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INTERIOR POINT METHODS AND KERNEL FUNCTIONS  
OF A LINEAR PROGRAMMING PROBLEM

by

LATRICEE Y. TANKSLEY

(Under the Direction of Goran Lesaja)

ABSTRACT

In this thesis the Interior – Point Method (IPM) for Linear Programming problem (LP) that is based on the generic kernel function is considered.

The complexity (in terms of iteration bounds) of the algorithm is first analyzed for a class of kernel functions defined by (3-1). This class is fairly general; it includes classical logarithmic kernel function, prototype self-regular kernel function as well as non-self-regular functions, thus it serves as a unifying frame for the analysis of IPM. Historically, most results in the theory of IPM are based on logarithmic kernel functions while other two classes are more recent. They were considered with the intention to improve theoretical and practical performance of IPMs. The complexity results that are obtained match the best known complexity results for these methods.

Next, the analysis of the IPM was summarized and performed for three more kernel functions. For two of them we again matched the best known complexity results.

The theoretical concepts of IPM were illustrated by basic implementation for the classical logarithmic kernel function and for the parametric kernel function both described in (3-1). Even this basic implementation shows potential for a good performance. Better implementation and more numerical testing would be necessary to draw more definite

conclusions. However, that was not the goal of the thesis, the goal was to show that IPM with kernel functions different than classical logarithmic kernel function can have best known theoretical complexity.

INDEX WORDS: Interior Point Methods, Linear Programming, Kernel Functions, Log Barrier Functions

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OF A LINEAR PROGRAMMING PROBLEM

by

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INTERIOR POINT METHODS AND KERNEL FUNCTIONS  
OF A LINEAR PROGRAMMING PROBLEM

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## NOMENCLATURE

$R^n$	Euclidean n dim space
$R^n_+$	All vectors of $R^n$ with nonnegative components
$x \in A$	Element $x$ belongs to set $A$
$\ x\ $	A Euclidian norm of vector $x \in R^n$ , $\ x\  = \sqrt{\sum_{i=1}^n x_i^2}$
$\mu \rightarrow 0$ ,	$\mu$ converges to 0
$f : R^n \rightarrow R$	A function with n variables
$\nabla f, \nabla^2 f$	A gradient and a Hessian of $f$
$F : R^n \rightarrow R^m$	A vector valued function
$\nabla F$	A Jacobian of $F$
$X = \text{diag}(x)$	A diagonal matrix that has components of the vector $x$ on the main diagonal and zeros everywhere else
$xs, \frac{x}{s}, x^{-1}$	Component wise operations (product, division, inverse) of vectors $x, s \in R^n$ . For example, $xs = (x_1s_1, \dots, x_ns_n)$ .
$o(f(n))$	There exist a constant $C$ and a function $g(n)$ such that $f(n) \geq Cg(n)$ ("small o" notation).

$O(f(n))$ 

There exist a constant  $C$  and a function  $g(n)$  such that

$$f(n) \leq Cg(n)$$

 $\Omega(f(n))$ 

There exist constants  $C, K$  and a function  $g(n)$  such that

$$Cg(n) \leq f(n) \leq Kg(n) \text{ ("big } \Omega \text{" notation).}$$

## CHAPTER 1

### INTRODUCTION

#### Linear Programming Model

The mathematical model of linear programming is useful in solving a wide range of problems in industry, business, science and government. This is by far the most used optimization model.

These problems and their linear programming models can often be complex as they consist of a huge number of variables and constraints ranging up to the hundred thousands. A linear programming model consists of an objective function, that is a linear function, and the constraints on that function that are also linear. Linear programming involves the planning of activities to obtain a result that reaches a specified goal among all feasible alternatives.

A Linear Program (LP) is a problem that can be expressed in standard form as follows:

$$\begin{array}{ll}
 \text{Minimum} & c^T x \\
 \text{subject to} & Ax = b \\
 & x \geq 0
 \end{array} \tag{1-1}$$

where  $x \in R^n$  is the vector of variables to be solved for, and matrix  $A \in R^{m \times n}$  and vectors  $c \in R^n$ ,  $b \in R^m$  are input data. . The linear function  $Z = c^T x$  is called the objective function, and the equations  $Ax = b$  are called functional constraints, while  $x \geq 0$  are called nonnegativity constraints. The set  $F = \{x \in R^n \mid Ax = b, x \geq 0\}$  is called a feasible set. Geometrically, the set represents a polyhedron in  $R^n$ .

Many practical problems can be modeled as LP models. To illustrate this fact we list the following simple example taken from the [Hillier, Lieberman, 2005].

**Example:** The Wyndor Glass Co. produces high-quality glass products, including windows and glass doors. It has three plants. Aluminum frames and hardware are made in Plant 1, wood frames are made in Plant 2, and Plant 3 produces the glass and assembles the products. Because of declining earnings, top management has decided to revamp the company's product line. Unprofitable products are being discontinued, releasing production capacity to launch two new products having large sales potential:

Product 1: An 8-foot glass door with aluminum framing

Product 2: A 4 x 6 foot double-hung wood-framed window

Product 1 requires some of the production capacity in Plants 1 and 3, but none in Plant 2.

Product 2 needs only Plants 2 and 3. The marketing division has concluded that the company could sell as much of either product as could be produced by these plants. However, because both products would be competing for the same production capacity in Plant 3, it is not clear which mix of these two products would be most profitable. Each product will be produced in batches of 20, so the production rate is defined as the number of batches produced per week.

Any combination of production rates that satisfies the restrictions is permitted, including producing none of one product and as much as possible of the other. Profit from selling one batch of Product 1 (glass doors) is \$3000 and profit from selling one batch of Product 2 (windows) is \$5000. We assume that all produced batches will be sold. Goal: To determine what the production rates should be for the two products in order to maximize the total profit, subject to the restrictions imposed by the limited production capacities available in the three plants.

Of course, this is a simplified real world situation but good enough to illustrate the usefulness of the model. Formulation of the Linear Programming (LP) Problem

Let  $x_1$  = number of batches of product 1 produced per week

$x_2$  = number of batches of product 2 produced per week

$Z$  = total profit per week in thousands of dollars from producing these two products

The following table summarizes the data gathered:

Wyndor Glass Company Data

Plant	Production Time Per Batch Hours		Production Time Available per Week, Hours
	Product		
	1	2	
1	1	0	4
2	0	2	12
3	3	2	18
Profit per Batch	3000	5000	

Table 1

The objective function is  $Z = 3x_1 + 5x_2$  and it represents a total profit measured in thousands of dollars. The objective function is subject to the restrictions imposed by the limited production capacities available in each of the plants, and they can be mathematically expressed by the following inequalities:

$$\begin{array}{ll}
 \text{Plant 1} & x_1 \leq 4 \\
 \text{Plant 2} & 2x_2 \leq 12 \\
 \text{Plant 3} & 3x_1 + 2x_2 \leq 18
 \end{array}$$

Thus, the overall linear programming model illustrating the Wyndor Glass Company is

$$\begin{array}{ll}
 \text{Maximize} & Z = 3x_1 + 5x_2 \\
 \\
 \text{subject to} & x_1 \leq 4 \\
 & 2x_2 \leq 12 \\
 & 3x_1 + 2x_2 \leq 18 \\
 & x_1 \geq 0, x_2 \geq 0
 \end{array}$$

By adding slack variables this problem can be transformed in the standard form (1-1).

$$\begin{array}{ll}
 \text{Maximize} & Z = 3x_1 + 5x_2 \\
 \text{subject to} & x_1 + x_3 = 4 \\
 & 2x_2 + x_4 = 12 \\
 & 3x_1 + 2x_2 + x_5 = 18 \\
 & x_1 \geq 0, x_2 \geq 0, x_3 \geq 0, x_4 \geq 0, x_5 \geq 0
 \end{array}$$

The similar procedure can be done for different inequality formulations of LP.

This example illustrates the applicability of the LP model. The number of problems that can be modeled as LP is huge and widespread to many areas of science, industry, business, finance, government, etc. For more examples see [HL] and other Operations Research textbooks. Therefore, the efficient methods to solve LP models are very important. In the following sections we will outline main methods that are used to solve LP models.

### Methods to Solve LP Models

The first successful general procedure, **Simplex Method**, for solving a LP problem was discovered by George Dantzig in 1947 although there were partial results discovered earlier. Theoretically the main idea of the simplex method (SM) is that it travels from vertex to vertex



on the boundary of the feasible region, repeatedly increasing or decreasing the objective function until either an optimal solution is found, or it is established that no solution exists. The number of iterations required in the worst case is an exponential function of the number of variables, as it was first discovered by Klee and Minti in 1972. However, the worst case behavior has not been observed in practice. On the contrary, the algorithm works very well in practice, typically requiring  $O(n)$  iterations. Highly sophisticated implementations are available (CPLEX, MOSEK, LINDO, EXCEL SOLVER) and have excellent codes for simplex algorithms. These codes are capable of solving huge problems with millions of variables and thousands of constraints. This discrepancy between exponential worst case complexity and good practical behavior of simplex method prompted the research in two directions. One direction was a search for the algorithm with the polynomial worst case complexity and the other direction was the analysis of average complexity of simplex method.

In 1979, Leonid Khachiyan showed that the **Ellipsoid Method**, created by A. Nemirovski and D. Yudin for nonlinear programming problems, solves any linear program in a number of steps which is a polynomial function of the amount of data defining the linear program. Unfortunately, in practice, the simplex method turned out to be far superior to the ellipsoid method. However, theoretical importance is significant because it provided a basis to prove that polynomial methods exist for many combinatorial problems.

In 1982 K. Borgward provided the first probabilistic analysis of the simplex method, showing that the expected number of iterations is polynomially bounded. Soon afterwards, other authors provided similar analysis. A relatively simple and complete analysis was provided by Adler and Megiddo in 1985. Using the clever probability model, they showed that

upper and lower bounds on an average number of iterations is a function of  $\Omega((\min\{m, n\})^2)$ , where  $m$  is the number of constraints and  $n$  is the number of variables.

In 1984, Narendra Karmarkar introduced an **Interior-Point Method** (IPM) for linear programming, combining the desirable theoretical properties of the ellipsoid method and practical advantages of the simplex method. Its success initiated an explosion in the development of interior-point methods that continue to this day.

These methods do not pass from vertex to vertex along the edges of the feasible region, which is the main feature of the simplex algorithm; they follow the central path in the interior of the feasible region. Though this property is easy to state, the analysis of interior-point methods is a subtle subject which is much less easily understood than the behavior of the simplex method. Interior-point methods are now generally considered competitive with the simplex method in most, though not all, applications, and sophisticated software packages implementing them are now available (CPLEX, MOSEK, LINDO, EXCEL SOLVER).

## CHAPTER 2

### INTERIOR POINT METHODS BASED ON KERNEL FUNCTIONS

The linear optimization problem in standard form is

$$(P) \quad \min \{c^T x : Ax = b, x \geq 0\},$$

where  $A \in R^{m \times n}$  ( $\text{rank}(A) = m$ ),  $b \in R^m$ ,  $c \in R^n$ . The dual problem of (P) is

$$(D) \quad \max \{b^T y : A^T y + s = c, s \geq 0\},$$

where  $s \in R^n$  is a dual slack variable.

We can assume that the Interior Point Condition (IPC) is satisfied without loss of generality; that is, there exists a point  $(x^0, s^0, y^0)$  such that  $Ax^0 = b$ ,  $x^0 > 0$  and  $A^T y^0 + s^0 = c$ ,  $s^0 > 0$ , which means that the interiors of the feasible regions of the primal(P) and dual(D), are not empty. If the problem doesn't satisfy the IPC, it can be modified so that it does and even in such a way that  $x^0 = s^0 = e$ , where  $e$  denotes a vector of all ones. The details can be found in [Roos, C *et. al.*, 1997].

Optimality conditions for (P) and (D) yield the following system of equations:

$$\begin{aligned} Ax &= b, x \geq 0 \\ A^T y + s &= c, s \geq 0 \\ xs &= 0, \end{aligned} \tag{2-1}$$

where the vector  $xs$  denotes the component wise product of vectors  $x$  and  $s$  which is also called Hadamard product.

The theory of interior point methods (IPMs) that is based on the use of Newton's Method suggests that the third equation in (2-1) has to be perturbed. The third equation is often called the **complementarity condition** for the primal and dual, and is replaced by  $xs = \mu e$ , where  $\mu$  is a positive parameter. The optimality conditions (2-1) are transformed to the following system:

$$\begin{aligned} Ax &= b, x > 0 \\ A^T y + S &= c, s > 0 \\ xs &= \mu e. \end{aligned} \tag{2-2}$$

Since  $\text{rank}(A) = m$ , this system has unique solution for each  $\mu > 0$ . We can write this solution as  $(x(\mu), y(\mu), s(\mu))$ , calling  $x(\mu)$  the  $\mu$ -center of (P) and  $(y(\mu), s(\mu))$  the  $\mu$ -center of (D). The set of all  $\mu$ -centers forms a homotopy path in the interior of the feasible region that is called the **central path**.

The main property of the central path can be summarized as follows: if  $\mu \rightarrow 0$ , then the limit of the central path exists, the limit points satisfy the complementarity condition, and the limit yields optimal solutions for (P) and (D). This limiting property of the central path leads to the main idea of the iterative methods for solving (P) and (D): Trace the central path while reducing  $\mu$  at each iteration. However, tracing the central path exactly would be too costly and inefficient. One of the main achievements of interior point methods was to show that tracing the central path approximately while still maintaining good properties of the algorithms is sufficient.

Tracing the central path means solving the system (2-2) using Newton Method on the function

$$F(x, y, s) = \begin{bmatrix} Ax - b \\ A^T y + s - c \\ xs - \mu e \end{bmatrix} = 0. \quad (2-3)$$

Tracing the central path approximately means that only one, or at most, a couple of iterations of a modified (damped) Newton's Method will be performed for a particular  $\mu$ . One iteration of a Newton Method for the function (2-3) and particular  $\mu$  is stated below.

$$\nabla F \begin{bmatrix} \Delta x \\ \Delta y \\ \Delta z \end{bmatrix} = -F(x, y, s) \quad (2-4)$$

where

$$\nabla F = \begin{bmatrix} A & 0 & 0 \\ 0 & A^T & I \\ S & 0 & X \end{bmatrix}$$

denotes the Jacobian of  $F$  and  $[\Delta x, \Delta y, \Delta z]^T$  is a Newton's search direction that we want to calculate.

Solving (2-4) reduces to solving this system.

$$\begin{aligned} A\Delta x &= 0 \\ A^T \Delta y + \Delta s &= 0 \\ s\Delta x + x\Delta s &= \mu e - xs \end{aligned} \quad (2-5)$$

Next we update the current iterates  $(x, y, s)$  by taking an appropriate step along the calculated direction.

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

The step size  $\alpha$  has to be chosen approximately, so that the new iterate is in a certain neighborhood of  $\mu$ -center. The choice of  $\alpha$  will be discussed later in the text.

The idea of the algorithm is illustrated in the Figure 1 blow.

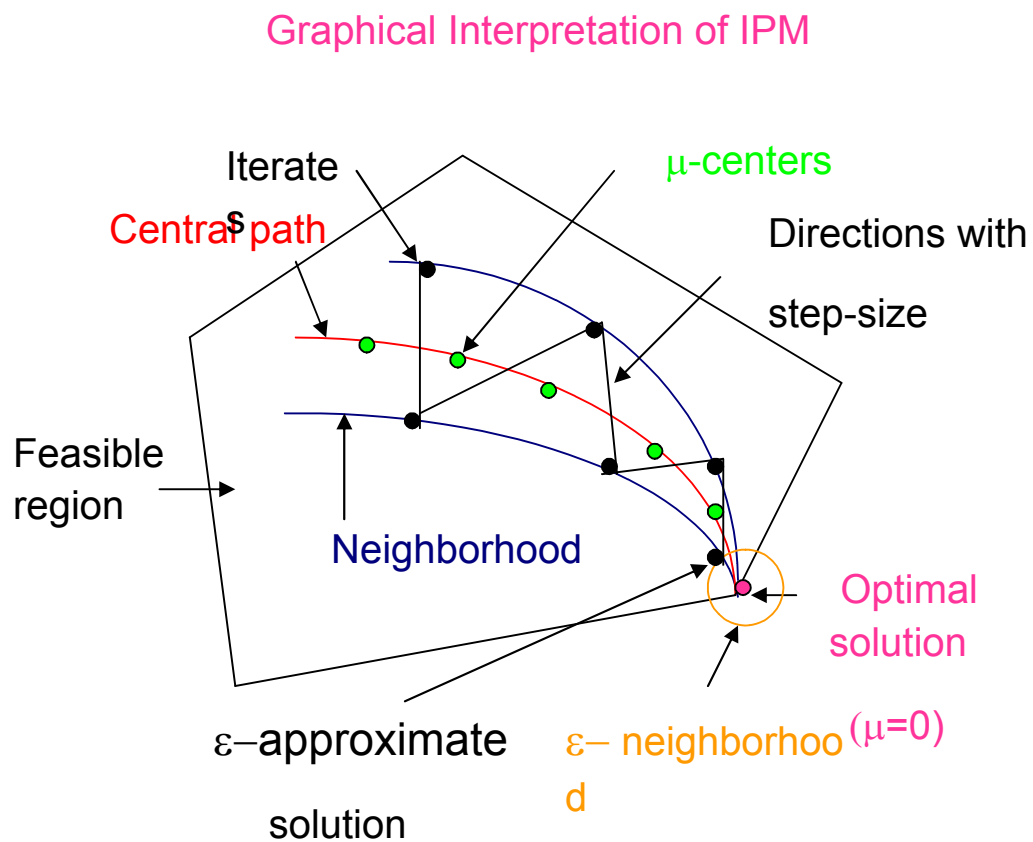


Figure 1

In order to generalize the algorithm outlined above, we introduce a new vector

$$\nu = \sqrt{\frac{x s}{\mu}} \quad (2-6)$$

which we use to define new scaled directions

$$d_x := \frac{\nu \Delta x}{x}, \quad d_s := \frac{\nu \Delta s}{s} \quad (2-7)$$

where the operations in (2-7) are component-wise product and division of vectors. Using the above definitions (2-6) and (2-7), the system (2-5) reduces to

$$\begin{aligned} A d_x &= 0 \\ A \Delta y + d_s &= 0 \\ d_x + d_s &= \nu^{-1} - \nu \end{aligned} \quad (2-8)$$

where  $\bar{A} = \frac{1}{\mu} A V^{-1} X$ ,  $V := \text{diag}(\nu)$ ,  $X = \text{diag}(x)$ . Note that  $X = \text{diag}(x)$  denotes a diagonal matrix that has components of the vector  $x$  on the main diagonal and zeros everywhere else.

