Bootcamp: Interior Point Algorithms I

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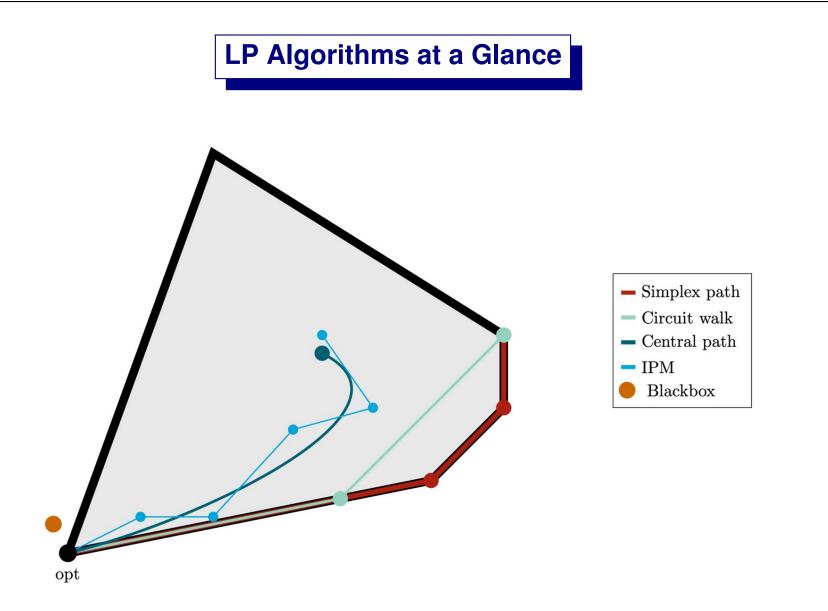


Figure 1: Slide from Daniel Dadush and Bento Natura 2023

Interior-Point Algorithms for LP

Consider linear program:

min
$$\mathbf{c}^T \mathbf{x}$$
 s.t. $A\mathbf{x} = \mathbf{b}, \ \mathbf{x} \ge \mathbf{0}.$

int
$$\mathcal{F}_p = \{\mathbf{x} : A\mathbf{x} = \mathbf{b}, \mathbf{x} > \mathbf{0}\} \neq \emptyset$$

int
$$\mathcal{F}_d = \{(\mathbf{y}, \mathbf{s}) : \mathbf{s} = \mathbf{c} - A^T \mathbf{y} > \mathbf{0}\} \neq \emptyset.$$

Let z^* denote the optimal value and

$$\mathcal{F} = \mathcal{F}_p \times \mathcal{F}_d.$$

We are interested in finding an ϵ -approximate solution for the LP problem:

$$\mathbf{c}^T \mathbf{x} - \mathbf{b}^T \mathbf{y} \le \epsilon.$$

For simplicity, we assume that an interior-point pair $(\mathbf{x}^0, \mathbf{y}^0, \mathbf{s}^0)$ is known, and we will use it as our initial point pair.

Karmarkar 1984, Renegar 1986...

Logarithmic Barrier Functions for LP

Consider the logarithmic barrier function optimization

$$(PB) \quad \begin{array}{ll} \text{minimize} & -\sum_{j=1}^n \log x_j \\ \text{s.t.} & \mathbf{x} \in \operatorname{int} \mathcal{F}_p \end{array}$$

and

$$(DB)$$
 maximize $\sum_{j=1}^{n} \log s_j$
s.t. $(\mathbf{y}, \mathbf{s}) \in \operatorname{int} \mathcal{F}_d$

They are linearly constrained convex programs (LCCP).

Much much earlier...

(1)

Analytic Center for the Primal Polytope

The maximizer $\bar{\mathbf{x}}$ of (PB) is called the analytic center of polytope \mathcal{F}_p . From the optimality condition theorem, we have

$$-(\bar{X})^{-1}\mathbf{e} - A^T\mathbf{y} = \mathbf{0}, \ A\bar{\mathbf{x}} = \mathbf{b}, \ \bar{\mathbf{x}} > \mathbf{0},$$

where e is the vector of all ones; or

$$X\mathbf{s} = \mathbf{e}$$
$$A\bar{\mathbf{x}} = \mathbf{b}$$
$$-A^T\mathbf{y} - \mathbf{s} = \mathbf{0}$$
$$\bar{\mathbf{x}} > \mathbf{0}$$

where X is the diagonal matrix generated from vector \mathbf{x} .

Sonnevend 1988, Bayer and Lagarias 1989, Megiddo 1989...

