

Algorithmics: Basic definitions and concepts

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Algorithms:
our context
Time
complexity
Asymptotic
notation
Graphs
Data structures
Traversals
Reductions
Divide and
conquer



The Algorithms course

Already known (nivell EDA)

- Algorithms cost and Asymptotic notation
- Sorting algorithms: Mergesort, Quicksort, ...
- Divide and conquer, recurrences, master theorem
- Complexity, P and NP, reductions
- Foundations on probability
- Basic data structures: Arrays, lists, stacks, queues, heaps, hashing ...
- Basics on graph theory, graph data structures
- Graph and digraph traversals (BFS, DFS) and applications.
- Backtracking algorithms

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The Algorithms course

Topics to cover:

- Divide and conquer: Linear Selection
- Sorting in linear time
- Greedy algorithms
- Dynamic programming
- Distances in graphs
- Flow networks: problems, algorithms and applications
- Linear Programming
- Approximation algorithms

Provide models to **solve** real problems

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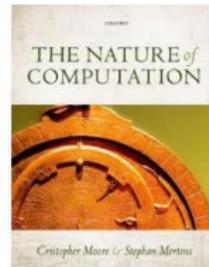
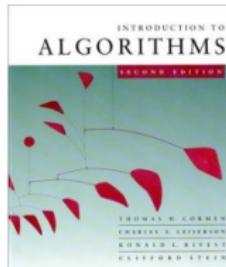
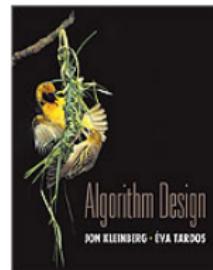
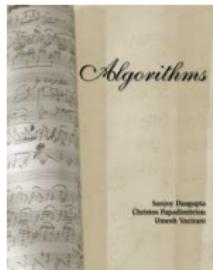
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References

Main references:



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"The algorithmic lenses: C. Papadimitriou"

- In 1936 Alan Turing demonstrated the universality of computational principles with his mathematical model of the Turing machine.
- Theoretical Computer Science views computation as a ubiquitous phenomenon, not one that it is limited to computers.
- Algorithms themselves have evolved into a complex set of techniques, for instances self-learning, Web services, concurrent, distributed or parallel, etc... Each of them with ad-hoc relevant computational limitations and social implications.
- However, this course will be a course on classical algorithms, which are the core needed to understand more advanced computational material.

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Algorithm: Precise recipe for a precise computational task.
Each step of the process must be clear and unambiguous, and
it should always yield a clear answer.

Sqrt (n)

$x_0 = 1$

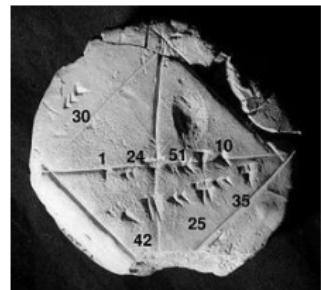
for $i = 1$ to 6 **do**

$x_i = (x_{i-1} + n/x_{i-1})/2$

end for

Babilònia (XVI BC)

For $n = 20$, x 's are 1 10.5 6.2023 4.7134 4.4783



Once we designed an algorithm: What do we want to know?

- Correctness, it always does what it should?
- Performance,
 - computing time,
 - memory use
 - communication cost, ...

For an algorithm \mathcal{A} , $t_{\mathcal{A}}(x)$ is the computing time on input x .

In this course, we use a **worst case analysis**: Given a problem, for which you designed an algorithm, you assume that your meanest adversary gives you the worst possible input.

We use as measure of **time complexity** or **cost** the function

$$T(n) = \max_{|x|=n} t_{\mathcal{A}}(x)$$

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Time complexity

The time complexity must be independent of the "used" machine

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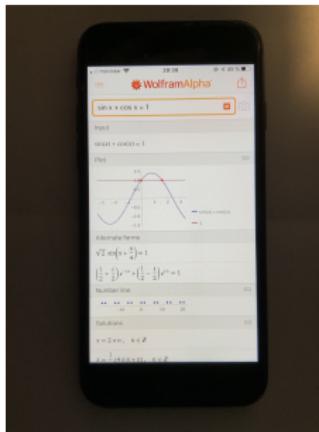
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We must consider carefully how operations scale with respect to size.

