



An Inexact Regularized Newton Framework with a Worst-Case Iteration Complexity of $\mathcal{O}(\epsilon^{-3/2})$ for Nonconvex Optimization

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COR@L Technical Report 17T-011-R2



An Inexact Regularized Newton Framework with a Worst-Case Iteration Complexity of $\mathcal{O}(\epsilon^{-3/2})$ for Nonconvex Optimization*

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Original Publication: September 27, 2017

Last Revised: March 14, 2018

Abstract

An algorithm for solving smooth nonconvex optimization problems is proposed that, in the worst-case, takes $\mathcal{O}(\epsilon^{-3/2})$ iterations to drive the norm of the gradient of the objective function below a prescribed positive real number ϵ and can take $\mathcal{O}(\epsilon^{-3})$ iterations to drive the leftmost eigenvalue of the Hessian of the objective above $-\epsilon$. The proposed algorithm is a general framework that covers a wide range of techniques including quadratically and cubically regularized Newton methods, such as the Adaptive Regularisation using Cubics (ARC) method and the recently proposed Trust-Region Algorithm with Contractions and Expansions (TRACE). The generality of our method is achieved through the introduction of generic conditions that each trial step is required to satisfy, which in particular allow for inexact regularized Newton steps to be used. These conditions center around a new subproblem that can be approximately solved to obtain trial steps that satisfy the conditions. A new instance of the framework, distinct from ARC and TRACE, is described that may be viewed as a hybrid between quadratically and cubically regularized Newton methods. Numerical results demonstrate that our hybrid algorithm outperforms a cubically regularized Newton method. unconstrained optimization, nonlinear optimization, nonconvex optimization, inexact Newton methods, worst-case iteration complexity, worst-case evaluation complexity

1 Introduction

This paper proposes an algorithm for solving

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

*This work was supported in part by the U.S. Department of Energy, Office of Science, Applied Mathematics, Early Career Research Program under Award Number DE-SC0010615 and by the U.S. National Science Foundation, Division of Mathematical Sciences, Computational Mathematics Program under Award Number DMS-1016291.

where the (possibly nonconvex) objective function $f : \mathbb{R}^n \rightarrow \mathbb{R}$ is assumed to be twice-continuously differentiable. The optimization problem (1) has been widely studied, as evidenced by its appearance as the focal point of numerous textbooks; e.g., see [1], [2], [10], [20], [23], and [24].

For many years, the most popular methods for solving (1) were in classes known as line search and trust region methods. Recently, however, cubic regularization methods have become popular, which are based on the pioneering work by [19] and [21]. Their rise in popularity is due to increased interest in algorithms with improved complexity properties, which stems from the impact of so-called optimal algorithms for solving convex optimization problems. For problem (1), by complexity properties, we mean a guaranteed bound on the number of iterations (or function evaluations or derivative evaluations) needed by an algorithm before the norm of the gradient of the objective must fall below a positive threshold $\epsilon > 0$. In other words, if x_k denotes the k th iteration of an algorithm, one seeks a bound on the number of iterations until it is guaranteed that

$$\|\nabla f(x_k)\| \leq \epsilon.$$

The complexity of a traditional trust region method (e.g., see Algorithm 6.1.1 in [10]) is $\mathcal{O}(\epsilon^{-2})$ (see [5]), which falls short of the $\mathcal{O}(\epsilon^{-3/2})$ complexity for cubic regularization methods (e.g., see the ARC method by [6, 7]). This latter complexity is optimal among a certain broad class of second-order methods when employed to minimize a broad class of objective functions; see [8]. That said, one can obtain even better complexity properties if higher-order derivatives are used; see [4] and [9].

The better complexity properties of regularization methods such as ARC have been a major point of motivation for discovering other methods that attain the same worst-case iteration complexity bounds. For example, the recently introduced (nontraditional) trust region method known as TRACE (see [11]) has the same optimal $\mathcal{O}(\epsilon^{-3/2})$ complexity, while at the same time allowing traditional trust region trial steps to be computed and used. A key aspect of the TRACE framework is that a solution to an implicit trust region problem is obtained by varying a regularization parameter instead of a trust region radius. This key idea has been adopted and advanced further by [3]; in particular, they propose an algorithm that has optimal iteration complexity by solving quadratic subproblems that have a carefully chosen quadratic regularization parameter.

Contributions The main contributions of this paper relate to advancing the understanding of optimal complexity algorithms for solving the smooth optimization problem (1). Our proposed framework is intentionally very general; it is not a trust region method, a quadratic regularization method, or a cubic regularization method. Rather, we propose a generic set of conditions that each trial step must satisfy that still allow us to establish an optimal first-order complexity result as well as a second-order complexity bound similar to the methods above. Our framework contains as special cases other optimal complexity algorithms such as ARC and TRACE. To highlight this generality of our contribution, we describe one particular instance of our framework that appears to be new to the literature.

During the final preparation of this article, we came across the work in [13] and [14]. This work shares certain commonalities with our own and appears to have been developed at the same time. Although there are numerous differences, we shall only point out three of them. First, the precise conditions that they require for each trial step are different from ours. In particular, the condition stated as (3.1c) in [14] requires that regularization is used to compute every trial step, a property not shared by our method (which can employ Newton steps). Second, they do not consider second-order convergence or complexity properties, although they might be able to do so by incorporating second-order conditions similar to ours. Third, they focus on strategies for identifying an appropriate value for the regularization parameter. An implementation of our method might consider their proposals, but could employ other strategies as well. In any case, overall, we believe that our papers are quite distinct, and in some ways are complementary.

Organization In §2, we present our general framework that is formally stated as Algorithm 1. In §3, we prove that our framework enjoys first-order convergence (see §3.1), an optimal first-order complexity (see §3.2), and certain second-order convergence and complexity guarantees (see §3.3). In §4, we show that ARC and TRACE can be viewed as special cases of our framework, and present yet another instance that is distinct

from these methods. In §5, we present details of implementations of a cubic regularization method and our newly proposed instance of our framework, and provide the results of numerical experiments with both. Finally, in §6, we present final comments.

Notation We use \mathbb{R}_+ to denote the set of nonnegative scalars, \mathbb{R}_{++} to denote the set of positive scalars, and \mathbb{N}_+ to denote the set of nonnegative integers. Given a real symmetric matrix A , we write $A \succeq 0$ (respectively, $A \succ 0$) to indicate that A is positive semidefinite (respectively, positive definite). Given a pair of scalars $(a, b) \in \mathbb{R} \times \mathbb{R}$, we write $a \perp b$ to indicate that $ab = 0$. Similarly, given such a pair, we denote their maximum as $\max\{a, b\}$ and their minimum as $\min\{a, b\}$. Given a vector v , we denote its (Euclidean) ℓ_2 -norm as $\|v\|$. Finally, given a discrete set \mathcal{S} , we denote its cardinality by $|\mathcal{S}|$.

Corresponding to the objective $f : \mathbb{R}^n \rightarrow \mathbb{R}$, we define the gradient function $g := \nabla f : \mathbb{R}^n \rightarrow \mathbb{R}^n$ and the Hessian function $H := \nabla^2 f : \mathbb{R}^n \rightarrow \mathbb{R}^{n \times n}$. Given an iterate x_k in an algorithm for solving (1), we define $f_k := f(x_k)$, $g_k := g(x_k) := \nabla f(x_k)$, and $H_k := H(x_k) := \nabla^2 f(x_k)$. Similarly, we apply a subscript to other algorithmic quantities whose definition depends on the iteration number k .

2 Algorithm Description

Our algorithm involves generic conditions that a trial step toward solving problem (1) must satisfy. One can obtain a step satisfying these conditions by computing—for appropriate positive lower and upper bounds σ_k^L and σ_k^U , respectively, on the ratio between a regularization variable $\lambda \geq 0$ and the norm of the trial step—an approximate solution of the subproblem

$$\begin{aligned} \mathcal{P}_k(\sigma_k^L, \sigma_k^U) &: \min_{(s, \lambda) \in \mathbb{R}^n \times \mathbb{R}_+} f_k + g_k^T s + \frac{1}{2} s^T (H_k + \lambda I) s \\ &\text{s.t. } (\sigma_k^L)^2 \|s\|^2 \leq \lambda^2 \leq (\sigma_k^U)^2 \|s\|^2. \end{aligned} \quad (2)$$

For a given value of the regularization variable λ , this problem involves a quadratic objective function and an upper bound on the norm of the trial step, just as in a trust region method. However, it also includes a lower bound on the norm of the trial step, and, in general, with λ as a variable, it encapsulates other types of subproblems as well, including those present in a cubic regularization framework. For additional details on the properties of this subproblem and its solutions, see Appendices A and B.

The conditions that the k th trial step and regularization pair, i.e., (s_k, λ_k) , must satisfy are stated in Assumption 2.1 below, wherein we invoke the following (unregularized) quadratic model of f at x_k :

$$q_k(s) := f_k + g_k^T s + \frac{1}{2} s^T H_k s.$$

Assumption 2.1. *The pair (s_k, λ_k) is computed such that it is feasible for problem (2) and, with*

$$\Delta_k(s_k, \lambda_k) := \begin{cases} \|s_k\| & \text{if } \lambda_k = 0 \\ \frac{1}{\sqrt{6}} \sqrt{\frac{\|g_k\| \|s_k\|}{\lambda_k}} & \text{if } \lambda_k > 0 \end{cases} \quad (3)$$

and constants $(\kappa_1, \kappa_2, \kappa_3) \in \mathbb{R}_{++} \times \mathbb{R}_{++} \times \mathbb{R}_{++}$, the following hold:

$$f_k - q_k(s_k) \geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + \|H_k\|}, \Delta_k(s_k, \lambda_k) \right\}; \quad (4a)$$

$$s_k^T (g_k + (H_k + \lambda_k I) s_k) \leq \min \{ \kappa_1 \|s_k\|^2, \frac{1}{2} s_k^T (H_k + \lambda_k I) s_k + \frac{1}{2} \kappa_2 \|s_k\|^3 \}; \quad \text{and} \quad (4b)$$

$$\|g_k + (H_k + \lambda_k I) s_k\| \leq \lambda_k \|s_k\| + \kappa_3 \|s_k\|^2. \quad (4c)$$

To see that Assumption 2.1 is well-posed and consistent with problem (2), we refer the reader to Theorem B.3 in Appendix B wherein we prove that any solution of problem (2) with s restricted to a sufficiently large dimensional subspace of \mathbb{R}^n satisfies all of the conditions in Assumption 2.1. We also claim that one can obtain a pair satisfying Assumption 2.1 in either of the following two ways:

- Choose $\sigma \in [\sigma_k^L, \sigma_k^U]$, compute s_k by minimizing the cubic function

$$c_k(s; \sigma) := q_k(s) + \frac{1}{2}\sigma\|s\|^3 = f_k + g_k^T s + \frac{1}{2}s^T H_k s + \frac{1}{2}\sigma\|s\|^3 \quad (5)$$

over a sufficiently large dimensional subspace of \mathbb{R}^n (assuming, when $\sigma = \sigma_k^L = 0$, that this function is not unbounded below), then set $\lambda_k \leftarrow \sigma\|s_k\|$. This is essentially the strategy employed in cubic regularization methods such as ARC.

- Choose $\lambda_k \geq 0$, then compute s_k by minimizing the objective of (2) with $\lambda = \lambda_k$ over a sufficiently large dimensional subspace of \mathbb{R}^n (assuming that the function is not unbounded below). The resulting pair (s_k, λ_k) satisfies Assumption 2.1 as long as it is feasible for (2). This is essentially the strategy employed in [3] and partly employed in TRACE.

One can imagine other approaches as well. Overall, we state problem (2) as a guide for various techniques for computing the pair (s_k, λ_k) . Our theory simply relies on the fact that any such computed pair satisfies the conditions in Assumption 2.1.

Our algorithm, stated as Algorithm 1, employs the following ratio (also employed, e.g., in TRACE) to determine whether a given trial step is accepted or rejected:

$$\rho_k := \frac{f_k - f(x_k + s_k)}{\|s_k\|^3}.$$

One potential drawback of employing this ratio is that the ratio is not invariant to scaling of the objective function. However, the use of this ratio can still be justified. For example, if one were to compute s_k by minimizing the cubic model (5) for some $\sigma > 0$, then the reduction in this model yielded by s_k is bounded below by a fraction of $\sigma\|s_k\|^3$ (see [7, Lemma 4.2]), meaning that $\rho_k \geq \eta$ holds when $\sigma \geq \eta$ and the actual reduction in f is proportional to the reduction in the cubic model. For further justification for this choice—such as how it allows the algorithm to accept Newton steps when the norm of the trial step is small (and, indeed, the norms of accepted steps vanish asymptotically as shown in Lemma 3.7 later on)—we refer the reader to [3] and [11].

3 Convergence Analysis

In this section, we prove global convergence guarantees for Algorithm 1. In particular, we prove under common assumptions that, from remote starting points, the algorithm converges to first-order stationarity, has a worst-case iteration complexity to approximate first-order stationarity that is on par with the methods in [7], [11], and [3], and—at least in a subspace determined by the search path of the algorithm—converges to second-order stationarity with a complexity on par with the methods in [7] and [11].

3.1 First-Order Global Convergence

Our goal in this subsection is to prove that the sequence of objective gradients vanishes. We make the following assumption about the objective function, which is assumed to hold throughout this section.

Assumption 3.1. *The objective function $f : \mathbb{R}^n \rightarrow \mathbb{R}$ is twice continuously differentiable and bounded below by a scalar $f_{\inf} \in \mathbb{R}$ on \mathbb{R}^n .*

We also make the following assumption related to the sequence of iterates.

