

WS 2011/12

Efficient Algorithms and Data Structures

Harald Räcke

Fakultät für Informatik
TU München

<http://www14.in.tum.de/lehre/2011WS/ea/>

Winter Term 2011/12

Part I

Organizational Matters

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- ▶ Modul: IN2003
- ▶ Name: “Efficient Algorithms and Data Structures”
“Effiziente Algorithmen und Datenstrukturen”
- ▶ ECTS: 8 Credit points
- ▶ Lectures:
 - ▶ 4 SWS
 - Mon 12:15–13:45 (Room 00.13.009A)
 - Thu 10:15–11:45 (Room 00.04.011, HS2)
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- ▶ IN0007
"Fundamentals of Algorithms and Data Structures"
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The Lecturer

- ▶ Harald RÄd'cke
- ▶ Email: raecke@in.tum.de
- ▶ Room: 03.09.044
- ▶ Office hours: (per appointment)

Tutorials

- ▶ Tutor:
 - ▶ Chintan Shah
 - ▶ chintan.shah@tum.de
 - ▶ Room: 03.09.059
 - ▶ Office hours: Wed 11:30–12:30
- ▶ Room: 00.08.038
- ▶ Time: Tue 14:14–15:45

Assessment

- ▶ In order to pass the module you need to
 1. pass an exam, and
 2. obtain at least 40% of the points in the assignment sheets.
- ▶ Exam:
 - Exam will be announced shortly.
 - There are no resources allowed, start from a blank sheet of paper (A4).
 - Answers should be given in English, but German is also accepted.

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 - ▶ Machine models
 - ▶ Efficiency measures
 - ▶ Asymptotic notation
 - ▶ Recursion
- ▶ Higher Data Structures
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 - ▶ Hashing
 - ▶ Priority queues
 - ▶ Union/Find data structures
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2 Literatur I

-  Alfred V. Aho, John E. Hopcroft, Jeffrey D. Ullman:
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-  Thomas H. Cormen, Charles E. Leiserson, Ron L. Rivest,
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-  Michael T. Goodrich, Roberto Tamassia:
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Grundlegende Algorithmen: Einführung in den Entwurf und die Analyse effizienter Algorithmen,

2. Auflage, Vieweg, 2003



Jon Kleinberg, Eva Tardos:

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3. Auflage, Addison-Wesley Publishing Company: Reading (MA), 1997

2 Literatur III



Donald E. Knuth:

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Christos H. Papadimitriou, Kenneth Steiglitz:

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2 Literatur IV



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The Algorithm Design Manual,

Springer, 1998

Part II

Foundations

3 Goals

- ▶ Gain knowledge about efficient algorithms for important problems, i.e., learn how to solve certain types of problems efficiently.
- ▶ Learn how to analyze and judge the efficiency of algorithms.
- ▶ Learn how to design efficient algorithms.

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What do you measure?

- ▶ Memory requirement
- ▶ Running time
- ▶ Number of comparisons
- ▶ Number of multiplications
- ▶ Number of hard-disc accesses
- ▶ Program size
- ▶ Power consumption
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How do you measure?

- ▶ Implementing and testing on representative inputs
 - ▶ How do you choose your inputs?
 - ▶ May be very time-consuming.
 - ▶ Very reliable results if done correctly.
 - ▶ Results only hold for a specific machine and for a specific set of inputs.
- ▶ Theoretical analysis in a specific model of computation.
 - ▶ Gives a worst case bound. (This algorithm always runs in $O(n^2)$.)
 - ▶ Usually ignores the details of the machine.
 - ▶ Can give a lower bound. (This algorithm needs at least $\Omega(n \log n)$ comparisons in the worst case.)

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- ▶ Theoretical analysis in a specific model of computation.
 - ▶ Gives a worst case bound. The algorithm always runs in $O(n^2)$.
 - ▶ Usually, we are interested in the average case performance.
 - ▶ Theoretical analysis is often used to compare algorithms in the worst case.

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Given a program P and an input I , does the algorithm always run in

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polynomial space?

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Input length

The theoretical bounds are usually given by a function $f : \mathbb{N} \rightarrow \mathbb{N}$ that maps the **input length** to the running time (or storage space, comparisons, multiplications, program size etc.).

The input length may e.g. be

the number of the input numbers

the length of the input string

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- ▶ the size of the input (number of bits)
- ▶ the number of arguments

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Suppose n numbers from the interval $\{1, \dots, N\}$ have to be sorted. In this case we usually say that the input length is n instead of e.g. $n \log N$, which would be the number of bits required to encode the input.

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1. Calculate running time and storage space etc. on a simplified, idealized model of computation, e.g. Random Access Machine (RAM), Turing Machine (TM), ...
2. Calculate number of certain basic operations: comparisons, multiplications, harddisc accesses, ...

Version 2. is often easier, but focusing on one type of operation makes it more difficult to obtain meaningful results.

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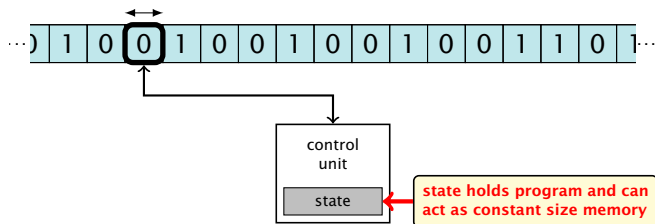
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Turing Machine

- ▶ Very simple model of computation.
- ▶ Only the “current” memory location can be altered.
- ▶ Very good model for discussing computability, or polynomial vs. exponential time.
- ▶ Some simple problems like recognizing whether input is of the form xx , where x is a string, have quadratic lower bound.

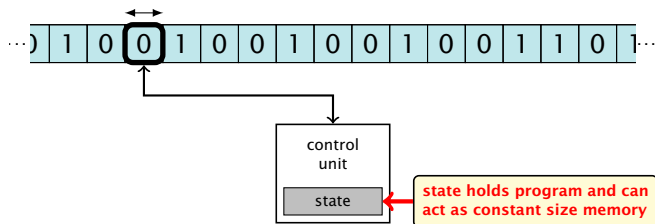
⇒ Not a good model for developing efficient algorithms.



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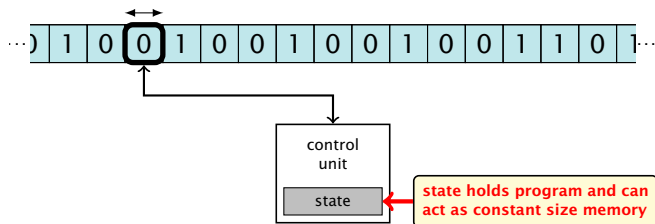
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Turing Machine

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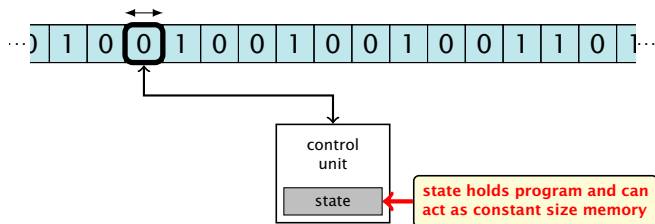
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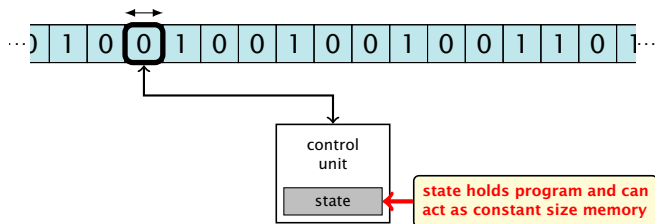
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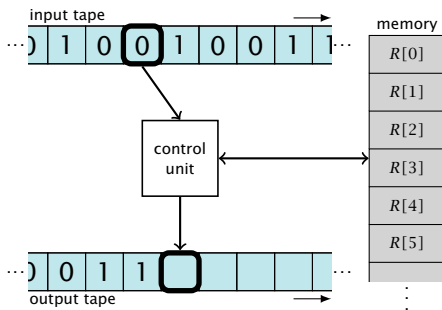
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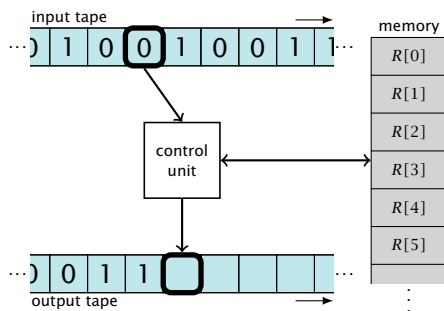
Random Access Machine (RAM)

- ▶ Input tape and output tape (sequences of zeros and ones; unbounded length).
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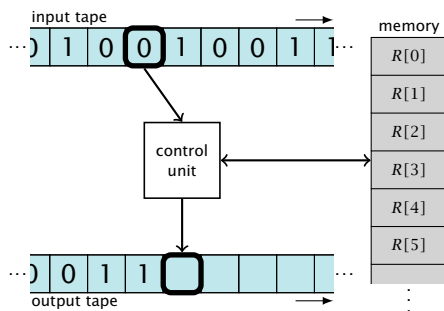
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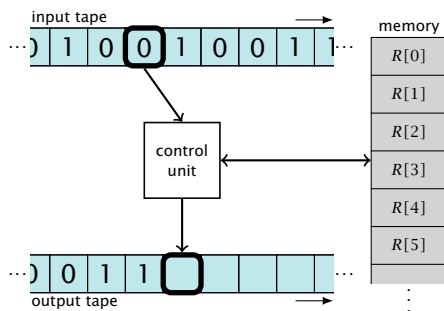
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Random Access Machine (RAM)

Operations

- ▶ input operations (input tape $\rightarrow R[i]$)
 - ▶ READ i
- ▶ output operations ($R[i] \rightarrow$ output tape)
 - ▶ WRITE i
- ▶ register-register transfers
 - ▶ $R[i] \leftarrow R[j]$
 - ▶ $R[i] \leftarrow 0$
- ▶ indirect addressing
 - ▶ $R[i] \leftarrow R[R[j]]$
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register i

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 - ▶ $R[j] := 4$
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jumps to position x in the program;
sets instruction counter to x ;
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 - ▶ jumpz $x R[i]$
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if not the instruction counter is increased by 1;
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jump to $R[i]$ (indirect jump);
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- ▶ **uniform** cost model

Every operation takes time 1.

- ▶ logarithmic cost model

The cost depends on the content of memory cells:

- ▶ The time for a step is equal to the largest operand involved.
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Bounded word RAM model: cost is uniform but the largest value stored in a register may not exceed w , where usually $w = \log_2 n$.

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Algorithm 1 RepeatedSquaring(n)

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Asymptotic Notation

Formal Definition

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Asymptotic Notation

There is an equivalent definition using limes notation. f and g are functions from \mathbb{N} to \mathbb{R}^+ .

$$\blacktriangleright g \in \mathcal{O}(f): 0 \leq \lim_{n \rightarrow \infty} \frac{g(n)}{f(n)} < \infty$$

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Let f, g be functions with the property

$\exists n_0 > 0 \forall n \geq n_0 : f(n) > 0$ (the same for g). Then

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Comments

- ▶ Do not use asymptotic notation within induction proofs.
- ▶ For any constants a, b we have $\log_a n = \Theta(\log_b n)$.
Therefore, we will usually ignore the base of a logarithm within asymptotic notation.
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6 Recurrences

Algorithm 2 mergesort(list L)

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1:  $s \leftarrow \text{size}(L)$ 
2: if  $s \leq 1$  return  $L$ 
3:  $L_1 \leftarrow L[1 \cdots \lfloor \frac{s}{2} \rfloor]$ 
4:  $L_2 \leftarrow L[\lceil \frac{s}{2} \rceil \cdots n]$ 
5: mergesort( $L_1$ )
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7:  $L \leftarrow \text{merge}(L_1, L_2)$ 
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This algorithm requires

$$T(n) \leq 2T\left(\left\lceil \frac{n}{2} \right\rceil\right) + \mathcal{O}(n)$$

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Methods for Solving Recurrences

1. **Guessing+Induction**

Guess the right solution and prove that it is correct via induction. It needs experience to make the right guess.

2. **Master Theorem**

For a lot of recurrences that appear in the analysis of algorithms this theorem can be used to obtain tight asymptotic bounds. It does not provide exact solutions.

3. **Characteristic Polynomial**

Linear homogenous recurrences can be solved via this method.

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First we need to get rid of the \mathcal{O} -notation in our recurrence:

$$T(n) \leq \begin{cases} 2T(\lceil \frac{n}{2} \rceil) + cn & n \geq 2 \\ 0 & \text{otherwise} \end{cases}$$

Assume that instead we had

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Formally one would make an induction proof, where the above is the induction step. The base case is usually trivial.

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Hence, statement is **true** if we choose $d \geq c$.

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Note that we can do this as for constant-sized inputs the running time is always some constant (b in the above case).

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$$\begin{aligned}T(n) &\leq 2T\left(\left\lceil \frac{n}{2} \right\rceil\right) + cn \\ &\leq 2\left(d\left\lceil \frac{n}{2} \right\rceil \log \left\lceil \frac{n}{2} \right\rceil\right) + cn\end{aligned}$$

$$\boxed{\left\lceil \frac{n}{2} \right\rceil \leq \frac{n}{2} + 1} \leq 2(d(n/2 + 1) \log(n/2 + 1)) + cn$$

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$$\left\lceil \frac{n}{2} \right\rceil \leq \frac{n}{2} + 1$$

$$\frac{n}{2} + 1 \leq \frac{9}{16}n$$

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$$\boxed{\left\lceil \frac{n}{2} \right\rceil \leq \frac{n}{2} + 1} \leq 2\left(d\left(\frac{n}{2} + 1\right) \log\left(\frac{n}{2} + 1\right)\right) + cn$$

$$\boxed{\frac{n}{2} + 1 \leq \frac{9}{16}n} \leq dn \log\left(\frac{9}{16}n\right) + 2d \log n + cn$$

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$$\left\lceil \frac{n}{2} \right\rceil \leq \frac{n}{2} + 1$$

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$$\log \frac{9}{16}n = \log n + (\log 9 - 4)$$

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$$\boxed{\log \frac{9}{16}n = \log n + (\log 9 - 4)} = dn \log n + (\log 9 - 4)dn + 2d \log n + cn$$

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$$\leq 2\left(d\left(\frac{n}{2} + 1\right) \log\left(\frac{n}{2} + 1\right)\right) + cn$$

$$\frac{n}{2} + 1 \leq \frac{9}{16}n$$

$$\leq dn \log\left(\frac{9}{16}n\right) + 2d \log n + cn$$

$$\log \frac{9}{16}n = \log n + (\log 9 - 4)$$

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$$\log n \leq \frac{n}{4}$$

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$$= dn \log n + (\log 9 - 4)dn + 2d \log n + cn$$

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$$\left\lceil \frac{n}{2} \right\rceil \leq \frac{n}{2} + 1$$

$$\leq 2\left(d\left(\frac{n}{2} + 1\right) \log\left(\frac{n}{2} + 1\right)\right) + cn$$

$$\frac{n}{2} + 1 \leq \frac{9}{16}n$$

$$\leq dn \log\left(\frac{9}{16}n\right) + 2d \log n + cn$$

$$\log \frac{9}{16}n = \log n + (\log 9 - 4)$$

$$= dn \log n + (\log 9 - 4)dn + 2d \log n + cn$$

$$\log n \leq \frac{n}{4}$$

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$$\leq dn \log n - 0.33dn + cn$$

6.1 Guessing+Induction

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$$\left\lceil \frac{n}{2} \right\rceil \leq \frac{n}{2} + 1$$

$$\leq 2\left(d\left(\frac{n}{2} + 1\right) \log\left(\frac{n}{2} + 1\right)\right) + cn$$

$$\frac{n}{2} + 1 \leq \frac{9}{16}n$$

$$\leq dn \log\left(\frac{9}{16}n\right) + 2d \log n + cn$$

$$\log \frac{9}{16}n = \log n + (\log 9 - 4)$$

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$$\log n \leq \frac{n}{4}$$

$$= dn \log n + (\log 9 - 3.5)dn + cn$$

$$\leq dn \log n - 0.33dn + cn$$

$$\leq dn \log n$$

for a suitable choice of d .

6.2 Master Theorem

Lemma 4

Let $a \geq 1$, $b \geq 1$ and $\epsilon > 0$ denote constants. Consider the recurrence

$$T(n) = aT\left(\frac{n}{b}\right) + f(n) .$$

Case 1.

If $f(n) = \mathcal{O}(n^{\log_b(a)-\epsilon})$ then $T(n) = \Theta(n^{\log_b a})$.

Case 2.

If $f(n) = \Theta(n^{\log_b(a)} \log^k n)$ then $T(n) = \Theta(n^{\log_b a} \log^{k+1} n)$.

Case 3.

If $f(n) = \Omega(n^{\log_b(a)+\epsilon})$ and for sufficiently large n
 $af\left(\frac{n}{b}\right) \leq cf(n)$ for some constant $c < 1$ then $T(n) = \Theta(f(n))$.

6.2 Master Theorem

We prove the Master Theorem for the case that n is of the form b^ℓ , and we assume that the non-recursive case occurs for problem size 1 and incurs cost 1.

The Recursion Tree

The running time of a recursive algorithm can be visualized by a recursion tree:

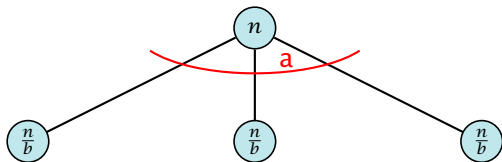
The Recursion Tree

The running time of a recursive algorithm can be visualized by a recursion tree:



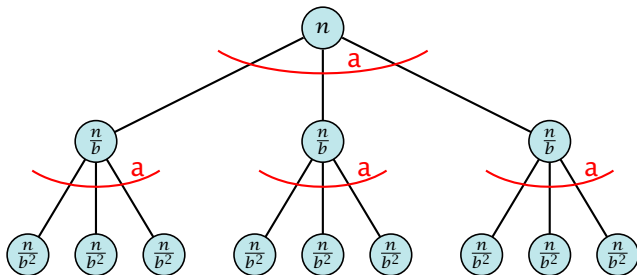
The Recursion Tree

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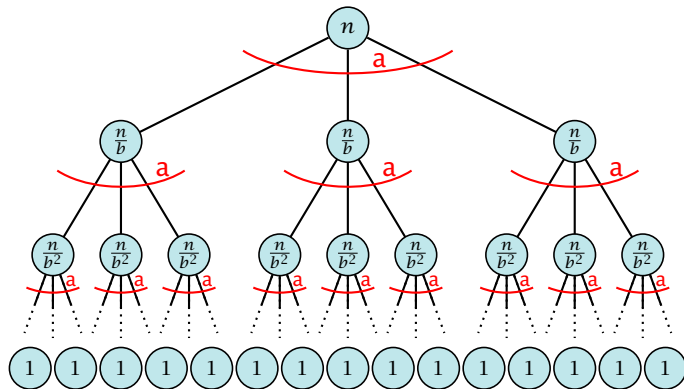
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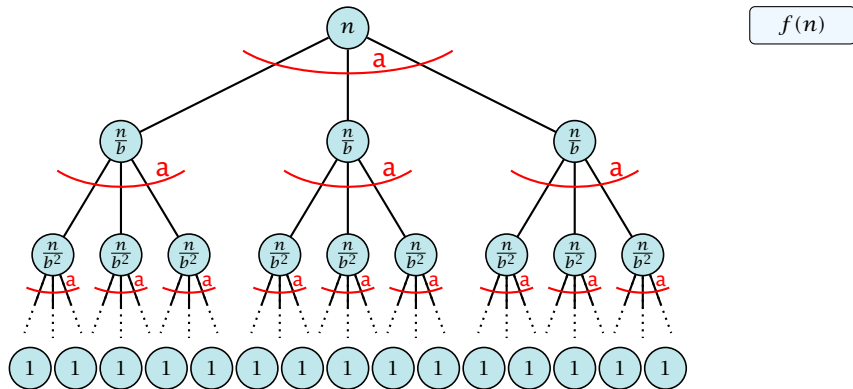
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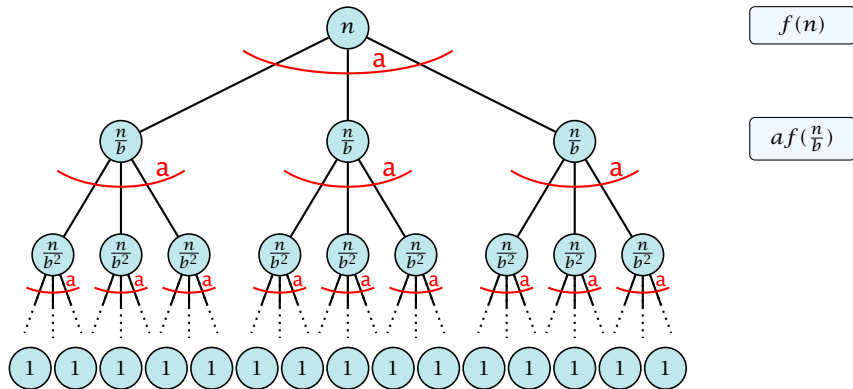
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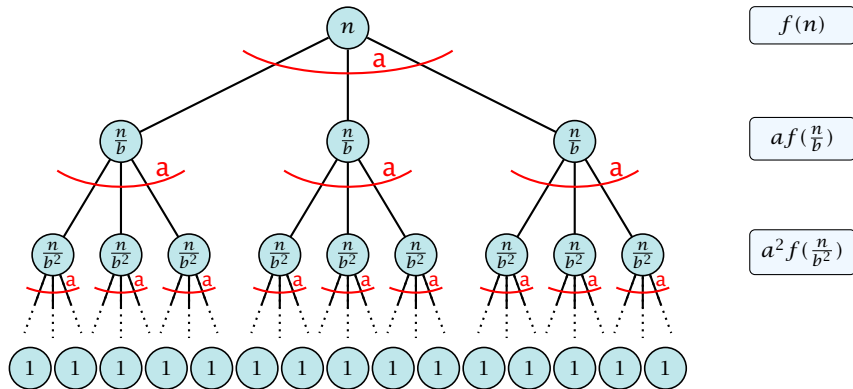
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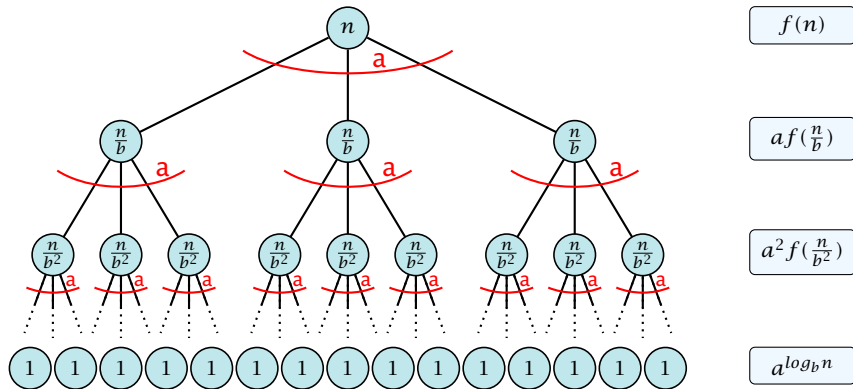
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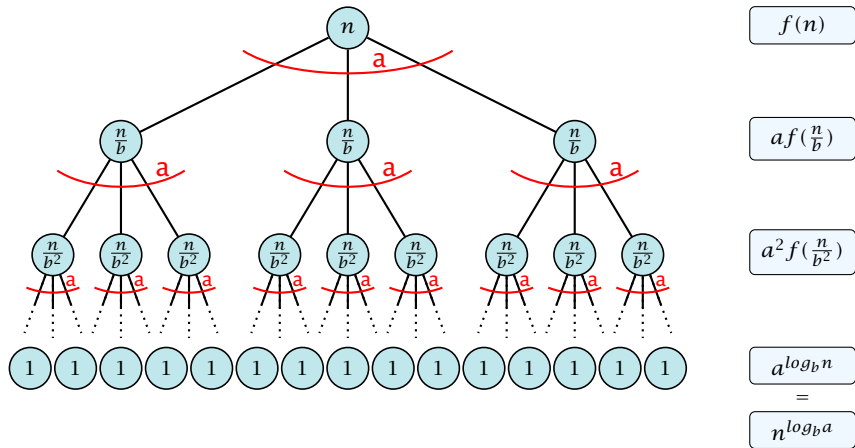
The Recursion Tree

The running time of a recursive algorithm can be visualized by a recursion tree:



The Recursion Tree

The running time of a recursive algorithm can be visualized by a recursion tree:



6.2 Master Theorem

This gives

$$T(n) = n^{\log_b a} + \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) .$$

Case 1. Now suppose that $f(n) \leq cn^{\log_b a - \epsilon}$.

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$$T(n) = n^{\log_b a}$$

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$$T(n) - n^{\log_b a} = \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right)$$

Case 1. Now suppose that $f(n) \leq cn^{\log_b a - \epsilon}$.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a - \epsilon} \end{aligned}$$

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$$b^{-i(\log_b a - \epsilon)} = b^{\epsilon i} (b^{\log_b a})^{-i} = b^{\epsilon i} a^{-i}$$

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$$\boxed{b^{-i(\log_b a - \epsilon)} = b^{\epsilon i} (b^{\log_b a})^{-i} = b^{\epsilon i} a^{-i}} = cn^{\log_b a - \epsilon} \sum_{i=0}^{\log_b n - 1} (b^{\epsilon})^i$$

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$$\boxed{\sum_{i=0}^k q^i = \frac{q^{k+1} - 1}{q - 1}}$$

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$$\boxed{\sum_{i=0}^k q^i = \frac{q^{k+1} - 1}{q - 1}} = cn^{\log_b a - \epsilon} (b^{\epsilon \log_b n} - 1) / (b^{\epsilon} - 1)$$

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$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a - \epsilon} \end{aligned}$$

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$$\begin{aligned} \boxed{\sum_{i=0}^k q^i = \frac{q^{k+1} - 1}{q - 1}} &= cn^{\log_b a - \epsilon} (b^{\epsilon \log_b n} - 1) / (b^{\epsilon} - 1) \\ &= cn^{\log_b a - \epsilon} (n^{\epsilon} - 1) / (b^{\epsilon} - 1) \end{aligned}$$

Case 1. Now suppose that $f(n) \leq cn^{\log_b a - \epsilon}$.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a - \epsilon} \end{aligned}$$

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$$\begin{aligned} \boxed{\sum_{i=0}^k q^i = \frac{q^{k+1} - 1}{q - 1}} &= cn^{\log_b a - \epsilon} (b^{\epsilon \log_b n} - 1) / (b^{\epsilon} - 1) \\ &= cn^{\log_b a - \epsilon} (n^{\epsilon} - 1) / (b^{\epsilon} - 1) \\ &= \frac{c}{b^{\epsilon} - 1} n^{\log_b a} (n^{\epsilon} - 1) / (n^{\epsilon}) \end{aligned}$$

Hence,

$$T(n) \leq \left(\frac{c}{b^{\epsilon} - 1} + 1 \right) n^{\log_b(a)}$$

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$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a - \epsilon} \end{aligned}$$

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$$\begin{aligned} \boxed{\sum_{i=0}^k q^i = \frac{q^{k+1} - 1}{q - 1}} &= cn^{\log_b a - \epsilon} (b^{\epsilon \log_b n} - 1) / (b^{\epsilon} - 1) \\ &= cn^{\log_b a - \epsilon} (n^{\epsilon} - 1) / (b^{\epsilon} - 1) \\ &= \frac{c}{b^{\epsilon} - 1} n^{\log_b a} (n^{\epsilon} - 1) / (n^{\epsilon}) \end{aligned}$$

Hence,

$$T(n) \leq \left(\frac{c}{b^{\epsilon} - 1} + 1 \right) n^{\log_b(a)} \quad \Rightarrow T(n) = \mathcal{O}(n^{\log_b a}).$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a}$.

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Hence,

$$T(n) = \mathcal{O}(n^{\log_b a} \log_b n)$$

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Hence,

$$T(n) = \mathcal{O}(n^{\log_b a} \log_b n) \quad \Rightarrow T(n) = \mathcal{O}(n^{\log_b a} \log n).$$

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Hence,

$$T(n) = \Omega(n^{\log_b a} \log_b n)$$

Case 2. Now suppose that $f(n) \geq cn^{\log_b a}$.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\geq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a} \\ &= cn^{\log_b a} \sum_{i=0}^{\log_b n - 1} 1 \\ &= cn^{\log_b a} \log_b n \end{aligned}$$

Hence,

$$T(n) = \Omega(n^{\log_b a} \log_b n) \quad \Rightarrow T(n) = \Omega(n^{\log_b a} \log n).$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

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$$T(n) = n^{\log_b a}$$

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$$T(n) - n^{\log_b a} = \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right)$$

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$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a} \cdot \left(\log_b\left(\frac{n}{b^i}\right)\right)^k \end{aligned}$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

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$$n = b^\ell \Rightarrow \ell = \log_b n$$

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$$\boxed{n = b^\ell \Rightarrow \ell = \log_b n} = cn^{\log_b a} \sum_{i=0}^{\ell-1} \left(\log_b \left(\frac{b^\ell}{b^i}\right)\right)^k$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a} \cdot \left(\log_b \left(\frac{n}{b^i}\right)\right)^k \end{aligned}$$

$n = b^\ell \Rightarrow \ell = \log_b n$
--

$$\begin{aligned} &= cn^{\log_b a} \sum_{i=0}^{\ell-1} \left(\log_b \left(\frac{b^\ell}{b^i}\right)\right)^k \\ &= cn^{\log_b a} \sum_{i=0}^{\ell-1} (\ell - i)^k \end{aligned}$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a} \cdot \left(\log_b\left(\frac{n}{b^i}\right)\right)^k \end{aligned}$$

$$\begin{aligned} \boxed{n = b^\ell \Rightarrow \ell = \log_b n} &= cn^{\log_b a} \sum_{i=0}^{\ell-1} \left(\log_b\left(\frac{b^\ell}{b^i}\right)\right)^k \\ &= cn^{\log_b a} \sum_{i=0}^{\ell-1} (\ell - i)^k \\ &= cn^{\log_b a} \sum_{i=1}^{\ell} i^k \end{aligned}$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

$$T(n) - n^{\log_b a} = \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ \leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a} \cdot \left(\log_b\left(\frac{n}{b^i}\right)\right)^k$$

$$n = b^\ell \Rightarrow \ell = \log_b n$$

$$= cn^{\log_b a} \sum_{i=0}^{\ell-1} \left(\log_b\left(\frac{b^\ell}{b^i}\right)\right)^k$$

$$= cn^{\log_b a} \sum_{i=0}^{\ell-1} (\ell - i)^k$$

$$= cn^{\log_b a} \sum_{i=1}^{\ell} i^k \approx \frac{1}{k} \ell^{k+1}$$

Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

$$\begin{aligned}
 T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\
 &\leq c \sum_{i=0}^{\log_b n - 1} a^i \left(\frac{n}{b^i}\right)^{\log_b a} \cdot \left(\log_b\left(\frac{n}{b^i}\right)\right)^k
 \end{aligned}$$

$$n = b^\ell \Rightarrow \ell = \log_b n$$

$$\begin{aligned}
 &= cn^{\log_b a} \sum_{i=0}^{\ell-1} \left(\log_b\left(\frac{b^\ell}{b^i}\right)\right)^k \\
 &= cn^{\log_b a} \sum_{i=0}^{\ell-1} (\ell - i)^k \\
 &= cn^{\log_b a} \sum_{i=1}^{\ell} i^k \\
 &\approx \frac{c}{k} n^{\log_b a} \ell^{k+1}
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Case 2. Now suppose that $f(n) \leq cn^{\log_b a} (\log_b(n))^k$.

