

A secant-based Nesterov method for convex functions

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A secant-based Nesterov method for convex functions.

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Abstract A simple secant-based fast gradient method is developed for problems whose objective function is convex and well-defined. The proposed algorithm extends the classical Nesterov gradient method by updating the estimate-sequence parameter with secant information whenever possible. This is achieved by imposing a secant condition on the choice of search point. Furthermore, the proposed algorithm embodies an "update rule with reset" that parallels the restart rule recently suggested in O'Donoghue and Candes (2013). The proposed algorithm applies to a large class of problems including logistic and least-square losses commonly found in the machine learning literature. Numerical results demonstrating the efficiency of the proposed algorithm are analyzed with the aid of performance profiles.

Keywords Convex optimization · Secant Methods · Fast gradient methods · Nesterov gradient method.

1 Introduction

This paper considers the unconstrained optimization of convex function f :

$$UCOP : \quad \min_{x \in \mathbb{R}^n} f(x) \tag{1}$$

where $f : \mathbb{R}^n \rightarrow \mathbb{R}$ is a continuously differentiable convex function. The domain of f , $\mathbf{dom} f$, is the convex set \mathbb{R}^n and x is a real vector. A necessary and sufficient condition for a point x^* to be a minimizer of f is $\nabla f(x^*) = 0$ [7]. Furthermore, it is assumed that f is bounded below and there exists a unique minimizer $x^* : f(x^*) = f^*$ with $f^* \leq f(x) \forall x \in \mathbb{R}^n$ [27].

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The idea of enhancing or accelerating gradient methods directly has been intensively researched [3, 12, 23, 24, 29, 32] since the pioneering works of Shah et. al. [33] and Polyak [31]. Accelerated gradient methods are easy to implement and offer much lower memory requirement as compared to higher-order methods such as Newton's method. Accelerated gradient methods compute future iterates by relying only on the local gradient and a history of past iterates. Accelerated gradient schemes can be thought of as momentum methods, in that the step taken at the current iteration depends on the previous iterations, and where the momentum grows from one iteration to the next [28]. Accelerated gradient methods, unlike gradient-descent methods, are not guaranteed to monotonically decrease the objective value. In other words, accelerated gradient methods are nonmonotone gradient methods that utilize the momentum from the previous iterates.

Particular schemes include the Barzilai-Borwein gradient method [2], the backpropagation method with momentum [5, 35] - a well-known algorithm in the neural network community - and a fast gradient method developed by Nesterov [24]. All the aforementioned accelerated gradient methods use only the previous iterate and as such they can be considered special cases of two-step iterative algorithms,

$$x_{k+1} = x_k - \alpha_k \nabla f(y_k) + \eta_k (x_k - x_{k-1}) \quad : \quad \alpha_k > 0, \eta_k \geq 0; \quad y_k = \sum_{i=0}^{i=k} \tau_i x_i, \tau_i \in \mathbb{R},$$

with appropriate choice of α_k , η_k and y_k . The nonmonotonicity of the accelerated gradient methods are beneficial and contribute to their increased convergence rate [1, 10, 15]. However they are susceptible to severe bumps in the objective values that may be detrimental and may lead to wasted iterations as noted in [28]. This is the case in the Nesterov gradient method when the momentum factor has exceeded a critical value. This can happen when the condition number (i.e. $q^{-1} = \frac{L}{\mu}$, where L , μ are as defined in § 2) is underestimated [28]. Moreover, accelerating the gradient method with the precise q in a well-conditioned region can also lead to wasted iterations [28]. The Lipschitz constant L can be estimated in a straightforward manner using backtracking (e.g. [22, pg. 162-163], [4, pg. 195]); however obtaining a nontrivial lower bound for the convexity parameter μ is much more challenging. In [25], a backtracking approach is taken to estimate a nontrivial convexity parameter. The use of fixed restart proportional to the condition number has also been considered (see e.g. [4, 13, 18]). A heuristic adaptive restart technique was recently introduced in [28] based on the idea of restarting the momentum factor to zero when a heuristic gradient condition is satisfied. The origin of momentum restart can in fact be traced back to the late 80's (see e.g. [36]). O'Donoghue and Candes [28] demonstrate dramatic speed up in the convergence rate of accelerated gradient methods by adaptively restarting the momentum factor with zero when a heuristic gradient condition is satisfied. They show that their restart scheme recovers the optimal complexity $\mathcal{O}(\sqrt{q^{-1}} \ln \frac{1}{\epsilon})$ for strongly-convex quadratic functions. A significant improved

performance of accelerated gradient methods combined with this heuristic adaptive restart of [28] was also reported in [11, 17, 19].

In this paper, we provide a theoretical justification for the heuristic restart condition of [28] by extending the Nesterov gradient method (NGM) to utilize available secant information. The proposed algorithm is based on updating the estimate-sequence parameter with secant information whenever possible. Furthermore, the proposed algorithm embodies an "update rule with reset" that parallels the restart rule suggested in [28]. Numerical examples indicate that the proposed algorithm significantly outperforms the adaptive restart [28]. The rest of this paper is organized as follows: In sections 3 and 4, the Nesterov gradient method and the quasi-Newton method are discussed. The proposed **Secant-Based-NGM** is described in section 5 and the global convergence for all convex functions will also be established therein.

2 Notation

A continuously differentiable function $h : \mathbb{R}^n \rightarrow \mathbb{R}$ has a Lipschitz continuous gradient on \mathbb{R}^n with constant L if there exist a constant $L > 0$ such that $\|\nabla h(x) - \nabla h(y)\| \leq L\|x - y\|$, $\forall x, y \in \mathbb{R}^n$. A continuously differentiable function $h : \mathbb{R}^n \rightarrow \mathbb{R}$ is convex with parameter μ if there exists a constant $\mu \geq 0$ such that $h(x) \geq h(y) + \nabla h(y)(x - y) + \frac{\mu}{2}\|x - y\|^2$, $\forall x, y \in \mathbb{R}^n$. Subsequently, the term trivial convexity parameter means a lower bound of convexity parameter while nontrivial convexity parameter refers to the greatest lower bound of the convexity parameter. Denote $\mathcal{F}_L^{1,1}(\mathbb{R}^n)$ as the class of convex functions with L -Lipschitz-continuous gradient. Let the function $f \in \mathcal{F}_L^{1,1} : \mathbb{R}^n \rightarrow \mathbb{R}$ denote a convex function bounded below and with a unique minimizer. If $\mu > 0$, then function f is strongly-convex, i.e. $f \in \mathcal{S}_{\mu,L}^{1,1}(\mathbb{R}^n) \subset \mathcal{F}_L^{1,1}(\mathbb{R}^n)$. The optimal values of $f(x)$ and $\Phi_k(x)$ are denoted by f^* and Φ_k^* respectively. Let ∇ denote the gradient operator and it is defined by $\nabla f(x) = \left[\frac{df(x)}{dx_1}, \dots, \frac{df(x)}{dx_n} \right]^T$.

Assumption 1. *A trivial convexity parameter $\mu \geq 0$ and the gradient's Lipschitz constant L are known.*

3 Nesterov Gradient Method

This section reviews the fast gradient method due to Nesterov [24]. Consider the following approximations of $f(x)$ at x_k :

$$\phi_k^1(x) = f(x_k) + \nabla f(x_k)^T(x - x_k) + \frac{1}{2\alpha}\|x - x_k\|^2. \quad (2)$$

$$\phi_k^2(x) = f(x_k) + \nabla f(x_k)^T(x - x_k) + \frac{1}{2}(x - x_k)^T \nabla^2 f(x_k)(x - x_k). \quad (3)$$

The Nesterov gradient method [24] attempts to use approximations which are better than $\phi_k^1(x)$ but less

expensive than $\phi_k^2(x)$ by defining an estimate-sequence (see Definition 1). Provided this estimate-sequence satisfies Nesterov's Principle (see below), then convergence to f^* is guaranteed (see Lemma 1).

Definition 1 ([24]). *A pair of sequences $\{\Phi_k(x)\}_{k=0}^{\infty}$, $\{\lambda_k\}_{k=0}^{\infty}$ is called an estimate-sequence of a function $f(x)$ if $\lambda_k \rightarrow 0$ and for any $x \in \mathbb{R}^n$ and for all $k \geq 0$, we have:*

$$\Phi_k(x) \leq (1 - \lambda_k)f(x) + \lambda_k\Phi_0(x), \quad (4)$$

where $\lambda_k > 0$ and $\Phi_k(x)$ is some local function. ■

The Nesterov gradient method is based on the principle of utilizing a sequence of local functions $\Phi_k(x)$ whose limit approaches the greatest global lower bound of $f(x)$.

Nesterov's Principle: This principle requires that the estimate-sequence (see Definition 1) defined by the local functions $\Phi_k(x)$ is constructed such that

$$f(x_k) \leq \Phi_k^*, \quad \Phi_k^* = \min_x \Phi_k(x). \quad \blacksquare \quad (5)$$

As graphically illustrated in Fig. 1, Nesterov's principle ensures that the local functions $\Phi_k(x)$ constituting the estimate-sequence have a continuum of minima that approaches the minimum of $f(x)$ as $\lambda_k \rightarrow 0$.

This convergence property of Nesterov's principle is made precise in Lemma 1.

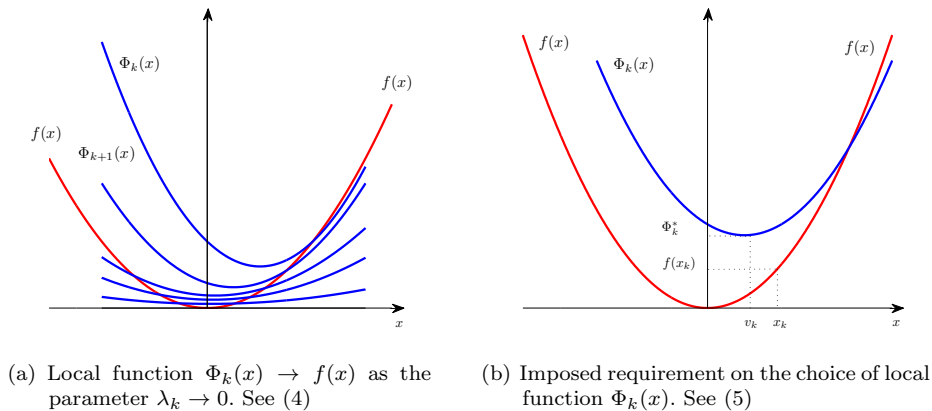


Fig. 1 Nesterov's optimal concept.

