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High-Order Evaluation Complexity for Convexly-Constrained Optimization with Non-Lipschitzian Group Sparsity Terms

X. Chen^{*} and Ph. L. Toint^{\dagger}

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Abstract

This paper studies high-order evaluation complexity for partially separable convexlyconstrained optimization involving non-Lipschitzian group sparsity terms in a nonconvex objective function. We propose a partially separable adaptive regularization algorithm using a *p*-th order Taylor model and show that the algorithm needs at most $O(e^{-(p+1)/(p-q+1)})$ evaluations of the objective function and its first *p* derivatives (whenever they exist) to produce an (ϵ, δ) -approximate *q*-th-order stationary point. Our algorithm uses the underlying rotational symmetry of the Euclidean norm function to build a Lipschitzian approximation for the non-Lipschitzian group sparsity terms, which are defined by the group ℓ_2 - ℓ_a norm with $a \in (0, 1)$. The new result shows that the partiallyseparable structure and non-Lipschitzian group sparsity terms in the objective function do not affect the worst-case evaluation complexity order.

Keywords: complexity theory, nonlinear optimization, non-Lipschitz functions, partially-separable problems, group sparsity, isotropic model.

AMS subject classifications:, 90C30, 90C46, 65K05

1 Introduction

Both applied mathematicians and computer scientists have, in recent years, made significant contributions to the fast-growing field of worst-case complexity analysis for nonconvex optimization (see [12] for a partial yet substantial bibliography). The purpose of this paper is to extend the available general theory in two distinct directions. The first is to cover the case where the problem involving non-Lipschitzian group sparsity terms. The second is to show that the ubiquitous partially-separable structure of the objective function (of which standard sparsity is a special case) can be exploited without affecting the complexity bounds derived in [12].

We consider the partially-separable convexly constrained nonlinear optimization problem:

$$\min_{x \in \mathcal{F}} f(x), \quad \text{where } f(x) = \sum_{i \in \mathcal{N}} f_i(U_i x) + \sum_{i \in \mathcal{H}} \|U_i x - b_i\|^a \stackrel{\text{def}}{=} \sum_{i \in \mathcal{N} \cup \mathcal{H}} f_i(U_i x), \qquad (1.1)$$

 $\mathcal{N} \cup \mathcal{H} \stackrel{\text{def}}{=} \mathcal{M}, \, \mathcal{N} \cap \mathcal{H} = \emptyset, \, f_i \text{ is a continuously } p \text{ times differentiable function from } \mathbb{R}^{n_i} \text{ into } \mathbb{R} \text{ for } i \in \mathcal{N}, \text{ and } f_i(U_i x) = \|U_i x - b_i\|^a \text{ for } i \in \mathcal{H}, \, a \in (0, 1), \, \|\cdot\| \text{ is the Euclidean norm,}$

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 $U_i \in \mathbb{R}^{n_i \times n}$ with $n_i \leq n$, and $b_i \in \mathbb{R}^{n_i}$. For simplicity, we assume that, for each $i \in \mathcal{M}$, U_i has full row rank and $||U_i|| = 1$, and that the ranges of the U_i^T for $i \in \mathcal{N}$ span \mathbb{R}^n so that the intersection of the nullspaces of the U_i is reduced to the origin. We also assume that the ranges of the U_i^T (for $i \in \mathcal{H}$) are orthogonal, that is

$$U_i U_j^T = 0 \text{ for } i \neq j, \, i, j \in \mathcal{H}.$$

$$(1.2)$$

Without loss of generality, we furthermore assume that the rows of U_i are orthonormal for $i \in \mathcal{H}$. Our final assumption, as in [17], is that the feasible set $\mathcal{F} \subseteq \mathbb{R}^n$ is non-empty closed and convex, and that it is "kernel-centered" in that is, if $P_{\mathcal{X}}[\cdot]$ is the orthogonal projection onto the convex set \mathcal{X} and [†] denotes the Moore-Penrose generalized inverse, then

$$U_i^{\dagger} b_i + P_{\ker(U_i)}[\mathcal{F}] \subseteq \mathcal{F} \text{ whenever } b_i \in U_i \mathcal{F}, \quad i \in \mathcal{H}.$$

$$(1.3)$$

These assumptions allow us to cover interesting applications. For example, consider the row sparse problem in multivariate regression [29, 30, 38]

$$\min_{X \in R^{\nu \times \gamma}} \|HX - B\|_F^2 + \lambda \|X\|_{\ell_a/\ell_2},$$
(1.4)

where $H \in \mathbb{R}^{\kappa \times \nu}, B \in \mathbb{R}^{\kappa \times \gamma}, \|\cdot\|_F$ is the Frobenius norm of a matrix,

$$||HX - B||_F^2 = \sum_{j=1}^{\gamma} \sum_{i=1}^{\kappa} (\sum_{\ell=1}^{\nu} H_{i\ell} X_{\ell,j} - B_{ij})^2 \text{ and } ||X||_{\ell_a/\ell_2} = \sum_{i=1}^{\nu} (\sum_{j=1}^{\gamma} X_{ij}^2)^{\frac{a}{2}}.$$

Let $n = \nu\gamma$, $\mathcal{F} = \mathbb{R}^n$, $b_i = 0$, $x = (x_{11}, x_{12}, \ldots, x_{\nu\gamma})^T \in \mathbb{R}^n$ and set $U_i \in \mathbb{R}^{\nu \times n}$ for $i \in \mathcal{N} = \{1, \ldots, \gamma\}$ be the projection whose entries are 0, or 1 such that $U_i x$ be the *i*th column of X and $U_i \in \mathbb{R}^{\gamma \times n}$ for $i \in \mathcal{H} = \{1, \ldots, \nu\}$ be the projection whose entries are 0, or 1 such that $U_i x$ be the *i*th row of X. Then problem (1.4) can be written in the form of (1.1). It is easy to see that the $\{U_i^T\}_{i\in\mathcal{N}}$ span \mathbb{R}^n . Hence, all assumptions mentioned above hold for problem (1.4).

Problem (1.1) encompasses the non-overlapping group sparse optimization problems. Let G_1, \ldots, G_m be subsets of $\{1, \ldots, n\}$ representing known groupings of the decision variable with size n_1, \ldots, n_m and $G_i \cap G_j = \emptyset, i \neq j$. In this case, problem (1.1) reduces to

$$\min_{x \in \mathcal{F}} f_1(x) + \lambda \sum_{i=1}^m \|U_i x\|^a,$$
(1.5)

where $f_1 : \mathbb{R}^n \to \mathbb{R}_+$ is a smooth loss function, $\lambda > 0$ is a positive number and $U_i \in \mathbb{R}^{n_i \times n}$ is defined in the following way

$$(U_i)_{kj} = \begin{cases} 1 & \text{if } j \in G_i \\ 0 & \text{otherwise} \end{cases} \quad \text{for} \quad k = 1, \dots, n_i.$$

Thus $U_i x = x_{G_i}$ is the *i*th group variable vector in \mathbb{R}^{n_i} with components $x_j, j \in G_i$. If $\mathcal{F} = \{x \mid \alpha_i \leq x_i \leq \beta_i, i = 1, ..., n\}$ with $\alpha_i < 0 < \beta_i$, then all assumptions mentioned above with $U_1 = I \in \mathbb{R}^{n \times n}$, for $1 \in \mathcal{N}$ hold for problem (1.5).

In problem (1.1), the decision variables have a group structure so that components in the same group tend to vanish simultaneously. Group sparse optimization problems have been

spherical harmonic representations of random fields on the sphere, $U_i x$ contains the coefficients of homogeneous harmonic polynomials of the degree *i*. Since $||U_i x||$ is rotationally invariant, the group sparse function $\sum_{i \in \mathcal{H}} ||U_i x||$ has been used in [31].

Note that problem (1.1) also covers some overlapping group sparse cases, provided (1.2) remains valid. A simple example is given by

$$f_1(U_1x) = [x_1^2 + x_2^2 + (x_3 - x_4)^2]^{\frac{a}{2}}$$
 and $f_2(U_2x) = [(x_3 + x_4)^2 + x_5^2)]^{\frac{a}{2}}$

with

$$U_1 = \begin{pmatrix} 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & -1 & 0 \end{pmatrix} \quad \text{and} \quad U_2 = \begin{pmatrix} 0 & 0 & 1 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{pmatrix}.$$

Problem (1.1) with $a \in (0, 1)$ and $n_i = 1, i \in \mathcal{H}$ has been studied in [5, 6, 14, 15, 16, 18]. Chen, Toint and Wang [17] show that an adaptive regularization algorithm using a p-th order Taylor model for p odd needs in general at most $O(\epsilon^{-(p+1)/p})$ evaluations of the objective function and its derivatives (at points where they are defined) to produce an ϵ -approximate first order stationary point. Since this complexity bound is identical in order to that already known for convexly constrained Lipschitzian minimization, the result in [17] shows that introducing non-Lipschitzian singularities in the objective function does not affect the worst-case evaluation complexity order.

The unconstrained optimization of smooth partially-separable was first considered in Griewank and Toint [28], studied by many researchers [25, 24, 13, 37, 19, 21] and extensively used in the popular CUTEst testing environment [26] as well as in the AMPL [23], LANCELOT [20] and FILTRANE [27] packages.

In problem (1.1), all these "element functions" f_i depend on $U_i x \in \mathbb{R}^{n_i}$ rather than on x, which is most useful when $n_i \ll n$. Letting

$$x_i = U_i x \in \mathbb{R}^{n_i} \text{ for } i \in \mathcal{M} \text{ and } f_{\mathcal{I}}(x) = \sum_{i \in \mathcal{I}} f_i(x) \text{ for any } \mathcal{I} \subseteq \mathcal{M},$$

we denote

$$f_{\mathcal{N}}(x) \stackrel{\text{def}}{=} \sum_{i \in \mathcal{N}} f_i(U_i x) = \sum_{i \in \mathcal{N}} f_i(x_i) \quad \text{and} \quad f_{\mathcal{H}}(x) \stackrel{\text{def}}{=} \sum_{i \in \mathcal{H}} f_i(U_i x) = \sum_{i \in \mathcal{H}} f_i(x_i).$$

The *p*-th degree Taylor series

$$T_{f_{\mathcal{N}},p}(x,s) = f_{\mathcal{N}}(x) + \sum_{j=1}^{p} \frac{1}{j!} \nabla_{x}^{j} f_{\mathcal{N}}(x)[s]^{j}, \text{ where } \nabla_{x}^{j} f_{\mathcal{N}}(x)[s]^{j} = \sum_{i \in \mathcal{N}} \nabla_{x_{i}}^{j} f_{i}(x_{i})[U_{i}s]^{j}, (1.6)$$

indicates that, for each j, only the $|\mathcal{N}|$ tensors $\{\nabla_{x_i}^j f_i(x_i)\}_{i\in\mathcal{N}}$ of dimension n_i^j needs to be computed and stored. Exploiting derivative tensors of order larger than 2 — and thus using the high-order Taylor series (1.6) as a local model of $f_{\mathcal{N}}(x+s)$ in the neighbourhood of x —

may therefore be practically feasible in our setting since n_i^j is typically orders of magnitude smaller than n. The same comment applies to $f_{\mathcal{H}}(x)$ whenever $||U_i x - b_i|| \neq 0$.

The main contribution of this paper is twofold.

- We propose a partially separable adaptive regularization algorithm with a *p*-th order Taylor model which uses the underlying rotational symmetry of the Euclidean norm function for $f_{\mathcal{H}}$ and the first *p* derivatives (whenever they exist) of the "element functions" f_i , for $i \in \mathcal{M}$.
- We show that the algorithm can produce an (ϵ, δ) -approximate q-th-order stationary point of problem (1.1) at most $O(\epsilon^{-(p+1)/(p-q+1)})$ evaluations of the objective function and its first p derivatives for any $q \in \{1, \ldots, p\}$.

Our results extend worst-case evaluation complexity bounds for smooth nonconvex optimization in [11, 12] which do not use the structure of partially separable functions and do not consider the non-Lipschitz continuity. Moreover, our results subsume the results for non-Lipschitz nonconvex optimization in [17] which only consider the complexity with q = 1 and $n_i = 1$ for $i \in \mathcal{H}$.

This paper is organized as follows. In Section 2, we define an (ϵ, δ) q-order necessary optimality conditions for local minimizers of problem (1.1). A Lipschitz continuous model to approximate f is proposed in Section 3. We then propose the partially separable adaptive regularization algorithm using the p-th order Taylor model in Section 4. In Section 5, we show that the algorithm produces an (ϵ, δ) -approximate q-th-order stationary point at most $O(\epsilon^{-(p+1)/(p-q+1)})$ evaluations of f and its first p derivatives.

