# COMPLEXITY OF PRIMAL-DUAL INTERIOR-POINT ALGORITHM FOR LINEAR PROGRAMMING BASED ON A NEW CLASS OF KERNEL FUNCTIONS

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In this paper, we first present a polynomial-time primal-dual interior-point method (IPM) for solving linear programming (LP) problems, based on a new kernel function (KF) with a hyperbolic-logarithmic barrier term. To improve the iteration bound, we propose a parameterized version of this function. We show that the complexity result meets the currently best iteration bound for large-update methods by choosing a special value of the parameter. Numerical experiments reveal that the new KFs have better results comparing with the existing KFs including  $\log t$  in their barrier term.

To the best of our knowledge, this is the first IPM based on a parameterized hyperbolic-logarithmic KF. Moreover, it contains the first hyperbolic-logarithmic KF (Touil and Chikouche in Filomat 34:3957-3969, 2020) as a special case up to a multiplicative constant, and improves significantly both its theoretical and practical results.

Keywords: linear programming, primal-dual interior-point methods, kernel function, complexity analysis, large and small-update methods

Classification: 90C05, 90C51

## 1. INTRODUCTION

LP also referred as "linear optimization" is a specific class of mathematical optimization problems. It represents optimization problems that involve linear relationships between the decision variables and the linear objective function, as well as linear constraints. LP has several applications in real life such as production planning, transportation and distribution, resource allocation, portfolio optimization, and many more. Due to the theoretical and practical importance of LPs, various methods were proposed to solve them. IPMs rank among the most efficient methods both theoretically and practically. IPMs were first introduced by Narendra Karmarkar in 1984 [32], the main property of these methods is that they follow the central path in the interior of the feasible region of the problem. The most important results on IPMs for solving LP problems are summarized in the monographs written by Roos et al. [48], Wright [56], Ye [57] and Peng et al. [41].

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Primal-dual IPMs based on KFs represent a fascinating subclass of IPMs. This approach combines two fundamental concepts: primal-dual IPMs and KFs. The choice of the KF plays an important role not only for the IPM analysis but also for the performance of the corresponding interior-point algorithm. Specifically, the KF is used to define an equivalent form of the central path of the IPM, to define the proximity measure, and to obtain search directions. In fact, the central path followed by IPMs is obtained by solving a parametric system that is characterized by a barrier function defined in terms of a KF. In addition, the gradient of the KF serves to define a measure of the distance between the iterates and the central path. Both the KF and the proximity measure effect the iteration bounds since some of their properties play an important role in the complexity analysis. That is why the complexity rate depends on the proposed KF.

The first primal-dual IPM based on the classical logarithmic barrier function was introduced by Roos et al. [48]. After that, Peng et al. introduced primal-dual IPMs based on the so-called self-regular barrier functions, in [39] for LP and semidefinite programming (SDP) and in [40] for second-order conic programming (SOCP). They significantly improved the theoretical complexity obtained for the classical logarithmic KF and obtained the currently best iteration bound for large-update IPMs namely,  $\mathcal{O}(\sqrt{n}\log n\log \frac{n}{\epsilon})$ , with n denotes the number of variables in the problem and  $\epsilon$  denotes the desired accuracy in terms of the objective value. This was one of the main motivations for considering other KFs as a substitute for the classical logarithmic KF. In this context, Bai et al. [7] introduced in 2002 an IPM based on an exponential barrier term which has a finite value at the boundary of the feasible region. The growth term of this finite KF was later parametrized by El Ghami et al. [23] and the approach was extended to solve different types of optimization problems including SDP [26], convex quadratic programming (CQP) [15],  $P_*(\kappa)$  horizontal linear complementarity problem [8] and Cartesian  $P_*(\kappa)$ -linear complementarity problem over symmetric cones (SCLCP) |54|.

In 2004, Bai et al. [6] proposed a new class of eligible KFs which are not necessarily self-regular. This class includes the classical logarithmic KF and self-regular KFs as special cases. They presented a unified analysis of primal-dual IPMs based on eligible KFs for LP. The results obtained in [6] were successfully extended to SDP, CQP, convex quadratic semidefinite programming (CQSDP), SOCP, convex quadratic programming over symmetric cone (SCCQP),  $P_*(\kappa)$ -linear complementarity problem (LCP), the Cartesian  $P_*(\kappa)$ -SCLCP and symmetric programming in [9, 13, 16, 20, 37, 52, 53, 55] respectively.

In [2, 3, 4], published in 2005 and 2007, Amini and his co-authors proposed parametrizations of the two exponential KFs proposed for the first time in [6]. They improved the iteration complexity for large-update methods from  $\mathcal{O}(\sqrt{n}\log^2 n\log\frac{n}{\epsilon})$  to  $\mathcal{O}(\sqrt{n}\log n\log\frac{n}{\epsilon})$ . The IPMs based on these parametric exponential KFs were extended to solve SDP [17, 42],  $P_*(\kappa)$ -LCP [5, 43] and SOCP [18].

Upon scrutiny of the works cited above that focus on IPMs based on KFs, it can be concluded that the theoretical complexity bounds remain unaffected by the type of problem when switching from one problem type to another using, of course, the same KF.

We cannot talk about KFs without mentioning trigonometric KFs. This type of functions has been extensively explored in the literature [14, 30, 31, 34, 36, 38, 44], starting with the work of El Ghami et al. [25] where the authors studied an IPM based on the first trigonometric KF introduced in [6]. They established that the complexity bounds for large- and small-update methods are  $\mathcal{O}(n^{\frac{3}{4}} \log \frac{n}{\epsilon})$  and  $\mathcal{O}(\sqrt{n} \log \frac{n}{\epsilon})$  respectively. This function was later generalized by Bouafia et al. [12] and El Ghami et al. [45] in 2016 for LP and SDP respectively. They obtained  $\mathcal{O}(\sqrt{n} \log n \log \frac{n}{\epsilon})$  iterations complexity for large-update methods which is a significant improvement from the results obtained by El Ghami et al. in [25].

Recently, Touil and Chikouche [51] introduced the first IPM based on a hyperboliclogarithmic KF for SDP. They proved that the corresponding interior-point algorithm meets  $\mathcal{O}(n^{\frac{2}{3}} \log \frac{n}{\epsilon})$  iterations as the worst case complexity bound for the large-update method. In another paper [50], they presented an IPM based on a pure hyperbolic barrier term. By parametrizing the latter, Guerdouh et al. in [29] improved the complexity bound for large-update IPMs from  $\mathcal{O}(n^{\frac{3}{4}} \log \frac{n}{\epsilon})$  to the currently best one. Remaining within the LP case, they proposed an IPM based on a different KF with an exponentialhyperbolic barrier term. The obtained complexity bound for large-update methods meets  $\mathcal{O}(\sqrt{n} \log^2 n \log \frac{n}{\epsilon})$  iterations [28].

Motivated by these works, we first introduce a primal-dual IPM for solving LP problems based on a new hyperbolic-logarithmic KF. Then, we parametrized the new function. The obtained class of KFs contains the hyperbolic-logarithmic KF proposed in [51] as a special case up to a multiplicative constant and improves its theoretical complexity bound to the currently best iteration bound for large-update methods namely,  $\mathcal{O}(\sqrt{n}\log n\log \frac{n}{\epsilon})$ .

We structure our paper as follows. In Section 2, we briefly recall the basics of IPMs for LP. In Section 3, we present the full complexity analysis for the basic KF followed by some numerical tests on eight different examples to illustrate the effectiveness of the proposed KF comparing to other available KFs. In section 4, we introduce a parametrization of the basic KF and derive its complexity bound with some additional numerical experiments. We finally end the paper by providing some concluding remarks.

Let us finish this introduction with some notations used in the whole paper :  $||x|| = \sqrt{x^T x}$  denotes the Euclidean norm of a vector  $x \in \mathbb{R}^n$ .  $\mathbb{R}^n_+$  and  $\mathbb{R}^n_{++}$  denote the nonnegative and the positive orthants respectively, e denotes the all-one vector. For given vectors  $x, s \in \mathbb{R}^n$ , X = diag(x) denotes the  $n \times n$  diagonal matrix whose diagonal entries are the components of x, and the vector xs indicate the component-wise product of x and s. Finally, if  $f(x), g(x) \ge 0$  are two real valued functions of a real nonnegative variable, the notation  $f(x) = \mathcal{O}(g(x))$  means that  $f(x) \le Cg(x)$  for some positive constant C and  $f(x) = \Theta(g(x))$  means that  $C_1g(x) \le f(x) \le C_2g(x)$  for two positive constants  $C_1$  and  $C_2$ .

### 2. PRELIMINARIES

In this section, we provide a brief description of the idea behind the IPMs based on KFs. We consider the standard form of LP problems

$$(P) \qquad \begin{cases} \min c^T x \\ Ax = b, \\ x \ge 0, \end{cases}$$

where  $A \in \mathbb{R}^{m \times n}$ ,  $c \in \mathbb{R}^n$  and  $b \in \mathbb{R}^m$  are given,  $\operatorname{rank}(A) = m \leq n$  and  $x \in \mathbb{R}^n$  is the vector of variables. The dual problem of (P) is given by

(D) 
$$\begin{cases} \max b^T y \\ A^T y + s = c, \\ s \ge 0, \end{cases}$$

where  $y \in \mathbb{R}^m$  and  $s \in \mathbb{R}^n$  are the vectors of variables.

Finding an optimal solution of (P) and (D) is equivalent to solving the non-linear system of equations

$$\begin{cases}
Ax = b, x \ge 0, \\
A^T y + s = c, s \ge 0, \\
xs = 0.
\end{cases}$$
(1)

The first and second pair of equations represent primal and dual feasibility. Replacing the last equation in (1), the so-called complementarity condition for (P) and (D), by the parameterized equation  $xs = \mu e$ , with parameter  $\mu > 0$ , implies that x, s > 0 and leads to the following system

$$\begin{cases}
Ax = b, x > 0, \\
A^T y + s = c, s > 0, \\
xs = \mu e.
\end{cases}$$
(2)

At this stage, we assume that the problems (P) and (D) satisfy the interior-point condition (IPC), i.e., there exists  $(x^0, y^0, s^0)$  such that

$$Ax^0 = b, \ x^0 > 0, \ A^T y^0 + s^0 = c, \ s^0 > 0.$$

This condition means that the interiors of the feasible regions of problems (P) and (D)are both non-empty. However, if the problem doesn't satisfy the IPC, it can be modified so that it does and even in such a way that  $x^0 = s^0 = e$ , the details can be found in [48]. Since the IPC holds, system (2) has a unique solution  $(x_{\mu}, y_{\mu}, s_{\mu})$  for each  $\mu > 0$ , (see [48, Theorem II.4]). The set of unique solutions  $\{(x_{\mu}, y_{\mu}, s_{\mu}) : \mu > 0\}$  forms the so-called central path of problems (P) and (D). If  $\mu \to 0$ , then the limit of the central path exists and yields optimal solutions for (P) and (D).

