# CS5321 Numerical Optimization

# 14 Linear Programming:The Interior Point Method



#### Interior point methods for LP

- There are many variations of IPM for LP
  - Primal IPM, dual IPM, primal-dual IPM
  - Potential Reduction Methods
  - Path Following Methods
    - Short-step methods
    - Long-step methods
    - Predictor-corrector methods
- We will cover the path following method and the long-step method



#### The dual problem of LP

 $\min_{x} z = c^{T} x$  subject to  $Ax = b, x \ge 0$ 

- The dual problem is (see chap 12)  $\max_{\lambda} b^{T} \lambda$  subject to  $A^{T} \lambda + s = c, s \ge 0$
- Weak duality

For any *x* and  $\lambda$  that are feasible in the primal and the dual problems,  $b^T \lambda \leq c^T x$ .

• Strong duality

If  $x^*$  and  $\lambda^*$  are solutions to the primal and the dual problems,  $b^T \lambda^* = c^T x^*$ .



### **Optimality conditions of LP**

 $\min_{x} z = c^{T} x$  subject to  $Ax = b, x \ge 0$ 

• The Lagrangian function

$$\mathcal{L}(x,\lambda,s) = c^{\mathrm{T}}x - \lambda^{\mathrm{T}}(Ax - b) - s^{\mathrm{T}}x$$

- The optimality conditions (KKT, see chap 12)  $A^T \lambda + s = c$  Ax = b  $x, s \ge 0$   $x_i s_i = 0$ 
  - The last one is the *complementarity* condition

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#### **Primal-dual method**

• The optimality conditions can be expressed as

$$F(x,\lambda,s) = \begin{pmatrix} A^T\lambda + s - c \\ Ax - b \\ XSe \end{pmatrix} = 0, \text{ for } x, s \ge 0$$

•  $X = \text{diag}(x), S = \text{diag}(s) \text{ and } e = (1, 1, ..., 1)^{T}$ .

- *F* is a *nonlinear* equation:  $\mathbb{R}^{2n+m} \to \mathbb{R}^{2n+m}$ .
  - Solved by the Newton's method: (*J* the Jacobian of *F*.)

$$(x, \lambda, s) += \alpha(\Delta x, \Delta \lambda, \Delta s), J(x, \lambda, s) \begin{pmatrix} \Delta x \\ \Delta \lambda \\ \Delta s \end{pmatrix} = -F(x, \lambda, s)$$



#### Jacobian of F

• Let 
$$F(x,\lambda,s) = \begin{pmatrix} A^T\lambda + s - c \\ Ax - b \\ XSe \end{pmatrix} = \begin{pmatrix} r_c \\ r_b \\ r_{XS} \end{pmatrix}$$

• The Jacobian  $J(x, \lambda, s)$  of  $F(x, \lambda, s)$  is

$$J = \begin{pmatrix} \nabla_x r_c & \nabla_\lambda r_c & \nabla_s r_c \\ \nabla_x r_b & \nabla_\lambda r_b & \nabla_s r_b \\ \nabla_x r_{XS} & \nabla_\lambda r_{XS} & \nabla_s r_{XS} \end{pmatrix} = \begin{pmatrix} 0 & A^T & I \\ A & 0 & 0 \\ S & 0 & X \end{pmatrix}$$

#### Solving the linear systems

- The Newton's direction is obtained by solving  $\begin{pmatrix} 0 & A^T & I \\ A & 0 & 0 \\ S & 0 & X \end{pmatrix} \begin{pmatrix} \Delta x \\ \Delta \lambda \\ \Delta s \end{pmatrix} = \begin{pmatrix} -r_c \\ -r_b \\ -r_{XS} \end{pmatrix}$ • S is an artificial variable. Let  $D = S^{-1/2}X^{1/2}$   $\begin{pmatrix} -D^{-2} & A^T \\ A & 0 \end{pmatrix} \begin{pmatrix} \Delta x \\ \Delta \lambda \end{pmatrix} = \begin{pmatrix} -r_c + X^{-1}r_{XS} \\ -r_b \end{pmatrix}$   $\Delta s = -X^{-1}r_{XS} - X^{-1}S\Delta x$ 
  - The left hand side is symmetric, but indefinite.

#### Centering

- In practice,  $r_{XS} = XSe \sigma \mu e$  rather than  $r_{XS} = XSe$ . Duality measure:  $\mu = \frac{1}{n} \sum_{i=1}^{n} x_i s_i = \frac{x^T s}{n}$
- Centering parameter  $\sigma \in [0,1]$ :
  - Decreasing as approaching solutions
  - This is related to the barrier method (chap 15)

Without Centered centering direction

## The path following algorithm

- 1. Given  $(x_0, \lambda_0, s_0)$  with  $(x_0, s_0) > 0$
- 2. For k = 0, 1, 2 ... until  $(x_k)^T s_k < \varepsilon$ 
  - a) Choose  $\sigma_k \in [0,1]$  and let  $\mu_k = (x_k)^T s_k / n$ .
  - b) Solve the Newton's direction  $(\Delta x_k, \Delta \lambda_k, \Delta s_k)$
  - c) Set  $(x_{k+1}, \lambda_{k+1}, s_{k+1}) = (x_k, \lambda_k, s_k) + \alpha_k (\Delta x_k, \Delta \lambda_k, \Delta s_k)$ , where  $\alpha_k$  is chosen to make  $(x_{k+1}, s_{k+1}) > 0$

#### **Central path and neighborhood**

- Feasible set  $\mathbf{F}^0 = \{(x,\lambda,s) \mid Ax = b, A^T\lambda + s = c, (x,s) > 0\}$
- Point  $(x_{\tau}, \lambda_{\tau}, s_{\tau}) \in \mathbf{F}^0$  is on the *central path C* if  $x_{\tau}(i)s_{\tau}(i) = \tau$  for all i=1,2,...,n.
- An example of the neighborhood of the central path

 $\mathbf{N}_{-\infty}(^{\circ}) = \{(x,\lambda,s) \in \mathbf{F}^0 | x_i s_i \ge {}^{\circ} \mu, i = 1, 2, \dots, n\}$ 

- μ is the duality measure (two slides before)
- Typically  $\gamma = 10^{-3}$ .



#### Long-step path-following

- 1. Given  $(x_0, \lambda_0, s_0) \in \mathbb{N}_{-\infty}$  and  $\gamma$ ,  $\sigma_{\min}$ ,  $\sigma_{\max}$ .
- 2. For k = 0, 1, 2 ... until  $(x_k)^T s_k < \varepsilon$ 
  - a) Choose  $\sigma_k \in [\sigma_{\min}, \sigma_{\max}]$  and let  $\mu_k = (x_k)^T s_k / n$ .
  - b) Solve the Newton's direction  $(\Delta x_k, \Delta \lambda_k, \Delta s_k)$
  - c) Set  $(x_{k+1}, \lambda_{k+1}, s_{k+1}) = (x_k, \lambda_k, s_k) + \alpha_k(\Delta x_k, \Delta \lambda_k, \Delta s_k)$ , where  $\alpha_k$  is chosen to make  $(x_{k+1}, \lambda_{k+1}, s_{k+1}) \in \mathbb{N}_{-\infty}$ .



#### Complexity



- The long step following method can converge in O(nlog(1/ε)) iterations (Theorem 14.4)
- The most effective interior point method for LP is the predictor-corrector algorithm (Mehrotra 1992)