# Notes for Operations Research & More

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To My Beloved Motherland China

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# Part I Preliminary Topics

# Chapter 1

# Introduction to Optimization

# 1.1 Optimization Model

The following is the basic forms of terminology:

$$(P) \quad \min \quad f(x) \tag{1.1}$$

s.t. 
$$g_i(x) \le 0, \quad i = 1, 2, ..., m$$
 (1.2)

$$h_j(x) = 0, \quad j = 1, 2, ..., l$$
 (1.3)

$$x \in X \tag{1.4}$$

We have

- $\bullet \ \ \textbf{-} \ x \in R^n \to X \subseteq R^m$
- $q_i(x)$  are called inequality constraints
- $h_j(x)$  are called equality constraints
- X is the domain of the variables (e.g. cone, polygon,  $\{0,1\}^n$ , etc.)
- Let F be the feasible region of (P):
  - $-x^0$  is a feasible solution iff  $x^0 \in F$
  - $-x^*$  is an optimized solution iff  $x^* \in F$  and  $f(x^*) \leq f(x^0), \forall x^0 \in F$  (for minimized problem)

**Notice:** Not every (P) has a feasible region, we can have  $F = \emptyset$ . Even if  $F \neq \emptyset$ , there might not be an solution to P, e.g. unbounded. If (P) has optimized solution(s), it could be 1) Unique, or 2) Infinite number of solution, or 3) Finite number of solution

Types of Optimization Problem

- $m = l = 0, x \in \mathbb{R}^n$ , unconstrained problem
- m+l>0, constrained problem
- $f(x), g_i(x), h_j(x)$  are linear, Linear Optimization
  - If  $X = \mathbb{R}^n$ , Linear Programming
  - If X is discrete, Discrete Optimization
  - If  $X \subseteq \mathbb{Z}^n$ , Integer Programming
  - If  $X \in \{0,1\}^n$ , Binary Programming
  - If  $X \in \mathbb{Z}^n \times \mathbb{R}^m$ , Mixed Integer Programming

#### Linear Programming Formulation Skills 1.2

#### Absolute Value 1.2.1

Consider the following model statement:

$$\min \quad \sum_{j \in J} c_j |x_j|, \quad c_j > 0 \tag{1.5}$$

s.t. 
$$\sum_{j \in J} a_{ij} x_j \gtrsim b_i, \quad \forall i \in I$$
 (1.6)

$$x_j$$
 unrestricted,  $\forall j \in J$  (1.7)

Modeling:

$$\min \sum_{j \in J} c_j(x_j^+ + x_j^-), \quad c_j > 0$$
(1.8)

s.t. 
$$\sum_{j \in J} a_{ij} (x_j^+ - x_j^-) \gtrsim b_i, \quad \forall i \in I$$
 (1.9)

$$x_j^+, x_j^- \ge 0, \quad \forall j \in J \tag{1.10}$$

#### 1.2.2 A Minimax Objective

Consider the following model statement:

$$\min \quad \max_{k \in K} \sum_{j \in J} c_{kj} x_j \tag{1.11}$$

s.t. 
$$\sum_{i \in I} a_{ij} x_j \gtrsim b_i, \quad \forall i \in I$$
 (1.12)

$$x_j \ge 0, \quad \forall j \in J \tag{1.13}$$

Modeling:

$$\min \quad z \tag{1.14}$$

s.t. 
$$\sum_{i \in J} a_{ij} x_j \gtrsim b_i, \quad \forall i \in I$$
 (1.15)

$$\sum_{j \in I} c_{kj} x_j \le z, \quad \forall k \in K \tag{1.16}$$

$$x_j \ge 0, \quad \forall j \in J$$
 (1.17)

#### A Fractional Objective 1.2.3

Consider the following model statement:

$$\min \quad \frac{\sum_{j \in J} c_j x_j + \alpha}{\sum_{j \in J} d_j x_j + \beta}$$
s.t. 
$$\sum_{j \in J} a_{ij} x_j \gtrsim b_i, \quad \forall i \in I$$

$$(1.18)$$

s.t. 
$$\sum_{j \in J} a_{ij} x_j \gtrsim b_i, \quad \forall i \in I$$
 (1.19)

$$x_i \ge 0, \quad \forall j \in J$$
 (1.20)

Modeling:

$$\min \quad \sum_{j \in J} c_j x_j t + \alpha t \tag{1.21}$$

s.t. 
$$\sum_{j \in J} a_{ij} x_j \gtrsim b_i, \quad \forall i \in J$$
 (1.22)

$$\sum_{j \in J} d_j x_j t + \beta t = 1 \tag{1.23}$$

$$t > 0 \tag{1.24}$$

$$x_j \ge 0, \quad \forall j \in J \tag{1.25}$$

$$(t = \frac{1}{\sum_{j \in J} d_j x_j + \beta}) \tag{1.26}$$

For the following statement:

$$\min \quad z^P = \frac{\mathbf{c}^\top \mathbf{x} + d}{\mathbf{e}^\top \mathbf{x} + f} \tag{1.27}$$

s.t. 
$$\mathbf{G}\mathbf{x} \le \mathbf{h}$$
 (1.28)

$$\mathbf{A}\mathbf{x} = \mathbf{b} \tag{1.29}$$

Modeling

$$\min \quad z^R = \mathbf{c}^\top \mathbf{y} + dz \tag{1.30}$$

s.t. 
$$\mathbf{G}\mathbf{y} - \mathbf{h}z \le 0$$
 (1.31)

$$\mathbf{A}\mathbf{y} - \mathbf{b}z = 0 \tag{1.32}$$

$$\mathbf{e}^{\mathsf{T}}\mathbf{y} + fz = 1 \tag{1.33}$$

$$z \ge 0 \tag{1.34}$$

# 1.2.4 A Range Constraint

Consider the following model statement:

$$\min \quad \sum_{j \in J} c_j x_j \tag{1.35}$$

s.t. 
$$d_i \le \sum_{j \in J} a_{ij} x_j \le e_i, \quad \forall i \in I$$
 (1.36)

$$x_j \ge 0, \quad \forall j \in J \tag{1.37}$$

Modeling:

$$\min \quad \sum_{j \in J} c_j x_j, \quad c_j > 0 \tag{1.38}$$

s.t. 
$$u_i + \sum_{j \in I} a_{ij} x_j = e_i, \quad \forall i \in I$$
 (1.39)

$$x_j \ge 0, \quad \forall j \in J$$
 (1.40)

$$0 \le u_i \le e_i - d_i, \quad \forall i \in I \tag{1.41}$$

# 1.3 Integer Programming Formulation Skills

# 1.3.1 A Variable Taking Discontinuous Values

In algebraic notation:

$$x = 0$$
, or  $l \le x \le u$  (1.42)

Modeling:

$$x \le uy \tag{1.43}$$

$$x \ge ly \tag{1.44}$$

$$y \in \{0, 1\} \tag{1.45}$$

where

$$y = \begin{cases} 0, & \text{if } x = 0\\ 1, & \text{if } l \le x \le u \end{cases}$$
 (1.46)

#### 1.3.2 Fixed Costs

In algebraic notation:

$$C(x) = \begin{cases} 0 & \text{for } x = 0\\ k + cx & \text{for } x > 0 \end{cases}$$
 (1.47)

Modeling:

$$C^*(x,y) = ky + cx \tag{1.48}$$

$$x \le My \tag{1.49}$$

$$x \ge 0 \tag{1.50}$$

$$y \in \{0, 1\} \tag{1.51}$$

where

$$y = \begin{cases} 0, & \text{if } x = 0\\ 1, & \text{if } x \ge 0 \end{cases}$$
 (1.52)

#### 1.3.3 Either-or Constraints

In algebraic notation:

$$\sum_{j \in J} a_{1j} x_j \le b_1 \text{ or } \sum_{j \in J} a_{2j} x_j \le b_2$$
 (1.53)

Modeling:

$$\sum_{j \in I} a_{1j} x_j \le b_1 + M_1 y \tag{1.54}$$

$$\sum_{j \in I} a_{2j} x_j \le b_2 + M_1 (1 - y) \tag{1.55}$$

$$y \in \{0, 1\} \tag{1.56}$$

where

$$y = \begin{cases} 0, & \text{if } \sum_{j \in J} a_{1j} x_j \le b_1 \\ 1, & \text{if } \sum_{j \in J} a_{2j} x_j \le b_2 \end{cases}$$
 (1.57)

Notice that the sign before M is determined by the inequality  $\geq$  or  $\leq$ , if it is " $\geq$ ", use "-", if it " $\leq$ ", use "+".

# 1.3.4 Conditional Constraints

If constraint A is satisfied, then constraint B must also be satisfied

If 
$$\sum_{j \in J} a_{1j} x_j \le b_1$$
 then  $\sum_{j \in J} a_{2j} x_j \le b_2$  (1.58)

The key part is to find the opposite of the first condition. We are using  $A \Rightarrow B \Leftrightarrow \neg B \Rightarrow \neg A$ Therefore it is equivalent to

$$\sum_{j \in J} a_{1j} x_j > b_1 \text{ or } \sum_{j \in J} a_{2j} x_j \le b_2$$
 (1.59)

Furthermore, it is equivalent to

$$\sum_{j \in J} a_{1j} x_j \ge b_1 + \epsilon \text{ or } \sum_{j \in J} a_{2j} x_j \le b_2$$
 (1.60)

Where  $\epsilon$  is a very small positive number. Modeling:

$$\sum_{j \in J} a_{1j} x_j \ge b_1 + \epsilon - M_2 y \tag{1.61}$$

$$\sum_{j \in J} a_{2j} x_j \le b_2 + M_2 (1 - y) \tag{1.62}$$

$$y \in \{0, 1\} \tag{1.63}$$

# 1.3.5 Special Ordered Sets

Out of a set of yes-no decisions, at most one decision variable can be yes. Also known as SOS1.

$$x_1 = 1, x_2 = x_3 = \dots = x_n = 0 \tag{1.64}$$

or 
$$(1.65)$$

$$x_2 = 1, x_1 = x_3 = \dots = x_n = 0 \tag{1.66}$$

or ... 
$$(1.67)$$

Modeling:

$$\sum_{i} x_i = 1, \quad i \in N \tag{1.68}$$

Out of a set of binary variables, at most two variables can be nonzero. In addition, the two variables must be adjacent to each other in a fixed order list. Also known as SOS2. Modeling: If  $x_1, x_2, ..., x_n$  is a SOS2, then

$$\sum_{i=1}^{n} x_i \le 2 \tag{1.69}$$

$$x_i + x_j \le 1, \forall i \in \{1, 2, ..., n\}, j \in \{i + 2, i + 3, ..., n\}$$
 (1.70)

$$x_i \in \{0, 1\} \tag{1.71}$$

There is another type of definition, that is out of a set of nonnegative variables **not binary here**, at most two variables can be nonzero. In addition, the two variables must be adjacent to each other in a fixed order list. All variables summing to 1.

This definition of SOS2 is used in Piecewise Linear Formulations.

#### 1.3.6 Piecewise Linear Formulations

The objective function is a sequence of line segments, e.g. y = f(x), consists k-1 linear segments going through k given points  $(x_1, y_1), (x_2, y_2), ..., (x_k, y_k)$ . Denote

$$d_i = \begin{cases} 1, & x \in (x_i, x_{i+1}) \\ 0, & \text{otherwise} \end{cases}$$
 (1.72)

Then the objective function is

$$\sum_{i \in \{1, 2, \dots, k-1\}} y = d_i f_i(x) \tag{1.73}$$

Modeling: Given that objective function as a piecewise linear formulation, we can have these constraints

$$\sum_{i \in \{1, 2, \dots, k-1\}} d_i = 1 \tag{1.74}$$

$$i \in \{1,2,...,k-1\}$$

$$d_i \in \{0, 1\}, i \in \{1, 2, ..., k - 1\}$$

$$(1.75)$$

$$x = \sum_{i \in \{1, 2, \dots, k\}} w_i x_i \tag{1.76}$$

$$y = \sum_{i \in \{1, 2, \dots, k\}} w_i y_i \tag{1.77}$$

$$w_1 \le d_1 \tag{1.78}$$

$$w_i \le d_{i-1} + di, i \in \{2, 3, ..., k - 1\}$$

$$(1.79)$$

$$w_k \le d_{k-1} \tag{1.80}$$

In this case,  $w_i \in SOS2$  (second definition)

#### 1.3.7Conditional Binary Variables

Choose at most n binary variable to be 1 out of  $x_1, x_2, ...x_m, m \ge n$ . If n = 1 then it is SOS1. Modeling:

$$\sum_{k \in \{1, 2, \dots, m\}} x_k \le n \tag{1.81}$$

Choose exactly n binary variable to be 1 out of  $x_1, x_2, ... x_m, m \ge n$ Modeling:

$$\sum_{k \in \{1, 2, \dots, m\}} x_k = n \tag{1.82}$$

Choose  $x_j$  only if  $x_k = 1$ 

Modeling:

$$x_j = x_k \tag{1.83}$$

"and" condition, iff  $x_1, x_2, ..., x_m = 1$  then y = 1Modeling:

$$y \le x_i, i \in \{1, 2, ..., m\} \tag{1.84}$$

$$y \ge \sum_{i \in \{1, 2, \dots, m\}} x_i - (m - 1) \tag{1.85}$$

#### Elimination of Products of Variables 1.3.8

For variables  $x_1$  and  $x_2$ ,

$$y = x_1 x_2 \tag{1.86}$$

Modeling: If  $x_1, x_2$  are binary, it is the same as "and" condition of binary variables.

If  $x_1$  is binary, while  $x_2$  is continuous and  $0 \le x_2 \le u$ , then

$$y \le ux_1 \tag{1.87}$$

$$y \le x_2 \tag{1.88}$$

$$y \ge x_2 - u(1 - x_1) \tag{1.89}$$

$$y \ge 0 \tag{1.90}$$

If both  $x_1$  and  $x_2$  are continuous, it is non-linear, we can use Piecewise linear formulation to simulate.

# Chapter 2

# Basic Concepts of Linear Algebra

# 2.1 Vector Spaces

#### 2.1.1 Field

**Definition 2.1.1** (Field). Let F denote either the set of real numbers or the set of complex numbers.

- Addition is commutative:  $x + y = y + x, \forall x, y \in F$
- Addition is associative:  $x + (y + z) = (x + y) + z, \forall x, y, z \in F$
- Element 0 exists and unique:  $\exists 0, x+0=x, \forall x \in F$
- To each  $x \in F$  there corresponds a unique element  $(-x) \in F$  such that x + (-x) = 0
- Multiplication is commutative:  $xy = yx, \forall x, y \in F$
- Multiplication is associative:  $x(yz) = (xy)z, \forall x, y, z \in F$
- Element 1 exists and unique:  $\exists 1, x1 = x, \forall x \in F$
- To each  $x \neq 0 \in F$  there corresponds a unique element  $x^{-1} \in F$  that  $xx^{-1} = 1$
- • Multiplication distributes over addition:  $x(y+z) = xy + xz, \forall x,y,z \in F$

Suppose one has a set F of objects  $x, y, z, \cdots$  and two operations on the elements of F as following:

- (Addition) associates with each pair of elements  $x, y \in F$  an element  $(x + y) \in F$ ,
- (Multiplication) associates with each pair x, y an element  $xy \in F$ ,

and these two operations satisfy all conditions above. The set F, together with these two operations, is then called a **field**.

**Definition 2.1.2** (Subfield). A subfield of the field C is a set F of complex numbers which itself is a field.

**Example.** The set of rational numbers is a field.

**Example.** The set of integers is **not** a field.

**Example.** The set of all complex numbers of the form  $x + y\sqrt{2}$  where x and y are rational, is a subfield of  $\mathbb{C}$ .

## 2.1.2 Vector Space and Subspace

**Definition 2.1.3** (Vector space). A **vector space** consists of the following:

- A field F of scalars;
- A set V of vectors;

• An addition operation, which associated with each pair of vectors  $\alpha, \beta \in V$  a vector  $\alpha + \beta$  in V called the **sum** of  $\alpha$  and  $\beta$ , in such a way that

$$-\alpha + \beta = \beta + \alpha;$$
  

$$-\alpha + (\beta + \gamma) = (\alpha + \beta) + \gamma;$$
  

$$-\forall \alpha \in V, \alpha + 0 = \alpha$$
  

$$-\forall \alpha \in V, \alpha + (-\alpha) = 0$$

• A multiplication operation, which associated with each scalar  $c \in F$  and a vector  $\alpha \in V$  a vector  $c\alpha \in V$  called the **product** of c and  $\alpha$ , in such a way that

$$-1\alpha = \alpha, \forall \alpha \in V$$

$$-(c_1c_2)\alpha = c_1(c_2\alpha)$$

$$-c(\alpha + \beta) = c\alpha + c\beta$$

$$-(c_1 + c_2)\alpha = c_1\alpha + c_2\alpha$$

we may simply denote the vector space as V, which is the same notation as the set of vectors, if the field F of scalars needs to be specified, we shall say V is a vector space over the field F.

**Definition 2.1.4** (Subspace). Let V be a vector space over the field F. A subspace of V is a subset H of V which is itself a vector space over F with the operations of vector addition and scalar multiplication on H

The definition of subspace implies  $\lambda x \in H, \forall \lambda \in F$  and if  $x, y \in V, x + y \in H$ . The following is equivalent:

- $H \subseteq \mathbb{R}^n$  is a subspace
- There is an  $m \times n$  matrix A such that  $H = \{x \in \mathbb{R}^n | Ax = 0\}$
- There is a  $k \times n$  matrix B such that  $H = \{x \in \mathbb{R}^n | x = uB, u \in \mathbb{R}^k\}$

Particularly,

**Definition 2.1.5** (Orthogonal subspace). For a subspace H, then  $\{x \in \mathbb{R}^n | xy = 0, y \in H\}$  is a **orthogonal** subspace and denoted by  $H^{\perp}$ 

**Proposition 1.** If  $H = \{x \in \mathbb{R}^n | Ax = 0\}$ , with A being an  $m \times n$  matrix, then  $H^{\perp} = \{x \in \mathbb{R}^n | x = A^{\top}u, u \in \mathbb{R}^m\}$ 

## 2.1.3 Linear, Conic, Affine, and Convex Combinations

**Proposition 2.** The following statements are equivalent:

- $x^1, x^2, \cdots, x^k \in \mathbb{R}^n$  are affinely independent
- $x^2 x^1, x^3 x^1, \dots, x^k x^1$  are linearly independent
- $\begin{bmatrix} x^1 \\ 1 \end{bmatrix}$ ,  $\begin{bmatrix} x^2 \\ 1 \end{bmatrix}$ ,  $\cdots$ ,  $\begin{bmatrix} x^k \\ 1 \end{bmatrix}$  are linearly independent

## 2.2 Determinants

#### 2.2.1

# 2.3 Inner Products

**Definition 2.3.1** (Inner Product). Let F be the field of real numbers or the field of complex numbers, and V a vector space over F. An **inner product** on V is a function which assigns to each ordered pair of vectors  $\alpha$ ,  $\beta$  in V a scalar  $< \alpha | \beta >$  in F in such a way that  $\forall \alpha, \beta, \gamma \in V, c \in \mathbb{R}$  that

• 
$$<\alpha + \beta | \gamma > = <\alpha | \gamma > + <\beta | \gamma >$$

2.4. NORMS 25

- $< c\alpha |\beta> = c < \alpha |\beta>$
- $<\alpha|\beta>=\overline{<\beta|\alpha>}$
- $\langle \alpha | \alpha \rangle \geq 0$ ,  $\langle \alpha | \alpha \rangle = 0$  iff  $\alpha = \mathbf{0}$

Furthermore, the above properties imply that

•  $<\alpha|c\beta+\gamma>=\bar{c}<\alpha|\beta>+<\alpha|\gamma>$ 

**Definition 2.3.2.** On  $F^n$  there is an inner product which we call the **standard inner product**. It is defined on  $\alpha = (x_1, x_2, ..., x_n)$  and  $\beta = (y_1, y_2, ..., y_n)$  by

$$<\alpha|\beta> = \sum_{j} x_{j} \bar{y_{j}}$$
 (2.1)

For  $F = \mathbb{R}^n$ 

$$<\alpha|\beta> = \sum_{j} x_j y_j$$
 (2.2)

In the real case, the standard inner product is often called the dot product and denoted by  $\alpha \cdot \beta$ 

**Example.** For  $\alpha = (x_1, x_2)$  and  $\beta = (y_1, y_2)$  in  $\mathbb{R}^2$ , the following is an inner product.

$$<\alpha|\beta> = x_1y_1 - x_2y_1 - x_1y_2 + 4x_2y_2$$
 (2.3)

**Example.** For  $\mathbb{C}^{n\times n}$ ,

$$\langle \mathbf{A}|\mathbf{B} \rangle = trace(\mathbf{B}^*\mathbf{A})$$
 (2.4)

is an inner product, where

$$\mathbf{A}_{ij}^* = \bar{\mathbf{A}}_{ji}$$
 (conjugate transpose) (2.5)

For  $\mathbb{R}^{n \times n}$ ,

$$\langle \mathbf{A} | \mathbf{B} \rangle = trace(\mathbf{B}^T \mathbf{A}) = \sum_{j} (AB^T)_{jj} = \sum_{j} \sum_{k} A_{jk} B_{jk}$$
 (2.6)

# 2.4 Norms

**Definition 2.4.1** (Norms). A **norm** on a vector space  $\mathcal{V}$  is a function  $\|\cdot\|:\mathcal{V}\to\mathbb{R}$  for which the following three properties hold for all point  $\mathbf{x},\mathbf{y}\in\mathcal{V}$  and scalars  $\lambda\in\mathbb{R}$ 

- (Absolute homogeneity)  $\|\lambda \mathbf{x}\| = |\lambda| \|\mathbf{x}\|$
- (Triangle inequality)  $\|\mathbf{x} + \mathbf{y}\| \le \|\mathbf{x}\| + \|\mathbf{y}\|$
- (Positivity) Equality  $\|\mathbf{x}\| = 0$  holds iff  $\mathbf{x} = 0$

**Definition 2.4.2** ( $L_p$ -norms). Let  $p \ge 1$  be a real number. We define the p-norm of vector  $\mathbf{v} \in \mathbb{R}^n$  as:

$$\|\mathbf{x}\|_{p} = \left(\sum_{i=1}^{n} |v_{i}|^{p}\right)^{\frac{1}{p}} \tag{2.7}$$

Particularly

$$\|\mathbf{v}\|_1 = \sum_{i=1}^n |v_i| \tag{2.8}$$

$$\|\mathbf{v}\|_2 = \sqrt{\sum_{i=1}^n v_i^2} \tag{2.9}$$

$$\|\mathbf{v}\|_{\infty} = \max_{i=1}^{n} |v_i| \tag{2.10}$$

**Definition 2.4.3** (Frobenius norm).  $\mathbf{X} \in \mathbb{R}^{m \times n}$ , the **Frobenius norm** is defined as

$$\|\mathbf{X}\|_F = \sqrt{trace(\mathbf{X}^\top \mathbf{X})} \tag{2.11}$$

**Definition 2.4.4** (Dual norm). For an arbitrary norm  $\|\cdot\|$  on Euclidean space **E**, the **dual norm**  $\|\cdot\|^*$  on **E** is defined by

$$\|\mathbf{v}\|^* = \max\{\langle \mathbf{v}|\mathbf{x} \rangle | \|\mathbf{x}\| \le 1\}$$
 (2.12)

For  $p, q \in [1, \infty]$ , the  $l_p$  and  $l_q$  norms on  $\mathbb{R}^n$  are dual to each other whenever  $\frac{1}{p} + \frac{1}{q} = 1$ .

# 2.5 Eigenvectors and Eigenvalues

**Definition 2.5.1.** If **A** is an  $n \times n$  matrix, then a nonzero vector  $\mathbf{x} \in \mathbb{R}^n$  is called an **eigenvector** of **A** if  $\mathbf{A}\mathbf{x}$  is a scalar multiple of  $\mathbf{x}$ , i.e.

$$\mathbf{A}\mathbf{x} = \lambda \mathbf{x} \tag{2.13}$$

for some scalar  $\lambda$ . The scalar  $\lambda$  is called **eigenvalue** of **A** and the vector **x** is said to be an **eigenvector corresponding to**  $\lambda$ 

**Theorem 2.1** (Characteristic Equation). If **A** is an  $n \times n$  matrix, then  $\lambda$  is an eigenvalue of **A** iff

$$\det(\lambda I - A) = 0 \tag{2.14}$$

Corollary 2.1.1.

$$\sum \lambda_A = tr(\mathbf{A}) \tag{2.15}$$

Corollary 2.1.2.

$$\prod \lambda_A = \det(\mathbf{A}) \tag{2.16}$$

**Notice:** Gaussian elimination changes the eigenvalues.

# 2.6 Decompositions

# Chapter 3

# Basic Concepts of Convex Analysis

# 3.1 Convex Sets

**Definition 3.1.1** (convex set, convex combination). A set X in  $\mathbb{R}^n$  is called a **convex set** if given any two points  $\mathbf{x}_1 \in X$  and  $\mathbf{x}_2 \in X$ , then  $\lambda \mathbf{x}_1 + (1 - \lambda)\mathbf{x}_2 \in X$ ,  $\forall \lambda \in [0, 1]$ . Any point of the form  $\lambda \mathbf{x}_1 + (1 - \lambda)\mathbf{x}_2$  where  $0 \le \lambda \le 1$  is called a **convex combination** of  $\mathbf{x}_1$  and  $\mathbf{x}_2$ . If  $\lambda \in (0, 1)$ , then the convex combination is called **strict**.

**Example.**  $S = \{(x_1, x_2) | 3x_1^2 + 4x_2^2 \le 1\}$ 

Example.  $X = \{\mathbf{x} | \mathbf{A}_{m \times n} \mathbf{x} = \mathbf{b}_m \}$ 

The following are some families of convex sets.

**Example.** Empty set is by convention considered as convex.

**Example.** Polyhedrons are convex sets.

**Example.** Let  $P = \{ \mathbf{x} \in \mathbb{R}^n | \mathbf{x}^\top \mathbf{A} \mathbf{x} \leq \mathbf{b} \}$  where  $\mathbf{A} \in \mathbb{S}_+^{n \times n}$  and  $\mathbf{b} \in \mathbb{R}_+$ . The set P is a convex subset of  $\mathbb{R}^n$ .

**Example.** Let  $\|.\|$  be any norm in  $\mathbb{R}^n$ . Then, the unit ball  $B = \{\mathbf{x} \in \mathbb{R}^n | \|\mathbf{x}\| \le b, b > 0\}$  is convex.

Let  $S_1, S_2$  be convex set, then:

- $S_1 \cap S_2$  is convex set
- $S_1 \oplus S_2$  (Minkowski addition) is convex set, where

$$S_1 \oplus S_2 = \{ \mathbf{x} \in \mathbb{R}^n | \mathbf{x} = \mathbf{x_1} + \mathbf{x_2}, \mathbf{x_1} \in S_1, \mathbf{x_2} \in S_2 \}$$
 (3.1)

•  $S_1 \ominus S_2$  is convex set, where

$$S_1 \oplus S_2 = \{ \mathbf{x} \in \mathbb{R}^n | \mathbf{x} = \mathbf{x_1} - \mathbf{x_2}, \mathbf{x_1} \in S_1, \mathbf{x_2} \in S_2 \}$$
 (3.2)

•  $f(S_1)$  is convex iff  $f(\mathbf{x}) = \mathbf{A}\mathbf{x} + \mathbf{b}, \mathbf{A} \in \mathbb{R}^{m \times n}, \mathbf{b} \in \mathbb{R}^m$ 

**Definition 3.1.2** (convex hull). Let S be an arbitrary set in  $\mathbb{R}^n$ . The **convex hull** of S, denoted by conv(S) is the collection of all convex combinations of elements in S

$$\mathbf{x} \in conv(S) \iff \mathbf{x} = \sum_{j=1}^{k} \lambda_j \mathbf{x}_j$$
 (3.3)

$$\sum_{j=1}^{k} \lambda_j = 1 \tag{3.4}$$

$$\lambda_j \le 0, \quad j = 1, \cdots, k \tag{3.5}$$

$$\mathbf{x}_1, \mathbf{x}_2, \cdots, \mathbf{x}_k \in S \tag{3.6}$$

**Lemma 3.1.** Let S be arbitrary set in  $\mathbb{R}^n$ . Then conv(S) is the smallest convex set containing S, which means conv(S) is the intersection of all convex sets containing S.

**Theorem 3.2** (Carathéodory's Theorem). Let  $S \subseteq \mathbb{R}^n$ . Then  $\forall \mathbf{x} \in conv(S)$ , there exists  $\mathbf{x}^1, \mathbf{x}^2, ..., \mathbf{x}^p \in S$ , where  $p \leq n+1$  such that  $\mathbf{x} \in conv\{\mathbf{x}^1, \mathbf{x}^2, ..., \mathbf{x}^p\}$ .

**Notice:** This theorem means, any point  $\mathbf{x} \in \mathbb{R}^n$  in a convex hull of S, i.e., conv(S), can be included in a convex subset  $S' \subseteq conv(S)$  that has n+1 extreme points.

**Theorem 3.3.** Let S be a convex set with nonempty interior. Let  $\mathbf{x}_1 \in cl(S)$  and  $\mathbf{x}_2 \in int(S)$ , then  $\mathbf{y} = \lambda \mathbf{x}_1 + (1 - \lambda)\mathbf{x}_2 \in int(S), \forall \lambda \in (0, 1)$ 

## 3.2 Convex Functions

**Definition 3.2.1.** Let  $C \in \mathbb{R}^n$  be a convex set. A function  $f: C \to \mathbf{R}$  is (resp. strictly) convex if

$$f(\lambda \mathbf{x}_1 + (1 - \lambda)\mathbf{x}_2) \le \lambda f(\mathbf{x}_1) + (1 - \lambda)f(\mathbf{x}_2) \tag{3.7}$$

$$\forall \mathbf{x}_1, \mathbf{x}_2 \in C, \forall \lambda \in (0, 1) \tag{3.8}$$

(resp.)

$$f(\lambda \mathbf{x}_1 + (1 - \lambda)\mathbf{x}_2) < \lambda f(\mathbf{x}_1) + (1 - \lambda)f(\mathbf{x}_2)$$
(3.9)

$$\forall \mathbf{x_1} \neq \mathbf{x}_2 \in C, \forall \lambda \in (0,1) \tag{3.10}$$

**Notice:** When calling a function convex, we imply that its domain is convex.

**Example.** Given any norm  $\|.\|$  on  $\mathbb{R}^n$ , the function  $f(x) = \|x\|$  is convex over  $\mathbb{R}^n$ .

**Definition 3.2.2.** Let S be a nonempty convex subset of  $\mathbb{R}^n$ ,  $f: S \to \mathbb{R}$  is (resp. strictly) **concave** if -f(x) is (resp. strictly) convex.

**Notice:** A function may be neither convex nor concave.

**Theorem 3.4.** Consider  $f: \mathbb{R}^n \to \mathbb{R}$ .  $\forall \bar{\mathbf{x}} \in \mathbb{R}^n$  and a nonzero direction  $\mathbf{d} \in \mathbb{R}^n$ . Define  $F_{\bar{\mathbf{x}},d}(\lambda) = f(\bar{\mathbf{x}} + \lambda \mathbf{d})$ . Then f is (resp. strictly) convex iff  $F_{\bar{\mathbf{x}},d}(\lambda)$  is (resp. strictly) convex for all  $\bar{\mathbf{x}} \in \mathbb{R}^n$ ,  $\forall \mathbf{d} \in \mathbb{R}^n \setminus \{0\}$ .

**Definition 3.2.3** (Level-set). Given a function  $f: \mathbb{R}^n \to \mathbb{R}$  and a scalar  $\alpha \in \mathbb{R}$ , we refer to the set  $S_{\alpha} = \{\mathbf{x} \in S | f(\mathbf{x}) \leq \alpha\} \subset \mathbb{R}^n$  as the  $\alpha$ -level-set of f.

**Lemma 3.5.** Let S be a nonempty convex set in  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}^n$  be a convex function, then the  $\alpha$ -level-set of f is a convex set for each value of  $\alpha \in \mathbb{R}$ .

**Notice:** The converse is not necessarily true.

**Definition 3.2.4** (Epigraphs, Hypographs). Let  $S \in \mathbb{R}^n$  be such that  $S \neq \emptyset$ . The **epigraph** of f, denoted by epi(f) is

$$epi(f) = \{(\mathbf{x}, y) \in S | \mathbf{x} \in S, y \in \mathbb{R}, y \ge f(x)\} \in \mathbb{R}^{n+1}$$

$$(3.11)$$

The **hypograph** of f, denoted by hypo(f) is

$$hypo(f) = \{(\mathbf{x}, y) \in S | \mathbf{x} \in S, y \in \mathbb{R}, y \le f(x)\} \in \mathbb{R}^{n+1}$$
(3.12)

**Theorem 3.6.** Let S be a nonempty convex subset in  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$ . Then f is convex iff epi(f) is convex.

**Theorem 3.7.** Let S be a nonempty convex subset in  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be a convex function on S. Then f is continuous in int(S).