The new search direction  $(\Delta x, \Delta y, \Delta s)$  is obtained by first solving the system (2-8). Once  $d_x$  and  $d_s$  are found we apply (2-6) to find  $\Delta x$  and  $\Delta s$ . This direction can also be obtained directly by solving the following system:

$$\begin{aligned} A \Delta x &= 0 \\ A^T \Delta y + \Delta s &= 0 \\ s \Delta x + x \Delta s &= -\mu \nu \nabla \Psi(\nu). \end{aligned} \quad (2-9)$$

This system can be reduced to

$$M\Delta y = r \quad (2-10)$$

where

$$\begin{aligned} M &= AS^{-1}XA^T \\ r &= \mu AS^{-1}v\nabla\psi(v) \end{aligned} \quad (2-11)$$

and  $X = \text{diag}(x)$ ,  $S = \text{diag}(s)$ ,  $V = \text{diag}(v)$ .

Once  $\Delta y$  is found,  $\Delta s$  and  $\Delta x$  are found by back substitutions

$$\Delta s = -A^T \Delta y \quad (2-12)$$

$$\Delta x = -s^{-1}(x\Delta s - \mu v\nabla\Psi(v)) \quad (2-13)$$

where products denote component-wise products of vectors.

The following observation is crucial for the generalization of the method. Observe that  $-(v^{-1} - v)$  is a gradient of the following function

$$\Psi_c(v) = \sum_{i=1}^n \left( \frac{v_i^2 - 1}{2} - \log v_i \right) = \left( \frac{v_1^2 - 1}{2} - \log v_1 \right) + \dots + \left( \frac{v_n^2 - 1}{2} - \log v_n \right). \quad (2-14).$$

This function is called **log-barrier** function.

One may easily verify that the Hessian  $\nabla^2\Psi_c(v) = \text{diag}(e + v^{-2})$ . Since this matrix is positive definite,  $\Psi_c(v)$  is strictly convex. Moreover, since  $\nabla\Psi_c(e) = 0$ , it follows that



$\Psi_c(v)$  attains its minimal value at  $v = e$ , with  $\Psi_c(e) = 0$ . Thus, it follows that  $\Psi_c(v)$  is nonnegative everywhere and vanishes if and only if  $v = e$ , that is, if and only if  $x = x(\mu)$  and  $s = s(\mu)$ . Hence, we see that the  $\mu$ -centers  $x(\mu)$  and  $s(\mu)$  can be characterized as the minimizers of the function  $\Psi_c(v)$ . Therefore, the function  $\Psi_c(v)$  serves as a **proximity measure** to the  $\mu$ -centers (central path.. The norm based proximity measure that is derived from  $\Psi_c(v)$  is defined as

$$\delta = \frac{1}{2} \|\nabla \Psi_c(v)\|. \quad (2-15)$$

Furthermore, the complimentary equation in (2-8) can be written as  $d_x + d_s = -\nabla \Psi_c(v)$ , which is also called the **scaled centering equation**. The importance of the equation arises from the fact that it essentially defines the search directions. Since  $d_x$  and  $d_s$  are orthogonal, we will still have  $d_x d_s = 0$  if and only if  $v = e$ . The same is true for  $\Delta x$  and  $\Delta s$ .

The main idea of the generalization of the method is to replace the log-barrier function (2-14) with some other barrier function that has the same properties as log barrier. The choice of this function will certainly affect the calculation of the search direction, the step size, and with that the rate of the convergence of the method. It is worth examining the classes of barrier functions that may lead to the improved behavior of the algorithm. In what follows, we will consider several such classes.

We will restrict ourselves to the case where the barrier function  $\Psi(v)$  is separable with identical coordinate functions  $\varphi(v_i)$ . Thus,

$$\Psi(v) = \sum_{i=1}^n \varphi(v_i), \quad (2-16)$$

where  $\varphi(t) : [0, \infty)$  is twice differentiable and attains its minimum at  $t = 1$ , with  $\varphi(1) = 0$ . The function  $\varphi(t)$  is called a **kernel function**. The log-barrier function belongs to this class.

The algorithm based on generic kernel function that was outlined above is summarized in the Figure 2 below. In principle, each barrier function gives rise to a different primal-dual algorithm. The parameters  $\tau, \theta$  and the step size  $\alpha$  in the algorithm should be tuned in such a way that the number of iterations required by the algorithm is as small as possible. The resulting iteration bound will depend on the kernel function, and our main task becomes to find a kernel functions that give a good and possibly best known iteration bound. The question of finding the kernel function that minimizes the iteration bound is still an open question.

**Generic Primal-Dual Interior-Point Algorithm for Linear Optimization**

Input:

An input data  $A, b, c$

A threshold parameter  $\tau \geq 1$ ;

An accuracy parameter  $\varepsilon > 0$ ;

A fixed barrier update parameter  $\theta < 1$ ;

Iteration:

begin

$x := e; s := e; \mu := 1$ ;

while  $n\mu > \varepsilon$  do

begin

calculate  $\mu := \mu(1 - \theta)$ ;

calculate  $\nu := \sqrt{\frac{xs}{\mu}}$ ;

while  $\Psi(\nu) > \tau$  do

begin

calculate the direction  $(\Delta x, \Delta y, \Delta s)$

using (2-10) – (2-13) :

calculate step size  $\alpha$  ;

update

$x := x + \alpha \Delta x, s := s + \alpha \Delta s, y := y + \alpha \Delta y$ ;

end

end

end

Figure 2 Generic IPM

## CHAPTER 3

### ANALYSIS OF THE ALGORITHM FOR A CLASS OF KERNEL FUNCTIONS

The following class of kernel functions will be used to analyze the algorithm, Generic IPM, discussed in the Chapter 2, Figure 2 .

$$\psi_{p,q}(t) = \begin{cases} \frac{t^{p+1} - 1}{p+1} + \frac{t^{1-q} - 1}{q-1}, & t > 0, p \in [0.1], q > 1 \\ \frac{t^{p+1} - 1}{p+1} - \log t, & t > 0, p \in [0.1] \end{cases} \quad (3-1)$$

where  $p$  is a growth parameter and  $q$  is a barrier parameter.

Notice that  $\frac{t^{1-q} - 1}{q-1} \rightarrow -\log t$  when  $q \rightarrow 1$ . This class of kernel functions is fairly general. As

we just explained, it includes log-kernel function as a special case. It also includes so-called self-regular kernel functions when  $p = 1$ . These functions have been extensively discussed in recent literature (Peng, J, *et. al*, 2002). Moreover, it also includes non self-regular functions when  $0 \leq p < 1$ . This class of functions was first discussed in (Lesaja G., *et.al.*, 2008). The results in this chapter follow the results presented in that paper. However, the proofs of several results that were omitted in the paper are outlined here.

#### Properties of Kernel Functions

The derivatives of  $\psi(t)$  in (3-1) play a crucial role in our analysis. Thus, we write down the first three derivatives:

$$\begin{aligned}
\psi'(t) &= t^p - t^{-q} \\
\psi''(t) &= pt^{p-1} + qt^{-q-1} \\
\psi'''(t) &= p(p-1)t^{p-2} - q(q+1)t^{-q-2}.
\end{aligned} \tag{3-2}$$

In the next several lemmas we will describe certain properties of kernel function and its derivatives and their relationships in terms of inequalities. These results will be used in the analysis of the Generic IPM.

The following lemma states the so-called exponential convexity of kernel function which is crucial in proving the polynomial complexity of the Generic IPM.

**Lemma 3.1** *If  $t_1 > 0$  and  $t_2 > 0$ , then  $\psi(\sqrt{t_1 t_2}) \leq \frac{1}{2}(\psi(t_1) + \psi(t_2))$ .*

Proof: It can be shown that the inequality in the lemma holds if and only if

$t\psi''(t) + \psi'(t) \geq 0$  for all  $t > 0$ . This result is beyond the scope of the thesis and can be found (Peng, J., et. al., 2002). Using (3-1) one can easily verify that

$$t\psi''(t) + \psi'(t) = t\left(pt^{p-1} + \frac{q}{t^{q+1}}\right) + t^p - \frac{1}{t^q} = (p+1)t^p + \frac{(q-1)}{t^q} > 0,$$

which completes the proof.

**Lemma 3.2** *If  $t \geq 1$ , then  $\frac{\psi'(t)}{2}(t-1) \leq \psi(t) \leq \frac{p+q}{2}(t-1)^2$ .*

Proof: If  $f(t) = 2\psi(t) - (t-1)\psi'(t)$ , then  $f'(t) = \psi'(t) - (t-1)\psi''(t)$  and

$f''(t) = -(t-1)\psi'''(t)$ . Also  $f(1) = 0$  and  $f'(1) = 0$ . Since  $\psi'''(t) < 0$  it follows that if

$t \geq 1$  then  $f''(t) \geq 0$  whence  $f'(t) \geq 0$  and  $f(t) \geq 0$ . This implies the first inequality. The second inequality follows from Taylor's theorem and the fact that  $\psi''(1) = p + q$ .

**Lemma 3.3** Suppose that  $\psi(t_1) = \psi(t_2)$  with  $t_1 \leq 1 \leq t_2$ . The following statements hold:

- i. One has  $\psi'(t_1) \leq 0$ ,  $\psi'(t_2) \geq 0$ , and  $-\psi'(t_1) \geq \psi'(t_2)$ .
- ii. If  $\beta \geq 1$ , then  $\psi(\beta t_1) \leq \psi(\beta t_2)$ ; equality holds if and only if  $\beta = 1$  or  $t_1 = t_2 = 1$ .

Proof: Proof of (i):

The statement is obvious if  $t_1 = 1$  or  $t_2 = 1$  because then  $\psi'(t_1) = \psi'(t_2) = 0$  implies  $t_1 = t_2 = 1$ .

Thus we may assume that  $t_1 < 1 < t_2$ . Suppose the opposite  $-\psi'(t_1) < \psi'(t_2)$ . By the mean value theorem we have

$$\psi'(t_1) = (1 - t_1)\psi''(\xi_2), \quad \text{for some } \xi_2 \in (1, t_2)$$

and

$$-\psi'(t_2) = (t_2 - 1)\psi''(\xi_1), \quad \text{for some } \xi_1 \in (t_1, 1).$$

Since  $\psi''(t)$  is monotonically decreasing one has  $\psi''(\xi_1) \geq \psi''(\xi_2)$ . Then we obtain

$$(t_2 - 1)\psi''(\xi_2) > (1 - t_1)\psi''(\xi_1) \geq (1 - t_1)\psi''(\xi_2),$$

Hence since  $\psi''(\xi_2) > 0$  it follows that  $t_2 - 1 > 1 - t_1$ . Using this and the fact that  $-\psi'(t)$  is convex, we may also write

$$\begin{aligned}
\psi(t_2) &= \\
&= \int_1^{t_2} \psi'(\xi) d\xi && \text{Since } \psi'(t) \text{ is concave,} \\
&\geq \frac{1}{2}(t_2 - 1)\psi'(t_2) && \text{Since } t_2 - 1 > 1 - t_1 \text{ and } \psi'(t_2) > 0, \\
&> \frac{1}{2}(1 - t_1)\psi'(t_2) && \text{Since } \psi'(t_1) < \psi'(t_2), \\
&> -\frac{1}{2}(1 - t_1)\psi'(1) && \text{Since } \psi'(t) \text{ is concave.} \\
&\geq -\int_{t_1}^1 \psi'(\xi) d\xi \\
&= \psi(t_1).
\end{aligned}$$

This contradiction proves the first part of the lemma.

Proof of (ii):

Consider,  $f(\beta) = \psi(\beta t_2) - \psi(\beta t_1)$ .

One has  $f(1) = 0$  and  $f'(\beta) = t_2 \psi'(\beta t_2) - t_1 \psi'(\beta t_1)$ .

Since  $\psi''(t) \geq 0$  for all  $t > 0$ ,  $\psi'(t)$  is monotonically increasing. Hence,  $\psi'(\beta t_1) \leq \psi'(\beta t_2)$ .

Substitution gives

$$f'(\beta) = t_2 \psi'(\beta t_2) - t_1 \psi'(\beta t_1) \geq t_2 \psi'(\beta t_2) - t_1 \psi'(\beta t_2) = \psi'(\beta t_2)(t_2 - t_1) \geq 0$$

The last inequality holds since  $t_2 \geq t_1$  and  $\psi'(t) \geq 0$  for  $t \geq 1$ . This proves that  $f(\beta) \geq 0$  for  $\beta \geq 1$ , and hence the inequality (ii) in the lemma follows.

If  $\beta = 1$  obviously we have equality. Otherwise, if  $\beta > 1$  and  $f(\beta) = 0$ , then the mean value theorem implies  $f'(\xi) = 0$  for some  $\xi \in (1, \beta)$ . But this implies  $\psi'(\xi t_2) = \psi'(\xi t_1)$ . Since  $\psi'(t)$

is strictly monotonic, this implies that  $\xi t_2 = \xi t_1$ , whence  $t_1 = t_2$ . Since also  $t_1 \leq 1 \leq t_2$  we obtain  $t_1 = t_2 = 1$ . This completes the proof of the second part of the lemma.

**Lemma 3.4** If  $t \geq 1$ , then  $\psi'(t)^2 \geq 2\psi(t)\psi''(t)$ .

Proof: Defining  $f(t) = \psi'(t)^2 - 2\psi(t)\psi''(t)$  one has  $f(1) = 0$  and

$$f'(t) = 2\psi'(t)\psi''(t) - 2\psi'(t)\psi''(t) - 2\psi(t)\psi'''(t) = -2\psi(t)\psi'''(t) > 0.$$

This proves the lemma.

**Lemma 3.5** Let  $\rho(s) : [0, \infty) \rightarrow (0, 1]$  be the inverse function of  $-\frac{1}{2}\psi'(t)$  for  $t \leq 1$ . The

following inequality holds:

$$\rho(s) \geq \frac{1}{(1+2s)^{\frac{1}{q}}}. \quad (3-3)$$

Proof: Since  $s = -\frac{1}{2}\psi'(t)$ , we have

$$-2s = t^p - t^{-q} \Rightarrow t^{-q} = t^p + 2s \leq 1 + 2s.$$

Since  $t = \rho(1)$ , this implies the lemma.

**Lemma 3.6** If  $t \geq 1$  and  $q \geq 2 - p$ , then  $t \leq 1 + \sqrt{t\psi(t)}$ .

Proof: Defining  $f(t) = t\psi(t) - (t-1)^2$  we have  $f(1) = 0$  and

$$f'(t) = \psi(t) + t\psi'(t) - 2(t-1).$$



Moreover, it is clear that  $f'(1) = 0$  and

$$f''(t) = 2\psi'(t) + t\psi''(t) - 2 = (2+p)t^p + (q-2)t^{-q} - 2 \geq pt^p + (q-2)t^q \geq p(t^p - t^{-q}) \geq 0.$$

The second inequality above is due to the fact that  $q \geq 2 - p$ . Thus we obtain

$$t\psi(t) \geq (t-1)^2,$$

which implies the lemma.

**Lemma 3.7** Let  $\varphi : [0, \infty) \rightarrow [1, \infty)$  be the inverse function of  $\psi(t)$  for  $t \geq 1$ . The following inequalities hold:

$$(1 + (p+1)s)^{\frac{1}{p+1}} \leq \varphi(s) \leq 1 + s + \sqrt{s^2 + 2s}. \quad (3-4)$$

If  $q \geq 2 - p$ , then

$$\varphi(s) \leq 1 + \sqrt{s + s^2 + s\sqrt{s^2 + 2s}} \quad (3-5)$$

Proof: Since  $q > 1$  and  $t \geq 1$ , we have

$$s = \psi(t) = \frac{t^{p+1} - 1}{p+1} + \frac{t^{1-q} - 1}{q-1} \leq \frac{t^{p+1} - 1}{p+1}$$

Hence, the first inequality in (3-4) follows.

The second inequality in (3-4) follows by using the first inequality of Lemma 3.2:

$$s = \psi(t) \geq \frac{1}{2}(t-1)\psi'(t) = \frac{1}{2}(t-1)(t^p - t^{-q}) \geq \frac{1}{2}(t-1)\left(1 - \frac{1}{t}\right) = \frac{1}{2}\left(t + \frac{1}{t} - 2\right).$$

Hence, solving the following inequality

$$t^2 - 2(1+s)^t + 1 \leq 0$$

leads to

$$t = \varphi(s) \leq 1 + s + \sqrt{s^2 + 2s}. \quad (3-6)$$

Finally, let  $q \geq 2 - p$ . By Lemma 3.6 one has  $t \leq 1 + \sqrt{t\psi(t)} \leq 1 + \sqrt{ts}$ .

Substitution of the upper bound for  $t$  given by (3-6) leads to

$$\varphi(s) \leq 1 + \sqrt{s + s^2 + s\sqrt{s^2 + 2s}}.$$

This completes the proof of the lemma.

Now we will derive a very important bound for normed proximity measure  $\delta(\nu) = \frac{1}{2} \|\nabla \Psi(\nu)\|$

in terms of the original proximity measure given by the barrier function  $\Psi(\nu)$ .

