

Generalized Stochastic Frank-Wolfe Algorithm with Stochastic “Substitute” Gradient for Structured Convex Optimization

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Abstract

The stochastic Frank-Wolfe method has recently attracted much general interest in the context of optimization for statistical and machine learning due to its ability to work with a more general feasible region. However, there has been a complexity gap in the guaranteed convergence rate for stochastic Frank-Wolfe compared to its deterministic counterpart. In this work, we present a new stochastic Frank-Wolfe method which closes this gap by introducing the notion of a “substitute” gradient that is a not-necessarily unbiased sample of the gradient. Moreover, we show that this new approach is equivalent to a randomized coordinate mirror descent algorithm applied to the dual problem, which in turn provides a new interpretation of dual coordinate descent method in the primal space. When the regularizer is furthermore strongly convex, we show that the generalized stochastic Frank-Wolfe method as well as the randomized dual coordinate descent present linear convergence. These new results are benefited from the understanding that first-order methods can inherently minimize the primal-dual gap.

1 Introduction

1.1 Problem set-up, examples, Algorithm 1

Our problem of interest is the following optimization problem:

$$\text{P :} \quad \min_{\beta} P(\beta) := \frac{1}{n} \sum_{j=1}^n l_j(x_j^T \beta) + R(\beta) , \quad (1)$$

where $\beta \in \mathbb{R}^p$, $l_j(\cdot) : \mathbb{R} \rightarrow \mathbb{R}$, $j = 1, \dots, n$, is a univariate function (the j^{th} loss function), $s_j = x_j^T \beta$ is the “fitted value” of the model β for the data sample x_j , and $R(\cdot)$ is some other function that can be used to model a regularizer and/or an indicator function of a feasible region Q , and/or a penalty term, coupling constraints, etc. Notice that the scalar variable s_j for $l_j(s_j)$ is a linear function of β , namely $s_j = x_j^T \beta$. We will shortly give several natural examples in statistical and machine learning where this structure arises quite naturally. Throughout this paper we assume the following regarding these functions:

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Assumption 1.1. *The following hold:*

1. *for $j = 1, \dots, n$, the univariate function $l_j(\cdot)$ is strictly convex and γ -smooth, namely $|\dot{l}_j(a) - \dot{l}_j(b)| \leq \gamma|a - b|$ for all a, b ,*
2. *$\text{dom}R(\cdot)$ is bounded, and the subproblem*

$$\min_{\beta} c^T \beta + R(\beta) \tag{2}$$

attains its optimum and can be easily solved for any c , and

3. *$0 \in \text{dom}R(\cdot)$.*

We note regarding (1.) above that the strict convexity (instead of simple convexity) of $l_j(\cdot)$ is only needed to guarantee that the conjugate function $l_j^*(\cdot)$ is differentiable, and that this can be relaxed both algorithmically as well as in the proofs herein, but at considerable notational and expositional expense. Regarding (2.), this is a generalization of a linear optimization oracle as follows: in the case when $R(\cdot)$ is the indicator function $\mathbf{I}_Q(\cdot)$ of a set $Q \subset \mathbb{R}^p$ (namely, $\mathbf{I}_Q(\beta) := 0$ if $\beta \in Q$, and $\mathbf{I}_Q(\beta) := +\infty$ otherwise), then Q is the feasible region of P , and (2.) states that the feasible region Q is bounded and that it is easy to solve linear optimization problems on Q . Also, (3.) above is for notational convenience, as we can always translate a given feasible point so that $0 \in \text{dom}R(\cdot)$.

Here we present several applications of our problem setup (1) in statistical and machine learning. (For other applications particularly amenable to solution by the Frank-Wolfe method, we refer the reader to [20].)

Example 1.1. LASSO[42], ridge regression[19], sparse logistic regression[35]. *Consider the least-squares regression problem where a set of training samples $\{(x_j, y_j)\}_{j=1}^n$ is given. The LASSO optimization problem (in constraint format) is:*

$$\begin{aligned} \min_{\beta} \quad & \frac{1}{2n} \sum_{j=1}^n (y_j - x_j^T \beta)^2 \\ \text{s.t.} \quad & \|\beta\|_1 \leq \delta, \end{aligned}$$

which is an instance of P by using the least squares loss function $l_j(\cdot) = \frac{1}{2}(y_j - \cdot)^2$ and using the indicator function of an ℓ_1 ball as the regularizer, namely $R(\beta) := \mathbf{I}_{\{\|\beta\|_1 \leq \delta\}}(\beta)$.

The ridge regression optimization problem adds the regularizer $\frac{\lambda}{2}\|\beta\|_2^2$ to the least squares objective function for the parameter $\lambda > 0$, and omits the ℓ_1 ball constraint. Notice that because $\beta = 0$ is a feasible solution it follows that the optimal objective value is bounded above by $\|y\|_2^2/(2n)$, and therefore we can model the regularizer using $R(\beta) = \frac{\lambda}{2}\|\beta\|_2^2 + \mathbf{I}_{\{\|\beta\|_2^2 \leq \|y\|_2^2/(n\lambda)\}}(\beta)$, which ensures that $\text{dom}R(\cdot)$ is bounded.

The ℓ_1 -regularized logistic regression optimization problem seeks a solution of:

$$\min_{\beta} P(\beta) = \frac{1}{n} \sum_{j=1}^n \ln(1 + \exp(-y_j x_j^T \beta)) + \lambda \|\beta\|_1,$$

for a given set of training samples $\{(x_j, y_j)\}_{j=1}^n$ where $y_j \in \{-1, 1\}$, and is an instance of P using the logistic loss functions $l_j(\cdot) = \ln(1 + \exp(-y_j \cdot))$ with the ℓ_1 -regularizer $R(\beta) = \lambda \|\beta\|_1 + \mathbf{I}_{\{\|\beta\|_1 \leq \ln(2)/\lambda\}}(\beta)$ where the indicator function term is structurally redundant but is added as in the previous example to ensure that $\text{dom}R(\cdot)$ is bounded.

Example 1.2. Matrix completion[10][7]. *In the matrix completion problem, we seek to compute a low-rank matrix that well-approximates a given matrix $M \in \mathbb{R}^{n \times p}$ on the set Ω of observed entries (i, j) . The convex relaxation of this problem is the following nuclear-norm optimization problem:*

$$\begin{aligned} \min_{\beta \in \mathbb{R}^{n \times p}} \quad & \frac{1}{2|\Omega|} \sum_{(i,j) \in \Omega} (M_{i,j} - \beta_{i,j})^2 \\ \text{s.t.} \quad & \|\beta\|_* \leq \delta, \end{aligned}$$

where $\|\cdot\|_*$ is the nuclear norm. In order to translate the matrix completion problem to the setting of P , we consider any index pair $(i, j) \in \Omega$ as a sample, and we have $l_{(i,j)}(\cdot) = \frac{1}{2}(\cdot - M_{i,j})^2$, and $R(\beta) = \mathbf{I}_{\{\|\beta\|_* \leq \delta\}}(\beta)$.

Example 1.3. Structured sparse matrix estimation with CUR factorization[29][31]. *We seek to compute an approximate factorization $M \approx CUR$ of a given data matrix $M \in \mathbb{R}^{n \times d}$ such that C contains a subset of c columns from M and R contains a subset of r rows from M . Mairal et al. [31] proposed the following convex relaxation of this problem:*

$$\begin{aligned} \min_{\beta} \quad & \frac{1}{2nd} \|M - M\beta M\|_F^2 \\ \text{s.t.} \quad & \sum_i \|\beta_{i,\cdot}\|_\infty \leq \delta \\ & \sum_j \|\beta_{\cdot,j}\|_\infty \leq \delta, \end{aligned}$$

which is an instance of P by modeling the $(i, j)^{\text{th}}$ loss term in P as $\frac{1}{2}(M_{i,j} - M_i^T \beta M_j)^2$ (which is a least squares loss of a particular linear function of the matrix variable β), and $R(\beta) = \mathbf{I}_{\{\sum_i \|\beta_{i,\cdot}\|_\infty \leq \delta, \sum_j \|\beta_{\cdot,j}\|_\infty \leq \delta\}}(\beta)$.

Let $X \in \mathbb{R}^{n \times p}$ denote the data matrix whose rows are comprised of the vectors $x_1, \dots, x_n$, i.e., the j^{th} row of X is the vector x_j , $j = 1, \dots, n$. Let us define $L(s) : \mathbb{R}^n \rightarrow \mathbb{R}$ by $L(s) := \sum_{j=1}^n l_j(s_j)$ which is the total losses associated with $s \in \mathbb{R}^n$. One can think of $s = X\beta$ as the “fitted values” in the context of linear or logistic regression.

Algorithm 1 presents the main algorithmic contribution of this paper, which is a first-order method for tackling the problem P . We call the method “Stochastic Generalized Frank-Wolfe method with Stochastic Substitute Gradient” for reasons which we will discuss as we walk through the structure of the method below.

We can write the first part of the objective function of P as $f(\beta) := \frac{1}{n}L(X\beta) = \frac{1}{n}L(s)$ with $s = X\beta$. We have $\nabla L(s) = (\dot{l}_1(s_1), \dots, \dot{l}_n(s_n))$ and the gradient of $f(\cdot)$ can be written as

$$\nabla f(\beta) = \frac{1}{n}X^T \nabla L(X\beta) = \frac{1}{n} \sum_{j=1}^n \dot{l}_j(x_j^T \beta) x_j, \quad (3)$$

which we can re-write as $\nabla f(\beta) = \frac{1}{n}X^T w$ where $w = \nabla L(s)$ and $s = X\beta$, and which can be alternatively stated as:

$$\nabla f(\beta) = \frac{1}{n} \sum_{j=1}^n w_j x_j \text{ where } w_j = \dot{l}_j(s_j) \text{ and } s_j = x_j^T \beta, \quad j = 1, \dots, n. \quad (4)$$

Algorithm 1 Stochastic Generalized Frank-Wolfe with Stochastic Substitute Gradient

Initialize. Initialize with $\bar{\beta}^{-1} = 0$, $s^0 = 0$, and substitute gradient $d^0 = \frac{1}{n}X^T \nabla L(s^0)$, with step-size sequences $\{\alpha_i\} \in (0, 1]$ and $\{\eta_i\} \in (0, 1]$.