For fixed  $\mu > 0$ , we apply Newton's method to the parameterized system (2). Thus, the search direction  $(\Delta x, \Delta y, \Delta s)$  is determined by the following system of linear equations

$$\begin{cases} A\Delta x = 0, \\ A^T \Delta y + \Delta s = 0, \\ s\Delta x + x\Delta s = \mu e - xs. \end{cases}$$
(3)

By taking a step along the search direction, one constructs a new iterate point

$$x_+ := x + \alpha \Delta x, \ y_+ := y + \alpha \Delta y, \ s_+ := s + \alpha \Delta s,$$

for some  $0 < \alpha \leq 1$  satisfying the positivity condition i.e.,  $(x_+, s_+) > 0$ . Now, we define the scaled vector v and the scaled search directions  $d_x$  and  $d_s$  as follows

$$v = \sqrt{\frac{xs}{\mu}}, \ d_x = \frac{v\Delta x}{x}, \ d_s = \frac{v\Delta s}{s}.$$
 (4)

System (3) can be rewritten as follows

$$\begin{cases} \overline{A}d_x = 0, \\ \overline{A}^T \Delta y + d_s = 0, \\ d_x + d_s = v^{-1} - v, \end{cases}$$
(5)

where  $\overline{A} = \frac{1}{\mu}AV^{-1}X$ , V = diag(v), X = diag(x).

Observe that the right-hand side in the last equation of (5) is equal to minus gradient of the classical logarithmic scaled barrier (proximity) function

$$\Psi_c(v) = \sum_{i=1}^n \psi_c(v_i),\tag{6}$$

with

$$\psi_c(t) = \frac{t^2 - 1}{2} - \log t.$$

Thus the third equation in (5), often called the scaled centering equation, can be rewritten as follows

$$d_x + d_s = -\nabla \Psi_c(v).$$

This indicates that  $\Psi_c$  essentially determines the search direction. In addition, it's easy to verify that

$$\nabla^2 \Psi_c(v) = \operatorname{diag}(e + v^{-2}),$$

and that the matrix diag $(e + v^{-2})$  is positive definite. This implies that the Hessian  $\nabla^2 \Psi_c(v)$  is positive definite which indicates that  $\Psi_c$  is strictly convex. Moreover,

$$\nabla \Psi_c(e) = 0.$$

It follows that  $\Psi_c(v)$  attains its minimal value at v = e, with  $\Psi_c(e) = 0$ . This means that

$$\nabla \Psi_c(v) = 0 \Leftrightarrow \Psi_c(v) = 0 \Leftrightarrow v = e \Leftrightarrow x = x_\mu \text{ and } s = s_\mu,$$

where  $x_{\mu}$  and  $s_{\mu}$  are the first and third components of the solution of system (2) respectively. Therefore,  $(x_{\mu}, s_{\mu})$  can be identified as the minimizers of the function  $\Psi_c$ . For this reason,  $\Psi_c$  serves mainly as a "proximity" measure of closeness for (x, s) with respect to the  $\mu$ -center. These observations concerning  $\Psi_c$  led to the creation of the concept of primal-dual IPMs based on KFs. The basic idea in these methods is to replace  $\psi_c$ by any strictly convex function  $\psi : ]0, +\infty[ \rightarrow [0, +\infty[$  which is minimal at t = 1 with  $\psi(1) = 0$ . The corresponding proximity function  $\Psi$  is then obtained by replacing  $\psi_c$  by  $\psi$  in (6). This explains the reason for calling  $\psi$  the KF of the barrier function  $\Psi$ . Thus,  $\Psi$  still serves as a proximity measure for closeness with respect to the  $\mu$ -center  $(x_{\mu}, s_{\mu})$ , and the inequality

$$\Psi(v) \le \tau,$$

defines a  $\tau$ -neighbourhood of the  $\mu$ -center.

In the literature, many KFs were proposed with a barrier term constructed by log t combined with another type of functions. In Table 1, we can find all these KFs (to our knowledge) and the complexity results for the corresponding algorithms, starting with the classical logarithmic KF  $\psi_c$  in [48] and its generalized version  $\psi_{1,p}$  proposed by El Ghami et al. in [22].

In this context, we consider in this paper a new univariate KF  $\psi$  defined as follows:

$$\psi(t) = \frac{t^2 - 1}{2} + \tanh^2(1) \big( \coth(t) - \log t \big) - \tanh(1), \ \forall t > 0, \tag{7}$$

and generalize the barrier term of  $\psi$  by applying a positive parameter p to obtain the following parametric KF

$$\psi_p(t) = \frac{t^2 - 1}{2} + \frac{\sinh^2(1)}{\sinh^2(1) + p \coth^{p-1}(1)} \Big( \coth^p(t) - \log t - \coth^p(1) \Big), \quad p \ge 2.$$
(8)

**Remark 2.1.** Taking p = 1 in (8) and using the hyperbolic identity

$$\cosh^2(t) - \sinh^2(t) = 1, \ \forall t \in \mathbb{R},$$

we obtain KF (7).

Coming back to system (5), we can convert it to

$$\begin{cases} \overline{A}d_x = 0, \\ \overline{A}^T \Delta y + d_s = 0, \\ d_x + d_s = -\nabla \Psi(v). \end{cases}$$
(9)

Since A has full row rank, system (9) has a unique solution. Furthermore, the vectors  $d_x$  and  $d_s$  are orthogonal and thus

$$d_x^T d_s = 0 \Leftrightarrow \nabla \Psi(v) = 0 \Leftrightarrow v = e \Leftrightarrow \Psi(v) = 0 \Leftrightarrow x = x_\mu \text{ and } s = s_\mu.$$

In the sequel, we will use the norm-based proximity measure  $\sigma(v)$  defined by

$$\sigma(v) = \frac{1}{2} \|d_x + d_s\| = \frac{1}{2} \|\nabla \Psi(v)\|.$$
(10)

Complexity	$egin{array}{l} \mathcal{O} & \left( n \log rac{n}{\epsilon}  ight) \ \mathcal{O} & \left( n \log rac{n}{\epsilon}  ight) \ \mathcal{O} & \left( n \log rac{n}{\epsilon}  ight) \ \mathcal{O} & \left( n rac{3}{\epsilon} \log rac{n}{\epsilon}  ight) \ \mathcal{O} & \left( p n rac{2r}{2r} \log rac{n}{\epsilon}  ight) \end{array}$	$ \begin{array}{c} \mathcal{O} \left( p \sqrt{n} (\log n)^{\frac{p+1}{p}} \log \frac{n}{\epsilon} \right) \\ \mathcal{O} \left( p^{-1} \sqrt{n} (\log n)^2 \log \frac{n}{\epsilon} \right) \end{array} $	$ \begin{array}{c} \mathcal{O} \left( n^{\frac{2}{3}} \log \frac{n}{\epsilon} \right) \\ \mathcal{O} \left( n^{\frac{2}{3}} \log \frac{n}{\epsilon} \right) \\ \mathcal{O} \left( (1+2k)n^{\frac{2p+1}{4p}} p^{\frac{2p+1}{2p}} \log \frac{n}{\epsilon} \right) \end{array} $	$egin{array}{l} \mathcal{O}\left(n^{rac{2}{3}}\lograc{n}{\epsilon} ight) \ \mathcal{O}\left(n^{rac{3}{4}}\lograc{n}{\epsilon}\lograc{p+2}{\epsilon} ight) \ \mathcal{O}\left(pn^{rac{2(p+1)}{2(p+1)}}\lograc{n}{\epsilon} ight) \end{array}$
Reference $\&$ Type of problem	[48, LP] [22, LP] [11, LP]	[19, LP] [10, LP]	$[46, LP] \\ [14, LP] \\ [30, LCP] ]$	[51, SDP] [new, LP] [new, LP]
Kernel functions	$\psi_{c}(t) = rac{t^{2}-1}{2} - \log t$ $\psi_{1,p}(t) = rac{t^{p+1}-1}{t^{p+1}} - \log t, \ \ 0  \psi_{2}(t) = 8t^{2} - 11t + rac{2}{\sqrt{t}} - 4\log t\psi_{3,p}(t) = rac{t^{2}-1-\log t}{2} + rac{t^{1-p-1}}{2(p-1)}, \ \ p \geq 2$	$\psi_{4,p}(t) = rac{t^2 - 1 - \log t}{2} + rac{e rac{1}{2} - 1 - 1}{2}, \ p \ge 1$ $\psi_{5,p}(t) = t^2 - 1 - \log t + rac{e^p (rac{1}{t} - 1) - 1}{p}, \ p \ge 1$	$\begin{split} \psi_{6}(t) &= \frac{t^{2}-1}{t^{2}} - \log t + \frac{1}{8} \tan^{2}(\pi \frac{1-t}{4t+2}) \\ \psi_{6,\lambda}(t) &= \frac{t^{2}-1}{2} - \log t + \lambda \tan^{2}(\pi \frac{1-t}{3t+2}),  0 < \lambda \leq \frac{8}{25\pi} \\ \psi_{7,p}(t) &= \frac{t^{2}-1}{2} - \log t - \int_{1}^{t} \frac{1}{2p(2+4x)^{2}} \tan^{2p}(\pi \frac{1-x}{4x+2}) dx,  p \in \mathbb{N} \end{split}$	$\begin{split} \psi_{\rm S}(t) &= \frac{1+2\coth(1)}{2}(t^2 - 1) + \coth^2(t) - \coth^2(1) - \log t\\ \psi_{\rm new}(t) &= \frac{t^2 - 1}{2} + \tanh^2(1) \left( \coth(t) - \log t \right) - \tanh(1)\\ \psi_{\rm new,p}(t) &= \frac{t^2 - 1}{2} + a \left( \coth^p(t) - \log t - \coth^p(1) \right),\\ \text{with } a &= \frac{t^2 - 1}{\sinh^2(1) + p\coth^{p-1}(1)} \text{ and } p \ge 2 \end{split}$
Type	əlqmiZ	Exponential	Trigonometric	Нурегьоліс

**Tab. 1.** Existing kernel functions with  $\log t$  in their barrier terms.

Now, we can describe the algorithm briefly. Given a strictly feasible point  $(x^0, y^0, s^0)$ which is in a  $\tau$ -neighbourhood of the given  $\mu$ -center, we reduce  $\mu$  to  $\mu_+ := (1 - \theta)\mu$  for some fixed  $0 < \theta < 1$  and then solve system (9) to obtain the search direction. The positivity condition of a new iterate is ensured by choosing an appropriate step size  $\alpha$ . This procedure is repeated until we find a new iterate  $(x_+, y_+, s_+)$  that again belongs to the  $\tau$ -neighbourhood of the current  $\mu$ -center, that is, until  $\Psi(v) \leq \tau$ . Then, we update the parameter  $\mu$  to  $\mu_+$  and we let  $(x, y, s) = (x_+, y_+, s_+)$ . This procedure is repeated until we find an iterate  $(x_+, y_+, s_+)$  such that  $x_+^T s_+ < \epsilon$ . In this case, an  $\epsilon$ -approximate optimal solution of problems (P) and (D) is found. The generic IPM is outlined in Algorithm 1.

