# Greedy algorithms 

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## Basic Idea

- We consider optimization problems. In $F$, the set of feasible solutions we must find an optimal element, i.e. an element where $c: F \rightarrow \mathbb{R}$, the objective function takes minimal/maximal value.
- Very often the feasible solutions are subsets of a universe.
- An easy algorithm paste the elements of $U$ (or the "surviving" elements of $U$ ). The MOST PROMISING element will be chosen as the next element of the output. After pasting all elements of $U$ we will have an output.


## Basic Idea: Why greedy?

- Taking the most promising element is a natural decision. The essence of greediness is that the above algorithm never overrules previous decisions. In spite of being a promising element at some point, later on we might realize that choosing that element is not a wise decision. A greedy algorithm do not step back.
- Greedy algorithms are very simple to implement. They are very fast. Unfortunately very often they are not able to guarantee that the output is optimal.
- The above description is not a mathematical definition. It is a scheme, that very often leads to good algorithms. Sometimes (rarely) greediness makes us to be able to find the optimal solution.
- Let us see an example.

Break


## The minimal cost spanning tree problem

## The problem

Given a connected graph, for each edge we have a positive cost (c: $E(G) \rightarrow \mathbb{R}_{++}$). This cost function can be naturally extended to subsets of $E(G)$ (the cost of an edge set is the sum of the costs of its elements).
Find a cheapest spanning tree of the input graph (the tree is considered as a set of edges).

The following algorithm, first described by Kruskal is a ,, prototype" of the greedy algorithm design.

## Kruskal's algorithm

## Kruskal's algorithm (1956)

(SORTING STEP) Sort the edges of the input graph in ascending order of cost. Let $E(G): e_{1}, e_{2}, \ldots, e_{m}$, i.e. $e_{1}$ is the cheapest edge, $e_{m}$ is the most expensive edge $\left(c\left(e_{1}\right) \leq c\left(e_{2}\right) \leq \ldots \leq c\left(e_{m}\right)\right)$.
(INITIALIZATION) Let $F$ be the set of edges selected so far. At the beginning of the algorithm $F=\emptyset$. // During the algorithm we only take care that $F$ is an edge set without cycle. (TESTS) In the $i^{\text {th }}$ step, we examine $e_{i}$. If $F \cup\left\{e_{i}\right\}$ is cycle-free, then we extend $F: F \leftarrow F \cup\left\{e_{i}\right\}$. If $F \cup\left\{e_{i}\right\}$ contains a cycle, then we do not change $F$.
(OUTPUT) After examining the last edge, we announce the current $F$ as the output.

## The main theorem

One can say: all our choices are the best decision at the time. Later some of the edges have been discarded. After that it is possible that we need to throw away an edge. This is a questionable choice. It was based on the fact that the previously selected edges form a part of the output. If our previous decisions are overruled, then we could have chosen to use the currently discarded edge (the cheapest edge of the remaining edge set). The cost of the calculated spanning tree cannot simply be compared to the tree calculated above.

Despite the huge question mark above, the calculated tree is optimal.

## Theorem (Kruskal's Theorem)

The output of the above algorithm is a minimum cost spanning tree of the input graph.

## Initial notations

- $n=|V(G)|$.
- Let $e_{1}, e_{2}, \ldots, e_{\ell}$ be the elements of the output in the order of their selection. Specifically

$$
c\left(e_{1}\right) \leq c\left(e_{2}\right) \leq \ldots \leq c\left(e_{\ell-1}\right) \leq c\left(e_{\ell}\right)
$$

- Let $F$ be an arbitrary spanning tree. We list the edges in increasing order of cost: $f_{1}, \ldots, f_{n-1}$, i.e.

$$
c\left(f_{1}\right) \leq c\left(f_{2}\right) \leq \ldots \leq c\left(f_{n-1}\right)
$$

- Based on the connectivity of the input graph it is easy to see that Kruskal's algorithm computes a spanning tree, i.e. $\ell=n-1$.