An illustration of Nesterov's principle ($f(x_k) \leq \Phi_k^*$ and $\Phi_k(x) \rightarrow f(x)$ as $\lambda_k \rightarrow 0$).

Lemma 1 ([24]). *If a local function $\Phi_k(x)$ is chosen such that Nesterov's principle is satisfied, then*

$$f(x_k) - f(x^*) \leq \lambda_k[\Phi_0(x^*) - f(x^*)], \quad \forall k > 0. \quad \blacksquare$$

Thus, for any scheme that satisfies Nesterov's principle [i.e. (4), (5)], the convergence rate of its minimization process is directly related to the rate of convergence of the λ_k sequence.

The Nesterov scheme is one approach that ensures satisfaction of the Nesterov's principle [i.e. (4), (5)].

The following Lemma 2, gives a recursive rule that satisfies (4) i.e. Definition 1.

Lemma 2 ([24]). *Let scalars $\lambda_0 = 1$, $\beta_k \in (0, 1)$, $\sum_{k=1}^{\infty} \beta_k = \infty$. The following recursive rules,*

$$\lambda_{k+1} = (1 - \beta_k)\lambda_k, \quad (6)$$

$$\Phi_{k+1}(x) \leq (1 - \beta_k)\Phi_k(x) + \beta_k f(x), \quad (7)$$

are sufficient to constitute an estimate-sequence $\{\Phi_k(x)\}_{k=0}^{\infty}$, $\{\lambda_k\}_{k=0}^{\infty}$ in the sense of Definition 1. ■

The Nesterov scheme uses Lemma 2 to construct the estimate-sequence defined in Lemma 3. Thereafter, the acceleration parameter β_k and search point y_k are carefully chosen such that (5) is satisfied.

3.1 NESTEROV'S CHOICE OF RECURSIVE RULE

The Nesterov choice of recursive rule for the local function $\Phi_k(x)$ satisfies the requirements of Lemma 2.

In this section, the recursive rule for $\Phi_k(x)$ used in the Nesterov scheme is given and illustrated in Fig. 2.

Definition 2 ([24]). *Define $\overline{\Phi_{k+1}}(x)$ as*

$$\overline{\Phi_{k+1}}(x) = (1 - \beta_k)\Phi_k(x) + \beta_k [f(y_k) + \nabla f(y_k)^T(x - y_k) + \frac{\mu_k}{2}\|x - y_k\|^2] \quad (8)$$

for a given sequence $\{y_k\}_{k=0}^{\infty}$ and with $\mu_k \in [0, \mu]$. ■

Remark 1. *The recursion $\Phi_{k+1}(x) = \overline{\Phi_{k+1}}(x)$ satisfies Lemma 2. In particular, Nesterov's choice of recursive rule corresponds to $\Phi_{k+1}(x) = \overline{\Phi_{k+1}}(x)$ with $\mu_k = \mu$.*

Consequently, Nesterov's choice of local function $\Phi_{k+1}(x)$ is a convex combination of the previous local function $\Phi_k(x)$ and the greatest global lower bound of $f(x)$. This choice is graphically illustrated in Fig. 2.

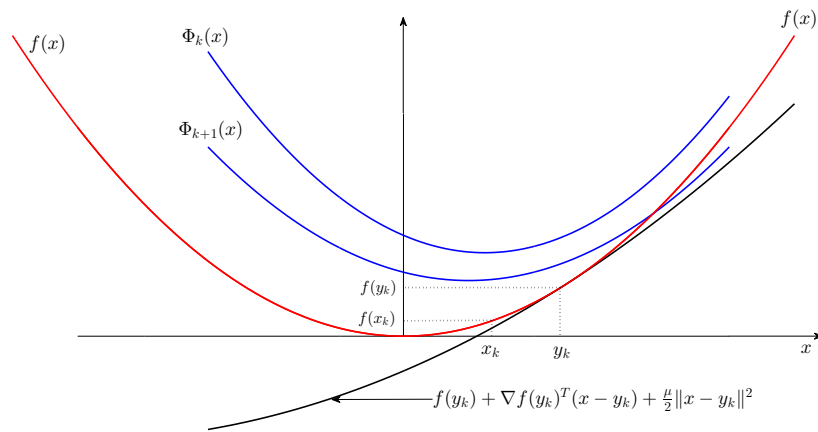


Fig. 2 Updating the local function $\Phi_k(x)$ for some y_k .

Local function $\Phi_{k+1}(x)$ is obtained as a convex combination of $\Phi_k(x)$ and the greatest global lower bound of $f(x)$ at y_k . (see Remark 1).

3.2 NESTEROV'S CHOICE OF ESTIMATE-SEQUENCE

In this section, a simple quadratic form is chosen as the initial local function $\Phi_0(x)$. This simple choice allows the requirements of Lemma 2 to be satisfied easily. Thus the recursion of the sequences defined in Lemma 3 defines an estimate-sequence that satisfies Definition 1.

Lemma 3 ([24]). *Let scalars $\beta_k \in (0, 1)$, $\gamma_k > 0$, $\mu \geq 0$, and $v_k, y_k \in \mathbb{R}^n$. Let $\Phi_0(x) = \Phi_0^* + \frac{\gamma_0}{2} \|x - v_0\|^2$. The recursive rules in Lemma 2 [i.e. (6), (7)] hold for*

$$\Phi_k(x) = \Phi_k^* + \frac{\gamma_k}{2} \|x - v_k\|^2 \quad (9)$$

provided the sequences $\{\gamma_k, v_k, \Phi_k^*\}_{k=0}^\infty$ are defined as

$$\gamma_{k+1} = (1 - \beta_k)\gamma_k + \beta_k\mu, \quad (10)$$

$$v_{k+1} = \frac{1}{\gamma_{k+1}} [(1 - \beta_k)\gamma_k v_k + \beta_k\mu y_k - \beta_k \nabla f(y_k)], \quad (11)$$

$$\begin{aligned} \Phi_{k+1}^* = & (1 - \beta_k)\Phi_k^* + \beta_k f(y_k) - \frac{\beta_k^2}{2\gamma_{k+1}} \|\nabla f(y_k)\|^2 + \\ & \frac{\beta_k(1 - \beta_k)\gamma_k}{\gamma_{k+1}} \left[\frac{\mu}{2} \|y_k - v_k\|^2 + \nabla f(y_k)^T (v_k - y_k) \right]. \end{aligned} \quad (12)$$

The variables γ_k and β_k shall subsequently be referred to as the estimate-sequence parameter and acceleration parameter respectively. The estimate-sequence is as desired but still the local condition (5) at the next iterate, $f(x_{k+1}) \leq \Phi_{k+1}^*$, needs to be ensured. This is subsequently achieved in § 3.3 by carefully choosing the accelerating parameter β_k and the search point y_k such that (5) is satisfied.

3.3 NESTEROV'S CHOICE OF β_k AND SEARCH POINT y_k

Suppose that $f(x_k) \leq \Phi_k^*$. Denote $\zeta(\beta_k) = \left[f(x_k) - f(y_k) + \frac{\beta_k\gamma_k}{\gamma_{k+1}} \left[\frac{\mu}{2} \|y_k - v_k\|^2 + \nabla f(y_k)^T (v_k - y_k) \right] \right]$.

Then Φ_{k+1}^* (12) can be written as

$$\Phi_{k+1}^* \geq f(y_k) - \frac{\beta_k^2}{2\gamma_{k+1}} \|\nabla f(y_k)\|^2 + (1 - \beta_k)\zeta(\beta_k). \quad (13)$$

The choice of y_k and β_k that satisfies Nesterov's Principle (i.e. Lemma 3 and $f(x_{k+1}) \leq \Phi_{k+1}^*$) is obtained as follows. Take $x_{k+1} = y_k - \alpha_k \nabla f(y_k)$, $\alpha_k = \frac{1}{L}$. Then, we have $f(x_{k+1}) \leq f(y_k) - \frac{1}{2L} \|\nabla f(y_k)\|^2$. Hence, to satisfy $f(x_{k+1}) \leq \Phi_{k+1}^*$, the search point y_k is chosen as

$$y_k = \frac{\beta_k\gamma_k v_k + \gamma_{k+1} x_k}{\gamma_k + \beta_k\mu} : \zeta(\beta_k) \geq \left[\frac{\beta_k\gamma_k}{\gamma_{k+1}} \left[\frac{\mu}{2} \|y_k - v_k\|^2 \right] \right], \quad (14)$$

the estimate-sequence parameter γ_{k+1} , **Algorithm 1a** simplifies to **Algorithm 1b**.

Algorithm 1b ([24]). The simplified Nesterov gradient method is outlined as follows:

Algorithm 1b Simplified Nesterov gradient method [24].

Given a starting point $x_0 \in \mathbf{dom} f$, $\beta_0 \in (0, 1)$, $y_0 = x_0$ and $q = \frac{\mu}{L}$.

repeat until stopping criterion is satisfied

1. Compute the Nesterov iterate: $x_{k+1} = y_k - \alpha_k \nabla f(y_k)$, with $\alpha_k = \frac{1}{L}$.
2. Compute $\beta_{k+1} \in (0, 1)$ from $\beta_{k+1}^2 = (1 - \beta_{k+1})\beta_k^2 + q\beta_{k+1}$.
3. Compute θ_{k+1} : $\theta_{k+1} = \frac{\beta_k(1 - \beta_k)}{\beta_{k+1} + \beta_k^2}$.
4. Compute $y_{k+1} = x_{k+1} + \theta_{k+1}(x_{k+1} - x_k)$.

end (repeat)

Remark 3. The choice of $\beta_0 = \sqrt{\frac{\mu}{L}}$ corresponds to $\gamma_0 = \mu$ while the corresponding β_0 for the case of $\gamma_0 = L$ can be obtained from (15). It is important to emphasize that $\beta_0 \neq 1$ since (15) cannot hold when $\beta_0 = 1$. Hence the choice $\beta_0 \in (0, 1)$. Furthermore, if $\beta_0 = \sqrt{\frac{\mu}{L}}$, then $\beta_k = \sqrt{\frac{\mu}{L}}$ and $\theta_k = \frac{\sqrt{L} - \sqrt{\mu}}{\sqrt{L} + \sqrt{\mu}}$ for all k . Were β_k be chosen as 0 for all $k \geq 0$, **Algorithm 1** would reduce to a fixed-step gradient-descent method. Subsequently, the variable θ_k shall be referred to as the momentum parameter.