(2)

Analytic Center for the Dual Polytope

The maximizer (\bar{y}, \bar{s}) of (DB) is called the analytic center of polytope \mathcal{F}_d , and we have

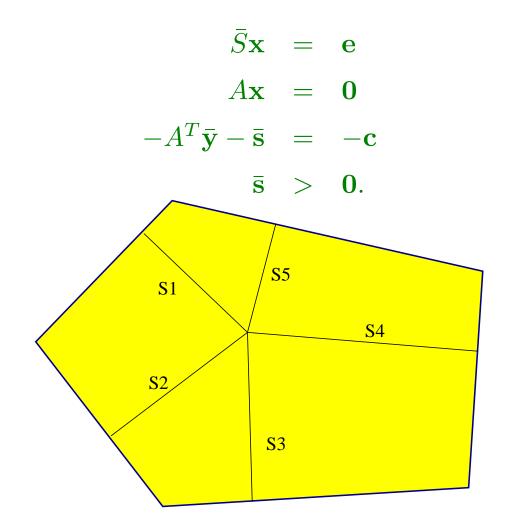


Figure 2: Analytic center maximizes the product of slacks.

Volumetric Center for the Dual Polytope

The maximizer (\bar{y}, \bar{s}) of the following problem is called the volumetric center of polytope \mathcal{F}_d , and we have

$$\begin{array}{ll} (DB) & \text{maximize} & -\log \det(AS^{-2}A^T) \\ & \text{s.t.} & (\mathbf{y},\mathbf{s}) \in \operatorname{int} \mathcal{F}_d \end{array}$$

More details see Vaidya 1996, Lee-Sidford 13-'19, van den Brand, Lee, Liu, Saranurak, Sidford, Song, Wang 21, etc.

Why Analytic

The analytic center of polytope \mathcal{F}_d is an analytic function of input data A, \mathbf{c} .

Consider $\Omega = \{y \in R : -y \le 0, y \le 1\}$, which is interval [0, 1]. The analytic center is $\overline{y} = 1/2$ with $\mathbf{x} = (2, 2)^T$.

Consider

$$\Omega' = \{ y \in R : \overbrace{-y \le 0, \cdots, -y \le 0}^{n \text{ times}}, y \le 1 \},\$$

which is, again, interval [0, 1] but " $-y \le 0$ " is copied n times. The analytic center for this system is $\bar{y} = n/(n+1)$ with $\mathbf{x} = ((n+1)/n, \dots, (n+1)/n, (n+1))^T$.

Analytic Volume of Polytope and Cutting Plane

$$AV(\mathcal{F}_d) := \prod_{j=1}^n \bar{s}_j = \prod_{j=1}^n (c_j - \mathbf{a}_j^T \bar{\mathbf{y}})$$

can be viewed as the analytic volume of polytope \mathcal{F}_d or simply \mathcal{F} in the rest of discussions.

If one inequality in \mathcal{F} , say the first one, needs to be translated, change $\mathbf{a}_1^T \mathbf{y} \leq c_1$ to $\mathbf{a}_1^T \mathbf{y} \leq \mathbf{a}_1^T \bar{\mathbf{y}}$; i.e., the first inequality is parallelly moved and it now cuts through $\bar{\mathbf{y}}$ and divides \mathcal{F} into two bodies. Analytically, c_1 is replaced by $\mathbf{a}_1^T \bar{\mathbf{y}}$ and the rest of data are unchanged. Let

$$\mathcal{F}^+ := \{ \mathbf{y} : \mathbf{a}_j^T \mathbf{y} \le c_j^+, \ j = 1, ..., n \},\$$

where $c_j^+ = c_j$ for j = 2, ..., n and $c_1^+ = \mathbf{a}_1^T \bar{\mathbf{y}}$.

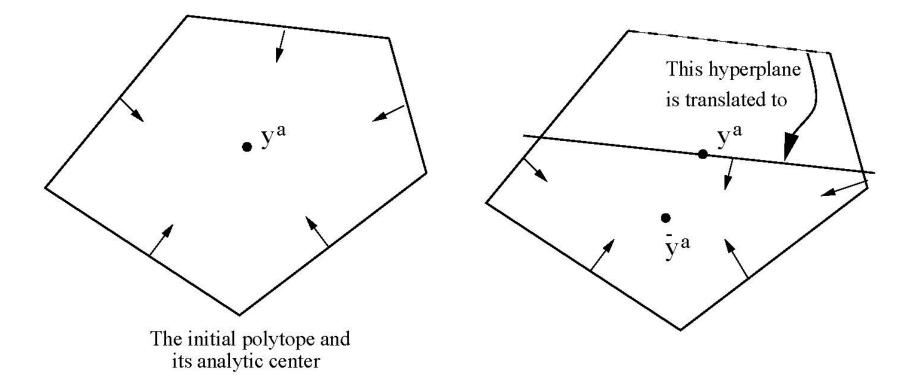


Figure 3: Translation of a hyperplane through the AC.

