

Adaptive cubic regularisation methods for unconstrained optimization. Part I: motivation, convergence and numerical results

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Abstract An Adaptive Regularisation algorithm using Cubics (ARC) is proposed for unconstrained optimization, generalizing at the same time an unpublished method due to Griewank (Technical Report NA/12, 1981, DAMTP, University of Cambridge), an algorithm by Nesterov and Polyak (Math Program 108(1):177–205, 2006) and a proposal by Weiser et al. (Optim Methods Softw 22(3):413–431, 2007). At each iteration of our approach, an approximate global minimizer of a local cubic regularisation of the objective function is determined, and this ensures a significant improvement in the objective so long as the Hessian of the objective is locally Lipschitz continuous. The new method uses an adaptive estimation of the local Lipschitz constant and approximations to the global model-minimizer which remain computationally-viable even for large-scale problems. We show that the excellent global and local convergence

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properties obtained by Nesterov and Polyak are retained, and sometimes extended to a wider class of problems, by our ARC approach. Preliminary numerical experiments with small-scale test problems from the CUTEr set show encouraging performance of the ARC algorithm when compared to a basic trust-region implementation.

Keywords Nonlinear optimization · Unconstrained optimization · Cubic regularization · Newton’s method · Trust-region methods · Global convergence · Local convergence

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1 Introduction

Trust-region [4] and line-search [8] methods are two commonly-used convergence schemes for unconstrained optimization and are often used to globalise Newton-like iterations. The work presented in this paper and in its companion [2] explores a third alternative: the use of a cubic overestimator of the objective function as a regularisation technique for the computation of the step from one iterate to the next. Specifically, suppose that we wish to find a local minimizer of $f : \mathbb{R}^n \rightarrow \mathbb{R}$, the smooth objective function of an unconstrained optimization problem, and that x_k is our current best estimate. Furthermore, suppose that the objective’s Hessian $\nabla_{xx}f(x)$ is globally Lipschitz continuous on \mathbb{R}^n with ℓ_2 -norm Lipschitz constant L . Then

$$\begin{aligned} f(x_k+s) &= f(x_k) + s^T g(x_k) + \frac{1}{2} s^T H(x_k) s + \int_0^1 (1-\tau) s^T [H(x_k+\tau s) - H(x_k)] s \, d\tau \\ &\leq f(x_k) + s^T g(x_k) + \frac{1}{2} s^T H(x_k) s + \frac{1}{6} L \|s\|_2^3 \stackrel{\text{def}}{=} m_k^C(s), \quad \text{for all } s \in \mathbb{R}^n, \end{aligned} \quad (1.1)$$

where we have defined $g(x) \stackrel{\text{def}}{=} \nabla_x f(x)$ and $H(x) \stackrel{\text{def}}{=} \nabla_{xx} f(x)$. Thus, so long as

$$m_k^C(s_k) < m_k^C(0) = f(x_k), \quad (1.2)$$

the new iterate $x_{k+1} = x_k + s_k$ improves $f(x)$. The minimization of $m_k^C(s)$ can therefore be used to generate the step s_k , forming the basis of new unconstrained minimization algorithms. Our work on such an algorithm draws on several independent lines of research, some recent and some more than twenty-five years old. We now briefly review the relevant contributions.

The bound (1.1) has been known for a long time, see for example [8, Lemma 4.1.14]. However, the use of the model m_k^C for computing a step was, as far as we know, first considered by Griewank (in an unpublished technical report [17]) as a means for constructing affine-invariant variants of Newton’s method which are globally convergent

to second-order critical points. Griewank introduces a variant of the model $m_k^C(s)$ with a variable weight replacing the constant L as a way to regularise Newton's quadratic model, especially in the presence of negative curvature. This variant is of the form

$$m_k^G(s) \stackrel{\text{def}}{=} f(x_k) + s^T g(x_k) + \frac{1}{2} s^T H(x_k) s + \frac{1}{3} \sigma_k \|s\|_{G_k}^3, \quad (1.3)$$

where $\sigma_k \|\cdot\|_{G_k}$ is iteratively chosen to ensure the overestimation property (1.1) while preserving affine invariance. He then proves global convergence to second-order critical points of a method where the step is computed by finding any second-order minimizer of the model which provides descent. All such minimizers (including the global one) are characterized; the convergence proofs are based on a global Hölder condition and the assumption that the matrices G_k stabilize over the iterations. Griewank also outlines a proof of quadratic local convergence. He finally suggests minimizing m_k^G approximately (using a variant of the nonlinear conjugate-gradients method), considers the rate of convergence of this modified method, and gives some preliminary numerical results.

More recently, Nesterov and Polyak [23] considered a similar idea and the model $m_k^C(s)$, although from a different perspective. They were able to show that, if the step is computed by globally minimizing the cubic model and if the objective's Hessian is globally Lipschitz continuous, then the resulting algorithm has a better global iteration-complexity bound than that achieved by the steepest descent method. They completed this remarkable result by showing that the model's global minimizer could be computed in a manner acceptable from the complexity point of view and proposed superior complexity bounds for the (star) convex and other special cases. Global convergence to second-order critical points and asymptotically quadratic rate of convergence were also proved for this method, but no numerical results were provided. Subsequently, Nesterov [22] has proposed more sophisticated methods which further improve the complexity bounds in the convex case.

Even more recently and again independently, Weiser et al. [26] also pursued the same line of thought, motivated (as Griewank) by the design of an affine-invariant version of Newton's method. Their approach, directly evolved from techniques for convex problems and elaborating on [9], makes use of the cubic model $m_k^G(s)$ with $G_k = \sigma_k G$, where G is positive definite and σ_k is an estimate of the global Lipschitz constant (the techniques for updating σ_k are similar to Griewank's). The proposed method does not consider global model minimization, but rather uses approximate techniques for finding a local minimizer, such as Krylov-space techniques and nonlinear conjugate-gradients. Again, global Lipschitz continuity is assumed, but no formal convergence or complexity analysis is presented. Limited but encouraging numerical experience is discussed.

Our purpose here and in [2] is to unify and extend these contributions into a coherent and numerically efficient algorithmic framework, for which global and asymptotic convergence results can be proved under weaker assumptions and with simpler proofs, while preserving the good complexity bound shown by Nesterov and Polyak [23]. Firstly, we relax the need to compute a global minimizer over \mathbb{R}^n , but show that a global minimizer on some suitable subspace is acceptable for obtaining the

desirable complexity bound. Secondly, though we follow our predecessors [17, 23, 26] by introducing a dynamic positive parameter σ_k instead of the scaled Lipschitz constant¹ $\frac{1}{2}L$ in (1.1), we do not insist that $H(x)$ be globally, or even locally, Lipschitz (or Hölder) continuous in general. Lastly, we allow for a symmetric approximation B_k to the local Hessian $H(x_k)$ in the cubic model on each iteration; this may be highly useful in practice. Thus, instead of (1.1), it is the model

$$m_k(s) \stackrel{\text{def}}{=} f(x_k) + s^T g_k + \frac{1}{2}s^T B_k s + \frac{1}{3}\sigma_k \|s\|^3, \quad (1.4)$$

that we employ as an approximation to f in each iteration of our Adaptive Regularisation with Cubics (ARC) method (the generic algorithmic framework is given in Sect. 2.1). Here, and for the remainder of the paper, for brevity we write $g_k = g(x_k)$ and $\|\cdot\| = \|\cdot\|_2$; our choice of the Euclidean norm for the cubic term is made for simplicity of exposition.

Note that whereas in the previous proposals discussed above, the cubic term of the model is rather closely linked to the size of the third derivative (provided the latter exists at all), in our approach, σ_k performs a double task. Namely, it may account not only for the discrepancy between the objective function and its second order Taylor expansion, but also for the difference between the exact and the approximate Hessian.

The rules for updating the parameter σ_k in the course of the ARC algorithm are justified by analogy to trust-region methods. In such a framework, σ_k might be regarded as the reciprocal of the trust-region radius (see our comments following the proof of Theorem 3.1 and the updating rules for the trust-region radius in [4]). Thus σ_k is increased if insufficient decrease is obtained in some measure of relative objective change, but decreased or unchanged otherwise.

Since finding a global minimizer of the model $m_k(s)$ may not be essential in practice, and as doing so might be prohibitively expensive from a computational point of view, we relax this requirement by letting s_k be an approximation to such a minimizer. Initially, we only require that s_k ensures that the decrease in the model is at least as good as that provided by a suitable Cauchy point obtained by globally minimizing (1.4) along the current negative gradient direction. Also, a milder condition than the inequality in (1.1) is required for the computed step s_k to be accepted. Provided the objective function is continuously differentiable and bounded below, and B_k , bounded above for all k , we show in Sect. 2.2 that the ARC iterates have at least one limit point that is first-order critical. Furthermore, when the gradient of f is uniformly continuous (on the iterates), the ARC algorithm is globally convergent to first-order critical points. The broad applicability of these convergence guarantees is particularly relevant when no (exact) second-order derivative information is available.

To improve on the performance and properties of the ARC algorithm, we further require that the step s_k globally minimizes the model (1.4) in a larger subspace. Suitable candidates include the Krylov subspaces generated by a Lanczos process or, in the limit, the whole of \mathbb{R}^n —recall that the Lanczos process is particularly appropriate for large-scale problems (see Sects. 6.2 and 7). Additional termination rules are

¹ The factor $\frac{1}{2}$ is for later convenience.

specified for the inner iterations, which guarantee that the steps s_k are not too short (see Lemmas 4.7 and 4.9). Any of these rules makes the ARC algorithm converge asymptotically at least Q-superlinearly (see Corollary 4.8 and the first remark following its proof), under appropriate assumptions but without assuming local or global Lipschitz continuity of the Hessian (Theorem 4.3). We also show that the well-known Dennis-Moré condition [7] on the Hessian approximation B_k is sufficient, and certain quasi-Newton formulae are thus appropriate. In the same context, we also show that the parameter σ_k stays bounded above and all steps s_k are eventually accepted (see Theorem 4.3). Under an asymptotic local Lipschitz assumption on $H(x)$, and slightly stronger agreement between B_k and $H(x_k)$ along s_k , Q-quadratic convergence of the iterates is shown when a specific termination criterion is employed (Corollary 4.10). We remark however that, in our numerical experiments, this rule is not the most efficient (see Sect. 7). Requiring asymptotic agreement between B_k and $H(x_k)$ (see (4.19)), without requiring Lipschitz continuity of the Hessian, we show, in a similar fashion to the analogous trust-region results, that the sequence of iterates $\{x_k\}$ is attracted to one of its limit points which is a local minimizer (Theorem 4.5). Without requiring local convexity of the objective as in the latter result, but assuming global Lipschitz continuity of the objective Hessian, we prove that any limit point of the sequence of iterates is weak second-order critical in the sense that the Hessian restricted to the subspaces of minimization is positive semidefinite in the limit (Theorem 5.4).

Worst-case iteration complexity bounds for the ARC family of methods are derived in [2]. Due to the direct correspondence between the count of (major) ARC iterations and the number of function- and derivative(s)-evaluations, these bounds also apply to the latter. When requiring the mild Cauchy condition on the step, and some stronger assumptions on the objective than in the global convergence results, we obtain an upper bound on the total number of iterations the ARC algorithm takes to drive the norm of the gradient of f below ϵ that is of order ϵ^{-2} , which is the same as for the steepest descent method [21, p. 29]. This to be expected since the Cauchy-point condition requires no more than a move in the negative gradient direction. The steepest-descent-like complexity bound can be improved when s_k is the global minimizer of the model (1.4) in a subspace containing the gradient g_k and an appropriate termination criterion is employed. In particular, assuming $H(x)$ to be globally Lipschitz continuous, and the approximation B_k “sufficiently close” to $H(x_k)$ along s_k , we show that the ARC algorithm has an overall worst-case iteration count of order $\epsilon^{-3/2}$ for generating $\|g_k\| \leq \epsilon$, and of order ϵ^{-3} for achieving approximate nonnegative curvature in a subspace containing s_k . These bounds match those proved by Nesterov and Polyak [23, Sect. 3] for their Algorithm 3.3. However, our framework, at least for the first-order results, is more general, as we allow more freedom in the choice of s_k and of B_k .

Despite the good convergence and complexity properties of the ARC algorithm, its practical efficiency ultimately relies on the ability to exactly or approximately minimize the cubic model m_k . Though m_k is non-convex, Theorem 3.1—first proved by different means in [17] and then, independently, in [23]—gives a powerful characterization of its global solutions over \mathbb{R}^n that can be exploited computationally as we show in Sect. 6.1. Our investigations suggest that the model can be globally minimized surprisingly efficiently, provided the factorization of the matrix B_k is (inexpensively) available. Since the latter may not be the case in large-scale optimization, we

also address computing cheaper and approximate minimizers of m_k , namely, global minimizers of m_k over certain subspaces, that do not involve explicit factorizations of B_k , only matrix-vector products (see Sect. 6.2). Our approach involves using the Lanczos process to build up an orthogonal basis for the Krylov subspace formed by successively applying B_k to g_k , and each direction s_k is the global minimizer of the model over the current Krylov subspace. It is easily shown that this technique of approximately minimizing the cubic model when employed with either of our termination criteria, is fully covered by our theoretical results. Furthermore, numerical experience with a Matlab implementation of this approach in the ARC algorithm shows this code to perform consistently better than a basic trust-region implementation when tested on all the small unconstrained problems from the CUTEr test set; see Sect. 7 and Fig. 4 for details. The exact Hessian was used as B_k in both the ARC and the trust-region implementations.

The outline of the paper (Part I) is as follows. Section 2.1 introduces the ARC algorithm, while Sect. 2.2 shows it to be globally convergent to first-order critical points. Section 3.1 gives a new proof to a known characterization of the global minimizer of the cubic model over \mathbb{R}^n , while Sect. 3.2 defines some more general properties that are satisfied by global minimizers of m_k over subspaces of \mathbb{R}^n . Then Sect. 3.3 prescribes some suitable termination criteria for the inner iterations employed to minimize the cubic model approximately. Using the results of Sect. 3, we show asymptotic convergence properties of the ARC algorithm in the presence of local convexity in Sect. 4.1, while we prove in Sect. 4.2 that the ARC algorithm then converges at least Q-superlinearly. Without assuming local convexity, Sect. 5 addresses conditions for the global convergence of the iterates to (weak) second-order critical limit points. Section 6 addresses ways of globally minimizing the cubic model both to high accuracy (Sect. 6.1) as well as approximately using Lanczos techniques (Sect. 6.2). We detail our numerical experiments in Sect. 7 and in Appendix A, and draw final conclusions in Sect. 8.

2 Cubic regularisation for unconstrained minimization

In this section, we assume that

$$\boxed{\text{AF.1}} \quad f \in C^1(\mathbb{R}^n). \tag{2.1}$$

2.1 The method

The iterative methods we shall consider for minimizing $f(x)$ fit into the generic Adaptive Regularisation framework using Cubics (ARC) summarized below.

Algorithm 2.1 Adaptive Regularisation using Cubics (ARC).

Given $x_0, \gamma_2 \geq \gamma_1 > 1, 1 > \eta_2 \geq \eta_1 > 0$, and $\sigma_0 > 0$, for $k = 0, 1, \dots$ until convergence,

1. Compute a step s_k for which

$$m_k(s_k) \leq m_k(s_k^c), \tag{2.2}$$

where the Cauchy point

$$s_k^c = -\alpha_k^c g_k \quad \text{and} \quad \alpha_k^c = \arg \min_{\alpha \in \mathbb{R}_+} m_k(-\alpha g_k). \tag{2.3}$$

2. Compute $f(x_k + s_k)$ and

$$\rho_k = \frac{f(x_k) - f(x_k + s_k)}{f(x_k) - m_k(s_k)}. \tag{2.4}$$

3. Set

$$x_{k+1} = \begin{cases} x_k + s_k & \text{if } \rho_k \geq \eta_1 \\ x_k & \text{otherwise.} \end{cases} \tag{2.5}$$

4. Set

$$\sigma_{k+1} \in \begin{cases} [0, \sigma_k] & \text{if } \rho_k > \eta_2 & \text{[very successful iteration]} \\ [\sigma_k, \gamma_1 \sigma_k] & \text{if } \eta_1 \leq \rho_k \leq \eta_2 & \text{[successful iteration]} \\ [\gamma_1 \sigma_k, \gamma_2 \sigma_k] & \text{otherwise.} & \text{[unsuccessful iteration]} \end{cases} \tag{2.6}$$

□

Given an estimate x_k of a critical point of f , a step s_k is computed that is only required to satisfy condition (2.2), and as such may be easily determined. Note that our ulterior interest is in the case when s_k is computed as an approximate (global) minimizer of the model $m_k(s)$ in (1.4), where B_k is a nontrivial approximation to the Hessian of f (provided the latter exists); these details, however, are not necessary at this point, as we first attempt to derive properties applicable to a wide class of problems and methods. The step s_k is accepted and the new iterate x_{k+1} set to $x_k + s_k$ whenever (a reasonable fraction of) the predicted model decrease $f(x_k) - m_k(s_k)$ is realized by the actual decrease in the objective, $f(x_k) - f(x_k + s_k)$. This is measured by computing the ratio ρ_k in (2.4) and requiring ρ_k to be greater than a prescribed positive constant η_1 (for example, $\eta_1 = 0.1$)—we shall shortly see (Lemma 2.1) that ρ_k is well-defined whenever $g_k \neq 0$. Since the current weight σ_k has resulted in a successful step, there is no pressing reason to increase it, and indeed there may be benefits in decreasing it if good agreement between model and function are observed. By contrast, if ρ_k is smaller than η_1 , we judge that the improvement in objective is insufficient—indeed there is no improvement if $\rho_k \leq 0$. If this happens, the step will be rejected and x_{k+1} left as x_k . Under these circumstances, the only recourse available is to increase the weight σ_k prior to the next iteration with the implicit intention of reducing the size of the step.

We note that, for Lipschitz-continuous Hessians, Griewank [17], Nesterov and Polyak [23] and Weiser et al. [26] all propose techniques for estimating the global Lipschitz constant L in (1.1). This is not our objective in the update (2.6)—even for the case of interest when $B_k \approx H(x_k)$ —since our only concern is local overestimation.

The connection between the construction of the ARC algorithm and of basic trust-region methods (see for example, [4, p. 116]) is superficially evident in the choice

of measure ρ_k and the criteria for step acceptance. At a deeper level, the parameter σ_k might be viewed as the reciprocal of the trust-region radius (see the remarks following the proof of Theorem 3.1). Thus the ways of updating σ_k in each iteration mimic those of changing the trust-region radius. Note that, as in the case of trust-region methods, finding the Cauchy point is computationally inexpensive as it is a one-dimensional minimization of a (two-piece) cubic polynomial; this involves finding roots of a quadratic polynomial and requires one matrix-vector and three vector products.

We remark that, due to the equivalence of norms on \mathbb{R}^n , the ℓ_2 -norm in the model $m_k(s)$ can be replaced by a more general, norm on \mathbb{R}^n of the form $\|x\| \stackrel{\text{def}}{=} \sqrt{x^\top Mx}$, $x \in \mathbb{R}^n$, where M is a given symmetric positive definite matrix. We may even allow for M to depend on k as long as it is uniformly positive definite and bounded as k increases, which may be relevant to preconditioning. It is easy to show that the convergence properties of the ARC algorithm established in what follows remain true in such a more general setting, although some of the constants involved change accordingly. The use of different norms may be viewed as an attempt to achieve affine invariance, an idea pursued by Griewank [17] and Weiser et al. [26]. Note also that regularisation terms of the form $\|s\|^\alpha$, for some $\alpha > 2$, may be employed in $m_k(s)$ instead of the cubic term (complexity aspects of such regularisations are discussed in [2]). Griewank [17] has considered just such extensions to cope with the possibility of Hölder rather than Lipschitz continuous Hessians.