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$$\boxed{n = b^\ell \Rightarrow \ell = \log_b n} = cn^{\log_b a} \sum_{i=0}^{\ell-1} \left(\log_b\left(\frac{b^\ell}{b^i}\right)\right)^k$$

$$= cn^{\log_b a} \sum_{i=0}^{\ell-1} (\ell - i)^k$$

$$= cn^{\log_b a} \sum_{i=1}^{\ell} i^k$$

$$\approx \frac{c}{k} n^{\log_b a} \ell^{k+1}$$

$$\Rightarrow T(n) = \mathcal{O}(n^{\log_b a} \log^{k+1} n).$$

Case 3. Now suppose that $f(n) \geq dn^{\log_b a + \epsilon}$, and that for sufficiently large n : $af(n/b) \leq cf(n)$, for $c < 1$.

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From this we get $a^i f(n/b^i) \leq c^i f(n)$, where we assume that $n/b^{i-1} \geq n_0$ is still sufficiently large.

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$$T(n) - n^{\log_b a} = \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right)$$

Case 3. Now suppose that $f(n) \geq dn^{\log_b a + \epsilon}$, and that for sufficiently large n : $af(n/b) \leq cf(n)$, for $c < 1$.

From this we get $a^i f(n/b^i) \leq c^i f(n)$, where we assume that $n/b^{i-1} \geq n_0$ is still sufficiently large.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &= \sum_{i=0}^{\log_b n - 1} c^i f(n) + \mathcal{O}(n^{\log_b a}) \end{aligned}$$

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$$q < 1 : \sum_{i=0}^n q^i = \frac{1 - q^{n+1}}{1 - q} \leq \frac{1}{1 - q}$$

Case 3. Now suppose that $f(n) \geq dn^{\log_b a + \epsilon}$, and that for sufficiently large n : $af(n/b) \leq cf(n)$, for $c < 1$.

From this we get $a^i f(n/b^i) \leq c^i f(n)$, where we assume that $n/b^{i-1} \geq n_0$ is still sufficiently large.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &= \sum_{i=0}^{\log_b n - 1} c^i f(n) + \mathcal{O}(n^{\log_b a}) \end{aligned}$$

$$\boxed{q < 1 : \sum_{i=0}^n q^i = \frac{1-q^{n+1}}{1-q} \leq \frac{1}{1-q}} \leq \frac{1}{1-c} f(n) + \mathcal{O}(n^{\log_b a})$$

Case 3. Now suppose that $f(n) \geq dn^{\log_b a + \epsilon}$, and that for sufficiently large n : $af(n/b) \leq cf(n)$, for $c < 1$.

From this we get $a^i f(n/b^i) \leq c^i f(n)$, where we assume that $n/b^{i-1} \geq n_0$ is still sufficiently large.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &= \sum_{i=0}^{\log_b n - 1} c^i f(n) + \mathcal{O}(n^{\log_b a}) \end{aligned}$$

$$\boxed{q < 1 : \sum_{i=0}^n q^i = \frac{1-q^{n+1}}{1-q} \leq \frac{1}{1-q}} \leq \frac{1}{1-c} f(n) + \mathcal{O}(n^{\log_b a})$$

Hence,

$$T(n) \leq \mathcal{O}(f(n))$$

Case 3. Now suppose that $f(n) \geq dn^{\log_b a + \epsilon}$, and that for sufficiently large n : $af(n/b) \leq cf(n)$, for $c < 1$.

From this we get $a^i f(n/b^i) \leq c^i f(n)$, where we assume that $n/b^{i-1} \geq n_0$ is still sufficiently large.

$$\begin{aligned} T(n) - n^{\log_b a} &= \sum_{i=0}^{\log_b n - 1} a^i f\left(\frac{n}{b^i}\right) \\ &= \sum_{i=0}^{\log_b n - 1} c^i f(n) + \mathcal{O}(n^{\log_b a}) \\ &\leq \frac{1}{1-c} f(n) + \mathcal{O}(n^{\log_b a}) \end{aligned}$$

$$q < 1 : \sum_{i=0}^n q^i = \frac{1-q^{n+1}}{1-q} \leq \frac{1}{1-q}$$

Hence,

$$T(n) \leq \mathcal{O}(f(n))$$

$$\Rightarrow T(n) = \Theta(f(n)).$$

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

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Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

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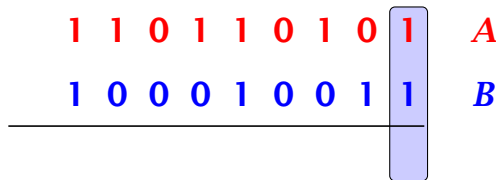
For this we first need to be able to add two integers A and B :

$$\begin{array}{r} 1\ 1\ 0\ 1\ 1\ 0\ 1\ 0\ 1\ A \\ 1\ 0\ 0\ 0\ 1\ 0\ 0\ 1\ 1\ B \\ \hline \end{array}$$

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1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									0

The diagram shows the addition of two 8-bit integers, A and B. The bits of A are 1 1 0 1 1 0 1 0 and the bits of B are 1 0 0 0 1 0 0 1. A horizontal line is drawn under the bits of B. A vertical bar on the right side of the diagram contains the bits 1, 1, and 0, corresponding to the bits of A, B, and the result of the addition, respectively. A small green '1' is written below the 8th bit of B, indicating a carry-in.

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1	1	0	1	1	0	1	0	1	<i>A</i>
1	0	0	0	1	0	0	1	1	<i>B</i>
<hr/>									
							1		
								0	

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For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
							1	1	
							0	0	

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For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									
						1	1	0	0

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

$$\begin{array}{rcccccccc} 1 & 1 & 0 & 1 & 1 & 0 & 1 & 0 & 1 & A \\ 1 & 0 & 0 & 0 & 1 & 0 & 0 & 1 & 1 & B \\ \hline & & & & & & 0 & 0 & 0 & \end{array}$$

The diagram illustrates the addition of two integers, A and B, using a register of constant size. The numbers are represented as binary strings. A vertical bar highlights the current bit position being processed, which is the 7th bit from the right (the least significant bit of the register). The carry bits are shown as small green '1's below the lines.

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									
					1		1	1	
						0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

$$\begin{array}{rcccccccc} 1 & 1 & 0 & 1 & 1 & 0 & 1 & 0 & 1 & A \\ 1 & 0 & 0 & 0 & 1 & 0 & 0 & 1 & 1 & B \\ \hline & & & & 0 & 1 & 1 & 1 & & \\ & & & & & 1 & 0 & 0 & 0 & \end{array}$$

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For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									
				0	1	1	1		
					1	0	0	0	

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Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
					1	0	1	1	1
					0	1	0	0	0

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For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									
				1	0	1	1	1	
				0	1	0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

$$\begin{array}{rcccccccc} 1 & 1 & 0 & 1 & 1 & 0 & 1 & 0 & 1 & A \\ 1 & 0 & 0 & 0 & 1 & 0 & 0 & 1 & 1 & B \\ \hline & & & 0 & 0 & 1 & 0 & 0 & 0 & \end{array}$$

The diagram illustrates the addition of two 9-bit integers, A and B, to produce a 9-bit result. The numbers are aligned by their least significant bits. A horizontal line is drawn under the second row. The result is shown below the line. A vertical blue box highlights the 4th bit position (index 3 from the right), which contains a 0. This bit is the result of adding the 4th bits of A and B (1 + 0) and the carry-in from the 3rd bit position (1). The carry-out from this position is 1, which is shown as a small green '1' below the 3rd bit of the second row. The carry-in for the 5th bit position is 0, shown as a small green '0' below the 5th bit of the second row. The carry-in for the 6th bit position is 1, shown as a small green '1' below the 6th bit of the second row. The carry-in for the 7th bit position is 1, shown as a small green '1' below the 7th bit of the second row. The carry-in for the 8th bit position is 1, shown as a small green '1' below the 8th bit of the second row.

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									
			1	1	0	1	1	1	
			0	0	1	0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

The diagram illustrates the addition of two integers, A and B , in binary. The numbers are aligned by their least significant bits. A horizontal line is drawn under the numbers. A vertical bar highlights the carry propagation from bit 2 to bit 3.

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
	0	1	1	0	1	1	1		
<hr/>									
		1	0	0	1	0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
<hr/>									
	0	1	1	0	1	1	1		
		1	0	0	1	0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

$$\begin{array}{r} 1\ 1\ 0\ 1\ 1\ 0\ 1\ 0\ 1\ A \\ 1\ 0\ 0\ 0\ 1\ 0\ 0\ 1\ 1\ B \\ \hline 1\ 1\ 0\ 0\ 1\ 0\ 0\ 0 \end{array}$$

The diagram illustrates the addition of two 9-bit integers, A and B, to produce a 9-bit result. The bits of A are 1, 1, 0, 1, 1, 0, 1, 0, 1. The bits of B are 1, 0, 0, 0, 1, 0, 0, 1, 1. The result is 1, 1, 0, 0, 1, 0, 0, 0. A vertical blue box highlights the first two bits of the result, 1 and 1, which correspond to the first two bits of the input integers, 1 and 1. Small green numbers (0, 0, 1, 1, 0, 1, 1, 1) are placed below the bits of B, representing carry bits from the previous addition steps.

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

1	1	0	1	1	0	1	0	1	A
1	0	0	0	1	0	0	1	1	B
0	0	1	1	0	1	1	1		
1	1	0	0	1	0	0	0		

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

	1	1	0	1	1	0	1	0	1	<i>A</i>
1	1	0	0	0	1	0	0	1	1	<i>B</i>
<hr/>										
	0	1	1	0	0	1	0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

	1	1	0	1	1	0	1	0	1	A
	1	0	0	0	1	0	0	1	1	B
	<hr/>									
	0	1	1	0	0	1	0	0	0	

Carry bits: 1 (left), 0 (right)

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For this we first need to be able to add two integers A and B :

	1	1	0	1	1	0	1	0	1	A
	1	0	0	0	1	0	0	1	1	B
	<hr/>									
1	0	1	1	0	0	1	0	0	0	

Example: Multiplying Two Integers

Suppose we want to multiply two n -bit Integers, but our registers can only perform operations on integers of constant size.

For this we first need to be able to add two integers A and B :

$$\begin{array}{r} 1\ 1\ 0\ 1\ 1\ 0\ 1\ 0\ 1\ A \\ 1\ 0\ 0\ 0\ 1\ 0\ 0\ 1\ 1\ B \\ \hline 1\ 0\ 1\ 1\ 0\ 0\ 1\ 0\ 0\ 0 \end{array}$$

This gives that two n -bit integers can be added in time $\mathcal{O}(n)$.

Example: Multiplying Two Integers

Suppose that we want to multiply an n -bit integer A and an m -bit integer B ($m \leq n$).

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$$\begin{array}{r} 10001 \times 1011 \\ \hline \end{array}$$

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$$\begin{array}{r} 10001 \times 1011 \\ \hline 10001 \end{array}$$

Example: Multiplying Two Integers

Suppose that we want to multiply an n -bit integer A and an m -bit integer B ($m \leq n$).

$$\begin{array}{r} 1\ 0\ 0\ 0\ 1 \times 1\ 0\ 1\ 1 \\ \hline 1\ 0\ 0\ 0\ 1 \end{array}$$

Example: Multiplying Two Integers

Suppose that we want to multiply an n -bit integer A and an m -bit integer B ($m \leq n$).

$$\begin{array}{r} 1\ 0\ 0\ 0\ 1 \times 1\ 0\ 1\ 1 \\ \hline 1\ 0\ 0\ 0\ 1 \\ 0 \\ 0 \end{array}$$

Example: Multiplying Two Integers

Suppose that we want to multiply an n -bit integer A and an m -bit integer B ($m \leq n$).

$$\begin{array}{r} 10001 \times 1011 \\ \hline 10001 \\ 100010 \\ \hline \end{array}$$

Example: Multiplying Two Integers

Suppose that we want to multiply an n -bit integer A and an m -bit integer B ($m \leq n$).

$$\begin{array}{r} 1\ 0\ 0\ 0\ 1 \times 1\ 0\ 1\ 1 \\ \hline 1\ 0\ 0\ 0\ 1 \\ 1\ 0\ 0\ 0\ 1\ 0 \end{array}$$

Example: Multiplying Two Integers

Suppose that we want to multiply an n -bit integer A and an m -bit integer B ($m \leq n$).

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Time requirement:

- ▶ Computing intermediate results: $\mathcal{O}(nm)$.
- ▶ Adding m numbers of length $\leq 2n$: $\mathcal{O}((m+n)m) = \mathcal{O}(nm)$.

Example: Multiplying Two Integers

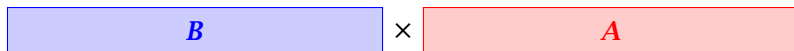
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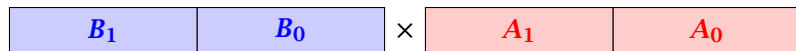
Suppose that integers A and B are of length $n = 2^k$, for some k .

$$\begin{array}{|c|c|} \hline B_1 & B_0 \\ \hline \end{array} \times \begin{array}{|c|c|} \hline A_1 & A_0 \\ \hline \end{array}$$

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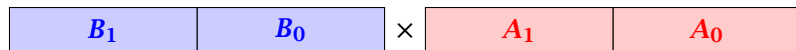
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$$A = A_1 \cdot 2^{\frac{n}{2}} + A_0 \text{ and } B = B_1 \cdot 2^{\frac{n}{2}} + B_0$$

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Hence,

$$A \cdot B = A_1 B_1 \cdot 2^n + (A_1 B_0 + A_0 B_1) \cdot 2^{\frac{n}{2}} + A_0 \cdot B_0$$

Example: Multiplying Two Integers

Algorithm 3 $\text{mult}(A, B)$

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1: if  $|A| = |B| = 1$  then  
2:   return  $a_0 \cdot b_0$   
3: split  $A$  into  $A_0$  and  $A_1$   
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We get the following recurrence:

$$T(n) = 4T\left(\frac{n}{2}\right) + \mathcal{O}(n) .$$

Example: Multiplying Two Integers

Master Theorem: Recurrence: $T[n] = aT(\frac{n}{b}) + f(n)$.

- ▶ Case 1: $f(n) = \mathcal{O}(n^{\log_b a - \epsilon})$ $T(n) = \Theta(n^{\log_b a})$
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In our case $a = 4$, $b = 2$, and $f(n) = \Theta(n)$. Hence, we are in Case 1, since $n = \mathcal{O}(n^{2-\epsilon}) = \mathcal{O}(n^{\log_b a - \epsilon})$.

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⇒ Not better than the “school method”.

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Example: Multiplying Two Integers

We can use the following identity to compute Z_1 :

$$\begin{aligned} Z_1 &= A_1B_0 + A_0B_1 && = Z_2 && = Z_0 \\ &= (A_0 + A_1) \cdot (B_0 + B_1) - \underbrace{A_1B_1}_{Z_2} - \underbrace{A_0B_0}_{Z_0} \end{aligned}$$

Hence,

Algorithm 4 mult(A, B)

1: **if** $|A| = |B| = 1$ **then**

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We get the following recurrence:

$$T(n) = 3T\left(\frac{n}{2}\right) + \mathcal{O}(n) .$$

Master Theorem: Recurrence: $T[n] = aT\left(\frac{n}{b}\right) + f(n)$.

- ▶ Case 1: $f(n) = \mathcal{O}(n^{\log_b a - \epsilon})$ $T(n) = \Theta(n^{\log_b a})$
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Again we are in Case 1. We get a running time of $\Theta(n^{\log_2 3}) \approx \Theta(n^{1.59})$.

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6.3 The Characteristic Polynomial

Consider the recurrence relation:

$$c_0T(n) + c_1T(n-1) + c_2T(n-2) + \dots + c_kT(n-k) = f(n)$$

This is the general form of a linear recurrence relation of order k with constant coefficients ($c_0, c_k \neq 0$).

The value of $T(n)$ only depends on the k preceding values. This means

the recurrence relation is of order k .

The recurrence is linear as there are no products of $T(n)$'s

and $T(n)$ only appears the first time in the relation. Therefore, a linear

recurrence relation is a linear recurrence relation of order k .

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Observations:

- ▶ The solution $T[0], T[1], T[2], \dots$ is completely determined by a set of boundary conditions that specify values for $T[0], \dots, T[k-1]$.
- ▶ In fact, any k consecutive values completely determine the solution.
- ▶ k non-consecutive values might not be an appropriate set of boundary conditions (depends on the problem).

Approach:

- ▶ First determine all solutions that satisfy recurrence relation.
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The Homogenous Case

The solution space

$$S = \left\{ T = T[0], T[1], T[2], \dots \mid T \text{ fulfills recurrence relation} \right\}$$

is a **vector space**. This means that if $T_1, T_2 \in S$, then also $\alpha T_1 + \beta T_2 \in S$, for arbitrary constants α, β .

How do we find a non-trivial solution?

We guess that the solution is of the form λ^n , $\lambda \neq 0$, and see what happens. In order for this guess to fulfill the recurrence we need

$$c_0 \lambda^n + c_1 \lambda^{n-1} + c_2 \cdot \lambda^{n-2} + \dots + c_k \cdot \lambda^{n-k} = 0$$

for all $n \geq k$.

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Dividing by λ^{n-k} gives that all these constraints are identical to

$$c_0\lambda^k + c_1\lambda^{k-1} + c_2 \cdot \lambda^{k-2} + \dots + c_k = 0$$

This means that if λ_i is a root (Nullstelle) of $P[\lambda]$ then $T[n] = \lambda_i^n$ is a solution to the recurrence relation.

Let $\lambda_1, \dots, \lambda_k$ be the k (complex) roots of $P[\lambda]$. Then, because of the vector space property

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Lemma 5

Assume that the characteristic polynomial has k *distinct* roots $\lambda_1, \dots, \lambda_k$. Then *all* solutions to the recurrence relation are of the form

$$\alpha_1 \lambda_1^n + \alpha_2 \lambda_2^n + \dots + \alpha_k \lambda_k^n .$$

Proof.

There is one solution for every possible choice of boundary conditions for $T[1], \dots, T[k]$.

We show that the above set of solutions contains one solution for every choice of boundary conditions.

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Suppose I am given boundary conditions $T[i]$ and I want to see whether I can choose the α'_i 's such that these conditions are met:

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We show that the column vectors are linearly independent. Then the above equation has a solution.

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This we show by induction:



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A vector (λ_i) , $\lambda_i \neq 0$ is linearly independent.
- ▶ **induction step** ($k \rightarrow k + 1$):
assume for contradiction that there exist α_i 's with

$$\alpha_1 \begin{pmatrix} \lambda_1 \\ \vdots \\ \lambda_1^{k-1} \\ \lambda_1^k \end{pmatrix} + \cdots + \alpha_k \begin{pmatrix} \lambda_k \\ \vdots \\ \lambda_k^{k-1} \\ \lambda_k^k \end{pmatrix} = 0$$

and not all $\alpha_i = 0$.



The Homogenous Case

Proof (cont.).

This we show by induction:

- ▶ **base case** ($k = 1$):
A vector (λ_i) , $\lambda_i \neq 0$ is linearly independent.
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and not all $\alpha_i = 0$. **Then all $\alpha_i \neq 0$!**



The Homogeneous Case

$$\alpha_1 \begin{pmatrix} \lambda_1 \\ \lambda_1^2 \\ \vdots \\ \lambda_1^{k-1} \\ \lambda_1^k \end{pmatrix} + \cdots + \alpha_k \begin{pmatrix} \lambda_k \\ \lambda_k^2 \\ \vdots \\ \lambda_k^{k-1} \\ \lambda_k^k \end{pmatrix} = 0$$

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$\lambda_1 v_1 =$ (bottom part of the first vector) $\lambda_k v_k =$ (bottom part of the second vector)

The Homogeneous Case

$$\alpha_1 \begin{pmatrix} \lambda_1 \\ \lambda_1^2 \\ \vdots \\ \lambda_1^{k-1} \\ \lambda_1^k \end{pmatrix} + \dots + \alpha_k \begin{pmatrix} \lambda_k \\ \lambda_k^2 \\ \vdots \\ \lambda_k^{k-1} \\ \lambda_k^k \end{pmatrix} = 0$$

$\lambda_1 v_1 =$ (left vector) $\lambda_k v_k =$ (right vector)

This means that

$$\sum_{i=1}^k \alpha_i v_i = 0 \quad \text{and} \quad \sum_{i=1}^k \lambda_i \alpha_i v_i = 0$$

The Homogeneous Case

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$$\sum_{i=1}^k \alpha_i v_i = 0 \quad \text{and} \quad \sum_{i=1}^k \lambda_i \alpha_i v_i = 0$$

Hence,

$$\sum_{i=1}^{k-1} \alpha_i v_i + \alpha_k v_k = 0 \quad \text{and} \quad -\frac{1}{\lambda_k} \sum_{i=1}^{k-1} \lambda_i \alpha_i v_i = \alpha_k v_k$$

The Homogeneous Case

This gives that

$$\sum_{i=1}^{k-1} \left(1 - \frac{\lambda_i}{\lambda_k}\right) \alpha_i \mathbf{v}_i = \mathbf{0} .$$

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This gives that

$$\sum_{i=1}^{k-1} \left(1 - \frac{\lambda_i}{\lambda_k}\right) \alpha_i v_i = 0 .$$

This is a contradiction as the v_i 's are linearly independent because of induction hypothesis.

The Homogeneous Case

What happens if the roots are not all distinct?

Suppose we have a root λ_i with multiplicity (Vielfachheit) at least 2. Then not only is λ_i^n a solution to the recurrence but also $n\lambda_i^n$.

To see this consider the polynomial

$$P(\lambda)\lambda^{n-k} = c_0\lambda^n + c_1\lambda^{n-1} + c_2\lambda^{n-2} + \dots + c_k\lambda^{n-k}$$

Since λ_i is a root we can write this as $Q(\lambda)(\lambda - \lambda_i)^2$. Calculating the derivative gives a polynomial that still has root λ_i .

This means

$$c_0n\lambda_i^{n-1} + c_1(n-1)\lambda_i^{n-2} + \dots + c_k(n-k)\lambda_i^{n-k-1} = 0$$

Hence,

$$\underbrace{c_0n\lambda_i^n}_{T[n]} + \underbrace{c_1(n-1)\lambda_i^{n-1}}_{T[n-1]} + \dots + \underbrace{c_k(n-k)\lambda_i^{n-k}}_{T[n-k]} = 0$$

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The Homogeneous Case

Suppose λ_i has multiplicity j . We know that

$$c_0 n \lambda_i^n + c_1 (n-1) \lambda_i^{n-1} + \dots + c_k (n-k) \lambda_i^{n-k} = 0$$

(after taking the derivative; multiplying with λ ; plugging in λ_i)

Doing this again gives

$$c_0 n^2 \lambda_i^n + c_1 (n-1)^2 \lambda_i^{n-1} + \dots + c_k (n-k)^2 \lambda_i^{n-k} = 0$$

We can continue $j-1$ times.

Hence, $n^\ell \lambda_i^n$ is a solution for $\ell \in 0, \dots, j-1$.

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The Homogeneous Case

Lemma 6

Let $P[\lambda]$ denote the characteristic polynomial to the recurrence

$$c_0T[n] + c_1T[n-1] + \dots + c_kT[n-k] = 0$$

Let λ_i , $i = 1, \dots, m$ be the (complex) roots of $P[\lambda]$ with multiplicities ℓ_i . Then the general solution to the recurrence is given by

$$T[n] = \sum_{i=1}^m \sum_{j=0}^{\ell_i-1} \alpha_{ij} \cdot (n^j \lambda_i^n) .$$

The full proof is omitted. We have only shown that any choice of α_{ij} 's is a solution to the recurrence.

Example: Fibonacci Sequence

$$T[0] = 0$$

$$T[1] = 1$$

$$T[n] = T[n - 1] + T[n - 2] \text{ for } n \geq 2$$

The characteristic polynomial is

$$\lambda^2 - \lambda - 1$$

Finding the roots, gives

$$\lambda_{1/2} = \frac{1}{2} \pm \sqrt{\frac{1}{4} + 1} = \frac{1}{2} (1 \pm \sqrt{5})$$

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Hence, the solution is of the form

$$\alpha \left(\frac{1 + \sqrt{5}}{2} \right)^n + \beta \left(\frac{1 - \sqrt{5}}{2} \right)^n$$

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$T[0] = 0$ gives $\alpha + \beta = 0$.

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$$\alpha \left(\frac{1 + \sqrt{5}}{2} \right) + \beta \left(\frac{1 - \sqrt{5}}{2} \right) = 1 \implies \alpha - \beta = \frac{2}{\sqrt{5}}$$

Example: Fibonacci Sequence

Hence, the solution is

$$\frac{1}{\sqrt{5}} \left[\left(\frac{1 + \sqrt{5}}{2} \right)^n - \left(\frac{1 - \sqrt{5}}{2} \right)^n \right]$$

The Inhomogeneous Case

Consider the recurrence relation:

$$c_0T(n) + c_1T(n - 1) + c_2T(n - 2) + \cdots + c_kT(n - k) = f(n)$$

with $f(n) \neq 0$.

While we have a fairly general technique for solving **homogeneous**, linear recurrence relations the inhomogeneous case is different.

The Inhomogeneous Case

The general solution of the recurrence relation is

$$T(n) = T_h(n) + T_p(n) ,$$

where T_h is **any** solution to the homogeneous equation, and T_p is **one** particular solution to the inhomogeneous equation.

There is no general method to find a particular solution.

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The Inhomogeneous Case

Example:

$$T[n] = T[n - 1] + 1 \quad T[0] = 1$$

Then,

$$T[n - 1] = T[n - 2] + 1 \quad (n \geq 2)$$

Subtracting the first from the second equation gives,

$$T[n] - T[n - 1] = T[n - 1] - T[n - 2] \quad (n \geq 2)$$

or

$$T[n] = 2T[n - 1] - T[n - 2] \quad (n \geq 2)$$

I get a completely determined recurrence if I add $T[0] = 1$ and $T[1] = 2$.

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$T[0] = 1$ gives $\alpha = 1$.

$T[1] = 2$ gives $1 + \beta = 2 \Rightarrow \beta = 1$.

The Inhomogeneous Case

If $f(n)$ is a polynomial of degree r this method can be applied $r + 1$ times to obtain a homogeneous equation:

$$T[n] = T[n - 1] + n^2$$

Shift:

$$T[n - 1] = T[n - 2] + (n - 1)^2$$

Difference:

$$T[n] - T[n - 1] = T[n - 1] - T[n - 2] + 2n - 1$$

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The Inhomogeneous Case

If $f(n)$ is a polynomial of degree r this method can be applied $r + 1$ times to obtain a homogeneous equation:

$$T[n] = T[n - 1] + n^2$$

Shift:

$$T[n - 1] = T[n - 2] + (n - 1)^2 = T[n - 2] + n^2 - 2n + 1$$

Difference:

$$T[n] - T[n - 1] = T[n - 1] - T[n - 2] + 2n - 1$$

$$T[n] = 2T[n - 1] - T[n - 2] + 2n - 1$$

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Shift:

$$T[n - 1] = 2T[n - 2] - T[n - 3] + 2(n - 1) - 1$$

$$T[n] = 2T[n - 1] - T[n - 2] + 2n - 1$$

Shift:

$$\begin{aligned} T[n - 1] &= 2T[n - 2] - T[n - 3] + 2(n - 1) - 1 \\ &= 2T[n - 2] - T[n - 3] + 2n - 3 \end{aligned}$$

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and so on...

6.4 Generating Functions

Definition 7 (Generating Function)

Let $(a_n)_{n \geq 0}$ be a sequence. The corresponding

- ▶ **generating function** (Erzeugendenfunktion) is

$$F(z) := \sum_{n=0}^{\infty} a_n z^n;$$

- ▶ exponential generating function (exponentielle Erzeugendenfunktion) is

$$F(z) = \sum_{n \geq 0} \frac{a_n}{n!} z^n.$$

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1. The generating function of the sequence $(1, 0, 0, \dots)$ is

$$F(z) = 1.$$

2. The generating function of the sequence $(1, 1, 1, \dots)$ is

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6.4 Generating Functions

There are two different views:

A generating function is a formal power series (formale Potenzreihe).

Then the generating function is an **algebraic object**.

Let $f = \sum_{n=0}^{\infty} a_n z^n$ and $g = \sum_{n=0}^{\infty} b_n z^n$.

- ▶ **Equality:** f and g are equal if $a_n = b_n$ for all n .
- ▶ **Addition:** $f + g := \sum_{n=0}^{\infty} (a_n + b_n) z^n$.
- ▶ **Multiplication:** $f \cdot g := \sum_{n=0}^{\infty} c_n z^n$ with $c = \sum_{p=0}^n a_p b_{n-p}$.

There are no convergence issues here.

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The arithmetic view:

We view a power series as a function $f : \mathbb{C} \rightarrow \mathbb{C}$.

Then, it is important to think about convergence/convergence radius etc.

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What does $\sum_{n=0}^{\infty} z^n = \frac{1}{1-z}$ mean in the **algebraic view**?

It means that the power series $1 - z$ and the power series $\sum_{n=0}^{\infty} z^n$ are invers, i.e.,

$$(1 - z) \cdot \left(\sum_{n=0}^{\infty} z^n \right) = 1 .$$

This is well-defined.

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Hence, the generating function of the sequence $a_n = n + 1$ is $1/(1-z)^2$.

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Hence, the generating function of the sequence $a_n = (n+1)(n+2)$ is $\frac{2}{(1-z)^2}$.

6.4 Generating Functions

Computing the k -th derivative of $\sum z^n$.

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The generating function of the sequence $a_n = \binom{n+k}{k}$ is $\frac{1}{(1-z)^{k+1}}$.

6.4 Generating Functions

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The generating function of the sequence $a_n = n$ is $\frac{z}{(1-z)^2}$.

6.4 Generating Functions

We know

$$\sum_{n \geq 0} y^n = \frac{1}{1-y}$$

Hence,

$$\sum_{n \geq 0} a^n z^n = \frac{1}{1-az}$$

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Hence, $a_n = n + 1$.

Some Generating Functions

n-th sequence element	generating function
1	$\frac{1}{1-z}$
$n+1$	$\frac{1}{(1-z)^2}$
$\binom{n+k}{n}$	$\frac{1}{(1-z)^{k+1}}$
n	$\frac{z}{(1-z)^2}$
a^n	$\frac{1}{1-az}$
n^2	$\frac{z(1+z)}{(1-z)^3}$
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n-th sequence element	generating function
cf_n	cF
$f_n + g_n$	$F + G$
$\sum_{i=0}^n f_i g_{n-i}$	$F \cdot G$
f_{n-k} ($n \geq k$); 0 otw.	$z^k F$
$\sum_{i=0}^n f_i$	$\frac{F(z)}{1-z}$
nf_n	$z \frac{dF(z)}{dz}$
$c^n f_n$	$F(cz)$

Some Generating Functions

n-th sequence element	generating function
cf_n	cF
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6. The coefficients of the resulting power series are the a_n .

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which gives

$$A = \frac{7}{4} \quad B = -\frac{1}{4} \quad C = -\frac{1}{2}$$

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6. This means $a_n = \frac{7}{4}3^n - \frac{1}{2}n - \frac{3}{4}$.

6.5 Transformation of the Recurrence

Example 9

$$f_0 = 1$$

$$f_1 = 2$$

$$f_n = f_{n-1} \cdot f_{n-2} \text{ for } n \geq 2 .$$

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Part III

Data Structures

Abstract Data Type

An abstract data type (ADT) is defined by an interface of operations or methods that can be performed and that have a defined behavior

The data types in this lecture all operate on objects that are represented by a **[key, value]** pair.