Analytic Volume Reduction of the New Polytope

Let $\bar{\mathbf{y}}^+$ be the analytic center of $\mathcal{F}^+.$ Then, the analytic volume of \mathcal{F}^+

$$AV(\mathcal{F}^{+}) = \prod_{j=1}^{n} (c_{j}^{+} - \mathbf{a}_{j}^{T} \bar{\mathbf{y}}^{+}) = (\mathbf{a}_{1}^{T} \bar{\mathbf{y}} - \mathbf{a}_{1}^{T} \bar{\mathbf{y}}^{+}) \prod_{j=2}^{n} (c_{j} - \mathbf{a}_{j}^{T} \bar{\mathbf{y}}^{+}).$$

We have the following volume reduction theorem:

Theorem 1

$$\frac{AV(\mathcal{F}^+)}{AV(\mathcal{F})} \le \exp(-1).$$



Since $\bar{\mathbf{y}}$ is the analytic center of $\mathcal F$, there exists $\bar{\mathbf{x}}>\mathbf{0}$ such that

$$\bar{X}\bar{\mathbf{s}} = \bar{X}(\mathbf{c} - A^T\bar{\mathbf{y}}) = \mathbf{e}$$
 and $A\bar{\mathbf{x}} = \mathbf{0}$.

Thus,

$$\bar{\mathbf{s}} = (\mathbf{c} - A^T \bar{\mathbf{y}}) = \bar{X}^{-1} \mathbf{e}$$
 and $\mathbf{c}^T \bar{\mathbf{x}} = (\mathbf{c} - A^T \bar{\mathbf{y}})^T \bar{\mathbf{x}} = \mathbf{e}^T \mathbf{e} = n.$

We have

$$\mathbf{e}^T \bar{X} \bar{\mathbf{s}}^+ = \mathbf{e}^T \bar{X} (\mathbf{c}^+ - A^T \bar{\mathbf{y}}^+) = \mathbf{e}^T \bar{X} \mathbf{c}^+$$
$$= \mathbf{c}^T \bar{\mathbf{x}} - \bar{x}_1 (c_1 - \mathbf{a}_1^T \bar{\mathbf{y}}) = n - 1.$$

$$\frac{AV(\mathcal{F}^+)}{AV(\mathcal{F})} = \prod_{j=1}^n \frac{\bar{s}_j^+}{\bar{s}_j}$$
$$= \prod_{j=1}^n \bar{x}_j \bar{s}_j^+$$
$$\leq \left(\frac{1}{n} \sum_{j=1}^n \bar{x}_j \bar{s}_j^+\right)^n$$
$$= \left(\frac{1}{n} \mathbf{e}^T \bar{X} \bar{\mathbf{s}}^+\right)^n$$
$$= \left(\frac{n-1}{n}\right)^n \leq \exp(-1).$$

Analytic Volume of Polytope and Multiple Cutting Planes

Now suppose we translate k(< n) hyperplanes, say 1, 2, ..., k, moved to cut the analytic center \bar{y} of \mathcal{F} , that is,

$$\mathcal{F}^+:=\{\mathbf{y}:\ \mathbf{a}_j^T\mathbf{y}\leq c_j^+,\ j=1,...,n\},$$
 where $c_j^+=c_j$ for $j=k+1,...,n$ and $c_j^+=\mathbf{a}_j^T\bar{\mathbf{y}}$ for $j=1,...,k.$
Corollary 1

$$\frac{AV(\mathcal{F}^+)}{AV(\mathcal{F})} \le \exp(-k).$$

The Analytic Center Cutting-Plane Method

Problem: Find a solution in the feasible set $\mathcal{F} := \{\mathbf{y} : \mathbf{a}_j^T \mathbf{y} \le c_j, j = 1, ..., n\}$. Start with the initial polytope

$$\mathcal{F}^{0} := \{ \mathbf{y} : \mathbf{a}_{j}^{T} \mathbf{y} \le c_{j}^{0} := c_{j} + R, \ j = 1, ..., n \}$$

where R is sufficiently large such that $\bar{\mathbf{y}}^0 = \mathbf{0}$ is an (approximate) analytic center of \mathcal{F}^0 .

Check if the (approximate) analytic center $\bar{\mathbf{y}}^k$ of \mathcal{F}^k is in \mathcal{F} or not. If not, define a new polytope \mathcal{F}^{k+1} by translating one or multiple violated constraint hyperplanes through $\bar{\mathbf{y}}^k$ as defined earlier, and compute an approximate analytic center $\bar{\mathbf{y}}^{k+1}$ of \mathcal{F}^{k+1} .

Continue this step till $\bar{\mathbf{y}}^k \in \mathcal{F}$.

Fair Pareto Optimal Solutions of Multiple Objectives

Problem: Find a solution in the feasible set $\mathcal{F} := \{\mathbf{y} : \mathbf{a}_j^T \mathbf{y} \leq c_j, j = 1, ..., n\}$ such that it is Pareto-Maximal for k objective function $\mathbf{b}_i^T \mathbf{y}, i = 1, ..., k$

- Weighted-Sum Objective Maximization: $\sum_i w_i \mathbf{b}_i^T \mathbf{y}$
- Alternating Cuting-Plane: Cuting the AC alternatively, or simultaneously with fixed proportions?