Theorem 1 ([24]). Let $\Phi_0(x) = \Phi_0^* + \frac{\gamma_0}{2}\|x - v_0\|^2$. Suppose $v_0 = x_0$. If a scheme satisfies Nesterov's principle [i.e. (4), (5)], then

$$f(x_k) - f(x^*) \leq \lambda_k \left[f(x_0) - f(x^*) + \frac{\gamma_0}{2}\|x_0 - x_0\|^2 \right], \quad \forall k > 0,$$

where $\lambda_0 = 1$ and $\lambda_k = \prod_{i=0}^{k-1} (1 - \beta_i)$. ■

Remark 4. Take $\gamma_0 = L$ in **Algorithm 1a** (or the corresponding β_0 in **Algorithm 1b**). Let $v_0 = x_0$. Then the Nesterov scheme satisfies the premises of Theorem 1. Since $\gamma_k > 0$ for all k , then the Nesterov gradient method **Algorithm 1** generates a sequence $\{x_k\}_{k=0}^{\infty}$ such that

$$f(x_k) - f^* \leq \frac{4L}{(k+2)^2} \times \|x_0 - x^*\|^2. \quad (17)$$

Furthermore, since $\gamma_k \geq \mu > 0$ for all k , then

$$f(x_k) - f^* \leq \min \left\{ L \left(1 - \sqrt{\frac{\mu}{L}}\right)^k, \frac{4L}{(k+2)^2} \right\} \times \|x_0 - x^*\|^2. \quad (18)$$

4 Quasi-Newton Method

In the quasi-Newton [27] method, the local function is a local quadratic model about x_k where B_k is

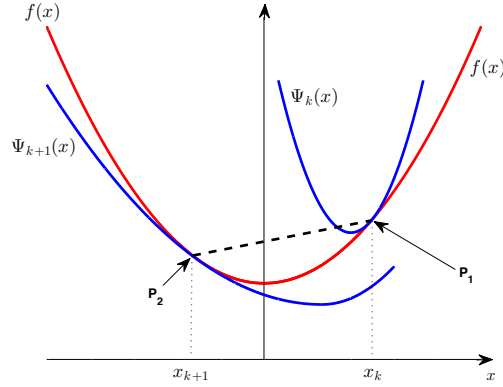


Fig. 4 Quasi-Newton method: Updating local function $\Psi_k(x)$.

Illustrating the secant line P_1 - P_2 . The imposed requirement on B_{k+1} is that $\nabla\Psi_{k+1}(x_k) = \nabla f(x_k)$.

the Hessian-approximate at x_k i.e.

$$\Psi_k(x) = f(x_k) + \nabla f(x_k)^T (x - x_k) + \frac{1}{2}(x - x_k)^T B_k (x - x_k). \quad (19)$$

Define the point where $\Psi_k(x_k) = f(x_k)$ and $\nabla\Psi_k(x_k) = \nabla f(x_k)$ as point P_1 (see Fig. 4). Suppose the new iterate x_{k+1} has been generated by minimizing $\Psi_k(x)$ i.e.

$$x_{k+1} = x_k + \alpha_k d_k = x_k - B_k^{-1} \nabla f(x_k) = x_k + s_k. \quad (20)$$

We wish to construct $\Psi_{k+1}(x)$ of the form

$$\Psi_{k+1}(x) = f(x_{k+1}) + \nabla f(x_{k+1})^T (x - x_{k+1}) + \frac{1}{2}(x - x_{k+1})^T B_{k+1} (x - x_{k+1}). \quad (21)$$

Similarly, define the point where $\Psi_{k+1}(x_{k+1}) = f(x_{k+1})$ and $\nabla\Psi_{k+1}(x_{k+1}) = \nabla f(x_{k+1})$ as P_2 (see Fig. 4). Requirements can be imposed on B_{k+1} based on our knowledge of the previous step. For a reliable Hessian estimate B_{k+1} , it is reasonable to expect that in addition to $\nabla\Psi_{k+1}(x_{k+1}) = \nabla f(x_{k+1})$, that it is desired to have $\nabla\Psi_{k+1}(x_k) = \nabla f(x_k)$. Therefore, it follows that

$$B_{k+1}(x_{k+1} - x_k) = \nabla f(x_{k+1}) - \nabla f(x_k). \quad (22)$$

This imposed condition (22) is known as the secant condition and can be written as:

$$B_{k+1} s_k = y_k. \quad (23)$$

Alternatively, the secant condition can be derived from the mean-value theorem for vector-valued func-

tions which implies that (23) is satisfied by the mean Hessian in the interval $[x_k, x_{k+1}]$ [6]. The pair (s_k, y_k) is said to be the secant pair associated with the secant condition (23). The matrix B_k is updated (see [27]) using symmetric rank-one updates (SR1) or symmetric rank-two updates (e.g. Powell-Symmetric-Broyden(PSB) and Davidon-Fletcher-Powell(DFP) updates).

5 A Secant-Based Nesterov Gradient Method

A careful study of (13) reveals that $f(x_k) \leq \Phi_k^*$ in the Nesterov scheme provided that $\beta_k^2 L - \gamma_{k+1} \leq 0$. This implies that the accelerating parameter β_k can be chosen such that $\beta_k \in (0, \beta'_k]$ where β'_k is the computed solution of step 1 of **Algorithm 1a**. A natural question that arises is "how should $\beta_k \in (0, \beta'_k]$ be chosen?". In this section, a new accelerated gradient method (**Secant-Based-NGM**) is proposed by extending the classical Nesterov gradient method to utilize available secant information whenever possible. **Secant-Based-NGM** is based on updating the estimate-sequence parameter γ_k by imposing a secant condition on the choice of search point y_k . This approach exploits the curvature information at the k^{th} iterate when determining the accelerating parameter β_k . As a result, the computed momentum parameter θ_k is varied in accordance to the changing curvature of the objective function $f(x)$. The global convergence of the proposed **Secant-Based-NGM** is also established for all convex functions.

5.1 RECURSIVE RULE REVISITED

Definition 3. Let $l_k(x) : l_k(x) = f(y_k) + \nabla f(y_k)^T(x - y_k) + \frac{\mu_k}{2} \|x - y_k\|^2$ for a given sequence $\{y_k\}_{k=0}^{\infty}$ and with $\mu_k \in [0, \mu]$.

Recall from Lemma 2 that it is sufficient for the recursive rule for $\Phi_k(x)$ to satisfy,

$$\Phi_{k+1}(x) \leq (1 - \beta_k)\Phi_k(x) + \beta_k f(x). \quad (24)$$

Recall from Remark 1 that $\overline{\Phi_{k+1}}(x) = \overline{\Phi_{k+1}}(x)$ satisfies (24) where,

$$\overline{\Phi_{k+1}}(x) = (1 - \beta_k)\Phi_k(x) + \beta_k l_k(x). \quad (25)$$

In the case $\mu_k \neq 0$, it then follows from (25) that

$$\overline{\Phi_{k+1}}(x) = (1 - \beta_k) \left[\Phi_k^* + \frac{\gamma_k}{2} \|x - v_k\|^2 \right] + \beta_k \left[l_k^* + \frac{\mu_k}{2} \|x - z_k\|^2 \right], \quad (26)$$

where $z_k = y_k - \frac{1}{\mu_k} \nabla f(y_k)$, $l_k^* = f(y_k) - \frac{1}{2\mu_k} \|\nabla f(y_k)\|^2$. In the case $\mu_k = 0$, it follows from (25) that,

$$\overline{\Phi_{k+1}}(x) = (1 - \beta_k) \left[\Phi_k^* + \frac{\gamma_k}{2} \|x - v_k\|^2 \right] + \beta_k \left[\hat{l}_k^* + \nabla f(y_k)^T x \right], \quad (27)$$

where $\hat{l}_k^* = f(y_k) - \nabla f(y_k)^T y_k$. Subsequently, we do not explore the flexibility in μ_k and take that $\mu_k = \mu$ as used in the original scheme of Nesterov. Furthermore, the proposed algorithm chooses $\Phi_{k+1}(x)$ such that $\Phi_{k+1}(x) \leq \overline{\Phi_{k+1}}(x)$ i.e.

$$\Phi_{k+1}(x) \leq (1 - \beta_k)\Phi_k(x) + \beta_k \left[f(y_k) + \nabla f(y_k)^T (x - y_k) + \frac{\mu}{2} \|x - y_k\|^2 \right]. \quad (28)$$

5.2 CONSTRUCTION OF **Secant-Based-NGM**

Basic Secant-Based-NGM extends the classical Nesterov scheme by utilizing the secant information in updating the estimate-sequence parameter γ_k . The two subsequent lemmas are used to arrive at an inequality (36) that gives an upper bound to γ_k .

