

A SIMPLE ITERATIVE ALGORITHM FOR MAXCUT*

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Abstract

We propose a simple iterative (SI) algorithm for the maxcut problem through fully using an equivalent continuous formulation. It does not need rounding at all and has advantages that all subproblems have explicit analytic solutions, the cut values are monotonically updated and the iteration points converge to a local optima in finite steps via an appropriate subgradient selection. Numerical experiments on G-set demonstrate the performance. In particular, the ratios between the best cut values achieved by SI and those by some advanced combinatorial algorithms in [Ann. Oper. Res., 248 (2017), 365–403] are at least 0.986 and can be further improved to at least 0.997 by a preliminary attempt to break out of local optima.

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1. Introduction

Given an undirected simple graph $G = (V, E)$ of order n with the vertex set V and the edge set E , a set pair (S, S') is called a cut of G if $S \cap S' = \emptyset$ and $S \cup S' = V$. The maxcut problem, one of Karp's 21 NP-complete problems [10], aims at finding a specific cut (S, S') of G to maximize the cut value

$$\text{cut}(S) = \sum_{\{i,j\} \in E(S,S')} w_{ij}, \quad (1.1)$$

where $E(S, S')$ collects all edges cross between S and S' in E , and w_{ij} denotes the nonnegative weight on the edge $\{i, j\} \in E$.

Due to its widespread applications in various areas [1, 2, 4], several maxcut algorithms have been proposed to search for approximate solutions and usually fall into two distinct categories: discrete algorithms and continuous ones. The former mainly refer to the combinatorial algorithms for maxcut, which directly deal with the discrete objective function (1.1) and usually adopt both complicated techniques to break out of local optima and advanced heuristics improving the solution quality, such as the scatter search [12], the tabu search [15] and hybrid

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strategies within the framework of evolutionary algorithms [11, 13]. In contrast, the objective functions for the latter, often obtained from the relaxation of the discrete objective function (1.1), are continuous, and thus standard continuous optimization algorithms can be applied into the relaxed problems in a straightforward manner, for instance, the Goemans-Williamson (GW) algorithm [9], the CirCut algorithm resulted from the rank-two relaxation heuristics [3], the spectral cut (SC) algorithm [8, 16] and its recursive implementation (RSC) [5, 14, 18]. For all these continuous algorithms, rounding is essential and indispensable in obtaining a cut from a solution of the corresponding relaxed problem. The GW algorithm rounds the solution of a relaxing semidefinite programming via randomly selecting the hyperplanes until it achieves an expected cut value. A deterministic strategy, named Procedure-CUT, is adopted by CirCut to round the angle-vector solution to get a best possible associated cut. The SC algorithm obtains a cut by rounding the maximal eigenvector of graph Laplacian by a threshold, while the RSC algorithm recursively distributes part of unabsorbed points into two sets corresponding to a cut where the selection and assignment are determined by rounding the approximate solution of the dual Cheeger cut problem.

In this work, we propose a novel continuous algorithm for the maxcut problem, i.e. a simple iterative (SI) algorithm. Compared with the above-mentioned continuous maxcut algorithms, the proposed SI algorithm has the following distinct advantages. First, our inner subproblem can be solved analytically (see Theorem 3.1), whereas no matter the GW algorithm or the RSC algorithm needs call other optimization solvers for the inner subproblems. This constitutes the main reason why we use the adjunct word simple for the proposed algorithm. Second, our continuous optimization problem is directly equivalent to the maxcut problem (see Theorem 2.1), and the corresponding cut is updated in an iterative manner and converges to the local maximum (see Theorem 3.4). That is, the SI algorithm does not need any rounding at all. Finally, as an iterative algorithm, SI may select the cut by SC to be the initial point (see Section 4). In other words, it can also be used to improve the quality of the solution obtained from any other algorithms.

The rest is organized as follows. Section 2 establishes an equivalent continuous formulation of the maxcut problem (1.1) and a Dinkelbach-type iterative algorithm with global convergence. However, the solvability of the related inner subproblem can not be assured due to both the NP-hardness and the lack of convexity. To this end, in Section 3, we propose our simple iterative algorithm with an analytical solution to the inner problem. Both cost analysis and quality check are performed through numerical experiments on G-set in Section 4. Besides, in order to further improve the quality, a preliminary attempt to break out of local optima is also implemented there. We are concluded with a few remarks in Section 5.

2. Equivalent Continuous Problems

Given an undirected graph $G = (V, E)$ with nonnegative weights, let

$$I(\mathbf{x}) = \sum_{\{i,j\} \in E} w_{ij} |x_i - x_j|, \quad (2.1)$$

$$\|\mathbf{x}\|_\infty = \max\{|x_1|, \dots, |x_n|\}, \quad (2.2)$$

$$F(\mathbf{x}) = \frac{I(\mathbf{x})}{\|\mathbf{x}\|_\infty}. \quad (2.3)$$

It can be readily verified that the nonnegative function $F(\mathbf{x})$ can achieve its maximum on $\mathbb{R}^n \setminus \{\mathbf{0}\}$ provided by its homogeneity of degree zero.

Let

$$S^\pm(\mathbf{x}) = \{i \in V : x_i = \pm \|\mathbf{x}\|_\infty\}, \quad (2.4)$$

$$S^<(\mathbf{x}) = \{i \in V : |x_i| < \|\mathbf{x}\|_\infty\}. \quad (2.5)$$

Then for any $\mathbf{x}^* \in \mathbb{R}^n \setminus \{\mathbf{0}\}$, we have

$$M(\mathbf{x}^*) := \{\mathbf{x} \mid \|\mathbf{x}\|_\infty = \|\mathbf{x}^*\|_\infty, S^\pm(\mathbf{x}^*) \subset S^\pm(\mathbf{x})\} \quad (2.6)$$

is a convex polytope. In fact, the convexity of $I(\mathbf{x})$ directly implies that, if $\mathbf{x}^* \in M(\mathbf{x}^*)$ is a maximizer of $F(\mathbf{x})$ on $\mathbb{R}^n \setminus \{\mathbf{0}\}$, so does any $\mathbf{x} \in M(\mathbf{x}^*)$.

For any nonempty subset $S \subset V$, we define an indicative vector $\mathbf{1}_S$

$$(\mathbf{1}_S)_i = \begin{cases} 1, & i \in S, \\ 0, & i \notin S, \end{cases}$$

and then

$$\mathbf{x} = \mathbf{1}_S - \mathbf{1}_{V \setminus S},$$

which satisfies

$$\frac{1}{2}F(\mathbf{x}) = \frac{1}{2}I(\mathbf{x}) = \sum_{\{i,j\} \in E} w_{ij} \frac{|x_i - x_j|}{2} = \text{cut}(S). \quad (2.7)$$

Theorem 2.1. *The maxcut problem (1.1) can be rewritten into*

$$\max_{S \subset V} \text{cut}(S) = \frac{1}{2} \max_{\mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}} F(\mathbf{x}), \quad (2.8)$$

and any vector \mathbf{x}^* reaching the maximum of $F(\mathbf{x})$ produces a maxcut (S, S') where the subset S satisfies $S^+(\mathbf{x}^*) \subset S \subset (S^-(\mathbf{x}^*))^c$.

Proof. Combining $\text{cut}(\emptyset) = 0$ and Eq. (2.7) for the nonempty set situation together leads directly to

$$\frac{1}{2} \max_{\mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}} F(\mathbf{x}) \geq \max_{S \subset V} \text{cut}(S). \quad (2.9)$$

On the other hand, suppose \mathbf{x}^* is a maximizer of $F(\mathbf{x})$ on $\mathbb{R}^n \setminus \{\mathbf{0}\}$, i.e.

$$\frac{1}{2}F(\mathbf{x}^*) = \frac{1}{2} \max_{\mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}} F(\mathbf{x}). \quad (2.10)$$

For any S^* satisfying $S^+(\mathbf{x}^*) \subset S^* \subset (S^-(\mathbf{x}^*))^c$, there exists $\hat{\mathbf{x}} \in M(\mathbf{x}^*)$ defined in Eq. (2.6) such that

$$\frac{\hat{\mathbf{x}}}{\|\hat{\mathbf{x}}\|_\infty} = \mathbf{1}_{S^*} - \mathbf{1}_{V \setminus S^*} \quad (2.11)$$

also maximizes $F(\mathbf{x})$ on $\mathbb{R}^n \setminus \{\mathbf{0}\}$ thanks to the zeroth-order homogeneousness of $F(\mathbf{x})$. That is, we have

$$\frac{1}{2} \max_{\mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}} F(\mathbf{x}) = \frac{1}{2}F\left(\frac{\hat{\mathbf{x}}}{\|\hat{\mathbf{x}}\|_\infty}\right) = \text{cut}(S^*) \leq \max_{S \subset V} \text{cut}(S), \quad (2.12)$$

where Eq. (2.7) has been applied. The proof is finished as a result of Eqs. (2.9) and (2.12). \square

Theorem 2.1 establishes an equivalent continuous optimization for the maxcut problem which will serve as the cornerstone of the subsequent design of an iterative algorithm. Now we only need to consider

$$r_{\max} = \max_{\mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}} F(\mathbf{x}). \quad (2.13)$$

As shown in Theorem 2.2, it can be solved via the following so-called Dinkelbach iterative scheme [7]:

$$\begin{cases} x^{k+1} = \arg \min_{\|\mathbf{x}\|_p=1} \{r^k \|\mathbf{x}\|_\infty - I(\mathbf{x})\}, & p \in [1, \infty], \\ r^{k+1} = F(\mathbf{x}^{k+1}), \end{cases} \quad (2.14a)$$

$$(2.14b)$$

where $\|\cdot\|_p$ denotes the standard p -norm in \mathbb{R}^n , i.e.

$$\|\mathbf{x}\|_p = (|x_1|^p + |x_2|^p + \cdots + |x_n|^p)^{\frac{1}{p}}. \quad (2.15)$$

Theorem 2.2 (Global Convergence). *The sequence $\{r^k\}$ generated by the iterative scheme (2.14) from any initial point $\mathbf{x}^0 \in \mathbb{R}^n \setminus \{\mathbf{0}\}$ increases monotonically to the global maximum r_{\max} .*

Proof. The definition of \mathbf{x}^{k+1} (see Eq. (2.14a)) implies

$$r^k \|\mathbf{x}\|_\infty - I(\mathbf{x}) \geq r^k \|\mathbf{x}^{k+1}\|_\infty - I(\mathbf{x}^{k+1}), \quad \forall \mathbf{x} \text{ s.t. } \|\mathbf{x}\|_p = 1,$$

and substituting $\mathbf{x} = \mathbf{x}^k$ into the above inequality yields

$$0 = r^k \|\mathbf{x}^k\|_\infty - I(\mathbf{x}^k) \geq r^k \|\mathbf{x}^{k+1}\|_\infty - I(\mathbf{x}^{k+1}),$$

which means

$$r^k \leq r^{k+1} \leq r_{\max}, \quad \forall k \in \mathbb{N}^+.$$

Therefore,

$$\exists r^* \in [0, r_{\max}] \text{ s.t. } \lim_{k \rightarrow +\infty} r^k = r^*,$$

and it suffices to show $r_{\max} \leq r^*$. To this end, we denote

$$f(r) = \min_{\|\mathbf{x}\|_p=1} (r \|\mathbf{x}\|_\infty - I(\mathbf{x})),$$

which must be continuous on \mathbb{R} by the compactness of the unit closed sphere

$$S_p = \{\mathbf{x} \in \mathbb{R}^n : \|\mathbf{x}\|_p = 1\}.$$

Note that

$$\begin{aligned} f(r^k) &= r^k \|\mathbf{x}^{k+1}\|_\infty - I(\mathbf{x}^{k+1}) \\ &= r^k \|\mathbf{x}^{k+1}\|_\infty - r^{k+1} \|\mathbf{x}^{k+1}\|_\infty \\ &= \|\mathbf{x}^{k+1}\|_\infty (r^k - r^{k+1}) \rightarrow 0 \text{ as } k \rightarrow +\infty, \end{aligned}$$

then we have

$$f(r^*) = \lim_{k \rightarrow +\infty} f(r^k) = 0,$$

which implies

$$r^* \|\mathbf{x}\|_\infty - I(\mathbf{x}) \geq 0, \quad \forall \mathbf{x} \text{ with } \|\mathbf{x}\|_p = 1.$$

Hence, $\forall \mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}$,

$$F(\mathbf{x}) = \frac{I(\mathbf{x})}{\|\mathbf{x}\|_\infty} = \frac{I(\mathbf{x})/\|\mathbf{x}\|_p}{\|\mathbf{x}\|_\infty/\|\mathbf{x}\|_p} = \frac{I(\hat{\mathbf{x}})}{\|\hat{\mathbf{x}}\|_\infty} \leq r^*,$$

where $\hat{\mathbf{x}} = \mathbf{x}/\|\mathbf{x}\|_p$. \square

Although the inner subproblem (2.14a) is not only non-convex but also non-solvable in polynomial time due to the NP-hardness of the maxcut problem, the Dinkelbach scheme (2.14) provides us a good starting point to a feasible iterative algorithm for the maxcut problem.

Remark 2.1. Obviously, the equivalent continuous optimization (2.13) has a fractional form, i.e. a ratio between two convex functions, but such kind of fractions have been hardly touched in the field of fractional programming [17], where concave optimization problems, like optimizing the ratio of a concave function to a convex one, are usually considered.