Typical computation times

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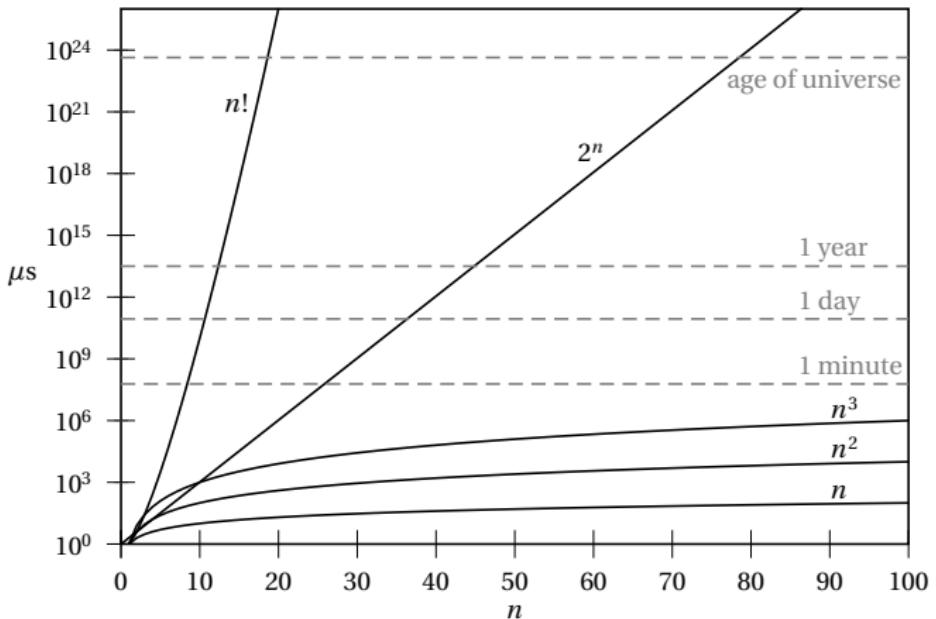
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We study the behavior of $T(n)$ when n can take very large values (i.e., $n \rightarrow \infty$)

- if $n = 10$, $n^2 = 100$ and $2^n = 1024$;
- if $n = 100$, $n^2 = 10000$ and
$$2^n = 12676506002282244014696703205376;$$
- if $n = 10^3$, $n^2 = 10^6$ and 2^n is a number with 302 digits.
- As a comparison, 10^{64} is estimated to be the number of atoms in earth ($< 2^{213}$).

Computation time assuming that an input with size $n = 1$ can be solved in 1 μ second:



From: Moore-Mertens, The Nature of Computation

Computation times as a function of input size n

	n	$n \lg n$	n^2	1.5^n	2^n
10	< 1s	< 1s	< 1s	< 1s	< 1s
50	< 1s	< 1s	< 1s	11m	36y
100	< 1s	< 1s	< 1s	12000y	$10^{17}y$
1000	< 1s	< 1s	< 1s	$> 10^{25}y$	$> 10^{25}y$
10^4	< 1s	< 1s	< 1s	$> 10^{25}y$	$> 10^{25}y$
10^5	< 1s	< 1s	< 1s	$> 10^{25}y$	$> 10^{25}y$
10^6	< 1s	20s	12d	$> 10^{25}y$	$> 10^{25}y$

From: Moore-Mertens, The Nature of Computation

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Efficient algorithms and practical algorithms

- When analyzing an algorithm, we say that it is **feasible** if its cost is polynomial.
 $n^{10^{10}}$ is a polynomial but this computing time could be prohibitive!
- In the same way, if we have cn^2 for constant $c = 10^{64}$, then c dominates inputs up to a size of $n > 10^{64}$.
- In this course, we will not enter in the analysis up to constants, but keep in mind that constants matter!!!!
- In practice, even for feasible algorithms with time complexity of for example n^4 , it could be too slow for $n \geq 1000$.

