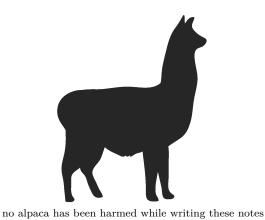
# Foundation of Operations Research

Lorenzo Rossi and everyone who kindly helped! 2022/2023

Last update: 2022-10-04

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# 1 Introduction

# 1.1 Algorithm

An algorithm for a problem is a sequence of instructions that allows to solve any of its instances. The execution time of an algorithm depends on various factors, most notably the instance and the computer.

#### Properties:

- An algorithm is **exact** if it provides an optimal solution for every instance.
  - otherwise is **heuristic**
- A **greedy algorithm** constructs a feasible solution iteratively, by making at each step a *locally optimal* choice, without reconsidering previous choices
  - for most discrete optimization problems, greedy type algorithms yield a feasible solution with no guarantee of optimality

# 1.2 Dynamic Programming

Proposed by *Richard Bellman* in 1950, **dynamic programming** (or DP) is a method for solving optimization problems, composed of a sequence of decisions, by solving a set of recursive equations.

*DP* is applicable to any sequential decision problem, for which the optimality property is satisfied; as such, it has a wide range of applications, including scheduling, transportation, and assignment problems.

# 1.3 Complexity of algorithms

In order analyze an algorithm, it's necessary to consider its complexity as a function of the size of the instance (the size of the input), independently of the computer; the complexity is defined as the number of elementary operations by assuming that each elementary operation takes a constant time.

Since it's hard to determine the exact number of elementary operations, an additional assumption is made: only the asymptotic number of elementary operations in the worst case (for the worst instances) is considered. The complexity evaluation is then performed by looking for the function f(n) that best approximates the upper bound on the number of elementary operations n for the worst instances.

# 1.3.1 Big-O notation

A function f if order of g, written  $f(n) = \mathcal{O}(g(n))$  if exists a constant c > 0 and a constant  $n_0 > 0$  such that  $f(n) \le c \cdot g(n)$  for all  $n \ge n_0$ .

An illustration of the big-O notation is shown in Figure 1.

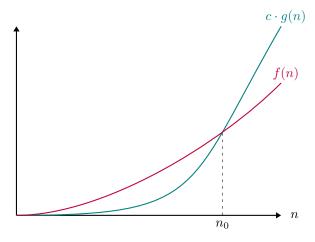


Figure 1: Big-O notation

n	$n^2$	$2^n$
1	$1 \mu s$	$1\mu s$
10	$100\mu s$	1.024ms
20	$400\mu s$	$\approx 1.04  s$
30	$900\mu s$	$\approx 18  m$
40	1.6ms	$\approx 13 d$
50	2.5ms	$\approx 36  y$
60	3.6ms	$\approx 36535  y$

Table 1: Complexity classes

# 1.4 Complexity classes

Two classes of algorithms are considered, according to their worst case order of complexity:

• Polynomial:  $\mathcal{O}\left(n^d\right)$  for a constant  $d>0, d\in\mathbb{R}$ • Exponential:  $\mathcal{O}\left(d^n\right)$  for a constant  $d>0, d\in\mathbb{R}$ 

Algorithms with a hight order Polynomial complexity are not considered efficient.

A comparison of the two classes, assuming that  $1 \mu s$  is needed for each elementary operation, is shown in Table 1.

# 2 Graph and Network Optimization

Many decision making problems can be formulated in terms of graphs and networks, such as:

- transportation and distribution problems
- network design problems
- location problem
- timetable scheduling
- ...

# 2.1 Graphs

A graph is a pair G = (N, E) with:

- N a set of **nodes** or **vertices**
- $E \subseteq N \times N$  a set of **edges** or arcs connecting them pairwise
  - an edge connecting nodes i and j is represented by  $\{i,j\}$  if the graph is **undirected**
  - an edge connecting nodes i and j is represented by (i,j) if the graph is **directed**

# Properties:

- Two **nodes** are **adjacent** if they are connected by an edge
- An edge e is incident in a node v if v is an endpoint of e
  - undirected graphs: the degree of a node is the number of incident edges
    - **directed** graphs: the in-degree (out-degree) of a node is the number of arcs that have it as successor (predecessor)
- A path from  $i \in N$  to  $j \in N$  is a sequence of edges

$$p = \langle \{v_1, v_2\}, \{v_2, v_3\}, \dots, \{v_{k-1}, v_k\} \rangle$$

connecting nodes  $v_1, \ldots, v_k$ , with  $\{v_i, v_{i+1} \in E\}$  for  $i = 1, \ldots, k-1$ 