## Strong form of Kruskal's Theorem

Theorem (Strong form of Kruskal's Theorem)
For $i=1,2, \ldots, n-1$ we have

$$
c\left(e_{i}\right) \leq c\left(f_{i}\right)
$$

## The Main Lemma

## Main Lemma

Let $F, F^{\prime}$ be two cycle-free edge sets over $V$. Suppose that $|F|<\left|F^{\prime}\right|$. Then there is an edge $e$ in $F^{\prime}-F$ such that $F \cup\{e\}$ is also cycle-free.

- Let $G_{F}$ be the cycle-free graph (forest) with vertex set $V$ and edge set $F$.
- We know that $G_{F}$ has $n-|F|$ components: $c\left(G_{F}\right)=n-|F|$.
- Similarly let $G_{F^{\prime}}$ be the cycle-free graph (forest) with vertex set $V$ and edge set $F^{\prime}$.
- We know that $G_{F^{\prime}}$ has $n-\left|F^{\prime}\right|$ components:
$c\left(G_{F^{\prime}}\right)=n-\left|F^{\prime}\right|<n-|F|=c\left(G_{F}\right)$.
- This can only be imagined if there is an edge $e \in F^{\prime}$ connecting two different components of $G_{F}$.
- Hence $e \notin F$ and $F \cup\{e\}$ is cycle-free.


## Strong form of Kruskal's Theorem: The proof

- Induction on $i$.
- $c\left(e_{1}\right) \leq c\left(f_{1}\right)$.
- Assume that for $i(<n-1)$

$$
c\left(e_{1}\right) \leq c\left(f_{1}\right), c\left(e_{2}\right) \leq c\left(f_{2}\right), \ldots, c\left(e_{i}\right) \leq c\left(f_{i}\right)
$$

- $F:=\left\{e_{1}, \ldots, e_{i}\right\}, F^{\prime}:=\left\{f_{1}, \ldots, f_{i}, f_{i+1}\right\}$. Apply the Main Lemma. $\rightarrow f \in F^{\prime}$.
- The Kruskal's algorithm tests $e_{i+1}$ not later than $f$.

$$
c\left(e_{i+1}\right) \leq c(f) \leq c\left(f_{i+1}\right)
$$

## The analysis of Kruskal's algorithm

- The cost of the sorting step is

$$
\mathcal{O}(|E| \log |E|)
$$

- The remaining part of algorithm is $|E|$ testing on an edge set. In each case the size of the edge set is is at most $|V|$. The cost of a test is $\mathcal{O}(|V|)$.
- Cost of the complete run

$$
\mathcal{O}(|E| \log |E|)+|E| \cdot \mathcal{O}(|V|)=\mathcal{O}(|E| \cdot|V|)
$$

- We have performed an analysis of a naive implementation. There are more clever solutions.

Break


## The basic question

## The computational problem

Given
(i) $\vec{G}$ directed graph,
(ii) $\ell: E(\vec{G}) \rightarrow \mathbb{R}_{++}$length function,
(iii) $s, t$ two distinguished.

Determine the distance of $s$ and $t$.

## Refreshing memory

## Definition: Walk in a directed graph

$\overrightarrow{u v}$-walk in $\vec{G}$ :

$$
\overrightarrow{\mathcal{S}}: u=w_{0}, \vec{e}_{1}, w_{1}, \vec{e}_{2}, \ldots, w_{L-1}, \vec{e}_{L}, w_{L}=v
$$

where $w_{i} \in V(i=0,1, \ldots, L), \vec{e}_{i} \in E(i=1, \ldots, L), e_{i}$ is an outgoing edge from $w_{i-1}$, and ingoing edge in $w_{i}(i=1, \ldots, L)$.

In the case of the existence of an $\overrightarrow{u v}$-walk we say that $v$ is reachable from $u$.
The graph theoretical length of $\overrightarrow{\mathcal{S}}$, an $\overrightarrow{u v}$-walk is $L$.