# 3.3 Subgradients and Subdifferentials

**Definition 3.3.1** (Subgradient). Let S be a nonempty convex set in  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be a convex function, then  $\xi$  is a subgradient of f at  $\bar{\mathbf{x}}$  if

$$f(\mathbf{x}) \ge f(\bar{\mathbf{x}}) + \xi^{\top}(\mathbf{x} - \bar{\mathbf{x}}), \forall \mathbf{x} \in S$$
 (3.13)

**Definition 3.3.2** (Subdifferential). The set of all subgradients of f at  $\bar{\mathbf{x}}$  is called **subdifferential** of f at  $\bar{\mathbf{x}}$ , denoted as  $\partial f(\bar{\mathbf{x}})$ 

**Theorem 3.8.** Let S be a nonempty convex set in  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be a convex function. Then for  $\bar{\mathbf{x}} \in int(S)$ , there exists a vector  $\xi$  such that

$$f(\mathbf{x}) \ge f(\bar{\mathbf{x}}) + \xi^{\top}(\mathbf{x} - \bar{\mathbf{x}}), \forall \mathbf{x} \in S$$
 (3.14)

In particular, the hyperplane

$$\mathcal{H} = \{ (\mathbf{x}, y) | y = f(\bar{\mathbf{x}}) + \xi^{\top} (\mathbf{x} - \bar{\mathbf{x}}) \}$$
(3.15)

is a supporting plane of epi(f) at  $(\bar{\mathbf{x}}, f(\bar{\mathbf{x}}))$ 

**Theorem 3.9.** Let S be a nonempty convex set in  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be a convex function. Suppose that for each  $\bar{\mathbf{x}} \in S$ , there exists  $\xi$  such that

$$f(\mathbf{x}) \ge f(\bar{\mathbf{x}}) + \xi^{\top}(\mathbf{x} - \bar{\mathbf{x}}), \forall \mathbf{x} \in S$$
 (3.16)

Then f is convex on int(S)

**Notice:** Not all convex functions are continuous, it has to be continuous in its interior, but it may not be continuous at the boundary.

# 3.4 Differentiable Functions

**Definition 3.4.1** (Differentiable Functions). Let S be a nonempty subset of  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$ . Then f is said to be **differentiable** at  $\bar{\mathbf{x}} \in int(S)$  if there exists a vector  $\nabla f(\bar{\mathbf{x}})$  and a function  $\alpha: \mathbb{R}^n \to \mathbb{R}$  such that

$$f(\mathbf{x}) = f(\bar{\mathbf{x}}) + \nabla f(\bar{\mathbf{x}})^{\top} (\mathbf{x} - \bar{\mathbf{x}}) + \alpha(\bar{\mathbf{x}}, \mathbf{x} - \bar{\mathbf{x}}) \|\mathbf{x} - \bar{\mathbf{x}}\|$$
(3.17)

for all  $\mathbf{x} \in S$  where  $\lim_{\mathbf{x} - \bar{\mathbf{x}}} \alpha(\bar{\mathbf{x}}, \mathbf{x} - \bar{\mathbf{x}}) = 0$ 

Remark. If function f is differentiable, then  $\nabla f(\bar{\mathbf{x}}) = (\frac{\partial f(\bar{\mathbf{x}})}{\partial \mathbf{x}_1}, \frac{\partial f(\bar{\mathbf{x}})}{\partial \mathbf{x}_2}, \cdots, \frac{\partial f(\bar{\mathbf{x}})}{\partial \mathbf{x}_n})$ , and the gradient is unique.

**Lemma 3.10.** Let  $S \neq \emptyset$  be a convex set of  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be convex. If f is differentiable at  $\bar{\mathbf{x}} \in int(S)$ , then the subdifferential of f at  $\bar{\mathbf{x}}$  is the singleton,  $\{\nabla f(\bar{\mathbf{x}})\}$ 

**Theorem 3.11.** Let S be a nonempty subset of  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be differentiable on S. Then f is (resp. strictly) convex on S iff  $\forall \bar{\mathbf{x}} \in S$ 

$$f(\mathbf{x}) \ge f(\bar{\mathbf{x}}) + \nabla f(\bar{\mathbf{x}})(\mathbf{x} - \bar{\mathbf{x}}), \forall \mathbf{x} \in S$$
 (3.18)

(resp.)

$$f(\mathbf{x}) > f(\bar{\mathbf{x}}) + \nabla f(\bar{\mathbf{x}})(\mathbf{x} - \bar{\mathbf{x}}), \forall \mathbf{x} \neq \bar{\mathbf{x}} \in S$$
 (3.19)

**Theorem 3.12** (Mean-value Theorem). Let S be a nonempty subset of  $\mathbb{R}^n$ . Let  $f: S \to \mathbb{R}$  be differentiable on S. Then for all  $\mathbf{x}_1, \mathbf{x}_2 \in S$ , there exists  $\lambda \in (0,1)$  such that

$$f(\mathbf{x}_2) = f(\mathbf{x}_2) + \nabla f(\hat{\mathbf{x}})(\mathbf{x}_2 - \mathbf{x}_1)$$
(3.20)

where

$$\hat{\mathbf{x}} = \lambda \mathbf{x}_1 + (1 - \lambda)\mathbf{x}_2 \tag{3.21}$$

# 3.5 Convex and Affine Hulls

# 3.6 Interior and Closure

# 3.7 Cones

# 3.8 Hyperplanes

# Chapter 4

# Basic Concepts of Polyhedral Theory

# 4.1 Extreme Points

**Definition 4.1.1** (extreme points). Let  $S \subseteq \mathbb{R}^n$ , an extreme point of S is a point  $x \in S$  such that

$$\begin{cases} x = \sum_{i=1}^{m} \lambda_i x^i \\ \sum_{i=1}^{m} \lambda_i = 1 \\ \lambda_i > 0, i = 1, \dots, m \\ x^i \in S, i = 1, \dots, m \end{cases}$$
 implies  $x^i = x, i = 1, \dots, m$  (4.1)

which means, if x is an extreme point, it cannot be written as a convex combination of points in S except the copies of x itself.

**Theorem 4.1.** Let P be a polyhedron in  $\mathbb{R}^n$  and  $v \in P$ . Then v is an extreme point of P iff  $dim(span\{a^i|(a^i)^topv = b_i\}) = n$ 

This theorem means, at an extreme point of a n-dimensional polyhedron, there will be n active inequalities. A direct result will be the equivalence between basic feasible solutions and extreme points in Linear Programming.

## 4.2 Polar Cones

# 4.3 Polyhedral and Dimension

## 4.3.1 Polyhedral

**Definition 4.3.1** (polyhedron). A **polyhedron** is a set of the form  $\{x \in \mathbb{R}^n | Ax \leq b\} = \{x \in \mathbb{R}^n | a^i x \leq b^i, \forall i \in M\}$ , where  $A \in \mathbb{R}^{m \times n}$  and  $b \in \mathbb{R}^m$ 

**Proposition 3.** A polyhedron is a convex set.

**Definition 4.3.2** (polytope). A polyhedron  $P \subset \mathbb{R}^n$  is **bounded** if there exists a constant K such that  $|x_i| < K, \forall x \in P, \forall i \in [1, n]$ , in this case the polyhedron is call **polytope**. The lower-bound of K is called **diagonal** denoted by d

**Definition 4.3.3** (cone).  $C \subseteq \mathbb{R}^n$  is a **cone** if  $x \in C$  implies  $\lambda x \in C, \forall \lambda \in \mathbb{R}_+$ 

# 4.3.2 Dimension of Polyhedral

**Definition 4.3.4** (dimension). A polyhedron P is **dimension** k, denoted dim(P) = k, if the maximum number of affinely independent points in P is k + 1

**Definition 4.3.5** (full-dimensional). A polyhedron  $P \subseteq \mathbb{R}^n$  is full-dimensional if dim(P) = n

**Proposition 4.** If  $P \subseteq \mathbb{R}^n$ , then  $dim(P) = n - rank(A^=, b^=)$ 

To proof a constraint  $(A^{=}, b^{=})$  is an equality constraint, we need to proof all point in the closure of P satisfied the constraint, to proof it is not an equality constraint, we need to find one point that is not in the hyperplane.

**Definition 4.3.6** (inner point, interior point).  $x \in P$  is called an **inner point** of P if  $a^i x < b_i, \forall i \in M^{\leq}, x \in P$  is called an **interior point** of P if  $a^i x < b_i, \forall i \in M$ 

Corollary 4.1.1. Every nonempty polyhedron has at least one inner point.

Corollary 4.1.2. A polyhedron has an interior point iff P is full-dimensional, i.e., there is no equality constraint.

## 4.4 Face and Facet

# 4.4.1 Valid Inequalities and Faces

The inequality denoted by  $(\pi, \pi_o)$  is called a **valid inequality** for P if  $\pi x \leq \pi_0, \forall x \in P$ . Note that  $(\pi, \pi_0)$  is a valid inequality iff P lies in the half-space  $\{x \in \mathbb{R}^n | Ax \leq b\}$ 

- If  $(\pi, \pi_0)$  is a valid inequality for P and  $F = \{x \in P | \pi x = x_0\}$ , F is called a **facet** of P and we say that  $(\pi, \pi_0)$ 

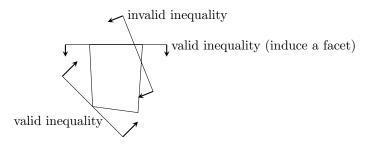


Figure 4.1: Example of valid/invalid inequality

#### represents or defines F

- A face is said to be **proper** if  $F \neq \emptyset$  and  $F \neq P$
- The face represented by  $(\pi, \pi_0)$  is nonempty iff  $\max\{\pi x | x \in P\} = \pi_0\}$
- If the face F is nonempty, we say it supports P
- Let P be a polyhedron with equality set  $M^{=}$ . If

$$F = \{x \in P | \pi^T x = \pi_0\} \tag{4.2}$$

is not empty, then F is a polyhedron. Let

$$M^{=} \subseteq M_F^{=}, M_F^{\leq} = M \setminus M_F^{=} \tag{4.3}$$

then

$$F = \{x | a_i^T x = b_i, \forall i \in M_F^=, a_i^T x \le b_i, \forall i \in M_f^{\le} \}$$
(4.4)

# 4.4.2 Facet

- A face F is said to be a **facet** of P if dim(F) = dim(P) 1
- Facets are all we need to describe polyhedral
- If F is a facet of P, then in any description of P, there exists some inequality representing F
- Every inequality that represents a face that is not a facet is unnecessary in the description of P Every full-dimensional polyhedron P has a unique (up to scalar multiplication) representation that consists of one inequality representing each facet of P
- If dim(P) = n k with k > 0, then P is described by a maximal set of linearly independent rows of  $(A^{=}, b^{=})$ , as well as one inequality representing each facet of P

# 4.4.3 Proving Facet

To prove an inequality  $\sum_i a_i x_i \le b_i$  is facet inducing for a D dimensional polyhedral, we need to prove there are D affinely independent vectors in  $\sum_i a_i x_i = b_i$ 

4.4. FACE AND FACET 33

# 4.4.4 Domination

 $\Pi x \leq \Pi_0$  dominates  $Mx \leq M_0$  if

$$\begin{cases}
\Pi \ge \mu M, \mu > 0 \\
\Pi_0 \le \mu M_0, \mu > 0 \\
(\Pi, \Pi_0) \ne (M, M_0)
\end{cases}$$
(4.5)

# Chapter 5

# Basic Concepts of Real Analysis

# 5.1 The Real Number System

# 5.2 Open Sets and Closed Sets

**Definition 5.2.1** (neighborhood).  $N_{\epsilon} = \{y \in \mathbb{R}^n | ||y - x|| < \epsilon \}$  as the **neighborhood** of  $x \in \mathbb{R}^n$ 

**Definition 5.2.2** (interior). Given  $S \subseteq \mathbb{R}^n$ , x belongs to the **interior** of S, denoted by int(S) if there is  $\epsilon > 0$  such that  $N_{\epsilon}(x) \subseteq S$ 

**Definition 5.2.3** (boundary). x belongs to the **boundary**  $\partial S$  if  $\forall \epsilon > 0$ ,  $N_{\epsilon}(x)$  contains at least one point in S and a point not in S

**Definition 5.2.4** (closure).  $x \in S$  belongs to the closure of S, denoted cl(s) if  $\forall \epsilon > 0$ ,  $N_{\epsilon}(x) \cap S = \emptyset$ 

**Definition 5.2.5** (metric space). A **metric space** is a set X where we have a notion of distance. That is, if  $x, y \in X$ , then d(x, y) is the distance between x and y. The particular distance function must satisfy the following conditions:

- $d(x,y) > 0, \forall x,y \in X$
- $d(x,y) = 0 \iff x = y$
- d(x,y) = d(y,x)
- $d(x,z) \leq d(x,y) + d(y,z)$

**Definition 5.2.6** (ball). Let X be a metric space. A ball B of radius r around a point  $x \in X$  is

$$B = \{ y \in X | d(x, y) < r \} \tag{5.1}$$

**Definition 5.2.7** (open set). A subset  $O \subseteq X$  is **open** if  $\forall x \in O, \exists r, B = \{x \in X | d(x,y) < r\} \subseteq O$ 

S is said to be an **open set** iff S = int(S)

**Theorem 5.1.** The union of any collection if open sets is open.

*Proof.* Sets  $S_1, S_2, ..., S_n$  are open sets, let  $S = \bigcup_{i=1}^n S_i$ , then  $\forall i, S_i \subseteq S$ .  $\forall x \in S, \exists i, x \in S_i$ . Given that  $S_i$  is an open set, then for x,  $\exists r$  that  $B = \{x \in S_i | d(x, y) < r\} \subseteq S_i \subseteq S$ , therefore S is an open set.  $\Box$ 

**Theorem 5.2.** The intersection of any finite number of open sets is open.

Proof. Sets  $S_1, S_2, ..., S_n$  are open sets, let  $S = \bigcap_{i=1}^n S_i$ , then  $\forall i, S \subseteq S_i$ .  $\forall x \in S, x \in S_i$ . For any i, we can define an  $r_i$ , such that  $B_i = \{x \in S_i | d(x,y) < r_i\} \subseteq S_i$ . Let  $r = \min_i \{r_i\}$ . Noticed that  $\forall i, B' = \{x \in S_i | d(x,y) < r\} \subseteq S_i$ . Therefore S is an open set.

Remark. The intersection of infinite number of open sets is not necessarily open.

Here we find an example that the intersection of infinite number of open sets can be closed.

**Example.** Let  $A_n \in \mathbb{R}$  and  $B_n \in \mathbb{R}$  be two infinite series, with the following properties. First,  $\forall n, A_n < a, \lim A_n = a$ , second,  $\forall n, B_n > b, \lim B_n = b$ , third a < b. Then we define infinite number of sets  $S_i$ , the *i*th set is defined as

$$S_i = (A_i, B_i) \subset \mathbb{R} \tag{5.2}$$

Then

$$S = \bigcap_{i=1}^{\infty} S_i = [a, b] \subset \mathbb{R}$$

$$(5.3)$$

and S is a closed set.

**Definition 5.2.8** (limit point). A point z is a **limit point** for a set A if every open set U that  $z \in U$  intersects A in a point other than z.

**Notice:** z is not necessarily in A.

**Definition 5.2.9** (closed set). A set C is **closed** iff it contains all of its limit points.

S is called **closed** iff S = cl(S)

**Theorem 5.3.**  $S \in \mathbb{R}^n$  is closed  $\iff \forall \{x_k\}_{k=1}^{\infty} \in S, \lim_{k \to \infty} \{x_k\}_{k=1}^{\infty} \in S$ 

**Theorem 5.4.** Every intersection of closed sets is closed.

**Theorem 5.5.** Every finite union of closed sets is closed.

Remark. The union of infinite number of closed sets is not necessarily closed.

**Theorem 5.6.** A set C is a closed set if  $X \setminus C$  is open

Proof. Let S be an open set,  $x \notin S$ , for any open set  $S_i$  that  $x \in S_i$ , we can find a correspond  $r_i > 0$ , such that  $B_i = \{x \in S_i | d(x,y) < r_i\}$ . Take  $r = \min_{\forall i} \{r_i\}$ , set  $B = \{x \notin S | d(x,y) < r\} \neq \emptyset$ . Which means for any  $x \notin S$ , we can find at least one point  $x' \in B$  that for all open set  $S_i$ ,  $x' \in S_i$ , which makes x a limit point of the complement of the open set. Notice that x is arbitrary, then the collection of x, i.e., the complement of S is a closed set.

Remark. The empty set is open and closed, the whole space X is open and closed.

- 5.3 Functions, Sequences, Limits and Continuity
- 5.4 Differentiation
- 5.5 Integration
- 5.6 Infinite Series of Constants
- 5.7 Power Series
- 5.8 Uniform Convergence
- 5.9 Arcs and Curves
- 5.10 Partial Differentiation
- 5.11 Multiple Integrals
- 5.12 Improper Integrals
- 5.13 Fourier Series

## Basic Concepts of Probability Theory

#### 6.1 Relationship between Some Random Variables

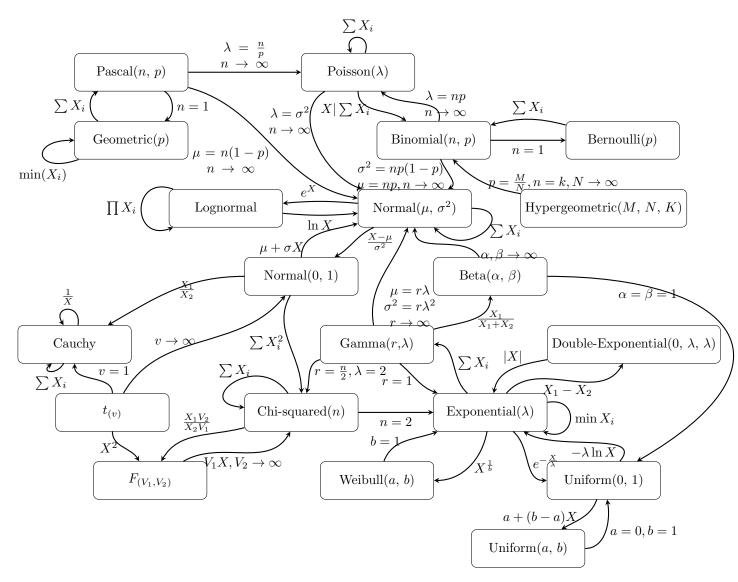


Figure 6.1: Relationship between Some Random Variables

#### 6.2 Discrete Random Variables

Distribution	PMF	$\operatorname{CDF}$	Exp.	Var.	MGF
$\operatorname{Uniform}(a,b)$	x = a, a+1,, b	x = a, a+1,, b	$\frac{b-a}{2}$	$\frac{(b-a+1)^2-1}{12}$	$ \frac{e^{at} - e^{(b+1)t}}{(b-a+1)(1-e^t)} \\ t \in \mathbb{R} $
Bernoulli(p)	$p^x(1-p)^{1-x}$ $x \in \{0,1\}$	$\begin{cases} 0, & x < 0 \\ 1 - p, & 0 \le x \le 1 \\ 1, & x > 1 \end{cases}$	p	p(1-p)	$1 - p + pe^t$ $t \in \mathbb{R}$
Binomial(n, p)	$\binom{n}{x} p^x (1-p)^{n-x}$ $x = 0, 1,, n$	$\sum_{k=0}^{x} \binom{n}{k} p^{k} (1-p)^{n-k}$ $x = 0, 1,, n$	np	np(1-p)	$ \left  \begin{array}{c} (1 - p + pe^t)^n \\ t \in \mathbb{R} \end{array} \right  $
$Poisson(\mu)$	x = 0, 1,, n $x = 0, 1,, n$ $x = 0, 1,, n,$	x = 0, 1,, n,	μ	$\mu$	$e^{\mu(e^t - 1)}$ $t \in \mathbb{R}$
Geometric(p)	$p(1-p)^{x}  x = 0, 1,, n,$	$1 - (1 - p)^{x+1}$ x = 0, 1,, n,	$\frac{1-p}{p}$	$\frac{1-p}{p^2}$	$\begin{array}{ c c }\hline & & & \\ \hline & & \hline & & \\ & & \hline & & \\ & & & \hline & & \\ & & & \hline & \\ & & & \\ & & & \hline & \\ & & & \\ & & & \hline & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & \\ & & \\ & $
$\operatorname{Pascal}(n,p)$	$\binom{n-1+x}{x} p^{n} (1-p)^{x}$ $x = 0, 1, 2,, n,$	$1 - I_p(k+1, n)$ x = 0, 1, 2,, n,	$\frac{n(1-p)}{p}$	$\frac{n(1-p)}{p^2}$	$ \begin{vmatrix} \left(\frac{p}{1 - (1 - p)e^t}\right)^n \\ t < -\ln(1 - p) \end{vmatrix} $
					(6.1

### 6.3 Continuous Random Variables

Distribution	PDF	CDF	Exp.	Var.	MGF	
Uniform $(a, b)$	$x = \begin{bmatrix} \frac{1}{b-a} \\ a, b \end{bmatrix}$	$x = \begin{bmatrix} \frac{x-a}{b-a} \\ x = [a, b] \end{bmatrix}$	$\frac{b-a}{2}$	$\frac{(b-a)^2}{12}$	$\begin{cases} 1, & t = 0\\ \frac{e^{bt} - e^{at}}{t(b-a)}, & t \neq 0 \end{cases}$	
$Normal(\mu, \sigma)$	$\frac{1}{\sqrt{2\pi}\sigma}e^{-\frac{(x-\mu)^2}{2\sigma^2}}$ $x \in \mathbb{R}$	$\int_{-\infty}^{x} \frac{1}{\sqrt{2\pi}\sigma} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$ $x \in \mathbb{R}$	$\mu$	$\sigma^2$	$e^{\frac{t(t\sigma^2+2\mu)}{2}}$ $t \in \mathbb{R}$	(6.2)
Exponential( $\lambda$ )	$\lambda e^{-\lambda x}$ $x > 0$	$1 - e^{\lambda x}$ $x > 0$	$\frac{1}{\lambda}$	$\frac{1}{\lambda^2}$	$t < \lambda$	
$\mathrm{Erlang}(n,\lambda)$	$\frac{\lambda^n x^{n-1} e^{-\lambda x}}{(n-1)!}$ $x > 0$	$1 - \sum_{i=0}^{n-1} \frac{\lambda^n x^n e^{-\lambda x}}{n!}$ $x > 0$	$\frac{n}{\lambda}$	$\frac{n}{\lambda^2}$	$\frac{\frac{1}{(1-\frac{t}{\lambda})^n}}{t < \lambda}$	

## Basic Concepts of Computational and Complexity Theory

- 7.1 Finite Automata and Regular Languages
- 7.2 Pushdown Automata and Context-Free Languages
- 7.3 Turing Machines and the Church-Turing Thesis
- 7.3.1 Turing Machine

7.3.2

#### 7.4 Computational Complexity

#### 7.4.1 Asymptotic Notation

**Definition 7.4.1** (asymptotically positive function).  $f: \mathbb{N} \to \mathbb{R}$  is an asymptotically positive function if  $\exists n_0 > 0$  such that  $\forall n > n_0$  we have f(n) > 0

**Definition 7.4.2** (O-Notation). For a function g(n),

$$O(g(n)) = \{f : \exists c > 0, n_0 > 0 \text{ such that } f(n) \le cg(n), \forall n \ge n_0\}$$

O-Notation also known as asymptotic upper bound.

**Definition 7.4.3** ( $\Omega$ -Notation). For a function g(n),

$$\Omega(g(n)) = \{ f : \exists c > 0, n_0 > 0 \text{ such that } f(n) \ge cg(n), \forall n \ge n_0 \}$$

 $\Omega$ -Notation also known as asymptotic lower bound.

**Definition 7.4.4** ( $\Theta$ -Notation). For a function g(n),

$$\Theta(g(n)) = \{ f : \exists c > 0, n_0 > 0 \text{ such that } c_1 g(n) \le f(n) \le c_2 g(n), \forall n \ge n_0 \}$$

 $\Omega$ -Notation and  $\Theta$ -Notation are not used very often when we talk about running times.

**Definition 7.4.5** (o-Notation). For a function g(n),

$$o(g(n)) = \{f : \exists c > 0, n_0 > 0 \text{ such that } f(n) < cg(n), \forall n \ge n_0 \}$$

**Definition 7.4.6** ( $\omega$ -Notation). For a function g(n),

$$\omega(g(n)) = \{ f : \exists c > 0, n_0 > 0 \text{ such that } f(n) > cg(n), \forall n \ge n_0 \}$$

**Notice:** O(g),  $\Omega(g)$  and  $\Theta(g)$  are sets, we use "=" to represent " $\in$ ". In here "=" is asymmetric. Equality such as  $O(n^3) = n^3 + n$  is incorrect.

#### **Example.** The following are some examples

f(n)	g(n)	O	Ω	Θ
$4n^2 + 3n$	$n^3 - 2n + 3$	Yes	No	No
$\lg^{10} n$	$n^{0.1}$	Yes	No	No
$\log_{10} n$	$\lg(n^3)$	Yes	Yes	Yes
$\lceil \sqrt{10n+100} \rceil$	n	Yes	No	No
$n^3 - 100n$	$10n^2 \lg n$	No	Yes	No
$2^n$	$2^{\frac{n}{2}}$	No	Yes	No
$\sqrt{n}$	$n^{\sin n}$	No	No	No

#### **Theorem 7.1.** Let f and g be two functions that

$$\lim_{n \to \infty} \frac{f(n)}{g(n)} = c > 0$$

Then  $f(n) = \Theta(g(n))$ 

*Proof.* Since  $\lim_{n\to\infty} \frac{f(n)}{g(n)}$  exists and positive, there is some  $n_0$  beyond which the ratio is always between  $\frac{1}{2}c$  and 2c. Thus,

$$\forall n > n_0, f(n) \le 2cg(n) \Rightarrow f(n) = O(g(n)) \tag{7.2}$$

$$\forall n > n_0, f(n) \ge \frac{1}{2}cg(n) \Rightarrow f(n) = \Omega(g(n))$$
(7.3)

A set of properties:

$$f(n) = O(g(n)) \iff g(n) = \Omega(f(n))$$
 (7.4)

$$f(n) = \Theta(g(n)) \iff f(n) = O(g(n)) \text{ and } g(n) = O(f(n))$$
 (7.5)

Another set of properties:

$$f = O(g), g = O(h) \Rightarrow f = O(h) \tag{7.6}$$

$$f = \Omega(g), g = \Omega(h) \Rightarrow f = \Omega(h)$$
 (7.7)

$$f = \Theta(g), g = \Theta(h) \Rightarrow f = \Theta(h)$$
 (7.8)

**Theorem 7.2.** If  $f_i = O(h)$ , for finite number of  $i \in K$ , then  $\sum_{i \in K} f_i = O(h)$ 

Proof is trivia.

**Notice:** Function f and g not necessarily have relation f = O(g) or g = O(f). E.g., for  $f = \sqrt{n}$  and  $g = n^{\sin n}$ ,  $f \notin O(g)$  and  $g \notin O(f)$ 

#### 7.4.2 Common Running Times

The following are examples of common running times.

Running Time	Examples			
O(n)	Scan through a list to find a element matching with input			
$O(\lg n)$	Binary search			
$O(n^2)$	Scan through every pair of elements, $\binom{n}{2}$			
$O(n^3)$	Matrix multiplication by definition			
$O(n \lg n)$	Many divide and conquer algorithm which in each step iteratively			
	divide the problem into two part and solve the subproblem, for			
	example mergesort.			
O(n!)	Enumerate all permutation, for example Hamiltonian Cycle Prob-			
	lem			
$O(c^n)$	Enumerate all elements in power set. $(O(2^n))$			
$O(n^n)$	Enumerate all combinations. (Can't find good example yet)			

Here is a comparison between running times:  $\lg n < \sqrt{n} < n < n \lg n < n^2 < n^{\sqrt{n}} < 2^n < e^n < n! < n^n$ 

#### 7.5 NP and Computational Intractability

#### 7.5.1 P, NP and Co-NP

**Definition 7.5.1** (Decision Problem). A problem X is a **decision problem** if the output is either 0 or 1 (yes/no).

The input of a problem can always be encoded into a binary string. The size of an input is the length of the encoded string s for the input.

**Notice:** For optimization problem X, we can always define a decision version X', by giving a threshold and ask if the objective function can satisfy that threshold or not. If that decision version X' can be solved in polynomial time, we can solve the original problem X in polynomial time.

**Definition 7.5.2** (Polynomial running time). An algorithm A has a **polynomial running time** if there is a polynomial function  $p(\cdot)$  such that  $\forall s$ , the algorithm A terminates on s in at most  $p(|\cdot|)$  steps.

**Definition 7.5.3** (P). The complexity class P is the set of decision problems X that can be solved in polynomial time.

**Example.** Shortest path, minimum spanning tree, determine if an integer is prime number, those are in P.

**Definition 7.5.4** (Certificate, Certifier). In order to check the an algorithm A, for a binary string s, such that  $s \in A$ , there is another separated algorithm B uses s and t, a separated string that contains the evidence that s is a "yes" instance of X, as input string and output (another) "yes". Then this string t is called **certificate**, this separated algorithm B(s,t) is called **certifier**.

**Notice:** Certificate is like a solution, and certifier is the algorithm to prove solution is correct. But both with fancy terminology

**Example.** For Independent set problem. The input (s) is a graph, the certificate (t) will be a set of size k, and the certifier will be an algorithm to check the given set is an independent set.

**Example.** For 3-SAT problem. The input s will be the 3-CNF (conjunctive normal form), the certificate t will be the assignment of true values for each terms in the 3-CNF. The certifier will be the algorithm calculates the true value of the clause given the true values in t.

**Definition 7.5.5** (Efficient Certifier). B is an efficient certifier for problem X if

• B runs in polynomial time with input s and t. (s is the input of origin algorithm and t is the certificate)

• There is a polynomial function  $p(|\cdot|)$  such that for every string s, we have  $s \in X$  (s is a "yes" solution for X) and B(s,t) = yes (the certifier returns "yes")

**Definition 7.5.6** (NP). The complexity class NP is the set of all problems for which there exists an efficient certifier.

From the definition of NP we can immediately know that  $P \subseteq NP$ . Because all the problems in P can satisfy the definition of NP. But is there a problem that  $X \in NP$  and  $X \notin P$ ? Solve that problem and you will win 100 million dollars and the chance to be remember forever.

#### Notice: NP stands for Non-deterministic Polynomial time

To prove if a problem is NP, we need to have a certifier that can output "yes" given the certificate in polynomial time. Otherwise it is not in NP.

**Example.** Given a graph G = (V, E) and an integer t > 0, whether the minimum vertex cover of G has size at least t. This problem is **unlikely** in NP. Reason is, the "yes" instance is "all minimum vertex cover has size at least t", that is not **likely to be** solvable in polynomial time for it need to find all minimum vertex covers.

**Definition 7.5.7** (Co-NP). For a problem X, the problem  $\bar{X}$  is the problem such that  $s \in \bar{X} \iff s \notin X$ . Then **Co-NP** is the set of decision problems X such that  $\bar{X} \in NP$ .

**Example** (Tautology Problem). Given a boolean formula, determine whether the formula is always evaluates to 1. This is a problem in Co-NP. Because we can have a polynomial time certifier to confirm that an instance is not a tautology.

**Example.** Given two boolean formulas, to determine whether or not they are equivalent. This is a problem in Co-NP. Because if we have one instance such that the output is "no", then we can easily prove there are counter examples in origin algorithm.

**Notice:** If the instance is "easy" to prove to be true, then it is in NP, if the counter instance is "easy" to prove to be false, then it is in Co-NP

Relation between P, NP and Co - NP is as following

- $P \subseteq NP$
- $P \subseteq Co NP$
- P = NP? is not known
- P = Co NP? is not known
- NP = Co NP? is not known
- If P = NP then P = Co NP

#### 7.5.2 Polynomial-Time Reductions

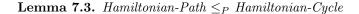
**Definition 7.5.8** (Polynomial-time reducible). Given an algorithm A that solves problem Y, if any instance of problem X can be solved using a polynomial number of standard computational steps, plus a polynomial number of calls to algorithm A, then we say X is **polynomial-time reducible** to Y, denoted as

$$X <_P Y \tag{7.9}$$

In a more intuitive way,  $X \leq_P Y$  means X can't be more difficult than Y. Solving X can be "transforming" X into an equivalent Y in polynomial number of steps and then solve it by calling Y polynomial number of times, usually one time. Thus, if  $X \leq_P Y$ ,

- If Y can be efficiently solved, X can be efficiently solved.
- If X cannot be efficiently solve, Y cannot be efficiently solved.

**Notice:** To prove  $X \leq_P Y$ , usually we already have an algorithm for Y, could be polynomial or not.



Proof. For graph G = (V, E), s and t are two vertices that in V, define a new graph  $G' = (V \cup \{v\}, E \cup \{(u, s)\} \cup \{(t, u)\}$ . To solve the Hamilton-Path from s to t in graph G, say  $p_{st}$ , is equivalent to solving Hamilton-Cycle problem in G', i.e., find  $sp_{st}te_{tu}ue_{us}$ .

#### Lemma 7.4. Hamiltonian- $Cycle \leq_P Hamiltonian$ -Path

*Proof.* For vertex s, make a copy and denote it as s', s' is connected to all the vertices that s connected. Solving the Hamiltonian-Cycle problem is equivalent to solving the Hamiltonian-Path problem from s to s'

#### **Lemma 7.5.** Hamiltonian-Path $\leq_P$ degree-3 spanning tree

*Proof.* In graph G, for vertex s and t, add vertices s', s'', t', t'' and edges (s, s'), (s, s''), (t, t'), (t, t''). For all the other vertex  $u \in V \setminus \{s\} \setminus \{t\}$ , add vertices u' and edge (u, u') to the graph. Then solving Hamiltonian-Path problem is equivalent to solve the degree spanning tree problem in this new G.

#### **Lemma 7.6.** Vertex Cover $\leq_P Set\ Cover$

Proof.  $\Box$ 

#### **Lemma 7.7.** Set $Cover \leq_P Vertex Cover$

Proof.

#### **Lemma 7.8.** Clique $\leq_P$ Independent Set

*Proof.* S is a clique in G = (V, E) iff S is an independent set in  $\overline{G} = (V, \overline{E})$ 

#### **Lemma 7.9.** Independent $Set \leq_P Clique$

*Proof.* S is an independent set in G = (V, E) iff S is a clique in  $\overline{G} = (V, \overline{E})$ 

#### **Lemma 7.10.** Vertex-Cover $\leq_P$ Independent Set

*Proof.* S is a vertex-cover of G = (V, E) iff  $V \setminus S$  is an independent set of G

#### **Lemma 7.11.** 3-Coloring $\leq_P$ 4-Coloring

*Proof.* For a graph G = (V, E), define a new graph  $G' = (V \cup \{u\}, E \cup \{(u, v) | \forall v \in V\})$ . Solving the 3-Coloring problem is equivalent to solving the 4-Coloring problem in G'

**Lemma 7.12.** Independent  $Set \leq_P Set \ Packing$ 

#### 7.5.3 NP-Completeness

**Definition 7.5.9** (NP-Completeness). A problem X is called **NP-Complete** if

- $X \in NP$ , and
- $Y \leq_P X, \forall Y \in NP$

An intuitive explanation will be, we can regard problems that is NP-Complete to be the most difficult problems in NP. If any of those can be solved in polynomial time, then all problems in NP can be solved in polynomial time.