**Theorem 3.1** The following inequality holds:

$$\delta(\nu) \geq \frac{1}{2} \psi'(\varphi(\Psi(\nu))). \quad (3-7)$$

The proof is beyond the scope of thesis and can be found in (Peng, J., et. al., 2002)

**Corollary 3.1** If  $\Psi(\nu) \geq \tau \geq 1$ , then  $\delta(\nu) \geq \frac{1}{6} (\Psi(\nu))^{\frac{p}{1+p}}$

Proof: Using Theorem 2.1, and the fact that  $\Psi(v) \geq \tau \geq 1$ , we have

$$\delta(v) \geq \frac{1}{2} \psi'(\varphi(\Psi(v))) = \frac{1}{2} (\varphi(\Psi(v)))^p - ((\varphi(\Psi(v)))^{-q}) \geq \frac{1}{2} (\varphi(\Psi(v)))^p - ((\varphi(\Psi(v)))^{-1}).$$
 Note that

$t^p - \frac{1}{t}$  is monotonically increasing in  $t$ . Thus, by using the first inequality in (3-4), we obtain

$$\begin{aligned} \delta(v) &\geq \frac{1}{2} \frac{(\varphi(\Psi(v)))^{p+1} - 1}{\varphi(\Psi(v))} \geq \frac{1}{2} \frac{(1 + (p+1)\Psi(v))^{\frac{p+1}{p+1}} - 1}{(1 + (p+1)\Psi(v))^{\frac{1}{p+1}}} \\ &= \frac{1}{2} \frac{(p+1)\Psi(v)}{(1 + (p+1)\Psi(v))^{\frac{1}{p+1}}} \geq \frac{1}{2} \frac{\Psi(v)}{(3\Psi(v))^{\frac{1}{p+1}}} \\ &\geq \frac{1}{6} (\Psi(v))^{\frac{p}{p+1}} \end{aligned}$$

Which proves the corollary.

### Analysis of the algorithm

The outline of the analysis of the algorithm is as follows.

1. Outer iteration estimates:

- Estimate of the increase of the barrier function after the  $\mu$  update

2. Inner iteration estimates:

- Estimator for default step size
- Estimate of the decrease of the barrier function during the inner iteration with the default step size

### Outer Iteration Estimates

At the start of each outer iteration of the algorithm, just before the update of the parameter  $\mu$  with the factor  $1 - \theta$ , we have  $\psi(v) \leq \tau$ . Since the  $\mu$  vector is updated to  $\mu^+ = (1 - \theta)\mu$ , with  $0 < \theta < 1$ , the vector  $v$  is updated to  $v^+ = \frac{v}{\sqrt{1 - \theta}}$ , which in general leads to an increase in the value of  $\Psi(v)$ . Then, during the subsequent inner iterations,  $\Psi(v)$  decreases until it passes the threshold  $\tau$  again. During the course of the algorithm the largest values of  $\Psi(v)$  occur just after the updates of  $\mu$ . That is why we need to derive an estimate for the effect of a  $\mu$ -update on the value of  $\Psi(v)$ .

**Theorem 3.2** Let  $\varphi : [0, \infty) \rightarrow [1, \infty)$  as defined in Lemma 3.7. Then for any positive vector  $v$  and any  $\beta \geq 1$  the following inequality holds:

$$\Psi(\beta v) \leq n\psi\left(\beta\varphi\left(\frac{\Psi(v)}{n}\right)\right).$$

The proof is beyond the scope of this thesis and can be found in [Peng, J., et. al., 2002].

**Corollary 3.2** Let  $0 \leq \theta \leq 1$  and  $v_+ = \frac{v}{\sqrt{1 - \theta}}$ . If  $\Psi(v) \leq \tau$ , then

$$\Psi(v_+) \leq n\psi\left(\frac{\varphi\left(\frac{\tau}{n}\right)}{\sqrt{1 - \theta}}\right) \leq \frac{(p + q)n}{2} \left(\frac{\varphi\left(\frac{\tau}{n}\right)}{\sqrt{1 - \theta}} - 1\right)^2. \quad (3-8)$$

Proof: With  $\beta \geq 1$  and  $\Psi(\nu) \leq 1$  the first inequality follows from Theorem 3.2. The second inequality follows by using Lemma 3.2 and  $\psi''(1) = \frac{p+q}{q}$ .

The following upper bounds on the value of  $\psi(\nu_+)$  after the  $\mu$ -update follow immediately

$$\psi(\nu_+) \leq L_1 := n\varphi \left( \frac{1 + \frac{\tau}{n} + \sqrt{\left(\frac{\tau}{n}\right)^2 + \frac{2\tau}{n}}}{\sqrt{1-\theta}} \right), \quad q > 1; \quad (3-9)$$

and

$$\psi(\nu_+) \leq L_2 := n\varphi \left( \frac{1 + \sqrt{\frac{\tau}{n} + \frac{\tau^2}{n^2} + \frac{\tau}{n} \sqrt{\frac{\tau^2}{n^2} + \frac{2\tau}{n}}}}{\sqrt{1-\theta}} \right), \quad q \geq 2 - p. \quad (3-10)$$

### Default Step-size

In this subsection, we will determine a default step size which not only keeps the iterations feasible but also gives rise to sufficiently large decrease of  $\Psi(\nu)$  in each inner iteration. During an inner iteration, the parameter  $\mu$  is fixed. After the step in the direction  $(\Delta x, \Delta y, \Delta s)$  with step size  $\alpha$ , the new iterate is

$$x_+ = x + \alpha \Delta x, \quad s_+ = s + \alpha \Delta s, \quad y_+ = y + \alpha \Delta y \quad (3-11)$$

And a new  $\nu$ -vector is given by

$$v_+ = \sqrt{\frac{x_+ s_+}{\mu}}. \quad (3-12)$$

Since

$$\begin{aligned} x_+ &= x \left( e + \alpha \frac{\Delta x}{x} \right) = x \left( e + \alpha \frac{d_x}{v} \right) = \frac{x}{v} (u + \alpha d_x), \\ s_+ &= s \left( e + \alpha \frac{\Delta s}{s} \right) = s \left( e + \alpha \frac{d_s}{v} \right) = \frac{s}{v} (u + \alpha d_s), \\ xs &= \mu v^2 \end{aligned}$$

we obtain

$$v_+ = \sqrt{(v + \alpha d_x)(v + \alpha d_s)}.$$

Next, we consider the decrease in  $\Psi$  as a function of  $\alpha$ . We define two functions

$$f(\alpha) = \Psi(v_+) - \Psi(v), \quad (3-13)$$

and

$$f_1(\alpha) := \frac{1}{2} (\Psi(v + \alpha d_x) + \Psi(v + \alpha d_s)) - \Psi(v). \quad (3-14)$$

Lemma 3.1 implies that

$$\Psi(v_+) = \Psi(\sqrt{(v + \alpha d_x)(v + \alpha d_s)}) \leq \frac{1}{2} (\Psi(v + \alpha d_x) + \Psi(v + \alpha d_s)).$$

The above inequality shows that  $f_1(\alpha)$  is an upper bound of  $f(\alpha)$ . Obviously,

$f(0) = f_1(0) = 0$ . Taking the derivative with respect to  $\alpha$ , we get

$$f'_1(\alpha) = \frac{1}{2} \sum_{i=1}^n (\psi'(v_i + \alpha d_{xi}) d_{xi} + \psi'(v_i + \alpha d_{si}) d_{si}).$$

From the above equation and using that  $d_x + d_s = -\nabla\Psi(v)$  we obtain

$$f'_1(0) = \frac{1}{2} \nabla\Psi(v)^T (d_x + d_s) = -\frac{1}{2} \nabla\Psi(v)^T \nabla\Psi(v) = -2\delta(v)^2. \quad (3-15)$$

Differentiating once again, we get

$$f''_1(\alpha) = \frac{1}{2} \sum_{i=1}^n (\psi''(v_i + \alpha d_{xi}) d_{xi}^2 + \psi''(v_i + \alpha d_{si}) d_{si}^2) > 0, \text{ unless } d_x = d_s = 0.$$

(3-16)

It is worthwhile to point out that during an inner iteration  $x$  and  $s$  are not both at the  $\mu$  center since  $\Psi(v) \geq \tau \geq 0$ , so we may conclude that  $f_1(\alpha)$  is strictly convex in  $\alpha$ .

**Lemma 3.8** The following inequality holds:

$$f_1''(\alpha) \leq 2\delta^2 \psi''(v_{\min} - 2\alpha\delta). \quad (3-17)$$

Proof: Since  $d_x \perp d_s$ , and  $d_x + d_s = -\nabla\Psi(v)$  it is easy to see that  $\|(d_x, d_s)\| = 2\delta$ . where

$|\delta = \frac{1}{2} \|\nabla\Psi(v)\|$  Therefore, we have  $\|d_x\| \leq 2\delta$  and  $\|d_s\| \leq 2\delta$ . Hence,

$$v_i + \alpha d_x \geq v_{\min} - 2\alpha\delta, \quad v_i + \alpha d_s \geq v_{\min} - 2\alpha\delta, \quad 1 \leq i \leq n.$$

Using (3-16) and definition of  $\delta$ , we get

$$f_1''(a) \leq \frac{1}{2} \psi''(v_{\min} - 2a\delta) \sum_{i=1}^n (d_{xi}^2 + d_{si}^2) = 2\delta^2 \psi''(v_{\min} - 2a\delta).$$

This proves the lemma.

**Lemma 3.9** If the step-size  $\alpha$  satisfies

$$-\psi'(v_{\min} - 2\alpha\delta) + \psi'(v_{\min}) \leq 2\delta, \quad (3-18)$$

then  $f_1'(\alpha) \leq 0$ .

Proof: Using the Lemma 3.8,(3-15) and (3-17), we write:

$$\begin{aligned} f_1'(a) &= f_1'(0) + \int_0^a f_1''(\zeta) d\zeta \\ &\leq -2\delta^2 + 2\delta^2 \int_0^a \psi''(v_{\min} - 2\zeta\delta) d\zeta \\ &= -2\delta^2 - \delta \int_0^a \psi''(v_{\min} - 2\zeta\delta) d(v_{\min} - 2\zeta\delta) \\ &= -2\delta^2 - \delta (\psi'(v_{\min} - 2\alpha\delta) - \psi'(v_{\min})) \\ &\leq -2\delta^2 + 2\delta^2 = 0. \end{aligned}$$

This proves the lemma.

**Lemma 3.10** The largest possible value of the step-size satisfying the condition of Lemma 3.9 is given by

$$\bar{\alpha} := \frac{1}{2\delta} (\rho(\delta) - \rho(2\delta)) \quad (3-19)$$

Proof: We want  $\alpha$  such that (3-18) holds with  $\alpha$  as large as possible. Let us denote  $v_{\min}$  as  $v_1$ . Since  $\psi''(t)$  is decreasing the derivative with respect to  $v_1$  of the expression which is to the left side of the inequality (3-18) (i.e.  $\psi''(v_1 - 2\alpha\delta) + \psi''(v_1)$ ) is negative. Hence, fixing  $\delta$ , the smaller  $v_1$  is, the smaller  $\alpha$  will be. We have



$$\delta = \frac{1}{2} \|\nabla\Psi(\nu)\| \geq \frac{1}{2} |\psi'(\nu_1)| \geq -\frac{1}{2} \psi'(\nu_1).$$

Equality holds if and only if  $\nu_1$  is the only coordinate in  $\nu$  that differs from 1 and  $\nu_1 \leq 1$  (in case  $\psi'(\nu_1) \leq 0$ ). Hence the worst situation for the step size occurs when  $\nu_1$  satisfies

$$-\frac{1}{2} \psi'(\nu_1) = \delta. \quad (3-20)$$

The derivative with respect to  $\alpha$  of the expression that is the left side of the inequality (3-18) equals  $2\delta\psi''(\nu_1 - 2\alpha\delta) \geq 0$  and hence this expression is increasing in  $\alpha$ . Thus, the largest possible value of  $\alpha$  satisfying (3-18), satisfies

$$-\frac{1}{2} \psi'(\nu_1 - 2\alpha\delta) = 2\delta. \quad (3-21)$$

Due to the definition of  $\rho$ , (3-20) and (3-21) can be written as

$$\nu_1 = \rho(\delta), \quad \nu_1 - 2\alpha\delta = \rho(2\delta).$$

This implies

$$\alpha = \frac{1}{2\delta} (\nu_1 - \rho(2\delta)) = \frac{1}{2\delta} (\rho(\delta) - \rho(2\delta))$$

and the lemma is proved.

**Lemma 3.11** Let  $\bar{\alpha}$  be defined by (3-19). The following inequality holds:

$$\bar{\alpha} \geq \frac{1}{\psi''(\rho(2\delta))}. \quad (3-22)$$

Proof: By definition of  $\rho$ ,

$$-\psi'(\rho(\delta)) = 2\delta$$

Taking the derivative with respect to  $\delta$ , we find

$$-\psi''(\rho(\delta))\rho'(\delta) = 2$$

which leads to

$$\rho'(\delta) = -\frac{2}{\psi''(\rho(\delta))} < 0. \quad (3-23)$$

Hence,  $\rho$  is monotonically decreasing. An immediate consequence of (3-19) is

$$\bar{\alpha} = \frac{1}{2\delta} \int_{2\delta}^{\delta} \rho'(\sigma) d\sigma = \frac{1}{\delta} \int_{\delta}^{2\delta} \frac{d\sigma}{\psi''(\rho(\sigma))} \quad (3-24)$$

where we also used (3-23). To obtain a lower bound for  $\bar{\alpha}$ , we want to replace the argument of the last integral by its minimal value. We would like to know when  $\psi''(\rho(\sigma))$  is a maximal for  $\sigma \in [\delta, 2\delta]$ . We know that  $\psi''$  is monotonically decreasing. Thus  $\psi''(\rho(\sigma))$  is maximal for  $\sigma \in [\delta, 2\delta]$  when  $\rho(\sigma)$  is minimal. Since  $\rho$  is monotonically decreasing this occurs when  $\sigma = 2\delta$ . Therefore

$$\bar{\alpha} = \frac{1}{\delta} \int_{\delta}^{2\delta} \frac{d\sigma}{\psi''(\rho(\sigma))} \geq \frac{1}{\delta} \frac{\delta}{\psi''(\rho(2\delta))} = \frac{1}{\psi''(\rho(2\delta))},$$

and the lemma is proved.

**Theorem 3.3** We have  $\bar{\alpha} \geq \tilde{\alpha} := \frac{1}{(p+q)(1+4\delta)^{\frac{1+q}{q}}}$  (3-25)

Proof: Using Lemma 3.11, and the fact that  $\psi''(t)$  is monotonically decreasing for  $t \in (0, +\infty)$  we have

$$\bar{\alpha} \geq \frac{1}{\psi''(p(2\delta))} \geq \frac{1}{p(1+4\delta)^{\frac{1+q}{q}} + q(1+4\delta)^{\frac{1+q}{q}}} \geq \frac{1}{(p+q)(1+4\delta)^{\frac{q+1}{q}}} = \tilde{\alpha}$$

and the theorem is proved.

Thus, we can define the following default step-size

$$\tilde{\alpha} := \frac{1}{(p+q)(1+4\delta)^{\frac{q+1}{q}}}, \quad (3-26)$$

### Inner Iteration Estimates

Using the lower bound on the step size obtained in (3-25), we can obtain results on the decrease of the barrier function during inner iteration.

**Lemma 3.12** If the step size is such that  $\alpha \leq \bar{\alpha}$ , then  $f(\alpha) \leq -\alpha\delta^2$ . (3-27)

Proof: Let the univariate function  $h$  be such that

$$h(0) = f_1(0) = 0, \quad h'(0) = f_1' = 2\delta^2, \quad h''(a) = 2\delta^2\psi''(v_{\min} - 2a\delta)$$

According to (3-17) we have  $f'''(\alpha) \leq h''(\alpha)$  and that implies  $f'(\alpha) \leq h'(\alpha)$  and  $f(\alpha) \leq h(\alpha)$ .

Taking  $\alpha \leq \tilde{\alpha}$ , with  $\tilde{\alpha}$  defined in (3-26)), we have

$$h'(\alpha) = -2\delta^2 + 2\delta^2 \int_0^{\alpha} \psi''(v_{\min} - 2\xi\delta) d\xi = -2\delta^2 - \delta(\psi'(v_{\min} - 2\alpha\delta) - \psi'(v_{\min})) \leq 0.$$

Since  $\psi'''(t)$  decreasing in  $t$ ,  $h''(\alpha)$  is increasing  $\alpha$ . Using Lemma 3.13 below, we get:

$$f_1(a) \leq h(a) \leq \frac{1}{2} ah'(0) = -\alpha\delta^2.$$

As we mentioned before,  $f_1(a)$  is an upper bound of  $f(a)$ , hence, the lemma is proved.

**Theorem 3.4** The following inequality holds:

$$f(\tilde{\alpha}) \leq -\frac{1}{60(p+q)} \Psi(v)^{\frac{p(q-1)}{q(p+1)}}. \quad (3-28)$$

Proof: According to Lemma 3.12, if the step-size is such that  $\alpha \leq \bar{\alpha}$ , then  $f(\alpha) \leq -\alpha\delta^2$ . By

(3-25) the default step-size  $\tilde{\alpha}$  satisfies  $\tilde{\alpha} \leq \alpha$ , hence, the following upper bound for  $f(\alpha)$  is

obtained  $f(\tilde{\alpha}) \leq -\tilde{\alpha}\delta^2$ . Using Corollary 3.2 and the fact that  $f(\tilde{\alpha})$  is monotonically

decreasing in  $\delta$ , we obtain

$$f(\tilde{\alpha}) \leq \frac{\delta^2}{(p+q)(1+4\delta)^{\frac{q+1}{q}}}$$

$$\leq \frac{\Psi(\nu)^{\frac{2p}{p+1}}}{36(p+q) \left(1 + \frac{2}{3} \Psi(\nu)^{\frac{p}{p+1}}\right)^{\frac{q+1}{q}}}$$

$$\leq -\frac{1}{60(p+q)} \Psi(\nu)^{\frac{p(q-1)}{q(p+1)}}$$

Thus, the proof is complete.

### Estimate of the total number of iterations

As we've already mentioned, there are two types of algorithms.

- The **Short-step Algorithms**, where the barrier update parameter  $\theta$  depends on the size of the problem; that is,  $\theta = O\left(\frac{1}{\sqrt{n}}\right)$ .
- The **Long-step Algorithms**, where the barrier update parameter  $\theta$  is fixed; that is,  $\theta \in (0,1)$ .

We will give the estimate of the total number of iterations needed for both types of algorithms.