Assumption 3.2. *The gradient function $g : \mathbb{R}^n \rightarrow \mathbb{R}^n$ is Lipschitz continuous with Lipschitz constant $g_{\text{Lip}} \in \mathbb{R}_{++}$ in an open convex set containing the sequences $\{x_k\}$ and $\{x_k + s_k\}$. Furthermore, the gradient sequence $\{g_k\}$ has $g_k \neq 0$ for all $k \in \mathbb{N}_+$ and is bounded in that there exists a scalar constant $g_{\max} \in \mathbb{R}_{++}$ such that $\|g_k\| \leq g_{\max}$ for all $k \in \mathbb{N}_+$.*

Algorithm 1 Inexact Regularized Newton Framework

Require: an acceptance constant $\eta \in \mathbb{R}_{++}$ with $0 < \eta < 1$

Require: bound update constants $\{\gamma_1, \gamma_2\} \subset \mathbb{R}_{++}$ with $1 < \gamma_1 \leq \gamma_2$

Require: ratio lower and upper bound constants $\{\underline{\sigma}, \bar{\sigma}\} \subset \mathbb{R}_{++}$ such that $\bar{\sigma} \geq \underline{\sigma}$

```

1: procedure INEXACT REGULARIZED NEWTON
2:   set  $x_0 \in \mathbb{R}^n$ 
3:   set  $\sigma_0^L \leftarrow 0$  and  $\sigma_0^U \in [\underline{\sigma}, \bar{\sigma}]$ 
4:   for  $k \in \mathbb{N}_+$  do
5:     set  $(s_k, \lambda_k)$  satisfying Assumption 2.1
6:     if  $\rho_k \geq \eta$  then [accept step]
7:       set  $x_{k+1} \leftarrow x_k + s_k$ 
8:       set  $\sigma_{k+1}^L \leftarrow 0$  and  $\sigma_{k+1}^U \leftarrow \sigma_k^U$ 
9:     else (i.e.,  $\rho_k < \eta$ ) [reject step]
10:      set  $x_{k+1} \leftarrow x_k$ 
11:      if  $\lambda_k < \underline{\sigma}\|s_k\|$  then
12:        set  $\sigma_{k+1}^L \in [\underline{\sigma}, \bar{\sigma}]$  and  $\sigma_{k+1}^U \in [\sigma_{k+1}^L, \bar{\sigma}]$ 
13:      else
14:        set  $\sigma_{k+1}^L \leftarrow \gamma_1 \frac{\lambda_k}{\|s_k\|}$  and  $\sigma_{k+1}^U \leftarrow \gamma_2 \frac{\lambda_k}{\|s_k\|}$ 
15:      end if
16:    end if
17:   end for
18: end procedure

```

It is worthwhile to note in passing that our complexity bounds for first- and second-order stationarity remain true even if one were to consider the possibility that $g_k = 0$ for some $k \in \mathbb{N}_+$, in which case one would have the algorithm terminate finitely or, if $H_k \not\succeq 0$, compute an improving direction of negative curvature for H_k . However, allowing this possibility—which is typically unlikely ever to occur in practice—would only serve to obscure certain aspects of our analysis. We refer the reader, e.g., to [7] (specifically, to the discussions at the ends of §2.1, §4, and §5 in that work) for commentary about why zero gradient values do not ruin complexity guarantees such as we present.

We begin with two lemmas each revealing an important consequence of Assumptions 3.1 and 3.2.

Lemma 3.1. *For all $k \in \mathbb{N}_+$, it follows that $s_k \neq 0$.*

Proof. The result follows by combining that $g_k \neq 0$ for all $k \in \mathbb{N}_+$ (see Assumption 3.2) with (4c). \square

Lemma 3.2. *The Hessian sequence $\{H_k\}$ is bounded in norm in that there exists a scalar constant $H_{max} \in \mathbb{R}_{++}$ such that $\|H_k\| \leq H_{max}$ for all $k \in \mathbb{N}_+$.*

Proof. The result follows by Assumption 3.1, the Lipschitz continuity of g in Assumption 3.2, and Lemma 1.2.2 in [22]. \square

In our next lemma, we prove an upper bound for the regularization variable λ_k .

Lemma 3.3. *For all $k \in \mathbb{N}_+$, the pair (s_k, λ_k) satisfies*

$$\lambda_k \leq 2 \frac{\|g_k\|}{\|s_k\|} + \frac{3}{2} H_{max} + \kappa_1.$$

Proof. Since (4a) ensures $q_k(s_k) - f_k \leq 0$, it follows with (4b) and Lemma 3.2 that

$$\begin{aligned} 0 &\geq q_k(s_k) - f_k = g_k^T s_k + \frac{1}{2} s_k^T H_k s_k \\ &\geq g_k^T s_k + \frac{1}{2} s_k^T H_k s_k + s_k^T (g_k + (H_k + \lambda_k I)s_k) - \kappa_1 \|s_k\|^2 \\ &= 2g_k^T s_k + \frac{3}{2} s_k^T H_k s_k + \lambda_k \|s_k\|^2 - \kappa_1 \|s_k\|^2 \\ &\geq -2\|g_k\| \|s_k\| - \frac{3}{2} H_{max} \|s_k\|^2 + \lambda_k \|s_k\|^2 - \kappa_1 \|s_k\|^2. \end{aligned}$$

After rearrangement and dividing by $\|s_k\|^2 \neq 0$ (see Lemma 3.1), the desired result follows. \square

Using Lemma 3.3, we now prove a lower bound for the reduction in q_k yielded by s_k .

Lemma 3.4. *For all $k \in \mathbb{N}_+$, the step s_k satisfies*

$$f_k - q_k(s_k) \geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + H_{max}}, \frac{\|s_k\|}{\sqrt{6}} \sqrt{\frac{\|g_k\|}{2\|g_k\| + \|s_k\|(\frac{3}{2}H_{max} + \kappa_1)}} \right\}.$$

Proof. If $\lambda_k = 0$, then by (4a) and Lemma 3.2 it follows that

$$f_k - q_k(s_k) \geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + \|H_k\|}, \|s_k\| \right\} \geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + H_{max}}, \|s_k\| \right\}.$$

On the other hand, if $\lambda_k > 0$, then (4a), Lemma 3.2, and Lemma 3.3 imply that

$$\begin{aligned} f_k - q_k(s_k) &\geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + \|H_k\|}, \frac{1}{\sqrt{6}} \sqrt{\frac{\|g_k\| \|s_k\|}{\lambda_k}} \right\} \\ &\geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + H_{max}}, \frac{\|s_k\|}{\sqrt{6}} \sqrt{\frac{\|g_k\|}{2\|g_k\| + \|s_k\|(\frac{3}{2}H_{max} + \kappa_1)}} \right\}. \end{aligned}$$

Combining the inequalities from these two cases, the desired result follows. \square

Going forward, for ease of reference, we respectively define sets of indices corresponding to accepted and rejected steps throughout a run of the algorithm as

$$\mathcal{A} := \{k \in \mathbb{N}_+ : \rho_k \geq \eta\} \quad \text{and} \quad \mathcal{R} := \{k \in \mathbb{N}_+ : \rho_k < \eta\}.$$

We now show that if the algorithm were only to compute rejected steps from some iteration onward, then the sequence $\{\lambda_k/\|s_k\|\}$ diverges to infinity.

Lemma 3.5. *If $k \in \mathcal{R}$ for all sufficiently large $k \in \mathbb{N}_+$, then $\{\lambda_k/\|s_k\|\} \rightarrow \infty$.*

Proof. Without loss of generality, assume that $\mathcal{R} = \mathbb{N}_+$. We now prove that the condition in Step 11 cannot be true more than once. Suppose, in iteration $\hat{k} \in \mathbb{N}_+$, Step 12 is reached, which means that $\lambda_{\hat{k}+1}/\|s_{\hat{k}+1}\| \geq \sigma$ since $(s_{\hat{k}+1}, \lambda_{\hat{k}+1})$ is required to be feasible for $\mathcal{P}_{\hat{k}+1}(\sigma_{\hat{k}+1}^L, \sigma_{\hat{k}+1}^U)$ in Step 5 where $\sigma_{\hat{k}+1}^L \geq \sigma$. Therefore, the condition in Step 11 tests false in iteration $(\hat{k}+1)$. Then, from Step 5, Step 14, and the fact that $\gamma_1 > 1$, it follows that $\{\lambda_k/\|s_k\|\}$ is monotonically increasing for all $k \geq \hat{k}$. Therefore, the condition in Step 11 cannot test true for any $k \geq \hat{k}+1$. Now, to see that the sequence diverges, notice from this fact, Step 5, and Step 14, it follows that for all $k \geq \hat{k}+1$ we have $\lambda_{k+1}/\|s_{k+1}\| \geq \gamma_1(\lambda_k/\|s_k\|)$ where $\gamma_1 > 1$. Thus, $\{\lambda_k/\|s_k\|\} \rightarrow \infty$, as claimed. \square

We now prove that if the gradients are bounded away from zero and the sequence of ratios $\{\lambda_k/\|s_k\|\}$ diverges, then $\rho_k \geq \eta$ for all sufficiently large $k \in \mathbb{N}_+$, meaning that the steps are accepted.

Lemma 3.6. Suppose that $\mathcal{I} \subseteq \mathbb{N}_+$ is an infinite index set such that for $\epsilon \in \mathbb{R}_{++}$ independent of k , one finds that $\|g_k\| \geq \epsilon$ for all $k \in \mathcal{I}$ and $\{\lambda_k/\|s_k\|\}_{k \in \mathcal{I}} \rightarrow \infty$. Then, for all sufficiently large $k \in \mathcal{I}$, it follows that $\rho_k \geq \eta$, meaning $k \in \mathcal{A}$.

Proof. From the Mean Value Theorem, there exists $\bar{x}_k \in [x_k, x_k + s_k]$ such that

$$\begin{aligned} q_k(s_k) - f(x_k + s_k) &= (g_k - g(\bar{x}_k))^T s_k + \frac{1}{2} s_k^T H_k s_k \\ &\geq -\|g_k - g(\bar{x}_k)\| \|s_k\| - \frac{1}{2} \|H_k\| \|s_k\|^2. \end{aligned} \quad (6)$$

From this, Lemma 3.4, and Assumption 3.2, it follows that, for all $k \in \mathcal{I}$,

$$\begin{aligned} f_k - f(x_k + s_k) &= f_k - q_k(s_k) + q_k(s_k) - f(x_k + s_k) \\ &\geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + H_{max}}, \frac{\|s_k\|}{\sqrt{6}} \sqrt{\frac{\|g_k\|}{2\|g_k\| + \|s_k\|(\frac{3}{2}H_{max} + \kappa_1)}} \right\} - (g_{Lip} + \frac{1}{2}H_{max}) \|s_k\|^2 \\ &\geq \frac{\epsilon}{6\sqrt{2}} \min \left\{ \frac{\epsilon}{1 + H_{max}}, \frac{\|s_k\|}{\sqrt{6}} \sqrt{\frac{\epsilon}{2g_{max} + \|s_k\|(\frac{3}{2}H_{max} + \kappa_1)}} \right\} - (g_{Lip} + \frac{1}{2}H_{max}) \|s_k\|^2. \end{aligned}$$

This shows that there exists a threshold $s_{thresh} > 0$ such that

$$f_k - f(x_k + s_k) \geq \eta \|s_k\|^3 \text{ whenever } k \in \mathcal{I} \text{ and } \|s_k\| \leq s_{thresh}.$$

We now claim that $\{\|s_k\|\}_{k \in \mathcal{I}} \rightarrow 0$. To prove this claim, suppose by contradiction that there exists an infinite subsequence $\mathcal{I}_s \subseteq \mathcal{I}$ and scalar $\epsilon_s \in \mathbb{R}_{++}$ such that $\|s_k\| \geq \epsilon_s$ for all $k \in \mathcal{I}_s$. It then follows from the boundedness of $\{\|g_k\|\}$ (see Assumption 3.2) and Lemma 3.3 that $\{\lambda_k\}_{k \in \mathcal{I}_s}$ is bounded. This allows us to conclude that $\{\lambda_k/\|s_k\|\}_{k \in \mathcal{I}_s}$ is bounded, which contradicts the assumptions of the lemma. Thus, $\{\|s_k\|\}_{k \in \mathcal{I}} \rightarrow 0$. Hence, there exists $k_s \in \mathcal{I}$ such that for all $k \in \mathcal{I}$ with $k \geq k_s$ one finds $\|s_k\| \leq s_{thresh}$. Therefore, for all $k \in \mathcal{I}$ with $k \geq k_s$, it follows that $\rho_k \geq \eta$, as claimed. \square

Next, we prove that the algorithm produces infinitely many accepted steps.

Lemma 3.7. It holds that $|\mathcal{A}| = \infty$ and $\{s_k\}_{k \in \mathcal{A}} \rightarrow 0$.

Proof. To derive a contradiction, suppose that $|\mathcal{A}| < \infty$. This implies that there exists k_0 such that, for all $k \geq k_0$, one has $k \in \mathcal{R}$ and $(x_k, g_k, H_k) = (x_{k_0}, g_{k_0}, H_{k_0})$. From this fact and Assumption 3.2, it follows that $\|g_k\| \geq \epsilon$ for all $k \geq k_0$ for some $\epsilon \in \mathbb{R}_{++}$. From the fact that $k \in \mathcal{R}$ for all $k \geq k_0$ and Lemma 3.5, it follows that $\{\lambda_k/\|s_k\|\} \rightarrow \infty$. This fact and $\|g_k\| \geq \epsilon$ for all $k \geq k_0$ imply that all the conditions of Lemma 3.6 are satisfied for $\mathcal{I} := \{k \in \mathbb{N}_+ : k \geq k_0\}$; therefore, Lemma 3.6 implies that for all sufficiently large $k \in \mathcal{I}$, one finds $\rho_k \geq \eta$ so that $k \in \mathcal{A}$, a contradiction.

To complete the proof, notice that the objective function values are monotonically decreasing. Combining this with the condition in Step 6, the fact that f is bounded below by f_{inf} (see Assumption 3.1), and $|\mathcal{A}| = \infty$, one deduces that $\{s_k\}_{k \in \mathcal{A}} \rightarrow 0$, as claimed. \square

We now prove that there exists an infinite subsequence of iterates such that the sequence of gradients computed at those points converges to zero.

Lemma 3.8. It holds that

$$\liminf_{k \in \mathbb{N}_+, k \rightarrow \infty} \|g_k\| = 0.$$

Proof. To derive a contradiction, suppose that $\liminf_{k \in \mathbb{N}_+, k \rightarrow \infty} \|g_k\| > 0$, which along with the fact that $g_{k+1} = g_k$ for any $k \in \mathbb{N}_+ \setminus \mathcal{A}$ means $\liminf_{k \in \mathcal{A}, k \rightarrow \infty} \|g_k\| > 0$. Thus, there exists $\epsilon \in \mathbb{R}_{++}$ such that

$$\|g_k\| \geq \epsilon \text{ for all sufficiently large } k \in \mathcal{A}. \quad (7)$$

Under (7), let us prove that $\{\lambda_k\}_{k \in \mathcal{A}} \rightarrow \infty$. To derive a contradiction, suppose there exists an infinite $\mathcal{A}_\lambda \subseteq \mathcal{A}$ such that $\lambda_k \leq \lambda_{max}$ for some $\lambda_{max} \in \mathbb{R}_{++}$. On the other hand, by $\{s_k\}_{k \in \mathcal{A}} \rightarrow 0$ (see Lemma 3.7) and (4c), it follows that $\{g_k + (H_k + \lambda_k I)s_k\}_{k \in \mathcal{A}_\lambda} \rightarrow 0$. Combining the upper bound on $\{\lambda_k\}_{k \in \mathcal{A}_\lambda}$, the fact that $\{s_k\}_{k \in \mathcal{A}} \rightarrow 0$, and $\|H_k\| \leq H_{max}$ (see Lemma 3.2), it follows that $\{g_k\}_{k \in \mathcal{A}_\lambda} \rightarrow 0$, which violates (7). Therefore, $\{\lambda_k\}_{k \in \mathcal{A}} \rightarrow \infty$.

