

An Optimal High-Order Tensor Method for Convex Optimization

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¹ Abstract

This paper is concerned with finding an *optimal* algorithm for minimizing a composite convex objective function. The basic setting is that the objective is the sum of two convex functions: the first function is smooth with up to the d -th order derivative information available, and the second function is possibly non-smooth, but its proximal tensor mappings can be computed approximately in an efficient manner. The problem is to find – in that setting – the best possible (optimal) iteration complexity for convex optimization. Along that line, for the smooth case (without the second non-smooth part in the objective), [Nesterov \(1983\)](#) proposed an optimal algorithm for the first-order methods ($d = 1$) with iteration complexity $O(1/k^2)$. A high-order tensor algorithm with iteration complexity of $O(1/k^{d+1})$ was proposed by [Baes \(2009\)](#) and [Nesterov \(2018\)](#). In this paper, we propose a new high-order tensor algorithm for the general composite case, with the iteration complexity of $O(1/k^{(3d+1)/2})$, which matches the lower bound for the d -th order methods as established in [Nesterov \(2018\)](#); [Arjevani et al. \(2018\)](#), and hence is optimal. Our approach is based on the *Accelerated Hybrid Proximal Extragradient* (A-HPE) framework proposed in [Monteiro and Svaiter \(2013\)](#), where a bisection procedure is installed for each A-HPE iteration. At each bisection step a proximal tensor subproblem is approximately solved, and the total number of bisection steps per A-HPE iteration is bounded by a logarithmic factor in the precision required.

Keywords: convex optimization; tensor method; acceleration; iteration complexity.

1. Introduction

In this paper, we consider the following composite unconstrained convex optimization:

$$\min_{x \in \mathbb{R}^n} F(x) := f(x) + h(x), \quad (1)$$

where f is differentiable and convex, and h is convex but possibly non-smooth. In this context, we assume that convex tensor (polynomial) proximal mappings regarding h can be approximately computed efficiently. Given that structure, a fundamental question is to find an *optimal* algorithm that solves the above problem, using the available derivative information of the smooth part f .

1. Extended abstract. Full version appears as arXiv reference, arXiv:1812.06557v2.

In case $F(x) = f(x)$, and only the gradient information of f is available, [Nesterov \(1983\)](#) proposed a gradient-type algorithm, which achieves the overall iteration complexity of $O(1/k^2)$, matching the lower bound on the iteration complexity of this class of solution methods, hence is known to be an *optimal* algorithm among all the first-order methods. When the Hessian information is available, [Nesterov \(2008\)](#) proposed an acceleration scheme for cubic regularized Newton’s method, and he showed that the iteration complexity bound improves from $O(1/k^2)$ to $O(1/k^3)$. A few years later, [Monteiro and Svaiter \(2013\)](#) proposed a different acceleration scheme, which they termed as *Accelerated Hybrid Proximal Extragradient Method* (A-HPE) framework, and they proved that if the second-order information is incorporated into the A-HPE framework then the corresponding accelerated Newton proximal extragradient method has a superior iteration complexity bound of $O(1/k^{7/2})$ over $O(1/k^3)$. In 2018, [Arjevani et al. \(2018\)](#) showed that $O(1/k^{7/2})$ is actually a lower bound for the oracle complexity of the second-order methods for convex smooth optimization. This shows that the accelerated Newton proximal extragradient method is an optimal second-order method.

As evidenced by the special cases $d = 1$ and $d = 2$, there is a clear tradeoff between the level of derivative information required and the overall iteration complexity improved. Therefore, a natural and important question arises:

What is the *exact* tradeoff relationship between d and the worst-case iteration complexity?

To answer such question, [Baes \(2009\)](#) extended the accelerated cubic regularized Newton method to the general high-order case with the iteration complexity of $O(1/k^{d+1})$, where d is the order of derivative information used in the algorithm. Similar extension was recently revisited by [Nesterov \(2018\)](#) with a discussion on the efficient implementation of the method when $d = 3$. [Jiang et al. \(2018\)](#) extended Nesterov’s approach to accommodate the composite optimization (1) and relaxed the requirement on the knowledge of problem parameters such as the Lipschitz constants and the requirement on the exact solutions of the subproblems while maintaining the same iteration bound as in [Nesterov \(2018\)](#). Along the line of bounding the worst case iteration complexity using up to the d -th order derivative information, there have also been significant progresses as well. [Arjevani et al. \(2018\)](#) showed that the worst case iteration complexity of any algorithm in that setting cannot be better than $O(1/k^{(3d+1)/2})$. A simplified analysis of the bound can be found in [Nesterov \(2018\)](#). So, there was a gap between the achieved iteration bound $O(1/k^{d+1})$ and the best possible bound of $O(1/k^{(3d+1)/2})$. Clearly at least one of the two bounds is improvable. In this paper, we aim to settle the above theoretical quest by providing a new implementable algorithm whose iteration complexity is precisely $O(1/k^{(3d+1)/2})$. As a result, the tradeoff relationship discussed above is pinned down to be exactly $O(1/k^{(3d+1)/2})$. We note the independent work [Gasnikov et al. \(2018\)](#) first in Russian and then translated to English in late December 2018, and two other independent works [Bubeck et al. \(2018\)](#) and [Bullins \(2018\)](#), which were posted on arxiv in December 2018 as well. They derive similar results to ours with a limitation on the smooth objective functions (or more specifically– quartics in [Bullins \(2018\)](#)), while we allow a composite objective.

Our algorithm is based on the A-HPE framework of [Monteiro and Svaiter \(2013\)](#), and it specifies a way to generate an approximate solution through the use of high order derivative information by Taylor expansion. In each iteration, such approximate solution is computed by means of a bisection process. At each bisection step, a regulated convex tensor (polynomial) optimization subproblem is approximately solved. Moreover, we show that, to implement one A-HPE iteration, the number

of bisection steps – each calling to solve a convex tensor subproblem – is upper bounded by a logarithmic factor in the inverse of the required precision. Our bisection procedure is similar to the one proposed in [Monteiro and Svaiter \(2013\)](#) for the case $d = 2$; however, a key modification is applied which enables the removal of the so-called “bracketing stage” used in [Monteiro and Svaiter \(2013\)](#).

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