3. A Simple Iterative Algorithm

The non-convex subproblem (2.14a) brings us a significant insight to deal with a relaxed subproblem alternatively, though it can not be solved in polynomial time.

Denote the subgradient of $I(\mathbf{x})$ by [6]

$$\partial I(\mathbf{x}) = \left\{ \mathbf{s} = (s_1, \dots, s_n) \mid s_i = \sum_{j:\{i,j\} \in E} w_{ij} z_{ij}, z_{ij} \in \text{Sgn}(x_i - x_j) \text{ and } z_{ij} = -z_{ji} \right\}, \quad (3.1)$$

where we have extended the sign function (note that $\text{sign}(0) = 1$ here)

$$\text{sign}(t) = \begin{cases} 1, & \text{if } t \geq 0, \\ -1, & \text{if } t < 0, \end{cases} \quad (3.2)$$

into a set-valued function

$$\text{Sgn}(t) = \begin{cases} \{1\}, & \text{if } t > 0, \\ \{-1\}, & \text{if } t < 0, \\ [-1, 1], & \text{if } t = 0. \end{cases} \quad (3.3)$$

For $\mathbf{x} = (x_1, x_2, \dots, x_n) \in \mathbb{R}^n$, we are able to define corresponding vectorized versions in an element-wise manner

$$\text{sign}(\mathbf{x}) = (s_1, s_2, \dots, s_n), \quad s_i = \text{sign}(x_i), \quad i = 1, 2, \dots, n, \quad (3.4)$$

$$\text{Sgn}(\mathbf{x}) = \{(s_1, s_2, \dots, s_n) \mid s_i \in \text{Sgn}(x_i), i = 1, 2, \dots, n\}. \quad (3.5)$$

Since the function $I(\cdot)$ is convex, it holds

$$I(\mathbf{x}) \geq I(\mathbf{y}) + (\mathbf{x} - \mathbf{y}, \mathbf{s}), \quad \forall \mathbf{s} \in \partial I(\mathbf{y}), \quad \forall \mathbf{x}, \mathbf{y} \in \mathbb{R}^n, \quad (3.6)$$

where (\cdot, \cdot) denotes the standard inner product in \mathbb{R}^n . Further considering the fact that $I(\cdot)$ is homogeneous of degree one, we have

$$I(\mathbf{y}) = (\mathbf{y}, \mathbf{s}), \quad \forall \mathbf{s} \in \partial I(\mathbf{y}), \quad \forall \mathbf{y} \in \mathbb{R}^n, \quad (3.7)$$

and

$$I(\mathbf{x}) \geq I(\mathbf{y}) + (\mathbf{x} - \mathbf{y}, \mathbf{s}) = (\mathbf{x}, \mathbf{s}), \quad \forall \mathbf{s} \in \partial I(\mathbf{y}), \quad \forall \mathbf{x}, \mathbf{y} \in \mathbb{R}^n. \quad (3.8)$$

Plugging the relaxation (3.8) into the subproblem (2.14a) modifies the two-step Dinkelbach iterative scheme into the following three-step one:

$$\begin{cases} \mathbf{x}^{k+1} = \arg \min_{\|\mathbf{x}\|_p=1} \{r^k \|\mathbf{x}\|_\infty - (\mathbf{x}, \mathbf{s}^k)\}, & p \in [1, \infty], & (3.9a) \\ r^{k+1} = F(\mathbf{x}^{k+1}), & & (3.9b) \\ \mathbf{s}^{k+1} \in \partial I(\mathbf{x}^{k+1}) & & (3.9c) \end{cases}$$

with an initial data $\mathbf{x}^0 \in \mathbb{R}^n \setminus \{\mathbf{0}\}$, $r^0 = F(\mathbf{x}^0)$ and $\mathbf{s}^0 \in \partial I(\mathbf{x}^0)$.

It can be readily verified that the subproblem (3.9a) is convex via the relaxation (3.8) from Eq. (2.14a). More importantly, we can write down a solution to the inner subproblem (3.9a) in an analytical manner (see Theorem 3.1). That is, no any other optimization solver is needed in (3.9) and so that a simple iterative algorithm we name it. Actually, we are able to prove that such iterative scheme (3.9) still keeps the monotonicity (see Theorem 3.2) and has local convergence (see Theorem 3.4).

3.1. Exact solution to the inner subproblem

Let

$$L(r, \mathbf{v}) := \min_{\|\mathbf{x}\|_p=1} \{r \|\mathbf{x}\|_\infty - (\mathbf{x}, \mathbf{v})\}, \quad r \in \mathbb{R}, \quad \mathbf{v} \in \mathbb{R}^n, \quad (3.10)$$

denote the minimal value in Eq. (3.9a). In order to obtain the exact solution, we need to show first a property of the simple iteration (3.9) (see Lemma 3.1), and use it to narrow the scope of discussion.

Lemma 3.1. *Suppose \mathbf{x}^k , r^k and \mathbf{s}^k are generated by the simple iteration (3.9). Then, for $k \geq 1$, we always have $r^k \leq \|\mathbf{s}^k\|_1$. In particular,*

- (1) $r^k = \|\mathbf{s}^k\|_1$ if and only if $\mathbf{x}^k / \|\mathbf{x}^k\|_\infty \in \text{Sgn}(\mathbf{s}^k)$.
- (2) $r^k < \|\mathbf{s}^k\|_1$ if and only if $L(r^k, \mathbf{s}^k) < 0$.

Proof. From Hölder's inequality,

$$r^k \|\mathbf{x}\|_\infty - (\mathbf{x}, \mathbf{s}^k) \geq r^k \|\mathbf{x}\|_\infty - \|\mathbf{s}^k\|_1 \|\mathbf{x}\|_\infty = (r^k - \|\mathbf{s}^k\|_1) \|\mathbf{x}\|_\infty, \quad (3.11)$$

where the equality holds if and only if $\mathbf{x} \in \|\mathbf{x}\|_\infty \text{Sgn}(\mathbf{s}^k)$. Plugging $\mathbf{x} = \mathbf{x}^k$ into Eq. (3.11) and using Eq. (3.7) lead to

$$(r^k - \|\mathbf{s}^k\|_1) \|\mathbf{x}^k\|_\infty \leq r^k \|\mathbf{x}^k\|_\infty - (\mathbf{x}^k, \mathbf{s}^k) = r^k \|\mathbf{x}^k\|_\infty - I(\mathbf{x}^k) = 0,$$

and thus $r^k \leq \|\mathbf{s}^k\|_1$.

Meanwhile, the statement (1) corresponds to the situation in which the equality holds.

Next we will show the statement (2) is true. On one hand, if $L(r^k, \mathbf{s}^k) < 0$, then there exists a vector \mathbf{x}^{k+1} satisfying

$$L(r^k, \mathbf{s}^k) = r^k \|\mathbf{x}^{k+1}\|_\infty - (\mathbf{x}^{k+1}, \mathbf{s}^k) < 0. \quad (3.12)$$

Substituting $\mathbf{x} = \mathbf{x}^{k+1}$ into Eq. (3.11) and using Eq. (3.12) yield

$$(r^k - \|\mathbf{s}^k\|_1)\|\mathbf{x}^{k+1}\|_\infty \leq r^k\|\mathbf{x}^{k+1}\|_\infty - (\mathbf{x}^{k+1}, \mathbf{s}^k) = L(r^k, \mathbf{s}^k) < 0,$$

and thus $r^k < \|\mathbf{s}^k\|_1$.

On the other hand, assume $r^k < \|\mathbf{s}^k\|_1$ holds. Choose $\mathbf{y} \in \text{Sgn}(\mathbf{s}^k)$ and let $\mathbf{x}^* = \mathbf{y}/\|\mathbf{y}\|_p$. It is obvious that $\|\mathbf{y}\|_\infty = 1$, $\|\mathbf{x}^*\|_p = 1$ and $(\mathbf{y}, \mathbf{s}^k) = \|\mathbf{s}^k\|_1$. In consequence, we have

$$\begin{aligned} L(r^k, \mathbf{s}^k) &\leq r^k\|\mathbf{x}^*\|_\infty - (\mathbf{x}^*, \mathbf{s}^k) \\ &= r^k\left\|\frac{\mathbf{y}}{\|\mathbf{y}\|_p}\right\|_\infty - \left(\frac{\mathbf{y}}{\|\mathbf{y}\|_p}, \mathbf{s}^k\right) \\ &= \frac{1}{\|\mathbf{y}\|_p}(r^k - \|\mathbf{s}^k\|_1) < 0. \end{aligned}$$

The proof is complete. \square

Given a real number $r > 0$ and a vector $\mathbf{v} = (v_1, \dots, v_n) \in \mathbb{R}^n$, in view of Lemma 3.1 and the subproblem (3.9a), we only need to consider the minimization problem

$$\mathbf{x}^* = \arg \min_{\|\mathbf{x}\|_p=1} \{r\|\mathbf{x}\|_\infty - (\mathbf{x}, \mathbf{v})\} \quad (3.13)$$

under the condition

$$0 < r \leq \|\mathbf{v}\|_1. \quad (3.14)$$

Without loss of generality, it suffices to discuss an ordered situation

$$|v_1| \geq |v_2| \geq \dots \geq |v_n| \geq |v_{n+1}| = 0, \quad (3.15)$$

where we have extended the index set into $\{1, \dots, n+1\}$ and introduced an auxiliary element $v_{n+1} = 0$ for convenience. Consider the accumulation of increment

$$A(m) = \sum_{j=1}^m (|v_j| - |v_{m+1}|), \quad m \in \{1, 2, \dots, n\}, \quad (3.16)$$

and then from Eq. (3.15) we have

$$0 = A(0) \leq A(1) \leq A(2) \leq \dots \leq A(n) = \|\mathbf{v}\|_1. \quad (3.17)$$

Meanwhile, we need a key step to increase the regularity for $p \in (1, \infty)$ through rewriting the minimization problem (3.13) into

$$\min_{\|\mathbf{x}\|_p=1} \{r\|\mathbf{x}\|_\infty - (\mathbf{x}, \mathbf{v})\} = \min_{\|\mathbf{u}\|_p=1} \{r\|\mathbf{u}\|_\infty - (\mathbf{u}, |\mathbf{v}|)\} \quad (3.18)$$

$$= \min_{\mathbf{w} \neq \mathbf{0}} \frac{r\|\mathbf{w}\|_\infty - (\mathbf{w}, |\mathbf{v}|)}{\|\mathbf{w}\|_p} \quad (3.19)$$

$$= \min_{\|\mathbf{z}\|_\infty=1} \frac{r - (\mathbf{z}, |\mathbf{v}|)}{\|\mathbf{z}\|_p} \quad (3.20)$$

$$= \min_{\mathbf{z} \in B_\infty} G(\mathbf{z}), \quad (3.21)$$

where

$$G(\mathbf{z}) = \frac{r - (\mathbf{z}, |\mathbf{v}|)}{\|\mathbf{z}\|_p}, \quad (3.22)$$

$$B_\infty = \{\mathbf{z} \in \mathbb{R}^n : \|\mathbf{z}\|_\infty \leq 1\}, \quad (3.23)$$

and the absolute value is taken in an element-wise manner, e.g., $|\mathbf{v}| = (|v_1|, |v_2|, \dots, |v_n|)$. Here we have used extensively the fact that $\|\cdot\|_p$, $\|\cdot\|_\infty$ and $(\cdot, |\mathbf{v}|)$ are all homogeneous of degree one, and in Eq. (3.21) the fact that $G(\mathbf{z})$ achieves its minimum on the boundary of the feasible region B_∞ due to

$$G(\lambda\mathbf{z}) - G(\mathbf{z}) = \left(\frac{1}{\lambda} - 1\right) \frac{r}{\|\mathbf{z}\|_p} < 0 \quad \text{for } \lambda > 1. \quad (3.24)$$

If one denotes the corresponding minimizers in Eqs. (3.18)-(3.21) by \mathbf{x}^* , \mathbf{u}^* , \mathbf{w}^* , and \mathbf{z}^* , respectively, then we have

$$\mathbf{u}^* = \text{sign}(\mathbf{v}) \cdot \mathbf{x}^*, \quad (3.25)$$

$$\mathbf{w}^* = \|\mathbf{w}^*\|_p \mathbf{u}^*, \quad (3.26)$$

$$\mathbf{z}^* = \frac{\mathbf{w}^*}{\|\mathbf{w}^*\|_\infty}, \quad (3.27)$$

where $\mathbf{x} \cdot \mathbf{y}$ in Eq. (3.25) denotes element-by-element multiplication of vectors \mathbf{x} and \mathbf{y} . It is obvious that we only need to search for the minimizer \mathbf{z}^* to the minimization problem (3.21), with which we are able to reach our target

$$\mathbf{x}^* = \frac{\text{sign}(\mathbf{v}) \cdot \mathbf{z}^*}{\|\mathbf{z}^*\|_p} \quad (3.28)$$

by virtue of the one-to-one mappings (3.25)-(3.27).

According to the condition (3.14), the rest of the discussion falls into the following three scenarios. Before that, we need the following lemma to characterize the minimizer of $G(\mathbf{z})$ on B_∞ , denoted by $\mathbf{z}^* = (z_1^*, z_2^*, \dots, z_n^*)$.