For iterations $i = 0, 1, \dots$

Solve l.o.o. subproblem: Compute $\tilde{\beta}^i \in \arg \min_{\beta} \left\{ (d^i)^T \beta + R(\beta) \right\}$

Choose random index: Choose $j_i \in \mathcal{U}[1, \dots, n]$

Update s value: $s_{j_i}^{i+1} \leftarrow (1 - \eta_i)s_{j_i}^i + \eta_i(x_{j_i}^T \tilde{\beta}^i)$, and $s_j^{i+1} \leftarrow s_j^i$ for $j \neq j_i$

Update substitute gradient: $d^{i+1} = \frac{1}{n}X^T \nabla L(s^{i+1}) = d^i + \frac{1}{n} \left(\dot{l}_{j_i}(s_{j_i}^{i+1}) - \dot{l}_{j_i}(s_{j_i}^i) \right) x_{j_i}$

Update primal variable: $\bar{\beta}^i \leftarrow (1 - \alpha_i)\bar{\beta}^{i-1} + \alpha_i \tilde{\beta}^i$.

(Optional Accounting:) $w^{i+1} \leftarrow \nabla L(s^{i+1})$

Here we emphasize that w is the vector of weights on the data values X in the composition of the gradient, and s is the vector of fitted values $X\beta$.

Especially in the context of “big data” applications of statistical and machine learning where n is huge, it can be extremely expensive to compute $\nabla f(\cdot)$. We therefore maintain a “substitute gradient” in Algorithm 1 that is constructed stochastically. This is accomplished as follows: let $\bar{\beta}^{i-1}$ be the value of β at the start of iteration i of the method, and we have a substitute gradient d^i that is the current proxy/substitute for the true gradient $\nabla f(\bar{\beta}^{i-1})$, where d^i is computed by:

$$d^i = \frac{1}{n} \sum_{j=1}^n w_j^i x_j \text{ where } w_j^i = \dot{l}_j(s_j^i), \quad j = 1, \dots, n, \quad (5)$$

for a given s^i that is the value of s at iteration i . But in contrast to (4) it will not necessarily hold that $s_j^i = x_j^T \bar{\beta}^i$ for $j = 1, \dots, n$ (equivalently $s^i = X\bar{\beta}^i$). (In fact, d^i will not necessarily be an unbiased estimate of $\nabla f(\bar{\beta}^i)$ as this will not be needed.) In the identical spirit as randomized coordinate descent, s^{i+1} will be determined by choosing a random index $j_i \in \mathcal{U}[1, \dots, n]$ and updating only the coordinate j_i of s^i , so that $s^{i+1} = s^i + \Delta^i e_{j_i}$ for some specific iteration-dependent scalar Δ^i (where e_ℓ denotes the ℓ^{th} unit coordinate vector in $\mathbb{R}^n$). This is accomplished in the “choose random index” step and the “update s value” step in Algorithm 1.

We now walk through the structure of Algorithm 1 in complete detail. The method is initialized with the initial decision variable β set to $\bar{\beta}^{-1} = 0$ and its fitted value $s^0 = X\bar{\beta}^{-1} = 0$ and initial substitute gradient $d^0 = \frac{1}{n}X^T \nabla L(s^0)$, which corresponds to the true fitted value and true gradient at $\bar{\beta}^{-1} = 0$. In iteration i , we use the substitute gradient d^i to compute $\tilde{\beta}^i$, which is a solution to the (generalized) linear optimization oracle (“l.o.o.”), where recall that this step specifies to solving a linear optimization problem over a set Q in the specific case when the $R(\cdot)$ is the indicator function of Q , namely $R(\cdot) = \mathbf{I}_Q(\cdot)$. Regarding updating the current fitted values s^i , we randomly choose a sample (a coordinate) j_i and only update s_{j_i} as a certain convex combination of the current fitted value $s_{j_i}^i$ and the fitted value for the j_i^{th} sample at $\tilde{\beta}^i$, namely $x_{j_i}^T \tilde{\beta}^i$, so that $s_{j_i}^{i+1} \leftarrow (1 - \eta_i)s_{j_i}^i + \eta_i(x_{j_i}^T \tilde{\beta}^i)$. Then we update the substitute gradient to make sure that $d^{i+1} = \frac{1}{n}X^T \nabla L(s^{i+1})$. The last step at iteration i is to take a Frank-Wolfe step to update $\bar{\beta}^i \leftarrow$

$(1 - \alpha_i)\bar{\beta}^{i-1} + \alpha_i\tilde{\beta}^i$ by taking a convex combination of the previous primal variable value $\bar{\beta}^{i-1}$ and the solution $\tilde{\beta}^i$ of the just-solved linear optimization oracle. Finally – and “optionally” since it does not affect future computations – we can perform an optional accounting step to update the dual variable $w^{i+1} \leftarrow \nabla L(s^{i+1})$ in order to compute a duality gap certificate. (The nature of this duality will be understood once we look at the dual problem of P in Section 2.)

Note that the computations in Algorithm 1 are minimally affected by the dimension n . Except for the initial computation of the gradient d^0 which is $O(np)$ operations, s^i and w^i are only updated by one coefficient at each iteration, and d^{i+1} is updated by adding a scalar multiple of x_{j_i} to d^i , which is $O(p)$ operations. The updates of $\bar{\beta}^i$ are $O(p)$ operations after solving for the optimal value $\tilde{\beta}^i$ in the linear optimization oracle, which is assumed to be easy to compute.

It is useful to place Algorithm 1 in the context of the Frank-Wolfe method. The Frank-Wolfe method is designed primarily to tackle the constrained convex optimization problem: $\min_{\beta \in Q} f(\beta)$ where $f(\cdot)$ is a smooth convex function and Q is a convex body, and it is assumed that linear optimization over Q is easy to compute. The optimization problem can of course be re-written as $\min_{\beta} f(\beta) + R(\beta)$ with $R(\cdot) = \mathbf{I}_Q(\cdot)$. The Frank-Wolfe update is:

$$\tilde{\beta}^i \in \arg \min_{\beta \in Q} \{\nabla f(\beta^i)^T \beta\} \quad \text{and} \quad \beta^{i+1} = (1 - \alpha_i)\beta^i + \alpha_i\tilde{\beta}^i. \quad (6)$$

It can be shown that with an appropriate choice of step-size sequence $\{\alpha_i\}$ that the Frank-Wolfe method computes an ε -optimal solution in $O(\frac{1}{\varepsilon})$ iterations, see [11], [14], and [12].

Due to its low iteration cost and convenient structural properties, the Frank-Wolfe method is especially applicable in several areas of statistical and machine learning and has thus received much renewed interest in recent years, see [20], [16], [13], [12], and the references therein. The Frank-Wolfe method can be generalized to deal with the more general problem $\min_{\beta} f(\beta) + R(\beta)$ where $R(\cdot)$ is any convex function with bounded domain and for which the “linear optimization problem” $\min_{\beta} c^T \beta + R(\beta)$ is easy to compute. The generalized Frank-Wolfe update then is:

$$\tilde{\beta}^i \in \arg \min_{\beta} \{\nabla f(\beta^i)^T \beta + R(\beta)\} \quad \text{and} \quad \beta^{i+1} = (1 - \alpha_i)\beta^i + \alpha_i\tilde{\beta}^i, \quad (7)$$

and notice that we recover the regular Frank-Wolfe update in the special case when $R(\cdot)$ is the indicator function $\mathbf{I}_Q(\cdot)$ of a feasible region Q , see [4] and [45] for a more detailed discussion on generalized Frank-Wolfe methods.

1.2 Related literature

Stochastic Frank-Wolfe methods. There have been several lines of research that investigate and develop stochastic Frank-Wolfe methods. Table 1 presents a summary comparison of the computational complexity of the most relevant stochastic Frank-Wolfe methods that we are aware of. The original Frank-Wolfe (FW) method [11] is a deterministic method. With an appropriate chosen step-size sequence, the method requires $O(\frac{1}{\varepsilon})$ iterations to attain ε -optimality; furthermore each iteration needs to make one exact gradient call and one linear optimization oracle call. A straightforward stochastic Frank-Wolfe (SFW) method randomly chooses an index j_i at iteration i and then computes and uses $\hat{l}_{j_i}(x_{j_i}^T \beta)x_{j_i}$ as an unbiased estimate of the full gradient $\frac{1}{n} \sum_{j=1}^n \hat{l}_j(x_j^T \beta)x_j$ and then uses this estimate in an otherwise standard Frank-Wolfe method. Hazan and Luo [18] showed

Table 1: Summary comparison of computational complexity of recent stochastic Frank-Wolfe methods to achieve an absolute ε -optimal solution.