- ▶ The **key** comes from a totally ordered set, and we assume that there is an efficient comparison function.
- ▶ The **value** can be anything; it usually carries satellite information important for the application that uses the ADT.

Dynamic Set Operations

- ▶ **S .search(k):** Returns pointer to object x from S with $\text{key}[x] = k$ or null.
- ▶ **S .insert(x):** Inserts object x into set S . $\text{key}[x]$ must not currently exist in the data-structure.
- ▶ **S .delete(x):** Given pointer to object x from S , delete x from the set.
- ▶ **S .minimum():** Return pointer to object with smallest key-value in S .
- ▶ **S .maximum():** Return pointer to object with largest key-value in S .
- ▶ **S .successor(x):** Return pointer to the next larger element in S or null if S is maximum.
- ▶ **S .predecessor(x):** Return pointer to the next smaller element in S or null if S is minimum.

Dynamic Set Operations

- ▶ **S. union(S'):** Sets $S := S \cup S'$. The set S' is destroyed.
- ▶ **S. merge(S'):** Sets $S := S \cup S'$. Requires $S \cap S' = \emptyset$.
- ▶ **S. split(k, S'):**
 $S := \{x \in S \mid \text{key}[x] \leq k\}$, $S' := \{x \in S \mid \text{key}[x] > k\}$.
- ▶ **S. concatenate(S'):** $S := S \cup S'$.
Requires $S.\text{maximum}() \leq S'.\text{minimum}()$.
- ▶ **S. decrease-key(x, k):** Replace $\text{key}[x]$ by $k \leq \text{key}[x]$.

Examples of ADTs

Stack:

- ▶ **$S.\text{push}(x)$** : Insert an element.
- ▶ **$S.\text{pop}()$** : Return the element from S that was inserted most recently; delete it from S .
- ▶ **$S.\text{empty}()$** : Tell if S contains any object.

Queue:

- ▶ **$S.\text{enqueue}(x)$** : Insert an element.
- ▶ **$S.\text{dequeue}()$** : Return the element that is longest in the structure; delete it from S .
- ▶ **$S.\text{empty}()$** : Tell if S contains any object.

Priority-Queue:

- ▶ **$S.\text{insert}(x)$** : Insert an element.
- ▶ **$S.\text{delete-min}()$** : Return the element with lowest key-value; delete it from S .

7 Dictionary

Dictionary:

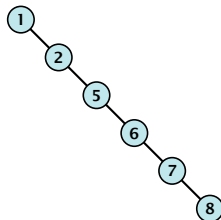
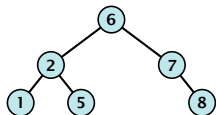
- ▶ **$S.insert(x)$** : Insert an element x .
- ▶ **$S.delete(x)$** : Delete the element pointed to by x .
- ▶ **$S.search(k)$** : Return a pointer to an element e with $key[e] = k$ in S if it exists; otherwise return null.

7.1 Binary Search Trees

An (**internal**) **binary search tree** stores the elements in a binary tree. Each tree-node corresponds to an element. All elements in the left sub-tree of a node v have a smaller key-value than $\text{key}[v]$ and elements in the right sub-tree have a larger-key value. We assume that all key-values are different.

(**External** Search Trees store objects only at leaf-vertices)

Examples:

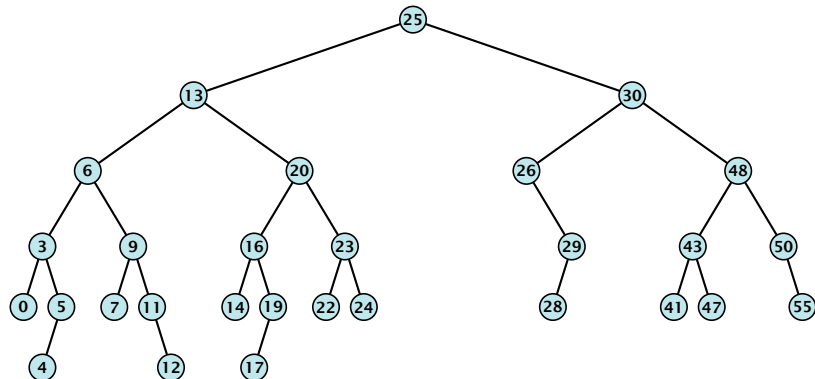


7.1 Binary Search Trees

We consider the following operations on binary search trees. Note that this is a super-set of the dictionary-operations.

- ▶ $T.\text{insert}(x)$
- ▶ $T.\text{delete}(x)$
- ▶ $T.\text{search}(k)$
- ▶ $T.\text{successor}(x)$
- ▶ $T.\text{predecessor}(x)$
- ▶ $T.\text{minimum}()$
- ▶ $T.\text{maximum}()$

Binary Search Trees: Searching

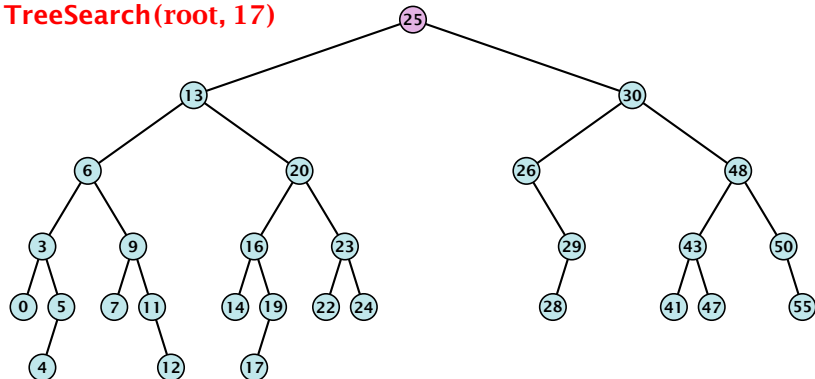


Algorithm 5 $\text{TreeSearch}(x, k)$

- 1: **if** $x = \text{null}$ **or** $k = \text{key}[x]$ **return** x
- 2: **if** $k < \text{key}[x]$ **return** $\text{TreeSearch}(\text{left}[x], k)$
- 3: **else return** $\text{TreeSearch}(\text{right}[x], k)$

Binary Search Trees: Searching

TreeSearch(root, 17)

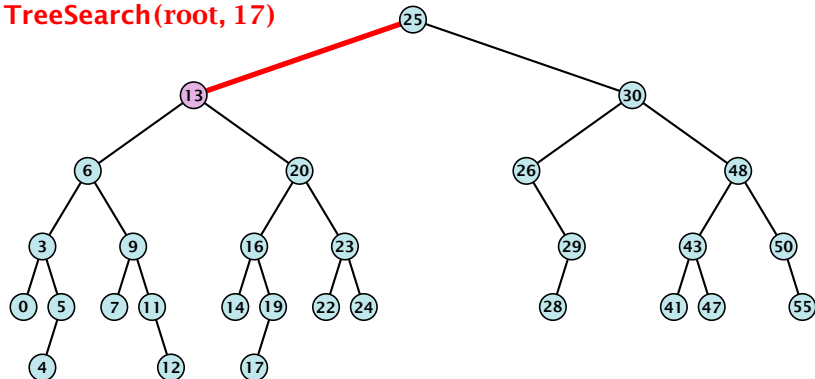


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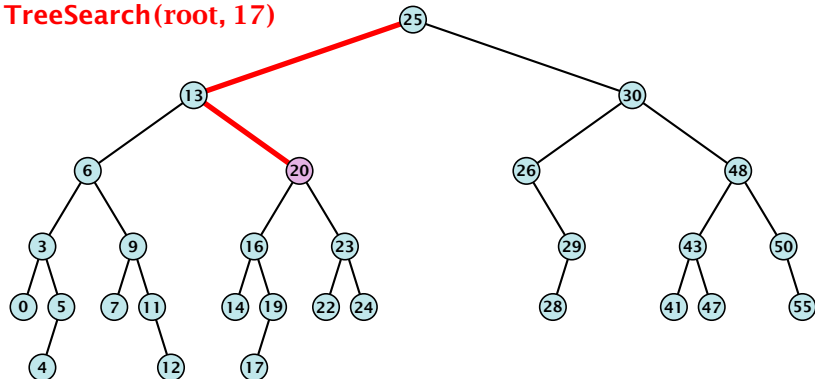


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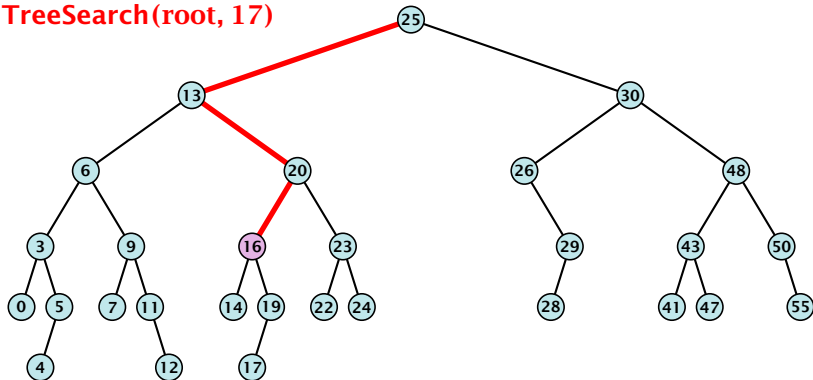


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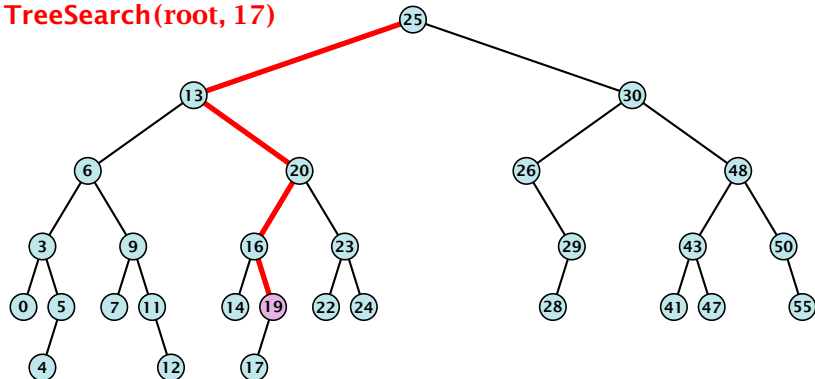


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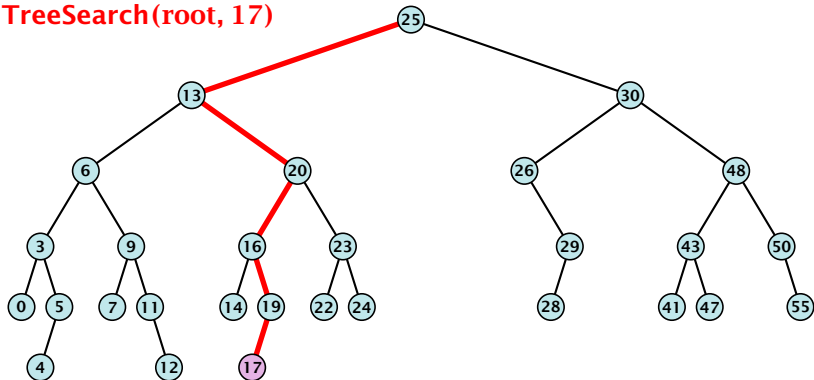


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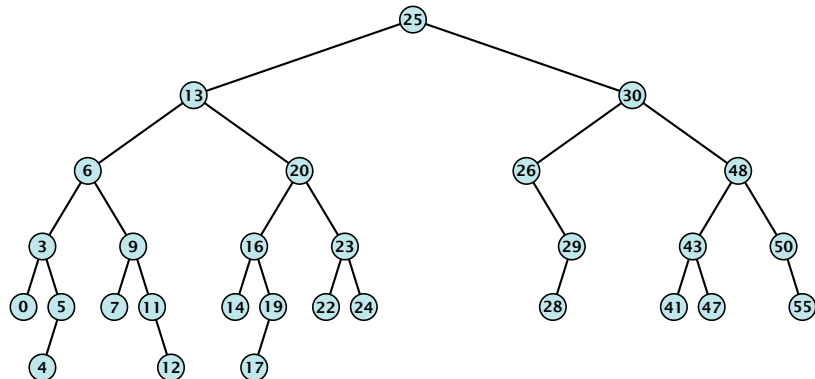
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Binary Search Trees: Searching

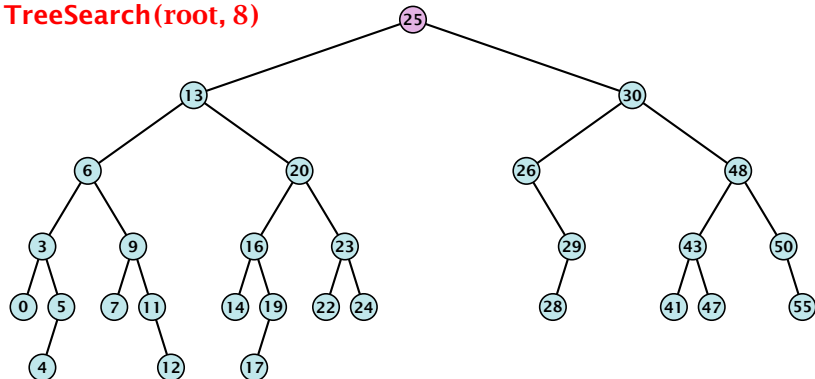


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Binary Search Trees: Searching

TreeSearch(root, 8)

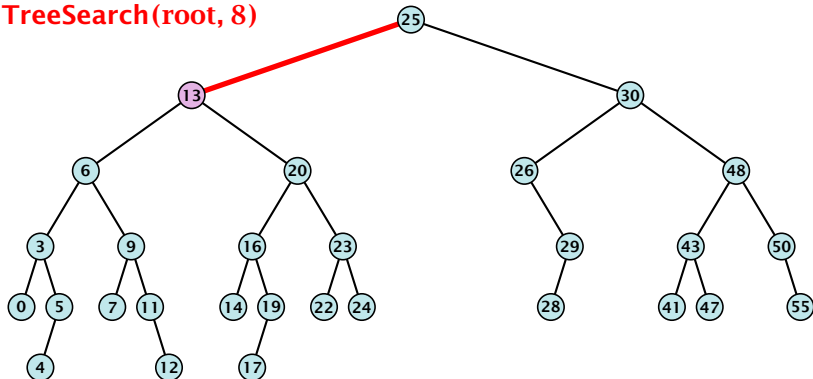


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Binary Search Trees: Searching

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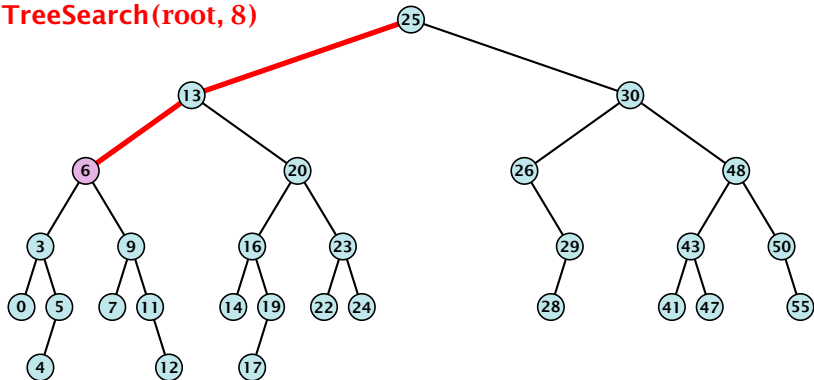


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Binary Search Trees: Searching

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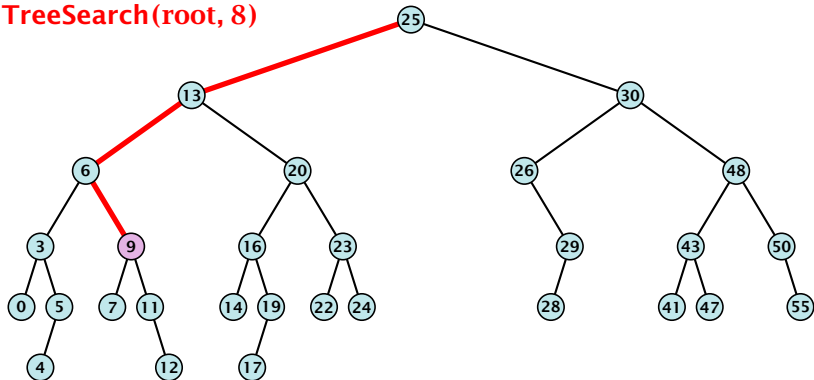


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Binary Search Trees: Searching

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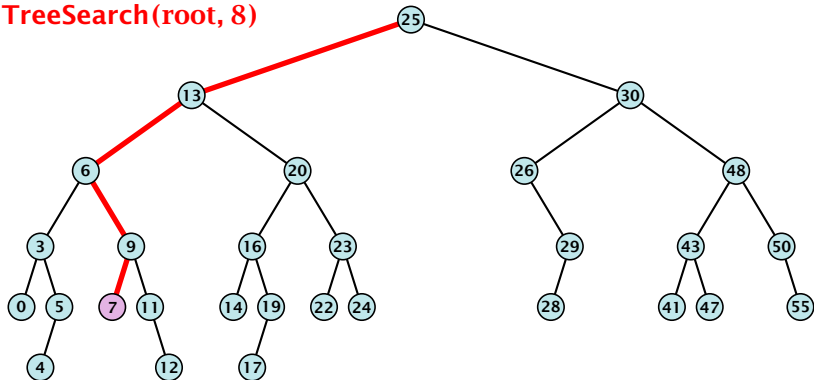


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Binary Search Trees: Searching

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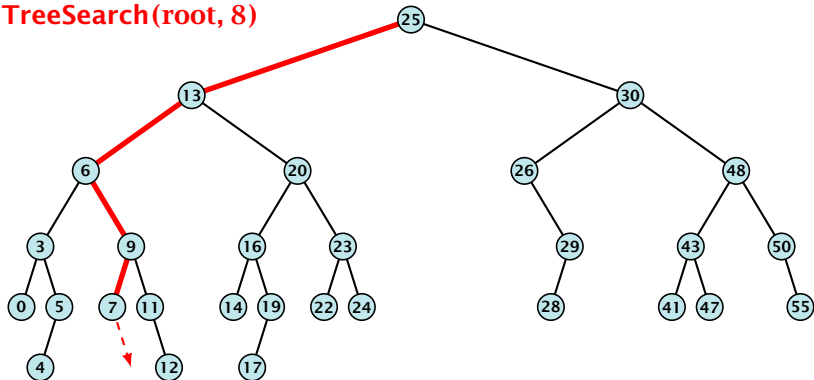


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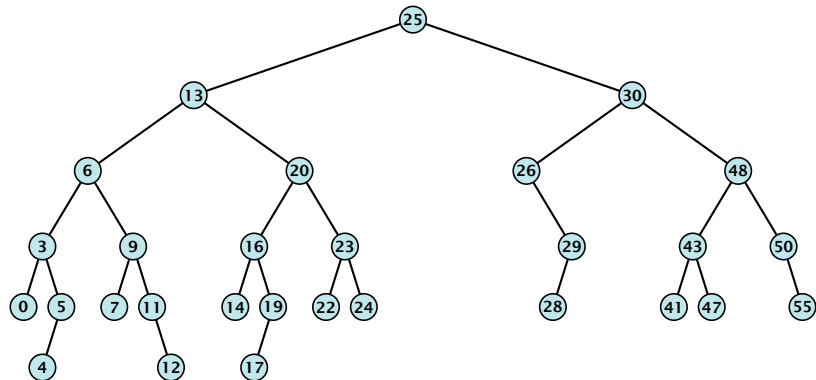
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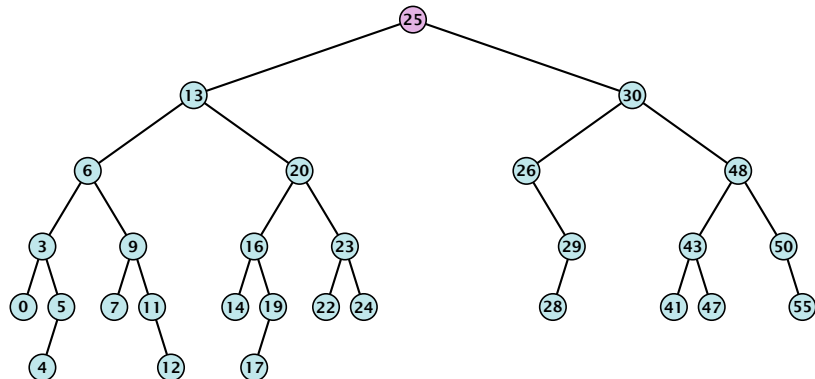
Binary Search Trees: Minimum



Algorithm 6 TreeMin(x)

- 1: **if** $x = \text{null}$ **or** $\text{left}[x] = \text{null}$ **return** x
- 2: **if** $k < \text{key}[x]$ **return** $\text{TreeSearch}(\text{left}[x], k)$
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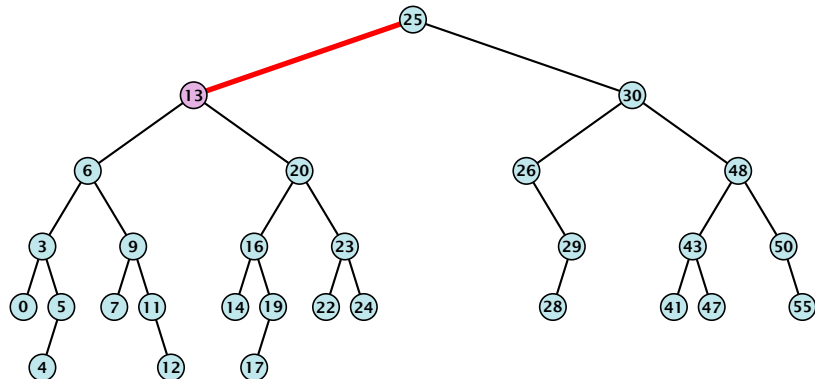
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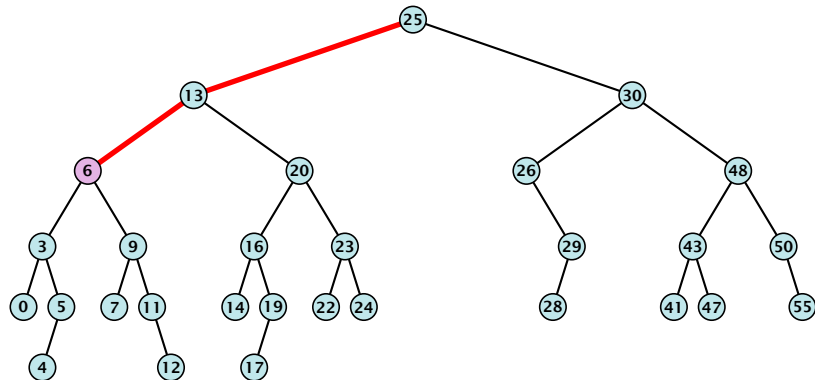
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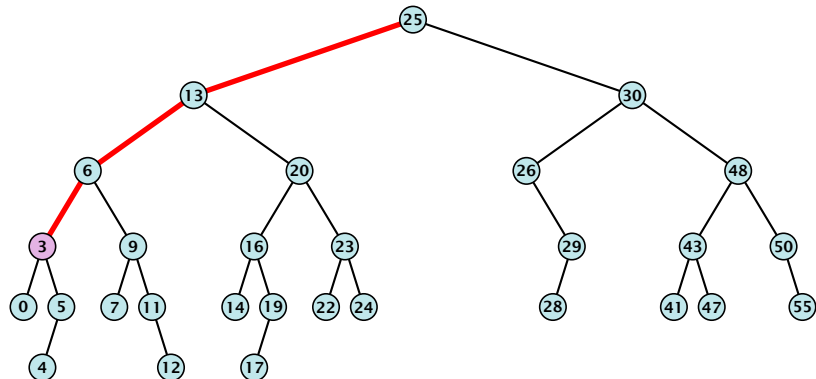
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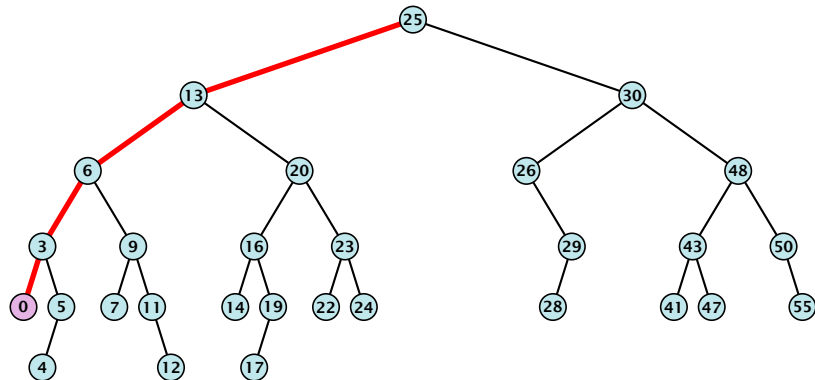
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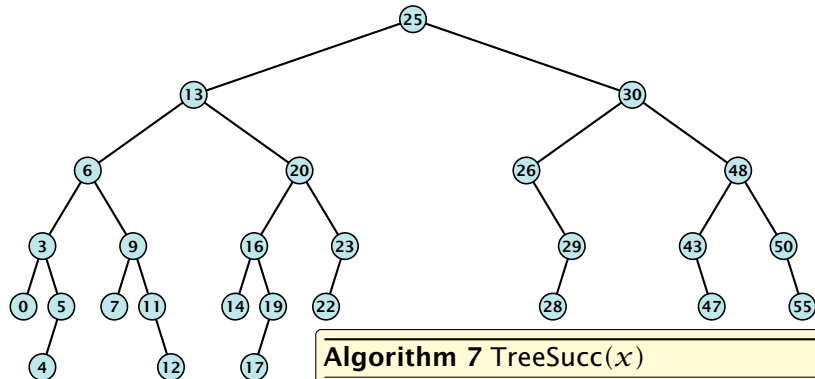
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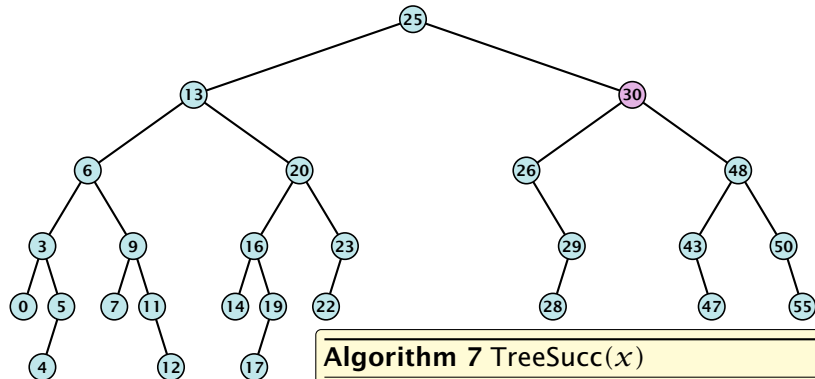
Binary Search Trees: Successor



Algorithm 7 TreeSucc(x)

- 1: **if** right[x] \neq null **return** TreeMin(right[x])
- 2: $y \leftarrow$ parent[x]
- 3: **while** $y \neq$ null **and** $x =$ right[y] **do**
- 4: $x \leftarrow y$; $y \leftarrow$ parent[x]
- 5: **return** y ;

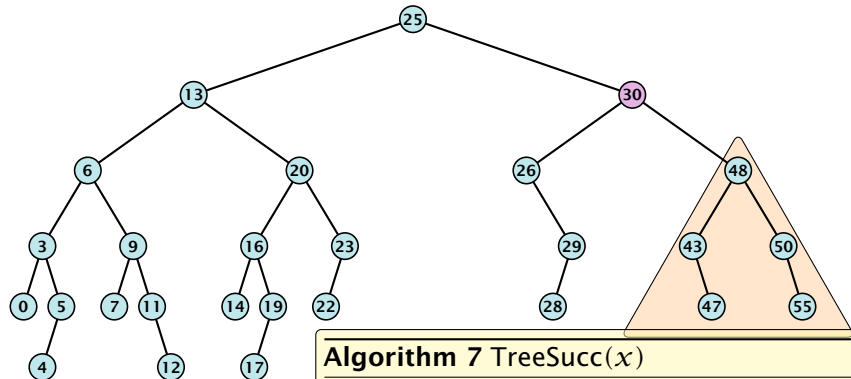
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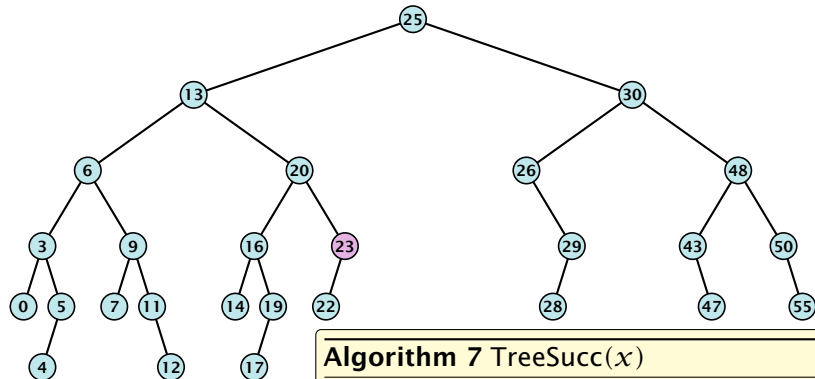
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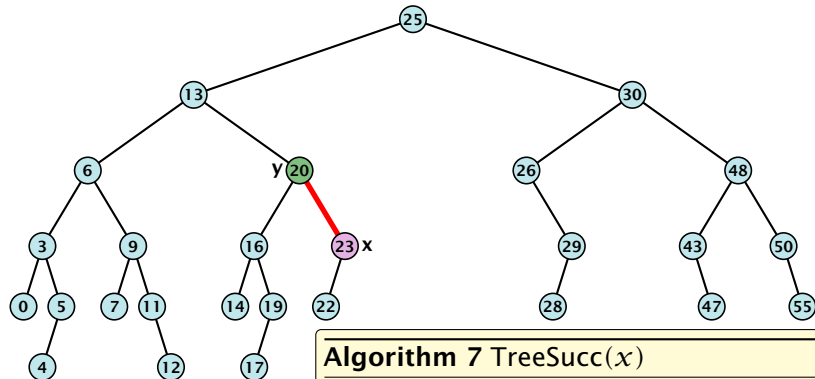
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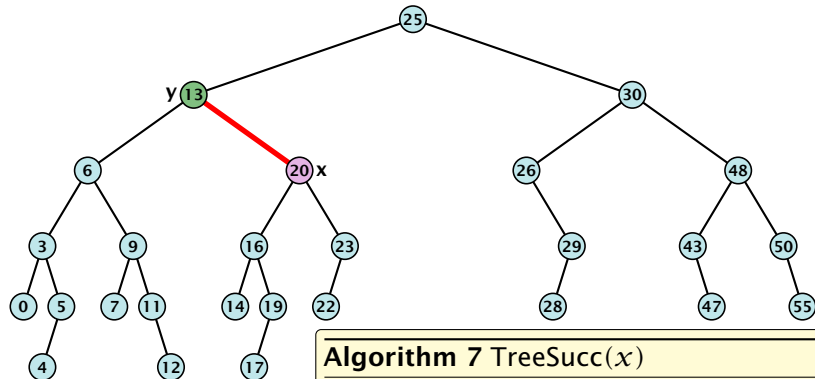
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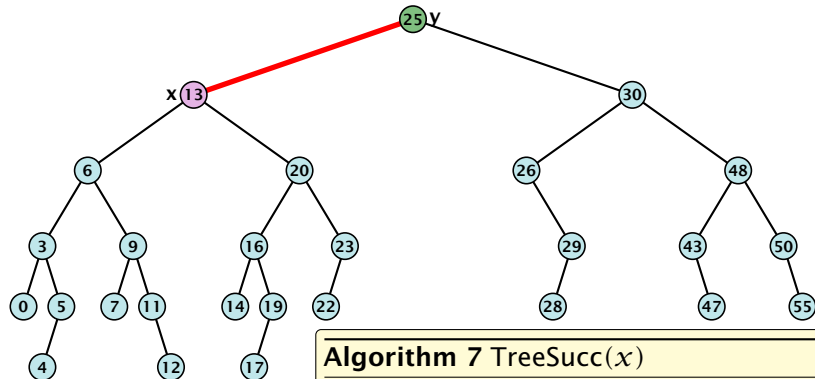
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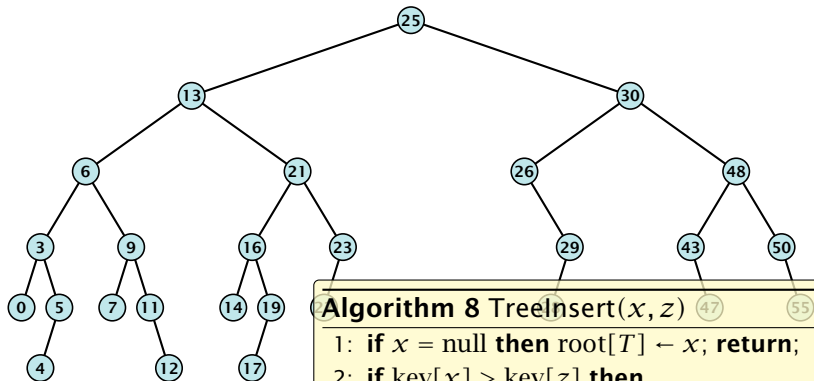
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Binary Search Trees: Insert

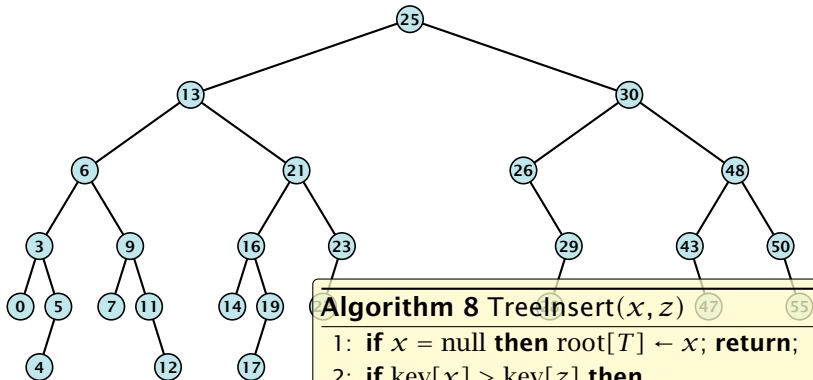


Algorithm 8 TreeInsert(x, z)

- 1: **if** $x = \text{null}$ **then** $\text{root}[T] \leftarrow x$; **return**;
- 2: **if** $\text{key}[x] > \text{key}[z]$ **then**
- 3: **if** $\text{left}[x] = \text{null}$ **then** $\text{left}[x] \leftarrow z$;
- 4: **else** TreeInsert($\text{left}[x], z$);
- 5: **else**
- 6: **if** $\text{right}[x] = \text{null}$ **then** $\text{right}[x] \leftarrow z$;
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Binary Search Trees: Insert

Insert element **not** in the tree.