**Theorem 7.13** (Cook's Theorem). 3-SAT is in NP-Complete.

Proof.

Theorem 7.14. 2-SAT is in P

#### Algorithm 1 Clear(S, x)

```
1: Initialize S' \leftarrow S
2: for Each clause c in S' do
       if c is a singleton x then
           Remove singleton clause x in S'
4:
5:
       else if c is a singleton \neg x then
           return null
6:
7:
       else if c is a 2-clause with x then
           Remove singleton clause c in S'
8:
       else if c is a 2-clause with \neg x then
9:
           Remove \neg x in clause c in S'
10:
11: return S'
```

#### **Algorithm 2** 2-SAT(S)

```
1: while Length of S is greater than 0 do
         if S has a singleton x (or \neg x) then
             S \leftarrow \text{Clear}(S, x) \text{ (or } S \leftarrow \text{Clear}(S, \neg x))
 3:
             if S == null then
 4:
                 return False
 5:
 6:
         else
             Denote S = (x_1 \wedge x_2) \vee S_r
 7:
 8:
             S_1 \leftarrow x_1 \vee S_r
             S_2 \leftarrow x_2 \vee S_r
 9:
             if 2\text{-SAT}(S_1) then
10:
                  return True
11:
             else if 2\text{-SAT}(S_2) then
12:
13:
                  return True
14:
             else
                  return False
15:
16: return True
```

*Proof.* The following process Clear(S, x) can be executed within O(n), in which S is a CNF with at least one clause being singleton x (or  $\neg x$ ), where n is the length of CNF S, if it returns S', S' will be a CNF without x (or  $\neg x$ ) in any of clause in S. Otherwise, if it returns null, then S is not satisfiable.

To determine a 2-SAT CNF satisfiability, the following process 2 - SAT(S) can be executed within O(n), where n is the length of CNF S. It returns True if S is satisfiable, False otherwise.

General idea is, if there are singletons in CNF S, if there are conflicts between singletons, the CNF is not satisfiable, otherwise, those singletons can be assigned True values, or False values, one singleton at a time. For the rest of clauses, either it becomes a singleton, or it remains a 2-clause. When we cannot find any singleton in CNF, we generate two sub-CNF, by separating the first 2-clause and keep the remaining of CNF. Then both two sub-CNF will have at least one singleton, we can perform the previous process to those two sub-CNF. If at least one of those sub-CNF is satisfiable, the original CNF is satisfiable, otherwise, if both sub-CNF are not satisfiable, the original CNF is not satisfiable. Therefore 2-SAT CNF Satisfiability is a problem in P.

**Theorem 7.15.** If X is NP-Complete and  $X \in P$ , then P = NP

Proof. Direct result from Cook's theorem.

#### 7.5.4 NP-Complete Problems

# Part II Linear Programming

## The Simplex Method - Basic

#### 8.1 Basic Feasible Solutions and Extreme Points

**Definition 8.1.1** (Basic Feasible Solutions). Consider the system  $\{\mathbf{A_{m \times n}x} = \mathbf{b_m}, \mathbf{b_m} \geq \mathbf{0}\}$ , suppose  $rank(\mathbf{A}, \mathbf{b}) = rank(\mathbf{A}) = m$ , we can rearrange the columns of  $\mathbf{A}$  so that we have a partition of  $\mathbf{A}$ . Let  $\mathbf{A} = \begin{bmatrix} \mathbf{B} & \mathbf{N} \end{bmatrix}$  where  $\mathbf{B}$  is an  $m \times m$  invertible matrix, and  $\mathbf{N}$  is an  $m \times (n - m)$  matrix. The solution  $\mathbf{x} = \begin{bmatrix} \mathbf{x_B} \\ \mathbf{x_N} \end{bmatrix}$  to the equation  $\mathbf{Ax} = \mathbf{b}$ , where  $\mathbf{x_B} = \mathbf{B}^{-1}\mathbf{b}$  and  $\mathbf{x_N} = \mathbf{0}$  is called **basic solution** of system. If  $\mathbf{x_B} \geq \mathbf{0}$ , it is called **basic feasible solution** or  $\mathbf{B}.\mathbf{F}.\mathbf{S}$ . If  $\mathbf{x_B} > \mathbf{0}$  it is called **non-degenerate basic feasible solution**. For  $\mathbf{x_B} \geq \mathbf{0}$ , if some  $x_j = 0$ , those components are called **degenerated basic feasible solution**.  $\mathbf{B}$  is called the **basic matrix**,  $\mathbf{N}$  is called **non-basic matrix** 

**Theorem 8.1.**  $\mathbf{x}$  is an extreme point  $\iff \mathbf{x}$  is a basic feasible solution.

*Proof.* This proof is lack of details. Denote S as feasible region.

FIXME

( $\Rightarrow$ ) First, Let  $\mathbf{x}$  be a B.F.S., Suppose  $\mathbf{x} = \lambda \mathbf{u} + (1 - \lambda)\mathbf{v}$ , for  $\mathbf{u}, \mathbf{v} \in \mathcal{S}, \lambda \in (0, 1)$ . Let  $I = \{i : x_i > 0\}$  be the set of index where the inequality constraint are not tight. Then for  $i \notin I$ ,  $x_i = 0$ , which implies  $u_i = v_i = 0$ .  $\mathbf{u}, \mathbf{v} \in \mathcal{S} \Rightarrow \mathbf{A}\mathbf{u} = \mathbf{A}\mathbf{v} = \mathbf{b} \Rightarrow \mathbf{A}(\mathbf{u} - \mathbf{v}) = \mathbf{0} \Rightarrow \sum_{i=1}^{n} (u_i - v_i)a_i = 0$ .

- if  $i \notin I$  then  $x_i = 0$ , which implies  $u_i = v_i = 0$  -  $\therefore$   $\mathbf{A}\mathbf{u} = \mathbf{A}\mathbf{v} = \mathbf{b}$ ,  $\therefore$   $\mathbf{A}(\mathbf{u} - \mathbf{v}) = \mathbf{0} \Rightarrow \sum_{i=1}^{n} (u_i - v_i)a_i = 0$ ,  $\therefore$   $u_i = v_i = 0$ , for  $i \notin I$ , it implies  $u_i = v_i$  for  $i \in I$ , Hence u = v, x is E.P.

( $\Leftarrow$ ) Second, suppose **x** is not B.F.S., i.e.  $\{a_i : i \in I\}$  are linearly dependent. Then there  $\exists \mathbf{u} \neq \mathbf{0}, u_i = 0, i \notin I$  such that  $\mathbf{A}\mathbf{u} = \mathbf{0}$ . Hence, for a small  $\epsilon$ ,  $\mathbf{x} = \frac{1}{2}(\mathbf{x} + \epsilon \mathbf{u}) + \frac{1}{2}(\mathbf{x} - \epsilon \mathbf{u})$ , **x** is not E.P.

#### 8.2 The Simplex Method

#### 8.2.1 Key to Simplex Method

#### Cost Coefficient

The cost coefficient can be derived from the following

$$z = cx \tag{8.1}$$

$$=c_B x_B + c_N x_N \tag{8.2}$$

$$= c_B(B^{-1}b - B^{-1}Nx_N) + c_N x_N (8.3)$$

$$= c_B B^{-1} b - \sum_{j \in N} (c_B B^{-1} a_j - c_j) x_j$$
(8.4)

$$= c_B B^{-1} b - \sum_{j \in N} (z_j - c_j) x_j$$
 (8.5)

We denote  $z_0=c_BB^{-1}b,\,z_j=c_B^{-1}a_j,\,\bar{b}=B^{-1}b$  and  $y_j=B^{-1}a_j$  for all nonbasic variables. The formulation can be transformed into

$$\min \quad z = z_0 - \sum_{j \in N} (z_j - c_j) x_j \tag{8.6}$$

s.t. 
$$\sum_{j \in N} y_j x_j + x_B = \bar{b}$$
 (8.7)

$$x_j \ge 0, j \in N \tag{8.8}$$

$$x_B \ge 0 \tag{8.9}$$

In the above formulation,  $z_j - c_j$  is the cost coefficient. If  $\exists j$  and  $z_j - c_j > 0$ , it means the objective function can still be optimized. (If  $\forall j, z_j - c_j \leq 0$ , then  $z \geq z_0$  for any feasible solution, z is the optimal solution)

#### Pivot

After finding the most violated  $z_j - c_j$ , we find a variable, say  $x_k$ , where  $z_k - c_k = \min\{z_j - c_j\}$  to be the variable leaving the basis.

If there are degenerated variables, we can perform different method to choose variable to enter basis.

#### Minimum Ratio

$$x_{B_i} = \bar{b}_i - y_{ik} x_k \ge 0 \tag{8.10}$$

Therefore we have the minimum ratio rule

$$x_k = \min_{i \in B} \{ \frac{\bar{b}_i}{y_{ik}}, y_{ik} > 0 \}$$
 (8.11)

If for the that column all  $y_{ik} \leq 0$ , unbounded.

#### 8.2.2 Simplex Method Algorithm

The pseudo-code of Simplex Method is given as following:

#### 8.3 Tableau Method for Simplex Method

The following is an example of using tableau to solve simplex method. Initial tableau:

		z	$x_1$	$x_2$	$x_3$	$x_4$	$x_5$	RHS
	z	1	1	3	0	0	0	0
Ì	$x_3$	0	1	-2	1	0	0	0
	$x_4$	0	-2	1	0	1	0	4
ĺ	$x_5$	0	5	3	0	0	1	15

Last tableau:

- The optimal basic variables are  $x_3$ ,  $x_2$ ,  $x_1$ . The optimal basis is the columns in the initial tableau with correspond columns

$$B = \begin{pmatrix} \frac{13}{11} & \frac{3}{11} & \frac{97}{11} \\ \frac{5}{11} & \frac{2}{11} & \frac{50}{11} \\ -\frac{3}{11} & \frac{1}{11} & \frac{3}{11} \end{pmatrix}$$
(8.14)

8.4. INITIAL BASIS 51

#### Algorithm 3 Simplex Method

```
Require: Given a basic feasible solution with basis B
Ensure: Optimal objective value min z = cx
 1: Set B for basic variables, N for nonbasic variables
 2: \mathbf{B} \leftarrow \text{all slack variables}
 3: N \leftarrow all variables excepts slack variables
 4: for \forall j do
            z_i = c_B B^{-1} a_i = 0
 6: while \exists z_j - c_j > 0 do
            z_j = wa_j - c_j = c_B B^{-1} a_j - c_j
           z_k - c_k = \max_{j \in \mathbf{N}} \{z_j - c_j\}y_k = B^{-1} a_k
 9:
            if \exists y_{ik} > 0 then
10:
                \theta_r = \min_{i \in \mathbf{B}} \{\theta_i = \frac{\bar{b}_i}{y_{ik}} : y_{ik} > 0\}
\mathbf{B} \leftarrow \mathbf{B} \setminus \{k\}
11:
12:
                 \mathbf{N} \leftarrow \mathbf{N} \cup \{k\}
13:
                 \mathbf{B} \leftarrow \mathbf{B} \cup \{r\}
14:
                 \mathbf{N} \leftarrow \mathbf{N} \setminus \{r\}
15:
16:
            else
                  Unbounded
17:
18: x_B^* = B^{-1}b = \bar{b}
19: x_N = 0
20: z^* = c_B B^{-1} b = c_B \bar{b} \mathbf{a_{B_k}}
```

- From the initial tableau, we can see the initial basis is built from slack variables  $x_3$ ,  $x_4$ ,  $x_5$ . The  $B^{-1}$  is the correspond columns in final tableau.

$$B = \begin{pmatrix} 1 & -2 & 1\\ 0 & 1 & -2\\ 0 & 3 & 5 \end{pmatrix} \tag{8.15}$$

- The optimal basic variables are  $x_3, x_2, x_1$ . Find  $c_B$  in the initial tableau.

$$c_B = \begin{pmatrix} 0\\3\\1 \end{pmatrix} \tag{8.16}$$

- Find  $w = c_B B^{-1}$  from the final tableau, correspond to the slack variable.

$$w = c_B B^{-1} = \begin{pmatrix} 0 \\ -\frac{12}{11} \\ -\frac{7}{11} \end{pmatrix}$$
 (8.17)

#### 8.4 Initial Basis

If some of the constraint is not in  $\sum_{i=1}^{n} a_i x_i \leq 0$  form, we cannot add a positive slack variable. In this case, we add an artificial variable other than slack variable.

$$\sum_{i=1}^{n} a_i x_i \ge (or =)0 \Rightarrow \sum_{j=1}^{n} a_i x_i + x_a = 0$$
(8.18)

Notice that in an optimal solution,  $x_a = 0$ , otherwise it is not valid. Artificial variables are only a tool to get the simplex method started.

#### 8.4.1 Two-Phase Method

#### Two-Phase Method

#### For **Phase I:**

Solve the following program start with a basic feasible solution  $x = 0, x_a = b$ , i.e., the artificial variable forms the basis.

$$\min \quad \mathbf{1}x_a \tag{8.19}$$

$$s.t. \quad Ax + x_a = b \tag{8.20}$$

$$x \ge 0 \tag{8.21}$$

$$x_a \ge 0 \tag{8.22}$$

If the optimal  $1x_a \neq 0$ , infeasible, stop. Otherwise proceed Phase II. For **Phase II**:

Remove the columns of artificial variables, replace the objective function with the original objective function, proceed to solve simplex method.

#### Discussion

Case A:  $x_a \neq 0$ 

Infeasible.

Case B.1:  $x_a = 0$  and all artificial variables are out of the basis

At the end of Phase I, we derive

$x_0$	$x_B$	$x_N$	$x_a$	RHS	
1	0	0	-1	0	(8.23)
0	I	$B^{-1}N$	$B^{-1}$	$B^{-1}b$	

We can discard  $x_a$  columns, (or we can leave it because it keeps track of  $B^{-1}$ ), and then we do the Phase II

z	ž	$x_B$	$x_N$	RHS
1		0	$c_B B^{-1} N - c_N$	$c_B B^{-1} b$
0	)	I	$B^{-1}N$	$B^{-1}b$

Case B.2: Some artificial variables are in the basis at zero values

This is because of degeneracy. We pivot on those artificial variables, once they leave the basis, eliminate them.

#### 8.4.2 Big M Method

#### 8.4.3 Single Artificial Variable

#### 8.5 Degeneracy and Cycling

#### 8.5.1 Degeneracy

#### Degeneracy in Simplex Method

If the basic variable  $x_B$  is not strictly > 0, i.e. if some basic variable equals to 0, we call it degenerate.

#### Degeneracy for Bounded Variables

If some basic variables are at their upper bound or lower bound, we call it degenerate.

#### 8.5.2 Cycling

In the degenerate case, pivoting by the simplex rule does not always give a strict decrease in the objective function value, because it may have  $b_r = 0$ . It is possible that the tableau may repeat if we use the simplex rule.

Geometrically speaking, it means that at the same point - extreme point - it corresponds to more than one feasible solutions, so when we are pivoting, we stays at the same place.

In computer algorithm, we rarely care about cycling because the data in computer is not precise, it is very hard to get into cycling.

#### 8.5.3 Cycling Prevent

#### Lexicographic Rule

- For entering variable, same as simplex rule
- For leaving variable, if there is a tie, choose the variable with the smallest  $\frac{y_{r1}}{y_{rk}}$ .

#### Bland's Rule

- For entering variable, choose the variable with smallest index where  $z_j c_j \leq 0$
- For leaving variable, if there is a tie, choose the variable with smallest index.

#### Successive Ratio Rule

- Select the pivot column as any column k where  $z_k-c_k\leq 0$
- Given k, select the pivot row r as the minimum successive ratio row associated with column k.

In other words, for pivot columns where there is no tie in the usual minimum ratio, the successive ratio rule reduces to the simplex rule

#### 8.6 As a Search Algorithm

#### 8.6.1 Improving Search Algorithm

A simplex method is a search algorithm, for each iteration it finds a not-worse solution, which can be represented as:

$$x^{t} = x^{t-1} + \lambda_{t-1}d^{t-1} \tag{8.25}$$

Where

- $x^t$  is the solution of the tth iteration
- $\lambda_t$  is the step length of tth iteration
- $d^t$  is the direction of the tth iteration

For each iteration, it contains three steps:

- Optimality test
- Find direction
- Find the step length

#### 8.6.2 Optimality Test

$$z = cx (8.26)$$

$$= \begin{bmatrix} c_B & c_N \end{bmatrix} \begin{bmatrix} x_B \\ x_N \end{bmatrix} \tag{8.27}$$

$$=c_B x_B + c_N x_N \tag{8.28}$$

and 
$$:Ax = b$$
 (8.29)

$$\therefore Bx_B + Nx_N = b, x_B \ge 0, x_N \ge 0 \tag{8.30}$$

$$\therefore x_B = B^{-1}b - B^{-1}Nx_N \tag{8.31}$$

$$z = c_B B^{-1} b - c_B B^{-1} N x_N + c_N x_N (8.32)$$

for current solution  $\hat{x} = \begin{bmatrix} \hat{x_B} \\ 0 \end{bmatrix}$ ,  $\hat{z} = c_B B^{-1} b$ , then

$$z - \hat{z} = \begin{bmatrix} 0 & c_N - c_B B^{-1} N \end{bmatrix} \begin{bmatrix} x_B \\ x_N \end{bmatrix}$$
 (8.33)

The  $c_N - c_B B^{-1} N$  is the reduced cost, for a minimized problem, if  $c_N - c_B B^{-1} N > 0$  means  $z - \hat{z} \ge 0$ , it reaches the optimality because we cannot find a solution less than  $\hat{z}$ .

#### 8.6.3 Find Direction

Suppose we choose  $x_k$  as a candidate to pivot into Basis

$$x = \begin{bmatrix} B^{-1}b - B^{-1}a_k x_k \\ 0 + e_k x_k \end{bmatrix} = \begin{bmatrix} B^{-1}b \\ 0 \end{bmatrix} + \begin{bmatrix} -B^{-1}a_k \\ e_k \end{bmatrix} x_k$$
 (8.34)

In this form, we can see: x is the result after tth iteration,  $\begin{bmatrix} B^{-1}b\\0 \end{bmatrix}$  is the result after (t-1)th iteration.  $\begin{bmatrix} -B^{-1}a_k\\e_k \end{bmatrix}$  is the iteration direction,  $x_k$  is the step length.

The only requirement of  $x_k$  is  $r_k < 0$  where  $r_k = c_k - z_k$  is reduce cost, which is the kth entry of  $c_N - c_B B^{-1} N$ . Generally speaking, we usually take  $r_k = \min\{c_j - z_j\}$  (which in fact can not guarantee the efficient of the algorithm.)

#### 8.6.4 Find the Step Length

We need to guarantee the non-negativity, so for each iteration, we need to make sure  $x \geq 0$ . Which means

$$B^{-1}b - B^{-1}a_k x_k \ge 0 (8.35)$$

Denote  $B^{-1}b$  as  $\bar{b}$ , denote  $B^{-1}a_k$  as  $y_k$ 

If  $y_k \leq 0$ , we can have  $x_k$  as large as infinite, which means unboundedness.

If  $y_k > 0$  now we can use the minimum ratio to guarantee non-negativity.

Remember hit the bound, basic variable leave the basis and become non-basic variable.

## The Simplex Method - Improved

#### 9.1 Revised Simplex Method

#### 9.1.1 Key to Revised Simplex Method

The procedure of Simplex Method is (almost) exactly the same as original simplex method. However, notice that we don't need to use N so for the revised simplex method, we don't calculate any matrix related to N The original matrix:

The revised matrix:

For each pivot iteration, calculate  $z_j - c_j = wa_j - c_j = c_B B^{-1} a_j - c_j$ ,  $\forall j \in N$ , pivot rules are the same as simplex method, each time find a variable  $x_k$  to enter basis

$$\begin{array}{c|cc}
B^{-1} & RHS \\
\hline
w & c_B \overline{b} \\
\hline
B^{-1} & \overline{b}
\end{array}
\qquad
\begin{array}{c|cc}
x_k \\
\hline
z_k - c_k \\
\hline
y_k
\end{array}$$
(9.3)

Do the minimum ratio rule to find the variable  $x_r$  to leave the basis

$$\begin{array}{c|cccc}
B^{-1} & RHS \\
\hline
w & c_B b \\
\hline
b_1 \\
\bar{b}_2 \\
B^{-1} & \bar{b}_2 \\
\vdots \\
B^{-1} & \bar{b}_r \\
\vdots \\
\bar{b}_m
\end{array}$$

$$\begin{array}{c|cccc}
x_k \\
z_k - c_k \\
\hline
y_{1k} \\
y_{2k} \\
\vdots \\
y_{rk} (\text{pivot at here}) \\
\vdots \\
y_{mk} \\
\end{array}$$

$$\begin{array}{c|cccc}
(9.4)
\end{array}$$

#### 9.1.2 Comparison between Simplex and Revised Simplex

#### Advantage of Revised Simplex

- Save storage memory
- Don't need to calculate N (including  $B^{-1}N$  and  $c_BB^{-1}N$ )
- More accurate because round up errors will not be accumulated

#### Disadvantage of Revised Simplex

- Need to calculate  $wa_i$  for all  $i \in N$  (in fact don't need to calculated it for the variable just left the basis)

#### **Computation Complexity**

Method	Type	Operations	
Simplex	×	(m+1)(n-m+1)	
	+	m(n-m+1)	(1
Revised Simplex	×	$(m+1)^2 + m(n-m)$	
	+	m(m+1) + m(n-m)	

#### When to use?

- When m >> n, do revise simplex method on the dual problem
- When  $m \simeq n$ , revise simplex method is not as good as simplex method
- When  $m \ll n$  perfect for revise simplex method.

#### 9.1.3 Decomposition of B inverse

Let  $B = \{a_{B_1}, a_{B_2}, ..., a_{B_r}, ..., a_{B_m}\}$  and  $B^{-1}$  is known. If  $a_{B_r}$  is replaced by  $a_{B_k}$ , then B becomes  $\bar{B}$ . Which means  $a_{B_r}$  enters the basis and  $a_{B_k}$  leaves the basis.

Then  $\bar{B}^{-1}$  can be represent by  $B^{-1}$ . Noting that  $a_k = By_k$  and  $a_{B_i} = Be_i$ , then

$$\bar{B} = (a_{B_1}, a_{B_2}, ..., a_{B_{r-1}}, a_k, a_{B_{r+1}}, a_m) \tag{9.6}$$

$$= (Be_1, Be_2, ..., Be_{r-1}, By_k, Be_{r+1}, ..., Be_m)$$

$$(9.7)$$

$$=BT\tag{9.8}$$

where T is

$$T = \begin{bmatrix} 1 & 0 & \dots & 0 & y_{1k} & 0 & \dots & 0 \\ 0 & 1 & \dots & 0 & y_{2k} & 0 & \dots & 0 \\ \vdots & \vdots & & \vdots & \vdots & & \vdots & \vdots \\ 0 & 0 & \dots & 1 & y_{r-1,k} & 0 & \dots & 0 \\ 0 & 0 & \dots & 0 & y_{rk} & 0 & \dots & 0 \\ 0 & 0 & \dots & 0 & y_{r+1,k} & 1 & \dots & 0 \\ \vdots & \vdots & & \vdots & \vdots & & \vdots \\ 0 & 0 & \dots & 0 & y_{mk} & 0 & \dots & 1 \end{bmatrix}$$
(9.9)

and

$$E = T^{-1} = \begin{bmatrix} 1 & 0 & \dots & 0 & \frac{-y_{1k}}{y_{rk}} & 0 & \dots & 0 \\ 0 & 1 & \dots & 0 & \frac{-y_{2k}}{y_{rk}} & 0 & \dots & 0 \\ \vdots & \vdots & & \vdots & \vdots & & \vdots \\ 0 & 0 & \dots & 1 & \frac{-y_{r-1,k}}{y_{rk}} & 0 & \dots & 0 \\ 0 & 0 & \dots & 0 & \frac{1}{y_{rk}} & 0 & \dots & 0 \\ 0 & 0 & \dots & 0 & \frac{-y_{r+1,k}}{y_{rk}} & 1 & \dots & 0 \\ \vdots & \vdots & & \vdots & \vdots & & \vdots \\ 0 & 0 & \dots & 0 & \frac{-y_{mk}}{y_{rk}} & 0 & \dots & 1 \end{bmatrix}$$

$$(9.10)$$

For each iteration, i.e. one variable enters the basis and one leaves the basis,  $\bar{B}^{-1} = T^{-1}B^{-1} = EB^{-1}$ . Given that the first iteration starts from slack variables, the first basis  $B_1$  is I, then we have

$$B_t^{-1} = E_{t-1}E_{t-2}\cdots E_2E_1I \tag{9.11}$$

Using E in calculation can simplify the product of matrix where

$$cE = c_1, c_2, ..., c_m \begin{bmatrix} 1 & 0 & ... & g_1 & ... & 0 \\ 0 & 1 & ... & g_2 & ... & 0 \\ \vdots & \vdots & & \vdots & & \vdots \\ 0 & 0 & ... & g_m & ... & 1 \end{bmatrix}$$

$$(9.12)$$

$$= (c_1, c_2, ..., c_{r-1}, cg, c_{r+1}, ..., c_m)$$

$$(9.13)$$

and

$$Ea = \begin{bmatrix} 1 & 0 & \dots & g_1 & \dots & 0 \\ 0 & 1 & \dots & g_2 & \dots & 0 \\ \vdots & \vdots & & \vdots & & \vdots \\ 0 & 0 & \dots & g_m & \dots & 1 \end{bmatrix} \begin{bmatrix} a_1 \\ a_2 \\ \vdots \\ a_m \end{bmatrix}$$
(9.14)

$$\begin{bmatrix}
a_1 \\
a_2 \\
\vdots \\
a_{r-1} \\
0 \\
a_{r+1} \\
\vdots \\
a_m
\end{bmatrix} + a_r \begin{bmatrix}
g_1 \\
g_2 \\
\vdots \\
g_{r-1} \\
g_r \\
g_r \\
g_{r+1} \\
\vdots \\
g_m
\end{bmatrix}$$
(9.15)

$$= \bar{a} + a_r g \tag{9.16}$$

Then we can calculate  $w, y_k$  and  $\bar{b}$ 

$$w = c_B B^{-1} = c_B E_{t-1} E_{t-2} \dots E_2 E_1 (9.17)$$

$$y_k = B^{-1}a_k = E_{t-1}E_{t-2}...E_2E_1a_k (9.18)$$

$$\bar{b} = B_{t+1}^{-1}b = E_t E_{t-1} E_{t-2} \dots E_2 E_1 b \tag{9.19}$$

#### 9.2 Dual Simplex Method

Maintain dual feasibility, i.e. primal optimality, and complementary slackness and work towards primal feasibility. **Tip:** The RHS become new  $z_j - c_j$ , the old  $z_j - c_j$  become new RHS. We are actually solving the dual problem.

#### 9.3 Simplex with Equations

#### 9.4 Simplex with Bounded Variables

#### 9.4.1 Bounded Variable Formulation

$$\min \quad cx \tag{9.20}$$

$$s.t. \quad Ax = b \tag{9.21}$$

$$l \le x \le b \tag{9.22}$$

Reason why we don't the following formulation

$$\min \quad cx \tag{9.23}$$

$$s.t. \quad Ax = b \tag{9.24}$$

$$x - Ix_l = l (9.25)$$

$$x + Ix_u = u (9.26)$$

$$x \ge 0 \tag{9.27}$$

$$x_l \ge 0 \tag{9.28}$$

$$x_u \ge 0 \tag{9.29}$$

is that this formulation increase the number of variable from n to 3n, and the number of constraint from m to m+2n, the size in increase significantly.

#### 9.4.2 Basic Feasible Solution

Consider the system Ax = b and  $l \le x \le b$ , where A is a  $m \times n$  matrix of rank m, the solution  $\bar{x}$  is a **basic feasible solution** if A can be partition into  $[B, N_l, N_u]$  where the solution x can be partition into  $x = (x_B, x_{N_l}, x_{N_u})$ , in which  $\bar{x}_{N_l} = l_{N_l}$  and  $\bar{x}_{N_u} = u_{N_u}$ , therefore

$$\bar{x}_B = B^{-1}b - B^{-1}N_l x_{N_l} - B^{-1}N_u x_{N_u} \tag{9.30}$$

Furthermore, similar to definition of nonnegative variables, if  $l_B \leq x_B \leq u_B$ ,  $x_B$  is a basic feasible solution, if  $l_B < x_B < u_B$ ,  $x_B$  is a non-degenerate basic feasible solution.

#### 9.4.3 Improving Basic Feasible Solution

The basic variables and the objective function can be derived as following:

$$x_B = B^{-1}b - B^{-1}N_l x_{N_l} - B^{-1}N_u x_{N_u} (9.31)$$

$$z = c_B x_B + c_{N_l} x_{N_l} + c_{N_u} x_{N_u} (9.32)$$

$$= c_B(B^{-1}b - B^{-1}N_l x_{N_l} - B^{-1}N_u x_{N_u}) (9.33)$$

$$+ c_B x_B + c_{N_l} x_{N_l} + c_{N_u} x_{N_u} \tag{9.34}$$

$$= c_B B^{-1} b + (c_{N_l} - c_B B^{-1} N_l) x_{N_l} (9.35)$$

$$+ (c_{N_u} - c_B B^{-1} N_u) x_{N_u} (9.36)$$

$$= c_B B^{-1} b - \sum_{j \in J_1} (z_j - c_j) x_j - \sum_{j \in J_2} (z_j - c_j) x_j$$
(9.37)

 $J_1$  is the set of variables at lower bound,  $J_2$  is the set of the variables at upper bound.

Notice that the right-hand-side no longer provide  $c_B B^{-1} b$  and  $B^{-1} b$ . For the variable entering the basis, find the variable with

$$\max\{\max_{j \in J_1} \{z_j - c_j\}, \max_{j \in J_2} \{c_j - z_j\}\}$$
(9.38)

to enter the basis

Tip: "Most violated rule"

The minimum ratio rule is revised for bounded simplex

$$\Delta = \min\{\gamma_1, \gamma_2, u_k - l_k\} \tag{9.39}$$

$$\gamma_{1} = \begin{cases} \min_{r \in J_{1}} \{ \frac{\bar{b}_{r} - l_{B_{r}}}{y_{rk}} : y_{rk} > 0 \} \\ \min_{r \in J_{2}} \{ \frac{\bar{b}_{r} - l_{B_{r}}}{-y_{rk}} : y_{rk} < 0 \} \\ \infty \end{cases}$$

$$(9.40)$$

$$\gamma_2 = \begin{cases} \min_{r \in J_1} \left\{ \frac{u_{B_r} - \bar{b}_r}{-y_{rk}} : y_{rk} < 0 \right\} \\ \min_{r \in J_2} \left\{ \frac{u_{B_r} - b_r}{y_{rk}} : y_{rk} > 0 \right\} \\ \infty \end{cases}$$

$$(9.41)$$

Tip:

Use  $l \le x + \Delta \le u$  to test the range of  $\delta$ , if it hits lower bound, it is called  $\gamma_1$ , if it hits upper bound, it is called  $\gamma_2$ .

## 9.5 Simplex with Unrestricted Variables

## **Duality and Sensitivity**

#### 10.1 Duality

#### 10.1.1 Dual Formulation

For any prime problem

$$\min \quad cx \tag{10.1}$$

s.t. 
$$Ax \ge b$$
 (10.2)

$$x \ge 0 \tag{10.3}$$

we can have a dual problem

$$\max \quad wb \tag{10.4}$$

s.t. 
$$wA \le c$$
 (10.5)

$$w \ge 0 \tag{10.6}$$

#### 10.1.2 Mixed Forms of Duality

For the following prime problem

P(or D) min 
$$c_1x_1 + c_2x_2 + c_3x_3$$
 (10.7)

s.t. 
$$A_{11}x_1 + A_{12}x_2 + A_{13}x_3 \ge b_1$$
 (10.8)

$$A_{21}x_1 + A_{22}x_2 + A_{23}x_3 \le b_2 \tag{10.9}$$

$$A_{31}x_1 + A_{32}x_2 + A_{33}x_3 = b_3 (10.10)$$

$$x_1 \ge 0 \tag{10.11}$$

$$x_2 \le 0 \tag{10.12}$$

$$x_3$$
 unrestricted (10.13)

The dual of the problem

$$D(\text{or P}) \quad \max \quad w_1 b_1 + w_2 b_2 + w_3 b_3 \tag{10.14}$$

s.t. 
$$w_1 A_{11} + w_2 A_{21} + w_3 A_{31} \le c_1$$
 (10.15)

$$w_1 A_{12} + w_2 A_{22} + w_3 A_{32} \ge c_2 (10.16)$$

$$w_1 A_{13} + w_2 A_{23} + w_3 A_{33} = c_3 (10.17)$$

$$w_1 \ge 0 \tag{10.18}$$

$$w_2 \le 0 \tag{10.19}$$

$$w_3$$
 unrestricted (10.20)

In sum,	the relation	between	primal	and	dual	problems	are	listed	as fo	llowing

	Minimization		Maximization	
	$\geq 0$	$\longleftrightarrow$	$\leq 0$	
Var	$\leq 0$	$\longleftrightarrow$	$\geq 0$	Cons
	Unrestricted	$\longleftrightarrow$	=	
	$\geq 0$	$\longleftrightarrow$	$\geq 0$	
Cons	$\leq 0$	$\longleftrightarrow$	$\leq 0$	Var
	=	$\longleftrightarrow$	Unrestricted	

#### 10.1.3 Dual of the Dual is the Primal

For a primal problem (P)

$$(P) \quad \min \quad cx \tag{10.21}$$

s.t. 
$$Ax \ge b$$
 (10.22)

$$x \ge 0 \tag{10.23}$$

The dual problem (D) is

$$(D) \quad \max \quad wb \tag{10.24}$$

s.t. 
$$wA \le c$$
 (10.25)

$$w \ge 0 \tag{10.26}$$

Rewrite the dual

$$\min \quad -b^{\mathsf{T}}w^{\mathsf{T}} \tag{10.27}$$

s.t. 
$$-A^{\mathsf{T}}w^{\mathsf{T}} \ge -c^{\mathsf{T}}$$
 (10.28)

$$w^{\top} \ge 0 \tag{10.29}$$

Find the dual of this problem

$$\max \quad x^{\top}(-c^{\top}) \tag{10.30}$$

s.t. 
$$x^{\top}(-A^{\top}) \le (-b^{\top})$$
 (10.31)

$$x^{\top} \ge 0 \tag{10.32}$$

(10.33)

Rewrite the dual of the dual

$$(P) \quad \min \quad cx \tag{10.34}$$

s.t. 
$$Ax \ge b$$
 (10.35)

$$x \ge 0 \tag{10.36}$$

#### 10.1.4 Primal-Dual Relationships

#### Weak Duality Property

Let  $x_0$  be any feasible solution of a primal minimization problem,

$$Ax_0 \ge b, \quad x_0 \ge 0 \tag{10.37}$$

Let  $x_0$  be any feasible solution of a dual maximization problem,

$$w_0 A \le c, \quad w_0 \ge 0 \tag{10.38}$$

Therefore, we have

$$cx_0 \ge w_0 A x_0 \ge w_0 b \tag{10.39}$$

which is called the weak duality property. This property is for any feasible solution in the primal and dual problem. Therefore, any feasible solution in the maximization problem gives the lower bound of its dual problem, which is a minimization problem, vice versa. We use this to give the bounds in using linear relaxation to solve IP problem.