Our aim now is to investigate the global convergence properties of the generic ARC algorithm.

2.2 Global convergence to first-order critical points

Throughout, we denote the index set of all successful iterations of the ARC algorithm by

$$S \stackrel{\text{def}}{=} \{k \geq 0 : k \text{ successful or very successful in the sense of (2.6)}\}. \tag{2.7}$$

We first obtain a guaranteed lower bound on the decrease in f predicted from the cubic model. This also shows that the analogue of (1.2) for m_k holds, provided $g_k \neq 0$.

Lemma 2.1 *Suppose that the step s_k satisfies (2.2). Then for $k \geq 0$, we have that*

$$\begin{aligned} f(x_k) - m_k(s_k) &\geq f(x_k) - m_k(s_k^c) \geq \frac{\|g_k\|^2}{6\sqrt{2} \max [1 + \|B_k\|, 2\sqrt{\sigma_k}\|g_k\|]} \\ &= \frac{\|g_k\|}{6\sqrt{2}} \min \left[\frac{\|g_k\|}{1 + \|B_k\|}, \frac{1}{2} \sqrt{\frac{\|g_k\|}{\sigma_k}} \right]. \end{aligned} \tag{2.8}$$

Proof Due to (2.2) and since the equality in (2.8) is straightforward, it remains to show the second inequality in (2.8). For any $\alpha \geq 0$, using the Cauchy-Schwarz inequality,

we have

$$\begin{aligned}
 m_k(s_k^c) - f(x_k) &\leq m_k(-\alpha g_k) - f(x_k) \\
 &= -\alpha \|g_k\|^2 + \frac{1}{2}\alpha^2 g_k^T B_k g_k + \frac{1}{3}\alpha^3 \sigma_k \|g_k\|^3 \\
 &\leq \alpha \|g_k\|^2 \left\{ -1 + \frac{1}{2}\alpha \|B_k\| + \frac{1}{3}\alpha^2 \sigma_k \|g_k\| \right\}. \tag{2.9}
 \end{aligned}$$

Now $m(s_k^c) \leq f(x_k)$ provided $-1 + \frac{1}{2}\alpha \|B_k\| + \frac{1}{3}\alpha^2 \sigma_k \|g_k\| \leq 0$ and $\alpha \geq 0$, the latter two inequalities being equivalent to

$$\alpha \in [0, \bar{\alpha}_k], \text{ where } \bar{\alpha}_k \stackrel{\text{def}}{=} \frac{3}{2\sigma_k \|g_k\|} \left[-\frac{1}{2}\|B_k\| + \sqrt{\frac{1}{4}\|B_k\|^2 + \frac{4}{3}\sigma_k \|g_k\|} \right].$$

Furthermore, we can express $\bar{\alpha}_k$ as

$$\bar{\alpha}_k = 2 \left[\frac{1}{2}\|B_k\| + \sqrt{\frac{1}{4}\|B_k\|^2 + \frac{4}{3}\sigma_k \|g_k\|} \right]^{-1}.$$

Letting

$$\theta_k \stackrel{\text{def}}{=} \left[\sqrt{2} \max \left(1 + \|B_k\|, 2\sqrt{\sigma_k \|g_k\|} \right) \right]^{-1}, \tag{2.10}$$

and employing the inequalities

$$\begin{aligned}
 \sqrt{\frac{1}{4}\|B_k\|^2 + \frac{4}{3}\sigma_k \|g_k\|} &\leq \frac{1}{2}\|B_k\| + \frac{2}{\sqrt{3}}\sqrt{\sigma_k \|g_k\|} \leq 2 \max \left(\frac{1}{2}\|B_k\|, \frac{2}{\sqrt{3}}\sqrt{\sigma_k \|g_k\|} \right) \\
 &\leq \sqrt{2} \max \left(1 + \|B_k\|, 2\sqrt{\sigma_k \|g_k\|} \right),
 \end{aligned}$$

and

$$\frac{1}{2}\|B_k\| \leq \max \left(1 + \|B_k\|, 2\sqrt{\sigma_k \|g_k\|} \right),$$

it follows that $0 < \theta_k \leq \bar{\alpha}_k$. Thus substituting the value of θ_k in the last inequality in (2.9), we obtain that

$$\begin{aligned}
 m_k(s_k^c) - f(x_k) &\leq \frac{\|g_k\|^2}{\sqrt{2} \max \left(1 + \|B_k\|, 2\sqrt{\sigma_k \|g_k\|} \right)} \\
 &\quad \times \left\{ -1 + \frac{1}{2}\theta_k \|B_k\| + \frac{1}{3}\theta_k^2 \sigma_k \|g_k\| \right\} \leq 0. \tag{2.11}
 \end{aligned}$$

It now follows from the definition (2.10) of θ_k that $\theta_k \|B_k\| \leq 1$ and $\theta_k^2 \sigma_k \|g_k\| \leq 1$, so that the expression in the curly brackets in (2.11) is bounded above by $(-1/6)$. This and (2.11) imply the second inequality in (2.8). \square

In the convergence theory of this section, the quantity $\sqrt{\|g_k\|/\sigma_k}$ plays a role similar to that of the trust-region radius in trust-region methods (compare (2.8) above with the bound (6.3.4) in [4]).

The following assumption will occur frequently in our results. For the model m_k , suppose

$$\boxed{\text{AM.1}} \quad \|B_k\| \leq \kappa_B, \quad \text{for all } k \geq 0, \text{ and some } \kappa_B \geq 0. \quad (2.12)$$

Next we obtain a useful bound on the step.

Lemma 2.2 *Suppose that AM.1 holds and that the step s_k satisfies (2.2). Then*

$$\|s_k\| \leq \frac{3}{\sigma_k} \max(\kappa_B, \sqrt{\sigma_k \|g_k\|}), \quad k \geq 0. \quad (2.13)$$

Proof Consider

$$\begin{aligned} m_k(s) - f(x_k) &= s^T g_k + \frac{1}{2} s^T B_k s + \frac{1}{3} \sigma_k \|s\|^3 \\ &\geq -\|s\| \|g_k\| - \frac{1}{2} \|s\|^2 \|B_k\| + \frac{1}{3} \sigma_k \|s\|^3 \\ &= (\frac{1}{3} \sigma_k \|s\|^3 - \|s\| \|g_k\|) + (\frac{2}{3} \sigma_k \|s\|^3 - \frac{1}{2} \|s\|^2 \|B_k\|). \end{aligned}$$

But then $\frac{1}{3} \sigma_k \|s\|^3 - \|s\| \|g_k\| > 0$ if $\|s\| > 3\sqrt{\|g_k\|/\sigma_k}$, while $\frac{2}{3} \sigma_k \|s\|^3 - \frac{1}{2} \|s\|^2 \|B_k\| > 0$ if $\|s\| > \frac{3}{4} \|B_k\|/\sigma_k$. Hence $m_k(s) > f(x_k)$ whenever

$$\|s\| > \frac{3}{\sigma_k} \max(\|B_k\|, \sqrt{\sigma_k \|g_k\|}).$$

But $m_k(s_k) \leq f(x_k)$ due to (2.8), and thus (2.13) holds, recalling also AM.1. □

For the proof of the next result, and some others to follow, we need to show that, under certain conditions, a step k is very successful in the sense of (2.6). Provided $f(x_k) > m_k(s_k)$, and recalling (2.4), we have

$$\rho_k > \eta_2 \iff r_k \stackrel{\text{def}}{=} f(x_k + s_k) - f(x_k) - \eta_2 [m_k(s_k) - f(x_k)] < 0. \quad (2.14)$$

Whenever $f(x_k) > m_k(s_k)$, we can express r_k as

$$r_k = f(x_k + s_k) - m_k(s_k) + (1 - \eta_2) [m_k(s_k) - f(x_k)], \quad k \geq 0. \quad (2.15)$$

An auxiliary lemma is given next.

Lemma 2.3 *Let AF.1 and AM.1 hold. Suppose that \mathcal{I} is an infinite index set such that*

$$\|g_k\| \geq \epsilon, \text{ for all } k \in \mathcal{I} \text{ and some } \epsilon > 0, \text{ and } \sqrt{\frac{\|g_k\|}{\sigma_k}} \rightarrow 0, \text{ as } k \rightarrow \infty, \quad k \in \mathcal{I}. \quad (2.16)$$

Then

$$\|s_k\| \leq 3\sqrt{\frac{\|g_k\|}{\sigma_k}}, \text{ for all } k \in \mathcal{I} \text{ sufficiently large.} \tag{2.17}$$

Additionally, if

$$x_k \rightarrow x_*, \text{ as } k \rightarrow \infty, \quad k \in \mathcal{I}, \text{ for some } x_* \in \mathbb{R}^n, \tag{2.18}$$

then each iteration $k \in \mathcal{I}$ that is sufficiently large is very successful, and

$$\sigma_{k+1} \leq \sigma_k, \text{ for all } k \in \mathcal{I} \text{ sufficiently large.} \tag{2.19}$$

Proof As (2.16) implies $\sqrt{\sigma_k \|g_k\|} \geq \epsilon \sqrt{\sigma_k / \|g_k\|} \rightarrow \infty$, as $k \rightarrow \infty, k \in \mathcal{I}$, the bound (2.17) now follows from (2.13). To prove (2.19), we use (2.14) and (2.15). Hence we first need to estimate the difference between the function and the model at $x_k + s_k$. A Taylor expansion of $f(x_k + s_k)$ around x_k gives

$$f(x_k + s_k) - m_k(s_k) = (g(\xi_k) - g_k)^T s_k - \frac{1}{2} s_k^T B_k s_k - \frac{\sigma_k}{3} \|s_k\|^3, \quad k \geq 0,$$

for some ξ_k on the line segment $(x_k, x_k + s_k)$, which, by employing AM.1 and (2.17), further gives

$$f(x_k + s_k) - m_k(s_k) \leq 3 \left\{ \|g(\xi_k) - g_k\| + \frac{3\kappa_B}{2} \sqrt{\frac{\|g_k\|}{\sigma_k}} \right\} \cdot \sqrt{\frac{\|g_k\|}{\sigma_k}}, \tag{2.20}$$

for some ξ_k on the line segment $(x_k, x_k + s_k)$ and for all $k \in \mathcal{I}$ sufficiently large. To bound the remaining term in (2.15), from (2.8), AM.1 and the inequality in (2.16), we obtain

$$f(x_k) - m_k(s_k) \geq \frac{\epsilon}{6\sqrt{2}} \min \left[\frac{\epsilon}{1 + \kappa_B}, \frac{1}{2} \sqrt{\frac{\|g_k\|}{\sigma_k}} \right], \text{ for all } k \in \mathcal{I},$$

which, by employing the limit in (2.16), further gives

$$f(x_k) - m_k(s_k) \geq \frac{\epsilon}{12\sqrt{2}} \sqrt{\frac{\|g_k\|}{\sigma_k}}, \text{ for all } k \in \mathcal{I} \text{ sufficiently large.}$$

This, (2.15) and (2.20) imply

$$r_k \leq \sqrt{\frac{\|g_k\|}{\sigma_k}} \left\{ 3\|g(\xi_k) - g_k\| + \frac{9\kappa_B}{2} \sqrt{\frac{\|g_k\|}{\sigma_k}} - \frac{(1 - \eta_2)\epsilon}{12\sqrt{2}} \right\},$$

for all $k \in \mathcal{I}$ sufficiently large. (2.21)

Since ξ_k belongs to the line segment $(x_k, x_k + s_k)$, we have $\|\xi_k - x_*\| \leq \|x_k - x_*\| + \|s_k\|$. Also, the limit in (2.16) and (2.17) imply $\|s_k\| \rightarrow 0$, as $k \rightarrow \infty, k \in \mathcal{I}$. It now follows from (2.18) that $\xi_k \rightarrow x_*$, as $k \rightarrow \infty, k \in \mathcal{I}$, and since g is continuous, we conclude that

$$\|g(\xi_k) - g_k\| \leq \|g(\xi_k) - g(x_*)\| + \|g_k - g(x_*)\| \rightarrow 0, \quad k \in \mathcal{I}, \quad k \rightarrow \infty.$$

This, the limit in (2.16) and (2.21) imply that $r_k < 0$, for all $k \in \mathcal{I}$ sufficiently large and thus, recalling (2.14), k is very successful. The inequality (2.19) now follows from (2.6). □

Note that if all iterations k sufficiently large belong to the iteration index set \mathcal{I} satisfying the conditions in Lemma 2.3, then a contradiction ensues from (2.19) and (2.16), implying that at least one of the two relations in (2.16) cannot hold. This is the crux of the convergence arguments in the proofs of Lemma 2.4 and Theorem 2.5, since, for example, whenever the first part of (2.16) cannot hold, it implies that $\liminf_{k \rightarrow \infty} \|g_k\| = 0$.

Next, we show that provided there are only finitely many successful iterations, all later iterates are first-order critical points.

Lemma 2.4 *Let AF.1 and AM.1 hold. Suppose furthermore that there are only finitely many successful iterations. Then $x_k = x_*$ for all sufficiently large k and $g(x_*) = 0$.*

Proof After the last successful iterate is computed, indexed by say k_0 , the construction of the algorithm implies that $x_{k_0+1} = x_{k_0+i} \stackrel{\text{def}}{=} x_*$, for all $i \geq 1$. If $\|g_{k_0+1}\| > 0$, then

$$\|g_k\| = \|g_{k_0+1}\| \stackrel{\text{def}}{=} \epsilon > 0, \quad \text{for all } k \geq k_0 + 1. \tag{2.22}$$

Thus, letting $I := \{k : k \geq k_0 + 1\}$ in Lemma 2.3, we have that (2.18) and the first part of (2.16) hold. Furthermore, since all iterations $k \geq k_0 + 1$ are unsuccessful, σ_k increases by at least a fraction γ_1 so that

$$\sigma_k \rightarrow \infty, \quad \text{as } k \rightarrow \infty,$$

which together with (2.22), implies that the second part of (2.16) is also achieved. Thus, Lemma 2.3 provides that each iteration $k \geq k_0 + 1$ sufficiently large is very successful. This contradicts the fact that $k \geq k_0 + 1$ is unsuccessful, and so $g_k = g(x_*) = 0, k \geq k_0 + 1$. □

We are now ready to prove the first convergence result for the ARC algorithm. In particular, we show that provided f is bounded from below, either we are in the above case and $g_k = 0$ for some finite k , or there is a subsequence of $\{g_k\}$ converging to zero.

Theorem 2.5 *Suppose that AF.1 and AM.1 hold. If $\{f(x_k)\}$ is bounded below, then*

$$\liminf_{k \rightarrow \infty} \|g_k\| = 0. \tag{2.23}$$

Proof Lemma 2.4 shows that the result is true when there are only finitely many successful iterations. Let us now assume infinitely many successful iterations occur, and recall the notation (2.7). We also assume that (2.23) does not hold, and so

$$\|g_k\| \geq \epsilon, \quad \text{for some } \epsilon > 0 \text{ and for all } k \geq 0. \tag{2.24}$$

Let us first prove that (2.24) implies that

$$\sum_{k \in \mathcal{S}} \sqrt{\frac{\|g_k\|}{\sigma_k}} < +\infty. \tag{2.25}$$

It follows from (2.4), (2.5), (2.8), AM.1 and (2.24) that

$$\begin{aligned} f(x_k) - f(x_{k+1}) &\geq \eta_1 [f(x_k) - m_k(s_k)] \\ &\geq \frac{\eta_1 \epsilon}{6\sqrt{2}} \cdot \min \left[\frac{\epsilon}{1 + \kappa_b}, \frac{1}{2} \sqrt{\frac{\|g_k\|}{\sigma_k}} \right], \quad \text{for all } k \in \mathcal{S}. \end{aligned} \tag{2.26}$$

Since $\{f(x_k)\}$ is monotonically decreasing and assumed to be bounded below, it is convergent, and hence, the minimum on the right-hand side of (2.26) will be attained at $\sqrt{\|g_k\|}/(2\sqrt{\sigma_k})$ as the left-hand side of (2.26) converges to zero. Thus we obtain

$$f(x_k) - f(x_{k+1}) \geq \frac{\eta_1 \epsilon}{12\sqrt{2}} \sqrt{\frac{\|g_k\|}{\sigma_k}}, \quad \text{for all } k \in \mathcal{S} \text{ sufficiently large,}$$

which summed up over all sufficiently large iterations provides

$$f(x_{k_0}) - f(x_{j+1}) = \sum_{k=k_0, k \in \mathcal{S}}^j [f(x_k) - f(x_{k+1})] \geq \frac{\eta_1 \epsilon}{12\sqrt{2}} \sum_{k=k_0, k \in \mathcal{S}}^j \sqrt{\frac{\|g_k\|}{\sigma_k}}, \tag{2.27}$$

for some iteration index k_0 sufficiently large and for any $j \in \mathcal{S}, j \geq k_0$. Thus, since $\{f(x_{j+1})\}$ is convergent, (2.25) follows by letting $j \rightarrow \infty$ in (2.27).

Now let us argue that the sequence of iterates $\{x_k\}, k \geq 0$, is a Cauchy sequence. Note that (2.25) implies

$$\sqrt{\|g_k\|/\sigma_k} \rightarrow 0, \quad k \rightarrow \infty, \quad k \in \mathcal{S}, \tag{2.28}$$

and so, also from (2.24), we deduce that (2.16) holds with $\mathcal{I} := \mathcal{S}$. It follows from (2.17) and the construction of the algorithm that

$$\begin{aligned} \|x_{l+r} - x_l\| &\leq \sum_{k=l}^{l+r-1} \|x_{k+1} - x_k\| = \sum_{k=l, k \in \mathcal{S}}^{l+r-1} \|s_k\| \\ &\leq 3 \sum_{k=l, k \in \mathcal{S}}^{l+r-1} \sqrt{\frac{\|g_k\|}{\sigma_k}}, \quad \text{for } l \geq 0 \text{ sufficiently large, } r \geq 0, \end{aligned}$$

whose right-hand side tends to zero as $l \rightarrow \infty$ due to (2.25). Thus $\{x_k\}$ is a Cauchy sequence, and

$$x_k \rightarrow x_*, \quad k \rightarrow \infty, \quad \text{for some } x_* \in \mathbb{R}^n. \tag{2.29}$$

From (2.24), (2.28) and (2.29), we have that (2.16) and (2.18) hold with $\mathcal{I} := \mathcal{S}$. Thus Lemma 2.3 provides that all $k \in \mathcal{S}$ sufficiently large are very successful, and (2.19) holds with $\mathcal{I} := \mathcal{S}$. Now, if all k sufficiently large belong to \mathcal{S} , namely, there are no unsuccessful iterations for k sufficiently large, then (2.19) implies that $\sigma_{k+1} \leq \sigma_k$, for all k sufficiently large, and so $\{\sigma_k\}$, $k \geq 0$, is bounded above. This however, contradicts $\sigma_k \rightarrow \infty$, which follows from (2.24) and (2.28). Thus (2.24) cannot hold.