We will need following technical results. The proofs can be found in (Peng, J., et. al., 2002)]

**Lemma 3.13** If  $\alpha \in [0,1]$  and  $t \geq -1$  then  $(1+t)^\alpha \leq 1 + \alpha t$ .

**Lemma 3.14** Let  $h(t)$  be twice differentiable convex function with  $h(0) = 0$ ,  $h'(0) < 0$ , and let  $h(t)$  attain its (global) minimum at  $t^* > 0$ . If  $h''(t)$  is monotonically increasing for  $t \in \{0, t^*\}$  then  $h(t) \leq \frac{th'(0)}{2}$ ,  $0 \leq t \leq t^*$ .

**Lemma 3.15** Let  $t_0, t_1, \dots, t_k$  be a sequence of positive numbers such that

$$t_{k+1} \leq t_k - \beta_k^{1-\gamma}, \quad k = 0, 1, \dots, K-1 \quad \text{where } \beta > 0 \text{ and } 0 < \gamma \leq 1. \text{ Then } K \leq \left\lceil \frac{t_0^\gamma}{\beta\gamma} \right\rceil.$$

### Long-step Algorithms

**Lemma 3.16** The total number of outer iterations in both cases is the same.

$$\frac{1}{\theta} \log \frac{n}{\varepsilon} \tag{3-29}$$

Proof: The number of outer iterations is the number of iterations  $K$  necessary to obtain  $n\mu \leq \varepsilon$ . Previous and new  $\mu$  are related as follows  $\mu := (1 - \theta)\mu$ . Thus,  $n\mu \leq \varepsilon$  can be written as  $\mu_0(1 - \theta)^K \leq \frac{\varepsilon}{n}$ . We can assume that  $\mu_0 = 1$  and by taking the logarithm of both sides of the inequality we obtain  $K \log(1 - \theta) \leq \log \frac{\varepsilon}{n}$ . Using the Taylor theorem for  $\log(1 - \theta)$  we obtain  $K \leq \frac{1}{\theta} \log \frac{n}{\varepsilon}$  proving the lemma.

Now we need to estimate the upper bound on the total number of inner iterations per one outer iteration for the large-step methods. That number is equal to the number of iterations necessary to return to the situation  $\Psi(\nu) \leq \tau$ . We denote the value of  $\Psi(\nu)$  after the  $\mu$  update

as  $\Psi_0$ . The subsequent values in the same outer iteration are denoted as  $\Psi_k$ ,  $k = 1, 2, \dots, K$  where  $K$  denotes the total number of inner iterations in the outer iteration. By using (3-9), we have

$$\Psi_0 \leq n\psi \left( \frac{1 + \frac{\tau}{n} + \sqrt{\left(\frac{\tau}{n}\right)^2 + \frac{2\tau}{n}}}{\sqrt{1 - \Theta}} \right).$$

Since  $\psi(t) \leq \frac{t^{p+1} - 1}{p+1}$  when  $t \geq 1$  and  $1 - (1 - \theta)^{\frac{p+1}{2}} \leq \theta$ , after some elementary reductions, we

obtain:

$$\Psi_0 \leq \frac{n\theta + (p+1)\tau + n(p+1)\sqrt{\left(\frac{\tau}{n}\right)^2 + \frac{2\tau}{n}}}{(p+1)(1-\theta)^{\frac{p+1}{2}}}. \quad (3-30)$$

Now Theorem 3.4 leads to

$$\Psi_{k+1} \leq \Psi_k - \beta(\Psi_k)^{1-\gamma}, \quad k = 0, 1, \dots, K-1, \quad (3-31)$$

where  $\beta = \frac{1}{60(p+q)}$  and  $\gamma = \frac{p+q}{q(p+1)}$ . Using Lemma 3.15 and (3-30) and (3-31) we obtain

the following upper bound on the number  $K$  of inner iterations.

$$K \leq 60q(p+1)(\Psi_0)^{\frac{p+q}{q(p+1)}} \leq 60q(p+1) \left( \frac{n\theta + (p+1)\tau + n(p+1)\sqrt{\left(\frac{\tau}{n}\right)^2 + \frac{2\tau}{n}}}{(p+1)(1-\theta)^{\frac{p+1}{2}}} \right)^{\frac{p+q}{q(p+1)}} \quad (3-$$

32)

Now we can derive an upper bound on the total number iterations needed by the large-update version of the Generic IPM in Figure 2.1. According to Lemma 3.16 the number of outer iterations is bounded above by

$$\frac{1}{\theta} \log \frac{n}{\varepsilon}$$

By multiplying the number of outer iterations and the number of inner iterations obtained in

(3-32) in the lemma above we get an upper bound for the total number of iterations

$$\frac{60q(p+1)}{\theta} \left( \frac{n\theta + (p+1)\tau + n(p+1)\sqrt{\left(\frac{\tau}{n}\right)^2 + \frac{2\tau}{n}}}{(p+1)(1-\theta)^{\frac{p+1}{2}}} \right)^{\frac{p+q}{q(p+1)}} \log \frac{n}{\varepsilon}.$$

For large update methods we know that  $\theta = \Theta(1)$ , and  $\tau = O(n)$ . After some elementary transformations the iteration bound reduces to the following bound

$$O \left( q n^{\frac{p+q}{q(p+q)}} \log \frac{n}{\varepsilon} \right) \quad (3-33)$$

This result is summarized in the theorem below



**Theorem 3.5:** Given that  $\theta = \Theta(1)$  and  $\tau = O(n)$  which are characteristics of the large-update methods the Generic IPM described in the Figure 2.1 will obtain  $\varepsilon$  - appropriate solutions of

(P) and (D) in at most  $O\left(q n^{\frac{p+q}{q(p+q)}} \log \frac{n}{\varepsilon}\right)$  iterations.

The obtained complexity result contains several previously known complexity results as special cases.

1. When  $p = 1$  and  $q > 1$ , the kernel function  $\psi(t)$  becomes the prototype self-regular function. If in addition,  $q = \log n$  the iteration bound reduces to the best known bound for self-regular function, which is  $O\left(\sqrt{n} \log n \log \frac{n}{\varepsilon}\right)$ .
2. Letting  $p = 1$  and  $q = 1$ , the iteration bound becomes  $O\left(n \log \frac{n}{\varepsilon}\right)$  and  $\psi(t)$  represents the classical logarithmic kernel function.
3. For  $q = 2$  and  $p = 0$ ,  $\psi(t)$  represents the simple kernel function  $\psi(t) = t - \frac{1}{t} - 2$  which is not self-regular. The iteration bound is the same as the one obtained for the logarithmic kernel function.

### Short-Step Algorithms

To get the best possible bound for short-step methods we need to use the bound described in (3-10).

$$\Psi_0 \leq n\psi \left( \frac{1 + \sqrt{\frac{\tau}{n} + \frac{\tau^2}{n^2} + \frac{\tau}{n} \sqrt{\frac{\tau^2}{n^2} + \frac{2\tau}{n}}}}{\sqrt{1-\theta}} \right) \leq \frac{(p+q)n}{2} \left( \frac{1 + \sqrt{\frac{\tau}{n} + \frac{\tau^2}{n^2} + \frac{\tau}{n} \sqrt{\frac{\tau^2}{n^2} + \frac{2\tau}{n}}}}{\sqrt{1-\theta}} - 1 \right)^2$$

Using  $1 - \sqrt{1-\theta} = \frac{\theta}{1 + \sqrt{1-\theta}} \leq \theta$  the above inequality can be simplified to

$$\Psi_0 \leq \frac{p+q}{2(1-\theta)} \left( \theta \sqrt{n + \sqrt{\tau + \frac{\tau^2}{n} + \tau \sqrt{\frac{\tau^2}{n^2} + \frac{2\tau}{n}}}} \right)^2 \quad (3-34)$$

Following the same line of arguments as in the above subsection 3.3.1 we conclude that the total number  $K$  of inner iterations is bounded above by

$$K \leq 60q(p+1)(\Psi_0)^{\frac{p+q}{q(p+1)}} \leq \frac{p+q}{(1-\theta)} \left( \theta \sqrt{n + \sqrt{\tau + \frac{\tau^2}{n} + \tau \sqrt{\frac{\tau^2}{n^2} + \frac{2\tau}{n}}}} \right)^{\frac{2(p+q)}{q(p+1)}} \quad (3-35)$$

Given the upper bound on the number of the outer iterations as mentioned in the previous subsection 3.3.1 the upper bound on the total number of iterations is

$$\frac{60q(p+q)}{\theta(1-\theta)} \left( \theta \sqrt{n + \sqrt{\tau \frac{\tau^2}{n} + \tau \sqrt{\frac{\tau^2}{n^2} + \frac{2\tau}{n}}}} \right)^{\frac{2(p+q)}{q(p+1)}} \log \frac{n}{\varepsilon}. \quad (3-36)$$

For small update methods it is well known that  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$  and  $\tau = O(1)$ . After some elementary reductions one easily obtains that the iteration bound is  $O\left(q^2 \sqrt{n} \log \frac{n}{\varepsilon}\right)$ . We summarize this result in the theorem below.

**Theorem 3.6:** Given that  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$  and  $\tau = O(1)$  which are characteristics of the small update methods the Generic IPM described in the Figure 2.1 we will obtain  $\varepsilon$  - appropriate solutions of (P) and (D) in at most  $O\left(q^2 \sqrt{n} \log \frac{n}{\varepsilon}\right)$  iterations.

## CHAPTER 4

### ANALYSIS OF THE ALGORITHM FOR ADDITIONAL KERNEL FUNCTIONS

#### Summary of the algorithm analysis

Looking carefully at the analysis of the Generic IPM described in Chapter 3 the procedure can be summarized in the following way.

Step 0: Input a kernel function  $\psi$  ; an update parameter  $\theta$ ,  $0 < \theta < 1$  ; a threshold parameter  $\tau$  ; and an accuracy parameter  $\varepsilon$  .

Step 1: Solve the equation  $-\frac{1}{2}\psi'=s$  to get  $\rho(s)$  the inverse function of  $-\frac{1}{2}\psi'(t), t \in (0,1]$  . If the equation is hard to solve, derive a lower bound for  $\rho(s)$  .

Step 2: Calculate the decrease of  $\Psi(t)$  in terms of  $\delta$  for the default step-size  $\tilde{\alpha}$  from

$$f(\tilde{\alpha}) \leq -\frac{\delta^2}{\psi''(\rho(2\delta))} .$$

Step 3: Solve the equation  $\psi(t) = s$  to get  $\varphi(s)$  , the inverse function of  $\psi(t), t \geq 1$  . If the equation is hard to solve, derive lower and upper bounds for  $\varphi(s)$  .

Step 4: Derive a lower bound for  $\delta$  in terms of  $\Psi(v)$  by using  $\delta(v) \geq \frac{1}{2}\psi'(\varphi(\Psi(v)))$  .

Step 5: Using the results of Step 3 and Step 4 find a valid inequality of the form

$$f(\tilde{\alpha}) \leq -\kappa\Psi(v)^{1-\gamma} \text{ for some positive constants } \kappa \text{ and } \gamma \in (0,1] .$$

Step 6: Calculate the upper bound of  $\Psi_0$  from

$$\Psi_0 \leq L_\psi(n, \theta, \tau) = n\psi\left(\frac{\varphi\left(\frac{\tau}{n}\right)}{\sqrt{1-\theta}}\right) \leq \frac{n}{2}\psi''(1)\left(\frac{\varphi\left(\frac{\tau}{n}\right)}{\sqrt{1-\theta}} - 1\right)^2.$$

Step 7: Derive an upper bound for the total number of iterations from

$$\frac{\Psi_0^\gamma}{\theta \kappa \gamma} \log \frac{n}{\varepsilon}.$$

Step 8: Set  $\tau = O(n)$  and  $\theta = \Theta(1)$  so as to calculate complexity bound for large-update methods, or set  $\tau = O(1)$  and  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$  so as to calculate the complexity bound for small update methods.

### Additional Kernel Functions

We will consider the following additional kernel functions.

$$\psi_1(t) = \frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q(q-1)} - \frac{q-1}{q}(t-1), \quad q > 1.$$

(4-1)

$$\psi_2(t) = \frac{t^2 - 1}{2} + \frac{e^{\frac{1}{t}} - e}{e} \tag{4-2}$$

$$\psi_3(t) = \frac{t^2 - 1}{2} - \int_1^t e^{\frac{1}{\zeta} - 1} d\zeta \quad (4-3)$$

The growth term in all of them is the same  $\psi_g(t) = \frac{t^2 - 1}{2}$  while the barrier term varies  $\psi_b(t)$ . The reason for considering this growth term is that according to the analysis above it seems to give the best complexity results and, thus, will give a more consistent view of the complexity analysis.

The following lemmas are useful for the above kernel functions. They are variations of the similar lemmas in the previous chapter and actually they are also valid for the class of kernel functions (3-1) used in that chapter. Their main purpose is to help facilitate the summary analysis described in the previous subsection.

**Lemma 4.1** When  $\psi(t) = \psi_i(t)$  for  $1 \leq i \leq 3$ , then  $\sqrt{1 + 2s} \leq \varphi(s) \leq 1 + \sqrt{2s}$ .

Proof: The inverse function of  $\psi(t)$  for  $t \in [1, \infty)$  is obtained by solving for  $t$  from the equation  $\psi(t) = s$ , for  $t \geq 1$ . In almost all cases it is hard to solve this equation explicitly.

However, we can easily find a lower and an upper bound for  $t$  and this suffices for our goal.

First one has

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

where  $\psi_b(t)$  denotes the barrier term. The inequality is due to the fact that  $\psi_b(1) = 0$  and  $\psi_b(t)$  is monotonically decreasing. It follows that

$$t = \varphi(s) \geq \sqrt{1 + 2s}$$

For the second inequality we use the fact that  $\psi_i''(t) \geq 1$  for  $1 \leq i \leq 3$ . Note that  $\psi_i(t)$  are nonnegative strictly convex functions such that  $\psi_i(1) = 0$ . This implies that  $\psi_i(t)$  is twice differentiable and therefore is completely determined by its second derivative

$$\psi_i(t) = \int_1^t \int_1^\xi \psi_i''(\zeta) d\zeta d\xi \quad (4-4)$$

Thus

$$s = \psi_i(t) = \int_1^t \int_1^\xi \psi_i''(\zeta) d\zeta d\xi \geq \int_1^t \int_1^\xi d\zeta d\xi = \frac{1}{2}(t-1)^2$$

which implies

$$t = \varphi(s) \leq 1 + \sqrt{2s}.$$

This completes the proof.

**Lemma 4.2** Let  $1 \leq i \leq 3$ . Then one has  $L_\psi(n, \theta, \tau) \leq \frac{\psi''(1) (\sqrt{2\tau} + \theta\sqrt{n})^2}{2(1-\theta)}$ . Hence, if

$\tau = O(1)$  and  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$ , then  $\Psi_0 = O(\psi''(1))$ .

Prof: By Lemma 4.1 we have  $\varphi(s) \leq 1 + \sqrt{2s}$ . Hence, by using Theorem 3.2 and first inequality in (3-8) we have

$$\Psi_0 \leq L_\psi(n, \theta, \tau) = n\psi \left( \frac{\varphi\left(\frac{\tau}{n}\right)}{\sqrt{1-\theta}} \right) \leq n\psi \left( \frac{1 + \sqrt{\frac{2\tau}{n}}}{\sqrt{1-\theta}} \right) \quad (4-5)$$

By using Taylor theorem and the fact that  $\psi(1) = \psi'(1) = 0$  we obtain

$$\psi(t) = \frac{1}{2}\psi''(1)(t-1)^2 + \frac{1}{3!}\psi'''(\xi)(\xi-1)^3$$

Given the fact that  $\psi'''(\xi) < 0$  we obtain for  $t > 1$  and  $1 < \xi < t$

$$\psi(t) \leq \frac{1}{2}\psi''(1)(t-1)^2 \quad (4-6)$$

Applying (4-6) to (4-5) with  $t = \frac{1 + \sqrt{\frac{2\tau}{n}}}{\sqrt{1-\theta}}$  we obtain

$$\psi_0 \leq \frac{n\psi''(1)}{2} \left( \frac{1 + \sqrt{\frac{2\tau}{n}}}{\sqrt{1-\theta}} - 1 \right)^2 \leq \frac{n\psi''(1)}{2} \left( \frac{\theta + \sqrt{\frac{2\tau}{n}}}{\sqrt{1-\theta}} \right)^2 = \frac{\psi''(1)}{2} \frac{(\sqrt{2\tau} + \theta\sqrt{n})^2}{1-\theta}$$

where we also used the fact that  $1 - \sqrt{1-\theta} = \frac{\theta}{1 + \sqrt{1-\theta}} \leq \theta$ . This proves the lemma.

**Lemma 4.3** Let  $\eta : [0, \infty) \rightarrow (0, 1]$  be the inverse function of the restriction of  $-\psi_b'(t)$  to the interval  $(0, 1]$  where  $\psi_b(t)$  is the barrier term in the kernel functions  $\psi_i(t)$ ,  $1 \leq i \leq 3$ . Then

$$\rho(s) \geq \eta(1 + 2s).$$



Proof: Let  $t = \rho(s)$ . Due to the definition of  $\rho$  as the inverse function of

$-\frac{1}{2}\psi'(t)$  for  $t \leq 1$  this means that

$$-2s = \psi'(t) = t + \psi_b'(t), \quad t \leq 1.$$

Since  $t \leq 1$  this implies

$$-\psi_b'(t) = t + 2s \leq 1 + 2s$$

Since  $-\psi_b'(t)$  is monotonically decreasing in all three cases, it follows that

$$t = \rho(s) \geq \eta(1 + 2s),$$

proving the lemma.

### **Analysis of the Generic IPM with Additional Kernel Functions**

In this subsection we will provide the analysis of the Generic IPM using additional kernel functions stated in the subsection. We will follow the steps described earlier.

#### **Example**

Consider the function

$$\psi_1(t) = \psi(t) = \frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q(q-1)} - \frac{q-1}{q}(t-1), \quad q > 1.$$

Step 1: The inverse function of  $-\psi_b'(t) = 1 + \frac{t-q-1}{q}$  is given by  $\eta(s) = -\frac{1}{(1+q(s-1))^{\frac{1}{q}}}$ .