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#### **Fundamental Theorem of Duality**

With regard to the primal and dual LP problems, one and only one of the following can be true.

- Both primal and dual has optimal solution  $x^*$  and  $w^*$ , where  $cx^* = w^*b$
- One problem has an unbounded optimal objective value, the other problem must be infeasible
- Both problems are infeasible.

#### **Strong Duality Property**

From KKT condition, we know that in order to make  $x^*$  the optimal solution, the following condition should be met.

- Primal Optimal:  $Ax^* \ge b, x^* \ge 0$
- Dual Optimal:  $w^*A \le c, w^* \ge 0$
- Complementary Slackness

$$\begin{cases} w^*(Ax^* - b) = 0\\ (c - w^*A)x^* = 0 \end{cases}$$
 (10.40)

The first condition means the primal has an optimal solution, the second condition means the dual has an optimal solution. The third condition means  $cx^* = w^*b$ , which is also called **strong duality property** 

**Notice:** w in the dual problem is the same as the  $w = c_B B^{-1}$  in primal problem.

#### Complementary Slackness Theorem

Let  $x^*$  and  $w^*$  be any feasible solutions, they are optimal iff

$$(c_j - \mathbf{w}^* \mathbf{a_j}) x_j^* = 0, \quad j = 1, ..., n$$
 (10.41)

$$w_i^*(\mathbf{a}^i \mathbf{x}^* - b_i) = 0, \quad i = 1, ..., m$$
 (10.42)

In particular

$$x_i^* > 0 \Rightarrow \mathbf{w}^* \mathbf{a}_j = c_j \tag{10.43}$$

$$\boldsymbol{w}^* \boldsymbol{a_j} < c_j \Rightarrow x_j^* = 0 \tag{10.44}$$

$$w_i^* > 0 \Rightarrow a^i x^* = b_i \tag{10.45}$$

$$\boldsymbol{a}^{i}\boldsymbol{x}^{*} > b_{i} \Rightarrow w_{i}^{*} = 0 \tag{10.46}$$

It means, if in optimal solution a variable is positive (has to be in the basis), the correspond constraint in the other problem is tight. If the constraint in one problem is not tight, the correspond variable in the other problem is zero. In the dual problem, we solved some w which is positive, we can know that the correspond constraint in primal is tight, furthermore we can solve the basic variables from those tight constraints, which becomes equality and we can solve it using Gaussian-Elimination.

#### 10.1.5 Shadow Price

#### Shadow Price under Non-degeneracy

Let B be an optimal basis for primal problem and the optimal solution  $x^*$  is non-degenerated.

$$z = c_B B^{-1} b - \sum_{j \in N} (z_j - c_j) x_j = w^* b - \sum_{j \in N} (z_j - c_j) x_j$$
(10.47)

therefore

$$\frac{\partial z^*}{\partial b_i} = c_B B_i^{-1} = w_i^* \tag{10.48}$$

 $w^*$  is the shadow prices for the right-hand-side vectors. We can also regard it as the **incremental cost** of producing one more unit of the *i*th product. Or  $w^*$  is the **fair price** we would pay to have an extra unit of the *i*th product.

#### Shadow Price under Degeneracy

For shadow price under degeneracy, the  $w^*$  may not be the true shadow price, for it may not be the right basis. In this case, the partial differentiation may not be valid, for component  $b_i$ , if  $x_i = 0$  and  $x_i$  is a basic variable, we can't find the differentiation.

#### 10.2 Sensitivity

- Change in the Cost Vector
  - Nonbasic Variable:  $c_B$  is not affected,  $z_j = c_B B^{-1} a_j$  is not changed, say nonbasic variable cost coefficient  $c_k$  changed into  $c'_k$ . For now  $z_k c_k \le 0$ , if  $z_k c'_k$  is positive,  $x_k$  must into the basis, the optimal value changed. Otherwise stays at the same.
  - Basic Variable: If  $c_{B_t}$  is replaced by  $c'_{B_t}$ , then  $z'_j c_j$  is

$$z'_{j} - c_{j} = c'_{B}B^{-1}a_{j} - c_{j} = (z_{j} - c_{j}) - (c'_{B_{t}} - c_{B_{t}})B^{-1}a_{B_{t}}$$
(10.49)

for j=k, it is a basic variable, therefore original  $z_k-c_k=0$ ,  $B^{-1}a_k=1$ . Hence  $z_k'-c_k=c_k'-c_k\Rightarrow z_k'-c_k'=0$ . The basis stays the same. The optimal solution updated as  $c_B'B^{-1}b=c_BB^{-1}b+(c_{B_t}'-c_{B_t})B^{-1}b_{B_t}$ .

- Change in the Right-Hand-Side: If b is replaced by b', then  $B^{-1}b$  is replaced by  $B^{-1}b'$ . If  $B^{-1}b' \geq 0$ , the basis remains optimal. Otherwise, we perform dual simplex method to continue.
- Change in the Matrix
  - Changes in Activity(Variable) Vectors for Nonbasic Columns: If a nonbasic column  $a_j$  is replaced by  $a'_j$ , then  $z_j = c_B B^{-1} a_j$  is replaced by  $z'_j = c_B B^{-1} a'_j$ , if new  $z'_j c_j \le 0$ , the basis stays optimal basis, the optimal value is the same because  $c_B$  stays the same.
  - Changes in Activity(Variable) Vectors for Basic Columns: If a basic columns changed, it means B and  $B^{-1}$  changed, and every column changed. We can do this in two steps:
    - \* step I: add a new column with  $a'_i$
    - \* step II: remove the original column  $a_j$

If in step I the new variable can enter basis, i.e.  $z'_j - c_j \leq 0$ , let it enter the basis and eliminate the original column directly (because at this time the original column leave the basis the nonbasic variable is 0); otherwise, if the new column can not form a new basis, treat  $x_j$ , the original variable as an artificial variable.

- Add a New Activity(Variable): Suppose we add a new variable  $x_{n+1}$  and  $c_{n+1}$  and  $a_{n+1}$  respectively. Calculate  $z_{n+1} c_{n+1}$  to determine if the new variable enters the basis, if not, remains the same optimal solution, otherwise, continue on to find a new optimal solution.
- Add a New Constraint: This is the basic of Branch-and-Cut/Bound, also, we can perform dual simplex method after we add a new constraint(cut).

## Column Generation

## Ellipsoid Methods

## Karmarkar's Algorithm

# Part III Integer Programming

# Branch and Bound

14.1 LP Relaxation Based Branch and Bound

# **Cutting Plane Methods**

- 15.1 Gomory Cuts
- 15.2 Kianfar's methods
- 15.3 Lift-and-project

# Lagrangian Relaxation

## 16.1 Basic Constructions

Consider an integer program

$$z_{IP} = \max$$
 **cx** s.t.  $\mathbf{A}\mathbf{x} \leq \mathbf{b}$   $\mathbf{x} \in \mathbb{Z}_+^n$ 

We can partial  $\bf A$  and  $\bf b$  into two sets of constraints, where  $\bf A = \begin{bmatrix} \bf A^1 \\ \bf A^2 \end{bmatrix}$  and  $\bf b = \begin{bmatrix} \bf b^1 \\ \bf b^2 \end{bmatrix}$ . So the IP model can be rewritten as

$$z_{IP} = \max$$
 **cx**  
s.t.  $\mathbf{A}^1 \mathbf{x} \leq \mathbf{b}^1$   
 $\mathbf{A}^2 \mathbf{x} \leq \mathbf{b}^2$   
 $\mathbf{x} \in \mathbb{Z}_+^n$ 

Usually,  $\mathbf{A}^1\mathbf{x} \leq \mathbf{b}^1$  are the "bad" constraints and  $\mathbf{A}^2\mathbf{x} \leq \mathbf{b}^2$  are the "nice" constraints.

We can derive an relaxation of the IP model by dropping  $\mathbf{A}^1\mathbf{x} \leq \mathbf{b}^1$ , and leave with the constraints that are easy, or relatively easier, to solve. This idea can be embedded into a more general framework called Lagrangian relaxation. Define and rewrite the IP model as

IP(Q) 
$$z_{IP} = \max$$
  $\mathbf{c}\mathbf{x}$   
s.t.  $\mathbf{A}^1\mathbf{x} \leq \mathbf{b}^1$   
 $\mathbf{x} \in Q, \quad Q = {\mathbf{A}^2\mathbf{x} \leq \mathbf{b}^2 | \mathbf{Z}_+^n}$ 

Notice that we have assumed the

# Benders Decomposition

# Total Unimodularity

# Semidefinite Programming

# **Examples: Classical Problems**

- 20.1 Packing and Matching
- 20.2 Knapsack Problem

# Part IV Graph and Network Theory

# Basic Structures

# 21.1 Graph

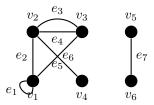
**Definition 21.1.1** (Graph). A graph G consists of a finite set V(G) on vertices, a finite set E(G) on edges and an **incident relation** than associates with any edge  $e \in E(G)$  an unordered pair of vertices not necessarily distinct called **ends**.

**Example.** The following graph can be represented as

$$V = V(G) = \{v_1, v_2, v_3, v_4, v_5, v_6\}$$
(21.1)

$$E = E(G) = \{e_1, e_2, e_3, e_4, e_5, e_6, e_7\}$$
(21.2)

$$e_1 = v_1 v_2, \quad e_2 = v_2 v_4, \quad \dots$$
 (21.3)



**Definition 21.1.2** (Loop, Parallel, Simple Graph). An edge with identical ends is called a **loop**, Two edges having the same ends are said to be **parallel**, A graph without loops or parallel edges is called **simple graph** 

**Definition 21.1.3** (Adjacent). Two edges of a graph are **adjacent** if they have a common end, two vertices are **adjacent** if they are jointed by an edge.

Saving a graph in computer program can be implemented in the following ways:

- Adjacency matrix:  $m \times n$  matrix, for A[u,v] = 1 if  $(u,v) \in E$  and A[u,v] = 0 otherwise
- Linked list: For every vertex v, there is a linked list containing all neighbors of v.

Assuming we are dealing with undirected graphs, n is the number of vertices, m is the number of edges,  $n-1 \le m \le n(n-1)/2$ ,  $d_v$  is the number of neighbors of v then

	Matrix	Linked lists
memory usage	$O(n^2)$	O(m)
time to check $(u, v) \in E$	O(1)	$O(d_u)$
time to list all neighbors of $v$	O(n)	$O(d_v)$

## 21.2 Subgraph

**Definition 21.2.1** (Subgraph). Given two graphs G and H, H is a **subgraph** of G if  $V(H) \subseteq V(G)$ ,  $E(H) \subseteq E(G)$  and an edge has the same ends in H as it does in G. Furthermore, if  $E(H) \neq E(G)$  then H is a proper subgraph.

**Definition 21.2.2** (Spanning). A subgraph H on G is spanning if V(H) = V(G)

**Definition 21.2.3** (Vertex-induced, Edge-induced). For a subset  $V' \subseteq V(G)$  we define an **vertex-induced** subgraph G[V'] to be the subgraph with vertices V' and those edges of G having both ends in V'. The **edge-induced** subgraph G[E'] has edges E' and those vertices of G that are ends to edges in E'.

**Notice:** If we combine node-induced or edge-induced subgraphs G(V') and G(V - V'), we cannot always get the entire graph.

## 21.3 Degree

**Definition 21.3.1** (Degree). Let  $v \in V(G)$ , then the **degree** of  $v \in V(G)$  denote by  $d_G(v)$  is defines to be the number of edges incident of v. Loops counted twice.

**Theorem 21.1.** For any graph G = (V, E)

$$\sum_{v \in V} d(v) = 2|E| \tag{21.4}$$

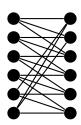
*Proof.*  $\forall$  edge e = uv with  $u \neq v$ , e is counted once for u and once for v, a total of two altogether. If e = uu, a loop, then it is counted twice for u.

**Problem 21.1.** Explain clearly, what is the largest possible number of vertices in a graph with 19 edges and all vertices of degree at least 3. Explain why this is the maximum value.

**Solution.** The maximum number is 12.

*Proof.* First we prove 12 vertices is possible, then we prove 13 vertices is not possible

• The following graph contains 12 vertices and 18 edges, each vertex has a degree of 3.



• For 13 vertices and each vertex has a degree of at least 3 will require at least

$$2|E| = \sum_{v \in V} d(v) \ge 3 \times |N| = 3 \times 13 \Rightarrow |E| \ge 19.5 > 19$$
(21.5)

edges, i.e., 13 vertices is not possible.

Corollary 21.1.1. Every graph has an even number of odd degree vertices.

Proof.

$$V = V_E \cup V_O \Rightarrow \sum_{v \in V} d(v) = \sum_{v \in V_E} d(v) + \sum_{v \in V_O} d(v) = 2|E|$$
 (21.6)

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## 21.4 Special Graphs

**Definition 21.4.1** (Complete Graph). A **complete** graph  $K_n (n \ge 1)$  is a simple graph with n vertices and with exactly one edge between each pair of distinct vertices.

**Definition 21.4.2** (Cycle). A cycle graph  $C_n (n \ge 3)$  consists of n vertices  $v_1, ... v_n$  and n edges  $\{v_1, v_2\}, \{v_2, v_3\}, ... \{v_{n-1}, v_n\}$ 

**Definition 21.4.3** (Wheel). A wheel graph  $W_n (n \ge 3)$  is a simple graph obtains by adding one vertex to the cycle graph  $C_n$ , and connecting this new vertex to all vertices of  $C_n$ 

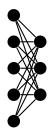
**Definition 21.4.4** (Bipartite Graph). A simple graph is said to be **bipartite** if the vertex set can be expressed as the union of two disjoint non-empty subsets  $V_1$  and  $V_2$  such that every edges has one end in  $V_1$  and another end in  $V_2$ 

**Example.** Here is an example for bipartite graph



**Definition 21.4.5** (Complete Bipartite Graph). The **complete bipartite graph**  $K_{mn}$  is the bipartite graph  $V_1$  containing m vertices and  $V_2$  containing n vertices such that each vertiex in  $V_1$  is adjacent to every vertex in  $V_2$ 

**Example.** Here is an example for  $K_{53}$ 



**Theorem 21.2.** A graph G is bipartite iff every cycle is even.

*Proof.* ( $\Rightarrow$ ) If the graph G is bipartite, by definition, the vertices of graph can be partition into two groups, that within the group there is no connection between vertices. Therefore, for each cycle, the odd index of vertices and even index of vertices has to be choose alternatively from each groups. Therefore the cycle has to be even.

(⇐) Prove by contradiction. A graph can be connected or not connected.

If G is connected and has at least two vertices, for an arbitrary vertex  $v \in V(G)$ , we can calculate the minimum number of edges between the other vertices v' and v (i.e., length, denoted by l(v', v)), for all the vertices that has odd length to v, assign them to set  $V_1$ , for the rest of vertices (and v), assign to set  $V_2$ . Assume that G is not bipartite, which means there are at least one edge between distinct vertices in set  $V_1$  or set  $V_2$ , without lost of generality, assume that edge is uw, u,  $w \in V_1$ . For all vertices in  $V_1$  there is an odd length of path between the vertex and v, therefore, there exists an odd l(u, v), and an odd l(w - v). The length of cycle l(u, w, v) = 1 + l(u, v) + l(w, v), which is an odd number, it contradict with the prerequisite that all cycles are even, which means the assumption that G is not bipartite is incorrect, G should be bipartite.

If G is not connected. Then G can be partition into a set of disjointed subgraphs which are connected with at least two vertices or contains only one vertex. For the component that has more that one vertices, we already proved that it has to be bipartite. For the subgraph  $G_i \subset G$ , i = 1, 2, ..., n, the vertices can be partition into  $V_{i1} \in V(G_i)$  and  $V_{i2} \in V(G_i)$ , where  $V_{i1} \cap V_{i2} = \emptyset$ , the union of those subgraphs are bipartite too because  $V_1 = \bigcup_{i=1}^n V_{i1} \in V(G)$  and  $V_2 = \bigcup_{i=1}^n V_{i2} \in V(G)$  satisfied the condition of bipartite. For the subgraph that has one one vertices, those vertices can be assigned into either  $V_1$  or  $V_2$ .

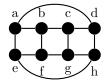
**Example.** The following graph is bipartite, it only contains even cycles.

We can rearrange the graph to be more clear as following

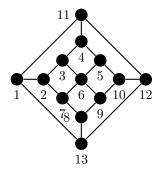
The vertices of graph G can be partition into two sets,  $\{a, c, f, h\}$  and  $\{b, d, e, g\}$ 

**Example.** The following graph is not bipartite

The cycle  $c = v_1 v_1 1 v_4 v_3 v_2$  have odd number of vertices.







# 21.5 Directed Graph

**Definition 21.5.1.** A graph G = (V, E) is called directed if for each edge  $e \in E$ , there is a **head**  $h(e) \in V$  and a **tail**  $t(e) \in V$  and the edges of e are precisely h(e) and t(e), denoted e = (t(e), h(e))

**Definition 21.5.2.** We call directed graphs **digraphs**, we call edges in a digraph are called **arcs**, and vertices in a digraph **nodes** 

**Definition 21.5.3.** Similar as in the undirected case we have walks, traces, paths and cycles in digraphs.

**Definition 21.5.4.** A vertex  $v \in V$  is **reachable** from a vertex  $u \in V$  if there is a (u, v)-dipath. If at the same time u is reachable from v, they are **strongly connected** 

**Definition 21.5.5.** A digraph is strongly connected if every pair of vertices are strongly connected.

**Definition 21.5.6.** A digraph is **strict** if it has no loops and whenever e and f are parallel, h(e) = t(f)

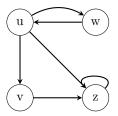
**Definition 21.5.7.** For a vertex v in a digraph D, the **indegree** of v in D, denoted by  $d^+(v)$  is the number of arcs of D having head V. The **outdegree** of v is denoted by  $d^-(v)$  is the number of arcs of D having tail v.

Let D = (V, A) be a digraph with no loops a vertex-arc **incident matrix** for D is a (0, 1, -1) matrix N with rows indexed by  $V = \{v_1, ..., v_n\}$  and column indexed by  $A = \{e_1, ..., e_m\}$  and where entry (i, j) in the matrix  $n_{ij}$  is

$$n_{ij} = \begin{cases} 1, & \text{if } v_i = h(e_j) \\ -1, & \text{if } v_i = t(e_j) \\ 0, & \text{otherwise} \end{cases}$$
 (21.7)

$$\begin{bmatrix} -1 & 0 & -1 & -1 & 1 \\ 1 & -1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & -1 \\ 0 & 1 & 1 & 0 & 0 \end{bmatrix}$$
 (21.8)

# 21.6 Sperner's Lemma



# Paths, Trees, and Cycles

## 22.1 Walk

**Definition 22.1.1** (walk). A walk in a graph G is a finite sequence  $w = v_0 e_1 v_1 e_2 ... e_k v_k$ , where for each  $e_i = v_{i-1} v_i$  the edge and its ends exists in G. We say that walk  $v_0$  to  $v_k$  on  $(v_0, v_k)$ -walk.

Example.

$$w = v_2 e_4 v_3 e_4 v_2 e_5 v_3 \tag{22.1}$$

is a walk, or  $(v_2, v_3)$ -walk

**Definition 22.1.2** (origin, terminal, internal, length). For  $(v_0, v_k)$ -walk, The vertices  $v_0$  and  $v_k$  are called the **origin** and the **terminal** of the walk w,  $v_1...v_{k-1}$  are called **internal** vertices. The integer k is the **length** of the walk. Length of w equals to the number of edges.

We can create a reverse walk  $w^{-1}$  by reversing w.

$$w^{-1} = v_k e_k v_{k-1} e_{k-1} \dots e_2 v_1 (22.2)$$

(The reverse walk is guaranteed to exist because it is an undirected graph)

Given two walks w and w' we can create a third walk denoted by ww' by concating w and w'. The new walk's origin is the same as terminal.

# 22.2 Path and Cycle

**Definition 22.2.1** (trail). A trail is a walk with no repeating edges. e.g.,  $v_3e_4v_2e_5v_3$ 

**Definition 22.2.2** (path). A path is a trail with no repeating vertices. e.g.,  $v_3e_4v_2$ 

Notice: Paths  $\subseteq$  Trails  $\subseteq$  Walks

**Definition 22.2.3** (closed, cycle). A path is **closed** if it has positive length and its origin and terminal are the same. e.g.,  $v_1e_2v_2e_4v_3e_3v_1$ . A closed trail where origin and internal vertices are distinct is called a **cycle** (The only time a vertex is repeated is the origin and terminal)

Definition 22.2.4 (even/odd cycle). A cycle is even if it has a even number of edges otherwise it is odd.

**Problem 22.1.** Prove that if  $C_1$  and  $C_2$  are cycles of a graph, then there exists cycles  $K_1, K_2, ..., K_m$  such that  $E(C_1)\Delta E(C_2) = E(K_1) \cup E(K_2) \cup ... \cup E(K_m)$  and  $E(K_i) \cap E(K_j) = \emptyset, \forall i \neq j$ . (For set X and Y,  $X\Delta Y = (X - Y) \cup (Y - X)$ , and is called the symmetric difference of X and Y)

*Proof.* Proof by constructing  $K_1, K_2, ... K_m$ . Denote

$$C_1 = v_{11}e_{11}v_{12}e_{12}v_{13}e_{13}...v_{1n}e_{1n}v_{11}$$
(22.3)

$$C_2 = v_{21}e_{21}v_{22}e_{22}v_{23}e_{23}...v_{2k}e_{2k}v_{21} (22.4)$$

Assume both cycle start at the same vertice,  $v_{11} = v_{12}$ . (If there is no intersected vertex for  $C_1$  and  $C_2$ , just simply set  $K_1 = C_1$  and  $K_2 = C_2$ )

The following algorithm can give us all  $K_j$ , j = 1, 2, ..., m by constructing  $E(C_1)\Delta E(C_2)$ . Also, the complexity is O(mn), which makes the proof doable.

Now we prove that  $K_i \cap K_j = \emptyset, \forall i \neq j$ . For each  $K_j$ , it is defined by two  $(v_o, v_t)$ -paths in the algorithm. From

### **Algorithm 4** Find $K_1, K_2, ...K_m$ by constructing $E(C_1)\Delta E(C_2)$

```
Require: Graph G, cycle C_1 and C_2
Ensure: K_1, K_2, ... K_m
 1: Initial, K \leftarrow \emptyset, j = 1
 2: Set temporary storage units, v_o \leftarrow v_{11}, v_t \leftarrow \emptyset
 3: for i = 1, 2, ..., n do
 4:
          if e_{1i} \in C_2 then
               if v_o \neq v_{1i} then
 5:
 6:
                    v_t \leftarrow v_{1i}
                    concate (v_o, v_t)-path \subset C_1 and (v_o, v_t)-path \subset C_2 to create a new K_j
 7:
 8:
                    Append K with K_j, K \leftarrow K \cup K_j
 9:
                    Reset temporary storage unit. v_o \leftarrow v_{1(i+1)} (or v_{11} if i = n), v_t \leftarrow \emptyset
10:
                    v_o \leftarrow v_{1(i+1)} \text{ (or } v_{11} \text{ if } i = n)
11:
```

the algorithm we know that all the edges in  $(v_o, v_t)$ -path in  $C_1$  are not intersecting with  $C_2$ , because if the edge in  $C_1$  is intersected with  $C_2$ , either we closed the cycle  $K_j$  before the edge, or we updated  $v_o$  after the edge (start a new  $K_j$  after that edge). By definition of cycle, all the  $(v_o, v_t)$ -path that are subset of  $C_1$  are not intersecting with each other, as well as all the  $(v_o, v_t)$ -path that are subset of  $C_2$ . Therefore,  $K_i \cap K_j = \emptyset, \forall i \neq j$ .

**Definition 22.2.5** (connected vertices). Two vertices u and v in a graph are said to be **connected** if there is a path between u and v.

**Definition 22.2.6** (component). Connectivity between vertices is an equivalence relation on V(G), if  $V_1, ... V_k$  are the corresponding equivalent classes then  $G[V_1]...G[V_k]$  are **components** of G. If graph has only one component, then we say the graph is connected. A graph is connected iff every pair of vertices in G are connected, i.e., there exists a path between every pair of vertices.

**Problem 22.2.** If G is a simple graph with at least two vertices, prove that G has two vertices with the same degree.

*Proof.* A simple graph can only be connected or not connected.

- If G is connected, i.e., for all vertices, the degree is greater than 0. Also the graph is simple, for a graph with |N| vertices, the degree of each vertex is less or equal to |N| 1 (cannot have loop or parallel edge). For |N| vertices, to make sure there is no two vertices that has same degree, it will need |N| options for degrees, however, we only have |N| 1 option. According to pigeon in holes principle, there has to be at least two vertices with the same degree.
- If G is not connected, i.e., the graph has more than one component. One of the following situation will happen:
  - For all components, each component contains only one vertex. Since we have at least two vertices, which means there are at least two component that has only one vertex. For those vertices, at least two vertices has the same degree as 0.
  - At least one component has more than one vertices. In this situation, we can find a component that has more than one vertices as a subgraph G' of the graph G. That G' is a connected simple graph by definition. We have already proved that a connected simple graph has two vertices with the same degree, which means G has two vertices with the same degree.

22.3. TREE AND FOREST

## 22.3 Tree and forest

**Definition 22.3.1** (acyclic graph). A graph is called **acyclic** if it has no cycles

Definition 22.3.2 (forest, tree). A acyclic graph is called a forest. A connected forest is called a tree.

**Theorem 22.1.** Prove that T is a tree, if T has exactly one more vertex than it has edges.

*Proof.* 1. First we prove for any tree T that has at least two vertices, there has to be at least one leaf, i.e., now we prove that we can find u with degree of 1. Proof by constructing algorithm. (In fact we can prove that there are at least two leaves.)

The above algorithm is guaranteed to have an end because a tree is acyclic by definition

#### **Algorithm 5** Find one leaf in a tree

```
Require: d(u) = 1
```

**Ensure:** A tree T has at least one vertex

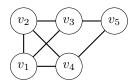
- 1: Let u and v be any distinct vertex in a tree T
- 2: Let p be the path between u and v
- 3: while  $d(u) \neq 1$  do
- 4: **if** d(u) > 1 **then**
- 5: Let n(u) be the set of neighboring vertices of u
- 6: In n(u), find a u' that the edge between u and u', denoted by  $e, e \notin p$
- 7:  $u \leftarrow u'$
- 8:  $p \leftarrow p \cup e$ 
  - 2. Then, if we remove one leaf in the tree, i.e., we remove an edge and a vertex, where that vertex only connects to the edge we removed. One of the following situations will happen:
    - (a) Situation 1: The remaining of T is one vertex. In this case, T has two vertices an one edge. (Exactly one more vertex than it has edges)
    - (b) Situation 2: The remaining of T is another tree T' (removal of edges will not change acyclic and connectivity), where |V(T)| = |V(T')| + 1 and |E(T)| = |E(V')| + 1. (one edge and one vertex has been removed)
  - 3. Do the leaf removal process recursively to T' if Situation 2 happens until Situation 1 happens.

# 22.4 Spanning tree

**Definition 22.4.1** (spanning tree). A subgraph T of G is a spanning tree if it is spanning (V(T) = V(G)) and it is a tree.

Example. In the following graph

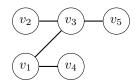
This is a spanning tree



**Problem 22.3.** Prove that if  $T_1$  and  $T_2$  are spanning trees of G and  $e \in E(T_1)$ , then there exists a  $f \in E(T_2)$ , such that  $T_1 - e + f$  and  $T_2 + e - f$  are both spanning trees of G.

*Proof.* One of the following situation has to happen:

1. If for given  $e \in E(T_1)$ ,  $\exists f = e \in E(T_2)$ , then  $T_1 - e + f = T_1$ ,  $T_2 + e - f = T_2$  are both spanning trees of G



- 2. If for given  $e \in E(T_1)$ ,  $e \notin E(T_2)$ , the following will find an edge f that  $T_1 e + f$  and  $T_2 + e f$  are both spanning trees of G.
  - (a)  $T_1$  is a spanning tree, removal of  $e \in E(T_1)$  will disconnect the spanning tree into two components (by definition of spanning tree), denoted by  $G_1 \subset G$  and  $G_2 \subset G$ , by definition,  $V(G_1)$  and  $V(G_2)$  is a partition of V(G).
  - (b) Add e into  $T_2$ . We can proof that by adding an edge into a tree will create exactly one cycle, denoted by C,  $e \in E(C)$ .
  - (c) For C, since it is a cycle and one end of e is in  $V(G_1)$ , the other end of e is in  $V(G_2)$ , there has to be at least two edges (can be more) that has one end in  $V(G_1)$  and the other end in  $V(G_2)$ , denote the set of those edges as  $E \subset E(C)$ , one of those edges is  $e \in E$
  - (d) Choose any  $f \in E$  and  $f \neq e$ , for that  $f, T_1 e + f$  and  $T_2 + e f$  are both spanning trees of G.
  - (e) Prove that  $T_1 e + f$  is a spanning tree
    - i.  $T_1 e + f$  have the same set of vertices as  $T_1$ , therefore it is spanning.
    - ii. It is connected both within  $G_1$  and  $G_2$ , for f, one end is in  $V(G_1)$ , the other end is in  $V(G_2)$  therefore  $T_1 e + f$  is connected.
    - iii.  $T_1 e + f$  have the same number of edges as  $T_1$ , which is  $|T_1| 1$ , therefore  $T_1 e + f$  is a tree. (We have proven the connectivity in the previous step.)
    - iv.  $T_1 e + f$  is spanning, connected, a tree, therefore it is a spanning tree.
  - (f) Prove that  $T_2 + e f$  is a spanning tree
    - i.  $T_2 + e f$  have the same set of vertices as  $T_2$ , therefore it is spanning.
    - ii.  $T_2$  is connected, adding an edge will not break connectivity, therefore  $T_2 + e$  is connected, removing an edge in a cycle will not break connectivity, therefore  $T_2 + e f$  is connected.
    - iii.  $T_2 e + f$  have the same number of edges as  $T_2$ , which is  $|T_2| 1$ , therefore  $T_2 + e f$  is a tree. (We have proven the connectivity in the previous step.)
    - iv.  $T_2 e + f$  is spanning, connected, a tree, therefore it is a spanning tree.

**Theorem 22.2.** Every connected graph has a spanning tree.

*Proof.* Prove by constructing algorithm:

Algorithm 6 Find a spanning tree for connected graph (Prim's Algorithm in unweighted graph)

**Require:** a connected graph G and an enumeration  $e_1, ... e_m$  of the edges of G

Ensure: a spanning tree T of G

- 1: Let T be the spanning subgraph of G with V(T) = V(G) and  $E(T) = \emptyset$
- $2: i \leftarrow 1$
- 3: while  $i \leq |E|$  do
- 4: **if**  $T + e_i$  is acyclic **then**
- 5:  $T \leftarrow T + e_i$
- 6:  $i \leftarrow i+1$

**Notice:** This algorithm can be improved, one idea is to make summation of edges in spanning subgraph less or equation to |V|-1

For the complexity of spanning tree algorithm:

- 1. Space complexity, 2|E|, which is O(|E|)
- 2. Time complexity
  - (a) How to check for acyclic?
    - i. At every stage T has certain components  $V_1,...V_t$ , (every time we add an edge, the number of components minus 1)
    - ii. So at the beginning t = |V| with  $|V_i| = 1 \forall i$  and at the end, t = 1.
  - (b) Count the amount of work for the algorithm.
    - i. Need to check for acyclic for each edge, which costs O(|E|)
    - ii. Need to flip the pointer for each vertex, for each vertex, at most will be flipped  $\log |V|$  times, altogether  $|V| \log |V|$  times.
    - iii. The time complexity is  $O(|E| + |V| \log |V|)$
- 3. First we need to input the data, create an array such that the first and the second entries are the ends of  $e_1$ , third and fourth are the ends of  $e_2$ , and so on.
- 4. The amount of storage needs in 2|E|, which is O(|E|)
- 5. The main work involved in the algorithm is for each edges  $e_i$  and the current T, to determine if  $T + e_i$  creates a cycle.
- 6. suppose we keep each component  $V_i$  by keeping for each vertex a pointer from the vertex to the name of the component containing it. Thus if  $\mu \in V_3$ , there will be a pointer from  $\mu$  to integer 3.
- 7. Then when edge  $e_i = \mu v$  is encountered in Step 2, we see that  $T + e_i$  contains a cycle if and only if  $\mu$  and v point to same integer which means they are in the same component
- 8. If they are not in the same component, we want to add the edge which means then I have to update the pointers.