Trajectory of Analytic Centers: Central Path for LP

Now consider the problem

maximize
$$\mathbf{b}^T \mathbf{y}$$

s.t. $A^T \mathbf{y} \leq \mathbf{c}$.

Assume that the feasible region is bounded, and the analytic center of the region is y^0 . Start with a polytope

$$\mathcal{F}(R) := \{ \mathbf{y} : A^T \mathbf{y} \le \mathbf{c}, \ \mathbf{b}^T \mathbf{y} \ge R, \cdots, \mathbf{b}^T \mathbf{y} \ge R \}$$

where R is so low such that \mathbf{y}^0 is also an (approximate) analytic center of $\mathcal{F}(R)$.

Define a family of polytopes $\mathcal{F}(R)$ by continuously increasing R toward the maximal value and consider its analytic center $\mathbf{y}(R)$: it forms a path of analytic centers from \mathbf{y}^0 toward the optimal solution set.

Better Parameterization: LP Regularized by the Barrier Function

An equivalent algebraic representation of the path is to consider the LP problem with the weighted barrier function

$$\begin{array}{ll} (LDB) & \text{maximize} & \mathbf{b}^T \mathbf{y} + \mu \sum_{j=1}^n \log s_j \\ & \text{s.t.} & (\mathbf{y}, \mathbf{s}) \in \operatorname{int} \mathcal{F}_d, \end{array}$$

and also

$$(LPB)$$
 minimize $\mathbf{c}^T \mathbf{x} - \mu \sum_{j=1}^n \log x_j$
s.t. $\mathbf{x} \in \operatorname{int} \mathcal{F}_p$

where μ is called the barrier (weight) parameter.

They are again linearly constrained convex programs (LCCP).

Self-Duality of LPB and LDB

They share the same first-order KKT conditions:

$$X\mathbf{s} = \mu \mathbf{e}$$
$$A\mathbf{x} = \mathbf{b}$$
$$-A^T \mathbf{y} - \mathbf{s} = -\mathbf{c};$$

where we have

$$\mu = \frac{\mathbf{x}^T \mathbf{s}}{n} = \frac{\mathbf{c}^T \mathbf{x} - \mathbf{b}^T \mathbf{y}}{n},$$

so that it's the average of complementarity or duality gap.

Denote by $(\mathbf{x}(\mu), \mathbf{y}(\mu), \mathbf{s}(\mu))$ the (unique) solution satisfying the conditions. As μ decreases to zero, $\mathbf{x}(\mu)$ form a path in the primal feasible region and $\mathbf{y}(\mu)$ form a path in the dual feasible region to-warding optimality respectively.

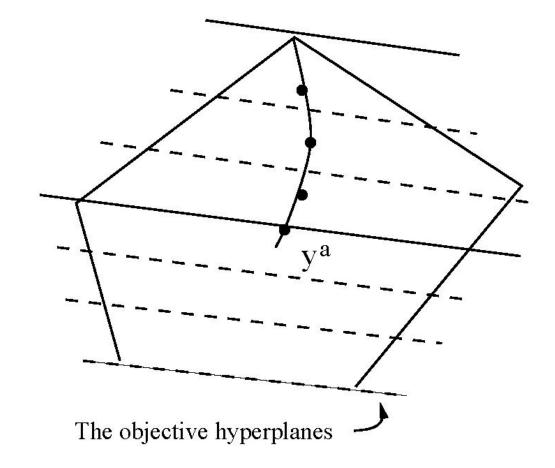


Figure 4: The central path of $\mathbf{y}(\mu)$ in a dual feasible region.

Central Path for Linear Programming Parametrized by μ

Theorem 2 Let $(\mathbf{x}(\mu), \mathbf{y}(\mu), \mathbf{s}(\mu))$ be on the central path of an linear program in standard form. i) The central path point $(\mathbf{x}(\mu), \mathbf{s}(\mu))$ is bounded for $0 < \mu \le \mu^0$ and any given $0 < \mu^0 < \infty$.

ii) For $0<\mu'<\mu$,

$$\mathbf{c}^T \mathbf{x}(\mu') < \mathbf{c}^T \mathbf{x}(\mu) \quad \textit{and} \quad \mathbf{b}^T \mathbf{y}(\mu') > \mathbf{b}^T \mathbf{y}(\mu)$$

if both primal and dual have nontrivial optimal solutions.

iii) $(\mathbf{x}(\mu), \mathbf{s}(\mu))$ converges to an optimal solution pair for (LP) and (LD). Moreover, the limit point $\mathbf{x}(0)_{P^*} > \mathbf{0}$ and the limit point $\mathbf{s}(0)_{Z^*} > \mathbf{0}$, where (P^*, Z^*) are the analytic centers on the primal and dual optimal faces, respectively (G[']uler and Y 1993).