Our next goal is to prove, still under (7), that $k \in \mathcal{A}$ for all sufficiently large $k \in \mathbb{N}_+$. To prove this, our strategy is to show that the sets of iterations involving a rejected step followed by an accepted step are finite. In particular, let us define the index sets

$$\begin{aligned}\mathcal{R}_1 &:= \{k \in \mathcal{R} : \text{the condition in Step 11 tests true and } (k+1) \in \mathcal{A}\} \text{ and} \\ \mathcal{R}_2 &:= \{k \in \mathcal{R} : \text{the condition in Step 11 tests false and } (k+1) \in \mathcal{A}\}.\end{aligned}$$

We aim to prove that these are finite. First, consider \mathcal{R}_1 . To derive a contradiction, suppose that $|\mathcal{R}_1| = \infty$. By definition, for all $k \in \mathcal{R}_1$, the condition in Step 11 tests true, meaning (s_{k+1}, λ_{k+1}) is found in Step 5 satisfying $\lambda_{k+1}/\|s_{k+1}\| \leq \bar{\sigma}$. On the other hand, since $(k+1) \in \mathcal{A}$ for all $k \in \mathcal{R}_1$, it follows from Lemma 3.7 that $\{s_{k+1}\}_{k \in \mathcal{R}_1} \rightarrow 0$. Combining the conclusions of these last two sentences shows that $\{\lambda_{k+1}\}_{k \in \mathcal{R}_1} \rightarrow 0$. However, this contradicts the conclusion of the previous paragraph, which showed that $\{\lambda_k\}_{k \in \mathcal{A}} \rightarrow \infty$. Hence, we may conclude that $|\mathcal{R}_1| < \infty$. Now consider \mathcal{R}_2 . To derive a contradiction, suppose that $|\mathcal{R}_2| = \infty$. The fact that the condition in Step 11 tests false for $k \in \mathcal{R}_2$ implies that (s_{k+1}, λ_{k+1}) is found in Step 5 satisfying $\lambda_{k+1}/\|s_{k+1}\| \leq \gamma_2 \lambda_k/\|s_k\|$. However, since $\{s_{k+1}\}_{k \in \mathcal{R}_2} \rightarrow 0$ (see Lemma 3.7) and $\{\lambda_{k+1}\}_{k \in \mathcal{R}_2} \rightarrow \infty$ (established in the previous paragraph), it follows that $\{\lambda_{k+1}/\|s_{k+1}\|\}_{k \in \mathcal{R}_2} \rightarrow \infty$, which combined with the previously established inequality $\lambda_{k+1}/\|s_{k+1}\| \leq \gamma_2 \lambda_k/\|s_k\|$ shows that $\{\lambda_k/\|s_k\|\}_{k \in \mathcal{R}_2} \rightarrow \infty$. Therefore, with (7), the conditions in Lemma 3.6 hold for $\mathcal{I} = \mathcal{R}_2$, meaning that, for all sufficiently large $k \in \mathcal{R}_2$, the inequality $\rho_k \geq \eta$ holds. This contradicts the fact that $\mathcal{R}_2 \subseteq \mathcal{R}$; hence, we conclude that \mathcal{R}_2 is finite. Since \mathcal{R}_1 and \mathcal{R}_2 are finite, it follows from the logic of Algorithm 1 that either $k \in \mathcal{A}$ for all sufficiently large k or $k \in \mathcal{R}$ for all sufficiently large k . By Lemma 3.7, it follows that $k \in \mathcal{A}$ for all sufficiently large k .

Thus far, we have proved under (7) that $\{\lambda_k\}_{k \in \mathcal{A}} \rightarrow \infty$ and that $k \in \mathcal{A}$ for all large $k \in \mathbb{N}_+$. From this latter fact, it follows that there exists k_σ such that $\sigma_k^U = \sigma_{k_\sigma}^U \in \mathbb{R}_{++}$ for all $k \geq k_\sigma$. In addition, from Step 5, it follows that for $k \geq k_\sigma$ one finds $\lambda_k/\|s_k\| \leq \sigma_k^U = \sigma_{k_\sigma}^U < \infty$. However, this leads to a contradiction to the facts that $\{\lambda_k\}_{k \in \mathcal{A}} \rightarrow \infty$ and $\{s_k\}_{k \in \mathcal{A}} \rightarrow 0$ (see Lemma 3.7). Overall, we have shown that (7) cannot be true, which proves the desired result. \square

We close with our main global convergence result of this subsection, the proof of which borrows much from that of Theorem 3.14 in [11].

Theorem 3.3. *Under Assumptions 2.1, 3.1, and 3.2, it follows that*

$$\lim_{k \in \mathbb{N}_+, k \rightarrow \infty} \|g_k\| = 0. \quad (8)$$

Proof. For the purpose of reaching a contradiction, suppose that (8) does not hold. Combining this with the fact that $|\mathcal{A}| = \infty$ (see Lemma 3.7), it follows that there exists an infinite subsequence $\{t_i\} \subseteq \mathcal{A}$ (indexed over $i \in \mathbb{N}_+$) and a scalar $\epsilon > 0$ such that, for all $i \in \mathbb{N}_+$, one finds $\|g_{t_i}\| \geq 2\epsilon > 0$. Also, the fact that $|\mathcal{A}| = \infty$ and Lemma 3.8 imply that there exists an infinite subsequence $\{\ell_i\} \subseteq \mathcal{A}$ (indexed over $i \in \mathbb{N}_+$) such that, for all $i \in \mathbb{N}_+$ and $k \in \mathbb{N}_+$ with $t_i \leq k < \ell_i$, one finds

$$\|g_k\| \geq \epsilon \text{ and } \|g_{\ell_i}\| < \epsilon. \quad (9)$$

Let us now restrict our attention to indices in the infinite index set

$$\mathcal{K} := \{k \in \mathcal{A} : t_i \leq k < \ell_i \text{ for some } i \in \mathbb{N}_+\}.$$

Observe from (9) that, for all $k \in \mathcal{K}$, it follows that $\|g_k\| \geq \epsilon$. Also, from the definition of \mathcal{A} ,

$$f_k - f_{k+1} \geq \eta \|s_k\|^3 \text{ for all } k \in \mathcal{K} \subseteq \mathcal{A}. \quad (10)$$

Since $\{f_k\}$ is monotonically decreasing and bounded below, one finds that $\{f_k\} \rightarrow \underline{f}$ for some $\underline{f} \in \mathbb{R}$, which when combined with (10) shows that

$$\lim_{k \in \mathcal{K}, k \rightarrow \infty} \|s_k\| = 0. \quad (11)$$

Using this fact, Lemma 3.4, Assumption 3.2, and the Mean Value Theorem (as it is used in the proof of Lemma 3.6 to yield (6)), it follows that for all sufficiently large $k \in \mathcal{K}$ one has

$$\begin{aligned} f_k - f_{k+1} &= f_k - q_k(s_k) + q_k(s_k) - f(x_k + s_k) \\ &\geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + H_{max}}, \frac{\|s_k\|}{\sqrt{6}} \sqrt{\frac{\|g_k\|}{2\|g_k\| + \|s_k\|(\frac{3}{2}H_{max} + \kappa_1)}} \right\} - (g_{Lip} + \frac{1}{2}H_{max})\|s_k\|^2 \\ &\geq \frac{\epsilon}{6\sqrt{2}} \frac{\|s_k\|}{\sqrt{6}} \sqrt{\frac{\|g_k\|}{2\|g_k\| + \|s_k\|(\frac{3}{2}H_{max} + \kappa_1)}} - (g_{Lip} + \frac{1}{2}H_{max})\|s_k\|^2. \end{aligned}$$

It now follows from (9) and (11) that, as $k \rightarrow \infty$ over $k \in \mathcal{K}$, the square root term in the previous inequality converges to $1/\sqrt{2}$. Since the second term in the previous inequality is of order $\|s_k\|^2$, the first term is of order $\|s_k\|$, and $1/\sqrt{2} > 1/\sqrt{3}$, one can thus conclude that $f_k - f_{k+1} \geq \epsilon\|s_k\|/36$ for all sufficiently large $k \in \mathcal{K}$. Consequently, it follows that for all sufficiently large $i \in \mathbb{N}_+$ one finds

$$\begin{aligned} \|x_{t_i} - x_{\ell_i}\| &\leq \sum_{k \in \mathcal{K}, k=t_i}^{\ell_i-1} \|x_k - x_{k+1}\| \\ &= \sum_{k \in \mathcal{K}, k=t_i}^{\ell_i-1} \|s_k\| \leq \sum_{k \in \mathcal{K}, k=t_i}^{\ell_i-1} \frac{36}{\epsilon} (f_k - f_{k+1}) = \frac{36}{\epsilon} (f_{t_i} - f_{\ell_i}). \end{aligned}$$

Since $\{f_{t_i} - f_{\ell_i}\} \rightarrow 0$ (recall that $\{f_k\} \rightarrow \underline{f}$ monotonically) this implies that $\{\|x_{t_i} - x_{\ell_i}\|\} \rightarrow 0$, which, in turn, implies that $\{\|g_{t_i} - g_{\ell_i}\|\} \rightarrow 0$ because of the continuity of g . However, this is a contradiction since, for any $i \in \mathbb{N}_+$, we have $\|g_{t_i} - g_{\ell_i}\| \geq \epsilon$ by the definitions of $\{t_i\}$ and $\{\ell_i\}$. Overall, we conclude that our initial supposition must be false, implying that (8) holds. \square

3.2 First-Order Complexity

Our next goal is to prove, with respect to a prescribed positive threshold, a worst-case upper bound on the number of iterations required for our algorithm to reduce the norm of the gradient below the threshold. In this subsection, along with Assumptions 2.1, 3.1, and 3.2, we add the following.

Assumption 3.4. *The Hessian function H is Lipschitz continuous on a path defined by the sequence of iterates and trial steps; in particular, it is Lipschitz continuous with a scalar Lipschitz constant $H_{Lip} > 0$ on the set $\{x_k + \tau s_k : k \in \mathbb{N}_+, \tau \in [0, 1]\}$.*

We begin our analysis in this subsection by providing a lemma that shows that successful steps always result if λ_k is sufficiently large relative to the size of the step.

Lemma 3.9. *For any $k \in \mathbb{N}_+$, if the pair (s_k, λ_k) satisfies*

$$\lambda_k \geq (H_{Lip} + \kappa_2 + 2\eta)\|s_k\|, \quad (12)$$

then $\rho_k \geq \eta$.

Proof. It follows from Assumption 3.4 and Taylor's expansion with Lagrange remainder that there exists \bar{x}_k on the line segment $[x_k, x_k + s_k]$ such that

$$q_k(s_k) - f(x_k + s_k) = \frac{1}{2}s_k^T(H_k - H(\bar{x}_k))s_k \geq -\frac{1}{2}H_{Lip}\|s_k\|^3. \quad (13)$$

Also, it follows from (4b) that

$$\begin{aligned}
f_k - q_k(s_k) &= -g_k^T s_k - \frac{1}{2} s_k^T H_k s_k \\
&= -s_k^T (g_k + (H_k + \lambda_k I)s_k) + \frac{1}{2} \lambda_k \|s_k\|^2 + \frac{1}{2} s_k^T (H_k + \lambda_k I)s_k \\
&\geq -\frac{1}{2} s_k^T (H_k + \lambda_k I)s_k - \frac{1}{2} \kappa_2 \|s_k\|^3 + \frac{1}{2} \lambda_k \|s_k\|^2 + \frac{1}{2} s_k^T (H_k + \lambda_k I)s_k \\
&= -\frac{1}{2} \kappa_2 \|s_k\|^3 + \frac{1}{2} \lambda_k \|s_k\|^2.
\end{aligned} \tag{14}$$

From (13) and (14), it follows that

$$\begin{aligned}
f_k - f(x_k + s_k) &= f_k - q_k(s_k) + q_k(s_k) - f(x_k + s_k) \\
&\geq \frac{1}{2} \lambda_k \|s_k\|^2 - \frac{1}{2} \kappa_2 \|s_k\|^3 - \frac{1}{2} H_{Lip} \|s_k\|^3,
\end{aligned}$$

which together with (12) implies that $\rho_k \geq \eta$, as claimed. \square

We now prove that the sequence $\{\sigma_k^U\}$ is bounded above.

Lemma 3.10. *There exists a scalar constant $\sigma_{max} \in \mathbb{R}_{++}$ such that, for all $k \in \mathbb{N}_+$,*

$$\sigma_k^U \leq \sigma_{max}.$$

Proof. Consider any $k \in \mathbb{N}_+$. If s_k is accepted (i.e., $k \in \mathcal{A}$), then $\sigma_{k+1}^U \leftarrow \sigma_k^U$. On the other hand, if s_k is rejected (i.e., $k \in \mathcal{R}$), then it follows from Step 12 and Step 14 that $\sigma_{k+1}^U \leq \max\{\bar{\sigma}, \gamma_2 \lambda_k / \|s_k\|\}$. Moreover, since $k \in \mathcal{R}$, meaning that $\rho_k < \eta$, it follows from Lemma 3.9 that $\lambda_k / \|s_k\|$ is bounded above by $(H_{Lip} + \kappa_2 + 2\eta)$. Thus, it follows that $\sigma_{k+1}^U \leq \max\{\bar{\sigma}, \gamma_2 (H_{Lip} + \kappa_2 + 2\eta)\}$ for all $k \in \mathcal{R}$. Overall, the desired result follows for any $\sigma_{max} \geq \max\{\bar{\sigma}, \gamma_2 (H_{Lip} + \kappa_2 + 2\eta)\}$. \square

We now establish a lower bound on the norm of any accepted trial step.