To prove algorithm we need to show the output is a spanning tree, which means three properties must hold:

- spanning (Step I)
- acyclic (We never add an edge that create a cycle)
- connected (Proof by contradiction)

So it is sufficient to show that the output will be connected.

Proof. (Proof by Contradiction) Suppose the output graph T of the algorithm is NOT connected. Let  $T_1$  be a component of T, let  $x \in T_1$  and  $y \notin T_1$ . But G is a connected graph (given from the beginning), so there must be a path in G that connects x and y. Let such a path in G be  $p = xe_1v_1e_2, ...v_{k-1}e_ky$ . Clearly,  $p \notin T_1$ . So there must be a first vertex in P that not in  $T_1$ . So  $e_i \notin E(T)$ , the only way this can happen when applying the algorithm is if  $T + e_i$  creates a cycle C, i.e.,  $e_i \in C$ , so  $C - e_i$  is a path connecting  $v_{i-1}$  and  $v_i$ . So  $c - e_i \in T$ , so  $v_{i-1}$  is connected to  $v_i \in T$ . Contradiction.

# 22.5 Cayley's Formula

# 22.6 Connectivity, DFS, BFS

## 22.6.1 Connectivity Problem

For connectivity problem, the input is a graph G = (V, E), with linked list representation, and two vertices  $s, t \in V$ . The problem is to find whether there is a path connecting s to t in G

There are two commonly use methods to solve connectivity problem: depth-first search and breath-first search.

## 22.6.2 Depth-First Search (DFS)

The idea of DFS is enumerating children before siblings when search in the graph/tree. It is a recursive algorithm. First, start with s, then travel through the fist edge leading out of the current vertex, when reached a "visited" vertex, go back and travel through next edge. If tried all edges leading out of the current vertex, go back. The algorithm is as following

#### Algorithm 7 Depth-First Search

- 1: Initialize, make all vertices as "unvisited"
- 2: RecursiveDFS(s)

#### **Algorithm 8** RecursiveDFS(v)

```
1: Mark v as "visited"

2: for \forall u \in d_s do:

3: if u is "unvisited" then

4: RecursiveDFS(u)
```

The running time of DFS is O(n+m).

If we only want to know if s and t are connected, the algorithm can be terminated if u = t.

## 22.6.3 Breadth-First Search (BFS)

The idea of BFS is enumerating siblings before children when search in the graph/tree. The general steps of BFS is as following: First, build layers  $L_0, L_1, \dots$ ; Next, set  $L_0 = \{s\}$ , where s is the starting vertex; Then,  $L_{j+1}$  contains all nodes that are not in  $\bigcup_{i=0}^{j} L_i$  and have an edge to a vertex in  $L_j$ 

The algorithm is as following

#### Algorithm 9 Breadth-First Search

```
1: Initialize, head \leftarrow 1, tail \leftarrow 1, queue[1] \leftarrow s, mark all vertices as "unvisited"

2: Mark s as "visited"

3: while head \geq tail do

4: v \leftarrow queue[tail], tail \leftarrow tail + 1

5: for \forall u \in d_v do

6: if u is "unvisited" then

7: head \leftarrow head + 1, queue[head] = u

8: Mark u as "visited"
```

The running time of BFS is O(n+m)

If we only want to know if s and t are connected, the algorithm can be terminated if u = t.

#### 22.6.4 Cycle detection

The following algorithm is for connected graph, if the graph is not connected, run the algorithm for each component until cycle is detected or all the components have been calculated. Since the complexity for running in connected graph is O(n+m), n as the number of vertices/nodes, and m as the number of edges/arcs, the running time of disconnected graph is the **summation** of running time in each component, where each component is connected. Therefore the complexity is the same in disconnected graph as in connected graph.

The main idea is starting with arbitrary vertex/node, using DFS or BFS to search on the graph try to revisit the vertex/node we start with. If succeed, a cycle is detected, otherwise if all the vertices/nodes has been visited, then no cycle exists. And in linked-list representation, the complexity is O(|V| + |E|), i.e. O(n+m). However, there is slightly different in undirected graph and directed graph, for undirected graph needs at least three vertices to form a cycle while directed graph needs at least two.

Here is the detail algorithm (DFS) for undirected graph:

#### Algorithm 10 Main algorithm

```
    For all nodes, labeled as "unvisited"
    Arbitrary choose a vertex v, add a dummy vertex w, add a dummy edge (w, v), label w as "visited"
    run DFS(w, v)
    Remove dummy vertex w and dummy edge (w, v)
    if DFS(w, v) returns "Cycle is found" then
    return "Cycle is found"
    else
    return "No cycle detected"
```

#### **Algorithm 11** DFS(w, v)

```
1: Label v as "visited"
2: if number of v 's neighbor is 1 then
3: return null
4: else
5: for all neighbor u in linked-list of v excepts w do
6: if u is labeled as "visited" then
7: return "Cycle is found"
8: else
9: run DFS(v, u)
```

Now check the complexity. Denote  $v \in N$  as a node in graph G, total number of nodes denoted by n, denote  $d_v$  as number of neighbors of node v. The complexity of DFS(w,v) is  $O(d_v)$  for each node v it visited (it should be O(v) because we need O(1) to check if a node is w), each node can only be visited, by "visited" it means DFS(w,v) is executed, at most once, which is controlled by the "visited" label. The total complexity is  $O(n + \sum_{v \in N} d_v) = O(n + m)$ 

For directed graph, we assume in the linked-list of each node, 1) it only contains the vertices where the node leading out to, or 2) we can distinguish the vertices it leading out to in O(1). Here is the detail algorithm (DFS) for directed graph:

#### Algorithm 12 Main algorithm

```
1: For all nodes, labeled as "unvisited"
2: Arbitrary choose a node v, run DFS(v)
3: if DFS(v) returns "Cycle is found" then
4: return "Cycle is found"
5: else
6: return "No cycle detected"
```

## **Algorithm 13** DFS(v)

```
1: Label v as "visited"
2: if number of vertices v leading out to is 0 then
3: return null
4: else
5: for all leading out vertices u in linked-list of v do
6: if u is labeled as "visited" then
7: return "Cycle is found"
8: else
9: run DFS(u)
```

Now check the complexity. Similar to undirected case, denote  $v \in N$  as a node in graph G, total number of nodes denoted by n, denote  $d_v$  as number of edges leading out from node v. The complexity of DFS(v) is  $O(d_v)$  for each

node v it visited, each node can only be visited, by "visited" it means DFS(v) is executed, at most once, which is controlled by the "visited" label. The total complexity is  $O(n + \sum_{v \in N} d_v) = O(n + m)$ 

Also, for undirected (connected) graph, if we don't need to find the cycle and only need to decide if there is a cycle or not, we can just check the number of vertices, n, and number of edges, m. If  $n \le m$  then G contains a cycle, otherwise no cycle, the complexity is O(1).

## 22.6.5 Test Bipartiteness

We have proved that a bipartite graph only has even cycles, and the graph with only even cycles are bipartite graph, however, that is not very continent to test if a graph is bipartite because it needs to enumerate all cycles.

The other idea to test bipartiteness is try to color the vertices of the graph, if it can be 2-colored, then the graph is bipartite, otherwise it is not.

The following is the algorithm (using BFS)

#### Algorithm 14 Test Bipartiteness

```
1: Initialize, head \leftarrow 1, tail \leftarrow 1, queue[1] \leftarrow s, mark all vertices as "unvisited"
 2: Mark s as "visited"
 3: color[s] \leftarrow 0
 4: while head \ge tail do
        v \leftarrow queue[tail], tail \leftarrow tail + 1
 5:
        for \forall u \in d_v do
 6:
             if u is "unvisited" then
 7:
                 head \leftarrow head + 1, queue[head] = u
 8:
                 Mark u as "visited"
 9:
                 color[u] \leftarrow 1 - color[v]
10:
             else
11:
                 if color[u] == color[v] then return False
12:
    return True
```

## 22.6.6 Topological Ordering

The topological ordering problem is given a directed acyclic graph G = (V, E), output a 1-to-1 function  $\pi : V \to \{1, 2, \dots, n\}$  so that if  $(u, v) \in E$ ,  $\pi(u) < \pi(v)$ 

The idea is each time take a vertex without incoming edges, then remove the vertex and all its outgoing edges. To make the algorithm efficient, we can 1) use linked lists of outgoing edges; 2) maintain the in degree  $d_v$  of vertices; 3) maintain a queue (or stack) of vertices v with  $d_v = 0$ 

The following is the algorithm

## **Algorithm 15** topological-sort(G)

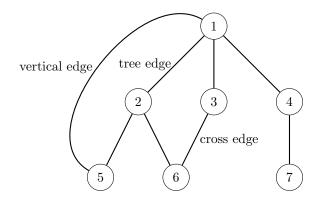
```
1: Initialize, let \forall v \in V, d_v = 0
 2: for v \in V do
          for u, (u, v) \in E do
 3:
               d_u \leftarrow d_u + 1
 5: S \leftarrow \{v : d_v = 0\}, i \leftarrow 0
 6: while S \neq \emptyset do
          v \leftarrow v \in S, S \leftarrow S \setminus \{v\}
 7:
          i \leftarrow i + 1, \pi(v) \leftarrow i
 8:
          for u, (u, v) \in E do
 9:
               d_u \leftarrow d_u - 1
10:
               if d_u = 0 then
11:
                    S \leftarrow S \cup \{u\}
13: if i < n then return Not directed acyclic graph
```

22.7. BLOCKS 103

## 22.6.7 Bridge

**Definition 22.6.1** (Tree Edge, Cross Edge, Vertical Edge). Given a graph G = (V, E) and a rooted tree T in G, edge in G can be classified by three types:

- Tree edges: edge in T
- Cross edges: (u, v), u and v do not have an ancestor-descendant relation
- Vertical edges: (u, v), u is an ancestor of v, or v is an ancestor of u



In a BFS tree T of a graph G, there can not be vertical edges, there cannot be cross edges (u, v) with u and v 2 levels apart. (Cross edge at most 1 level apart)

In a DFS tree T of a graph G, there can not be cross edges, there can only be tree edges and vertical edges.

**Definition 22.6.2** (bridge). Given a connected graph G = (V, E), an edge  $e \in E$  is called a **bridge** if the graph  $G = (V, E \setminus \{e\})$  is disconnected.

The idea to find bridge is through a DFS tree. Notice that there are only tree edges and vertical edges in DFS tree. Vertical edges are not bridges, a tree edge (u, v) is not a bridge if some vertical edge jumping from below u to above v. Other tree edges are bridges.

Define level(v) as the level of vertex v in DFS tree.  $T_v$  as the sub tree rooted at v, h(v) as the smallest level that can be reached using a vertical edge from vertices in  $T_v$ . (parent(u), u) is a bridge if  $h(u) \ge level(u)$ . The algorithm is as following:

## Algorithm 16 FindBridge(G)

- 1: Mark all vertices as "unvisited"
- 2: for  $v \in V$  do
- 3: **if** v is "unvisited" **then**
- 4:  $level(v) \leftarrow 0$
- 5: RecursiveDFS(v)

## 22.7 Blocks

## Algorithm 17 RecursiveDFS(v)

```
1: mark v as "visited"
 2: h(v) \leftarrow \infty
 3: for u \in d_v do
         \mathbf{if}\ u \ \mathrm{is}\ \text{``unvisited''}\ \mathbf{then}
             level(u) \leftarrow level(v) + 1
 5:
             RecursiveDFS(u)
 6:
             if h(u) \ge level(u) then
 7:
                  (u, v) is a bridge
 8:
             if h(u) < h(v) then
 9:
10:
                  h(v) \leftarrow h(u)
         else
11:
12:
             if level(u) < level(v) - 1 and level(u) < h(v) then
                  h(v) \leftarrow level(u)
13:
```

# **Euler Tours and Hamilton Cycles**

- 23.1 Euler Tours
- 23.2 Hamilton Cycles
- 23.3 The Chinese Postman Problem
- 23.4 The Traveling Salesman Problem

# Matriod, Planarity

#### Plane and Planar Graphs 24.1

#### 24.2 **Dual Graphs**

#### 24.3 Matroids

**Definition 24.3.1** (Matroids). Let S be a finite set of **elements** and let d be a collection of subsets of S satisfying the property

If 
$$x \le y, y \in d, \Rightarrow x \in d$$
 (24.1)

The pair (S,d) is called an **independent system** and the members of d are called **independent sets**.

**Example.** Let G be a graph and let  $S \in E(G)$  define  $M \subseteq S$  to be independent if M is a matching

$$S = \{(1,2), (2,3), (2,4), (3,5), (4,6), (5,6)\}$$
(24.2)

$$d = \{\emptyset, \{(1,2)\}, \{(2,3)\}, \dots, \{\text{other matching...}\}\}$$
(24.3)

**Example.** Let G be a graph and let S = V(G) define  $X \subseteq S$  to be independent if no two member of x are adjacent.

$$S = \{1, 2, 3, 4\} \tag{24.4}$$

$$d = \{\emptyset, 1, 2, 3, 4, (1, 3), (1, 4), (3, 4), (1, 3, 4)\}$$
(24.5)

**Example.** Let G be a connected graph and let S = E(G), define  $X \subseteq S$  to be independent if G[X] contains cycles.

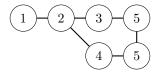
Given any independent system, there is a natural combinatorial optimization problem of finding the maximum cardinality independent set.

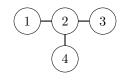
Let's try the following: Greedy algorithm

Step 1: Set  $I = \emptyset$ 

Step 2: If there exists  $e \in S \setminus I$  such that I + e is independent, set  $I \leftarrow I + e$  and go to Step 1, otherwise stop. Those Independent systems for which the greedy algorithm guarantee to find a maximum cardinality independent

set are very special called matroids





- 24.4 Independent Sets
- 24.5 Ramsey's Theorem
- 24.6 Turán's Theorem
- 24.7 Schur's Theorem
- 24.8 Euler's Formula
- 24.9 Bridges
- 24.10 Kuratowski's Theorem
- 24.11 Four-Color Theorem
- 24.12 Graphs on other surfaces

## Matchings

#### 25.1 Maximum Matching

**Definition 25.1.1** (Matching). Let G = (V, E) be a graph, a **matching** is a subset of edges  $M \subseteq E$  such that no two elements of M are adjacent. The two ends of an edge in M are said to be **matched under** M. A matching M saturates a vertex v, and v is said to be **M-saturated** or **M-covered**, if some edge of M is incident with v. Otherwise, v is **M-unsaturated** or **M-exposed**.

**Definition 25.1.2** (Perfect matching, Maximum matching). If every vertex of G is M-saturated, then the matching is said to be **perfect matching**. M is a **maximum matching** if G has no matching M' with |M'| > |M|. Every perfect matching is maximum. The maximum matching does not necessarily to be perfect. Perfect matching and maximum matching may not be unique.

**Definition 25.1.3** (M-alternating). An **M-alternating** path in G is a path whose edges are alternately in  $E \setminus M$  and M.

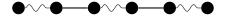
**Definition 25.1.4** (M-augmenting). An **M-augmenting** path in G is an M-alternating path whose origin and terminus are M-unsaturated.

**Lemma 25.1.** Every augmenting path P has property that let  $M' = P\Delta M = (M \cup P) \setminus (M \cap P)$  then M' contains one more edge then M

The following path is an M-augmenting path



The following path is  $M' = P\Delta M(M \cup P) \setminus (M \cap P)$  and all the vertices are M-saturated.



**Theorem 25.2** (Berge, 1957). A matching M in a graph G is maximum iff G has no M-augmenting path.

*Proof.* ( $\Rightarrow$ ) It is clear that if M is maximum, it has no augmenting paths since otherwise by problem claim we can increase by one.

( $\Leftarrow$ ) Suppose M is not maximum and let M' be a bigger matching. Let  $A = M\Delta M'$  now no vertex of G is incident to more than two members of A. For otherwise either two members of M or two members of M' would be adjacent. Contradict the definition of matching. It follows that every component of the edges incident subgraph G[A] is either an even cycle with edge augmenting in  $M\Delta M'$  or else A path with edges alternating between M and M'. Since  $|M'| \geq |M|$  then the even cycle cannot help because exchanging M and M' will have same cardinality. The path case implies that p is alternating in M and since |M'| > |M| the end arc exposed so that p is augmenting.

**Definition 25.1.5** (Vertex-cover). The **vertex-cover** is a subset of vertices X such that every edge of G is incident to some member of X.

**Lemma 25.3.** The cardinality of any matching is less than or equal to the cardinality of any vertex cover.

*Proof.* Consider any matching. Any vertex cover must have nodes that at least incident to the edges in the matching. Since all the edges in the matching are disjointed, so for a single node can at most cover one edge in the matching. If the matching is not perfect, for the edges that not in the tree, they may or may not be possible to be covered by the nodes incident to the edges in the matching, with an easy triangle graph example, we can prove this lemma.  $\Box$ 

**Theorem 25.4** (König Theorem). If G is bipartite, the cardinality of the maximum matching is equal to the cardinality of the minimum vertex cover.

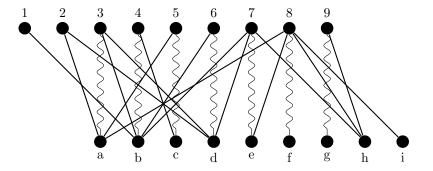
*Proof.* Let G be a bipartite graph, G = (V, E) where  $V = X \cup Y$  as X and Y are two disjointed sets of vertices. Let M be a maximum matching on G. For each edge in M, denoted by  $e_i = a_i b_i$  where  $e_i \in M$ ,  $a_i \in A$  and  $b_i \in B$  and  $A = \{a_i : e_i \in M\} \subseteq X$ , and  $B = \{b_i : e_i \in M\} \subseteq Y$ . Therefore, we can partition X by A and  $U = X \setminus A$ , partition Y by B and  $W = Y \setminus B$ .

We can further partition the matching M into  $M_1$  and  $M_2$ . For all the edges in  $M_1$  can be included into an M-alternating path starts from a vertex in U (which includes the edges directly linked to vertices in U), and  $M_2 = M \setminus M_1$ . For edges in  $M_1$ , we take the ends of edges in  $M_2$  in the vertex cover, denoted by  $M_1$ , take the ends of edges in  $M_2$ , we take the ends of edges in  $M_3$  in the vertex cover, denoted by  $M_3$ , and the ends of edges in  $M_3$  as a subset denoted by  $M_3$ , and the ends of edges in  $M_3$  as a subset denoted by  $M_3$ .

We claim that all the vertices in U can only be connected to vertices in  $B_1$  and vertices in W can only be connected to vertices in  $A_2$ .

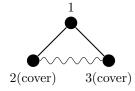
 $U \subset X$  connects to vertices in  $B_1$  by definition. If vertices in  $W \subset Y$  is connected to vertices in  $A_1$ , then we will have M-augmenting path which is contradicted to the assumption that M is maximum matching.

The following is an example. Where the edge in the matching that accessible from members of  $U = \{1, 2\}$  in an M-alternating path is edge 3a, 4b, 5c, 6d.



In which  $U = \{1, 2\}$ ,  $M_1 = \{3a, 4b, 5c, 6d\}$ .  $U = \{1, 2, \}$ ,  $A_1 = \{3, 4, 5, 6\}$ ,  $A_2 = \{7, 8, 9\}$ ,  $W = \{h, i\}$ ,  $B_1 = \{a, b, c, d\}$ ,  $B_2 = \{e, f, g\}$ . The vertex cover is  $\{a, b, c, d, 7, 8, 9\}$ .

The above theorem does not apply to non-bipartite graph. The following is an example



The maximum matching has one edge, where the minimum cover has two vertices.

## 25.2 Maximum Matching Algorithm

**Definition 25.2.1** (M-alternating tree). An M-alternating tree T is a rooted tree satisfied the following condition:

- The root r is M-unsaturated
- The unique path from r to any vertex T is M-alternating
- Every vertex in T, except r is incident to a matching edge of T

A vertex x of T is called **inner** if (r, s)-path in T has an odd number of edges. Otherwise x is called **outer**.

**Lemma 25.5.** Let M be a matching in G and let T be an M-alternating tree with root r, then the following conclusion hold

- If  $v \neq r$  is an outer vertex and p is the unique (r, v)-path in T then the edge of p incident to v is in M
- The number of inner vertices in T equals the number of matching edges in T.

**Definition 25.2.2** (Alternating forest). An alternating forest is a forest of G where every components is an alternating tree. An alternating forest (F, e) is an alternating forest to be then with an edge e = M, V where M and V are outer vertices contained in two distinct components of F.

**Example** (Hungarian forest). A **Hungarian forest** F is an alternating forest containing all exposed vertices of G and such that the outer vertices of F are adjacent in G only to inner vertices of F

**Definition 25.2.3** (Augmenting forest). An augmenting forest (F, e) where F is an augmenting forest and e = uv connects two outer vertices in distinct components of F.

The plan is to grow an alternating forest that eventually become augmenting or Hungarian. Augmenting forest will increase the cardinality of the match, Hungarian implies that you have found optimal maximum cardinality matching.

**Theorem 25.6.** Let (F, e) be an augmenting forest. And let  $T_1$  and  $T_2$  be the two components of F containing an end of e. Let  $p_i$  be the unique path in  $T_i$  from the root to the end of e, then  $p_1ep_2^{-1}$  is an augmenting path.

**Theorem 25.7.** If F is a Hungarian forest for some matching M then M is a maximum match.

The above theorems suggest a method for computing maximum matching. Let M be a matching of G and let F be an alternating forest in G made up of all M-exposing vertices. If F happens to be Hungarian, stop with the max matching M. If (F, e) for some e is augmenting then we increase our matching by 1 and start process.

Suppose F is neither Hungarian nor augmenting, by definition, there must be an edge e incident to an outer vertex of F to no inner vertex of F. But e cannot be incident to two outer vertices of F in distinct components, since (F, e) is not augmenting. Hence there are only two cases:

Case 1: e = uv where u is outer in F and v is not outer in F. The only way its possible if v is covered. Augmenting F to M will increase the matching.

Case 2: e = uv where u and v are outer vertices in F. Let r be the root of the component of F containing u and v, let  $p_u$  and  $p_v$  be (r, u)-path and (r, v)-path in F. Let b be the last vertex these two paths have in common and let p be the (u, v)-path in F, let  $p'_u$  be (b - u)-path and  $p'_v$  be the (b, v)-path respectively. Let n be the length of  $p_u$ , let m be the length of  $p_v$ , let  $p'_v$  be the length of  $p'_v$  be the cycle  $p'_v$  be the cycle  $p'_v$ . Number of vertices in  $p'_v$  is  $p'_v$  and  $p'_v$  be the length cycle. If  $p'_v$  has no odd cycles, we call those graph bipartite. To this case can not happen in bipartite graph so algorithm without case 2 will solve bipartite matching. If a graph has no odd cycles, i.e., bipartite, then we have an algorithm using augmenting forest and Hungarian forest and case 1.

We now have to deal with odd cycles. The idea is to "shrink" add cycles to a super node Let  $S \subseteq E(G)$ , denote by G : S the subgraph with edge set S

$$G: S = G \setminus (E(G) - S) \tag{25.1}$$

The contraction of S to be the  $G \setminus S$  with E(G/S) = E(G) - S. V(G/S) to be the components of G : S and if  $e \in E(G/S)$  then the ends of e in G/S are in components of G : S containing both ends in G. Let network to general case 2. b is outer vertex. Let B be the set of edges of cycle C, we call B a **blossom**. We propose to replace M by M - B, G by G/B and F by F/B

## 25.3 Edmonds's Blossom Algorithm

 $O(|V|^4)$  - Non-bipartite matching is one of very few problems in P, for which LP relaxation will not provide optimal solution.

- 25.4 Hall's Marriage Theorem
- 25.5 Transversal Theory
- 25.6 Menger's Theorem
- 25.7 The Hungarian Algorithm

## Colorings

26.1	Edge Chromatic Number
26.2	Vizing's Theorem
26.3	The Timetabling Problem
26.4	Vertex Chromatic Number
26.5	Brooks' Theorem
26.6	Hajós' Theorem
26.7	Chromatic Polynomials
26.8	Girth and Chromatic Number

## Minimum Spanning Tree Problem

#### 27.1 Basic Concepts

**Example.** A company wants to build a communication network for their offices. For a link between office v and office w, there is a cost  $c_{vw}$ . If an office is connected to another office, then they are connected to with all its neighbors. Company wants to minimize the cost of communication networks.

**Definition 27.1.1** (Cut vertex). A vertex v of a connected graph G is a **cut vertex** if  $G \setminus v$  is not connected.

**Definition 27.1.2** (Connection problem). Given a connected graph G and a positive cost  $C_e$  for each  $e \in E$ , find a minimum-cost spanning connected subgraph of G. (Cycles all allowed)

**Lemma 27.1.** An edge  $e = uv \in G$  is an edge of a cycle of G iff there is a path  $G \setminus e$  from u to v.

**Definition 27.1.3** (Minumum spanning tree problem). Given a connected graph graph G, and a cost  $C_e$ ,  $\forall e \in E$ , find a minimum cost spanning tree of G

The only way a connection problem will be different than MSP is if we relax the restriction on  $C_e > 0$  in the connection problem.

### 27.2 Kroskal's Algorithm

**Algorithm 18** Kroskal's Algorithm,  $O(m \log m)$ 

Require: A connected graph

Ensure: A MST

Keep a spanning forest H = (V, F) of G, with  $F = \emptyset$ 

**while** |F| < |V| - 1 **do** 

add to F a least-cost edge  $e \notin F$  such that H remains a forest.

### 27.3 Prim's Algorithm

**Algorithm 19** Prim's Algorithm, O(nm)

Require: A connected graph

Ensure: A MST

Keep H = (V(H), T) with  $V(H) = \{v\}$ , where  $r \in V(G)$  and  $T = \emptyset$ 

while |V(T)| < |V| do

Add to T a least-cost edge  $e \notin T$  such that H remains a tree.

#### 27.4 Extensible MST

**Definition 27.4.1** (cut). For a graph G = (V, E) and  $A \subseteq V$  we denote  $\delta(A) = \{e \in E : e \text{ has an end in } A \text{ and an end in } V \setminus A\}$ . A set of the form  $\delta(A)$  for some A is called a **cut** of G.

**Definition 27.4.2.** We also define  $\gamma(A) = \{e \in E : \text{both ends of } e \text{ are in } A\}$ 

**Theorem 27.2.** A graph G = (V, E) is connected iff there is no  $A \subseteq V$  such that  $\emptyset \neq A \neq V$  with  $\delta(A) = \emptyset$ 

**Definition 27.4.3.** Let us call a subset  $A \in E$  extensible to a minimum spanning tree problem if A is contained in the edge set of some MST of G

**Theorem 27.3.** Suppose  $B \subseteq E$ , that B is extensible to an MST and that e is a minimum cost edge of some cut D satisfying  $D \cap B = \emptyset$ , then  $B \cup \{e\}$  is extensible to an MST.

#### Solve MST in LP 27.5

Given a connected graph G = (V, E) and a cost on the edges  $C_e$  for all  $e \in E$ , Then we can formulate the following LP

$$X_e = \begin{cases} 1, & \text{if edge } e \text{ is in the optimal solution} \\ 0, & \text{otherwise} \end{cases}$$
 (27.1)

The formulation is as following

min 
$$\sum_{e \in E} c_e x_e$$
 (27.2)  
s.t. 
$$\sum_{e \in E} x_e = |V| - 1$$
 (27.3)

s.t. 
$$\sum_{e \in E} x_e = |V| - 1$$
 (27.3)

$$x_e \ge 0 \tag{27.4}$$

$$e \in E \tag{27.5}$$

$$\sum_{e \in E(S)} x_e = |S| - 1, \forall S \subseteq V, S \neq \emptyset$$
(27.6)

(27.7)

## Shortest-Path Problem

#### 28.1 Basic Concepts

All Shortest-Path methods are based on the same concept, suppose we know there exists a dipath from r to v of a cost  $y_v$ . For each vertex  $v \in V$  and we find an arc  $(v, w) \in E$  satisfying  $y_v + v_{vw} < y_w$ . Since appending (v, w) to the dipath to v takes a cheaper dipath to w then we can update  $y_w$  to a lower cost dipath.

**Definition 28.1.1** (feasible potential). We call  $y = (y_v : v \in V)$  a **feasible potential** if it satisfies

$$y_v + c_{vw} \ge y_w \quad \forall (v, w) \in E \tag{28.1}$$

and  $y_r = 0$ 

**Proposition 5.** Feasible potential provides lower bound for the shortest path cost.

*Proof.* Suppose that you have a dipath  $P = v_0 e_1 v_1, ..., e_k v_k$  where  $v_0 = r$  and  $v_k = v$ , then

$$C(P) = \sum_{i=1}^{k} C_{e_i} \ge \sum_{i=1}^{k} (y_{v_i} - y_{v_{i-1}}) = y_{v_k} - y_{v_0}$$
(28.2)

28.2 Breadth-First Search Algorithm

#### 28.3 Ford's Method

Define a predecessor function  $P(w), \forall w \in V$  and set P(w) to v whenever  $y_w$  is set to  $y_v + c_{vw}$ 

#### Algorithm 20 Ford's Method

**Ensure:** Shortest Paths from r to all other nodes in V Require: A digraph with arc costs, starting node r

Initialize,  $y_r = 0$  and  $y_v = \infty, v \in V \setminus r$ 

Initialize,  $P(r) = 0, P(v) = -1, \forall v \in V \setminus r$ 

while y is not a feasible potential do

Let  $e = (v, w) \in E$  (this could be problematic)

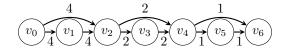
if  $y_v + c_{vw} < y_w$  (incorrect) then

 $y_w \leftarrow y_v + c_{vw} \text{ (correct it)}$ 

P(w) = v (set v as predecessor)

**Notice:** Technically speaking, this is not an algorithm, for the following reasons: 1) We did not specify how to pick e, 2) This procedure might not stop given some situations, e.g., if there is a cycle with minus total weight

**Notice:** This method can be really bad. Here is another example that could take  $O(2^n)$  to solve.



#### 28.4 Ford-Bellman Algorithm

#### Algorithm 21 Ford-Bellman Algorithm

```
Ensure: Shortest Paths from r to all other nodes in V
Require: A digraph with arc costs, starting node r
Initialize y and p
for i = 0; i < N; i + + do
for \forall e = (v, w) \in E do
if y_v + c_{vw} < y_w (incorrect) then
y_w \leftarrow y_v + c_{vw} (correct it)
P(w) = v \text{ (set } v \text{ as predecessor)}
for \forall e = (v, w) \in E do
if y_v + c_{vw} < y_w (incorrect) then
Return error, negative cycle
```

Notice: Only correct the node that comes from a node that has been corrected.

A usual representation of a digraph is to store all the arcs having tail v in a list  $L_v$  to scan v means the following:

• For  $(v, w) \in L_v$ , if (v, w) is incorrect, then correct (v, w)

For Bellman, will either terminate with shortest path from r to all  $v \in V \setminus r$  or it will terminate stating that there is a negative cycle. In O(mn)

In the algorithm if i = n and there exists a feasible potential, the problem has a negative cycle.

Suppose that the nodes of G can be ordered from left to right so that all arcs go from left to right. That is suppose there is an ordering  $v_1, v_2, ..., v_n \in V$  so that  $(v_i, v_j) \in V$  implies i < j. We call such an ordering **topological** sort. If we order E in the sequence that  $v_i v_j$  precedes  $v_k v_i$  if i < k based on topological order then ford algorithm will terminate in one pass.