Proof of (i)

 $(\mathbf{x}(\mu^0) - \mathbf{x}(\mu))^T (\mathbf{s}(\mu^0) - \mathbf{s}(\mu)) = 0,$ since $(\mathbf{x}(\mu^0) - \mathbf{x}(\mu)) \in \mathcal{N}(A)$ and $(\mathbf{s}(\mu^0) - \mathbf{s}(\mu)) \in \mathcal{R}(A^T)$. This can be rewritten as

$$\sum_{j=1}^{n} \left(s(\mu^{0})_{j} x(\mu)_{j} + x(\mu^{0})_{j} s(\mu)_{j} \right) = n(\mu^{0} + \mu) \le 2n\mu^{0},$$

or

$$\sum_{j}^{n} \left(\frac{x(\mu)_j}{x(\mu^0)_j} + \frac{s(\mu)_j}{s(\mu^0)_j} \right) \le 2n.$$

Thus, $\mathbf{x}(\mu)$ and $\mathbf{s}(\mu)$ are bounded, which proves (i).

The Path-Following Algorithms

In general, one can start from an (approximate) central path point $\mathbf{x}(\mu^0)$, $(\mathbf{y}(\mu^0), \mathbf{s}(\mu^0))$, or $(\mathbf{x}(\mu^0), \mathbf{y}(\mu^0), \mathbf{s}(\mu^0))$ where μ^0 is sufficiently large.

Then, let μ^1 be a slightly smaller parameter than μ^0 . Then, we compute an (approximate) central path point $\mathbf{x}(\mu^1)$, $(\mathbf{y}(\mu^1), \mathbf{s}(\mu^1))$, or $(\mathbf{x}(\mu^1), \mathbf{y}(\mu^1), \mathbf{s}(\mu^1))$. They can be updated from the previous point at μ^0 using the Newton method.

 μ might be reduced at each stage by a specific factor, giving $\mu^{k+1} = \gamma \mu^k$ where γ is at most $1 - \frac{1}{3\sqrt{n}}$, where k is the iteration count.

This is called the primal, dual, or primal-dual path-following method; see Renegar 1988, Gonzagar 1989, Kojima et al. 1989, Monteiro and Adler 1989,...

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The Newton Method of Primal-Dual Path-Following

Given a pair $(\mathbf{x}^k, \mathbf{y}^k, \mathbf{s}^k) \in \operatorname{int} \mathcal{F}$ closely to the central path, that is,

$$\|X^k S^k \mathbf{e} - \mu^k \mathbf{e}\| \le \eta \mu^k$$

for a small positive constant η , we compute direction vectors \mathbf{d}_x , \mathbf{d}_y and \mathbf{d}_s from the system equations:

$$S^{k}\mathbf{d}_{x} + X^{k}\mathbf{d}_{s} = \gamma \mu^{k}\mathbf{e} - X^{k}S^{k}\mathbf{e},$$

$$A\mathbf{d}_{x} = \mathbf{0},$$

$$-A^{T}\mathbf{d}_{y} - \mathbf{d}_{s} = \mathbf{0}.$$
(3)

where $\gamma = (1 - \frac{1}{3\sqrt{n}})$. Then we update

$$\mathbf{x}^{k+1} = \mathbf{x}^k + \mathbf{d}_x > \mathbf{0}, \ \mathbf{y}^{k+1} = \mathbf{y}^k + \mathbf{d}_y, \ \mathbf{s}^{k+1} = \mathbf{s}^k + \mathbf{d}_s > \mathbf{0}.$$

Then one can prove

$$||X^{k+1}S^{k+1}\mathbf{e} - \mu^{k+1}\mathbf{e}|| \le \eta(1 - \frac{1}{3\sqrt{n}})\mu^k = \eta\mu^{k+1}.$$

This leads to \sqrt{n} iteration complexity.

Adaptive Path-Following Algorithms

Here we describe and analyze the Predictor-Corrector interior-point algorithm (Mizuno-Todd-Y 1993, Mehrotra 1993). Consider the neighborhood

$$\mathcal{N}_2(\eta) = \left\{ (\mathbf{x}, \mathbf{s}) \in \operatorname{int} \mathcal{F} : \| X\mathbf{s} - \mu \mathbf{e} \| \le \eta \mu \quad \text{where} \quad \mu = \frac{\mathbf{x}^T \mathbf{s}}{n} \right\} \text{ for some } \eta \in (0, 1).$$

Given $(\mathbf{x}^0, \mathbf{s}^0) \in \mathcal{N}_2(\eta)$ with $\eta = 1/4$. Set $k := 0$.

While $(\mathbf{x}^k)^T \mathbf{s}^k > \epsilon$ do:

1. Predictor step: set $(\mathbf{x}, \mathbf{s}) = (\mathbf{x}^k, \mathbf{s}^k)$ and compute $\mathbf{d} = \mathbf{d}(\mathbf{x}, \mathbf{s}, 0)$ from (3); compute the largest $\overline{\theta}$ so that

$$(\mathbf{x}(\theta), \mathbf{s}(\theta)) \in \mathcal{N}_2(2\eta) \text{ for } \theta \in [0, \overline{\theta}].$$

- 2. Corrector step: set $(\mathbf{x}', \mathbf{s}') = (\mathbf{x}(\bar{\theta}), \mathbf{s}(\bar{\theta}))$ and compute $\mathbf{d}' = \mathbf{d}(\mathbf{x}', \mathbf{s}', 1)$ from (3); set $(\mathbf{x}^{k+1}, \mathbf{s}^{k+1}) = (\mathbf{x}' + \mathbf{d}'_x, \mathbf{s}' + \mathbf{d}'_s).$
- 3. Let k := k + 1 and return to Step 1.