It remains to show that all sufficiently large iterations belong, indeed, to \mathcal{S} . Let us assume the contrary: recalling that \mathcal{S} is assumed to be infinite (and hence, not all iterations can be consecutively unsuccessful for all k sufficiently large), let $\{k_i\}$ denote an (infinite) subsequence of very successful iterations such that $k_i - 1$ is unsuccessful for all $i \geq 0$ (as all $k \in \mathcal{S}$ sufficiently large are very successful, without loss of generality, we can disregard successful iterations). Then (2.6) implies $\sigma_{k_i} \leq \gamma_2 \sigma_{k_i-1}$, for all i , and $k_i - 1$ being unsuccessful gives $g_{k_i} = g_{k_i-1}$, for all i . Thus, also from (2.28), we deduce

$$\sqrt{\|g_{k_i-1}\|/\sigma_{k_i-1}} \rightarrow 0, \quad \text{as } i \rightarrow \infty. \tag{2.30}$$

It follows from (2.24), (2.29) and (2.30) that (2.16) and (2.18) are satisfied with $\mathcal{I} := \{k_i - 1 : i \geq 0\}$, and so Lemma 2.3 provides that $k_i - 1$ is very successful for all i sufficiently large. This contradicts our assumption that $k_i - 1$ is unsuccessful for all i . □

To show that the whole sequence of gradients $\{g_k\}$ converges to zero, we employ the additional assumption that the gradient g is uniformly continuous on the sequence of iterates $\{x_k\}$, namely,

AF.2

$$\|g_{t_i} - g_{l_i}\| \rightarrow 0 \text{ whenever } \|x_{t_i} - x_{l_i}\| \rightarrow 0, \quad i \rightarrow \infty, \tag{2.31}$$

where $\{x_{t_i}\}$ and $\{x_{l_i}\}$ are subsequences of $\{x_k\}$. Clearly, AF.2 is satisfied if g is uniformly continuous on \mathbb{R}^n . It is also achieved if g is globally Lipschitz continuous on $\{x_k\}$ (see (4.21)); if $f \in \mathcal{C}^2(\mathbb{R}^n)$, then the latter, and hence AF.2, holds if the Hessian of f is bounded above on the convex hull of all the iterates.

Corollary 2.6 *Let AF.1–AF.2 and AM.1 hold. If $\{f(x_k)\}$ is bounded below, then*

$$\lim_{k \rightarrow \infty} \|g_k\| = 0. \tag{2.32}$$

Proof If there are finitely many successful iterations, then Lemma 2.4 implies that (2.32) holds. Now let \mathcal{S} be infinite, and assume that there is an infinite subsequence $\{t_i\} \subseteq \mathcal{S}$ such that

$$\|g_{t_i}\| \geq 2\epsilon, \quad \text{for some } \epsilon > 0 \text{ and for all } i. \tag{2.33}$$

Note that only successful iterates need to be considered since the gradient remains constant on all the other iterates (due to the construction of the algorithm). Theorem 2.5 implies that for each t_i , there is a first successful iteration $l_i > t_i$ such that $\|g_{l_i}\| < \epsilon$. Thus $\{l_i\} \subseteq \mathcal{S}$ and for all i , we have

$$\|g_k\| \geq \epsilon, \quad \text{for all } k \text{ with } t_i \leq k < l_i, \quad \text{and} \quad \|g_{l_i}\| < \epsilon. \tag{2.34}$$

Let $\mathcal{K} \stackrel{\text{def}}{=} \{k \in \mathcal{S} : t_i \leq k < l_i\}$, where the subsequences $\{t_i\}$ and $\{l_i\}$ were defined above; note that \mathcal{K} is also infinite. Since $\mathcal{K} \subseteq \mathcal{S}$, it follows from (2.4), (2.5), (2.8), AM.1 and (2.34) that

$$f(x_k) - f(x_{k+1}) \geq \frac{\eta_1 \epsilon}{6\sqrt{2}} \cdot \min \left[\frac{\epsilon}{1 + \kappa_B}, \frac{1}{2} \sqrt{\frac{\|g_k\|}{\sigma_k}} \right], \quad k \in \mathcal{K}. \tag{2.35}$$

Since $\{f(x_k)\}$ is monotonically decreasing and bounded from below, it is convergent, and hence the left-hand side of (2.35) converges to zero as $k \rightarrow \infty$. Thus (2.35) implies

$$\sqrt{\|g_k\|/\sigma_k} \rightarrow 0, \quad k \rightarrow \infty, \quad k \in \mathcal{K}. \tag{2.36}$$

Due to (2.36), the bound (2.35) asymptotically becomes

$$f(x_k) - f(x_{k+1}) \geq \frac{\eta_1 \epsilon}{12\sqrt{2}} \sqrt{\frac{\|g_k\|}{\sigma_k}}, \quad \text{for all } k \in \mathcal{K} \text{ sufficiently large.} \tag{2.37}$$

From (2.36) and (2.34), we have that (2.16) is satisfied with $\mathcal{I} := \mathcal{K}$, and thus (2.17) holds for this choice of \mathcal{I} , which together with (2.37) and the definition of \mathcal{K} , provides the bound

$$f(x_k) - f(x_{k+1}) \geq \frac{\eta_1 \epsilon}{36\sqrt{2}} \|s_k\|, \quad \text{for all } t_i \leq k < l_i, \quad k \in \mathcal{S}, \quad i \text{ sufficiently large.} \tag{2.38}$$

Summing up (2.38) over k with $t_i \leq k < l_i$, and employing (2.5) and the triangle inequality, we obtain

$$\frac{36\sqrt{2}}{\eta_1\epsilon} [f(x_{t_i}) - f(x_{l_i})] \geq \sum_{k=t_i, k \in \mathcal{S}}^{l_i-1} \|s_k\| = \sum_{k=t_i}^{l_i-1} \|x_{k+1} - x_k\| \geq \|x_{t_i} - x_{l_i}\|, \quad (2.39)$$

for all i sufficiently large. Since $\{f(x_k)\}$ is convergent, $\{f(x_{t_i}) - f(x_{l_i})\}$ converges to zero as $i \rightarrow \infty$. Thus (2.39) implies that $\|x_{t_i} - x_{l_i}\|$ converges to zero as $i \rightarrow \infty$, and by AF.2, $\|g_{l_i} - g_{t_i}\|$ tends to zero. We have reached a contradiction, since (2.33) and (2.34) imply $\|g_{l_i} - g_{t_i}\| \geq \|g_{t_i}\| - \|g_{l_i}\| \geq \epsilon$ for all $i \geq 0$. \square

From now on, we assume throughout that

$$g_k \neq 0, \quad \text{for all } k \geq 0; \quad (2.40)$$

we will discuss separately the case when $g_l = 0$ for some l (see our remarks at the end of Sect. 3.2 and Sect. 5). It follows from (2.8) and (2.40) that

$$f(x_k) > m_k(s_k), \quad k \geq 0. \quad (2.41)$$

A comparison of the above results to those in [20] outlines the similarities of the two approaches, as well as the differences; see also [25]. Compare for example, Lemma 2.4, Theorem 2.5 and Corollary 2.6 to Theorems 4.10 and 4.14 in [20]. In particular, note that we extended the results in [20] as no lower bound was imposed in ARC on the rate of decrease of σ_k on very successful iterations, while the change to the trust-region radius on the same type of iterations is required to be bounded (above) in [20].

3 On approximate minimizers of the model

3.1 Optimality conditions for the minimizer of m_k over \mathbb{R}^n

In this first subsection, we give a new proof of a result concerning necessary and sufficient optimality conditions for the *global* minimizer of the cubic model, for which different proofs are provided by Griewank [17] and Nesterov and Polyak [23, Sect. 5.1]. Our approach is closer in spirit to trust-region techniques, thus offering new insight into this surprising result, as well as a proper fit in the context of our paper.

We may express the derivatives of the cubic model $m_k(s)$ in (1.4) as

$$\nabla_s m_k(s) = g_k + B_k s + \lambda s \quad \text{and} \quad \nabla_{ss} m_k(s) = B_k + \lambda I + \lambda \left(\frac{s}{\|s\|} \right) \left(\frac{s}{\|s\|} \right)^T, \quad (3.1)$$

where $\lambda = \sigma_k \|s\|$ and I is the n by n identity matrix.

We have the following global optimality result.

Theorem 3.1 Any s_k^* is a global minimizer of $m_k(s)$ over \mathbb{R}^n if and only if it satisfies the system of equations

$$(B_k + \lambda_k^* I) s_k^* = -g_k, \tag{3.2}$$

where $\lambda_k^* = \sigma_k \|s_k^*\|$ and $B_k + \lambda_k^* I$ is positive semidefinite. If $B_k + \lambda_k^* I$ is positive definite, s_k^* is unique.

Proof In this proof, we drop the iteration subscript k for simplicity. Firstly, let s^* be a global minimizer of $m(s)$ over \mathbb{R}^n . It follows from (3.1) and the first- and second-order necessary optimality conditions at s^* that

$$g + (B + \lambda^* I) s^* = 0,$$

and hence that (3.2) holds, and that

$$w^T \left(B + \lambda^* I + \lambda^* \left(\frac{s^*}{\|s^*\|} \right) \left(\frac{s^*}{\|s^*\|} \right)^T \right) w \geq 0 \tag{3.3}$$

for all vectors w .

If $s^* = 0$, (3.3) is equivalent to $\lambda^* = 0$ and B being positive semi-definite, which immediately gives the required result. Thus we need only consider $s^* \neq 0$.

There are two cases to consider. Firstly, suppose that $w^T s^* = 0$. In this case, it immediately follows from (3.3) that

$$w^T (B + \lambda^* I) w \geq 0 \quad \text{for all } w \text{ for which } w^T s^* = 0. \tag{3.4}$$

It thus remains to consider vectors w for which $w^T s^* \neq 0$. Since w and s^* are not orthogonal, the line $s^* + \alpha w$ intersects the ball of radius $\|s^*\|$ at two points, s^* and $u^* \neq s^*$, say, and thus

$$\|u^*\| = \|s^*\|. \tag{3.5}$$

We let $w^* = u^* - s^*$, and note that w^* is parallel to w .

Since s^* is a global minimizer, we immediately have that

$$\begin{aligned} 0 &\leq m(u^*) - m(s^*) \\ &= g^T (u^* - s^*) + \frac{1}{2} (u^*)^T B u^* - \frac{1}{2} (s^*)^T B s^* + \frac{\sigma}{3} (\|u^*\|^3 - \|s^*\|^3) \\ &= g^T (u^* - s^*) + \frac{1}{2} (u^*)^T B u^* - \frac{1}{2} (s^*)^T B s^*, \end{aligned} \tag{3.6}$$

where the last equality follows from (3.5). But (3.2) gives that

$$g^T (u^* - s^*) = (s^* - u^*)^T B s^* + \lambda^* (s^* - u^*)^T s^*. \tag{3.7}$$

In addition, (3.5) shows that

$$(s^* - u^*)^T s^* = \frac{1}{2}(s^*)^T s^* + \frac{1}{2}(u^*)^T u^* - (u^*)^T s^* = \frac{1}{2}(w^*)^T w^*. \tag{3.8}$$

Thus combining (3.6) and (3.7), we find that

$$\begin{aligned} 0 &\leq \frac{1}{2}\lambda^*(w^*)^T w^* + \frac{1}{2}(u^*)^T B u^* - \frac{1}{2}(s^*)^T B s^* + (s^*)^T B s^* - (u^*)^T B s^* \\ &= \frac{1}{2}(w^*)^T (B + \lambda^* I) w^* \end{aligned} \tag{3.9}$$

from which we deduce that

$$w^T (B + \lambda^* I) w \geq 0 \quad \text{for all } w \text{ for which } w^T s^* \neq 0. \tag{3.10}$$

Hence (3.4) and (3.10) together show that $B + \lambda^* I$ is positive semidefinite. The uniqueness of s^* when $B + \lambda^* I$ is positive definite follows immediately from (3.2). For the sufficiency implication, let s^* satisfy (3.2) and note that since $m_k(s)$ is bounded below, a global minimizer, say u^* , exists and also satisfies (3.2). Then, whenever B is positive semidefinite, or indefinite with g not orthogonal to the eigenvector of B corresponding to its leftmost eigenvalue, say λ_1 , the argument in the second and third paragraphs of Sect. 6.1 implies that there exists a unique solution (s, λ) to the system $(B + \lambda I)s = -g$ and $\lambda = \sigma \|s\|$, and so $s^* = u^*$. Else, $\lambda^* = -\lambda_1$ and s^* and u^* are solutions of (6.6) for some values of α . It follows that $\|u^*\| = \|s^*\|$ and the equalities in (3.6), and (3.8)–(3.10) give $m(u^*) \geq m(s^*)$, which implies $m(u^*) = m(s^*)$ since u^* is a global minimizer. Thus so is s^* . \square

Note how similar this result and its proof are to those for the trust-region subproblem (see [4, Theorem 7.2.1]), for which we aim to minimize $g_k^T s + \frac{1}{2}s^T B_k s$ within an ℓ_2 -norm trust region $\|s\| \leq \Delta_k$ for some “radius” $\Delta_k > 0$. Often, the global solution s_k^* of this subproblem satisfies $\|s_k^*\| = \Delta_k$. Then, recalling that s_k^* would also satisfy (3.2), we have from Theorem 3.1 that $\sigma_k = \lambda_k^*/\Delta_k$. Hence one might interpret the parameter σ_k in the ARC algorithm as inversely proportional to the trust-region radius.

In Sect. 6.1, we discuss ways of computing the global minimizer s_k^* .

3.2 Minimizing the cubic model in a subspace

The only requirement on the step s_k computed by the ARC algorithm has been that it satisfies the Cauchy condition (2.2). As we showed in Sect. 2.2, this is enough for the algorithm to converge to first-order critical points. To be able to guarantee stronger convergence properties for the ARC algorithm, further requirements need to be placed on s_k . The strongest such conditions are, of course, the first and second order (necessary) optimality conditions that s_k satisfies provided it is the (exact) global minimizer of $m_k(s)$ over \mathbb{R}^n (see Theorem 3.1). This choice of s_k , however, may be in general prohibitively expensive from a computational point of view, and thus, for most (large-scale) practical purposes, (highly) inefficient (see Sect. 6.1). As in the case of trust-region methods, a much more useful approach in practice is to compute

an approximate global minimizer of $m_k(s)$ by (globally) minimizing the model over a sequence of (nested) subspaces, in which each such subproblem is computationally quite inexpensive (see Sect. 6.2). Thus the conditions we require on s_k in what follows, are some derivations of first- and second-order optimality when s_k is the global minimizer of m_k over a subspace (see (3.11), (3.12) and Lemma 3.2). Then, provided each subspace includes g_k , not only do the previous results still hold, but we can prove further convergence properties (see Sect. 4.1) and deduce good complexity bounds (see [2]) for the ARC algorithm. Furthermore, our approach and results widen the scope of the convergence and complexity analysis in [23] which addresses solely the case of the exact global minimizer of m_k over \mathbb{R}^n .

In what follows, we require that s_k satisfies

$$g_k^\top s_k + s_k^\top B_k s_k + \sigma_k \|s_k\|^3 = 0, \quad k \geq 0, \tag{3.11}$$

and

$$s_k^\top B_k s_k + \sigma_k \|s_k\|^3 \geq 0, \quad k \geq 0. \tag{3.12}$$

Note that (3.11) is equivalent to $\nabla_s m_k(s_k)^\top s_k = 0$, due to (3.1).

The next lemma presents some suitable choices for s_k that achieve (3.11) and (3.12).

Lemma 3.2 *Suppose that s_k is the global minimizer of $m_k(s)$, for $s \in \mathcal{L}_k$, where \mathcal{L}_k is a subspace of \mathbb{R}^n . Then s_k satisfies (3.11) and (3.12). Furthermore, letting Q_k denote any orthogonal matrix whose columns form a basis of \mathcal{L}_k , we have that*

$$Q_k^\top B_k Q_k + \sigma_k \|s_k\| I \text{ is positive semidefinite.} \tag{3.13}$$

In particular, if s_k^ is the global minimizer of $m_k(s)$, $s \in \mathbb{R}^n$, then s_k^* achieves (3.11) and (3.12).*

Proof Let s_k be the global minimizer of m_k over some \mathcal{L}_k , i.e., s_k solves

$$\min_{s \in \mathcal{L}_k} m_k(s). \tag{3.14}$$

Let l denote the dimension of the subspace \mathcal{L}_k . Let Q_k be an orthogonal $n \times l$ matrix whose columns form a basis of \mathcal{L}_k . Thus $Q_k^\top Q_k = I$ and for all $s \in \mathcal{L}_k$, we have $s = Q_k u$, for some $u \in \mathbb{R}^l$. Recalling that s_k solves (3.14), and letting

$$s_k = Q_k u_k, \tag{3.15}$$

we have that u_k is the global minimizer of

$$\min_{u \in \mathbb{R}^l} m_{k,r}(u) \stackrel{\text{def}}{=} f(x_k) + (Q_k^\top g_k)^\top u + \frac{1}{2} u^\top Q_k^\top B_k Q_k u + \frac{1}{3} \sigma_k \|u\|^3, \tag{3.16}$$

where we have used the following property of the Euclidean norm when applied to orthogonal matrices,

$$\|Q_k u\| = \|u\|, \quad \text{for all } u. \tag{3.17}$$

Applying Theorem 3.1 to the reduced model $m_{k,r}$ and u_k , it follows that

$$Q_k^\top B_k Q_k u_k + \sigma_k \|u_k\| u_k = -Q_k^\top g_k,$$

and multiplying by u_k , we have

$$u_k^\top Q_k^\top B_k Q_k u_k + \sigma_k \|u_k\|^3 = -g_k^\top Q_k u_k,$$

which is the same as (3.11), due to (3.15) and (3.17). Moreover, Theorem 3.1 implies that $Q_k^\top B_k Q_k + \sigma_k \|u_k\| I$ is positive semidefinite. Due to (3.15) and (3.17), this is (3.13), and also implies

$$u_k^\top Q_k^\top B_k Q_k u_k + \sigma_k \|u_k\|^3 \geq 0,$$

which is (3.12). □

Note that the Cauchy point (2.3) satisfies (3.11) and (3.12) since it globally minimizes m_k over the subspace generated by $-g_k$. To improve the properties and performance of ARC, however, it may be necessary to minimize m_k over (increasingly) larger subspaces.

The next lemma gives a lower bound on the model decrease when (3.11) and (3.12) are satisfied.

Lemma 3.3 *Suppose that s_k satisfies (3.11). Then*

$$f(x_k) - m_k(s_k) = \frac{1}{2} s_k^\top B_k s_k + \frac{2}{3} \sigma_k \|s_k\|^3. \tag{3.18}$$

Additionally, if s_k also satisfies (3.12), then

$$f(x_k) - m_k(s_k) \geq \frac{1}{6} \sigma_k \|s_k\|^3. \tag{3.19}$$

Proof Relation (3.18) can be obtained by eliminating the term $s_k^\top g_k$ from (1.4) and (3.11). It follows from (3.12) that $s_k^\top B_k s_k \geq -\sigma_k \|s_k\|^3$, which we then substitute into (3.18) and obtain (3.19). □

Requiring that s_k satisfies (3.11) may not necessarily imply (2.2), unless $s_k = -g_k$. Nevertheless, when minimizing m_k globally over successive subspaces, condition (2.2) can be easily ensured by including g_k in each of the subspaces. This is the approach we take in our implementation of the ARC algorithm, where the subspaces generated by

Lanczos method naturally include the gradient (see Sects. 6 and 7). Thus, throughout, we assume the Cauchy condition (2.2) still holds.

The assumption (2.40) provides the implication

$$s_k \text{ satisfies (3.11)} \implies s_k \neq 0. \tag{3.20}$$

To see this, assume $s_k = 0$. Then (3.18) gives $f(x_k) = m_k(s_k)$. This, however, contradicts (2.41).

