

# Heuristic Algorithms

Master's Degree in Computer Science/Mathematics

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# Effectiveness of a heuristic algorithm

A heuristic algorithm is useful if it is

- ① **efficient**: it “costs” much less than an exact algorithm
- ② **effective**: it “frequently” returns a solution “close to” an exact one

In particular, the **effectiveness** of heuristic algorithms is given by:

- closeness of the solution obtained to an optimal one
- frequency of hitting optimal or nearly optimal solutions

These features can be combined into a

- **frequency distribution of solutions more or less close to the optimum**

The effectiveness of a heuristic algorithm can be investigated with a

- **theoretical analysis (*a priori*)**, proving that the algorithm finds always or with a given frequency **solutions** with a given **guarantee of quality**
- **experimental analysis (*a posteriori*)**, measuring the performance of the algorithm on sampled **benchmark instances** to show that a **guarantee of quality** is respected in practice

We here discuss the theoretical analysis of effectiveness

# Indices of effectiveness

The **effectiveness of a heuristic optimisation algorithm**  $A$  is measured by the **difference** between the heuristic value  $f_A(I)$  and the optimum  $f^*(I)$

- **absolute difference:**

$$\tilde{\delta}_A(I) = |f_A(I) - f^*(I)| \geq 0$$

rarely used, and **only when the objective is a pure number**

- **relative difference:**

$$\delta_A(I) = \frac{|f_A(I) - f^*(I)|}{f^*(I)} \geq 0$$

frequent in **experimental analysis** (*usually as a percent ratio*)

- **approximation ratio:**

$$\rho_A(I) = \max \left[ \frac{f_A(I)}{f^*(I)}, \frac{f^*(I)}{f_A(I)} \right] \geq 1$$

frequent in **theoretical analysis**: the first form is used for minimisation problems, the second one for maximisation problems

# Theoretical analysis (in the worst case)

To obtain a compact measure, independent from  $I$ , find the worst case  
(as for efficiency, that is complexity)

The difference between  $f_A(I)$  and  $f^*(I)$  is in general unlimited, but for some algorithms it is limited:

- absolute approximation:

$$\exists \tilde{\alpha}_A \in \mathbb{N} : \tilde{\delta}_A(I) \leq \tilde{\alpha}_A \text{ for each } I \in \mathcal{I}$$

A (rare) example is Vizing's algorithm for *Edge Coloring* ( $\tilde{\alpha}_A = 1$ )

- relative approximation:

$$\exists \alpha_A \in \mathbb{R}^+ : \rho_A(I) \leq \alpha_A \text{ for each } I \in \mathcal{I}$$

Factor  $\alpha_A$  ( $\tilde{\alpha}_A$ ) is the relative (absolute) approximation guarantee

For other algorithms, the guarantee depends on the instance size

$$\rho_A(I) \leq \alpha_A(n) \text{ for each } I \in \mathcal{I}_n, n \in \mathbb{N}$$

Effectiveness can be independent from size (contrary to efficiency)

# How to achieve an approximation guarantee?

For a minimisation problem, the aim is to prove that

$$\exists \alpha_A \in \mathbb{R} : f_A(I) \leq \alpha_A f^*(I) \text{ for each } I \in \mathcal{I}$$

- ① find a way to build an **underestimate**  $LB(I)$

$$LB(I) \leq f^*(I) \quad I \in \mathcal{I}$$

- ② find a way to build an **overestimate**  $UB(I)$ ,  
related to  $LB(I)$  by a coefficient  $\alpha_A$

$$UB(I) = \alpha_A LB(I) \quad I \in \mathcal{I}$$

- ③ find an **algorithm**  $A$  whose solution is not worse than  $UB(I)$

$$f_A(I) \leq UB(I) \quad I \in \mathcal{I}$$

Then  $f_A(I) \leq UB(I) = \alpha_A LB(I) \leq \alpha_A f^*(I)$ , for each  $I \in \mathcal{I}$

$$f_A(I) \leq \alpha_A f^*(I) \text{ for each } I \in \mathcal{I}$$

# A 2-approximated algorithm for the *VCP*

Given a undirected graph  $G = (V, E)$  find the minimum cardinality vertex subset such that each edge of graph is incident to it

A **matching** is a set of nonadjacent edges

**Maximal matching** is a matching such that any other edge of the graph is adjacent to one of its edges  
*(it cannot be enlarged)*