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Symbol	$L = \lim_{n \rightarrow \infty} \frac{f(n)}{g(n)}$	intuition
$f(n) = O(g(n))$	$L < \infty$	$f \leq g$
$f(n) = \Omega(g(n))$	$L > 0$	$f \geq g$
$f(n) = \Theta(g(n))$	$0 < L < \infty$	$f = g$
$f(n) = o(g(n))$	$L = 0$	$f < g$
$f(n) = \omega(g(n))$	$L = \infty$	$f > g$

Names used for specific function classes

name	definition
polylogarithmic	$f = O(\log^c n)$ (c constant)
polynomial	$f = O(n^c)$ (c constant) or $n^{O(1)}$
subexponential	$f = o(2^{n^\epsilon})$ ($0 < \epsilon < 1$)
exponential	$f = 2^{\text{poly}(n)}$
double exponential	$f = 2^{\exp(n)}$

Notation:

$\lg \equiv \log_2$; $\ln \equiv \log_e$; $\log \equiv \log_{10}$.

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Some math. you should remember

Given an integer $n > 0$ and a real $a > 1$ and $a \neq 0$:

- Arithmetic summation: $\sum_{i=0}^n i = \frac{n(n+1)}{2}$.
- Geometric summation: $\sum_{i=0}^n a^i = \frac{1-a^{n+1}}{1-a}$.

Logarithms and Exponents: For $a, b, c \in \mathbb{R}^+$,

- $\log_b a = c \Leftrightarrow a = b^c \Rightarrow \log_b 1 = 0$
- $\log_b ac = \log_b a + \log_b c$, $\log_b a/c = \log_b a - \log_b c$.
- $\log_b a^c = c \log_b a \Rightarrow c^{\log_b a} = a^{\log_b c} \Rightarrow 2^{\log_2 n} = n$.
- $\log_b a = \log_c a / \log_c b \Rightarrow \log_b a = \Theta(\log_c a)$

Stirling: $n! = \sqrt{2\pi n}(n/e)^n + O(1/n) + \gamma \Rightarrow n! + \omega((n/2)^n)$.

n -Harmonic: $H_n = \sum_{i=1}^n 1/i \sim \ln n$.

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Graphs

See for ex. Chapter 3 of Dasgupta, Papadimitriou, Vazirani (DPV).

Graph: $G = (V, E)$, where V is the set of vertices, $n = |V|$, and $E \subset V \times V$ is the set of edges, $m = |E|$,

- Graphs: *undirected graphs (graphs)* and *directed graphs (digraphs)*
- The **degree** of v ($d(v)$) is the number of edges which are incident to v .
- A **clique** on n vertices (K_n) is a **complete graph** (with $m = n(n - 1)/2$).
- A undirected G is said to be connected if there is a path between any two distinct vertices.
- If G is connected, then $m \geq n - 1$.

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Directed graphs

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- Edges are directed.
- The connectivity concept in digraphs is the so called **strong connectivity**: There is a directed path between any two vertices.
- In a digraph $m \leq n(n - 1)$.

Density of a graph

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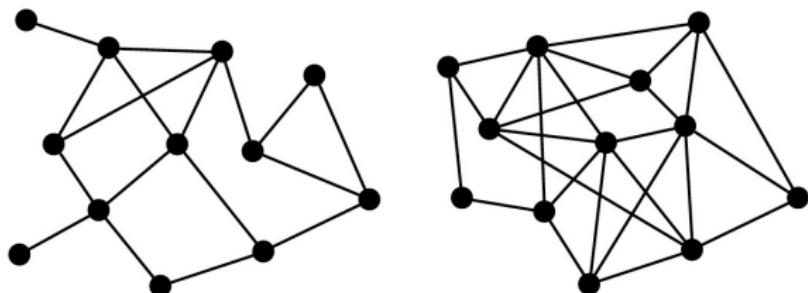
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A G with n vertices is said to be **dense** when $m = \Theta(n^2)$.
When $m = o(n^2)$, G is said to be **sparse**.



Common graph's data structures

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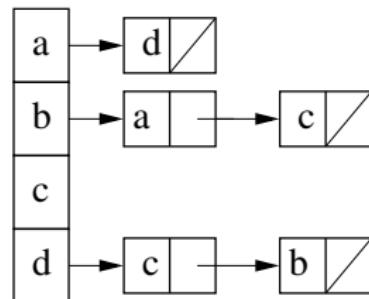
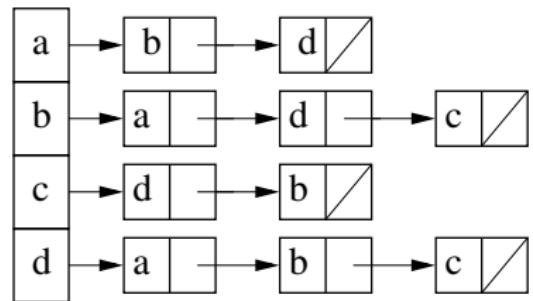
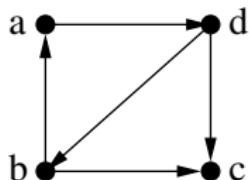
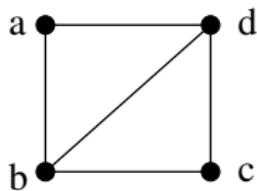
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Let G be a graph with $V = \{1, 2, \dots, n\}$.

