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#### Abstract

The aggregate subgradient method is developed for solving unconstrained nonsmooth difference of convex (DC) optimization problems. The proposed method shares some similarities with both the subgradient and the bundle methods. Aggregate subgradients are defined as a convex combination of subgradients computed at null steps between two serious steps. At each iteration search directions are found using only two subgradients: the aggregate subgradient and a subgradient computed at the current null step. It is proved that the proposed method converges to a critical point of the DC optimization problem and also that the number of null steps between two serious steps is finite. The new method is tested using some academic test problems and compared with several other nonsmooth DC optimization solvers.

**Keywords:** Nonsmooth optimization, Nonconvex optimization, DC optimization, Subgradient method

**TUCS Laboratory** Turku Optimization Group (TOpGroup)

## **1** Introduction

Consider the unconstrained DC minimization problem

$$\begin{cases} \text{minimize} & f(x) \\ \text{subject to} & x \in \mathbb{R}^n, \end{cases}$$
(1)

where  $f(x) = f_1(x) - f_2(x)$ , functions  $f_1$  and  $f_2$  are convex and in general, nonsmooth. Here  $\mathbb{R}^n$  is the *n*-dimensional Euclidean space.

To date, several methods have been developed to locally solve Problem (1) [9, 15, 19, 21, 24, 25, 33, 34, 35]. Some of these methods are extensions of the Difference of Convex Algorithm (DCA) [24, 25, 33, 34] while others can be considered as the extension of the bundle method for convex optimization [9, 15, 19, 21]. In this paper, a different approach is proposed to develop an algorithm for solving Problem (1). The key element in this approach is the concept of *the aggregate subgradient*. Aggregate subgradients have been used to design nonsmooth optimization algorithms [3, 16, 23]. Although there are different definitions of these subgradients in all cases they are convex combinations of subgradients from the current bundle.

The subgradient method is widely used in nonsmooth optimization due to its simple structure. Its convergence has been proved only for convex problems [1, 10, 11, 29, 30, 31, 32]. This method uses one subgradient and one function evaluation at each iteration and involves no line search procedure. It is applicable to large-scale problems due to its small storage requirements.

Bundle methods are efficient methods in nonsmooth optimization (see, [2, 4, 13, 14, 18, 23, 27, 28]). The problem of finding the search direction in these methods is reduced to a certain quadratic programming subproblem whose size may increase significantly as the number of variables increases. Various approaches, in particular based on the notion of the aggregate subgradient, to reduce the size of this subproblem has been proposed in [16, 17, 26]. These approaches allow one to design bundle methods which are applicable to large-scale problems.

The method introduced in this paper shares some similarities with both the subgradient and the bundle methods. It is a descent method, uses more than one subgradient at each iteration and involves the line search procedure (similar to bundle methods). On the other hand, it does not involve any time consuming quadratic programming subproblem to find search directions (similar to the subgradient methods). The proposed method has a simple structure, and it is easy to implement. We prove the convergence of this method and demonstrate its performance using some nonsmooth unconstrained DC optimization test problems.

The paper is structured as follows. In Section 2, the new aggregate subgradient method is introduced and its convergence is studied. Results of numerical experiments are reported in Section 3. Section 4 concludes the paper. Throughout this paper, we use following notations and definitions. The inner product in  $\mathbb{R}^n$  is  $\langle u, v \rangle = \sum_{i=1}^n u_i v_i$ , and  $\|\cdot\|$  is the associated norm.  $S_1 = \{x \in \mathbb{R}^n : \|x\| = 1\}$  is the unit sphere, and  $B(x; \varepsilon) = \{y \in \mathbb{R}^n : \|x - y\| < \varepsilon\}$  is the open ball centered at  $x \in \mathbb{R}^n$ . The convex hull of a set is denoted by "conv". For a convex function  $f : \mathbb{R}^n \to \mathbb{R}$ , its subdifferential at a point  $x \in \mathbb{R}^n$  is [5]

$$\partial f(x) = \{ \xi \in \mathbb{R}^n : f(y) - f(x) \ge \langle \xi, y - x \rangle \ \forall y \in \mathbb{R}^n \}$$

and its  $\varepsilon$ -subdifferential is

$$\partial_{\varepsilon} f(x) = \{ \xi \in \mathbb{R}^n : f(y) - f(x) \ge \langle \xi, y - x \rangle - \varepsilon \ \forall \ y \in \mathbb{R}^n \}.$$

A point  $x^* \in \mathbb{R}^n$  is called a *critical point* of Problem (1) if

$$\partial f_1(x^*) \cap \partial f_2(x^*) \neq \emptyset.$$

# 2 The aggregate subgradient method

In this section, we introduce an aggregate subgradient method for solving Problem (1). For a given number  $\tau > 0$  consider the set

$$Q_{1\tau}(x) = \operatorname{conv} \bigcup_{d \in S_1} \partial f_1(x + \tau d), \ x \in \mathbb{R}^n.$$