In the case when $g_k = 0$ for some $k \geq 0$ and thus assumption (2.40) is not satisfied, we need to be more careful. If s_k minimizes m_k over a subspace \mathcal{L}_k generated by the columns of some orthogonal matrix Q_k , we have

$$(3.13) \text{ holds and } \lambda_{\min}(Q_k^\top B_k Q_k) < 0 \implies s_k \neq 0, \tag{3.21}$$

since Lemma 3.2 holds even when $g_k = 0$. But if $\lambda_{\min}(Q_k^\top B_k Q_k) \geq 0$ and $g_k = 0$, then $s_k = 0$ and the ARC algorithm will terminate. Hence, if our intention is to identify whether B_k is indefinite, it will be necessary to build Q_k so that $Q_k^\top B_k Q_k$ predicts negative eigenvalues of B_k . This will ultimately be the case with probability one if Q_k is built as the Lanczos basis of the Krylov space $\{B_k^l v\}_{l \geq 0}$ for some random initial vector $v \neq 0$. Note that we have the implication

$$(3.19), (3.21) \text{ and } \sigma_k > 0 \implies (2.41), \tag{3.22}$$

and thus the step will reduce the model.

3.3 Termination criteria for the approximate minimization of m_k

In the previous section, the bound (3.19) on the model decrease was deduced. However, for this to be useful for investigating rates of convergence (and later on in [2], complexity bounds) for the ARC algorithm, we must ensure that s_k does not become too small compared to the size of the gradient. To deduce a lower bound on $\|s_k\|$, we need to be more specific about the ARC algorithm. In particular, suitable termination criteria for the method used to minimize $m_k(s)$ need to be made precise.

Let us assume that some iterative solver is used on each (major) iteration k to approximately minimize $m_k(s)$. Let us set the termination criterion for its inner iterations i to be

$$\|\nabla_s m_k(s_{i,k})\| \leq \theta_{i,k} \|g_k\|, \tag{3.23}$$

where

$$\theta_{i,k} \stackrel{\text{def}}{=} \kappa_\theta \min(1, h_{i,k}), \tag{3.24}$$

where $s_{i,k}$ are the inner iterates generated by the solver, κ_θ is any constant in $(0, 1)$, and

$$h_{i,k} \stackrel{\text{def}}{=} h_{i,k}(\|s_{i,k}\|, \|g_k\|)$$

are positive parameters. In particular, we are interested in two choices for $h_{i,k}$, namely,

$$h_{i,k} = \|s_{i,k}\|, \quad i \geq 0, \quad k \geq 0, \tag{3.25}$$

and

$$h_{i,k} = \|g_k\|^{1/2}, \quad i \geq 0, \quad k \geq 0. \tag{3.26}$$

The first choice gives improved complexity for the ARC algorithm (see [2]), while the second yields the best numerical performance of the algorithm in our experiments (see Sect. 7). Note that $g_k = \nabla_s m_k(0)$.

The condition (3.23) is always satisfied by any minimizer $s_{i,k}$ of m_k , since then $\nabla_s m_k(s_{i,k}) = 0$. Thus condition (3.23) can always be achieved by an iterative solver, the worst that could happen is to iterate until an exact minimizer of m_k is found. We hope in practice to terminate well before this inevitable outcome.

It follows from (3.23) and (3.24) that

$$\boxed{\text{TC.h}} \quad \|\nabla_s m_k(s_k)\| \leq \theta_k \|g_k\|, \quad \text{where } \theta_k = \kappa_\theta \min(1, h_k), \quad k \geq 0, \tag{3.27}$$

where $h_k \stackrel{\text{def}}{=} h_{i,k} > 0$ with i being the last inner iteration. In particular, for the choice (3.25), we have

$$\boxed{\text{TC.s}} \quad \|\nabla_s m_k(s_k)\| \leq \theta_k \|g_k\|, \quad \text{where } \theta_k = \kappa_\theta \min(1, \|s_k\|), \quad k \geq 0, \tag{3.28}$$

while for the choice (3.26), we obtain

$$\boxed{\text{TC.g}} \quad \|\nabla_s m_k(s_k)\| \leq \theta_k \|g_k\|, \quad \text{where } \theta_k = \kappa_\theta \min(1, \|g_k\|^{1/2}), \quad k \geq 0. \tag{3.29}$$

The lower bounds on s_k that the criteria TC.h, TC.s and TC.g provide are given in Lemmas 4.7 and 4.9.

4 Local convergence properties

For the remainder of the paper, we assume that

$$\boxed{\text{AF3}} \quad f \in C^2(\mathbb{R}^n). \tag{4.1}$$

Provided AF.3 holds, a Taylor expansion of $f(x_k + s_k)$ and its agreement with the model to first-order give the following estimate of the difference between the function

and the model at $x_k + s_k$, namely,

$$f(x_k + s_k) - m_k(s_k) = \frac{1}{2} s_k^\top [H(\xi_k) - B_k] s_k - \frac{\sigma_k}{3} \|s_k\|^3, \quad k \geq 0, \quad (4.2)$$

for some ξ_k on the line segment $(x_k, x_k + s_k)$. The expression (4.2) will be useful in some proofs.

4.1 Locally convex models

In this section, we investigate the convergence properties of the ARC algorithm in the case when the approximate Hessians B_k become positive definite asymptotically, at least along the direction s_k . Some results in this section follow closely those of Sect. 6.5 in [4].

Our main assumption in this section is that s_k satisfies (3.11). We remark that condition (3.12) is automatically achieved when B_k is positive semidefinite. Thus at present, we do not assume explicitly that s_k satisfies (3.12). Furthermore, no requirement of a termination criterion for the inner iterations is made (thus none of the definitions in Sect. 3.3 are employed in this section). Significantly, none of the results in this section requires the Hessian of the objective to be globally or locally Lipschitz continuous.

Let

$$R_k(s_k) \stackrel{\text{def}}{=} \frac{s_k^\top B_k s_k}{\|s_k\|^2}, \quad k \geq 0, \quad (4.3)$$

denote the Rayleigh quotient of s_k with respect to B_k , representing the curvature of the quadratic part of the model m_k along the step. We show that if (3.11) holds, we can guarantee stronger lower bounds on the model decrease than (3.19).

Lemma 4.1 *Let s_k satisfy (3.11). Then*

$$f(x_k) - m_k(s_k) \geq \frac{1}{2} R_k(s_k) \|s_k\|^2, \quad (4.4)$$

where $R_k(s_k)$ is the Rayleigh quotient (4.3). In particular,

$$f(x_k) - m_k(s_k) \geq \frac{1}{2} \lambda_{\min}(B_k) \|s_k\|^2, \quad (4.5)$$

where $\lambda_{\min}(B_k)$ denotes the leftmost eigenvalue of B_k .

Proof The bound (4.4) follows straightforwardly from (3.18) and (4.3), while for (4.5), we also employed the Rayleigh quotient inequality [4, p.19]. \square

When the Rayleigh quotient (4.3) is uniformly positive, the size of s_k is of order $\|g_k\|$, as we show next.

Lemma 4.2 *Suppose that s_k satisfies (3.11). If the Rayleigh quotient (4.3) is positive, then*

$$\|s_k\| \leq \frac{1}{R_k(s_k)} \|g_k\|. \tag{4.6}$$

Furthermore, if B_k is positive definite, then

$$\|s_k\| \leq \frac{1}{\lambda_{\min}(B_k)} \|g_k\|. \tag{4.7}$$

Proof The following relations are derived from (3.11) and the Cauchy-Schwarz inequality

$$R_k(s_k)\|s_k\|^2 \leq s_k^\top B_k s_k + \sigma_k \|s_k\|^3 = -g_k^\top s_k \leq \|g_k\| \cdot \|s_k\|.$$

The first and the last terms above give (4.7) since $s_k \neq 0$ because of (3.20), and $R_k(s_k) > 0$. The bound (4.7) follows from (4.6) and the Rayleigh quotient inequality. \square

The next theorem shows that all iterations are ultimately very successful provided some further assumption on the level of resemblance between the approximate Hessians B_k and the true Hessians $H(x_k)$ holds as the iterates converge to a local minimizer. In particular, we require

$$\boxed{\text{AM.2}} \quad \frac{\|(B_k - H(x_k))s_k\|}{\|s_k\|} \rightarrow 0, \quad \text{whenever } \|g_k\| \rightarrow 0. \tag{4.8}$$

The first limit in (4.8) is known as the Dennis–Moré condition [7]. A number of quasi-Newton techniques for updating B_k achieve this condition provided some additional properties hold [24, Sect. 3.3, Chapter 8]. For example, for Powell symmetric Broyden updates, it was shown in [18] that the Dennis–Moré condition is satisfied “on average” when the iterates converge to a stationary point x_* with $H(x_*)$ positive definite and locally Lipschitz continuous, provided also that the distances to the solution—or equivalently the steps—are square summable. It follows from (2.8) that the latter condition holds in our case over successful iterations.

Theorem 4.3 *Let AF.2–AF.3 and AM.1–AM.2 hold, and also let s_k satisfy (3.11), and*

$$x_k \rightarrow x_*, \quad \text{as } k \rightarrow \infty, \tag{4.9}$$

where $H(x_*)$ is positive definite. Then there exists $R_{\min} > 0$ such that

$$R_k(s_k) \geq R_{\min}, \quad \text{for all } k \text{ sufficiently large.} \tag{4.10}$$

Also, we have

$$\|s_k\| \leq \frac{1}{R_{\min}} \|g_k\|, \quad \text{for all } k \text{ sufficiently large.} \tag{4.11}$$

Furthermore, all iterations are eventually very successful, and σ_k is bounded from above.

Proof Since f is continuous, the limit (4.9) implies $\{f(x_k)\}$ is bounded below. Thus Corollary 2.6 provides that x_* is a first-order critical point and $\|g_k\| \rightarrow 0$. The latter limit and AM.2 imply

$$\frac{\|(H(x_k) - B_k)s_k\|}{\|s_k\|} \rightarrow 0, \quad k \rightarrow \infty, \tag{4.12}$$

i.e., the Dennis–Moré condition holds. Since $H(x_*)$ is positive definite, so is $H(x_k)$ for all k sufficiently large. In particular, there exists a constant R_{\min} such that

$$\frac{s_k^\top H(x_k)s_k}{\|s_k\|^2} > 2R_{\min} > 0, \quad \text{for all } k \text{ sufficiently large.} \tag{4.13}$$

From (4.3), (4.12) and (4.13), we obtain that for all sufficiently large k ,

$$2R_{\min}\|s_k\|^2 \leq s_k^\top H(x_k)s_k = s_k^\top [H(x_k) - B_k]s_k + s_k^\top B_k s_k \leq [R_{\min} + R_k(s_k)]\|s_k\|^2,$$

which gives (4.10). The bound (4.11) now follows from (4.6) and (4.10).

It follows from (2.41) that the equivalence (2.14) holds. We are going to derive an upper bound on the expression (2.15) of r_k and show that it is negative for all k sufficiently large. From (4.2), we have, also since $\sigma_k \geq 0$,

$$f(x_k + s_k) - m_k(s_k) \leq \frac{1}{2}\|(H(\xi_k) - B_k)s_k\| \cdot \|s_k\|, \tag{4.14}$$

where ξ_k belongs to the line segment $(x_k, x_k + s_k)$. Relation (4.4) in Lemma 4.1, and (4.10), imply

$$f(x_k) - m_k(s_k) \geq \frac{1}{2}R_{\min}\|s_k\|^2, \quad \text{for all } k \text{ sufficiently large.} \tag{4.15}$$

It follows from (2.15), (4.14) and (4.15) that

$$r_k \leq \frac{1}{2}\|s_k\|^2 \left\{ \frac{\|(H(\xi_k) - B_k)s_k\|}{\|s_k\|} - (1 - \eta_2)R_{\min} \right\}, \quad \text{for all } k \text{ sufficiently large.} \tag{4.16}$$

We have

$$\frac{\|(H(\xi_k) - B_k)s_k\|}{\|s_k\|} \leq \|H(x_k) - H(\xi_k)\| + \frac{\|(H(x_k) - B_k)s_k\|}{\|s_k\|}, \quad k \geq 0. \tag{4.17}$$

Since $\xi_k \in (x_k, x_k + s_k)$, we have $\|\xi_k - x_k\| \leq \|s_k\|$, which together with (4.11) and $\|g_k\| \rightarrow 0$, gives $\|\xi_k - x_k\| \rightarrow 0$. This, (4.9) and $H(x)$ continuous, give $\|H(x_k) -$

$H(\xi_k)\| \rightarrow 0$, as $k \rightarrow \infty$. It now follows from (4.12) and (4.17) that

$$\frac{\|(H(\xi_k) - B_k)s_k\|}{\|s_k\|} \rightarrow 0, \quad k \rightarrow \infty.$$

This, (3.20) and (4.16) imply $r_k < 0$, for all k sufficiently large. Since σ_k is not allowed to increase on the very successful steps of the ARC algorithm, and every k sufficiently large is very successful, σ_k is bounded from above. \square

The next two theorems address conditions under which the assumption (4.9) holds.

Theorem 4.4 *Suppose that AF2–AF.3, AM.1 and (3.11) hold, and that $\{f(x_k)\}$ is bounded below. Also, assume that $\{x_{k_i}\}$ is a subsequence of iterates converging to some x_* and that there exists $\underline{\lambda} > 0$ such that*

$$\lambda_{\min}(B_k) \geq \underline{\lambda}, \tag{4.18}$$

whenever x_k is sufficiently close to x_ . Let $H(x_*)$ be nonsingular. Then $x_k \rightarrow x_*$, as $k \rightarrow \infty$.*

Proof The conditions of Corollary 2.6 are satisfied, and thus $\|g_k\| \rightarrow 0, k \rightarrow \infty$. We deduce that $g(x_*) = 0$ and x_* is a first-order critical point. By employing (4.7) in Lemma 4.2, the proof now follows similarly to that of [4, Theorem 6.5.2]. \square

We remark that the sequence of iterates $\{x_k\}$ has a converging subsequence provided, for example, the level set of $f(x_0)$ is bounded.

The above theorem does not prevent the situation when the iterates converge to a critical point that is not a local minimizer. In the next theorem, besides assuming that x_* is a strict local minimizer, we require the approximate Hessians B_k to resemble the true Hessians $H(x_k)$ whenever the iterates approach a first-order critical point, namely,

AM.3

$$\|H(x_k) - B_k\| \rightarrow 0, \quad k \rightarrow \infty, \quad \text{whenever } \|g_k\| \rightarrow 0, \quad k \rightarrow \infty. \tag{4.19}$$

This condition is ensured, at least from a theoretical point of view, when B_k is set to the approximation of $H(x_k)$ computed by finite differences [8, 24]. It is also satisfied when using the symmetric rank one approximation to update B_k and the steps are linearly independent [1, 3].

Theorem 4.5 *Let AF2–AF.3, AM.1, AM.3 and (3.11) hold. Let also $\{f(x_k)\}$ be bounded below. Furthermore, suppose that $\{x_{k_i}\}$ is a subsequence of iterates converging to some x_* with $H(x_*)$ positive definite. Then the whole sequence of iterates $\{x_k\}$ converges to x_* , all iterations are eventually very successful, and σ_k stays bounded above.*

Proof Corollary 2.6 and f bounded below provide that x_* is a first-order critical point and $\|g_k\| \rightarrow 0$. The latter limit and AM.3 imply

$$\|H(x_k) - B_k\| \rightarrow 0, \quad k \rightarrow \infty. \tag{4.20}$$

Let $\{k_i\}$ index all the successful iterates x_{k_i} that converge to x_* (recall that the iterates remain constant on unsuccessful iterations). Since $H(x_*)$ is positive definite and $x_{k_i} \rightarrow x_*$, it follows from (4.20) that B_{k_i} is positive definite for all sufficiently large i , and thus there exists $\underline{\lambda} > 0$ such that (4.18) holds. Theorem 4.4 now provides that the whole sequence $\{x_k\}$ converges to x_* .

The conditions of Theorem 4.3 now hold since AM.3 implies AM.2. Thus the latter part of Theorem 4.5 follows from Theorem 4.3. \square

We remark that in the conditions of Theorem 4.5, B_k is positive definite asymptotically.

4.2 Asymptotic rate of convergence

In this section, the termination criteria in Sect. 3.3 are employed to show that the steps s_k do not become too small compared to the size of g_k (Lemmas 4.7 and 4.9), which then implies, in the context of Theorems 4.3 and 4.5, that the ARC algorithm is at least Q-superlinearly convergent (Corollaries 4.8 and 4.10).

Firstly, a technical result is deduced from the termination criterion TC.h, which requires that g is Lipschitz continuous on an open convex set X containing all the iterates $\{x_k\}$, namely,

$$\boxed{\text{AF.4}} \quad \|g(x) - g(y)\| \leq \kappa_H \|x - y\|, \quad \text{for all } x, y \in X, \text{ and some } \kappa_H \geq 0. \tag{4.21}$$

If AF.3 holds, AF.4 is satisfied if the Hessian $H(x)$ is bounded above on X . Note that AF.4 implies AF.2.

Lemma 4.6 *Let AF.3–AF.4 and TC.h hold. Then, for each $k \in \mathcal{S}$, with \mathcal{S} defined in (2.7), we have*

$$\begin{aligned} (1 - \kappa_\theta) \|g_{k+1}\| &\leq \left\| \int_0^1 H(x_k + \tau s_k) d\tau - H(x_k) \right\| \cdot \|s_k\| + \|(H(x_k) - B_k)s_k\| \\ &\quad + \kappa_\theta \kappa_H h_k \|s_k\| + \sigma_k \|s_k\|^2, \end{aligned} \tag{4.22}$$

where $\kappa_\theta \in (0, 1)$ occurs in TC.h.

Proof Let $k \in \mathcal{S}$, and so $g_{k+1} = g(x_k + s_k)$. Then

$$\begin{aligned} \|g_{k+1}\| &\leq \|g(x_k + s_k) - \nabla_s m_k(s_k)\| + \|\nabla_s m_k(s_k)\| \\ &\leq \|g(x_k + s_k) - \nabla_s m_k(s_k)\| + \theta_k \|g_k\|, \end{aligned} \tag{4.23}$$

where we used TC.h to derive the last inequality. We also have from Taylor’s theorem and (3.1)

$$\|g(x_k + s_k) - \nabla_s m_k(s_k)\| \leq \left\| \int_0^1 [H(x_k + \tau s_k) - B_k] s_k d\tau \right\| + \sigma_k \|s_k\|^2. \tag{4.24}$$

From AF.4 and the triangle inequality, we obtain

$$\|g_k\| \leq \|g_{k+1}\| + \|g_{k+1} - g_k\| \leq \|g_{k+1}\| + \kappa_H \|s_k\|. \tag{4.25}$$

Substituting (4.25) and (4.24) into (4.23), we deduce

$$(1 - \theta_k) \|g_{k+1}\| \leq \left\| \int_0^1 [H(x_k + \tau s_k) - B_k] s_k d\tau \right\| + \theta_k \kappa_H \|s_k\| + \sigma_k \|s_k\|^2. \tag{4.26}$$

It follows from the definition of θ_k in (3.27) that $\theta_k \leq \kappa_\theta h_k$ and $\theta_k \leq \kappa_\theta$, and (4.26) becomes

$$(1 - \kappa_\theta) \|g_{k+1}\| \leq \left\| \int_0^1 [H(x_k + \tau s_k) - B_k] s_k d\tau \right\| + \kappa_\theta \kappa_H h_k \|s_k\| + \sigma_k \|s_k\|^2. \tag{4.27}$$

The triangle inequality provides

$$\left\| \int_0^1 [H(x_k + \tau s_k) - B_k] s_k d\tau \right\| \leq \left\| \int_0^1 H(x_k + \tau s_k) d\tau - H(x_k) \right\| \cdot \|s_k\| + \|(H(x_k) - B_k) s_k\|, \tag{4.28}$$

and so (4.22) follows from (4.27). □

The next lemma establishes conditions under which the TC.h criterion provides a lower bound on s_k .