Adjacency list

Adjacency matrix

Adjacency list



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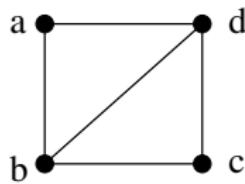
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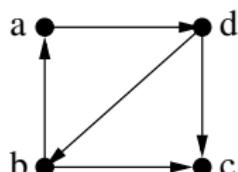
Adjacency matrix

Given G with $|V| = n$ define its **adjacency matrix** as the $n \times n$ matrix:

$$A[i,j] = \begin{cases} 1 & \text{if } (i,j) \in E, \\ 0 & \text{if } (i,j) \notin E. \end{cases}$$



$$\begin{matrix} & \begin{matrix} a & b & c & d \end{matrix} \\ \begin{matrix} a \\ b \\ c \\ d \end{matrix} & \begin{pmatrix} 0 & 1 & 0 & 1 \\ 1 & 0 & 1 & 1 \\ 0 & 1 & 0 & 1 \\ 1 & 1 & 1 & 0 \end{pmatrix} \end{matrix}$$



$$\begin{matrix} & \begin{matrix} a & b & c & d \end{matrix} \\ \begin{matrix} a \\ b \\ c \\ d \end{matrix} & \begin{pmatrix} 0 & 0 & 0 & 1 \\ 1 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 1 & 1 & 0 \end{pmatrix} \end{matrix}$$

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Adjacency matrix

- If G is undirected, its adjacency matrix $A(G)$ is symmetric.
- If A is the adjacency matrix of G , then A^2 gives, for $i, j \in V$, whether there is a path between i and j in G , with length 2.
For $k > 0$, A^k indicates if there is a path with length k in G .
- If G has weights on edges, i.e. $w_{i,j}$ for each $(i,j) \in E$, $A(G)$ has w_{ij} in $a_{i,j}$.
- Adjacency matrix allows the use of tools from linear algebra.

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Comparison between the matrix and the list DS

- The adjacency list uses a register per vertex and two per edge. As each register needs 64 bits, then the space to represent a graph is $\Theta(n + m)$.
- The use of the adjacency matrix needs n^2 bits ($\{0, 1\}$), so for an unweighted graph G , we need $\Theta(n^2)$ bits.
- For weighted G , we need $64n^2$ bits (assuming weights are reasonably “small”).
- In general, for unweighted dense graphs, the adjacency matrix is better, otherwise the adjacency list is a shorter representation.

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Complexity issues between matrix and list DS

- Adding a new edge to G : In both data structures we need $\Theta(1)$.
- Edge query, for u and v in $V(G)$, $(u, v) \in E(V)$?:
 - For matrix representation: $\Theta(1)$.
 - For list representation: $O(n)$.
- Explore all neighbours of vertex v :
 - For matrix representation: $\Theta(n)$
 - For list representation: $\Theta(|d(v)|)$
- Erase an edge in G : The same as Edge query.
- Erase a vertex in G :
 - For matrix representation: $\Theta(n)$.
 - For list representation: $O(m)$.

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Searching a graph: Breadth First Search

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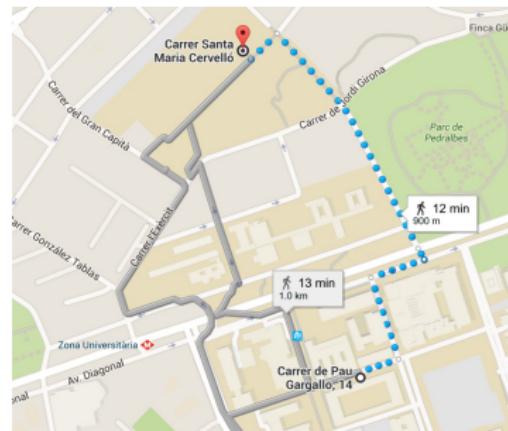
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- 1 Start with vertex v , visit v and all its neighbors.
- 2 Then, the non-visited neighbors of visited ones.
- 3 Repeat until all vertices are visited.