Algorithm and Reference	Number of Exact Gradient Calls	Number of Stochastic Gradient Calls	Number of Linear Optimization Oracle Calls
FW, Frank and Wolfe [11]	$O(\frac{1}{\varepsilon})$	0	$O(\frac{1}{\varepsilon})$
SFW, Hazan and Luo [18]	0	$O(\frac{1}{\varepsilon^3})$	$O(\frac{1}{\varepsilon})$
Online-FW, Hazan and Kale [17]	0	$O(\frac{1}{\varepsilon^4})$	$O(\frac{1}{\varepsilon^4})$
SCGS, Lan and Zhou [24]	0	$O(\frac{1}{\varepsilon^2})$	$O(\frac{1}{\varepsilon})$
SVRFW, Hazan and Luo [18]	$O(\ln \frac{1}{\varepsilon})$	$O(\frac{1}{\varepsilon^2})$	$O(\frac{1}{\varepsilon})$
STORC, Hazan and Luo [18]	$O(\ln \frac{1}{\varepsilon})$	$O(\frac{1}{\varepsilon^{1.5}})$	$O(\frac{1}{\varepsilon})$
GSFW, this work	1	$O(\frac{1}{\varepsilon})$	$O(\frac{1}{\varepsilon})$

that this method requires $O(\frac{1}{\varepsilon^3})$ stochastic gradient calls and $O(\frac{1}{\varepsilon})$ linear optimization oracle calls to compute an ε -optimal solution. Hazan and Kale [17] proposed an online Frank-Wolfe (Online-FW) method, which requires $O(\frac{1}{\varepsilon^4})$ stochastic gradient calls and $O(\frac{1}{\varepsilon^4})$ linear optimization oracle calls. Lan and Zhou [24] proposed a new technique – the Stochastic Conditional Gradient Sliding (SCGS) – which combines Nesterov’s acceleration techniques and the Frank-Wolfe method, and requires $O(\frac{1}{\varepsilon^2})$ stochastic gradient calls and $O(\frac{1}{\varepsilon})$ linear optimization oracle calls. In [18], Hazan and Luo developed two different types of stochastic Frank-Wolfe methods by utilizing a variance-reduction technique, namely Stochastic Variance-Reduced Frank-Wolfe (SVRFW) and STOchastic variance-Reduced Conditional gradient sliding (STORC). Both methods require $O(\frac{1}{\varepsilon})$ linear optimization oracle calls. While SVRF requires $O(\frac{1}{\varepsilon^2})$ stochastic gradient calls, STORC requires $O(\frac{1}{\varepsilon^{1.5}})$ stochastic gradient calls. In the present work, which we call GSFW in Table 1, we show that Algorithm 1 requires $O(\frac{1}{\varepsilon})$ stochastic gradient calls and $O(\frac{1}{\varepsilon})$ linear optimization oracle calls, which is the same as for deterministic Frank-Wolfe. (Just before submission, we learned of a concurrent paper of Shah [38] which also claims the same computational complexity. However, despite reasonable efforts, we are not able to verify the proofs in that paper.)

Randomized Dual Coordinate Descent. Dual coordinate descent methods have been widely used in statistical and machine learning applications. For example, sequence minimization optimization (SMO) (a variant of dual greedy coordinate descent) is known as one of the best solvers for kernel SVMs [21] and is implemented in LIBSVM [8]. Randomized dual coordinate descent for solving the P was first proposed in [40]. There are many follow-up works on randomized dual coordinate descent, for example, accelerated proximal randomized dual coordinate, see [41], [25], using a non-uniform distribution to choose the coordinate [34], and a primal-dual coordinate method [47], among others. All of these dual methods (or primal-dual methods) require the regularizer $R(\cdot)$ to be a strongly convex function or require adding a dummy strongly convex regularizer to the objective function. In the standard Frank-Wolfe set-up, $R(\cdot)$ is an indicator function and so is not strongly convex. (We will further discuss the connections and differences between the above methods and our method in Appendix A.4.) Another issue for dual coordinate methods is that even though one can rewrite the dual coordinate method entirely in the primal space [39], there are still explicit dual variables which lack intuition or interpretation in the primal space. In contrast, we show here in Lemma 2.1 that Algorithm 1 can be interpreted in the optional dual variables as a randomized coordinate mirror descent algorithm in these variables. One can also apply the analysis

of randomized coordinate descent algorithms in [36] [32] directly to the dual problem (8), but that only generates a dual convergence guarantee and is insufficient if one is interested in the primal problem. In contrast, here we will show how a primal (and/or dual) first-order method naturally implies a primal-dual guarantee without any strong convexity assumptions. Moreover, when $R(\cdot)$ is not strongly convex, the objective function in the dual problem (8) is not differentiable, which is outside of the standard set-up for randomized coordinate descent [36] [32]. But as a byproduct of our analysis, we will present convergence guarantees for randomized coordinate descent for a non-differentiable function, see Appendix A.5.

Variance Reduction Techniques for Stochastic Optimization. There have been many recent algorithmic developments designed to directly tackle the optimization problem P. In order to obtain improved convergence guarantees over the standard Stochastic Gradient Descent (SGD) method, variance reduction techniques have been proposed and extensively studied in recent years. SAG [37] is the first variance reduction method in the literature that we are aware of. In contrast to the sublinear convergence rate of SGD, SAG and several concurrent and/or subsequent works – such as SVRG [22], MISO [30], and SAGA [9] – obtain linear convergence when the objective function is both smooth and strongly convex. Variance reduction techniques can also be applied to non-strongly convex optimization [37], [30], [9], [2], which leads to improved convergence guarantees as well. More recently, Allen-Zhu [1] has proposed an accelerated stochastic method for directly solving P. We recommend [1] for a more detailed discussion on variance reduction techniques overall. It is also worth mentioning that the dual coordinate method [40] also corresponds to a variant of a variance reduction technique in the primal space [39].

1.3 Contributions

Algorithm 1, which we call the Generalized Stochastic Frank-Wolfe (GSFW) method, and its analysis in Lemma 2.1, Theorems 3.1 and 3.2, makes the following contributions to the research on first-order methods for solving loss minimization problems in statistical and machine learning:

1. GSFW is a new primal stochastic Frank-Wolfe method that improves on the computational complexity of stochastic Frank-Wolfe methods – including the new variance reduction methods; and indeed its complexity is on par with that of deterministic Frank-Wolfe. GSFW requires $O(\frac{1}{\varepsilon})$ stochastic gradient oracle calls and $O(\frac{1}{\varepsilon})$ linear optimization oracle calls to compute an absolute ε -optimal solution of P (Theorem 3.1); and in the case when $R(\cdot)$ is strongly convex GSFW requires $O(\ln(\frac{1}{\varepsilon}))$ stochastic gradient oracle calls and $O(\ln(\frac{1}{\varepsilon}))$ linear optimization oracle calls to compute an absolute ε -optimal solution of P (Theorem 3.2).
2. We show that GSFW is equivalent to a randomized coordinate mirror descent algorithm applied to the dual problem (Algorithm 2), which in turn provides a natural interpretation of a dual coordinate descent method in the primal space. This is discussed in Section 2.
3. We show that first-order methods inherently minimize the primal-dual gap with no need for extra conditions. In particular, we show that randomized coordinate mirror descent for the dual problem does not require $R(\cdot)$ to be strongly convex, in contrast with the current literature in this context.
4. As a byproduct, we present a convergence bound for randomized coordinate mirror descent

for minimizing non-smooth functions. This is shown in Section A.5.

1.4 Notation

We use e_j to denote the j^{th} unit coordinate vector in $\mathbb{R}^p$. The ℓ_p norm is denoted $\|\cdot\|_p$. We use $\dot{l}_j(\cdot)$ to denote the first derivative of $l_j(\cdot)$. The Bregman distance function associated with a convex function $h(\cdot)$ is defined as $D_h(y, x) := h(y) - h(x) - \nabla h(x)^T(y - x)$. We use $\mathbb{E}$ to denote expectation and $\mathbb{E}_{j_i}$ to denote expectation conditional on the randomly chosen index j_i . For indicator functions, we use $\mathbf{I}_Q(\cdot)$ to denote the indicator function for the set Q , namely $\mathbf{I}_Q(\beta) := 0$ if $\beta \in Q$, and $\mathbf{I}_Q(\beta) := +\infty$ otherwise; and we use $\mathbf{I}_{\{\text{constraint}\}}(\beta)$ to denote the indicator function of a particular constraint (or condition), namely $\mathbf{I}_{\{\text{constraint}\}}(\beta) := 0$ if the constraint is true at β , and $\mathbf{I}_{\{\text{constraint}\}}(\beta) := +\infty$ otherwise. In a slight abuse of terminology we refer to the “subgradient” of a concave function when it is perhaps more technically accurate to refer to this as a sup-gradient. A differentiable function $f(\cdot)$ is μ -strongly convex with respect to a norm $\|\cdot\|$ if it holds that $f(y) \geq f(x) + \nabla f(x)^T(y - x) + \frac{\mu}{2}\|y - x\|^2$ for all $x, y \in \text{dom} f(\cdot)$. A differentiable function $f(\cdot)$ is μ -strongly convex with respect to a reference function $h(\cdot)$ if it holds that $f(y) \geq f(x) + \nabla f(x)^T(y - x) + \mu D_h(y, x)$ for all $x, y \in \text{dom} f(\cdot)$.