We end this section by introducing notations used in the next four sections. **Notations.** For a symmetric tensor S of order p, $S[v]^p$ is the result of applying S to p copies of the vector v and

$$\|S\|_{[p]} \stackrel{\text{def}}{=} \max_{\|v\|=1} |S[v]^p| \tag{1.7}$$

is the associated induced norm for such tensors. If S_1 and S_2 are tensors, $S_1 \otimes S_2$ is their tensor product and $S_1^{k\otimes}$ is the product of S_1 with itself k times. For any set \mathcal{X} , $|\mathcal{X}|$ denotes its cardinality.

Because the notion of partial separability hinges on geometric interpretation of the problem, it is useful to introduce the various subspaces of interest for our analysis. We will extensively use the following definitions. As will become clear in Section 2, we will need to identify

$$\mathcal{C}(x,\epsilon) \stackrel{\text{def}}{=} \{ i \in \mathcal{H} \mid \|U_i x - b_i\| \le \epsilon \} \text{ and } \mathcal{A}(x,\epsilon) \stackrel{\text{def}}{=} \mathcal{H} \setminus \mathcal{C}(x,\epsilon),$$
(1.8)

the collection of hard elements which are close to non-Lipschitz continuity for a given x and its complement (the "active" elements), and

$$\mathcal{R}(x,\epsilon) \stackrel{\text{def}}{=} \bigcap_{i \in \mathcal{C}(x,\epsilon)} \ker(U_i) = \left[\underset{i \in \mathcal{C}(x,\epsilon)}{\operatorname{span}} (U_i^T) \right]^{\perp}$$
(1.9)

the subspace in which those nearly singular elements are invariant. (When $\mathcal{C}(x,\epsilon) = \emptyset$, we set $\mathcal{R}(x,\epsilon) = \mathbb{R}^n$.) For convenience, if $\epsilon = 0$, we denote $\mathcal{C}(x) \stackrel{\text{def}}{=} \mathcal{C}(x,0)$, $\mathcal{A}(x) \stackrel{\text{def}}{=} \mathcal{A}(x,0)$ and $\mathcal{R}(x) \stackrel{\text{def}}{=} \mathcal{R}(x,0)$. From these definitions, we have

$$U_i d = 0$$
, for $i \in \mathcal{C}(x)$, $d \in \mathcal{R}(x)$. (1.10)

Also denote by

$$\mathcal{R}_{\{i\}} \stackrel{\text{def}}{=} \operatorname{span}(U_i^T) \tag{1.11}$$

and observe that (1.2) implies that the $\mathcal{R}_{\{i\}}$ are orthogonal for $i \in \mathcal{H}$. Hence $\mathcal{R}_{\{i\}}$ is also the subspace in which all singular elements are invariant but the *i*-th. We also denote the "working" collection of elements from \mathcal{N} and those in \mathcal{H} not close to non-Lipschitz continuity by

$$\mathcal{W}(x,\epsilon) \stackrel{\text{def}}{=} \mathcal{N} \cup \mathcal{A}(x,\epsilon). \tag{1.12}$$

We denote $\mathcal{W}(x) \stackrel{\text{def}}{=} \mathcal{W}(x,0)$ when $\epsilon = 0$.

If $\{x_k\}$ is a sequence of iterates in \mathbb{R}^n , we also use the shorthands

$$C_k = C(x_k, \epsilon), \quad A_k = A(x_k, \epsilon), \quad \mathcal{R}_k = \mathcal{R}(x_k, \epsilon) \text{ and } \mathcal{W}_k = \mathcal{W}(x_k, \epsilon).$$
 (1.13)

We will make frequent use of

$$f_{\mathcal{W}_k}(x) \stackrel{\text{def}}{=} \sum_{i \in \mathcal{W}_k} f_i(x), \tag{1.14}$$

which is objective function "reduced" to the elements "away from non-Lipschitz continuity" at x_k .

For some $x, s \in \mathbb{R}^n$, we often use the notations $r_i = U_i x - b_i$ and $s_i = U_i s$.

2 Necessary optimality conditions

At variance with the theory developed in [17], which solely covers convergence to ϵ -approximate first-order stationary points, we now consider arbitrary orders of optimality. To this aim, we follow [11] and define, for a sufficiently smooth function $h : \mathbb{R}^n \to \mathbb{R}$ and a convex set $\mathcal{F} \subseteq \mathbb{R}^n$, the vector x to be an (ϵ, δ) -approximate q-th-order stationary point $(\epsilon > 0, \delta > 0,$ $q \in \{1, \ldots, p\}$) of $\min_{x \in \mathcal{F}} h(x)$ if, for some $\delta \in (0, 1]$

$$\phi_{h,q}^{\delta}(x) \le \epsilon \chi_q(\delta) \tag{2.1}$$

where

$$\phi_{h,q}^{\delta}(x) \stackrel{\text{def}}{=} h(x) - \min_{\substack{x+d \in \mathcal{F} \\ \|d\| \le \delta}} T_{h,q}(x,d), \tag{2.2}$$

and

$$\chi_q(\delta) \stackrel{\text{def}}{=} \sum_{\ell=1}^q \frac{\delta^\ell}{\ell!}.$$
(2.3)

In other words, we declare x to be an (ϵ, δ) -approximate q-th-order stationary point if the scaled maximal decrease that can be obtained on the q-th order Taylor series for h in a neighbourhood of x of radius δ is at most ϵ . We refer the reader to [11] for a detailed motivation and discussion of this measure. For our present purpose, it is enough to observe that $\phi_{h,q}^{\delta}(x)$ is a continuous function of x and δ for any q. Moreover, for q = 1 and q = 2, δ can be chosen equal to one and $\phi_{h,1}^1(x)$ and $\phi_{h,2}^1(x)$ are easy to compute. In the unconstrained case,

$$\phi_{h,1}^{1}(x) = \|\nabla_{x}^{1}h(x)\|$$

and computing $\phi_{h,2}^1$ reduces to solving the standard trust-region problem

$$\phi_{h,2}^{1}(x) = \left| \min_{\|d\| \le 1} \nabla_{x}^{1} h(x)[d] + \frac{1}{2} \nabla_{x}^{2} h(x)[d]^{2} \right|$$

In the constrained case,

$$\phi_{h,1}^1(x) = \left| \min_{\substack{x+d \in \mathcal{F} \\ \|d\| \le 1}} \nabla_x^1 h(x)[d] \right|,$$

which is the optimality measure used in [9] or [17] among others. However, given the potential difficulty of solving the global optimization problem in (2.2) for q > 2, our approach remains, for now, conceptual for such high optimality orders.

We now claim that we can extend the definition (2.1) to cover problem (1.1) as well. The key observation is that, by the definition of $\mathcal{W}(x,\epsilon)$ and $\mathcal{R}(x,\epsilon)$,

$$f_{\mathcal{A}(x,\epsilon)}(x) = f_{\mathcal{A}(x,\epsilon)}(x+d) \le f_{\mathcal{H}}(x+d) \le f_{\mathcal{A}(x,\epsilon)}(x+d) + \epsilon^a |\mathcal{H}| \text{ for all } d \in \mathcal{R}(x,\epsilon).$$
(2.4)

Note now that $f_{\mathcal{W}(x,\epsilon)}$ is smooth around x because it only contains elements which are away from non-Lipschitz continuity, and hence that $T_{f_{\mathcal{W}(x,\epsilon)},p}(x,s)$ is well-defined. We may therefore define x to be an (ϵ, δ) -approximate q-th-order stationary point for (1.1) if, for some $\delta \in (0, 1]$

$$\psi_{f,q}^{\epsilon,\delta}(x) \le \epsilon \chi_q(\delta), \tag{2.5}$$

where we define

$$\psi_{f,q}^{\epsilon,\delta}(x) \stackrel{\text{def}}{=} f(x) - \min_{\substack{x+d\in\mathcal{F}\\\|d\|\leq\delta,\,d\in\mathcal{R}(x,\epsilon)}} T_{f_{\mathcal{W}(x,\epsilon)},q}(x,d).$$
(2.6)

By the definition of $\mathcal{W}(x,\epsilon)$, we have $f_{\mathcal{W}(x,\epsilon)}(x) \leq f(x)$ and thus

$$f_{\mathcal{W}(x,\epsilon)}(x) - \min_{\substack{x+d\in\mathcal{F}\\ \|d\| \le \delta, d\in\mathcal{R}(x,\epsilon)}} T_{f_{\mathcal{W}(x,\epsilon)},q}(x,d) \le \psi_{f,q}^{\epsilon,\delta}(x).$$

Taking $\epsilon = 0$, x is q-th-order stationary point if $\psi_{f_{\mathcal{W}(x)},q}^{0,\delta} = 0$.

The optimality measure (2.5) may give the impression (in particular in its use of $\mathcal{R}(x, \epsilon)$) that the "singular" and "smooth" parts of the problem are merely separated, and that one could possibly apply the existing theory for smooth problems to the latter. Unfortunately, this is not true, because the "separation" implied by (2.5) does depend on ϵ , and one therefore needs to show that the complexity of minimizing the "non-singular" part does not explode (in particular with the unbounded growth of the Lispchitz constant) when ϵ tends to zero. Designing a suitable algorithm and proving an associated complexity result comparable to what is known for smooth problems is the main challenge in what follows.

Theorem 2.1 If x_* is a local minimizer of (1.1), then there is $\delta \in (0, 1]$ such that

$$\psi_{f,q}^{0,\delta}(x_*) = 0. \tag{2.7}$$

Proof. Suppose first that $\mathcal{R}(x_*) = \{0\}$ (which happens if there exists $x_* \in \mathcal{F}$ such that $f_{\mathcal{H}}(x_*) = 0$ and $\operatorname{span}_{i \in \mathcal{H}} \{U_i^T\} = \mathbb{R}^n$). Then (2.7) holds vacuously with any $\delta \in (0, 1]$. Now suppose that $\mathcal{R}(x_*) \neq \{0\}$. Let

$$\delta_1 = \min\left[1, \min_{i \in \mathcal{A}(x_*)} \|U_i x_* - b_i\|\right] \in (0, 1].$$

Since x_* is a local minimizer of (1.1), there exists $\delta_2 > 0$ such that

$$f(x_{*}) = \min_{\substack{x_{*}+d\in\mathcal{F}\\\|d\|\leq\delta_{2}}} f_{\mathcal{N}}(x_{*}+d) + \sum_{i\in\mathcal{H}} \|U_{i}(x_{*}+d) - b_{i}\|^{a}$$

$$\leq \min_{\substack{x_{*}+d\in\mathcal{F}\\\|d\|\leq\delta_{2}, d\in\mathcal{R}(x_{*})}} f_{\mathcal{N}}(x_{*}+d) + \sum_{i\in\mathcal{H}} \|U_{i}(x_{*}+d) - b_{i}\|^{a}$$

$$= \min_{\substack{x_{*}+d\in\mathcal{F}\\\|d\|\leq\delta_{2}, d\in\mathcal{R}(x_{*})}} f_{\mathcal{N}}(x_{*}+d) + \sum_{i\in\mathcal{A}(x_{*})} \|U_{i}(x_{*}+d) - b_{i}\|^{a}$$

$$= \min_{\substack{x_{*}+d\in\mathcal{F}\\\|d\|\leq\delta_{2}, d\in\mathcal{R}(x_{*})}} f_{\mathcal{W}(x_{*})}(x_{*}+d),$$

where we used (1.10) and (1.12) to derive the last two equalities, respectively. Now we consider the reduced problem

$$\min_{\substack{x_*+d\in\mathcal{F}\\\|d\|\leq\delta_2,\,d\in\mathcal{R}(x_*)}} f_{\mathcal{W}(x_*)}(x_*+d).$$
(2.8)

Since we have that

$$f_{\mathcal{W}(x_*)}(x_*) = f_{\mathcal{N}}(x_*) + \sum_{i \in \mathcal{A}(x_*)} \|U_i x_* - b_i\|^a = f_{\mathcal{N}}(x_*) + \sum_{i \in \mathcal{H}} \|U_i x_* - b_i\|^a = f(x_*)$$

we obtain that

$$f_{\mathcal{W}(x_{*})}(x_{*}) \leq \min_{\substack{x_{*}+d \in \mathcal{F} \\ \|d\| \leq \delta_{2}, d \in \mathcal{R}(x_{*})}} f_{\mathcal{W}(x_{*})}(x_{*}+d)$$

and x_* is a local minimizer of problem (2.8).

Note that there is $\delta_3 \in (0, \delta_1)$ such that for any $x_* + d$ in the ball $B(x_*, \delta_3)$, we have

$$||U_i(x_*+d) - b_i|| \ge ||U_ix_* - b_i|| - ||U_id|| \ge \delta_1 - ||U_i|| ||d|| = \delta_1 - \delta_3 > 0, \quad i \in \mathcal{A}(x_*).$$

Hence $f_{\mathcal{W}(x_*)}(x_* + d)$ is q-times continuously differentiable, and has Lipschitz continuous derivatives of orders 1 to q in $B(x^*, \delta_3)$. By Theorem 3.1 in [10], there is a $\delta \in (0, \min[\delta_2, \delta_3]]$, such that

$$\psi_{f_{\mathcal{W}(x_*)},q}^{0,\delta}(x_*) = f_{\mathcal{W}(x_*)}(x_*) - \min_{\substack{x_* + d \in \mathcal{F} \\ \|d\| \le \delta, \ d \in \mathcal{R}(x_*)}} T_{f_{\mathcal{W}(x_*)},q}(x_*,d) = 0.$$

This and the identity $f(x_*) = f_{\mathcal{W}(x_*)}(x_*)$, give the desired result (2.7).