From the above, we can summarize the role of KFs in what follows

- The KF, more specifically the gradient of the barrier function is used to solve system (9) and hence to define new search directions.
- The proximity measure  $\sigma$ , which is used to measure the closeness of the triplet (x, y, s) to the  $\mu$  center, is defined in term of the KF.
- The KF serves to define a τ-neighbourhood of the μ-center and thus it intervenes in the stopping criterion of the inner iteration of the algorithm.

# Algorithm 1 : Generic Interior-Point Algorithm for LP

### Input

a threshold parameter  $\tau \geq 1$ ; an accuracy parameter  $\epsilon > 0$ ; a fixed barrier update parameter  $\theta \in [0, 1]$ ;  $(x^0, y^0, s^0)$  satisfy the IPC and  $\mu^0 = 1$  such that  $\Phi\left(x^0, s^0; \mu^0\right) := \Psi(v^0) \le \tau.$ begin  $x: = x^0; s: = s^0; \mu: = \mu^0;$ while  $n\mu \ge \epsilon$ **begin** (outer iteration)  $\mu := (1 - \theta)\mu;$ while  $\Phi(x, s; \mu) := \Psi(v) > \tau$ **begin** (inner iteration) Solve system (9) and use (4) to obtain  $(\Delta x, \Delta y, \Delta s)$ ; Choose a suitable step size  $\alpha$ ;  $x := x + \alpha \Delta x; y := y + \alpha \Delta y; s := s + \alpha \Delta s; v := \sqrt{\frac{xs}{\mu}};$ end while (inner iteration) end while (outer iteration)

## 3. THE NEW KERNEL FUNCTION

In this section, we give some properties of our new KF and derive the complexity analysis.

#### 3.1. Some technical results

In the analysis of the algorithm based on  $\psi$ , we need its first three derivatives with respect to t which are given for all t > 0 by

$$\psi'(t) = t - \tanh^2(1) \left(\frac{1}{\sinh^2(t)} + \frac{1}{t}\right),$$
(11)

$$\psi''(t) = 1 + \tanh^2(1) \left( 2 \frac{\coth(t)}{\sinh^2(t)} + \frac{1}{t^2} \right),\tag{12}$$

and

$$\psi^{\prime\prime\prime}(t) = -\tanh^2(1)\left(\frac{2}{\sinh^4(t)} + 4\frac{\coth^2(t)}{\sinh^2(t)} + \frac{2}{t^3}\right) < 0.$$
(13)

Clearly,  $\psi'(1) = \psi(1) = 0$ ,  $\lim_{t \to 0^+} \psi(t) = \lim_{t \to +\infty} \psi(t) = +\infty$  and  $\psi''(t) > 0$ ,  $\forall t > 0$ . This implies that  $\psi$  is a KF.

Furthermore, the two functions  $\psi : [1, +\infty[ \rightarrow [0, +\infty[ \text{ and } \psi' :]0, 1] \rightarrow] -\infty, 0]$  are both monotonically increasing. So, we can define the following functions  $\varrho : [0, +\infty[ \longrightarrow [1, +\infty[ \text{ the inverse function of } \psi(t) \text{ for } t \ge 1.$  $\rho : [0, +\infty[ \longrightarrow ]0, 1]$  the inverse function of  $-\frac{1}{2}\psi'(t)$  for  $0 < t \le 1$ .

In what follows, we develop some technical lemmas. From (12), we see that

$$\psi''(t) \ge 1, \ \forall t > 0,$$

thus we have the following lemma.

**Lemma 3.1.** (Bai at al. [7, Lemma 2.1]) Let  $\psi$  be defined as in (7). Then

$$\frac{1}{2}(t-1)^2 \le \psi(t) \le \frac{1}{2}(\psi'(t))^2, \quad \forall t > 0$$

**Corollary 3.2.** (Bai at al. [7, Corollary 2.2]) Let  $\sigma(v)$  be defined by (10). Then, for any v > 0, we have

$$\sigma(v) \ge \sqrt{\frac{\Psi(v)}{2}}.$$

**Remark 3.3.** Throughout the paper, we assume that  $\tau \ge 1$ . Using Corollary 3.2 and the assumption that  $\Psi(v) \ge \tau$ , we have

$$\sigma(v) \ge \sqrt{\frac{1}{2}}.$$

Furthermore, we can write  $\psi(t)$  as

$$\psi(t) = \frac{t^2 - 1}{2} + \psi_b(t),$$

where  $\frac{t^2-1}{2}$  is the growth term and  $\psi_b(t)$  is the barrier term of our KF. Moreover,  $\psi_b(1) = 0$  and  $\psi_b$  is monotonically decreasing. Hence, we can take advantage of the following lemma.

**Lemma 3.4.** (Bai at al. [6, Lemma 6.2]) One has

$$\sqrt{1+2s} \le \varrho\left(s\right) \le 1 + \sqrt{2s}, \ \forall s \in [0, +\infty[s])$$

Now, we give an implicit lower bound for  $\rho(s)$ .

**Lemma 3.5.** For all  $(z,t) \in [0, +\infty[\times]0, 1]$  such that  $z = -\frac{1}{2}\psi'(t)$ , one has

$$\operatorname{coth}(t) \le \operatorname{coth}(1) (2z+2)^{\frac{1}{2}}.$$

Proof. Let  $z \ge 0$  and  $t \in [0, 1]$  such that  $z = -\frac{1}{2}\psi'(t)$ , then  $\rho(z) = t$ . Since  $\frac{1}{\sinh^2(t)} = \coth^2(t) - 1$ , using (11), we have

$$2z = -\psi'(t)$$
  
=  $-t + \tanh^2(1)\left(\coth^2(t) - 1 + \frac{1}{t}\right),$ 

which implies that

$$\operatorname{coth}^{2}(t) = \operatorname{coth}^{2}(1)(2z+t) + 1 - \frac{1}{t}$$
  
 $\leq \operatorname{coth}^{2}(1)(2z+2).$ 

This proves the lemma.

Lemma 3.6. (Touil and Chikouche [50, Lemma 1]) One has

$$2t \operatorname{coth}(t) - 1 > 0, \ \forall t > 0.$$
 (14)

The following lemma reveals some key properties of the new KF.

**Lemma 3.7.** Let  $\psi$  be as defined in (7). Then,

(i) 
$$t\psi''(t) - \psi'(t) > 0, \ \forall t > 0.$$

- (ii)  $t\psi''(t) + \psi'(t) > 0, \ \forall t > 0.$
- (iii)  $\psi''$  is monotonically decreasing on  $]0, +\infty[$ .

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Proof. For the first and second item, using (11) and (12), we have

$$t\psi''(t) - \psi'(t) = \tanh^2(1) \left(\frac{2t\coth(t) + 1}{\sinh^2(t)} + \frac{2}{t}\right) > 0,$$

and

$$t\psi''(t) + \psi'(t) = 2t + \tanh^2(1)\left(\frac{2t\coth(t) - 1}{\sinh^2(t)}\right) > 0,$$

by taking into acount (14) of Lemma 3.6.

For the third item, using (13), we have  $\psi'''(t) < 0$ ,  $\forall t > 0$ , then  $\psi''(t)$  decreases monotonically. This completes the proof.

**Lemma 3.8.** (Peng et al. [41, Lemma 2.1.2]) A twice differentiable function  $\psi : \mathbb{R}_{++} \to \mathbb{R}$  verifies property (ii) of Lemma 3.7 if and only if it verifies one of the following equivalent properties

- $\psi(\sqrt{t_1 t_2}) \leq \frac{1}{2} (\psi(t_1) + \psi(t_2)), \ \forall t_1, t_2 > 0.$
- the function  $\xi \mapsto \psi(e^{\xi})$  is convex.

We say that  $\psi$  is exponentially convex or shortly e-convex.

# 3.2. Growth behaviour

Before updating  $\mu$  in the generic interior-point algorithm, we have  $\Psi(v) \leq \tau$ . After updating  $\mu$  in an outer iteration, the vector v is divided by the factor  $(1 - \theta)$ , which generally leads to an increase of the value of  $\Psi(v)$ . Thus, during the inner iterations, the value of  $\Psi(v)$  decreases until it passes the threshold  $\tau$ . We proceed by studying the effect of updating the barrier parameter  $\mu$  on the value of  $\Psi(v)$ .

**Theorem 3.9.** (Bai at al. [6, Theorem 3.2]) For any v > 0 and  $\beta > 1$ , we have

$$\Psi\left(\beta v\right) \le n\psi\left(\beta \varrho\left(\frac{\Psi(v)}{n}\right)\right)$$

Applying this theorem with  $\beta = \frac{1}{1-\theta}$ ,  $0 < \theta < 1$ , and taking into account the increasing of  $\psi$  on  $[1, +\infty]$ , we get

**Corollary 3.10.** By the assumption  $\Psi(v) \leq \tau$  just before the  $\mu$ -update to  $(1 - \theta)\mu$ , we have

$$\Psi(v_{+}) \le n\psi\left(\frac{\varrho\left(\frac{\tau}{n}\right)}{\sqrt{1-\theta}}\right),\tag{15}$$

where  $v_+ = \frac{v}{\sqrt{1-\theta}}$ .