Lemma 3.11. *For all $k \in \mathcal{A}$, it follows that*

$$\|s_k\| \geq \left(\frac{1}{2} H_{Lip} + 2\sigma_{max} + \kappa_3 \right)^{-1/2} \|g_{k+1}\|^{1/2}.$$

Proof. Let $k \in \mathcal{A}$. It follows that

$$\begin{aligned}
\|g_{k+1}\| &\leq \|g_{k+1} - (g_k + (H_k + \lambda_k I)s_k)\| + \|g_k + (H_k + \lambda_k I)s_k\| \\
&\leq \|g_{k+1} - (g_k + H_k s_k)\| + \lambda_k \|s_k\| + \|g_k + (H_k + \lambda_k I)s_k\|.
\end{aligned} \tag{15}$$

By Taylor's theorem and Assumption 3.4, the first term on the right-hand side of this inequality satisfies

$$\begin{aligned}
\|g_{k+1} - (g_k + H_k s_k)\| &\leq \left\| \int_0^1 (H(x_k + \tau s_k) - H_k) s_k d\tau \right\| \\
&\leq \int_0^1 \|H(x_k + \tau s_k) - H_k\| d\tau \cdot \|s_k\| \\
&\leq \int_0^1 \tau d\tau \cdot H_{Lip} \|s_k\|^2 = \frac{1}{2} H_{Lip} \|s_k\|^2.
\end{aligned}$$

Combining this with (15) and observing Step 5, (4c), and Lemma 3.10, it follows that

$$\begin{aligned}
\|g_{k+1}\| &\leq \frac{1}{2} H_{Lip} \|s_k\|^2 + 2 \frac{\lambda_k}{\|s_k\|} \|s_k\|^2 + \kappa_3 \|s_k\|^2 \\
&\leq \frac{1}{2} H_{Lip} \|s_k\|^2 + 2\sigma_{max} \|s_k\|^2 + \kappa_3 \|s_k\|^2,
\end{aligned}$$

which, after rearrangement, completes the proof. \square

We are now prepared to prove a worst-case upper bound on the total number of accepted steps that may occur for iterations in which the norm of the gradient of the objective is above a positive threshold.

Lemma 3.12. *For any $\epsilon \in \mathbb{R}_{++}$, the total number of elements in the index set*

$$\mathcal{K}_\epsilon := \{k \in \mathbb{N}_+ : k \geq 1, (k-1) \in \mathcal{A}, \|g_k\| > \epsilon\}$$

is at most

$$\left\lfloor \left(\frac{f_0 - f_{\inf}}{\eta(\frac{1}{2}H_{Lip} + 2\sigma_{max} + \kappa_3)^{-3/2}} \right) \epsilon^{-3/2} \right\rfloor =: N_{\mathcal{A}}(\epsilon) \geq 0. \quad (16)$$

Proof. The proof follows in a similar manner as that of Lemma 3.20 in [11]. By Lemma 3.11, it follows that, for all $k \in \mathcal{K}_\epsilon$, one finds

$$\begin{aligned} f_{k-1} - f_k &\geq \eta \|s_{k-1}\|^3 \\ &\geq \eta (\frac{1}{2}H_{Lip} + 2\sigma_{max} + \kappa_3)^{-3/2} \|g_k\|^{3/2} \\ &\geq \eta (\frac{1}{2}H_{Lip} + 2\sigma_{max} + \kappa_3)^{-3/2} \epsilon^{3/2}. \end{aligned}$$

In addition, it follows from Theorem 3.3 that $|\mathcal{K}_\epsilon| < \infty$. Hence, the reduction in f obtained up to the largest index in \mathcal{K}_ϵ , call it \bar{k}_ϵ , satisfies

$$f_0 - f_{\bar{k}_\epsilon} = \sum_{k=1}^{\bar{k}_\epsilon} (f_{k-1} - f_k) \geq \sum_{k \in \mathcal{K}_\epsilon} (f_{k-1} - f_k) \geq |\mathcal{K}_\epsilon| \eta (\frac{1}{2}H_{Lip} + 2\sigma_{max} + \kappa_3)^{-3/2} \epsilon^{3/2}.$$

Rearranging this inequality to yield an upper bound for $|\mathcal{K}_\epsilon|$ and using the fact that $f_0 - f_{\inf} \geq f_0 - f_{\bar{k}_\epsilon}$, one obtains the desired result. \square

In order to prove a result similar to Lemma 3.12 for the *total* number of iterations with $\|g_k\| > \epsilon$, we require an upper bound on the total number of trial steps that may be rejected between accepted steps. To this end, let us define, for a given $\hat{k} \in \mathcal{A} \cup \{0\}$, the iteration number and corresponding set

$$\begin{aligned} k_{\mathcal{A}}(\hat{k}) &:= \min\{k \in \mathcal{A} : k > \hat{k}\} \\ \text{and } \mathcal{I}(\hat{k}) &:= \{k \in \mathbb{N}_+ : \hat{k} < k < k_{\mathcal{A}}(\hat{k})\}, \end{aligned}$$

i.e., we let $k_{\mathcal{A}}(\hat{k})$ be the smallest of all iteration numbers in \mathcal{A} that is strictly larger than \hat{k} , and we let $\mathcal{I}(\hat{k})$ be the set of iteration numbers between \hat{k} and $k_{\mathcal{A}}(\hat{k})$.

We now show that the number of rejected steps between the first iteration and the first accepted step, or between consecutive accepted steps, is bounded above.

Lemma 3.13. *For any $\hat{k} \in \mathcal{A} \cup \{0\}$, it follows that*

$$|\mathcal{I}(\hat{k})| \leq 1 + \left\lfloor \frac{1}{\log(\gamma_1)} \log \left(\frac{\sigma_{max}}{\underline{\sigma}} \right) \right\rfloor =: N_{\mathcal{R}} \geq 0.$$

Proof. The proof follows in a similar manner as for Lemma 3.24 in [11]. First, the result holds trivially if $|\mathcal{I}(\hat{k})| = 0$. Thus, we may assume that $|\mathcal{I}(\hat{k})| \geq 1$. Since $(\hat{k}+1) \in \mathcal{R}$ by construction, it follows from Steps 11–14 and Step 5 that $\lambda_{\hat{k}+2}/\|s_{\hat{k}+2}\| \geq \underline{\sigma}$, which, due to the lower bound on $\lambda_{k+1}/\|s_{k+1}\|$ in Step 14 and Step 5, leads to

$$\lambda_{k_{\mathcal{A}}(\hat{k})} \geq \underline{\sigma} (\gamma_1)^{k_{\mathcal{A}}(\hat{k})-\hat{k}-2} \|s_{k_{\mathcal{A}}(\hat{k})}\|.$$

Combining this with Step 5 and Lemma 3.10 shows that

$$\sigma_{max} \geq \sigma_{k_{\mathcal{A}}(\hat{k})}^{\mathbb{U}} \geq \lambda_{k_{\mathcal{A}}(\hat{k})}/\|s_{k_{\mathcal{A}}(\hat{k})}\| \geq \underline{\sigma} (\gamma_1)^{k_{\mathcal{A}}(\hat{k})-\hat{k}-2}.$$

After rearrangement, it now follows that

$$k_{\mathcal{A}}(\hat{k}) - \hat{k} - 2 \leq \frac{1}{\log(\gamma_1)} \log \left(\frac{\sigma_{max}}{\underline{\sigma}} \right).$$

The desired result follows from this inequality since $|\mathcal{I}(\hat{k})| = k_{\mathcal{A}}(\hat{k}) - \hat{k} - 1$. \square

We are now prepared to prove our main complexity result of this subsection.

Theorem 3.5. *Under Assumptions 2.1, 3.1, 3.2, and 3.4, for a scalar $\epsilon \in \mathbb{R}_{++}$, the total number of elements in the index set $\{k \in \mathbb{N}_+ : \|g_k\| > \epsilon\}$ is at most*

$$N(\epsilon) := 1 + N_{\mathcal{R}} N_{\mathcal{A}}(\epsilon), \quad (17)$$

where $N_{\mathcal{A}}(\epsilon)$ and $N_{\mathcal{R}}$ are defined in Lemmas 3.12 and 3.13, respectively. Consequently, for any $\bar{\epsilon} \in \mathbb{R}_{++}$, it follows that $N(\epsilon) = \mathcal{O}(\epsilon^{-3/2})$ for all $\epsilon \in (0, \bar{\epsilon}]$.

Proof. Without loss of generality, we may assume that at least one iteration is performed. Lemma 3.12 guarantees that the total number of elements in the index set $\{k \in \mathcal{A} : k \geq 1, \|g_k\| > \epsilon\}$ is at most $N_{\mathcal{A}}(\epsilon)$, where, immediately prior to each of the corresponding accepted steps, Lemma 3.13 guarantees that at most $N_{\mathcal{R}}$ trial steps are rejected. Accounting for the first iteration, the desired result follows. \square

3.3 Second-Order Global Convergence and Complexity

Our goal in this subsection is to prove results showing that, in some sense, the algorithm converges to second-order stationarity and does so with a worst-case iteration complexity on par with the methods in [7] and [11]. In particular, our results show that if the algorithm computes each search direction to satisfy a curvature condition over a subspace, then second-order stationarity is reached in a manner that depends on the subspaces.

In this subsection, we make the following additional assumption about the subproblem solver.

Assumption 3.6. *For all $k \in \mathbb{N}_+$, let $\mathcal{L}_k \subseteq \mathbb{R}^n$ denote a subspace with an orthonormal basis formed from the columns of a matrix R_k . The step s_k satisfies*

$$\xi(R_k^T H_k R_k) \geq -\kappa_4 \|s_k\| \quad (18)$$

for some $\kappa_4 \in \mathbb{R}_+$, where $\xi(R_k^T H_k R_k)$ indicates the smallest eigenvalue of $R_k^T H_k R_k$.

This assumption is reasonable, e.g., in cases when s_k is computed by solving problem 2 with the component s restricted to a subspace of \mathbb{R}^n . We refer the reader to Theorem B.3 for a proof of this fact, which also reveals that this assumption is congruous with Assumption 2.1.

Under this assumption, we have the following second-order convergence result.

Theorem 3.7. *Suppose Assumptions 2.1, 3.1, 3.2, 3.4, and 3.6 hold. It follows that*

$$\liminf_{k \in \mathcal{A}, k \rightarrow \infty} \xi(R_k^T H_k R_k) \geq 0.$$

Proof. The result follows from (18) since $\{s_k\}_{k \in \mathcal{A}} \rightarrow 0$ (see Lemma 3.7). \square

As a consequence of Theorem 3.7, if the sequence $\{R_k\}_{k \in \mathcal{A}}$ tends toward full-dimensionality as $k \rightarrow \infty$, then any limit point x_* of $\{x_k\}$ must have $H(x_*) \succeq 0$.

Our next goal is to prove a worst-case iteration complexity result for achieving second-order stationarity in a sense similar to that in Theorem 3.7. Toward this end, we first prove the following lemma, which is similar to Lemma 3.12.

Lemma 3.14. *For any $\epsilon \in \mathbb{R}_{++}$, the total number of elements in the index set*

$$\mathcal{K}_{\epsilon,\xi} := \{k \in \mathbb{N}_+ : k \geq 1, (k-1) \in \mathcal{A}, \xi(R_k^T H_k R_k) < -\epsilon\}$$

is at most

$$\left\lfloor \left(\frac{f_0 - f_{\inf}}{\eta \kappa_4^{-3}} \right) \epsilon^{-3} \right\rfloor =: N_{\mathcal{A},\xi}(\epsilon) \geq 0. \quad (19)$$

Proof. Under Assumption 3.6, it follows that, for all $k \in \mathcal{K}_{\epsilon,\xi}$, one finds

$$f_{k-1} - f_k \geq \eta \|s_{k-1}\|^3 \geq \eta \left(\frac{-\xi(R_k^T H_k R_k)}{\kappa_4} \right)^3 \geq \eta \kappa_4^{-3} \epsilon^3.$$

It follows from this inequality, the fact that f is monotonically decreasing over the sequence of iterates, and Assumption 3.1 that

$$f_0 - f_{\inf} \geq \sum_{k \in \mathcal{K}_{\epsilon,\xi}} (f_{k-1} - f_k) \geq |\mathcal{K}_{\epsilon,\xi}| \eta \kappa_4^{-3} \epsilon^3.$$

Rearranging this inequality to yield an upper bound for $|\mathcal{K}_{\epsilon,\xi}|$ gives the result. \square

We close with the following second-order complexity result.

Theorem 3.8. *Under Assumptions 2.1, 3.1, 3.2, 3.4, and 3.6, for any pair of scalars $(\epsilon_1, \epsilon_2) \in \mathbb{R}_{++} \times \mathbb{R}_{++}$, the number of elements in the index set*

$$\{k \in \mathbb{N}_+ : \|g_k\| > \epsilon_1 \vee \xi(R_k^T H_k R_k) < -\epsilon_2\}$$

is at most

$$N(\epsilon_1, \epsilon_2) := 1 + N_{\mathcal{R}} \max\{N_{\mathcal{A}}(\epsilon_1), N_{\mathcal{A},\xi}(\epsilon_2)\}, \quad (20)$$

where $N_{\mathcal{A}}(\cdot)$, $N_{\mathcal{R}}$, and $N_{\mathcal{A},\xi}(\cdot)$ are defined in Lemmas 3.12, 3.13, and 3.14, respectively. Consequently, for any pair of scalars $(\bar{\epsilon}_1, \bar{\epsilon}_2) \in \mathbb{R}_{++} \times \mathbb{R}_{++}$, it follows that

$$N(\epsilon_1, \epsilon_2) = \mathcal{O}(\max\{\epsilon_1^{-3/2}, \epsilon_2^{-3}\}) \text{ for all } (\epsilon_1, \epsilon_2) \in (0, \bar{\epsilon}_1] \times (0, \bar{\epsilon}_2].$$

Proof. The proof follows in a similar manner as that of Theorem 3.5 by additionally incorporating the bound proved in Lemma 3.14. \square

4 Algorithm Instances

Algorithm 1 is a broad framework containing, amongst other algorithms, ARC and TRACE. Indeed, the proposed framework and its supporting analyses cover a wide range of algorithms as long as the pairs in the sequence $\{(s_k, \lambda_k)\}$ satisfy Assumption 2.1.

In this section, we show that ARC and TRACE are special cases of our proposed framework in that the steps these algorithms accept would also be acceptable for our framework, and that the procedures followed by these methods after a step is rejected are consistent with our framework. We then introduce an instance of our framework that is new to the literature. (If desired for the guarantees in §3.3, one could also mind whether the elements in the sequence $\{(s_k, \lambda_k)\}$ satisfy Assumption 3.6. However, for brevity in this section, let us suppose that one is interested only in Assumption 2.1.)