Lemma 3.2. *Let $\mathbf{z}^* = (z_1^*, z_2^*, \dots, z_n^*)$ be the minimizer of $G(\mathbf{z})$ on B_∞ . If $r < \|\mathbf{v}\|_1$ and $1 \leq p < \infty$, then \mathbf{z}^* has not only nonnegative elements, but also the same order as $|\mathbf{v}|$.*

Proof. Since $\mathbf{z}^* = (z_1^*, z_2^*, \dots, z_n^*)$ is the minimizer of $G(\mathbf{z})$ on B_∞ , it holds

$$G(\mathbf{z}^*) \leq G(\mathbf{z}), \quad \forall \mathbf{z} \in B_\infty.$$

In the following, the proof is by contradiction and split into three steps.

First, the condition $r < \|\mathbf{v}\|_1$ directly implies that \mathbf{z}^* satisfies

$$r - (\mathbf{z}^*, |\mathbf{v}|) = G(\mathbf{z}^*) \|\mathbf{z}^*\|_p \leq G(\mathbf{1}) \|\mathbf{z}^*\|_p = \frac{\|\mathbf{z}^*\|_p}{\|\mathbf{1}\|_p} (r - \|\mathbf{v}\|_1) < 0, \quad (3.29)$$

where $\mathbf{1} \in B_\infty$.

Second, if there exists some $i \in \{1, \dots, n\}$ such that $z_i^* < 0$, and let $\hat{\mathbf{z}} = \{\hat{z}_1, \dots, \hat{z}_n\}$ be a vector satisfying: $\hat{z}_i = 0$ and $\hat{z}_j = z_j^*$ for $j \neq i$, then we have $\hat{\mathbf{z}} \in B_\infty$ and

$$G(\hat{\mathbf{z}}) = \frac{r - (\hat{\mathbf{z}}, |\mathbf{v}|)}{\|\hat{\mathbf{z}}\|_p} \leq \frac{r - (\mathbf{z}^*, |\mathbf{v}|)}{\|\hat{\mathbf{z}}\|_p} < \frac{r - (\mathbf{z}^*, |\mathbf{v}|)}{\|\mathbf{z}^*\|_p} = G(\mathbf{z}^*) \quad (3.30)$$

from (3.29). This contradicts to the fact that \mathbf{z}^* is the minimizer of $G(\mathbf{z})$ on B_∞ .

Third, if there exists $i, j \in \{1, \dots, n\}$ such that $(z_i^* - z_j^*)(|v_i| - |v_j|) < 0$, let $\hat{\mathbf{z}}$ be a vector obtained by exchanging the i -th and the j -th elements of \mathbf{z} , then $\hat{\mathbf{z}} \in B_\infty$, $\|\hat{\mathbf{z}}\|_p = \|\mathbf{z}^*\|_p$, and

$$G(\hat{\mathbf{z}}) - G(\mathbf{z}^*) = \frac{(z_i^* - z_j^*)(|v_i| - |v_j|)}{\|\mathbf{z}\|_p} < 0. \quad (3.31)$$

This also contradicts with the fact that \mathbf{z}^* achieves the minimum of $G(\mathbf{z})$ on B_∞ . \square

More importantly, under the conditions of Lemma 3.2, the objective function $G(\mathbf{z})$ is now differentiable with the i -th partial derivative being

$$\frac{\partial G(\mathbf{z}^*)}{\partial z_i} = -\frac{|v_i| \|\mathbf{z}^*\|_p + (r - (\mathbf{z}^*, |\mathbf{v}|)) \|\mathbf{z}^*\|_p^{1-p} (z_i^*)^{p-1}}{\|\mathbf{z}^*\|_p^2}. \quad (3.32)$$

• **Scenario 1:**

$$r < \|\mathbf{v}\|_1 \quad \text{and} \quad 1 < p < \infty. \quad (3.33)$$

In view of Lemma 3.2 and the decreasing order of $|\mathbf{v}|$ described in Eq. (3.15), we may assume that there exists an integer $m_0 \in \{1, 2, \dots, n\}$ such that \mathbf{z}^* satisfies

$$1 = z_1^* = z_2^* = \dots = z_{m_0}^* > z_{m_0+1}^* \geq \dots \geq z_n^* \geq 0. \quad (3.34)$$

Let

$$T = \{\mathbf{t} \in \mathbb{R}^n : \mathbf{z}^* + \epsilon \mathbf{t} \in B_\infty \text{ for sufficiently small } \epsilon > 0\} \quad (3.35)$$

denote the tangent cone of B_∞ at \mathbf{z}^* .

From Eq. (3.34), the tangent cone can be readily rewritten into

$$T = \{\mathbf{t} = (t_1, t_2, \dots, t_n) \in \mathbb{R}^n : t_i \leq 0, \quad i = 1, 2, \dots, m_0\}. \quad (3.36)$$

Since the minimizer \mathbf{z}^* achieves a local minimum of $G(\mathbf{z})$ on B_∞ , according to Eq. (3.36), we have

$$\frac{\partial G(\mathbf{z}^*)}{\partial \mathbf{t}} \geq 0, \quad \forall \mathbf{t} \in T \quad (3.37)$$

$$\Leftrightarrow \begin{cases} \frac{\partial G(\mathbf{z}^*)}{\partial z_i} \leq 0, & i = 1, 2, \dots, m_0, \\ \frac{\partial G(\mathbf{z}^*)}{\partial z_i} = 0, & i = m_0 + 1, \dots, n. \end{cases} \quad (3.38)$$

For the sake of subsequent discussion, it is convenient to introduce the following three auxiliary quantities:

$$\alpha = \sum_{i=1}^{m_0} |v_i| - r, \quad (3.39)$$

$$\beta = \sum_{i=m_0+1}^n z_i^* |v_i|, \quad (3.40)$$

$$\gamma = \sum_{i=m_0+1}^n (z_i^*)^p, \quad (3.41)$$

and then

$$\alpha + \beta = (\mathbf{z}^*, |\mathbf{v}|) - r > 0, \quad (3.42)$$

$$m_0 + \gamma = \|\mathbf{z}^*\|_p^p > 0. \quad (3.43)$$

Therefore, the partial derivative (3.32) becomes

$$\frac{\partial G(\mathbf{z}^*)}{\partial z_i} = \frac{1}{\|\mathbf{z}^*\|_p} \left(\frac{\alpha + \beta}{m_0 + \gamma} (z_i^*)^{p-1} - |v_i| \right), \quad (3.44)$$

and then Eq. (3.38) directly implies

$$\begin{cases} (z_i^*)^{p-1} \leq \frac{m_0 + \gamma}{\alpha + \beta} |v_i|, & i = 1, 2, \dots, m_0, \\ (z_i^*)^{p-1} = \frac{m_0 + \gamma}{\alpha + \beta} |v_i|, & i = m_0 + 1, \dots, n. \end{cases} \quad (3.45)$$

Substituting Eq. (3.45) into Eq. (3.41) and using Eq. (3.40) yields

$$\gamma = \sum_{i=m_0+1}^n (z_i^*)^{p-1} z_i^* = \sum_{i=m_0+1}^n \frac{m_0 + \gamma}{\alpha + \beta} |v_i| z_i^* = \frac{m_0 + \gamma}{\alpha + \beta} \beta, \quad (3.46)$$

and then, from $m_0 > 0$, we have

$$\frac{m_0 + \gamma}{\alpha + \beta} = \frac{m_0}{\alpha}, \quad \alpha > 0. \quad (3.47)$$

That is, the condition for local minimizers, Eq. (3.38) or Eq. (3.45), can be further simplified into

$$\begin{cases} (z_i^*)^{p-1} \leq a_i, & i = 1, 2, \dots, m_0, \\ (z_i^*)^{p-1} = a_i, & i = m_0 + 1, \dots, n, \end{cases} \quad (3.48)$$

where

$$a_i = \frac{m_0}{\alpha} |v_i| = \frac{m_0 |v_i|}{\sum_{j=1}^{m_0} |v_j| - r}, \quad i = 1, 2, \dots, n. \quad (3.49)$$

Combining Eqs. (3.34) and (3.48) determines the minimizer $\mathbf{z}^* = (z_1^*, z_2^*, \dots, z_n^*)$

$$z_i^* = \min \left\{ 1, a_i^{\frac{1}{p-1}} \right\}, \quad i = 1, 2, \dots, n, \quad (3.50)$$

and then we are able to reach our target \mathbf{x}^* by Eq. (3.28).

Therefore the exact solution to the inner subproblem for Scenario 1 has been obtained. The only remaining thing is how to determine m_0 efficiently used in Eq. (3.49). This can be achieved with the help of the accumulation function $A(m)$ defined in Eq. (3.16). Namely, m_0 is the smallest integer m satisfying $A(m) > r$

$$m_0 = \min \{ m \in \{1, 2, \dots, n\} : A(m) > r \}, \quad (3.51)$$

which can be readily verified as follows:

$$\begin{aligned} a_{m_0} &\geq (z_{m_0}^*)^{p-1} = 1 > (z_{m_0+1}^*)^{p-1} = a_{m_0+1} \\ \Leftrightarrow \frac{m_0 |v_{m_0}|}{\sum_{j=1}^{m_0} |v_j| - r} &\geq 1 > \frac{m_0 |v_{m_0+1}|}{\sum_{j=1}^{m_0} |v_j| - r} \\ \Leftrightarrow \sum_{j=1}^{m_0} (|v_j| - |v_{m_0+1}|) &> r \geq \sum_{j=1}^{m_0-1} (|v_j| - |v_{m_0}|) \\ \Leftrightarrow A(m_0) &> r \geq A(m_0 - 1). \end{aligned}$$

• **Scenario 2:**

$$r < \|\mathbf{v}\|_1 \quad \text{and} \quad p = 1. \quad (3.52)$$

In this scenario, we still have the minimizer $\mathbf{z}^* = (z_1^*, z_2^*, \dots, z_n^*)$ has nonnegative elements according to Lemma 3.2, namely, $\forall i \in \{1, 2, \dots, n\}, 0 \leq z_i^* \leq 1$.

Let m_1 be the largest integer satisfying $A(m-1) < r$, i.e.

$$m_1 = \max \{m \in \{1, 2, \dots, n\} : A(m-1) < r\}, \quad (3.53)$$

where $A(m)$ is the accumulation function defined by Eq. (3.16). Comparing Eq. (3.53) with Eq. (3.51), we have $m_1 \leq m_0$ where m_0 is defined by Eq. (3.51), and specifically

$$A(m_1 - 1) < r = A(m_1) = A(m_1 + 1) = \dots = A(m_0 - 1) < A(m_0), \quad (3.54)$$

thereby indicating

$$|v_{m_1}| > |v_{m_1+1}| = \dots = |v_{m_0}| > |v_{m_0+1}|. \quad (3.55)$$

For $i \in \{1, 2, \dots, n\}$, consider a continuous function on $[0, 1]$

$$G_i(t) = G(z_1^*, \dots, z_{i-1}^*, t, z_{i+1}^*, \dots, z_n^*) = -|v_i| + \frac{r - \sum_{j \neq i} z_j^* (|v_j| - |v_i|)}{\sum_{j \neq i} |z_j^*| + t}. \quad (3.56)$$

Since \mathbf{z}^* is a minimizer of $G(\mathbf{z})$ on B_∞ , it can be readily verified that $\min_{t \in [0, 1]} G_i(t) = G(\mathbf{z}^*)$, i.e.

$$z_i^* \in \arg \min_{t \in [0, 1]} G_i(t), \quad \forall i \in \{1, 2, \dots, n\}, \quad (3.57)$$

from which we are able to determine z_i^* . The related discussion needs to split the index set into the following three cases:

- (1) For $1 \leq i \leq m_1$, we claim that $G_i(t)$ is a strictly monotonically decreasing function due to $r - \sum_{j \neq i} z_j^* (|v_j| - |v_i|) > 0$ and thus $z_i^* = 1$. The verification shown below is in a straightforward manner

$$\begin{aligned} r - \sum_{j \neq i} z_j^* (|v_j| - |v_i|) &\geq r - \sum_{j=1}^{i-1} z_j^* (|v_j| - |v_i|) \\ &\geq r - \sum_{j=1}^{i-1} (|v_j| - |v_i|) \\ &= r - A(i-1) > 0. \end{aligned}$$

- (2) For $m_0 < i \leq n$, we claim that $z_i^* = 0$. Suppose the contrary that $z_{i_0}^* > 0$ is the positive element with the largest index and $i_0 > m_0$. Then the order given in Eqs. (3.15) and (3.55) directly implies

$$\begin{aligned} r - \sum_{j \neq i_0} z_j^* (|v_j| - |v_{i_0}|) &= r - \sum_{j=1}^{m_1} (|v_j| - |v_{i_0}|) - \sum_{j=m_1+1}^{i_0-1} z_j^* (|v_j| - |v_{i_0}|) \\ &\leq r - \sum_{j=1}^{m_1} (|v_j| - |v_{i_0}|) \end{aligned}$$

$$\begin{cases} < r - \sum_{j=1}^{m_1} (|v_j| - |v_{m_1+1}|) = r - A(m_1) = 0, & \text{if } m_1 < m_0, \\ \leq r - \sum_{j=1}^{m_1} (|v_j| - |v_{m_1+1}|) = r - A(m_0) < 0, & \text{if } m_1 = m_0, \end{cases}$$

and thus we have $G_{i_0}(t)$ is a strictly monotonically increasing function. That is, the minimizer of $G_{i_0}(t)$ on $[0, 1]$ is 0, i.e. $z_{i_0}^* = 0$, which is obviously a contradiction.