Lemma 4.7 *Let AF.3–AF.4, AM.2 and the limit $x_k \rightarrow x_*$, $k \rightarrow \infty$, hold. Let TC.h be achieved with*

$$h_k \rightarrow 0, \text{ as } k \rightarrow \infty, k \in \mathcal{S}. \tag{4.29}$$

Then s_k satisfies

$$\|s_k\| (d_k + \sigma_k \|s_k\|) \geq (1 - \kappa_\theta) \|g_{k+1}\| \text{ for all } k \in \mathcal{S}, \tag{4.30}$$

where $d_k > 0$ for all $k \geq 0$, and

$$d_k \rightarrow 0, \text{ as } k \rightarrow \infty, k \in \mathcal{S}. \tag{4.31}$$

Proof The inequality (4.22) can be expressed as

$$(1 - \kappa_\theta) \|g_{k+1}\| \leq \left[\left\| \int_0^1 [H(x_k + \tau s_k) - H(x_k)] d\tau \right\| + \frac{\|(H(x_k) - B_k)s_k\|}{\|s_k\|} + \kappa_\theta \kappa_H h_k \right] \cdot \|s_k\| + \sigma_k \|s_k\|^2,$$

where $k \in \mathcal{S}$. Let d_k denote the term in the curly brackets multiplying $\|s_k\|$. Then $d_k > 0$ since $h_k > 0$. Furthermore, $x_k + \tau s_k \in (x_k, x_{k+1})$, for all $\tau \in (0, 1)$, and $x_k \rightarrow x_*$, imply

$$\left\| \int_0^1 [H(x_k + \tau s_k) - H(x_k)] d\tau \right\| \rightarrow 0, \text{ as } k \rightarrow \infty, \tag{4.32}$$

since the Hessian of f is continuous. Since AF.4 implies AF.2, Corollary 2.6 provides $\|g_k\| \rightarrow 0$. It now follows from AM.2, (4.29) and (4.32) that $d_k \rightarrow 0$, as the index k of successful iterations increases. \square

By employing Lemma 4.7 in the context of Theorem 4.3, we show that the ARC algorithm is asymptotically Q-superlinearly convergent.

Corollary 4.8 *In addition to the conditions of Theorem 4.3, assume that AF.4 holds and TC.h is satisfied with $h_k \rightarrow 0, k \rightarrow \infty, k \in \mathcal{S}$. Then*

$$\frac{\|g_{k+1}\|}{\|g_k\|} \rightarrow 0, \text{ as } k \rightarrow \infty, \tag{4.33}$$

and

$$\frac{\|x_{k+1} - x_*\|}{\|x_k - x_*\|} \rightarrow 0, \text{ as } k \rightarrow \infty. \tag{4.34}$$

In particular, the limits (4.33) and (4.34) hold when $h_k = \|s_k\|, k \geq 0$, or $h_k = \|g_k\|^{1/2}, k \geq 0$, i.e., in the case of the termination criteria TC.s and TC.g, respectively.

Proof Note that AF.4 implies that AF.2 holds in the conditions of Theorem 4.3. Since the conditions of Lemma 4.7 hold, so does the bound (4.30). Moreover, as Theorem 4.3 gives that all iterates are eventually very successful and σ_k is bounded above, say by some σ_{sup} , (4.30) holds for all k sufficiently large and thus

$$\|s_k\|(d_k + \sigma_{sup}\|s_k\|) \geq (1 - \kappa_\theta)\|g_{k+1}\| \tag{4.35}$$

for all k sufficiently large, where $d_k > 0$ and $\kappa_\theta \in (0, 1)$. Employing the upper bound (4.11) on s_k , (4.35) becomes

$$\frac{1}{R_{\min}} \left(d_k + \frac{\sigma_{sup}}{R_{\min}} \|g_k\| \right) \|g_k\| \geq \|s_k\|(d_k + \sigma_{sup}\|s_k\|) \geq (1 - \kappa_\theta)\|g_{k+1}\|, \tag{4.36}$$

for all k sufficiently large, and further, because of (2.40),

$$\frac{\|g_{k+1}\|}{\|g_k\|} \leq \frac{R_{\min}d_k + \sigma_{sup}\|g_k\|}{R_{\min}^2(1 - \kappa_\theta)}, \tag{4.37}$$

for all k sufficiently large. The right-hand side of (4.37) tends to zero as $k \rightarrow \infty$, since

$$d_k \rightarrow 0 \quad \text{and} \quad \|g_k\| \rightarrow 0, \quad \text{as } k \rightarrow \infty; \tag{4.38}$$

the first limit above comes from (4.31) and all k sufficiently large being successful, while the second limit follows from Corollary 2.6. Thus (4.33) holds. The limit (4.34) is obtained from standard Taylor expansions of g_k and g_{k+1} around x_* , and from $g(x_*) = 0$ with positive definite $H(x_*)$.

The bound (4.11) and the second limit in (4.38) imply that the choices of h_k in TC.s and TC.g converge to zero, and thus the limits (4.33) and (4.34) hold for these choices of h_k . □

Note that the limits (4.33) and (4.34) also hold if we let $h_k = \|s_k\|/\sigma_k, k \geq 0$, in TC.h, provided the conditions of Theorem 4.3 hold (since then, σ_k is bounded above asymptotically). See (7.3) in Sect. 7.

Note also that no assumption on the Hessian of f being globally or locally Lipschitz continuous has been imposed in Lemma 4.7 or in Corollary 4.8. Our next results, however, make a *local* Lipschitz continuity assumption on the Hessian of f in a neighbourhood of a given point x_* , i.e.,

AF.5

$$\|H(x) - H(y)\| \leq L_*\|x - y\|, \quad \text{for all } x, y \text{ sufficiently close to } x_*,$$

and some $L_* > 0$,

and show a tighter bound on s_k than (4.30) (see Lemma 4.9), and further, Q-quadratic asymptotic convergence of the iterates (Corollary 4.10). In this context, we also slightly strengthen the condition AM.2, by requiring that B_k satisfies

AM.4

$$\|(H(x_k) - B_k)s_k\| \leq C\|s_k\|^2, \text{ for all } k \geq 0, \text{ and some constant } C > 0. \tag{4.39}$$

We remark that if the inequality in AM.4 holds for sufficiently large k , it also holds for all $k \geq 0$. The condition AM.4 is trivially satisfied with $C = 0$ when we set $B_k = H(x_k)$ for all $k \geq 0$. Quasi-Newton methods may still satisfy AM.4 in practice, though theoretically, only condition AM.2 can be ensured.

Lemma 4.9 *Let AF.3–AF.5, AM.4 and TC.s hold. Suppose also that $x_k \rightarrow x_*$, as $k \rightarrow \infty$. If*

$$\sigma_k \leq \sigma_{\max}, \quad \text{for all } k \geq 0, \tag{4.40}$$

for some $\sigma_{\max} > 0$, then s_k satisfies

$$\|s_k\| \geq \kappa_g^* \sqrt{\|g_{k+1}\|} \quad \text{for all sufficiently large } k \in \mathcal{S}, \tag{4.41}$$

where κ_g^* is the positive constant

$$\kappa_g^* \stackrel{\text{def}}{=} \sqrt{\frac{1 - \kappa_\theta}{\frac{1}{2}L_* + C + \sigma_{\max} + \kappa_\theta \kappa_H}}. \tag{4.42}$$

Proof The conditions of Lemma 4.6 are satisfied with $h_k = \|s_k\|$. Thus, for any $k \in \mathcal{S}$ sufficiently large, (4.22) becomes, due also to AM.4 and (4.40),

$$(1 - \kappa_\theta)\|g_{k+1}\| \leq \left\| \int_0^1 [H(x_k + \tau s_k) - H(x_k)]d\tau \right\| \cdot \|s_k\| + C\|s_k\|^2 + (\sigma_{\max} + \kappa_\theta \kappa_H)\|s_k\|^2. \tag{4.43}$$

Since $x_k \rightarrow x_*$, AF.5 and $x_k + \tau s_k$ being on the line segment (x_k, x_{k+1}) for any $\tau \in (0, 1)$, imply

$$\left\| \int_0^1 [H(x_k + \tau s_k) - H(x_k)]d\tau \right\| \leq \int_0^1 \|H(x_k + \tau s_k) - H(x_k)\|d\tau \leq \frac{1}{2}L_*\|s_k\|, \tag{4.44}$$

for all sufficiently large $k \in \mathcal{S}$. Thus (4.43) becomes

$$(1 - \kappa_\theta)\|g_{k+1}\| \leq \left(\frac{1}{2}L_* + C + \sigma_{\max} + \kappa_\theta \kappa_H\right) \|s_k\|^2, \tag{4.45}$$

which together with (4.42) provides (4.41). □

Our next result employs Lemma 4.9 to show Q-quadratic asymptotic convergence of the ARC algorithm.

Corollary 4.10 *In addition to the conditions of Theorem 4.3, assume that AF.4–AF.5, AM.4 and TC.s hold. Then g_k converges to zero, and x_k , to x_* , Q-quadratically, as $k \rightarrow \infty$.*

Proof Note that AF.4 implies that AF.2 holds in the conditions of Theorem 4.3. Also, AM.4 implies AM.2, since (4.10) and $\|g_k\| \rightarrow 0$ give $\|s_k\| \rightarrow 0$, as $k \rightarrow \infty$. Theorem 4.3 implies that σ_k is bounded above and thus (4.40) holds. Recalling that all the iterates are eventually very successful, Lemma 4.9 now implies that

$$\|s_k\| \geq \kappa_g^* \sqrt{\|g_{k+1}\|}, \quad \text{for all } k \text{ sufficiently large,} \tag{4.46}$$

where $\kappa_g^* > 0$. It follows from (4.10) and (4.46) that

$$\frac{1}{R_{\min}} \|g_k\| \geq \|s_k\| \geq \kappa_g^* \sqrt{\|g_{k+1}\|}, \quad \text{for all } k \text{ sufficiently large,} \tag{4.47}$$

and thus

$$\frac{\|g_{k+1}\|}{\|g_k\|^2} \leq \frac{1}{R_{\min}^2 (\kappa_g^*)^2}, \quad \text{for all } k \text{ sufficiently large,} \tag{4.48}$$

and g_k converges Q-quadratically. The Q-quadratic rate of convergence of the iterates follows in a standard way, using Taylor’s theorem. □

Analoguees of Corollaries 4.8 and 4.10 hold in the case when the stronger conditions of Theorem 4.5 are satisfied. In particular, we require the stronger condition AM.3, instead of AM.2 or AM.4, to be achieved by B_k ; then, the limit $x_k \rightarrow x_*$ is guaranteed to hold. The weaker assumption AM.2, however, makes Corollary 4.8 applicable to quasi-Newton methods (see our remarks following (4.8)).

Note that no positive lower bound on σ_k was required for the convergence results in Sects. 2.2, 4.1 and 4.2 to hold. In particular, asymptotically, it may be desirable in implementations to let σ_k to go to zero, possibly at the same rate as $\|g_k\|$. This feature is included in our implementation of the ARC algorithm (see Sect. 7).

We remark that assumption AF.4 was only employed in Sect. 4.2 and will not be used again for the remainder of the paper (Part I). Thus note that (4.21) could have been replaced by the weaker condition that g is Lipschitz continuous only on the iterates $\{x_k\}$; due to connections to Part II [2], we prefer the formulation in (4.21).

5 Global convergence to second-order critical points

This section addresses the convergence of the sequence of iterates to second-order critical points in a framework that does not require global or local convexity of the model or the function f at the iterates or their limit points. Then, however, we shall see that other conditions such as $H(x)$ being globally Lipschitz continuous, need to be imposed. A common assumption in this section is that

$$\sigma_k \geq \sigma_{\min}, \quad k \geq 0, \tag{5.1}$$

for some $\sigma_{\min} > 0$. The first lemma gives a useful property of the steps s_k , derived from (3.19).

Lemma 5.1 *Let $\{f(x_k)\}$ be bounded below by f_{low} . Also, assume that s_k satisfies (3.11) and (3.12), and σ_k , the bound (5.1). Then, recalling (2.7), we have*

$$\|s_k\| \rightarrow 0, \quad \text{as } k \rightarrow \infty, \quad k \in \mathcal{S}. \tag{5.2}$$

Proof The construction of the ARC algorithm, (2.41), the model decrease (3.19) and (5.1) give

$$f(x_k) - f(x_{k+1}) \geq \eta_1[f(x_k) - m_k(s_k)] \geq \frac{1}{6}\eta_1\sigma_{\min}\|s_k\|^3, \quad k \in \mathcal{S}. \quad (5.3)$$

Summing up over all iterates from 0 to k , we obtain from (5.3)

$$f(x_0) - f(x_{k+1}) \geq \frac{\eta_1}{6}\sigma_{\min} \sum_{j=0, j \in \mathcal{S}}^k \|s_j\|^3, \quad k \geq 0,$$

which further gives, together with $\{f(x_k)\}$ being bounded below,

$$\frac{6}{\eta_1\sigma_{\min}}[f(x_0) - f_{\text{low}}] \geq \sum_{j=0, j \in \mathcal{S}}^k \|s_j\|^3, \quad k \geq 0. \quad (5.4)$$

Thus the series $\sum_{j=0, j \in \mathcal{S}}^{\infty} \|s_j\|^3$ is convergent, and (5.2) holds. \square

The next lemma shows that σ_k cannot blow up provided the objective f has a globally Lipschitz continuous Hessian, namely,

$$\boxed{\text{AF.6}} \quad \|H(x) - H(y)\| \leq L\|x - y\|, \quad \text{for all } x, y \in \mathbb{R}^n, \text{ where } L > 0, \quad (5.5)$$

and B_k and $H(x_k)$ agree along s_k in the sense of AM.4.

Lemma 5.2 *Let AF.3, AF.6 and AM.4 hold. Then*

$$\sigma_k \leq \max\left(\sigma_0, \frac{3}{2}\gamma_2(C + L)\right) \stackrel{\text{def}}{=} L_0, \quad \text{for all } k \geq 0. \quad (5.6)$$

Proof Let $L_1 \stackrel{\text{def}}{=} 3(C + L)/2$. To prove (5.6), it is enough to show the implication

$$\sigma_k \geq L_1 \implies k \text{ very successful}, \quad (5.7)$$

which further gives $\sigma_{k+1} \leq \sigma_k$. We allow the factor γ_2 in L_0 for the case when σ_k is only slightly less than L_1 and k is not very successful, while the term σ_0 in (5.6) accounts for choices at start-up.

To show (5.7), we deduce from (4.2) that for each $k \geq 0$,

$$\begin{aligned} f(x_k + s_k) - m_k(s_k) &\leq \frac{1}{2}\|H(\xi_k) - H(x_k)\| \cdot \|s_k\|^2 + \frac{1}{2}\|(H(x_k) - B_k)s_k\| \cdot \|s_k\| \\ &\quad - \frac{\sigma_k}{3}\|s_k\|^3 \leq \left(\frac{C + L}{2} - \frac{\sigma_k}{3}\right)\|s_k\|^3, \end{aligned} \quad (5.8)$$

where to obtain the second inequality, we employed AF.6, $\|\xi_k - x_k\| \leq \|s_k\|$ and AM.4. It follows from (5.8) that

$$\sigma_k \geq L_1 \implies f(x_k + s_k) \leq m_k(s_k). \tag{5.9}$$

The second inequality in (5.9) and (2.41) imply that the ratio (2.4) satisfies $\rho_k \geq 1$ and so $\rho_k > \eta_2$, for any $\eta_2 \in (0, 1)$. Thus the step k is very successful. \square

Next, we show that all the limit points of the sequence of Rayleigh quotients of B_k and of $H(x_k)$ at successful steps s_k are nonnegative.

Theorem 5.3 *Let AF.3, AF.6 and AM.4 hold, and $\{f(x_k)\}$ be bounded below by f_{low} . Also, assume that s_k satisfies (3.11) and (3.12), and σ_k , (5.1). Then, recalling (4.3), we have*

$$\liminf_{k \in \mathcal{S}, k \rightarrow \infty} R_k(s_k) \geq 0 \quad \text{and} \quad \liminf_{k \in \mathcal{S}, k \rightarrow \infty} \frac{s_k^\top H(x_k)s_k}{\|s_k\|^2} \geq 0. \tag{5.10}$$

Proof For all $k \geq 0$ such that $R_k(s_k) < 0$, (3.12), (3.20) and (5.6) imply

$$L_0 \|s_k\| \geq \sigma_k \|s_k\| \geq -R_k(s_k) = |R_k(s_k)|.$$

Now (5.2) implies the limit $R_k(s_k) \rightarrow 0, k \in \mathcal{S}$ and $k \rightarrow \infty$, for all $R_k(s_k) < 0$. This proves the first limit in (5.10). The second inequality in (5.10) now follows from the first, (5.2) and the inequalities

$$R_k(s_k) \leq \frac{\|[H(x_k) - B_k]s_k\|}{\|s_k\|} + \frac{s_k^\top H(x_k)s_k}{\|s_k\|^2} \leq C \|s_k\| + \frac{s_k^\top H(x_k)s_k}{\|s_k\|^2}, \quad k \geq 0,$$

where we employed AM.4 to obtain the last inequality. \square

The next theorem gives conditions which ensure that the limit points of the sequence of iterates are second order critical points. Beforehand, we remark a useful property concerning the Hessian $H(x_k)$ and its approximation B_k . Given any matrix Q_k with orthogonal columns, [13, Corollary 8.1.6] provides the first inequality below

$$\begin{aligned} |\lambda_{\min}(Q_k^\top H(x_k) Q_k) - \lambda_{\min}(Q_k^\top B_k Q_k)| &\leq \|Q_k^\top [H(x_k) - B_k] Q_k\| \\ &\leq \sqrt{n} \|H(x_k) - B_k\|, \quad k \geq 0, \end{aligned} \tag{5.11}$$

while the second inequality above employs $\|Q_k^\top\| \leq \sqrt{n}$ and $\|Q_k\| = 1$.

Theorem 5.4 *Let AF.3, AF.6 and AM.4 hold. Assume that $\{f(x_k)\}$ is bounded below by f_{low} , and that σ_k satisfies (5.1). Also, let s_k be a global minimizer of m_k over a subspace \mathcal{L}_k , and let Q_k be any orthogonal matrix whose columns form a basis of \mathcal{L}_k . Then any subsequence of negative leftmost eigenvalues $\{\lambda_{\min}(Q_k^\top B_k Q_k)\}$ converges to zero as $k \rightarrow \infty, k \in \mathcal{S}$, and thus*

$$\liminf_{k \in \mathcal{S}, k \rightarrow \infty} \lambda_{\min}(Q_k^\top B_k Q_k) \geq 0. \tag{5.12}$$

Additionally, assume that AF.2, AM.1 and AM.3 hold. Then any subsequence of negative leftmost eigenvalues $\{\lambda_{\min}(Q_k^\top H(x_k) Q_k)\}$ converges to zero as $k \rightarrow \infty, k \in \mathcal{S}$, and thus

$$\liminf_{k \in \mathcal{S}, k \rightarrow \infty} \lambda_{\min}(Q_k^\top H(x_k) Q_k) \geq 0. \tag{5.13}$$

Furthermore, if Q_k becomes a full orthogonal basis of \mathbb{R}^n as $k \rightarrow \infty, k \in \mathcal{S}$, then, provided it exists, any limit point of the sequence of iterates $\{x_k\}$ is second-order critical.

Proof For all $k \geq 0$ such that $\lambda_{\min}(Q_k^\top B_k Q_k) < 0$, we employ Lemma 3.2, in particular, (3.13), and also (5.6), to obtain

$$L_0 \|s_k\| \geq \sigma_k \|s_k\| \geq -\lambda_{\min}(Q_k^\top B_k Q_k) = |\lambda_{\min}(Q_k^\top B_k Q_k)|.$$

Now (5.2) implies the limit

$$\lambda_{\min}(Q_k^\top B_k Q_k) \rightarrow 0, k \in \mathcal{S} \text{ and } k \rightarrow \infty, \text{ for all } \lambda_{\min}(Q_k^\top B_k Q_k) < 0, \tag{5.14}$$

which gives (5.12).