4.1 ARC as a Special Case

The ARC method, which was inspired by the work in [19] and [21], was first proposed and analyzed in [6, 7]. In these papers, various sets of step computation conditions are considered involving exact and inexact subproblem solutions yielding different types of convergence and worst-case complexity guarantees. For our purposes here, we consider the more recent variant of ARC stated and analyzed as “AR p ” with $p = 2$ in [4]. (For ease of comparison, we consider this algorithm when their regularization parameter update—see Step 4 in their algorithm—uses $\eta_1 = \eta_2$. Our algorithm is easily extended to employ a two-tier acceptance condition, involving two thresholds η_1 and η_2 , as is used in [4] and [6, 7].)

Suppose that a trial step s_k is computed by this version of ARC. In particular, let us make the reasonable assumption that the subproblem for which s_k is an approximate solution is defined by some regularization value $\sigma_k \in [\sigma_k^L, \sigma_k^U]$ (with $\sigma_k^L \geq \sigma_{min}$ since ARC ensures that $\sigma_k \geq \sigma_{min} \in \mathbb{R}_{++}$ for all $k \in \mathbb{N}$) and that this subproblem is minimized over a subspace \mathcal{L}_k such that $g_k \in \mathcal{L}_k$ (see Appendix B). As is shown using a similar argument as in the proof of our Theorem B.3(b), one can show under these conditions that (s_k, λ_k) with $\lambda_k = \sigma_k \|s_k\|$ satisfies (4a). In addition, considering the algorithm statement in [4], but using our notation, one is required to have

$$g_k^T s_k + \frac{1}{2} s_k^T H_k s_k + \lambda_k \|s_k\|^2 < 0 \quad \text{and} \quad \|g_k + (H_k + \lambda_k I)s_k\| \leq \theta \|s_k\|^2 \quad \text{for some } \theta \in \mathbb{R}_{++}.$$

It is easily seen that (s_k, λ_k) satisfying these conditions also satisfies (4b)–(4c) for any $(\kappa_1, \kappa_2, \kappa_3)$ such that $\kappa_1 \geq \frac{1}{2} H_{max}$ and $\kappa_3 \geq \theta$. Overall, we have shown that a trial step s_k computed by this version of ARC satisfies Assumption 2.1, meaning that it satisfies the condition in Step 5 in Algorithm 1. If this trial step is accepted by ARC, then this means that $f_k - f(x_k + s_k) \geq \eta_1(f_k - q_k(s_k))$. Along with [4, Lemma 2.1], this implies that $f_k - f(x_k + s_k) \geq \frac{1}{3} \eta \sigma_k \|s_k\|^3$, meaning that $\rho_k \geq \frac{1}{3} \eta_1 \sigma_{min}$. Hence, this trial step would also be accepted in Algorithm 1 under the assumption that $\eta \in (0, \frac{1}{3} \eta_1 \sigma_{min}]$.

Finally, if a trial step is rejected in this version of ARC, then σ_{k+1} is set to a positive multiple of σ_k . This is consistent with the procedure after a step rejection in Algorithm 1, where it is clear that, with appropriate parameter choices, one would find $\sigma_{k+1} \in [\sigma_{k+1}^L, \sigma_{k+1}^U]$.

4.2 TRACE as a Special Case

TRACE is proposed and analyzed in [11]. Our goal in this subsection is to show that, with certain parameter settings, a trial step that is computed and accepted by TRACE could also be one that is computed and accepted by Algorithm 1, and that the procedures for rejecting a step in TRACE are consistent with those in Algorithm 1. Amongst other procedures, TRACE involves dynamic updates for two sequences, $\{\delta_k\}$ and $\{\Delta_k\}$. The elements of $\{\delta_k\}$ are the trust region radii while $\{\Delta_k\}$ is a monotonically nondecreasing sequence of upper bounds for the trust region radii; consequently, $\|s_k\| \leq \delta_k \leq \Delta_k$ with $\Delta_{k+1} \geq \Delta_k$ for all $k \in \mathbb{N}$. For simplicity in our discussion here, let us assume that $\|s_k\| < \Delta_k$ for all $k \in \mathbb{N}$. This is a fair assumption since, as shown in [11, Lemma 3.11], the manner in which $\{\Delta_k\}$ is set ensures that $\|s_k\| = \Delta_k$ only a finite number of times in any run.

In TRACE, during iteration $k \in \mathbb{N}$, a trust region radius $\delta_k \in \mathbb{R}_{++}$ is given and a trial step s_k and regularization value λ_k are computed satisfying the standard trust region subproblem optimality conditions

$$g_k + (H_k + \lambda_k I)s_k = 0, \quad H_k + \lambda_k I \succeq 0, \quad \text{and} \quad \lambda_k(\delta_k - \|s_k\|) = 0, \quad \text{where } (\lambda_k, \delta_k - \|s_k\|) \geq 0.$$

By the first of these conditions, the pair (s_k, λ_k) clearly satisfies (4b)–(4c). In addition, one can use standard trust region theory, in particular related to Cauchy decrease (see [10] or [23]), to show that the pair also satisfies (4a). Overall, assuming that the pair (σ_k^L, σ_k^U) is set such that $\lambda_k / \|s_k\| \in [\sigma_k^L, \sigma_k^U]$, it follows that Assumption 2.1 is satisfied, meaning that TRACE offers the condition in Step 5 in Algorithm 1. If the trial step s_k is subsequently accepted by TRACE, then it would also be accepted by Algorithm 1 since both algorithms use the same step acceptance condition.

Now suppose that a trial step is not accepted in TRACE. This can occur in two circumstances. It can occur if $\rho_k \geq \eta$ while $\lambda_k > \sigma_k \|s_k\|$, in which case the trust region radius is *expanded* and a new subproblem

is solved. By the proof of [11, Lemma 3.7], the solution of this new subproblem yields (in iteration $k + 1$ in TRACE) the relationship that $\lambda_{k+1}/\|s_{k+1}\| \leq \sigma_{k+1} = \sigma_k$. Hence, under the same assumption as above that the pair (σ_k^L, σ_k^U) is set such that $\lambda_k/\|s_k\| \in [\sigma_k^L, \sigma_k^U]$, this shows that the procedure in TRACE involving *an expansion of the trust region radius and the computation of the subsequent trial step* yields a trial step that would be offered in a *single iteration* in Algorithm 1. The other circumstance in which a trial step is rejected in TRACE is when $\rho_k < \eta$, in which case the trust region radius is contracted. In this case, one can see that the outcome of the CONTRACT subroutine in TRACE is consistent with Steps 11–14 of Algorithm 1 in the sense that the solution of the subsequent subproblem in TRACE will have $\lambda_{k+1}/\|s_{k+1}\| \in [\underline{\sigma}, \bar{\sigma}]$ (if $\lambda_k < \underline{\sigma}\|s_k\|$) or $\lambda_{k+1}/\|s_{k+1}\|$ within a range defined by positive multiples of $\lambda_k/\|s_k\|$; see Lemmas 3.17 and 3.23 in [11].

4.3 A Hybrid Algorithm

The primary distinguishing feature of our algorithm instance is the manner in which we compute the pair (s_k, λ_k) in Step 5 of Algorithm 1. Our newly proposed hybrid algorithm considers two cases.

Case 1: $\sigma_k^L > 0$. In this case, we find a pair (s_k, λ_k) by solving problem (A.3) over a sequence of increasingly higher dimensional Krylov subspaces as described in [6] until (4) and (18) are satisfied. The reason we know that (4) and (18) will eventually be satisfied can be seen as follows. Solving problem (A.3) over a Krylov subspace is equivalent to solving problem (B.3) with an appropriate choice of R_k as a basis for that Krylov subspace, then setting $s_k = R_k v_k$. Then, it follows from Theorem B.2(i) that solving (B.3) is equivalent to solving (B.2), which in turn is equivalent to solving (B.1) in the sense that if $(v_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ is a first-order primal-dual solution of problem (B.2), then $(s_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ with $s_k = R_k v_k$ is a solution of problem (B.1). Finally, we need only note from Theorem B.3 that solutions to problem (B.1) satisfy (4a) for all Krylov subspaces \mathcal{L}_k (recall that g_k is contained in all Krylov subspaces), (4b) for all Krylov subspaces, (4c) if the Krylov subspace \mathcal{L}_k includes enough of the space (in the worst case, $\mathcal{L}_k = \mathbb{R}^n$), and (18) for all Krylov subspaces.

Case 2: $\sigma_k^L = 0$. In this case, we begin by applying the linear CG method in an attempt to solve the linear system $H_k s = -g_k$, which iteratively solves

$$\min_{s \in \mathbb{R}^n} q_k(s) \quad (21)$$

over a sequence of expanding Krylov subspaces. One of two outcomes is possible. First, the CG algorithm may ultimately identify a vector s_k such that (s_k, λ_k) with $\lambda_k = 0$ satisfies (4) and (18). Second, the CG algorithm may never identify a vector s_k such that (s_k, λ_k) with $\lambda_k = 0$ satisfies (4) and (18). Indeed, this might occur if CG encounters a direction of negative curvature—in which case we terminate CG immediately—or if CG solves (21) accurately or reaches an iteration limit, and yet at least one condition in (4)/(18) is not satisfied. In such a case, we choose to reset $\sigma_k^L \in (0, \sigma_k^U]$, then solve problem (A.3) over a sequence of expanding Krylov subspaces as described in Case 1. In this manner, we are guaranteed to identify a pair (s_k, λ_k) satisfying (4) and (18) as required.

5 Implementation and Numerical Results

We implemented two algorithms in MATLAB, one following the strategy in §4.3 and, for comparison purposes, one following the ARC algorithm in [7] with ideas from [4]. We refer to our implementation of the former as **iR_Newton**, for inexact Regularized Newton, and to our implementation of the latter as **iARC**, for inexact ARC. In this section, we describe our approach for computing the pairs $\{(s_k, \lambda_k)\}$ in **iR_Newton** and **iARC**, as well as other implementation details, and discuss the results of numerical experiments on a standard set of nonlinear optimization test problems.

5.1 Implementation Details

Let us begin by noting that the implemented algorithms terminate in iteration $k \in \mathbb{N}_+$ if

$$\|g_k\|_\infty \leq 10^{-6} \max\{\|g_0\|_\infty, 1\}.$$

We chose not to employ a termination test based on a second-order stationarity condition. Correspondingly, neither of the algorithms check a second-order condition when computing a trial step; e.g., in **iR_Newton**, we are satisfied with a step satisfying (4) and do not check (18). In addition, for practical purposes, we set a maximum iteration limit of 10^6 , a time limit of four hours, and a minimum step norm limit of 10^{-20} . For reference, the input parameter values we used are given in Table 1. We chose these values as ones that worked well on our test set for both implemented algorithms.

Table 1: Input parameters for **iARC** and **iR_Newton**

η_1	$1.0e-16$	γ_0	$2.0e-01$	κ_1	$1.0e+00$	$\underline{\sigma}$	$1.0e-10$
η_2	$1.0e-01$	γ_1	$1.0e+01$	κ_2	$1.0e+00$	$\bar{\sigma}$	$1.0e+20$
		γ_2	$2.0e+02$	κ_3	$1.0e+00$		

For both implemented algorithms, we employ a sequence $\{\sigma_k\}$ that is updated dynamically. In **iARC**, this sequence is handled as described in [7], namely,

$$\sigma_{k+1} \leftarrow \begin{cases} \max\{\underline{\sigma}, \gamma_0 \sigma_k\} & \text{if } \frac{f_k - f(x_k + s_k)}{f_k - c_k(s_k; \sigma_k)} \geq \eta_2 \\ \sigma_k & \text{if } \frac{f_k - f(x_k + s_k)}{f_k - c_k(s_k; \sigma_k)} \in [\eta_1, \eta_2) \\ \gamma_1 \sigma_k & \text{if } \frac{f_k - f(x_k + s_k)}{f_k - c_k(s_k; \sigma_k)} < \eta_1 \end{cases}$$

The value σ_k is used in defining $c_k(\cdot; \sigma_k)$ (recall (5)) that is minimized approximately to compute the trial step s_k for all $k \in \mathbb{N}_+$. In particular, the implementation iteratively constructs Krylov subspaces of increasing dimension using the Lanczos process, where for each subspace we employ the RQS function from the GALAHAD software library (see [17] and [18]) to minimize $c_k(\cdot; \sigma_k)$ over the subspace. If the subspace is full-dimensional or the resulting step s_k satisfies

$$\|g_k + (H_k + \sigma_k \|s_k\| I)s_k\| \leq \kappa_3 \|s_k\|^2, \quad (22)$$

then it is used as the trial step. Otherwise, the process continues with a larger subspace. We remark that condition (22) is more restrictive than our condition (4c), but we use it since it is one that has been proposed for cubic regularization methods; e.g., see (2.13) in [4].

One could employ more sophisticated techniques for setting the elements of the sequence $\{\sigma_k\}$ in **iARC** that attempt to reduce the number of rejected steps; e.g., see [16]. Such improvements might aid **iR_Newton** as well. However, for simplicity and to avoid the need for additional parameter tuning, we did not include such enhancements in our implemented algorithms.

As for **iR_Newton**, for consistency between the two implementations, we do not explicitly compute the sequence $\{\lambda_k\}$, but rather employ $\{\sigma_k^L \|s_k\|\}$ in its place. For example, whenever an acceptable step is computed with $\sigma_k^L = 0$, then, as described in **Case 2** in §4.3, we effectively use $\lambda_k = 0$. On the other hand, when $\sigma_k^L > 0$, we employ the same iterative approach as used for **iARC** to compute the trial step s_k as an approximate minimizer of $c_k(\cdot; \sigma_k^L)$, where in place of λ_k in (4) we employ $\sigma_k^L \|s_k\|$. Then, in either case, in the remainder of iteration $k \in \mathbb{N}_+$, specifically for setting σ_{k+1}^L and σ_{k+1}^U , we use $\sigma_k^L \|s_k\|$ in place of λ_k in Steps 11 and 14. We also define an auxiliary sequence $\{\sigma_k\}$ using the update

$$\sigma_{k+1} \leftarrow \begin{cases} \max\{\underline{\sigma}, \gamma_0 \sigma_k\} & \text{if } \rho_k \geq \eta_1 \text{ and } \sigma_k^L > 0 \\ \sigma_k & \text{if } \sigma_k^L = 0 \\ \min\{\gamma_1 \sigma_k, \bar{\sigma}\} & \text{if } \rho_k < \eta_1 \text{ and } \sigma_k^L > 0. \end{cases}$$

This update is similar to the one employed for **iARC** with the added assurance that $\{\sigma_k\} \subset [\underline{\sigma}, \bar{\sigma}]$. The elements of this sequence are used in two circumstances. First, if, as described in **Case 2** in §4.3, CG fails to produce a trial step s_k satisfying (4) (with $\lambda_k = 0$), then we reset $\sigma_k^L \leftarrow \sigma_k$ and revert to the same scheme as above to compute the trial step when $\sigma_k^L > 0$. Second, if a step is rejected and $\sigma_k^L < \underline{\sigma}$ (equivalently, $\lambda_k < \underline{\sigma}\|s_k\|_2$ as in Step 12 in Algorithm 1), then we set $\sigma_{k+1}^L \leftarrow \sigma_{k+1}$. Lastly, we note that if CG ever performs n iterations and the resulting solution (due to numerical error) does not satisfy (4) and no negative curvature is detected, then the resulting approximate solution s_k is used as the trial step.