(3) For $m_1 < i \leq m_0$, using the results for above two cases and the order (3.55) yields

$$\begin{aligned} r - \sum_{j \neq i} z_j^* (|v_j| - |v_i|) &= r - \sum_{j=1}^{m_1} (|v_j| - |v_i|) - \sum_{j=m_1+1}^{m_0} z_j^* (|v_j| - |v_i|) \\ &= r - \sum_{j=1}^{m_1} (|v_j| - |v_{m_1+1}|) = r - A(m_1) = 0, \end{aligned}$$

and thus we have $G_i(t)$ is a constant function on $[0, 1]$, i.e. $G_i(t) \equiv -|v_i|$. That is, the minimizer z_i^* can take any value in $[0, 1]$.

In a word, the minimizers $\mathbf{z}^* = (z_1^*, \dots, z_n^*)$ for Scenario 2 constitute the following set:

$$\{(z_1^*, \dots, z_n^*) \in [0, 1]^n \mid z_i^* = 1, i = 1, \dots, m_1, \text{ and } z_i^* = 0, i = m_0 + 1, \dots, n\}, \quad (3.58)$$

and thus \mathbf{x}^* can be also obtained through Eq. (3.28).

• **Scenario 3:**

$$r = \|\mathbf{v}\|_1 \quad \text{or} \quad p = \infty. \quad (3.59)$$

For $r = \|\mathbf{v}\|_1$, according to Lemma 3.1, we have that \mathbf{x}^* is a minimizer of problem (3.13) and only if

$$\frac{\mathbf{x}^*}{\|\mathbf{x}^*\|_\infty} \in \text{Sgn}(\mathbf{v}), \quad \|\mathbf{x}^*\|_p = 1. \quad (3.60)$$

For $p = \infty$, the minimization problem Eq. (3.13) becomes a linear optimization problem

$$\mathbf{x}^* = \arg \min_{\|\mathbf{x}\|_\infty=1} \{r - (\mathbf{x}, \mathbf{v})\}, \quad (3.61)$$

the solution of which can be still represented by Eq. (3.60). Hence, the solution Eq. (3.60) solves the minimization problem (3.13) for Scenario 3.

Summarizing the above analysis for three scenarios, we have figured out the exact solution to the inner subproblem (3.9a), as stated in the following theorem.

Theorem 3.1 (Exact Solution). *The solution of the minimization problem (3.13) under the condition (3.14) can be expressed analytically in Eqs. (3.28) and (3.50) for $r < \|\mathbf{v}\|_1$, and $1 < p < \infty$, Eqs. (3.28) and (3.58) for $r < \|\mathbf{v}\|_1$, and $p = 1$, and Eq. (3.60) otherwise.*

Hereafter we use a set denoted by X_p^{k+1} to collect all those analytical solutions of the inner subproblem (3.9a), i.e. $\mathbf{x}^{k+1} \in X_p^{k+1}$. Obviously, X_p^{k+1} is closed.

Using the above analytical solution, we are able to get a lower bound for $F(\mathbf{x})$ below, which will be useful in the subsequent convergence analysis.

Corollary 3.1. *Under the condition (3.14) the minimizer \mathbf{x}^* to the problem (3.13) satisfies*

$$\frac{(\mathbf{x}^*, \mathbf{v})}{\|\mathbf{x}^*\|_\infty} \geq \sum_{i=1}^m |v_i|, \quad (3.62)$$

where $m = m_0$ for both Scenario 1 and Scenario 2, and $m = n$ for Scenario 3.

Proof. For both Scenario 1 and Scenario 2, there exists a minimizer $\mathbf{z}^* = (z_1^*, z_2^*, \dots, z_n^*) \in B_\infty \cap \mathbb{R}_+^n$ to the equivalent problem (3.21), which satisfies

$$z_1^* = \dots = z_{m_0}^* = 1,$$

according to Eqs. (3.50) and (3.58), respectively. Then, from Eq. (3.28), we have

$$\frac{(\mathbf{x}^*, \mathbf{v})}{\|\mathbf{x}^*\|_\infty} = (\text{sign}(\mathbf{v}) \cdot \mathbf{z}^*, \mathbf{v}) = (\mathbf{z}^*, |\mathbf{v}|) \geq \sum_{i=1}^{m_0} |v_i|. \quad (3.63)$$

As for Scenario 3, using Eq. (3.60), it can be easily verified that

$$\frac{(\mathbf{x}^*, \mathbf{v})}{\|\mathbf{x}^*\|_\infty} = \left(\frac{\mathbf{x}^*}{\|\mathbf{x}^*\|_\infty}, \mathbf{v} \right) = (\text{Sgn}(\mathbf{v}), \mathbf{v}) = \|\mathbf{v}\|_1 = \sum_{i=1}^n |v_i|. \quad (3.64)$$

Thus the proof is finished. \square

3.2. Subgradient selection and convergence analysis

Besides the inner subproblem (3.9a), another key issue to implement the simple iteration (3.9) is how to choose the subgradient (see Eq. (3.9c)). For a general selection, according to Eqs. (3.7)-(3.9), we are able to obtain $r^k \leq r^{k+1}$, because

$$\begin{aligned} 0 &= r^k \|\mathbf{x}^k\|_\infty - I(\mathbf{x}^k) = r^k \|\mathbf{x}^k\|_\infty - (\mathbf{x}^k, \mathbf{s}^k) \\ &\geq r^k \|\mathbf{x}^{k+1}\|_\infty - (\mathbf{x}^{k+1}, \mathbf{s}^k) \geq r^k \|\mathbf{x}^{k+1}\|_\infty - I(\mathbf{x}^{k+1}) \\ &= \|\mathbf{x}^{k+1}\|_\infty (r^k - r^{k+1}). \end{aligned}$$

Theorem 3.2 (Monotonicity). *The sequence $\{r^k\}$ generated by the iterative scheme (3.9) from any initial point $\mathbf{x}^0 \in \mathbb{R}^n \setminus \{\mathbf{0}\}$ increases monotonically.*

The monotone increasing in Theorem 3.2 is also a direct consequence of Lemma 3.1 and Corollary 3.1 because there exists an index set $\iota \subset \{1, 2, \dots, n\}$, the size of which should be m in Eq. (3.62) (the upper bound of summation index) such that

$$r^{k+1} = \frac{I(\mathbf{x}^{k+1})}{\|\mathbf{x}^{k+1}\|_\infty} \geq \frac{(\mathbf{x}^{k+1}, \mathbf{s}^k)}{\|\mathbf{x}^{k+1}\|_\infty} \geq \sum_{i \in \iota} |s_i^k| \geq r^k, \quad (3.65)$$

where the subgradient \mathbf{s}^k is not required to be ordered like Eq. (3.15). In particular, such monotone increasing could be strict, i.e. $r^{k+1} > r^k$ via improving the last “ \geq ” in Eq. (3.65) into “ $>$ ”, if $\|\mathbf{s}^k\|_1 > r^k$ holds in each step. To this end, we only need to determine a subgradient $\mathbf{s}^\sigma \in \partial I(\mathbf{x}^k)$ such that $\|\mathbf{s}^\sigma\|_1 > r^k$ if there exists $\mathbf{s} \in \partial I(\mathbf{x}^k)$ such that $\|\mathbf{s}\|_1 > r^k$.

Suppose $\mathbf{x}^k = (x_1^k, \dots, x_n^k)$, $\mathbf{s} = (s_1, \dots, s_n)$, and then we have

$$\lambda_i(t) = |t| - \frac{x_i^k}{\|\mathbf{x}^k\|_\infty} t \geq 0, \quad \forall i \in \{1, \dots, n\}, \quad (3.66)$$

$$\|\mathbf{s}\|_1 - r^k = \|\mathbf{s}\|_1 - \frac{(\mathbf{x}^k, \mathbf{s})}{\|\mathbf{x}^k\|_\infty} = \sum_{i=1}^n \lambda_i(s_i), \quad (3.67)$$

which yields

$$\|\mathbf{s}\|_1 - r^k > 0 \Leftrightarrow \exists i \in \{1, \dots, n\}, \quad \text{s.t.} \quad \lambda_i(s_i) > 0. \quad (3.68)$$

Given $\mathbf{x} \in \mathbb{R}^n \setminus \{\mathbf{0}\}$, let

$$\mathbf{q} = (q_1, \dots, q_n) \in (\mathbb{R}^+)^n, \quad q_i = \sum_{j:\{i,j\} \in E} w_{ij} \mathbf{1}_{x_i = x_j}, \quad (3.69)$$

$$\mathbf{p} = (p_1, \dots, p_n), \quad p_i = \sum_{j:\{i,j\} \in E} w_{ij} \text{sign}(x_i - x_j) - q_i. \quad (3.70)$$

It can be readily verified that $\mathbf{p} \in \partial I(\mathbf{x})$, and

$$s_i \in (\partial I(\mathbf{x}))_i = [p_i - q_i, p_i + q_i], \quad \forall \mathbf{s} = (s_1, \dots, s_n) \in \partial I(\mathbf{x}). \quad (3.71)$$

Here $(\partial I(\mathbf{x}))_i$ denotes the projected interval of the convex domain $\partial I(\mathbf{x})$ onto the i -th coordinate.

Denote $\bar{\mathbf{p}} = (\bar{p}_1, \dots, \bar{p}_n)$ by

$$\bar{p}_i = \begin{cases} p_i \mp q_i, & \text{if } i \in S^\pm(\mathbf{x}), \\ p_i + \text{sign}(p_i)q_i, & \text{if } i \in S^<(\mathbf{x}), \end{cases} \quad (3.72)$$

which is located on the boundary of $(\partial I(\mathbf{x}))_i = [p_i - q_i, p_i + q_i]$.

Accordingly, combining Eqs. (3.68), (3.71) and the convexity of $\lambda_i(t)$ given in Eq. (3.66) leads to

$$\begin{aligned} & \exists \mathbf{s} \in \partial I(\mathbf{x}^k), \quad \text{s.t.} \quad \|\mathbf{s}\|_1 > r^k \\ & \Leftrightarrow \exists i \in \{1, \dots, n\}, \quad \text{s.t.} \quad \max_{s_i \in (\partial I(\mathbf{x}))_i} \lambda_i(s_i) > 0 \\ & \Leftrightarrow \exists i \in \{1, \dots, n\}, \quad \text{s.t.} \quad \lambda_i(\bar{p}_i) > 0. \end{aligned}$$

That is, if there exists $i \in \{1, \dots, n\}$ such that $\lambda_i(\bar{p}_i) > 0$, then there should exist a subgradient $\mathbf{s} \in \partial I(\mathbf{x}^k)$ satisfying $s_i = \bar{p}_i$ and $\|\mathbf{s}\|_1 > r^k$. Hence, if $\bar{\mathbf{p}} \in \partial I(\mathbf{x})$ (but hardly holds in general), then we may directly select $\bar{\mathbf{p}}$; otherwise we are able to use $\bar{\mathbf{p}}$ as an indicator for the subgradient selection.

Define a partial order relation “ \leq ” on \mathbb{R}^2 by $(x_1, y_1) \leq (x_2, y_2)$ if and only if either $x_1 < x_2$ or $x_1 = x_2, y_1 \leq y_2$ holds. Let $\Sigma(\mathbf{x})$ be a collection of permutations of $\{1, 2, \dots, n\}$ such that for any $\sigma \in \Sigma(\mathbf{x})$, it holds

$$(x_{\sigma(1)}, \bar{p}_{\sigma(1)}) \leq (x_{\sigma(2)}, \bar{p}_{\sigma(2)}) \leq \dots \leq (x_{\sigma(n)}, \bar{p}_{\sigma(n)}). \quad (3.73)$$

For any $\sigma \in \Sigma(\mathbf{x})$, we select

$$\mathbf{s}^\sigma = (s_1^\sigma, s_2^\sigma, \dots, s_n^\sigma) \in \partial I(\mathbf{x}), \quad (3.74)$$

$$s_i^\sigma = \sum_{j:\{i,j\} \in E} w_{ij} z_{ij}, \quad i = 1, 2, \dots, n, \quad (3.75)$$

$$z_{ij} = \text{sign}(\sigma^{-1}(i) - \sigma^{-1}(j)) \in \text{Sgn}(x_i - x_j), \quad i, j = 1, 2, \dots, n. \quad (3.76)$$

The following theorem demonstrates that such subgradient selection is sufficient to guarantee the strict monotonicity $r^{k+1} > r^k$ (if any) with the help of the statement (1) of Lemma 3.1.