Assume now that AF.2, AM.1 and AM.3 holds. Then AF.2, AM.1, $\{f(x_k)\}$ bounded below, (2.40) and Corollary 2.6 give that

$$\|g_k\| \rightarrow 0, \quad k \rightarrow \infty, \tag{5.15}$$

so that the first limit in AM.3 holds, i.e.,

$$\|H(x_k) - B_k\| \rightarrow 0, \quad k \rightarrow \infty. \tag{5.16}$$

We deduce from (5.11) that for all $k \geq 0$ with $\lambda_{\min}(Q_k^\top H(x_k) Q_k) < 0$, we have

$$\begin{aligned} 0 \leq -\lambda_{\min}(Q_k^\top H(x_k) Q_k) &\leq \sqrt{n} \|H(x_k) - B_k\| - \lambda_{\min}(Q_k^\top B_k Q_k) \\ &\leq \sqrt{n} \|H(x_k) - B_k\| + \max(0, -\lambda_{\min}(Q_k^\top B_k Q_k)). \end{aligned} \tag{5.17}$$

It follows from (5.14) and (5.16) that the right-hand side of (5.17) converges to zero as $k \in \mathcal{S} \rightarrow \infty$, and so $\lambda_{\min}(Q_k^\top H(x_k) Q_k) \rightarrow 0, k \in \mathcal{S} \rightarrow \infty$ with $\lambda_{\min}(Q_k^\top H(x_k) Q_k) < 0$. This gives (5.13).

Assume now that there exists x_* such that $x_k \rightarrow x_*, k \in \mathcal{K}$ and $k \rightarrow \infty$. Then (5.15) and the arguments that give (5.15) imply $\|g(x_*)\| = 0$ and $\mathcal{K} \subseteq \mathcal{S}$, where the set inclusion also uses the fact that the iterates remain constant on unsuccessful iterations. Employing AF.3, we have $H(x_k) \rightarrow H(x_*), k \in \mathcal{K}, k \rightarrow \infty$. Since the set of orthogonal matrices is compact and Q_k becomes full-dimensional as $k \rightarrow \infty, k \in \mathcal{K}$, any limit point, say Q , of $\{Q_k\}_{k \in \mathcal{K}}$ is a full orthogonal basis of \mathbb{R}^n . Due to similarity and from (5.13), we now have $\lambda_{\min}(H(x_*)) = \lambda_{\min}(Q^\top H(x_*) Q) \geq 0$, and so x_* is second-order critical. □

In our implementation of the ARC algorithm, though we minimize m_k only in certain subspaces, our particular approach (see Sect. 6.2) implies that ever more accurate Ritz approximations to the extreme eigenvalues of B_k are computed provided the gradient is not orthogonal to any eigenvector of B_k [14]. In other words, except for the latter case, we expect that the orthogonal bases of the generated subspaces become full-dimensional asymptotically, and so Theorem 5.4 implies that the solutions we obtain will be second-order critical in the limit.

When $Q_k = I$ and $B_k = H(x_k)$ for all $k \geq 0$, the last part of Theorem 5.4 is essentially [23, Theorem 2].

Encountering zero gradient values. Recall the discussion in the last paragraph of Sect. 3.2, where we assume that there exists $k \geq 0$ such that $g_k = 0$ and thus (2.40) does not hold. Then (3.21) provides $s_k \neq 0$ and (2.41) holds. These two relations imply that Lemmas 5.1 and 5.2 continue to hold even when some of the iterates have zero gradients (and the ARC algorithm continues to iterate to try to attain second-order criticality in the subspaces). Employing these Lemmas and the conditions of Corollary 5.4, the limit (5.12) can be shown as before since the value of the gradient was irrelevant in its derivation. To ensure (5.13), we further assume, in addition to the requirements of Corollary 5.4, that

$$B_k = H_k \quad \text{for all } k \text{ for which } g_k = 0. \quad (5.18)$$

The proof of (5.13) follows as before, except that if there are infinitely many k_l such that

$$g_{k_l} = 0 \quad \text{and} \quad \lambda_{\min}(Q_{k_l}^\top H(x_{k_l}) Q_{k_l}) < 0, \quad (5.19)$$

then (5.13) and (5.18) give $\liminf_{k_l \rightarrow \infty, k_l \in \mathcal{S}} \lambda_{\min}(Q_{k_l}^\top H(x_{k_l}) Q_{k_l}) \geq 0$. Note that the ARC algorithm ultimately moves away from iterates satisfying (5.19): since σ_k is bounded above as in (5.6), the ARC algorithm cannot take an infinite number of unsuccessful steps at x_{k_l} (when σ_k is increased by a fixed fraction).

The last statement of Corollary 5.4 also holds in this case provided Q_k is full-dimensional also when $g_k = 0$.

6 Methods for approximately minimizing the cubic model

While the ARC algorithm provides a powerful general framework for unconstrained minimization, the practicality and efficiency of this algorithm is obviously related to our ability to find a suitable (approximate) minimizer of the cubic model at each iteration. In this section we consider this issue in some detail. The optimality conditions in Theorem 3.1 for the global minimizer of $m_k(s)$ over \mathbb{R}^n are highly suggestive of efficient algorithms in many cases, as we discuss in the first subsection. We then concentrate on one way in which this minimizer may be approximated in practice, while retaining both the convergence and complexity properties of the true model minimizer. Most especially, the method we propose is “matrix-free”—that is, it only

requires Hessian-vector products rather than access to the Hessian itself—and thus may be used in principle for large, unstructured problems.

Throughout this section, we drop the (major) iteration subscript k for convenience.

6.1 Computing the global solution

To start with, we suppose that we wish to compute the global model minimizer of $m(s)$ over \mathbb{R}^n . Theorem 3.1 shows that we seek a pair (s, λ) for which

$$(B + \lambda I)s = -g \quad \text{and} \quad \lambda^2 = \sigma^2 s^T s \quad (6.1)$$

and for which $B + \lambda I$ is positive semidefinite. Just as in the trust-region case² [4, Sect. 7.3.1], suppose that B has an eigendecomposition

$$B = U^T \Lambda U, \quad (6.2)$$

where Λ is a diagonal matrix of eigenvalues $\lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$ and U is an orthonormal matrix of associated eigenvectors. Then

$$B(\lambda) = U^T (\Lambda + \lambda I) U.$$

We deduce immediately from Theorem 3.1 that the value of λ we seek must satisfy $\lambda \geq -\lambda_1$ as only then is $B(\lambda)$ positive semidefinite.

Suppose that $\lambda > -\lambda_1$. Then $B + \lambda I$ is positive definite, and let

$$s(\lambda) = -(B + \lambda I)^{-1} g = -U^T (\Lambda + \lambda I)^{-1} U g. \quad (6.3)$$

But, of course, the solution we are looking for depends upon the nonlinear equality $\|s(\lambda)\| = \lambda/\sigma$. To say more, we need to examine $\|s(\lambda)\|_2$ in detail. For convenience, define $\psi(\lambda) \stackrel{\text{def}}{=} \|s(\lambda)\|^2$. We have that

$$\psi(\lambda) = \|U^T (\Lambda + \lambda I)^{-1} U g\|^2 = \|(\Lambda + \lambda I)^{-1} U g\|^2 = \sum_{i=1}^n \frac{\gamma_i^2}{(\lambda_i + \lambda)^2}, \quad (6.4)$$

where γ_i is the i th component of Ug .

If B is positive semidefinite, the required solution is given by the single positive root to either of the equivalent univariate nonlinear equations

$$\theta_2(\lambda) \stackrel{\text{def}}{=} \psi(\lambda) - \frac{\lambda^2}{\sigma^2} = 0 \quad \text{or} \quad \theta_1(\lambda) \stackrel{\text{def}}{=} \sqrt{\psi(\lambda)} - \frac{\lambda}{\sigma} = 0, \quad (6.5)$$

² Griewank also comments in [17] on the similarity between the two problems, but does not use it in practice.

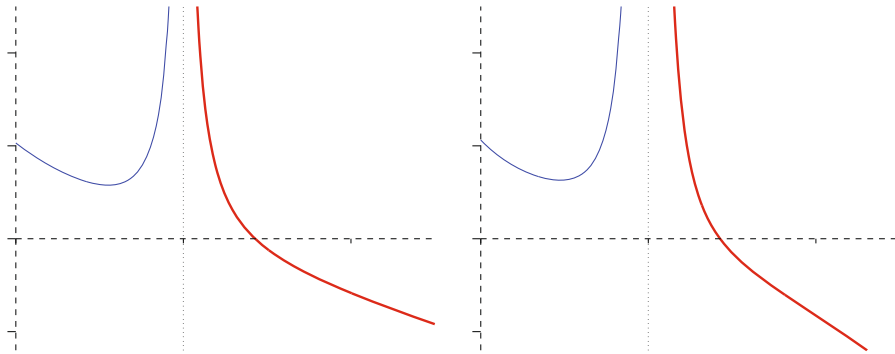


Fig. 1 Graphs of the functions $\theta_1(\lambda)$ (left) and $\theta_2(\lambda)$ (right) from (6.5) when $g = (0.25 \ 1)^T$, $H = \text{diag}(-1 \ 1)$ and $\sigma = 2$

as the left-hand sides in (6.5) are strictly decreasing functions of λ , for $\lambda \geq \max(0, -\lambda_1) = 0$, and range between some positive value and $-\infty$. If B is indefinite and $\gamma_1 \neq 0$, the solution is likewise the root larger than $-\lambda_1$ of the same equations. Hence in both these cases, the model $m(s)$ has a unique global minimizer. But if B is indefinite and $\gamma_1 = 0$, we have difficulties analogous to those for the hard case [4, Sect. 7.3.1.3] for the trust-region subproblem in which the required solution s_* is made up as a suitable linear combination of u_1 and $\lim_{s \rightarrow -\lambda_1} s(\lambda)$. To determine the values of the coefficients of this linear combination, in place of the trust-region constraint, we employ (6.1), and find a value for $\alpha \in \mathbb{R}$ such that

$$-\lambda_1 = \sigma \|s(-\lambda_1) + \alpha u_1\|. \tag{6.6}$$

See Fig. 1 for an illustration of the graphs of θ_1 and θ_2 .

In practice, just as in the trust-region case, it may be preferable to solve one of

$$\begin{aligned} \phi_2(\lambda) \stackrel{\text{def}}{=} \frac{1}{\psi(\lambda)} - \frac{\sigma^2}{\lambda^2} = 0, \quad \phi_1(\lambda) \stackrel{\text{def}}{=} \frac{1}{\sqrt{\psi(\lambda)}} - \frac{\sigma}{\lambda} = 0, \\ \beta_2(\lambda) \stackrel{\text{def}}{=} \frac{\lambda^2}{\psi(\lambda)} - \sigma^2 = 0 \quad \text{or} \quad \beta_1(\lambda) \stackrel{\text{def}}{=} \frac{\lambda}{\sqrt{\psi(\lambda)}} - \sigma = 0 \end{aligned} \tag{6.7}$$

instead of (6.5). Figures 2 and 3 illustrate these alternatives.

In any event, a safeguarded univariate Newton iteration to find the required root of whichever of the functions (6.5) or (6.7) we prefer is recommended, but in all cases this requires the solution of a sequence of linear equations

$$B(\lambda^{(k)})s^{(k)} \equiv (B + \lambda^{(k)}I)s^{(k)} = -g$$

for selected $\lambda^{(k)} > \max(0, -\lambda_1)$. Clearly to use Newton’s method, derivatives of (simple functions of) $\psi(\lambda)$ will be required, but fortunately these are easily obtained once a factorization of $B + \lambda^{(k)}I$ is known. In particular, we have the result below.

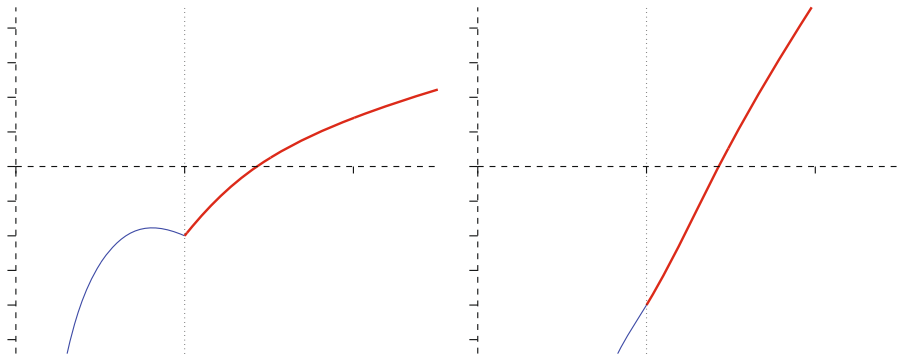


Fig. 2 Graphs of the functions $\phi_1(\lambda)$ (left) and $\phi_2(\lambda)$ (right) from (6.7) when $g = (0.25 \ 1)^T$, $H = \text{diag}(-1 \ 1)$ and $\sigma = 2$

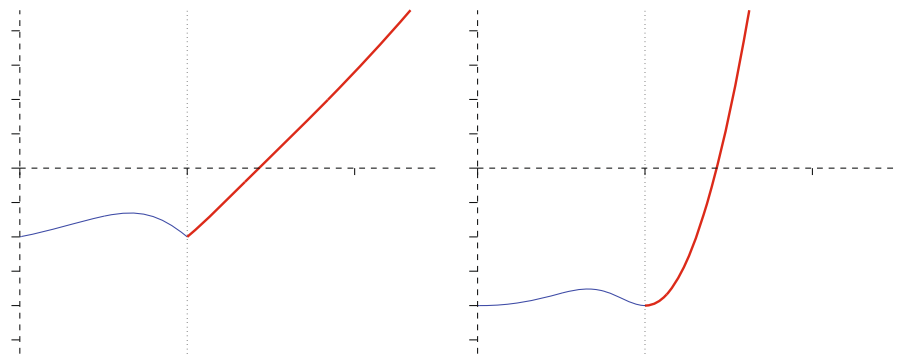


Fig. 3 Graphs of the functions $\beta_1(\lambda)$ (left) and $\beta_2(\lambda)$ (right) from (6.7) when $g = (0.25 \ 1)^T$, $H = \text{diag}(-1 \ 1)$ and $\sigma = 2$

Lemma 6.1 Suppose that $s(\lambda)$ satisfies (6.3), $\psi(\lambda) = \|s(\lambda)\|^2$ and $\lambda > -\lambda_1$. Then

$$\psi'(\lambda) = 2s(\lambda)^T \nabla_\lambda s(\lambda) \quad \text{and} \quad \psi''(\lambda) = 6\|\nabla_\lambda s(\lambda)\|^2, \tag{6.8}$$

where

$$\nabla_\lambda s(\lambda) = -B(\lambda)^{-1}s(\lambda). \tag{6.9}$$

Moreover, given the Cholesky factorization $B(\lambda) = L(\lambda)L^T(\lambda)$, it follows that

$$s(\lambda)^T \nabla_\lambda s(\lambda) = -\|w(\lambda)\|^2,$$

where $L(\lambda)L^T(\lambda)s(\lambda) = -g$ and $L(\lambda)w(\lambda) = s(\lambda)$.

Proof See the proof of [4, Lemma 7.3.1]. □

Then, for example, for $\phi_1(\lambda)$ from (6.7), we obtain the following expressions.

Corollary 6.2 *Suppose $g \neq 0$. Then the function $\phi_1(\lambda)$ is strictly increasing, when $\lambda > \max(0, -\lambda_1)$, and concave. Its first two derivatives are*

$$\phi'_1(\lambda) = -\frac{s(\lambda)^T \nabla_\lambda s(\lambda)}{\|s(\lambda)\|^3} + \frac{\sigma}{\lambda^2} > 0 \tag{6.10}$$

and

$$\phi''_1(\lambda) = \frac{3(s(\lambda)^T \nabla_\lambda s(\lambda))^2 - \|s(\lambda)\|_2^2 \|\nabla_\lambda s(\lambda)\|^2}{\|s(\lambda)\|^5} - \frac{2\sigma}{\lambda^3} < 0. \tag{6.11}$$

Proof Again, see the proof of [4, Lemma 7.3.1]. □

In this case, the basic Newton iteration is as follows.

Algorithm 6.1: Newton’s method to solve $\phi_1(\lambda) = 0$

Let $\lambda > \max(0, -\lambda_1)$ be given.

Step 1. Factorize $B(\lambda) = LL^T$.

Step 2. Solve $LL^T s = -g$.

Step 3. Solve $Lw = s$.

Step 4. Compute the Newton correction $\Delta\lambda^N = \frac{\lambda\left(\|s\| - \frac{\lambda}{\sigma}\right)}{\|s\| + \frac{\lambda}{\sigma}\left(\frac{\lambda\|w\|^2}{\|s\|^2}\right)}$.

Step 5. Replace λ by $\lambda + \Delta\lambda^N$. □

In practice, various safeguards of the kinds described for the trust-region case in [4, Sects. 7.3.4–7.3.8] should be added to ensure convergence of Algorithm 6.1 from an arbitrary initial λ . Numerical experience has indicated that the speed of (global) convergence may be improved by only linearizing the term $\omega(\lambda) \stackrel{\text{def}}{=} 1/\sqrt{\psi(\lambda)}$ of ϕ_1 in (6.7)—and not the σ/λ term as does Newton’s method—when computing a correction $\Delta\lambda^c$ to the estimate λ of the required root of ϕ_1 . The resulting correction thus satisfies the equation

$$\omega(\lambda) + \omega'(\lambda)\Delta\lambda^c \equiv \frac{1}{\psi^{\frac{1}{2}}(\lambda)} - \frac{1}{2}\frac{\psi'(\lambda)}{\psi^{\frac{3}{2}}(\lambda)}\Delta\lambda^c = \frac{\sigma}{\lambda + \Delta\lambda^c}, \tag{6.12}$$

which may be rewritten as a quadratic equation for $\Delta\lambda^c$.

Although Algorithm 6.1 and the variant using (6.12) are not generally globally convergent, there is one very important case in which they will be.

Theorem 6.3 *Suppose $\lambda > -\lambda_1$ and $\phi_1(\lambda) < 0$. Then both the Newton iterate $\lambda + \Delta\lambda^N$ and alternative $\lambda + \Delta\lambda^c$ inherit these properties and converge monotonically towards the required root, λ_* . The convergence is globally Q -linear with factor at least*

$$\gamma_\lambda \stackrel{\text{def}}{=} 1 - \frac{\phi'_1(\lambda_*)}{\phi'_1(\lambda)} < 1$$

and is ultimately Q -quadratic. Moreover $\lambda + \Delta\lambda^N \leq \lambda + \Delta\lambda^c \leq \lambda_*$.

Proof The proof in the case of the Newton iterate is essentially identical to that of [4, Lemma 7.3.2]. Since $\omega(\lambda)$ is a concave function of λ , (6.7) and (6.12) give that

$$\phi_1(\lambda + \Delta\lambda^c) = \omega(\lambda + \Delta\lambda^c) - \frac{\sigma}{\lambda + \Delta\lambda^c} \leq \omega(\lambda) + \omega'(\lambda)\Delta\lambda^c - \frac{\sigma}{\lambda + \Delta\lambda^c} = 0.$$

The Newton correction satisfies the linearized equation

$$\omega(\lambda) + \omega'(\lambda)\Delta\lambda^N = \frac{\sigma}{\lambda} - \frac{\sigma}{\lambda^2}\Delta\lambda^N. \tag{6.13}$$

But, as σ/λ is a convex function of λ ,

$$\frac{\sigma}{\lambda + \Delta\lambda^c} \geq \frac{\sigma}{\lambda} - \frac{\sigma}{\lambda^2}\Delta\lambda^c,$$

and hence

$$\omega(\lambda) + \omega'(\lambda)\Delta\lambda^c \geq \frac{\sigma}{\lambda} - \frac{\sigma}{\lambda^2}\Delta\lambda^c,$$

from (6.12). Combining this with (6.13), we obtain

$$\phi'_1(\lambda)(\Delta\lambda^c - \Delta\lambda^N) = \left(\omega'(\lambda) + \frac{\sigma}{\lambda^2}\right)(\Delta\lambda^c - \Delta\lambda^N) \geq 0$$

and hence $\Delta\lambda^c \geq \Delta\lambda^N > 0$ since Corollary 6.2 gives $\phi'_1(\lambda) > 0$. Thus the alternative iterate improves on the Newton one. \square

Similar results are easily derived for the other root functions defined in (6.5) and (6.7).