**Lemma 3.11.** (Bai at al. [6, Lemma 6.3]) Let  $\theta$  be such that  $0 < \theta < 1$ . If  $\Psi(v) \le \tau$ , then

$$\Psi(v_{+}) \leq \frac{\psi''(1)}{2} \frac{\left(\theta\sqrt{n} + \sqrt{2\tau}\right)^2}{1-\theta}.$$

**Corollary 3.12.** Let  $\theta$  be such that  $0 < \theta < 1$ . If  $\Psi(v) \leq \tau$ , then

$$\Psi(v_{+}) \leq \frac{3\coth(1)+1}{(1-\theta)} \left(\theta\sqrt{n} + \sqrt{2\tau}\right)^{2} := \Psi_{0}.$$

 $\Psi_0$  is an upper bound for  $\Psi(v_+)$  during the process of the algorithm.

### 3.3. An estimation of the step size

The purpose of this subsection is to compute a default step size  $\alpha$  such that the iterates  $(x_+, y_+, s_+)$  defined in Algorithm 1 are feasible and the proximity function decreases sufficiently. Note that during an inner iteration the parameter  $\mu$  is fixed. Hence, using (4), we have

$$x_{+} = \frac{x}{v}(v + \alpha d_{x}), \quad s_{+} = \frac{s}{v}(v + \alpha d_{s}).$$

It follows that the new v-vector is given by

$$v_+ = \sqrt{\frac{x_+s_+}{\mu}} = \sqrt{(v + \alpha d_x)(v + \alpha d_s)}.$$

The e-convexity property of  $\psi$  implies that

$$\Psi(v_{+}) \leq \frac{1}{2} \left( \Psi(v + \alpha d_{x}) + \Psi(v + \alpha d_{s}) \right).$$

Now, we consider the decrease in  $\Psi$  as a function of  $\alpha$  noted f defined by

$$f(\alpha) = \Psi(v_{+}) - \Psi(v).$$

In the remainder of this section, we put for simplicity  $\sigma := \sigma(v)$ . From Lemmas 4.3, 4.4, 4.5 and Theorem 4.6 in [6], we have the following theorem.

**Theorem 3.13.** (Bai at al. [6]) The largest possible value of the step size  $\alpha^*$  is

$$\alpha^* = \frac{\rho\left(\sigma\right) - \rho\left(2\sigma\right)}{2\sigma}.$$

Furthermore

$$\alpha^* \ge \frac{1}{\psi''\left(\rho\left(2\sigma\right)\right)},$$

and we have for all  $\alpha \in [0, \alpha^*]$ :

$$f(\alpha) \le -\alpha \sigma^2.$$

**Corollary 3.14.** Let us set  $\bar{\alpha} = \frac{1}{\psi''(\rho(2\sigma))}$ , as the default step size. Then

$$f(\bar{\alpha}) \le -\frac{\sigma^2}{\psi''(\rho(2\sigma))}.$$
(16)

We can obtain the upper bound for the decreasing value of the proximity in the inner iteration by the following theorem.

**Theorem 3.15.** If  $\bar{\alpha}$  is the default step size and  $\sigma \geq 1$ , then we have

$$f\left(\bar{\alpha}\right) \le -\frac{\Psi(v)^{\frac{1}{4}}}{273}.\tag{17}$$

Proof. Using (12) and (14) we have

$$\psi''(t) \leq 1 + \tanh^2(1) \left( 2 \coth(t) (\coth^2(t) - 1) + 4 \coth^2(t) \right)$$
  
$$\leq 1 + \tanh^2(1) \left( 2 \coth^3(t) + 4 \coth^2(t) \right)$$
  
$$\leq 1 + 6 \tanh^2(1) \coth^3(t).$$

Hence, putting  $t = \rho \left( 2\sigma \right)$  and using Lemma 3.5 we get

$$\psi''(\rho(2\sigma)) \le (1 + 6 \tanh^2(1)) \coth^3(\rho(2\sigma))$$
  
$$\le \coth^3(1) (1 + 6 \tanh^2(1)) (4\sigma + 2)^{\frac{3}{2}}$$
  
$$\le \coth^3(1) (1 + 6 \tanh^2(1)) (4\sigma + 2(2\sigma))^{\frac{3}{2}}$$

Moreover,

$$f(\bar{\alpha}) \leq -\frac{\sigma^2}{\psi''\left(\rho\left(2\sigma\right)\right)}$$

This implies that

$$f(\bar{\alpha}) \leq -\frac{\sigma^2}{\coth^3(1)\left(1+6\tanh^2(1)\right)(8\sigma)^{\frac{3}{2}}} \\ = -\frac{\sqrt{\sigma}}{8^{\frac{3}{2}}\coth^3(1)\left(1+6\tanh^2(1)\right)}.$$

Using Corollary 3.2, we get

$$\begin{split} f\left(\bar{\alpha}\right) &\leq -\frac{\Psi(v)^{\frac{1}{4}}}{2^{\frac{1}{4}}8^{\frac{3}{2}}\coth^{3}(1)\left(1+6\tanh^{2}(1)\right)} \\ &\leq -\frac{\Psi(v)^{\frac{1}{4}}}{273}. \end{split}$$

Hence the theorem is proved.

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#### 3.4. Iteration complexity

We need to compute how many inner iterations are required to return to the situation where  $\Psi(v) \leq \tau$  after  $\mu$ -update. Let us define the value of  $\Psi(v)$  after  $\mu$ -update as  $\Psi_0$ , and the subsequent values in the same outer iteration as  $\Psi_k, k = 1, \ldots, K$ , where K stands for the total number of inner iterations in the outer iteration. By the definition of  $f(\alpha)$  and according to (17), for  $k = 1, \ldots, K - 1$ , we obtain

$$\Psi_{k+1} \le \Psi_k - \frac{\Psi_k^{\frac{1}{4}}}{273}.$$

**Lemma 3.16.** (Peng et al. [39, Proposition 2.2]) Suppose  $t_0, t_1, \ldots, t_k$  be a sequence of positive numbers such that

$$t_{k+1} \le t_k - \beta t_k^{1-\gamma}, k = 0, 1, \dots, K-1,$$

where  $\beta > 0$  and  $0 < \gamma \leq 1$ . Then

$$K \le \left[\frac{t_0^{\gamma}}{\beta \gamma}\right].$$

As a consequence, by taking  $t_k = \Psi_k$ ,  $\beta = \frac{1}{273}$  and  $\gamma = \frac{3}{4}$ , we get the following lemma.

**Lemma 3.17.** Let K be the total number of inner iterations in the outer iteration. Then, we have

$$K \le 364 \, \Psi_0^{\frac{3}{4}}.$$

Now we derive the complexity bounds for large and small-update methods.

**Theorem 3.18.** The total number of iterations to obtain an approximate solution with  $n\mu \leq \epsilon$  is bounded by

$$\left(364\,\Psi_0^{\frac{3}{4}}\frac{\log\frac{n}{\epsilon}}{\theta}\right).\tag{18}$$

Proof. It is well known that the number of outer iterations for the situation  $n\mu \leq \epsilon$  is bounded above by  $\frac{1}{\theta}(\log \frac{n}{\epsilon})$  (see [48, Lemma II.17, page 116]).

Knowing that an upper bound for the total number of iterations is obtained by multiplying the number of inner and outer iterations, we obtain the result thanks to the above lemma.  $\Box$ 

For large-update IPMs with  $\tau = \mathcal{O}(n)$  and  $\theta = \Theta(1)$ , we have  $\mathcal{O}\left(n^{\frac{3}{4}}\log\frac{n}{\epsilon}\right)$  iteration complexity for LP problems based on our new KF.

For small-update IPMs with  $\tau = \mathcal{O}(1)$  and  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$ , we get the currently best known iteration bound, namely  $\mathcal{O}\left(\sqrt{n}\log\frac{n}{\epsilon}\right)$  iterations.

#### 3.5. Numerical tests

In this subsection, we carry out numerical experiments of the interior-point algorithm based on different KFs. Our experiments are implemented in MATLAB R2012b and performed on Supermicro dual-2.80 GHz Intel Core i5 server with 4.00 Go RAM.

We have taken  $\epsilon = 10^{-8}$  and  $\tau = n$ . We chose a practical step size  $\alpha$  as in [33] i.e.,  $\alpha = \min(\alpha_x, \alpha_s)$ , with

$$\alpha_x = \min_{i=1..n} \begin{cases} -\frac{x_i}{\Delta x_i} & \text{if } \Delta x_i < 0, \\ 1 & \text{elsewhere,} \end{cases}$$

and

$$\alpha_s = \min_{i=1..n} \begin{cases} -\frac{s_i}{\Delta s_i} & \text{if } \Delta s_i < 0, \\ 1 & \text{elsewhere.} \end{cases}$$

This choice of  $\alpha$  guarantees the strict positivity of the new point. Moreover, we increase the step size by a fixed factor  $0 < \beta < 1$  (in our case we chose  $\beta = 0.9$ ).

Our purpose is to show that the generic interior-point algorithm based on our hyperboliclogarithmic KF can be very efficient in solving LP problems. As in [24], we conducted comparative numerical tests between the KFs provided in Table 1 on the following eight test problems with different sizes, ranging from very small to big size problems.

#### 3.5.1. Examples with fixed size

All examples with fixed size are taken from [1].