We recall x_* a q-th-order stationary point of (1.1) if there is $\delta \in (0, 1]$ such that (2.7) holds.

Theorem 2.2 For each k, let x_k be an (ϵ_k, δ_k) -approximate q-th-order stationary point of (1.1) with $1 \ge \delta_k \ge \overline{\delta} > 0$ and $\epsilon_k \to 0$. Then any cluster point of $\{x_k\}$ is a q-th-order stationary point of (1.1).

Proof. Let x_* be a cluster point of $\{x_k\}$. Without loss of generality, we assume that $x_* = \lim_{k\to\infty} x_k$. From $0 < \chi_q(\delta) \le 2$ and $\psi_{f,q}^{\epsilon,\delta}(x) \ge 0$ for any $\delta \in (0,1)$, we have from (2.5) that $\lim_{k\to\infty} \psi_{f,q}^{\epsilon_k,\delta_k}(x_k) = 0$. We now need to prove that $\psi_{f,q}^{0,\overline{\delta}}(x_*) = 0$.

If $\mathcal{R}(x_*) = \{0\}$, (2.7) holds vacuously with any $\delta > 0$, and hence x_* is a *q*th-order-stationary point of (1.1). Suppose now that $\mathcal{R}(x_*) \neq \{0\}$. We first claim that there exists a $k_* \geq 0$ such that

$$\mathcal{C}(x_k, \epsilon_k) \subseteq \mathcal{C}(x_*) \quad \text{for} \quad k \ge k_*.$$
 (2.9)

To prove this inclusion, we choose k_* sufficiently large to ensure that

$$||x_k - x_*|| + \epsilon_k < \min_{j \in \mathcal{A}(x_*)} ||U_j x_* - b_j||, \text{ for } k \ge k_*.$$
 (2.10)

Such a k_* must exist, since the right-hand side of this inequality is strictly positive by definition of $\mathcal{A}(x_*)$. For an arbitrary $k \geq k_*$ and an index $i \in \mathcal{C}(x_k, \epsilon_k)$, using the definition of $\mathcal{C}(x, \epsilon)$, the identity $||U_i|| = 1$ and (2.10), we obtain that

$$||U_i x_* - b_i|| \le ||U_i (x_* - x_k)|| + ||U_i x_k - b_i|| \le ||x_* - x_k|| + \epsilon_k < \min_{j \in \mathcal{A}(x_*)} ||U_j x_* - b_i||.$$

This implies that $||U_i x_* - b_i|| = 0$ and $i \in \mathcal{C}(x_*)$. Hence (2.9) holds. By the definition of $\mathcal{R}(x, \epsilon)$ and $\mathcal{W}(x, \epsilon)$, (2.9) implies that, for all k,

$$\mathcal{R}(x_*) \subseteq \mathcal{R}(x_k, \epsilon_k) \quad \text{and} \quad \mathcal{W}(x_*) \subseteq \mathcal{W}(x_k, \epsilon_k).$$
 (2.11)

For any fixed $k \ge k_*$, consider now the following three minimization problems:

$$(A,k) \quad \begin{cases} \min_d & T_{f_{\mathcal{W}(x_k,\epsilon_k)},q}(x_k,d) \\ \text{s.t.} & x_k + d \in \mathcal{F}, \ d \in \mathcal{R}(x_k,\epsilon_k), \ \|d\| \le \delta_k, \end{cases}$$
(2.12)

$$(B,k) \quad \begin{cases} \min_d & T_{f_{\mathcal{W}(x_k,\epsilon_k)},q}(x_k,d) \\ \text{s.t.} & x_k + d \in \mathcal{F}, \ d \in \mathcal{R}(x_*), \ \|d\| \le \delta_k, \end{cases}$$
(2.13)

and

$$(C,k) \quad \begin{cases} \min_d & T_{f_{\mathcal{W}(x_*)},q}(x_k,d) \\ \text{s.t.} & x_k + d \in \mathcal{F}, \ d \in \mathcal{R}(x_*), \ \|d\| \le \delta_k. \end{cases}$$
(2.14)

Since d = 0 is a feasible point of these three problems, their minimum values, which we respectively denote by $\vartheta_{A,k}$, $\vartheta_{B,k}$ and $\vartheta_{C,k}$, are all smaller than $f(x_k)$. Moreover, it follows from the first part of (2.11) that, for each k,

$$\vartheta_{B,k} \ge \vartheta_{A,k}.\tag{2.15}$$

It also follows from (2.9) and (1.2) that

$$T_{f_{\mathcal{W}(x_{k},\epsilon_{k})},q}(x_{k},d) = T_{f_{\mathcal{W}(x_{*})},q}(x_{k},d) - f_{\mathcal{W}(x_{*})}(x_{k}) + f_{\mathcal{W}(x_{k},\epsilon_{k})}(x_{k}) \le T_{f_{\mathcal{W}(x_{*})},q}(x_{k},d) + |\mathcal{H}|\epsilon_{k}^{a}$$

for all $d \in \mathcal{R}(x_*)$, and thus (2.15) becomes

$$\vartheta_{A,k} \le \vartheta_{B,k} \le \vartheta_{C,k} + |\mathcal{H}|\epsilon_k^a \quad \text{for all } k \ge k_*.$$
(2.16)

The assumption that x_k is an (ϵ_k, δ_k) -approximate qth-order stationary point of (1.1) implies that

$$f(x_k) - \vartheta_{C,k} - |\mathcal{H}|\epsilon_k^a \le f(x_k) - \vartheta_{A,k} \le \epsilon_k \chi_q(\delta_k), \text{ for all } k \ge k_*.$$
(2.17)

Now (2.11) implies that $T_{f_{\mathcal{W}(x_*)},q}(x_k,d) \leq T_{f_{\mathcal{W}(x_k)},q}(x_k,d)$. Hence

$$f(x_k) - \vartheta_{C,k} = f(x_k) - \min_{\substack{x_k + d \in \mathcal{F} \\ \|d\| \le \delta_k, \ d \in \mathcal{R}(x_*)}} T_{f_{\mathcal{W}(x_*)},q}(x_k,d) \ge \psi_{f,q}^{0,\delta_k}(x_k)$$

As a consequence, (2.17) implies that

$$\psi_{f,q}^{0,\delta_k}(x_k) \le \epsilon_k \chi_q(\delta_k) + |\mathcal{H}|\epsilon_k^a.$$
(2.18)

In addition, the feasible sets of the three problems (2.12)-(2.14) are convex, and the objectives functions are polynomials with degree q. By the perturbation theory for optimization problems [21, Theorem 3.2.8], we can claim that

$$\lim_{k \to \infty} \vartheta_{C,k} = \min_{\substack{x_* + d \in \mathcal{F} \\ \|d\| \le \delta_*, \ d \in \mathcal{R}(x_*)}} T_{f_{\mathcal{W}(x_*)},q}(x_*,d), \tag{2.19}$$

where $\delta_* = \liminf_{k \to \infty} \delta_k \ge \overline{\delta}$. This implies that letting $k \to \infty$ in (2.18) gives

$$\psi_{f,q}^{0,\delta}(x_*) = f(x_*) - \min_{\substack{x_* + d \in \mathcal{F} \\ \|d\| \le \delta, d \in \mathcal{R}(x_*)}} T_{f_{\mathcal{W}(x_*)},q}(x_*,d) = 0.$$

3 A Lipschitz continuous model of $f_{\mathcal{W}_k}(x+s)$

Our minimization algorithm, described in the next section, involves the approximate minimization of a model $m(x_k, s)$ of $f_{\mathcal{W}_k}$ in the intersection of a neighbourhood of x_k with \mathcal{R}_k . This model, depending on function and derivatives values computed at x_k , should be able to predict values and derivatives of f at some neighbouring point $x_k + s$ reasonably accurately. This is potentially difficult if the current point happens to be near a point at which the function is not Lipschitz continuous.

Before describing our proposal, we need to state a useful technical result.

Lemma 3.1 Let a be a positive number and $r \neq 0$. Define, for a positive integer j,

$$\pi(a-j) \stackrel{\text{def}}{=} a \prod_{i=1}^{j-1} (a-i).$$
(3.1)

Then, if $\nabla^j_{\cdot} \|r\|^a$ is the value of the *j*-th derivative tensor of the function $\|\cdot\|^a$ with respect to its argument, evaluated at *r*, we have that,

$$\nabla^{j}_{\cdot} \|r\|^{a} = \sum_{i=1}^{j} \phi_{i,j} \|r\|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes}$$
(3.2)

for some scalars $\{\phi_{i,j}\}_{i=1}^j$ such that $\sum_{i=1}^j \phi_{i,j} = \pi(a-j)$, and that

$$\left\| \nabla_{\cdot}^{j} \| r \|^{a} \right\|_{[j]} = |\pi(a-j)| \| r \|^{a-j}.$$
(3.3)

Moreover, if β_1, β_2 are positive reals and ||r|| = 1, then

$$\left\| \nabla_{\cdot}^{j} \| \beta_{1} r \|^{a} - \nabla_{\cdot}^{j} \| \beta_{2} r \|^{a} \right\|_{[j]} = \left| \pi (a - j) \right| \left| \beta_{1}^{a - j} - \beta_{2}^{a - j} \right|.$$
(3.4)

Proof. See appendix.

Consider now the elements f_i for $i \in \mathcal{N}$. Each such element is p times continuously differentiable and, if we assume that its p-th derivative tensor $\nabla_x^p f_i$ is globally Lipschitz continuous with constant $L_i \geq 0$ in the sense that, for all $x_i, y_i \in \mathbb{R}^{n_i}$

$$\|\nabla_{x_i}^p f_i(x_i) - \nabla_{x_i}^p f_i(y_i)\|_{[p]} \le L_i \|x_i - y_i\|,$$
(3.5)

then it can be shown (see [11, Lemma 2.1]) that

$$f_i(x_i + s_i) = T_{f_i, p}(x_i, s_i) + \frac{1}{(p+1)!} \tau_i L_i \|s_i\|^{p+1} \quad \text{with} \quad |\tau_i| \le 1.$$
(3.6)

Because L_i in (3.6) is usually unknown in practice, it may not be possible to use (3.6) directly to model f_i in a neighbourhood of x. However, we may replace this term with an adaptive parameter σ_i , which yields the following (p + 1)-rst order model for the *i*-th smooth element f_i ,

$$m_i(x_i, s_i) = T_{f_i, p}(x_i, s_i) + \frac{1}{(p+1)!} \sigma_i ||s_i||^{p+1}, \quad (i \in \mathcal{N}).$$
(3.7)

Now we consider the elements f_i for $i \in \mathcal{H}$. Let $r_i = U_i x - b_i$. Using the associated Taylor's expansion would indeed ignore the non-Lipschitzian continuity occurring for $r_i = 0$ and this would restrict the validity of the model to a possibly very small neighbourhood of x_k whenever r_i is small for some $i \in \mathcal{A}(x_k, \epsilon)$. Our proposal is to use the underlying rotational symmetry of the Euclidean norm function to build a better Lipschtzian model. Suppose that $r_i \neq 0 \neq r_i + s_i$ and let

$$u_i = \frac{r_i}{\|r_i\|}, \quad r_i^+ = r_i(x+s) = r_i + s_i \text{ and } u_i^+ = \frac{r_i^+}{\|r_i^+\|}.$$
 (3.8)

Moreover, let R_i be the rotation in the (u_i, u_i^+) plane⁽¹⁾ such that

$$R_i u_i^+ = u_i. aga{3.9}$$

⁽¹⁾If $u_i = u_i^+$, $R_i = I$. If $n_i = 1$ and $r_i r_i^+ < 0$, this rotation is just the mapping from \mathbb{R}_+ to \mathbb{R}_- , defined by a simple sign change, as in the two-sided model of [17].