**Proposition 1.** The set  $Q_{1\tau}(x)$  is convex and compact for any finite  $\tau > 0$ .

*Proof.* The convexity of the set  $Q_{1\tau}(x)$  follows from its definition. Since the subdifferential is bounded over any bounded set the set  $Q_{1\tau}(x)$  is bounded. To prove its closedness it is sufficient to prove the closedness of the set

$$\bar{Q}_{1\tau}(x) = \bigcup_{d \in S_1} \partial f_1(x + \tau d).$$

Take any sequence  $\{u_k\} \subset \bar{Q}_{1\tau}(x)$  such that  $u_k \to u$  as  $k \to \infty$ . This means that there exists the sequence  $\{d_k\} \subset S_1$  such that  $u_k \in \partial f_1(x + \tau d_k)$ . Since the set  $S_1$ is compact all limit points of the sequence  $\{d_k\}$  belong to this set. Without loss of generality, assume that  $d_k \to \bar{d} \in S_1$  as  $k \to \infty$ . Then the upper semicontinuity of the subdifferential mapping implies that  $u \in \partial f_1(x + \tau \bar{d})$  and therefore,  $u \in \bar{Q}_{1\tau}(x)$ . Since a convex hull of a closed set is also closed it follows that the set  $Q_{1\tau}(x)$  is closed.

**Remark 1.** Note that the set  $Q_{1\tau}(x)$  was introduced in [6], where it is called the spherical subdifferential.

The proof of the following theorem can be found in [5] (see Theorem 2.33).

**Theorem 1.** Let  $f : \mathbb{R}^n \to \mathbb{R}$  be convex with a Lipschitz constant L > 0 at  $x \in \mathbb{R}^n$ . Then for all  $\varepsilon \ge 0$  we have

$$\partial f(y) \subseteq \partial_{\varepsilon} f(x) \quad \forall \ y \in B\left(x; \frac{\varepsilon}{2L}\right).$$
 (2)

**Proposition 2.** For the set  $Q_{1\tau}(x)$ ,  $x \in \mathbb{R}^n$ ,  $\tau > 0$  the following holds:

$$Q_{1\tau}(x) \subseteq \partial_{\varepsilon} f_1(x)$$

for all  $\varepsilon > 2L\tau$ .

*Proof.* It is clear that  $x + \tau d \in B\left(x; \frac{\varepsilon}{2L}\right)$  for any  $\varepsilon > 2L\tau$  and  $d \in S_1$ . Then the proof follows from Theorem 1 and the convexity of  $\partial_{\varepsilon} f_1(x)$ .

Now take any subgradient  $v \in \partial f_2(x)$  and construct the set

$$Q_{\tau}(x) = \{\xi \in \mathbb{R}^n : \xi = u - v, u \in Q_{1\tau}(x)\}$$

The convexity of functions  $f_1$  and  $f_2$  and the subgradient inequality imply that

$$f(x+\tau d) - f(x) \le \tau \langle u - v, d \rangle \quad \forall \ u \in \partial f_1(x+\tau d)$$
(3)

and therefore,

$$f(x+\tau d) - f(x) \le \tau \max_{\xi \in Q_{\tau}(x)} \langle \xi, d \rangle.$$

It follows from Proposition 1 that the set  $Q_{\tau}(x)$  is convex and compact for any finite  $\tau > 0$ . If  $0_n \notin Q_{\tau}(x)$  at  $x \in \mathbb{R}^n$ , then one can use the set  $Q_{\tau}(x)$  to compute descent directions of the function f [8]. However, it is not always easy to compute the whole set  $Q_{\tau}(x)$ .

Next we describe the proposed algorithm, Algorithm 1. At each iteration of this algorithm, we use only two elements of the set  $Q_{\tau}(x)$  to compute search directions. The algorithm consists of inner and outer iterations. The inner iteration contains null steps whereas the outer iteration involves serious steps. Aggregate subgradients are calculated in the inner iterations.

In the next proposition, we prove that for each outer iteration  $k \ge 0$  the number of inner iterations, null steps, in Algorithm 1 is finite.

#### **Proposition 3.** Let

$$C_k = \max\{\|\xi\| : \xi \in Q_{\tau_k}(x_k)\} < +\infty.$$
(5)

Assume that  $\delta_k < C_k$ . Then the inner loop at the *k*-th iteration of Algorithm 1 stops after  $m_k > 0$  iterations, where

$$m_k \leq \left\lceil \frac{2\log_2(\delta_k/C_k)}{\log_2 C_{1k}} \right\rceil, \ C_{1k} = 1 - (1 - c_1)^2 (4C_k^2)^{-1} \delta_k^2.$$

Algorithm 1: Aggregate subgradient method.

