{p_{\min}} \mathbb{E}[F(x(k)) - F^*] + \frac{L_i^2}{2} \sum_{i=1}^{n} \alpha_i p_i \mathbb{E}[N_i(k)^{-1}].
$$

(43)

By recalling that $0 < \alpha_i \leq \frac{2\beta - 1}{L_i (1 + \beta)}$, we get $\frac{L_i (1 + \beta)}{2} - \frac{2\beta - 1}{2\alpha_i} \leq 0$. Thus, by subtracting $F^*$ from both sides of (43), we obtain (16). \qed