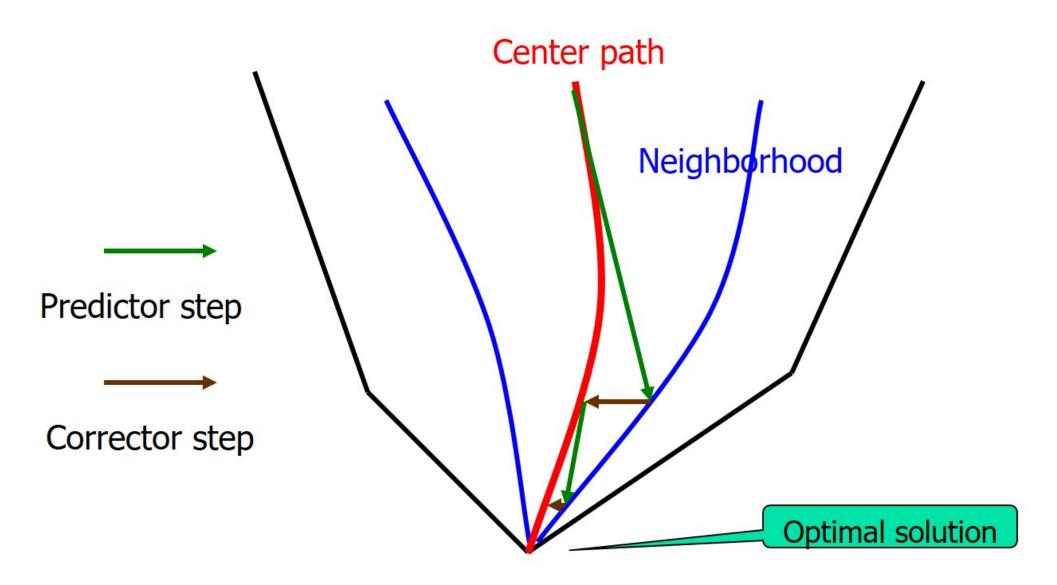


Figure 5: Illustration of the predictor-corrector algorithm.

Reduce the Analytical Volume Directly: Potential Reduction

Problem: Find a solution in the feasible set $\mathcal{F} := \{\mathbf{y} : \mathbf{a}_j^T \mathbf{y} \le c_j, j = 1, ..., n\}.$

For $\mathbf{x} \in \operatorname{int} \mathcal{F}_p$, Karmarkar's primal potential function is defined by

$$\psi_n(\mathbf{x}) := n \log(\mathbf{c}^T \mathbf{x} - z^*) - \sum_{j=1}^n \log(x_j),$$

where z^* is the optimal objective value of the LP problem.

This leads to n iteration complexity.

Primal-Dual Potential Function for LP

Typically, a single merit-function driven algorithm is preferred since it can adaptively take large step sizes as long as the merit value is sufficiently reduced, comparing to check and balance of hyper-parameters/measures of the path-following type of algorithms.

For $x \in int \mathcal{F}_p$ and $(y, s) \in int \mathcal{F}_d$, the joint Tanabe-Todd-Ye primal-dual potential function is defined by

$$\psi_{n+\rho}(\mathbf{x}, \mathbf{s}) := (n+\rho)\log(\mathbf{x}^T\mathbf{s}) - \sum_{j=1}^n \log(x_j s_j),$$

m

where $\rho \geq 0$ and it is fixed.

$$\psi_{n+\rho}(\mathbf{x}, \mathbf{s}) = \rho \log(\mathbf{x}^T \mathbf{s}) + \psi_n(\mathbf{x}, \mathbf{s}) \ge \rho \log(\mathbf{x}^T \mathbf{s}) + n \log n,$$

then, for $\rho > 0$, $\psi_{n+\rho}(\mathbf{x}, \mathbf{s}) \to -\infty$ implies that $\mathbf{x}^T \mathbf{s} \to 0$. More precisely, we have

$$\mathbf{x}^T \mathbf{s} \le \exp(\frac{\psi_{n+\rho}(\mathbf{x},\mathbf{s}) - n\log n}{\rho}).$$

Choosing $\rho=\sqrt{n}$ leads to \sqrt{n} iteration complexity.