• A directed path  $i \in N$  to  $j \in N$  is a sequence of arcs

$$p = \langle (v_1, v_2), (v_2, v_3), \dots, (v_{k-1}, v_k) \rangle$$

connecting nodes, with  $(v_i, v_{i+1} \in E)$  for  $i = 1, \dots, k-1$ 

- Nodes u and v are **connected** if exists a path connecting them
- A graph (N, E) is **connecting** if u, v are connecting  $\forall u, v \in N$
- A graph (N, E) is **strongly connected** if u, v are connected by a directed path  $\forall u, v \in N$
- A graph is **bipartite** if there is a partition  $N = N_1 \cup N_2$ ,  $N_1 \cap N_2 = \emptyset$  such that  $\forall (u, v) \in E, u \in N_1$  and  $v \in N_2$
- A graph is **complete** if  $E = \{\{v_i, v_j\} \mid v_i, v_j \in N \land i \leq j\}$
- Given a directed graph G = (N, A) and  $S \subseteq N$ , the **outgoing cut** induced by S is the set of arcs:

$$\delta^+(S) = \{(u, v) \in A \mid u \in S \land v \in N \setminus S\}$$

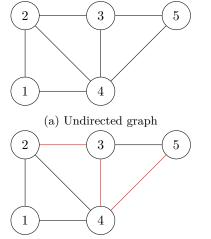
the **incoming cut** induced by S is the set of arcs:

$$\delta^{-}(S) = \{(u, v) \in A \mid v \in S \land u \in N \setminus S\}$$

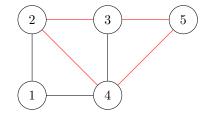
Some examples are shown in Figure 2.

# Properties of graphs:

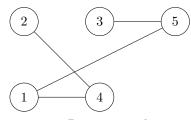
- A graph with n nodes has at most  $m = \frac{n(n-1)}{2}$  edges
- A directed graph with n nodes has at most m = n(n-1) arcs
  - a graph is **dense** if  $m \approx n^2$
  - a graph is sparse if  $m \ll n$



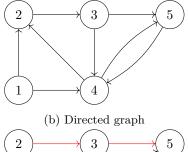
Connected graph, nodes 2 and 5 are connected (c)  $\langle \{2,3\}, \{3,4\}, \{4,5\} \rangle \text{ is a path}$ 

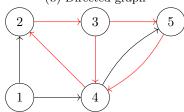


 $\begin{array}{c} {\rm Cycle} \\ {\rm (e)}\ \langle \{2,3\}, \{3,5\}, \{5,4\}, \{4,2\} \rangle \ {\rm is\ a\ cycle} \end{array}$ 

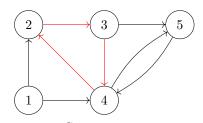


Bipartite graph (g)  $N_1 = \{1, 2, 3\}, N_2 = \{4, 5\}$ 

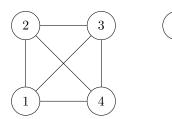




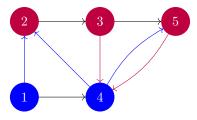
Not strongly connected graph (d)  $\langle \{3,5\}, \{5,4\}, \{4,2\}, \{2,3\}, \{3,4\} \rangle$  is a directed path



Circuit (f)  $\langle (2,3), (3,4), (4,2) \rangle$  is a circuit

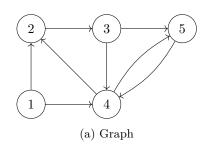


 $\begin{array}{c} \text{Complete graph} \\ \text{(h)} \ \ N = \{1,2,3,4\} \end{array}$ 



(i) incoming  $(\delta^+)$  and outgoing  $(\delta^-)$  cuts of two sets of nodes (purple and blue)

Figure 2: Examples of graphs



$$A = \begin{bmatrix} 0 & 1 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 & 1 \\ 0 & 1 & 0 & 0 & 1 \\ 0 & 0 & 0 & 1 & 0 \end{bmatrix}$$

$$S(1) = \{2, 4\}$$

$$S(2) = \{3\}$$

$$S(3) = \{4, 5\}$$

$$S(4) = \{2, 5\}$$

$$S(5) = \{4\}$$

(b) Adjacency matrix

(c) Adjacency list

Figure 3: Graph representation

## 2.1.1 Graphs representation

Graphs are represented by:

• Adjacency **matrix A** of size  $n \times n$  if the graph is dense:

$$a_{ij} = \begin{cases} 1 & \text{if } (i,j) \in A \\ a_{ij} & \text{otherwise} \end{cases}$$

• Adjacency **list**  $\mathbf{A}$  of size n if the graph is sparse

The same representation can be used for both directed and undirected graphs; the adjacency matrix for an undirected graph is **symmetric**.

An example of a graph representation is shown in Figure 3.

## 2.1.2 Graph reachability problem

**Definition**: Given a directed graph G = (N, A) and a node  $s \in N$ , find all nodes reachable from s

## Goal

 $\rightarrow$  Input: graph G=(N,A), described via successor lists, and a node  $s\in N$   $\rightarrow$  Output: subset  $M\subseteq N$  of nodes of G reachable from s

The goal is reached by an efficient algorithm to solve the problem, with the following properties:

- $\bullet$  a queue Q of nodes not yet processed is kept by the algorithm
- the queue uses a FIFO policy
- the nodes exploration is performed in a breadth-first manner

**Algorithm** The algorithm pseudocode is shown in Code 1.

```
Q := \{s\}
    M := { } { }
    while Q is not empty do
 4
       u := node \in Q
 5
       Q := Q \setminus \{u\}
 6
       M := M \cup \{u\}
 7
       for (u, v) \in \delta^+(u) do
          if v \notin M and v \notin Q then
 8
9
            Q := Q \cup \{v\}
10
          end
       end
11
12
    end
```

Code 1: Graph reachability

The algorithm stops when  $\delta^+(M) = \emptyset$  (when the outgoing cut of the set of nodes M is empty);  $\delta^-(M)$  is the set of arcs with head node in M and tail in  $N \setminus M$ .