**Example 3.19.** m = 2, n = 4.

$$A = \begin{pmatrix} 1 & 1 & 1 & 1 \\ 1 & 1 & 0 & -3 \end{pmatrix}, \quad b = (1, 0.5)^T, \quad c = (1, 2, 3, 4)^T,$$

where the initial feasible solutions are defined as follows

$$x^{0} = (0.5, 0.27, 0.14, 0.09)^{T}, \quad y^{0} = (0, 0)^{T}, \quad s^{0} = (1, 2, 3, 4)^{T}.$$

The optimal solution is  $x^* = (0.87500, 0, 0, 0.12500)^T,$   $y^* = (1.75000, -0.75000)^T$  $s^* = (0.00000, 1.00000, 1.25000, 0.00000)^T.$ 

**Example 3.20.** m = 3, n = 5.

$$A = \begin{pmatrix} 2 & 1 & 1 & 0 & 0 \\ 1 & 2 & 0 & 1 & 0 \\ 0 & 1 & 0 & 0 & 1 \end{pmatrix}, \quad b = (8,7,3)^T, \quad c = (-4, -5, 0, 0, 0)^T,$$

where the initial feasible solutions are defined as follows  $x^0 = (2.85, 1.9, 0.4, 0.35, 1.1)^T$ ,  $y^0 = (-1.2, -1.8, -0.5)^T$ ,  $s^0 = (0.2, 0.3, 1.2, 1.8, 0.5)^T$ . The optimal solution is

 $\begin{aligned} x^* &= (3.0000, 2.0000, 0.0000, 0.0000, 1.0000)^T, \\ y^* &= (-1.0000, -2.0000, -0.0000)^T, \\ s^* &= (0.0000, 0.0000, 1.0000, 2.0000, 0.0000)^T. \end{aligned}$ 

# **Example 3.21.** m = 3, n = 6.

$$A = \begin{pmatrix} 2 & 1 & 0 & -1 & 0 & 0 \\ 0 & 0 & 1 & 0 & 1 & -1 \\ 1 & 1 & 1 & 1 & 1 & 1 \end{pmatrix}, \quad b = (0,0,1)^T, \quad c = (3,-1,1,0,0,0)^T,$$

where the initial feasible solutions are defined as follows  $x^0 = (0.06757, 0.13258, 0.13302, 0.26774, 0.13302, 0.2664)^T, y^0 = (-2, -2, -3)^T,$   $s^0 = (10, 4, 6, 1, 5, 1)^T.$ The optimal solution is  $x^* = (0.0000, 0.5000, 0.0000, 0.5000, 0.0000, 0.0004)^T,$   $y^* = (-0.5000, -0.5000, -0.5000)^T,$  $s^* = (4.5000, 0.0000, 2.0000, 0.0000, 1.0000, 0.0000)^T.$ 

## **Example 3.22.** m = 5, n = 9.

$$A = \begin{pmatrix} 0 & 1 & 2 & -1 & 1 & 1 & 0 & 0 & 0 \\ 1 & 2 & 3 & 4 & -1 & 0 & 1 & 0 & 0 \\ -1 & 0 & -2 & 1 & 2 & 0 & 0 & 1 & 0 \\ 1 & 2 & 0 & -1 & -2 & 0 & 0 & 0 & 1 \\ 1 & 3 & 4 & 2 & 1 & 0 & 0 & 0 & 0 \end{pmatrix}, \quad b = \begin{pmatrix} 1 \\ 2 \\ 3 \\ 2 \\ 1 \end{pmatrix}, \quad c = (1, 0, -2, 1, 1, 0, 0, 00)^T,$$

where the initial feasible solutions are defined as follows  $\begin{aligned} x^0 &= (0.1819, 0.0699, 0.063, 0.1105, 0.2012, 0.6732, 1.1885, 2.835, 2.1912)^T, \\ y^0 &= (-1.3843, -0.8751, -0.4241, -0.4463, -3.0424)^T, \\ s^0 &= (4.9398, 13.1544, 14.7156, 9.1788, 4.5072, 1.3843, 0.8751, 0.4241, 0.4463)^T. \\ \text{The optimal solution is} \\ x^* &= (0.0000, 0.0000, 0.2664, 0.0000, 0.0000, 0.4269, 1.1406, 3.5729, 2.0000)^T, \\ y^* &= (-0.0000, -0.0000, -0.0000, -0.0000, -0.4999)^T, \\ s^* &= (1.5000, 1.4999, 0.0000, 1.9999, 1.4999, 0.0000, 0.0000, 0.0000), 0.0000)^T. \end{aligned}$ 

# 3.5.2. Examples with variable size

**Example 3.23.** (Bouafia et al. [12]) The matrix A is defined as

$$A(i,j) = \begin{cases} 1 & \text{if } i = j \text{ or } j = i + m, \\ 0 & \text{otherwise,} \end{cases}$$

c(i) = -1, c(i + m) = 0 and b(i) = 2, for i = 1, ..., m. We start by an initial point  $(x^0, y^0, s^0)$  such that  $x^0 = e, y^0(i) = -2$  and  $s^0(i) = 1, s^0(i + m) = 2$ , for i = 1, ..., m. The optimal solution is  $x^*(i) = 2, x^*(i + m) = 0, y^*(i) = -1, s^*(i) = 0, s^*(i + m) = 1$ , for i = 1, ..., m. **Example 3.24.** (Anane [1, Example 5]) The matrix A is defined as

$$A(i,j) = \begin{cases} 1 & \text{if } i = j \text{ or } j = i + m, \\ 0 & \text{otherwise,} \end{cases}$$

c = -e and b(i) = 2, i = 1, ..., m. We start by an initial point  $(x^0, y^0, s^0)$  such that  $x^0 = s^0 = e$  and  $y^0(i) = -2$ , for i = 1, ..., m. The optimal solution is  $x^* = e, s^* = 0$ , and  $y^*(i) = -1$ , for i = 1, ..., m.

**Example 3.25.** (Touil et al. [49, Example 4]) The matrix A is defined as

$$A(i,j) = \begin{cases} 1 & \text{if } i = j \text{ or } j = i + m, \\ 0 & \text{otherwise,} \end{cases}$$

c = -e and b(i) = 2, i = 1, ..., m. We start by an initial point  $(x^0, y^0, s^0)$  such that  $x^0(i) = 1.5, x^0(i+m) = 0.5, y^0(i) = -2$ , for i = 1, ..., m, and  $s^0 = e$ . The optimal solution is  $x^*(i) = 1.4793, x^*(i+m) = 0.5207, y^*(i) = -1$ , for i = 1, ..., m, and  $s^* = 0$ .

**Example 3.26.** The matrix A is defined as

$$A(i,j) = \begin{cases} 1 & \text{if } i = j \text{ or } j = i + m, \\ 0 & \text{otherwise,} \end{cases}$$

 $\begin{array}{l} c(i) = \frac{1}{4}, \, c(i+m) = \frac{3}{4} \text{ and } b\,(i) = 2, \, i = 1, \ldots, m. \\ \text{We start by an initial point}\,\,(x^0, y^0, s^0) \text{ such that} \\ x^0(i) = \frac{5}{4}, \, x^0(i+m) = \frac{3}{4}, \, y^0\,(i) = \frac{-7}{12}, \text{ and } s^0(i) = \frac{5}{6}, \, s^0(i+m) = \frac{4}{3}, \, \text{for } i = 1, \ldots, m. \\ \text{The optimal solution is} \\ x^*(i) = 2, \, x^*(i+m) = 0, \, y^*\,(i) = \frac{1}{4}, \, \text{and } s^*(i) = 0, \, s^*(i+m) = \frac{1}{2}, \, \text{for } i = 1, \ldots, m. \end{array}$ 

Problems with fixed size were tested for multiple values of  $\theta$ ,  $\theta \in \{0.1, 0.3, 0.5, 0.7, 0.9\}$ . The table corresponding to fixed size problems (Table 2) shows that, in most cases, larger  $\theta$  gives better iteration numbers. Thus, for variable size problems, we only choose  $\theta \in \{0.7, 0.9, 0.99\}$ . This reduces the number of experiments since we perform Algorithm 1 with the considered KFs on four test problems for seven different sizes n = 2m where  $m \in \{5, 25, 50, 100, 200, 400, 1000\}$ . This left us with 104 experiments for each KF. The summary of results is given in tables below.

	Exa	amp	ole	4		Exa	amp	ole	3		Exa	amp	ole	2		Exa	amp	ole	1	Ex
0.9	0.7	0.5	0.3	0.1	0.9	0.7	0.5	0.3	0.1	0.9	0.7	0.5	0.3	0.1	0.9	0.7	0.5	0.3	0.1	θ
24	28	31	<b>5</b> 8	196	33	17	30	57	192	11	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_c$
67	45	38	58	196	73	44	41	57	192	45	36	36	57	191	57	41	38	56	188	$\psi_{1,rac{1}{2}}$
41	22	34	112	240	34	17	32	61		12	19	31	63	243	18	36	47	104		$\psi_2$
21	<b>21</b>	30	58	196	28	18	30	57	192	10	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_{3,2}$
24	28	31	58	196	33	17	30	57	192	11	17	<b>29</b>	57	191	33	17	<b>29</b>	56	188	$\psi_{3,rac{\log n}{2}}$
24	<b>21</b>	33	58	196	19	20	30	57	197	10	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_{4,1}$
24	24	30	<b>5</b> 8	196	21	20	30	62	197	10	17	<b>29</b>	57	191	13	17	32	56	188	$\psi_{4,2}$
22	28	36	66	220	22	23	37	69	206	9	25	35	75	215	13	17	31	60	190	$\psi_{5,2}$
23	28	36	74	214	22	21	30	62	195	9	21	33	57	191	12	17	<b>29</b>	56	188	$\psi_{5,\log n}$
23	22	31	58	196	39	17	30	57	192	11	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_6$
23	22	31	58	196	26	17	30	57	192	11	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_{6,0.05}$
23	22	31	58	196	42	17	30	57	192	11	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_{7,2}$
23	22	31	58	196		17	30	57	192	11	17	29	57	191	11	17	29	56	188	$\psi_{7,[\log n]}$
20	24	42	78	212	25	20	33	66	204	9	21	39	75	215	14	23	44	70	200	$\psi_8$
21	<b>21</b>	မို	58 58	196	21	18	30	57	192	10	17	<b>29</b>	57	191	11	17	<b>29</b>	56	188	$\psi_{ m new}$

Tab. 2. Number of inner iterations for fixed size examples.