Description of Algorithm

Given $(\mathbf{x}^0, \mathbf{y}^0, \mathbf{s}^0) \in \operatorname{int} \mathcal{F}$. Set $\rho \ge \sqrt{n}$ and k := 0. While $(\mathbf{x}^k)^T \mathbf{s}^k \ge \epsilon$ do

1. Set $(\mathbf{x}, \mathbf{s}) = (\mathbf{x}^k, \mathbf{s}^k)$ and $\gamma = n/(n+\rho)$ and compute $(\mathbf{d}_x, \mathbf{d}_y, \mathbf{d}_s)$ from (3).

2. Let
$$\mathbf{x}^{k+1} = \mathbf{x}^k + \bar{\alpha} \mathbf{d}_x$$
, $\mathbf{y}^{k+1} = \mathbf{y}^k + \bar{\alpha} \mathbf{d}_y$, and $\mathbf{s}^{k+1} = \mathbf{s}^k + \bar{\alpha} \mathbf{d}_s$ where
 $\bar{\alpha} = \arg\min_{\alpha \ge 0} \psi_{n+\rho} (\mathbf{x}^k + \alpha \mathbf{d}_x, \mathbf{s}^k + \alpha \mathbf{d}_s).$

3. Let k := k + 1 and return to Step 1.

Allow to take longer step-size α !

Alternating Direction Method

Recall that for $x \in int \mathcal{F}_p$ and $(y, s) \in int \mathcal{F}_d$, the joint primal-dual potential function is defined as

$$\psi_{n+\rho}(\mathbf{x}, \mathbf{s}) := (n+\rho) \log(\mathbf{x}^T \mathbf{s}) - \sum_{j=1}^n \log(x_j s_j)$$

= $(n+\rho) \log(\mathbf{c}^T \mathbf{x} - \mathbf{b}^T \mathbf{y}) - \sum_{j=1}^n \log(x_j) - \sum_{j=1}^n \log(s_j).$

Alternate Updating of primal \mathbf{x} and (\mathbf{y}, \mathbf{s}) : at the *k*th step, fix $(\mathbf{y}^k, \mathbf{s}^k)$ and reduce the potential function by a constant via updating from \mathbf{x}^k to \mathbf{x}^{k+1} while keep $(\mathbf{y}^{k+1}, \mathbf{s}^{k+1}) = (\mathbf{y}^k, \mathbf{s}^k)$:

$$\psi_{n+\rho}(\mathbf{x}^{k+1},\mathbf{s}^{k+1}) - \psi_{n+\rho}(\mathbf{x}^k,\mathbf{s}^k) \le -\delta.$$

Once can prove that, if by updating primal only one cannot reduce the potential function by a constant anymore, then one must be able to update the dual from $(\mathbf{y}^k, \mathbf{s}^k)$ to $(\mathbf{y}^{k+1}, \mathbf{s}^{k+1})$ (while keep $\mathbf{x}^{k+1} = \mathbf{x}^k$) and reduce the potential function by a constant; see Y 1989. The sample complexity result holds and it was the first one extended to solving SDP by Alizadeh 1992.

First-Order Potential Reduction

At the *k*th iteration, we compute the direction vectors $(\mathbf{d}_x, \mathbf{d}_y, \mathbf{d}_s)$ using the steepest descent direction:

$$\min \quad \nabla_x \phi(\mathbf{x}^k, \mathbf{s}^k)^T \mathbf{d}_x + \nabla_s \phi(\mathbf{x}^k, \mathbf{s}^k)^T \mathbf{d}_s$$
s.t.
$$A \mathbf{d}_x = \mathbf{0}$$

$$A^T \mathbf{d}_y + \mathbf{d}_s = \mathbf{0}.$$

Thus,

$$\begin{aligned} \mathbf{d}_x &= -(I - A^T (A A^T)^{-1} A) \nabla_x \phi(\mathbf{x}^k, \mathbf{s}^k), \\ \mathbf{d}_y &= A \nabla_s \phi(\mathbf{x}^k, \mathbf{s}^k), \\ \mathbf{d}_s &= -A^T A \nabla_s \phi(\mathbf{x}^k, \mathbf{s}^k). \end{aligned}$$

First-Order Potential Reduction as a Presolver

- First-order method solves to 1e-02 accuracy and then switch to second-order
- An average solution reduction of 30%

Accuracy	1e-04	1e-06	1e-08	1e-10
First-order	7.5	798.0	>1200	>1200
Second-order	33.0	56.7	89.3	93.3
First + Second	5.4	12.1	14.1	15.2

Figure 6: Speed-Up on QAP-LP

Potential Reduction for General Linear Complementarity

Given $M \in R^{n \times n}$ and $\mathbf{q} \in R^n$, find (\mathbf{x}, \mathbf{s})

$$\mathbf{s} = M\mathbf{x} + \mathbf{q}, \ \mathbf{x} \ge \mathbf{0}, \ \mathbf{s} \ge \mathbf{0}, \ \text{ and } \ \mathbf{x}^T\mathbf{s} = 0$$