**Data:**  $\sigma_1 \in (0,1)$  - decrease parameter for  $\tau$ ,  $\sigma_2 \in (0,1]$  - decrease parameter for the inner iteration tolerance  $\delta$ ,  $c_1 \in (0,1)$  - search control parameter,  $c_2 \in (0,c_1]$  - line search parameter,  $\varepsilon > 0$  - optimality tolerance.

**Result:** An approximate critical point of Problem (1).

*Step 0.* (Initialization). Choose a starting point  $x_0 \in \mathbb{R}^n$  and numbers  $\tau_0 > 0$ ,  $\delta_0 > 0$ . Set k = 0.

#### **Outer iteration**

#### Inner iteration

Step 1. (Initialization). Select any  $d_k^0 \in S_1$ . Compute  $u_k^0 \in \partial f_1(x_k + \tau_k d_k^0)$  and  $v_k \in \partial f_2(x_k)$ . Set  $\xi_k^0 = u_k^0 - v_k$ ,  $\bar{\xi}_k^0 = \xi_k^0$  and l = 0.

Step 2. Solve the problem

 $\begin{cases} \text{minimize} & \varphi^l(\lambda) = \|\lambda \xi_k^l + (1-\lambda) \bar{\xi}_k^l\|^2 \\ \text{subject to} & \lambda \in [0,1]. \end{cases}$ 

Let  $\bar{\lambda}_l$  be a solution to this problem.

Step 3. Calculate  $\bar{\xi}_{k}^{l+1} = \bar{\lambda}_{l}\xi_{k}^{l} + (1 - \bar{\lambda}_{l})\bar{\xi}_{k}^{l}$ . If  $\|\bar{\xi}_{k}^{l+1}\| \leq \delta_{k},$  (4)

then EXIT inner iterations and go to Step 6.

Step 4. (Search direction). Compute  $d_k^{l+1} = -\|\bar{\xi}_k^{l+1}\|^{-1}\bar{\xi}_k^{l+1}$ .

Step 5. If

$$f(x_k + \tau_k d_k^{l+1}) - f(x_k) > -c_1 \tau_k \|\bar{\xi}_k^{l+1}\|$$

then compute the subgradient  $u_k^{l+1} \in \partial f_1(x_k + \tau_k d_k^{l+1})$ , set  $\xi_k^{l+1} = u_k^{l+1} - v_k$ , l = l+1 and go to Step 2. Otherwise EXIT inner iterations and go to Step 7.

*Step 6.* (Stopping criterion). If  $\tau_k \leq \varepsilon$ , then STOP since an approximate critical point has been found. Otherwise, set  $\tau_{k+1} = \sigma_1 \tau_k$ ,  $\delta_{k+1} = \sigma_2 \delta_k$ ,  $x_{k+1} = x_k$ , k = k + 1 and go to Step 1.

*Step* 7. (Line search). Calculate the step-length  $\alpha_k \ge \tau_k$  as follows:

$$\alpha_k = \operatorname{argmax} \left\{ \alpha \geq 0 : f(x_k + \alpha d_k^{l+1}) - f(x_k) \leq -c_2 \alpha \| \bar{\xi}_k^{l+1} \| \right\}.$$

*Step 8.* Set  $x_{k+1} = x_k + \alpha_k d_k^{l+1}$ ,  $\tau_{k+1} = \tau_k$ ,  $\delta_{k+1} = \delta_k$ , k = k+1 and go to Step 1.

*Proof.* The inner loop in Algorithm 1 terminates when either the condition (4) or the condition of the sufficient descent

$$f(x_k + \tau_k d_k^{l+1}) - f(x_k) \le -c_1 \tau_k \|\bar{\xi}_k^{l+1}\|$$
(6)

is met. Let  $S_k^l = \{\xi \in \mathbb{R}^n : \xi = \lambda \xi_k^l + (1-\lambda) \overline{\xi_k^l}, \lambda \in [0,1]\}$ . In order to prove this proposition it is sufficient to estimate the upper bound of the number of iterations  $m_k$  when the condition (4) occurs. If none of the conditions (4) and (6) are satisfied at the *l*-th inner iteration, then the approximated subgradient  $\xi_k^{l+1} = u_k^{l+1} - v_k$  computed in Step 5 does not belong to the set  $S_k^l$ . Indeed, according to the definition of  $\overline{\xi}_k^{l+1}$  and the necessary and sufficient condition for a minimum (see Lemma 5.2.6 in [27]) we have

$$\langle \boldsymbol{\xi}, \bar{\boldsymbol{\xi}}_k^{l+1} \rangle \ge \| \bar{\boldsymbol{\xi}}_k^{l+1} \|^2 \quad \forall \; \boldsymbol{\xi} \in S_k^l.$$

$$\tag{7}$$

On the other hand, since the condition (6) does not hold we get

$$f(x_k + \tau_k d_k^{l+1}) - f(x_k) > -c_1 \tau_k \|\bar{\xi}_k^{l+1}\|.$$

In addition, it follows from (3) that

$$f(x_k + \tau_k d_k^{l+1}) - f(x_k) \le \tau_k \langle \xi_k^{l+1}, d_k^{l+1} \rangle,$$

therefore, we have

$$-c_1\tau_k\|\bar{\xi}_k^{l+1}\| < \tau_k\langle \xi_k^{l+1}, d_k^{l+1}\rangle,$$

which means that

$$\langle \xi_k^{l+1}, \bar{\xi}_k^{l+1} \rangle < c_1 \| \bar{\xi}_k^{l+1} \|^2.$$
 (8)