Theorem 3.3. *For any permutation $\sigma \in \Sigma(\mathbf{x}^k)$, we have $\|\mathbf{s}^\sigma\| > r^k$ if and only if there exists $\mathbf{s} \in \partial I(\mathbf{x}^k)$ such that $\|\mathbf{s}\| > r^k$.*

Proof. The necessity is obvious, we only need to prove the sufficiency. Suppose $\mathbf{s} = (s_1, \dots, s_n) \in \partial I(\mathbf{x}^k)$ satisfies $\|\mathbf{s}\| > r^k$. Then it can be derived that $\mathbf{x}^k / \|\mathbf{x}^k\|_\infty \notin \text{Sgn}(\mathbf{s})$ using the statement (1) of Lemma 3.1, i.e. there exists an $i_0 \in \{1, 2, \dots, n\}$ such that

$$\frac{x_{i_0}^k}{\|\mathbf{x}^k\|_\infty} \notin \text{Sgn}(s_{i_0}). \quad (3.77)$$

Denote $J = \{j \mid x_j^k = x_{i_0}^k\}$. Let $j_1, j_2 \in J$ be the indexes minimizing and maximizing function $\sigma^{-1}(\cdot)$ over J , respectively. Then we have

$$\begin{aligned} s_{j_1}^\sigma &= \sum_{t: \{t, j_1\} \in E} w_{j_1 t} \text{sign}(\sigma^{-1}(j_1) - \sigma^{-1}(t)) = p_{j_1} - q_{j_1}, \\ s_{j_2}^\sigma &= \sum_{t: \{t, j_2\} \in E} w_{j_2 t} \text{sign}(\sigma^{-1}(j_2) - \sigma^{-1}(t)) = p_{j_2} + q_{j_2}. \end{aligned}$$

We claim that either j_1 or j_2 is an integer j such that $x_j^k \notin \|\mathbf{x}^k\|_\infty \text{Sgn}(s_j^\sigma)$. This directly yields $\|\mathbf{s}^\sigma\| > r^k$ according to the statement (1) of Lemma 3.1 and thus completes the proof. Suppose the contrary

$$x_{i_0}^k = x_j^k \in \|\mathbf{x}^k\|_\infty \text{Sgn}(s_j^\sigma), \quad j = j_1, j_2, \quad (3.78)$$

and split the discussion into the following two cases:

(1) $i_0 \in S^<(\mathbf{x}^k)$. Eq. (3.78) implies that

$$s_{j_1}^\sigma = s_{j_2}^\sigma = 0 \Leftrightarrow p_{j_1} = q_{j_1} \geq 0, \quad p_{j_2} = -q_{j_2} \leq 0,$$

and thus, from Eq. (3.72), we have

$$\begin{aligned} \bar{p}_{j_1} &= p_{j_1} + \text{sign}(p_{j_1})q_{j_1} = p_{j_1} + \text{sign}(p_{j_1})p_{j_1} \geq 0, \\ \bar{p}_{j_2} &= p_{j_2} + \text{sign}(p_{j_2})q_{j_2} = p_{j_2} - \text{sign}(p_{j_2})p_{j_2} \leq 0. \end{aligned}$$

Since

$$(x_{j_1}, \bar{p}_{j_1}) \leq (x_{i_0}, \bar{p}_{i_0}) \leq (x_{j_2}, \bar{p}_{j_2}), \quad x_{j_1} = x_{i_0} = x_{j_2},$$

it yields

$$0 \geq \bar{p}_{j_2} \geq \bar{p}_{i_0} \geq \bar{p}_{j_1} \geq 0 \Rightarrow \bar{p}_{i_0} = 0 \Rightarrow p_{i_0} = q_{i_0} = 0 \Rightarrow s_{i_0} = 0,$$

which contradicts Eq. (3.77).

(2) $i_0 \in S^\pm(\mathbf{x}^k)$. Eq. (3.78) implies that

$$\begin{cases} s_{i_0} \geq p_{i_0} - q_{i_0} = \bar{p}_{i_0} \geq \bar{p}_{j_1} = s_{j_1}^\sigma \geq 0, & \text{if } i_0 \in S^+(\mathbf{x}^k), \\ s_{i_0} \leq p_{i_0} + q_{i_0} = \bar{p}_{i_0} \leq \bar{p}_{j_2} = s_{j_2}^\sigma \leq 0, & \text{if } i_0 \in S^-(\mathbf{x}^k), \end{cases} \quad (3.79)$$

both of which contradict Eq. (3.77). \square

In a word, we choose $\mathbf{s}^k = \mathbf{s}^\sigma \in \partial I(\mathbf{x}^k)$ with $\sigma \in \Sigma(\mathbf{x}^k)$ in the simple iterative scheme (3.9). Besides the above-mentioned strict increasing, we are able to show below that such subgradient selection assures finite-step local convergence. Before that, we would like to further specify the choice of \mathbf{x}^{k+1} from the closed solution set X_p^{k+1} at the first step. There is a natural isomorphism h by a central projection between S_p and S_∞ which are the unit spheres in norms p and ∞ , respectively. Denote $\partial X_p^{k+1} \in X_p^{k+1}$ be the collection of points corresponding to the vertices of $h(X_p^{k+1})$ which is a convex polytope, in view of the fact that the convex function $I(\mathbf{x})$ achieves its maximum values on vertices among $h(X_p^{k+1})$. Hence, the three-step iterative scheme (3.9) can be crystallized into

$$\begin{cases} \mathbf{x}^{k+1} \in \partial X_p^{k+1}, & (3.80a) \\ r^{k+1} = F(\mathbf{x}^{k+1}), & (3.80b) \\ \mathbf{s}^{k+1} = \mathbf{s}^\sigma, \quad \sigma \in \Sigma(\mathbf{x}^{k+1}). & (3.80c) \end{cases}$$

Let

$$C = \left\{ \mathbf{x} \in \mathbb{R}^n \mid F(\mathbf{y}) \leq F(\mathbf{x}), \forall \mathbf{y} \in \{T_i \mathbf{x} : i \in \{1, \dots, n\}\} \right\}, \quad (3.81)$$

where $T_i \mathbf{x}$ is defined as

$$(T_i \mathbf{x})_j = \begin{cases} x_j, & j \neq i, \\ -x_j, & j = i. \end{cases} \quad (3.82)$$

Theorem 3.4 (Finite-Step Local Convergence). *Assume the sequences $\{\mathbf{x}^k\}$ and $\{r^k\}$ are generated by the simple iterative scheme (3.80) from any initial point $\mathbf{x}^0 \in \mathbb{R}^n \setminus \{\mathbf{0}\}$. There must exist $N \in \mathbb{Z}^+$ and $r^* \in \mathbb{R}$ such that, for any $k > N$, $r^k = r^*$, and $\mathbf{x}^{k+1} \in C$ are local maximizers.*

Proof. According to Eq. (3.65), for any integer $k > 0$, there exists $\iota \subset \{1, \dots, n\}$ such that $r^k \leq \sum_{i \in \iota} |s_i^k| \leq r^{k+1}$, which means the sequence $\{r^k\}$ can only take finite values because the set

$$\left\{ \sum_{i \in \iota} |s_i| \mid \mathbf{s}^\sigma = (s_1, \dots, s_n), \forall \sigma \in \Sigma(\mathbf{x}), \forall \mathbf{x}, \iota \right\}$$

is finite. Thus there exist $N \in \mathbb{Z}^+$ and $r^* \in \mathbb{R}$ such that $r^k = r^*$ for any $k > N$, thereby implying that

$$\begin{aligned} r^k &= \|\mathbf{s}^k\|_1, \\ X_p^{k+1} &= \left\{ \mathbf{x} \mid \frac{\mathbf{x}}{\|\mathbf{x}\|_\infty} \in \text{Sgn}(\mathbf{s}^k), \|\mathbf{x}\|_p = 1 \right\} \end{aligned}$$

by virtue of Lemma 3.1. That is, $\forall \mathbf{x}^{k+1} \in \partial X_p^{k+1}$, we have

$$\frac{\mathbf{x}^{k+1}}{\|\mathbf{x}^{k+1}\|_\infty} \in \text{Sgn}(\mathbf{s}), \quad \forall \mathbf{s} \in \partial I(\mathbf{x}^{k+1}), \quad (3.83)$$

$$S^<(\mathbf{x}^{k+1}) = \emptyset. \quad (3.84)$$

Now we will show $\mathbf{x}^{k+1} \in C$ and neglect the superscript $k+1$ hereafter for simplicity. Suppose the contrary that there exists $i \in S^\pm(\mathbf{x})$ satisfying $F(T_i \mathbf{x}) > F(\mathbf{x})$, and then we have

$$\|T_i \mathbf{x}\|_\infty = \|\mathbf{x}\|_\infty, \quad I(T_i \mathbf{x}) - I(\mathbf{x}) = \pm \sum_{j: \{j, i\} \in E} w_{ij} 2x_j > 0. \quad (3.85)$$

Let $\mathbf{s} = (s_1, \dots, s_n) \in \partial I(\mathbf{x})$ be generated by

$$z_{ij} = -\frac{x_j}{\|\mathbf{x}\|_\infty} \in \text{Sgn}(x_i - x_j),$$

and thus

$$x_i s_i = x_i \sum_{j:\{j,i\} \in E} w_{ij} z_{ij} = - \sum_{j:\{j,i\} \in E} \frac{w_{ij} x_i x_j}{\|\mathbf{x}\|_\infty} = -\frac{1}{2}(I(T_i \mathbf{x}) - I(\mathbf{x})) < 0,$$

which contradicts Eq. (3.83).

Finally, we want to show \mathbf{x} is a local maximizer of $F(\cdot)$ on $\mathbb{R}^n \setminus \{\mathbf{0}\}$. Let U be a neighborhood of \mathbf{x} such that

$$\begin{aligned} \left\| \frac{\mathbf{y}}{\|\mathbf{y}\|_\infty} - \frac{\mathbf{x}}{\|\mathbf{x}\|_\infty} \right\|_\infty &< \frac{1}{2}, & \forall \mathbf{y} \in U, \\ \mathbf{y}' = \frac{\|\mathbf{x}\|_\infty}{\|\mathbf{y}\|_\infty} \mathbf{y}, \quad g(t) &= I(t(\mathbf{y}' - \mathbf{x}) + \mathbf{x}), & \forall \mathbf{y} \in U. \end{aligned}$$

We claim

$$g(t) - g(0) \leq 0, \quad \forall t \in [0, 1], \quad (3.86)$$

with which we are able to verify \mathbf{x} is a local maximizer as follows:

$$F(\mathbf{y}) - F(\mathbf{x}) = F(\mathbf{y}') - F(\mathbf{x}) = \frac{1}{\|\mathbf{x}\|_\infty} (I(\mathbf{y}') - I(\mathbf{x})) = \frac{1}{\|\mathbf{x}\|_\infty} (g(1) - g(0)) \leq 0, \quad \forall \mathbf{y} \in U.$$

The only remaining thing is to prove Eq. (3.86). First, it can be easily shown that $g(t)$ is linear on $[0, 1]$, and there exists $\mathbf{s} = (s_1, \dots, s_n) \in \partial I(\mathbf{x})$ such that its slope can be determined by

$$g(t) - g(0) = t(\mathbf{s}, \mathbf{y}' - \mathbf{x}).$$

Then, the verification of Eq. (3.86) can be completed by

$$x_j (y'_j - x_j) \leq 0 \Rightarrow s_j (y'_j - x_j) \leq 0, \quad j = 1, 2, \dots, n,$$

where we have used $\|\mathbf{y}'\|_\infty = \|\mathbf{x}\|_\infty$ as well as Eqs. (3.83) and (3.84). \square

Remark 3.1. Notice that $\{T_i\}_{i=1}^n$ can generate a commutative group \mathcal{T} on \mathbb{R}^n . If we further restrict its domain to be $\{\mathbf{x} \mid S^<(\mathbf{x}) = \emptyset\}$, then \mathbf{x} is a global maximizer of $F(\cdot)$ on $\mathbb{R}^n \setminus \{\mathbf{0}\}$ if and only if $F(\mathbf{y}) \leq F(\mathbf{x})$ holds for any $\mathbf{y} = T\mathbf{x}$, $\forall T \in \mathcal{T}$. In such sense, the set C given in (3.81) is the first order approximation to global maximizers.

4. Numerical Experiments

In this section, we conduct performance evaluation of the proposed SI algorithm (3.80) on the graphs with positive weight in G-set (<https://web.stanford.edu/yye/yye/Gset/>), and always set the initial data \mathbf{x}^0 to be the maximal eigenvector of the graph Laplacian [8, 16]. The three bipartite graphs $G48$, $G49$, $G50$ will not be considered because their optima cuts can be achieved at the initial step. The recently updated cut values achieved by a multiple search operator heuristic are chosen to be the reference [13]. SI is capable of producing approximate cuts with high quality: the ratios between the resulting cut values and the reference ones are at least 0.986 (see Table 4.2), and can be improved to 0.997 (see Table 4.3) after introducing a straightforward perturbation to break out of local optima.

4.1. Implementation and cost analysis

We begin with the algorithm implementation plus a preliminary cost analysis. Algorithm 4.1 gives the pseudo-code of the SI algorithm (3.80), where \mathbf{x}^{k+1} , r^{k+1} and \mathbf{s}^{k+1} are generated in Line 4, Line 5 and Lines 6-9, respectively. In Line 4, the function EXACT_SOLUTION randomly selects the iteration point $\mathbf{x}^{k+1} \in \partial X_p^{k+1}$ where the exact solution set X_p^{k+1} is given in Theorem 3.1. After $\bar{\mathbf{p}}$ is obtained in Line 8 using Eq. (3.72), we are able to arrive at the partial order (3.73) through the Bubble Sort procedure during which the requested subgradient \mathbf{s}^{k+1} can be automatically updated in an iterative manner (see Lines 17 and 18). This constitutes the subroutine named by SUBGRADIENT which starts from Line 12. It should be pointed out that there is randomness in determining both \mathbf{x}^{k+1} and \mathbf{s}^{k+1} , so that the performance evaluation below is conducted in the sense of average by re-running SI (3.80) 100 times from the same initial data. A preliminary estimate of the cost for three iteration steps $T = 50, 500, 2000$ is presented in Table 4.1 where we have set $p = 2$ for instance.