Lemma 4. *Given $a, b > 0$ and $x_1, x_2 \in \mathbb{R}^n$, there exists $x_3 \in \mathbb{R}^n$ and $d \geq 0$ such that $a\|x - x_1\|^2 + b\|x - x_2\|^2 = (a + b)\|x - x_3\|^2 + d$ for all $x \in \mathbb{R}^n$.*

Proof:

Take

$$x_3 = \frac{ax_1 + bx_2}{a + b} \quad \text{and} \quad d = \frac{ab(x_1 - x_2)^T(x_1 - x_2)}{a + b}. \quad (29)$$

It then follows that

$$a\|x - x_1\|^2 + b\|x - x_2\|^2 = (a + b)\|x - x_3\|^2 + d. \quad (30)$$

Hence, $d = 0$ if and only if $x_1 = x_2$. Thus $d \geq 0$ if $a, b > 0$. ■

Lemma 5. *Given $a, b > 0$ and $x_1, x_2 \in \mathbb{R}^n$, there exists $x_3 \in \mathbb{R}^n$ and $\hat{d} \in \mathbb{R}$ such that $a\|x - x_1\|^2 + bx_2^T x = a\|x - x_3\|^2 + \hat{d}$ for all $x \in \mathbb{R}^n$.*

Proof:

Take

$$x_3 = \frac{2ax_1 - bx_2}{2a} \quad \text{and} \quad \hat{d} = \frac{4abx_1^T x_2 - b^2 x_2^T x_2}{4a}. \quad (31)$$

It then follows that $a\|x - x_1\|^2 + bx_2^T x = a\|x - x_3\|^2 + \hat{d}$. It is clear from (31) that $\hat{d} \in \mathbb{R}$ if $a, b > 0$. ■

The subsequent Lemma 6 gives the needed freedom in updating the γ_k sequence and also gives the upper bound to the acceptable value of γ_k .

Lemma 6. *Let scalars $\beta_k \in (0, 1)$, $\gamma_k > 0$, $\mu \geq 0$, and $v_k, y_k \in \mathbb{R}^n$. Let $\Phi_0(x) = \Phi_0^* + \frac{\gamma_0}{2}\|x - v_0\|^2$. The recursive rules in Lemma 2 [i.e. (6), (7)] hold for*

$$\Phi_k(x) = \Phi_k^* + \frac{\gamma_k}{2}\|x - v_k\|^2 \quad (32)$$

provided the sequences $\{\gamma_k, v_k, \Phi_k^*\}_{k=0}^\infty$ are defined as

$$\gamma_{k+1}^F = (1 - \beta_k)\gamma_k + \beta_k\mu, \quad (33)$$

$$v_{k+1} = \frac{1}{\gamma_{k+1}^F}[(1 - \beta_k)\gamma_k v_k + \beta_k\mu y_k - \beta_k \nabla f(y_k)], \quad (34)$$

$$\begin{aligned} \Phi_{k+1}^* &= (1 - \beta_k)\Phi_k^* + \beta_k f(y_k) - \frac{\beta_k^2}{2\gamma_{k+1}^F} \|\nabla f(y_k)\|^2 + \\ &\quad \frac{\beta_k(1 - \beta_k)\gamma_k}{\gamma_{k+1}^F} \left[\frac{\mu}{2} \|y_k - v_k\|^2 + \nabla f(y_k)^T (v_k - y_k) \right], \end{aligned} \quad (35)$$

$$\gamma_{k+1} \leq \gamma_{k+1}^F. \quad (36)$$

Proof:

From (25), $\overline{\Phi_{k+1}}(x)$ is defined as $\overline{\Phi_{k+1}}(x) = (1 - \beta_k)\Phi_k(x) + \beta_k l_k(x)$. Furthermore, choose v_{k+1} as the unconstrained minimum of $\overline{\Phi_{k+1}}(x)$. Since $\gamma_{k+1}^F = (1 - \beta_k)\gamma_k + \beta_k\mu$, we can obtain v_{k+1} as,

$$v_{k+1} = \frac{1}{\gamma_{k+1}^F}[(1 - \beta_k)\gamma_k v_k + \beta_k\mu y_k - \beta_k \nabla f(y_k)]. \quad (37)$$

Next, the inequality (36) is established for the cases when the convexity parameter $\mu \neq 0$ and $\mu = 0$.

Case 1: $\mu \neq 0$

By Lemma 4, it follows from (26) that $\overline{\Phi_{k+1}}(x) = (1 - \beta_k)\Phi_k^* + \beta_k l_k^* + d + \frac{\gamma_{k+1}^F}{2}\|x - v_{k+1}\|^2$ for some $d \geq 0$. The case of $d = 0$ occurs if and only if we have coincident minimizers i.e. $v_{k+1} = v_k = z_k$ (see [12]).

Nesterov's choice then corresponds to $\Phi_{k+1}(x) = \overline{\Phi_{k+1}}(x)$ with $\gamma_{k+1} = \gamma_{k+1}^F$ and $\Phi_{k+1}^* = (1 - \beta_k)\Phi_k^* + \beta_k l_k^* + d$. However, we choose

$$\gamma_{k+1} \leq \gamma_{k+1}^F. \quad (38)$$

Consequently, we then have $\Phi_{k+1}(x) \leq \overline{\Phi_{k+1}}(x)$ which still satisfies (24) and more importantly we still have $\Phi_{k+1}(v_{k+1}) = \overline{\Phi_{k+1}}(v_{k+1})$. Also note that the unconstrained minimum of this choice of $\Phi_{k+1}(x)$ still remains as v_{k+1} (34). This inequality (36) is crucial in the sense that it gives an upper bound to γ_k for all k and allows the use of the secant information in updating the γ_k sequence whenever possible.

Case 2: $\mu = 0$

By Lemma 5, it follows from (27) that $\overline{\Phi_{k+1}}(x) = (1 - \beta_k)\Phi_k^* + \beta_k \hat{l}_k^* + \hat{d} + \frac{\gamma_{k+1}^F}{2}\|x - v_{k+1}\|^2$ for some

$\hat{d} \in \mathbb{R}$. Thus the inequality (36) still holds for this case using similar arguments as in the case of $\mu \neq 0$.

It remains to compute Φ_{k+1}^* . It follows from (32) that at $x = v_{k+1}$, we have that $\Phi_{k+1}^* = \Phi_{k+1}(v_{k+1})$.

Since $\Phi_{k+1}(v_{k+1}) = \overline{\Phi_{k+1}(v_{k+1})}$, then by (25) it follows that Φ_{k+1}^* can be computed as

$$\Phi_{k+1}^* = (1 - \beta_k)\Phi_k(v_{k+1}) + \beta_k l_k(v_{k+1}). \quad (39)$$

Substitute for $\Phi_k(v_{k+1})$ in (39) using (32), then (39) becomes

$$\begin{aligned} \Phi_{k+1}^* &= (1 - \beta_k)\Phi_k^* + \frac{(1 - \beta_k)\gamma_k}{2}\|v_{k+1} - v_k\|^2 + \\ &\quad \beta_k f(y_k) + \beta_k \nabla f(y_k)(v_{k+1} - y_k) + \frac{\beta_k \mu}{2}\|v_{k+1} - y_k\|^2. \end{aligned} \quad (40)$$

It follows from (37) that

$$v_{k+1} - y_k = \frac{(1 - \beta_k)\gamma_k}{\gamma_{k+1}^F}(v_k - y_k) - \frac{\beta_k}{\gamma_{k+1}^F}\nabla f(y_k), \quad (41)$$

and that

$$v_{k+1} - v_k = \frac{\beta_k \mu}{\gamma_{k+1}^F}(y_k - v_k) - \frac{\beta_k}{\gamma_{k+1}^F}\nabla f(y_k). \quad (42)$$

Substituting (41) and (42) into (40), then the result for Φ_{k+1}^* follows,

$$\begin{aligned} \Phi_{k+1}^* &= (1 - \beta_k)\Phi_k^* + \beta_k f(y_k) - \frac{\beta_k^2}{2\gamma_{k+1}^F}\|\nabla f(y_k)\|^2 + \\ &\quad \frac{\beta_k(1 - \beta_k)\gamma_k}{\gamma_{k+1}^F} \left[\frac{\mu}{2}\|y_k - v_k\|^2 + \nabla f(y_k)^T(v_k - y_k) \right]. \quad \blacksquare \end{aligned} \quad (43)$$

Remark 5. Lemma 3 is a special case of Lemma 6 if γ_{k+1} is chosen as $\gamma_{k+1} = \gamma_{k+1}^F$.

The secant information is used in updating γ_k by requiring that $\nabla \Phi_{k+1}(y_k) = \nabla f(y_k)$. It then follows that $\hat{\gamma}_{k+1}(y_k - v_{k+1}) = \nabla f(y_k)$ where $\hat{\gamma}_{k+1}$ is a possible update of γ_k . Using a symmetric rank-1 update, it then follows that

$$\hat{\gamma}_{k+1} = \frac{\nabla f(y_k)^T(y_k - v_{k+1})}{(y_k - v_{k+1})^T(y_k - v_{k+1})}. \quad (44)$$

However, this γ_k updating is subject to the constraint (36). An effective way of enforcing (36) is that larger values of $\hat{\gamma}_{k+1}$ correspond to updates closer to γ_{k+1}^F and vice-versa. To this end, we define γ_{k+1}^E ,

$$\gamma_{k+1}^E = \frac{\hat{\gamma}_{k+1}}{\hat{\gamma}_{k+1} + \gamma_{k+1}^F} \times \gamma_{k+1}^F. \quad (45)$$

where the effective curvature γ_{k+1}^E is to be used in defining an update rule for γ_k in the proposed algorithm.

The search point y_k still has to satisfy (16) and the accelerating parameter β_k computed from $\beta_k^2 L = \gamma_{k+1}^F$

as in the classical Nesterov gradient method. However, updating γ_k based on (44) and (45) ensures that the computed $(\beta_k \in (0, \beta'_k])$ satisfies $\beta_k^2 \mathbf{L} - \gamma_{k+1}^{FF} \leq 0$ (where $\gamma_{k+1}^{FF} = (1 - \beta_k)\gamma_k^F + \beta_k\mu$, see introduction of § 5). Moreover, β_k is determined by the curvature information at the k^{th} iterate (see step 1 of **Algorithm 2a**).

This computation of $\hat{\gamma}_{k+1}$ comes at an extra cost of 2 vector-vector multiplication. However, as shown in the numerical results, the benefits of computing γ_{k+1} outweigh the extra cost of its computation. The update γ_{k+1} can now be appended to the classical Nesterov gradient method in a straight-forward manner as shown below in step 6 of **Basic Secant-Based-NGM** below.