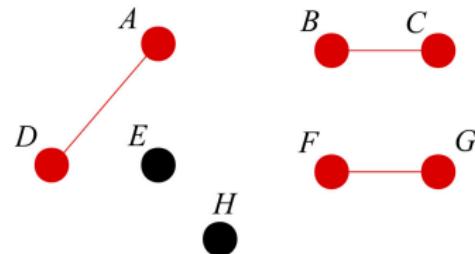
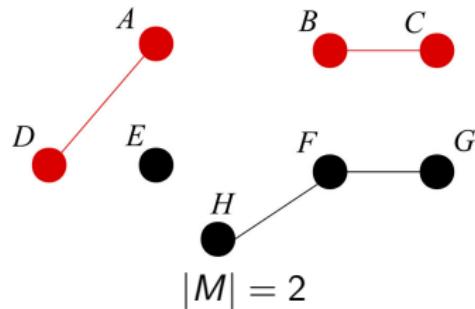
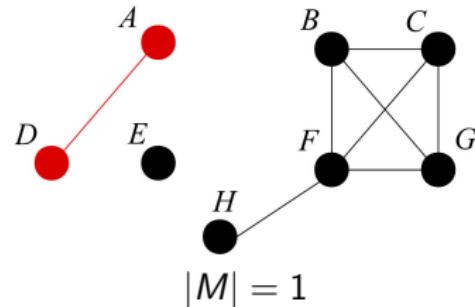
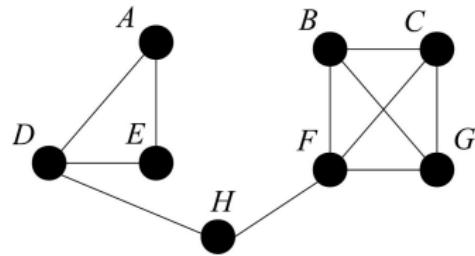
Matching algorithm:

- ① Build a **maximal matching**  $M \subseteq E$  scanning the edges of  $E$  and including in  $M$  those not adjacent to  $M$   
*(now every edge of  $E \setminus M$  is adjacent to an edge of  $M$ )*
- ② The **set of extreme vertices of the matching edges** is a *VCP* solution

$$x_A := \bigcup_{(u,v) \in M} \{u, v\}$$

and it can be improved removing the redundant vertices

# Example



The optimum is  $f^* = 5$

The matching algorithm is 2-approximated

- ① The cardinality of matching  $M$  is an underestimate  $LB(I)$ 
  - the cardinality of an optimal covering for any subset of edges  $E' \subseteq E$  does not exceed that of an optimal covering for  $E$

$$|x_{E'}^*| \leq |x_E^*|$$

*(it costs more to cover all edges than only the matching)*

- the optimal covering of a matching  $M$  has cardinality  $|M|$   
*(each edge of the matching requires exactly one different vertex)*

- ② Including both the extremes of each edge of the matching yields
  - an overestimate *(it covers both the matching and the adjacent edges)*
  - of value  $UB(I) = 2LB(I)$  *(two different vertices for each edge)*
- ③ The matching algorithm returns solutions of value  $f_A(I) \leq UB(I)$   
*(possibly removing redundant vertices)*

This implies  $f_A(I) \leq 2f^*(I)$  for each  $I \in \mathcal{I}$ , that is  $\alpha_A = 2$

... and the bound is tight!

Since  $\alpha_A$  relates  $UB(I)$  and  $LB(I)$ ,  $f_A(I)$  and  $f^*(I)$  could be closer

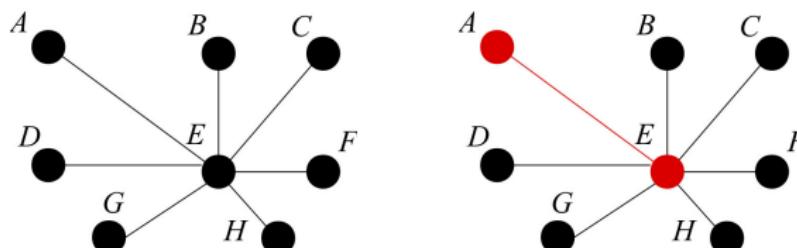
Actually, for many instances  $\rho_A(I)$  is much better than  $\alpha_A$

Are there instances  $\bar{I}$  for which  $f_A(\bar{I}) = \alpha_A f^*(\bar{I})$ ? How are they like?