2 Dual problem, and equivalence of Algorithm 1 in the dual with Randomized Coordinate Mirror Descent

Recall the definition of the conjugate of a function $f(\cdot)$:

$$f^*(y) := \sup_{x \in \text{dom} f(\cdot)} \{y^T x - f(x)\} .$$

We will also be interested in the following dual problem of (1) that is constructed using the conjugate functions of the component functions of (1):

$$\text{D :} \quad \max_w D(w) := -R^* \left(-\frac{1}{n} X^T w \right) - \frac{1}{n} \sum_{j=1}^n l_j^*(w_j) . \quad (8)$$

Notice that we can write:

$$R^* \left(-\frac{1}{n} X^T w \right) = - \min_{\beta} \left\{ \frac{1}{n} w^T X \beta + R(\beta) \right\} . \quad (9)$$

Also, defining the convex/concave saddle-function $\phi(\cdot, \cdot)$:

$$\phi(\beta, w) := \frac{1}{n} w^T X \beta - \frac{1}{n} \sum_{i=1}^n l_i^*(w_i) + R(\beta) , \quad (10)$$

we can write P and D in saddlepoint minimax format as:

$$\text{P :} \quad \min_{\beta} \max_w \phi(\beta, w) \quad \text{and} \quad \text{D :} \quad \max_w \min_{\beta} \phi(\beta, w) . \quad (11)$$

Another standard first-order method for convex optimization is the mirror descent algorithm (also called primal gradient method with Bregman distance) [43], [27], [26], [6], which we now briefly review in the context of solving the dual problem D in (8), which is a concave maximization problem. The Bregman distance of a differentiable “prox” function $h(\cdot)$ is defined to be:

$$D_h(w_1, w_2) := h(w_1) - h(w_2) - \langle \nabla h(w_2), w_1 - w_2 \rangle.$$

The (deterministic) mirror descent algorithm for solving D has the following update:

$$w^{i+1} \leftarrow \arg \min_w \{-g(w^i)^T(w - w^i) + \eta_i D_h(w, w^i)\},$$

where $g(\cdot)$ is a subgradient of the objective function $D(\cdot)$ at w (which we call a subgradient even though $D(\cdot)$ is concave), and $\{\eta_i\}$ is the step-size sequence. It is shown in Bach [4] that the generalized Frank-Wolfe method for the primal (1) is equivalent to mirror descent algorithm for the dual (8).

Algorithm 2 presents a Randomized Coordinate Mirror Descent method applied to solve the dual problem D. The algorithm uses the average of the conjugate functions $l_i^*(\cdot)$ as the prox function, namely $h(\cdot) = \frac{1}{n} \sum_{i=1}^n l_i^*(w_i)$, and it initializes the dual variable w^0 to be the prox-center (which is the point that minimizes the prox function). At the start of the i^{th} iteration, the algorithm randomly chooses a coordinate j_i and computes the j_i^{th} coordinate of a subgradient of the dual objective function $D(w)$ at $w = w^i$, since indeed it is straightforward to verify that $\frac{1}{n}(X\tilde{\beta}^i - \nabla L^*(w^i))$ is a subgradient of $D(w)$ at $w = w^i$. The algorithm then performs a coordinate mirror descent step to update the dual variable w^i . Last of all – and optionally since it does not affect future computations – the algorithm updates the primal variable $\tilde{\beta}^i$ in order to compute a primal-dual optimality gap certificate.

Algorithm 2 Randomized Coordinate Mirror Descent applied to the dual problem (8)

Initialize. Define the prox function $h(w) := \frac{1}{n} \sum_{i=1}^n l_i^*(w_i)$. Initialize with $w^0 = \arg \min_w \frac{1}{n} \sum_{i=1}^n l_i^*(w_i)$ and step-size sequences $\{\alpha_i\} \in (0, 1]$ and $\{\eta_i\} \in (0, 1]$. (Optional: set $\tilde{\beta}^{-1} = 0$.)

For iterations $i = 0, 1, \dots$

Compute Randomized Coordinate of Subgradient of $D(\cdot)$ at w^i

Compute $\tilde{\beta}^i \in \arg \min_{\beta} \left\{ \left(\frac{1}{n} (w^i)^T X \beta + R(\beta) \right) \right\}$

Choose random index. Choose $j_i \in \mathcal{U}[1, \dots, n]$

Compute subgradient coordinate vector: $\tilde{g}^i \leftarrow \frac{1}{n} \left(x_{j_i}^T \tilde{\beta}^i - l_{j_i}^*(w_{j_i}^i) \right) e_{j_i}$

Update dual variable: Compute $w^{i+1} = \arg \min_w \left\{ \langle -\eta_i \tilde{g}^i, w - w^i \rangle + D_h(w, w^i) \right\}$

(Optional Accounting:) $\tilde{\beta}^i \leftarrow (1 - \alpha_i) \tilde{\beta}^{i-1} + \alpha_i \tilde{\beta}^i$.

The main result of this section is the following lemma concerning the equivalence of Algorithm 1 and Algorithm 2.

Lemma 2.1. (Equivalence Lemma) *Algorithm 1 and Algorithm 2 are equivalent as follows: the iterate sequence of either algorithm exactly corresponds to an iterate sequences of the other.* $\square$

As a means to proving the lemma, we first reinterpret the update of w_{j_i} at iteration i of Algorithm 2 in the following proposition:

Proposition 2.1. *At iteration i of Algorithm 2 it holds that:*

- (1.) $\dot{l}_{j_i}^*(w_{j_i}^{i+1}) = (1 - \eta_i)\dot{l}_{j_i}^*(w_{j_i}^i) + \eta_i x_{j_i}^T \tilde{\beta}^i$, and
- (2.) $w_{j_i}^{i+1} = \dot{l}_{j_i} \left((1 - \eta_i)\dot{l}_{j_i}^*(w_{j_i}^i) + \eta_i x_{j_i}^T \tilde{\beta}^i \right)$.

Proof: Because $h(w)$ is a coordinate-wise separable function, we can rewrite the update for $w_{j_i}^{i+1}$ as

$$\begin{aligned} w_{j_i}^{i+1} &= \arg \min_{w_{j_i}} \left\langle -\frac{\eta_i}{n} \left(x_{j_i}^T \tilde{\beta}^i - \dot{l}_{j_i}^*(w_{j_i}^i) \right), w_{j_i} \right\rangle + D_{\frac{1}{n} \dot{l}_{j_i}^*}(w_{j_i}, w_{j_i}^i) \\ &= \arg \min_{w_{j_i}} \left\langle -\eta_i \left(x_{j_i}^T \tilde{\beta}^i - \dot{l}_{j_i}^*(w_{j_i}^i) \right), w_{j_i} \right\rangle + D_{\dot{l}_{j_i}^*}(w_{j_i}, w_{j_i}^i) \\ &= \arg \min_{w_{j_i}} \left\langle -\eta_i x_{j_i}^T \tilde{\beta}^i - (1 - \eta_i)\dot{l}_{j_i}^*(w_{j_i}^i), w_{j_i} \right\rangle + \dot{l}_{j_i}^*(w_{j_i}^i) . \end{aligned}$$

From the first-order optimality condition of the above 1-dimensional problem we have $\dot{l}_{j_i}^*(w_{j_i}^{i+1}) = \eta_i x_{j_i}^T \tilde{\beta}^i + (1 - \eta_i)\dot{l}_{j_i}^*(w_{j_i}^i)$, which shows (1.); and (2.) follows directly from (1.) by the properties of the conjugate function in Proposition A.1. $\square$

Proof of Lemma 2.1 We show that the iterate sequence of Algorithm 2 corresponds exactly to an iterate sequence of Algorithm 1. The $\{s^i\}$ sequence is not formally defined in Algorithm 2, so let us define $s^i := \nabla L^*(w^i)$ for all $i = 0, 1, \dots$, which is consistent through conjugacy with the relationship $w^i = \nabla L(s^i)$ in the Optional Accounting step of Algorithm 1 (see Proposition A.1). In order to show the correspondence we proceed by induction on the iteration counter i . For $i = 0$ we have from conjugacy that $s^0 := \nabla L^*(w^0) = 0$ from the definition w^0 in the initialization of Algorithm 2. We also need to show that $\tilde{\beta}^0$ is a solution to the linear optimization oracle problem in Algorithm 1. We have for all $i = 0, \dots$, that:

$$\begin{aligned} \tilde{\beta}^i \in \arg \min_{\beta} \left\{ \frac{1}{n} (w^i)^T X \beta + R(\beta) \right\} &= \arg \min_{\beta} \left\{ \frac{1}{n} (\nabla L(s^i))^T X \beta + R(\beta) \right\} \\ &= \arg \min_{\beta} \left\{ (d^i)^T \beta + R(\beta) \right\} , \end{aligned}$$

thus showing that β^i corresponds to a linear optimization oracle solution at iteration i in Algorithm 1 for all $i = 0, \dots$. Now suppose that the correspondence holds for some iteration counter i , and let us examine $s^{i+1} := \nabla L^*(w^{i+1})$. We have from Proposition 2.1 that:

$$w_{j_i}^{i+1} = \dot{l}_{j_i} \left((1 - \eta_i)\dot{l}_{j_i}^*(w_{j_i}^i) + \eta_i x_{j_i}^T \tilde{\beta}^i \right) = \dot{l}_{j_i} \left((1 - \eta_i)s_{j_i}^i + \eta_i x_{j_i}^T \tilde{\beta}^i \right) , \quad (12)$$

where the first equality is from Proposition 2.1 and the second equality uses induction. This then implies that

$$(1 - \eta_i)s_{j_i}^i + \eta_i x_{j_i}^T \tilde{\beta}^i = \dot{l}(w_{j_i}^{i+1}) = s^{i+1} .$$

And for all coefficient indices $j \neq i$ we have

$$s_j^{i+1} = \dot{l}^*(w_j^{i+1}) = \dot{l}(w_j^i) = s^i ,$$

where the second equality follows from conjugacy, whereby s^{i+1} satisfies the update rule as stated in Algorithm 1, thus demonstrating that the iterate sequence of Algorithm 2 corresponds exactly to an iterate sequence of Algorithm 1. The same type of analysis as above can be used to prove that the iterate sequence of Algorithm 1 corresponds exactly to an iterate sequence of Algorithm 2. $\square$

3 Convergence Guarantees

In this section we develop computational guarantees for Algorithm 2, which automatically provide computational guarantees for Algorithm 1 due to the equivalence shown in Theorem 2.1. Our first – and main – result is Theorem 3.1, which is an expected $O(1/k)$ guaranteed decrease in the duality gap between P and D. Secondly, in the case when $R(\cdot)$ is a strongly convex function, we present a linear convergence result on the duality gap in Theorem 3.2. We start by defining two measures – M and $D_{\max}$ – associated with P and whose values will enter our computational bounds.