Algorithm 2a (Basic secant-based algorithm). The outline is as follows:

Basic Secant-Based-NGM.

Given a starting point $x_0 \in \mathbf{dom} f$, $\gamma_0 > 0$ and $v_0 = x_0$.

repeat until stopping criterion is satisfied

1. Compute $\beta_k \in (0, 1)$ from $\beta_k^2 \mathbf{L} + \beta_k(\gamma_k - \mu) - \gamma_k = 0$.
2. Compute γ_{k+1}^F : $\gamma_{k+1}^F = (1 - \beta_k)\gamma_k + \beta_k\mu$. (33)
3. Compute search point: $y_k = x_k - \rho_k\gamma_k(x_k - v_k)$.
4. Compute the Nesterov iterate: $x_{k+1} = y_k - \alpha_k \nabla f(y_k)$, with $\alpha_k = \frac{1}{\mathbf{L}}$.
5. Compute v_{k+1} : $v_{k+1} = \frac{1}{\gamma_{k+1}^F} [(1 - \beta_k)\gamma_k v_k + \beta_k\mu y_k - \beta_k \nabla f(y_k)]$. (44)
6. Compute $\hat{\gamma}_{k+1}$: $\hat{\gamma}_{k+1} = \frac{\nabla f(y_k)^T (y_k - v_{k+1})}{(y_k - v_{k+1})^T (y_k - v_{k+1})}$. (34)
7. Compute γ_{k+1} : $\gamma_{k+1} = \widehat{\min}_\mu (\gamma_{k+1}^E, \gamma_{k+1}^F)$ update rule
8. If $\hat{\gamma}_{k+1} < 0$, then set $\gamma_{k+1} = \beta_k^2 \mu$ reset rule

end (repeat)

Remark 6. The $\widehat{\min}_\mu$ operator rule in step 7 is given by

$$c = \widehat{\min}_\mu (a, b) : \begin{cases} c = a & \text{if } \gamma_{k+1}^F > \mu, \\ c = \min(\hat{\gamma}_{k+1}, b) & \text{if } \gamma_{k+1}^F < \mu \text{ and } \hat{\gamma}_{k+1} > \hat{\gamma}_k \\ c = b & \text{if otherwise.} \end{cases}$$

The extra condition in the case of $\gamma_{k+1}^F < \mu$ serves to penalize oscillation in the trajectory of $\hat{\gamma}_{k+1}$.

Remark 7. In the case $\mu \neq 0$, see Remark 8 for the appropriate reset rule.

Just as with the classical Nesterov gradient method, $v_{k+1} = x_{k+1} + \frac{1 - \beta_k}{\beta_k}(x_{k+1} - x_k)$ and the variable v_k can therefore be eliminated. With this elimination of v_{k+1} , **Basic Secant-Based-NGM** simplifies

to **Simplified Secant-Based-NGM**.

Algorithm 2b (Simplified secant-based algorithm). The outline is as follows:

Simplified Secant-Based-NGM.

Given a starting point $x_0 \in \mathbf{dom} f$, $\beta_0 \in (0, 1)$ and $y_0 = x_0$.

repeat until stopping criterion is satisfied

1. Compute Nesterov iterate: $x_{k+1} = y_k - \alpha_k \nabla f(y_k)$, with $\alpha_k = \frac{1}{L}$.
2. Compute $\gamma_{k+1}^F = \beta_k^2 L$; $\tau_k = \frac{1 - \beta_k}{\beta_k}$.
3. Compute $y_v = [\alpha_k \nabla f(y_k) - \tau_k(x_{k+1} - x_k)]$; $\hat{\gamma}_{k+1} = \frac{y_v^T \nabla f(y_k)}{y_v^T y_v}$.
4. Compute γ_{k+1} : $\gamma_{k+1} = \widehat{\min}_\mu(\gamma_{k+1}^E, \gamma_{k+1}^F)$ update rule
5. If $\hat{\gamma}_{k+1} < 0$, then set $\gamma_{k+1} = \beta_k^2 \mu$ reset rule
6. Compute $\beta_{k+1} \in (0, 1)$ from $\beta_{k+1}^2 L + \beta_{k+1}(\gamma_{k+1} - \mu) - \gamma_{k+1} = 0$.
7. Compute θ_{k+1} : $\theta_{k+1} = \rho_{k+1} \gamma_{k+1} \tau_k$, where $\rho_{k+1} = \frac{\beta_{k+1}}{\gamma_{k+1} + \beta_{k+1} \mu}$.
8. Compute $y_{k+1} = x_{k+1} + \theta_{k+1}(x_{k+1} - x_k)$.

end (repeat)

Fact: Let γ_{k+1} , γ_{k+1}^F , β_k and θ_{k+1} be as previously defined. If $\mu = 0$ and $\gamma_{k+1} = \frac{\gamma_{k+1}^F}{n}$ for some $n \gg 1$, then the ratio $\frac{\beta_{k+1}}{\beta_k} \approx \frac{1}{\sqrt{n}}$ and $\theta_{k+1} \approx \frac{1 - \beta_k}{\sqrt{n}}$. The proof follows from the definitions of γ_{k+1}^F , β_{k+1} and θ_{k+1} in steps (2, 6, 7) respectively.

Remark 8. In the case $\mu = 0$, then the reset rule in step 5 can be replaced with $\gamma_{k+1} = \frac{\gamma_{k+1}^F}{L^2}$ and this corresponds to a momentum factor $\theta_{k+1} \approx \frac{1}{L}$. However, to account for the case where $L \gg 1$, we have used $\gamma_{k+1} = \min(\frac{\gamma_{k+1}^F}{L^2}, \epsilon \gamma_{k+1}^F)$ where $\epsilon = 10^{-6}$.

In what follows, the gradient restart condition and restart rule of [28] is contrasted with a gradient condition (46) and the proposed "update rule with reset" respectively. With the substitution of v_{k+1} in **Basic Secant-Based-NGM**, the reset condition $\hat{\gamma}_{k+1} < 0$ in step 5 of **Simplified Secant-Based-NGM** is then equivalent to

$$\alpha_k \|\nabla f(y_k)\|^2 - \tau_k \nabla f(y_k)^T (x_{k+1} - x_k) < 0. \quad (46)$$

This gradient condition (46) is more conservative than the gradient-scheme restart condition suggested in [28] especially when the iterates are far away from the optimum point. Thus the gradient condition (46) is less frequently satisfied. The advantage of the conservativeness of (46) is reinforced by the observation

in [28] that ”... restarting far from the optimum can slow down the early convergence slightly, until the quadratic region is reached and the algorithm enters the rapid linear convergence phase.”.

The restart rule of [28] is given as

1. setting β_{k+1} as 1.
2. setting the momentum factor θ_{k+1} as 0.

Firstly, it should be noted that the reset β_{k+1} should be in the interval $(0, 1)$ since $\beta_k \in (0, 1)$ for all $k \geq 0$ (see Remark 3). Moreover, an arbitrary choice of $\beta_{k+1} \in (0, 1)$ may correspond to a γ_{k+1} that violates the inequality $\gamma_{k+1} \leq \gamma_{k+1}^F$ (36). Furthermore, the proposed **Simplified Secant-Based-NGM** proceeds with the computed β_{k+1} unlike the restart rule 1 of [28]. It is emphasized that the proposed **Simplified Secant-Based-NGM** does not reset the momentum factor to zero (see Remark 8) unlike the restart rule 2 of [28]. Thus the **Secant-Based-NGM** is not a momentum-restart algorithm, and the proposed ”update rule with reset” (see Remarks 6, 7, 8) in the **Secant-Based-NGM** satisfies inequality (36). Moreover, restarting the momentum factor θ_{k+1} with 0 as stated in [28] inhibits fast convergence since it discards the entire accumulated information from previous iterates. In the next sub-section, we establish the global convergence of the proposed **Secant-Based-NGM** for all convex functions.

5.3 GLOBAL CONVERGENCE OF PROPOSED SCHEME

The scheme construction of the proposed algorithm (**Secant-Based-NGM**) satisfies Nesterov’s principle [i.e. (4), (5)]. Thus the proposed scheme satisfies the premises of Theorem 1 and therefore the **Secant-Based-NGM** is globally convergent with

$$f(x_k) - f(x^*) \leq \left(\prod_{i=0}^{k-1} (1 - \beta_i) \right) \times \left[f(x_0) - f(x^*) + \frac{\gamma_0}{2} \|x - x_0\|^2 \right], \quad \forall k \geq 1, \quad (47)$$

where $\beta_i \in (0, 1)$ and $\gamma_0 > 0$. If the gradient’s Lipschitz constant L is known, then (47) reduces to

$$f(x_k) - f(x^*) \leq \left(\prod_{i=0}^{k-1} (1 - \beta_i) \right) \times \left[\frac{\gamma_0 + L}{2} \|x - x_0\|^2 \right], \quad \forall k \geq 1. \quad (48)$$

6 Numerical Results

Consider test examples of the form

$$UCOP : \quad \min_{x \in \mathbb{R}^n} f(x), \quad (49)$$

where $x \in \mathbb{R}^n$ is the unknown variable. The numerical tests investigate the effects of increasing the

dimension and condition number respectively on the performance of the proposed algorithm.

Computational Setup: All numerical tests were coded in 64-bit MATLAB on a Dell-Optilex-780 PC with Intel dual-core CPU of 2.93 GHz, RAM of 16GB and a 160GB-free Hard-Disk. All Matlab sessions were single-threaded, *feature('accel','off')* and process-Priority set to "High". All Matlab sessions were executed on the PC running Windows 7 in "Safe Mode". The stopping criterion was $\|\nabla f(x)\| \leq 10^{-9}$ and $\|\nabla f(x)\| \leq 10^{-6}$ for examples 1,3 and 2,4 respectively. All matrices are random square matrices such that the Hessian has eigenvalues in the interval $[1+\mu, L+\mu]$. This was achieved using singular value decomposition to allocate the desired eigenvalues. All matrices and vectors were randomly generated in MATLAB with seed *rng(1234,'twister')*. Except for Ex. 4, the computations were also repeated with seed *rng(5678,'twister')*. Except for Ex. 2, the first run-time effects were accounted for by ignoring the run-times of the first run while averaging over the subsequent 3 runs of each algorithm.