Then (7) and the condition  $c_1 \in (0, 1]$  imply that  $\xi_k^{l+1} \notin S_k^l$ . Therefore, it is proved that the newly calculated subgradient  $\xi_k^{l+1}$  is not on the segment between  $\overline{\xi}_k^l$  and  $\xi_k^l$ , and by updating subgradients in this manner, we change the set  $S_k^l$  at each inner iteration of Algorithm 1. This means that  $\overline{\xi}_k^{l+2} \neq \overline{\xi}_k^{l+1}$ . Indeed, if  $\overline{\xi}_k^{l+2} = \overline{\xi}_k^{l+1}$ , then the optimality condition (7) would hold also for all  $\xi \in S_k^{l+1}$  and since  $\xi_k^{l+1} \in S_k^{l+1}$  we get

$$\langle \boldsymbol{\xi}_k^{l+1}, \bar{\boldsymbol{\xi}}_k^{l+1} \rangle \ge \| \bar{\boldsymbol{\xi}}_k^{l+1} \|^2$$

contradicting (8).

Next, we prove that if none of the stopping criteria is satisfied, then the new aggregated subgradient  $\bar{\xi}_k^{l+2}$  computed at Step 3 is better than  $\bar{\xi}_k^{l+1}$  in the sense that  $\varphi^{l+1}(\bar{\lambda}_{l+1}) < \varphi^l(\bar{\lambda}_l)$ . Since the function  $\varphi^{l+1}$  is strictly convex  $\bar{\lambda}_{l+1}$  is its only minimizer. Furthermore, it follows from

$$\bar{\xi}_{k}^{l+2} = \bar{\lambda}_{l+1}\xi_{k}^{l+1} + (1 - \bar{\lambda}_{l+1})\bar{\xi}_{k}^{l+1}$$

and  $\bar{\xi}_k^{l+2} \neq \bar{\xi}_k^{l+1}$  that  $\bar{\lambda}_{l+1} > 0$ . Then we have

$$\varphi^{l+1}(\bar{\lambda}_{l+1}) = \|\bar{\xi}_k^{l+2}\|^2 < \varphi^{l+1}(0) = \|\bar{\xi}_k^{l+1}\|^2 = \varphi^l(\bar{\lambda}_l).$$

In addition, it is clear that

$$\|\bar{\xi}_k^{l+2}\|^2 \le \|t\xi_k^{l+1} + (1-t)\bar{\xi}_k^{l+1}\|^2 \quad \forall t \in [0,1].$$

This means that for all  $t \in [0, 1]$ 

$$\|\bar{\xi}_{k}^{l+2}\|^{2} \leq t^{2} \|\xi_{k}^{l+1} - \bar{\xi}_{k}^{l+1}\|^{2} + 2t\langle\xi_{k}^{l+1} - \bar{\xi}_{k}^{l+1}, \bar{\xi}_{k}^{l+1}\rangle + \|\bar{\xi}_{k}^{l+1}\|^{2}.$$

It follows from (5) that

$$\|\boldsymbol{\xi}_k^{l+1} - \bar{\boldsymbol{\xi}}_k^{l+1}\| \le 2C_k.$$

This together with the inequality (8) implies that

$$\varphi^{l+1}(\bar{\lambda}_{l+1}) = \|\bar{\xi}_k^{l+2}\|^2 \leq 4t^2 C_k^2 + (1 - 2t(1 - c_1))\|\bar{\xi}_k^{l+1}\|^2$$

for all  $t \in [0,1]$ . Let  $t_0 = (1-c_1) \|\bar{\xi}_k^{l+1}\|^2 (4C_k^2)^{-1}$ . It is clear that  $t_0 \in (0,1)$ . Therefore, for  $t = t_0$  we get

$$\|\bar{\xi}_k^{l+2}\|^2 \le \left(1 - (1 - c_1)^2 (4C_k^2)^{-1} \|\bar{\xi}_k^{l+1}\|^2\right) \|\bar{\xi}_k^{l+1}\|^2.$$

At any *l*-th iteration, l > 0, where the stopping condition (4) is not met, we have  $\|\bar{\xi}_k^{l+1}\| > \delta_k$ . Then it follows that

$$\|\bar{\xi}_k^{l+2}\|^2 < \left(1 - (1 - c_1)^2 (4C_k^2)^{-1} \delta_k^2\right) \|\bar{\xi}_k^{l+1}\|^2.$$