Let $M := \max_{\beta \in \text{dom}R(\cdot)} \max_{j=1, \dots, n} \{|x_j^T \beta|\}$, and note that $M < +\infty$ since $\text{dom}R(\cdot)$ is bounded by Assumption 1.1.

Let $\mathcal{W} \subset \mathbb{R}^n$ be the set of “optimal w responses” to values $\beta \in \text{dom}R(\cdot)$ in the saddle-function $\phi(\beta, w)$, namely:

$$\mathcal{W} := \{\hat{w} \in \mathbb{R}^n : \hat{w} \in \arg \max_w \phi(\hat{\beta}, w) \text{ for some } \hat{\beta} \in \text{dom}R(\cdot)\} ,$$

and let $D_{\max}$ be any upper bound on $D_h(\hat{w}, w^0)$ as $\hat{w}$ ranges over all values in $\mathcal{W}$, so that

$$D_h(\hat{w}, w^0) \leq D_{\max} \text{ for all } \hat{w} \in \mathcal{W} .$$

Note at the moment that there is no guarantee that $D_{\max} < +\infty$, but this will be remedied below in Proposition 3.4.

Remark 3.1. A suitable value of $D_{\max}$ can often be easily derived based on the structure of $l_j(\cdot)$. For example, in logistic regression where the loss function is $l_j(s_j) := \log(1 + \exp(-y_j s_j))$ for the given label $y_j \in \{-1, 1\}$, we have $l_j^*(w_j) = -y_j w_j \ln(-y_j w_j) + (1 + y_j w_j) \ln(1 + y_j w_j)$ with $\text{dom}l_j^*(\cdot) = \{w_j : 0 \leq -y_j w_j \leq 1\}$ (where a $\ln(a) := 0$ for $a = 0$). Therefore for all $\hat{w} \in \mathcal{W}$ it holds that

$$D_h(\hat{w}, w^0) \leq \max_{0 \leq -Y\hat{w} \leq e} D_h(w, w^0) = \frac{1}{n} \left(\max_{0 \leq -Y\hat{w} \leq e} L^*(w) - L^*(w^0) \right) = \ln(2) ,$$

where Y is the diagonal matrix whose diagonal coefficients correspond to y and $e = [1, \dots, 1]^T$, so we may set $D_{\max} = \ln(2)$.

Notice in Algorithm 1 and Algorithm 2 that j_i is a random variable; and that s^i , d^i , w^i , etc., are random variables that depend on all previous random variable values $j_0, j_1, \dots, j_{i-1}$, and we denote this string of random variables by

$$\xi_i = \{j_0, j_1, \dots, j_{i-1}\} .$$

We now state our main computational guarantee for Algorithm 2 (and hence for Algorithm 1 as well).

Theorem 3.1. *Consider the Stochastic Generalized Frank-Wolfe method (Algorithm 1) or the Randomized Dual Coordinate Mirror Descent method (Algorithm 2), with step-size sequences $\alpha_i = \frac{2(2n+i)}{(i+1)(4n+i)}$ and $\eta_i = \frac{2n}{2n+i+1}$ for $i = 0, 1, \dots$. Denote*

$$\bar{w}^k = \frac{2}{(4n+k)(k+1)} \sum_{i=0}^k (2n+i)w^i.$$

Under Assumption 1.1, it holds for all $k \geq 0$ that

$$\mathbb{E}_{\xi_k} \left[P(\bar{\beta}^k) - D(\bar{w}^k) \right] \leq \frac{8n\gamma M^2}{(4n+k)} + \frac{2n(2n-1)D_{\max}}{(4n+k)(k+1)} \leq \frac{8n\gamma M^2}{(4n+k)} + \frac{2n(2n-1)\gamma M^2}{(4n+k)(k+1)}. \quad \square$$

Remark 3.2. *Algorithm 1 and Algorithm 2 as well as their analysis can be directly extended to the mini-batch setting. The only difference would be that each index j_i in the statement of Algorithm 1 and Algorithm 2 would be replaced by a random subset of the indices; and in the analysis one analyzes a mini-batch of samples instead of a single sample. Moreover, the updates of the two Algorithms in the mini-batch setting can be implemented in parallel as a result of the separability of samples in Algorithm 1 and of coordinates in Algorithm 2.*

The following string of propositions will be needed for the proof of Theorem 3.1.

Proposition 3.1. *For all iterates i and any $j \in \{1, \dots, n\}$ it holds that $\left| \dot{l}_j^*(w_j^i) \right| \leq M$.*

Proof. We prove this by induction on i . The proposition is true for $i = 0$ because $\dot{l}_j^*(w_j^0) = 0$ for all j by the definition of w^0 . Next suppose that $\left| \dot{l}_j^*(w_j^i) \right| \leq M$ for a given iterate i and for all $j = 1, \dots, n$. Then at iteration $i+1$ and any $j \neq j_i$ we have $w_j^{i+1} = w_j^i$, whereby $\left| \dot{l}_j^*(w_j^{i+1}) \right| = \left| \dot{l}_j^*(w_j^i) \right| \leq M$. And it follows from Proposition 2.1 that

$$\left| \dot{l}_{j_i}^*(w_{j_i}^{i+1}) \right| = \left| (1 - \eta_i) \dot{l}_{j_i}^*(w_{j_i}^i) + \eta_i x_{j_i}^T \tilde{\beta}^i \right| \leq (1 - \eta_i)M + \eta_i M = M,$$

and therefore for any $j = 1, \dots, n$, we have $\left| \dot{l}_j^*(w_j^{i+1}) \right| \leq M$, which completes the proof by induction. $\square$

As a simple corollary we obtain an upper bound on $\|\tilde{g}^i\|_2$ as follows:

Corollary 3.1. $\|\tilde{g}^i\|_2 = \frac{1}{n} \left| x_{j_i}^T \tilde{\beta}^i - \dot{l}_{j_i}^*(w_{j_i}^i) \right| \leq \frac{2M}{n}$.

Proposition 3.2. $h(\cdot)$ is $\frac{1}{n\gamma}$ -strongly convex with respect to the norm $\|\cdot\|_2$.

Proof. Recall that $h(w) = \frac{1}{n} \sum_{j=1}^n l_j^*(w_j)$. It follows from Assumption 1.1 and Proposition A.1 that $\dot{l}_j^*(\cdot)$ is $\frac{1}{\gamma}$ -strongly convex. Therefore for any $w^1, w^2 \in \text{dom}h(\cdot)$ it holds that:

$$\begin{aligned} h(w^1) &= \frac{1}{n} \sum_{j=1}^n l_j^*(w_j^1) \\ &\geq \frac{1}{n} \sum_{j=1}^n \left(l_j^*(w_j^2) + \dot{l}_j^*(w_j^2)(w_j^1 - w_j^2) + \frac{1}{2\gamma} |w_j^2 - w_j^1|^2 \right) \\ &= h(w^2) + \langle \nabla h(w^2), w^1 - w^2 \rangle + \frac{1}{2n\gamma} \|w^2 - w^1\|_2^2. \quad \square \end{aligned}$$

Proposition 3.3. $\phi(\tilde{\beta}^i, w) = D(w^i) + \left\langle \nabla_w \phi(\tilde{\beta}^i, w^i), w - w^i \right\rangle - D_h(w, w^i)$.

Proof. The proof follows from straightforward substitution using $\phi(\tilde{\beta}^i, w) = \frac{1}{n} \left(w^T X \tilde{\beta}^i - \sum_{j=1}^n l_j^*(w_j) \right) + R(\tilde{\beta}^i)$ and noticing from the construction of $\tilde{\beta}^i$ that $D(w^i) = \phi(\tilde{\beta}^i, w^i)$. $\square$

We have the following proposition which establishes an upper bound on $D_{\max}$:

Proposition 3.4. *Under Assumption 1.1 it holds that $D_{\max} \leq \gamma M^2$.*

Before proving this proposition, we first show that there is a natural boundedness constraint for the dual problem:

Proposition 3.5. *Let $T := \{w \in \mathbb{R}^n : \|w - w^0\|_\infty \leq \gamma M\}$. Then:*

1. *for any $\hat{\beta} \in \text{dom} R(\cdot)$ it holds that $\arg \max_w \phi(\hat{\beta}, w) \in T$, and*
2. *for all w^i generated in Algorithm 2, it holds that $w^i \in T$.*

Proof. We first prove (1.). Notice that $w_0 = \nabla L(0)$ and $\arg \max_w \phi(\hat{\beta}, w) = \nabla L(X\hat{\beta})$ (from conjugacy via Proposition A.1), whereby the γ -smoothness of $l_j(\cdot)$ implies that

$$\left\| \arg \max_w \phi(\hat{\beta}, w) - w_0 \right\|_\infty = \left\| \nabla L(X\hat{\beta}) - \nabla L(0) \right\|_\infty = \max_j \left| \dot{l}_j(x_j^T \hat{\beta}) - \dot{l}_j(0) \right| \leq \gamma \max_j |x_j^T \hat{\beta}| \leq \gamma M,$$

which proves (1.). It follows from Proposition 3.1 that for any coordinate j and iterate i it holds that $\left| \dot{l}_j^*(w_j^i) \right| \leq M$. Together with $\dot{l}_j^*(w_j^0) = 0$, we have

$$\frac{1}{\gamma} |w_j^i - w_j^0| \leq \left| \dot{l}_j^*(w_j^i) - \dot{l}_j^*(w_j^0) \right| \leq M,$$

(where the first inequality is from the $\frac{1}{\gamma}$ -strong convexity of $l_j^*(w_j)$), from which it follows that $\|w^i - w^0\|_\infty \leq \gamma M$, which proves (2.). $\square$