## Refreshing memory (cont'd)

Length of a walk in a weighted graph
The (weighted) length of $\overrightarrow{\mathcal{S}}$, an $\overrightarrow{u v}$-walk is

$$
\sum_{i=1}^{L} \ell\left(\vec{e}_{i}\right)
$$

## Definition: The distance of two vertices in a weighted graph

$d(u, v)$ denotes the distance of two vertices $u$ and $v$, that is the minimal length among the $\overrightarrow{u v}$-walks. // $d(u, v)=\infty$ is possible!

## Observation

The shortest $\overrightarrow{u v}$-walk will be a path.

## Initial remarks

- We assume the our graph has no loop. We assume that for any two vertices there is at most one edge from $u$ to $v$. I.e. we assume that $\vec{G}$ is simple (in directed sense).


## The case of graph theoretical distance: unweighted case

- We solve a harder problem. The input will be ( $\vec{G}, s$ ). We determine $S$, the set of vertices that are reachable from $s$.
- The set of reachable vertices will be $S=S_{0} \dot{\cup} S_{1} \dot{\cup} \ldots \dot{\cup} S_{L}$, where $S_{i}$ contains exactly those vertices that are at distance $i$ from $s$.
- L denotes the length of longest path starting at $s$.
- Each edge between $S$ and $\bar{S}=V(G)-S$ are oriented from $\bar{S}$ to $S$. This property will be a proof of the fact that the elements of $\bar{S}$ are not reachable from $s$.


## Breadth first search algorithm

## Breadth first search

(I) // Inicialization // Let $S_{0}=\{s\}$.
(E) // Extension cycle // While $S_{i} \neq \emptyset$

$$
\begin{array}{r}
S_{i+1}=\left\{x \in V(G)-\left(S_{0} \cup \ldots \cup S_{i}\right): \text { there is } \sigma \in S_{i},\right. \\
\text { such that } \overrightarrow{\sigma x} \in E(G)\}
\end{array}
$$

- $i \leftarrow i+1$.
end-while


## The correctness of breadth first search algorithm

## Theorem

The above algorithm is correct. I.e.
(i) In the case of $x \in S_{i}$ the graph theoretical distance of $s$ and $x$ is $i$.
(ii) If $x \notin S$, then $x$ is not reachable from $s$.

## Breadth first search tree

See the following modification of the algorithm:

## Breadth first search tree (modification)

$\left(I^{*}\right) F:=\emptyset$
$\left(\mathrm{E}^{+}\right)$In cycle $(\mathrm{E})$, when we insert $x$ into $S_{i+1}(i \geq 0)$, then the algorithm search for a suitable $\sigma \in S_{i}$ and find one: $\sigma_{x} \in S_{i}$.
// Suitable means that $\overrightarrow{\sigma_{x} x}$ is an edge.
(Extending $F$ ) $F \leftarrow F \cup\left\{\overrightarrow{\sigma_{x} x}\right\}$

## Theorem

In the graph $\left.G\right|_{S}$ the edge set $F$ will be the edge set of a spanning tree $\mathcal{F}$. $\mathcal{F}, s)$ is a rooted tree.
For each vertex $x \in S$ there is exactly one $\overrightarrow{s x}$-path in $\mathcal{F}$. This one of the (graph theoretically) shortest $\overrightarrow{s x}$-path.

## Weighted case: The basic idea

- We also assume that all vertices are reachable from $s$.
- Assume that for a vertex set $S \subset V(\vec{G})$ we have
(o) $s \in S$,
(i) for each $v \in S$ we have the shortest $\overrightarrow{s v}$ path, and that is inside $S$.
- The greedy algorithm chooses the most promising vertex from $\bar{S}:=V(G) \backslash S$. For this we assume that the following information is also computed:
(ii) For each $v \in \bar{S}$ we have the shortest $\overrightarrow{s v}$ path, and that is inside $S$ except the last $\overrightarrow{v^{-} v}$ step $\left(v^{-} \in S\right)$.
- One should think about the promised information as a labeling, $c: V(\vec{G}) \rightarrow \mathbb{R} \cup\{\infty\}$, of the vertices.