5.2 Results on the CUTEst Test Set

We employed our implemented algorithms, **iARC** and **iR_Newton**, to solve unconstrained problems in the CUTEst test set; see [15]. Among 171 unconstrained problems in the set, one (**FLETCBV2**) was removed since the algorithms terminated at the initial point, five (**ARGLINC**, **DECONVU**, **FLETCHBV**, **INDEFM**, and **POWER**) were removed due to a function evaluation error or our memory limitation of 8GB, and nine (**EIGENBLS**, **EIGENCLS**, **FMINSURF**, **NONMSQRT**, **SBRYBND**, **SCURLY10**, **SCURLY20**, **SCURLY30**, and **SSCOSINE**) were removed since neither algorithm terminated within our time limit. In addition, four were removed since neither of the algorithms terminated successfully: for **HIELOW**, **iARC** reached our maximum iteration limit; for **CURLY20** and **SCOSINE**, **iARC** reached the time limit; for **INDEF**, **iARC** terminated due to a subproblem solver error; and for all of these four problems, **iR_Newton** terminated due to our minimum step norm limit. The remaining set consisted of 152 test problems with number of variables ranging from 2 to 100,000. For additional details on the problems used and their sizes, see Appendix C.

To compare the performance of the implemented algorithms, we generated performance profiles for the number of iterations and number of Hessian-vector products required before termination. These are shown in Figure 1. A performance profile graph of an algorithm at point α shows the fraction of the test set for which the algorithm is able to solve within a factor of 2^α of the best algorithm for the given measure; see [12]. When generating the profiles, we did not include three of the test problems—**CURLY10**, **CURLY30**, and **MODBEALE**—on which **iARC** was unsuccessful while **iR_Newton** was successful. (In particular, **iARC** reached the time limit for all problems.) We feel that this gives a fairer comparison with respect to the problems on which both algorithms were successful.

As seen in Figure 1, the algorithms performed relatively comparably when it came to the number of iterations required, though clearly **iR_Newton** had an edge in terms of requiring fewer iterations on various problems. The difference in terms of numbers of Hessian-vector products required was more drastic, and indeed we point to this as the main measure of improved performance for **iR_Newton** versus **iARC**. One reason for this discrepancy is that **iR_Newton** required fewer iterations on some problems. However, more significantly, the difference was due in part to **iR_Newton**'s ability to employ and accept inexact Newton steps (with $\lambda_k = 0$) on many iterations. This is due to the fact that, in CG, one is able to compute the Hessian-vector product $H_k s_k$, needed to check the termination conditions for the computation of s_k , by taking a linear combination of Hessian-vector products already computed in CG; i.e., if $\{p_{k,i}\}$ are the search directions computed in CG such that $s_k = \sum_i \alpha_{k,i} p_{k,i}$, then CG involves computing $H_k p_{k,i}$ for each i and can compute $H_k s_k = \sum_i \alpha_{k,i} (H_k p_{k,i})$. By contrast, one is unable to retrieve this product via a linear combination when the step is computed from the minimization of a cubic function, as is needed in **iARC** and in **iR_Newton** whenever $\sigma_k^L > 0$. Overall, we claim that the primary strength of **iR_Newton** as compared to **iARC** is its ability to employ inexact Newton steps.

For further details of our numerical results, see Appendix C. In these results, we also indicate the number of tridiagonal factorizations required; at least one is needed involving a tridiagonal matrix of size $m \times m$ every time an algorithm solves a cubic subproblem over an m -dimensional subspace.

6 Conclusion

We have proposed a general framework for solving smooth nonconvex optimization problems and proceeded to prove worst-case iteration complexity bounds for it. In fact, for a certain class of second-order methods em-

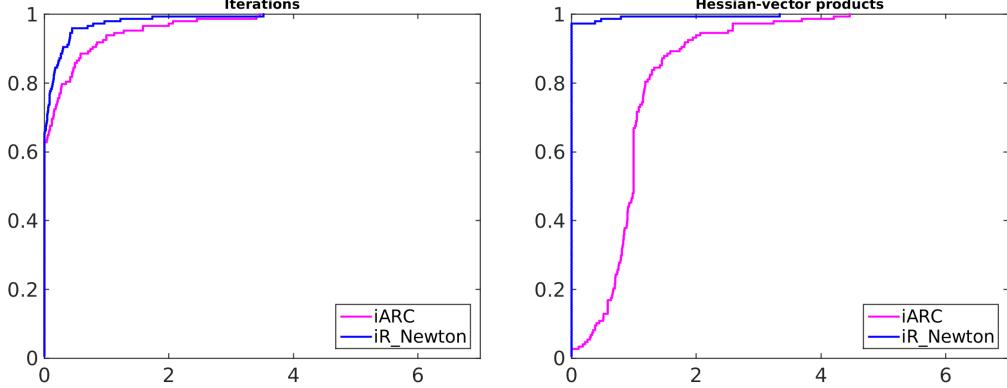


Figure 1: Performance profiles for **iARC** and **iR_Newton**.

ployed to minimize a certain class of nonconvex functions, our first-order complexity result for our method is known to be optimal; see [8]. Our framework is flexible enough to cover a wide range of popular algorithms, an achievement made possible by the use of generic conditions that each trial step is required to satisfy. The use of such conditions allows for the calculation of inexact Newton steps, for example by performing minimization over expanding Krylov subspaces. Although we have presented a particular instance of our framework motivated by subproblem (2), additional instances can easily be derived by applying other optimization strategies for solving (2). Numerical experiments with an instance of our algorithm showed that it can lead to improved performance on a broad test set as compared to an implementation of a straightforward cubic regularization approach.

Acknowledgements

We thank the anonymous referees for their valuable comments, suggestions, and corrections which helped to improve the paper. We are also grateful to the Associate Editor for handling the paper.

A Subproblem Solution Properties

In this appendix, we explore properties of any first-order stationary solution of problem $\mathcal{P}_k(\sigma_k^L, \sigma_k^U)$ defined as (2). Let us define a Lagrangian function for (2) as

$$\begin{aligned} \mathcal{L}(s, \lambda, \beta^L, \beta^U, \beta^N) = & f_k + g_k^T s + \frac{1}{2} s^T (H_k + \lambda I) s \\ & - \frac{\beta^L}{2} (\lambda^2 - (\sigma_k^L)^2 \|s\|^2) + \frac{\beta^U}{2} (\lambda^2 - (\sigma_k^U)^2 \|s\|^2) - \beta^N \lambda, \end{aligned}$$

where $(\beta^L, \beta^U) \in \mathbb{R}_+ \times \mathbb{R}_+$ are the dual variables associated with the left-hand and right-hand constraints on λ , respectively, and $\beta^N \in \mathbb{R}_+$ is the dual variable associated with the nonnegativity constraint on λ . The tuple $(s_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ is a first-order primal-dual stationary solution of $\mathcal{P}_k(\sigma_k^L, \sigma_k^U)$ if it satisfies the following conditions:

$$g_k + (H_k + \lambda_k I)s_k + \beta_k^L(\sigma_k^L)^2 s_k - \beta_k^U(\sigma_k^U)^2 s_k = 0, \quad (\text{A.1a})$$

$$\frac{1}{2}\|s_k\|^2 - \lambda_k(\beta_k^L - \beta_k^U) - \beta_k^N = 0, \quad (\text{A.1b})$$

$$0 \leq \beta_k^L \perp (\lambda_k^2 - (\sigma_k^L)^2 \|s_k\|^2) \geq 0, \quad (\text{A.1c})$$

$$0 \leq \beta_k^U \perp (\lambda_k^2 - (\sigma_k^U)^2 \|s_k\|^2) \leq 0, \quad \text{and} \quad (\text{A.1d})$$

$$0 \leq \beta_k^N \perp \lambda_k \geq 0. \quad (\text{A.1e})$$

We make the following assumption throughout this appendix.

Assumption A.1. *The vector g_k is nonzero.*

Under this assumption, the following lemma is a simple consequence of (A.1a).

Lemma A.1. *Any solution of (2) has $s_k \neq 0$.*

We now establish conditions that must hold depending on the value of $\sigma_k^L \in \mathbb{R}_+$.

Lemma A.2. *The following hold true for any solution of (A.1).*

(i) *If $\sigma_k^L > 0$, then $\lambda_k > 0$, $\beta_k^N = 0$, $\beta_k^L > 0$, and $\lambda_k = \sigma_k^L \|s_k\|$.*

(ii) *If $\sigma_k^L = 0$, then $\lambda_k = 0$.*

Proof. Consider part (i). For the sake of deriving a contradiction, suppose $\sigma_k^L > 0$ and $\lambda_k = 0$. These, along with Lemma A.1, imply that $0 = \lambda_k^2 < (\sigma_k^L)^2 \|s_k\|^2$, which contradicts (A.1c). Hence, $\lambda_k > 0$, as claimed. Then, it follows from (A.1e) that $\beta_k^N = 0$, as claimed. Next, observe that from (A.1b), Lemma A.1, $\beta_k^N = 0$, $\lambda_k > 0$, and $(\beta_k^L, \beta_k^U) \geq 0$, it follows that $\beta_k^L > 0$, as claimed. This, along with (A.1c), implies that $\lambda_k^2 = (\sigma_k^L)^2 \|s_k\|^2$. This implies that $\lambda_k = \pm(\sigma_k^L) \|s_k\|$, which combined with $\lambda_k \in \mathbb{R}^+$ means that $\lambda_k = \sigma_k^L \|s_k\|$, as claimed.

Now consider part (ii). For the sake of deriving a contradiction, suppose that $\sigma_k^L = 0$ and $\lambda_k > 0$. Then, it follows from (A.1e) that $\beta_k^N = 0$. Moreover, combining $\sigma_k^L = 0$ and $\lambda_k > 0$, it follows from (A.1c) that $\beta_k^L = 0$. It now follows from $\beta_k^L = 0$, $\beta_k^N = 0$, and (A.1b) that

$$\frac{1}{2} \|s_k\|^2 = -\lambda_k \beta_k^U \leq 0, \quad (\text{A.2})$$

where the inequality follows from $\lambda_k > 0$ and $\beta_k^U \geq 0$. This contradicts Lemma A.1. \square

Our main result is the following. In part (i) with $\sigma_k^L > 0$, we show that solving (2) is equivalent to solving what may be referred to as an ARC subproblem [6]. In part (ii) with $\sigma_k^L = 0$, we show that it is equivalent to minimizing a quadratic, if a minimizer exists.

Theorem A.2. *The following hold true.*

(i) *Suppose $\sigma_k^L > 0$. Then, (2) has a solution (s_k, λ_k) , which can be obtained as*

$$s_k \in \arg \min_{s \in \mathbb{R}^n} (f_k + g_k^T s + \frac{1}{2} s^T H_k s + \frac{1}{2} \sigma_k^L \|s\|^3), \quad (\text{A.3})$$

then setting $\lambda_k = \sigma_k^L \|s_k\| > 0$.

(ii) *If $\sigma_k^L = 0$, then a solution of problem (2) exists if and only if $H_k \succeq 0$ and $g_k^T u = 0$ for all $u \in \text{Null}(H_k)$.*

In such cases, computing a solution (s_k, λ_k) of problem (2) is equivalent to computing a solution s_k of problem (21) and setting $\lambda_k = 0$.

Proof. Consider part (i). Since $\sigma_k^L > 0$, it follows from Lemma A.2 that problem (2) is equivalent to

$$\begin{aligned} \min_{(s, \lambda) \in \mathbb{R}^n \times \mathbb{R}_+} & f_k + g_k^T s + \frac{1}{2} s^T (H_k + \lambda I) s \\ \text{s.t. } & \sigma_k^L \|s\| = \lambda, \end{aligned} \quad (\text{A.4})$$

where, by Lemma A.1, it follows that the solution has $\lambda_k > 0$, as desired. Substituting the constraint of (A.4) into the objective of (A.4), one finds that solving it is equivalent to solving (A.3) for s_k , then setting $\lambda_k = \sigma_k^L \|s_k\|$, as claimed. Since $\sigma_k^L > 0$, a minimizer of problem (A.3) exists because it involves the minimization of a coercive function.

Now consider part (ii). Since $\sigma_k^L = 0$, it follows from Lemma A.2 that $\lambda_k = 0$, meaning that problem (2) is equivalent to (21). This problem has a solution if and only if the objective is bounded below, which is the case if and only if $H_k \succeq 0$ and $g_k^T u = 0$ for all $u \in \text{Null}(H_k)$. \square

B Subproblem Solution Properties Over Subspaces

In this appendix, we explore properties of any first-order stationary solution (when one exists) of problem $\mathcal{P}_k(\sigma_k^L, \sigma_k^U)$ defined as (2) when the search space for s is restricted to a subspace of \mathbb{R}^n . Specifically, for some m -dimensional subspace $\mathcal{L}_k \subseteq \mathbb{R}^n$, consider the problem

$$\begin{aligned} & \min_{(s, \lambda) \in \mathcal{L}_k \times \mathbb{R}_+} f_k + g_k^T s + \frac{1}{2} s^T (H_k + \lambda I) s \\ & \text{s.t. } (\sigma_k^L)^2 \|s\|^2 \leq \lambda^2 \leq (\sigma_k^U)^2 \|s\|^2. \end{aligned} \quad (\text{B.1})$$

Given an orthogonal basis R_k for \mathcal{L}_k , a solution of (B.1) can be obtained from that of

$$\begin{aligned} & \min_{(v, \lambda) \in \mathbb{R}^m \times \mathbb{R}_+} f_k + g_k^T R_k v + \frac{1}{2} (R_k v)^T (H_k + \lambda I) R_k v \\ & \text{s.t. } (\sigma_k^L)^2 \|v\|^2 \leq \lambda^2 \leq (\sigma_k^U)^2 \|v\|^2. \end{aligned} \quad (\text{B.2})$$

Specifically, if $(v_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ is a first-order primal-dual solution of problem (B.2), then the tuple $(s_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ with $s_k = R_k v_k$ is such a solution of problem (B.1).