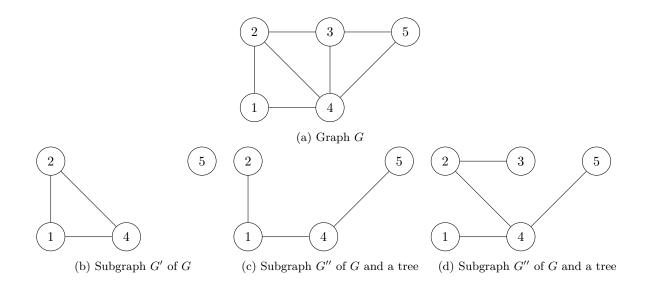


Figure 4: Subgraphs and trees

## 2.1.2.1 Complexity analysis

At each iteration of the while loop:

- 1. A node u is **removed** from the queue Q and **added** to the set M
- 2. For all nodes v directly reachable from u and not already in M or Q, v is added to Q

Since each node u is inserted in Q at most once and each arch (u, v) is considered at most once, the overall complexity is:

$$\mathcal{O}(n+m)$$
  $n=|N|, m=|A|$ 

For dense graphs, this value converges to  $\mathcal{O}(n^2)$ .

# 2.2 Subgraphs and Trees

Let G = (N, E) be a graph. Then:

- G' = (N', E') is a **subgraph** of G if  $N' \subseteq N$  and  $E' \subseteq E$
- A tree  $G_T = (N', T)$  of G is a connected, acyclic, subgraph of G
- $G_T = (N', T)$  is a spanning tree of G if it contains all the nodes (N' = N)
- The leaves of a tree are the nodes with degree 1

A representation of these concepts is shown in Figure 4.

# 2.3 Properties of trees

# 2.3.1 Property 1 - number of edges

Every tree with n nodes has n-1 edges.

#### 2.3.1.1 Proof

- Base case: the claim holds for n = 1 (a tree with a single node has no edges)
- Inductive steps: show that the claim is valid for for any tree with n+1 nodes
  - let  $T_1$  be a tree with n+1 and recall with any tree with  $n\geq 2$  nodes has at least 2 leaves
  - by deleting one of the leaves and its incident edge, a tree  $T_2$  with n nodes is obtained
  - by induction hypothesis,  $T_2$  has n-1 edges; therefore,  $T_1$  has n-1+1=n edges

# 2.3.2 Property 2 - number of paths

Any pair of nodes in a tree is connected via a unique path. Otherwise, the tree would contain a cycle.

#### 2.3.3 Property 3 - new cycles

By adding a new edge to a tree, a new unique cycle is created. This cycle consists of the path in Property 2 - number of paths and the new edge.

# 2.3.4 Property 4 - exchange property

Let  $G_T = (N, T)$  be a spanning tree of G = (N, E). Consider an edge  $e \notin T$  and the unique cycle C of  $T \cup \{e\}$ . For each edge  $f \in C \setminus \{e\}$ , the subgraph  $T \cup \{e\} \setminus \{f\}$  is a spanning tree of G.

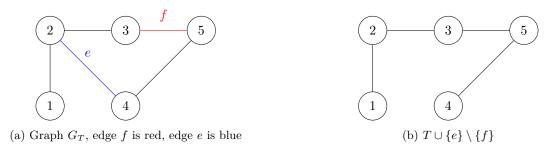


Figure 5: Exchange property

## 2.3.5 Property 5 - cut property

Let F be a partial tree (spanning nodes in  $S \subseteq N$ ) contained in a optimal spanning tree of G = (N, E). Consider  $e = \{u, v\} \in \delta(S)$  of minimum cost, then there exists a minimum cost spanning tree of G containing e.

# 2.3.5.1 Proof

By contradiction, assume  $T^* \subseteq E$  is a minimum cost spanning tree with  $F \subseteq T^*$  and  $e \notin T^*$ . Adding an edge e to  $T^*$  creates the cycle C. Let  $f \in \delta(S) \cap C$ :

- If  $c_e = c_f$ , then  $T^* \cup \{e\} \setminus \{f\}$  is a minimum cost spanning tree of G as it has the same cost as  $T^*$
- If  $c_e < c_f$ , then  $c(T^* \cup \{e\} \setminus \{f\}) < c(T^*)$ , hence  $T^*$  is not optimal

# 2.4 Optimal cost spanning tree

Spanning trees have a number of applications:

- network design
- IP network protocols
- compact memory storage
- ..