Let  $C_{1k} = 1 - (1 - c_1)^2 (4C_k^2)^{-1} \delta_k^2$ . It is obvious that  $C_{1k} \in (0, 1)$ . Then we have

$$\varphi^{l+1}(\bar{\lambda}_{l+1}) < C_{1k}\varphi^l(\bar{\lambda}_l).$$

Since  $\varphi^0(\bar{\lambda}_0) = \|\xi_k^0\|^2 \le C_k^2$  we get

$$\varphi^l(\bar{\lambda}_l) < C_{1k}^l C_k^2.$$

Thus, the inequality (4) is satisfied if

$$C_{1k}^{m_k}C_k^2 \leq \delta_k^2,$$

which must happen after at most  $m_k$  iterations, where

$$m_k \leq \left\lceil \frac{2\log_2(\delta_k/C_k)}{\log_2 C_{1k}} \right\rceil.$$

This completes the proof.

Next, we prove that Algorithm 1 generates a sequence which converges to a critical point of Problem (1). For a given  $x_0 \in \mathbb{R}^n$  consider the level set

$$\mathscr{L}(x_0) = \{ x \in \mathbb{R}^n : f(x) \le f(x_0) \}.$$

**Theorem 2.** Assume that the set  $\mathscr{L}(x_0)$  is bounded for any starting point  $x_0 \in \mathbb{R}^n$ , and  $\varepsilon = 0$  in Algorithm 1. Then any limit point of the sequence  $\{x_k\}$  generated by this algorithm is a critical point of Problem (1).

*Proof.* Algorithm 1 consists of inner and outer iterations. In all inner iterations, the values of  $\tau_k$  and  $\delta_k$  remain unchanged. Furthermore, these parameters remain unchanged in Steps 7 and 8 of outer iterations. Their values are updated in Step 6 when the condition (4) is satisfied. According to Proposition 3 the number of steps in inner iterations is finite for the given  $\tau_k$  and  $\delta_k$ . As an outcome of the inner loop either Step 6 or Steps 7 and 8 are executed.

First, we show that for each unchanged values of  $\tau_k$  and  $\delta_k$  the number of executions of Steps 7 and 8 is finite. Indeed, in this case the line search is carried out in Step 7 and we have

$$f(x_{k+1}) - f(x_k) \le -c_2 \tau_k \|\bar{\xi}_k^{l+1}\|, \|\bar{\xi}_k^{l+1}\| > \delta_k$$

Therefore, we get

$$f(x_{k+1}) - f(x_k) \le -c_2 \tau_k \delta_k.$$
(9)

Since the set  $\mathscr{L}(x_0)$  is compact and the function *f* is continuous it follows that

$$f_* := \inf \{ f(x) : x \in \mathbb{R}^n \} > -\infty.$$
(10)

If the values of  $\tau_k$  and  $\delta_k$  are unchanged infinitely many times then (9) implies that  $f(x_k) \to -\infty$  as  $k \to \infty$ . This contradicts the condition (10). Thus,  $\tau_k$  and  $\delta_k$ are updated in Step 6 after a finite number of outer iterations, and we denote those iterations by  $k_p$ , p = 1, 2... During the iteration  $k_p$ , the condition (4) is satisfied meaning that

$$\|\bar{\boldsymbol{\xi}}_{k_p}\| \leq \boldsymbol{\delta}_{k_p} \quad \forall \ p = 1, 2, \dots,$$

where  $\bar{\xi}_{k_p} = \bar{\xi}_{k_p}^l$  for some l > 0. It is obvious that  $\bar{\xi}_{k_p} \in Q_{\tau_{k_p}}(x_{k_p})$  which, in turn, implies that

$$\min_{\boldsymbol{\xi}\in\mathcal{Q}_{\tau_{k_p}}(x_{k_p})}\|\boldsymbol{\xi}\|\leq\delta_{k_p},$$

or

$$\min_{u \in Q_{1\tau_{k_p}}(x_{k_p})} \|u - v_{k_p}\| \le \delta_{k_p}.$$
 (11)

In addition, it follows from Proposition 2 that

$$Q_{1\tau_{k_p}}(x_{k_p}) \subseteq \partial_{\bar{\varepsilon}} f_1(x_{k_p}) \tag{12}$$

for all  $\bar{\epsilon} > 2L\tau_{k_p}$ , where L > 0 is the Lipschitz constant of the function  $f_1$  over the set  $\mathscr{L}(x_0)$ . Take any  $c_0 > 1$  and consider  $\bar{\epsilon}_{k_p} = 2c_0L\tau_{k_p}$ . Then the inequality (11) and the inclusion (12) imply that

$$\min_{u\in\partial_{\bar{e}_{k_p}}f_1(x_{k_p})}\|u-v_{k_p}\|\leq \delta_{k_p},$$

which means that

$$\partial_{\bar{\boldsymbol{\varepsilon}}_{k_p}} f_1(x_{k_p}) \cap B(v_{k_p}; \boldsymbol{\delta}_{k_p}) \neq \boldsymbol{\emptyset},$$

or

$$\partial_{\bar{\varepsilon}_{k_p}} f_1(x_{k_p}) \cap B(\partial f_2(x_{k_p}); \delta_{k_p}) \neq \emptyset.$$
(13)