We observe that, given the isotropic nature of the Euclidean norm and the value $||r_i||^a$, the derivatives of $||r_i + s_i||^a$ with respect to s_i can be deduced from those $|| ||r_i||u_i^+||^a$. More precisely, for any $d \in \mathbb{R}^{n_i}$,

$$\left\| \|r_i\|u_i^+\|^a = \|r_i\|^a \text{ and } \nabla^{\ell}_{\cdot} \| \|r_i\|u_i^+\|^a[d]^{\ell} = \nabla^{\ell}_{\cdot} \|r_i\|^a[R_id]^{\ell}.$$
(3.10)

For example, when $\ell = 1$, one verifies that

$$\nabla^{1}_{\cdot} \|r_{i}\|^{a} [R_{i}d] = a \|r_{i}\|^{a-2} r_{i}^{T} R_{i}d$$

$$= a \|r_{i}\|^{a-2} \|r_{i}\| (R_{i}^{T} u_{i})^{T}d$$

$$= a \|\|r_{i}\|u_{i}^{+}\|^{a-2} (\|r_{i}\|u_{i}^{+})^{T}d$$

$$= \nabla^{1}_{\cdot} \|\|r_{i}\|u_{i}^{+}\|^{a} [d].$$

We may then choose to compute the Taylor's expansion for the function $\|\cdot\|^a$ around $\|r_i\|u_i^+$, that is $\|r_i^+\|^a = \|\|r_i^+\|u_i^+\|^a$

$$\begin{aligned} r_{i} \parallel^{\infty} &= \| \|r_{i}\| u_{i} \| \\ &= \| \|r_{i}\| u_{i}^{+} \|^{a} + \sum_{\ell=1}^{\infty} \frac{1}{\ell!} \nabla^{\ell} \| \|r_{i}\| u_{i}^{+} \|^{a} [(\|r_{i}^{+}\| - \|r_{i}\|) u_{i}^{+}]^{\ell} \\ &= \|r_{i}\|^{a} + \sum_{\ell=1}^{\infty} \frac{(\|r_{i}^{+}\| - \|r_{i}\|)^{\ell}}{\ell!} \nabla^{\ell} \|r_{i}\|^{a} [R_{i} u_{i}^{+}]^{\ell} \\ &= \|r_{i}\|^{a} + \sum_{\ell=1}^{\infty} \frac{(\|r_{i}^{+}\| - \|r_{i}\|)^{\ell}}{\ell!} \nabla^{\ell} \|r_{i}\|^{a} [u_{i}]^{\ell}. \end{aligned}$$

Using the expression (3.2) applied to ℓ copies of the unit vector u_i and the fact that $r_i^{j\otimes}[u_i]^j = (r_i^T u_i)^j = ||r_i||^j$ for all $j \in \mathbb{N}$, we now deduce that, for $\zeta_i = ||r_i + s_i|| - ||r_i|| \ge -||r_i||$,

$$||r_i + s_i||^a = ||r_i||^a + \sum_{\ell=1}^{\infty} \frac{\pi(a-\ell)}{\ell!} \zeta_i^{\ell} ||r_i||^{a-\ell},$$
(3.11)

which is nothing else than the Taylor's expansion of $||r_i + \zeta_i u_i||^a$ (or, equivalently, of $||||r_i||u_i^+ + \zeta_i u_i^+||^a)$ expressed as a function of the scalar variable $\zeta_i \geq -||r_i||$. As can be expected from the isotropic nature of the Euclidean norm, the value of $||r_i^+||^a$ (and of its derivatives after a suitable rotation) only depend(s) on the distance of r_i^+ to the non-Lipschitz continuity at zero. Thus, limiting the development (3.11) to degree p (as in (3.7)), it is natural to define

$$\mu(\|r_i\|,\zeta_i) \stackrel{\text{def}}{=} \|r_i\|^a + \sum_{\ell=1}^p \frac{\pi(a-\ell)}{\ell!} \zeta_i^\ell \|r_i\|^{a-\ell}, \quad (i \in \mathcal{A}(x,\epsilon)),$$
(3.12)

which is a unidimensional model of $||r_i + \zeta_i u_i||^a$ based of the residual value $||r_i||$. Note that $\mu(||r_i||, \zeta_i)$ is Lipschitz continuous as a function of ζ_i as long as $||r_i||$ remains uniformly bounded away from zero, with a Lipschitz constant depending on the lower bound. We then define the isotropic model

$$m_i(x_i, s_i) \stackrel{\text{def}}{=} \mu(\|r_i\|, \zeta_i) = \mu_i(\|r_i\|, \|r_i + s_i\| - \|r_i\|) \quad \text{for} \quad i \in \mathcal{A}(x, \epsilon),$$
(3.13)

so that, abusing notation slightly,

$$m_i(x_i, s_i) = T_{m_i, p}(x_i, s_i) \stackrel{\text{def}}{=} ||r_i||^a + \sum_{\ell=1}^p \frac{\pi(a-\ell)}{\ell!} \zeta_i^\ell ||r_i||^{a-\ell}, \quad (i \in \mathcal{A}(x, \epsilon)).$$

We now state some useful properties of this model.

Lemma 3.2 Suppose that p is odd and that $\mathcal{A}(x) \neq \emptyset$ for some $x \in \mathcal{F}$. Then, for $i \in \mathcal{A}(x)$,

$$m_i(x_i, s_i) \ge \|r_i + s_i\|^a,$$
(3.14)

and, whenever $||r_i + s_i|| \le ||r_i(x)||$,

$$\nabla_{\zeta}^{\ell}\mu(\|r_i\|,\zeta_i)) \ge \nabla_{\zeta}^{\ell}\mu(\|r_i\|,0) = \pi(a-\ell)\|r_i\|^{a-\ell} \quad (\ell \ odd), \tag{3.15}$$

and

$$\nabla_{\zeta}^{\ell}\mu(\|r_i\|,\zeta_i) \le \nabla_{\zeta}^{\ell}\mu(\|r_i\|,0) = \pi(a-\ell)\|r_i\|^{a-\ell} \quad (\ell \ even).$$
(3.16)

As a consequence, $m_i(x_i, \zeta_i)$ is a concave function of ζ_i on the interval $[-||r_i||, 0]$.

Proof. Let $i \in \mathcal{A}(x)$. From the mean-value theorem and (3.11), we have that, for some $\nu \in (0, 1)$,

$$\|r_i + s_i\|^a = \|r_i\|^a + \sum_{\ell=1}^p \frac{\pi(a-\ell)}{\ell!} \zeta_i^\ell \|r_i\|^{a-\ell} + \frac{\pi(a-p-1)}{(p+1)!} \zeta_i^{p+1} \|r_i + \nu\zeta_i u_i\|^{a-p-1}.$$
 (3.17)

Since p is odd, we obtain that $\pi(a-p-1) < 0$ and $\zeta_i^{p+1} \ge 0$. Thus (3.14) directly results from (3.17), (3.12) and (3.13). Now (3.12) and (3.13) together imply that

$$\nabla_{\zeta}^{\ell}\mu(\|r_i\|,\zeta_i) = \nabla_{\zeta}^{\ell}\mu(\|r_i\|,0) + \sum_{j=\ell+1}^{p} \frac{\pi(a-j)}{(j-\ell)!} \zeta_i^{j-\ell} \|r_i\|^{a-j}$$
(3.18)

for $\zeta_i = \|r_i(x) + s_i\| - \|r_i(x)\| \le 0$. Suppose first that ℓ is odd. Then we have that $\pi(a-j)$ is negative for even j, that is exactly when $\zeta_i^{j-\ell}$ is non-positive. Hence every term in the sum of the right-hand side of (3.18) is non-negative and (3.15) follows. Suppose now that ℓ is even. Then $\pi(a-j)$ is negative for odd j, which is exactly when $\zeta_i^{j-\ell}$ is non-negative. Hence every term in the sum of the right-hand side of (3.18) is non-positive and (3.16) follows. The last conclusion of the lemma is then deduced by considering $\ell = 2$ in (3.16) and observing that $\pi(a-\ell)\|r_i(x)\|^{a-\ell} = a(a-1)\|r_i(x)\|^{a-2} < 0$.

Thus the isotropic model $m_i(x_i, s_i)$ overestimates the true function $||r_i(x) + s_i||^a$ and correctly reflects its concavity in the direction of its non-Lipschitz continuity. But $m_i(x_i, s_i) = \mu(||r_i||, \zeta_i)$ is now Lipschitz continuous as a function of s_i , while $||r_i(x) + s_i||^a$ is not.

Combining (3.7) and (3.13) now allows us to define a model for the complete f on $\mathcal{R}(x,\epsilon)$ as

$$m(x,s) \stackrel{\text{def}}{=} \sum_{i \in \mathcal{W}(x,\epsilon)} m_i(x_i, s_i).$$
(3.19)

Since $r_i(x) \neq 0$ for $i \in \mathcal{A}(x, \epsilon)$, this model is in turn well defined.

We conclude this section by observing that writing the problem in the partially-separable form (1.1) is the key to expose the singular parts of the objective function, which then allows exploiting their rotational symmetry.

4 The adaptive regularization algorithm

Having defined a model of f, we may then use this model within a regularization minimization method inspired from the ARp algorithm in [11]. In such an algorithm, the step from an iterate x_k is obtained by attempting to (approximately) minimize the model ((3.19) in our case). If an (ϵ, δ) -approximate q-th-order-stationary point is sought, this minimization is terminated as soon as the step s_k is long enough, in that

$$\|s_k\| \ge \varpi \epsilon^{\frac{1}{p-q+1}} \tag{4.1}$$

for some constant $\varpi \in (0, 1]$, or as soon as the trial point $x_k + s_k$ is an approximate q-thorder-stationary point of the model, in the sense that

$$\psi_{m,q}^{\epsilon,\delta_k}(x_k,s_k) \le \min\left[\frac{\theta \|s_k\|^{p-q+1}}{(p-q+1)!}, a\min_{i\in\mathcal{A}(x_k+s_k,\epsilon)} \|r_i(x_k+s_k)\|\right] \chi_q(\delta_k)$$
(4.2)

for some $\theta, \delta_k \in (0, 1]$, where $\psi_{m,q}^{\epsilon, \delta_k}(x_k + s_k)$ is the optimality measure (2.6) computed for the model m(x, s), that is

$$\psi_{m,q}^{\epsilon,\delta}(x,s) \stackrel{\text{def}}{=} m(x,s) - \min_{\substack{x+s+d\in\mathcal{F}\\ \|d\| \le \delta, d\in\mathcal{R}(x,\epsilon)}} T_{m,q}(x,s+d).$$
(4.3)

When q = 1, $\delta_k = 1$ and the feasible set $\mathcal{F} = \mathbb{R}^n$, we have $\chi_q(\delta_k) = 1$, and (4.2) holds if $\theta \|s_k\| \ge a\epsilon$ and

$$\psi_{m,1}^{\epsilon,1}(x,s_k) = \|\nabla f_{\mathcal{N}}(x) + \nabla f_{\mathcal{A}(x,\epsilon)}(x)\| \le a\epsilon.$$

In view of the optimality condition (2.5), we also require that, if $||r_i(x+s)|| \leq \epsilon$ occurs for some $i \in \mathcal{H}$ in the course of the model minimization, the value of $r_i(x+s)$ is fixed, implying that the remaining minimization is carried out on $\mathcal{R}(x_k + s, \epsilon)$. As a consequence, the dimension of $\mathcal{R}(x_k + s, \epsilon)$ (and thus of \mathcal{R}_k) is monotonically non-increasing during the step computation and across all iterations. It was shown in [11, Lemma 2.5] that, unless x_k is an $(\epsilon, 1)$ -approximate *p*-th-order-stationary point (which is obviously enough for the whole algorithm to terminate), a step satisfying (4.2) can always be found. The fact that this condition must hold on a subspace of potentially diminishing dimension clearly does not affect the result, and indicates that (4.2) is well-defined. This model minimization is in principle simpler than the original problem because the general nonlinear f_i have been replaced by locally accurate polynomial approximations and also because the model is now Lipschitz continuous, albeit still non-smooth. Importantly, the model minimization does not involve any evaluation of the objective function or its derivatives, and model evaluations within this calculation therefore do not affect the overall evaluation complexity of the algorithm.

We now introduce some useful notation for describing our algorithm. Define

$$x_{i,k} \stackrel{\text{def}}{=} U_i x_k, \quad r_{i,k} \stackrel{\text{def}}{=} U_i x_k - b_i, \quad s_{i,k} \stackrel{\text{def}}{=} U_i s_k, \quad u_{i,k} \stackrel{\text{def}}{=} \frac{r_{i,k}}{\|r_{i,k}\|}$$

and

$$\mathcal{A}_{k}^{+} \stackrel{\text{def}}{=} \mathcal{A}(x_{k} + s_{k}, \epsilon), \quad \mathcal{R}_{k}^{+} \stackrel{\text{def}}{=} \mathcal{R}(x_{k} + s_{k}, \epsilon) \text{ and } \mathcal{W}_{k}^{+} \stackrel{\text{def}}{=} \mathcal{W}(x_{k} + s_{k}, \epsilon).$$

Also let

$$\Delta f_{i,k} \stackrel{\text{def}}{=} f_i(x_{i,k}) - f_i(x_{i,k} + s_{i,k}), \quad \Delta f_k \stackrel{\text{def}}{=} f_{\mathcal{W}_k^+}(x_k) - f_{\mathcal{W}_k^+}(x_k + s_k) = \sum_{i \in \mathcal{W}_k^+} \Delta f_{i,k},$$

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$$\Delta m_{i,k} \stackrel{\text{def}}{=} m_i(x_{i,k}, 0) - m_i(x_{i,k}, s_{i,k}), \quad \Delta m_k = \sum_{i \in \mathcal{W}_k^+} \Delta m_{i,k},$$

and

$$\Delta T_{k} \stackrel{\text{def}}{=} T_{f_{\mathcal{W}_{k}^{+}, p}}(x_{k}, 0) - T_{f_{\mathcal{W}_{k}^{+}, p}}(x_{k}, s_{k})$$

$$= [T_{f_{\mathcal{N}, p}}(x_{k}, 0) - T_{f_{\mathcal{N}, p}}(x_{k}, s_{k})] + [m_{\mathcal{A}_{k}^{+}}(x_{k}, 0) - m_{\mathcal{A}_{k}^{+}}(x_{k}, s_{k})]$$

$$= \Delta m_{k} + \frac{1}{(p+1)!} \sum_{i \in \mathcal{N}} \sigma_{i, k} \|s_{i, k}\|^{p+1}.$$
(4.4)

Our partially-separable adaptive regularization degree-p algorithm PSARp is then given by Algorithm 4.1 on the following page.