BFS use a QUEUE, (FIFO) to keep the neighbors of visited vertices.

Recall that vertices are labeled to avoid visiting them more than once.

Searching a graph: Depth First Search

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explore

- 1 From current vertex, move to a neighbor.
- 2 Until you get stuck.
- 3 Then backtrack till new place to explore.



DFS use a STACK, (LIFO)

Time Complexity of DFS and BFS

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For graphs given by adjacency lists:

$$O(|V| + |E|)$$

For graphs given by adjacency matrix:

$$O(|V|^2)$$

Therefore, both procedures can be implemented in linear time
with respect to the size of the input graph.

Connected components in undirected graphs

A connected component is a maximal connected subgraph of G .

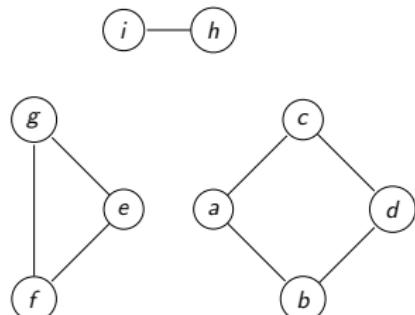
A connected graph has a unique connected component.

Connected Components

INPUT: undirected graph G

QUESTION: Find all the connected components of G .

To find connected components in G use DFS and keep track of the set of vertices visited in each **explore** call.



The problem can be solved in $O(|V| + |E|)$.

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Strongly connected components in a digraph

A digraph $G = (V, E)$, is *strongly connected*, if for all $u, v \in V$, there are paths $u \rightarrow v$ and $v \rightarrow u$.

A strongly connected component is a maximal strongly connected graph.

Strongly Connected Components

INPUT: digraph G

QUESTION: Find the strongly connected components of G .

Kosaraju-Sharir's algorithm: Uses BFS (twice). Complexity $T(n) = O(|V| + |E|)$

Tarjan's algorithm: Based in using DFS. Complexity

$T(n) = O(|V| + |E|)$

Both algorithms are optimal, i.e. linear time, but in practice Tarjan's algorithm is easier to implement.

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Strongly connected components in a digraph

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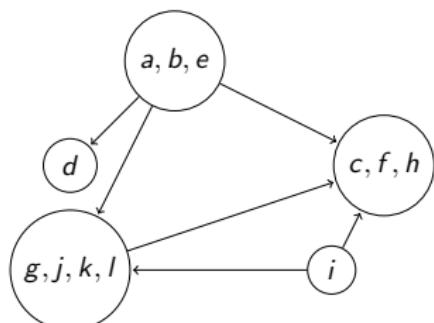
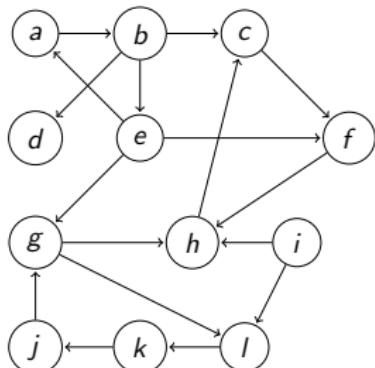
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A nice property: For every digraph, the graph on its strongly connected components is *acyclic*.



The classes P and NP

- Recall that a problem belongs to the class P if there exists an algorithm that is polynomial in the worst-case analysis, (for the worst input given by a malicious adversary)
- A problem given in **decisional form** belongs to the class NP **non-deterministic polynomial time** if given a certificate of a solution we can verify in polynomial time that indeed the certificate is a valid solution to the problem in decisional form and those certificates have polynomial size
- It is easy to see that $P \subseteq NP$, but it is an open problem to prove that $P=NP$ or that $P \neq NP$.
- The class NP-complete is the class of most difficult problems in decisional form that are in NP. Most difficult in the sense that if one of them is proved to be in P then $P=NP$.