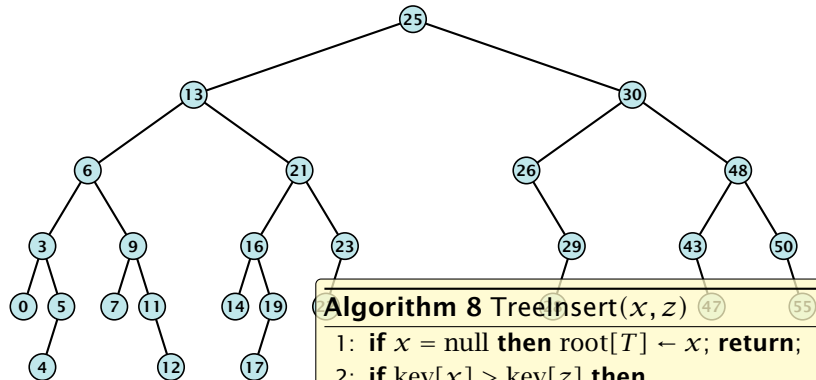


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Search for z . At some point the search stops at a null-pointer. This is the place to insert z .

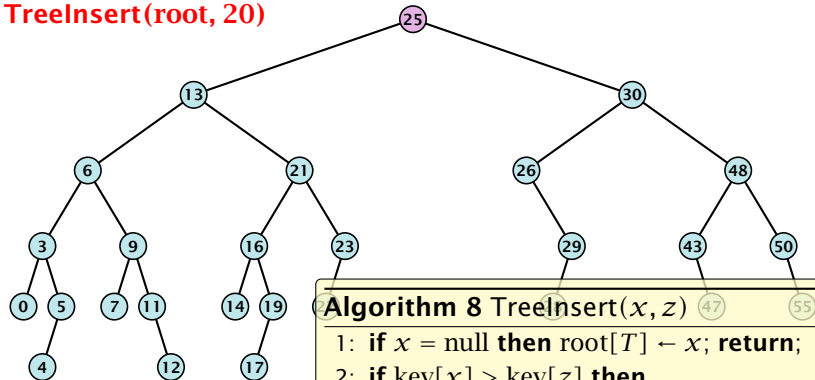
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Binary Search Trees: Insert

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TreeInsert(root, 20)



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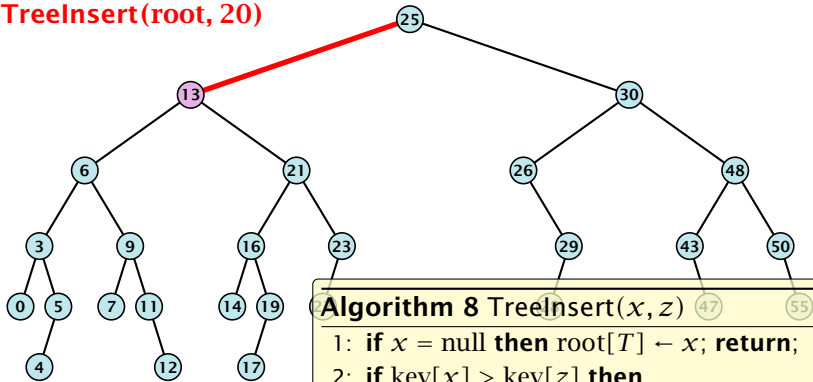
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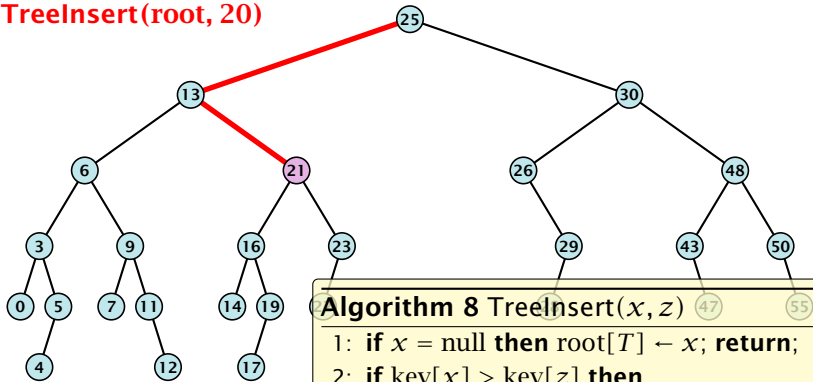
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Search for z . At some point the search stops at a null-pointer. This is the place to insert z .

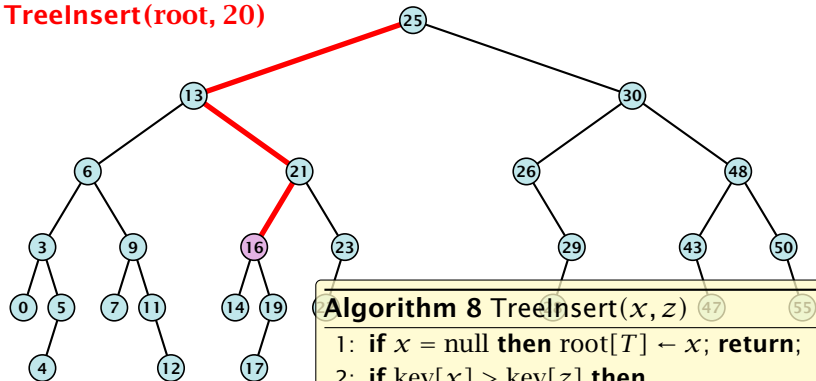
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- 1: **if** $x = \text{null}$ **then** $\text{root}[T] \leftarrow x$; **return**;
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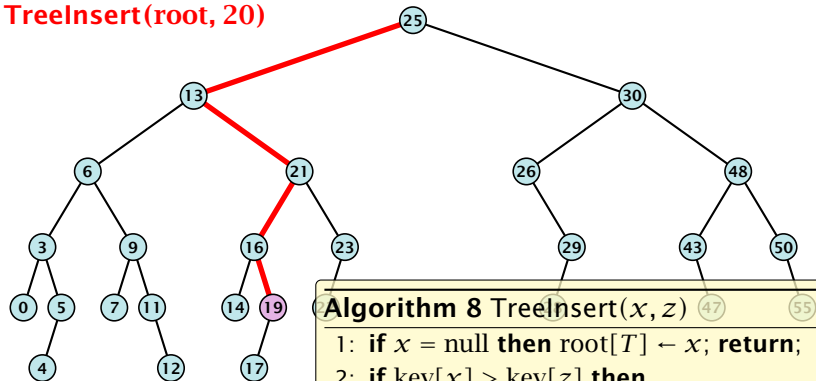
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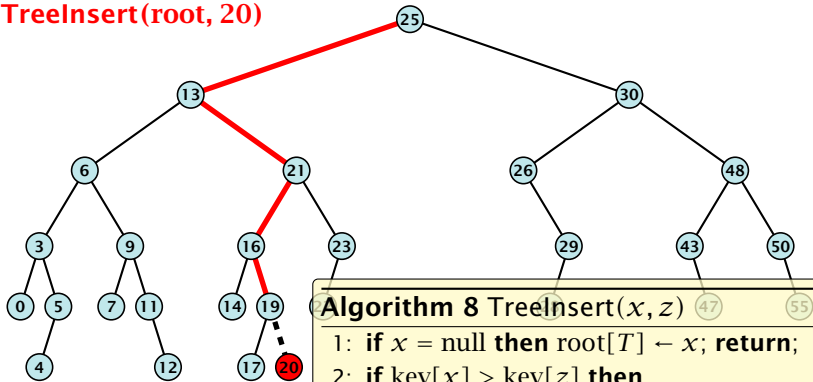
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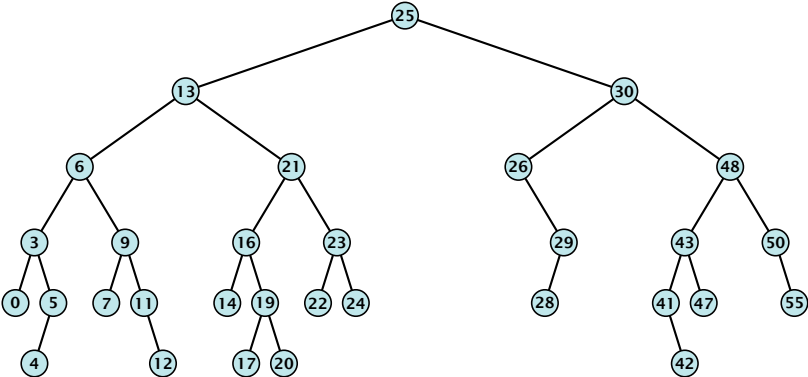


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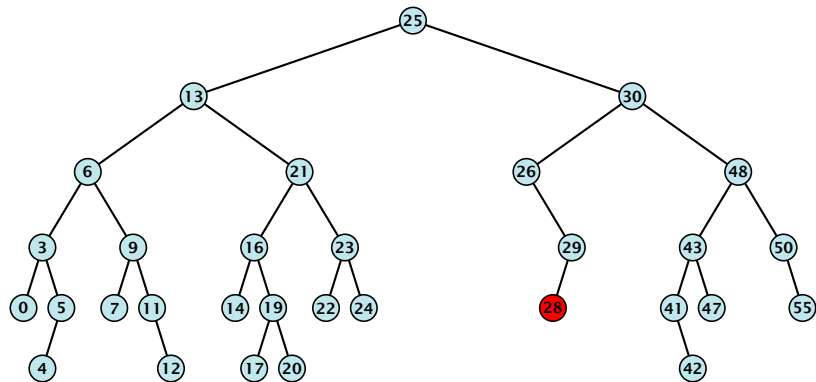
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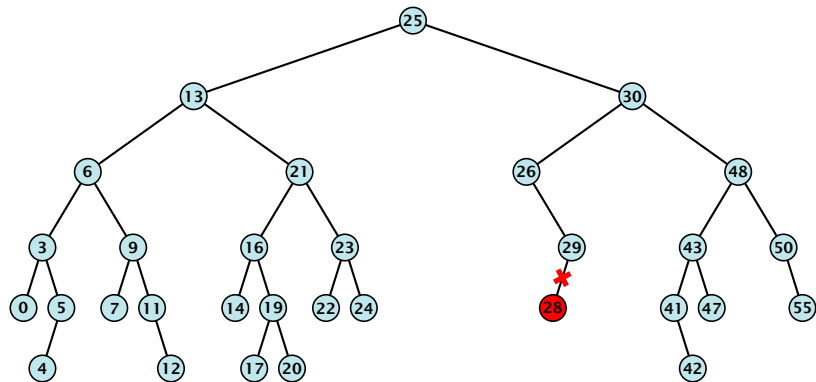


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Binary Search Trees: Delete

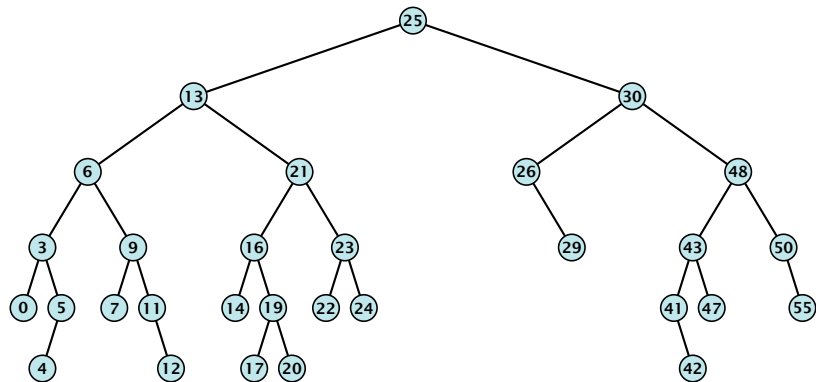


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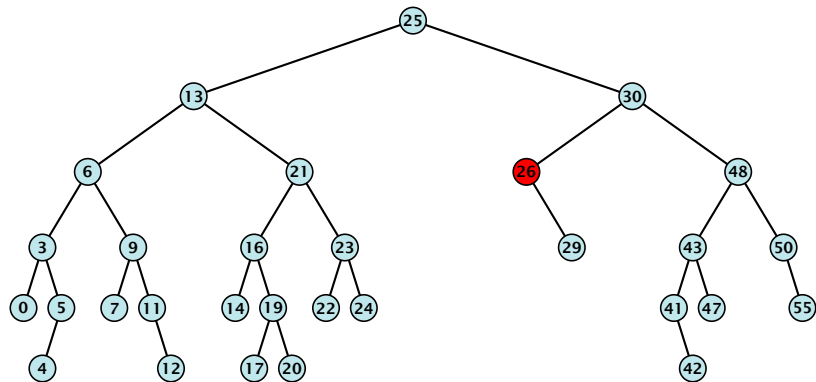


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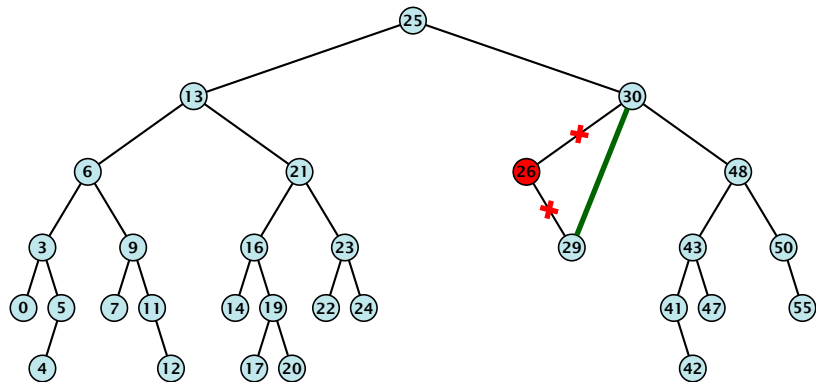


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Element has exactly one child

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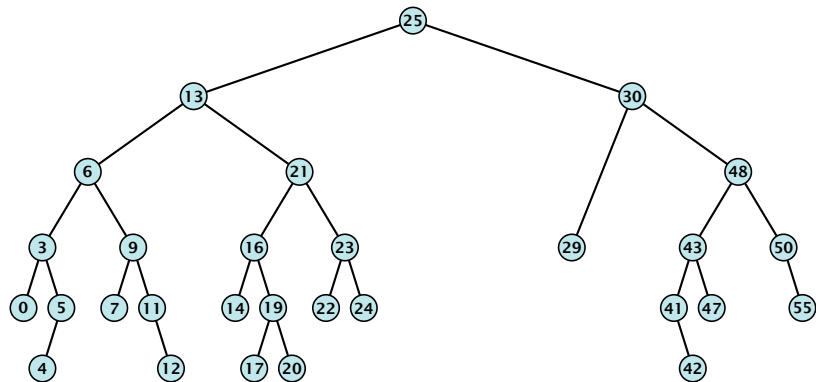


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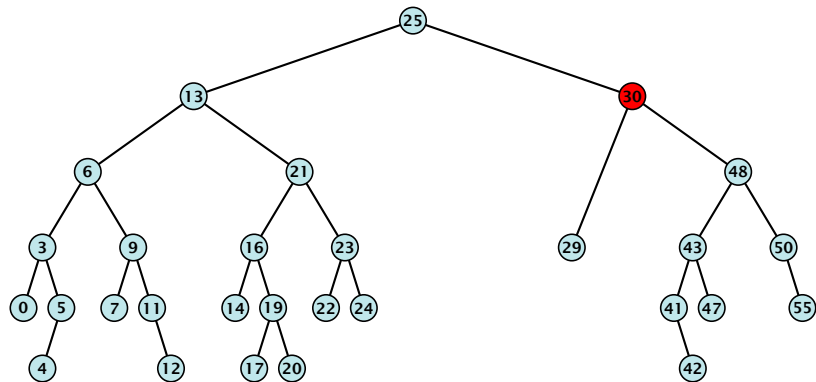


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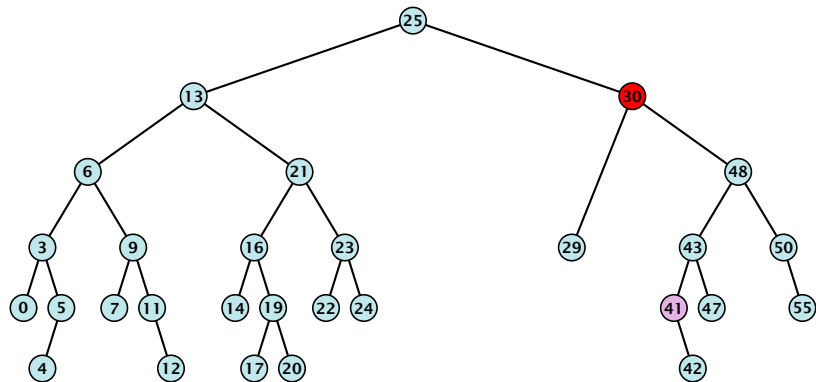


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Element has two children

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- ▶ Splice successor out of the tree
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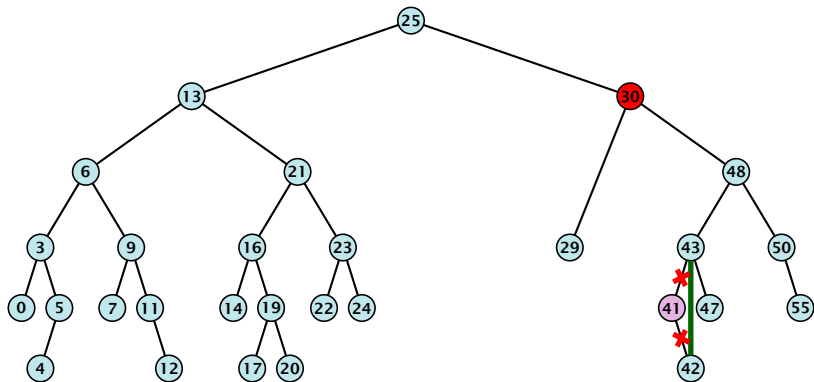


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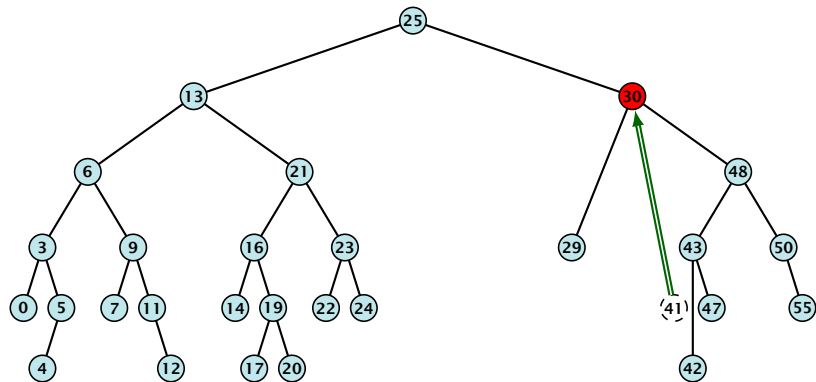


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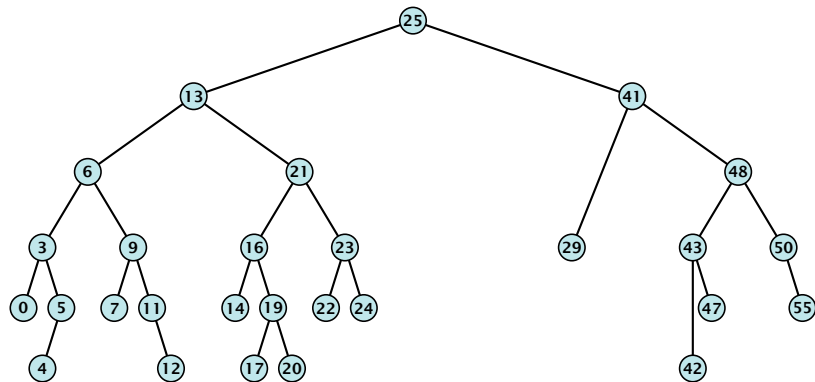


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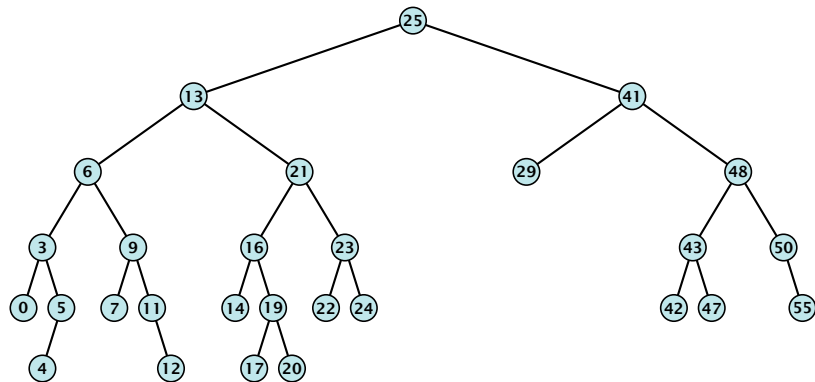


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Algorithm 9 TreeDelete(z)

```
1: if left[ $z$ ] = null or right[ $z$ ] = null
2:   then  $y \leftarrow z$  else  $y \leftarrow \text{TreeSucc}(z)$ ;   select  $y$  to splice out
3:   if left[ $y$ ]  $\neq$  null
4:     then  $x \leftarrow \text{left}[y]$  else  $x \leftarrow \text{right}[y]$ ;  $x$  is child of  $y$  (or null)
5:   if  $x \neq \text{null}$  then parent[ $x$ ]  $\leftarrow$  parent[ $y$ ];   parent[ $x$ ] is correct
6:   if parent[ $y$ ] = null then
7:     root[ $T$ ]  $\leftarrow x$ 
8:   else
9:     if  $y = \text{left}[\text{parent}[x]]$  then
10:      left[parent[ $y$ ]]  $\leftarrow x$ 
11:     else
12:      right[parent[ $y$ ]]  $\leftarrow x$ 
13:   if  $y \neq z$  then copy  $y$ -data to  $z$ 
```

} fix pointer to x

Balanced Binary Search Trees

All operations on a binary search tree can be performed in time $\mathcal{O}(h)$, where h denotes the height of the tree.

However the height of the tree may become as large as $\Theta(n)$.

Balanced Binary Search Trees

With each insert- and delete-operation perform **local** adjustments to guarantee a height of $\mathcal{O}(\log n)$.

AVL-trees, Red-black trees, Scapegoat trees, 2-3 trees, B-trees, AA trees, Treaps

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7.2 Red Black Trees

Definition 11

A red black tree is a balanced binary search tree in which each internal node has two children. Each internal node has a colour, such that

1. The root is black.
2. All leaf nodes are black.
3. For each node, all paths to descendant leaves contain the same number of black nodes.
4. If a node is red then both its children are black.

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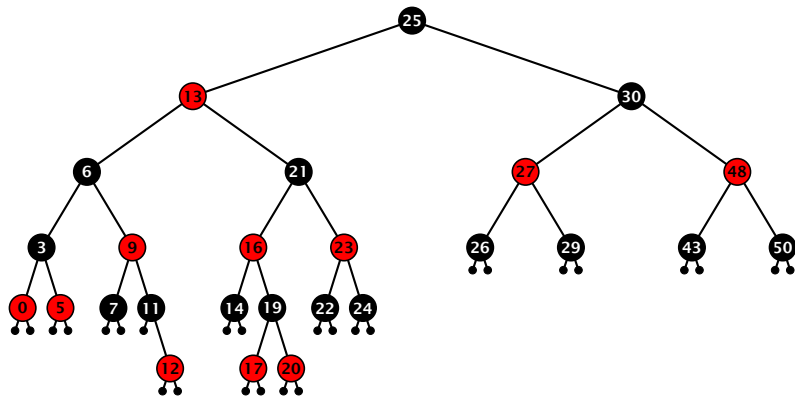
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Red Black Trees: Example



7.2 Red Black Trees

Lemma 12

A red-black tree with n internal nodes has height at most $\mathcal{O}(\log n)$.

Definition 13

The black height $\text{bh}(v)$ of a node v in a red black tree is the number of black nodes on a path from v to a leaf vertex (not counting v).

We first show:

Lemma 14

A sub-tree of black height $\text{bh}(v)$ in a red black tree contains at least $2^{\text{bh}(v)} - 1$ internal vertices.

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Induction on the height of v .

base case ($\text{height}(v) = 0$)

- if $\text{height}(v)$ (maximum distance from v and a node in the subtree rooted at v) is 0 then v is a leaf.
- The black height of v is 0.
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Proof (cont.)

induction step

- Suppose v is a node with height $|v| > 0$.
- v has two children with strictly smaller height.
- These children (v_l, v_r) either have $bal(v_l) = bal(v_r) = 0$ or $bal(v_l) = bal(v_r) = \pm 1$.
- By induction hypothesis both subtrees contain at least $\frac{1}{2} \cdot 2^{|v_l|} - 1$ internal nodes.
- Thus T_v contains at least $2(\frac{1}{2} \cdot 2^{|v_l|} - 1) + 1 = 2^{|v_l|} - 1$



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- ▶ Suppose v is a node with $\text{height}(v) > 0$.
- ▶ v has two children with strictly smaller height.
- ▶ These children (c_1, c_2) either have $\text{bh}(c_i) = \text{bh}(v)$ or $\text{bh}(c_i) = \text{bh}(v) - 1$.
- ▶ By induction hypothesis both sub-trees contain at least $2^{\text{bh}(v)-1} - 1$ internal vertices.
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Let h denote the height of the red-black tree, and let p denote a path from the root to the furthest leaf.

At least half of the nodes on p must be black, since a red node must be followed by a black node.

Hence, the black height of the root is at least $h/2$.

The tree contains at least $2^{h/2} - 1$ internal vertices. Hence,
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Hence, $h \leq 2 \log n + 1 = \mathcal{O}(\log n)$. □

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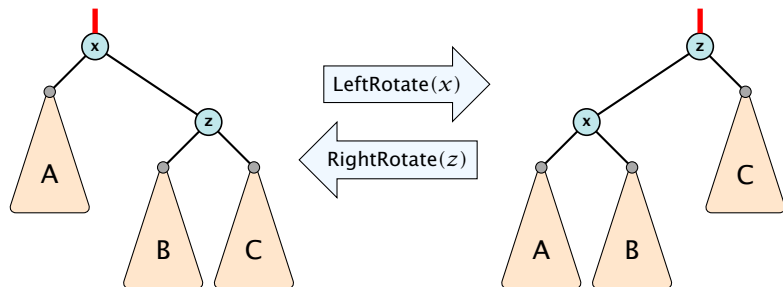
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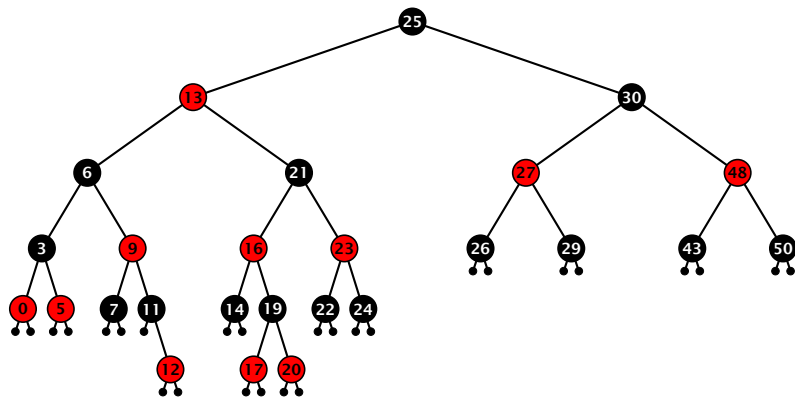
We need to adapt the insert and delete operations so that the red black properties are maintained.