Note that an $x_0 \in \mathcal{F}$ can always be computed by projecting an infeasible starting point onto \mathcal{F} . The motivation for the second and third parts of (4.9) and (4.10) is to identify cases where the isotropic model m_i overestimates the element function f_i to an excessive extent, leaving some room for reducing the regularization and hence allowing longer steps. The requirement that $\rho_k \geq \eta$ in both (4.9) and (4.10) is intended to prevent a situation where a particular regularization parameter is increased and another decreased at a given unsuccessful iteration, followed by the opposite situation at the next iteration, potentially leading to cycling.

It is worthwhile noting the differences between the PSARp algorithm and the algorithm discussed in [17]. The first and most important is that the new algorithm is intended to find an (ϵ, δ) -approximate q-th-order stationary point for problem (1.1), rather than a first-order stationary point. This is made possible by using the q-th-order termination criterion (4.5) instead of a criterion only involving the first-order model decrease, and by simultaneously using the step termination criteria (4.1) and (4.2) which again replace a simpler version based solely on first-order information. The second is that the PSARp algorithm applies to the more general problem (1.1), in particular using the isotropic model (3.12) to allow $n_i > 1$ for $i \in \mathcal{H}$.

As alluded to above and discussed in [12] and [4], the potential termination of the algorithm in Step 2 can only happen whenever q > 2 and $x_k = x_{\epsilon}$ is an $(\epsilon, 1)$ -approximate *p*-th-orderstationary point within \mathcal{R}_k , which, together with (2.5), imply that the same property holds for problem (1.1). This is a significantly stronger optimality condition than what is required by (4.5). Also note that the potentially costly calculation of (4.2) may be avoided if (4.1) holds.

Let the index set of the "successful" and "unsuccessful" iterations be given by

$$\mathcal{S} \stackrel{\text{def}}{=} \{k \ge 0 \mid \rho_k \ge \eta\} \text{ and } \mathcal{U} \stackrel{\text{def}}{=} \{k \ge 0 \mid \rho_k < \eta\}.$$

Also define

$$\mathcal{S}_k \stackrel{\text{def}}{=} \mathcal{S} \cap \{0, \dots, k\} \text{ and } \mathcal{U}_k \stackrel{\text{def}}{=} \{0, \dots, k\} \setminus \mathcal{S}_k$$

We then state a bound on $|\mathcal{U}_k|$ as a function of $|\mathcal{S}_k|$. This is a standard result for nonpartially-separable problems (see [7, Theorem 2.4] for instance), but needs careful handling of the model's overestimation properties to apply to our present context.

Algorithm 4.1: Partially-Separable Adaptive Regularization (PSAR_p)

Step 0: Initialization: $x_0 \in \mathcal{F}$ and $\{\sigma_{i,0}\}_{i\in\mathcal{N}} > 0$ are given as well as the accuracy $\epsilon \in (0,1]$ and constants $0 < \gamma_0 < 1 < \gamma_1 \leq \gamma_2, \eta \in (0,1), \theta \geq 0, \delta_{-1} = 1, \sigma_{\min} \in (0, \min_{i\in\mathcal{N}} \sigma_{i,0}]$ and $\kappa_{\text{big}} > 1$. Set k = 0.

Step 1: Termination: Evaluate $f(x_k)$ and $\{\nabla_x^j f_{\mathcal{W}_k}(x_k)\}_{j=1}^q$. If

$$\psi_{f,q}^{\epsilon,\delta_{k-1}}(x_k) \le \epsilon \chi_q(\delta_{k-1}) \tag{4.5}$$

return $x^{\epsilon} = x_k$ and terminate. Otherwise evaluate $\{\nabla_x^j f_{\mathcal{W}_k}(x_k)\}_{j=q+1}^p$.

- Step 2: Step computation: Attempt to compute a step $s_k \in \mathcal{R}_k$ such that $x_k + s_k \in \mathcal{F}$, $m(x_k, s_k) < m(x_k, 0)$ and either (4.1) holds or (4.2) holds for some $\delta_k \in (0, 1]$. If no such step exists, return $x^{\epsilon} = x_k$ and terminate.
- Step 3: Step acceptance: Compute

$$\rho_k = \frac{\Delta f_k}{\Delta T_k} \tag{4.6}$$

and set $x_{k+1} = x_k$ if $\rho_k < \eta$, or $x_{k+1} = x_k + s_k$ if $\rho_k \ge \eta$.

Step 4: Update the "nice" regularization parameters: For $i \in \mathcal{N}$, if

$$f_i(x_{i,k} + s_{i,k}) > m_i(x_{i,k}, s_{i,k}) \tag{4.7}$$

 set

$$\sigma_{i,k+1} \in [\gamma_1 \sigma_{i,k}, \gamma_2 \sigma_{i,k}]. \tag{4.8}$$

Otherwise, if either

$$\rho_k \ge \eta \quad \text{and} \quad \Delta f_{i,k} \le 0 \quad \text{and} \quad \Delta f_{i,k} < \Delta m_{i,k} - \kappa_{\text{big}} |\Delta f_k| \tag{4.9}$$

or

$$\rho_k \ge \eta \quad \text{and} \quad \Delta f_{i,k} > 0 \quad \text{and} \quad \Delta f_{i,k} > \Delta m_{i,k} + \kappa_{\text{big}} |\Delta f_k| \tag{4.10}$$

then set

$$\sigma_{i,k+1} \in [\max[\sigma_{\min}, \gamma_0 \sigma_{i,k}], \sigma_{i,k}], \tag{4.11}$$

else set

$$\sigma_{i,k+1} = \sigma_{i,k}.\tag{4.12}$$

Increment k by one and go to Step 1.

Lemma 4.1 Suppose that AS.2 and AS.3 hold and that $\sigma_{i,k} \leq \sigma_{\max}$ for all $i \in \mathcal{M}$ and all $k \geq 0$. Then, for all $k \geq 0$, $k \leq \kappa^a |\mathcal{S}_k| + \kappa^b$,

where

$$\kappa^{a} \stackrel{\text{def}}{=} 1 + \frac{|\mathcal{N}| |\log \gamma_{0}|}{\log \gamma_{1}} \quad and \quad \kappa^{b} \stackrel{\text{def}}{=} \frac{|\mathcal{N}|}{\log \gamma_{1}} \log \left(\frac{\sigma_{\max}}{\sigma_{\min}}\right).$$

Proof. Note that ρ_k is defined by Δf_k and ΔT_k which are computed on $\mathcal{W}_k^+ = \mathcal{N} \cup \mathcal{A}_k^+$. Following the proof of [17, Lemma 4.11] and (3.14), we have $\Delta f_k \geq \Delta m_k$ if (4.7) fails for all $i \in \mathcal{N}$. Hence, from (4.6), we have $\rho_k \geq 1 > \eta$ and $k \in \mathcal{S}_k$. If $k \in \mathcal{U}_k$, $\sigma_{i,k}$ is increased with (4.8) for at least one $i \in \mathcal{N}$. Let $\mathcal{J}_{i,k}$ be the set of iterations where $\sigma_{i,j}$ $(j \leq k)$ is increased. Then $|\mathcal{U}_k| \leq |\mathcal{N}| \max_{i \in \mathcal{N}} |\mathcal{J}_{i,k}|$. Using $k = |\mathcal{S}_k| + |\mathcal{U}_k| - 1$ completes the proof.

5 Evaluation complexity analysis

We are now ready for a formal analysis of the evaluation complexity of the PSARp algorithm for problem (1.1), under the following assumptions.

AS.1The feasible set
$$\mathcal{F}$$
 is closed, convex, non-empty and kernel-centered (in the sense of (1.3)).AS.2Each element function f_i ($i \in \mathcal{N}$) is p times continuously differentiable in an open set containing \mathcal{F} , where p is odd whenever $\mathcal{H} \neq \emptyset$.AS.3The p -th derivative of each f_i ($i \in \mathcal{N}$) is Lipschitz continuous on \mathcal{F} with associated Lipschitz constant L_i (in the sense of (3.5)).AS.4There exists a constant f_{low} such that $f_{\mathcal{N}}(x) \geq f_{\text{low}}$ for all $x \in \mathcal{F}$.AS.5If $\mathcal{H} \neq \emptyset$, there exists a constant $\kappa_{\mathcal{N}} \geq 0$ such that $\|\nabla_x^j f_i(U_i x)\| \leq \kappa_{\mathcal{N}}$ for all $x \in \mathcal{V}, i \in \mathcal{N}$ and $j \in \{1, \dots, p\}$, where

$$\mathcal{V} \stackrel{\text{def}}{=} \left\{ x \in \mathcal{F} \mid \text{ there exists } i \in \mathcal{H} \text{ with } \|r_i(x)\| \le \frac{a}{16} \right\}.$$
(5.1)

Note that AS.4 is necessary for problem (1.1) to be well-defined. Also note that, because of AS.2, AS.5 automatically holds if \mathcal{F} or \mathcal{V} are bounded or if the iterates $\{x_k\}$ remain in a bounded set. It is possible to weaken AS.2 and AS.3 by replacing \mathcal{F} with the level set $\mathcal{L} = \{x \in \mathcal{F} \mid f(x) \leq f(x_0)\}$ without affecting the results below. Finally observe that \mathcal{V} need not to be bounded, in particular if $\operatorname{span}_{i \in \mathcal{H}}(U_i)$ is a proper subspace of \mathbb{R}^n . The motivation for the particular choice of $\frac{1}{16}a$ in (5.1) will become clear in Lemma 5.5 below.

We first recall a result providing useful bounds.

Lemma 5.1 There exist a constant $\varsigma > 0$ such that, for all $s \in \mathbb{R}^m$ and all $v \ge 1$,

$$\varsigma^{v} \|s\|^{v} \le \sum_{i \in \mathcal{N}} \|s_{i}\|^{v} \le |\mathcal{N}| \|s\|^{v}.$$

$$(5.2)$$

Proof. See [17, Lemma 4.1].

This lemma states that $\sum_{i \in \mathcal{N}} \|\cdot\|$ is a norm on \mathbb{R}^n whose equivalence constants with respect to the Euclidean one are ς and $|\mathcal{N}|$.

Our next step is to specify under which conditions the standard ϵ -independent overestimation and derivative accuracy bounds typical of the Lipschitz case (see [11, Lemma 2.1] for instance) can be obtained for the elements functions of (1.1). We define, for a given $k \ge 0$ and a given constant $\mu > 0$ independent of ϵ ,

$$\mathcal{O}_{k,\mu} \stackrel{\text{def}}{=} \{ i \in \mathcal{A}_k^+ \mid \min[\|r_{i,k}\|, \|r_{i,k} + s_{i,k}\|] \ge \mu \}.$$
(5.3)

Observe that if, for some $i \in \mathcal{H}$ and $b_i \notin U_i \mathcal{F}$, then both $||r_{i,k}||$ and $||r_{i,k} + s_{i,k}||$ are bounded away from zero, so $i \in \mathcal{O}_{k,\mu}$ for all k and all μ such that $\mu \leq \min_{x \in \mathcal{F}} ||U_i x - b_i||$. Thus we assume, without loss of generality, that

$$b_i \in U_i \mathcal{F}$$
 for all $i \in \mathcal{H}$. (5.4)

We then obtain the following crucial error bounds.