Since the set  $\mathscr{L}(x_0)$  is compact and  $x_{k_p} \in \mathscr{L}(x_0)$  for any p > 0 the sequence  $\{x_{k_p}\}$  has at least one limit point. For the sake of simplicity, assume that  $x_{k_p} \to \bar{x}$  as  $p \to \infty$ . The upper semicontinuity of the subdifferential mapping and the fact that  $\tau_{k_p}, \delta_{k_p} \downarrow 0$  as  $p \to \infty$  imply that for  $\gamma > 0$  there exists  $p_{\gamma} > 0$  such that

$$\partial_{\bar{\varepsilon}_{k_p}} f_1(x_{k_p}) \subset \partial f_1(\bar{x}) + B(0_n; \gamma), \quad \partial f_2(x_{k_p}) \subset \partial f_2(\bar{x}) + B(0_n; \gamma) \text{ and } \delta_{k_p} < \gamma$$

for all  $p > p_{\gamma}$ . Then for all  $p > p_{\gamma}$  we have

$$B(\partial f_2(x_{k_p}); \delta_{k_p}) \subseteq B(\partial f_2(x_{k_p}); \gamma) \subseteq \partial f_2(\bar{x}) + B(0_n; 2\gamma).$$

Taking into account (13) we get

$$\left(\partial f_1(\bar{x}) + B(0_n; \gamma)\right) \cap \left(\partial f_2(\bar{x}) + B(0_n; 2\gamma)\right) \neq \emptyset.$$

Since  $\gamma > 0$  is arbitrary we have

$$\partial f_1(\bar{x}) \cap \partial f_2(\bar{x}) \neq \emptyset.$$

This completes the proof.

#### 

## **3** Numerical results

In this section, we compare the proposed aggregate subgradient algorithm with three other nonsmooth optimization methods. The collection of test problems used in our experiments consists of nonsmooth unconstrained DC optimization problems which are formulated by modifying various convex problems introduced in [16, 19, 20]. The problems are given in Appendix. We present only the results of those problems where all the solvers converge to the same critical point. Overall, this leaves us 150 test problems. The problems can be formulated with different number of variables: we use 100 (extra small), 200 (small), 500 (medium) and 1000 (large)<sup>1</sup> variables in our tests.

<sup>&</sup>lt;sup>1</sup>We include 148 large test problems since none of four methods succeeded in solving two other problems.

**Solvers and Parameters.** We use the following solvers in our experiments:

- DCA is an implementation of the well-known difference of convex algorithm (DCA) [25]. The DCA, in general, finds critical solutions to the DC optimization problems. In the implementation of the DCA, we apply the proximal bundle method MPBNGC [27] to solve the convex subproblem.
- PBDC [19] is a proximal bundle method for DC optimization. Similar to the DCA, the PBDC is guaranteed to find critical solutions to the DC optimization problems.
- SolvOpt [22] is an implementation of Shor's *r*-algorithm [32]. Unlike the other methods considered in this paper, the SolvOpt is not utilizing the DC structure of the problem.
- AggSub is an implementation of the proposed algorithm. The following parameters were used in our experiments:

$$\begin{cases} \sigma_1 = 0.2, \ \sigma_2 = 1.0, \ \delta_0 = 10^{-7}, \\ \varepsilon = 10^{-5}, \ c_1 = 0.2, \ c_2 = 0.05, \end{cases} \quad \text{and} \quad \tau_0 = \begin{cases} 10.0, & n < 200, \\ 50.0, & n \ge 200. \end{cases}$$

The Fortran 95 source code of AggSub is available at http://napsu. karmitsa.fi/aggsub/. All experiments are performed on an Intel<sup>®</sup> Core<sup>TM</sup> i5-2400 CPU (3.10GHz, 3.10GHz) running under Windows 7. To compile the codes, we use gfortran, the GNU Fortran compiler. All the solvers are used with the default settings of the codes.

We say that a solver finds a solution with respect to a tolerance  $\varepsilon > 0$  if

$$\frac{f_{\text{best}} - f_{\text{opt}}}{1 + |f_{\text{opt}}|} \le \varepsilon,$$

where  $f_{\text{best}}$  is the best value of the objective function obtained by the solver and  $f_{\text{opt}}$  is the best known (or optimal) solution. Otherwise, the solver *fails*. We set  $\varepsilon = 10^{-3}$ . In addition to the usual stopping criteria of the solvers, we terminate the experiment if the elapsed CPU time exceeds half an hour per problem.