The study of these instances is useful to

- evaluate whether they are rare or frequent
- introduce *ad hoc* modifications to improve the algorithm

In the literature the typical expression “*and the bound is tight*” introduces the description of instances exhibiting the worst case



If all worst cases are patched, the approximation guarantee improves

# The *TSP* under the triangle inequality

Consider the *TSP* with the additional (rather common) assumptions that

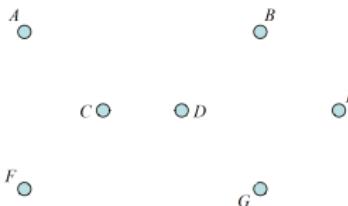
- graph  $G = (N, A)$  is **complete**
- cost  $c$  is **nonnegative, symmetric** and satisfies the **triangle inequality**

$$c_{ij} = c_{ji} \geq 0 \quad \forall i, j \in N \quad \text{and} \quad c_{ij} + c_{jk} \geq c_{ik} \quad \forall i, j, k \in N$$

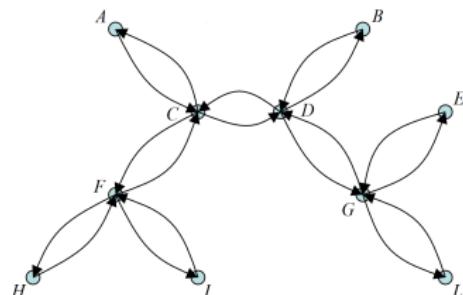
Double-tree algorithm

- ① Consider the **complete undirected graph** corresponding to  $G$
- ② Build a **minimum cost spanning tree**  $T^* = (N, X^*)$
- ③ Make a **pre-order visit** of  $T^*$  and build two lists of arcs:
  - a  $x'$  lists the arcs used both by the visit and the backtracking:  
this is a **circuit** visiting each node, possibly several times
  - b  $x$  lists the arcs linking the nodes in pre-order ending with the first:  
this is a **circuit** visiting each node exactly once

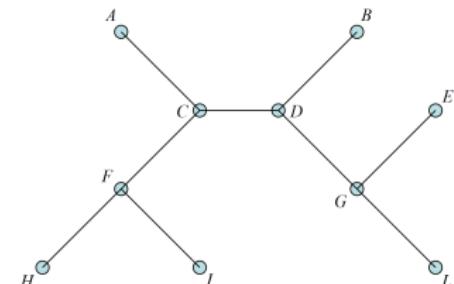
# Example



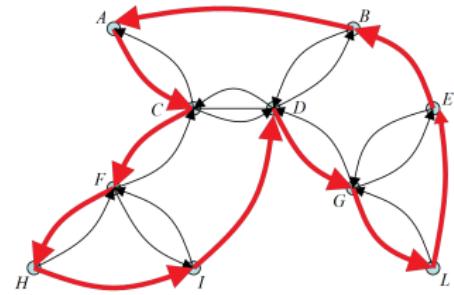
1) Complete graph  $G$  (arcs omitted)



3.a)  $x' = (A, C, F, H, \textcolor{blue}{F}, I, \textcolor{blue}{F}, \textcolor{blue}{C}, D, G, L, \textcolor{blue}{G}, E, \textcolor{blue}{G}, \textcolor{blue}{D}, B, \textcolor{blue}{D}, \textcolor{blue}{C}, A)$



2. Minimum spanning tree  $T^*$



3.b)  $x = (A, C, F, H, I, D, G, L, E, B, A)$

The double-tree algorithm is 2-approximated

- ① the cost of the minimum spanning tree is an underestimate  $LB(I)$ 
  - deleting an arc from a Hamiltonian circuit yields a Hamiltonian path that is cheaper
  - a Hamiltonian path is a spanning tree (usually not of minimum cost)
- ② the cost of circuit  $x'$  is
  - an overestimate  $UB(I)$  *(it is a nonminimum Hamiltonian circuit)*
  - equal to  $2LB(I)$  *(two arcs correspond to each edge)*
- ③ the cost of circuit  $x$  is  $f_A(I) \leq UB(I)$   
*(a single direct arc replacing a sequence decreases the cost)*

This implies that  $f_A(I) \leq 2f^*(I)$  for each  $I \in \mathcal{I}$ , that is  $\alpha_A = 2$

Notice:  $x'$  is used in the approximation proof, but needs not be computed

# Inapproximability

For an **inapproximable problem**, **approximated algorithms** would be exact

Consider this family of *TSP* instances on complete graphs:

- $c_{ij} = 0$  for  $(i, j) \in A_0$
- $c_{ij} = 1$  for  $(i, j) \in (N \times N) \setminus A_0$  (*the triangle inequality is violated!*)