**Model**: an undirected graph G = (N, E), n = |N|, m = |E| and a cost function  $c : E \to \mathbb{R}$ , that assigns a cost to each edge, with  $e = \{u, v\} \in E$ .

## Required properties

1. Each pair of nodes must be in a path  $\Rightarrow$  the output must be a connected subgraph containing all the nodes N of G

2. The subgraph must have no cycles  $\Rightarrow$  the output must be a tree

**Formalized problem**: given an undirected graph G = (N, E) and a cost function  $c : E \to \mathbb{R}$ , find a spanning tree  $G_T(N, T)$  of G of minimum, total cost.

The objective is finding:

$$\min_{T \in X} \sum_{e \in T} c_e \qquad X = \text{ set of all spanning trees of } G$$

#### 2.4.1 Theorem 1 - number of nodes in spanning trees

The **Theorem 1**, formulated by Cayley in 1889, states that: A complete graph with n nodes  $(n \ge 1)$  has  $n^{n-2}$  spanning trees.

## 2.4.2 Prim's algorithm

Idea: iteratively build a spanning tree.

#### Method

- 1. Start from initial tree (S,T) with  $S = \{u\}, S \subseteq N$  and  $T = \emptyset$
- 2. At each ste, add to the current partial tree (S,T) an edge of minimum cost among those which connect a node in S to a node in  $N \setminus S$

#### Goal

- $\rightarrow$  Input: connected graph G = (N, E) with edge costs.
- $\rightarrow$  Output: subset  $T \subseteq N$  of edges of G such that  $G_T = (N, T)$  is a minimum cost spanning tree of G.

Complexity if all edges are scanned at each iteration, the complexity order is  $\mathcal{O}(nm)$ 

**Algorithm** the pseudocode of the algorithm is shown in Code 2.

Code 2: Prim's algorithm

Prim's algorithm is **greedy**: at each step a minimum cost edge is selected among those in the cut  $\delta(S)$  induced by the current set of nodes S.

## 2.4.2.1 Correcteness of Prim's algorithm

**Proposition**: Prim's algorithm is exact.

The exactness does not depend on the choice of the first node nor on the selected edge of minimum cost  $\delta(S)$ . Each selected edge is part of the optimal solution as it belongs to a minimum spanning tree.

## 2.4.2.2 Optimality test

The optimality condition allows to verify whether a spanning tree T is optimal or not; it suffices to check that each  $e \in E \setminus T$  is not a cost decreasing edge.

#### 2.4.2.3 Implementation in quadratic time

The Prim's algorithm can be implemented in quadratic time  $(\mathcal{O}(n^2))$ .

#### Data structure

- ullet number of edges selected so far
- Subset  $S \subseteq N$  of nodes incident to the selected edges
- Subset  $T \subseteq E$  of selected edges

• 
$$C_j = \begin{cases} \min\{c_{ij} \mid i \in S\} & j \notin S \\ +\infty & \text{otherwise} \end{cases}$$

• 
$$closest_j = \begin{cases} arg \min\{c_{ij} \mid i \in S\} & j \notin S \\ predecessor of j in the minimum spanning tree & j \in S \end{cases}$$

An example of a step is shown in Figure 6.

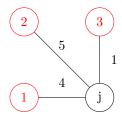


Figure 6: Data structure nodes  $1, 2, 3 \in S$ , node  $j \notin S$  closest<sub>j</sub> = 3  $c_{\text{closest}_{j}j} = 1$ 

The spanning tree is built by selecting the node j with minimum cost  $C_j$  and adding the edge  $\{j, closest_j\}$  to the spanning tree.

The code for this algorithm is shown in Code 3.

```
T := {}
   S := \{u\}
   // initialization
   for j \notin N \setminus S do
5
      if \{u, j\} \in E then
6
        C_{j} := c_{u, j}
7
      else
        C_j := + \inf y
8
9
      end
10
      closest_j := u
11
   end
12
   for k := 1 to n - 1 do
      \min := +\infty // \text{ selection of min cost edge}
13
14
      for j := 1, \ldots, n do
        if j \notin S and C_j < min then
15
16
           min := C_j
17
           v := j
18
        end
19
      S := S \cup \{v\} // \text{ extend } S
20
      T := T \cup \{\{v, closest_v\}\} // extend T
21
22
      for j := 1 to n do
```

Code 3: Prim's algorithm in quadratic time

The complexity of this algorithm is  $\mathcal{O}(n^2)$ . For sparse graphs, where  $m \ll \frac{n(n-1)}{2}$ , a more efficient implementation  $(\mathcal{O}(m \log(2)))$  (using priority queues) is possible.

#### 2.4.3 Optimality condition

Given a spanning tree T, an edge  $e \notin T$  is cost decreasing if when added to T, it creates a cycle C with  $C \subseteq T \cup \{e\}$  and  $\exists f \in C \setminus \{e\}$  such that  $c_e < c_f$ .

# 2.4.4 Theorem 2 - Tree optimality condition

A tree T is of minimum total cost if and only if no cost decreasing edge exists.