Rotations

The properties will be maintained through rotations:



Red Black Trees: Insert

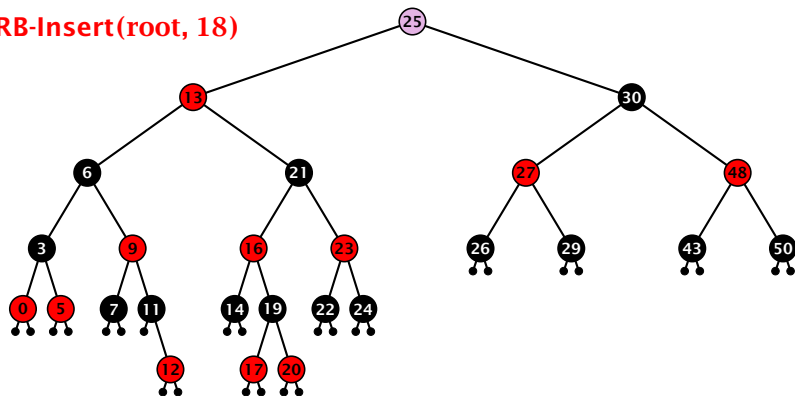


Insert:

- ▶ first make a normal insert into a binary search tree
- ▶ then fix red-black properties

Red Black Trees: Insert

RB-Insert(root, 18)

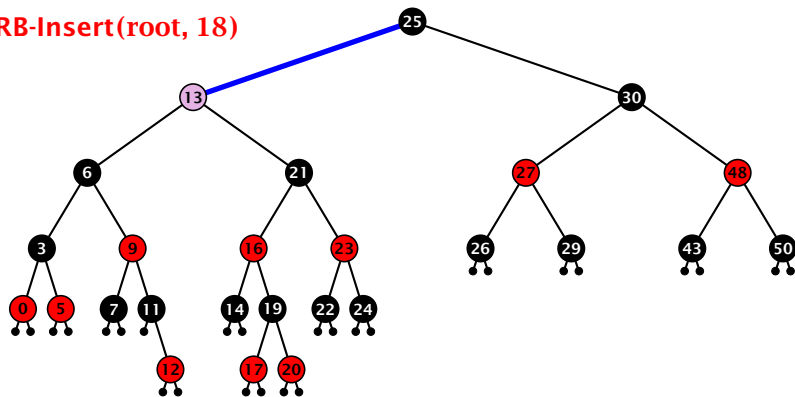


Insert:

- ▶ first make a normal insert into a binary search tree
- ▶ then fix red-black properties

Red Black Trees: Insert

RB-Insert(root, 18)

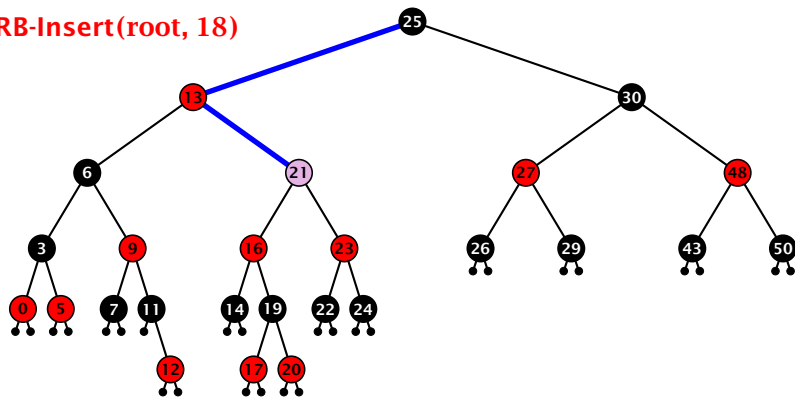


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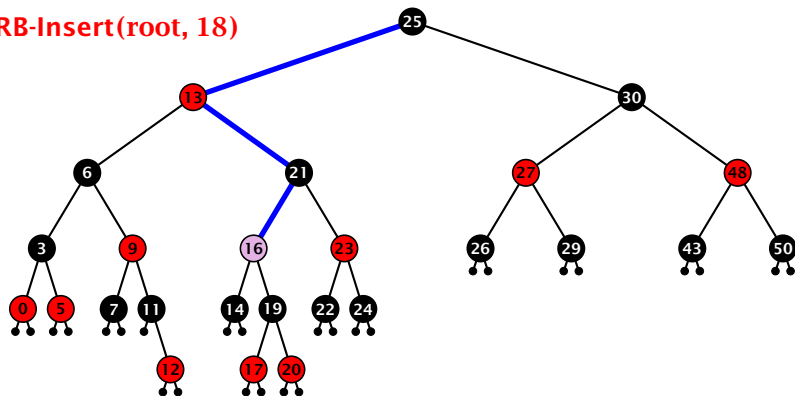


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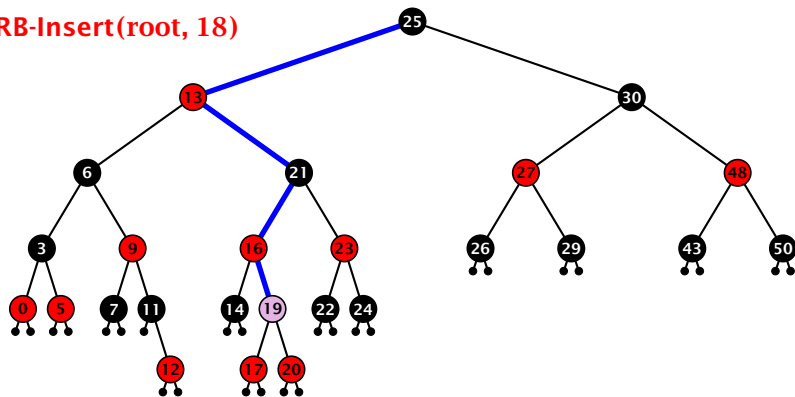


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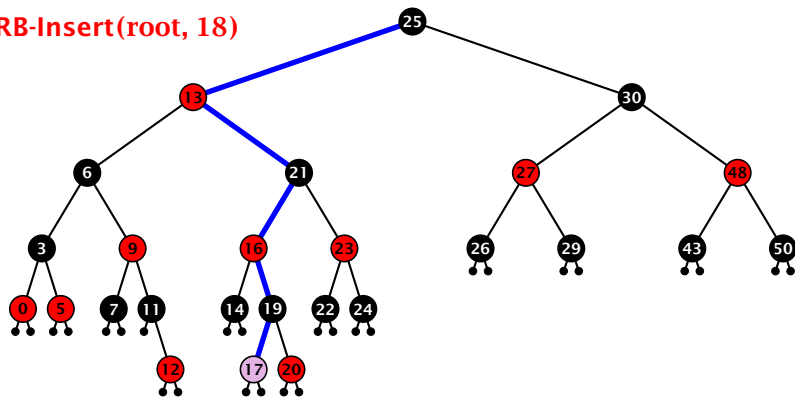


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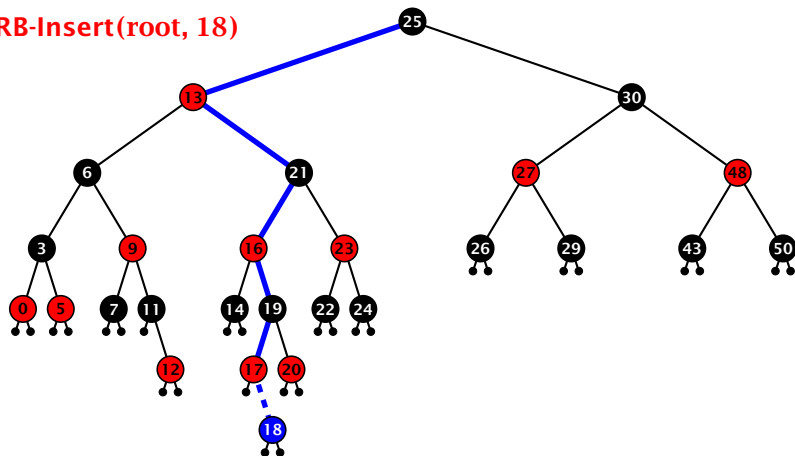


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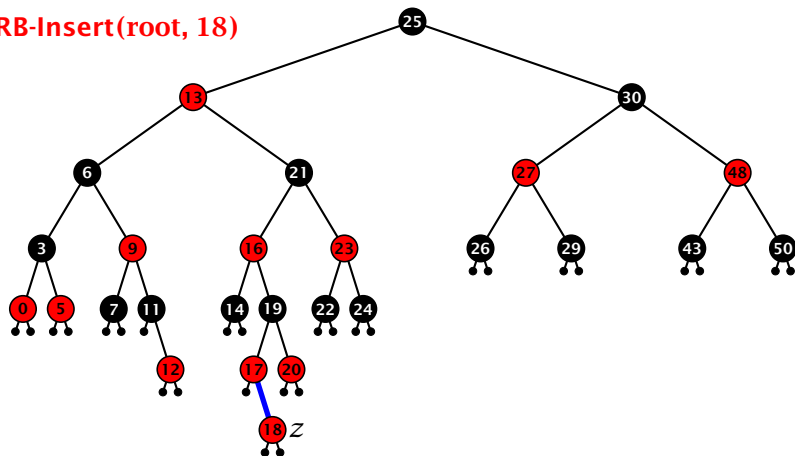


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Red Black Trees: Insert

RB-Insert(root, 18)



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- ▶ then fix red-black properties

Red Black Trees: Insert

Invariant of the fix-up algorithm:

- ▶ z is a red node
- ▶ the black-height property is fulfilled at every node
- ▶ the only violation of red-black properties occurs at z and $\text{parent}[z]$
 - either both of them are red (most important case)
 - or the parent does not exist (violation does not need to be fixed)

If z has a parent but no grand-parent we could simply color the parent/root black; however this case never happens.

Red Black Trees: Insert

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- ▶ z is a red node
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- ▶ the only violation of red-black properties occurs at z and $\text{parent}[z]$

(either both of them are red

or both are black, important case)

(if the parent is red, not color

violations since root must be black)

If z has a parent but no grand-parent we could simply color the parent/root black; however this case never happens.

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Red Black Trees: Insert

Algorithm 10 InsertFix(z)

```
1: while parent[ $z$ ]  $\neq$  null and col[parent[ $z$ ]] = red do
2:   if parent[ $z$ ] = left[gp[ $z$ ]] then
3:      $uncle \leftarrow$  right[grandparent[ $z$ ]]
4:     if col[ $uncle$ ] = red then
5:       col[p[ $z$ ]]  $\leftarrow$  black; col[ $u$ ]  $\leftarrow$  black;
6:       col[gp[ $z$ ]]  $\leftarrow$  red;  $z \leftarrow$  grandparent[ $z$ ];
7:     else
8:       if  $z$  = right[parent[ $z$ ]] then
9:          $z \leftarrow$  p[ $z$ ]; LeftRotate( $z$ );
10:      col[p[ $z$ ]]  $\leftarrow$  black; col[gp[ $z$ ]]  $\leftarrow$  red;
11:      RightRotate(gp[ $z$ ]);
12:     else same as then-clause but right and left exchanged
13: col(root[ $T$ ])  $\leftarrow$  black;
```

Red Black Trees: Insert

Algorithm 10 InsertFix(z)

```
1: while parent[ $z$ ]  $\neq$  null and col[parent[ $z$ ]] = red do
2:   if parent[ $z$ ] = left[gp[ $z$ ]] then  $z$  in left subtree of grandparent
3:      $uncle \leftarrow$  right[grandparent[ $z$ ]]
4:     if col[ $uncle$ ] = red then
5:       col[p[ $z$ ]]  $\leftarrow$  black; col[ $u$ ]  $\leftarrow$  black;
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1: while parent[ $z$ ]  $\neq$  null and col[parent[ $z$ ]] = red do
2:   if parent[ $z$ ] = left[gp[ $z$ ]] then
3:     uncle  $\leftarrow$  right[grandparent[ $z$ ]]
4:     if col[uncle] = red then Case 1: uncle red
5:       col[p[ $z$ ]]  $\leftarrow$  black; col[u]  $\leftarrow$  black;
6:       col[gp[ $z$ ]]  $\leftarrow$  red;  $z$   $\leftarrow$  grandparent[ $z$ ];
7:     else
8:       if  $z$  = right[parent[ $z$ ]] then
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6:       col[gp[ $z$ ]]  $\leftarrow$  red;  $z \leftarrow$  grandparent[ $z$ ];
7:   else Case 2: uncle black
8:     if  $z$  = right[parent[ $z$ ]] then
9:        $z \leftarrow$  p[ $z$ ]; LeftRotate( $z$ );
10:    col[p[ $z$ ]]  $\leftarrow$  black; col[gp[ $z$ ]]  $\leftarrow$  red;
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Red Black Trees: Insert

Algorithm 10 InsertFix(z)

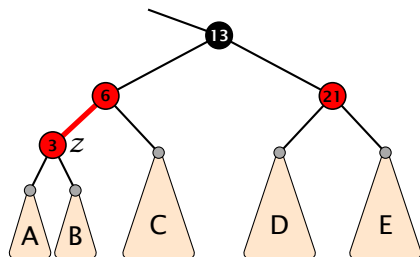
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1: while parent[ $z$ ]  $\neq$  null and col[parent[ $z$ ]] = red do
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7:     else
8:       if  $z =$  right[parent[ $z$ ]] then 2a:  $z$  right child
9:          $z \leftarrow$  p[ $z$ ]; LeftRotate( $z$ );
10:        col[p[ $z$ ]]  $\leftarrow$  black; col[gp[ $z$ ]]  $\leftarrow$  red;
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3:     uncle  $\leftarrow$  right[grandparent[ $z$ ]]
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10:      col[p[ $z$ ]]  $\leftarrow$  black; col[gp[ $z$ ]]  $\leftarrow$  red; 2b:  $z$  left child
11:      RightRotate(gp[ $z$ ]);
12:   else same as then-clause but right and left exchanged
13: col(root[ $T$ ])  $\leftarrow$  black;
```

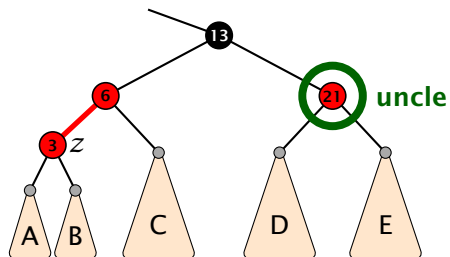
Case 1: Red Uncle



1. recolour
2. move z to grand-parent
3. invariant is fulfilled for new z
4. you made progress



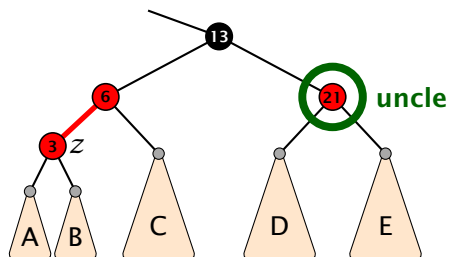
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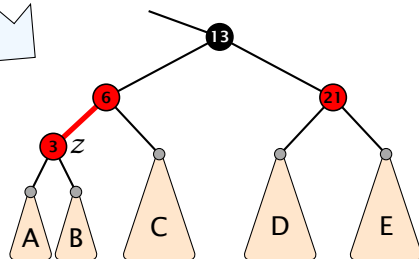
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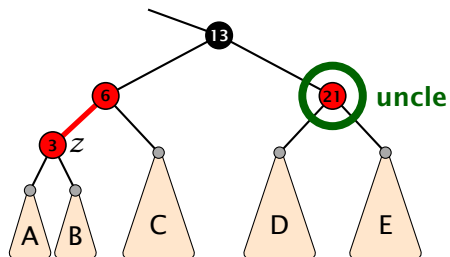
Case 1: Red Uncle



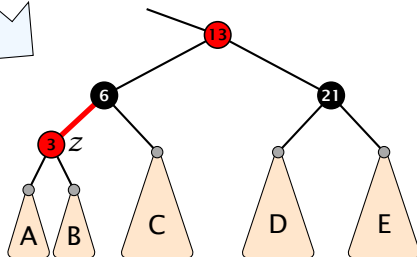
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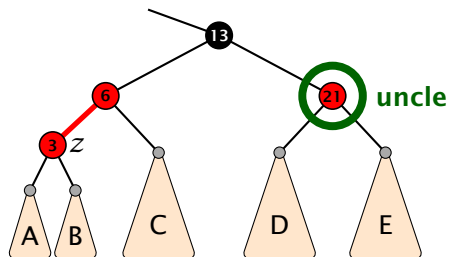
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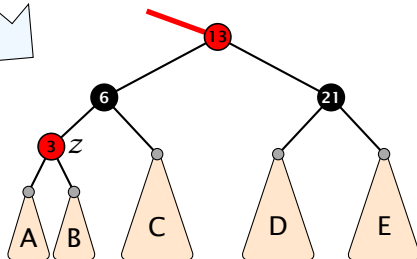
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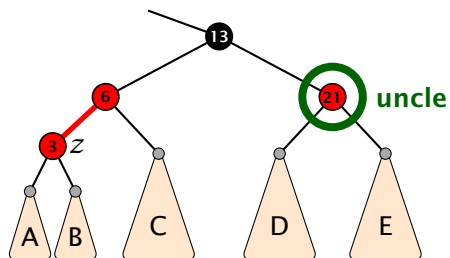
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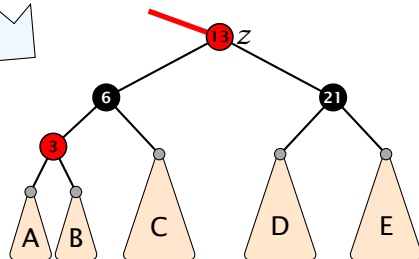
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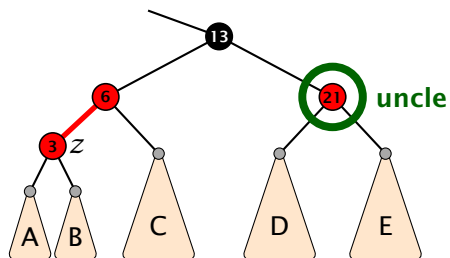
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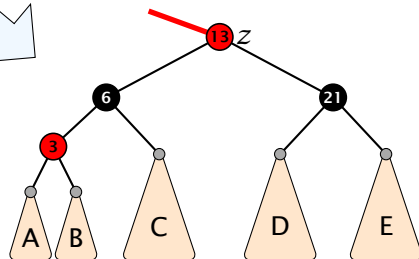
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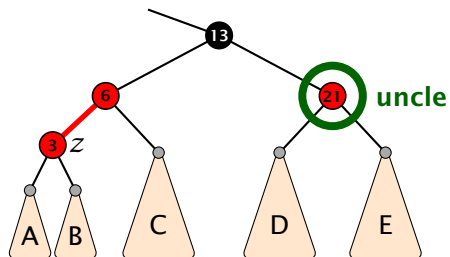
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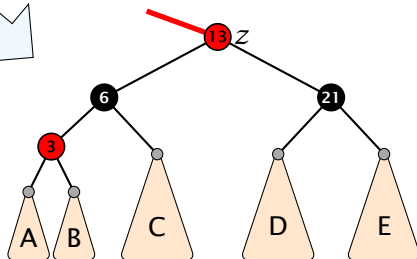
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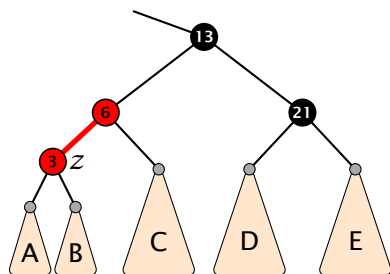


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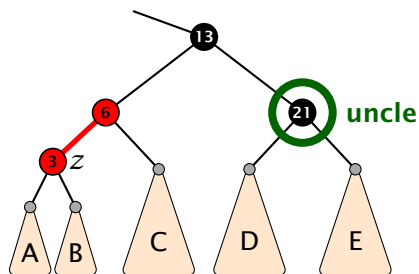
Case 2b: Black uncle and z is left child

1. rotate around grandparent
2. re-colour to ensure that black height property holds
3. you have a red black tree



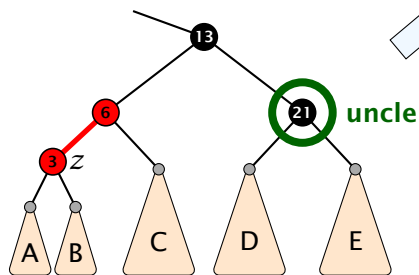
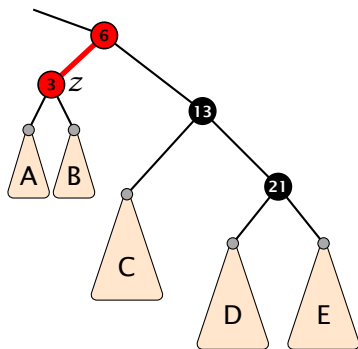
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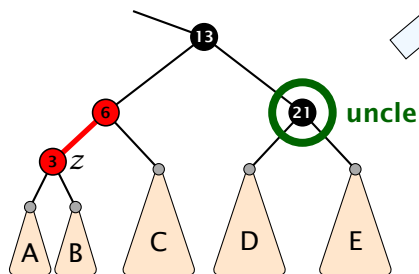
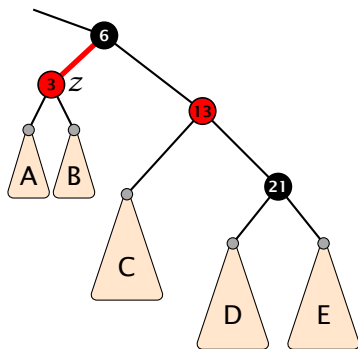
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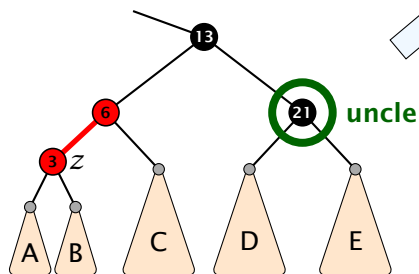
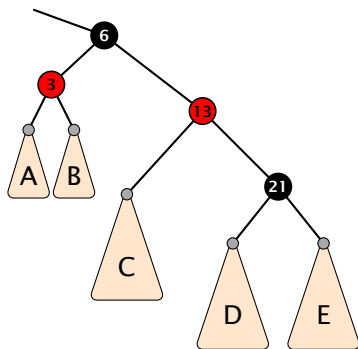
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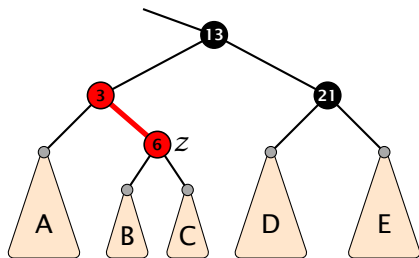
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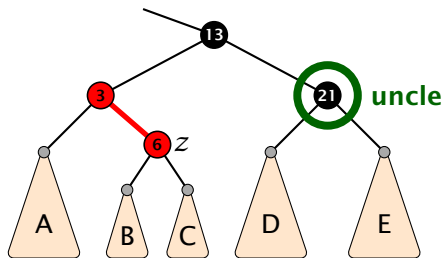
Case 2a: Black uncle and z is right child

1. rotate around parent
2. move z downwards
3. you have case 2b.



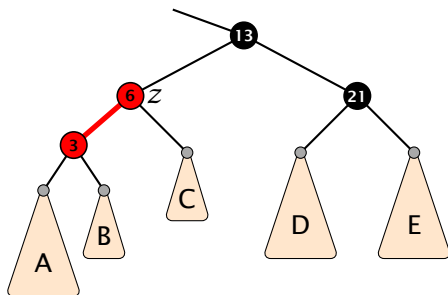
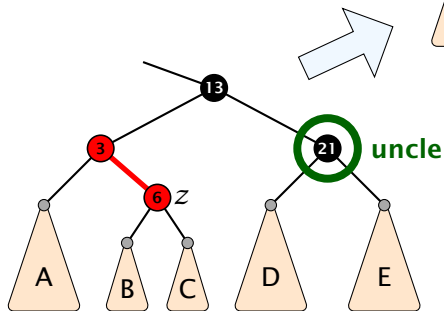
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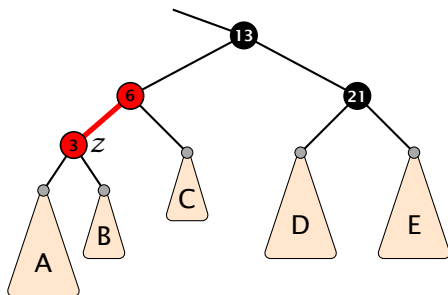
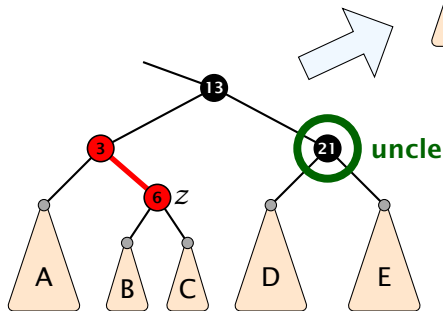
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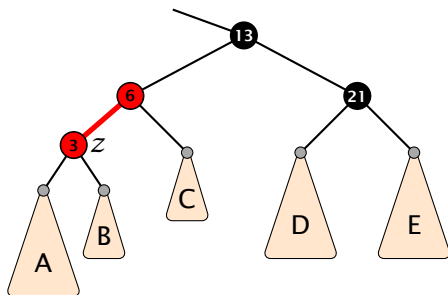
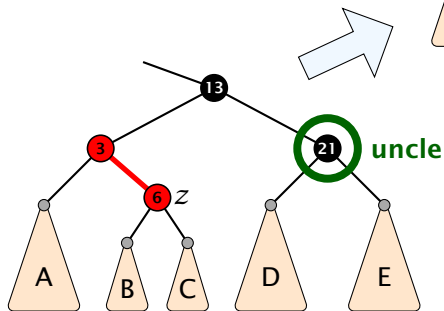
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Red Black Trees: Insert

Running time:

- ▶ Only Case 1 may repeat; but only $h/2$ many steps, where h is the height of the tree.
- ▶ Case 2a → Case 2b → red-black tree
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Performing step one $\mathcal{O}(\log n)$ times and every other step at most once, we get a red-black tree. Hence $\mathcal{O}(\log n)$ re-colourings and at most 2 rotations.

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First do a standard delete.

If the spliced out node x was red everything is fine.

If it was black there may be the following problems.

1. Parent and child of x were red; two adjacent red vertices.

2. If you delete the root, the root may now be red.

3. Every path from an ancestor of x to a descendant leaf of x changes the number of black nodes. Black height property might be violated.

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If it was black there may be the following problems.

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2. If x was the root, the root may now be red.

3. Every path from an ancestor of x to a descendant leaf of x changes the number of black nodes. Black height property may not be violated.

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• Parent and child of x were red, two adjacent red vertices.

• x was the root of the tree, the root may now be red.

• x was the left or right child of a red vertex, the balance

changes, the number of black nodes, Black Height, property

is not preserved.

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- ▶ Every path from an ancestor of x to a descendant leaf of x changes the number of black nodes. Black height property might be violated.

Red Black Trees: Delete

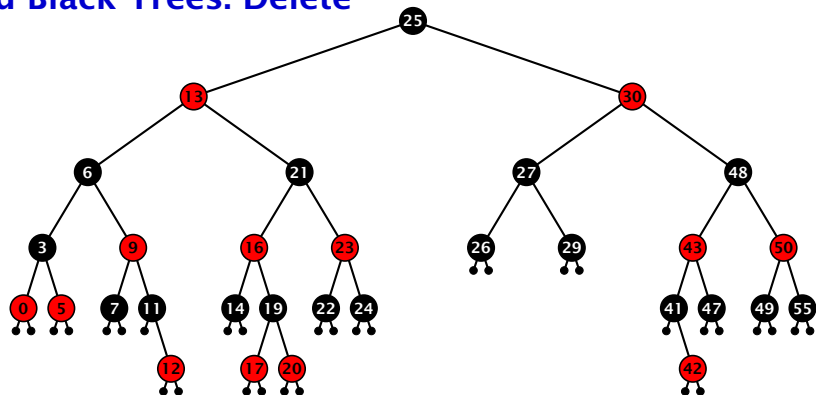
First do a standard delete.

If the spliced out node x was red everything is fine.

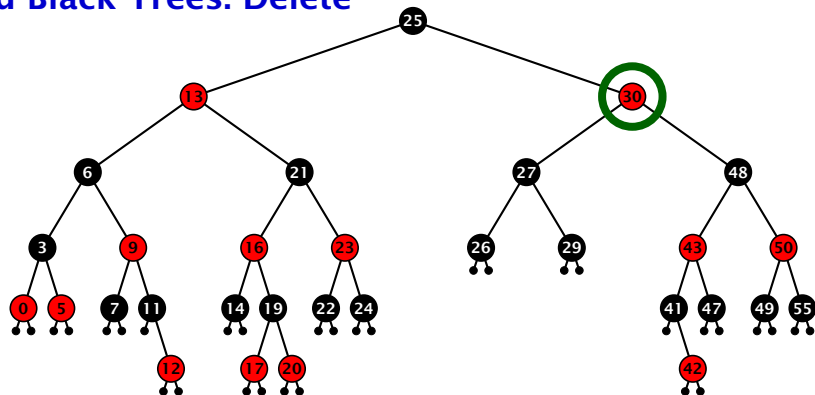
If it was black there may be the following problems.

- ▶ Parent and child of x were red; two adjacent red vertices.
- ▶ If you delete the root, the root may now be red.
- ▶ Every path from an ancestor of x to a descendant leaf of x changes the number of black nodes. Black height property might be violated.