$\psi_{ m new}$	18	19	20	20	21	21	22	14	15	15	17	17	17	18	28	28	30	30	30	30	30
$\psi_8 \psi$	35		37		40	40	41	12	ŝ	ŝ	14	4	4	Ŋ	H	П	ŝ	ŝ	23	ŝ	23
	с. 	<u></u>	<u></u>	<u></u>	4	4	4	-	-	-	-	-	-	-	2	2	2	2	2	2	5
$\psi_{7,[\log n]}$	18	19	20	20	21	21	22	15	20	22	17	17	18	20	37	45	57	57	57	55	54
$\psi_{7,2}$	18	19	20	20	21	21	22	15	17	17	18	18	18	19	37	37	40	40	40	40	40
$\psi_{6,0.05}$	18	19	20	20	<b>21</b>	<b>21</b>	22	16	17	17	18	18	18	20	30	30	33	33	33	33	33
$\psi_6$	18	19	20	20	21	21	22	16	17	17	19	19	19	20	28	28	31	31	31	31	31
$\psi_{5,\log n}$	32	36	39	46	62	49	77	26	I	I	I	I	I	I	32	41	42	42	46	I	I
$\psi_{5,2}$	26	27	28	28	31	31	34	31	32	32	33	33	33	34	15	15	17	17	17	17	17
$\psi_{4,2}$	21	22	23	23	24	24	25	23	24	25	25	25	25	26	40	40	42	42	42	42	42
$\psi_{4,1}$	18	19	20	20	21	21	22	13	14	14	16	16	16	17	39	39	41	41	41	41	41
$\psi_{3,\frac{\log n}{2}}$	18	19	20	20	<b>21</b>	<b>21</b>	22	25	15	15	16	15	16	17	21	30	29	27	30	24	31
$\psi_{3,2}$	18	19	20	20	<b>21</b>	21	22	14	15	15	16	16	16	18	29	29	31	31	31	31	31
$\psi_2$	_	I	I		I	l	l	I	I	I		I	I		20	20	22	22	22	22	22
$\psi_{1,\frac{1}{2}}$	44	46	49	49	52	52	54	58	63	63	60	60	60	60	89	89	104	104	104	104	104
$\psi_c$	18	19	20	20	<b>21</b>	21	22	21	23	23	24	24	24	25	21	21	24	24	24	24	24
m	5	25	50	100	200	400	1000	5	25	50	100	200	400	1000	5	25	50	100	200	400	1000
θ			7.(	) =	θ					6.0	) =	θ					66	0 =	θ		

**Tab. 3.** Number of inner iterations for Example 5 with different sizes n = 2m.

		$\theta$	= 0	.99					θ	= (	).9					θ	= (	).7			θ
1000	400	200	100	50	25	сл	1000	400	200	100	50	25	თ	1000	400	200	100	50	25	თ	m
13	13	13	13	13	11	11	15	13	13	13	12	12	11	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_c$
73	73	73	73	73	60	60	55	51	51	51	45	45	40	48	46	46	44	44	41	39	$\psi_{1,\frac{1}{2}}$
12	12	12	12	12	10	10	12	11	11	11	10	10	9				I			54	$\psi_2$
13	13	13	13	13	10	10	14	12	12	12	11	11	10	22	21	21	20	20	19	<b>18</b>	$\psi_{3,2}$
13	13	13	13	13	10	11	13	12	12	12	11	11	11	22	<b>21</b>	<b>21</b>	20	20	19	<b>18</b>	$\psi_{3,rac{\log n}{2}}$
13	13	13	13	13	10	10	13	12	12	12	11	11	10	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_{4,1}$
13	13	13	13	13	10	10	13	12	12	12	11	11	10	<b>22</b>	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_{4,2}$
12	12	12	12	12	10	10	12	11	11	11	10	10	9	42	41	41	40	40	37	34	$\psi_{5,2}$
12	12	12	12	12	10	10	12	11	11	11	10	10	9		I	I	32	45	28	30	$\psi_{5,\log n}$
13	13	13	13	13	11	11	14	13	13	13	12	12	11	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_6$
13	13	13	13	13	11	11	14	13	13	13	12	12	11	22	<b>21</b>	<b>21</b>	20	20	19	<b>18</b>	$\psi_{6,0.05}$
13	13	13	13	13	11	11	15	13	13	13	12	12	11	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_{7,2}$
13	13	13	13	13	11	11	15	13	13	13	12	12	11	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_{7,[\log n]}$
12	12	12	12	12	10	10	12	11	11	11	10	10	9	34	33 33	33 33	32	32	29	28	$\psi_8$
13	13	13	13	13	10	10	14	12	12	12	11	11	10	22	21	21	20	20	19	18	$\psi_{\rm new}$

$\psi_{\mathrm{new}}$	18	19	20	20	21	21	22	10	11	11	12	12	12	14	10	10	13	13	13	13	13
$\psi_8$		I	I	1				6	10	10	11	11	11	12	10	10	12	12	12	12	12
$\psi_{7,[\log n]}$	18	19	20	20	21	21	22	11	12	12	13	13	13	15	11	11	13	13	13	13	13
$\psi_{7,2}$	18	19	20	20	21	21	22	11	12	12	13	13	13	15	11	11	13	13	13	13	13
$\psi_{6,0.05}$	18	19	20	<b>20</b>	<b>21</b>	<b>21</b>	22	11	12	12	13	13	13	15	11	11	13	13	13	13	13
$\psi_6$	18	19	20	20	21	21	22	11	12	12	13	13	13	15	11	11	13	13	13	13	13
$\psi_{5,\log n}$		I	I					6	10	10	11	11	11	14	10	10	12	12	12	14	14
$\psi_{5,2}$		I	I	I	I	I	I	6	10	10	11	11	11	12	10	10	12	12	12	12	12
$\psi_{4,2}$	18	19	20	20	21	21	22	10	11	11	12	12	12	13	10	10	13	13	13	13	13
$\psi_{4,1}$	18	19	20	<b>20</b>	<b>21</b>	<b>21</b>	22	10	11	11	12	12	12	13	10	10	13	13	13	13	13
$\psi_{3,\frac{\log n}{2}}$	18	19	20	20	<b>21</b>	<b>21</b>	22	11	11	11	12	12	12	13	11	10	13	13	13	13	13
$\psi_{3,2}$	18	19	20	20	<b>21</b>	<b>21</b>	22	10	11	11	12	12	12	14	10	10	13	13	13	13	13
$\psi_2$	_	I	Ι	I	I	I	I	6	10	10	11	11	11	12	10	10	12	12	12	12	12
$\psi_{1,\frac{1}{2}}$	42	44	47	47	49	49	52	44	49	49	54	54	54	60	64	64	80	80	80	80	80
$\psi_c$	18	19	20	20	21	21	22	11	12	12	13	13	13	15	11	11	13	13	13	13	13
m	5	25	50	100	200	400	1000	5	25	50	100	200	400	1000	5	25	50	100	200	400	1000
θ			7.0	) =	θ					6.0	) =	θ					66	0 =	θ		

**Tab. 5.** Number of inner iterations for Example 7 with different sizes n = 2m.

		θ	= 0	.99					θ	= (	).9					θ	= (	).7			θ
1000	400	200	100	50	25	сл	1000	400	200	100	50	25	сл	1000	400	200	100	50	25	თ	m
23	23	23	23	23	21	21	16	15	15	15	14	14	12	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_c$
96	96	96	96	96	82	82	72	67	67	67	61	61	56	52	50	50	47	47	45	42	$\psi_{1,\frac{1}{2}}$
25	25	25	25	25	23	23	35	34	34	34	33	33 33	32	62	59	59	56	56	53	50	$\psi_2$
28	28	28	28	28	26	26	15	14	14	14	13	13	12	22	21	21	20	20	19	18	$\psi_{3,2}$
28	24	23	22	26	26	22	16	<b>14</b>	14	<b>14</b>	13	13	12	<b>22</b>	<b>21</b>	<b>21</b>	<b>20</b>	<b>20</b>	19	18	$\psi_{3,rac{\log n}{2}}$
30	30	30	30	30	28	28	16	15	15	15	14	14	13	<b>22</b>	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_{4,1}$
38	38	38	38	38	36	36	20	19	19	19	17	17	16	22	<b>21</b>	<b>21</b>	20	20	19	<b>18</b>	$\psi_{4,2}$
30	30	30	30	30	28	28	19	18	18	18	17	17	16	31	30	30	29	29	28	27	$\psi_{5,2}$
43	45	38	100	36	31	33	29	32	92	50	25	49	16				59	52	31	29	$\psi_{5,\log n}$
26	26	26	26	26	24	24	16	15	15	15	14	14	12	22	<b>21</b>	<b>21</b>	<b>20</b>	20	19	18	$\psi_6$
26	26	26	26	26	24	24	16	15	15	15	14	14	12	<b>22</b>	<b>21</b>	<b>21</b>	<b>20</b>	<b>20</b>	19	18	$\psi_{6,0.05}$
29	29	29	29	29	27	27	16	15	15	15	14	14	12	22	<b>21</b>	<b>21</b>	20	20	19	<b>18</b>	$\psi_{7,2}$
34	33	40	40	40	32	27	18	16	16	16	15	14	12	22	<b>21</b>	<b>21</b>	20	20	19	18	$\psi_{7,[\log n]}$
21	<b>21</b>	<b>21</b>	<b>21</b>	<b>21</b>	19	19	15	<b>14</b>	<b>14</b>	<b>14</b>	13	13	12	36	33	33	32	32	31	28	$\psi_8$
26	26	26	26	26	23	23	15	14	14	14	13	13	12	22	21	21	<b>20</b>	20	19	18	$\psi_{\mathrm{new}}$

Tab. 6. Number of inner iterations for Example 8 with different sizes

n = 2m.

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## 3.5.3. Comments

Recall that the numerical results were obtained by performing Algorithm 1 with the KFs defined in Table 1 on eight test problems.

When there are parameters  $p \ge 1$  involved in the definition of a KF, we used two values of these parameters: the parameter that gives the best theoretical iteration bound and a common value p = 2, except for  $\psi_{1,p}$  and  $\psi_{6,\lambda}$  where we chose  $p = \frac{1}{2}$  and  $\lambda = 0.05$  since the parameters p and  $\lambda$  take their values in the intervals ]0,1] and  $]0,\frac{8}{25\pi}]$  respectively. This left us with 15 different KFs.

For each example, we used **bold** font to highlight the best, i.e., the smallest, iteration number.