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Beyond worst-case analysis

- Under the hypothesis that $P \neq NP$, if the decision version of a problem is NP-complete, then the optimization problem will require at least exponential time, for some inputs.
- The classification of a problem as NP-complete is a case of worst-case analysis, and for many problems the "expensive inputs" are few, and far from practical typical inputs. We will see some examples through the course.
- Therefore, there are alternative ways to get in practice, solutions for NP-complete problems, with the use of alternative algorithmic techniques, as approximation (we will see some examples), heuristics and self-learning algorithms, that are deferred to other courses.

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A powerful tool to solve problems: Reductions

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You have been introduced in previous courses to the concept of reduction between decision problems, to define the class NP-complete.

We have to extend the concept to function problems.

Reductions

Given problems A and B , assume we have an algorithm \mathcal{A}_B to solve the problem B on any input y .

A polynomial time **reduction** $A \leq B$ is a pair of polynomial time functions (f, g) such that

- f maps any input x to A , in polynomial time to an input $f(x)$ to problem B in such a way that x has a valid solution for A iff $f(x)$ has a valid solution for B .
- g maps solutions to $f(x)$ into solutions to x .

Therefore if we have that $A \leq B$, as there is an algorithm \mathcal{A}_B to solve problem B in polynomial time, then we have an algorithm \mathcal{A}_A , for any input x of A : Compute $g(\mathcal{A}_B(f(x)))$, that runs in polynomial time.

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The VERTEX COVER problem

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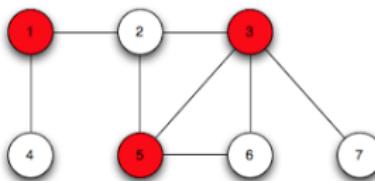
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VERTEX COVER: Given a graph $G = (V, E)$ with $|V| = n, |E| = m$, find the minimum set of vertices $S \subseteq V$ such that it covers every edge of G .

Example:



The VERTEX COVER problem is known to be in NP-hard.

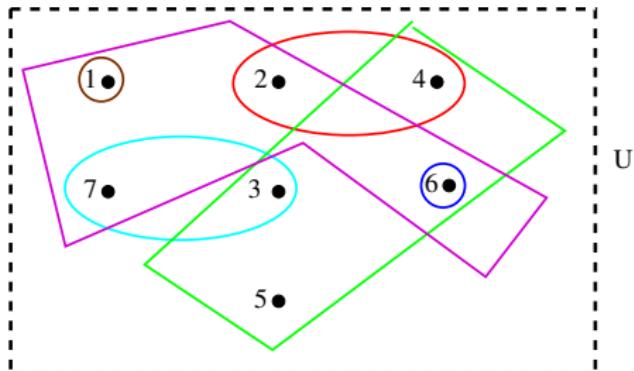
The SET COVER problem

SET COVER: Given a set U of m elements, a collection $S = \{S_1, \dots, S_n\}$ where each $S_i \subseteq U$, select the minimum number of subsets in such a way that their union is equal to U .

There is a weighted version of the problem, but this simpler version already is NP-hard.

Example: Given $U = \{1, 2, 3, 4, 5, 6, 7\}$ ($m = 7$), with $S_a = \{3, 7\}$, $S_b = \{2, 4\}$, $S_c = \{3, 4, 5, 6\}$, $S_d = \{6\}$, $S_e = \{1\}$, $S_f = \{1, 2, 6, 7\}$.

Solution: $\{S_c, S_f\}$



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SET COVER

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The VERTEX COVER problem is a special case of the SET COVER problem. As a model, the SET COVER has important practical applications.

To understand the computational complexity of SET COVER it is important to understand first the complexity of special cases as VERTEX COVER.

VERTEX COVER \leq SET COVER

Given an input to VERTEX COVER $G = (V, E)$, of size $|V| + |E| = n + m$ we want to construct in polynomial time on $n + m$ a specific input $f(G) = (U, S)$ to SET COVER such that if there exist a polynomial algorithm \mathcal{A} to find a min. vertex cover in G , then $\mathcal{A}(f(G))$ is an efficient algorithm to find an optimal solution to set cover.