The main task for reaching the exact solution set X_p^{k+1} by Theorem 3.1 is to determine m_0 with which m_1 can be automatically obtained by Eq. (3.55). In our implementation, EXACT_SOLUTION uses the Bubble Sort to arrange \mathbf{s}^k in an ascending order such that the accumulation of increment $A(n), A(n-1), \dots, A(1)$ is calculated, successively, until $m_0(k)$

Algorithm 4.1: Pseudo-Code of the Simple Iterative (SI) Algorithm.

```

1 Initialize  $\mathbf{x}^0, r^0, \mathbf{q}^0, \mathbf{p}^0, \sigma^0, \mathbf{s}^0, T$ .
2  $k \leftarrow 0$ .
3 while  $k < T$  do
4    $\mathbf{x}^{k+1} \leftarrow \text{EXACT\_SOLUTION}(\mathbf{s}^k, r^k)$ ;
5    $r^{k+1} \leftarrow r^k + \delta_F(\mathbf{x}^{k+1}, \mathbf{x}^k)$ ;
6    $\mathbf{q}^{k+1} \leftarrow \mathbf{q}^k + \delta_{\mathbf{q}}(\mathbf{x}^{k+1}, \mathbf{x}^k)$ ;
7    $\mathbf{p}^{k+1} \leftarrow \mathbf{p}^k + \delta_{\mathbf{p}}(\mathbf{x}^{k+1}, \mathbf{x}^k)$ ;
8    $\bar{\mathbf{p}}^{k+1} \leftarrow \text{P\_BAR}(\mathbf{x}^{k+1}, \mathbf{q}^{k+1}, \mathbf{p}^{k+1})$ ;
9    $(\mathbf{s}^{k+1}, \sigma^{k+1}) \leftarrow \text{SUBGRADIENT}(\mathbf{x}^{k+1}, \bar{\mathbf{p}}^{k+1}, \mathbf{s}^k, \sigma^k)$ ;
10   $k \leftarrow k + 1$ .
11 end
12 function  $(\mathbf{s}, \sigma) = \text{SUBGRADIENT}(\mathbf{x}, \bar{\mathbf{p}}, \mathbf{s}, \sigma)$ 
13 for  $i = 1$  to  $n - 1$  do
14   for  $j = i + 1$  to  $n$  do
15     if  $(\mathbf{x}(\sigma(j)), \bar{\mathbf{p}}(\sigma(j))) \leq (\mathbf{x}(\sigma(j-1)), \bar{\mathbf{p}}(\sigma(j-1)))$  then
16        $\sigma \leftarrow \text{SWAP}(\sigma, j-1, j)$ ;
17        $\mathbf{s}(\sigma(j-1)) \leftarrow \mathbf{s}(\sigma(j-1)) - 2w_{\sigma(j)\sigma(j-1)}$ ;
18        $\mathbf{s}(\sigma(j)) \leftarrow \mathbf{s}(\sigma(j)) + 2w_{\sigma(j)\sigma(j-1)}$ .
19     end
20   else
21     break.
22   end
23 end
24 end function
25 end function

```

Table 4.1: Cost analysis: Mean values of $(n - m_0^k)/n$, $\delta_\sigma(k)/n$, and $|V(k)|/n$ obtained from 100 runs of the simple iterative (SI) algorithm (3.80) with $p = 2$. Here $(n - m_0^k)$, $\delta_\sigma(k)$, and $|V(k)|$ are given in Eq. (3.51), Eq. (4.1), and Eq. (4.5), respectively, n denotes the size of graph and T gives the total step of iterations. The time complexity of SI is $\mathcal{O}(\text{mean}(c(k))nT)$ with $c(k) = (n - m_0(k)) + \delta_\sigma(k) + |V(k)|$ depending on the underlying graph. The values of $\text{mean}(c(k))/n$ are at most 0.095, 0.077, 0.077 for $T = 50, 500, 2000$, respectively, thereby indicating $\text{mean}(c(k))$ is much smaller than n .

Graph	n	$\text{mean}(n - m_0^k)/n$			$\text{mean}(\delta_\sigma(k))/n$			$\text{mean}(V(k))/n$		
		$T = 50$	$T = 500$	$T = 2000$	$T = 50$	$T = 500$	$T = 2000$	$T = 50$	$T = 500$	$T = 2000$
G1	800	0.0073	0.0046	0.0045	0.0091	0.0082	0.0083	0.0010	0.0000	0.0000
G2	800	0.0069	0.0043	0.0042	0.0090	0.0084	0.0082	0.0008	0.0000	0.0000
G3	800	0.0072	0.0048	0.0046	0.0093	0.0081	0.0081	0.0009	0.0000	0.0000
G4	800	0.0061	0.0042	0.0040	0.0090	0.0082	0.0081	0.0061	0.0042	0.0040
G5	800	0.0059	0.0045	0.0043	0.0089	0.0081	0.0079	0.0007	0.0000	0.0000
G14	800	0.0510	0.0423	0.0412	0.0246	0.0243	0.0251	0.0034	0.0000	0.0000
G15	800	0.0550	0.0444	0.0440	0.0248	0.0253	0.0247	0.0068	0.0000	0.0000
G16	800	0.0532	0.0445	0.0428	0.0229	0.0237	0.0237	0.0532	0.0445	0.0428
G17	800	0.0567	0.0471	0.0461	0.0244	0.0244	0.0244	0.0020	0.0000	0.0000
G22	2000	0.0146	0.0106	0.0102	0.0120	0.0125	0.0120	0.0018	0.0001	0.0001
G23	2000	0.0119	0.0100	0.0099	0.0123	0.0117	0.0118	0.0021	0.0000	0.0000
G24	2000	0.0135	0.0100	0.0097	0.0125	0.0117	0.0120	0.0135	0.0100	0.0097
G25	2000	0.0161	0.0115	0.0111	0.0132	0.0124	0.0122	0.0161	0.0115	0.0111
G26	2000	0.0140	0.0101	0.0097	0.0126	0.0127	0.0127	0.0038	0.0000	0.0000
G35	2000	0.0475	0.0380	0.0370	0.0274	0.0259	0.0255	0.0475	0.0380	0.0370
G36	2000	0.0566	0.0484	0.0466	0.0252	0.0286	0.0278	0.0054	0.0005	0.0000
G37	2000	0.0516	0.0427	0.0414	0.0260	0.0272	0.0271	0.0076	0.0000	0.0000
G38	2000	0.0595	0.0478	0.0466	0.0260	0.0292	0.0302	0.0091	0.0000	0.0000
G43	1000	0.0140	0.0102	0.0100	0.0131	0.0125	0.0122	0.0024	0.0000	0.0000
G44	1000	0.0169	0.0134	0.0133	0.0126	0.0135	0.0129	0.0013	0.0000	0.0000
G45	1000	0.0152	0.0137	0.0135	0.0126	0.0130	0.0132	0.0152	0.0137	0.0135
G46	1000	0.0112	0.0093	0.0091	0.0121	0.0121	0.0121	0.0011	0.0000	0.0000
G47	1000	0.0125	0.0103	0.0100	0.0125	0.0125	0.0122	0.0125	0.0103	0.0100
G51	1000	0.0527	0.0440	0.0427	0.0228	0.0241	0.0239	0.0016	0.0004	0.0000
G52	1000	0.0485	0.0406	0.0391	0.0203	0.0239	0.0235	0.0485	0.0406	0.0391
G53	1000	0.0544	0.0462	0.0453	0.0212	0.0242	0.0245	0.0031	0.0000	0.0000
G54	1000	0.0616	0.0544	0.0520	0.0215	0.0241	0.0235	0.0017	0.0000	0.0000

(i.e. m_0 at the k -step iteration) can be determined via Eq. (3.51). The cost of this procedure is $\mathcal{O}((n - m_0(k))n)$. Table 4.1 shows that the mean values of $(n - m_0(k))/n$ are far less than 1, which are at most 0.060, 0.048, 0.047 for $T = 50, 500, 2000$, respectively. In particular, EXACT_SOLUTION reduces to Scenario 3 and thus only costs $\mathcal{O}(n)$ for $p = \infty$.

The subroutine SUBGRADIENT is also a Bubble Sort procedure and costs $\mathcal{O}(n^2)$ in worst cases. However, it should be pointed out that the efficiency of Bubble Sort depends on the

initial order and actually costs $\mathcal{O}((\delta_\sigma(k) + 1)n)$, where

$$\delta_\sigma(k) = \frac{1}{2n} \sum_{i=1}^n |\sigma^{k+1}(i) - \sigma^k(i)| \quad (4.1)$$

denotes an average displacement of the permutation σ used in Eq. (3.73) for the k -th iteration. Table 4.1 shows that the ratios between $\text{mean}(\delta_\sigma(k))$ and n are at most 0.03.

It remains to estimate the cost for updating $r, \mathbf{q}, \mathbf{p}$ in Lines 5-7, where we prefer to only calculate the increment

$$\delta_\Gamma(\mathbf{x}^{k+1}, \mathbf{x}^k) := \Gamma(\mathbf{x}^{k+1}) - \Gamma(\mathbf{x}^k), \quad \Gamma \in \{F, \mathbf{q}, \mathbf{p}\}. \quad (4.2)$$

Let

$$\mathbf{z}^k = \frac{\mathbf{x}^k}{\|\mathbf{x}^k\|_\infty}, \quad \mathbf{z}^{k,i} = (z_1^{k,i}, \dots, z_n^{k,i}) = (z_1^{k+1}, \dots, z_i^{k+1}, z_{i+1}^k, \dots, z_n^k). \quad (4.3)$$

Then we have

$$\begin{aligned} \delta_\Gamma(\mathbf{x}^{k+1}, \mathbf{x}^k) &= \delta_\Gamma(\mathbf{z}^{k+1}, \mathbf{z}^k) = \sum_{i=1}^n \delta_\Gamma(\mathbf{z}^{k,i}, \mathbf{z}^{k,i-1}) \\ &= \sum_{i \in V(k)} \delta_\Gamma(\mathbf{z}^{k,i}, \mathbf{z}^{k,i-1}), \quad \Gamma \in \{F, \mathbf{q}, \mathbf{p}\}, \end{aligned} \quad (4.4)$$

where

$$V(k) = \{i \mid z_i^{k+1} \neq z_i^k\}. \quad (4.5)$$

Consequently, the complexity of calculating $\delta_\Gamma(\mathbf{x}^{k+1}, \mathbf{x}^k)$ is $\mathcal{O}(|V(k)|n)$ since we need $\mathcal{O}(n)$ to produce each component $\delta_\Gamma(\mathbf{z}^{k,i}, \mathbf{z}^{k,i-1})$. Table 4.1 reveals that $\text{mean}(|V(k)|)$ is much smaller than n and sometimes vanishes as T increases.

In total, the SI algorithm costs $\mathcal{O}(c(k)n)$ in the k -th step and

$$c(k) := (n - m_0(k)) + \delta_\sigma(k) + |V(k)|$$

depends on the underlying graph like its weights and order. Actual numerical experiments in Table 4.1 show that the mean value of $c(k)$ is much smaller than n .

4.2. Quality check

We are now ready for quality check of numerical solutions achieved by the simple algorithm (3.80). The numerical results for the RSC algorithm based on graph Laplacian (Δ_2 -RSC) [14] and graph 1-Laplacian (Δ_1 -RSC) [5], as well as the GW algorithm [14] are adopted for comparison.

The quality check is performed based on numerical solutions until $T = 2000$. Table 4.2 shows the minimum, mean and maximum cut values during 100 runs for $p = 1, 2, \infty$. It can be easily seen there that the results for different $p (= 1, 2, \infty)$ are comparable and are all very close to the reference values. Actually, the ratios between the best cut values by SI (chosen from the maximum cut values over $p = 1, 2, \infty$) and the reference ones are at least 0.986 (see the results for G36), while the numerical lower bound for such ratios is about 0.946, 0.933 and 0.949 for the GW, Δ_2 -RSC and Δ_1 -RSC algorithms, respectively. In particular, for the case of $p = \infty$, the ratios between the minimum, mean, maximum cut values and the reference ones are at

Table 4.2: Quality check: Cut values for 27 problem instances in G-set obtained from 100 runs of the simple iterative (SI) algorithm (3.80) with $p = 1, 2, \infty$. Each run starts from the maximal eigenvector of the graph Laplacian and undergoes $T = 2000$ iterations. The minimum, mean and maximum cut values are recorded and compared to the reference ones obtained by combinatorial algorithms [13], while the initial cut values are listed in the second column. The ratios between the best cut values (in italics) by SI and the reference values are at least 0.986 (see G36).