Proof of Proposition 3.4: Let $L(s) := \sum_{j=1}^n l_j(s_j)$ and $L^*(w) := \sum_{j=1}^n l_j^*(w_j)$, and note that $L(\cdot)$ and $L^*(\cdot)$ are a conjugate pair. Let $\hat{w} \in \mathcal{W}$ and let $\hat{\beta}$ be such that $\hat{w} \in \arg \max_w \phi(\hat{\beta}, w)$. Then

$$\begin{aligned}
D_h(\hat{w}, w^0) &= \frac{1}{n} (L^*(\hat{w}) - L^*(w^0)) \\
&= \frac{1}{n} \left((\hat{w})^T X \hat{\beta} - L(X \hat{\beta}) - L^*(w^0) \right) \\
&\leq \frac{1}{n} \left(\max_{w \in T, \beta \in \text{dom} R(\cdot)} \{w^T X \beta - L(X \beta)\} - L^*(w^0) \right) \\
&= \frac{1}{n} \left(\max_{w \in T, \beta \in \text{dom} R(\cdot)} \left\{ (w - w^0)^T X \beta + (w^0)^T X \beta - L(X \beta) \right\} - L^*(w^0) \right) \\
&\leq \frac{1}{n} \left(\max_{w \in T, \beta \in \text{dom} R(\cdot)} \{ (w - w^0)^T X \beta \} + \max_{\beta \in \text{dom} R(\cdot)} \{ (w^0)^T X \beta - L(X \beta) \} - L^*(w^0) \right) \\
&\leq \frac{1}{n} \left(n \max_{w \in T, \beta \in \text{dom} R(\cdot)} \|w - w^0\|_\infty \|X \beta\|_\infty + L^*(w^0) - L^*(w^0) \right) \\
&\leq \gamma M^2,
\end{aligned}$$

where the second equality follows from Proposition A.1, the first inequality uses $\hat{\beta} \in \text{dom} R(\cdot)$ and $\hat{w} \in T$ (from Proposition 3.5), and the last inequality uses $\max_{\beta \in \text{dom} R(\cdot)} \|X \beta\|_\infty \leq M$. $\square$

Proposition 3.6. Consider the series $\{\alpha_i\}$ defined by $\alpha_i = \frac{2(2n+i)}{(4n+i)(i+1)}$ for $i \geq 0$ and define the series $\{\bar{\beta}^i\}$ by $\bar{\beta}^{-1} = 0$ and $\bar{\beta}^i = (1 - \alpha_i)\bar{\beta}^{i-1} + \alpha_i \tilde{\beta}^i$ for $i \geq 0$. Also define $\gamma_i = 2n + i$ for $i \geq 0$. Then

$$\bar{\beta}^k = \frac{\sum_{i=0}^k \gamma_i}{\sum_{i=0}^k \gamma_i} \text{ for all } k \geq 0.$$

Proof: The proof follows easily by induction and using $\sum_{i=0}^k \gamma_i = \frac{(4n+k)(k+1)}{2}$. $\square$

Proof of Theorem 3.1. Denote $g^i := \frac{1}{n} (X \tilde{\beta}^i - \nabla L^*(w^i))$, whereby g^i is a subgradient of $D(w)$ at w^i , and $\tilde{g}^i$ is an unbiased estimator of g^i up to the scalar n , namely $\mathbb{E}_{j_i}[\tilde{g}^i] = \frac{1}{n} g^i$. Therefore we have for any i and any $w \in \mathcal{W}$ that:

$$\begin{aligned}
\langle -g^i, w - w^i \rangle &= n \mathbb{E}_{j_i} [\langle -\tilde{g}^i, w - w^i \rangle] \\
&\geq n \mathbb{E}_{j_i} \left[\langle -\tilde{g}^i, w^{i+1} - w^i \rangle + \frac{1}{\eta_i} D_h(w^{i+1}, w^i) + \frac{1}{\eta_i} D_h(w, w^{i+1}) - \frac{1}{\eta_i} D_h(w, w^i) \right] \\
&\geq n \mathbb{E}_{j_i} \left[\langle -\tilde{g}^i, w^{i+1} - w^i \rangle + \frac{1}{2n\gamma\eta_i} \|w^{i+1} - w^i\|_2^2 + \frac{1}{\eta_i} D_h(w, w^{i+1}) - \frac{1}{\eta_i} D_h(w, w^i) \right] \\
&\geq n \mathbb{E}_{j_i} \left[-\frac{1}{2} n \gamma \eta_i \|\tilde{g}^i\|_2^2 + \frac{1}{\eta_i} D_h(w, w^{i+1}) - \frac{1}{\eta_i} D_h(w, w^i) \right] \\
&\geq -2\gamma M^2 \eta_i + \frac{n}{\eta_i} \mathbb{E}_{j_i} [D_h(w, w^{i+1})] - \frac{n}{\eta_i} D_h(w, w^i),
\end{aligned} \tag{13}$$

where the first inequality is from the “three point property” of Tseng (Lemma A.1 in the Appendix), the second inequality is due to the fact that $h(w)$ is $\frac{1}{n\gamma}$ -strongly convex with respect to the norm

$\|\cdot\|_2$ (Proposition 3.2), and the third inequality is an application of the Cauchy-Schwarz inequality, and the last inequality uses Corollary 3.1.

On the other hand, we have from Proposition 3.3 that

$$\langle -g^i, w - w^i \rangle = \langle -\nabla_w \phi(\tilde{\beta}^i, w^i), w - w^i \rangle = D(w^i) - \phi(\tilde{\beta}^i, w) - D_h(w, w^i). \quad (14)$$

Combining (13) and (14) and rearranging yields

$$-(\phi(\tilde{\beta}^i, w) - D(w^i)) \geq -2\gamma M^2 \eta_i + \frac{n}{\eta_i} \mathbb{E}_{j_i}[D_h(w, w^{i+1})] - \left(\frac{n}{\eta_i} - 1\right) D_h(w, w^i).$$

Substituting $\eta_i = \frac{2n}{2n+i+1}$ and multiplying by $2n+i$ results, we arrive at the following inequality after rearranging terms:

$$\begin{aligned} & (2n+i)(\phi(\tilde{\beta}^i, w) - D(w^i)) \\ & \leq 4n\gamma M^2 \left(\frac{2n+i}{2n+i+1}\right) + \frac{1}{2} ((2n+i)(2n+i-1)D_h(w, w^i) - (2n+i)(2n+i+1)\mathbb{E}_{j_i}[D_h(w, w^{i+1})]) . \end{aligned}$$

Summing the above inequality for $i = 0, \dots, k$ and recalling from Proposition 3.6 that $\bar{\beta}^k := \frac{2}{(4n+k)(k+1)} \sum_{i=0}^k (2n+i)\tilde{\beta}^i$, and taking the unconditional expectation, we arrive at:

$$\begin{aligned} \frac{(4n+k)(k+1)}{2} \mathbb{E}_{\xi_k}[\phi(\bar{\beta}^k, w) - D(\bar{w}^k)] &= \left(\sum_{i=0}^k 2n+i\right) \mathbb{E}_{\xi_k}[\phi(\bar{\beta}^k, w) - D(\bar{w}^k)] \\ &\leq \mathbb{E}_{\xi_k} \left[\sum_{i=0}^k (2n+i)(\phi(\tilde{\beta}^i, w) - D(w^i)) \right] \\ &\leq 4(k+1)n\gamma M^2 + \frac{1}{2}(2n)(2n-1)D_h(w, w^0) \\ &\leq 4(k+1)n\gamma M^2 + n(2n-1)D_{\max}, \end{aligned}$$

where the first inequality uses the convexity of $\phi(\beta, w)$ over β and the concavity of $D(w)$, the second inequality follows from the summation and canceling terms in the telescoping series, and the third inequality uses $w \in \mathcal{W}$. Choosing $\hat{w} = \arg \max_w \phi(\bar{\beta}^k, w)$, we have $P(\bar{\beta}^k) = \phi(\bar{\beta}^k, \hat{w})$, which yields:

$$\mathbb{E}_{\xi_k}[P(\bar{\beta}^k) - D(\bar{w}^k)] \leq \frac{8n\gamma M^2}{(4n+k)} + \frac{2n(2n-1)D_{\max}}{(4n+k)(k+1)},$$

thus showing the first inequality in the statement of the theorem. The second inequality in the statement of the theorem then follows as a simple application of Proposition 3.4. $\square$

3.1 Linear Convergence when $R(\cdot)$ is Strongly Convex

In this section, we further assume $R(\cdot)$ is a μ -strongly convex function, and we develop a linear convergence guarantee for Algorithms 1 and 2. We first formally define a separable function.

Definition 3.1. The function $h : R^n \rightarrow R$ is separable if

$$h(x) = \sum_{i=1}^n h_i(x_i),$$

where x_i is the i^{th} coordinate of x and h_i is a univariate function.

Next we introduce the notation of relative smoothness and relative strong convexity developed recently in [27],[26],[5],[15]. We adapt a simplified version of the coordinate-wise relative smoothness condition as in [15].

Definition 3.2. $f(\cdot)$ is coordinate-wise σ -smooth relative to a separable reference function $h(\cdot)$ if for any x , scalar t and coordinate j it holds that:

$$f(x + te_j) \leq f(x) + \langle \nabla f(x), te_j \rangle + \sigma D_h(x + te_j, x). \quad (15)$$

We also adapt the notion of relative strong convexity developed in [27].

Definition 3.3. $f(\cdot)$ is μ -strongly convex relative to $h(\cdot)$ if for any x, y , it holds that

$$f(y) \geq f(x) + \langle \nabla f(x), y - x \rangle + \mu D_h(y, x). \quad (16)$$

The next proposition states that the dual function $D(w)$ is both coordinate-wise smooth and strongly concave relative to the reference function $h(w) := \frac{1}{n} \sum_{j=1}^n l_j^*(w_j)$. In the proposition, recall that x_j is the j^{th} row of the matrix X .

Proposition 3.7.