Test Functions and Solvers: We consider four sets of problems with varied number of instances. The test function in Ex. 1 is a well-known convex quadratic function used for benchmarking convex solvers while the test functions in Ex. 2 and Ex. 3 are convex non-quadratic functions usually encountered in machine-learning literature [8, 14, 26, 34, 37]. The test problems in Ex. 4 is the well-known Maros and Mészáros's collection used for benchmarking convex quadratic program solvers [16, 20]. A first-order oracle [24] is used in all numerical computations. The set of solvers \mathcal{S} considered is:

- Classical Nesterov gradient method(NGM)¹,
- Adaptive restart [28],
- Fixed restart² after $k = \sqrt{\frac{8L}{\mu}}$,
- Proposed algorithm: **(Simplified) Secant-Based-NGM**.³

In all cases, $\beta_0 = \sqrt{\mu/L}$ if $\mu \neq 0$ and $\beta_0 = \frac{\sqrt{5}-1}{2}$ if otherwise. These chosen β_0 correspond to $\gamma_0 = \mu$ and $\gamma_0 = L$ respectively. The performance profile in the sense of Dolan and Moré [9] is adopted to analyze the performance data of the above set of solvers \mathcal{S} on each problem set \mathcal{P} . The percentage of the test problems for which a method is the fastest is given on the left axis of the profile plot. The right side of the profile plot gives the percentage of the test problems that were successfully solved by each of the methods. In other words, the right side of the profile plot is a measure of an algorithm's robustness.

Ex. 1: Ridge regression problem [21]

This is a linear least squares problem with Tikhonov regularization. Given $A \in \mathbb{R}^{m \times n}$, $b \in \mathbb{R}^n$ and

¹ The simplified Nesterov gradient method **Algorithm 1b** is used for computations.

² In the case of $\mu = 0$ (Ex. 2 and Ex. 4), fixed restart is done after $k = \max\{N, \sqrt{L}\}$ iterations.

³ In the case of $\mu = 0$ (Ex. 2 and Ex. 4), the algorithm resets with $\gamma_{k+1} = \min\left(\frac{\gamma_k+1}{L^2}, 10^{-6}\gamma_{k+1}^F\right)$, see Remark 8.

the convexity parameter $\mu = 0.1$.

$$f(x) = \frac{\mu}{2}\|x\|_2^2 + \frac{1}{2}\|Ax - b\|_2^2. \quad (50)$$

The objective function $f(x) \in \mathcal{S}_{\mu,L}^{1,1}$ is a positive-definite quadratic convex function with Lipschitz gradient of $L = \lambda_{\max}(A^T A) + \mu$ and trivial convexity parameter of $\mu = 0.1$. All algorithms use the trivial convexity parameter $\mu = 0.1$ except algorithm NGM1 which used a nontrivial convexity parameter of $\mu = 1.1$.

The plots in Fig. 5(a) and Fig. 5(b) show the effect of increasing the problem size (N) and condition number (L/μ) respectively. As expected the NGM with trivial convexity parameter $\mu = 0.1$ (NGM2) is

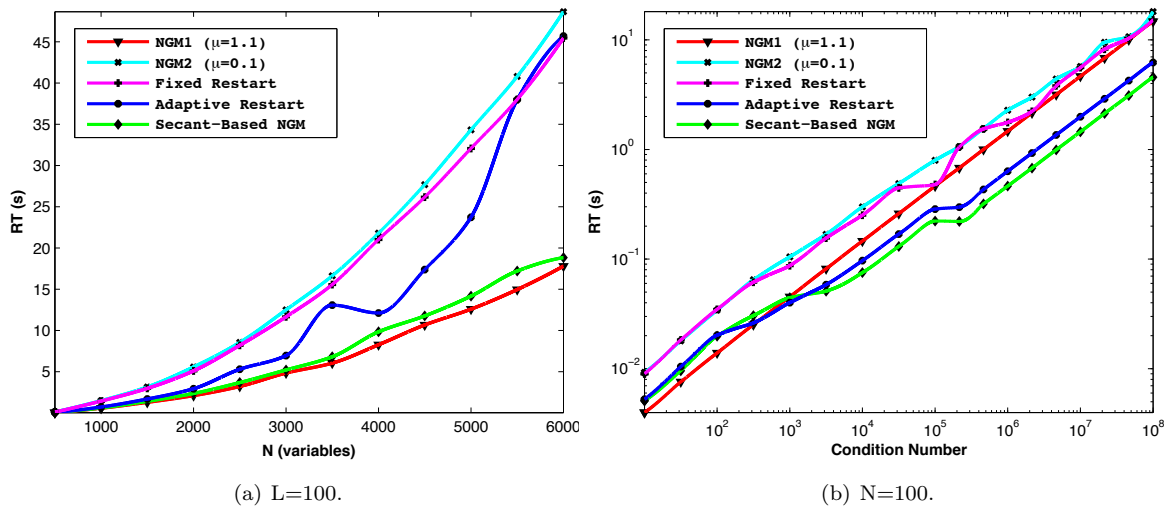


Fig. 5 Effect of problem size (N) and condition number (L/μ) on the run-time.
NGM1 - NGM with nontrivial μ ; NGM2 - NGM with trivial μ .

slowest with the largest run-time (RT) as observed in Fig. 5(a). Using a fixed restart shows slightly improved performance with increasing dimension. However, the adaptive restart and **Secant-Based-NGM** perform significantly better than NGM2 while the **Secant-Based-NGM** performs comparable with NGM1 as the dimension number increases. It can be noted in Fig. 5(b) that the adaptive restart and **Secant-Based-NGM** perform better than even the NGM1. Moreover, the **Secant-Based-NGM** outperforms the adaptive restart as the condition number becomes high.

The computations were repeated for randomly-generated matrices and vectors with seed `rng(5678, 'twister')`.

The plots in Fig. 6(a) and Fig. 6(b) show the effect of increasing the problem size (N) and the condition

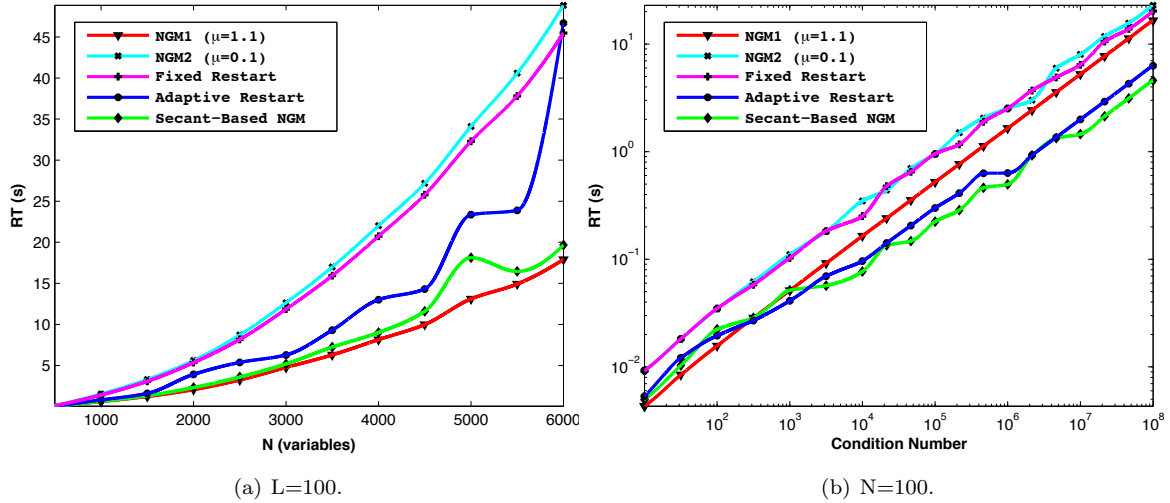


Fig. 6 Effect of problem size (N) and condition number (L/μ) on the run-time. NGM1 - NGM with nontrivial μ ; NGM2 - NGM with trivial μ .

number (L/μ) respectively. Similar conclusions can be drawn from the plots in Fig. 6(a) and Fig. 6(b). The performance profiles [9] for all problem instances (i.e. with seeds $rng(1234, 'twister')$ and $rng(5678, 'twister')$) is shown in Fig. 7(a). As expected, It is clear from Fig. 7(a) that the NGM with nontrivial convexity has the most wins (i.e. the highest probability of being the optimal solver). However, it can also be observed

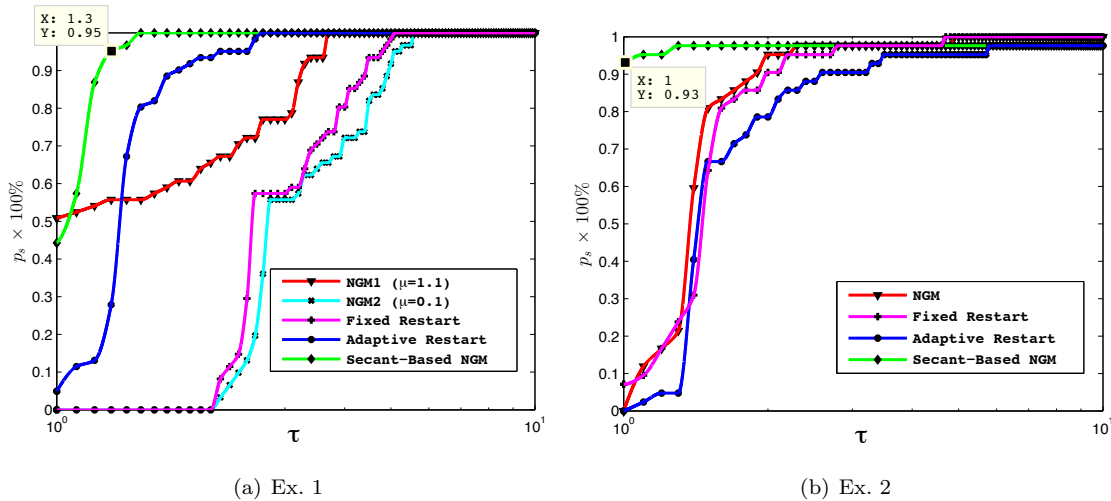


Fig. 7 The performance profiles of the set of solvers S for Ex. 1 and Ex. 2.

from Fig. 7(a) that the performance of NGM with a trivial convexity parameter is improved when a fixed restart or adaptive restart [28] is used. Moreover, **Secant-Based-NGM** performs better than when restarts are used with NGM. In general, the proposed **Secant-Based-NGM** has the highest

probability ($p_s(\tau)=0.95$) of being the fastest solver within a factor $\tau = 1.3$ of the best solver.