#### 2.4.4.1 Proof

- $\Rightarrow$  If a cost decreasing edge exists, then T is not of minimum total cost
- $\Leftarrow$  If no cost decreasing edge exists, then T is of minimum total cost
  - let  $T^*$  be a minimum cost spanning tree of graph G, found via by Prim's algorithm
  - it can be verified that  $T^*$  can be iteratively (changing one edge at a time) transformed into T without changing the total cost
  - thus, T is also optimal

# 2.5 Optimal paths

Optimal (shortest, longest, ...) paths have a wide range of applications:

- Google Maps, GPS navigators
- Planning and management of transportation, electrical, and telecommunication networks
- Problem planning
- ...

**Model**: Given a directed graph G = (N, A) with a cost  $c_{ij} \in \mathbb{R}$  associated to each arc  $(i, j) \in A$ , and two nodes s and t, determine a minimum cost (shortest) path from s to t.

- Each value  $c_{i,j}$  represents the cost (or length, travel time, ...) of arc  $(i,j) \in A$
- Node s is the origin (or source), node t is the destination (or sink)

## **Properties** of optimal paths:

• A path  $\langle (i_1, i_2), (i_2, i_3), \dots, (i_{k-1}, i_k) \rangle$  is simple if no node is visited more than once

# 2.5.1 Property 6 - shortest path

If  $c_{ij} \geq 0$  for all  $(i, j) \in A$ , there is at least one shortest path that is simple.

#### 2.5.2Dijkstra's algorithm

**Idea**: consider the nodes in increasing order of length (cost) of the shortest path from s to any one of the other nodes.

#### Method

- To each **node**  $j \in N$ , a **label**  $L_j$  is associated
  - $\Rightarrow$  at the end of the algorithm, this label will be the cost of the minimum cost path from s to j
- Another label  $predecessor_j$  is associated to each node  $j \in N$ 
  - $\Rightarrow$  at the end of the algorithm, this label will be the node that precedes j on the minimum cost path from s to i
- Make a **greedy** choice with respect to the paths from s to j
- A set of shortest paths from s to any node  $j \notin s$  can be retrieved backwards from t to s iterating over

#### Goal

- $\rightarrow$  Input: graph G = (N, A), cost  $c_{ij} \ge 0 \,\forall i, j$ , origin  $s \in N$
- $\rightarrow$  Output: shortest path from s to all other nodes in G

#### Data structure

- $S \subseteq N$ : subset of nodes whose labels are permanent

• 
$$X \subseteq N$$
: subset of nodes with temporary labels  
•  $L_j = \begin{cases} \text{cost of a shortest path from } s \text{ to } j & j \in S \\ \min\{L_i + c_{ij} \mid (i,j) \in \delta^+(S) & j \notin S \} \end{cases}$ 

- $\rightarrow$  given a directed graph G and the current subset of nodes  $S \subset N$ , consider the outgoing cut  $\delta^+(S)$ and select  $(u, v) \in \delta^+(S)$  such that:  $L_u + c_{uv} = \min\{L_i + c_{ij} \mid (i, j) \in \delta^+(S)\}$
- $\rightarrow$  thus:  $L_u + c_{uv} \le L_i + c_{ij}, \forall (i,j) \in \delta^+(S)$
- predecessor<sub>j</sub> =  $\begin{cases} \text{predecessor of } j \text{ in the shortest path from } s \text{ to } j & j \in S \\ u \text{ such that } L_u + c_{uj} = \min\{L_i + c_{ij} \mid i \in S\} & j \notin S \end{cases}$

Complexity: the complexity of the algorithm depends on the how the arc (u, v) is selected among those of the current cut  $\delta^+(u)$ .

- If all m arcs are scanned, the overall complexity would be  $\mathcal{O}(nm)$ , hence  $\mathcal{O}(n^3)$
- If all labels  $L_i$  are determined by appropriate updates (as in Prim's algorithm), only a single arc of  $\delta^+(j)$ is scanned, hence the complexity is  $\mathcal{O}(n^2)$

# Notes:

- A set of shortest paths from s to all the nodes  $j \in N$  can be retrieved backwards from t to s iterating over the predecessors
- The union of a set of shortest paths from node s to all the other nodes of G is an arborescence rooted at s
- Dijkstra's algorithm does not work when there are arcs with negative cost: if G contains a circuit of negative cost, the shortest path problem may not be well defined

The code for this algorithm is shown in Code 4.

```
1 \quad S := \{ \}
   X := \{s\}
   for u \in N do
4
      L_u := \infty
5
   end
   L_s := 0
6
7
   while |S| < |N| do
      u := argmin\{L_i \mid i \in X\}
8
      X := X \setminus \{u\}
9
10
      S := S \cup \{u\}
      for (u, v) \in \delta^+(u) do
11
12
         if L v > L u + c uv then
           L_v := L_u + c_uv
13
14
           predecessor_v := u
15
           X := X \cup \{v\}
16
         end
17
      end
18
   end
```

Code 4: Dijkstra's algorithm

## 2.5.2.1 Correcteness of Dijkstra's algorithm

Dijkstra's algorithm is correct.