Hence, by Lemma 6.3,  $\rho(s) = \frac{1}{(1+2qs)^{\frac{1}{q}}}$ .

Step 2: It follows that

$$f(\alpha) \leq -\frac{\delta^2}{\psi''(\eta(2\delta))} = -\frac{\delta^2}{1 + \frac{1}{\rho(2\delta)^{q+1}}} \leq -\frac{\delta^2}{1 + (1+4q\delta)^{\frac{q}{q+1}}}. \quad (4-7)$$

Step 3: By Lemma 6.1 the inverse function of  $\psi(t)$  for  $t \in [1, \infty)$  satisfies

$$\sqrt{1+2s} \leq \varphi(s) \leq 1 + \sqrt{2s}$$

Omitting the argument  $\nu$ , we therefore have

$$\varphi(\Psi(\nu)) \geq \sqrt{1+2\Psi}.$$

Step 4: Now using the fact that  $\delta(\nu) \geq \frac{1}{2}\psi'(\varphi(\Psi(\nu)))$ , and assuming  $\Psi \geq \tau \geq 1$ , we obtain

$$\delta \geq \frac{1}{2}(\sqrt{1+2\Psi} - 1) + \frac{1}{q} \left( 1 - \frac{1}{(1+2\Psi)^q} \right) \geq \frac{1}{2}(\sqrt{1+2\Psi} - 1) = \frac{\Psi}{1 + \sqrt{1+2\Psi}}. \quad (4-8)$$

Step 5: Combining (4-7) and (4-8) after some elementary reductions, we obtain

$$f(\tilde{\alpha}) \leq -\frac{\delta^2}{1 + (1+4q\delta)^{\frac{q}{q+1}}} \leq -\frac{1}{53} \Psi^{\frac{q-1}{2q}}. \quad (4-9)$$

Thus, it follows that

$$\Psi_{k+1} \leq \Psi_k - \kappa(\Psi_k)^{1-\gamma}, \quad k = 0, 1, \dots, K-1$$

with  $\kappa = \frac{1}{53q}$  and  $\gamma = \frac{q+1}{2q}$ , and  $K$  denotes the number of inner iterations. Hence, by Lemma

3.15 the number  $K$  of inner iterations is bounded above by

$$K \leq \frac{\Psi_0}{k\gamma} = \frac{106q^2}{q+1} \Psi_0 \leq 106q \Psi_0^{\frac{q+1}{2q}}. \quad (4-10)$$

Step 6: To estimate  $\Psi_0$  we use Lemma 6.2, with  $\psi''(1) = 2$ . Thus, we obtain

$$\Psi_0 \leq \frac{(\theta\sqrt{n} + \sqrt{2\tau})^2}{1-\theta}.$$

Step 7: Thus, using Lemma 3.16 the total number of iterations is bounded above by

$$\frac{K}{\theta} \log \frac{n}{\varepsilon} \leq \frac{106q}{\theta} \left( \frac{(\theta\sqrt{n} + \sqrt{2\tau})^2}{1-\theta} \right)^{\frac{q+1}{2q}} \log \frac{n}{\varepsilon}. \quad (4-11)$$

Step 8: For large update methods ( with  $\tau = O(n)$  and  $\theta = \Theta(1)$  ) the right hand side expression becomes

$$O \left( qn^{\frac{q+1}{2q}} \log \frac{n}{\varepsilon} \right). \quad (4-12)$$

For small update methods ( with  $\tau = O(1)$  and  $\theta = \Theta \left( \frac{1}{\sqrt{n}} \right)$  ) the right hand side expression

becomes

$$O \left( q\sqrt{n} \log \frac{n}{\varepsilon} \right). \quad (4-13)$$

### Example

Consider the kernel function

$$\psi_2(t) = \psi(t) = \frac{t^2 - 1}{2} + \frac{e^{\frac{1}{t}} - e}{e}.$$

Step 1: The inverse function of  $-\psi_b'(t) = \frac{e^{\frac{1}{t}} - 1}{t^2}$  is such that  $\eta(s) = t \Leftrightarrow \frac{e^{\frac{1}{t}} - 1}{t^2} = s, t \leq 1$ . It

follows that  $\frac{e^{\frac{1}{t}} - 1}{t^2} = st^2 \leq s$  whence we obtain  $\eta(s) = t \geq \frac{1}{1 + \log s}$ . Hence, by Lemma 4.3,

$$\rho(s) \geq \eta(1 + 2s).$$

Step 2: Since  $\psi''(t)$  is monotonically decreasing we have

$$f(\tilde{\alpha}) \leq -\frac{\delta^2}{\psi''(\rho(2\delta))} \leq \frac{\delta^2}{\psi''(\eta(1 + 4q\delta))}.$$

Now putting  $t = \eta(1 + 4q\delta)$  we have  $t \leq 1$  and we may write

$$f(\alpha) \leq -\frac{\delta^2}{\psi''(t)} = -\frac{\delta^2}{1 + \frac{1+2t}{t^4} e^{\frac{1}{t}-1}} \leq \frac{\delta^2}{1 + \frac{3}{t^4} e^{\frac{1}{t}-1}} = \frac{\delta^2}{1 + \frac{3(1+4\delta)}{t^2}} \leq \frac{\delta^2}{1 + \frac{15\delta}{t^2}}.$$

Since  $\frac{1}{t^2} = \frac{1}{(\eta(1 + 4\delta))^2} \leq (1 + \log(1 + 4\delta))^2$  we finally get

$$f(\alpha) = \frac{\delta^2}{1 + 15\delta(1 + \log(1 + 4\delta))^2}. \quad (4-14)$$

Step 3: By Lemma 6.1 the inverse function of  $\psi(t)$  for  $t \in [1, \infty)$  satisfies

$$\sqrt{1+2s} \leq \varphi(s) \leq 1 + \sqrt{2s}. \text{ Omitting the argument } \nu, \text{ we thus have } \varphi(\Psi(\nu)) \geq \sqrt{1+2\Psi}.$$

Step 4: Now using that  $\delta(\nu) \geq \frac{1}{2}\psi'(\varphi(\Psi(\nu)))$ , we obtain

$$\delta \geq \frac{1}{2} \left( \sqrt{1+2\Psi} - \frac{e^{\frac{1}{\sqrt{1+2\Psi}}-1}}{1+2\Psi} \right) \geq \frac{1}{2} (\sqrt{1+2\Psi} - 1) = \frac{\Psi}{1+\sqrt{1+2\Psi}}. \quad (4-15)$$

Step 5: Substitution of ((4-15) into the (4-14) gives, after some elementary reductions, while assuming  $\Psi_0 \geq \Psi \geq \tau \geq 1$

$$f(\alpha) \leq -\frac{\Psi^{\frac{1}{2}}}{44(1+\log(1+\sqrt{2\Psi}))^2} \leq -\frac{\Psi^{\frac{1}{2}}}{44(1+\log(1+\sqrt{2\Psi_0}))^2}.$$

Thus, it follows that

$$\Psi_{k+1} \leq \Psi_k - \kappa(\Psi_k)^{1-\gamma}, \quad k = 0, 1, \dots, K-1$$

with  $\kappa = \frac{1}{44(1+\log(1+\sqrt{2\Psi_0}))^2}$  and  $\gamma = \frac{1}{2}$ , and  $K$  denotes the number of inner iterations.

Hence, by Lemma 3.15 the number  $K$  of inner iterations is bounded above by

$$K \leq \frac{\Psi_0}{k\gamma} = 88(1+\log(1+\sqrt{2\Psi_0}))^2 \Psi_0^{\frac{1}{2}}.$$

Step 6: We use Lemma 4.2 with  $\psi''(1) = 4$ , to estimate  $\Psi_0$ .

$$\Psi_0 \leq 2 \frac{(\theta\sqrt{n} + \sqrt{2\tau})^2}{1-\theta}. \quad (4-16)$$

Substitution of (4-16) in the expression for  $K$  gives

$$K \leq 88\sqrt{2} \left( 1 + \log \left( 1 + 2 \frac{\theta\sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}} \right) \right)^2 \frac{\theta\sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}}. \quad (4-17)$$

Step 7: Thus the total number of iterations is bounded above by

$$\frac{K}{\theta} \log \frac{n}{\varepsilon} \leq 88\sqrt{2} \left( 1 + \log \left( 1 + 2 \frac{\theta\sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}} \right) \right)^2 \frac{\theta\sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}} \log \frac{n}{\varepsilon}. \quad (4-18)$$

Step 8: For large update methods, when  $\tau = O(n)$  and  $\theta = \Theta(1)$  the right hand side expression

(4-18) becomes

$$O\left(\sqrt{n}(\log n)^2 \log \frac{n}{\varepsilon}\right). \quad (4-19)$$

For small update methods, when  $\tau = O(1)$  and  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$ , the right hand side expression

(4-18) becomes

$$O\left(\sqrt{n} \log \frac{n}{\varepsilon}\right) \quad (4-20)$$

### Example

Consider the kernel function

$$\psi_3(t) = \psi(t) = \frac{t^2 - 1}{2} - \int_1^t e^{\frac{1}{s}-1} d\zeta .$$

Step 1: The inverse function of  $-\psi'_b(t) = e^{\frac{1}{t}-1}$  is given by  $\eta(s) = \frac{1}{1 + \log s}$ . Hence, by Lemma

6.3,

$$\rho(s) \geq \frac{1}{1 + \log(1 + 2s)} .$$

Step 2: It follows that

$$f(\alpha) \leq -\frac{\delta^2}{\psi''(\rho(2\delta))} = -\frac{\delta^2}{1 + \frac{1}{e^{\frac{1}{\rho(2\delta)}-1}}} \leq -\frac{\delta^2}{1 + (1 + 4q\delta)(1 + \log(1 + 4\delta))^2} .$$

(4-21)

Step 3: By Lemma 4.1 the inverse function of  $\psi(t)$  for  $t \in [1, \infty)$  satisfies

$$\sqrt{1 + 2s} \leq \varphi(s) \leq 1 + \sqrt{2s} .$$

Thus we have, omitting the argument  $v$ ,

$$\varphi(\Psi(v)) \geq \sqrt{1 + 2\Psi} .$$

Step 4: Now using that  $\delta \geq \frac{1}{2}\psi'(\varphi(\Psi(v)))$  we obtain

$$\delta \geq \frac{1}{2} \left( \sqrt{1 + 2\Psi} - e^{\frac{1}{\sqrt{1 + 2\Psi}}-1} \right) \geq \frac{1}{2} (\sqrt{1 + 2\Psi} - 1) = \frac{\Psi}{1 + \sqrt{1 + 2\Psi}} . \quad (4-22)$$

Step 5: Substitution of (4-22) into the (4-21) gives, after some elementary reductions, while assuming  $\Psi_0 \geq \Psi \geq \tau \geq 1$

$$f(\alpha) \leq -\frac{\Psi^{\frac{1}{2}}}{21(1 + \log(1 + \sqrt{\Psi}))^2} \leq -\frac{\Psi^{\frac{1}{2}}}{21(1 + \log(1 + \sqrt{\Psi_0}))^2}.$$

Thus, it follows that

$$\Psi_{k+1} \leq \Psi_k - \kappa(\Psi_k)^{1-\gamma}, \quad k = 0, 1, \dots, K-1$$

with  $\kappa = \frac{1}{21(1 + \log(1 + \sqrt{2\Psi_0}))^2}$  and  $\gamma = \frac{1}{2}$ , and  $K$  denotes the number of inner iterations.

Hence, by Lemma 3.15 the number  $K$  of inner iterations is bounded above by

$$K \leq \frac{\Psi_0}{k\gamma} = 42(1 + \log(1 + \sqrt{2\Psi_0}))^2 \Psi_0^{\frac{1}{2}}. \quad (4-23)$$

We use Lemma 4.2, with  $\psi''(1) = 2$  to estimate  $\Psi_0$ . This gives

$$\Psi_0 \leq \frac{(\theta\sqrt{n} + \sqrt{2\tau})^2}{1-\theta}. \quad (4-24)$$

Substitution of (4-24) into (4-23) leads to the following estimate for number  $K$  of inner iterations

$$K \leq 42 \left( 1 + \log \left( 1 + \frac{\theta\sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}} \right) \right)^2 \frac{\theta\sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}}. \quad (4-25)$$

Step 7: By Lemma 3.16 the total number of iterations is bounded above by



$$\frac{K}{\theta} \log \frac{n}{\varepsilon} \leq 42 \left( 1 + \log \left( 1 + \frac{\theta \sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}} \right) \right)^2 \frac{\theta \sqrt{n} + \sqrt{2\tau}}{\sqrt{1-\theta}} \log \frac{n}{\varepsilon} \quad (4-26)$$

Step 8: For large-update methods when  $\tau = O(n)$  and  $\theta = \Theta(1)$  the right hand side expression

(4-26) becomes

$$O\left(\sqrt{n}(\log n)^2 \log \frac{n}{\varepsilon}\right). \quad (4-27)$$

For small update methods, when  $\tau = O(1)$  and  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$ , the right hand side expression

(4-26) becomes

$$O\left(\sqrt{n} \log \frac{n}{\varepsilon}\right). \quad (4-28)$$

### Summary of complexity results

The complexity results for Generic IPM with kernel functions defined in (3-1) and in (4-1)-(4-3) are summarized in the Table for large-step methods and in the table for small-step methods.

For the class of kernel functions (3-1) we consider three special cases,

- the logarithmic kernel function:  $\psi(t) = \frac{t^2 - 1}{2} - \log t$
- the classical self-regular function when  $p = 1$ :  $\psi(t) = \frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q - 1}$

- the linear non-self-regular function when  $p = 0$ :  $\psi(t) = t - 1 + \frac{t^{1-q} - 1}{q - 1}$

### Complexity of large-update methods

The complexity results for large-step methods are summarized below. They are obtained by taking into the account that  $\tau = O(n)$  and  $\theta = \Theta(1)$ .

#### Complexity results for long-step methods

i	Kernel Function $\psi_i(t)$	Iteration Bound
1	$\frac{t^2 - 1}{2} - \log t$	$O(n) \log \frac{n}{\varepsilon}$
2	$\frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q - 1}$	$O(qn^{\frac{q+1}{2q}}) \log \frac{n}{\varepsilon}$
3	$t - 1 + \frac{t^{1-q} - 1}{q - 1}$	$O(qn) \log \frac{n}{\varepsilon}$
4	$\frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q(q-1)} - \frac{q-1}{q}(t-1)$	$O(qn^{\frac{q+1}{2q}}) \log \frac{n}{\varepsilon}$
5	$\frac{t^2 - 1}{2} + \frac{e^t - e}{e}$	$O(\sqrt{n} \log^2 n) \log \frac{n}{\varepsilon}$
6	$\frac{t^2 - 1}{2} - \int_1^t e^{\xi-1} d\xi$	$O(\sqrt{n} \log^2 n) \log \frac{n}{\varepsilon}$

Table 2

Notice that the best bound is obtained in case of 3 and 4 by taking  $q = \frac{1}{2} \log n$  which gives the

iteration bound of

$$O\left(\sqrt{n} \log n \log \frac{n}{\varepsilon}\right), \quad (4-29)$$

which is currently the best known bound for large-update methods.

### Complexity of small update methods

The complexity results for small-update methods are summarized below. They are obtained by

taking into the account that  $\theta = \Theta\left(\frac{1}{\sqrt{n}}\right)$  and  $\tau = O(1)$ .

#### Complexity results for short-step methods

i	Kernel Function $\psi_i(t)$	Iteration Bound
1	$\frac{t^2 - 1}{2} - \log t$	$O(n) \log \frac{n}{\varepsilon}$
2	$\frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q-1}$	$O(q^2 \sqrt{n}) \log \frac{n}{\varepsilon}$
3	$t - 1 + \frac{t^{1-q-1}}{q-1}$	$O(q^2 \sqrt{n}) \log \frac{n}{\varepsilon}$
4	$\frac{t^2 - 1}{2} + \frac{t^{1-q} - 1}{q(q-1)} - \frac{q-1}{q}(t-1)$	$O(q \sqrt{n}) \log \frac{n}{\varepsilon}$
5	$\frac{t^2 - 1}{2} + \frac{e^t - e}{e}$	$O(q \sqrt{n}) \log \frac{n}{\varepsilon}$
6	$\frac{t^2 - 1}{2} - \int_1^t e^{\xi} d\xi$	$O(q \sqrt{n}) \log \frac{n}{\varepsilon}$

Table 3

The above table shows that the small-update methods based on listed kernel functions all have the same complexity, namely

$$O\left(\sqrt{n} \log \frac{n}{\varepsilon}\right). \quad (4-30)$$

This is up till now the best iteration bound for IPMs solving LP problems.

Historically most of the IPMs were based on the logarithmic kernel function. Notice that the gap between theoretical complexity of short-step methods and large-step methods is significant; the short step methods have much better theoretical complexity. However, in practical implementations large-step methods work better. This discrepancy was one of the motivations to consider other kernel functions in hopes to find the kernel function which would not have a gap or the gap would be smaller. As we can see, this goal has been achieved; for cases 2 and 4 the gap is much smaller because for these kernel functions large step method has much better complexity than for the classical logarithmic kernel function. This is one of the main achievements of considering different classes of kernel functions.

## CHAPTER 5

### NUMERICAL RESULTS

The Generic IPM described in the Figure 1 was implemented in MATLAB 7.6.0 with the class of kernel functions described by formula (3-1)

$$\psi_{p,q}(t) = \begin{cases} \frac{t^{p+1} - 1}{p+1} + \frac{t^{1-q} - 1}{q-1}, & t > 0, p \in [0,1], q > 1 \\ \frac{t^{p+1} - 1}{p+1} - \log t, & t > 0, p \in [0,1] \end{cases}$$

This imply that there are two implementations of the algorithm, one for the classical logarithmic kernel function

$$\psi(t) = \frac{t^2 - 1}{2} - \log t \quad (5-1)$$

and one for the kernel function with parameters  $p$  and  $q$

$$\psi_{p,q}(t) = \frac{t^{p+1} - 1}{p+1} + \frac{t^{1-q} - 1}{q-1}, \quad t > 0, p \in [0,1], q > 1 \quad (5-2)$$

We call the first implementation “Classical Method” and the second implementation “New Method”. Both codes are listed in the Appendix A.