From Tables 2-6, we may draw a few conclusions:

- For  $\psi_{1,p}$  (the classical logarithmic barrier function  $\psi_c$  occurs if p = 1), although the theoretical iteration bound of the algorithm is independent of the parameter p, numerical tests show that p influences the iteration count. In both types of problems (fixed or variable size), p = 1 gives better results which is in accordance with the analysis carried out by [22].
- The function  $\psi_{1,\frac{1}{2}}$  never gives the smallest iteration number in examples with variable size, even for examples with fixed size it gives the smallest iteration number only for the values 0.1 and 0.3 of  $\theta$ .
- For KFs  $\psi_2, \psi_{5,2}, \psi_{5,\log n}, \psi_{7,[\log n]}$  and  $\psi_8$ , the dashes in the corresponding columns of Tables 2–6 indicate that the algorithms require more than 10<sup>4</sup> iterations to obtain an optimal solution. Despite this,  $\psi_2, \psi_{5,2}, \psi_{5,\log n}$  and  $\psi_8$ , (and exclusively  $\psi_8$ ) are the only ones to give the smallest iteration number for  $\theta = 0.9$  in Examples 6 and 7 (in Example 5), while for  $\theta = 0.99$  in Example 5 (resp. in Example 8),  $\psi_{5,2}$  (resp.  $\psi_8$ ) is the only function to achieve the best iteration number.
- The iteration numbers of the algorithm based on our KF  $\psi_{\text{new}}$  depend on the values of the parameter  $\theta$ . In fact, the value 0.9 of  $\theta$  gives better iteration numbers in general.
- In all examples with variable size for  $\theta = 0.7$ , the KF  $\psi_{\text{new}}$  has the smallest iterations number with the KFs  $\psi_c, \psi_{3,p}, \psi_{4,p}, \psi_6, \psi_{6,0.05}, \psi_{7,2}, \psi_{7,[\log n]}$  except  $\psi_{4,2}$  in Example 5.
- For our KF, the obtained iteration numbers coincided with, or at worst was close to the best ones with a slackness of at most 4 iterations, except the case  $\theta = 0.99$  in Example 5, where the slackness attain thirty.

To confirm the superiority of our algorithm in terms of the total number of iterations, we compute, for each KF the percentage of cases where the KF gives the best iteration number. As an illustration, we plot a histogram that we use as a statistical tool to compare the performance of the algorithms.

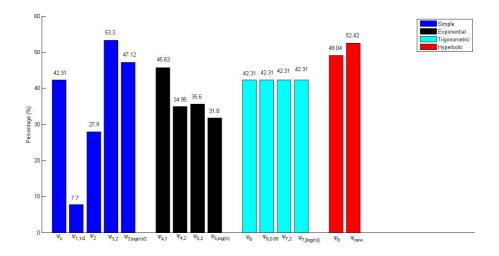


Fig. 1. Performance comparison between KFs in Table 1.

Supported by the performance bar graph, We may conclude some remarks:

- For  $\psi_{4,p}$ , the parameter which gives the best theoretical complexity bound has better percentage than p = 2. But the concordance between the theoretical and the numerical results is not always satisfied, as we can see in the case of  $\psi_{3,p}$ ,  $\psi_{5,p}$ and  $\psi_{7,p}$ .
- We can easily see that the algorithms based on hyperbolic KFs attain the most wins, on average, among the considered algorithms based on other types.
- Although the theoretical complexity obtained for  $\psi_8$  is better than the one of  $\psi_{new}$ , but numerical tests reveal that  $\psi_{new}$  has better percentage. In fact, a thorough analysis shows that in the examples with fixed size, it's  $\psi_{new}$  that gives the smallest iteration number with a slackness which can amount up to 24 iterations. As for examples with variable size for  $\theta = 0.7$ ,  $\psi_{new}$  performs better than  $\psi_8$  in Examples 5, 6 and 8 while in Examples 7,  $\psi_8$  doesn't have the ability to complete the run successfully. In contrast, for the values 0.9 and 0.99 of  $\theta$ ,  $\psi_8$  meets or exceeds  $\psi_{new}$ with a slackness of no more than 7 iterations.

The numerical effectiveness of  $\psi_{\text{new}}$  comparing with all KFs with logarithmic barrier term motivates us to propose a generalization of this function in order to achieve a better theoretical complexity.

#### 4. PARAMETRIZED VERSION

In the previous section, we performed a well-detailed complexity analysis for  $\psi_{\text{new}}$ . Thus, we discuss more briefly the analysis for the generalized KF  $\psi_{\text{new},p}$ .

$$\psi_{\text{new},p}(t) = \frac{t^2 - 1}{2} + \psi_{b,p}(t), \qquad (19)$$

with 
$$\psi_{b,p}(t) = a \Big( \coth^p(t) - \log t - \coth^p(1) \Big), a = \frac{\sinh^2(1)}{\sinh^2(1) + p \coth^{p-1}(1)}$$
 and  $p \ge 2$ .

**Remark 4.1.** This function can also be considered as a generalization, up to the multiplicative constant 1/a, of the KF  $\psi_8$  introduced in [51] (see Table 1).

Using simple calculations, we can easily prove that  $\psi_{\text{new},p}$  is indeed a KF and we have:

$\psi_{\text{new},p}'(t) = t - a \left(\frac{p \coth^{p-1}(t)}{\sinh^2(t)} + \frac{1}{t}\right)$
$\psi_{\text{new},p}''(t) = 1 + a \left( \frac{p(p-1) \coth^{p-2}(t)}{\sinh^4(t)} + \frac{2p \coth^p(t)}{\sinh^2(t)} + \frac{1}{t^2} \right) > 1$
$\psi_{\text{new},p}^{\prime\prime\prime}(t) = -a \Big( \frac{p(p-1)(p-2) \operatorname{coth}^{p-3}(t)}{\sinh^6(t)} + \frac{(6p^2 - 4p) \operatorname{coth}^{p-1}(t)}{\sinh^4(t)} + \frac{4p \operatorname{coth}^{p+1}(t)}{\sinh^2(t)} + \frac{2}{t^3} \Big) < 0$
$t\psi_{\text{new},p}'' - \psi_{\text{new},p}' = a\Big(\frac{p(p-1)t\coth^{p-2}(t)}{\sinh^4(t)} + \frac{2pt\coth^p(t)}{\sinh^2(t)} + \frac{2}{t} + \frac{p\coth^{p-1}(t)}{\sinh^2(t)}\Big)$
$t\psi_{\text{new},p}'' + \psi_{\text{new},p}' = 2t + a\left(\frac{p(p-1)t \coth^{p-2}(t)}{\sinh^4(t)} + \frac{p \coth^{p-1}(t)}{\sinh^2(t)}\left(2t \coth(t) - 1\right)\right)$

**Tab. 7.** The first three derivatives of  $\psi_{\text{new},p}$  with  $t\psi''_{\text{new},p} \pm \psi'_{\text{new},p}$ .

Thus,  $\psi_{\text{new},p}$  satisfies all the results of Lemma 3.1, Corollary 3.2, Lemma 3.4 and Lemma 3.7. Whereas, the equivalent of Lemma 3.5 for  $\psi_{\text{new},p}$  is

**Lemma 4.2.** For all  $(z,t) \in [0,+\infty) \times (0,1]$  such that  $z = -\frac{1}{2}\psi'_{\text{new},p}(t)$ , we have

$$\operatorname{coth}(t) \le \left( (\sinh^2(1) + \coth^{p-1}(1)) \operatorname{coth}^2(1) \right)^{\frac{1}{p+1}} (2z+1)^{\frac{1}{p+1}}.$$

# 4.1. Complexity analysis

The complexity analysis for the parameterized KF  $\psi_{\text{new},p}$  proceeds in the same way as in the previous section for  $\psi_{\text{new}}$ . To avoid repetition, we do not present all the details of the computations. We only present the outcome of each step of our computational scheme in Table 8.

	$f(\bar{\alpha}) \le -\frac{\sqrt{2}\Psi(v)^{\frac{p}{2(p+1)}}}{72\left(1+2\sinh^2(1)\coth^5(1)(p+5)\right)}$	$\Psi(v_+) \le \frac{1}{2}$	$\frac{\coth(1)(p+2)}{(1-\theta)} \left(\theta\sqrt{n} + \right.$	$\sqrt{2\tau}\Big)^2 := \Psi_0$
β =	$= \frac{\sqrt{2}}{72\left(1+2\sinh^2(1)\coth^5(1)(p+5)\right)} \text{ and } \gamma = \frac{p+2}{2(p+1)}$	$K \leq$	$\frac{72\sqrt{2}(5+4p)(p+1)}{(p+2)}$	$\Psi_0^{\frac{p+2}{2(p+1)}}$

Tab. 8. Outcomes of the computational scheme.

For small-update method with  $\tau = \mathcal{O}(1)$  and  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$ , the complexity of the primal-dual interior-point algorithm for LP problems based on the new parametric KF

is  $\mathcal{O}(p^2 \sqrt{n} \log \frac{n}{\epsilon})$  iterations complexity.

As for large-update method i. e.,  $\tau = \mathcal{O}(n)$  and  $\theta = \Theta(1)$ , the substitution of these values into

$$K \le \left[\frac{72\sqrt{2}(5+4p)(p+1)}{(p+2)}\right] \Psi_0^{\frac{p+2}{2(p+1)}}$$

does not give the best possible bound. A better bound is obtained using the following lemma

**Lemma 4.3.** (El Ghami et al. [21]) Let  $0 \le \theta < 1$ ,  $v_+ = \frac{v}{\sqrt{1-\theta}}$ . If  $\Psi(v) \le \tau$ , then we have

$$\Psi(v_+) \le \frac{\theta n + 2\tau + 2\sqrt{2\tau n}}{2(1-\theta)} := \Psi_0.$$

 $\Psi_0$  is an upper bound for  $\Psi(v_+)$  during the process of the algorithm.

Proof. For simplicity, we denote

$$\psi(t) := \psi_{\text{new},p}(t).$$

We can easily verify that  $\psi_{b,p}(1) = 0$ ,  $\psi'_{b,p}(1) = -1$ , and  $\psi_{b,p}$  is monotonically decreasing on  $\mathbb{R}_{++}$ . Hence, we have

$$\psi(t) \le \frac{t^2 - 1}{2}, \ \forall t \ge 1.$$

Putting  $t = \frac{\varrho(\frac{\tau}{n})}{\sqrt{1-\theta}} \ge 1$  and using (15), we get

$$\Psi(v_{+}) \leq \frac{n}{2} \left( \frac{\varrho\left(\frac{\tau}{n}\right)^{2}}{1-\theta} - 1 \right).$$

Moreover, using Lemma 3.4, we obtain

$$\Psi(v_{+}) \leq \frac{n}{2} \left[ \frac{\left(1 + \sqrt{\frac{2\tau}{n}}\right)^2}{1 - \theta} - 1 \right]$$
$$= \frac{\theta n + 2\tau + 2\sqrt{2\tau n}}{2(1 - \theta)}.$$

Using the upper bound defined in Lemma 4.3, we get  $\mathcal{O}\left((p+1)n^{\frac{p+2}{2(p+1)}}\log\frac{n}{\epsilon}\right)$  iterations complexity for large-update methods. This expression is minimal at  $p = \frac{\log n}{2} - 1$  and then is equal to  $\mathcal{O}\left(\sqrt{n}\log n\log\frac{n}{\epsilon}\right)$ .