- $\bullet \ \operatorname{LP:} M \text{ skew symmetric} \\$
- Convex QP: M symmetric and monotone: $(\mathbf{x} \mathbf{y})^T M(\mathbf{x} \mathbf{y}) \ge 0$
- Monotone LCP: $M + M^T$ is Positive Semidefinite. The PRA terminates in $O(\sqrt{n} \log(\epsilon^{-1}))$ iterations for above three problems.
- P-matrix: $0 < \gamma = \max_{j} \min \frac{\mathbf{x}^{T} M^{T} \mathbf{x}}{\|\mathbf{x}\|^{2}}$. The PRA terminates in $O(n^{2} \max(|\lambda|/(\gamma n), 1) \log(\epsilon^{-1}))$ iterations, where λ (is the least eigenvalue of $(M + M^{T})/2$; see Kojima et al. 1992
- General QP: M symmetric but non-convex: The PRA terminates in $O(n^2 \epsilon^{-1} \log(\epsilon^{-1}) + n \log(n))$ iterations (no condition-numbers!) with a solution that is ϵ accurate on both the first and second order

optimality conditions; see Y 1998.

Initialization

- Combining the primal and dual into a single linear feasibility problem, then applying LP algorithms to find a feasible point of the problem. Theoretically, this approach can retain the currently best complexity result.
- The big M method, i.e., add one or more artificial column(s) and/or row(s) and a huge penalty parameter M to force solutions to become feasible during the algorithm.
- Phase I-then-Phase II method, i.e., first try to find a feasible point (and possibly one for the dual problem), and then start to look for an optimal solution if the problem is feasible and bounded.
- Combined Phase I-Phase II method, i.e., approach feasibility and optimality simultaneously. To our knowledge, the "best" complexity of this approach is $O(n \log(R/\epsilon))$.

Homogeneous and Self-Dual Algorithm

- It solves the linear programming problem without any regularity assumption concerning the existence of optimal, feasible, or interior feasible solutions, while it retains the currently best complexity result
- It can start at any positive primal-dual pair, feasible or infeasible, near the central ray of the positive orthant (cone), and it does not use any big M penalty parameter or lower bound.
- Each iteration solves a system of linear equations whose dimension is almost the same as that solved in the standard (primal-dual) interior-point algorithms.
- If the LP problem has a solution, the algorithm generates a sequence that approaches feasibility and optimality simultaneously; if the problem is infeasible or unbounded, the algorithm will produce an infeasibility certificate for at least one of the primal and dual problems.

Primal-Dual Alternative Systems

A pair of LP has two alternatives

An Integrated Homogeneous System

The two alternative systems can be homogenized as one:

$$\begin{array}{ll} (HP) & A\mathbf{x} - \mathbf{b}\tau &= \mathbf{0} \\ & -A^T\mathbf{y} + \mathbf{c}\tau &= \mathbf{s} \ge \mathbf{0}, \\ & \mathbf{b}^T\mathbf{y} - \mathbf{c}^T\mathbf{x} &= \kappa \ge 0, \\ & \mathbf{y} \text{ free}, \ (\mathbf{x}; \tau) &\ge \mathbf{0} \end{array}$$

where the two alternatives are

 $(\text{Solvable}): \ (\tau>0, \kappa=0) \quad \text{or} \quad (\text{Infeasible}): \ (\tau=0, \kappa>0)$

The Homogeneous System is Self-Dual

$$\begin{array}{lll} (HP) & A\mathbf{x} - \mathbf{b}\tau &= \mathbf{0}, \ (\mathbf{y}') & (HD) & A\mathbf{x}' - \mathbf{b}\tau' &= \mathbf{0}, \\ & -A^T\mathbf{y} + \mathbf{c}\tau &= \mathbf{s} \geq \mathbf{0}, \ (\mathbf{x}') & A^T\mathbf{y}' - \mathbf{c}\tau' &\leq \mathbf{0}, \\ & \mathbf{b}^T\mathbf{y} - \mathbf{c}^T\mathbf{x} &= \kappa \geq 0, \ (\tau') & -\mathbf{b}^T\mathbf{y}' + \mathbf{c}^T\mathbf{x}' &\leq 0, \\ & \mathbf{y} \text{ free}, \ (\mathbf{x};\tau) &\geq \mathbf{0} & \mathbf{y}' \text{ free}, \ (\mathbf{x}';\tau') &\geq \mathbf{0} \end{array}$$

Theorem 3 System (HP) is feasible (e.g. all zeros) and any feasible solution $(\mathbf{y}, \mathbf{x}, \tau, \mathbf{s}, \kappa)$ is *self-complementary*:

$$\mathbf{x}^T \mathbf{s} + \tau \kappa = 0.$$

Furthermore, it has a strictly self-complementary feasible solution

$$\left(egin{array}{c} \mathbf{x} + \mathbf{s} \ au + \kappa \end{array}
ight) > \mathbf{0},$$

Start from any infeasible but interior-solution pair in the primal and dual cones, and apply IPM to solve the Phase I problem; see Y-Todd-Mizuno 1994.