**Results.** The results are analyzed using the performance profiles (Figures 1 - 2) introduced in [12]. We compare the performance of the solvers both in terms of the computational time (CPU time) and the number of the subgradient evaluations (evaluations for short) — we use the sum of the number of the subgradient evaluations of the DC components in the AggSub, the PBDC and the DCA, while for the SolvOpt we use the number of the subgradient evaluations of the solvert we use the number of the subgradient evaluations.

number of function evaluations follows the similar trends with the number of the subgradient evaluations with all the solvers, and thus, we omit these results.

In the performance profiles, the value of  $\rho_s(\tau)$  at  $\tau = 0$  shows the ratio of the test problems for which the solver *s* is the best — that is, the solver *s* uses the least computational time or evaluations — while the value of  $\rho_s(\tau)$  at the rightmost abscissa gives the ratio of the test problems that the solver *s* can solve — that is, the robustness of the solver *s*. In addition, the higher is a particular curve, the better is the corresponding solver.

It is clear from Figure 1 that the AggSub is the most efficient and accurate with all sizes of the test problems — it is superior to other three methods in the larger problems. With the number of evaluations, as usual for any subgradient method, the AggSub has rather high numbers as can be seen in Figure 2 (a) and (b). However, this observation is faded out with the larger problems (see Figure 2 (c) and (d)). The reason for most of the failures with the SolvOpt is its inaccuracy while the PBDC usually stops far away from the solution due to the time limit. From these results, we conclude that, overall, the AggSub is an efficient and robust method.

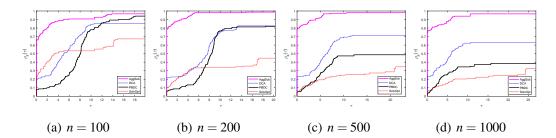


Figure 1: CPU time.

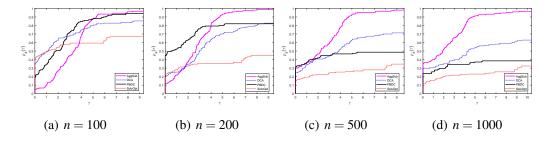


Figure 2: Number of subgradient evaluations.

# 4 Concluding remarks

In this paper, an aggregated subgradient method is developed to solve unconstrained nonsmooth difference of convex (DC) optimization problems. The method combines the strengths of the well-known subgradient and bundle methods. It is a descent method, easy to implement and does not involve solving of any timeconsuming quadratic programming subproblem, since only two subgradients are utilized to calculate search directions. In addition, the proposed method is globally convergent to a critical point.

The aggregated subgradient method was tested using large-scale nonsmooth unconstrained DC programming problems and compared with several other nonsmooth DC programming solvers. Results demonstrate that the method is efficient and robust for solving nonsmooth DC optimization problems, although in some cases it may require a large number of function and subgradient evaluations. The proposed method outperforms other methods when the evaluation of the objective function and its subgradient is not expensive.

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# **Appendix: New DC Test Problems**

All the test problems are unconstrained DC optimization problems, where objective functions are presented as DC functions:

$$f(x) = f_1(x) - f_2(x).$$

In what follows we first present some convex functions. These functions are used as  $f_1$  and  $f_2$  to obtain a DC problem. Different combinations used with suitable starting points are given in Table 1. The functions and starting points are combined such that the resulting problem satisfies the following criteria:

- the problem is bounded from below,
- the function  $f = f_1 f_2$  is nonconvex,
- the starting point for the problem is not critical,
- the problem has not too many local solution. That is, algorithms converge mostly to same (or only couple of) minima from different starting points,
- the computational time with 1000 variables is more than 1 sec. with AggSub (i.e. the problem is not too simple),

#### **Convex functions**

1. 
$$f(x) = (n+1) \max \{x_i^2 : i = 1, ..., n\}$$
  
2.  $f(x) = \max \{x_i^2 : i = 1, ..., n\}$  (Generalization of MAXQ)  
3.  $f(x) = \sum_{i=1}^n x_i^2$   
4.  $f(x) = n \max \{|x_i| : i = 1, ..., n\}$   
5.  $f(x) = \sum_{i=1}^n |x_i|$   
6.  $f(x) = 20 \max \{ \left| \sum_{i=1}^n (x_i - x_i^*) t_j^{i-1} \right| : j = 1, ..., 20 \}, t_j = 0.05j, j = 1, ..., 20$   
7.  $f(x) = \sum_{j=1}^{20} \left| \sum_{i=1}^n (x_i - x_i^*) t_j^{i-1} \right|, t_j = 0.05j, j = 1, ..., 20$   
8.  $f(x) = \sum_{i=2}^n |x_i - x_{i-1}|$   
9.  $f(x) = \max \{ \left| \sum_{j=1}^n \frac{x_j}{i+j-1} \right| : i = 1, ..., n \}$  (Generalization of MXHILB)