Graph	Initial	$p = 1$			$p = 2$			$p = \infty$			Reference
		min	mean	max	min	mean	max	min	mean	max	
G1	11221	11477	11524.52	<i>11554</i>	11505	11527.27	11548	11499	11523.3	11553	11624
G2	11283	11483	11525.91	<i>11583</i>	11484	11505.6	11535	11473	11517.42	11557	11620
G3	11298	11497	11542.53	11599	11564	11586.14	11602	11557	11586.99	<i>11605</i>	11622
G4	11278	11520	11561.1	11597	11545	11564.49	<i>11606</i>	11539	11568.26	11598	11646
G5	11370	11530	11568.89	11602	11561	11579.89	<i>11607</i>	11558	11577.22	11602	11631
G14	2889	2998	3016.89	<i>3036</i>	3015	3022.74	3030	3012	3026.18	3033	3064
G15	2771	2968	2986.36	3005	2989	2995.56	3006	2985	2996.68	<i>3008</i>	3050
G16	2841	2973	2995.32	3009	3005	3010.14	3014	3005	3011.76	<i>3017</i>	3052
G17	2866	2963	2981.65	3004	2996	3002.27	<i>3012</i>	2995	3002.7	3010	3047
G22	12876	13198	13243.55	<i>13302</i>	13237	13267.7	13286	13248	13266.04	13290	13359
G23	12817	13173	13220.99	13267	13238	13252.85	<i>13271</i>	13232	13246.72	13264	13344
G24	12826	13194	13221.16	13244	13235	13260.34	<i>13295</i>	13233	13262.4	13287	13337
G25	12781	13155	13189.61	<i>13245</i>	13185	13221	13240	13187	13213.86	13238	13340
G26	12752	13140	13176.81	<i>13222</i>	13185	13201.6	13221	13183	13197.21	13210	13328
G35	7194	7512	7540.39	7572	7558	7576.68	<i>7595</i>	7560	7575.4	7586	7687
G36	7124	7502	7529.77	7557	7535	7548.22	7566	7531	7553.36	<i>7576</i>	7680
G37	7162	7505	7541.03	7563	7565	7581.08	<i>7603</i>	7575	7590.59	7602	7691
G38	7122	7513	7542.91	7568	7561	7573.63	7589	7544	7565.93	<i>7593</i>	7688
G43	6395	6570	6609.44	6635	6599	6625.47	<i>6645</i>	6604	6625.8	6644	6660
G44	6439	6575	6598.42	6618	6593	6608.63	6617	6590	6609.27	<i>6621</i>	6650
G45	6364	6574	6599.49	<i>6624</i>	6587	6595.39	6612	6586	6593.07	6599	6654
G46	6389	6557	6586.68	<i>6612</i>	6562	6583.31	6597	6569	6583.93	6602	6649
G47	6353	6552	6583.5	<i>6614</i>	6584	6598.8	6609	6584	6597.24	6610	6657
G51	3645	3769	3787.88	3810	3788	3797.76	<i>3811</i>	3785	3794.12	3806	3848
G52	3645	3766	3787.31	3807	3808	3811.55	3816	3806	3811.85	<i>3820</i>	3851
G53	3630	3778	3793.98	3812	3795	3804.37	<i>3818</i>	3797	3808.45	3817	3850
G54	3655	3767	3789.34	3805	3800	3809.05	3818	3808	3813.69	<i>3821</i>	3852

least 0.979, 0.982, 0.986, respectively, all of which are larger than the average ratios over these 27 graphs for the Δ_1 -RSC ($\simeq 0.971$), GW ($\simeq 0.960$), Δ_2 -RSC ($\simeq 0.958$), and SC ($\simeq 0.951$) algorithms. The SC cuts are obtained by rounding the initial data with a threshold of 0 and the cut values are shown in the second column of Table 4.2. In order to further show the overall performance in achieving high ratio by the SI algorithm, we plot the histogram of the ratios for all 3×2700 runs in Fig. 4.1. We are able to clearly observe there that:

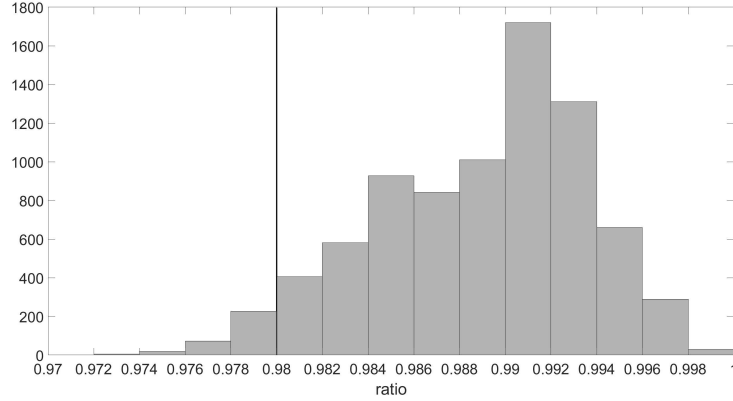


Fig. 4.1. Quality check: Histogram of the ratios for all 2700×3 runs depicted in Table 4.2 (more explanations are referred to Table 4.2). The percent of runs obtained a ratio larger than 0.980, which lie on the right of the black vertical line, exceeds 95%.

- (1) More than 95% of runs achieve ratios exceeding 0.986 (see the black vertical line of Fig. 4.1).
- (2) The percent of runs obtained a ratio larger than 0.986 exceeds 72%.

4.3. Breaking out of local optima

Within the SI algorithm, we are allowed to plug into local breakout techniques to further improve the solution quality. A preliminary attempt is to generate a new point $\tilde{\mathbf{x}}^{k+1} = (\tilde{x}_1, \dots, \tilde{x}_n)$ in a stochastic manner

$$\tilde{x}_i = \begin{cases} -x_i^k & \text{with the probability of } e^{-\beta|\bar{p}_i^k|}, \\ x_i^k & \text{with the probability of } 1 - e^{-\beta|\bar{p}_i^k|}, \end{cases} \quad (4.6)$$

when SI is stuck at $\mathbf{x}^k = (x_1^k, \dots, x_n^k)$, namely, it cannot make the function $F(\cdot)$ increase. Here we choose β to be a controllable parameter, and $\bar{\mathbf{p}}^k = (\bar{p}_1^k, \dots, \bar{p}_n^k)$ to generate $\tilde{\mathbf{x}}^{k+1}$ in order to decrease $F(\cdot)$ (if any) as little as possible in view of the following fact:

$$F(T_i \mathbf{x}^k) - F(\mathbf{x}^k) = -|\bar{p}_i^k|, \quad i = 1, \dots, n, \quad (4.7)$$

where T_i is defined in Eq. (3.82). In such sense, we regard the manner (4.6) as a special kind of perturbation, and the resulting algorithm is named by the simple iteration with perturbation (SI-P).

Algorithm 4.2 presents the pseudo-code of SI-P. In Lines 17-39, SI_PERTURB undergoes the simple iteration equipped with jumping out of local optima until a given final time T during which the perturbation (4.6) triggers with a prescribed β if the function values remain unchanged for t steps (see Lines 29-32). It should be noted that SI_PERTURB returns the maximal cut value r_{opt} and the corresponding point \mathbf{x}_{opt} during the period of T . We are now left only to choose β which should not only depend on the iteration time but also have a specific range because $\tilde{\mathbf{x}}^{k+1} = -\mathbf{x}^k$ when $\beta = 0$ and $\tilde{\mathbf{x}}^{k+1} = \mathbf{x}^k$ when $\beta \rightarrow \infty$. Here Algorithm 4.2 adopts

Algorithm 4.2: Pseudo-Code of the Simple Iteration with Perturbation (SI-P).

```

Require: initial  $\mathbf{x}^0, L, T$ .
Ensure :  $count, \mathbf{x}^{count}, r^{count}$ .
1  $r^0 \leftarrow 0, count \leftarrow 0$ .
2 while True do
3   for  $ipert = 1$  to  $L$  do
4      $\beta$ : chosen from  $(0, 1)$  randomly;
5      $(\tilde{\mathbf{x}}^{ipert}, \tilde{r}^{ipert}) \leftarrow \text{SI\_PERTURB}(\mathbf{x}^{count}, T, t, \beta)$ .
6   end
7    $ipert \leftarrow \operatorname{argmax}_{ipert \in \{1, 2, \dots, L\}} \tilde{r}^{ipert}$ ;
8    $count \leftarrow count + 1$ ;
9   if  $\tilde{r}^{ipert} > r^{count-1}$  then
10     $\mathbf{x}^{count} \leftarrow \tilde{\mathbf{x}}^{ipert}$ ;
11     $r^{count} \leftarrow \tilde{r}^{ipert}$ .
12  end
13  else
14    break.
15  end
16 end
17 function  $(\mathbf{x}^{opt}, r^{opt}) = \text{SI\_PERTURB}(\mathbf{x}^0, T, t, \beta)$ 
18 initial  $r^0, \mathbf{q}^0, \mathbf{p}^0, \sigma^0, \mathbf{s}^0$ .
19  $k \leftarrow 0$ .
20  $r^{opt} \leftarrow 0$ .
21  $\bar{\mathbf{p}}^k \leftarrow \text{P\_BAR}(\mathbf{x}^k, \mathbf{q}^k, \mathbf{p}^k)$ .
22 while  $k < T$  do
23    $\mathbf{x}^{k+1} \leftarrow \text{EXACT\_SOLUTION}(\mathbf{s}^k, r^k)$ ;
24    $r^{k+1} \leftarrow r^k + \delta_F(\mathbf{x}^{k+1}, \mathbf{x}^k)$ ;
25   if  $r^{k+1} > r^{opt}$  then
26      $r^{opt} \leftarrow r^{k+1}$ ;
27      $\mathbf{x}^{opt} \leftarrow \mathbf{x}^{k+1}$ .
28   end
29   if  $r^{k+1} = r^k = \dots = r^{k-t}$  then
30      $\mathbf{x}^{k+1} \leftarrow \text{PERTURB}(\bar{\mathbf{p}}^k, \beta)$ ;
31      $r^{k+1} \leftarrow r^k + \delta_F(\mathbf{x}^{k+1}, \mathbf{x}^k)$ .
32   end
33    $\mathbf{q}^{k+1} \leftarrow \mathbf{q}^k + \delta \mathbf{q}(\mathbf{x}^{k+1}, \mathbf{x}^k)$ ;
34    $\mathbf{p}^{k+1} \leftarrow \mathbf{p}^k + \delta \mathbf{p}(\mathbf{x}^{k+1}, \mathbf{x}^k)$ ;
35    $\bar{\mathbf{p}}^{k+1} \leftarrow \text{P\_BAR}(\mathbf{x}^{k+1}, \mathbf{q}^{k+1}, \mathbf{p}^{k+1})$ ;
36    $(\mathbf{s}^{k+1}, \sigma^{k+1}) \leftarrow \text{SUBGRADIENT}(\mathbf{x}^{k+1}, \bar{\mathbf{p}}^{k+1}, \mathbf{s}^k, \sigma^k)$ ;
37    $k \leftarrow k + 1$ .
38 end
39 end function

```

a naive way via an outer loop in Lines 2-16, and takes L runs of `SI_PERTURB` with different β randomly chosen from $(0, 1)$ within each turn of loop. This outer loop continues until the cut value stops increasing (see Lines 9-15) and the variable *count* records the total number of turns (see Line 8).

Table 4.3 shows the numerical results by SI-P with $t = 3, T = 2000, L = 20$, and $p = \infty$. Now the cut values are all increased for those 27 problem instances in G-set and the ratios between the best cut values and the reference ones become at least 0.997. Moreover, the complexity of the function `SI_PERTURB` is almost the same as `SI`. Therefore, SI-P calls `SI_PERTURB` $count \times L$ times and performs $count \times L \times T$ iterations in total. This is the price we should pay for the improved cut values, which is at most $17 \times 20 \times 2000 = 680000$ iteration steps in the numerical experiments on G-set (see the last column of Table 4.3).

Table 4.3: Improved cut values achieved by the simple iteration with perturbation (SI-P) depicted in Algorithm 4.2 with $t = 3, L = 20, T = 2000$, and $p = \infty$. The ratios between the best cut values by SI-P and the reference ones by combinatorial algorithms [13] are at least 0.997 (see G37). The number of turns of outer loop, recorded by *count* in Line 8 of Algorithm 4.2, is at most 17, thereby meaning that the SI-P algorithm runs at most $count \times L \times T = 17 \times 20 \times 2000 = 680000$ iterations.