Lemma 5.2 Suppose that AS.2 and AS.3 hold. Then, for $k \ge 0$ and $L_{\max} \stackrel{\text{def}}{=} \max_{i \in \mathcal{N}} L_i$,

$$f_i(x_{i,k} + s_{i,k}) = m_i(x_{i,k}, s_{i,k}) + \frac{1}{(p+1)!} \Big[\tau_{i,k} L_{\max} - \sigma_{i,k} \Big] \|s_{i,k}\|^{p+1} \quad with \quad |\tau_{i,k}| \le 1, \quad (5.5)$$

for all $i \in \mathcal{N}$. If, in addition, $\mu > 0$ is given and independent of ϵ , then there exists a constant $L(\mu)$ independent of ϵ such that, for $\ell \in \{1, \ldots, p\}$,

$$\|\nabla_{x}^{\ell} f_{\mathcal{N}\cup\mathcal{O}_{k,\mu}}(x_{k}+s_{k}) - \nabla_{s}^{\ell} T_{f_{\mathcal{N}\cup\mathcal{O}_{k,\mu}},p}(x_{k},s_{k})\| \leq \frac{L(\mu)}{(p-\ell+1)!} \|s_{k}\|^{p-\ell+1}.$$
 (5.6)

Proof. First note that, if f_i has a Lipschitz continuous *p*-th derivative as a function of $x_i = U_i x$, then (1.6) shows that it also has a Lipschitz continuous *p*-th derivative as a function of *x*. It is therefore enough to consider the element functions as functions of x_i .

Observe now that, for each k and $i \in \mathcal{N}$, AS.2 and AS.3 ensure (5.5), and the inequality

$$\|\nabla_{x_i}^{\ell} f_i(x_{i,k} + s_{i,k}) - \nabla_{s_i}^{\ell} T_{f_i,p}(x_{i,k}, s_{i,k})\| \le \frac{L_i}{(p - \ell + 1)!} \|s_{i,k}\|^{p - \ell + 1}$$
(5.7)

immediately follows from the known bounds for p times continuously differentiable functions with Lipschitz continuous p-th derivative (see [11, Lemma 2.1]). Consider now $i \in \mathcal{O}_{k,\mu}$ for some k and some fixed $\mu > 0$, implying that $\min[||r_{i,k}||, ||r_{i,k} + s_{i,k}||] \ge \mu > 0$. Then

$$\nabla^{\ell}_{\cdot} \|r_{i,k} + s_{i,k}\|^{a} [d]^{\ell} = \nabla^{\ell}_{\cdot} \|\|r_{i,k} + s_{i,k}\|u^{+}_{i,k}\|^{a} [d]^{\ell} = \nabla^{\ell}_{\cdot} \|\|r_{i,k} + s_{i,k}\|u_{i,k}\|^{a} [R_{i,k}d]^{\ell}$$
(5.8)

where $R_{i,k}$ is the rotation such that $R_{i,k}u_{i,k}^+ = u_{i,k}$. We also have from (3.10) with x replaced by $x_k + s_k$ that

$$\nabla_{s_i}^{\ell} T_{m_i,p}(x_{i,k}, s_{i,k})[d]^{\ell} = \nabla_{\cdot}^{\ell} \left\| \|r_{i,k}\| u_{i,k} \right\|^a [R_{i,k}d]^{\ell}.$$
(5.9)

Taking the difference between (5.8) and (5.9), we obtain, successively using the definition of the tensor norm, the fact that $R_{i,k}$ is orthonormal and (3.4) in Lemma 3.1, that

$$\begin{split} \left\| \nabla^{\ell}_{\cdot} \| r_{i,k} + s_{i,k} \|^{a} - \nabla^{\ell}_{s_{i}} T_{m_{i},p}(x_{i,k}, s_{i,k}) \right\|_{[\ell]} \\ &= \max_{\||d\|=1} \left| \nabla^{\ell}_{\cdot} \| r_{i,k} + s_{i,k} \|^{a} [d]^{\ell} - \nabla^{\ell}_{s_{i}} T_{m_{i},p}(x_{i,k}, s_{i,k}) [d]^{\ell} \right| \\ &= \max_{\|d\|=1} \left| \nabla^{\ell}_{\cdot} \| \| r_{i,k} + s_{i,k} \| u_{i,k} \|^{a} [R_{i,k}d]^{\ell} - \nabla^{\ell}_{\cdot} \| \| r_{i,k} \| u_{i,k} \|^{a} [R_{i,k}d]^{\ell} \right| \\ &= \left\| \nabla^{\ell}_{\cdot} \| \| r_{i,k} + s_{i,k} \| u_{i,k} \|^{a} - \nabla^{\ell}_{\cdot} \| \| r_{i,k} \| u_{i,k} \|^{a} \|_{[\ell]} \\ &= \left| \pi (a - \ell) \right| \left\| r_{i,k} + s_{i,k} \|^{a - \ell} - \| r_{i,k} \|^{a - \ell} \right|. \end{split}$$

Now the univariate function $\nu(t) \stackrel{\text{def}}{=} t^a$ is (more than) p+1 times continuously differentiable with bounded (p+1)-rst derivative on the interval $[t_1, t_2]$ and thus, from Lemma 5.2, we have that

$$\pi(a-\ell)\left|t_1^{a-\ell} - t_2^{a-\ell}\right| = \left|\frac{d^\ell\nu}{dt^\ell}(t_1) - \frac{d^\ell\nu}{dt^\ell}(t_2)\right| \le \frac{L_\nu}{(p-\ell+1)!}|t_1 - t_2|^{p-\ell+1},$$

where L_{ν} is the upper bound on the (p+1)-rst derivative of $\nu(t)$ on interval $[t_1, t_2]$, that is $L_{\nu} = |\pi(a-p-1)| \min[t_1, t_2]^{a-p-1}$. As a consequence, we obtain that

$$\left\|\nabla_{\cdot}^{\ell}\|r_{i,k}+s_{i,k}\|^{a}-\nabla_{s_{i}}^{\ell}T_{m_{i},p}(x_{i,k},s_{i,k})\right\|_{[\ell]} \leq \frac{L(\mu)}{(p-\ell+1)!}\left|\left\|r_{i,k}+s_{i,k}\right\|-\left\|r_{i,k}\right\|\right|^{p-\ell+1},$$

where we use the fact that $\min[||r_{i,k}||, ||r_{i,k} + s_{i,k}||] \ge \mu$ to define

$$L(\mu) \stackrel{\text{def}}{=} \max \left| \pi(a-p-1) | \mu^{a-p-1}, L_{\max} \right].$$

We then observe that $||s_{i,k}|| = ||r_{i,k} + s_{i,k} - r_{i,k}|| \ge ||r_{i,k} + s_{i,k}|| - ||r_{i,k}|||$ which finally yields that

$$\left\|\nabla^{\ell}_{\cdot} \|r_{i,k} + s_{i,k}\|^{a} - \nabla^{\ell}_{s_{i}} T_{m_{i},p}(x_{i,k}, s_{i,k})\right\|_{[\ell]} \leq \frac{L(\mu)}{(p-\ell+1)!} \|s_{i,k}\|^{p-\ell+1}$$

Combining this last inequality with (5.7) and the fact that $\nabla_{x_i}^{\ell} \|r_{i,k} + s_{i,k}\|^a = \nabla_{\cdot}^{\ell} \|r_{i,k} + s_{i,k}\|^a$ then ensures that (5.6) holds.

Observe that the Lipschitz constant L is independent of ϕ whenever $\mathcal{H} = \emptyset$. Our model definition also implies the following bound.

Lemma 5.3 For all $k \ge 0$ before termination, $s_k \ne 0$, (4.6) is well-defined and

$$\Delta T_k \ge \frac{\sigma_{\min} \varsigma^{p+1}}{(p+1)!} \, \|s_k\|^{p+1}.$$
(5.10)

Proof. We immediately deduce that

$$\Delta T_k \ge \frac{\sigma_{\min}}{(p+1)!} \sum_{i \in \mathcal{N}} \|s_{i,k}\|^{p+1}$$
(5.11)

from (4.4), the observation that, at successful iterations, the algorithm enforces $\Delta m_k > 0$ and (4.11). As a consequence, $s_k \neq 0$. Hence at least one $||s_{i,k}||$ is strictly positive because of (5.2), and (5.11) therefore implies that (4.6) is well-defined. The inequality (5.10) then follows from Lemma 5.1. Following a now well-oiled track in convergence proofs for regularization methods, we derive an upper bound on the regularization parameters.

Lemma 5.4 [17, Lemma 4.6] Suppose that AS.2 and AS.3 hold. Then, for all $i \in \mathcal{N}$ and all $k \geq 0$,

$$\sigma_{i,k} \in [\sigma_{\min}, \sigma_{\max}], \tag{5.12}$$

where $\sigma_{\max} \stackrel{\text{def}}{=} \gamma_2 L_{\max}$.

Proof. Assume that, for some $i \in \mathcal{N}$ and $k \geq 0$, $\sigma_{i,k} \geq L_i$. Then (5.5) gives that (4.7) must fail, ensuring (5.12) because of the mechanism of the algorithm.

It is important to note that σ_{\max} is independent of ϵ . We now verify that the trial step produced by Step 2 of the PSAR*p* Algorithm either essentially fixes the residuals r_i to zero (their values being then fixed for the rest of the calculation), or the step is long enough (i.e. (4.1) holds), or it maintains these residuals safely away from zero in the sense that their norm exceeds an ϵ -independent constant.

Lemma 5.5 Suppose that AS.1, AS.2, AS.3 and AS.5 hold, that $\mathcal{H} \neq \emptyset$ and that (4.1) fails. Let

$$\omega \stackrel{\text{def}}{=} \min\left[\frac{a}{16}, \left(\frac{a}{12|\mathcal{N}|\left(\kappa_{\mathcal{N}} + \frac{\sigma_{\max}}{(p-q+1)!}\right)}\right)^{\frac{1-a}{2}}\right].$$
(5.13)

Then, if, for some $i \in \mathcal{H}$,

$$\|r_{i,k}\| < \omega, \tag{5.14}$$

we have that

$$|r_{i,k} + s_{i,k}|| \le \epsilon \quad or \quad ||r_{i,k} + s_{i,k}|| \ge \omega.$$
 (5.15)

Proof. The conclusion is obvious if $i \in C_k^+ = \mathcal{H} \setminus \mathcal{A}_k^+$. Consider now $i \in \mathcal{A}_k^+$ and suppose, for the purpose of deriving a contradiction, that

$$||r_{i,k} + s_{i,k}|| \in (\epsilon, \omega) \quad \text{for some} \quad i \in \mathcal{A}_k^+, \tag{5.16}$$

and immediately note that the failure of (4.1) and the orthonormality of the rows of U_i imply that

$$\|s_{i,k}\| \le \|s_k\| < \varpi \epsilon^{\frac{1}{p-q+1}} \le 1$$
(5.17)

and also that (4.2) must hold. As a consequence, for some $\delta_k \in (0, 1]$,

$$a \| r_{i,k} + s_{i,k} \| \chi_q(\delta_k) \ge \psi_{m,q}^{\epsilon,\delta_k}(x_k, s_k).$$
(5.18)

Consider now the vector

$$d_{k} = -\min\left[\delta_{k}, \|r_{i,k} + s_{i,k}\|\right] v_{i,k}^{+} \text{ with } v_{i,k}^{+} = U_{i}^{\dagger} u_{i,k}^{+} \stackrel{\text{def}}{=} U_{i}^{\dagger} \frac{r_{i,k} + s_{i,k}}{\|r_{i,k} + s_{i,k}\|}.$$
(5.19)

We now verify that d_k is admissible for problem (4.3). Clearly $||d_k|| = \delta_k$ because the rows of U_i are orthonormal. We also see that (1.2) and (1.11) imply that, since $i \in \mathcal{A}_k^+$,

$$d_k \in \mathcal{R}_{\{i\}} \subseteq \mathcal{R}_k^+. \tag{5.20}$$

Moreover, we have that

$$x_k + s_k + d_k \in [\![x_k + s_k, x_k + s_k - U_i^{\dagger}(r_{i,k} + s_{i,k})]\!],$$
(5.21)

where [v, w] denotes the line segment joining the vectors v and w. But

$$\begin{aligned} x_k + s_k - U_i^{\dagger}(r_{i,k} + s_{i,k}) &= x_k + s_k - U_i^{\dagger}U_i(x_k + s_k) + U_i^{\dagger}b_i \\ &= (I - U_i^{\dagger}U_i)(x_k + s_k) + U_i^{\dagger}b_i \\ &= P_{\ker(U_i)}[x_k + s_k] + U_i^{\dagger}b_i \\ &\in \mathcal{F}, \end{aligned}$$

where we have used (1.3) to deduce the last inclusion. Since \mathcal{F} is convex and $x_k + s_k \in \mathcal{F}$, we deduce from (5.21) that $x_k + s_k + d_k \in \mathcal{F}$. As a consequence, d_k is admissible for problem (4.3) and hence, using (5.18),

$$a \| r_{i,k} + s_{i,k} \| \chi_q(\delta_k) \ge \psi_{m,q}^{\epsilon,\delta_k}(x_k, s_k) \ge \max\left[0, m(x_k, s_k) - T_{m,q}(x_k, s_k - d_k)\right].$$
(5.22)