(1.) Suppose $R(\cdot)$ is a μ -strongly convex function with respect to $\|\cdot\|_2$, then $-D(\cdot)$ is coordinate-wise $\left(\frac{\gamma \max_j \|x_j\|_2^2}{n\mu} + 1\right)$ -smooth relative to $h(\cdot)$, and

(2.) $-D(\cdot)$ is 1-strongly convex relative to $h(\cdot)$.

Proof. (1.) Consider w_1 and w_2 such that $w_2 = w_1 + te_j$ for some coordinate j , namely w_1 and w_2 only differ in one coordinate. It follows from Proposition A.1 that $R^*(\cdot)$ is $\frac{1}{\mu}$ -smooth with respect to $\|\cdot\|_2$, thus we have

$$\begin{aligned} R^*\left(-\frac{1}{n}X^T w_2\right) &\leq R^*\left(-\frac{1}{n}X^T w_1\right) + \left\langle \nabla R^*\left(-\frac{1}{n}X^T w_1\right), -\frac{1}{n}X^T(w_2 - w_1) \right\rangle + \frac{1}{2\mu} \left\| \frac{1}{n}X^T(w_2 - w_1) \right\|_2^2 \\ &= R^*\left(-\frac{1}{n}X^T w_1\right) + \left\langle -\frac{1}{n}X \nabla R^*\left(-\frac{1}{n}X^T w_1\right), w_2 - w_1 \right\rangle + \frac{t^2}{2n^2\mu} \|x_j\|_2^2 \\ &\leq R^*\left(-\frac{1}{n}X^T w_1\right) + \left\langle -\frac{1}{n}X \nabla R^*\left(-\frac{1}{n}X^T w_1\right), w_2 - w_1 \right\rangle + \frac{\gamma \|x_j\|_2^2}{n\mu} D_h(w_2, w_1), \end{aligned}$$

where the first inequality follows from smoothness, the equality is from $w_2 = w_1 + te_j$, and the last inequality utilizes the fact that $h(\cdot)$ is $(\frac{1}{n\gamma})$ -strongly convexity with respect to $\|\cdot\|_2$. Therefore it holds that $\hat{f}(w) := R^*\left(-\frac{1}{n}X^T w\right)$ is coordinate-wise $(\frac{\gamma \max_j \|x_j\|_2^2}{n\mu})$ -smooth relative to $h(\cdot)$. The proof is completed by noticing that $-D(w) = R^*\left(-\frac{1}{n}X^T w\right) + h(w)$.

(2.) This follows from the additivity property of relative strong convexity (Proposition 1.2 in [27]), whereby $D(\cdot)$ is 1-strongly concave relative to $h(w)$. $\square$

The following theorem states a linear convergence guarantee in the case when $R(\cdot)$ is strongly convex.

Theorem 3.2. *Suppose $D(\cdot)$ is coordinate-wise σ -smooth relative to $h(\cdot)$. Consider the Stochastic Generalized Frank-Wolfe method (Algorithm 1) or the Randomized Dual Coordinate Mirror Descent method (Algorithm 2), with step-size sequences $\eta_i = \frac{1}{\sigma}$ and $\alpha_i = \frac{n^{-1}\sigma^i}{\sigma^{i+1} - (\sigma-1/n)^{i+1}}$. Under Assumption 1.1 it holds for all $k \geq 1$ that*

$$\mathbb{E}_{\xi_k} \left[P(\bar{\beta}^{k-1}) - D(w^k) \right] \leq \frac{D_{\max}}{\left(1 + \frac{1}{n\sigma-1}\right)^k - 1} \leq \frac{\gamma M^2}{\left(1 + \frac{1}{n\sigma-1}\right)^k - 1}. \quad (17)$$

$\square$

Notice that the first inequality in (17) shows linear convergence; indeed, in this case it holds that

$$\frac{1}{\left(1 + \frac{1}{n\sigma-1}\right)^k - 1} \leq n\sigma \left(1 - \frac{1}{n\sigma}\right)^k. \quad (18)$$

(This inequality holds trivially for $k = 1$, and induction on k establishes the result for $k \geq 2$.) Furthermore, when k is large the -1 term in the denominator of the left-hand side can be ignored which yields the asymptotic bound $(1 - \frac{1}{n\sigma})^k D_{\max}$. The next corollary states the implication of this linear convergence bound in terms of the values γ and μ of the γ -smoothness of $l_1(\cdot), \dots, l_n(\cdot)$ and the μ -strong convexity of $R(\cdot)$.

Corollary 3.2. *Choose $\sigma = \frac{\gamma \max_j \|x_j\|_2^2}{n\mu} + 1$ as per Proposition 3.7. Then Theorem 3.2 and (18) imply*

$$\mathbb{E}_{\xi_k} \left[P(\bar{\beta}^{k-1}) - D(w^k) \right] \leq \frac{D_{\max}}{\left(1 + \frac{1}{\frac{\gamma \max_j \|x_j\|_2^2}{\mu} + n - 1}\right)^k - 1} \leq D_{\max} \left(\frac{\gamma \max_j \|x_j\|_2^2}{\mu} + n \right) \left(1 - \frac{1}{n + \frac{\gamma \max_j \|x_j\|_2^2}{\mu}} \right)^k.$$

Before proving Theorem 3.2, we first present an elementary proposition for a separable reference function $h(\cdot)$, whose proof is given in Appendix A.3.

Proposition 3.8. *Suppose $h(\cdot) : \mathbb{R}^n \rightarrow \mathbb{R}$ is a separable function. Let $j \sim \mathcal{U}[1, \dots, n]$. For given $x, a, y \in \mathbb{R}^n$, define the random variable $b \in \mathbb{R}^n$ such that $b_j = a_j$, and $b_i = x_i$ for all $i \neq j$. Then:*

$$D_h(y, a) - D_h(y, x) = n \mathbb{E}_j (D_h(y, b) - D_h(y, x)). \quad \square$$

We also will use the following proposition whose proof follows easily by induction on k .

Proposition 3.9. Consider the series $\{\alpha_i\}$ defined by $\alpha_i = \frac{n^{-1}\sigma^i}{\sigma^{i+1} - (\sigma-1/n)^{i+1}}$ for $i \geq 0$, and define the series $\{\bar{\beta}^i\}$ by $\bar{\beta}^{-1} = 0$ and $\bar{\beta}^i = (1 - \alpha_i)\bar{\beta}^{i-1} + \alpha_i\tilde{\beta}^i$ for $i \geq 0$. Also define $\gamma_i = \left(\frac{n\sigma}{n\sigma-1}\right)^i$ for $i \geq 0$. Then

$$\bar{\beta}^k = \frac{\sum_{i=0}^k \gamma_i \tilde{\beta}^i}{\sum_{i=0}^k \gamma_i} \text{ for all } k \geq 0.$$

□

Proof of Theorem 3.2.

Notice that $\tilde{g}^i = \nabla_{j_i} D(w^i) e_{j_i}$, and w^{i+1} is a coordinate update from w^i , whereby we have

$$-D(w^{i+1}) \leq -D(w^i) - \langle \tilde{g}^i, w^{i+1} - w^i \rangle + \sigma D_h(w^{i+1}, w^i) \leq -D(w^i),$$

and hence the dual function value sequence $\{D(w^i)\}$ is non-decreasing.

Define $r^{i+1} := \arg \min_w \{\langle -\nabla D(w^i), w - w^i \rangle + \sigma D_h(w, w^i)\}$, then we have

$$\begin{aligned} \mathbb{E}_{j_i}[-D(w^{i+1})] &\leq \mathbb{E}_{j_i}[-D(w^i) - \langle \nabla D(w^i), w^{i+1} - w^i \rangle + \sigma D_h(w^{i+1}, w^i)] \\ &= \mathbb{E}_{j_i}[-D(w^i) - \frac{1}{n} (\langle \nabla D(w^i), r^{i+1} - w^i \rangle + \sigma D_h(r^{i+1}, w^i))] \\ &\leq \mathbb{E}_{j_i}[-D(w^i) - \frac{1}{n} (\langle \nabla D(w^i), w - w^i \rangle + \sigma D_h(w, w^i) - \sigma D_h(w, r^{i+1}))] \\ &= \mathbb{E}_{j_i}[-D(w^i) - \frac{1}{n} \langle \nabla D(w^i), w - w^i \rangle + \sigma D_h(w, w^i) - \sigma D_h(w, w^{i+1})] \\ &= \mathbb{E}_{j_i}[-\frac{n-1}{n} D(w^i) - \frac{1}{n} (D(w^i) + \langle \nabla D(w^i), w - w^i \rangle) + \sigma D_h(w, w^i) - \sigma D_h(w, w^{i+1})], \end{aligned} \tag{19}$$

where the first inequality is from the coordinate-wise σ -smoothness of $D(w)$ relative to $h(w)$ and the fact that w^{i+1} is a coordinate update from w^i , the first equality is due to expectation and the separability of $h(\cdot)$, the second inequality uses the three-point property (Lemma A.1), the second equality uses Proposition 3.8, and the third equality is just arithmetic rearrangement.