- 28.5 SPFA Algorithm
- 28.6 Dijkstra Algorithm
- 28.7 A\* Algorithm

## 28.8 Floyd-Warshall Algorithm

If all weights/distances in the graph are nonnegative then we could use Dijkstra within starting nodes being any one of the vertices of the graph. This method will take  $O(n^3)$ 

If weight/distances are arbitrary and we would like to find shortest path between all pairs of vertices or detect a negative cycle we could use Bellman-Ford Algorithm with  $O(n^4)$ 

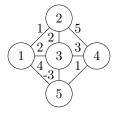
We would like an algorithm to find shortest path between any two pairs in a graph for arbitrary weights (determined, negative, cycles) in  $O(n^3)$ 

#### Algorithm 22 Dijkstra Algorithm

Ensure: Shortest Paths from r to all other nodes in V Require: A digraph with arc costs, starting node r Initialize y and p  $S \leftarrow V$  while  $S \neq \emptyset$  do

Choose  $v \in S$  with minimum  $y_v$   $S \leftarrow S \setminus v$  for  $\forall w, (v, w) \in E$  do

if  $y_v + c_{vw} < y_w$  (incorrect) then  $y_w \leftarrow y_v + c_{vw}$  (correct it) P(w) = v (set v as predecessor)



Let  $d_{ij}^k$  denote the length of the shortest path from i to j such that all intermediate vertices are contained in the set  $\{1, ..., k\}$ 

In this case  $d_{14}^5 = 5$ 

If the vertex k is not an intermediate vertex on p, then  $d_{ij}^k = d_{ij}^{k-1}$ , notice that  $d_{15}^4 = -1$ , node 4 is not intermediate, so  $d_{15}^3 = -1$ 

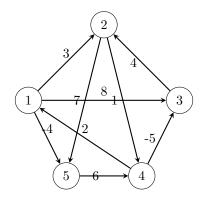
If the vertex k is an intermediate on p, then  $d_{ij}^k = d_{ik}^{k-1} + d_{kj}^{k-1}$ ,  $d_{14}^5 = 0$   $(p = 1 \rightarrow 3 \rightarrow 5 \rightarrow 4)$ , i.e.,  $d_{14}^5 = d_{15}^4 + d_{54}^4 = 0$ Therefore  $d_{ij}^k = \min\{d_{i-1}^{k-1}, d_{i-1}^{k-1} + d_{i-1}^{k-1}\}$ 

Therefore  $d_{ij}^k = \min\{d_{ij}^{k-1}, d_{ik}^{k-1} + d_{kj}^{k-1}\}$ Input: graph G = (V, E) with weight on edges Output: Shortest path between all pairs of vertices on existence of a negative cycle Step 1: Initialize

$$d_{ij}^{0} = \begin{cases} c_{ij} & \text{distance from } i \text{ to } j \text{ if } (i,j) \in E \\ 0 & \text{if } i = j \\ \infty & \text{if } (i,j) \notin E \end{cases}$$
(28.3)

Step: For k = 1 to n For i = 1 to n For j = 1 to n  $d_{ij}^k = \min\{d_{ij}^{k-1}, d_{ik}^{k-1} + d_{kj}^{k-1}\}$  Next j Next i Next k Between optimal matrix  $D^n$ 

$$D^{0} = \begin{bmatrix} 0 & 3 & 8 & \infty & -4 \\ \infty & 0 & \infty & 1 & 7 \\ \infty & 4 & 0 & \infty & \infty \\ 2 & \infty & -5 & 0 & \infty \\ \infty & \infty & \infty & 6 & 0 \end{bmatrix}$$
 (28.4)



$$\Pi^{0} = \begin{bmatrix}
1 & 1 & 1 & 1 \\
 & 2 & 2 \\
 & 3 & \\
 & 4 & 4 & \\
 & & 5 & 
\end{bmatrix}$$
(28.5)

$$D^{1} = \begin{bmatrix} 0 & 3 & 8 & \infty & -4 \\ \infty & 0 & \infty & 1 & 7 \\ \infty & 4 & 0 & \infty & \infty \\ 2 & \mathbf{5} & -5 & 0 & -\mathbf{2} \\ \infty & \infty & \infty & 6 & 0 \end{bmatrix}$$
 (28.6)

$$\Pi^{1} = \begin{bmatrix}
1 & 1 & 1 \\
 & 2 & 2 \\
3 & & & \\
4 & 1 & 4 & 1 \\
 & & 5
\end{bmatrix}$$
(28.7)

$$D^{2} = \begin{bmatrix} 0 & 3 & 8 & 4 & -4 \\ \infty & 0 & \infty & 1 & 7 \\ \infty & 4 & 0 & 5 & 11 \\ 2 & 5 & -5 & 0 & -2 \\ \infty & \infty & \infty & 6 & 0 \end{bmatrix}$$
 (28.8)

$$\Pi^{2} = \begin{bmatrix}
1 & 1 & \mathbf{2} & 1 \\
 & 2 & 2 \\
3 & \mathbf{2} & \mathbf{2} \\
4 & 1 & 4 & 1 \\
 & 5
\end{bmatrix}$$
(28.9)

$$D^{3} = \begin{bmatrix} 0 & 3 & 8 & 4 & -4 \\ \infty & 0 & \infty & 1 & 7 \\ \infty & 4 & 0 & 5 & 11 \\ 2 & -1 & -5 & 0 & -2 \\ \infty & \infty & \infty & 6 & 0 \end{bmatrix}$$
 (28.10)

$$\Pi^{3} = \begin{bmatrix}
1 & 1 & 2 & 1 \\
 & 2 & 2 \\
 & 3 & 2 & 2 \\
 & 4 & 3 & 4 & 1 \\
 & & 5
\end{bmatrix}$$
(28.11)

$$D^{4} = \begin{bmatrix} 0 & 3 & -1 & 4 & -4 \\ 3 & 0 & -4 & 1 & -1 \\ 7 & 4 & 0 & 5 & 3 \\ 2 & -1 & -5 & 0 & -2 \\ 8 & 5 & 1 & 6 & 0 \end{bmatrix}$$
(28.12)

$$\Pi^{4} = \begin{bmatrix}
1 & 4 & 2 & 1 \\
4 & 4 & 2 & 1 \\
4 & 3 & 2 & 1 \\
4 & 3 & 4 & 1 \\
4 & 3 & 4 & 5
\end{bmatrix}$$
(28.13)

$$D^{5} = \begin{bmatrix} 0 & \mathbf{1} & -\mathbf{3} & \mathbf{2} & -4 \\ 3 & 0 & -4 & 1 & -1 \\ 7 & 4 & 0 & 5 & 3 \\ 2 & -1 & -5 & 0 & -2 \\ 8 & 5 & 1 & 6 & 0 \end{bmatrix}$$
 (28.14)

$$\Pi^{5} = \begin{bmatrix}
3 & 4 & 5 & 1 \\
4 & 4 & 2 & 1 \\
4 & 3 & 2 & 1 \\
4 & 3 & 4 & 5 \\
4 & 3 & 4 & 5
\end{bmatrix}$$
(28.15)

Time complexity  $O(n^3)$ 

If during the previous processes, there exist an element of negative value in the diagonal, it means there exists negative cycle.

## 28.9 Johnson's Algorithm

## Maximum Flow Problem

#### 29.1 Basic Concept

Let D = (V, A) be a strict diagraph with distinguished vertices s and t. We call s the source and t the sink, let  $u = \{u_e : e \in A\}$  be a nonnegative integer-valued capacity function defined on the arcs of D. The maximum flow problem on (D, s, t, u) is the following Linear program.

$$\max \quad v \tag{29.1}$$

s.t. 
$$\sum_{h(e)=i} x_e - \sum_{t(e)=i} x_e = \begin{cases} -v, & \text{if } i = s \\ v, & \text{if } i = t \\ 0, & \text{otherwise} \end{cases}$$
 (29.2)

$$0 \le x_e \le u_e, \quad \forall e \in A \tag{29.3}$$

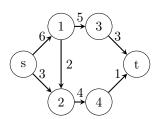
We think of  $x_e$  as being the flow on arc e. Constraint says that for  $i \neq s, t$  the flow into a vertex has to be equal to the flow out of vertex. That is, flow is conceded at vertex i for i = s and for i = t the net flow in the entire digraph must be equal to v. A  $\mathbf{x_e}$  that satisfied the above constraints is an (s, t)-flow of value v. If in addition it satisfies the bounding constraints, then it is a feasible (s, t)-flow. A feasible (s, t)-flow that has maximum v is optimal on maximum.

## 29.2 Solving Maximum Flow Problem in LP

**Theorem 29.1.** For  $S \subseteq V$  we define  $(S, \bar{S})$  to be a (s, t)-cut if  $s \in S$  and  $t \in \bar{S} = V - S$ , the capacity of the cut, denoted  $u(S, \bar{S})$  as  $\sum \{u_e : e \in \delta^-(S)\}$  where  $\delta^-(S) = \{e \in A : t(e) \in S \text{ and } h(e) \in \bar{S}\}$ 

**Example.** For the following graph:

Let 
$$S = \{1, 2, 3, s\}, \bar{S} = \{4, t\}$$

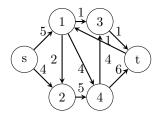


then  $\delta^-(S) = \{(2,4), (3,t)\} \Rightarrow u(S,\bar{S}) = 7$ 

**Definition 29.2.1.** If (S,S) has minimum capacity of all (s,t)-cuts, then it is called **minimum cut**.

**Definition 29.2.2.** Let  $\delta^{+}(S) = \delta^{-}(V - S)$ 

**Example.** Let  $S = \{s, 1, 2, 3\}$ ,  $\bar{S} = \{4, t\}$ ,  $u(S, \bar{S}) = u_{14} + u_{24} + u_{3t} = 10$ ,  $\delta^{-}(S) = \{(1, 4), (2, 4), (3, t)\}$ ,  $\delta^{+}S = \{(4, 3), (t, 1)\}$ 



**Lemma 29.2.** If x is a (s,t) flow of value v and  $(S,\bar{S})$  is a (s,t)-cut, then

$$v = \sum_{e \in \delta^{-}(S)} x_e - \sum_{e \in \delta^{+}(S)} x_e$$
 (29.4)

*Proof.* Summing the first set of constraints over the vertices of S,

$$\sum_{i \in S} \left( \sum_{h(e)=i} x_e - \sum_{t(e)=i} x_e \right) = -v \tag{29.5}$$

Now for an arc e with both ends in S,  $x_e$  will occur twice once with a positive and once with negative so they cancel and the above sum is reduced to

$$\sum_{e \in \delta^+(S)} x_e - \sum_{e \in \delta^-(S)} x_e = -v \tag{29.6}$$

Notice: Flow is the prime variable, capacity is the dual variable.

**Corollary 29.2.1.** If x is a feasible flow of value v, and  $(S, \bar{S})$  is an (s, t)-cut, then

$$v \le u(S, \bar{S}) \quad (Weak \ duality)$$
 (29.7)

**Definition 29.2.3.** Define an arc e to be saturated if  $x_e = u_e$ , and to be flowless if  $x_e = 0$ 

**Corollary 29.2.2.** Let x be a feasible flow and  $(S, \bar{S})$  be a (s, t)-cut, if  $\forall e \in \delta^-(S)$  is saturated, and  $\forall e \in \delta^+(S)$  is flowless, then x is a maximum flow and  $(S, \bar{S})$  is a minimum cut. (Strong duality)

*Proof.* If every arc of  $\delta^-(S)$  is saturated then

$$\sum_{e \in \delta^{-}(S)} x_e = \sum_{e \in \delta^{-}(S)} u_e \tag{29.8}$$

If every arc of  $\delta^+(S)$  is flowless then

$$\sum_{e \in \delta^+(S)} x_e = 0 \tag{29.9}$$

 $\Rightarrow x$  is as large as it can get when as  $u(S, \bar{S})$  is as small as it can get.

#### 29.3 Prime and Dual of Maximum Network Flow Problem

The LP of maximum flow can be modeled as following, WLOG, we let  $s = v_1 \in V, t = v_{|V|} \in V$ .

$$\max \quad f = \begin{bmatrix} 0 & 0 & \cdots & 0 & 1 \end{bmatrix} \begin{bmatrix} \mathbf{x} \\ f \end{bmatrix} \tag{29.10}$$

s.t. 
$$\begin{bmatrix} \mathbf{A} & \mathbf{F} \end{bmatrix} \begin{bmatrix} \mathbf{x} \\ f \end{bmatrix} = \mathbf{0}$$
 (29.11)

$$\mathbf{Ix} \le \mathbf{u} \tag{29.12}$$

In which **A** is the vertex-arc incident matrix and **F** is a column vector where the first row is -1, last row is 1 and all other rows are 0s, which is because we denote the first vertex as source s and the last vertex as the sink t. **u** is the column vector of upper bound of each arcs.

$$\mathbf{A} = \mathbf{A}_{|E| \times |V|} = [a_{ij}], \text{ where } a_{ij} = \begin{cases} 1, & \text{if } v_i = h(e_j) \\ -1, & \text{if } v_i = t(e_j) \\ 0, & \text{otherwise} \end{cases}$$
 (29.14)

$$\mathbf{F} = \begin{bmatrix} -1 & \cdots & 0 & \cdots & 1 \end{bmatrix}^{\mathsf{T}} \tag{29.15}$$

$$\mathbf{u} = \begin{bmatrix} u_1 & u_2 & \cdots & u_{|E|} \end{bmatrix}^{\top} \tag{29.16}$$

Then, we take the dual of LP

$$\min \quad \mathbf{u}\mathbf{w}_{\mathbf{E}} \tag{29.17}$$

s.t. 
$$\begin{bmatrix} \mathbf{w}_{\mathbf{V}} & \mathbf{w}_{\mathbf{E}} \end{bmatrix} \begin{bmatrix} \mathbf{A} \\ \mathbf{I} \end{bmatrix} \ge 0$$
 (29.18)

$$\begin{bmatrix} \mathbf{w}_{\mathbf{V}} & \mathbf{w}_{\mathbf{E}} \end{bmatrix} \begin{bmatrix} \mathbf{F} \\ \mathbf{0} \end{bmatrix} = 1 \tag{29.19}$$

$$\mathbf{w}_{\mathbf{V}}$$
 unrestricted (29.20)

$$\mathbf{w_E} \ge \mathbf{0} \tag{29.21}$$

In which  $\mathbf{w_V}$  is "whether or not" vertex v is in S where  $(S, \bar{S})$  represents a cut,  $\mathbf{w_E}$  is "whether or not" an arc in in  $\delta^+(S)$ .  $\mathbf{u}, \mathbf{E}, \mathbf{F}$  have the same meaning as in prime.

$$\mathbf{w}_{\mathbf{V}} = \begin{bmatrix} w_1 & w_2 & \cdots & w_{|V|} \end{bmatrix}^{\top} \tag{29.22}$$

$$\mathbf{w_E} = \begin{bmatrix} w_{|V|+1} & w_{|V|+2} & \cdots & w_{|V|+|E|} \end{bmatrix}^{\top}$$
 (29.23)

To make it more clear, it can be rewritten as following

$$\min \quad \sum_{e \in E} u_e w_e \tag{29.24}$$

s.t. 
$$w_i - w_j + w_{|V|+e} \ge 0, \forall e = (i, j) \in E$$
 (29.25)

$$-w_1 + w_{|V|} = 1 (29.26)$$

$$\mathbf{w}_{\mathbf{V}}$$
 unrestricted (29.27)

$$\mathbf{w_E} \ge \mathbf{0} \tag{29.28}$$

The meaning for the first set of constraint is to decide whether or not an arc is in  $\delta^+(S)$  of a  $(S, \bar{S})$ , which is decided by  $w_V$ . The  $w_1 - w_{|V|} = 1$ , which is the second set of constraint means the source  $s = v_1$  and the sink  $t = v_{|V|}$  has to be in S and  $\bar{S}$  respectively.

#### 29.4 Maximum Flow Minimum Cut Theorem

**Definition 29.4.1.** Let P be a path, (not necessarily a dipath), P is called **unsaturated** if every **forward** arc is unsaturated  $(x_e < u_e)$  and ever **reverse** arc has positive flow  $(x_e > 0)$ . If in addition P is an (s, t)-path, then P is called an **x-augmenting path** 

**Theorem 29.3.** A feasible flow x in a digraph D is maximum iff D has no augmenting paths.

*Proof.* (Prove by contradiction)

 $(\Rightarrow)$  Let x be a maximum flow of value v and suppose D has an augmenting path. Define in P (augmenting path):

$$D_1 = \min\{u_e - x_e : e \text{ forward in } P\}$$
(29.29)

$$D_2 = \min\{x_e : e \text{ backward in } P\} \tag{29.30}$$

$$D = \min\{D_1, D_2\} \tag{29.31}$$

Since P is augmenting, then D > 0, let

$$\hat{x_e} = \begin{cases} x_e + D & \text{If } e \text{ is forward in } P \\ x_e - D & \text{If } e \text{ is backward in } P \\ x_e & otherwise \end{cases}$$
 (29.32)

It is easy to see that  $\hat{x}$  is feasible flow and that the value is V+D, a contradiction.

( $\Leftarrow$ ) Suppose D admits no x-augmenting path, Let S be the set of vertices reachable from s by x-unsaturated path clearly  $s \in S$  and  $t \notin S$  (because otherwise there would be an augmenting path). Thus,  $(S, \bar{S})$  is a (s, t)-cut.

Let  $e \in \delta^-(S)$  then e must be saturated. For otherwise we could add the h(e) to S

Let  $e \in \delta^+(S)$  then e must be flow less. For otherwise we could add the t(e) to S.

According to previous corollary, that x is maximum.

**Theorem 29.4.** (Max-flow = Minimum-cut) For any digraph, the value of a maximum (s,t)-flow is equal to the capacity of a minimum (s,t)-cut

#### 29.5 Ford-Fulkerson Method

Finding augmenting paths is the key of max-flow algorithm, we need to describe two functions, labeling and scanning a vertex.

A vertex is first labeled if we can find x-unsaturated path from s, i.e., (s, v)-unsaturated path.

The vertex v is scanned after we attempted to extend the x-unsaturated path.

This algorithm is incomplete/incorrect, needs to be fixed

#### Algorithm 23 Labeling algorithm

**Ensure:** Max-flow x with value v

**Require:** Digraph with source s and sink t, a capacity function u and a feasible flow (could be  $x_e = 0$ )

Initialize,  $v \leftarrow x$ 

Designate all vertices as unlabeled and unscanned

Label s

while There exists vertex unlabeled or unscanned do

Let i be such a vertex, for each arc e with  $t(e) = i, x_e < u_e$  and h(e) unlabeled, label h(e)

For each arc e with  $h(e) = i, x_e > 0$  and t(e) unlabeled, label t(e), designate i as scanned.

If t is not label

x is the maximum.

#### Algorithm 24 Ford-Fulkerson algorithm

**Ensure:** Max-flow x with value v

**Require:** Digraph with source s and sink t, a capacity function u and a feasible flow (could be  $x_e = 0$ )

Initialize,  $v \leftarrow x$ 

Designate all vertices as unlabeled and unscanned

Label s

while There exists vertex unlabeled or unscanned do

Let i be such a vertex, for each arc e with  $t(e) = i, x_e < u_e$  and h(e) unlabeled, label h(e)

For each arc e with  $h(e) = i, x_e > 0$  and t(e) unlabeled, label t(e), designate i as scanned.

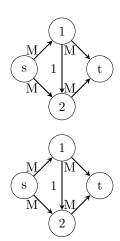
If t is not label

x is the maximum.

Labeling algorithm can be exponential, the following is an example

### 29.6 Polynomial Algorithm for max flow

Let (D, s, t, u) be a max flow problem and let x be a feasible flow for D, the **x-layers** of D are define be the following algorithm



Layer algorithm (Dinic 1977) Input: A network (D, s, t, u) and a feasible flow x Output: The **x-layers**  $V_0, V_1, ..., V_l$  where  $V_i \cap V_j = \emptyset \forall i \neq j$ 

Step 1: Set  $V_0 = \{s\}, i \leftarrow 0$  and I(x) = 0 Step 2: Let I(x) = 0 be the set of vertices I(x) = 0 such that there is an arc I(x) = 0 step 1: Set I(x) = 0 step 2: Let I(x) = 0 step 3: Let I(x) = 0 step 4: Let I(x) = 0 step 4: Let I(x) = 0 step 5: Let I(x) = 0 step 6: Let I(x) =

- $t(e) \in V_i, h(e) = w, x_e < u_e \text{ or }$
- $h(e) \in V_j, t(e) = w, x_e > 0$

Step 3: If  $t \in R$ , set  $V_{i+1} = \{t\}$ , l(t) = i+1 and stop. Set  $V_{i+1} \leftarrow R \setminus \bigcup_{0 \le j \le i} V_j$ ,  $l \leftarrow i+1$ , l(x) = i, goto Step 2. If  $V_{i+1} = \emptyset$ , set l(x) = i and Stop.

**Example.** For the following graph

Second iteration 
$$(29.33)$$

$$V_0 = \{s\}, i = 0, l(x) = 0 \tag{29.34}$$

$$R = \{1, 2\} \tag{29.35}$$

$$V_1 \leftarrow \{1, 2\}, i = 1, l(x) = 1$$
 (29.36)

$$R = \{3, 4, 5\} \tag{29.37}$$

$$V_2 \leftarrow \{3, 4\}, i = 2, l(x) = 2$$
 (29.38)

$$R = \{1, 5, 6, 3\} \tag{29.39}$$

$$V_3 \leftarrow \{5, 6\}, i = 3, l(x) = 3$$
 (29.40)

$$R = \{4, t\} \tag{29.41}$$

$$V_4 = \{t\} \tag{29.42}$$

$$A_1 = \{(s,1), (s,2)\} \tag{29.43}$$

$$A_2 = \{(1,3), (2,4)\}\tag{29.44}$$

$$A_3 = \{(3,5), (4,6)\} \tag{29.45}$$

$$A_4 = \{(5, t), (6, t)\}\tag{29.46}$$

The layer network  $D_x$  is defined by  $V(D_x) = V_0 \cup V_1 \cup V_2 \cdots \cup V_{l(x)}$ 

Suppose we have computed the layers of D and  $t \in V_l(x)$ , the last layer (last layer I am goin to  $V_e$ ) For each  $i, 1 \le i \le l$ , define a set of arcs  $A_i$  and a function  $\hat{u}$  on  $A_i$  as following. For each  $e \in A(D)$ 

- If  $t(e) \in V_{i-1}$ ,  $h(e) \in V_i$  and  $x_e < u_e$  then add arc e to  $A_i$  and define  $\hat{u}_e = u_e x_e$
- If  $h(e) \leftarrow V_{i-1}, t(e) \in V_i$  and  $x_e > 0$  then add arc e' = (h(e), t(e)) to  $A_i$  with  $\hat{u}_e x_e$

Let  $\hat{u}$  be the capacity function on  $D_x$  and let the source and sink of  $D_x$  be s and t

We can think of  $D_x$  as being make of arc shortest (in terms of numbers of arcs) x-augmenting paths.

A feasible flow in a network is said to be maximal (does not means maximum) if every (s,t)-directed path contains at least one saturated arc.

For layered algorithm  $V_0, V_1, ..., V_L$ 

Arcs:

- If  $t(e) \in V_{i-1}$ ,  $h(e) \in V_i$  and  $x_e < u_e$ , then add e to  $A_i$  with  $\hat{u}_e = u_e x_e$
- If  $h(e) \in V_{i-1}$ ,  $t(e) \in V_i$  and  $x_e > 0$ , then add arc e' = (h(e), t(e)) to  $A_i$  and define  $\hat{u_e} = x_e$

Maximal Flow: If every directed (s, t)-path has at least one saturated arc.

Computing maximal flow is easier than computing maximum flow, since we never need to consider canceling flows on reverse arcs,

Let  $\hat{x}$  be a maximal flow on the layered network  $D_x$ , we can define new flows in D(x') by

$$x'_e = x_e + \hat{x_e}, \quad \text{If } t(e) \in V_{i-1}, h(e) \in V_i$$

$$x'_e = x_e - \hat{x_e}, \quad \text{If } h(e) \in V_{i-1}, t(e) \in V_i$$
(29.47)

$$x'_e = x_e - \hat{x_e}, \quad \text{If } h(e) \in V_{i-1}, t(e) \in V_i$$
 (29.48)

#### 29.7Dinic Algorithm

Input: A layered network  $(D_x, s, t, \hat{u})$  and a feasible flow x Output: A maximal flow  $\hat{x}$  from  $D_x$ Step 1: Set  $H \leftarrow D_x$  and  $i \leftarrow S$  Step 2: If there is no arc e with t(e) = i, goto Step 4, otherwise let e be such an arc Step 3: Set  $T(h(e)) \leftarrow i$  and  $i \leftarrow h(e)$ , if i = t goto Step 5, otherwise goto Step 2. Step 4: If i=s, Stop, Otherwise delete i and all incident arcs with H, set  $i\leftarrow T(i)$  and goto Step 2 Step 5: Construct the directed path,  $s = i_0 e_1 i_1 e_2 \dots e_k i_k = t$  where  $i_{j-1} = T(i_j), 1 \le j \le k$ . Set  $D = \min\{\hat{u_{e_j}} - \hat{x_{e_j}} : i \le j \le k\}$ , set  $\hat{x_{e_j}} \leftarrow \hat{x_{e_j}} + D, i \leq j \leq k$ . Delete from H all saturated arcs on this path, set  $i \leftarrow 1$  and goto Step 2.

#### Algorithm 25 Dinic Algorithm

**Ensure:** A maximal flow  $\hat{x}$  from  $D_x$ 

**Require:** A layered network  $(D_x, s, t, \hat{u})$  and a feasible flow x

Initialize  $H \leftarrow D_x$  and  $i \leftarrow S$ 

**Theorem 29.5.** Dinic algorithm runs in  $O(|E||V|^2)$ 

*Proof.* Step 1 is O(|E||V|) Step 2 runs Step 1 for O(|V|) times

## Minimum Cost Flow Problem

#### 30.1 Transshipment Problem

Transshipment Problem (D, b, w) is a linear program of the form

$$\min \quad wx \tag{30.1}$$

$$s.t. \quad Nx = b \tag{30.2}$$

$$x \ge 0 \tag{30.3}$$

Where N is a vertex-arc incident matrix. For a feasible solution to LP to exist, the sum of all bs must be zero. Since the summation of rows of N is zero. The interpretation of the LP is as follows.

The variables are defined on the edges of the digraph and that  $x_e$  denote the amount of flow of some commodity from the tail of e to the head of e

Each constraints

$$\sum_{h(e)=i} x_e - \sum_{t(e)=i} x_e = b_i \tag{30.4}$$

represents consequential of flow of all edges into k vertex that have a demand of  $b_i > 0$ , or a supply of  $b_i < 0$ . If  $b_i = 0$  we call that vertex a transshipment vertex.

## 30.2 Network Simplex Method

**Lemma 30.1.** Let  $C_1$  and  $C_2$  be distinct cycles in a graph G and let  $e \in C_1 \cup C_2$ . Then  $(C_1 \cup C_2) \setminus e$  contains a cycle.

*Proof.* Case 1:  $C_1 \cap C_2 = \emptyset$ . Trivia.

Case 2:  $C_1 \cap C_2 \neq \emptyset$ . Let  $e \in C_2$  and  $f = uv \in C_1 \setminus C_2$ . Starting at v traverse  $C_1$  in the direction away from u until the first vertex of  $C_2$ , say x. Denote the (v, x)-path as P. Starting at u traverse  $C_1$  in the direction away from v until the first vertex of  $C_2$ , say y. Denote the (u, y)-path as Q.  $C_2$  is a cycle, there are two (x, y)-path in  $C_2$ . Denote the (x, y)-path without e as R. Then  $vPxRyQ^{-1}uf$  is a cycle.

**Theorem 30.2.** Let T be a spanning tree of G. And let  $e \in E \setminus T$  then T + e contains a unique cycle C and for any edge  $f \in C$ , T + e - f is a spanning tree of G

Let (D, b, w) be a transshipment problem. A feasible solution x is a **feasible tree solution** if there is a spanning tree T such that  $||x|| = \{e \in A, x_e \neq 0\} \subseteq T$ .

The strategy of network simplex algorithm is to generate negative cycles, if negative cycle exists, it means the solution can be improved.

For any tree T of D and for  $e \in A \setminus T$ , it follows from above theorem that T + e contains a unique cycle. Denote that cycle C(T, e) and orient it in the direction of e, define

$$w(T, e) = \sum \{w_e : e \text{ forward in } C(T, e)\}$$
$$-\sum \{w_e : e \text{ reverse in } C(T, e)\}$$
(30.5)

We think of w(T, e) as the weight of C(T, e).

## 30.2.1 Network Simplex Method

```
Algorithm 26 Network Simplex Method Algorithm

Ensure: An optimal solution or the conclusion that (D,b,w) is unbounded

Require: A transshipment problem (D,b,w) and a feasible tree solution x containing to a spanning tree T

while \exists e \in A \setminus T, w(T,e) < 0 do

let e \in A \setminus T be such that w(T,e) < 0.

if C(T,e) has no reverse arcs then

Return unboundedness

else

Set \theta = \min\{x_f : f \text{ reverse in } C(T,e)\} and set f = \{f \in C(T,e) : f \text{ reverse in } C(T,e), x_f = \theta\}

if f forward in C(T,e) then

x_f \leftarrow x_f + \theta

else

x_f \leftarrow x_f - \theta

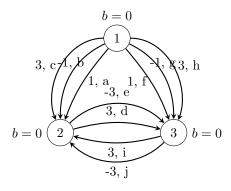
Let f \in F and T \leftarrow T + e - f

Return x as optimal
```

#### 30.2.2 Example for cycling

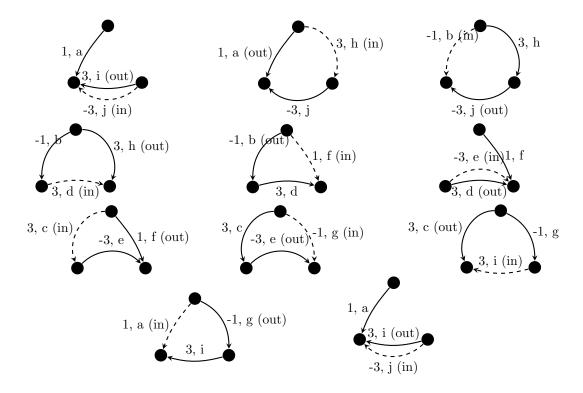
**Notice:** Similar to Simplex Method in LP, even though in worst case may be inefficient. In most cases it is simple and empirically efficient. Also, similarly, there will be cycling problems.

The following is an example of cycling



```
Then for the following steps we can detect cycling: w(T,j) = w_j - w_i = -3 - 3 = -6, therefore j in entering basis, i is leaving basis. w(T,h) = w_h + w_j - w_a = 3 - 3 - 1 = -1, therefore h is entering basis, a is leaving basis. w(T,b) = w_b - w_j - w_h = -1 + 3 - 3 = -1, therefore b is entering basis, j is leaving basis. w(T,d) = w_d - w_h + w_b = 3 - 3 - 1 = -1, therefore d is entering basis, h is leaving basis. w(T,f) = w_f - w_d - w_b = 1 - 3 + 1 = -1, therefore f is entering basis, h is leaving basis. w(T,e) = w_e - w_d = -3 - 3 = -6, therefore e is entering basis, e is leaving basis. w(T,c) = w_c + w_e - w_f = 3 - 3 - 1 = -1, therefore e is entering basis, e is leaving basis. w(T,g) = w_g - w_e - w_c = -1 + 3 - 3 = -1, therefore e is entering basis, e is leaving basis. e is leaving basis. e is entering basis, e is leaving basis. e is entering basis, e is leaving basis. e is entering basis, e is leaving basis.
```

The last graph is the same as the first graph, i.e., cycling detected.



#### 30.2.3 Cycling prevention

To Avoid cycling we will introduce the Modified Network Simplex Method. Let T be a **rooted** spanning tree. Let f be an arc in T, we say f is **away** from the root r if t(f) is the component of T - f. Otherwise we say f is **towards** r.

Let x be a feasible tree solution associated with T, then we say T is a **strong feasible tree** if for every arc  $f \in T$  with  $x_f = 0$  then f is away from  $r \in T$ . Modification to NSM:

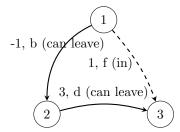
- The algorithm is initialed with a strong feasible tree.
- f in pivot phase is chosen to be the first reverse arc of C(T, e) having  $x_f = \theta$ . By "first", we mean the first arc encountered in traversing C(T, e) in the direction of e, starting at the vertex i of C(T, e) that minimizes the number of arcs in the unique (r, i)-path in T.

**Notice:** In the second rule above, r could also be in the cycle, in that case, i is r.

Continue the previous example. Now should how we can avoid cycling:

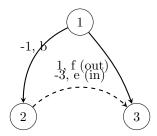
The first few (four) steps are the same as previous example, starting from

 $w(T,f) = w_f - w_d - w_b = 1 - 3 + 1 = -1$ . f is entering basis, both b and d can leave the basis, according to



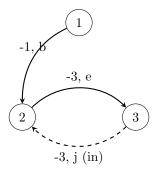
the modified pivot rule, we choose the "first" arc encountered in traversing C(T, e), which is d to leave the basis, instead of b.

 $w(T,e) = w_e - w_f + w_b = -5$ , e is entering basis, f is leaving basis. Now the only arc to enter basis and maintain



negative w is j.

 $w(T,j) = w_j + w_e = -6$ , but in C(T,j) there is no reversing arc, therefore we detect unboundedness.

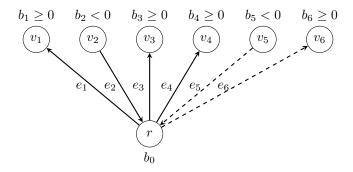


#### 30.2.4 Finding Initial Strong Feasible Tree

Pick a vertex in D to be root r. The tree T has an arc e with the t(e) = r and h(e) = v. For each  $v \in V \setminus r$  with  $b_v \ge 0$  and has an arc e with h(e) = r and t(e) = v for each  $v \in V \setminus r$  for which  $b_v < 0$ . Wherever possible the arcs of T are chosen from A, where an appropriate arc doesn't exist. We create an **artificial arc** and give its weight  $|V|(\max\{w_e : e \in A\}) + 1$ . This is similar to Big-M method and if optimal solution contains artificial arcs ongoing arc problem is infeasible.

Here is an example after adding artificial arcs:

Where  $e_5$  and  $e_6$  are artificial arcs, the weight of those arcs are  $|V|(\max\{w_e:e\in\mathcal{A}\})+1$ . And the above tree is



a basic feasible solution.

We need to prove that such artificial arc has sufficiently large weight to guarantee

- It will leave the basis, and
- It will not enter the basis again (for this, just delete the artificial arc after it leaves the basis, then it will never enter the basis again)

*Proof.* Now prove that such arcs will always leave the basis. Before the prove we give some notation.

- Define set E as the set of arcs which is not artificial arc, in the above example,  $E = \{e_1, e_2, e_3, e_4\}$ .
- Define set A as the set of arcs which are artificial arcs, in the above example,  $A = \{e_5, e_6\}$ . Noticed that  $E \cap A = \emptyset$ .

- Define set M as the vertices in the spanning tree that is reachable from r by E, in the above example,  $M = \{v_1, v_2, v_3, v_4\}.$
- Define  $M' = (V \setminus r) \setminus M$  in the tree that can only be reached from r by A, i.e., artificial arcs, in the above example,  $M' = \{v_5, v_6\}$ .

Then the initial basic feasible solution is a graph

$$G_0 = \langle M \cup M' \cup \{r\}, E \cup A \rangle$$
 (30.6)

Denote the origin graph

$$G = \langle V, \mathcal{A} \rangle \tag{30.7}$$

Notice that with the artificial arcs,  $G_0$  is not a subgraph of G.

Let  $(M \cup \{r\}, M')$  be a cut in the origin graph G. For the vertices in M', one of the following cases will happen:

- case 1:  $\sum_{v \in M'} b_v \ge 0$
- case 2:  $\sum_{v \in M'} b_v < 0$

For case 1, we claim that at least one of the vertices  $v_{M'} \in M'$  with  $b_{v_{M'}} \geq 0$  linked by an arc, say f, such that  $h(f) = v_{M'}$  and  $t(f) = v_M \in M$ . Otherwise the balance of flow cannot hold in the origin graph G. Furthermore, denote the artificial arc from r to  $v_{M'}$  by  $e_{rv_{M'}}$ .