In Appendix A, we proved properties of a solution (if one exists) of a problem of the form (B.2). Let us now translate the results of that appendix to the present setting, for which we require the following assumption on the reduced gradient $R_k^T g_k$.

Assumption B.1. *The vector $R_k^T g_k$ is nonzero.*

Lemma B.1. *Any solution of (B.2) has $v_k \neq 0$.*

Lemma B.2. *The following hold for any first-order primal-dual solution of (B.1).*

- (i) *If $\sigma_k^L > 0$, then $\lambda_k > 0$, $\beta_k^N = 0$, $\beta_k^L > 0$, and $\lambda_k = \sigma_k^L \|v_k\|$.*
- (ii) *If $\sigma_k^L = 0$, then $\lambda_k = 0$.*

Theorem B.2. *The following hold true.*

- (i) *Suppose $\sigma_k^L > 0$. Then, (B.2) has a solution (v_k, λ_k) , which can be obtained as*

$$v_k \in \arg \min_{v \in \mathbb{R}^m} (f_k + g_k^T R_k v + \frac{1}{2} v^T R_k^T H_k R_k v + \frac{1}{2} \sigma_k^L \|v\|^3), \quad (\text{B.3})$$

then setting $\lambda_k = \sigma_k^L \|v_k\| > 0$.

- (ii) *If $\sigma_k^L = 0$, then a solution of (B.2) exists if and only if $R_k^T H_k R_k \succeq 0$ and $g_k^T R_k u = 0$ for all $u \in \text{Null}(R_k^T H_k R_k)$. In such cases, computing a solution (v_k, λ_k) of problem (B.2) is equivalent to computing a solution v_k of*

$$\min_{v \in \mathbb{R}^m} f_k + g_k^T R_k v + \frac{1}{2} v^T R_k^T H_k R_k v \quad (\text{B.4})$$

and setting $\lambda_k = 0$.

Considering problem (B.3), we obtain the following result from [6, Lemma 3.2].

Lemma B.3. *If $\sigma_k^L > 0$, then v_k from (B.3) satisfies*

$$g_k^T R_k v_k + v_k^T R_k^T H_k R_k v_k + \frac{3}{2} \sigma_k^L \|v_k\|^3 = 0 \quad (\text{B.5a})$$

$$v_k^T R_k^T H_k R_k v_k + \frac{3}{2} \sigma_k^L \|v_k\|^3 \geq 0 \quad (\text{B.5b})$$

$$R_k^T H_k R_k + \frac{3}{2} \sigma_k^L \|v_k\| I \succeq 0. \quad (\text{B.5c})$$

We now show that, under certain reasonable assumptions, solutions of the primal-dual reduced-space subproblem (B.1) satisfy the conditions required by Assumptions 2.1 and 3.6.

Theorem B.3. *The following hold true.*

- (a) Any solution of problem (B.1) satisfies (4b).
- (b) Any solution of problem (B.1) satisfies (4a) provided $g_k \in \mathcal{L}_k$.
- (c) Any solution of problem (B.1) satisfies (4c) provided $\mathcal{L}_k = \mathbb{R}^n$.
- (d) Any solution of problem (B.1) satisfies (18) for any $\kappa_4 \geq \frac{3}{2} \sup_{k \in \mathbb{N}_+} \{\sigma_k^L\}$.

Proof. Any first-order primal-dual solution $(s_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ of problem (B.1) corresponds to such a solution $(v_k, \lambda_k, \beta_k^L, \beta_k^U, \beta_k^N)$ of problem (B.2) where $s_k = R_k v_k$. Hence, throughout this proof, for any solution vector s_k for problem (B.1), we may let $s_k = R_k v_k$ where v_k satisfies the properties in Lemmas B.1–B.3.

First, suppose $\sigma_k^L > 0$, which by Theorem B.2(i) implies that problem (B.1) has a solution. Then, it follows from (B.5a), $s_k = R_k v_k$, and Lemma B.2(i) that

$$0 = g_k^T s_k + s_k^T H_k s_k + \frac{3}{2} \sigma_k^L \|s_k\|^3 = g_k^T s_k + s_k^T H_k s_k + \frac{3}{2} \lambda_k \|s_k\|^2,$$

which means that

$$s_k^T (g_k + (H_k + \lambda_k I) s_k) = -\frac{1}{2} \lambda_k \|s_k\|^2. \quad (\text{B.6})$$

Meanwhile, from (B.5b), $s_k = R_k v_k$, and Lemma B.2(i), it follows that

$$0 \leq s_k^T H_k s_k + \frac{3}{2} \sigma_k^L \|s_k\|^3 = s_k^T H_k s_k + \frac{3}{2} \lambda_k \|s_k\|^2 = s_k^T (H_k + \lambda_k I) s_k + \frac{1}{2} \lambda_k \|s_k\|^2,$$

which means that

$$-\frac{1}{4} \lambda_k \|s_k\|^2 \leq \frac{1}{2} s_k^T (H_k + \lambda_k I) s_k. \quad (\text{B.7})$$

It follows from (B.6), (B.7), $\lambda_k > 0$ (by Lemma B.2(i)), and $(\kappa_1, \kappa_2) \in \mathbb{R}_{++} \times \mathbb{R}_{++}$ that

$$\begin{aligned} s_k^T (g_k + (H_k + \lambda_k I) s_k) &= -\frac{1}{2} \lambda_k \|s_k\|^2 \leq \min\{\frac{1}{2} \kappa_1 \|s_k\|^2, \frac{1}{2} s_k^T (H_k + \lambda_k I) s_k - \frac{1}{4} \lambda_k \|s_k\|^2\} \\ &\leq \min\{\frac{1}{2} \kappa_1 \|s_k\|^2, \frac{1}{2} s_k^T (H_k + \lambda_k I) s_k + \frac{1}{2} \kappa_2 \|s_k\|^3\}, \end{aligned}$$

which implies (4b). This establishes that part (a) is true. Now consider part (b). From Theorem B.2, [6, Lemma 2.1], and $s_k = R_k v_k$, it follows that

$$f_k - q_k(s_k) - \frac{1}{2} \sigma_k^L \|s_k\|^3 \geq \frac{\|R_k^T g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|R_k^T g_k\|}{1 + \|R_k^T H_k R_k\|}, \frac{1}{\sqrt{6}} \sqrt{\frac{\|R_k^T g_k\|}{\sigma_k^L}} \right\}.$$

Since, under assumption, $g_k \in \mathcal{L}_k$ so that $g_k = R_k y$ for some $y \in \mathbb{R}^m$, it follows that

$$\|R_k^T g_k\| = \|R_k^T R_k y\| = \|y\| = \|R_k y\| = \|g_k\|.$$

Combining this with $\|R_k^T H_k R_k\| \leq \|H_k\|$ and the previous displayed inequality shows

$$f_k - q_k(s_k) - \frac{1}{2} \sigma_k^L \|s_k\|^3 \geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + \|H_k\|}, \frac{1}{\sqrt{6}} \sqrt{\frac{\|g_k\|}{\sigma_k^L}} \right\}.$$

This may now be combined with Theorem B.2 (specifically $\lambda_k = \sigma_k^L \|s_k\| > 0$) to obtain

$$f_k - q_k(s_k) \geq f_k - q_k(s_k) - \frac{1}{2} \sigma_k^L \|s_k\|^3 \geq \frac{\|g_k\|}{6\sqrt{2}} \min \left\{ \frac{\|g_k\|}{1 + \|H_k\|}, \frac{1}{\sqrt{6}} \sqrt{\frac{\|g_k\| \|s_k\|}{\lambda_k}} \right\},$$

which means that (s_k, λ_k) satisfies (4a), proving part (b). Now consider part (c). It follows from Theorem A.2(i) and the optimality conditions for problem (A.3) that

$$0 = g_k + H_k s_k + \frac{3}{2} \sigma_k^L \|s_k\| s_k = g_k + H_k s_k + \frac{3}{2} \lambda_k s_k = g_k + (H_k + \lambda_k I) s_k + \frac{1}{2} \lambda_k s_k.$$

This and the fact that $\kappa_3 > 0$ imply that

$$\|g_k + (H_k + \lambda_k I)s_k\| = \frac{1}{2}\lambda_k\|s_k\| \leq \lambda_k\|s_k\| + \kappa_3\|s_k\|^2,$$

which completes the proof of part (c). Finally, consider part (d). From (B.5c), the fact that $\|s_k\| = \|v_k\|$, and $\kappa_4 \geq \frac{3}{2} \sup_{k \in \mathbb{N}_+} \{\sigma_k^L\}$, it follows that

$$\xi(R_k^T H_k R_k) \geq -\frac{3}{2}\sigma_k^L\|s_k\| \geq -\kappa_4\|s_k\|,$$

as desired to prove part (d).

Now suppose that $\sigma_k^L = 0$. From Theorem B.2(ii), a solution of problem (B.1) exists if and only if $R_k^T H_k R_k \succeq 0$ and $g_k^T R_k u = 0$ for all $u \in \text{Null}(R_k^T H_k R_k)$. If this is not the case, then there is nothing left to prove; hence, let us assume that these conditions hold. From these conditions, Theorem B.2(ii), the optimality conditions of problem (B.4), the fact that $\lambda_k = 0$, and $s_k = R_k v_k$, it follows that

$$g_k^T s_k + s_k^T H_k s_k = 0 \quad \text{and} \quad s_k^T H_k s_k \geq 0.$$

This shows that (4b) holds, proving part (a) for this case. Next, since v_k is given by the solution of problem (B.4), it follows that the reduction in the objective yielded by v_k is at least as large as the reduction obtained by minimizing the objective over the span of $-R_k^T g_k$. Hence, from standard theory on Cauchy decrease (see [10] or [23]), one can conclude that

$$f_k - q_k(s_k) \geq \frac{\|R_k^T g_k\|}{2} \min \left\{ \frac{\|R_k^T g_k\|}{1 + \|R_k^T H_k R_k\|}, \|s_k\| \right\}.$$

Thus, using the arguments in the previous paragraph under the assumption that $g_k \in \mathcal{L}_k$, one is led to the conclusion that (4a) holds, which proves part (b) for this case. Next, when $\mathcal{L}_k = \mathbb{R}^n$, the optimality conditions for problem (B.4) imply that $g_k + H_k s_k = 0$, which, since $\lambda_k = 0$, implies that (4c) holds, proving part (c). Finally, since $R_k^T H_k R_k \succeq 0$, it follows that (18) holds, proving part (d). \square

C Detailed Numerical Results

Further details of the results of our numerical experiments are shown in Table 2. In the table, #Var indicates the number of variables, #Iter indicates the number of iterations required (with %Newton indicating the percentage that were inexact Newton steps with $\lambda_k = 0$), #Acc indicates the number of accepted steps (again with %Newton indicating the percentage that were inexact Newton steps), #Hv-prod indicates the number of Hessian-vector products required, and #T-fact indicates the number of tridiagonal matrix factorizations required.

Table 2: Numerical results for iARC and iR_Newton.

Prob	#Var	Alg	#Iter (%Newton)	#Acc (%Newton)	#Hv-prod	#T-fact
AKIVA	2	iARC iR_Newton	5 5 (%100)	5 5 (%100)	15 10	20 0
ALLINITU	4	iARC iR_Newton	11 8 (%50)	8 6 (%67)	56 25	61 21
ARGLINA	200	iARC iR_Newton	3 1 (%100)	3 1 (%100)	6 1	3 0
ARGLINB	200	iARC iR_Newton	2 1 (%100)	2 1 (%100)	4 1	2 0
ARWHEAD	5000	iARC iR_Newton	4 4 (%100)	4 4 (%100)	10 5	6 0
BARD	3	iARC iR_Newton	11 11 (%91)	8 10 (%90)	50 28	52 6
BDQRTIC	5000	iARC iR_Newton	9 9 (%100)	9 9 (%100)	34 17	33 0
BEALE	2	iARC iR_Newton	11 12 (%42)	8 8 (%62)	33 29	47 29

Table 2 -- continued from previous page

Prob	#Var	Alg	#Iter (%Newton)	#Acc (%Newton)	#Hv-prod	#T-fact
BIGGS6	6	iARC iR.Newton	457 419 (%96)	383 406 (%97)	3446 1557	3679 183
BOX	10000	iARC iR.Newton	3 4 (%75)	3 3 (%67)	14 9	12 4
BOX3	3	iARC iR.Newton	7 7 (%100)	7 7 (%100)	30 16	32 0
BOXPPOWER	20000	iARC iR.Newton	3 7 (%100)	3 7 (%100)	10 13	9 0
BRKMCC	2	iARC iR.Newton	2 2 (%100)	2 2 (%100)	6 4	7 0
BROWNAL	200	iARC iR.Newton	2 1 (%100)	2 1 (%100)	6 1	4 0
BROWNBS	2	iARC iR.Newton	53 5 (%80)	38 5 (%80)	142 11	191 5
BROWNDEN	4	iARC iR.Newton	8 9 (%100)	8 9 (%100)	35 20	35 0
BROYDN7D	5000	iARC iR.Newton	472 812 (%29)	279 346 (%2)	7202 5033	12598 10022
BRYBND	5000	iARC iR.Newton	19 17 (%53)	10 11 (%82)	240 206	367 81
CHAINWOO	4000	iARC iR.Newton	81 70 (%81)	57 59 (%88)	798 409	942 185
CHNROSNB	50	iARC iR.Newton	64 53 (%75)	40 40 (%82)	1126 499	1456 294
CHNRSNBM	50	iARC iR.Newton	96 101 (%58)	58 59 (%59)	1708 899	2320 960
CLIFF	2	iARC iR.Newton	14 14 (%100)	14 14 (%100)	28 14	14 0
COSINE	10000	iARC iR.Newton	12 11 (%55)	7 7 (%71)	108 45	140 29
CRAGGLVY	5000	iARC iR.Newton	30 31 (%100)	30 31 (%100)	228 108	208 0
CUBE	2	iARC iR.Newton	42 35 (%69)	27 25 (%76)	126 72	170 51
CURLY10	10000	iARC iR.Newton	---	---	271881	19957
CURLY30	10000	iARC iR.Newton	---	---	125639	630
DENSCHNA	2	iARC iR.Newton	5 5 (%100)	5 5 (%100)	15 10	15 0
DENSCHNB	2	iARC iR.Newton	7 7 (%71)	6 5 (%80)	20 12	22 8
DENSCHNC	2	iARC iR.Newton	13 11 (%82)	9 9 (%89)	38 21	46 8
DENSCHND	3	iARC iR.Newton	61 44 (%86)	57 40 (%95)	206 82	172 27
DENSCHNE	3	iARC iR.Newton	24 21 (%52)	15 16 (%69)	68 41	62 58
DENSCHNF	2	iARC iR.Newton	5 5 (%100)	5 5 (%100)	15 10	15 0
DIXMAANA	3000	iARC iR.Newton	6 6 (%100)	6 6 (%100)	14 7	8 0
DIXMAANB	3000	iARC iR.Newton	7 7 (%100)	7 7 (%100)	16 8	9 0
DIXMAANC	3000	iARC iR.Newton	8 8 (%100)	8 8 (%100)	18 9	9 0
DIXMAAND	3000	iARC iR.Newton	9 9 (%100)	9 9 (%100)	20 10	10 0
DIXMAANE	3000	iARC iR.Newton	59 60 (%100)	59 60 (%100)	670 331	622 0
DIXMAANF	3000	iARC iR.Newton	38 37 (%100)	37 37 (%100)	510 249	487 0
DIXMAANG	3000	iARC iR.Newton	39 40 (%100)	39 40 (%100)	532 288	514 0
DIXMAANH	3000	iARC iR.Newton	41 41 (%100)	41 41 (%100)	448 224	421 0
DIXMAANI	3000	iARC iR.Newton	193 249 (%100)	193 249 (%100)	3456 2814	3464 0
DIXMAANJ	3000	iARC iR.Newton	34 34 (%100)	34 34 (%100)	324 162	296 0
DIXMAANK	3000	iARC iR.Newton	30 30 (%100)	30 30 (%100)	248 124	225 0
DIXMAANL	3000	iARC	29	29	180	148