The algorithm was tested on several examples with different sizes ranging from very small to moderate size problems. The data was entered in some cases “by hand” and for the others they were generated randomly.

**Example:** Consider the following simple LP model.

$$\begin{aligned}
 \max \quad & 2x_1 + 2x_2 \\
 \text{s.t.} \quad & x_1 + x_2 \leq 3 \\
 & x_1 \geq 0, x_2 \geq 0.
 \end{aligned}$$

It is easy to see that this problem has infinitely many optimal solutions, they are all the points on the segment  $[(1,0);(0,1)]$  and the optimal value is  $Z = 6$ . The problem was solved by New Method with  $p = 0.8$ ,  $q = 1.2$  with accuracy parameter  $\varepsilon = 0.001$ . It took unusually many iterations (57), however algorithm steadily converged to the expected result  $(x_1, x_2) = (1.5, 1.5)$ .

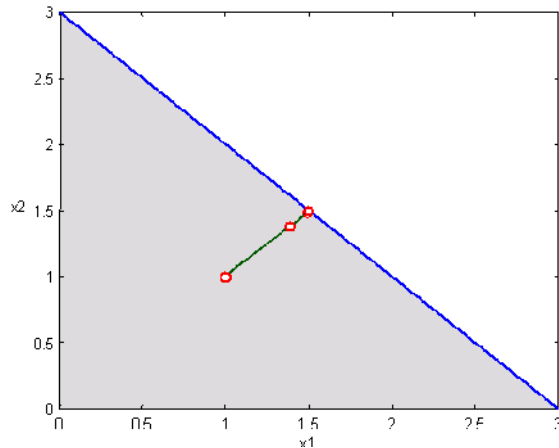
Numerical results for Example

x			y	s		
1	1	1	0	1	1	1
1.383	1.383	0.234	- 0.823 4	0.1	0.1	1.248 9
1.488 3	1.488 3	0.023 4	- 1.602 3	0.065 6	0.065 6	1.756 7
....	....	.....	.....	.....	.....	..... ...
1.499 9	1.499 9	0.000 2	- 2.000 2	0.000 2	0.000 2	2.000 2

Table 4

Objective function value  $Z = -5.9997$

This example also illustrates an important feature of the interior-point methods that distinguish them from simplex-type methods; that in the case of infinitely many optimal solutions they converge to the center of the optimal set rather than to the vertex. The graphical illustration of the above example with several first iterations is given below.



IPM Based on generic kernel function

Figure 3

Next, the algorithm was examined on the set of 200 randomly generated “small” problems for with sizes less than 10. The average number of iteration and CPU time is given in the table below.

Classical method		New method	
Average Number Of Iterations	Time	Average Number Of Iterations	Time
24.6	0.0362	40.7	0.0448

Numerical results for randomly generated “small problems” with dimension less than 10

Table 5

Next, the algorithm was examined on the set of 200 randomly generated “moderate size” problems of the size 200x300. The average number of iteration and CPU time is given in the table below.

Classical method		New method	
Average Number Of Iterations	Time	Average Number Of Iterations	Time
35.81	2.6626	40	2.8812

Numerical results for randomly generated problems with dimension 200x300

Table 6

The algorithm was then applied to the randomly generated problems of the bigger size 400x700. The result is given in the table below

Classical method		New method	
Average Number Of Iterations	Time	Average Number Of Iterations	Time
57	35.8048	41	25.4030

Numerical results for randomly generated problem with dimension 400x700

Table 7

Results summarized in tables seem to suggest that Classical Methods works slightly better for the problems of the smaller size while, as the dimension of the problem increases, the New Method becomes better. Another feature of the IPM is also visible from these examples and



that is that the number of iterations does not increase significantly with the increase in the size of the problem.

In the sequel the New Method was examined on the set of 200 randomly generated “small” problems with sizes less than 10 and for different values of parameters  $p$  and  $q$ . The results are given in the Table below.

p=0.7,q=1.3		p=0.6,q=1.4	
Iteration	Time	Iteration	Time
40.6	0.0673	41.6	0.0648
p=0.9,q=1.1		p=0.8,q=1.2	
Iteration	Time	Iteration	Time
35.9	0.0526	37.2	0.0603

Table 8: More Numerical Results

Next, the New Method was examined on the set of 20 randomly generated problems of the size  $200 \times 300$  and for different values of parameters  $p$  and  $q$ . The results are given in the Table below.

Table 9: More Numerical Results

The table seems to suggest that for the problems of the smaller size the New Method works the best when  $p = 0.9$ ,  $q = 1.1$ , which is in line with theoretical expectation. However, for the problems of the higher dimension it is hard to make conclusion which combination of parameters works the best. Theory suggests that  $p = 1$ ,  $q = \log n$  where  $n$  is the number of variables gives the best complexity. However for the particular set of problems in the previous table, it seems that combination  $p = 0.7$ ,  $q = 1.3$  works the best.

Better implementation and more testing is necessary for more definite conclusions. However, that was not the intention of the thesis. The goal was to make the basic implementations that illustrates theoretical concepts discussed in the thesis. Even on this basic level the implementation of the New Method shows the potential to work well. Of course, with more sophisticated implementation the performance can be further improved.

## CHAPTER 6

### CONCLUSION

In this thesis the Interior – Point Method (IPM) for Linear Programming problem (LP) that is based on the generic kernel function is considered. The algorithm is described in Chapter 2.

In Chapter 3 the complexity (in terms of iteration bounds) of the algorithm is analyzed for a class of kernel functions defined by (3-1). This class is fairly general; it includes classical logarithmic kernel function, prototype self-regular kernel function as well as non-self-regular functions, thus it serves as a unifying frame for the analysis of IPMs. Two versions of the IPMs are considered, the short-step algorithms where barrier parameter  $\theta$  depends on the size of the problem and long-step algorithms where barrier parameter is a fixed constant  $\theta \in (0,1)$ .

Historically most of the IPMs were based on the logarithmic kernel function. Notice that the gap between theoretical complexity of short-step methods and large-step methods is significant; the short step methods have much better theoretical complexity. However, in practical implementations large-step methods work better. This discrepancy was one of the motivations to consider other kernel functions in hopes to find the kernel function which would not have a gap or the gap would be smaller. As we can see this goal has been achieved; for kernel functions 2 and 4, the gap is much smaller than for the classical logarithmic kernel function. In addition, the complexity results that are obtained match the best known complexity results for these methods. This chapter is mostly based on the paper (Bai, Y., et al., 2008) with the addition of most of the proofs that were omitted in the paper.

The main contribution of the thesis is contained in Chapter 4. The detailed complexity analysis of the IPM that was provided in Chapter 3 for kernel function (3-1) is summarized and the analysis of the algorithm was performed for three additional kernel functions (4-1) – (4-3). For one of them we again matched the best known complexity results for the large-step methods and for the other two the complexity is slightly weaker, however still significantly improved in comparison with classical logarithmic kernel function.

The IPM that is theoretically analyzed in Chapter 3 is implemented in Chapter 5 for the classical logarithmic kernel function (Classical Method) and for the parametric kernel function (New Method) both described in (3-1). Although the implementation is on the basic level, it shows potential for a good performance of IPM based on kernel function different than classical logarithmic kernel function on which most of the commercial codes are based. The preliminary calculations seem to indicate that IPM with classical kernel logarithmic function perform better on problems of the smaller size while for larger problems the New Method seems to work slightly better. Also, based on the preliminary numerical tests it is hard to make conclusion which combination of parameters  $p$  and  $q$  in (3-1) works the best. Better implementation and more numerical testing would be necessary to draw more definite conclusions.

However, that was not the goal of the thesis, the goal was to show that IPM with kernel functions different than classical logarithmic kernel function can have best known theoretical complexity and to show that they have potential for practical implementations.

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## APPENDICES

### APPENDIX A

#### MatLab codes for the Classical Method

```

function [i,xx,yy,ss,z,d]=IpmClassical(A, b, c, epsilon)
%
% input tau, epsilon, theta
%
% sizes: A--m*n, b--m, s--n, x--n, y--m, c--n
%
% To call this function, please set up the problem by defining A, b and
% c. Or load the example problem.

[m,n]=size(A);
x=1*ones(n,1); s=1*ones(n,1); y=zeros(m,1); mu=x'*s/n;
rd=c-A'*y-s;
rp=b-A*x;

i=1;
xx(i,:)=x;
yy(i,:)=y;
ss(i,:)=s;

while norm(rd)>epsilon||norm(rp)>epsilon||n*mu > epsilon
    i=i+1;
    [dx,ds,dy]=SolvesystemClassical(A,b,c,x,y,s,mu);
    inds=find(ds<0);indx=find(dx<0);
    alpha=0.9*min(abs([1;s(inds)]./ds(inds);x(indx)]./dx(indx)));
    x=x+alpha*dx;
    y=y+alpha*dy;
    s=s+alpha*ds;
    mu=min(x'*s/n,0.9*mu);
    rd=c-A'*y-s;
    rp=b-A*x;

    xx(i,:)=x;
    yy(i,:)=y;
    ss(i,:)=s;
    d(i,:)=dx';
    if i>200 || alpha<1e-11
        i;
        break;
    end
end
z=x'*c;
%[i,z]
end

```

```

function [dx,ds,dy]=SolvesystemClassical(A,b,c,x,y,s,mu)
% This function solves the following system
%      A*dx = 0
%      A^T*dy + ds = 0
%      S*dx + x*ds = - mu*v.*grad(Psi(v))
%
% Psi(v)=sum((v.^(p+1)-1)/(p+1)+(v.^(1-q)-1)/(q-1));
% grad(v)=v.^p-v.^q;

    gama=.1;
    X=diag(x);
    S=diag(s);
    S_inv=diag(1./s);
    rd=c-A'*y-s;
    rp=b-A*x;

    M=A*S_inv*X*A';
    r=b+A*S_inv*(X*rd-gama*mu*ones(size(A,2),1));

    dy=M\r;
    ds=rd-A'*dy;
    dx=-x+S_inv*(gama*mu*ones(size(A,2),1)-X*ds);

end

```

### MatLab codes for the New Method

```

function [i,xx,yy,ss,z,d]=IpmNew(A, b, c, epsilon)
%
%
% input tau, epsilon, theta
% v>0, 0<=p<=1, q>1, tau>1
% sizes: A--m*n, b--m, s--n, x--n, y--m, c--n
%
% To call this function, please set up the problem by defining A, b and
% c. Or load the example problem.

p=1-0.2;
q=1+0.2;
theta=0.1;
%tau=1.5;
    [m,n]=size(A);
    x=ones(n,1); s=ones(n,1); y=zeros(m,1); mu=x'*s/n;
    rd=c-A'*y-s;
    rp=b-A*x;
    i=1;
    xx(i,:)=x;
    yy(i,:)=y;
    ss(i,:)=s;
    while norm(rd)>epsilon||norm(rp)>epsilon||n*mu > epsilon
        i=i+1;
        v=sqrt(x.*s./mu);
        [dx,ds,dy]=SolvesystemNew(A,b,c,x,y,s,mu,v,p,q);
    end

```

```

delta=1/2*sqrt(sum((v.^p-1./v.^q).^2));
alpha=1/((p+q)*(1-4*abs(delta))^(1+1/q));
alpha=abs(alpha);
inds=find(ds<0);indx=find(dx<0);
alpha=0.9*min(abs([1;s(inds)./ds(inds);x(indx)./dx(indx)]));
x=x+alpha*dx;
y=y+alpha*dy;
s=s+alpha*ds;

rd=c-A'*y-s;
rp=b-A*x;
mu=(1-theta)*mu;
xx(i,:)=x';
yy(i,:)=y';
ss(i,:)=s';
d(i,:)=dx';
if i>200 || alpha<1e-10
    i;
    break;
end
end
end
z=x'*c;
%[i,z]
end

```

```

function [dx,ds,dy]=SolvesystemNew(A,b,c,x,y,s,mu,v,p,q)
% This function solves the following system
%      A*dx = 0
%      A^T*dy + ds = 0
%      S*dx + x*ds = - mu*v.*grad(Psi(v))
%
% Psi(v)=sum((v.^(p+1)-1)/(p+1)+(v.^(1-q)-1)/(q-1));
% grad(v)=v.^p-v.^q;
gama=.1;
X=diag(x);
S=diag(s);
S_inv=diag(1./s);
rd=c-A'*y-s;
rp=b-A*x;

M=A*S_inv*X*A';
r=A*S_inv*(X*rd+mu*v.*(v.^p-gama*v.^(-q)))+rp;

dy=M\r;
ds=rd-A'*dy;
dx=-S_inv*(X*ds+mu*v.*(v.^p-gama*v.^(-q)));
end

```

### Problem Generator

```

function R=mytest(NO,m,n)
% This function solve the random systems of dimention m*n

```



```

% input m,n are the dimension of A
% input NO is the number of the iterations
%
% MYTEST(); will apply both ipm methods to a random matrix with random
% dimension m*n, where m,n are positive integers less than 10
%
% MYTEST(N) will iterate 'MYTEST()' N times.
%
% MYTEST(N,m,n) will iterately apply the methods to N random problems
% with dimension m*n.

    if ~exist('NO')
        NO=1;
    end
    m_ne=0;n_ne=0;
    if ~exist('m')
        m_ne=1;
    end
    if ~exist('n')
        n_ne=1;
    end

    dim=zeros(NO,2);
    k=1;
    while k<=NO
        if m_ne
            m=fix(rand(1)*10);
        end
        if n_ne
            n=fix(rand(1)*10);
        end
        A=rand(m,n);b=rand(m,1);c=rand(n,1);
        if rank(A)<min(m,n) || min(m,n)==0
            %fprintf('rank(A)<min(m,)\n');
            if m==0 && ~m_ne
                display('STOP, m=0');
                return;
            elseif n==0 && ~n_ne
                display('STOP, n=0');
                return;
            end
            continue;
        end
        tic;
        [i1(k),x1,yy1,ss1,z1(k)]=IpmNew(A, b, c, 0.01);
        t1(k)=toc;
        tic;
        [i2(k),xx2,yy2,ss2,z2(k)]=IpmClassical(A, b, c, 0.01);
        t2(k)=toc;
        dim(k,1)=m;dim(k,2)=n;
        k=k+1;
    end
    R = [dim(:,1),dim(:,2),i1',z1',t1',i2',z2',t2'];
    myfun@;
end

```

## Organization of output

```
function myfun@

    fprintf('          New method          Classical method \n');
    fprintf('  idx dim(m*n)  iteration Opt      time      iteration Opt
time\n');
    fprintf('-----\n');
    --\n';
    for k=1:size(R,1)
        fprintf('   %3d %3d%3d  %3d  %6.4f  %4.4f      %3d  %6.4f
%4.4f\n', ...
                k,R(k,:));
    end
    R_ave=sum(R)/size(R,1);
    fprintf('-----\n');
    --\n';
    fprintf('          New method          Classical method \n');
    fprintf('          dim(m*n)  iteration time      iteration time\n');
    fprintf(' average %3.1f %3.1f      %3.1f  %4.4f      %3.1f
%4.4f\n', ...
            R_ave([1,2,3,5,6,8]));
    fprintf('-----\n');
    --\n';

end
```

**APENDIX B****MatLab code Example 5.1**

```
max 3 * x_1 + 5 * x_2
```

```
s.t.
```

```
    x_1 +      <=4
```

```
        2 * x_2 <=12
```

```
    3 * x_1 + 2 * x_2 <=18
```

```
        x_1, x_2 >= 0
```

```
A =
```

```
    1  0  1  0  0
```

```
    0  2  0  1  0
```

```
    3  1  0  0  1
```

```
(change to min-problem)
```

```
c =
```

```
   -3
```

```
   -5
```

```
    0
```

```
    0
```

```
    0
```

```
b =
```

```
    4
```

```
   12
```

```
   18
```

```
>> [i,x,y,s,z]=lpmClassical(A, b, c, epsilon)
```

```
i =
```

```
7
```

```
z =
```

```
-41.9805
```

```
>> [i,x,y,s,z]=lpmNew(A, b, c, epsilon)
```

```
i =
```

```
60
```

```
z =
```

```
-41.9993
```

```
>> [i,x,y,s,z]=lpmClassical(A, b, c, epsilon)
```

```
i =
```

```
7
```

```
x =
```

```
1.0000 1.0000 1.0000 1.0000 1.0000
```

```
1.4437 1.7530 0.8831 0.9645 1.0398
```

```
2.0334 3.0422 0.7370 0.3825 0.8654
```

```
2.9815 4.6039 0.4101 0.0542 0.4968
```

```
3.8206 5.7689 0.0779 0.0054 0.1095
```

```
3.9803 5.9762 0.0095 0.0020 0.0169
```

```
3.9977 5.9975 0.0013 0.0005 0.0028
```

y =

```
    0    0    0
-0.1333 -0.0519  0.0235
-0.2783 -0.5925 -0.1361
-0.5626 -1.3849 -0.4117
-0.8317 -2.0210 -0.6562
-0.9001 -2.1384 -0.6935
-0.9363 -2.1549 -0.6873
```

s =

```
1.0000  1.0000  1.0000  1.0000  1.0000
0.4093  0.1000  0.9699  0.8885  0.8131
0.1458  0.0100  0.8931  1.2073  0.7509
0.0146  0.0068  0.8669  1.6891  0.7159
0.0032  0.0026  0.8825  2.0717  0.7069
0.0011  0.0007  0.9052  2.1434  0.6986
0.0003  0.0002  0.9368  2.1554  0.6878
```

z =

```
-41.9805
```

```
>> [i,x,y,s,z]=lpmNew(A, b, c, epsilon)
```

i =

```
60
```

z =

```
-41.9993
```

x =

1.0000	1.0000	1.0000	1.0000	1.0000
1.4437	1.7530	0.8831	0.9645	1.0398
2.0409	3.0601	0.7365	0.3780	0.8702
3.0192	4.6730	0.4016	0.0474	0.5044
3.8797	5.8542	0.0624	0.0308	0.1302
3.9559	5.9723	0.0383	0.0293	0.1223
3.9623	5.9859	0.0372	0.0256	0.1236
3.9669	5.9883	0.0330	0.0232	0.1105
3.9702	5.9896	0.0297	0.0208	0.0997
3.9732	5.9906	0.0268	0.0188	0.0897
3.9759	5.9916	0.0241	0.0169	0.0808
3.9783	5.9924	0.0217	0.0152	0.0727
3.9805	5.9932	0.0195	0.0137	0.0655
3.9824	5.9938	0.0176	0.0123	0.0589
3.9842	5.9945	0.0158	0.0111	0.0530
3.9857	5.9950	0.0143	0.0100	0.0478
3.9872	5.9955	0.0128	0.0090	0.0430
3.9884	5.9960	0.0116	0.0081	0.0387
3.9896	5.9964	0.0104	0.0073	0.0348
3.9906	5.9967	0.0094	0.0065	0.0314
3.9916	5.9971	0.0084	0.0059	0.0282
3.9924	5.9973	0.0076	0.0053	0.0254
3.9932	5.9976	0.0068	0.0048	0.0229
3.9939	5.9979	0.0061	0.0043	0.0206
3.9945	5.9981	0.0055	0.0039	0.0185
3.9950	5.9983	0.0050	0.0035	0.0167