We conclude this section with an interesting observation concerning the global minimizer $s(\sigma)$ of the cubic model $m(s, \sigma) \equiv m(s)$ in (1.4), where we now make clear the explicit dependence on the parameter σ .

Theorem 6.4 *Suppose that $s(\sigma) \neq 0$ is a global minimizer of the model $m(s, \sigma) \equiv m(s)$ in (1.4). Then the length of the minimizer $v(\sigma) \stackrel{\text{def}}{=} \|s(\sigma)\|$ is a non-increasing function of σ .*

Proof We have from Theorem 3.1 that

$$(B + \sigma \|s(\sigma)\|I)s(\sigma) = -g \tag{6.14}$$

and that $B + \sigma \|s(\sigma)\|I$ and thus $B + \sigma \|s(\sigma)\|I + \sigma \|s(\sigma)\|^{-1}s(\sigma)s^T(\sigma)$ are positive semidefinite. We consider the derivative $v'(\sigma) = \|s(\sigma)\|^{-1}s^T(\sigma)\nabla_\sigma s(\sigma)$. Differentiating (6.14) with respect to σ reveals that

$$(B + \sigma \|s(\sigma)\|I)\nabla_\sigma s(\sigma) + \|s(\sigma)\|s(\sigma) + \sigma \|s(\sigma)\|^{-1}s(\sigma)s^T(\sigma)\nabla_\sigma s(\sigma) = 0$$

and thus that

$$\left(B + \sigma \|s(\sigma)\| I + \sigma \|s(\sigma)\|^{-1} s(\sigma) s^T(\sigma) \right) \nabla_{\sigma} s(\sigma) = -\|s(\sigma)\| s(\sigma). \tag{6.15}$$

Premultiplying (6.15) by $s^T(\sigma)$ and dividing by $\|s(\sigma)\|$ gives that

$$v'(\sigma) = -\frac{\nabla_{\sigma} s^T(\sigma) (B + \sigma \|s(\sigma)\| I + \sigma \|s(\sigma)\|^{-1} s(\sigma) s^T(\sigma)) \nabla_{\sigma} s(\sigma)}{\|s(\sigma)\|^2} \leq 0$$

since we have seen that $B + \sigma \|s(\sigma)\| I + \sigma \|s(\sigma)\|^{-1} s(\sigma) s^T(\sigma)$ is positive semidefinite. Thus $v'(\sigma)$ is a non-increasing function of σ . □

Griewank [17] provides a more complicated proof of the same result.

6.2 Computing an approximate solution

Of course, the dominant cost of the methods we have just discussed is that of factorizing $B + \lambda I$ for various λ , and this may be prohibitive for large n ; indeed factorization may be impossible. An obvious alternative is to use a Krylov method to approximate the solution. This was first proposed in [14] for trust-region methods.

The Lanczos method may be used to build up an orthogonal basis $\{q_0, \dots, q_j\}$ for the Krylov space $\{g, Bg, B^2g, \dots, B^jg\}$, formed by applying successively B to g . Letting $Q_j = (q_0, \dots, q_j)$, the key relationships are

$$Q_j^T Q_j = I, \quad Q_j^T B Q_j = T_j \quad \text{and} \quad Q_j^T g = \gamma_0 e_1, \tag{6.16}$$

where e_1 is the first unit vector of appropriate length and T_j is a symmetric, tridiagonal matrix.

We shall consider vectors of the form

$$s \in \mathcal{S}_j = \{s \in \mathbb{R}^n \mid s = Q_j u\}$$

and seek

$$s_j = Q_j u_j, \tag{6.17}$$

where s_j solves the problem

$$\min_{s \in \mathcal{S}_j} m(s). \tag{6.18}$$

It then follows directly from (6.16) that u_j solves the problem

$$\min_{u \in \mathbb{R}^{j+1}} m_j(u) \stackrel{\text{def}}{=} f + \gamma_0 u^T e_1 + \frac{1}{2} u^T T_j u + \frac{1}{3} \sigma \|u\|_2^3. \tag{6.19}$$

There are a number of crucial observations to be made here. Firstly, as T_j is tridiagonal, it is feasible to use the method broadly described in Sect. 6.1 to compute the solution to (6.18) even when n is large. Secondly, having found u_j , the matrix Q_j is needed to recover s_j , and thus the Lanczos vectors q_j will either need to be saved on backing store or regenerated when required. We only need Q_j once we are satisfied that continuing the Lanczos process will give little extra benefit. Thirdly, one would hope that as a sequence of such problems may be solved, and as T_j only changes by the addition of an extra diagonal and superdiagonal entry, solution data from one subproblem may be useful for starting the next. Lastly, this is a clear extension of the GLTR method for the solution of the trust-region problem [14], and many of the implementation issues and details follow directly from there.

Furthermore, employing this approach within the ARC algorithm benefits from the theoretical guarantees of convergence in Sects. 2.2–5, and satisfies the complexity bounds developed in [2]. To see this, let $\mathcal{L}_k = \mathcal{S}_j$ in Lemma 3.2 and note that the current gradient is included in all subspaces \mathcal{S}_j .

6.3 Scaled regularisation

The preceding development can trivially be generalized if the regularisation $\frac{1}{3}\sigma\|s\|_2^3$ is replaced by the scaled variant $\frac{1}{3}\sigma\|s\|_M^3$, where we define $\|s\|_M = s^T M s$ for some symmetric positive definite M . All that changes is that the key second-derivative matrix is $B(\lambda) = B + \lambda M$ in Theorem 3.1 and its successors, and that M -orthogonal vectors are generated using the preconditioned Lanczos method; the regularisation in the tridiagonal problem (6.19) is not altered.

6.4 A further possibility

Another possibility is suppose that the optimal $\|s\|_2$ is known to be of size Δ . In this case, the required value of $m(s)$ is $f + s^T g + \frac{1}{2}s^T B s + \frac{1}{3}\sigma\Delta^3$, where s solves the trust-region problem

$$q(\Delta) = \min_{s \in \mathbb{R}^n} s^T g + \frac{1}{2}s^T B s \text{ subject to } \|s\|_2 = \Delta.$$

Hence

$$\min_{s \in \mathbb{R}^n} m(s) = \min_{\Delta \in \mathbb{R}_+} q(\Delta) + \frac{1}{3}\sigma\Delta^3$$

and we may use known algorithms for the trust-region problem to accomplish the univariate minimization of $q(\Delta) + \frac{1}{3}\sigma\Delta^3$. We have not considered this possibility further at this stage.

7 Numerical results

We now turn our attention to investigating how the cubic regularisation method performs in practice. We have implemented the ARC algorithm with B set to the true Hessian H , together with both the exact and inexact subproblem solvers described in Sects. 6.1 and 6.2. To be specific, when solving the subproblem, we compute the required root λ of $\phi_1(\lambda)$; we use Algorithm 6.1 to find the required root, replacing the correction $\Delta\lambda$ in Step 4 by the improvement given by (6.12). To simplify matters, we start the root finding from $\max(0, -\lambda_1) + \epsilon$ for some tiny ϵ —this of course entails that we find the eigenvalue λ_1 of B (Sect. 6.1 and the built-in MATLAB function `eigs`) or T_j (Sect. 6.2 and the specially-designed algorithm given in [14]), which is potentially expensive in the case of B , but far less so for T_j , especially since we have that of T_{j-1} as a starting estimate—in which case Theorem 6.3 will ensure global (and ultimately) rapid convergence.

In view of its suitability for large-scale problems, in the results we shall present, we used the Lanczos-based inexact solver described in Sect. 6.2. Using the exact solver gives similar results for the small-scale problems that we tested, since the values of the parameters in the stopping rules we have chosen to use require that we solve to reasonably high relative accuracy in the inexact case (see (7.1)–(7.3)). We considered three stopping rules for the inexact inner iteration, all derived from the TC.h criteria in Sect. 3.3. In the first, recalling TC.g, we stop as soon as the approximate solution in Step 2 of the ARC algorithm satisfies

$$\|\nabla m_k(s_k)\| \leq \min(0.0001, \|\nabla m_k(0)\|^{\frac{1}{2}}) \|\nabla m_k(0)\|; \tag{7.1}$$

the aim is to try to encourage rapid ultimate convergence [5, 6, 19] without the expense of “over-solving” when far from optimality—we refer to (7.1) as the “ g rule” (see Corollary 4.8 where we show the ARC algorithm with such a termination criterion converges Q-superlinearly). The remaining rules are geared more towards ensuring the best overall complexity bound we have obtained (see [2]). Thus our second “ s rule” comes from TC.s, and it is to stop as soon as

$$\|\nabla m_k(s_k)\| \leq \min(0.0001, \|s_k\|) \|\nabla m_k(0)\|. \tag{7.2}$$

However since we were concerned that this might be overly stringent when s_k is small when σ_k is large rather than because $\nabla m_k(0)$ is small, our final “ s/σ rule” is to stop as soon as

$$\|\nabla m_k(s_k)\| \leq \min\left(0.0001, \frac{\|s_k\|}{\max(1, \sigma_k)}\right) \|\nabla m_k(0)\|. \tag{7.3}$$

The ARC algorithm converges at least Q-superlinearly also when (7.2) and (7.3) are employed (see Corollaries 4.8 and 4.10, and the remarks inbetween).

The other crucial ingredient is the management of the regularisation parameter σ_k in Step 4 of the ARC algorithm. Here, on very successful iterations, we set $\sigma_{k+1} = \max(\min(\sigma_k, \|g_k\|), \epsilon_M)$, where $\epsilon_M \approx 10^{-16}$ is the relative machine precision—the

intention is to try to reduce the model rapidly to Newton ($\sigma = 0$) model once convergence sets in, while maintaining some regularisation before the asymptotics are entered. For other successful steps we leave σ_k unchanged, while for unsuccessful steps we increase σ_k by 2 (the choice of the “2” factor is for simplicity, and it is likely that better values are possible in a similar vein to [15]). We start with $\sigma_0 = 1$, and use $\eta_1 = 0.1$ and $\eta_2 = 0.9$, similar performance being observed for other choices of initial parameters.

By way of a comparison, we have also implemented a standard trust-region method [4, Alg. 6.1.1.]. Here we have used the GLTR method [14] to find an approximate solution of the trust-region problem, stopping as above as soon as (7.1) is satisfied. The trust-region radius Δ_{k+1} following a very successful iteration is $\min(\max(2\|s_k\|, \Delta_k), 10^{10})$, it is left unchanged if the iteration is merely successful, while an unsuccessful step results in a halving of the radius. The initial radius is always set to 1.

We give the results obtained by applying both Algorithms to all of the unconstrained problems from the CUTEr collection [16]; for those whose dimensions may be adjusted, we chose small variants simply so as not to overload our (Matlab) computing environment, most particularly the CUTEr interface to Matlab. All of our experiments were performed on a single processor of a 3.06 GHz Dell Precision 650 Workstation. Both our new algorithm, and the trust-region algorithm were implemented as Matlab M-files, and the tests performed using Matlab 7.2.

We give the complete set of results in Appendix A. The algorithm is stopped as soon as the norm of the gradient $\|g(x_k)\|$ is smaller than 10^{-5} . An upper limit of 10000 iterations was set, and any run exceeding this is flagged as a failure. In Fig. 4, we present the iteration-count performance profile [10] for the methods. We would have liked to have given a similar figure for CPU times, but the Matlab CPU timer proved too inaccurate for this purpose—we defer such a comparison until we have produced a careful (Fortran) implementation and tested it on larger examples in a controlled computing environment.

While we should regard our results as tentative, since improvements to both algorithms are most likely, we are delighted with the overall performance of our new algorithm. Trust-region methods are generally considered to be reliable and effective methods for unconstrained minimization, and thus it is gratifying to be able to improve upon them in so many cases. Of course, there are cases where the new algorithm isn't as effective, but in general the algorithm appears able to move to the asymptotic region more rapidly. Whether this is a consequence of a provably good complexity bound or for other reasons of robustness is not clear. For the three cases where the new algorithm fails, slow (but sure) progress is made towards the solution, but these problems are well known to be very hard. Of the three variants of the new method, that which is less concerned with provably superior worst-case complexity, appears to be more promising. This is perhaps not so surprising since the more conservative acceptance rules (7.2) and (7.3) aim for guaranteed rather than speculative reduction. There is very little to choose between the latter pair, indicating that our concern that (7.2) may be over-stringent may have been misplaced.

Using quasi-Newton updates rather than the exact Hessian in computations will require careful implementation and thought. Clearly, we do not need positive definiteness of the updates due to the regularisation term of the cubic model, and hence

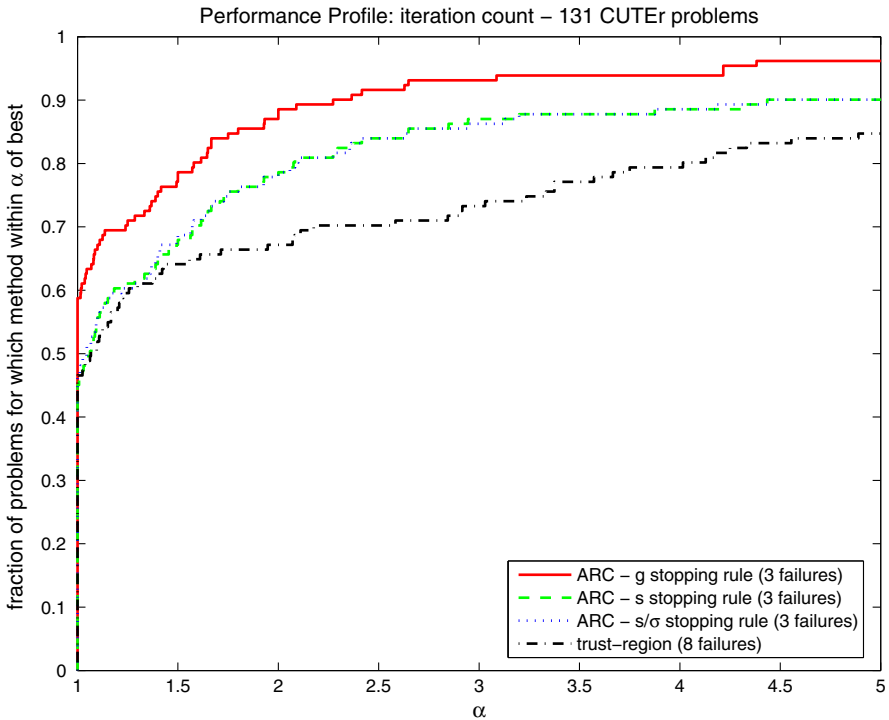


Fig. 4 Performance profile, $p(\alpha)$: Iteration counts for the 131 problems under consideration

indefinite update formulae, such as the symmetric rank-one formula, may be used. Indeed, choices that ultimately aim to imitate the true Hessian (such as the symmetric rank-one and Powell symmetric-Broyden formulae) should be used so as to reveal problem negative curvature and hence allow convergence to weak second-order critical points (see Sect. 5). Recalling our remarks following (4.8) and (4.19), and also some potentially useful linear algebra details in [11, 12], the Powell symmetric-Broyden and symmetric rank-one updates may prove to be a reasonable choice in this context, for medium-scale problems.

8 Conclusions

In this paper we have considered the convergence properties of a new general cubic regularisation framework for unconstrained optimization which has roots in earlier algorithmic proposals by Griewank [17], Nesterov and Polyak [23] and Weiser et al. [26]. The framework allows for the approximate solution of the key step calculation, and is suitable for large-scale problems. We presented a Lanczos-based approximation method which is covered by our theoretical developments. Encouraging preliminary numerical experiments with small-scale problems were reported.

Our next goal is to implement and test these ideas carefully in the large-scale case, and also to investigate the computational efficiency of using quasi-Newton updates in

place of the exact Hessian. The latter approach fits within the theoretical framework of this paper.

Extensions to these ideas are obvious and far-reaching. In particular, since the use of trust-region models is widespread in optimization, it is worth investigating where cubic models might be employed in their place. Projection methods for bound constrained minimization and penalty/barrier/augmented Lagrangian methods for constrained optimization are obvious candidates. Note that in the case of linear equality constraints, the (semi-)norm will only need to regularise in the null-space of the constraints, and solving the subproblem is likewise easy so long as the Krylov subspace is projected onto the constraint manifold [4]. More generally, difficulties resulting from the incompatibility of the intersection of linearized constraints with trust-region bounds has been a perennial problem in constrained optimization; (cubic) regularisation offers an easy way to avoid this difficulty.

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Appendix A

Here we give the complete set of results from our tests. For each problem, in Table 1 we report its number of variables (n), along with the number of iterations (= the number of function evaluations) required for convergence (iter), the number of gradient evaluations ($\#g$), and the best objective function value found (f ; the subscript gives the base-10 exponent) for the four rival methods. The symbol $>$ indicates that the iteration limit was exceeded.