Red Black Trees: Delete



Red Black Trees: Delete

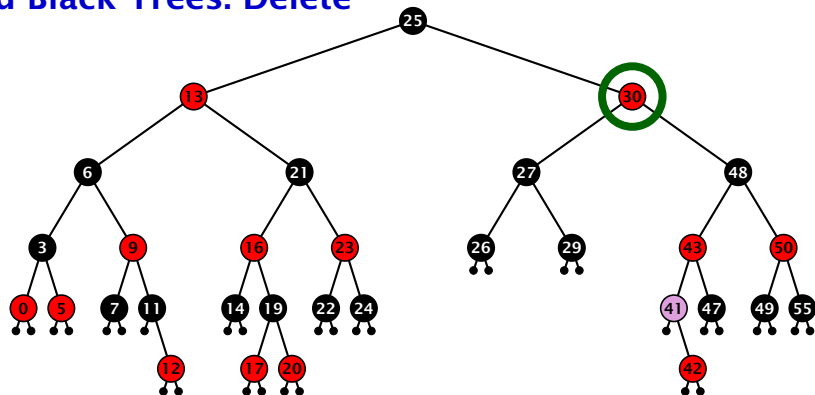


Case 3:

Element has two children

- ▶ do normal delete
- ▶ when replacing content by content of successor, don't change color of node

Red Black Trees: Delete

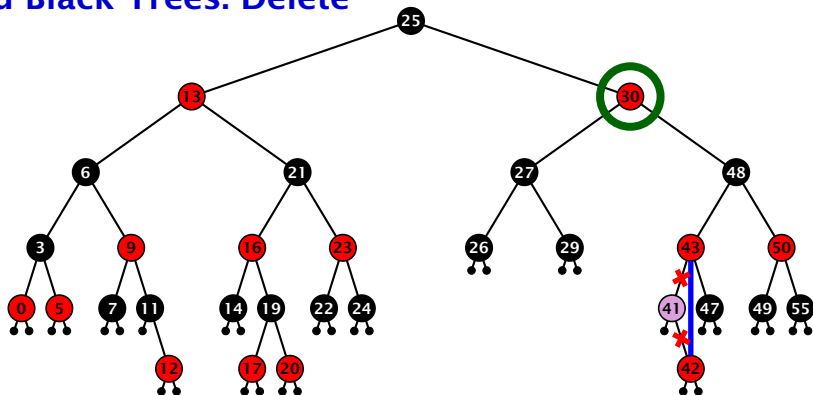


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Red Black Trees: Delete

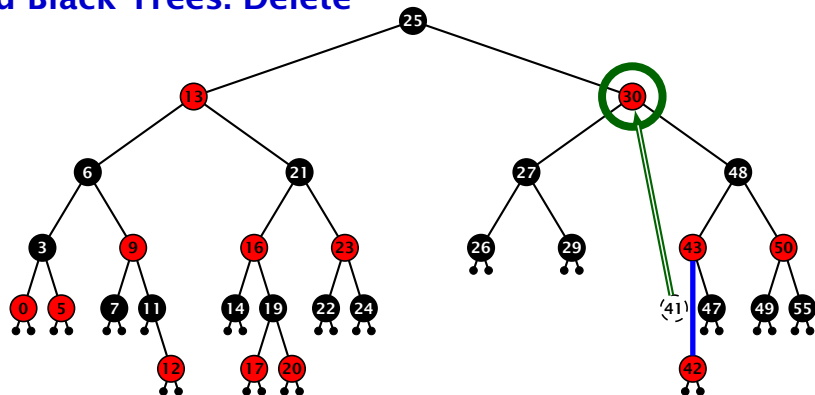


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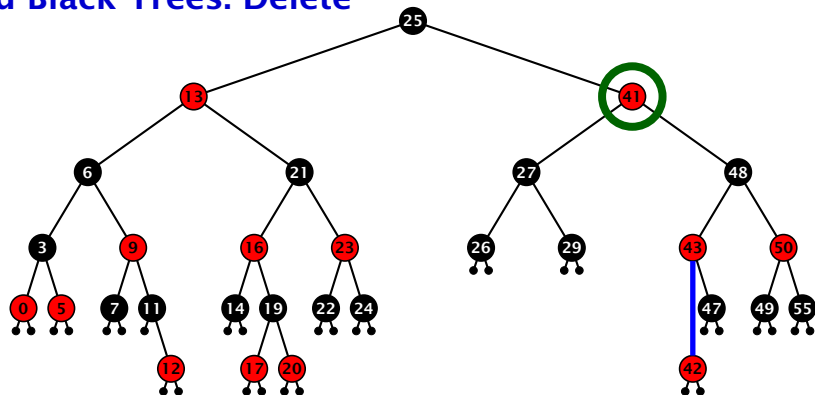


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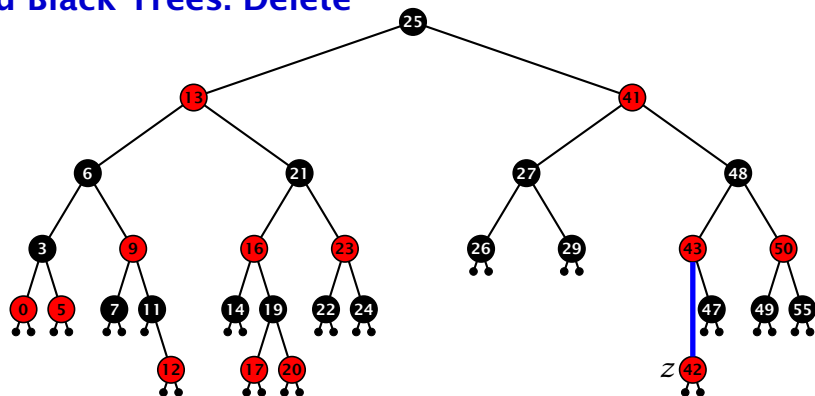


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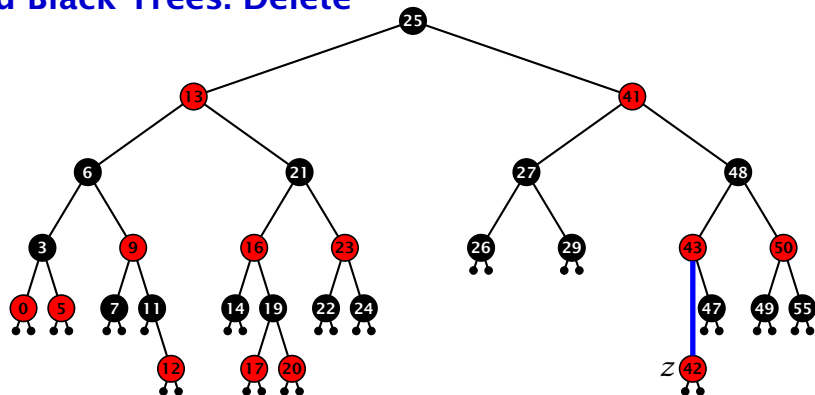
Red Black Trees: Delete



Delete:

- ▶ deleting black node messes up black-height property
- ▶ if z is red, we can simply color it black and everything is fine
- ▶ the problem is if z is black (e.g. a dummy-leaf); we call a fix-up procedure to fix the problem.

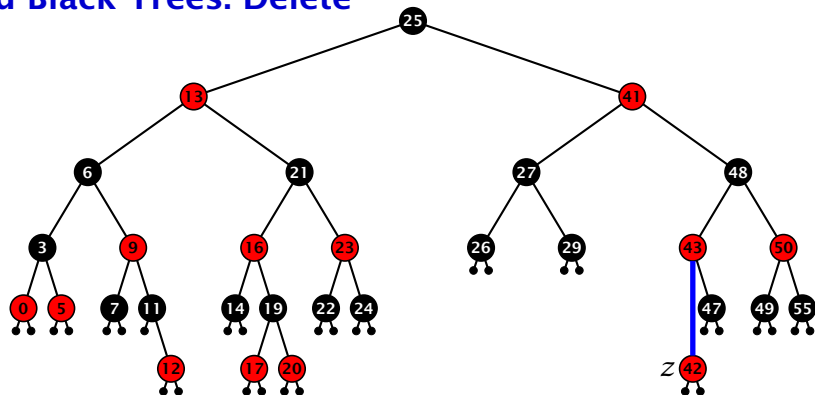
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Red Black Trees: Delete

Invariant of the fix-up algorithm

- ▶ the node z is black
- ▶ if we “assign” a fake black unit to the edge from z to its parent then the black-height property is fulfilled

Goal: make rotations in such a way that you at some point can remove the fake black unit from the edge.

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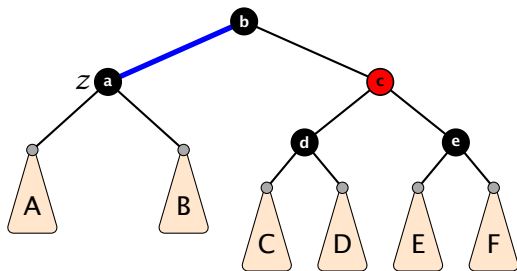
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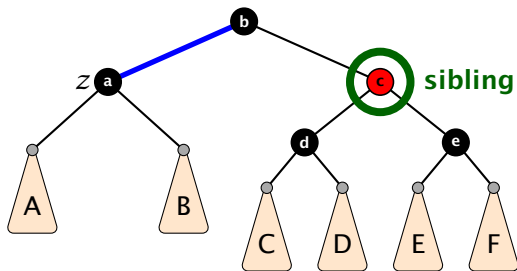
Case 1: Sibling of z is red



1. left-rotate around parent of z
2. recolor nodes b and c
3. the new sibling is black
(and parent of z is red)
4. Case 2 (special),
or Case 3, or Case 4



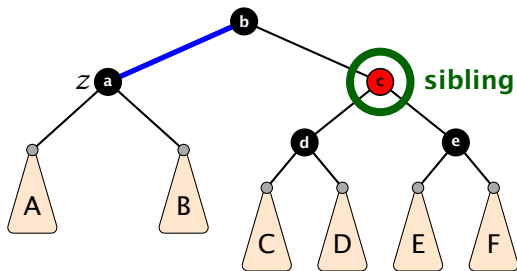
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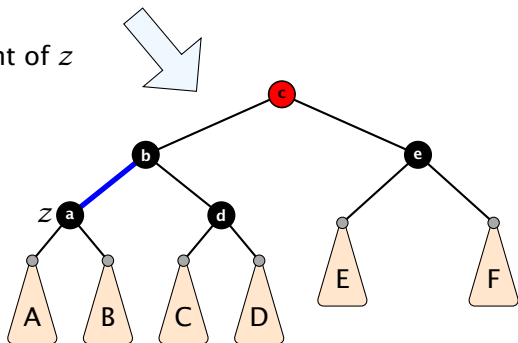
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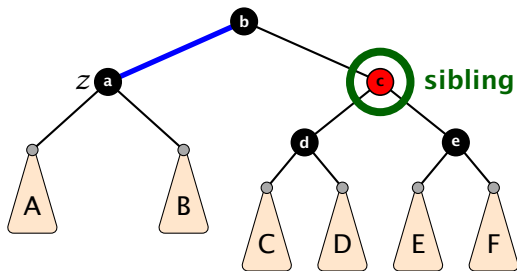
Case 1: Sibling of z is red



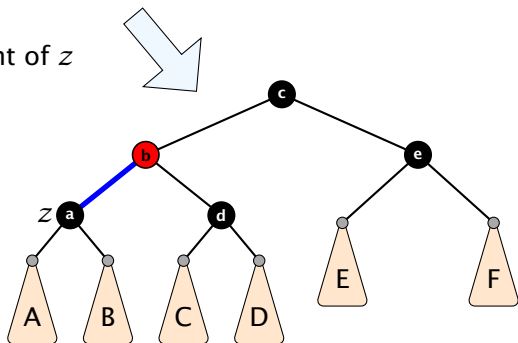
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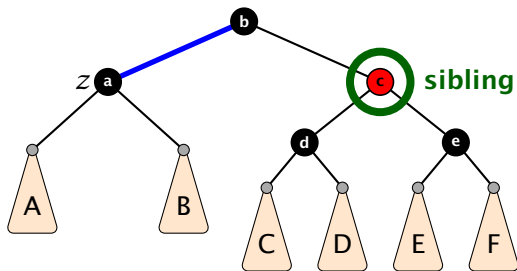
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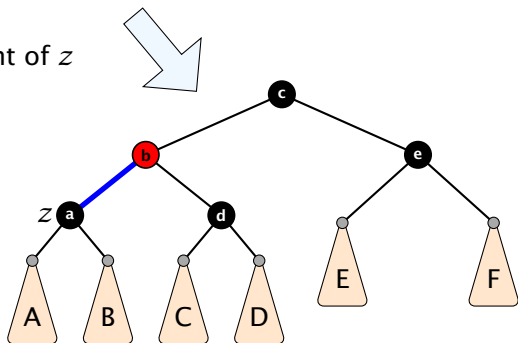
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4. Case 2 (special), or Case 3, or Case 4



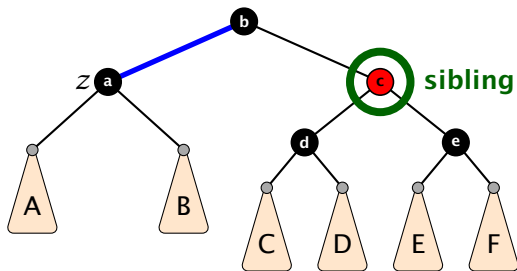
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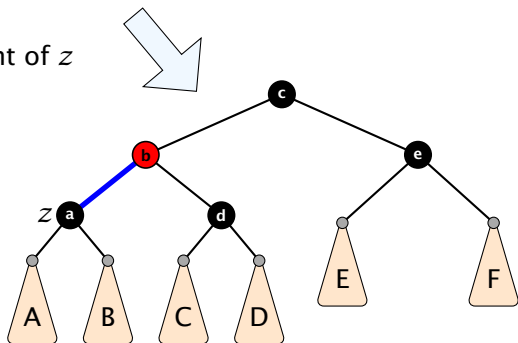
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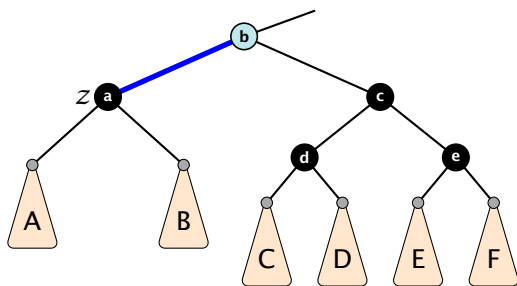
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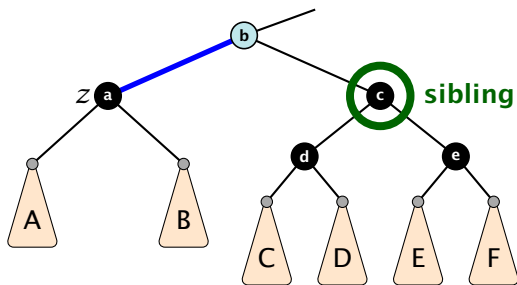
Case 2: Sibling is black with two black children



1. re-color node c
2. move fake black unit upwards
3. move z upwards
4. we made progress
5. if b is red we color it black and are done



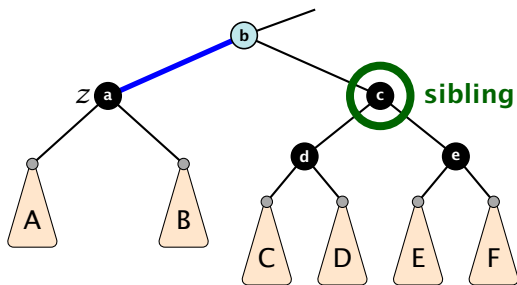
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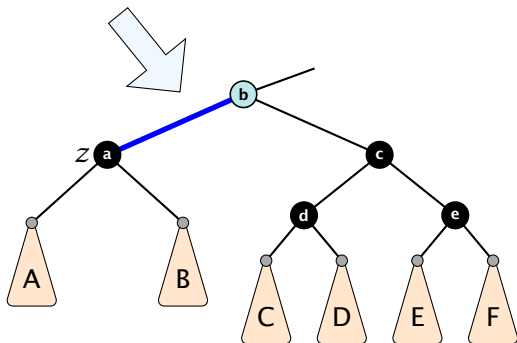
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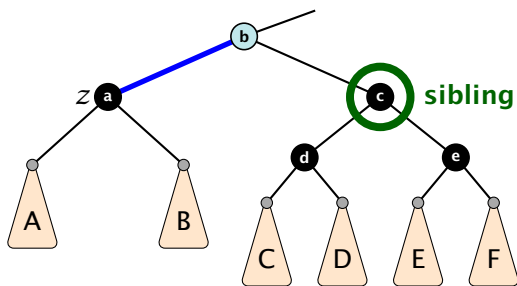
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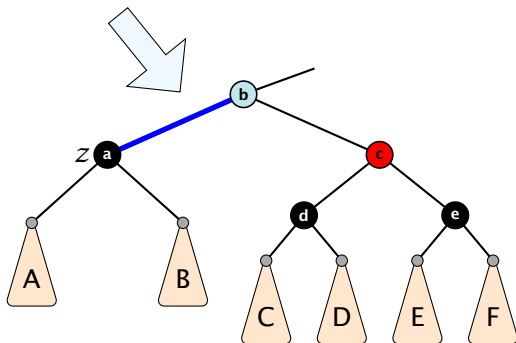
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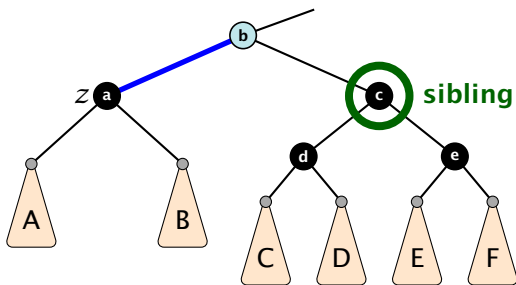
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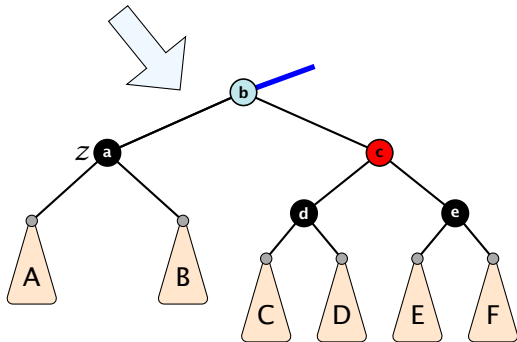
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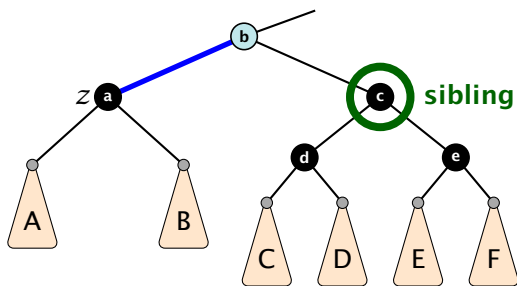
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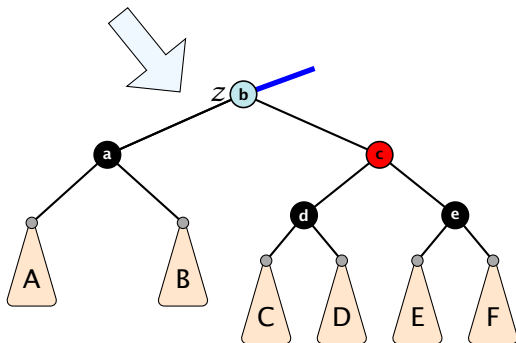
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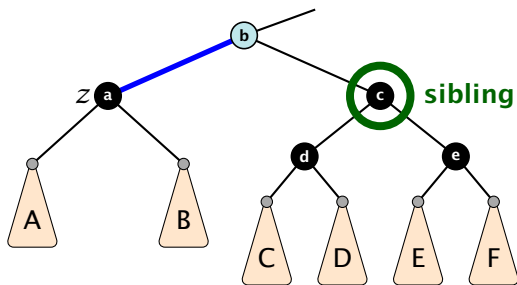
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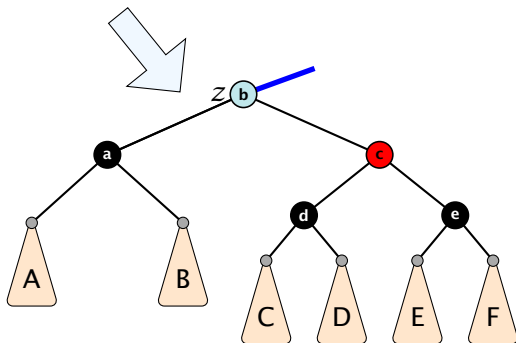
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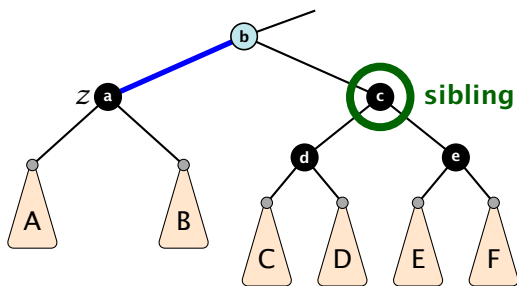
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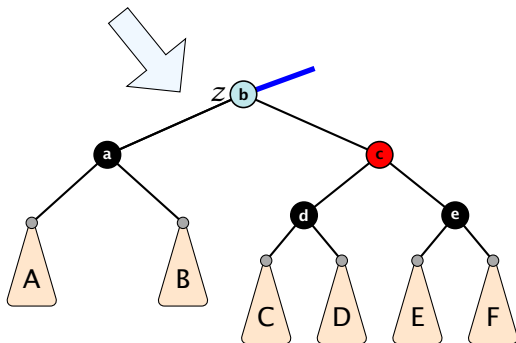
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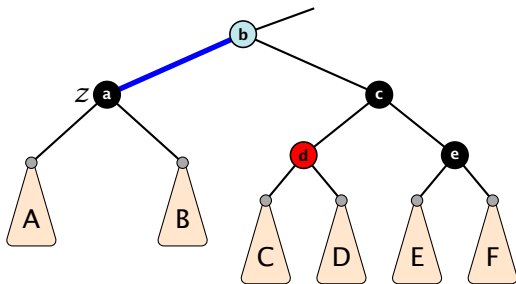


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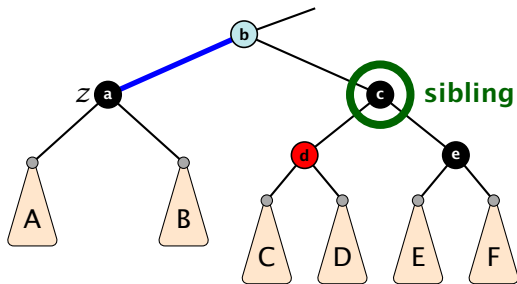
Case 3: Sibling black with one black child to the right

1. do a right-rotation at sibling
2. recolor c and d
3. new sibling is black with red right child (Case 4)



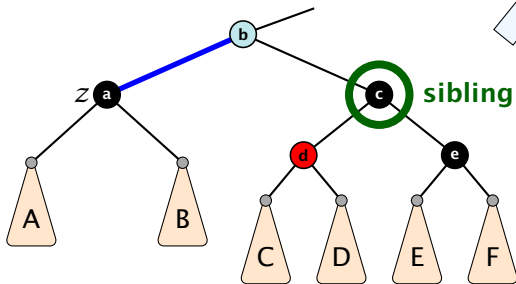
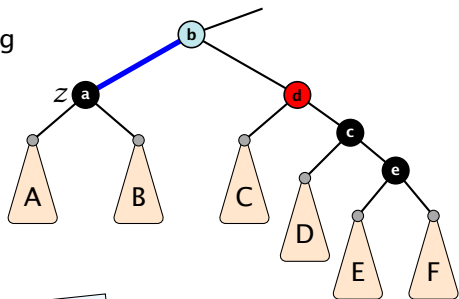
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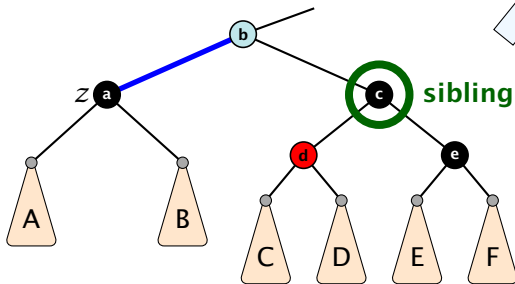
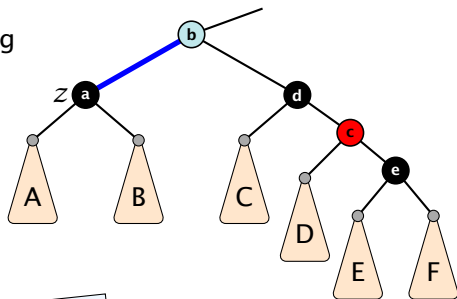
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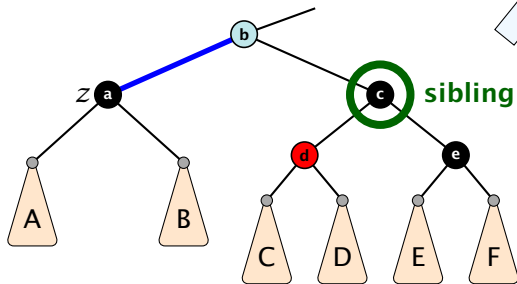
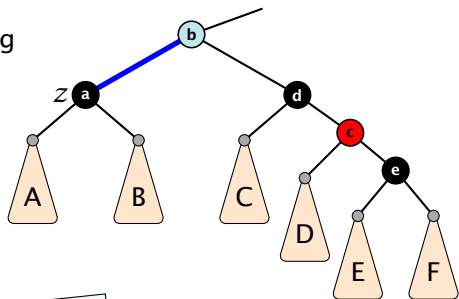
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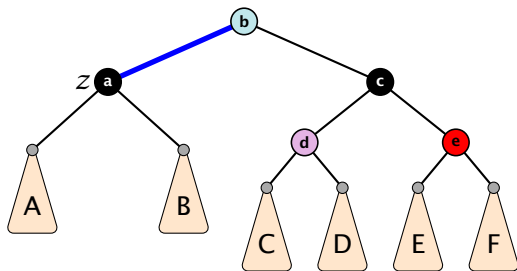


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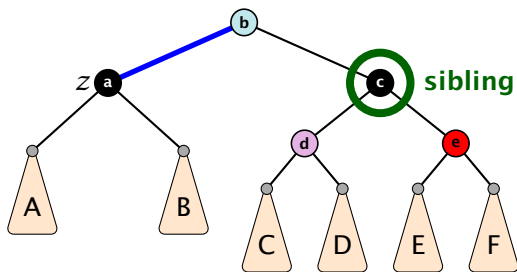
Case 4: Sibling is black with red right child



1. left-rotate around b
2. recolor nodes b , c , and e
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4. you have a valid red black tree



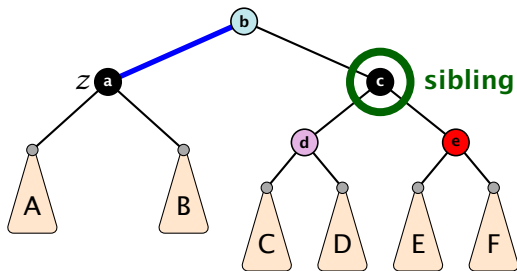
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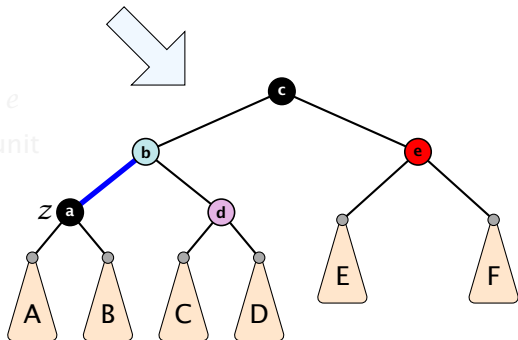
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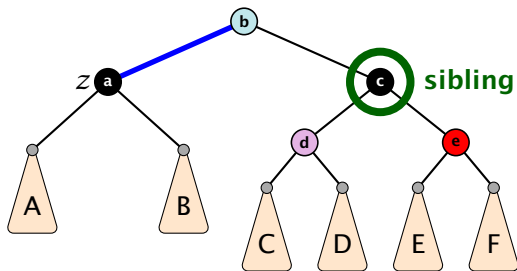
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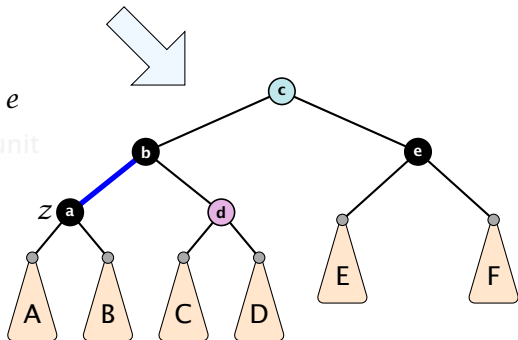
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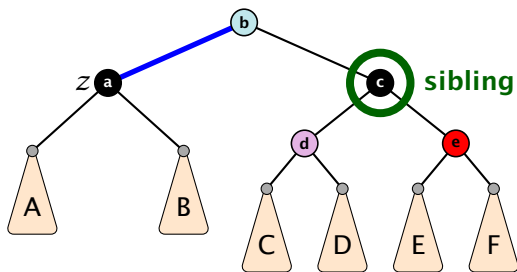
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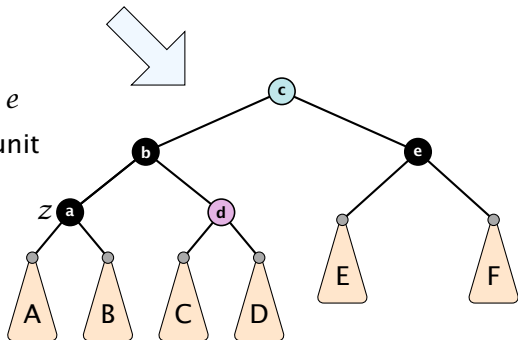
1. left-rotate around *b*
2. recolor nodes *b*, *c*, and *e*
3. remove the fake black unit
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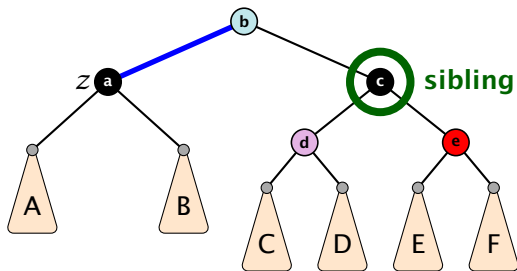
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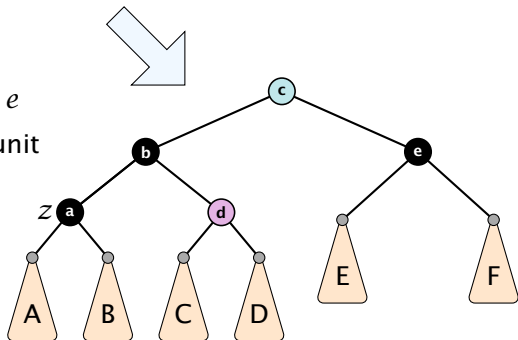
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Case 4: Sibling is black with red right child



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2. recolor nodes b , c , and e
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4. you have a valid red black tree



Running time:

- ▶ only Case 2 can repeat; but only h many steps, where h is the height of the tree
- ▶ Case 1 → Case 2 (special) → red black tree
- ▶ Case 1 → Case 3 → Case 4 → red black tree
- ▶ Case 1 → Case 4 → red black tree
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Performing Case 2 $O(\log n)$ times and every other step at most once, we get a red black tree. Hence, $O(\log n)$ re-colourings and at most 3 rotations.

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7.3 AVL-Trees

Definition 15

AVL-trees are binary search trees that fulfill the following balance condition. For every node v

$$|\text{height}(\text{left sub-tree}(v)) - \text{height}(\text{right sub-tree}(v))| \leq 1 .$$

Lemma 16

An AVL-tree of height h contains at least $F_{h+2} - 1$ and at most $2^h - 1$ internal nodes, where F_n is the n -th Fibonacci number ($F_0 = 0, F_1 = 1$), and the height is the maximal number of edges from the root to an (empty) dummy leaf.

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Proof.

The upper bound is clear, as a binary tree of height h can only contain

$$\sum_{j=0}^{h-1} 2^j = 2^h - 1$$

internal nodes.

Proof (cont.)

Induction (base cases):

1. an AVL-tree of height $h = 1$ contains at least one internal node, $1 \geq F_3 - 1 = 2 - 1 = 1$.
2. an AVL tree of height $h = 2$ contains at least two internal nodes, $2 \geq F_4 - 1 = 3 - 1 = 2$



Proof (cont.)

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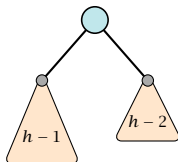


Induction step:

An AVL-tree of height $h \geq 2$ of minimal size has a root with sub-trees of height $h - 1$ and $h - 2$, respectively. Both, sub-trees have minimal node number.