Table 2 -- continued from previous page

Prob	#Var	Alg	#Iter (%Newton)	#Acc (%Newton)	#Hv-prod	#T-fact
		iR_Newton	29 (%100)	29 (%100)	90	0
DIXMAANM	3000	iARC iR_Newton	375 398 (%100)	375 398 (%100)	10902 6126	11542 0
DIXMAANN	3000	iARC iR_Newton	82 87 (%100)	82 87 (%100)	1368	1358
DIXMAANO	3000	iARC iR_Newton	63 59 (%100)	63 59 (%100)	908 371	893 0
DIXMAANP	3000	iARC iR_Newton	51 51 (%100)	51 51 (%100)	476 238	432 0
DIXON3DQ	10000	iARC iR_Newton	2257 2476 (%100)	2256 2476 (%100)	143968 81042	164858 0
DJTL	2	iARC iR_Newton	215 204 (%32)	120 81 (%5)	642 404	866 512
DQDRITC	5000	iARC iR_Newton	6 4 (%100)	6 4 (%100)	34 10	32 0
DQRTIC	5000	iARC iR_Newton	15 11 (%100)	15 11 (%100)	30 11	15 0
EDENSCH	2000	iARC iR_Newton	15 15 (%100)	15 15 (%100)	44 22	25 0
EG2	1000	iARC iR_Newton	3 3 (%100)	3 3 (%100)	6 3	3 0
EIGENALS	2550	iARC iR_Newton	179 173 (%84)	134 150 (%89)	15548 7871	20388 1999
ENGVAL1	5000	iARC iR_Newton	9 9 (%100)	9 9 (%100)	64 32	54 0
ENGVAL2	3	iARC iR_Newton	21 21 (%57)	15 15 (%80)	100 56	139 34
ERRINROS	50	iARC iR_Newton	131 108 (%94)	122 103 (%97)	1202 504	1106 37
ERRINRSM	50	iARC iR_Newton	404 167 (%98)	396 163 (%99)	6566 1154	7225 27
EXPFIT	2	iARC iR_Newton	14 11 (%27)	9 6 (%50)	42 26	62 38
EXTROSNB	1000	iARC iR_Newton	179 185 (%62)	107 114 (%64)	2978 1576	3586 1553
FLETBV3M	5000	iARC iR_Newton	41 56 (%43)	34 32 (%41)	86 65	43 32
FLETCHCR	1000	iARC iR_Newton	2437 2187 (%66)	1450 1438 (%69)	66056 29012	90373 23819
FMINSRF2	5625	iARC iR_Newton	875 905 (%50)	567 448 (%40)	6528 2666	7378 1989
FREUROTH	5000	iARC iR_Newton	17 18 (%39)	11 10 (%60)	102 51	120 35
GENHUMPS	5000	iARC iR_Newton	14931 3567 (%2)	11710 2077 (%1)	477824 25952	1724919 85744
GENROSE	500	iARC iR_Newton	593 690 (%19)	350 341 (%4)	24494 11862	54811 30583
GROWTHLS	3	iARC iR_Newton	8 8 (%100)	8 8 (%100)	32 16	32 0
GULF	3	iARC iR_Newton	46 40 (%62)	31 29 (%62)	196 101	249 78
HAIRY	2	iARC iR_Newton	28 19 (%21)	15 10 (%40)	84 45	133 64
HATFLDD	3	iARC iR_Newton	23 23 (%65)	19 18 (%78)	108 64	140 33
HATFLDE	3	iARC iR_Newton	23 24 (%62)	18 18 (%83)	112 69	153 45
HATFLDFL	3	iARC iR_Newton	851 1127 (%84)	711 961 (%85)	4255 3429	5134 1313
HEART6LS	6	iARC iR_Newton	1502 1071 (%52)	895 620 (%51)	15002 5164	27063 6637
HEART8LS	8	iARC iR_Newton	115 186 (%31)	69 97 (%23)	1407 1178	2400 2185
HELIX	3	iARC iR_Newton	11 15 (%33)	7 9 (%44)	52 48	72 57
HILBERTA	2	iARC iR_Newton	5 3 (%100)	5 3 (%100)	12 4	9 0
HILBERTB	10	iARC iR_Newton	4 3 (%100)	4 3 (%100)	16 6	12 0
HIMMELBB	2	iARC iR_Newton	10 11 (%27)	6 6 (%33)	26 20	36 27
HIMMELBF	4	iARC iR_Newton	53 70 (%71)	36 62 (%77)	310 244	408 185

Table 2 -- continued from previous page

Prob	#Var	Alg	#Iter (%Newton)	#Acc (%Newton)	#Hv-prod	#T-fact
HIMMELBG	2	iARC iR_Newton	7 7 (%57)	6 6 (%67)	21 13	25 3
HIMMELBH	2	iARC iR_Newton	5 6 (%67)	4 4 (%75)	13 7	12 2
HUMPS	2	iARC iR_Newton	125 89 (%13)	80 49 (%10)	370 218	655 291
HYDC2OLS	99	iARC iR_Newton	11 11 (%73)	9 9 (%89)	402 215	539 165
JENSMP	2	iARC iR_Newton	8 8 (%100)	8 8 (%100)	24 16	26 0
JIMACK	3549	iARC iR_Newton	54 52 (%100)	54 52 (%100)	36564 16267	45769 0
KOWOSB	4	iARC iR_Newton	20 18 (%94)	20 18 (%94)	114 55	121 4
LIARWHD	5000	iARC iR_Newton	12 11 (%100)	12 11 (%100)	46 21	45 0
LOGHAIRY	2	iARC iR_Newton	167 326 (%39)	116 233 (%49)	390 542	414 460
MANCINO	100	iARC iR_Newton	6 4 (%100)	6 4 (%100)	14 5	8 0
MARATOSB	2	iARC iR_Newton	3 3 (%100)	3 3 (%100)	7 4	5 0
MEXHAT	2	iARC iR_Newton	11 11 (%100)	11 11 (%100)	33 22	43 0
MEYER3	3	iARC iR_Newton	12 16 (%75)	12 15 (%73)	48 38	51 20
MODBEALE	20000	iARC iR_Newton	---	---	---	---
			3317 (%99)	3304 (%100)	65293	351
MOREBV	5000	iARC iR_Newton	4 1 (%100)	4 1 (%100)	1102 401	2064 0
MSQRTALS	1024	iARC iR_Newton	39 44 (%73)	33 36 (%83)	9830 4743	12602 149
MSQRTBLS	1024	iARC iR_Newton	32 39 (%69)	26 31 (%81)	5822 3131	7090 156
NCB20	5010	iARC iR_Newton	106 65 (%32)	70 43 (%42)	2888 688	4664 614
NCB20B	5000	iARC iR_Newton	29 38 (%47)	18 19 (%42)	4286 3297	9958 8386
NONCVXU2	5000	iARC iR_Newton	10302 11094 (%100)	10302 11094 (%100)	20604 11094	10302 0
NONCVXUN	5000	iARC iR_Newton	23771 20913 (%100)	23771 20913 (%100)	47542 20913	23771 0
NONDIA	5000	iARC iR_Newton	2 2 (%100)	2 2 (%100)	4 2	2 0
NONDQUAR	5000	iARC iR_Newton	45 38 (%95)	37 36 (%97)	156 70	126 2
OSBORNEA	5	iARC iR_Newton	36 21 (%43)	28 12 (%67)	289 75	412 73
OSBORNEB	11	iARC iR_Newton	25 28 (%75)	19 23 (%78)	396 225	539 105
OSCIGRAD	100000	iARC iR_Newton	13 15 (%40)	10 9 (%56)	190 92	220 61
OSCI PATH	10	iARC iR_Newton	222974 227426 (%59)	131997 134306 (%59)	4233806 2273151	5617938 2521354
PALMER1C	8	iARC iR_Newton	161 74 (%100)	161 74 (%100)	482 145	362 0
PALMER1D	7	iARC iR_Newton	1069 196 (%100)	1069 196 (%100)	3586 379	2567 0
PALMER2C	8	iARC iR_Newton	109 76 (%100)	109 76 (%100)	326 147	245 0
PALMER3C	8	iARC iR_Newton	64 36 (%100)	64 36 (%100)	252 69	201 0
PALMER4C	8	iARC iR_Newton	27 90 (%100)	27 90 (%100)	102 177	84 0
PALMER5C	6	iARC iR_Newton	9 10 (%100)	9 10 (%100)	22 13	13 0
PALMER6C	8	iARC iR_Newton	238 252 (%100)	238 252 (%100)	870 503	654 0
PALMER7C	8	iARC iR_Newton	65 65 (%100)	65 65 (%100)	196 120	143 0
PALMER8C	8	iARC iR_Newton	76 90 (%100)	76 90 (%100)	300 174	229 0
PARKCH	15	iARC	31	22	478	685

Table 2 -- continued from previous page

Prob	#Var	Alg	#Iter (%Newton)	#Acc (%Newton)	#Hv-prod	#T-fact
		iR_Newton	33 (%61)	23 (%74)	270	209
PENALTY1	1000	iARC	14	14	28	14
		iR_Newton	12 (%100)	12 (%100)	12	0
PENALTY2	200	iARC	22	22	314	319
		iR_Newton	22 (%100)	22 (%100)	157	0
PENALTY3	200	iARC	24	20	168	161
		iR_Newton	25 (%60)	18 (%72)	90	44
POWELLSG	5000	iARC	17	17	98	91
		iR_Newton	17 (%100)	17 (%100)	49	0
QUARTC	5000	iARC	15	15	30	15
		iR_Newton	11 (%100)	11 (%100)	11	0
ROSENBR	2	iARC	29	20	87	116
		iR_Newton	32 (%62)	20 (%70)	64	51
S308	2	iARC	12	9	36	41
		iR_Newton	10 (%80)	8 (%88)	20	8
SCHMVETT	5000	iARC	5	5	142	166
		iR_Newton	6 (%100)	6 (%100)	89	0
SENSORS	100	iARC	17	12	146	263
		iR_Newton	21 (%19)	12 (%33)	66	85
SINEVAL	2	iARC	66	42	194	256
		iR_Newton	63 (%63)	41 (%68)	123	81
SINQUAD	5000	iARC	16	11	64	63
		iR_Newton	15 (%33)	9 (%44)	32	26
SISSER	2	iARC	12	12	24	12
		iR_Newton	12 (%100)	12 (%100)	12	0
SNAIL	2	iARC	103	63	290	364
		iR_Newton	107 (%55)	63 (%56)	203	182
SPARSINE	5000	iARC	153	143	15246	18485
		iR_Newton	188 (%88)	174 (%94)	10745	183
SPARSQUR	10000	iARC	15	15	64	49
		iR_Newton	15 (%100)	15 (%100)	32	0
SPMSRTLS	4999	iARC	17	15	582	761
		iR_Newton	17 (%76)	15 (%87)	275	4
SROSENBR	5000	iARC	9	7	36	35
		iR_Newton	10 (%70)	7 (%86)	20	12
SSBRYBND	5000	iARC	75	45	77075	177454
		iR_Newton	39 (%38)	23 (%52)	22010	11269
STRATEC	10	iARC	74	65	886	1069
		iR_Newton	67 (%90)	61 (%95)	413	87
TESTQUAD	5000	iARC	162	162	16908	19812
		iR_Newton	163 (%100)	163 (%100)	8271	0
TOINTGOR	50	iARC	11	11	234	275
		iR_Newton	11 (%100)	11 (%100)	117	0
TOINTGSS	5000	iARC	4	4	14	10
		iR_Newton	3 (%100)	3 (%100)	7	0
TOINTPSP	50	iARC	35	22	254	335
		iR_Newton	41 (%49)	20 (%40)	156	66
TOINTQOR	50	iARC	7	7	104	103
		iR_Newton	7 (%100)	7 (%100)	52	0
TQUARTIC	5000	iARC	11	11	44	50
		iR_Newton	1 (%100)	1 (%100)	2	0
TRIDIA	5000	iARC	16	16	2128	2630
		iR_Newton	17 (%100)	17 (%100)	1310	0
VARDIM	200	iARC	12	12	24	12
		iR_Newton	12 (%100)	12 (%100)	12	0
VAREIGVL	50	iARC	5	5	42	38
		iR_Newton	5 (%100)	5 (%100)	21	0
VIBRBEAM	8	iARC	70	41	644	1131
		iR_Newton	39 (%31)	25 (%48)	183	226
WATSON	12	iARC	14	14	174	214
		iR_Newton	14 (%100)	14 (%100)	88	0
WOODS	4000	iARC	15	15	40	26
		iR_Newton	172 (%87)	157 (%92)	404	144
YFITU	3	iARC	54	38	270	348
		iR_Newton	55 (%65)	39 (%64)	172	130
ZANGWIL2	2	iARC	3	3	6	3
		iR_Newton	1 (%100)	1 (%100)	1	0

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