Graph	Reference	Result	Ratio	Ratio without perturbation	<i>count</i>
G1	11624	11624	1.0000	0.9939	9
G2	11620	11620	1.0000	0.9946	4
G3	11622	11622	1.0000	0.9985	3
G4	11646	11646	1.0000	0.9959	17
G5	11631	11630	0.9999	0.9975	9
G14	3064	3063	0.9997	0.9899	5
G15	3050	3050	1.0000	0.9862	12
G16	3052	3052	1.0000	0.9885	6
G17	3047	3046	0.9997	0.9879	14
G22	13359	13358	0.9999	0.9948	8
G23	13344	13339	0.9996	0.9940	7
G24	13337	13335	0.9999	0.9963	7
G25	13340	13337	0.9998	0.9924	4
G26	13328	13318	0.9992	0.9911	3
G35	7687	7663	0.9969	0.9869	11
G36	7680	7656	0.9969	0.9865	13
G37	7691	7665	0.9966	0.9884	7
G38	7688	7673	0.9980	0.9876	12
G43	6660	6660	1.0000	0.9976	2
G44	6650	6650	1.0000	0.9956	4
G45	6654	6654	1.0000	0.9917	2
G46	6649	6646	0.9995	0.9929	5
G47	6657	6657	1.0000	0.9929	3
G51	3848	3841	0.9982	0.9891	4
G52	3851	3849	0.9995	0.9920	8
G53	3850	3846	0.9990	0.9914	10
G54	3852	3845	0.9982	0.9920	9

4.4. Comparison with CirCut

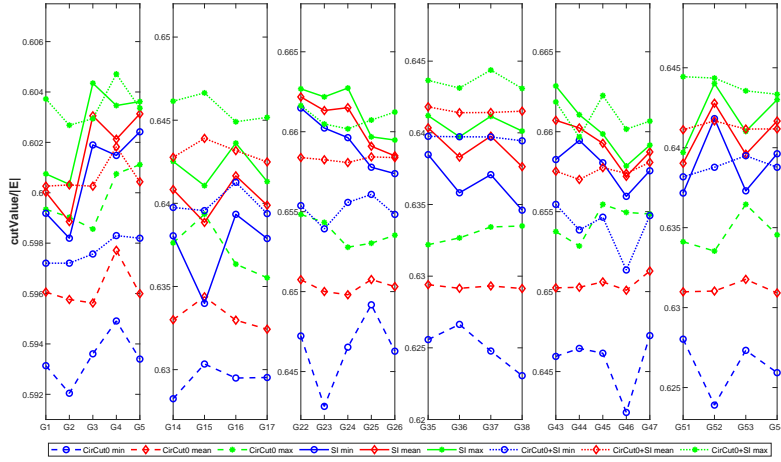
Finally, we compare SI in Algorithm 4.1 and SI_PERTURB in Algorithm 4.2 with the primal CirCut algorithm given in [3, Algorithm 1] which does not invoke any local search techniques. For convenience, we adopt the same notations as those used in [3] unless otherwise specified. It should be noted that CirCut do need a simple gradient algorithm with a backtracking Armijo line-search to minimize the nonconvex objective function $f(\theta)$ from θ^0 instead of calling an external solver. The stopping condition for the line-search is, either the relative change in $f(\theta)$ is less than ϵ_f or the relative change in its gradient is less than ϵ_g . We use T_f (resp. T_g) to count the total number of times $f(\theta)$ (resp. $\nabla f(\theta)$) has been evaluated during the line-search. CirCut stops until N consecutive random perturbations cannot improve the approximate cut, and let T_P count the total number of perturbations. For a fair comparison, we extract the first two steps: line-search and Procedure-CUT, to form a pure rank-two relaxation for maxcut, abbreviated as CirCut0, which excludes the perturbation step, and set it against SI. When the random perturbation is invoked, CirCut is compared with SI_PERTURB. For a given graph, one usually runs the algorithm M times with multiple random starting points: $\theta^0 \sim \mathcal{U}(0, 2\pi)$, and \bar{T}_γ gives the average value of T_γ over these M times for $\gamma \in \{f, g, P\}$. The iteration of SI is not stopped until the cut values remain unchanged for t consecutive steps and we run SI_PERTURB \bar{T}_P times starting from a single initial data where T_S counts the total number of iterations. Both SI and SI_PERTURB are re-run M times from the same initial data given by the maximal eigenvector of the graph Laplacian where \bar{T}_S denotes the average value of T_S .

We set $N = 10, M = 20, t = 3$ and use the implementation of CirCut available at Github¹⁾ where the parameters for the line-search are: the maximum number of rounds $n_{\max} = 200$, and the tolerances $\epsilon_f = \epsilon_g = 1e-4$. Fig. 4.2 plots the minimum, mean and maximum cut values (normalized by the number of edges $|E|$) from the simulations with the M starting points. We are able to observe there that the quality of SI solutions is much better than CirCut0, and the approximate cuts produced by SI_PERTURB are of comparable quality to CirCut with the same number of perturbations.

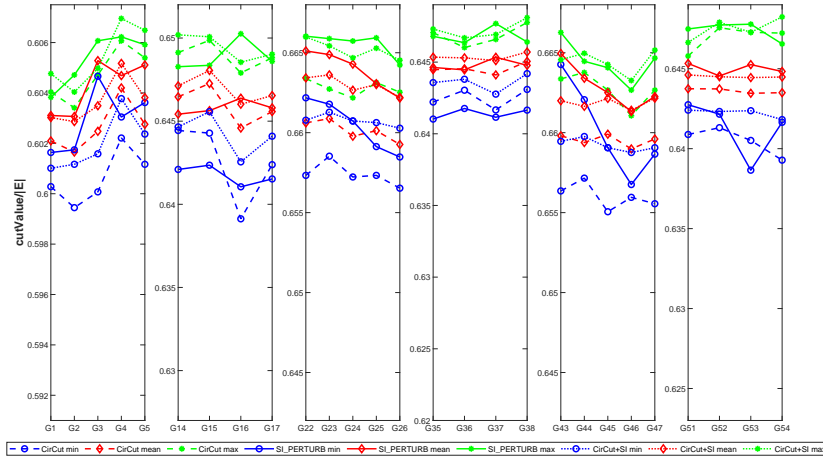
According to Theorems 3.2 and 3.4, the cut values obtained by SI are monotonically updated, and the iterative solution converges to a local optimum from any given initial data. So it will be interesting to see whether SI can improve the output cuts of CirCut0 and the results are displayed in Fig. 4.2(a) with dot lines and legend CirCut0+SI where the maximum number of SI iterations is set to be 100. It is shown there that SI does improve the quality of solutions obtained by CirCut0, which clearly indicates that CirCut0 is not guaranteed to get local optimum. Almost the same story happens with CirCut+SI (see Fig. 4.2(b)). On the other hand, we should point out that CirCut0 cannot improve the quality of the solution produced by SI. In fact, the solution obtained by SI must be a cut, while, by [3, Theorem 3.4], any cut corresponds to a critical point of $f(\theta)$, which means that the stopping condition $\nabla f(\theta) \approx \mathbf{0}$ in the line-search is immediately satisfied, and there is obviously no way to improve it further.

We run all above-mentioned algorithms in Matlab (r2019b) on a High-Performance Computing Platform: 2*Intel Xeon E5-2650-v4 (2.2 GHz, 30 MB Cache, 9.6 GT/s QPI Speed, 12 Cores, 24 Threads) with 128 GB Memory. In Table 4.4, we list the values of \bar{T}_γ with $\gamma \in \{f, g, S, P\}$ as well as the average wall-clock time in seconds needed by only one thread for each run. In each step of the line-search, it is required to calculate $f(\theta)$ and its gradient the complexity of which

¹⁾ see <https://github.com/MQLib/MQLib/tree/master/src/heuristics/maxcut>



(a) SI vs CirCut0



(b) SI_PERTURB vs CirCut

Fig. 4.2. Comparison with CirCut: The minimum, mean, and maximum cut values (normalized by the number of edges $|E|$) produced by CirCut0, SI, and CirCut0+SI from multiple starting points are displayed in (a), while those obtained by CirCut, SI_PERTURB, and CirCut+SI are presented in (b). CirCut0 refers to the pure rank-two relaxation which consists of the line-search and Procedure-CUT, only the first two steps of [3, Algorithm 1]. CirCut0+SI means the output of CirCut0 serves as the input to SI for possible solution quality improvements and so does CirCut+SI.

is $\mathcal{O}(n^2)$ and does not change significantly as the search proceeds. Then, it can be deduced that the run time of both CirCut0 and CirCut should be roughly proportional to $\bar{T}_f + \bar{T}_g$, which can be readily verified in Table 4.4. By comparison, thanks to its monotonicity in Theorem 3.2 and local adjustability in Theorem 3.4, the complexity of one iteration step decreases as SI goes on, and its average over all steps is $\mathcal{O}(\text{mean}(c(k))n)$ and usually much less than $\mathcal{O}(n^2)$ as already shown in Table 4.1. There are two important implications. One is that the time ratio between SI and CirCut0 (or CirCut) should be roughly proportional to $\bar{T}_S/(\bar{T}_f + \bar{T}_g)$ where \bar{T}_S for SI is not so large (less than 50). This can be easily checked in Table 4.4. The other is the run time of

Table 4.4: Comparison with CirCut: The wall-clock time in seconds, the numbers of calculating $f(\theta)$ and its gradient \bar{T}_f and \bar{T}_g in CirCut0 and CirCut, the number of iterations \bar{T}_S in SI and SI_PERTURB, and the same number of perturbations \bar{T}_P in CirCut and SI_PERTURB. It is clearly shown that SI and SI_PERTURB require much fewer iterations and thus run much faster than CirCut0 and CirCut, respectively. CirCut0 refers to the pure rank-two relaxation which only contains the first two steps of [3, Algorithm 1].

Graph	CirCut0			SI		CirCut			SI_PERTURB		
	Time	\bar{T}_g	\bar{T}_f	Time	\bar{T}_s	Time	\bar{T}_g	\bar{T}_f	Time	\bar{T}_s	\bar{T}_p
G1	1.73	74.45	126	0.34	32.65	39.37	1589.05	2817.9	0.61	643.55	30
G2	1.76	74.75	125.55	0.37	33.15	38.33	1544	2763.15	0.55	652.9	30
G3	1.75	73.85	124.35	0.31	32.1	41.82	1745.25	3103.15	0.50	637.25	33
G4	2.09	92.95	156.85	0.32	34.65	37.95	1585.1	2788.6	0.54	617.7	30
G5	1.88	77	130	0.33	29	39.18	1657.25	2909.15	0.56	677.9	31
G14	1.09	75	118.3	0.10	24.15	27.09	2070.7	3183.85	0.60	791.1	34
G15	1.09	75.5	118	0.11	28.25	26.14	2038.9	3137.1	0.58	780.15	32
G16	1.17	81.4	127.5	0.11	25	22.18	1695.25	2628.1	0.46	660.8	28
G17	1.00	76.25	119.35	0.10	23.3	22.75	1752	2694.95	0.45	658.35	29
G22	4.26	65.6	101.9	0.90	31.05	90.58	1344.4	2076.35	3.81	801.15	31
G23	3.51	54.4	84.7	0.91	31.25	104.28	1538.15	2395.05	3.74	795.7	34
G24	3.94	62.35	96.3	0.91	32.9	94.22	1404.1	2175.95	3.69	805.25	31
G25	3.95	63	97.7	0.90	32.45	87.42	1297.55	2007.25	3.50	732.9	29
G26	4.02	63.8	99.9	0.86	32.45	83.02	1244.25	1916.05	3.64	750.6	28
G35	4.50	86.15	136.7	0.58	27.6	155.32	2968.25	4606.1	5.57	1020	40
G36	4.28	83.6	131.2	0.64	30.45	176.06	3405.65	5276.25	6.67	1072.1	42
G37	4.67	88.15	141.5	0.64	31.1	165.54	3206.85	4959.4	6.44	1042.5	41
G38	4.87	92.75	145.7	0.64	29.9	164.64	3217.9	4957.6	7.35	1170.75	43
G43	1.39	62.55	97.45	0.24	33.35	29.69	1275.85	1998.3	0.62	598.75	29
G44	1.41	63.25	99.1	0.24	29.9	34.89	1517.6	2402.25	0.72	746.2	34
G45	1.47	65.55	102.6	0.24	28.2	33.33	1458.75	2319.2	0.69	739.2	32
G46	1.37	60.95	95.15	0.24	27.35	29.63	1262.55	2002.3	0.55	606.75	28
G47	1.37	64.25	99.95	0.23	27.55	25.59	1127.9	1777.85	0.58	624.6	25
G51	1.56	82.5	129.95	0.15	25.1	41.71	2281.45	3516.85	1.02	793.35	35
G52	1.50	79.85	125.55	0.15	24.3	45.90	2506.95	3875	1.22	922.45	38
G53	1.66	88.85	139.55	0.17	25.4	36.32	1959.1	3017.85	0.93	753.15	32
G54	1.49	79.65	124.25	0.15	23.25	35.04	1873.75	2885.05	0.88	690.9	30

SI_PERTURB should grow much more slowly as the iteration goes on where \bar{T}_S is always larger than 500, which can be seen by comparing the time between SI and SI_PERTURB in Table 4.4. Hence, we are able to tell that SI and SI_PERTURB require much fewer iterations and thus run much faster than CirCut0 and CirCut, respectively. That is, although both SI and CirCut do not need call any external solver, quickly locating and checking stationary points of $f(\theta)$ as many as possible in CirCut is not that simple.

5. Conclusion and Outlook

An equivalent continuous fractional optimization problem and a simple iterative (SI) algorithm going from one cut to another in a monotonic and rounding-free way for the maxcut problem were proposed. “Simple” means SI utilizes the exact solutions of the inner subproblems. Numerical experiments on G-set demonstrated that the continuous SI algorithm can produce more qualified solutions than all other existing continuous algorithms. The underlying guiding thought is to build a firm bridge between discrete data world and continuous math field and then use it to design more efficient algorithms. Introducing more advanced combinatorial heuristics into SI and further improving the quality of solutions are on the way. Our attempts on the maxcut problem may provide a valuable reference for other combinatorial problems and fractional programming problems.

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