## Dijkstra's algorithm

## Dijkstra's algorithm

(I) // Inicialization
$S=\{s\}, c(s)=0$.
In the case of $v \notin S, \overrightarrow{s v} \in E$ we have $c(v)=\ell(\overrightarrow{s v})$.
In the case of $v \notin S, \overrightarrow{s v} \notin E$ we have $c(v)=\infty$.
(E) // Extension (until $\bar{S} \neq \emptyset$ )
$m \in \bar{S}$ with smallest label in $\vec{S}$.
$S \leftarrow S \cup\{m\}$.
$/ / \bar{S} \leftarrow \bar{S} \backslash\{m\}$.
(U) // Update

We might change the label of $n \in \bar{S}$ in the case of $\overrightarrow{m h} \in E$ :

$$
c_{\text {new }}(n)=\min \left\{c_{\text {old }}(n), c(m)+\ell(\overrightarrow{m h})\right\}
$$

If $t \notin S$ (or $\bar{S} \neq \emptyset$ ) back to (E).

## Correctness of the algorithm

## Theorem

In each update step Dijkstra's algorithm computes a correct labeling. I.e.
(i) The label of $v \in S$ is the length of shortest $\overrightarrow{s v}$ path. Furthermore this shortest length can be realized by a $\overrightarrow{s v}$-path inside $S$.
(ii) The label of $v \in \bar{S}$ is the length of shortest $\overrightarrow{s v}$ path, that leaves $S$ only when it makes the last step.

- We prove by mathematical induction on $|S|$. Case $|S|=1$ is straight forward.
- For the induction step we need to prove two claims:
(i) The label of $m$ is correct,
(ii) The label of $x \in \bar{S}$ is correct.


## Analysis of the algorithm

- $\bar{S}_{\text {fin }}$ denotes the set of nodes from $\bar{S}$ with finite label.
- Extension step: Delete the node with minimum label from $\bar{S}_{\text {fin }}$. The number of extension steps is at most $|V|-1$.
- Label updates steps: Some vertex with label " $\infty$ " enters $\bar{S}_{\text {fin }}$. Some labels in $\bar{S}_{\text {fin }}$ will be decreased by a value $\delta(>0)$. Computation for label updates is needed for at most $|E|$ times.
- The total number of steps is

$$
\mathcal{O}\left(|V|^{2}+|E|\right)
$$

- The above argument followed a naive implementation. There are cleverer ways to implement Dijkstra's high level description.


## Final remarks

- Every label has a corresponding edge, that is responsible for its value.
- Changing the value of the label/updating: changing the "responsible" node too.
- If we keep track of these edges responsible for the actual value we will obtain a rooted, directed spanning tree of the original graph (see the computation of breadth first search tree). This tree maintains/contains for each vertex a shortest path leading to that vertex.

Break


## Coding texts

## Definition of Text

Let $\Sigma$ be a finite alphabet. The elements of $\Sigma$ are called characters. A text is a finite sequence of characters $\left(\in \Sigma^{\star}\right)$.

The length of a text is the number of characters in the text.

## Coding texts

An algorithm/function $c: \Sigma^{\star} \rightarrow\{0,1\}^{\star}$

$$
+
$$

decoding algorithm $d:\{0,1\}^{\star} \rightarrow \Sigma^{\star}$.

Character based coding: $c_{0}: \Sigma \rightarrow\{0,1\}^{\star}$ coding of characters. The code of a text is obtained by "putting together" the codes of its characters.