3.9955	5.9984	0.0045	0.0031	0.0150
3.9960	5.9986	0.0040	0.0028	0.0135
3.9964	5.9987	0.0036	0.0025	0.0122
3.9967	5.9989	0.0033	0.0023	0.0109
3.9971	5.9990	0.0029	0.0021	0.0098
3.9974	5.9991	0.0026	0.0018	0.0089
3.9976	5.9992	0.0024	0.0017	0.0080
3.9979	5.9993	0.0021	0.0015	0.0072
3.9981	5.9993	0.0019	0.0013	0.0065
3.9983	5.9994	0.0017	0.0012	0.0058
3.9984	5.9995	0.0016	0.0011	0.0052
3.9986	5.9995	0.0014	0.0010	0.0047
3.9987	5.9996	0.0013	0.0009	0.0042
3.9989	5.9996	0.0011	0.0008	0.0038
3.9990	5.9996	0.0010	0.0007	0.0034
3.9991	5.9997	0.0009	0.0006	0.0031
3.9992	5.9997	0.0008	0.0006	0.0028
3.9993	5.9997	0.0007	0.0005	0.0025
3.9993	5.9998	0.0007	0.0005	0.0023
3.9994	5.9998	0.0006	0.0004	0.0020
3.9995	5.9998	0.0005	0.0004	0.0018
3.9995	5.9998	0.0005	0.0003	0.0016
3.9996	5.9998	0.0004	0.0003	0.0015
3.9996	5.9999	0.0004	0.0003	0.0013
3.9996	5.9999	0.0004	0.0002	0.0012
3.9997	5.9999	0.0003	0.0002	0.0011
3.9997	5.9999	0.0003	0.0002	0.0010

3.9997	5.9999	0.0003	0.0002	0.0009
3.9998	5.9999	0.0002	0.0002	0.0008
3.9998	5.9999	0.0002	0.0001	0.0007
3.9998	5.9999	0.0002	0.0001	0.0006
3.9998	5.9999	0.0002	0.0001	0.0006
3.9998	5.9999	0.0002	0.0001	0.0005
3.9999	6.0000	0.0001	0.0001	0.0005

y =

0	0	0
-0.1333	-0.0519	0.0235
-0.2846	-0.6023	-0.1375
-0.5916	-1.4288	-0.4214
-0.9196	-2.0876	-0.6622
-1.4299	-2.2344	-0.5248
-1.5993	-2.2683	-0.4714
-1.5888	-2.2669	-0.4748
-1.5890	-2.2668	-0.4743
-1.5884	-2.2665	-0.4741
-1.5878	-2.2662	-0.4739
-1.5873	-2.2660	-0.4738
-1.5868	-2.2658	-0.4737
-1.5864	-2.2656	-0.4735
-1.5860	-2.2654	-0.4734
-1.5857	-2.2652	-0.4733
-1.5854	-2.2651	-0.4732
-1.5851	-2.2649	-0.4732



-1.5848 -2.2648 -0.4731  
-1.5846 -2.2647 -0.4730  
-1.5844 -2.2646 -0.4730  
-1.5842 -2.2645 -0.4729  
-1.5841 -2.2645 -0.4729  
-1.5839 -2.2644 -0.4728  
-1.5838 -2.2643 -0.4728  
-1.5837 -2.2643 -0.4728  
-1.5835 -2.2642 -0.4727  
-1.5834 -2.2642 -0.4727  
-1.5834 -2.2641 -0.4727  
-1.5833 -2.2641 -0.4727  
-1.5832 -2.2641 -0.4727  
-1.5831 -2.2640 -0.4726  
-1.5831 -2.2640 -0.4726  
-1.5830 -2.2640 -0.4726  
-1.5830 -2.2640 -0.4726  
-1.5830 -2.2639 -0.4726  
-1.5829 -2.2639 -0.4726  
-1.5829 -2.2639 -0.4726  
-1.5828 -2.2639 -0.4726  
-1.5828 -2.2639 -0.4725  
-1.5828 -2.2639 -0.4725  
-1.5828 -2.2639 -0.4725  
-1.5828 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725

-1.5827 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725  
-1.5827 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725  
-1.5826 -2.2638 -0.4725

s =

1.0000	1.0000	1.0000	1.0000	1.0000
0.4093	0.1000	0.9699	0.8885	0.8131
0.1425	0.0100	0.8960	1.2136	0.7489
0.0143	0.0168	0.8812	1.7184	0.7110
0.0220	0.0112	0.9486	2.1166	0.6911
0.0160	0.0110	1.4328	2.2373	0.5277
0.0148	0.0097	1.5996	2.2686	0.4717
0.0132	0.0088	1.5888	2.2669	0.4748
0.0119	0.0079	1.5890	2.2668	0.4743
0.0107	0.0071	1.5884	2.2665	0.4741

0.0096	0.0064	1.5878	2.2662	0.4739
0.0087	0.0057	1.5873	2.2660	0.4738
0.0078	0.0052	1.5868	2.2658	0.4737
0.0070	0.0047	1.5864	2.2656	0.4735
0.0063	0.0042	1.5860	2.2654	0.4734
0.0057	0.0038	1.5857	2.2652	0.4733
0.0051	0.0034	1.5854	2.2651	0.4732
0.0046	0.0031	1.5851	2.2649	0.4732
0.0041	0.0027	1.5848	2.2648	0.4731
0.0037	0.0025	1.5846	2.2647	0.4730
0.0033	0.0022	1.5844	2.2646	0.4730
0.0030	0.0020	1.5842	2.2645	0.4729
0.0027	0.0018	1.5841	2.2645	0.4729
0.0024	0.0016	1.5839	2.2644	0.4728
0.0022	0.0015	1.5838	2.2643	0.4728
0.0020	0.0013	1.5837	2.2643	0.4728
0.0018	0.0012	1.5835	2.2642	0.4727
0.0016	0.0011	1.5834	2.2642	0.4727
0.0014	0.0010	1.5834	2.2641	0.4727
0.0013	0.0009	1.5833	2.2641	0.4727
0.0012	0.0008	1.5832	2.2641	0.4727
0.0010	0.0007	1.5831	2.2640	0.4726
0.0009	0.0006	1.5831	2.2640	0.4726
0.0008	0.0006	1.5830	2.2640	0.4726
0.0008	0.0005	1.5830	2.2640	0.4726
0.0007	0.0005	1.5830	2.2639	0.4726
0.0006	0.0004	1.5829	2.2639	0.4726

0.0006	0.0004	1.5829	2.2639	0.4726
0.0005	0.0003	1.5828	2.2639	0.4726
0.0005	0.0003	1.5828	2.2639	0.4725
0.0004	0.0003	1.5828	2.2639	0.4725
0.0004	0.0002	1.5828	2.2639	0.4725
0.0003	0.0002	1.5828	2.2638	0.4725
0.0003	0.0002	1.5827	2.2638	0.4725
0.0003	0.0002	1.5827	2.2638	0.4725
0.0002	0.0002	1.5827	2.2638	0.4725
0.0002	0.0001	1.5827	2.2638	0.4725
0.0002	0.0001	1.5827	2.2638	0.4725
0.0002	0.0001	1.5827	2.2638	0.4725
0.0002	0.0001	1.5827	2.2638	0.4725
0.0001	0.0001	1.5827	2.2638	0.4725
0.0001	0.0001	1.5826	2.2638	0.4725
0.0001	0.0001	1.5826	2.2638	0.4725
0.0001	0.0001	1.5826	2.2638	0.4725
0.0001	0.0001	1.5826	2.2638	0.4725
0.0001	0.0001	1.5826	2.2638	0.4725
0.0001	0.0001	1.5826	2.2638	0.4725
0.0001	0.0000	1.5826	2.2638	0.4725
0.0001	0.0000	1.5826	2.2638	0.4725
0.0001	0.0000	1.5826	2.2638	0.4725

z =

-41.9993

>>

### MatLab code for Table 5.2

random problems with random dimentions less that 10

	New method				Classical method			
idx	dim(m*n)	iteration	Opt	time	iteration	Opt	time	
-----								
36	6 5 12	1.4119	0.0056		13	1.4164	0.0396	
37	6 7 15	0.9627	0.0065		16	0.9627	0.0046	
38	6 8 17	0.7974	0.0072		24	0.7974	0.0073	
39	4 8 17	0.7864	0.0336		21	0.7864	0.0061	
40	8 3 15	0.4944	0.0070		38	0.3653	0.0140	
41	4 9 72	1.0654	0.0757		27	1.0668	0.0086	
42	5 8 18	0.4801	0.0076		24	0.4801	0.0072	
43	8 9 14	1.3682	0.0070		19	1.3682	0.0064	
44	2 2 13	0.7773	0.0042		14	0.7773	0.0033	
45	6 5 13	0.2225	0.0059		14	0.2229	0.0049	
46	8 6 14	1.6662	0.0068		20	1.6114	0.0268	
47	3 6 17	0.3840	0.0069		30	NaN	0.0101	
48	3 7 64	0.5853	0.0260		7	0.5886	0.0018	
49	3 3 12	0.8750	0.0044		14	0.8750	0.0038	
50	2 7 64	0.6310	0.0458		7	0.6367	0.0017	
51	7 7 12	1.7618	0.0047		13	1.7618	0.0036	
52	2 7 64	0.1590	0.0244		7	0.1647	0.0018	

	New method		Classical method	
dim(m*n)	iteration	time	iteration	time
-----				
average	4.9	5.4	40.7	0.0448
			24.6	0.0362
-----				

	New method				Classical method			
idx	dim(m*n)	iteration	Opt	time	iteration	Opt	time	
1	9 1 45	NaN	0.0681	30	NaN	0.0326		
2	9 1 45	NaN	0.0674	45	NaN	0.7326		
3	7 7 12	3.1727	0.0053	14	3.1727	0.0042		
4	7 7 12	2.0364	0.0050	13	2.0364	0.0037		
5	5 4 14	0.7954	0.0675	16	0.7257	0.0059		
6	6 3 49	0.4286	0.0243	92	0.4560	0.1107		
7	6 1 45	NaN	0.1613	45	NaN	0.0612		
8	3 9 66	0.2055	0.0284	9	0.1994	0.0024		
9	3 1 45	NaN	0.1071	36	NaN	0.0133		
10	7 8 13	2.3593	0.0057	14	2.3593	0.0042		
11	2 7 64	0.3550	0.0246	8	0.3596	0.0019		
12	6 7 14	1.7947	0.0055	15	1.7947	0.0043		
13	2 6 62	NaN	0.0637	27	NaN	0.0096		
14	3 2 201	0.4458	0.3814	71	0.1206	0.1475		
15	7 8 14	2.3068	0.0059	16	2.3068	0.0258		
16	7 5 28	0.4972	0.0138	19	0.5465	0.0067		
17	4 9 17	1.2697	0.0076	18	1.2697	0.0053		
18	2 7 64	NaN	0.0814	17	NaN	0.0044		
19	1 4 58	0.0069	0.0650	6	0.0106	0.0015		
20	5 5 12	2.0070	0.0045	13	2.0070	0.0036		
21	1 5 60	0.0321	0.0245	6	0.0387	0.0026		
22	1 2 52	0.2947	0.0739	5	0.2976	0.0012		
23	9 6 23	0.7775	0.0118	52	0.5957	0.0617		

24	5	3	18	1.7080	0.0088	33	1.2630	0.0476
25	4	3	31	0.3020	0.0528	28	0.3146	0.0426
26	5	8	16	2.4185	0.0069	16	2.4185	0.0046
27	4	7	16	0.2408	0.0064	20	0.2408	0.0059
28	9	8	17	1.8185	0.0091	24	1.2417	0.0096
29	1	2	52	0.2448	0.0208	5	0.2478	0.0012
30	3	4	14	1.0822	0.0051	17	1.0822	0.0046
31	4	5	14	1.4672	0.0054	16	1.4672	0.0044
32	4	2	201	0.0760	0.3930	59	0.0470	0.1529
33	5	8	25	0.5901	0.0115	28	0.5901	0.0087
34	1	9	66	0.0042	0.0293	7	0.0082	0.0018
35	3	7	15	0.8637	0.0060	201	0.2481	0.3810
36	6	5	12	1.4119	0.0056	13	1.4164	0.0396
37	6	7	15	0.9627	0.0065	16	0.9627	0.0046
38	6	8	17	0.7974	0.0072	24	0.7974	0.0073
39	4	8	17	0.7864	0.0336	21	0.7864	0.0061
40	8	3	15	0.4944	0.0070	38	0.3653	0.0140
41	4	9	72	1.0654	0.0757	27	1.0668	0.0086
42	5	8	18	0.4801	0.0076	24	0.4801	0.0072
43	8	9	14	1.3682	0.0070	19	1.3682	0.0064
44	2	2	13	0.7773	0.0042	14	0.7773	0.0033
45	6	5	13	0.2225	0.0059	14	0.2229	0.0049
46	8	6	14	1.6662	0.0068	20	1.6114	0.0268
47	3	6	17	0.3840	0.0069	30	NaN	0.0101
48	3	7	64	0.5853	0.0260	7	0.5886	0.0018
49	3	3	12	0.8750	0.0044	14	0.8750	0.0038
50	2	7	64	0.6310	0.0458	7	0.6367	0.0017

51	7	7	12	1.7618	0.0047	13	1.7618	0.0036
52	2	7	64	0.1590	0.0244	7	0.1647	0.0018
53	4	5	15	3.6371	0.0059	16	3.6371	0.0455
54	9	3	42	0.4919	0.0216	41	0.7683	0.0158
55	7	1	45	NaN	0.0833	28	NaN	0.0224
56	5	4	20	1.4544	0.0331	14	1.6696	0.0063
57	2	1	45	NaN	0.0891	45	NaN	0.0296
58	7	5	41	1.6545	0.0388	33	0.8089	0.0497
59	1	7	64	0.1189	0.0272	8	0.1203	0.0020
60	9	9	12	3.7354	0.0071	15	3.7354	0.0049
61	6	8	16	1.3473	0.0067	17	1.3473	0.0051
62	2	8	65	0.1011	0.0256	7	0.1045	0.0018
63	2	6	16	0.5000	0.0055	19	0.5000	0.0048
64	2	6	62	0.4073	0.0233	6	0.4123	0.0014
65	4	6	15	1.0340	0.0374	19	1.0340	0.0054
66	3	9	66	0.9751	0.0292	8	0.9772	0.0021
67	5	5	13	0.9247	0.0050	14	0.9247	0.0039
68	6	6	13	1.6887	0.0051	14	1.6887	0.0452
69	2	2	13	0.1492	0.0045	15	0.1492	0.0036
70	3	2	171	NaN	0.1243	54	2.9584	0.0724
71	7	2	201	0.2807	0.2557	66	0.0886	0.0771
72	5	2	201	0.6778	0.4181	75	1.0557	0.1723
73	2	9	66	0.0872	0.0264	7	0.0909	0.0018
74	5	7	14	1.4273	0.0056	15	1.4273	0.0040
75	9	9	12	4.2331	0.0055	14	4.2331	0.0048
76	4	4	13	0.7073	0.0047	14	0.7073	0.0040
77	2	9	66	0.4177	0.0272	7	0.4248	0.0019



78	7	9	15	2.3273	0.0066	16	2.3273	0.0050
79	7	5	15	0.6883	0.0382	13	0.8011	0.0524
80	3	5	17	0.7023	0.0066	19	0.7023	0.0053
81	8	5	17	1.6375	0.0082	87	0.5440	0.1686
82	5	6	17	0.5266	0.0070	22	0.5266	0.0065
83	6	9	16	1.1541	0.0073	19	1.1541	0.0059
84	1	2	52	0.9460	0.0199	5	0.9489	0.0012
85	3	9	20	1.4204	0.0087	39	NaN	0.0153
86	9	4	29	0.1871	0.0378	32	0.6971	0.0615
87	9	5	37	0.9494	0.0764	32	1.3586	0.0558
88	4	9	18	0.5192	0.0078	23	0.5192	0.0411
89	9	5	41	0.5489	0.0475	18	1.5730	0.0528
90	6	4	58	NaN	0.1482	58	NaN	0.1727
91	3	5	15	0.6959	0.0058	20	0.6959	0.0055
92	7	5	14	2.9338	0.0065	14	2.3208	0.0053
93	3	5	60	0.4861	0.0749	7	0.4866	0.0020
94	3	2	196	4.5525	0.3504	146	1.8557	0.3399
95	2	8	65	0.6134	0.0258	7	0.6152	0.0018
96	6	8	15	1.6926	0.0062	19	1.6926	0.0058
97	3	6	15	0.8336	0.0057	17	0.8336	0.0047
98	4	3	56	NaN	0.0830	14	1.4102	0.0045
99	7	2	201	0.0365	0.4461	83	1.0238	0.2068
100	6	1	45	NaN	0.1986	45	NaN	0.0328
101	1	4	58	0.6234	0.0448	6	0.6307	0.0014
102	7	5	14	1.1884	0.0070	38	0.4592	0.1011
103	2	5	60	0.5702	0.0634	6	0.5733	0.0014
104	8	5	38	1.2510	0.0198	19	1.3811	0.0390