Notice that

$$\begin{aligned} &-D(w^i) - \langle \nabla D(w^i), w - w^i \rangle + n\sigma D_h(w, w^i) - n\sigma D_h(w, w^{i+1}) \\ &= -D(w^i) - \langle \nabla_w \phi(\tilde{\beta}^i, w^i), w - w^i \rangle + n\sigma D_h(w, w^i) - n\sigma D_h(w, w^{i+1}) \\ &= -\phi(\tilde{\beta}^i, w) + (n\sigma - 1)D_h(w, w^i) - n\sigma D_h(w, w^{i+1}), \end{aligned}$$

where the last equality utilizes Proposition 3.3. We can then rewrite (19) (after multiplying by n on both sides) as

$$\mathbb{E}_{j_i}[-nD(w^{i+1})] \leq \mathbb{E}_{j_i} \left[-(n-1)D(w^i) - \phi(\tilde{\beta}^i, w) + (n\sigma - 1)D_h(w, w^i) - n\sigma D_h(w, w^{i+1}) \right]. \tag{20}$$

Multiplying (20) by $\left(\frac{n\sigma}{n\sigma-1}\right)^{i+1}$ and summing over $i = 0, \dots, k-1$, we obtain:

$$\begin{aligned}
& \mathbb{E}_{\xi_k} \left[- \sum_{i=1}^k n \left(\frac{n\sigma}{n\sigma-1} \right)^i D(w^i) \right] \\
& \leq \mathbb{E}_{\xi_k} \left[- \sum_{i=1}^k (n-1) \left(\frac{n\sigma}{n\sigma-1} \right)^i D(w^{i-1}) - \sum_{i=1}^k \left(\frac{n\sigma}{n\sigma-1} \right)^i \phi(\tilde{\beta}^{i-1}, w) + n\sigma D_h(w, w^0) \right] \\
& \leq \mathbb{E}_{\xi_k} \left[- \sum_{i=1}^k (n-1) \left(\frac{n\sigma}{n\sigma-1} \right)^i D(w^{i-1}) - \left(\sum_{i=1}^k \left(\frac{n\sigma}{n\sigma-1} \right)^i \right) \phi(\bar{\beta}^{k-1}, w) + n\sigma D_h(w, w^0) \right],
\end{aligned}$$

where the last inequality is from Proposition 3.9 and the convexity of $\phi(\beta, w)$ in β . Since the sequence $\{D(w^i)\}$ is non-decreasing in i it follows that:

$$\mathbb{E}_{\xi_k} \left[- \left(\sum_{i=1}^k \left(\frac{n\sigma}{n\sigma-1} \right)^i \right) D(w^k) \right] \leq \mathbb{E}_{\xi_k} \left[- \left(\sum_{i=1}^k \left(\frac{n\sigma}{n\sigma-1} \right)^i \right) \phi(\bar{\beta}^{k-1}, w) \right] + n\sigma D_h(w, w^0). \quad (21)$$

Let us substitute the following value of w in (21): $w \leftarrow \hat{w}^{k-1} := \arg \max_w \{\phi(\bar{\beta}^{k-1}, w)\}$, which yields:

$$\left(\sum_{i=1}^k \left(\frac{n\sigma}{n\sigma-1} \right)^i \right) \mathbb{E}_{\xi_k} [\phi(\bar{\beta}^{k-1}, \hat{w}^{k-1}) - D(w^k)] \leq n\sigma D_h(\hat{w}^{k-1}, w^0) \leq n\sigma D_{\max},$$

where the last inequality above comes from the definition of $D_{\max}$. Therefore we have

$$\mathbb{E}_{\xi_k} [P(\bar{\beta}^{k-1}) - D(w^k)] \leq \frac{n\sigma}{\left(\sum_{i=1}^k \left(\frac{n\sigma}{n\sigma-1} \right)^i \right)} D_{\max} = \frac{D_{\max}}{\left(1 + \frac{1}{n\sigma-1} \right)^k - 1},$$

which furnishes the proof by utilizing Proposition 3.9. $\square$

Remark 3.3. A natural question to ask next is whether one can achieve an accelerated convergence rate when $R(\cdot)$ is strongly convex, similar to that in [41], [25]. The answer actually is yes, as one can utilize similar proof techniques as those developed in [25]. However, the accelerated version may not have a natural interpretation in the primal variables.

A Appendix

A.1 Properties of Conjugate Functions

Recall the definition of the conjugate of a function $f(\cdot)$:

$$f^*(y) := \sup_{x \in \text{dom } f} \{y^T x - f(x)\}.$$

The following properties of conjugate functions are used in this paper:

Proposition A.1. (see [3], [46], [23]) If $f(\cdot)$ is a closed convex function, then $f^{**}(\cdot) = f(\cdot)$. Furthermore:

1. $f(\cdot)$ is γ -smooth with domain $\mathbb{R}^p$ with respect to the norm $\|\cdot\|$ if and only if $f^*(\cdot)$ is $1/\gamma$ -strongly convex with respect to the (dual) norm $\|\cdot\|^*$.
2. If $f(\cdot)$ is differentiable and strictly convex, then the following three conditions are equivalent:
 - (a) $y = \nabla f(x)$
 - (b) $x = \nabla f^*(y)$, and
 - (c) $x^T y = f(x) + f^*(y)$.

A.2 Three-Point Property

We state here the “three-point property” as memorialized by Tseng [44]:

Lemma A.1. (Three-Point Property [44]) Let $\phi(x)$ be a convex function, and let $D_h(\cdot, \cdot)$ be the Bregman distance for $h(\cdot)$. For a given vector z , let

$$z^+ := \arg \min_{x \in Q} \{\phi(x) + D_h(x, z)\}.$$

Then

$$\phi(x) + D_h(x, z) \geq \phi(z^+) + D_h(z^+, z) + D_h(x, z^+) \quad \text{for all } x \in Q. \quad \square$$

A.3 Proof of Proposition 3.8

Note that

$$\begin{aligned} \langle \nabla h(a) - \nabla h(x), y \rangle &= \sum_{i=1}^n \langle \nabla h_i(a_i) - \nabla h_i(x_i), y_i \rangle \\ &= n \mathbb{E}_j \langle \nabla h_j(a_j) - \nabla h_j(x_j), y_j \rangle \\ &= n \mathbb{E}_j \langle \nabla h_j(b_j) - \nabla h_j(x_j), y_j \rangle \\ &= n \mathbb{E}_j \sum_{i=1}^n \langle \nabla h_i(b_i) - \nabla h_i(x_i), y_i \rangle \\ &= n \mathbb{E}_j \langle \nabla h(b) - \nabla h(x), y \rangle, \end{aligned}$$

where the second equation is from expectation, and the third and fourth equation follow because $b_j = a_j$ and $b_i = x_i$ for all $i \neq j$. Using similar logic it also holds that

$$\langle \nabla h(a), a \rangle - \langle \nabla h(x), x \rangle = n \mathbb{E}_j (\langle \nabla h(b), b \rangle - \langle \nabla h(x), x \rangle),$$

and

$$h(a) - h(x) = n \mathbb{E}_j (h(b) - h(x)).$$

Therefore,

$$\begin{aligned}
D_h(y, a) - D_h(y, x) &= \langle \nabla h(a), y - a \rangle - \langle \nabla h(x), y - x \rangle - (h(a) - h(x)) \\
&= \langle \nabla h(a) - \nabla h(x), y \rangle - (\langle \nabla h(a), a \rangle - \langle \nabla h(x), x \rangle) - (h(a) - h(x)) \\
&= n\mathbb{E}_j [\langle \nabla h(b) - \nabla h(x), y \rangle - (\langle \nabla h(b), b \rangle - \langle \nabla h(x), x \rangle) - (h(b) - h(x))] \\
&= n\mathbb{E}_j [\langle \nabla h(b), y - b \rangle - \langle \nabla h(x), y - x \rangle - (h(b) - h(x))] \\
&= n\mathbb{E}_j [D_h(y, b) - D_h(y, x)] \quad . \quad \square
\end{aligned}$$

A.4 Connections and Comparisons to Stochastic Dual Coordinate Descent Methods

Compared with the stochastic dual coordinate descent methods developed in [40][25][41], our analysis does not require $R(\cdot)$ to be strongly convex. And in the case when $R(\cdot)$ is strongly convex, our coordinate update at each iteration requires the solution of a univariate problem of the following form for a suitably given scalar c_{j_i} :

$$\min_{w_{j_i}} c_{j_i} w_{j_i} + l_{j_i}^*(w_{j_i}) \quad , \quad (22)$$

in comparison with the algorithms in [40] or [25] for which the coordinate update at each iteration requires the solution of the following univariate problem for suitably given scalars c_{j_i} and b_{j_i} :

$$\min_{w_{j_i}} c_{j_i} w_{j_i} + l_{j_i}^*(w_{j_i}) + b_{j_i} w_{j_i}^2 \quad . \quad (23)$$

There are many cases where (22) has a closed-form solution whereas (23) does not, for example when the loss $l_{j_i}(\cdot)$ comes from logistic regression. This makes our method more efficient in practice for these types of loss functions.

The unaccelerated version of [25] is a randomized coordinate method with a composite function, and has the following update :

$$w_{j_i}^{i+1} = \arg \min_{w_{j_i}} \left\{ c^i w_{j_i} + \frac{1}{2\eta} (w_{j_i} - w_{j_i}^i)^2 + \frac{1}{n} l_{j_i}^*(w_{j_i}) \right\} \quad ,$$

where c^i is one coordinate of the gradient. The above update can be viewed as a coordinate mirror descent method update with reference function $h(w) := \frac{1}{2\eta} \|w\|^2 + \frac{1}{n} L^*(w)$. This is similar to the equivalence of composite optimization and mirror descent in the deterministic case discussed in Section 3.3 of [27].

A.5 Regarding Randomized Coordinate Mirror Descent with Non-smooth Functions

Since the seminal work of Nesterov [32], there have been many research results on randomized coordinate descent for minimizing a smooth objective function. However, there has not been

much research on randomized coordinate descent for minimizing a non-smooth objective function; in fact the only work as far as we know is [33] which considers a special non-smooth objective function – the maximum of some convex functions. On the other hand, the analysis in [32] and the many following-up papers (including [36], [28], and others) cannot be applied directly to analyze the non-smooth case so far as we can tell. In our analysis in Theorem 3.1, we use (in the dual) randomized coordinate mirror descent for a non-smooth objective function. Indeed, one can consider a randomized coordinate of a subgradient as an unbiased stochastic subgradient descent (up to a scalar), and then use the analysis of stochastic mirror descent algorithm (developed in [26]) to develop convergence guarantees.

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