Fixed-length codes, Example: ASCII (1972) (source: wiki)

| $b_{7} b_{6} b_{5}$ |  |  |  |  | ${ }^{0} \mathrm{O}$ | $0^{0}$ | ${ }^{0} 1_{0}$ | $0^{0} 1$ | ${ }^{1} 0$ | ${ }^{1} 0$ | $10$ | 11. |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\underbrace{1} \underbrace{b_{4}}_{\downarrow}$ | $\left[\begin{array}{c} b_{3} \\ d \end{array}\right.$ | $\begin{gathered} \mathrm{b}_{2} \\ \downarrow \end{gathered}$ | $\begin{gathered} \mathrm{b}_{1} \\ \downarrow \end{gathered}$ |  | 0 | 1 | 2 | 3 | 4 | 5 | 6 | 7 |
| 0 | 0 | 0 | 0 | 0 | NUL | DLE | SP | 0 | @ | P | , | p |
| 0 | 0 | 0 | 1 | 1 | SOH | DCI | ! | 1 | A | Q | a | q |
| 0 | 0 | 1 | 0 | 2 | STX | DC2 | " | 2 | B | R | b | $r$ |
| 0 | 0 | 1 | 1 | 3 | ETX | DC3 | \# | 3 | C | S | C | S |
| 0 | 1 | 0 | 0 | 4 | EOT | DC4 | \$ | 4 | D | T | d | $\dagger$ |
| 0 | 1 | 0 | 1 | 5 | ENQ | NAK | \% | 5 | E | U | e | $u$ |
| 0 | 1 | 1 | 0 | 6 | ACK | SYN | Q | 6 | F | V | $f$ | $v$ |
| 0 | 1 | 1 | 1 | 7 | BEL | ETB | , | 7 | G | W | g | W |
| 1 | 0 | 0 | 0 | 8 | BS | CAN | $($ | 8 | H | X | h | x |
| 1 | 0 | 0 | 1 | 9 | HT | EM | $)$ | 9 | I | Y | i | y |
| 1 | 0 | 1 | 0 | 10 | LF | SUB | * | : | $J$ | Z | j | 2 |
| 1 | 0 | 1 | 1 | 11 | VT | ESC | + | ; | K | [ | k |  |
| 1 | 1 | 0 | 0 | 12 | FF | FS | , | < | L | 1 | 1 | 1 |
| 1 | 1 | 0 | 1 | 13 | CR | GS | - | = | M | ] | m | , |
| 1 | 1 | 1 | 0 | 14 | SO | RS | - | $>$ | N | ヘ | $n$ | $\sim$ |
| 1 | 1 | 1 | 1 | 15 | SI | US | / | ? | 0 | - | 0 | DEL |

Variable-length codes, Example: Morse code (1837-44) (source:wiki)

International Morse Code

1. The length of a dot is one unit.
2. A dash is three units.
3. The space between parts of the same letter is one unit.
4. The space between letters is three units.

5 . The space between words is seven units.


## Variable-length codes without comma: Prefix codes

## Definition

Rooted binary plane tree.
Leaf of a rooted binary plane tree.

## Definition: Prefix tree for $\Sigma$

Let $(T, r)$ be a rooted binary plane tree. Let $L$ be the set of leaves of $(T, r)$. pause $(T, r, \ell)$ is a prefix tree for $\Sigma$, iff $\ell: \Sigma \rightarrow L$ is a bijection.

The coding of the characters based on a prefix tree

$$
k(\in \Sigma) \mapsto \text { labels of the } r-\ell(k) \text { path in } T
$$

## The fundamental question

## Problem

Given $\Sigma$ alphabet and a text $\tau(\rightarrow$ probability distribution over $\Sigma /$ frequency table $\left(\in \mathbb{N}^{\Sigma}\right)$ ). Find a prefix tree over $\Sigma$, that minimize the lenth of the code of $\tau$.