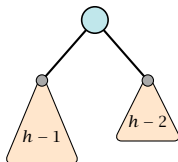
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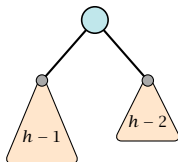


Let

$$f_h := 1 + \text{minimal size of AVL-tree of height } h .$$

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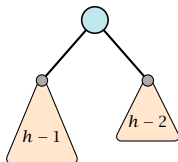
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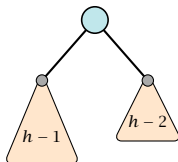
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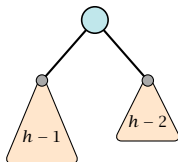
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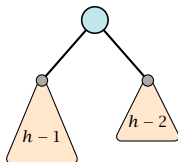
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$$f_h = f_{h-1} + f_{h-2} \qquad = F_{h+2}$$

7.3 AVL-Trees

Since

$$F(k) \approx \frac{1}{\sqrt{5}} \left(\frac{1 + \sqrt{5}}{2} \right)^k ,$$

an AVL-tree with n internal nodes has height $\Theta(\log n)$.

7.3 AVL-Trees

We need to maintain the balance condition through rotations.

For this we store in every internal tree-node v the balance of the node. Let v denote a tree node with left child c_ℓ and right child c_r .

$$\text{balance}[v] := \text{height}(T_{c_\ell}) - \text{height}(T_{c_r}) ,$$

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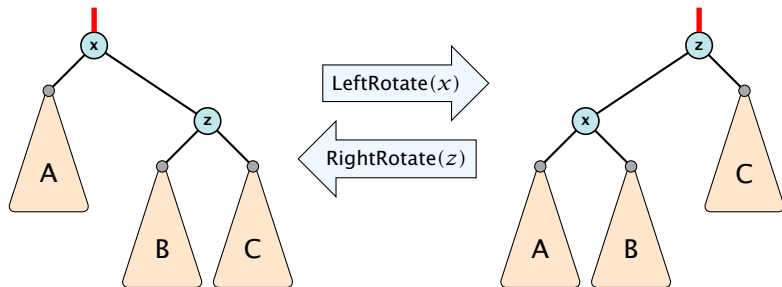
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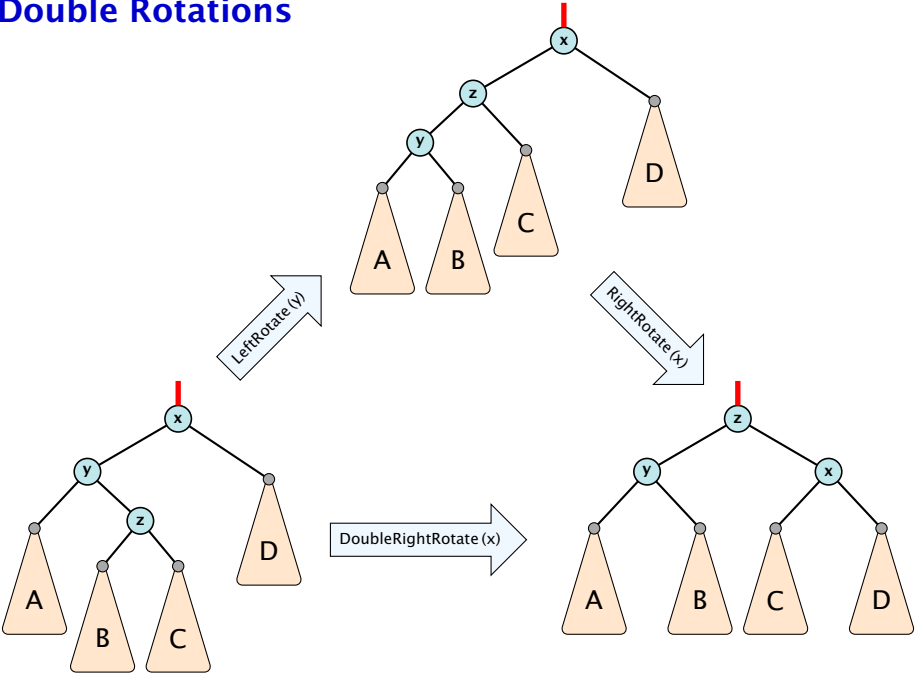
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Rotations

The properties will be maintained through rotations:



Double Rotations



AVL-trees: Insert

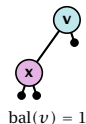
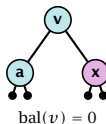
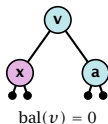
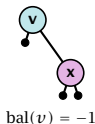
- ▶ Insert like in a binary search tree.

AVL-trees: Insert

- ▶ Insert like in a binary search tree.
- ▶ Let v denote the parent of the newly inserted node x .

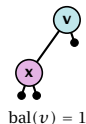
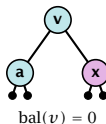
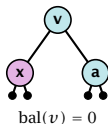
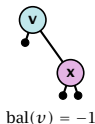
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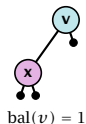
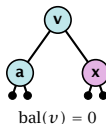
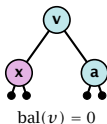
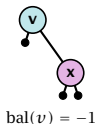
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- ▶ If $\text{bal}[v] \neq 0$, T_v has changed height; the balance-constraint may be violated at ancestors of v .
- ▶ Call $\text{fix-up}(\text{parent}[v])$ to restore the balance-condition.

Invariant at the beginning $\text{fix-up}(v)$:

1. The balance constraints holds at all descendants of v .
2. A node has been inserted into T_c , where c is either the right or left child of v .
3. T_c has increased its height by one (otw. we would already have aborted the fix-up procedure).
4. The balance at the node c fulfills $\text{balance}[c] \in \{-1, 1\}$. This holds because if the balance of c is 0, then T_c did not change its height, and the whole procedure will have been aborted in the previous step.

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AVL-trees: Insert

Algorithm 11 AVL-fix-up-insert(v)

- 1: **if** $\text{balance}[v] \in \{-2, 2\}$ **then** DoRotationInsert(v);
- 2: **if** $\text{balance}[v] \in \{0\}$ **return**;
- 3: AVL-fix-up-insert(parent[v]);

We will show that the above procedure is correct, and that it will do at most one rotation.

AVL-trees: Insert

Algorithm 12 DoRotationInsert(v)

```
1: if balance[ $v$ ] = -2 then
2:     if balance[right[ $v$ ]] = -1 then
3:         LeftRotate( $v$ );
4:     else
5:         DoubleLeftRotate( $v$ );
6: else
7:     if balance[left[ $v$ ]] = 1 then
8:         RightRotate( $v$ );
9:     else
10:        DoubleRightRotate( $v$ );
```

AVL-trees: Insert

It is clear that the invariant for the fix-up routine holds as long as no rotations have been done.

We have to show that after doing one rotation all balance constraints are fulfilled.

We show that after doing a rotation at v :

- ▶ v fulfills balance condition.
- ▶ All children of v still fulfill the balance condition.
- ▶ The height of T_v is the same as before the insert-operation took place.

We only look at the case where the insert happened into the right sub-tree of v . The other case is symmetric.

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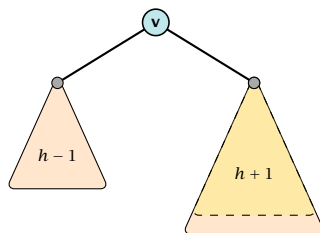
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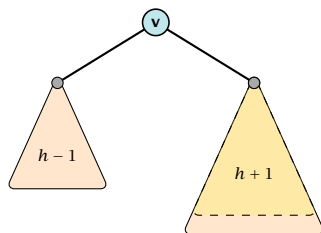


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Before the insertion the height of T_v was $h + 1$.

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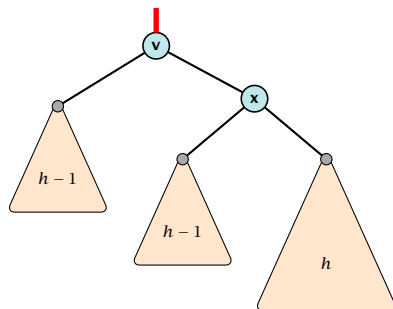
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Case 1: $\text{balance}[\text{right}[v]] = -1$

We do a left rotation at v

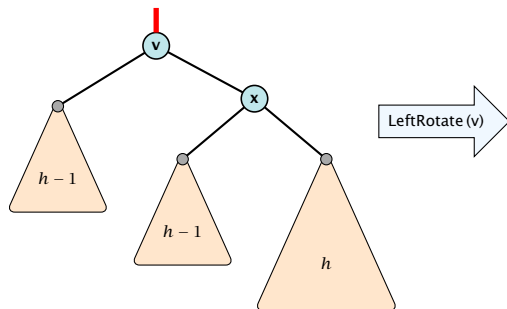
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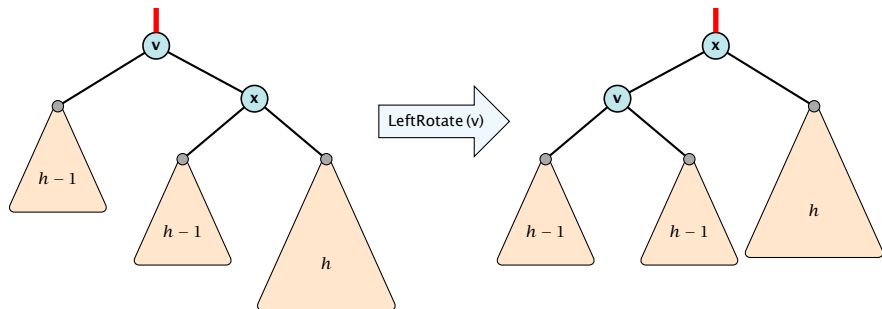
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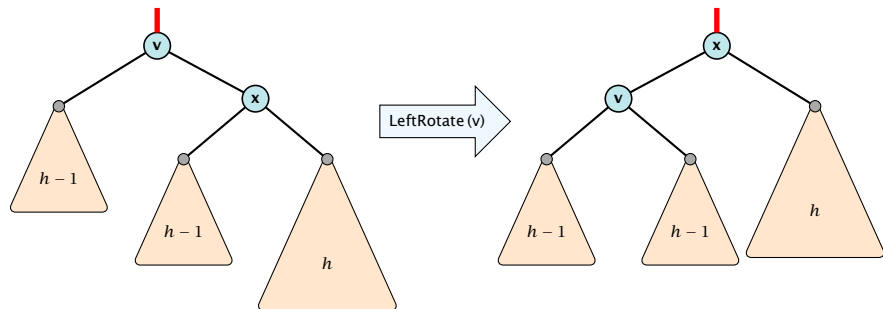
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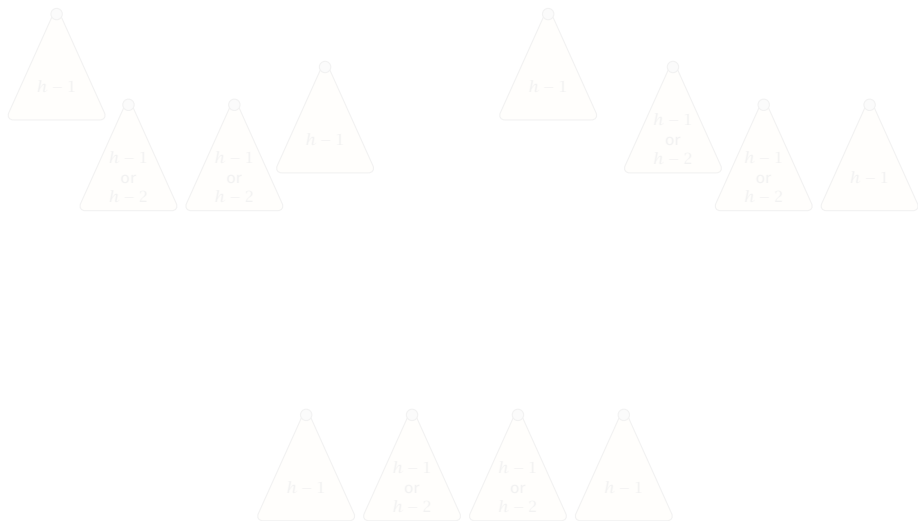
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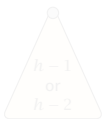
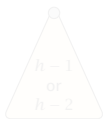
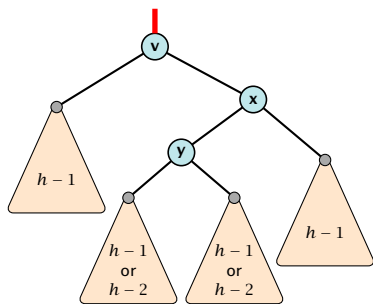


Now, T_v has height $h + 1$ as before the insertion. Hence, we do not need to continue.

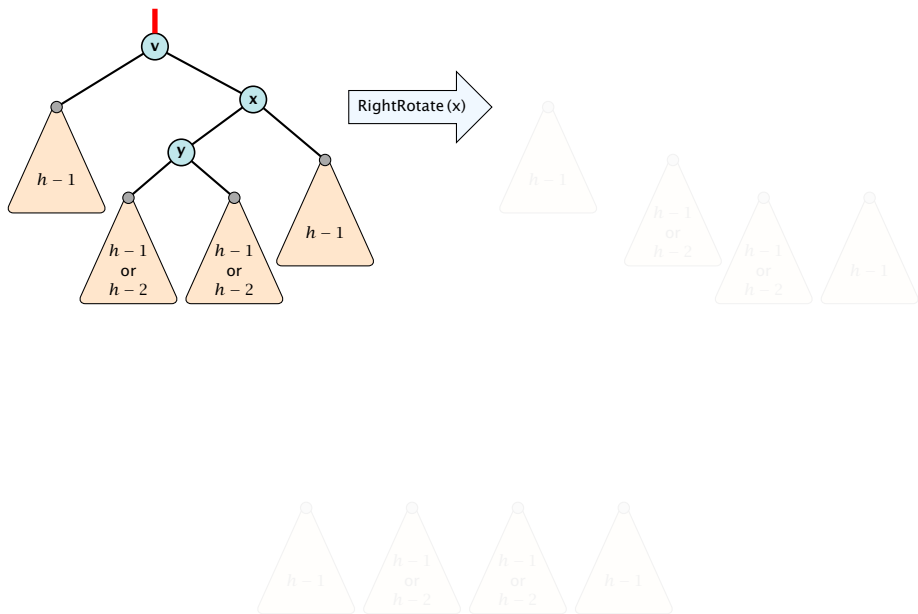
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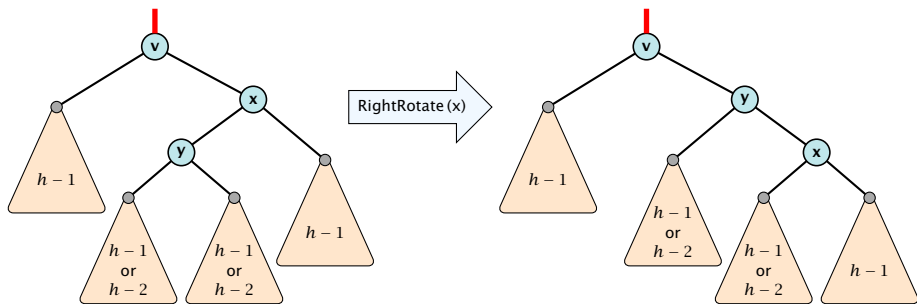
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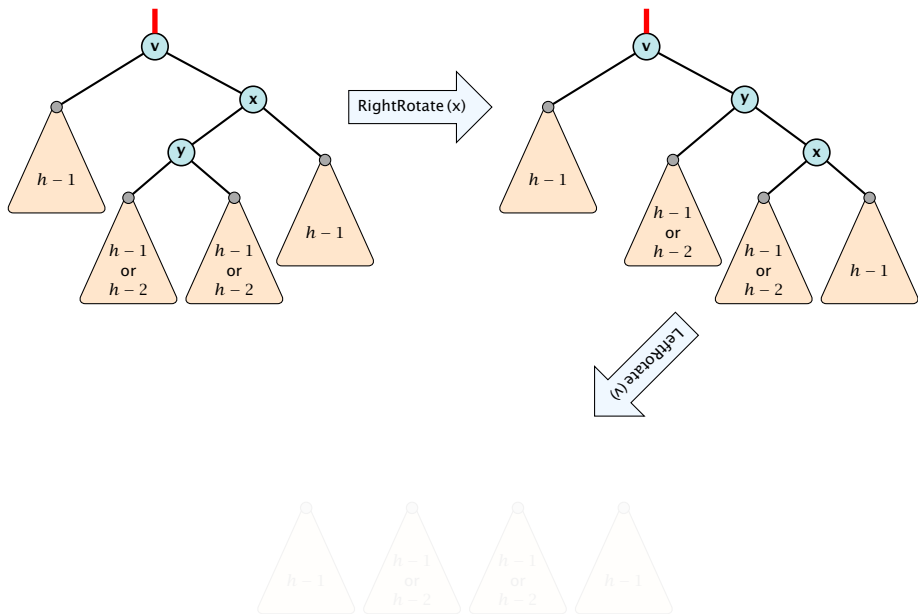
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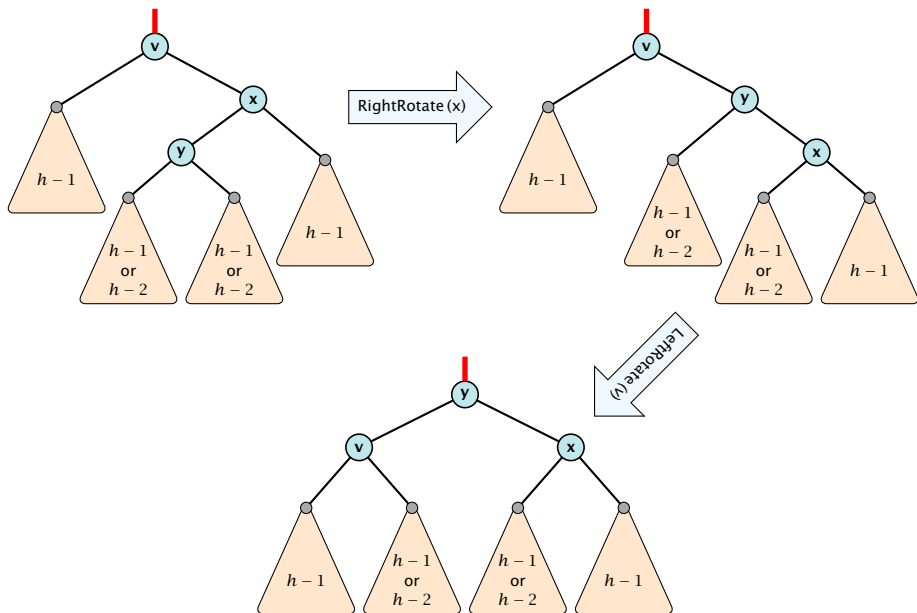
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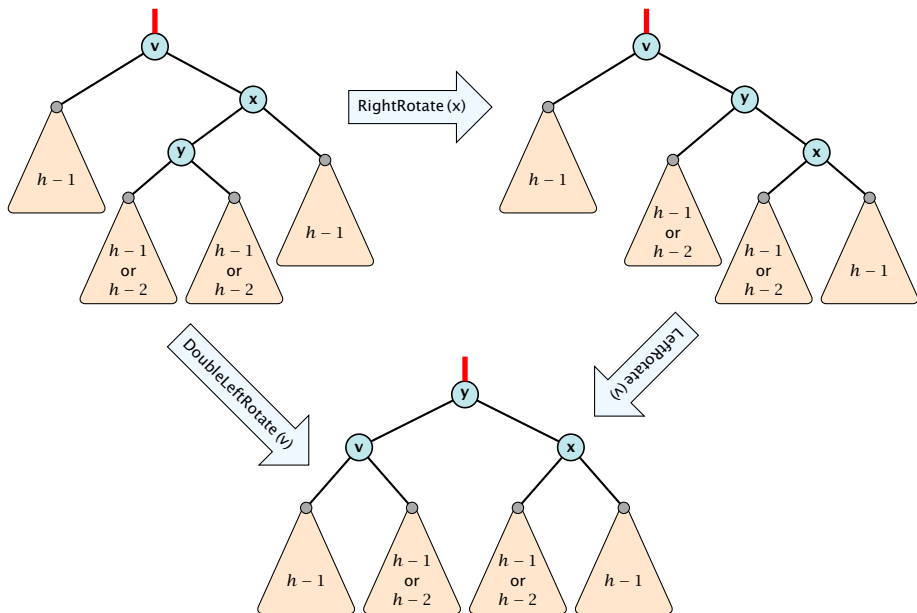
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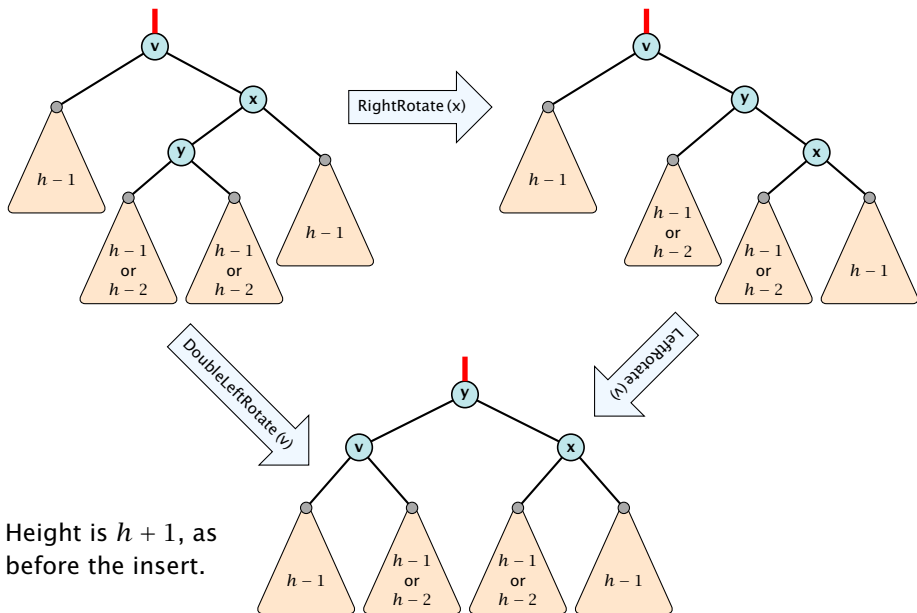
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AVL-trees: Delete

- ▶ Delete like in a binary search tree.
- ▶ Let v denote the parent of the node that has been spliced out.
- ▶ The balance-constraint may be violated at v , or at ancestors of v , as a sub-tree of a child of v has reduced its height.
- ▶ Initially, the node c —the new root in the sub-tree that has changed—is either a dummy leaf or a node with two dummy leaves as children.



Case 1



Case 2

In both cases $\text{bal}[c] = 0$.

- ▶ Call $\text{fix-up}(v)$ to restore the balance-condition.

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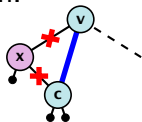
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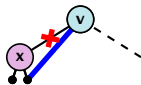
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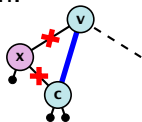
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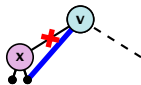
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Algorithm 13 AVL-fix-up-delete(v)

- 1: **if** $\text{balance}[v] \in \{-2, 2\}$ **then** DoRotationDelete(v);
- 2: **if** $\text{balance}[v] \in \{-1, 1\}$ **return**;
- 3: AVL-fix-up-delete($\text{parent}[v]$);

We will show that the above procedure is correct. However, for the case of a delete there may be a logarithmic number of rotations.

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AVL-trees: Delete

Algorithm 14 DoRotationDelete(v)

```
1: if balance[ $v$ ] = -2 then
2:     if balance[right[ $v$ ]] = -1 then
3:         LeftRotate( $v$ );
4:     else
5:         DoubleLeftRotate( $v$ );
6: else
7:     if balance[left[ $v$ ]] = {0, 1} then
8:         RightRotate( $v$ );
9:     else
10:        DoubleRightRotate( $v$ );
```

AVL-trees: Delete

It is clear that the invariant for the fix-up routine holds as long as no rotations have been done.

We show that after doing a rotation at v :

- ▶ v fulfills balance condition.
- ▶ All children of v still fulfill the balance condition.
- ▶ If now $\text{balance}[v] \in \{-1, 1\}$ we can stop as the height of T_v is the same as before the deletion.

We only look at the case where the deleted node was in the right sub-tree of v . The other case is symmetric.

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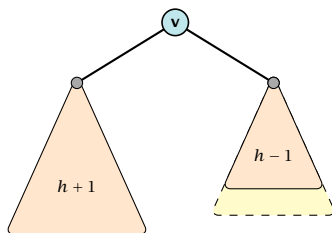
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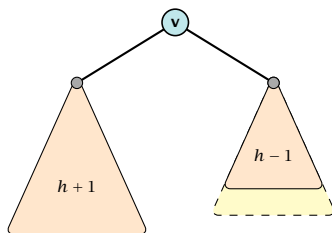


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Before the insertion the height of T_v was $h + 2$.

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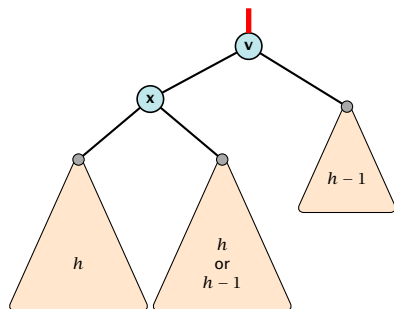
Case 1: $\text{balance}[\text{left}[v]] \in \{0, 1\}$



If the middle subtree has height h the whole tree has height $h + 2$ as before the deletion. The iteration stops as the balance at the root is non-zero.