Notice that for  $v_M$  there is not necessarily be an arc between r and  $v_M$ , but there must exists an  $(r, v_M)$ -path denoted by P, for M is the set of vertices that reachable from r by arcs in E.

Take that arc f as entering arc to the basis. Then

$$C(T,f) = re_{rv_M} v_{M'} f v_M P r \tag{30.8}$$

For

$$w(T,f) = w_f - w_{e_{rv_{M'}}} + \sum_{e \in P} d_e w_e$$
(30.9)

where  $d_e = 1$  if  $w_e$  is forward in P and  $d_e = -1$  otherwise.

Now that  $w_{e_{rv_{M'}}} = |V|(\max\{w_e : e \in A\}) + 1$ , it guarantees that

$$w(T,f) = w_f + \sum_{e \in P} d_e w_e - w_{e_{rv_{M'}}}$$
(30.10)

$$\leq w_f + \sum_{e \in P} w_e - w_{e_{rv_{M'}}} \tag{30.11}$$

$$\leq \sum_{e \in A} w_e - w_{e_{rv_{M'}}}$$
(30.12)

$$\leq |V|(\max\{w_e : e \in A\}) - w_{e_{rv}}$$
 (30.13)

$$\leq -1 < 0 \tag{30.14}$$

So f can enter the basis, and the artificial variable  $e_{rv_{M'}}$  will leave the basis, for it is the most violated reverse arc in the C(T, f). When we put f into the basis, update  $G_0$ , such that  $M \leftarrow M \cup \{v_{M'}\}$  and  $M' \leftarrow M' \setminus \{v_{M'}\}$ . For case 2, it is similar. At least one of the vertices  $v_{M'} \in M'$  with  $b_{v_{M'}} < 0$  linked by an arc, say f', such that  $t(f') = v_{M'}$  and  $h(f') = v_M \in M$ . Otherwise the balance of flow cannot hold in the origin graph G. Furthermore, denote the artificial arc from  $v_{M'}$  to r by  $e_{v_{M'}r}$ .

Similarly we can find a cycle  $C(T, f') = rP'v_M f'v_{M'} e_{v_{M'}r} r$ .  $w(T, f') = w_{f'} - w_{e_{rv_{M'}}} + \sum_{e \in P'} d_e w_e$ , where  $d_e = 1$  if  $w_e$  is forward in P' and  $d_e = -1$ . We can prove  $w(T, f') \leq -1 < 0$ . That that f' as entering arc to the basis, similarly move  $v_{M'}$  form set M' to M.

The above case can be dealt with iteratively until set M' become  $\emptyset$ , at which stage there is no artificial arc in the basic feasible solution. Which means all the artificial variable can leave the basis.

**Notice:** This algorithm can be really bad, its mimic of Simplex Method of LP, which means we can run into exponential operations

#### 30.3 Transshipment Problem and Circulation Problem

**Definition 30.3.1.** The minimum weight circulation problem is defined as follows:

$$\min \quad wx \tag{30.15}$$

s.t. 
$$Nx = 0$$
 (30.16)

$$l \le x \le u \tag{30.17}$$

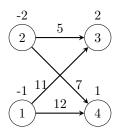
It turns out that the circulation problem is equivalent with transshipment problem.

We will show how to transform any transshipment into circulation.

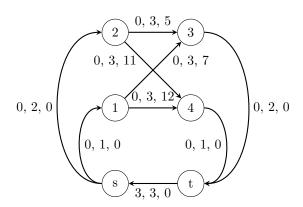
Let (D, b, w) be a transshipment problem and define two new vertices s and t.

- For each supply vertex x add the arc (s,x) to D with  $l_x = 0, u_x = -b_x, w_x = 0$ .
- Similarly, for each demand vertex x, add the arc (x,t) to D with  $l_x=0, u_x=b_x, w_x=0$ .
- Finally, add an arc (t, s) having  $w_{ts} = 0, l_{ts} = u_{ts} = \sum \{b_x : \forall x, x \text{ is demand vertex}\}.$
- Each original arc is given a  $l_x = 0, u_x = \sum \{b_x : \forall x, x \text{ is a demand vertex}\}, w_x \text{ remains unchanged.}$

The following is a graph for transshipment problem.



After the above procedures, it is now transformed into a circulation problem.



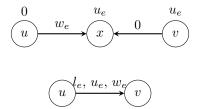
Then, we will show how to transform any circulation problem into transshipment problem.

If the lower bound of the arc is zero, i.e.,  $l_e = 0$ , then for each such arc (u, v), introduce a vertex in between u and v, replace the arc e = (u, v) by  $e_1 = (u, x)$  and  $e_2 = (x, v)$ . Both arcs are uncapacitated. Let  $w_1(u, x) = w(u, v)$  and  $w_2(x, v) = 0$  be the new weights for the arcs. Let  $u_e$  be the demands of newly added vertex x and add  $u_e$  to the supplies of vertex v (in v the supplies is the summation from all arcs that go to v).

$$u$$
  $0, u_e, w_e$   $v$ 

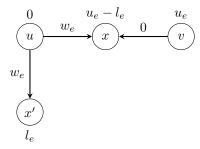
Will be transform into

If the lower bound of the arc is not zero, i.e.,  $l_e \neq 0$ , then for each such arc (u, v), introduce two new vertices, one vertex is in between u and v, similar to the previous case, the difference is the demands of this new vertex is



 $u_e - l_e$ , the others stays the same. Then add the other vertex, this vertex, denoted as x' is added along with an arc between u and x', the weight of this arc is w(u, v), the demands on this new vertex is  $l_e$ .

Will be transform into



Perform the above procedures to all the arcs in a circulation problem. In O(E) (polynomially times) transformation, such problem can be transformed into transshipment problem.

### 30.4 Out-of-Kilter algorithm

This algorithm is a Primal-dual method and is applied to the minimum weight circulation problem.

For LP optimality conditions we need primal feasibility, dual feasibility and complementary slackness, i.e., KKT conditions. Primal and dual feasibility are obvious so we need to show complementary slackness through following theorem.

**Theorem 30.3.** Let x be a feasible circulation flow for (D, l, u, w). And suppose there exists a real value vector  $\{y_i : i \in V\}$  which we called **vertex-numbers**. For all edges  $e \in A$ 

$$y_{h(e)} - y_{t(e)} > w_e \text{ implies } x_e = u_e$$
 (30.18)

$$y_{h(e)} - y_{t(e)} < w_e \text{ implies } x_e = l_e$$
 (30.19)

Then x is optimal to the circulation problem.

*Proof.* For each  $e \in A$  define

$$\gamma_e = \max\{y_{h(e)} - y_{t(e)} - w_e, 0\}$$
(30.20)

$$\mu_e = \max\{w_e - y_{h(e)} + y_{t(e)}, 0\}$$
(30.21)

Then

$$\gamma_e - \mu_e = y_{h(e)} - y_{t(e)} - w_e \tag{30.22}$$

Furthermore

$$\sum_{e \in A} (\mu_e l_e - \gamma_e u_e) \tag{30.23}$$

$$= \sum_{e \in A} (\mu_e l_e - \gamma_e u_e) + \sum_{i \in V} y_i (\sum_{h(e)=i} x_e - \sum_{t(e)=i} x_e)$$
(30.24)

$$= \sum_{e \in A} (\mu_e l_e - \gamma_e u_e + x_e (y_{h(e)} - y_{t(e)}))$$
(30.25)

$$= \sum_{e \in A} (\mu_e l_e - \gamma_e u_e + x_e (\gamma_e - \mu_e + w_e))$$
 (30.26)

$$= \sum_{e \in A} (\gamma_e(x_e - u_e) + \mu_e(l_e - x_e) + x_e w_e)$$
(30.27)

$$\leq \sum_{e \in A} x_e w_e \tag{30.28}$$

The last inequality will be satisfied as equality iff the first two hold.

The following is the formulation of circulation problem

$$(P) \quad \min \quad wx \tag{30.29}$$

s.t. 
$$Nx = 0$$
  $y$  (30.30)

$$x \ge l \quad z^l \tag{30.31}$$

$$-x \le -u \quad z^u \tag{30.32}$$

(D) 
$$\max lz^l - uz^u \tag{30.33}$$

(s.t.) 
$$yN^{-1} + z^l - z^u \le w$$
 (30.34)

$$y free$$
 (30.35)

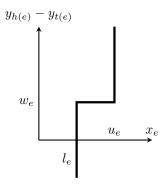
$$z^l, z^u \ge 0 \tag{30.36}$$

(CS) 
$$y_{h(e)} - y_{t(e)} > w_e \Rightarrow x_e = u_e$$
 (30.37)

$$y_{h(e)} - y_{t(e)} < w_e \Rightarrow x_e = l_e$$
 (30.38)

There is an alternative way of circulation optimality for a circulation problem. We define a **kilter-diagram** as follows.

For every edge construct the following:

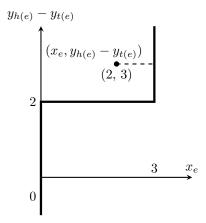


For each point  $(x_e, y_{h(e)} - y_{t(e)})$  we define a **kilter-number**  $k_e$ , be the minimum positive distance change in  $x_e$  required to put in on the kilter line.

**Example.** For edge  $e: w_e = 2, l_e = 0, u_e = 3$ , assume  $x_e = 2, y_{h(e)} - y_{t(e)} = 3$ , then  $k_e = 1$ 

**Lemma 30.4.** If for every circulation x and vertex number y we have  $\sum_{e \in A} k_e = 0$ , then x is optimal.

*Proof.* Since  $k_e$  is a nonnegative number, then the only way that  $\sum_{e \in A} k_e = 0$  is  $k_e = 0, \forall e \in A$ , which means  $\forall e \in A, l_e \leq x_e \leq u_e$ . Furthermore, the complementary slackness are satisfied.



General idea of algorithm follows. Suppose we are given a circulation x and vertex-numbers y (we do not require feasibility). Usually we pick x = 0, y = 0. If every edge is in kilter-line then we are optimal.

Otherwise there is at least one edge  $e^*$  that is out-of kilter. The algorithm consist of two phases, one called **flow-change** phase (horizontally), then other **number-change** phase (vertically).

In the flow-change phase, we want to find a new circulation for an out-of-kilter edge  $e^*$  say  $\hat{e}$  such that we reduce the kilter number  $k_{e^*}$ , without increasing any other kilter number for other edges.

To do this, denote the edges of  $e^*$  to be s and t, where such that  $k_{e^*}$  will be decreased by increasing the flow from s to t on  $e^*$ .

If  $e^* = (s, t)$  this will accomplished by increasing  $x_{e^*}$  and if e = (t, s) it is accomplished by decreasing  $x_{e^*}$ . To do this we look for an (s, t)-path p of the following edges.

- If e is forward in p, then increasing  $x_e$  does not increase  $k_e$  and
- If e is reversed in p, then decreasing  $x_e$  dose not increase  $k_e$

In terms of kilter diagram, an arc satisfies "forward" if it is forward and in left side of kilter line, and it satisfies "reversed" if it is reverse and in right side of kilter line.

Suppose we can not find such a path. From s to t, let x be the vertices that can decrease by an augmenting path. Then either we can change the vertex numbers y so that  $\sum_{e \in A} k_e$  does not increase but x does, or we can show that problem is infeasible.

INPUT a minimum circulation problem (D, l, u, w) a circulation x and vertex-numbers y

OUTPUT conclusion that (D, l, u, w) is infeasible or an minimum weighted flow.

Step 1: If every arc is in kilter  $(k_e = 0, \forall e \in A)$ . Stop with x is optimal. Otherwise let  $e^*$  be an out-of-kilter arc. If increasing  $x_{e^*}$  decreases  $k_{e^*}$  set  $s = h(e^*)$  and  $t(e^*)$  otherwise set  $s = t(e^*)$  and  $t = h(e^*)$ 

Step 2: If there exists an (s,t) augmenting path p then goto Step 3, otherwise goto Step 4.

STEP 3: Set  $y_e = y_{h(e)} - y_{t(e)}, e \in A$  Set  $\Delta_1 = \min\{u_e - x_e : e \text{ is forward and } y_e \geq w_e\}$  Set  $\Delta_2 = \min\{l_e - x_e : e \text{ is forward and } y_e < w_e\}$  Set  $\Delta_3 = \min\{x_e - l_e : e \text{ is reverse and } y_e \leq w_e\}$  Set  $\Delta_4 = \min\{x_e - u_e : e \text{ is reverse and } y_e > w_e\}$   $\Delta = \min\{\Delta_i, i = 1, 2, 3, 4\}$ 

Increase  $x_e$  by  $\Delta$  on each forward arc in p, decrease  $x_e$  by  $\Delta$  on each reverse arc in p.

If  $e^* = (s, t)$  decrease  $x_{e^*}$  by  $\Delta$ , otherwise increase  $x_{e^*}$  by  $\Delta$ 

If  $k_{e^*} > 0$  goto Step 2. otherwise goto Step 1.

Step 4: Let X be the set of vertices reachable from s by augmenting paths, then  $t \notin X$ , if every arc e with  $h(e) \in X$  has  $x_e \leq l_e$  and every arc e with  $t(e) \in X$  has  $x_e \geq u_e$ , and at least one of the above inequality is strict, then Stop with problem infeasible

Otherwise set  $\delta_1 = \min\{w_e - y_e : t(e) \in X, y_e < w_e, x_e \le u_e \ne l_e\}$   $\delta_2 = \min\{y_e - w_e : h(e) \in X, y_e > w_e, x_e \ge u_e \ne l_e\}$   $\delta = \min\{\delta_1, \delta_2\}$ 

Set  $y_i = y_i + \delta$  for  $i \notin X$ 

If  $k_{e^*} > 0$ , goto Step 2, otherwise goto Step 1.

Out-of-kilter takes O(|E||V|K) where  $K = \sum_{e \in A} k_e$ . However, there is an algorithm called **scaling algorithm** that uses out-of-kilter as subroutine that runs in  $O(R|E|^2|V|)$  where  $R = \lceil \max\{\log_2 u_e : e \in A\} \rceil$ 

## 30.5 Complexity of Different Minimum Weighted Flow Algorithms

Let arc capacities between 1 and U, costs between -C and C

Year	Discoverer	Method	Big O
1951	Dantzig	Network Simplex	$O(E^2V^2U)$
1960	Minty, Fulkerson	Out-of-Kilter	O(EVU)
1958	$_{ m Jewell}$	Successive Shortest Path	O(EVU)
1962	Ford-Fulkerson	Primal Dual	$O(EV^2U)$
1967	Klein	Cycle Canceling	$O(E^2CU)$
1972	Edmonds-Karp, Dinitz	Capacity Scaling	$O(E^2 \log U)$
1973	Dinitz-Gabow	Improved Capacity Scaling	$O(EV \log U)$
1980	Rock, Bland-Jensen	Cost Scaling	$O(EV^2 \log C)$
1985	Tardos	$\epsilon$ -optimality	poly(E, V)
1988	Orlin	Enhanced Capacity Scaling	$O(E^2)$

## Social Network Analysis

# Part V

## Heuristic and Metaheuristic Optimization

## Search Algorithm

Special Topic: Vehicle Routing Problem

### The Traveling Salesman Problem

#### 33.1 **Formulations**

In this section, we are going to compare between different formulations of TSP. Generally speaking, let G = (V, A)be a graph where V is a set of n vertices, and A is a set of arcs (or edges). Let  $C = c_{ij}$  be a cost (distance) matrix associated with A. The TSP consists of determining a minimum cost (distance) Hamiltonian circle (or cycle) that visits each vertex once and only once. If for all  $i, j \in V, c_{ij} = c_{ji}$ , then the TSP is symmetrical, otherwise is asymmetrical.

Define the decision variable  $x_{ij}$  as the following

$$x_{ij} = \begin{cases} 1, & \text{if goes from } i \text{ to } j \\ 0, & \text{otherwise} \end{cases}, \quad (i,j) \in A$$
 (33.1)

The objective function will be

$$\min \sum_{(i,j)\in A} c_{ij} x_{ij} \tag{33.2}$$

#### 33.1.1Dantzig-Fulkerson-Johnson (DFJ) Formulation

The first famous formulations for TSP is the Dantzig-Fulkerson-Johnson (DFJ) formulation:

$$\sum_{j \in V, (i,j) \in A} x_{ij} = 1, \quad \forall i \in V$$

$$\sum_{i \in V, (i,j) \in A} x_{ij} = 1, \quad \forall j \in V$$

$$(33.3)$$

$$\sum_{i \in V, (i,j) \in A} x_{ij} = 1, \quad \forall j \in V \tag{33.4}$$

$$\sum_{j \notin S, i \in S, (i,j) \in A} x_{ij} \ge 1, \quad \forall S \subset V, 2 \le |S| \le n - 1$$

$$(33.5)$$

In the formulation, constraints (33.3) and constraints (33.4) are degree constraints, which specify that every vertex is entered exactly once. Constraints (33.5) is the sub-tour constraints, they prohibit the formation of sub-tours. S is a non-empty subset of V, and has at least 2 vertices. 33.5 can be replaced by

$$\sum_{i,j \in S, (i,j) \in A} x_{ij} \le |S| - 1, \quad \forall S \subset V, 2 \le |S| \le n - 1$$
(33.6)

If we list all sub-tour constraints in DFJ, there will be  $O(2^n)$  constraints and  $O(n^2)$  binary variables. The exponential number of constraints makes it impractical to solve directly. Instead, lazy constraints are usually implemented for the sub-tour elimination constraints (33.5 or 33.6).

#### 33.1.2 Miller-Tucker-Zemlin (MTZ) Formulation

We can also formulate TSP using sequential formulations, namely, Miller-Tucker-Zemlin (MTZ) formulation. In the MTZ formulation, the degree constraints (33.3 and 33.4) are the same as in DFJ formulation.

Define a new set of integer decision variables  $u_i$ ,  $u_i$  defined as the sequence in which node i is visited,  $u_1 = 1$ . The sub-tour constraints (33.5 or 33.6) are replaced by the following:

$$u_i - u_j + (n-1)x_{ij} \le n-2, \quad i, j = 2, \dots, n \in V, (i,j) \in A$$
 (33.7)

$$1 \le u_i \le n - 1, \quad i \in 2, \dots, n \in V \tag{33.8}$$

In MTZ formulation, there are  $O(n^2)$  constraints,  $O(n^2)$  binary variables, and O(n) continuous variables.

#### 33.1.3 Quadratic Formulation (QAP)

In this section, we are going to go over a TSP formulation are super bad. However, it still has some value for further study.

The idea is to transform TSP into an assignment problem. Assuming we have n boxes, which represents n steps in the path. Define  $x_{ij}$  as

$$x_{ij} = \begin{cases} 1, & \text{Vertex } i \text{ is assigned to box } j \\ 0, & \text{Otherwise} \end{cases}$$
 (33.9)

The constraints are simple as an assignment problem as following

$$\sum_{j=1}^{n} x_{ij} = 1, \quad \forall i \in V \tag{33.10}$$

$$\sum_{i \in V}^{n} x_{ij} = 1, \quad j = 1, \dots, n$$
(33.11)

However, the tricky part is in the objective function

$$\min \sum_{i \in V} \sum_{j \in V \setminus \{i\}} \sum_{k=1}^{n-1} c_{ij} x_{ik} x_{j,k+1} + \sum_{i \in V} \sum_{j \in V \setminus \{i\}} c_{ij} x_{in} x_{j1}$$
(33.12)

Notice that the objective function is not linear function, with the multiplications of decision variables. Now we are going to linearize them. The linearized version is as following

$$\min \sum_{i \in V} \sum_{j \in V \setminus \{i\}} \sum_{k=1}^{n-1} c_{ij} w_{ij}^k + \sum_{i \in V} \sum_{j \in V \setminus \{i\}} c_{ij} w_{ij}^n$$
(33.13)

s.t. 
$$\sum_{j=1}^{n} x_{ij} = 1$$
,  $\forall i \in V$  (33.14)

$$\sum_{i \in V}^{n} x_{ij} = 1, \quad j = 1, \dots, n$$
(33.15)

$$w_{ij}^k \ge x_{ik} + x_{j,k+1} - 1, \quad i \in V, j \in V \setminus \{i\}, k = 1, \dots, n-1$$
 (33.16)

$$w_{ij}^k \ge x_{ik} + x_{j1} - 1, \quad i \in V, j \in V \setminus \{i\}, k = n$$
 (33.17)

$$w_{ij}^k \in \{0,1\}, \quad i \in V, j \in V \setminus \{i\}, k = 1, \dots, n$$
 (33.18)

$$x_{ij} \in \{0, 1\}, \quad i \in V, j \in V \setminus \{i\}$$
 (33.19)

We can prove that this is very very bad.

#### 33.1.4 Flow Based Formulations

In this section, flow based formulations are discussed, which includes **Single Commodity Flow**, **Two Commodity Flow** and **Multi-Commodity Flow**. In these formulations, continuous variables are introduced to represent the flow on the arcs.

In Single Commodity Flow formulation, define  $y_{ij}$  as the flow in an arc  $(i,j) \in A$ .

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Degree constraints 33.3 and 33.4 are retained. The following constraints are introduced:

$$y_{ij} \le (n-1)x_{ij}, \quad \forall i, j \in V, (i,j) \in A$$
 (33.20)

$$\sum_{i \in V(1,1) \in A} y_{1j} = n - 1 \tag{33.21}$$

$$\sum_{j \in V, (1,j) \in A} y_{1j} = n - 1$$

$$\sum_{i \in V, (i,j) \in A} y_{ij} - \sum_{k \in V, (j,k) \in A} y_{jk} = 1, \quad \forall j \in V \setminus \{1\}$$
(33.21)

(33.23)

Constraints (33.20) can be tighten by the following:

$$y_{ij} \le (n-1)x_{ij}, \quad i = 1, j \in V \setminus \{1\}, (i,j) \in A$$
 (33.24)

$$y_{ij} \le (n-2)x_{ij}, \quad i, j \in V \setminus \{1\}, (i, j) \in A$$
 (33.25)

(33.26)

In SCM formulation, there are  $O(n^2)$  constraints,  $O(n^2)$  binary variables and  $O(n^2)$  continuous variables. In Two Commodity Flow formulation, define  $y_{ij}$  as the flow in an arc  $(i,j) \in A$ , for commodity type 1, and define  $z_{ij}$  as the flow in an arc  $(i,j) \in A$ , for commodity type 2.

Besides degree constraints, the other constraints are as following

$$y_{ij} + z_{ij} = (n-1)x_{ij}, \quad \forall i, j \in V, (i,j) \in A$$
 (33.27)

$$\sum_{j \in V \setminus \{1\}} (y_{1j} - y_{j1}) = n - 1, \quad (1, j) \in A$$
(33.28)

$$\sum_{i \in V} (y_{ij} - y_{ji}) = 1, \quad \forall i \in V \setminus \{1\}, (i, j) \in A$$
 (33.29)

$$\sum_{j \in V \setminus \{1\}} (z_{1j} - z_{j1}) = 1 - n, \quad (1, j) \in A$$
(33.30)

$$\sum_{j \in V} (z_{ij} - z_{ji}) = -1, \quad \forall i \in V \setminus \{1\}, (i, j) \in A$$
 (33.31)

$$\sum_{i \in V} (y_{ij} + z_{ij}) = n - 1, \quad \forall i \in V$$
(33.32)

(33.33)

In TCM formulation, constraints (33.27) only allow flow in an arc if present. Constraints (33.28) and (33.29) forces (n-1) units of commodity type 1 to flow in at node 1 and 1 unit to flow out at every other nodes. Constraints (33.30) and (33.31) are similar, those forces (n-1) units of commodity type 2 to flow out at node 1 and 1 unit to flow in at every other nodes. Constraints (33.32) forces exactly (n-1) units of combined commodity in each arc. In TCM formulation, there are  $O(n^2)$  constraints,  $O(n^2)$  binary variables and  $O(n^2)$  continuous variables.

The SCM and the TCM can be generalized into Multi-Commodity Flow formulation. As usual, degree constraints are retained. The following continuous variables are introduced. Define  $y_{ij}^k$  as the flow of commodity type k in arc  $(i, j) \in A$ .

The other constraints are

$$y_{ij}^k \le x_{ij}, \quad \forall i, j, k \in \mathbb{N}, k \ne 1$$
(33.34)

$$\sum_{i \in V} y_{1i}^k = 1, \quad \forall k \in V \setminus \{1\}$$
(33.35)

$$\sum_{i \in V} y_{i1}^k = 0, \quad \forall k \in V \setminus \{1\}$$
(33.36)

$$\sum_{i \in V} y_{ik}^k = 1, \quad \forall k \in V \setminus \{1\}$$
 (33.37)

$$\sum_{j \in V} y_{jk}^k = 0, \quad \forall k \in V \setminus \{1\}$$
(33.38)

$$\sum_{i \in V} y_{ij}^k - \sum_{i \in V} y_{ji}^k = 0, \quad \forall j, k \in V \setminus \{1\}, (i, j) \in A$$
 (33.39)

Constraints (33.34) only allow flow in an arc which is present. Constraints (33.35) forces exactly one unit of each type of commodity to flow in at node 1. Constraints (33.36) prevent any commodity flow out at node 1. Constraints (33.37), and Constraints (33.38), forces exactly one unit of type k commodity to flow out, and in, at every node except node 1. Constraints (33.39) forces balance of all types of commodities at every node except node 1.

This formulation has  $O(n^3)$  constraints,  $O(n^2)$  binary variables, and  $O(n^3)$  continuous variables.

#### 33.1.5 Shortest Path Formulation

#### A graph for timed staged shortest path

In this section, we are going to introduce another form of formulation with different definition of decision variable and objective function.

Assuming for a completed graph G = (V, A). Define  $x_{ij}^t$  as the following

$$x_{ij}^{t} = \begin{cases} 1, & \text{If path crosses arc } (i,t) \text{ and } (j,t+1) \\ 0, & \text{Otherwise} \end{cases}, \quad i \in V, j \in V \setminus \{i\}, t = 1, \dots, n$$
 (33.40)

The objective function will be

$$\min \sum_{i \in V} \sum_{j \in V \setminus \{i\}} c_{ij} \sum_{t=1}^{n} x_{ij}^{t}$$
(33.41)

The constraints are as following

$$\sum_{j \in V \setminus \{1\}} x_{1j}^1 = 1 \tag{33.42}$$

$$\sum_{j \in V \setminus \{1, i\}} x_{ij}^2 - x_{1i}^1 = 0, \quad \forall i \in V \setminus \{1\}$$
(33.43)

$$\sum_{j \in V \setminus \{1, i\}} x_{ij}^t - \sum_{j \in V \setminus \{1, i\}} x_{ji}^{t-1} = 0, \quad \forall i \in V \setminus \{i\}, t \in \{2, \dots, n-1\}$$
(33.44)

$$x_{i1}^{n} - \sum_{j \in V \setminus \{1, i\}} x_{ji}^{n-1} = 0, \quad \forall i \in V \setminus \{1\}$$
(33.45)

$$\sum_{i \in V \setminus \{1\}} x_{i1}^n = 1 \tag{33.46}$$

$$\sum_{t=2}^{n-1} \sum_{j \in V \setminus \{1, i\}} x_{ij}^t + x_{i1}^n \le 1, \quad \forall i \in V \setminus \{1\}$$
 (33.47)

Notice that constraint (33.47) can be replaced by

$$x_{1i}^{1} + \sum_{t=2}^{n-1} \sum_{j \in V \setminus \{1, i\}} x_{ji}^{t} \le 1, \quad \forall i \in V \setminus \{1\}$$
 (33.48)

33.2	$\mathbf{NP}$	Completeness	of	<b>TSP</b>
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- 33.2.1 Proof of  $TSP \in NPC$
- 33.2.2 Polynomially Solvable Special Cases of TSP
- 33.3 Lower Bounds of TSP
- 33.3.1 The Assignment Lower Bound
- 33.3.2 The Minimum Spanning Tree (Arborescence) Bound
- 33.3.3 The 2-match Problem
- 33.4 Exact Algorithm
- 33.5 Constructive Heuristic
- 33.5.1 Nearest Neighborhood Algorithm
- 33.5.2 Insertion Algorithm
- 33.5.3 Sweep Algorithm
- 33.5.4 Christofides Algorithm
- 33.6 Local Search Heuristic
- 33.6.1 Lin-Kernighan Algorithm

The general idea of Lin-Kernighan Algorithm is based on a substantial generalization of the interchange transformation.

- 33.7 Metaheuristic
- 33.7.1 Genetic Algorithm for TSP

# The Vehicle Routing Problem

# The Capacitate Vehicle Routing Problem

The Vehicle Routing Problem with Time Windows

## Pickup-and-Delivery Problem

#### 37.1 Problem Formulation

Let N be the set of transportation requests. For each transportation request  $i \in N$ , a load of size  $\bar{q}_i \in \mathbb{N}$  has to be transported from a set of origins  $N_i^+$  to a set of destinations  $N_i^-$ . Each load is subdivided as follows

$$\bar{q}_i = \sum_{j \in N_i^+} q_j = -\sum_{j \in N_i^-} q_j \tag{37.1}$$

Define  $N^+ = \bigcup_{i \in N} N_i^+$  as the set of all origins and  $N^- = \bigcup_{i \in N} N_i^-$  as the set of all destinations. Let  $V = N^+ \cup N^-$ . Furthermore, the M be the set of vehicles. Each vehicle  $k \in M$  has a capacity  $Q_k \in \mathbb{N}$ , a start location  $k^+$ , and an end location  $k^-$ . Define  $M^+ = \{k^+ | k \in M\}$  as the set of start locations and  $M^- = \{k^- | k \in M\}$  as the set of end locations. Let  $W = M^+ \cup M^-$ .

For all  $i, j \in V \cup W$  let  $d_{ij}$  denote the travel distance,  $t_{ij}$  denote the travel time, and  $c_{ij}$  denote the travel cost. Note that the dwell time at origins and destinations can be easily incorporated in the travel time and therefore will not be considered explicitly.

**Definition 37.1.1.** A pickup and delivery route  $R_k$  for vehicle k is a directed route through a subset  $V_k \in V$  such that:

- $R_k$  starts in  $k^+$
- $\forall i \in N, (N_i^+ \cup N_i^-) \cap V_k = \emptyset \text{ or } N_i^+ \cup N_i^-$
- If  $N_i^+ \cup N_i^- \subseteq V_k$ , then all locations in  $N_i^+$  are visited before locations in  $N_i^-$
- Vehicle k visits each location in  $V_k$  exactly once
- The vehicle load never exceeds  $Q_k$
- $R_k$  ends in  $k^-$

**Definition 37.1.2.** A pickup and delivery plan is a set of routes  $\mathcal{R} = \{R_k | k \in M\}$  such that:

- $R_k$  is the pickup and delivery route for vehicle k, for each  $k \in M$
- $\{V_k|k\in M\}$  is a partition of V,

Here are the special cases of the General Pickup and Delivery Problem

**Example. The pickup and delivery problem**, where |W| = 1,  $|N_i^+| = |N_i^-| = 1$ ,  $\forall i \in N$ . In this case we define  $i^+$  as the unique element of  $|N_i^+|$  and  $i^-$  as the unique element of  $|N_i^-|$ .

**Example. The dial-a-ride problem**, where |W| = 1 and  $|N_i^+| = |N_i^-| = 1, \bar{q}_i = 1, \forall i \in N$ 

**Example.** The vehicle routing problem, where |W| = 1,  $|N_i^+| = |N_i^-| = 1 \forall i \in \mathbb{N}$  and  $N^+ = W$  or  $N^- = W$ .

**Notice:** Generally speaking, we usually have problem with  $|N_i^+| = |N_i^-| = 1$ . In the cases where  $|N_i^+| > 1$  or  $|N_i^-| > 1$ , the transportation requests can be decomposed into several independent requests with  $|N_i^+| = |N_i^-| = 1$ , unless it has to be served by the same vehicle.

Notice: We are not aware of any real-life applications where both  $|N_i^+| > 1$  and  $|N_i^-| > 1$  at the same time yet.

The following is the formulation for General Pickup and Delivery Problem.

k travels from location i to location j and 0 otherwise.

For  $i \in V \cup W$ , specifying the departure time at vertex i.

and 0 otherwise.

 $D_i$ 

Table 37.1: Decision variable notation

For  $i \in N, k \in M$  Equals to 1 if transportation request i is assigned to vehicle k

For  $(i,j) \in (V \times V) \cup \{(k^+,j)|j \in V\} \cup \{(j,k^-)|j \in V\}, k \in \text{Equals to 1 if vehicle}$ 

For  $i \in V \cup W$ , specifying the load of vehicle arriving at vertex i. Define  $q_{k+} =$ 

91	$0, \forall k \in M.$	or veniere arriving at verter in Beline q <sub>k</sub> +	
min	f(x)		(37.2)
s.t.	$\sum_{k \in M} z_i^k = 1$	$orall i \in N$	(37.3)
	$\sum_{j \in V \cup W} x_{lj}^k = z_i^k$	$\forall i \in N, l \in N_i^+ \cup N_i^-, k \in M$	(37.4)
	$\sum_{j \in V \cup W} x_{jl}^k = z_i^k$	$\forall i \in N, l \in N_i^+ \cup N_i^-, k \in M$	(37.5)
	$\sum_{j \in V \cup \{k^-\}} x_{k^+j}^k = 1$	$\forall k \in M$	(37.6)
	$\sum_{j \in V \cup \{k^+\}} x_{ik^-}^k = 1$	$\forall k \in M$	(37.7)
	$D_{k+} = 0$	$\forall k \in M$	(37.8)
	$D_p \le D_q$	$\forall i \in N, p \in N_i^+, q \in N_i^-$	(37.9)
	$D_i + t_{ij} \le D_j + M(1 - x_{ij}^k)$	$\forall i,j \in V \cup W, k \in M$	(37.10)
	$y_{k^+} = 0$	$\forall k \in M$	(37.11)
	$y_l \le \sum_{k \in M} Q_k z_i^k$	$\forall i \in N, l \in N_i^+ \cup N_i^-$	(37.12)
	$y_j \ge y_l + q_i x_{lj}^k - M(1 - x_{lj}^x)$	$\forall i \in N, l \in N_i^+, j \in V \cup \{k^+\}, l \neq j$	(37.13)
	$y_j \le y_l + q_i x_{lj}^k + M(1 - x_{lj}^x)$	$\forall i \in N, l \in N_i^+, j \in V \cup \{k^+\}, l \neq j$	(37.14)
	$y_j \ge y_l - q_i x_{lj}^k - M(1 - x_{lj}^x)$	$\forall i \in N, l \in N_i^-, j \in V \cup \{k^-\}, l \neq j$	(37.15)
	$y_j \le y_l - q_i x_{lj}^k + M(1 - x_{lj}^x)$	$\forall i \in N, l \in N_i^-, j \in V \cup \{k^-\}, l \neq j$	(37.16)
	$x_{ij}^k \in \{0,1\}$	$\forall i,j \in V \cup W, k \in M$	(37.17)
	$z_i^k \in \{0, 1\}$	$\forall i \in N, k \in M$	(37.18)
	$D_i \ge 0$	$\forall i \in V \cup W$	(37.19)
	$y_i \ge 0$	$orall i \in V \cup W$	(37.20)

Explanation of constraints:

• Constraint (37.3) - Each transportation request is assigned to exactly one vehicle.