#### **Proof**:

- 1. A the k-th step:
  - $\begin{array}{l} \bullet \;\; S = \{s, i_1, \dots, i_{k-1}\} \\ \bullet \;\; \begin{cases} \text{cost of a minimum cost path from } s \; \text{to} \;\; j \; \in S \\ \text{cost of a minimum cost path with all intermediate nodes in } S \;\; \quad j \notin S \\ \end{array}$
- 2. By induction on the number k of steps:
  - base case: for k = 1 the statement holds, since

$$S = \{s\}, \quad L_s = 0, \quad L_j = +\infty, \quad \forall j \notin S$$

- inductive step: assume that the statement holds for k+1
  - let  $u \notin S$  be the node that is inserted in S and  $\phi$  the path from s to u such that:

$$L_v + c_{vu} \le L_i + c_{iu}, \quad \forall (i, v) \in \delta^+(S)$$

• every path  $\pi$  from s to u has  $c(\pi) \geq c(\phi)$ , as there exists  $i \in S$  and  $j \notin S$  such that:

$$\pi = \pi_1 \cup \{(i, j)\} \cup \pi_2$$

where (i, j) is the first arc in  $\pi \cap \delta^+(S)$ 

• it holds that

$$c(\pi) = c(\pi_1) + c_{ij} + c(\pi_2) \ge L_i + c_{ij}$$

because  $c_{ij} \geq 0 \Rightarrow c(\pi_2) \geq 0$  and by the choice of  $(v, u), c(\pi_1) \geq L_i$ 

• finally, by induction assumption:

$$L_i + c_{ij} \ge L_v + c_{vu} = c(\phi)$$

• a visualization of this step of the proof is shown in Figure 7

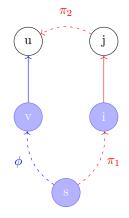
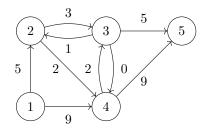


Figure 7: Proof of the induction step; nodes s, v, i are in cut S



(a) Sample graph, with the cost of each arc

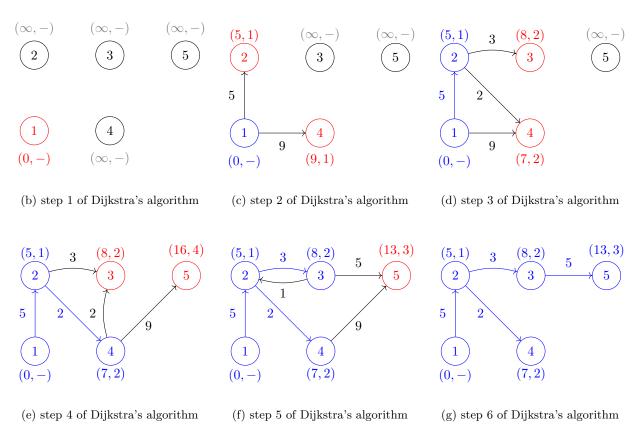


Figure 8: Example of Dijkstra's algorithm

# 2.5.2.2 Example of Dijkstra's algorithm

An example of Dijkstra's algorithm is shown in Figure 8.

## 2.5.3 Floyd-Warshall's algorithm

#### Goal

```
\rightarrow Input: a directed graph G = (N, A) with an n \times n cost matrix C = [c_{ij}] \rightarrow Output: for each pair of nodes i, j \in N, the cost c_{ij} of the shortest path from i to j
```

# Data structure

 $\bullet\,$  Two  $n\times n$  matrices D,P whose elements correspond, at the end of the algorithm, to:

```
d_{ij} the cost of the shortest path from i to j p_{ij} the predecessor of j on the shortest path from i to j
```

#### Method

1. Initialization of D and P:

$$p_{ij} = i \quad \forall i$$

$$d_{ij} = \begin{cases} 0 & i = j \\ c_{ij} & i \neq j \land (i,j) \in A \\ +\infty & \text{otherwise} \end{cases}$$

2. Triangular operation: for each pair of nodes i, j, where  $i \neq u, j \neq u$ , check whether the path from i to j is shorter by going through u (i.e.  $d_{iu} + d_{uj} < d_{ij}$ )

#### Complexity

• Since in the worst case the triangular operation is executed for all nodes u ad for each pair of nodes i, j, the complexity is  $\mathcal{O}(n^3)$ 

The code for this algorithm is shown in Code 5.

```
1
    for i := 1 to n do
 2
      for j := 1 to n do
 3
         p_id := i
 4
         if i = j then
 5
           d_{ij} := 0
         else if (i, j) in A then
 7
8
           d_ij := c_ij
9
           d_{ij} := +\infty
10
11
         end
12
      end
13
   for u \in N do
14
      for i \in N \setminus \{u \} do
15
         for j \in N \setminus \{u\}
16
           if d_iu + d_uj < d_ij then</pre>
17
18
              p_ij := p_uj
              d_ij := d_iu + d_uj
19
20
           end
21
         end
```

```
22     for i ∈ N do
23     if d_ij < 0 then
24         error "negative cycle"
25     end
26     end</pre>
```

Code 5: Floyd-Warshall's algorithm

#### 2.5.3.1 Correctness of Floyd-Warshall's algorithm

Floyd-Warshall's algorithm is correct.