Ex. 2: Binary classification problem [8,26]

This a logistic regression problem with l_2 -regularization. Given $z_i \in \mathbb{R}^n$, $y_i \in \{-1, 1\}$, $\mu = 0$ and $N \geq 1$.

$$f(x) = \frac{\mu}{2} \|x\|_2^2 + \sum_{i=1}^N \log(1 + e^{-y_i z_i^T x}). \quad (51)$$

The objective function $f(x) \in \mathcal{F}_L^{1,1}$ is a non-quadratic convex function with Lipschitz gradient of $L = 0.25\lambda_{\max}(F^T F)$ and convexity parameter of $\mu = 0$, where the design matrix $F = [z_1 \cdots z_N]^T \in \mathbb{R}^{N \times n}$. The explained variable $y = \text{sign}(w^T F^T)$ was generated as described in [8,30] except that $w = [1; 1; 1 \cdots 1]$. The choice of w means that each feature has an equal effect on the explained variable y .

The plots in Fig. 8(a) and Fig. 8(b) show the effect of increasing the problem size (N) and condition number (L/μ) respectively. The computations were repeated for randomly-generated matrices and vectors

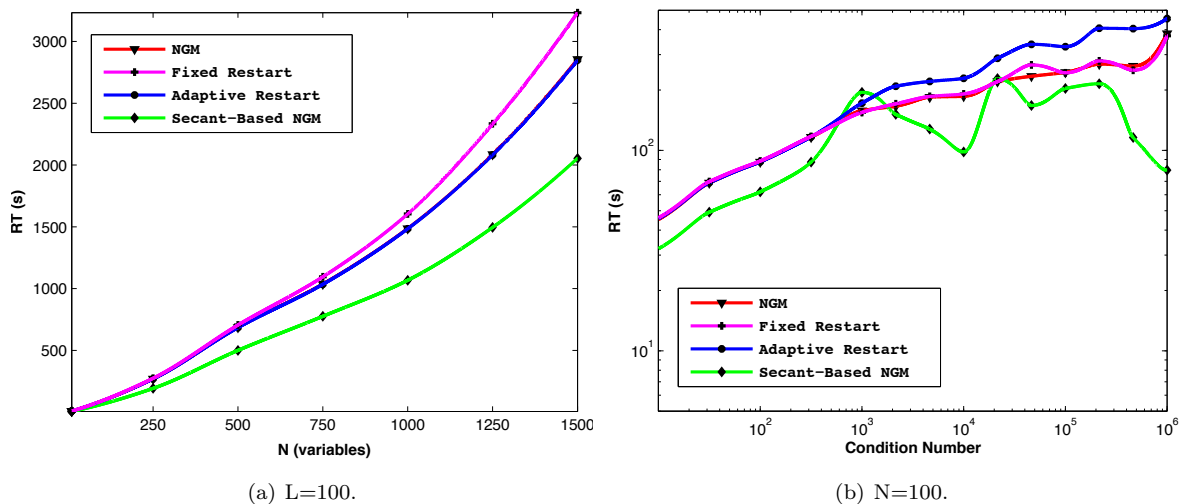


Fig. 8 Effect of problem size (N) and condition number (L/μ) on the run-time.

with seed *rng(5678, 'twister')*. The plots in Fig. 9(a) and Fig. 9(b) show the effect of increasing the problem size (N) and condition number (L/μ) respectively. Similar conclusions can be drawn from the plots in both Fig. 8 and Fig. 9. It can be observed in these figures that all the algorithms have similar performance except for **Secant-Based-NGM** that clearly outperforms the rest as the dimension number increases. In Fig. 9(b), we observe a rather strange behaviour of the Fixed Restart and NGM at the end of the plot. Nevertheless, the proposed **Secant-Based-NGM** significantly outperforms all other

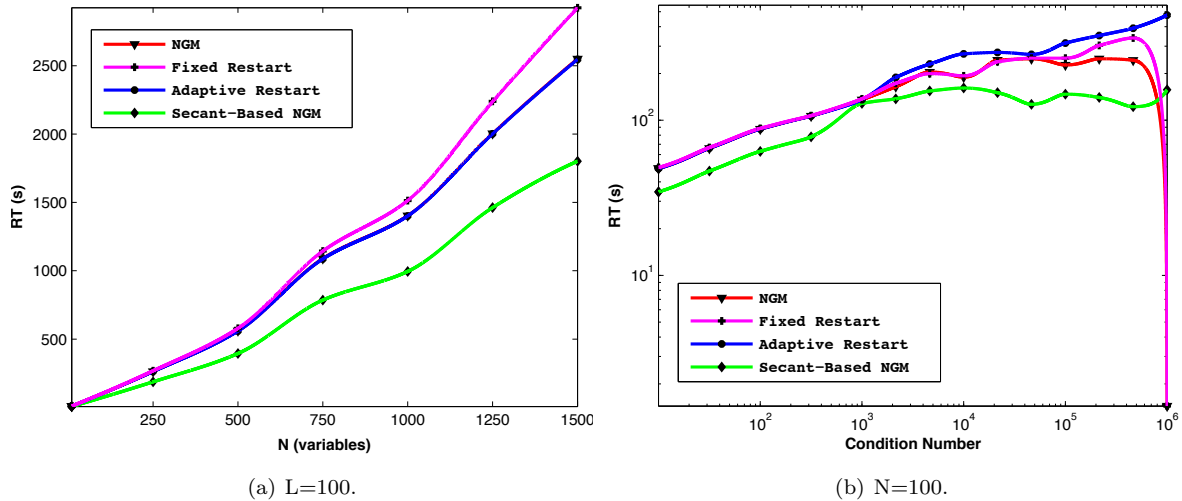


Fig. 9 Effect of problem size (N) and condition number (L/μ) on the run-time.

algorithms as the condition number increases while the adaptive restart performs worse than the NGM.

The performance profiles [9] for all problem instances is shown in Fig. 7(b). It is clear from Fig. 7(b) that **Secant-Based-NGM** is the most efficient of the considered solvers. It solved 93% of the problems significantly faster than the other solvers. It can also be observed from Fig. 7(b) that the performance of adaptive restart [28] is worse than the fixed restart or the classical NGM. In general, the proposed **Secant-Based-NGM** significantly improves over and above the classical NGM and the adaptive restart suggested in [28].

Ex. 3: Approximate Huber loss [14, 34, 37]

Given $z_i \in \mathbb{R}^n$, $y_i \in \{-1, 1\}$, $\mu = 0.1$ and $N \geq 1$.

$$f(x) = \frac{\mu}{2} \|x\|_2^2 + \sum_{i=1}^N \log(\cosh(y_i - z_i^T x)). \quad (52)$$

The objective function $f(x) \in \mathcal{S}_{\mu, L}^{1,1}$ is a non-quadratic convex function with Lipschitz gradient of $L = \lambda_{\max}(F^T F)$ and convexity parameter of $\mu = 0.1$, where the design matrix $F = [z_1 \cdots z_N]^T \in \mathbb{R}^{N \times n}$. The explained variable $y = \text{sign}(w^T F^T)$ was generated as described in Ex. 2.

The plots in Fig. 10(a) and Fig. 10(b) show the effect of increasing the problem size (N) and condition number (L/μ) respectively. The computations were repeated for randomly-generated matrices and vectors with seed `rng(5678, 'twister')`. The plots in Fig. 11(a) and Fig. 11(b) show the effect of increasing the problem size (N) and condition number (L/μ) respectively. Similar conclusions can be drawn from the

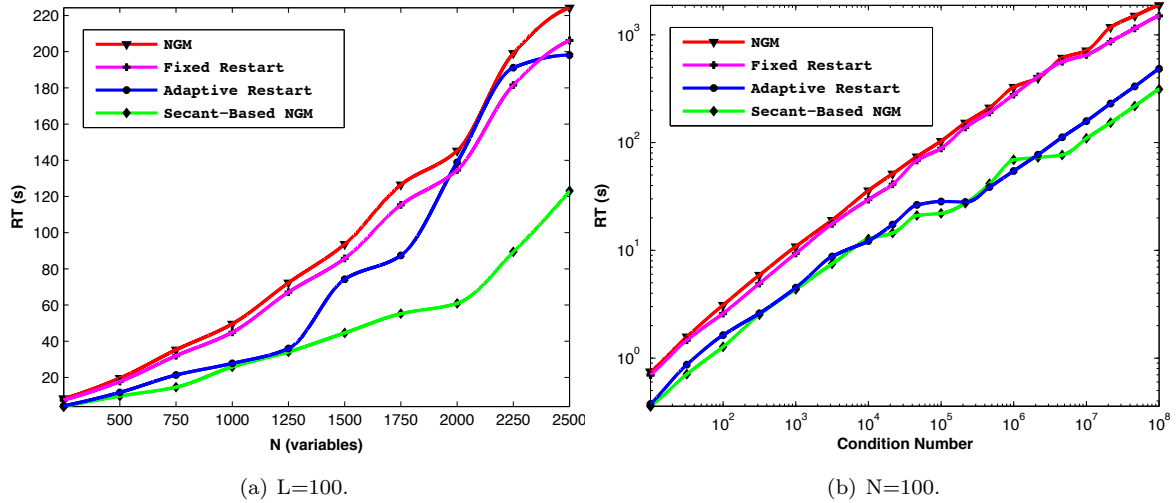


Fig. 10 Effect of problem size (N) and condition number (L/μ) on the run-time.