Table 1 Comparison between the trust-region and ARC algorithms

Name	n	Trust-region			ARC with g -rule			ARC with s -rule			ARC with s/σ -rule		
		iter	$\#g$	f	iter	$\#g$	f	iter	$\#g$	f	iter	$\#g$	f
ALLINITU	4	8	8	5.74	16	9	5.74	16	9	5.74	16	9	5.74
ARGLINA	200	30	30	2.00 ₊₂	8	8	2.00 ₊₂	8	8	2.00 ₊₂	8	8	2.00 ₊₂
ARWHEAD	100	6	6	6.59 ₋₁₄	6	6	8.79 ₋₁₄	6	6	8.79 ₋₁₄	6	6	8.79 ₋₁₄
BARD	3	8	8	8.21 ₋₃	8	8	8.21 ₋₃	8	8	8.21 ₋₃	8	8	8.21 ₋₃
BDQRTIC	100	14	14	3.79 ₊₂	10	10	3.79 ₊₂	10	10	3.79 ₊₂	10	10	3.79 ₊₂
BEALE	2	9	7	7.55 ₋₁₄	10	7	9.89 ₋₁₂	10	7	9.89 ₋₁₂	10	7	9.89 ₋₁₂
BIGGS6	6	200	196	2.43 ₋₁	66	47	1.66 ₋₁₀	76	52	1.35 ₋₁₁	76	52	1.35 ₋₁₁
BOX3	3	8	8	2.03 ₋₁₁	9	9	5.84 ₋₁₆	9	9	5.84 ₋₁₆	9	9	5.84 ₋₁₆
BRKMCC	2	3	3	1.69 ₋₁	4	4	1.69 ₋₁	4	4	1.69 ₋₁	4	4	1.69 ₋₁
BROWNAL	200	11	11	5.49 ₋₂₀	5	5	8.29 ₋₁₇	3	3	1.47 ₋₉	3	3	1.47 ₋₉
BROWNBS	2	>	>	9.80 ₊₁₁	28	27	2.17 ₋₂₉	28	27	1.78 ₋₂₉	28	27	1.78 ₋₂₉
BROWNDEN	4	44	44	8.58 ₊₄	9	9	8.58 ₊₄	9	9	8.58 ₊₄	9	9	8.58 ₊₄
BROYDN7D	100	22	22	3.24 ₊₁	25	17	3.01 ₊₁	25	17	3.01 ₊₁	25	17	3.01 ₊₁
BRYBND	100	18	15	5.54 ₋₁₇	16	10	5.32 ₋₁₇	16	10	5.34 ₋₁₇	16	10	5.32 ₋₁₇
CHAINWOO	100	255	253	3.22 ₊₁	60	42	1.00	60	41	1.00	60	42	1.00

Table 1 continued

Name	n	Trust-region			ARC with <i>g</i> -rule			ARC with <i>s</i> -rule			ARC with <i>s/σ</i> -rule		
		iter	# <i>g</i>	<i>f</i>	iter	# <i>g</i>	<i>f</i>	iter	# <i>g</i>	<i>f</i>	iter	# <i>g</i>	<i>f</i>
CHNROSNB	50	63	61	1.28 ₋₁₈	68	42	1.82 ₋₁₆	68	42	1.80 ₋₁₆	68	42	1.80 ₋₁₆
CLIFF	2	28	28	2.00 ₋₁	28	28	2.00 ₋₁	28	28	2.00 ₋₁	28	28	2.00 ₋₁
CRAGGLVY	202	29	29	6.67 ₊₁	14	14	6.67 ₊₁	14	14	6.67 ₊₁	14	14	6.67 ₊₁
CUBE	2	32	28	1.00 ₋₁₉	56	27	7.73 ₋₂₁	56	27	7.73 ₋₂₁	56	27	7.73 ₋₂₁
CURLY10	50	15	15	-5.02 ₊₃	27	16	-5.02 ₊₃	27	16	-5.02 ₊₃	27	16	-5.02 ₊₃
CURLY20	50	12	12	-5.02 ₊₃	29	15	-5.02 ₊₃	29	15	-5.02 ₊₃	29	15	-5.02 ₊₃
CURLY30	50	19	18	-5.02 ₊₃	30	15	-5.02 ₊₃	30	15	-5.02 ₊₃	30	15	-5.02 ₊₃
DECONVU	61	46	38	1.40 ₋₉	142	56	3.92 ₋₁₁	131	55	1.27 ₋₁₀	144	56	1.41 ₋₁₀
DENSCHNA	2	6	6	2.21 ₋₁₂	6	6	1.83 ₋₁₂	6	6	1.83 ₋₁₂	6	6	1.83 ₋₁₂
DENSCHNB	2	6	6	1.04 ₋₁₃	6	6	2.74 ₋₁₅	6	6	2.74 ₋₁₅	6	6	2.74 ₋₁₅
DENSCHNC	2	11	11	2.18 ₋₂₀	11	11	5.50 ₋₁₉	11	11	5.50 ₋₁₉	11	11	5.50 ₋₁₉
DENSCHND	3	36	36	1.77 ₋₈	50	33	8.74 ₋₉	50	33	8.74 ₋₉	50	33	8.74 ₋₉
DENSCHNE	3	16	16	1.10 ₋₁₈	24	14	9.92 ₋₁₄	25	15	1.84 ₋₁₁	25	15	1.84 ₋₁₁
DENSCHNF	2	7	7	6.51 ₋₂₂	7	7	6.93 ₋₂₂	7	7	6.93 ₋₂₂	7	7	6.93 ₋₂₂
DIXMAANA	150	27	27	1.00	8	8	1.00	8	8	1.00	8	8	1.00
DIXMAANB	150	27	27	1.00	8	8	1.00	8	8	1.00	8	8	1.00
DIXMAANC	150	27	27	1.00	22	15	1.00	23	16	1.00	23	16	1.00
DIXMAAND	150	27	27	1.00	26	18	1.00	26	18	1.00	26	18	1.00
DIXMAANE	150	31	31	1.00	12	12	1.00	12	12	1.00	12	12	1.00
DIXMAANF	150	29	29	1.00	24	19	1.00	24	19	1.00	24	19	1.00
DIXMAANG	150	29	29	1.00	29	23	1.00	29	23	1.00	29	23	1.00
DIXMAANH	150	29	29	1.00	31	23	1.00	31	23	1.00	31	23	1.00
DIXMAANI	150	37	37	1.00	17	17	1.00	17	17	1.00	17	17	1.00
DIXMAANJ	150	38	36	1.00	33	28	1.00	33	28	1.00	33	28	1.00
DIXMAANK	150	35	34	1.00	35	28	1.00	35	28	1.00	35	28	1.00
DIXMAANL	150	41	39	1.00	37	29	1.00	37	29	1.00	37	29	1.00
DJTL	2	81	72	-8.95 ₊₃	1,655	1,581	-8.95 ₊₃	1,655	1,581	-8.95 ₊₃	1,655	1,581	-8.95 ₊₃
DQRTIC	100	594	594	1.20 ₋₇	25	25	2.36 ₋₈	25	25	2.36 ₋₈	25	25	2.36 ₋₈
EDENSCH	100	76	76	6.03 ₊₂	12	12	6.03 ₊₂	12	12	6.03 ₊₂	12	12	6.03 ₊₂
EG2	100	4	4	-9.89 ₊₁	4	4	-9.89 ₊₁	4	4	-9.89 ₊₁	4	4	-9.89 ₊₁
EIGENALS	110	23	23	2.83 ₋₁₅	25	20	3.38 ₋₁₅	25	20	4.56 ₋₁₅	25	20	6.56 ₋₁₆
EIGENBLS	110	88	83	8.84 ₋₁₅	170	71	7.46 ₋₁₁	202	78	1.40 ₋₁₂	208	81	2.02 ₋₁₂
EIGENCLS	132	46	44	1.02 ₋₁₅	57	35	9.47 ₋₁₃	57	35	1.82 ₋₁₁	56	35	4.34 ₋₁₂
ENGVAL1	100	19	19	1.09 ₊₂	9	9	1.09 ₊₂	9	9	1.09 ₊₂	9	9	1.09 ₊₂
ENGVAL2	3	14	14	9.71 ₋₁₇	27	16	3.08 ₋₂₀	27	16	3.08 ₋₂₀	27	16	3.08 ₋₂₀
ERRINROS	50	127	123	4.39 ₊₁	66	39	4.39 ₊₁	61	41	3.99 ₊₁	61	41	3.99 ₊₁
EXPFIT	2	10	8	2.41 ₋₁	14	7	2.41 ₋₁	14	7	2.41 ₋₁	14	7	2.41 ₋₁
EXTROSNB	100	1,619	1,591	7.96 ₋₉	6,826	1,197	1.67 ₋₈	6,984	1,219	1.58 ₋₈	6,768	1,187	1.71 ₋₈
FLETGBV2	100	3	3	-5.14 ₋₁	5	5	-5.14 ₋₁	5	5	-5.14 ₋₁	5	5	-5.14 ₋₁
FLETGBV3	50	>	>	-3.43	>	>	-1.17 ₊₂	>	>	-1.17 ₊₂	>	>	-1.18 ₊₂
FLETCHBV	10	5,012	5,012	-1.99 ₊₆	341	299	-2.17 ₊₆	496	419	-2.28 ₊₆	465	388	-2.22 ₊₆

Table 1 continued

Name	n	Trust-region			ARC with g -rule			ARC with s -rule			ARC with s/σ -rule		
		iter	# g	f	iter	# g	f	iter	# g	f	iter	# g	f
FLETCHCR	100	154	151	1.00 ₋₁₅	230	154	6.20 ₋₁₇	230	154	6.20 ₋₁₇	230	154	6.20 ₋₁₇
FMINSRF2	121	128	126	1.00	45	38	1.00	45	38	1.00	45	38	1.00
FMINSURF	121	67	66	1.00	43	36	1.00	43	36	1.00	43	36	1.00
FREUROTH	100	55	30	1.20 ₊₄	17	12	1.20 ₊₄	17	12	1.20 ₊₄	17	12	1.20 ₊₄
GENHUMPS	10	>	>	1.42 ₊₃	7,599	3,977	1.32 ₋₁₀	7,442	3,831	3.66 ₋₁₇	7,596	3,983	1.96 ₋₁₀
GENROSE	100	79	77	1.00	130	62	1.00	135	63	1.00	133	63	1.00
GENROSEB	500	663	660	1.00	647	290	1.00	652	294	1.00	657	294	1.00
GROWTHLS	3	142	134	1.00	115	69	1.00	113	71	1.00	113	71	1.00
GULF	3	248	246	2.96 ₋₁₁	45	33	2.89 ₋₁₁	45	33	2.90 ₋₁₁	45	33	2.89 ₋₁₁
HAIRY	2	103	100	2.00 ₊₁	75	31	2.00 ₊₁	81	33	2.00 ₊₁	80	33	2.00 ₊₁
HATFLDD	3	19	18	6.62 ₋₈	26	23	6.62 ₋₈	26	23	6.62 ₋₈	26	23	6.62 ₋₈
HATFLDE	3	20	19	5.12 ₋₇	22	17	5.12 ₋₇	22	17	5.12 ₋₇	22	17	5.12 ₋₇
HEART6LS	6	3,369	3,363	4.09 ₋₁₆	407	310	4.02 ₋₁₄	413	300	2.55 ₋₁₆	426	304	1.32 ₋₁₄
HEART8LS	8	103	99	1.42 ₋₁₉	94	49	1.99 ₋₁₇	98	48	8.26 ₋₁₈	95	50	7.57 ₋₁₅
HELIX	3	13	12	1.16 ₋₁₅	21	12	3.13 ₋₁₃	21	12	3.13 ₋₁₃	21	12	3.13 ₋₁₃
HIMMELBB	2	14	10	2.52 ₋₂₁	21	7	9.71 ₋₂₆	21	7	9.71 ₋₂₆	21	7	9.71 ₋₂₆
HUMPS	2	6,015	6,007	2.39 ₋₁₈	1,552	636	2.02 ₋₁₂	1,591	663	1.64 ₋₁₀	1,498	659	3.68 ₋₁₀
HYDC20LS	99	>	>	5.37 ₋₁	>	>	7.67 ₋₁	>	>	7.75 ₋₁	>	>	7.68 ₋₁
JENSMP	2	10	10	1.24 ₊₂	10	10	1.24 ₊₂	10	10	1.24 ₊₂	10	10	1.24 ₊₂
KOWOSB	4	13	10	3.08 ₋₄	10	8	3.08 ₋₄	10	8	3.08 ₋₄	10	8	3.08 ₋₄
LIARWHD	100	35	35	3.87 ₋₁₄	12	12	4.44 ₋₂₆	12	12	4.44 ₋₂₆	12	12	4.44 ₋₂₆
LOGHAIRY	2	9,349	9,346	1.82 ₋₁	2,737	1,208	1.82 ₋₁	2,696	1,198	1.82 ₋₁	2,617	1,145	1.82 ₋₁
MANCINO	100	754	754	1.27 ₋₂₁	27	11	1.35 ₋₂₁	27	11	1.35 ₋₂₁	27	11	1.16 ₋₂₁
MEXHAT	2	37	28	-4.00 ₋₂	98	27	-4.00 ₋₂	98	27	-4.00 ₋₂	98	27	-4.00 ₋₂
MEYER3	3	>	>	9.02 ₊₁	361	234	8.79 ₊₁	422	269	8.79 ₊₁	420	268	8.79 ₊₁
MOREBV	100	95	95	2.87 ₋₇	59	59	4.16 ₋₇	102	102	2.86 ₋₇	102	102	2.86 ₋₇
MSQRTALS	100	22	19	1.42 ₋₁₄	18	15	5.84 ₋₁₅	18	15	4.73 ₋₁₅	18	15	5.77 ₋₁₅
MSQRTBLS	100	20	17	3.32 ₋₁₂	18	15	2.07 ₋₁₅	18	15	1.99 ₋₁₅	18	15	2.06 ₋₁₅
NONCVXU2	100	714	714	2.33 ₊₂	51	42	2.32 ₊₂	49	41	2.32 ₊₂	49	41	2.32 ₊₂
NONCVXUN	100	689	689	2.35 ₊₂	43	33	2.33 ₊₂	47	38	2.33 ₊₂	47	38	2.33 ₊₂
NONDIA	100	7	7	1.50 ₋₁₈	11	8	2.90 ₋₁₈	11	8	2.90 ₋₁₈	11	8	2.90 ₋₁₈
NONDQUAR	100	17	17	2.58 ₋₁₅	16	16	1.42 ₋₈	62	38	1.81 ₋₆	62	38	1.81 ₋₆
NONMSQRT	100	3,838	3,838	1.81 ₊₁	2,700	2,535	1.81 ₊₁	2,959	2,926	1.81 ₊₁	3,809	3,777	1.81 ₊₁
OSBORNEA	5	111	104	5.46 ₋₅	468	266	4.69 ₋₂	697	465	4.70 ₋₂	991	607	4.70 ₋₂
OSBORNEB	11	21	20	4.01 ₋₂	21	18	4.01 ₋₂	21	18	4.01 ₋₂	21	18	4.01 ₋₂
OSCIPTH	8	3,910	3,829	5.61 ₋₈	8,172	2,474	8.67 ₋₈	8,115	2,463	8.75 ₋₈	8,209	2,485	8.22 ₋₈
PALMER5C	6	64	64	2.13	8	8	2.13	8	8	2.13	8	8	2.13
PALMER6C	8	512	512	1.64 ₋₂	124	124	1.64 ₋₂	239	239	1.64 ₋₂	239	239	1.64 ₋₂
PALMER7C	8	1,243	1,243	6.02 ₋₁	55	55	6.02 ₋₁	162	162	6.02 ₋₁	162	162	6.02 ₋₁
PALMER8C	8	590	590	1.60 ₋₁	303	303	1.60 ₋₁	311	311	1.60 ₋₁	311	311	1.60 ₋₁
PARKCH	15	1	1	4.73 ₋₇	1	1	4.73 ₋₇	42	25	1.62 ₊₃	56	26	1.62 ₊₃

Table 1 continued

Name	n	Trust-region			ARC with g -rule			ARC with s -rule			ARC with s/σ -rule		
		iter	# g	f	iter	# g	f	iter	# g	f	iter	# g	f
PENALTY1	100	610	602	9.02 ₋₄	85	35	9.02 ₋₄	85	35	9.02 ₋₄	85	35	9.02 ₋₄
PENALTY2	200	2	2	4.71 ₊₁₃	4	4	4.71 ₊₁₃	11	11	4.71 ₊₁₃	11	11	4.71 ₊₁₃
PENALTY3	200	1	1	1.16 ₊₆	1	1	1.16 ₊₆	29	14	1.00 ₋₃	24	14	9.97 ₋₄
PFIT1LS	3	331	315	1.66 ₋₁₂	870	233	1.07 ₋₁₃	870	233	1.07 ₋₁₃	870	233	1.07 ₋₁₃
PFIT2LS	3	104	92	2.71 ₋₁₃	246	69	1.46 ₋₁₆	247	69	1.16 ₋₁₆	245	69	2.55 ₋₁₆
PFIT3LS	3	131	119	3.44 ₋₁₄	574	168	7.03 ₋₁₈	581	166	1.61 ₋₁₃	581	166	4.21 ₋₁₃
PFIT4LS	3	227	212	4.62 ₋₁₆	1,319	417	2.06 ₋₁₄	1,319	419	1.01 ₋₁₃	1,316	419	9.86 ₋₁₄
POWELLSG	4	5	5	1.93 ₋₃₀	5	5	1.81 ₋₁₂	16	16	4.54 ₋₉	16	16	4.54 ₋₉
POWER	100	28	28	1.25 ₋₉	24	24	1.61 ₋₉	24	24	1.61 ₋₉	24	24	1.61 ₋₉
QUARTC	100	566	566	0.00	12	12	1.42 ₋₂₅	25	25	2.36 ₋₈	25	25	2.36 ₋₈
ROSENBR	2	4	4	1.71 ₋₃₂	5	5	1.07 ₋₁₅	37	20	1.80 ₋₁₂	37	20	1.80 ₋₁₂
S308	2	1	1	0.00	1	1	0.00	10	10	7.73 ₋₁	10	10	7.73 ₋₁
SBRYBND	100	>	>	6.76 ₊₆	>	>	6.76 ₊₆	>	>	6.76 ₊₆	>	>	2.44 ₊₂
SCHMVEIT	100	6	6	-2.94 ₊₂	5	5	-2.94 ₊₂	5	5	-2.94 ₊₂	5	5	-2.94 ₊₂
SENSORS	100	25	23	-1.97 ₊₃	34	23	-1.94 ₊₃	41	25	-1.94 ₊₃	34	23	-1.94 ₊₃
SINEVAL	2	57	53	1.58 ₋₂₅	94	43	4.77 ₋₁₄	94	43	4.80 ₋₁₄	94	43	4.77 ₋₁₄
SINQUAD	100	16	16	-4.01 ₊₃	15	9	-4.01 ₊₃	15	9	-4.01 ₊₃	15	9	-4.01 ₊₃
SISSER	2	13	13	1.07 ₋₈	13	13	1.14 ₋₈	13	13	1.14 ₋₈	13	13	1.14 ₋₈
SNAIL	2	12	12	2.54 ₋₁₅	7	7	8.46 ₋₁₆	100	63	2.16 ₋₁₇	100	63	2.16 ₋₁₇
SPARSINE	100	11	11	3.63 ₋₁₇	25	17	1.84 ₋₁₇	25	17	2.17 ₋₁₇	25	17	1.86 ₋₁₇
SPARSQR	100	17	17	2.32 ₋₈	17	17	7.78 ₋₉	17	17	7.78 ₋₉	17	17	7.78 ₋₉
SPMSRTL	100	12	12	9.34 ₋₁₄	17	16	3.90 ₋₁₃	17	16	4.53 ₋₁₃	17	16	4.53 ₋₁₃
SROSENBR	100	7	7	9.08 ₋₁₃	9	9	6.52 ₋₁₆	9	9	6.52 ₋₁₆	9	9	6.52 ₋₁₆
STREG	4	>	>	1.00 ₊₂₀	104	52	2.95 ₋₁₃	214	147	9.88 ₋₂₀	214	147	9.88 ₋₂₀
TOINTGOR	50	41	41	1.37 ₊₃	9	9	1.37 ₊₃	9	9	1.37 ₊₃	9	9	1.37 ₊₃
TOINTGSS	100	32	32	1.01 ₊₁	11	10	1.02 ₊₁	11	10	1.02 ₊₁	11	10	1.02 ₊₁
TOINTPSP	50	35	35	2.26 ₊₂	30	19	2.26 ₊₂	30	19	2.26 ₊₂	30	19	2.26 ₊₂
TQUARTIC	100	78	75	1.09 ₋₂₀	18	13	1.62 ₋₂₃	18	13	1.62 ₋₂₃	18	13	1.62 ₋₂₃
VARDIM	200	10	10	1.05 ₋₂₅	3	3	7.19 ₋₂₆	30	30	6.79 ₋₂₅	30	30	6.79 ₋₂₅
VAREIGVL	50	20	17	7.98 ₋₁₀	14	14	8.39 ₋₁₀	14	14	7.97 ₋₁₀	14	14	7.99 ₋₁₀
VIBRBEAM	8	>	>	4.54	261	144	1.56 ₋₁	304	188	1.56 ₋₁	257	140	1.56 ₋₁
WATSON	12	11	11	9.13 ₋₉	10	10	1.86 ₋₉	10	10	1.83 ₋₉	10	10	1.83 ₋₉
WOODS	4	71	68	2.18 ₋₂₁	69	39	6.99 ₋₁₉	69	39	6.99 ₋₁₉	69	39	6.99 ₋₁₉
YFITU	3	246	245	6.79 ₋₁₃	59	46	6.82 ₋₁₃	59	46	6.82 ₋₁₃	59	46	6.82 ₋₁₃

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