The optimum of any such instance  $\bar{I}$  is:

$$\begin{cases} f^*(\bar{I}) = 0 & \text{if } A_0 \text{ contains a Hamiltonian circuit} \\ f^*(\bar{I}) \geq 1 & \text{otherwise} \end{cases}$$

(*in the latter case, the optimal solution contains at least an arc  $\notin A_0$* )

Assume that a polynomial algorithm  $A$  provide a guarantee  $\alpha_A$

$$f^*(I) \leq f_A(I) \leq \alpha_A f^*(I) \quad \forall I \in \mathcal{I}$$

Then  $f^*(\bar{I}) = 0 \Leftrightarrow f_A(\bar{I}) = 0$

Whenever the subgraph  $G(N, A_0)$  has a Hamiltonian circuit,  $A$  finds it, solving an  $\mathcal{NP}$ -complete problem in polynomial time ( $\mathcal{P} = \mathcal{NP}$ )

For hard problems

- exact algorithms provide the best approximation guarantee ( $\alpha_A = 1$ ), but require exponential time  $T_A$
- approximated algorithms provide a worse guarantee ( $\alpha_A > 1$ ), but could require polynomial time  $T_A$

Some problems admit a family of algorithms providing a whole range of **compromises between efficiency and effectiveness**

- better and better approximation guarantees:  $\alpha_{A_1} > \dots > \alpha_{A_r}$
- worse and worse computational complexities:  $T_{A_1} < \dots < T_{A_r}$

**Approximation scheme** is a parametric algorithm  $A_\alpha$  allowing to choose  $\alpha$   
(Example: the *KP*)

# Beyond the worst case

As usual, the worst-case approach is rough:  
some algorithms often have a good performance, though sometimes bad

The alternative approaches are similar to the ones used for complexity

- **parametrisation**: prove an approximation guarantee that depends on other parameters of the instances besides the size  $n$
- **average-case**: assume a probability distribution on the instances and evaluate the expected value of the approximation factor  
(*the algorithm could have a bad performance only on rare instances*)

but there is at least another approach

- **randomisation**: the operations of the algorithm depend not only on the instance, but also on pseudorandom numbers, so that the solution becomes a **random variable** which can be investigated  
(*the time complexity could also be random, but usually is not*)

# Randomised approximation algorithms

For a randomised algorithm  $A$ ,  $f_A(I, \omega)$  and  $\rho_A(I, \omega)$  are random variables depending on the pseudorandom number seed  $\omega$

A **randomised approximation algorithm** has an **approximation ratio** whose expected value is limited by a constant

$$E[\rho_A(I, \omega)] \leq \alpha_A \text{ for each } I \in \mathcal{I}$$

Max-SAT problem: given a CNF, find a truth assignment to the logical variables that satisfy a maximum weight subset of formulae

Purely random algorithm:

Assign to each variable  $x_j$  ( $j = 1, \dots, n$ )

- value *False* with probability  $1/2$
- value *True* with probability  $1/2$

*What is the expected value of the solution?*

# Randomised approximation for the MAX-SAT

The objective  $f(x) = f_A(I, \omega)$  is the total weight of the satisfied clauses and its expected value is

$$E[f_A(I, \omega)] = E \left[ \sum_{i \in \mathcal{C}} \delta_i(x(\omega)) w_i \right] = \sum_{i \in \mathcal{C}} (w_i \cdot \Pr[\delta_i(x(\omega)) = 1])$$

where  $\delta_i(x(\omega)) = 1$  if solution  $x(\omega)$  satisfies clause  $i$ , 0 otherwise

Let  $k_i$  be the number of literals of formula  $i \in \mathcal{C}$  and  $k_{\min} = \min_{i \in \mathcal{C}} k_i$

$$\Pr[\delta_i(x(\omega)) = 1] = 1 - \left(\frac{1}{2}\right)^{k_i} \geq 1 - \left(\frac{1}{2}\right)^{k_{\min}} \text{ for each } i \in \mathcal{C}$$

$$\Rightarrow E[f_A(I, \omega)] \geq \sum_{i \in \mathcal{C}} w_i \cdot \left[1 - \left(\frac{1}{2}\right)^{k_{\min}}\right] = \left[1 - \left(\frac{1}{2}\right)^{k_{\min}}\right] \sum_{i \in \mathcal{C}} w_i$$

and since  $\sum_{i \in \mathcal{C}} w_i \geq f^*(I)$  for each  $I \in \mathcal{I}$  one obtains

$$\frac{E[f_A(I, \omega)]}{f^*(I)} \geq \left[1 - \left(\frac{1}{2}\right)^{k_{\min}}\right] \geq \frac{1}{2}$$