plots in both Fig. 10 and Fig. 11. It can be observed in these figures that all the algorithms have similar performance except for **Secant-Based-NGM** that clearly outperforms the rest as the dimension number increases. Also in Fig. 10(b), the proposed **Secant-Based-NGM** significantly outperforms other

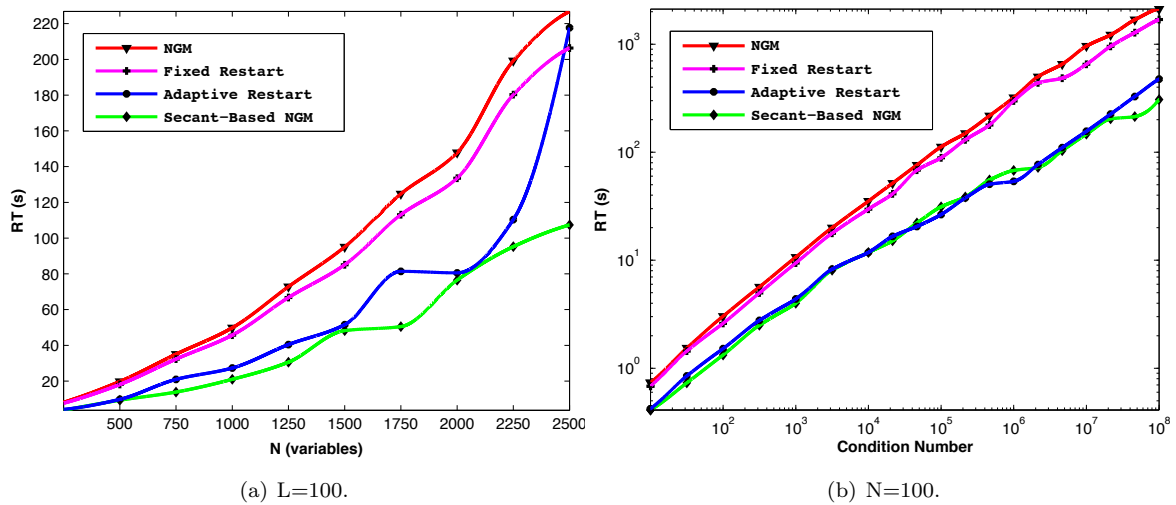


Fig. 11 Effect of problem size (N) and condition number (L/μ) on the run-time.

algorithms as the condition number increases while the NGM has the worst performance.

The performance profiles [9] for all problem instances is shown in Fig. 12(a). It is clear from Fig. 12(a) that **Secant-Based-NGM** is the most efficient of the considered solvers. It solved about 87% of the problems significantly faster than the other solvers. It can also be observed from Fig. 12(a) that the adaptive restart [28] has better run-time performance than the classical NGM. In general,

the proposed **Secant-Based-NGM** significantly improves over and above the classical NGM and the adaptive restart suggested in [28].

Ex. 4: Maros and Mészáros Test Problems [16, 20]

This is a test suite of 138 convex quadratic programs by Maros and Mészáros. This test suite contains selected problems from the BRUNEL and CUTE collections. The objective functions were regularized with $10^{-9}\|x\|_2^2$ to ensure that the Hessians are positive-definite. All considered algorithms use the trivial convexity parameter $\mu = 0$. The initial vectors x_0 were randomly-generated with seed $rng(1234, 'twister')$.

In Table 1 (see Appendix), we report the running time (RT) for each of the solvers. We report only on those test problems in which the run-time (RT) of the classical NGM is within the interval [1s, 500s]. We also ensured that the objective functions of the considered test problems are not in repetition.

As expected the classical NGM had the slowest run-times. The run-time is slightly improved using a fixed-restart and further improved using the adaptive restart. However, as shown in Fig. 12(b), our proposed algorithm **Secant-Based-NGM** performs significantly better than the adaptive restart in about 65% of the test problems and no worse than 1.65 times slower than the adaptive restart.

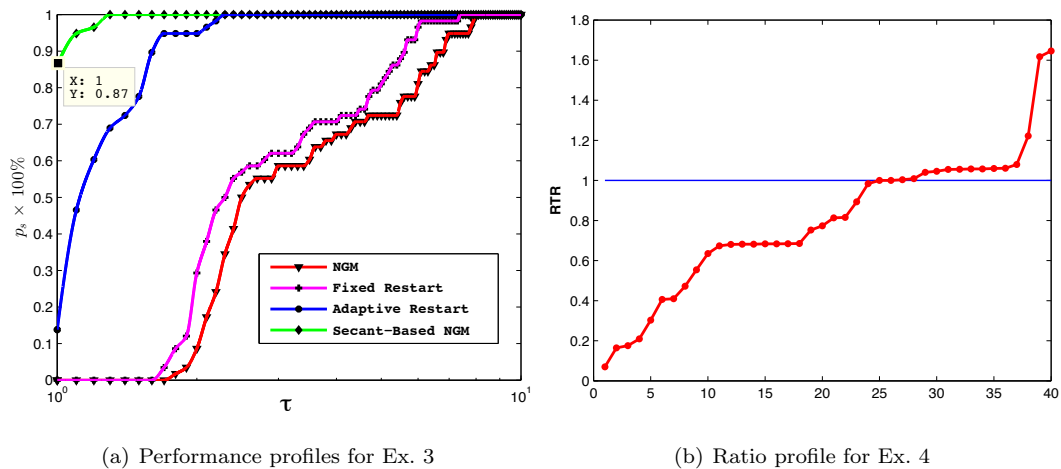


Fig. 12 The performance profiles and ratio profile of the set of solvers \mathcal{S} for Ex. 3 and Ex. 4 respectively. RTR is the ratio of the Secant-Based NGM's run-time to the Adaptive Restart's run-time.

7 Conclusion

The Nesterov gradient method (NGM) needs to be restarted (e.g. [18, 28]) when a nontrivial convexity parameter is not available. This paper introduces a new secant-based Nesterov gradient method (**Secant-Based-NGM**) and also establishes that it is globally convergent for all convex functions. The algorithm only requires a trivial lower bound of the convexity parameter $\mu \geq 0$ and the gradient's Lip-

schitz constant. The efficiency of the proposed algorithm derives from updating the estimate-sequence parameter γ_k by imposing a secant condition on the choice of search point y_k . Furthermore, the proposed **Secant-Based-NGM** embodies an "update rule with reset" that parallels the restart rule suggested in [28]. The effectiveness of the proposed algorithm is confirmed in numerical computations involving large dataset of test problems with varying dimension and condition number. The proposed **Secant-Based-NGM** significantly improves over and above the classical NGM and when restarts (e.g. [18, 28]) are used with the Nesterov gradient method.

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8 Appendix

See Table 1.

Table 1 Performance of solvers S on Ex. 4. All algorithms were stopped after 500s.

S/N	Test Problem	Secant-Based-NGM RT(s)	Adaptive Restart RT(s)	Fixed Restart RT(s)	Classical NGM RT(s)
1	009 'BOYD1'	4.50	5.04	66.81	67.03
2	017 'CVXQP1.L'	234.41	142.38	169.84	415.49
3	018 'CVXQP1.M'	5.16	5.14	13.87	18.89
4	028 'DUAL1'	0.17	0.10	0.89	1.11
5	032 'DUALC1'	0.61	0.79	26.58	25.95
6	035 'DUALC8'	0.17	0.23	1.81	2.48
7	039 'GOULDQP3'	0.22	0.23	0.28	0.99
8	042 'HS268'	0.53	0.65	5.95	5.86
9	069 'PRIMAL1'	10.09	14.76	16.78	16.53
10	070 'PRIMAL2'	10.96	16.09	18.26	17.96
11	071 'PRIMAL3'	11.28	16.50	18.81	18.50
12	072 'PRIMAL4'	13.44	19.61	22.38	22.14
13	073 'PRIMALC1'	9.78	14.51	16.52	16.26
14	074 'PRIMALC2'	9.73	14.27	16.19	15.95
15	075 'PRIMALC5'	10.01	14.63	16.60	16.40
16	076 'PRIMALC8'	10.62	15.58	17.71	17.41
17	079 'QAFIRO'	32.65	58.93	54.56	53.63
18	080 'QBANDM'	47.61	44.87	43.45	42.75
19	081 'QBEACONF'	500.00	500.00	227.99	224.73
20	083 'QBRANDY'	58.08	55.81	53.99	53.35
21	084 'QCAPRI'	78.65	74.56	72.25	71.33
22	085 'QE226'	82.19	500.00	388.40	383.64
23	087 'QFFFFFF80'	333.21	318.67	355.52	345.00
24	088 'QFORPLAN'	160.76	152.01	147.27	145.41
25	090 'QGROW15'	203.03	500.00	340.63	335.75
26	091 'QGROW22'	264.73	250.82	242.93	240.55
27	092 'QGROW7'	121.38	114.55	110.63	109.59
28	098 'QPILOTNO'	253.61	251.34	254.42	225.82
29	100 'QRECIPE'	34.76	500.00	58.01	57.20
30	104 'QSCFXM1'	151.47	500.00	255.69	252.71
31	105 'QSCFXM2'	204.93	500.00	344.88	341.66
32	106 'QSCFXM3'	236.10	500.00	397.89	394.56
33	109 'QSCSD1'	87.64	500.00	147.92	145.46
34	110 'QSCSD6'	104.74	500.00	176.15	174.71
35	111 'QSCSD8'	189.45	179.14	173.31	171.69
36	114 'QSCTAP3'	500.00	500.00	500.00	408.50
37	117 'QSHARE2B'	88.83	82.29	79.56	78.61
38	128 'S268'	0.53	0.65	5.94	5.84
39	132 'STCQP1'	0.17	0.14	1.20	1.46
40	138 'ZECEVIC2'	18.17	28.61	30.20	29.82