## Huffman's algorithm: Basic idea

- We consider the the characters as a one-node prefix trees. So we start with $|\Sigma|$ prefix trees.
- We choose two trees and merge them.Merging: We introduce a new root-vertex. Its two children are the roots of the two merged tree.
- Using merging steps we compute the output (equiv greediness).
- The initial trees have frequency values. During the algorithm each tree has an f-vaule: The SUM of the frequencies assigned to its leaves.
- ??? HOW TO CHOOSE THE TWO TREES TO MERGE ???
- The natural idea: choose the two trees with lowest frequencies.


## Huffman's algorithm

## Huffman's algorithm (1951)

(INITIALIZATION) We construct a rooted binary plane tree for each character. Each tree has an $f$-value, the frequency of the corresponding character.
// During the algorithm we always have a set of prefix trees with $f$-values: $\mathcal{T}$.
(CHOICE) (until $|\mathcal{T}|>1$ ) Take the two trees with the lowest frequencies: $T_{1}, T_{2}$.
(Merge) Merge the two chosen trees $\rightarrow M\left(T_{1}, T_{2}\right)$.

$$
\mathcal{T} \leftarrow \mathcal{T}-\left\{T_{1}, T_{2}\right\} \cup\left\{M\left(T_{1}, T_{2}\right)\right\}
$$

The $f$-value of $M\left(T_{1}, T_{2}\right)$ is the some of the $f$-values of the merged two trees. Back to (CHOICE).

## The correctness of Huffman's algorithm: the main idea

## Huffman's theorem

The output of the Huffman's algorithm is a prefix tree the produces the shortest prefix coding of the input text.

Let $T$ be a prefix tree. Choose two leaves $\left(\ell, \ell^{\prime}\right)$ that are siblings on the lowest level of $T$. Let $k$ and $k^{\prime}$ be the two characters that are in the first merge step when we run Huffman's coding. Modify $T$ by changing the labels of the leaves. With at most two transpositions we obtain $T^{\prime}$ where the label of $\ell$ is $k$, and the label of $\ell^{\prime}$ is $k^{\prime}$.
Let $\beta^{\prime}$ be the code of $\tau$ based on $T^{\prime}$. Let $\beta$ be the code of $\tau$ based on $T$.

## Observation

The length of $\beta^{\prime}$ is at most the length of $\beta$.

## "this is an example of a huffman tree"



Source: wikipedia

Break


## The greedy algorithm for matchings

## Greedy algorithm for finding large matchings

(Initialization) Start with a matching $M$.
WHILE there is an edge $e \in E(G)-M$ such that $M \cup\{e\}$ is a
matching too, do
(Greedy extension step) $M \leftarrow M \cup\{e\}$.
(Halting) The actual matching is the output.
// At halting we have that each edge of $E(G) \backslash M$ is neighbors of an edge from $M$.

- There is no fear of infinite loop.
- We know that int the case of halting the output can't be augmented by extensions (by adding further edges to the output).


## Example



Figure: Our graph has four disjoint levels of equal sized vertex sets (let $n$ be the size of the levels, in our example $n=4$ ). Between two adjacent levels all possible edges are present and there are no further edges. It is possible that the greedy algorithm first chooses the yellow edges, matching the two middle levels. Then it halts. The green edges form a perfect matching.

## Analysis

## Theorem

Let $\nu_{\text {greedy }}(G)$ denote the size of the output of the greedy algorithm. Then

$$
\frac{\nu(G)}{2} \leq \nu_{\text {greedy }}(G) \leq \nu(G)
$$

The second inequality is obvious since our algorithm computes a matching.

## Analysis: the proof

- Let $M_{\text {greedy }}$ denote the output matching of the greedy algorithm.
- $L=V\left(M_{\text {greedy }}\right)$ is the set of matched vertices.
- It is obvious that $L$ is a covering vertex set, and $|L|=2 \nu_{\text {greedy }}(G)$.
- The size of $L$ gives an upper bound on the size of an arbitrary matching, hence $\nu(G) \leq|L|=2 \nu_{\text {greedy }}(G)$.


## Thank you for your attention!