Moreover (5.20) and (3.13) imply that

$$m(x_{k}, s_{k}) - T_{m,q}(x_{k}, s_{k} - d_{k})$$

$$= m_{\mathcal{N}}(x_{k}, s_{k}) - T_{m_{\mathcal{N}},q}(x_{k}, s_{k} - d_{k}) + m_{i}(x_{i,k}, s_{i,k}) - m_{i}(x_{i,k}, s_{i,k} - U_{i}d_{k})$$

$$\geq -|m_{\mathcal{N}}(x_{k}, s_{k}) - T_{m_{\mathcal{N}},q}(x_{k}, s_{k} - d_{k})| + m_{i}(x_{i,k}, s_{i,k}) - m_{i}(x_{i,k}, s_{i,k} - U_{i}d_{k}).$$
(5.23)

We start by considering the first term in the right-hand side of this inequality. Observe now that (5.14) ensures that $x_k \in \mathcal{V}$ (as defined in (5.1)). Hence AS.5, (5.14) and (5.17) together imply that, for each $i \in \mathcal{N}$,

$$\begin{aligned} |m_{i}(x_{k},s_{k}) - T_{m_{i},q}(x_{k},s_{k} - d_{k})| \\ &\leq \left| \sum_{\ell=1}^{q} \frac{1}{\ell!} \nabla_{x}^{\ell} T_{m_{i},q}(x_{i,k},s_{i,k}) [-U_{i}d_{k}]^{\ell} \right| \\ &= \left| \sum_{\ell=1}^{q} \frac{1}{\ell!} \left(\sum_{t=\ell}^{p} \frac{1}{(t-\ell)!} \nabla_{x}^{t} f_{i}(x_{i,k}) [s_{i,k}]^{t-\ell} + \frac{\sigma_{i,k}}{(p+1)!} \|\nabla_{\cdot}^{\ell}\| s_{i,k}\|^{p+1} \| \right) [-U_{i}d_{k}]^{\ell} \right|. \end{aligned}$$

$$(5.24)$$

Using now the identity $||U_i d_k|| = ||d_k|| = \delta_k$ and the fact that

$$\sum_{t=\ell}^{p} \frac{1}{(t-\ell)!} \le 1 + \chi_{p-\ell}(1) < 3,$$

we obtain from (5.24), the triangle inequality and (5.17) that

$$|m_i(x_k, s_k) - T_{m_i, q}(x_k, s_k - d_k)| < \sum_{\ell=1}^q \frac{1}{\ell!} \left(3 \|\nabla_x^t f_i(x_{i,k})\|_{[t]} + \frac{\sigma_{i,k}}{(p+1)!} \|\nabla_{\cdot}^{\ell}\|_{s_{i,k}} \|^{p+1} \| \right) \delta_k^{\ell}.$$
(5.25)

But we have from Lemma 3.1 and (5.17) that, for $\ell \in \{1, \ldots, q\}$,

$$\left\|\nabla^{\ell}_{\cdot}\|s_{i,k}\|^{p+1}\right\| = \left|\pi(p-\ell+1)\right| \|s_{i,k}\|^{p+1-\ell} \le \frac{(p+1)!}{(p-q+1)!},\tag{5.26}$$

and therefore that

$$\begin{aligned} \left| m_{\mathcal{N}}(x_{k},s_{k}) - T_{m_{\mathcal{N}},q}(x_{k},s_{k}-\delta_{k}\|r_{i,k}+s_{i,k}\|v_{i,k}^{+}) \right| &< 3|\mathcal{N}| \left(\kappa_{\mathcal{N}} + \frac{1}{(p-q+1)!}\sigma_{\max}\right)\chi_{q}(\delta_{k}) \\ &\leq \frac{1}{4}a\omega^{a-1}\chi_{q}(\delta_{k}), \end{aligned}$$

$$(5.27)$$

where we have used (5.13) to derive the last inequality. Let us now consider the second term in the right-hand side of (5.23). Applying Lemma 3.2, we obtain that $\mu(||r_{i,k} + s_{i,k}||, \cdot)$ is concave between 0 and $-||r_{i,k} + s_{i,k}||$ and $\mu(||r_{i,k}||, \cdot)$ is concave between 0 and $-||r_{i,k}||$. Therefore, because of (5.14), we may deduce that

$$\begin{split} m_i(x_{i,k}, s_{i,k}) - m_i(x_{i,k}, s_{i,k} - U_i d_k) &= \mu(\|r_{i,k} + s_{i,k}\|, 0) - \mu(\|r_{i,k} + s_{i,k}\|, \|U_i d_k\|) \\ &\geq \nabla_{\zeta}^1 \mu(\|r_{i,k} + s_{i,k}\|, 0) \|U_i d_k\| \\ &\geq \nabla_{\zeta}^1 \mu(\|r_{i,k}\|, \|r_{i,k} + s_{i,k}\| - \|r_{i,k}\|) \|U_i d_k\| \\ &\geq a \|r_{i,k}\|^{a-1} \delta_k \\ &\geq \frac{1}{2} a \omega^{a-1} \chi_q(\delta_k), \end{split}$$

where the second and third inequalities result from (3.15). Combining now this inequality with (5.22), (5.23) and (5.27), we deduce that

$$a \| r_{i,k} + s_{i,k} \| \chi_q(\delta_k) > \frac{1}{2} a \omega^{a-1} \chi_q(\delta_k) - \frac{1}{4} a \omega^{a-1} \chi_q(\delta_k) = \frac{1}{4} a \omega^{a-1} \chi_q(\delta_k).$$

Finally, we obtain using (5.16) that

$$\omega > \frac{1}{4}\omega^{a-1},$$

which is impossible in view of (5.13). Hence (5.16) cannot hold and the proof is complete. \Box

This last result is crucial in that it shows that there is a "forbidden" interval (ϵ, ω) for the residual's norms $||r_i(x_k + s_k)||$, where ω only depends on the problem and it is independent of ϵ . This in turn allows to partition the successful iterates into subsets, distinguishing iterates which "fix" a residual to a near zero value, iterates with long steps and iterates with possibly short steps in regions where the considered objective function's *p*-th derivative tensor is safely bounded independently of ϵ . Our analysis now follows the broad outline of [17] while simplifying some arguments. Focusing on the case where $\mathcal{H} \neq \emptyset$, we first isolate the set of successful iterations which "deactivate" a residual, that is

$$\mathcal{S}_{\epsilon} \stackrel{\text{def}}{=} \{k \in \mathcal{S} \mid ||r_{i,k} + s_{i,k}|| \le \epsilon \text{ and } ||r_{i,k}|| > \epsilon \text{ for some } i \in \mathcal{H}\},\$$

and notice that, by construction

$$|\mathcal{S}_{\epsilon}| \le |\mathcal{H}|. \tag{5.28}$$

We next define the ϵ -independent constant

$$\alpha = \frac{3}{4}\omega$$

and

$$\mathcal{S}_{\omega} \stackrel{\text{def}}{=} \{ k \in \mathcal{S} \mid \|s_k\| \ge \frac{1}{4}\omega \}.$$
(5.29)

Moreover, for an iteration $k \in S \setminus (S_{\epsilon} \cup S_{\omega})$, we verify that \mathcal{A}_k can be partitioned into

$$\begin{aligned} \mathcal{I}_{\heartsuit,k} &\stackrel{\text{def}}{=} & \{i \in \mathcal{A}_k \mid \|r_{i,k}\| \in [\alpha, +\infty) \text{ and } \|r_{i,k} + s_{i,k}\| \in [\alpha, +\infty) \} \\ \mathcal{I}_{\diamondsuit,k} \stackrel{\text{def}}{=} & \{i \in \mathcal{A}_k \mid \left(\|r_{i,k}\| \in [\omega, +\infty) \text{ and } \|r_{i,k} + s_{i,k}\| \in (\epsilon, \alpha)\right) \\ & \text{or } & \left(\|r_{i,k}\| \in (\epsilon, \alpha) \text{ and } \|r_{i,k} + s_{i,k}\| \in [\omega, \infty)\right) \} \\ \mathcal{I}_{\clubsuit,k} \stackrel{\text{def}}{=} & \{i \in \mathcal{A}_k \mid \|r_{i,k}\| \in (\epsilon, \omega) \text{ and } \|r_{i,k} + s_{i,k}\| \in (\epsilon, \omega) \}. \end{aligned}$$

Morever, Lemma 5.5 shows that $\mathcal{I}_{\mathbf{a},k}$ is always empty and one additionally has that, if $i \in \mathcal{I}_{\diamond,k}$, then

$$||s_k|| \ge ||s_{i,k}|| \ge |||r_{i,k} + s_{i,k}|| - ||r_{i,k}||| \ge \omega - \alpha = \frac{1}{4}\omega,$$

implies that $k \in \mathcal{S}_{\|s\|}$. Hence $\mathcal{I}_{\diamondsuit,k}$ is also empty and

$$\mathcal{A}_k = \mathcal{I}_{\heartsuit,k} \quad \text{for} \quad k \in \mathcal{S} \setminus (\mathcal{S}_\epsilon \cup \mathcal{S}_\omega) \stackrel{\text{def}}{=} \mathcal{S}_\heartsuit.$$
 (5.30)

The next important result shows that steps at iteration belonging to S_{\heartsuit} are long enough, because they are taken over regions where a good ϵ -independent Lipschitz bound holds. Indeed, if $\mathcal{H} \neq \emptyset$ and assuming that $\epsilon \leq \alpha$, we have, for $k \in S_{\heartsuit}$, that $\mathcal{A}_k^+ = \mathcal{A}_k$ and thus that $\mathcal{W}_k^+ = \mathcal{W}_k$ and $\mathcal{R}_k^+ = \mathcal{R}_k$. Moreover, the definition of $I_{\heartsuit,k} = \mathcal{A}_k$ ensures that $\mathcal{A}_k \subseteq \mathcal{O}_{k,\alpha}$ and thus that Lemma 5.2 (and in particular (5.6)) guarantees that $f_{\mathcal{W}_k}$ satisfies standard derivative error bounds for functions with Lipschitz continuous *p*-th derivative (with corresponding Lipschitz constant $L(\alpha)$). We may therefore apply known results for such functions to $f_{\mathcal{W}_k}$. The following lemma is extracted from [11], by specializing Lemma 3.3 in that reference to the optimization of $f_{\mathcal{W}_k}$ over \mathcal{R}_k for functions with Lipschitz continuous *p*-th derivative (i.e. $\beta = 1$ in [11]).

Lemma 5.6 Suppose that AS.1 – AS.3 and AS.5 hold, that

$$\epsilon \le \alpha \quad if \quad \mathcal{H} \neq \emptyset \tag{5.31}$$

and consider $k \in S_{\heartsuit}$ such that the PSARp Algorithm does not terminate at iteration k + 1. Then

$$\|s_k\| \ge \kappa_{\heartsuit} \epsilon^{\frac{1}{p-q+1}} \quad with \quad \kappa_{\heartsuit} \stackrel{\text{def}}{=} \left(\frac{(p-q+1)!}{L(\alpha) + \theta + \sigma_{\max}}\right)^{\frac{1}{p-q+1}}.$$
(5.32)

We may finally establish our final evaluation complexity bound by combining our results so far.

Theorem 5.7 Suppose that AS.1-AS.5 and (5.31) hold. Then the PSARp Algorithm requires at most

$$\left\lfloor \kappa_{\mathcal{S}}(f(x_0) - f_{\text{low}}) \epsilon^{-\frac{p+1}{p-q+1}} \right\rfloor + |\mathcal{H}|$$
(5.33)

successful iterations and at most

$$\left\lfloor \left\lfloor \kappa_S(f(x_0) - f_{\text{low}}) \left(e^{-\frac{p+1}{p-q+1}} \right) \right\rfloor \left(1 + \frac{|\log \gamma_1|}{\log \gamma_2} \right) + \frac{1}{\log \gamma_2} \log \left(\frac{\sigma_{\max}}{\sigma_0} \right) \right\rfloor + |\mathcal{H}| + 1 \quad (5.34)$$

evaluations of f and its p first derivatives to return an (ϵ, δ) -approximate q-th-order-stationary point for problem (1.1), where

$$\kappa_S \stackrel{\text{def}}{=} \frac{(p+1)!}{\eta \sigma_{\min} \varsigma^{p+1}} \left(\frac{(p-q+1)!}{L(\alpha) + \theta + \sigma_{\max}} \right)^{-\frac{1}{p-q+1}}.$$
(5.35)

Proof. Consider $k \in S$ before termination. Because the iteration is successful, we obtain from AS.4, Step 3 of the algorithm and Lemma 5.3 that

$$f(x_0) - f_{\text{low}} \ge f(x_0) - f(x_{k+1}) = \sum_{j \in \mathcal{S}_k} \Delta f_k \ge \eta \sum_{j \in \mathcal{S}_k} \Delta T_k \ge \frac{\eta \sigma_{\min} \varsigma^{p+1}}{(p+1)!} \sum_{j \in \mathcal{S}_k} \|s_k\|^{p+1}.$$
 (5.36)