105	5	7	16	0.7932	0.0068	19	0.7932	0.0267
106	2	3	56	0.6463	0.0198	5	0.6492	0.0012
107	7	9	18	1.2119	0.0081	24	1.2119	0.0078
108	4	2	116	0.4112	0.1974	116	0.3480	0.2755
109	2	4	58	0.3450	0.0323	6	0.3478	0.0014
110	8	6	47	0.8233	0.0480	25	0.8203	0.0931
111	1	5	60	0.0042	0.0245	7	0.0052	0.0017
112	5	7	19	2.2536	0.0077	25	2.2536	0.0074
113	7	7	13	1.4697	0.0053	14	1.4697	0.0044
114	9	2	120	0.4624	0.2295	201	0.8555	0.4208
115	9	8	12	1.9521	0.0063	13	1.9450	0.0872
116	1	1	45	0.4377	0.0147	5	0.4378	0.0011
117	6	3	40	1.3974	0.0493	60	0.0616	0.1541
118	2	5	60	0.0523	0.0223	6	0.0577	0.0014
119	4	5	16	1.3757	0.0060	17	1.3757	0.0046
120	9	8	13	2.8652	0.0067	15	3.0675	0.0062
121	8	8	11	5.0975	0.0048	15	5.0975	0.0047
122	2	2	13	0.3116	0.0042	39	NaN	0.0116
123	9	5	18	0.4133	0.0092	21	0.5460	0.0381
124	5	4	17	1.3629	0.0079	18	1.3382	0.0401
125	9	5	17	1.4464	0.0085	16	1.9442	0.0707
126	7	6	15	2.1528	0.0073	13	1.5332	0.0388
127	1	2	52	0.4030	0.0202	5	0.4060	0.0012
128	9	8	17	1.5577	0.0091	16	1.3553	0.0158
129	7	1	45	NaN	0.1823	36	NaN	0.0295
130	8	9	17	2.5650	0.0080	21	2.5650	0.0070
131	3	4	15	0.3709	0.0055	17	0.3709	0.0062

132	8	7	13	1.5864	0.0063	16	2.2049	0.0061
133	9	9	12	1.4834	0.0053	14	1.4834	0.0045
134	8	4	88	0.1456	0.1307	58	0.4504	0.1475
135	6	2	27	0.0678	0.0600	16	NaN	0.0094
136	2	9	66	0.4524	0.0270	8	0.4538	0.0019
137	1	7	64	0.1021	0.0279	7	0.1043	0.0017
138	1	6	62	0.0028	0.0263	7	0.0051	0.0017
139	7	7	12	3.4885	0.0052	13	3.4885	0.0037
140	3	4	15	0.2784	0.0056	17	0.2784	0.0048
141	7	9	14	3.1083	0.0062	15	3.1083	0.0047
142	1	9	66	0.0150	0.0301	7	0.0228	0.0018
143	8	8	12	3.1648	0.0050	13	3.1648	0.0041
144	3	8	16	0.6641	0.0651	17	0.6641	0.0048
145	6	2	127	NaN	0.1905	29	0.0231	0.0666
146	7	7	11	3.2350	0.0046	13	3.2350	0.0043
147	2	4	58	0.7658	0.0200	6	0.7670	0.0013
148	8	8	12	2.1261	0.0050	13	2.1261	0.0038
149	2	5	60	0.2748	0.0218	6	0.2799	0.0526
150	1	3	56	0.1017	0.0223	6	0.1030	0.0014
151	8	9	15	2.2372	0.0071	17	2.2372	0.0058
152	7	3	195	0.4054	0.3624	48	0.3538	0.1150
153	3	9	17	0.7720	0.0073	20	0.7720	0.0058
154	1	5	60	0.0085	0.0624	7	0.0096	0.0017
155	5	7	16	0.6305	0.0068	16	0.6305	0.0046
156	7	8	16	1.9168	0.0070	25	1.9168	0.0083
157	4	5	14	0.7501	0.0052	15	0.7501	0.0045
158	7	1	45	NaN	0.1248	16	NaN	0.0114

159	3	3	11	0.9672	0.0264	12	0.9672	0.0031
160	1	1	45	0.4010	0.0133	4	0.4008	0.0008
161	4	8	18	0.6044	0.0076	19	0.6044	0.0053
162	6	5	14	1.2959	0.0066	16	1.2533	0.0058
163	9	6	15	1.5190	0.0076	23	1.2027	0.0529
164	9	9	12	2.4810	0.0054	13	2.4810	0.0041
165	7	2	54	3.0844	0.0905	26	1.1829	0.0092
166	9	5	11	0.6774	0.0054	20	0.8400	0.0390
167	3	9	66	0.1178	0.0305	7	0.1240	0.0019
168	2	3	56	1.2827	0.0201	5	1.2850	0.0012
169	8	5	14	3.1274	0.0066	21	2.3811	0.0387
170	2	4	58	0.2706	0.0206	6	0.2715	0.0014
171	4	3	79	1.8914	0.1462	159	0.6958	0.3458
172	1	5	60	0.0769	0.0245	6	0.0821	0.0015
173	1	3	56	0.1561	0.0220	5	0.1636	0.0012
174	7	4	55	0.3958	0.0958	31	0.8488	0.0474
175	7	6	22	0.4917	0.0108	23	0.5037	0.0401
176	1	4	58	0.0114	0.0541	6	0.0139	0.0014
177	6	7	15	1.2088	0.0063	19	1.2088	0.0057
178	8	7	22	2.3937	0.0109	31	3.0084	0.1104
179	9	8	12	2.3249	0.0063	23	2.0587	0.0430
180	1	8	65	0.0225	0.0277	8	0.0236	0.0329
181	8	3	21	0.2064	0.0111	52	0.1950	0.0961
182	8	1	45	NaN	0.1603	45	NaN	0.0903
183	3	1	45	NaN	0.0892	45	NaN	0.0170
184	4	9	16	0.7594	0.0070	25	0.7594	0.0080
185	7	6	12	1.7165	0.0058	16	1.2626	0.0795

186	9	9	12	4.2814	0.0057	13	4.2814	0.0042
187	7	9	15	1.0139	0.0067	17	1.0139	0.0051
188	2	3	56	0.0452	0.0192	5	0.0469	0.0012
189	8	4	48	0.4489	0.0365	14	1.0610	0.0049
190	1	2	52	0.0018	0.0198	5	0.0047	0.0012
191	4	9	18	1.5071	0.0082	23	1.5071	0.0282
192	4	3	15	0.5991	0.0068	16	0.6997	0.0056
193	6	3	123	0.0110	0.2378	67	0.0438	0.1039
194	1	6	62	0.1704	0.0566	7	0.1724	0.0017
195	3	2	53	NaN	0.0634	17	NaN	0.0057
196	2	4	58	0.0721	0.0200	6	0.0768	0.0015
197	3	8	16	2.0542	0.0065	20	2.0542	0.0056
198	5	1	45	NaN	0.1366	45	NaN	0.0839
199	4	5	13	1.3755	0.0054	15	1.3755	0.0041
200	3	3	14	1.8529	0.0051	15	1.8529	0.0039

-----

	New method		Classical method			
	dim(m*n)	iteration	time	iteration	time	
average	4.9	5.4	40.7	0.0448	24.6	0.0362

-----

### MatLab code for Table 5.3

solve the random system with dimension 200\*300

```
m=200;n=300;
```

```
A=rand(m,n);
```

```
b=rand(m,1);
```

```
c=rand(n,1);
```

```
epsilon=0.01;
```

in new mehtod, p=0.8,q=1.2

New method				Classical method			
idx	iteration	Opt	time	iteration	Opt	time	
-----							
34	31	13.3783	2.1870	34	13.3783	2.4084	
35	33	13.5477	2.3660	38	13.5477	2.7975	
36	35	12.2685	2.4543	47	12.2685	3.6504	
37	34	12.1387	2.4273	42	12.1387	3.1664	
38	31	14.3202	2.1804	42	14.3202	3.1886	
39	33	11.7495	2.4815	30	11.7495	2.0452	
40	30	15.0938	2.1312	39	15.0938	2.8654	
41	40	15.4534	3.0469	44	15.4534	3.3165	
42	38	12.4233	2.8703	44	12.4233	3.2338	
43	66	12.7004	5.6847	40	12.7022	2.8478	
44	39	12.4386	2.9631	40	12.4386	2.8733	
45	33	13.5479	2.3851	42	13.5479	3.0937	
46	36	13.7716	2.6777	43	13.7716	3.1809	
47	33	12.7877	2.5047	35	12.7877	2.5006	
48	36	16.7568	2.7086	36	16.7568	2.5545	
49	38	15.3569	2.9702	34	15.3569	2.4201	

average

New method		Classical method	
iteration	time	iteration	Opt time
-----			
35.8100	2.6626	40.0000	2.8812

	New method		Classical method				
idx	iteration	Opt	time	iteration	Opt	time	
-----							
1	33	14.1325	2.2903	36	14.1325	2.4028	
2	34	11.4602	2.3497	40	11.4602	2.6396	
3	34	14.8578	2.3979	43	14.8578	3.4298	
4	32	14.0615	2.3182	42	14.0615	2.9494	
5	37	13.8361	2.6571	40	13.8361	2.7531	
6	27	12.5830	1.8774	31	12.5830	2.0912	
7	35	15.0891	2.4443	46	15.0891	3.1579	
8	42	15.4758	3.0902	51	15.4758	3.5747	
9	33	13.9757	2.3580	36	13.9755	2.3991	
10	33	14.4968	2.3421	34	14.4968	2.3027	
11	31	11.4472	2.1374	65	11.4295	4.6683	
12	29	13.3044	1.9933	36	13.3044	2.4292	
13	34	13.6009	2.3862	50	13.6009	3.4991	
14	34	11.4376	2.4010	41	11.4376	2.8593	
15	31	14.8721	2.1535	35	14.8721	2.3473	
16	38	14.0051	2.7457	35	14.0051	2.3498	
17	29	13.8349	2.0148	35	13.8349	2.3490	
18	35	12.8241	2.4992	42	12.8241	2.9333	
19	33	13.0498	2.3489	42	13.0499	2.9874	
20	33	14.5358	2.3046	35	14.5358	2.3323	
21	33	13.7607	2.3485	38	13.7607	2.6101	
22	31	13.3368	2.1436	36	13.3368	2.4539	
23	31	12.6427	2.1378	44	12.6427	3.0703	
24	33	13.1727	2.3204	47	13.1727	3.3550	

25	38	12.4437	2.7154	45	12.4437	3.1244
26	30	16.4101	2.0944	39	16.4101	2.6710
27	43	14.5286	3.1413	39	14.5286	2.6813
28	32	14.2258	2.2594	31	14.2258	2.1112
29	34	11.9038	2.4134	51	11.9038	3.6221
30	29	12.9792	1.9981	33	12.9792	2.1890
31	30	13.5566	2.0404	39	13.5566	2.6850
32	30	13.9748	2.1090	31	13.9748	2.0360
33	38	13.9658	2.6906	41	13.9658	2.7484
34	31	13.3783	2.1870	34	13.3783	2.4084
35	33	13.5477	2.3660	38	13.5477	2.7975
36	35	12.2685	2.4543	47	12.2685	3.6504
37	34	12.1387	2.4273	42	12.1387	3.1664
38	31	14.3202	2.1804	42	14.3202	3.1886
39	33	11.7495	2.4815	30	11.7495	2.0452
40	30	15.0938	2.1312	39	15.0938	2.8654
41	40	15.4534	3.0469	44	15.4534	3.3165
42	38	12.4233	2.8703	44	12.4233	3.2338
43	66	12.7004	5.6847	40	12.7022	2.8478
44	39	12.4386	2.9631	40	12.4386	2.8733
45	33	13.5479	2.3851	42	13.5479	3.0937
46	36	13.7716	2.6777	43	13.7716	3.1809
47	33	12.7877	2.5047	35	12.7877	2.5006
48	36	16.7568	2.7086	36	16.7568	2.5545
49	38	15.3569	2.9702	34	15.3569	2.4201
50	35	10.1950	2.5621	41	10.1950	2.9928
51	39	14.4130	3.0114	47	14.4130	3.6163



52	35	14.5251	2.6399	40	14.5251	2.9441
53	72	14.9484	6.2311	45	14.9489	3.2947
54	35	16.3867	2.5885	40	16.3867	2.8813
55	38	12.7772	2.9128	42	12.7772	3.1070
56	34	13.0523	2.5302	38	13.0523	2.8084
57	35	13.0530	2.5921	43	13.0530	3.2378
58	34	13.0528	2.5316	35	13.0528	2.4582
59	31	15.8548	2.2426	35	15.8548	2.5212
60	29	14.2370	2.0127	39	14.2370	2.8138
61	43	13.5499	3.3413	38	13.5499	2.7139
62	35	12.8535	2.5790	35	12.8535	2.4133
63	33	12.9314	2.4403	37	12.9314	2.6749
64	95	13.7239	9.0412	44	13.7796	3.2587
65	31	15.8651	2.2830	34	15.8651	2.3814
66	37	13.8844	2.7863	38	13.8844	2.6984
67	38	14.4416	2.9258	46	14.4415	3.5383
68	35	13.3879	2.5361	44	13.3879	3.1893
69	37	14.7301	2.6624	47	14.7301	3.5063
70	34	14.0756	2.5296	40	14.0755	2.8794
71	37	13.8434	2.7015	41	13.8434	2.9964
72	32	14.9581	2.3169	32	14.9581	2.1915
73	31	13.0970	2.2517	40	13.0970	2.9671
74	32	12.1322	2.3486	42	12.1322	3.1600
75	33	14.5811	2.4676	36	14.5811	2.5803
76	36	15.1297	2.7490	40	15.1297	2.9157
77	28	16.3501	2.0110	37	16.3501	2.6951
78	37	11.7357	2.7682	38	11.7357	2.6857

79	47	13.4661	3.8289	38	13.4661	2.7929
80	30	14.1476	2.1423	38	14.1476	2.7840
81	33	14.0924	2.4596	38	14.0924	2.7350
82	37	14.7447	2.7495	39	14.7447	2.7690
83	40	13.4680	2.9901	46	13.4680	3.4051
84	37	11.5136	2.7620	42	11.5136	3.0184
85	36	14.5218	2.7568	36	14.5218	2.6255
86	54	13.8599	3.9193	36	13.8574	2.6522
87	32	12.6705	2.4086	41	12.6705	3.1666
88	34	12.5076	2.6738	38	12.5076	2.8091
89	40	14.6875	3.2020	34	14.6875	2.5030
90	30	11.3641	2.2120	33	11.3641	2.3333
91	34	14.8216	2.5382	40	14.8216	2.9601
92	34	16.0217	2.5202	48	16.0217	3.8445
93	33	14.0469	2.4813	41	14.0469	3.1175
94	40	15.8152	3.0568	41	15.8152	2.9229
95	30	14.3876	2.2207	34	14.3876	2.4590
96	30	15.8316	2.2985	40	15.8316	3.0683
97	33	14.2434	2.4493	38	14.2434	2.8031
98	37	13.7590	2.8397	46	13.7590	3.5587
99	39	15.5669	2.8991	52	15.5669	3.9323
100	35	15.0051	2.6977	48	15.0051	3.7811

-----

average

New method		Classical method	
iteration	time	iteration Opt	time
35.8100	2.6626	40.0000	2.8812

### MatLab code for Table 5.4

solve the random system with dimention 400\*700

```
m=400;n=700;
```

```
A=rand(m,n);
```

```
b=rand(m,1);
```

```
c=rand(n,1);
```

```
epsilon=0.01;
```

in Dr. Lesaja's mehtod,  $p=0.8, q=1.2$

	New method			Classical method		
idx	iteration	Opt	time	iteration	Opt	time
-----						
1	41	16.8015	25.4030	57	16.8018	35.8048
-----						

### MatLab code for Table 5.4

```
impNew
```

```
p=1-lambda;
```

```
q=1+lambda;
```

average of 200 iterations

```

-----
          0.1          0.2
dim(m*n) iteration time iteration time
average 5.0 5.1 35.9 0.0526 37.2 0.0603

```

```

-----
          0.3          0.4
dim(m*n) iteration time iteration time
average 4.8 5.1 40.6 0.0673 41.6 0.0648

```

### MatLab code for Table 5.6

20 iterations

```

-----
          0.3          0.4
dim(m*n) iteration time iteration time
average 200.0 300.0 33.8 2.5982 37.8 2.9957

```

```

-----
          0.1          0.2
dim(m*n) iteration time iteration time
average 200.0 300.0 36.6 2.8611 34.7 2.6867

```

---

		0.1		0.2			
idx	dim(m*n)	iteration	Opt	time	iteration	Opt	time
<hr/>							
1	200300	36	15.3765	2.7472	35	15.3765	2.6816
2	200300	33	13.6721	2.5727	33	13.6721	2.5643
3	200300	39	13.1033	3.0115	41	13.1033	3.2237
4	200300	38	14.9661	3.0094	40	14.9661	3.2188
5	200300	32	12.9225	2.4362	28	12.9225	2.0493
6	200300	32	13.3005	2.3987	34	13.3005	2.5932
7	200300	38	11.4939	2.9336	36	11.4939	2.8237
8	200300	32	13.6853	2.4423	37	13.6853	2.9811
9	200300	33	12.6659	2.5796	28	12.6659	2.0762
10	200300	79	13.1793	6.9372	43	13.1802	3.4659
11	200300	33	12.4495	2.4903	33	12.4495	2.4935
12	200300	33	14.7143	2.5612	31	14.7143	2.3323
13	200300	32	13.0591	2.3790	34	13.0591	2.6057
14	200300	41	14.7339	3.2123	33	14.7339	2.4766
15	200300	34	15.4971	2.6664	33	15.4971	2.6206
16	200300	37	12.8786	2.9451	40	12.8786	3.0927
17	200300	36	12.6252	2.7991	37	12.6252	2.9537
18	200300	33	14.2547	2.5594	32	14.2547	2.4313
19	200300	32	15.6560	2.4247	32	15.6560	2.4553
20	200300	29	14.5694	2.1154	34	14.5694	2.5937

---