**Proof**: assume that the nodes of G are numbered from 1 to n. Verify that, if the node index order is followed, after the u-th cycle the value  $d_{ij}$  (for any i, j) corresponds to the cost of a shortest path from i to j with at most u intermediate nodes  $(\{1, \ldots, u\})$ 

# 2.6 Optimal paths in directed, acyclic graphs

A directed graph G = (N, A) is **acyclic** if it does not contain any circuit. A directed acyclic graph G is then referred to as a  $\mathbf{DAG}$ .

Property of DAGs: the nodes of any directed acyclic graph G can be ordered topologically, i.e. indexed so that for each arc  $(i, j) \in A$  the index of i is less than the index of j  $(i \le j)$ . The topological order can be exploited by dynamic programming algorithms to compute efficiently the shortest paths in a DAG.

**Problem**: given a  $DAG\ G = (N, A)$  with a cost  $c_{ij} \in \mathbb{R}$  and nodes s, t, determine the shortest (or longest) path from s to t.

## 2.6.1 Topological ordering method

The method requires G = (N, A) to be a DAG represented via the list of predecessors  $\delta^-(v)$  and the list of successors  $\delta^+(v)$  of each node  $v \in N$ . Then, it works as follows:

- 1. Assign the smallest positive integer not yet assigned to a node  $v \in N$  with  $\delta^-(v) = \emptyset$ 
  - $\rightarrow$  such node always exists because G does not contain circuits
- 2. Delete the node v with all its incident arcs
- 3. Go to step (1) until all nodes have been assigned a number

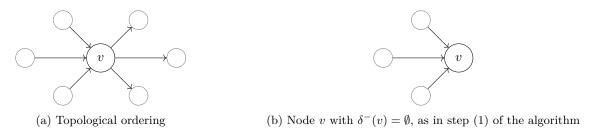


Figure 9: Topological ordering method

This algorithm has complexity  $\mathcal{O}(|A|)$ , because each node is assigned a number only once. Furthermore, all arcs incident to a node are deleted only once.

# 2.6.2 Dynamic programming for shortest path in DAGs

Any shortest path from 1 to t, called  $\pi_t$ , with at least 2 arcs can be subdivided into two parts:

- $\pi_i$ , the shortest subpath from s to i
- (i,t), the remaining part

This decomposition is called the optimality principle of shortest paths in DAGs. An illustration of this decomposition is shown in Figure 10.



Figure 10: Shortest path from 1 to t

The strategy to find the shortest path is:

- 1. For each node i = 1, ..., t let  $L_i$  be the cost of a shortest path from 1 to i
  - $\rightarrow L_t = \min_{(i,t) \in \delta^-(t)} \left\{ L_i + c_{it} \right\}$
  - $\rightarrow$  the minimum is taken over all possible predecessors i of t
- 2. If G is topologically ordered DAG, then the only possible predecessors of t in a shortest path  $\pi_t$  from 1 to t are those with index i < t

  - $\rightarrow L_t = \min_{i < t} \{L_i + c_{it}\}$   $\rightarrow$  in a graph with circuits, any node i can be a predecessor of t if  $i \neq t$

For DAGs whose nodes are topologically ordered  $L_{t-1}, \ldots, L_1$  satisfy the same type of recursive relations:

$$L_{t-1} = \min_{i < t-1} \{L_i + c_{i,t-1}\}; \dots; L_2 = \min_{i=1} \{L_i + c_{i2}\} = L_1 + c_{12}; L_1 = 0$$

which can be solved in reversed order

$$L_1 = 0; L_2 = L_1 + c_{12}; \dots; L_t = \min_{i < t-1} \{L_i + c_t\}$$

**Algorithm**: finally, the algorithm is shown in pseudocode in Algorithm 6.

```
1 sort the nodes of G topologically
2
  L_1 := 0
   for j := 2 to n do
       L_j := \min\{L_i + c_{ij} \mid (i, j) \in \delta^-(j) \land i < j\}
4
       \texttt{pred\_j} := \texttt{v} \texttt{ such that (v, j)} = \texttt{argmin}\{\texttt{L\_i} + \texttt{c\_\{ij\}} \ | \ (\texttt{i, j}) \in \delta^-(\texttt{j}) \land \texttt{i} < \texttt{j}\}
                                                   Code 6: Shortest path in DAG
```

Complexity of the algorithm is  $\mathcal{O}(|A|)$ :

- Topological ordering of the nodes:  $\mathcal{O}(m)$  with m = |A| (number of arcs)
- Each node/arc is processed only once:  $\mathcal{O}(n+m)$

In order to find the longest path, the algorithm can be adapted as follows:

$$L_t = \max_{i < t} \left\{ L_i + c_{it} \right\}$$

# 2.6.2.1 Optimality of the algorithm

The Dynamic Programming algorithm for finding shortest or longest paths in DAGs is exact. This is due to the optimality principle, already explored in the previous section.