Defining now

$$\mathcal{S}_{\epsilon,k} \stackrel{\text{def}}{=} \mathcal{S}_{\epsilon} \cap \{0, \dots, k\}, \quad \mathcal{S}_{\omega,k} \stackrel{\text{def}}{=} \mathcal{S}_{\omega} \cap \{0, \dots, k\} \text{ and } \mathcal{S}_{\heartsuit,k} \stackrel{\text{def}}{=} \mathcal{S}_{\heartsuit} \cap \{0, \dots, k\},$$

we verify that $\mathcal{S}_{\omega,k}$ and $\mathcal{S}_{\heartsuit,k}$ form a partition of $\mathcal{S}_k \setminus \mathcal{S}_{\epsilon,k}$. As a consequence, we have that

$$f(x_{0}) - f_{\text{low}} \geq \frac{\eta \sigma_{\min}\varsigma^{p+1}}{(p+1)!} \left\{ |\mathcal{S}_{\omega,k}| \min_{j \in \mathcal{S}_{\|s\|,k}} \|s_{k}\|^{p+1} + |\mathcal{S}_{\heartsuit,k}| \min_{j \in \mathcal{S}_{\heartsuit,k}} \|s_{k}\|^{p+1} \right\} \\ \geq \frac{\eta \sigma_{\min}\varsigma^{p+1}}{(p+1)!} \left\{ |\mathcal{S}_{\omega,k}| (\frac{1}{4}\omega)^{p+1} + |\mathcal{S}_{\heartsuit,k}| \left(\kappa_{\heartsuit}\epsilon^{\frac{1}{p-q+1}}\right)^{p+1} \right\} \\ \geq \frac{\eta \sigma_{\min}\varsigma^{p+1}}{(p+1)!} \left\{ |\mathcal{S}_{\omega,k}| + |\mathcal{S}_{\heartsuit,k}| \right\} \min\left[(\frac{1}{4}\omega)^{p+1}, \left(\kappa_{\heartsuit}\epsilon^{\frac{1}{p-q+1}}\right)^{p+1} \right] \\ \geq \frac{\eta \sigma_{\min}\varsigma^{p+1}}{(p+1)!} |\mathcal{S}_{k} \setminus \mathcal{S}_{\epsilon,k}| \kappa_{\heartsuit}^{\frac{1}{p-q+1}} \epsilon^{\frac{1}{p-q+1}},$$

where we have used (5.36), Lemma 5.1, (5.29) and (5.32) to deduce the second inequality, and the assumption that (without loss of generality in view of (5.32)) $\kappa_{\heartsuit} \leq \frac{1}{4}\omega$ to deduce the last. The above inequality yields that

$$|\mathcal{S}_k| = |\mathcal{S}_k \setminus \mathcal{S}_{\epsilon,k}| + |\mathcal{S}_{\epsilon,k}| \le \kappa_S (f(x_0) - f_{\text{low}}) \epsilon^{-\frac{p+1}{p-q+1}} + |\mathcal{S}_{\epsilon,k}|,$$

where κ_S is given by (5.35). Since $|\mathcal{S}_{\epsilon,k}| \leq |\mathcal{S}_{\epsilon}| \leq |\mathcal{H}|$, we finally deduce that the bound (5.33) holds. The bound (5.34) then follows by applying Lemma 4.1 and observing that f and its first p derivatives are evaluated at most once per iteration, plus once at termination. \Box

We conclude our development by recalling that the above result is valid for $\mathcal{H} = \emptyset$, in which case the problem is a smooth convexly-constrained partially-separable problem. Note that the norm-equivalence constant ς occurs in (5.35), which indicate that the underlying geometry of the problem's invariants subspaces ker (U_i) may have a significant impact on complexity.

6 Conclusions

We have shown that an (ϵ, δ) -approximate q-th-order stationary point of partially-separable convexly-constrained optimization with non-Lipschitzian singularities can be found at most $O(\epsilon^{-(p+1)/(p-q+1)})$ evaluations of the objective function and its first p derivatives for any $q \in \{1, 2, \ldots, p\}$ whenever the smooth element functions $f_i, i \in \mathcal{N}$ of the objective function are p times differentiable. This worst-case complexity is obtained via our Algorithm 4.1 (PSAR_p) with a p-th order Taylor model which uses the underlying rotational symmetry of the Euclidean norm function for $f_{\mathcal{H}}$ and the first p derivatives (whenever they exist) of the "element functions" f_i , for $i \in \mathcal{M}$.

Several observations are of interest. A first one is that the results remain valid if Lipschitz continuity is not assumed on the whole of the feasible set, but restricted to the segments of

the "path of iterates", that is $\bigcup_k [\![x_k, x_{k+1}]\!]$. While this might in general be difficult to ensure a priori, there may be cases where problem structure could help. A second observation is that convexity of the feasible set is only used on the segments $\bigcup_{i,k} [\![x_{i,k}, U_i^{\dagger}b_i]\!]$. Again this might be exploitable in some cases. The third observation is that, in line with [11], it is possible to replace the Lipschitz continuity assumption by a weaker Hölder continuity.

We have seen in the introduction that our framework covers general non-overlapping as well as special cases of overlapping group sparsity, where (1.2) still holds. While it would obviously be interesting to relax this orthogonality condition, this seems very challenging with our present conceptual tools. Indeed, (1.2) is crucial for obtaining the second inclusion of (5.20) in the proof of Lemma 5.5, itself a central piece of our argumentation. The main difficulty, if (1.2) is not assumed, is to reconcile the analysis of the isotropic model described in Section 3 in the segment between the current point and singularity (as assumed in the definition of μ in (3.12) and Lemma 3.2) with the requirement that a step along the same segment leaves the (nearly) singular elements invariant (as requested in the definition of the optimality measure (2.6)). Further progress covering the general overlapping group sparsity case is therefore likely to require a new proof technique.

While it may be possible to handle non-kernel-centered feasible sets (maybe along the lines of the discussion in [17]), this remains open at this stage. Another interesting perspective is a more generic exploitation of geometric symmetries inherent to optimization problems: our treatment here focuses on a specific case of rotational symmetry, but this should not, one hopes, be limitative.

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Appendix

Proof Lemma 3.1 The proof of (3.3) is essentially borrowed from [11, Lemma 2.4], although details differ because the present version covers $a \in (0, 1)$. We first observe that $\nabla^{j}_{\cdot} ||r||^{a}$ is a *j*-th order tensor, whose norm is defined using (1.7). Moreover, using the relationships

$$\nabla^{1}_{\cdot} \|r\|^{\tau} = \tau \, \|r\|^{\tau-2} r \quad \text{and} \quad \nabla^{1}_{\cdot} (r^{\tau \otimes}) = \tau \, r^{(\tau-1) \otimes} \otimes I, \quad (\tau \in \mathbb{R}), \tag{A.1}$$

defining

$$\nu_0 \stackrel{\text{def}}{=} 1, \quad \text{and} \quad \nu_i \stackrel{\text{def}}{=} \prod_{\ell=1}^i (a+2-2\ell),$$
(A.2)

and proceeding by induction, we obtain that, for some $\mu_{j,i} \ge 0$ with $\mu_{1,1} = 1$,

$$\begin{split} \nabla_{\cdot}^{1} \left[\nabla_{\cdot}^{j-1} \| r \|^{a} \right] \\ &= \nabla_{\cdot}^{1} \left[\sum_{i=2}^{j} \mu_{j-1,i-1} \nu_{i-1} \| r \|^{a-2(i-1)} r^{(2(i-1)-(j-1))\otimes} \otimes I^{((j-1)-(i-1))\otimes} \right] \\ &= \sum_{i=2}^{j} \mu_{j-1,i-1} \nu_{i-1} \left[(a-2(i-1)) \| r \|^{a-2(i-1)-2} r^{(2(i-1)-(j-1)+1)\otimes} \otimes I^{(j-i)\otimes} \right. \\ &\quad + ((2(i-1)-(j-1))) \| r \|^{a-2(i-1)} r^{(2(i-1)-(j-1)-1)\otimes} \otimes I^{(j-1)-(i-1)+1)\otimes} \right] \\ &= \sum_{i=2}^{j} \mu_{j-1,i-1} \nu_{i-1} \left[(a+2-2i) \| r \|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \right. \\ &\quad + (2(i-1)-j+1) \| r \|^{a-2(i-1)} r^{(2(i-1)-j)\otimes} \otimes I^{(j-(i-1))\otimes} \right] \\ &= \sum_{i=2}^{j} \mu_{j-1,i-1} \nu_{i-1} (a+2-2i) \| r \|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \\ &\quad + \sum_{i=1}^{j-1} (2i-j+1) \mu_{j-1,i} \nu_{i} \| r \|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \\ &= \sum_{i=1}^{j} \left((a+2-2i) \mu_{j-1,i-1} \nu_{i-1} + (2i-j+1) \mu_{j-1,i} \nu_{i} \right) \| r \|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \end{split}$$

where the last equation uses the convention that $\mu_{j,0} = 0$ and $\mu_{j-1,j} = 0$ for all j. Thus we may write

$$\nabla^{j}_{\cdot} \|r\|^{a} = \nabla^{1}_{\cdot} \left[\nabla^{j-1}_{\cdot} \|r\|^{a} \right] = \sum_{i=1}^{j} \mu_{j,i} \nu_{i} \|r\|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes}$$
(A.3)

with

$$\mu_{j,i}\nu_i = (a+2-2i)\mu_{j-1,i-1}\nu_{i-1} + (2i-j+1)\mu_{j-1,i}\nu_i$$

= $[\mu_{j-1,i-1} + (2i-j+1)\mu_{j-1,i}]\nu_i,$ (A.4)

where we used the identity

$$\nu_i = (a+2-2i)\nu_{i-1}$$
 for $i = 1, \dots, j$ (A.5)

to deduce the second equality. Now (A.3) gives that

$$\nabla^{j}_{\cdot} \|r\|^{a} [v]^{j} = \sum_{i=1}^{j} \mu_{j,i} \nu_{i} \|r\|^{a-j} \left(\frac{r^{T} v}{\|r\|}\right)^{2i-j} (v^{T} v)^{j-i}.$$

It is then easy to see that the maximum in (1.7) is achieved for v = r/||r||, so that

$$\|\nabla_{\cdot}^{j}\|r\|^{a}\|_{[j]} = \left|\sum_{i=1}^{j} \mu_{j,i}\nu_{i}\right|\|r\|^{a-j} = |\pi_{j}|\|r\|^{a-j}$$
(A.6)

with

$$\pi_j \stackrel{\text{def}}{=} \sum_{i=1}^j \mu_{j,i} \nu_i. \tag{A.7}$$

Successively using this definition, (A.4), (A.5) (twice), the identity $\mu_{j-1,j} = 0$ and (A.7) again, we then deduce that

$$\pi_{j} = \sum_{\substack{i=1\\j=1}}^{j} \mu_{j-1,i-1}\nu_{i} + \sum_{\substack{i=1\\j=1\\j=1}}^{j} (2i-j+1)\mu_{j-1,i}\nu_{i}$$

$$= \sum_{\substack{i=1\\j=1\\j=1}}^{j-1} \mu_{j-1,i} [\nu_{i+1} + (2i-j+1)\nu_{i}]$$

$$= \sum_{\substack{i=1\\j=1}}^{j-1} \mu_{j-1,i} [(a+2-2(i+1))\nu_{i} + (2i-j+1)\nu_{i}]$$

$$= (a+1-j)\sum_{\substack{i=1\\i=1}}^{j-1} \mu_{j-1,i}\nu_{i}$$

$$= (a+1-j)\pi_{j-1}.$$
(A.8)

Since $\pi_1 = a$ from the first part of (A.1), we obtain from (A.8) that

$$\pi_j = \pi(a-j),\tag{A.9}$$

which, combined with (A.6) and (A.7), gives (3.3). Moreover, (A.9), (A.7) and (A.3) give (3.2) with $\phi_{i,j} = \mu_{j,i} \nu_i$. In order to prove (3.4) (where now ||r|| = 1), we use (A.3), (A.7), (A.9) and obtain that

$$\begin{split} \nabla_{\cdot}^{j} \|\beta_{1}r\|^{a} - \nabla_{\cdot}^{j}\|\beta_{2}r\|^{a} &= \sum_{i=1}^{j} \mu_{j,i}\nu_{i} \|\beta_{1}r\|^{a-2i} \beta_{1}^{(2i-j)}r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \\ &- \sum_{i=1}^{j} \mu_{j,i}\nu_{i} \|\beta_{2}r\|^{a-2i} \beta_{2}^{(2i-j)}r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \\ &= \pi(a-j) \left[\beta_{1}^{a-j} - \beta_{2}^{a-j}\right] \|r\|^{a-2i} r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes} \\ &= \pi(a-j) \left[\beta_{1}^{a-j} - \beta_{2}^{a-j}\right] r^{(2i-j)\otimes} \otimes I^{(j-i)\otimes}. \end{split}$$

Using (1.7) again, it is easy to verify that the maximum defining the norm is achieved for v = r and (3.4) then follows from ||r|| = 1.