REDUCTION f :

- Consider U as the set E of edges.
- For each vertex $i \in V$, S_i is the set of edges incident to i .
Therefore $|S| = n$ and for each S_i , $|S_i| \leq m$.
- The cost of the reduction from G to (U, S) is $O(n + m)$

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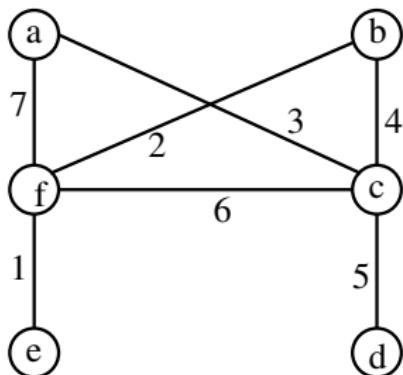
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Example for the reduction

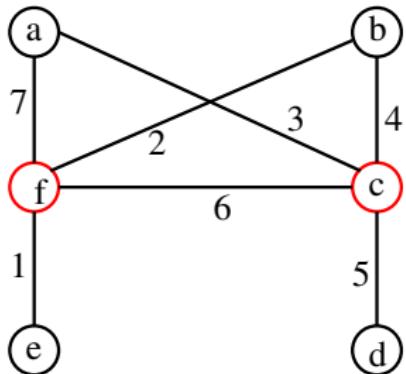
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$$\begin{aligned} U &= \{1, 2, 3, 4, 5, 6, 7\} \\ S &= \{S_a, S_b, S_c, S_d, S_e, S_f\} \\ \overset{f}{\overbrace{S_a}} &= \{3, 7\}, S_b = \{3, 7\}, \\ \overset{f}{\overbrace{S_c}} &= \{3, 4, 5, 6\}, S_d = \{5\}, \\ S_e &= \{1\}, S_f = \{1, 2, 6, 7\}. \end{aligned}$$

Example for the reduction

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$$\begin{aligned}U &= \{1, 2, 3, 4, 5, 6, 7\} \\S &= \{S_a, S_b, S_c, S_d, S_e, S_f\} \\S_a &= \{3, 7\}, S_b = \{3, 7\}, \\&\stackrel{f}{\Rightarrow} S_c = \{3, 4, 5, 6\}, S_d = \{5\}, \\S_e &= \{1\}, S_f = \{1, 2, 6, 7\}.\end{aligned}$$

If there is an algorithm to solve the SET COVER for G , the same algorithm apply to $(U, S) = f(G)$ will yield a solution for VERTEX COVER on input (U, V) .

But both VERTEX COVER and SET COVER are known to be NP-hard.

The divide-and-conquer strategy.

- 1 Break the problem into smaller subproblems,
- 2 recursively solve each problem,
- 3 appropriately combine their answers.



Julius Caesar (I-BC)
"Divide et impera"

Known Examples:

- Binary search
- Merge-sort
- Quicksort
- Strassen matrix multiplication

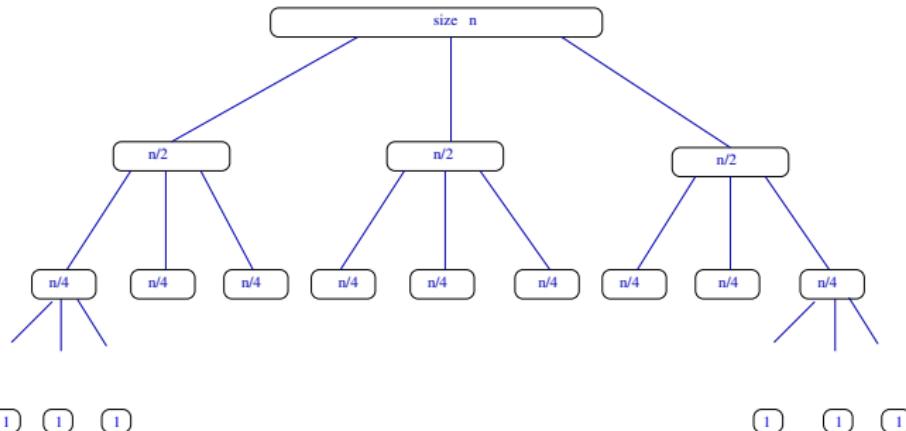


J. von Neumann
(1903-57) Merge sort

Recurrences Divide and Conquer

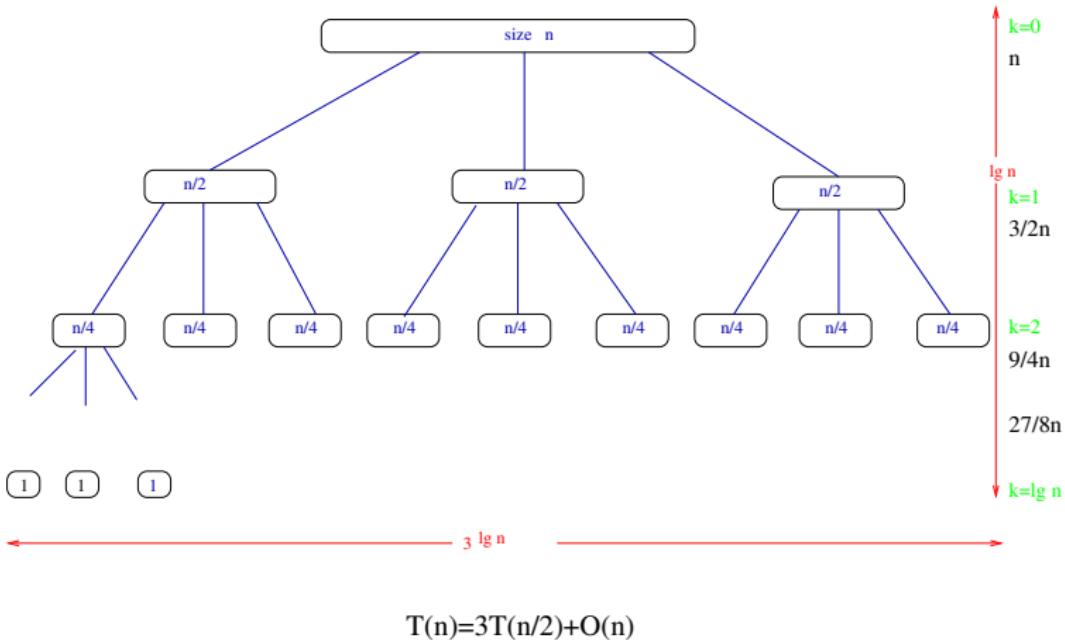
$$T(n) = 3T(n/2) + O(n)$$

The algorithm under analysis divides input of size n into 3 subproblems, each of size $n/2$, at a cost (of dividing and joining the solutions) of $O(n)$



$$T(n) = 3T(n/2) + O(n).$$

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At depth k of the tree there are 3^k subproblems, each of size $n/2^k$.

For each of those problems we need $O(n/2^k)$ (splitting time + combination time).

Therefore, for some constant c , the cost at depth k is:

$$3^k \times \left(\frac{n}{2^k}\right) = \left(\frac{3}{2}\right)^k \times c n.$$

with max. depth $k = \lg n$, so $T(n)$ is

$$\left(1 + \frac{3}{2} + \left(\frac{3}{2}\right)^2 + \left(\frac{3}{2}\right)^3 + \cdots + \left(\frac{3}{2}\right)^{\lg n}\right) c n$$

From $T(n) = c n \left(\sum_{k=0}^{\lg n} \left(\frac{3}{2}\right)^k \right)$,

We have a **geometric series** of ratio $3/2$, starting at 1 and ending at $\left(\left(\frac{3}{2}\right)^{\lg n}\right) = \frac{n^{\lg 3}}{n^{\lg 2}} = \frac{n^{1.58}}{n} = n^{0.58}$.

As the series is increasing, $T(n)$ is dominated by the last term:

$$T(n) = c n \left(\frac{n^{\lg 3}}{n} \right) = O(n^{1.58}).$$

A basic Master Theorem

There are several versions of the Master Theorem to solve D&C recurrences. The one presented below is taken from DPV's book.

Theorem (DPV-2.2)

If $T(n) = aT(\lceil n/b \rceil) + O(n^d)$ for constants $a \geq 1, b > 1, d \geq 0$, then has asymptotic solution:

$$T(n) = \begin{cases} O(n^d), & \text{if } d > \log_b a, \\ O(n^d \lg n), & \text{if } d = \log_b a, \\ O(n^{\log_b a}), & \text{if } d < \log_b a. \end{cases}$$

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Master Theorems

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This basic Master Theorem does not provide always exact bounds

A different one can be found in CLRS's book providing exact bounds but leaving cases outside.

For stronger versions:

Akra-Bazi Theorem: <https://courses.csail.mit.edu/6.046/spring04/handouts/akrabazzi.pdf>

Salvador Roura Theorems

<http://www.lsi.upc.edu/~diaz/RouraMT.pdf>