10. 
$$f(x) = \sum_{i=1}^{n-1} \max\left\{-x_i - x_{i+1}, -x_i - x_{i+1} + (x_i^2 + x_{i+1}^2 - 1)\right\}$$
 (Chained LQ)

- 11.  $f(x) = \sum_{i=1}^{n-1} \max\left\{x_i^4 + x_{i+1}^2, (2-x_i)^2 + (2-x_{i+1})^2, 2e^{-x_i + x_{i+1}}\right\}$  Chained CB3 I
- 12.  $f(x) = \max\left\{\sum_{i=1}^{n-1} (x_i^4 + x_{i+1}^2), \sum_{i=1}^{n-1} ((2 x_i)^2 + (2 x_{i+1})^2), \sum_{i=1}^{n-1} (2e^{-x_i + x_{i+1}})\right\}$  Chained CB3 II
- 13.  $f(x) = \sum_{i=1}^{n} |x_i| + 10 \sum_{i=1}^{n} \max \left\{ 2(x_i^2 x_i 1), 0 \right\}$
- 14.  $f(x) = n \max \left\{ \left| \sum_{j=1}^{n} \frac{x_j}{i+j-1} \right| : i = 1, \dots, n \right\}$
- 15.  $f(x) = \sum_{i=1}^{n} \left| \sum_{j=1}^{n} \frac{x_j}{i+j-1} \right|$

16. 
$$f(x) = (n-1) \max \left\{ \max \left\{ x_i^4 + x_{i+1}^2, (2-x_i)^2 + (2-x_{i+1})^2, 2e^{-x_i + x_{i+1}} \right\} : i = 1, \dots, n-1 \right\}$$
  
17. 
$$f(x) = \max \left\{ 2\sum_{i=1}^{n-1} \left( x_i^2 + (x_{i+1} - 1)^2 + x_{i+1} - 1 \right), 0 \right\}$$

18. 
$$f(x) = \sum_{i=1}^{n-1} (x_i^2 + (x_{i+1} - 1)^2 + x_{i+1} - 1)$$

### **Starting points**

- 1.  $x_0 = (1, 1, ..., 1)^T \in \mathbb{R}^n$ 2.  $x_0 = (i, i = 1, ..., ]n/2[, -i, i = ]n/2[+1, ..., n)^T$ 3.  $x_{0,i} = 1$  for odd *i* and  $x_{0,i} = -1$  for even *i*, i = 1, ..., n
- 4.  $x_{0,i} = -1.5$  for odd *i* and  $x_{0,i} = 2$  for even *i*, i = 1, ..., n
- 5.  $x_{0,i} = i, \quad i = 1, \dots, n$

# New DC problems (152 different).

Table 1. Combinations of problems.		
$f_1$	$f_2$	<i>x</i> <sub>0</sub>
1	3	1, 2, 3, 4, 5
1	5	1, 2, 3, 4, 5
1	6	1, 2, 3, 4, 5
1	9	1, 2, 3, 4, 5
2	5	1, 2, 3, 4, 5
3	2	1, 2, 3, 4, 5
3	4	1, 2, 3, 4, 5
3	5	1, 2, 3, 4, 5
4	9	1, 2, 3, 4, 5
5	9	1,2,3,4,5
10	9	1, 2, 3, 4, 5
10	15	1, 2, 3, 4, 5
11	7	1, 2, 3, 4, 5
11	9	2, 3, 4, 5
11	14	1, 2, 3, 4, 5 <sup>(1)</sup>
11	15	2, 3, 4, 5
12	9	2, 3, 4, 5
12	14	2, 3, 4, 5 <sup>(2)</sup>
12	15	1, 2, 3, 4, 5
13	14	2, 3, 4
13	15	1, 2, 3, 4, 5
16	2	2, 3, 4, 5
16	3	2, 3, 4, 5
16	4	2, 3, 4, 5
16	5	2, 3, 4, 5
16	6	1, 3, 4, 5 <sup>(3)</sup>
16	7	2, 3, 4, 5
16	8	2, 3, 4, 5
16	9	2, 3, 4, 5
16	10	2, 3, 4, 5
16	14	2, 3, 4, 5
16	15	2, 3, 4, 5
16	18	2, 3, 4, 5
17	15	1, 2, 3, 4, 5

Table 1: Combinations of problems.

<sup>1</sup> Omitted also  $x_0 = 2$  with n = 200,500 and 1000, and  $x_0 = 5$  with n = 200,500.

<sup>2</sup> Omitted also  $x_0 = 4$  with n = 100 and  $x_0 = 5$  with n = 200, 500.

<sup>3</sup> Omitted also  $x_0 = 5$  with n = 100.



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