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• Constraints (37.4), (37.5) - A vehicle only enters or leaves a location *l* if it is an origin or a destination of a transportation request assigned to that vehicle.

- Constraints (37.6), (37.7) Make sure the vehicle is leaving from /arriving at the correct place.
- Constraint (37.8) Initial starting time for all vehicles is 0.
- Constraint (37.9) For each item to be delivered, it should be picked up before delivery
- Constraint (37.10) Traveling distance from location i to j
- Constraint (37.11) All vehicles leaving the initial location with no loading.
- Constraint (37.12) Load capacity limit for vehicles.
- Constraints (37.13), (37.14), (37.15), (37.16) Load / unload item when arriving pickup / delivery location
- Constraints (37.17), (37.18), (37.19), (37.20) Binary variable definition and non-negativity constraints.

#### 37.2 Heuristic Method

# Stochastic Vehicle Routing Problem

# Dynamic Vehicle Routing Problem

# Part VI Nonlinear Programming

# **Optimality Conditions and Duality**

- 40.1 The Fritz John Optimality Conditions
- 40.2 The Karush-Kuhn-Tucker Optimality Conditions
- 40.3 Constraint Qualification
- 40.4 Lagrangian Duality and Saddle Point Optimality Condition

# **Unconstrained Optimization**

# Quadratic Programming

# Penalty and Barrier Functions

# $\begin{array}{c} {\bf Part~VII} \\ {\bf Algorithms} \end{array}$

### General Paradigms

This chapter we are going to discussion three types of strategies for algorithm design, greedy algorithm, divide-and-conquer, and dynamic programming.

- Greedy algorithm
  - Make a greedy choice
  - Prove that the greedy choice is safe
  - Reduce the problem to a sub-problem and solve it iteratively
  - Usually for optimization problems.
- Divide-and-Conquer
  - Break a problem into many independent sub-problems
  - Solve each sub-problem separately
  - Combine solutions for sub-problems to form a solution for the origin one
  - Usually used to design more efficient algorithm
- Dynamic Programming
  - Break up a problem into many overlapping sub-problems
  - Build solutions for larger and larger sub-problems
  - Use a table to store solutions for sub-problems for reuse

#### 44.1 Greedy Algorithms

#### 44.1.1 Introduction

Greedy algorithm solve the problem incrementally. Its often for optimization problems. Solving optimization problem typically requires a sequences of steps, at each step, an irrevocable decision will be made, and makes the choice looks the best at the moment. Based on that, a small instance will be added to the problem and we solve it again.

Greedy algorithm do not always yield optimal solution, but it usually can provide a relatively acceptable computational complexity. They often run in polynomial time due to the incrementally of instances. Sometimes even if we cannot guarantee the solution is optimal, we still use it in optimization, because of its cheap computation burden. One of the examples will be the constructive heuristic of VRP problems.

Greedy algorithm usually gives polynomial time complexity, but that is not all the cases. In simplex method, each pivot is greedily searching through for the extreme point. Although simplex method usually gives us traceable computation running time, however, it is not a polynomial algorithm.

The following is a very rough sketch of generic greedy algorithm

If we can prove the following, we can claim the greedy algorithm. First, it need to prove that for "current moment", the strategy is safe, i.e., there is always an optimum solution that agrees with the decision made according to the

#### Algorithm 27 Generic Greedy Algorithm

- 1: while The instance is non-trivial do
- 2: Make the choice using the greedy strategy
- 3: Reduce the instance

strategy, this is usually difficult. Then, it need to show that the remaining task after applying the strategy is to solve a/many smaller instance(s) of the same problem.

Although greedy algorithm is intuitive and usually leads to satisfactory complexity, however, for most of the problems there is **no** natural greedy algorithm that works.

#### 44.1.2 Examples

#### **Interval Scheduling**

For n jobs, job i starts at time  $s_i$ , and finishes at time  $f_i$ . i and j are compatible if  $[s_i, f_i)$  and  $[s_j, f_j)$  are disjoint. Find the maximum-size of mutually compatible jobs.

The following is the greedy algorithm to solve this problem

#### **Algorithm 28** Interval Scheduling, S(s, f, n)

```
1: Sort jobs by f

2: t \leftarrow 0, S \leftarrow \emptyset

3: for every j \in [n] according to non-decreasing order of f_j do

4: if s_j \geq t then

5: S \leftarrow S \cup \{j\}

6: t \leftarrow f_j

return S
```

Now we proof the that it is safe to schedule the job j with the earliest finish time, i.e., there is an optimum solution where the job j with the earliest finish time is scheduled.

*Proof.* For arbitary optimum solution S, one of the following cases will happen:

Case 1: S contains j, done

Case 2: S does not contain j. Then the first job in S can be replaced by j to obtain another optimum schedule S'.

#### Unit-length interval covering

Given a set of n points  $X = \{x_1, x_2, \dots, x_n\}$  on the real line, WLOG, assuming the points have already been sorted. We want to use the smallest number of unit-length closed intervals to cover all the points in X.

The following is a greedy algorithm to find the set of unit-length intervals that cover all the points in real line.

#### Algorithm 29 Cover points with unit-length intervals

```
1: Initial, S \leftarrow \emptyset

2: for i = 1, 2, ..., n do

3: if x_i is not covered by any unit-length intervals then

4: Add one unit-length interval starting from x_i, i.e., S \leftarrow S \cup \{x_i\}
```

This strategy (algorithm) is a greedy algorithm because it build up the solution in steps, in each iteration it considers one more point to be covered, i.e., it is optimal for each step in the iteration. Now we prove this algorithm is "safe".

*Proof.* First we consider the case where there is only one point on the real line. Then the optimal number of unit-length interval will be 1, according to the algorithm, that interval will be started at that point.

Then assuming for the case where there are k points from left to right, i.e.,  $X = \{x_1, x_2, \dots, x_k\}$ , and p unit-intervals is already the minimal number and placed by the algorithm, then the  $(k+1)^{th}$  point can only be one of the following cases:

Case 1: The  $(k+1)^{th}$  point is covered by the  $p^{th}$  unit-length interval. According to the strategy, no new unit-length interval will be needed, the number of unit-length interval for k+1 points will be the same as when there are k points. Therefore in this case for k+1 points, p is the minimal (optimal) number. So the strategy is "safe" in this case.

Case 2: The  $(k+1)^{th}$  point is not covered by the  $p^{th}$  unit-length interval. According to the strategy, there will be one new unit-length interval added. Notice that p unit-length intervals will not be feasible to cover k+1 points in this case, because if we move the  $p^{th}$  unit-length interval to the right, it will not be able to cover at least one point which overlapped with the starting point of that unit-length interval. Since p is infeasible and p+1 unit-length interval is feasible, then p+1 is the minimal (optimal) number. So the strategy is "safe" in this case.

Notice that for the k we have mentioned above, k can start from 1 to infinite number of integer. So this strategy is "safe" in every cases.

## 44.2 Divide and Conquer

#### 44.2.1 Introduction

The divide-and-conquer algorithm contains three steps: divide, conquer and combine. Step one, divide instances into many smaller instances. Step two, conquer small instance by solving each of smaller instances recursively and separately. Step three, combine solutions to small instances to obtain a solution for the origin large instance.

Divide and conquer can sometimes solve the problems that greedy algorithm cannot solve, but they often not strong enough to reduce exponential brute-force search down to polynomial time. What usually happen is that they reduce a running time that unnecessarily large, but already polynomial, down to a faster running time.

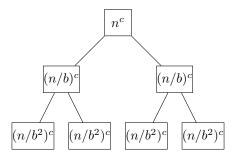
### 44.2.2 Master Theorem

The Master Theorem is useful in analyzing the running time of divide and conquer algorithm. Assume at each step we divide the origin problem of size n into subproblems of size n/b, run for a times of "itself" to conquer those subproblems with a combine operation of  $O(n^c)$ , then the total running time can be derived by the Master Theorem as following:

**Theorem 44.1** (Master Theorem). For running time in forms of  $T(n) = aT(n/b) + O(n^c)$ , where  $a \ge 1$ , b > 1,  $c \ge 0$  are constants. Then

$$T(n) = \begin{cases} O(n^{\lg_b a}), & c < \lg_b a \\ O(n^c \lg n), & c = \lg_b a \\ O(n^c), & c > \lg_b a \end{cases}$$
(44.1)

Proof of Master theorem using recursion tree



*Proof.* The  $i^{th}$  level has  $a^{i-1}$  nodes. For the following cases, we can derive the time complexity:

Case 1:  $c < \lg_b a$  bottom-level dominates  $(\frac{a}{b^c} \lg_b n) n^c = O(n^{\lg_b a})$ 

Case 2:  $c = \lg_b a$  all levels have same time  $n^c \lg_b n = O(n^c \lg n)$ 

Case 3:  $c > \lg_b a$  top-level dominates  $O(n^c)$ 

### 44.2.3 Examples

Counting Inversions

## 44.3 Dynamic Programming

#### 44.3.1 Introduction

The principal of dynamic programming is essentially opposite of the greedy algorithm. Dynamic programming implicitly explores the space of all possible solutions, by carefully decomposing the origin problem into subproblems, and store the solution of those subproblems. Base on those subproblems, then build up larger and larger problem until the origin problem is solved.

### 44.3.2 Examples

### Weighted Interval Scheduling

Consider n jobs, job i starts at time  $s_i$  and finishes at time  $f_i$ , each job has a weight of  $v_i > 0$ . Job i and job j are compatible if  $[s_i, f_i)$  and  $[s_j, f_j)$  are disjoint. Find the maximum-size subset of mutually compatible jobs. A special case of this problem is when the values of all jobs are equal, which will be the interval scheduling problem as discussed in the greedy algorithm examples.

Define p(j) as for a job j, the largest index i < j such that job j is disjoint with job i. Define opt(i) as the optimal value for instance only containing jobs  $\{1, 2, \dots, i\}$ .

Before the algorithm for weighted interval scheduling, sort the jobs by non-decreasing order of finishing time first, in  $O(n \lg n)$  The dynamic programming algorithm is as following:

### Algorithm 30 ComputeOpt(i)

```
1: if i == 0 then

2: return 0

3: else

4: return \max\{v_i + ComputeOpt(p(i)), ComputeOpt(i-1)\}
```

For finding p(i) for one job, it takes  $O(\lg n)$  by binary search. For n jobs the complexity will be  $O(n \lg n)$ . The running time of this algorithm can be exponential in n, if each time ComputeOpt(i) is computed repeatedly. However, if we store the value of each ComputeOpt(i) and reuse it, we can reduce the running time to O(n). We can recover the set of jobs for given (valid) opt(i) by the following algorithm, assuming the jobs has been sorted by non-decreasing order of finishing time.

### Algorithm 31 RecoverJobs()

```
1: Compute p_1, p_2, \cdots, p_n
 2: opt(0) \leftarrow 0
 3: for i \leftarrow 1 to n do
          if opt(i-1) \ge v_i + opt(p_i) then
 4:
 5:
               opt(i) \leftarrow opt(i-1)
 6:
               b[i] \leftarrow N
 7:
          else
               opt(i) \leftarrow v_i + opt(p_i)
 8:
               b[i] \leftarrow Y
10: i \leftarrow n, S \leftarrow \emptyset
     while i \neq 0 do
11:
          if b[i] == N then
12:
               i \leftarrow i - 1
13:
          else
14:
               S \leftarrow S \cup \{i\}
15:
               i \leftarrow p_i
16:
17: return S
```

The above algorithm is using memorized recursion to solve the problem, there is a second efficient algorithm to solve the Weighted Interval Scheduling Problem.

#### Subset Sum Problem

Given an integer bound W>0 and a set of n items, each with an integer weight  $w_i>0$ , find a subset S of items that

$$\max \sum_{i \in S} w_i \tag{44.2}$$

$$\max \sum_{i \in S} w_i$$
s.t. 
$$\sum_{i \in S} w_i \le W$$

$$(44.2)$$

Consider this instance, i items,  $(w_1, w_2, \dots, w_i)$ , budget is W'. For opt[i, W'] there can only be one of the following

Case 1: The value of optimum solution does not contain  $w_i$ , then opt[i, W'] = opt[i-1, W'], else

Case 2: The value of optimum solution contains  $w_i$ , then  $opt[i, W'] = opt[i-1, W'-w_i] + w_i$ 

The algorithm is as following

### Algorithm 32 Optimum Subset

```
1: for W' \leftarrow 0 to W do
 2:
         opt[0, W'] \leftarrow 0
 3: for i \leftarrow 1 to n do
 4:
         for W' \leftarrow 0 to W do
             opt[i, W'] \leftarrow opt[i-1, W']
 5:
             b[i,W'] \xleftarrow{\cdot} N
 6:
             if w_i \leq W' and opt[i-1, W'-w_i] + w_i \geq opt[i, W'] then
 7:
                  opt[i, W'] \leftarrow opt[i-1, W'-w_i] + w_i
 8:
                  b[i, W'] \leftarrow Y
 9:
10: return opt[n, W]
```

### **Algorithm 33** Recover the Optimum Set

```
1: i \leftarrow n, W' \leftarrow W, S \leftarrow 0
2: while i > 0 do
3:
        if b[i, W'] == Y then
             W' \leftarrow W' - w_i
4:
             S \leftarrow \cup \{i\}
5:
        i \leftarrow i-1
7: return S
```

### Optimum Binary Search Tree

Given n elements  $e_1 < e_2 < \cdots < e_n$ ,  $e_i$  has frequency  $f_i$ , the goal is to build a binary search tree for  $\{e_1, e_2, \cdots, e_n\}$ with the minimum accessing cost

$$\sum_{i=1}^{n} f_i d_i \tag{44.4}$$

Where  $d_i$  is the depth of  $e_i$  in the tree.

Suppose we choose  $e_k$  to be the root, than  $e_1, e_2, \dots, e_{k-1}$  are in left-sub tree, and  $e_{k+1}, \dots, e_n$  will be in right-sub

tree, then, denote the cost for the tree and subtrees to be  $C, C_L, C_R$  respectively

$$C = \sum_{j=1}^{n} f_j d_j \tag{44.5}$$

$$=\sum_{j=1}^{n} f_j + \sum_{j=1}^{n} f_j (d_j - 1)$$
(44.6)

$$= \sum_{j=1}^{n} f_j + \sum_{j=1}^{k-1} f_j(d_j - 1) + \sum_{j=k+1}^{n} f_j(d_j - 1)$$
(44.7)

$$=\sum_{j=1}^{n} f_j + C_L + C_R \tag{44.8}$$

Denote opt(i,j) to be the optimal value for the instance of  $(e_i,e_{i+1},\cdots,e_j)$ , then, for every i,j such that  $1 \leq i \leq j \leq n$ 

$$opt(i,j) = \sum_{k=i}^{j} f_k + \min_{k:i \le k \le j} \{ opt(i,k-1) + opt(k+1,j) \}$$
(44.9)

Here is an example

**Example.** Consider the following optimum binary search tree instance. We have 5 elements  $e_1, e_2, e_3, e_4$  and  $e_5$  with  $e_1 < e_2 < e_3 < e_4 < e_5$  and their frequencies are  $f_1 = 5, f_2 = 25, f_3 = 15, f_4 = 10$  and  $f_5 = 30$ . Recall that the goal is to find a binary search tree for the 5 elements so as to minimize  $\sum_{i=1}^{5} \operatorname{depth}(e_i) f_i$ , where  $\operatorname{depth}(e_i)$  is the depth of the element  $e_i$  in the tree. You need to output the best tree as well as its cost. You can try to complete the following tables and show the steps. In the two tables,  $\operatorname{opt}(i,j)$  is the cost of the best tree for the instance containing  $e_i, e_{i+1}, ..., e_j$  and  $\pi(i,j)$  is the root of the best tree.

$ opt(i,j) \setminus j $ $i $	1	2	3	4	5
1	5	35	65	95	170
2		25	55	85	155
3			15	35	90
4				10	50
5					30

$\pi(i,j) \setminus j$	1	2	3	4	5
1	1	2	2	2	3
2		2	2	2 (or 3)	3
3			3	3	5
4				4	5
5					5

Table 44.1: opt and  $\pi$  tables for the optimum binary search tree instance. For cleanness of the table, we assume opt(i, j) = 0 if j < i and there are not shown in the left table.

```
opt(1,2) = min\{0 + opt(2,2), opt(1,1) + 0\} + (f_1 + f_2) = min\{25,5\} + 5 + 25 = 35
opt(2,3) = min\{0 + opt(3,3), opt(2,2) + 0\} + (f_2 + f_3) = min\{15,25\} + 25 + 15 = 55
opt(3,4) = min\{0 + opt(4,4), opt(3,3) + 0\} + (f_3 + f_4) = min\{10,15\} + 15 + 10 = 35
opt(4,5) = min\{0 + opt(5,5), opt(4,4) + 0\} + (f_4 + f_5) = min\{30,10\} + 10 + 30 = 50
opt(1,3) = min\{0 + f(2,3), f(1,1) + f(3,3), f(1,2) + 0\} + (f_1 + f_2 + f_3)
         = \min\{55, 20, 35\} + 5 + 25 + 15 = 65
opt(2,4) = min\{0 + f(3,4), f(2,2) + f(4,4), f(2,3) + 0\} + (f_2 + f_3 + f_4)
         = \min\{35, 35, 55\} + 25 + 15 + 10 = 85
opt(3,5) = min\{0 + f(4,5), f(3,3) + f(5,5), f(3,4) + 0\} + (f_3 + f_4 + f_5)
         = \min\{50, 45, 35\} + 15 + 10 + 30 = 90
opt(1,4) = min\{0 + f(2,4), f(1,1) + f(3,4), f(1,2) + f(4,4), f(1,3) + 0\}
             +(f_1+f_2+f_3+f_4)=95
         = \min\{85, 40, 45, 65\} + 5 + 25 + 15 + 10
opt(2,5) = min\{0 + f(3,5), f(2,2) + f(4,5), f(2,3) + f(5,5), f(2,4) + 0\}
             +(f_2+f_3+f_4+f_5)=155
         =\min\{90,75,85,85\}+25+15+10+30
opt(1,5) = min\{0 + f(2,5), f(1,1) + f(3,5), f(1,2) + f(4,5), f(1,3) + f(5,5) + f(1,4) + 0\}
             +(f_1+f_2+f_3+f_4+f_5)
         = \min\{155, 95, 85, 95, 95\} + 5 + 25 + 15 + 10 + 30 = 170
```

### Matrix Chain Multiplication

Given n matrices  $A_1, A_2, \dots, A_n$  of sizes  $r_1 \times c_1, r_2 \times c_2, \dots, r_n \times c_n$  where  $c_i = r_{i+1}$  for every  $i = 1, 2, \dots, n-1$ . The Matrix Chain Multiplication finds the order of computing  $A_1A_1 \cdots A_n$  with the minimum number of multiplications. The idea is as following. Assume the last step in the multiplication is  $(A_1A_2 \cdots A_i)(A_{i+1}A_{i+2} \cdots A_n)$ . The cost of this step will be  $r_i \times c_i \times c_n$ . So we need to optimally solve two sub-instances, i.e.,  $(A_1A_2 \cdots A_i)$  and  $(A_{i+1}A_{i+2} \cdots A_n)$ .

Denote opt[i, j] as the minimum cost of computing  $A_i A_{i+1} \cdots A_j$ , then

$$opt[i,j] = \begin{cases} 0, & i = j \\ \min_{k:i \le k < j} (opt[i,k] + opt[k+1,j] + r_i \times c_k \times c_j), & j < j \end{cases}$$
(44.10)

The algorithm is as following

#### Algorithm 34 MatrixChainMultiplication

```
1: opt[i,i] \leftarrow 0 \quad \forall i=1,2,\cdots,n
 2: for l \leftarrow 2 to n do
         for i \leftarrow 1 to n - l + 1 do
 3:
              j \leftarrow i + l - 1
 4:
              opt[i,j] \leftarrow \infty
 5:
              for k \leftarrow i to j-1 do
 6:
                   if opt[i,k] + opt[k+1,j+r_ic_kc_i < opt[i,j] then
 7:
 8:
                       opt[i,j] \leftarrow opt[i,k] + opt[k+1,j] + r_i c_k c_j
                       \pi[i,j] \leftarrow k
 9:
10: return opt[1, n]
```

With above algorithm, to construct the optimal solution, the follow algorithm is needed

## 44.4 Compare between three paradigms

## $\overline{\textbf{Algorithm 35} \; \text{PrintOptimalOrder(i, j)}}$

```
1: if i == j then

2: \operatorname{Print}(A_i)

3: else

4: \operatorname{Print}("("))

5: \operatorname{PrintOptimalOrder}(i, \pi[i, j])

6: \operatorname{PrintOptimalOrder}(\pi[i, j] + 1, j)

7: \operatorname{Print}(")")
```

## Sorting

- 45.1 Exchange Sorts
- 45.1.1 Bubble Sort
- 45.1.2 Cocktail Shaker Sort
- 45.1.3 Odd-even Sort
- 45.1.4 Comb Sort
- 45.1.5 Gnome Sort
- 45.1.6 Quicksort

The idea of quicksort is to recursively divide an array into two subarrays, one with smaller number and one with large number, and concatenate the subarrays after all subarrays are singletons.

The most ideal way is to divide the subarrays equally, which requires an algorithm to find the median of an array of size n in O(n) time.

The quicksort algorithm is as following

#### Algorithm 36 Quicksort(A, n)

```
1: if n = 1 then return A

2: Initial, A_L \leftarrow \emptyset, A_H \leftarrow \emptyset

3: x \leftarrow \text{Median}(A)

4: for element \in A do

5: if element \leq x then

6: A_L \leftarrow A_L \cup \{element\}

7: else

8: A_H \leftarrow A_H \cup \{element\}

9: B_L \leftarrow \text{Quicksort}(A_L, A_L.\text{size})

10: B_H \leftarrow \text{Quicksort}(A_H, A_H.\text{size})

11: t \leftarrow \text{number of times } element \text{ appear in } A \text{ return } B_L + element^t + B_H
```

Running time T(n) = 2T(n/2) + O(n),  $T(n) = O(n \lg n)$ For the median finding algorithm

### Algorithm 37 Median(A)

```
1: (To be finished)
```

If we don't use the median finding algorithm, we can modify the Quicksort(A, n) to be a random algorithm by replacing line 3 by  $x \leftarrow \texttt{RandomElement(A)}$ . This modified algorithm has an expected running time of  $O(n \lg n)$ . The worse case running time is  $O(n^2)$ .

Based on Quicksort algorithm, we can define an O(n) algorithm to find the *i*th smallest number in A, given that we have an O(n) algorithm to find median of array.

The selection algorithm is as follows

### Algorithm 38 Selection(A, n, i)

```
1: if n = 1 then
 2:
        return A
 3: else
 4:
        x \leftarrow \texttt{Median(A)}
        for element \in A do
 5:
            if element \leq x then
 6:
                A_L \leftarrow A_L \cup \{element\}
 7:
 8:
                 A_H \leftarrow A_H \cup \{element\}
 9:
        if i \leq A_L.size then
10:
            return Selection (A_L, A_L.size, i)
11:
        else if i > n - A_R.size then
12:
13:
            return Selection (A_R, A_R.size, i - (n - A_R.size))
        else
14:
15:
            \mathbf{return}\ x
```

Similarly, without Median(A), we can replace line 4 by  $x \leftarrow \texttt{RandomElement(A)}$ . Then the expected running time will be O(n)

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### 45.2 Selection Sorts

- 45.2.1 Selection Sort
- 45.2.2 Heapsort
- 45.2.3 Smoothsort
- 45.2.4 Cartesian Tree Sort
- 45.2.5 Tournament Sort
- 45.2.6 Cycle Sort
- 45.2.7 Weak-heap Sort
- 45.3 Insertion Sorts
- 45.3.1 Insertion Sort
- 45.3.2 Shell Sort
- 45.3.3 Splaysort
- 45.3.4 Tree Sort
- 45.3.5 Library Sort
- 45.3.6 Patience Sorting
- 45.4 Merge Sorts

## 45.4.1 Merge Sort

Merge sort is a typical divide and conquer algorithm. It recursively separate an array into two subarrays, and sort while merging them. The algorithm is as following

### **Algorithm 39** MergeSort(A, n)

- 1: if n = 1 then return A
- 2: **else**
- 3:  $B \leftarrow \texttt{MergeSort}(A[0..\lfloor n/2 \rfloor], \lfloor n/2 \rfloor)$
- 4:  $C \leftarrow \texttt{MergeSort}(A[\lceil n/2 \rceil..n], \lceil n/2 \rceil)$ return  $\texttt{Merge}(B, C, \lfloor n/2 \rfloor, \lceil n/2 \rceil)$

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45.4.2	Cascade Merge Sort
45.4.3	Oscillating Merge Sort
45.4.4	Polyphase Merge Sort
45.5	Distribution Sorts
45.5.1	American Flag Sort
45.5.2	Bead Sort
45.5.3	Bucket Sort
45.5.4	Burstsort
45.5.5	Counting Sort
45.5.6	interpolation Sort
45.5.7	Pigenhole Sort
45.5.8	Proxmap Sort
45.5.9	Radix Sort
45.5.10	Flashsort
45.6	Concurrent Sorts
45.6.1	Bitonic Sorter
45.6.2	Batcher Odd-even Mergesort
45.6.3	Pairwise Sorting Network
45.6.4	Samplesort
45.7	Hybird Sorts
45.7.1	Block Merge Sort
45.7.2	Timsort
45.7.3	Spreadsort
45.7.4	Merge-insertion Sort
45.8	Please Don't Do that Sorts

45.8.1 Slowsort

45.8.2 Bogosort

45.8.3 Stooge Sort

## Mathematical Algorithm

#### Polynomial Multiplication 46.1

For given two polynomials of degree n-1, the algorithm outputs the product of two polynomials.

### Example.

$$(3x^3 + 2x^2 - 6x + 9) \times (-2x^3 + 7x^2 - 8x + 4) \tag{46.1}$$

$$= -6x^{6} + 17x^{5} + 24x^{4} - 60x^{3} + 119x^{2} - 96x + 36$$

$$(46.2)$$

Then for input as (3, 2, -6, 9) and (-2, 7, -8, 4), the output will be (-6, 17, 2, -60, 119, -96, 36)

A naive algorithm to solve this problem will be  $O(n^2)$ 

#### Algorithm 40 PolyMultNaive(A, B, n)

- 1: Let C[k] = 0 for every  $k = 0, 1, \dots, 2n 2$
- 2: for  $i \leftarrow 0$  to n-1 do
- for  $j \leftarrow 0$  to n-1 do  $C[i+j] \leftarrow C[i+j] + A[i] \times B[j]$

Use divide and conquer can reduce the running time. The idea is to divide the polynomial with degree of n-1(WLOG, let n be even number) by two polynomials, i.e.

$$p(x) = p_H(x)x^{\frac{n}{2}} + p_L(x)$$
(46.3)

Both  $p_H$  and  $p_L$  are polynomials with degree of  $\frac{n}{2}-1$ , then

$$p(x)q(x) = (p_H(x)x^{\frac{n}{2}} + p_L(x)) \times (q_H(x)x^{\frac{n}{2}} + q_L(x))$$
(46.4)

$$= p_H q_H x^n + (p_H q_L + p_L q_H) x^{\frac{n}{2}} + p_L q_L$$
(46.5)

Therefore

$$multiply(p,q) = multiply(p_H, q_H)x^n$$
 (46.6)

+ 
$$(multiply(p_H, q_L) + multiply(p_L, q_H))x^{\frac{n}{2}} + multiply(p_L, q_L)$$
 (46.7)

$$= multiply(p_H, q_H)x^n \tag{46.8}$$

+ 
$$(multiply(p_H + p_L, q_H + q_L) - multiply(p_H, q_H) - multiply(p_L, q_L))x^{\frac{n}{2}}$$
 (46.9)

$$+ multiply(p_L, q_L)$$
 (46.10)

(46.11)

The algorithm is as following

The running time  $T(n) = 3T(n/2) + O(n), T(n) = O(n^{\log_2 3})$ 

### Algorithm 41 PolyMultiDC(A, B, n)

```
1: if n = 1 then return A[0]B[0]

2: A_L \leftarrow A[0..n/2 - 1], A_H \leftarrow A[n/2..n - 1]

3: B_L \leftarrow B[0..n/2 - 1], B_H \leftarrow B[n/2..n - 1]

4: C_L \leftarrow \text{PolyMultiDC}(A_L, B_L, n/2)

5: C_H \leftarrow \text{PolyMultiDC}(A_H, B_H, n/2)

6: C_M \leftarrow \text{PolyMultiDC}(A_H + A_L, B_H + B_L, n/2)

7: C \leftarrow 0 array of length 2n - 1

8: for i \leftarrow 0 to n - 2 do

9: C[i] \leftarrow C[i] + C_L[i]

10: C[i + n] \leftarrow C[i + n] + C_H[i]

11: C[i + n/2] \leftarrow C[i + n/2] + C_M[i] - C_L[i] - C_H[i]

return C
```

## 46.2 Matrices Multiplication

## 46.3 Gaussian Elimination

## 46.4 Curve Fitting

## 46.5 Integration

Searching

# String

- 48.1 String Searching
- 48.2 Pattern Matching
- 48.3 Longest Common Subsequence
- 48.4 Parse
- 48.5 Optimal Caching
- 48.6 File Compression
- 48.7 Cryptology

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## **Data Structures**

- 49.1 Elementary Data Structures
- 49.2 Hash Tables
- 49.3 Binary Search Trees
- 49.4 Red-Black Trees
- **49.5** B-Trees
- 49.6 Fibonacci Heaps
- 49.7 van Emde Boas Trees

Special Topic: Computational Geometry

## Convex Hull

- 50.1 Computing Slope Statistics
- 50.2 Convexity
- 50.3 Graham's Scan
- 50.4 Turning and orientations

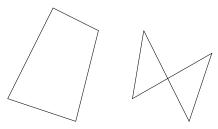
## Intersections

## Triangulation and Partitioning

## 52.1 Polygon Triangulation

### 52.1.1 Types of Polygons

**Definition 52.1.1** (simple polygon). A simple polygon is a closed polygonal curve without self-intersection.

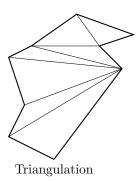


Simple Polygon Non-simple Polygon

Polygons are basic building blocks in most geometric applications. It can model arbitrarily complex shapes, and apply simple algorithms and algebraic representation/manipulation.

### 52.1.2 Triangulation

**Definition 52.1.2** (Triangulation). **Triangulation** is to partition polygon P into non-overlapping triangles using diagonals only. It reduces complex shapes to collection of simpler shapes. Every simple n-gon admits a triangulation which has n-2 triangles.



Theorem 52.1. Every polygon has a triangulation

Lemma 52.2. Every polygon with more than three vertices has a diagonal.

*Proof.* (by Meisters, 1975) Let P be a polygon with more than three vertices. Every vertex of a P is either *convex* or *concave*. W.L.O.G.(any polygon must has convex corner) Assume p is a convex vertex. Denote the neighbors of p as q and r. If  $\bar{q}r$  is a diagonal, done, and we call  $\triangle pqr$  is an ear. If  $\triangle pqr$  is not an ear, it means at least one vertex is inside  $\triangle pqr$ , assume among those vertexes inside  $\triangle pqr$ , s is a vertex closest to s, then s is a diagonal.

### 52.1.3 Art Gallery Theorem

**Theorem 52.3.** Every n-gon can be guarded with  $\lfloor \frac{n}{3} \rfloor$  vertex guards

**Lemma 52.4.** Triangulation graph can be 3-colored.

**Problem 52.1.** The floor plan of an art gallery modeled as a simple polygon with n vertices, there are guards which is stationed at fixed positions with 360 degree vision but cannot see through the walls. How many guards does the art gallery need for the security? (Fun fact: This problem was posted to Vasek Chvatal by Victor Klee in 1973).

*Proof.* - P plus triangulation is a planar graph

- 3-coloring means there exist a 3-partition for vertices that no edge or diagonal has both endpoints within the same set of vertices.
- Proof by Induction:
- Remove an ear (there will always exist ear)
- Inductively 3-color the rest
- Put ear back, coloring new vertex with the label not used by the boundary diagonal.

## 52.1.4 Triangulation Algorithms

# Voronoi Diagrams

# Arrangement and Duality

# **Delaunay Triangulations**

Search

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# **Motion Planning**

## Quadtrees

### Visibility Graphs

## Part VIII Stochastic Methods

### Markov Chain

Queueing Theory

# Part IX Inventory Theory

# Part X Game Theory

## Part XI Simulation

### Random Numbers

### Monte Carlo Simulation

#### Discrete Event Simulation