# 4.2. Numerical tests

Now, we would like to investigate the influence of parameterizing the KF (7) on the computational behavior of the generic primal-dual algorithm for LP presented in Algorithm 1. For this purpose, we present some numerical simulations for implementing the algorithm based on the new parametric KF (8) on the aforementioned set of problems with the same values of the parameters  $\theta, \mu$  and  $\tau$  and for different values of the parameter p. We also compare with the KF that gave the best performance in the previous section  $\psi_{3,2}$  and with  $\psi_8$  which is of the same hyperbolic-logarithmic type. The obtained results are listed in tables below.

Ex	$\theta$	$\psi_{\text{new}}$	$\psi_{\text{new},2}$	$\psi_{\text{new},3}$	$\psi_{\text{new},4}$	$\psi_{\mathrm{new},\frac{\log n}{2}-1}$	$\psi_{3,2}$	$\psi_8$
	0.1	188	188	188	188	$18\tilde{8}$	188	200
ole	0.3	<b>56</b>	56	56	56	<b>56</b>	<b>56</b>	70
lm	0.5	29	29	29	29	29	<b>29</b>	44
Example	0.7	17	17	17	17	17	17	23
щ	0.9	11	11	10	12	24	11	14
5	0.1	191	191	191	191	191	191	215
ole	0.3	57	57	57	57	57	57	75
Example	0.5	29	29	29	29	<b>29</b>	<b>29</b>	39
Jxa	0.7	17	17	17	17	17	17	21
щ	0.9	10	9	9	9	11	10	9
3	0.1	192	192	192	192	192	192	204
ole	0.3	57	57	57	57	57	57	66
Example	0.5	30	30	30	30	30	30	33
Jxa	0.7	18	19	18	19	17	18	20
щ	0.9	21	19	19	22	33	28	25
4	0.1	196	196	196	196	196	196	212
ole	0.3	<b>58</b>	58	58	<b>58</b>	58	<b>58</b>	78
Example	0.5	33	30	30	30	31	30	42
Dxa	0.7	21	20	22	24	28	21	24
Щ	0.9	21	18	22	20	24	21	20

Tab. 9. Number of inner iterations for fixed size examples.

$\theta = 0.99$	$\theta = 0.9$	$\theta = 0.7$	θ
5 25 200 200 400 1000	$5 \\ 25 \\ 50 \\ 100 \\ 200 \\ 400 \\ 1000$	-	m
28 28 30 30 30	14 15 15 17 17 17 17 18	18     19     20     20     21     21     21     22	$\psi_{ m new}$
24 24 27 27 27 27 27 27 27	14 15 16 16 16 16 16	18 19 20 20 20 21 21 21 21 22	$\psi_{ m new,2}$
29 29 31 31 31 31 31	16 17 17 18 18 18 18 18	21 22 23 23 23 24 24 24 24 25	] $\psi_{ m new,3}$
30 32 32 32 32 32	16 17 17 18 18 18 18 18	22 23 24 24 24 25 25 25 26	Exampl $\psi_{\text{new},4}$
<b>21</b> 28 31 27 25 31	22 15 16 15 16 15	18 19 20 20 21 21 21 21	e 5 $\psi_{\text{new},\frac{\log n}{2}-1}$
29 29 31 31 31 31 31	$14 \\ 15 \\ 15 \\ 16 \\ 16 \\ 16 \\ 16 \\ 18 \\ 18$	18 19 20 20 20 20 21 21 21 21	$\psi_{3,2}$
21 21 23 23 23 23 23 23	12     13     13     14     14     14     14     15     15     15     1	35363737404041	$\psi_8$
10 13 13 13 13 13	10     11     11     11     12     1	18     19     20     20     21     21     22     22     22	$\psi_{ m new}$
10     10     12     1	<b>9</b> 111 112 12 12 12 12 13	18     19     20     20     21     21     22     22     2	$\psi_{ m new,2}$
$   \begin{array}{c}     10 \\     12 \\     12 \\     12 \\     12 \\     12 \\     12 \\     12 \\     12 \\   \end{array} $	9 10 10 11 11 11 11	18     19     20     20     21     21     22     22	] $\psi_{ m new,3}$
10 10 12 12 12 12 12 12	9 10 10 11 11 11 11 11	18     19     20     20     21     21     22     22     22	Example $\psi_{\text{new},4}$
$   \begin{array}{c}     11 \\     10 \\     12 \\     12 \\     12 \\     12 \\     12 \\     12 \\     12 \\   \end{array} $	11 11 11 12 12 12 12 12	18     19     20     20     21     21     22     22	le 6 $\psi_{\text{new},\frac{\log n}{2}-1}$
10 10 13 13 13 13 13 13	10     11     11     11     12     12     12     12     12     14     14	18     19     20     20     21     21     22     22     2	$\psi_{3,2}$
10 10 12 12 12 12 12 12 12	9 10 10 11 11 11 11 11	28 29 32 32 33 33 33 34	$\psi_8$

**Tab. 10.** Number of inner iterations for Example 5 and 6 with different sizes n = 2m.

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	$\psi_8$	28	31	32	32	33	33	36	12	13	13	14	14	14	15	19	19	21	21	21	21	21	
	$\psi_{3,2}$	18	19	20	20	<b>21</b>	<b>21</b>	22	12	<b>13</b>	<b>13</b>	14	14	14	15	26	26	28	28	28	28	28	
8 0	$\psi_{\mathrm{new},\frac{\log n}{2}-1}$	18	19	20	20	<b>21</b>	<b>21</b>	22	12	13	13	14	14	15	16	21	26	27	25	23	24	82	
Example	$\psi_{\rm new,4}$	18	19	20	20	21	21	22	12	13	13	14	14	15	16	22	22	24	24	24	24	24	
I	$\psi_{\mathrm{new,3}}$	18	19	20	20	21	21	22	12	13	13	14	14	15	16	24	24	26	26	26	26	26	
	$\psi_{\mathrm{new,2}}$	18	19	20	20	21	<b>21</b>	22	12	13	13	14	14	14	15	21	21	23	23	23	23	23	
	$\psi_{\rm new}$	18	19	20	20	21	21	22	12	13	13	14	14	14	15	23	23	26	26	26	26	26	
	$\psi_8$	I	I	I	I		I	I	6	10	10	11	11	11	12	10	10	12	12	12	12	12	
	$\psi_{3,2}$	18	19	20	<b>20</b>	<b>21</b>	<b>21</b>	22	10	11	11	12	12	12	14	10	10	13	13	13	13	13	
e 7	$\psi_{\mathrm{new}, rac{\log n}{2} - 1}$	18	19	20	20	<b>21</b>	21	22	11	11	11	12	12	12	13	11	10	12	12	12	12	12	
Example 7	$\psi_{\rm new,4}$	18	19	20	20	21	21	22	6	10	10	11	11	11	12	10	10	12	12	12	12	12	
I	$\psi_{\mathrm{new,3}}$	18	19	20	20	21	21	22	6	10	10	11	11	11	12	10	10	12	12	12	12	12	
	$\psi_{\mathrm{new},2}$	18	19	20	<b>20</b>	21	<b>21</b>	22	6	11	11	12	12	12	13	10	10	12	12	12	12	12	
	$\psi_{\mathrm{new}}$	18	19	20	20	21	21	22	10	11	11	12	12	12	14	10	10	13	13	13	13	13	
		5	25	50	100	200	400	1000	ъ	25	50	100	200	400	1000	5	25	50	100	200	400	1000	
U U	2			7.0	) =	θ					6.0	) =	θ					66	0 =	θ			

**Tab. 11.** Number of inner iterations for Example 7 and 8 with different sizes n = 2m.

$\psi_i$	$\psi_{ m new}$	$\psi_{ m new,2}$	$\psi_{ m new,3}$	$\psi_{ m new,4}$	$\psi_{\mathrm{new},\frac{\log n}{2}-1}$	$\psi_{3,2}$	$\psi_8$
%	50.48	66.99	68.93	66.99	58.25	51.45	55.33

**Tab. 12.** The percentage of cases where the KF gives the bestiteration number.

# 4.2.1. Comments

By comparing the results in Tables 9-12, we notice that:

- The number of iterations clearly depends on the value of the parameter p as the gap between two different values of p can amount up to 59. It should also be noted that  $\psi_{\text{new},3}$  significantly reduces the number of iterations although  $\psi_{\text{new},\frac{\log n}{2}-1}$  has the best complexity bound theoretically.
- Comparing with  $\psi_{\text{new}}$ , the KFs  $\psi_{\text{new},2}$ ,  $\psi_{\text{new},3}$ ,  $\psi_{\text{new},4}$  and  $\psi_{\text{new},\frac{\log n}{2}-1}$  were able to produce even better iteration numbers especially in Examples 6 and 7 for  $\theta = 0.9$  and  $\theta = 0.99$ , while maintaining similar performance for  $\theta = 0.7$  in Examples 6, 7 and 8.
- $\psi_{\text{new},p}$ , for all tested parameters, outperformed  $\psi_{3,2}$  which had the best performance in the previous section. This confirms that the parametrization has effected the number of iterations of the algorithm positively.
- $\psi_{\text{new},2}$  far outperformed  $\psi_8$  despite having the same complexity bounds and more than, the same expression up to a multiplicative constant (see Remark 4.1).

# 5. CONCLUSIONS AND REMARKS

In this paper, we investigate an interior-point algorithm for solving LP problems based on a new KF which is a combination of the classic KF and a hyperbolic barrier term. The complexity analysis for large-update primal-dual IPMs based on this KF yields an  $\mathcal{O}\left(n^{\frac{3}{4}}\log\frac{n}{\epsilon}\right)$  iteration bound which improves the classical iteration bound with a factor  $n^{\frac{1}{4}}$ . For small-update methods, the proposed algorithm enjoys the favourable iteration bound. To achieve the best iteration complexity, we introduce a class of KFs by parametrizing the new hyperbolic-logarithmic KF. The simulation results illustrate that the proposed algorithm is efficient and robust.

As a further research we would like to extend our approach to more general optimization problems. Another question of interest is whether the new KF can be used to design an efficient infeasible interior-point algorithm as in [27, 35, 47].

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