# 2.7 Project planning

A project consists of a set of m activities with their (estimated) duration: activity  $A_i$  has duration  $d_i \geq 0, i = 1, \ldots, m$ . Some pair of activities allow a precedent constraint:  $A_i \propto A_j$  indicated that  $A_i$  must be performed before  $A_i$ .

**Model**: a project can be represented by a directed graph G = (N, A) where:

- each arc corresponds to an activity
- the arc length represent the duration of the corresponding activity

In order to account for precedence constraints, the arcs must be positioned such that for activities  $A_i \propto A_j$  there exists a directed path where the arc associated to  $A_i$  precedes the arc associated to  $A_j$ . Such notation is shown in Figure 11.



Figure 11: Precedence relation in project planning

Therefore, a nove v marks an event corresponding to the end fo all the activities  $(i, v) \in \delta^-(v)$  and the (possible) start of all the activities  $(v, j) \in \delta^+(v)$ .

# 2.7.1 Property 7

The directed graph G representing a project is acyclic (is a DAG).

Proof: by contradiction, if  $A_{i1} \propto A_{12} \propto \ldots \propto A_{jk} \propto A_{kj}$  there would be a logical inconsistency.

#### 2.7.2 Optimal paths

A graph G can be simplified by contracting some arcs, but it's important to not introduce unwanted precedence constraints. Artificial nodes or artificial arcs are introduced so that graph G:

- Contains a unique initial node s corresponding to the event "beginning of the project"
- Contains a unique final node t corresponding to the event "end of the project"
- does not contain multiple arcs with the same origin and destination

**Problem**: given a project (set of activities with duration and precedence constraints), schedule the activities in order to minimize the overall project duration (the time needed to complete all the activities).

# 2.7.2.1 Property 8

The minimum overall project duration is the length of a longest path from s to t in the graph G.

**Proof**: since any s-t path represents a sequence of activities that must be executed in the specified order, its length provides a lower bound on the minimum overall project duration.

## 2.7.2.2 Critical path method - CPM

The critical path method (*CPM*) determines:

• A schedule (a plan for executing the activities specifying the order and the assigned time) that minimizes the overall project duration

• The slack of each activity (the amount of time by which its execution can be delayed without affecting the overall minimum project duration)

**Initialization**: construct the graph G representing the project.

#### Method:

- 1. Find a topological order of the nodes
- 2. Consider the nodes by increasing indices and for each  $h \in N$  find the earliest time  $T_{min_h}$  at which the event associated to node h can occur
  - $\rightarrow T_{min_h}$  corresponds to the minimum project duration
- 3. Consider the nodes by decreasing indices and for each  $h \in N$  find the latest time  $T_{max_h}$  at which the event associated to node h can occur without delaying the project completion date beyond  $T_{min_n}$
- 4. For each activity  $(i, j) \in A$  find the slack
  - $\rightarrow$  the slack is calculated as  $\sigma ij T_{max_j} T_{min_i} d_{ij}$

```
Input: graph G = (N, A) with n = |N| and the duration d_{ij} associated to each (i, j) \in A/ Output: (T_{min_i}, T_{max_i}), i = 1, ..., n
```

**Algorithm**: finally, the algorithm is shown in pseudocode in Algorithm 7.

```
1 Sort the nodes topologically 2 T_min_i := 0 3 for j = 2 to n do 4 T_min_j := \max\{T_min_i + d_ij \mid (i, j) \in \delta^-(j)\} 5 end for 6 T_max_n := T_min_n // \min project duration 7 for <math>i = n-1 to 1 do 8 T_max_i := \min\{T_max_j - d_ij \mid (i, j) \in \delta^+(i)\} 9 end for
```

Code 7: Critical path method

**Complexity**: the overall complexity is  $\mathcal{O}((n+m)) \approx \mathcal{O}((m))$ , due to the sum of

- complexity of the topological sort  $\mathcal{O}((n+m))$
- complexity of the first loop  $\mathcal{O}((n+m))$
- complexity of the second loop  $\mathcal{O}((n+m))$

#### 2.7.2.3 Critical paths

An activity (i, j) with zero slack  $\sigma_{ij} = T_{max_j} = T_{min_i} = d_{ij} = 0$  is called **critical**. A critical path is a path in a s - t composed uniquely by critical activities. At least one always exists.

# 2.7.2.4 Gantt charts

A Gantt chart is a graphical representation of a project schedule. It was introduced in 1896 by Henry Gantt, an American mechanical engineer and management consultant.

There are two types of Gantt charts:

- Gantt chart at earliest each activity (i,j) starts at  $T_{min_i}$  and ends at  $T_{min_i} + d_{ij}$
- Gantt chart at latest each activity (i,j) starts at  $T_{max_i}$  and ends at  $T_{max_i} + d_{ij}$