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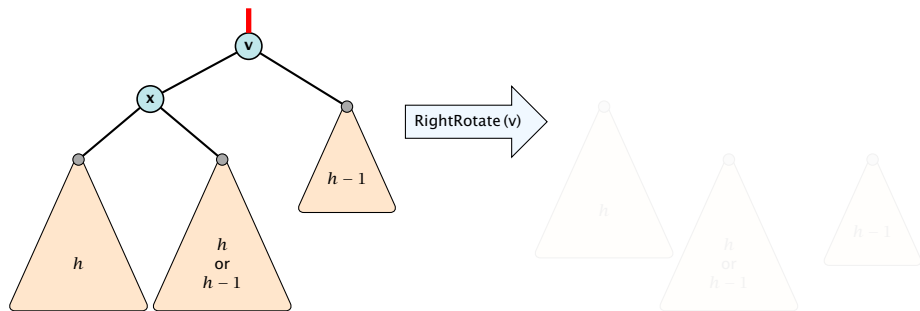
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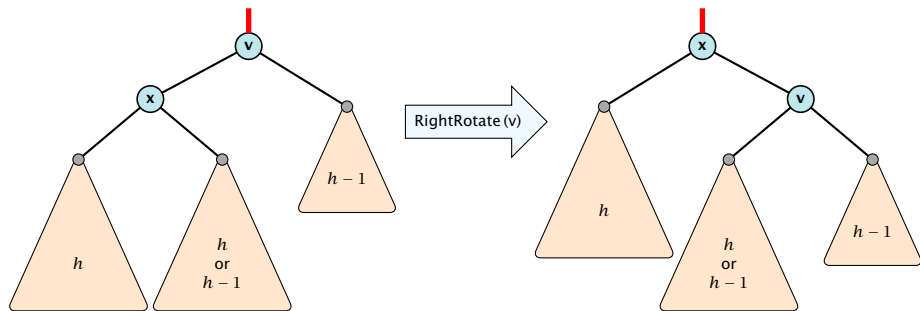
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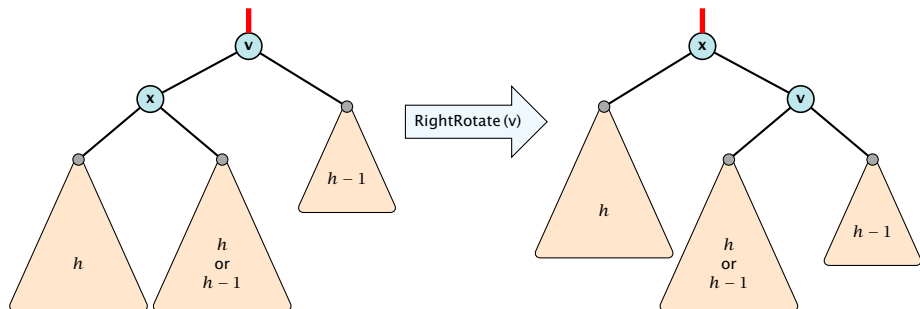
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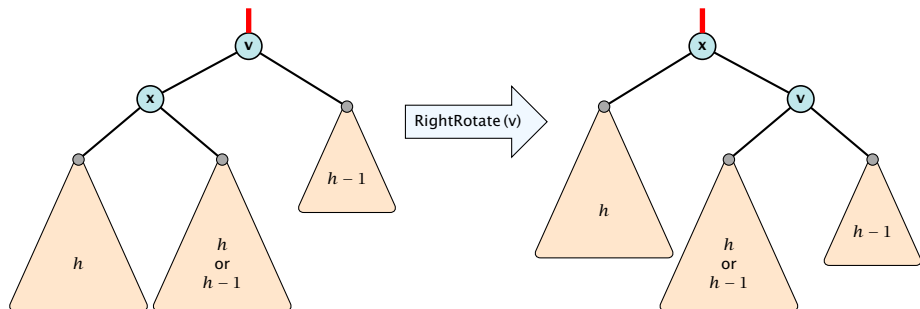
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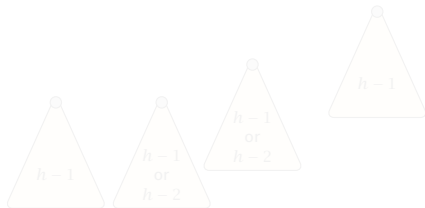
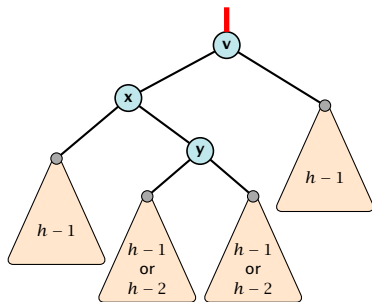
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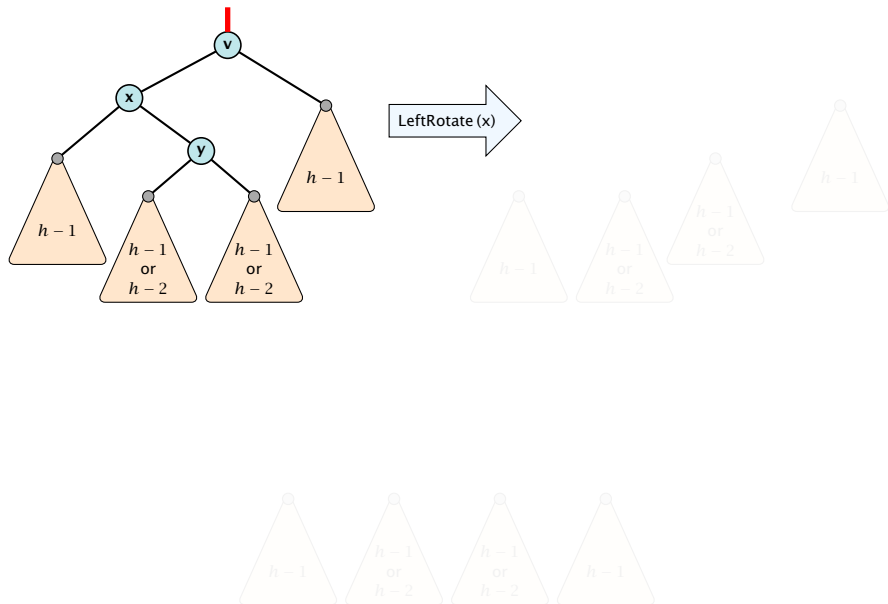
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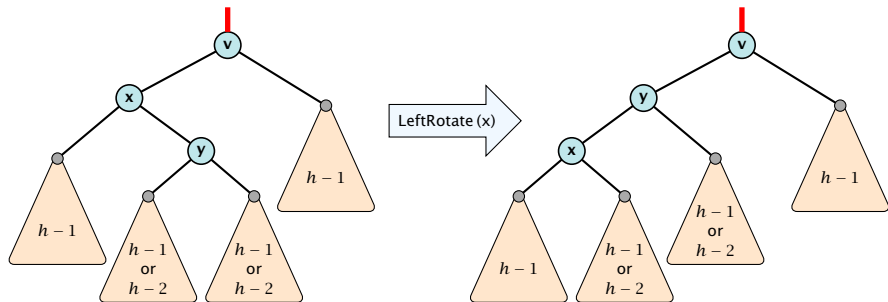
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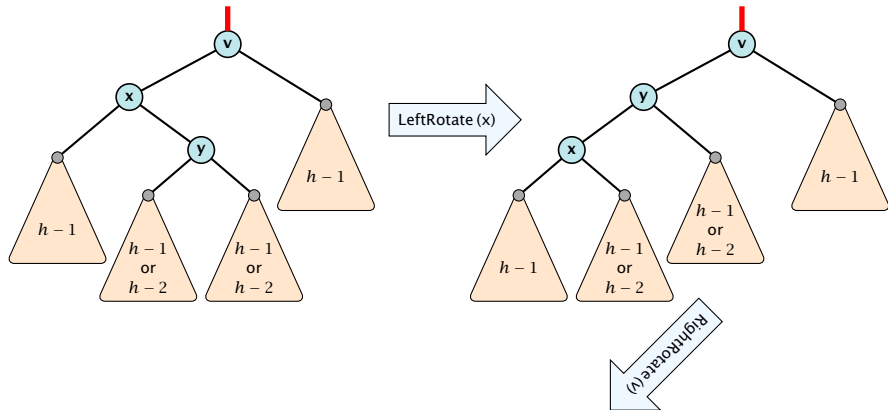
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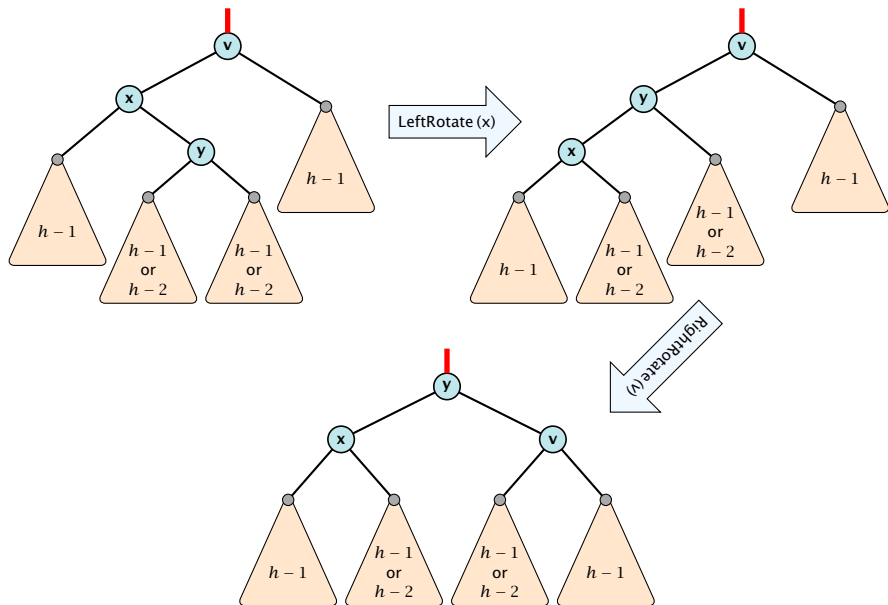
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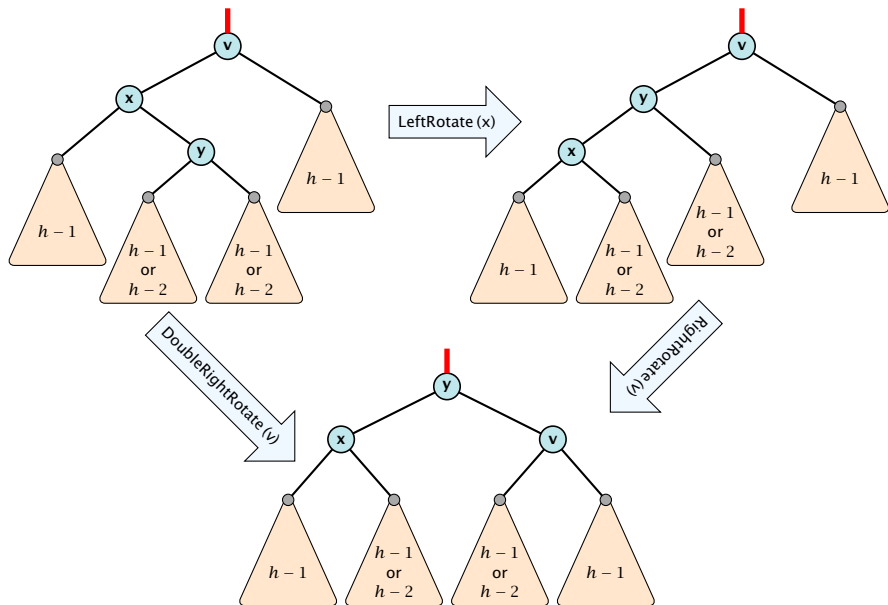
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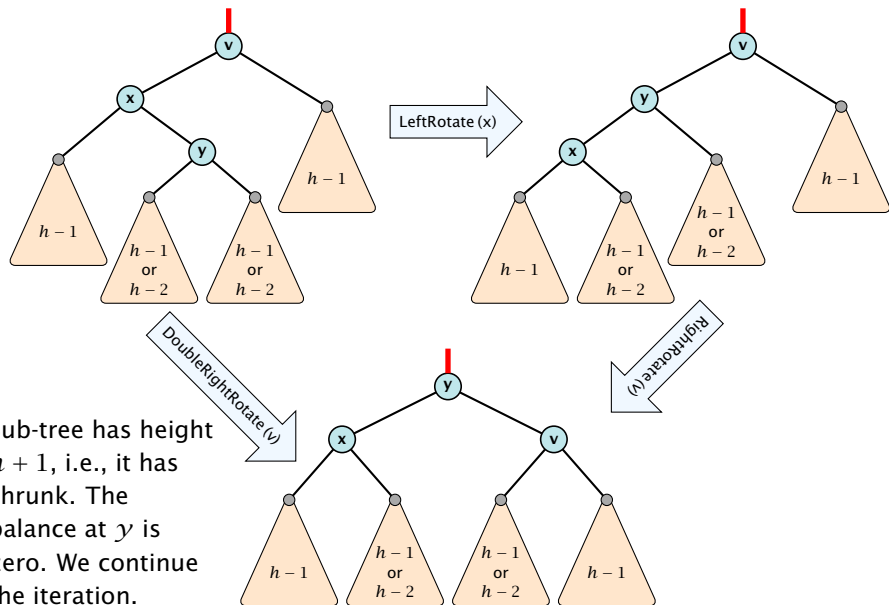
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7.4 (a, b)-trees

Definition 17

For $b \geq 2a - 1$ an (a, b) -tree is a search tree with the following properties

1. all leaves have the same distance to the root
2. every internal non-root vertex v has at least a and at most b children
3. the root has degree at least 2 if the tree is non-empty
4. the internal vertices do not contain data, but only keys (external search tree)
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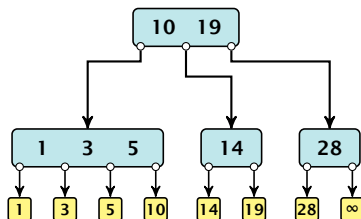
Each internal node v with $d(v)$ children stores $d - 1$ keys k_1, \dots, k_{d-1} . The i -th subtree of v fulfills

$$k_{i-1} < \text{key in } i\text{-th sub-tree} \leq k_i ,$$

where we use $k_0 = -\infty$ and $k_d = \infty$.

7.4 (a, b)-trees

Example 18



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Variants

- ▶ The dummy leaf element may not exist; this only makes implementation more convenient.
- ▶ Variants in which $b = 2a$ are commonly referred to as B -trees.
- ▶ A B -tree usually refers to the variant in which keys and data are stored at internal nodes.
- ▶ A B^+ tree stores the data only at leaf nodes as in our definition. Sometimes the leaf nodes are also connected in a linear list data structure to speed up the computation of successors and predecessors.
- ▶ A B^* tree requires that a node is at least $2/3$ -full as only $1/2$ -full (the requirement of a B -tree).

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Let T be an (a, b) -tree for $n > 0$ elements (i.e., $n + 1$ leaf nodes) and height h (number of edges from root to a leaf vertex). Then

1. $2a^{h-1} \leq n + 1 \leq b^h$
2. $\log_b(n + 1) \leq h \leq \log_a\left(\frac{n+1}{2}\right)$

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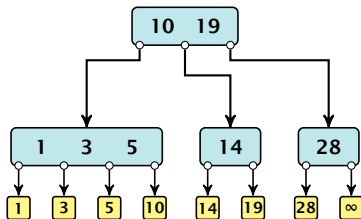
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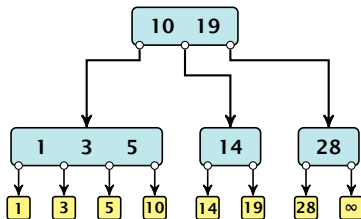


Search



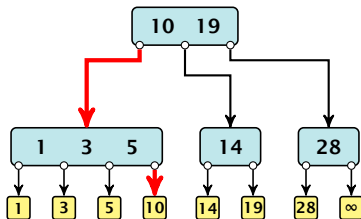
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Search(8)



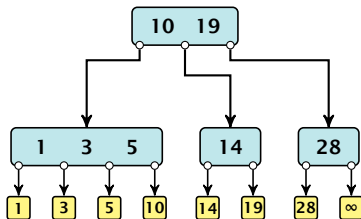
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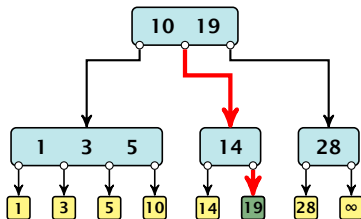
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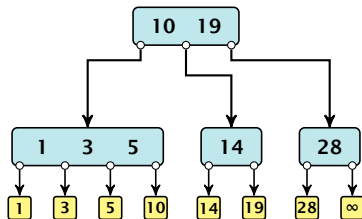


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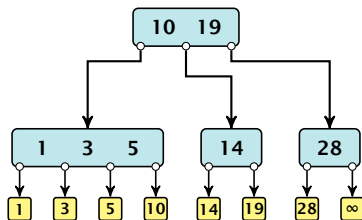


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Time: $\mathcal{O}(b \cdot h) = \mathcal{O}(b \cdot \log n)$, if the individual nodes are organized as linear lists.

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Insert element x :

- ▶ Follow the path as if searching for $\text{key}[x]$.
- ▶ If this search ends in leaf ℓ , insert x **before** this leaf.
- ▶ For this add $\text{key}[x]$ to the key-list of the last internal node v on the path.
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Rebalance(v):

- ▶ Let k_i , $i = 1, \dots, b$ denote the keys stored in v .
- ▶ Let $j := \lfloor \frac{b+1}{2} \rfloor$ be the middle element.
- ▶ Create two nodes v_1 , and v_2 . v_1 gets all keys k_1, \dots, k_{j-1} and v_2 gets keys k_{j+1}, \dots, k_b .
- ▶ Both nodes get at least $\lfloor \frac{b-1}{2} \rfloor$ keys, and have therefore degree at least $\lfloor \frac{b-1}{2} \rfloor + 1 \geq a$ since $b \geq 2a - 1$.
- ▶ They get at most $\lceil \frac{b-1}{2} \rceil$ keys, and have therefore degree at most $\lceil \frac{b-1}{2} \rceil + 1 \leq b$ (since $b \geq 2$).
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Rebalance(v):

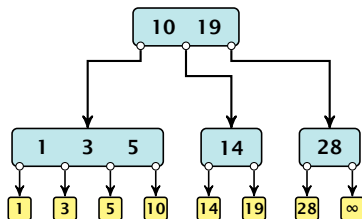
- ▶ Let k_i , $i = 1, \dots, b$ denote the keys stored in v .
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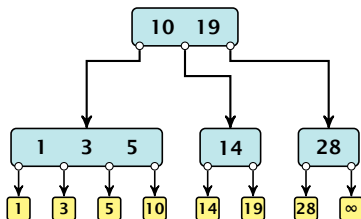
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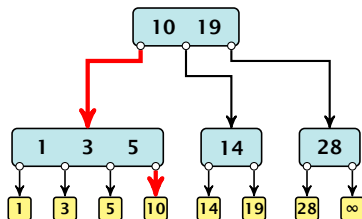
Insert

Insert(8)



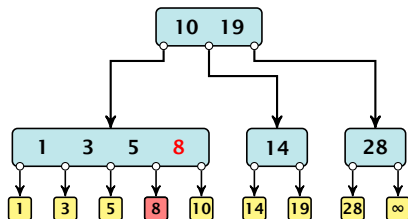
Insert

Insert(8)



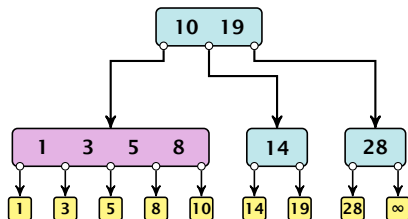
Insert

Insert(8)



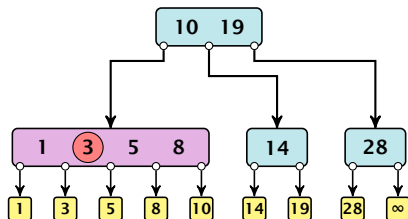
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Insert(8)

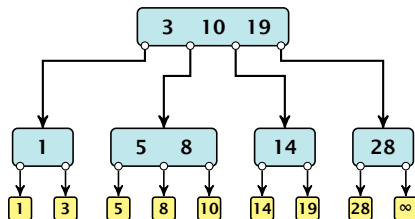


Insert

Insert(8)

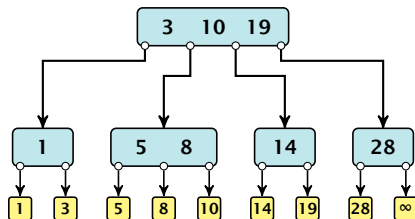


Insert



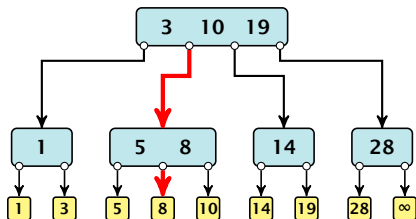
Insert

Insert(6)



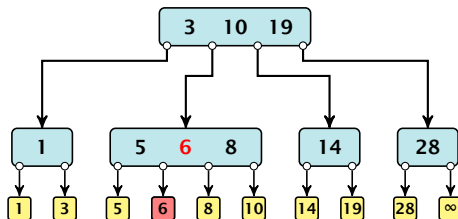
Insert

Insert(6)



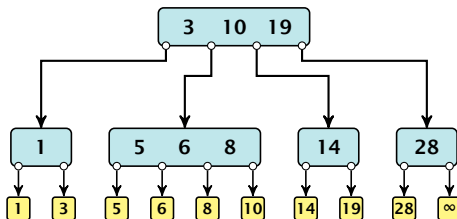
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Insert(6)



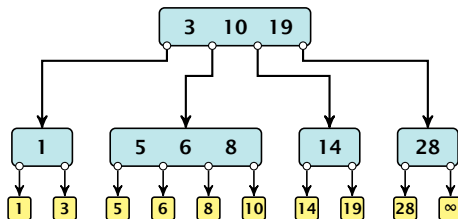
Insert

Insert(6)



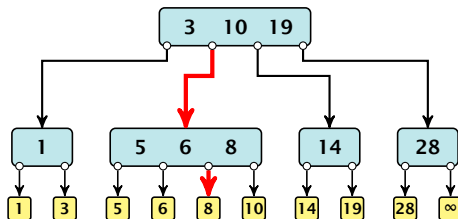
Insert

Insert(7)



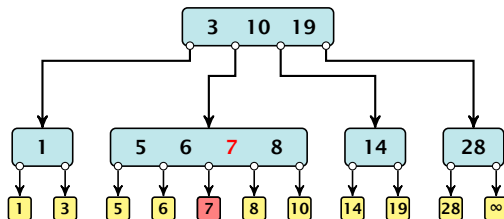
Insert

Insert(7)



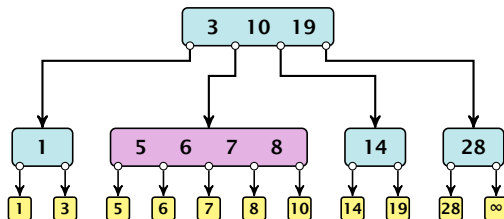
Insert

Insert(7)



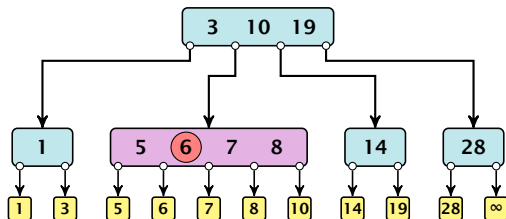
Insert

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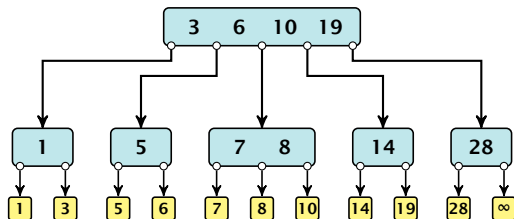
Insert

Insert(7)



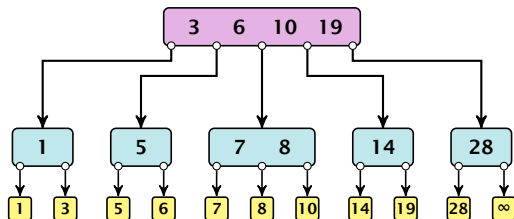
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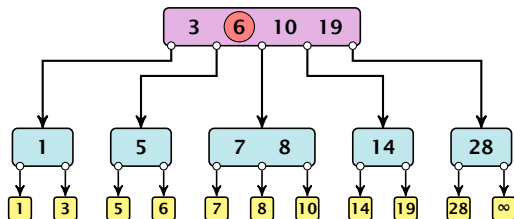
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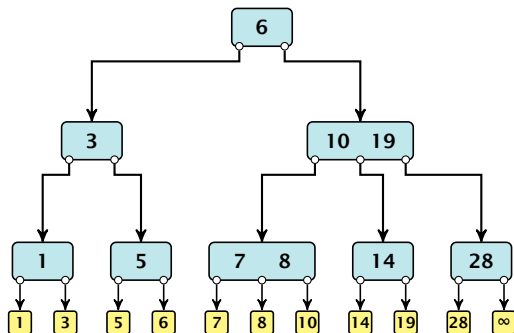
Insert

Insert(7)



Insert

Insert(7)



Delete

Delete element x (pointer to leaf vertex):

- ▶ Let v denote the parent of x . If $\text{key}[x]$ is contained in v , remove the key from v , and delete the leaf vertex.
- ▶ Otherwise delete the key of the predecessor of x from v ; delete the leaf vertex; and replace the occurrence of $\text{key}[x]$ in internal nodes by the predecessor key. (Note that it appears in exactly one internal vertex).
- ▶ If now the number of keys in v is below $a - 1$ perform $\text{Rebalance}'(v)$.

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Rebalance'(v):

- ▶ If there is a neighbour of v that has at least a keys take over the largest (if right neighbor) or smallest (if left neighbour) and the corresponding sub-tree.
- ▶ If not: merge v with one of its neighbours.
- ▶ The merged node contains at most $(a - 2) + (a - 1) + 1$ keys, and has therefore at most $2a - 1 \leq b$ successors.
- ▶ Then rebalance the parent.
- ▶ During this process the root may become empty. In this case the root is deleted and the height of the tree decreases.

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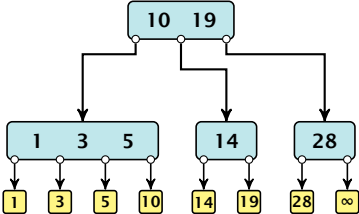
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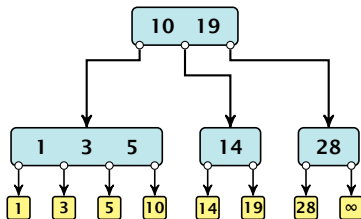
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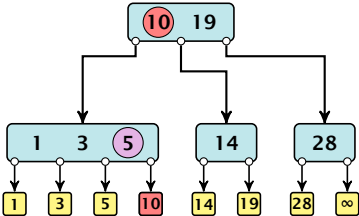
Delete

Delete(10)



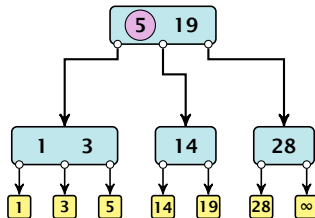
Delete

Delete(10)

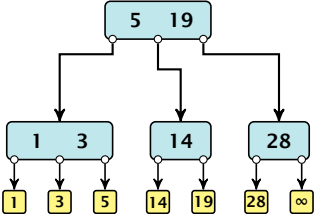


Delete

Delete(10)

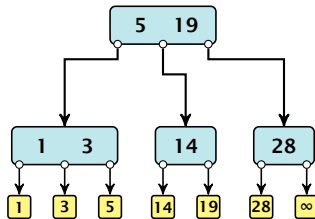


Delete



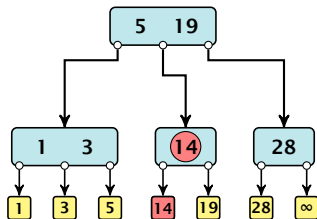
Delete

Delete(14)



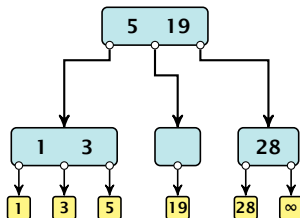
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Delete(14)



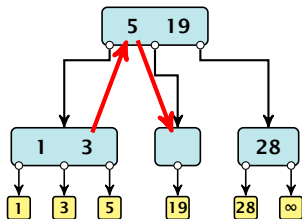
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Delete(14)



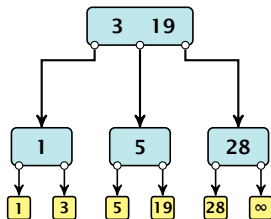
Delete

Delete(14)

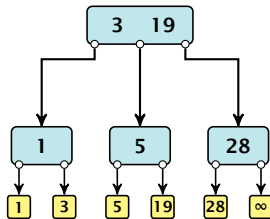


Delete

Delete(14)

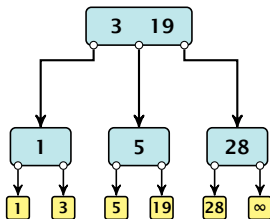


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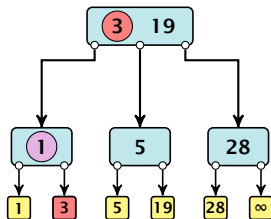
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Delete(3)



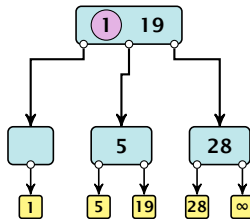
Delete

Delete(3)



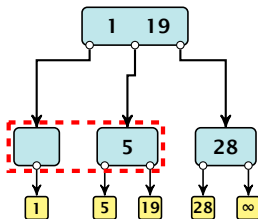
Delete

Delete(3)



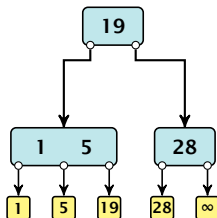
Delete

Delete(3)

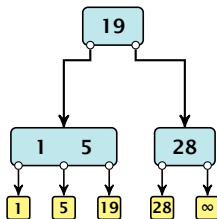


Delete

Delete(3)

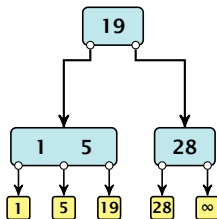


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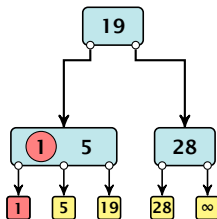
Delete

Delete(1)



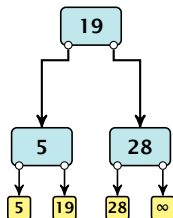
Delete

Delete(1)

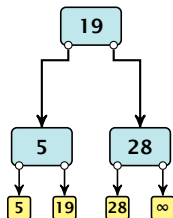


Delete

Delete(1)

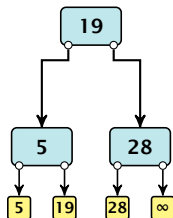


Delete



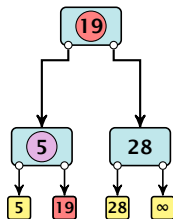
Delete

Delete(19)



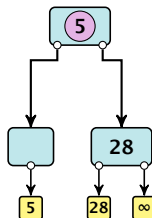
Delete

Delete(19)



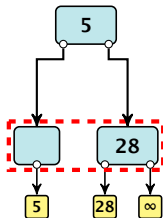
Delete

Delete(19)



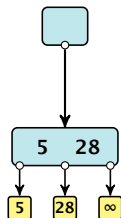
Delete

Delete(19)



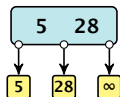
Delete

Delete(19)



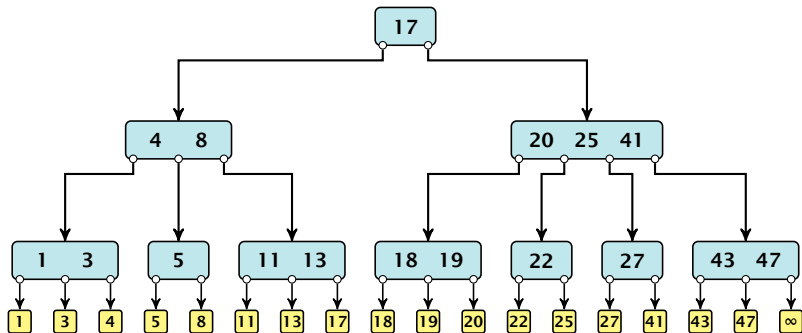
Delete

Delete(19)



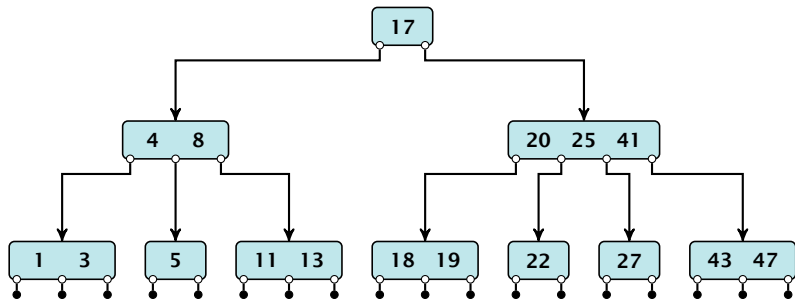
(2, 4)-trees and red black trees

There is a close relation between red-black trees and (2, 4)-trees:



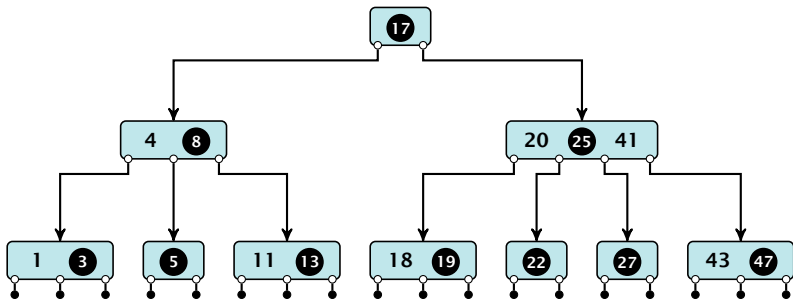
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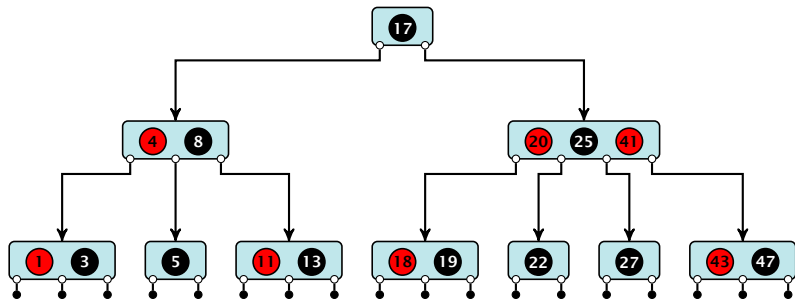
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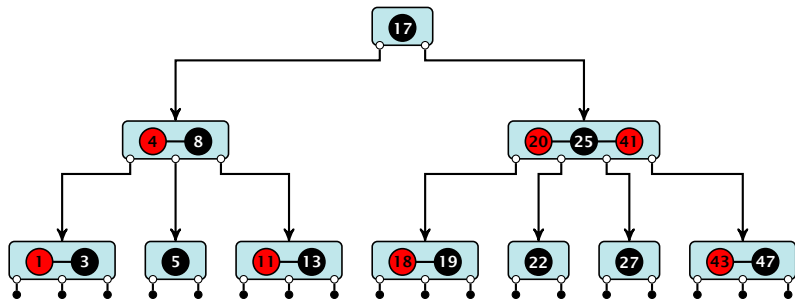
(2, 4)-trees and red black trees

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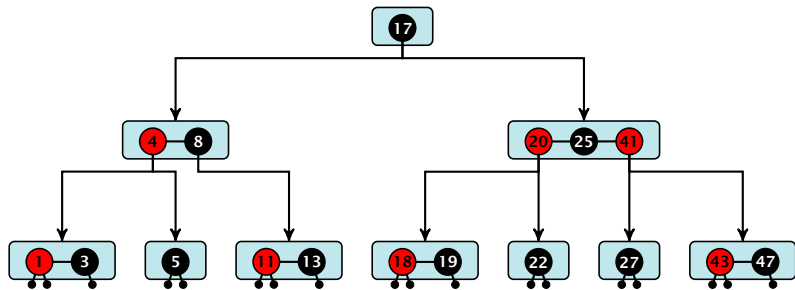
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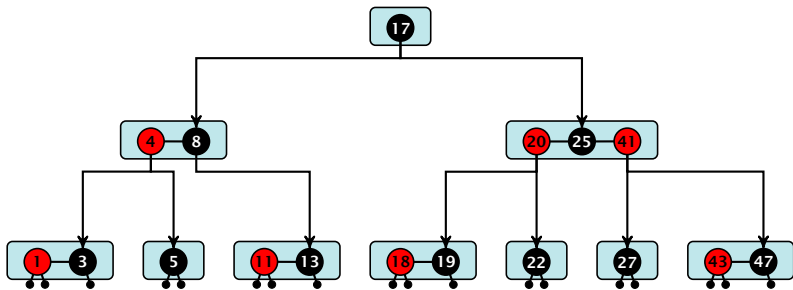
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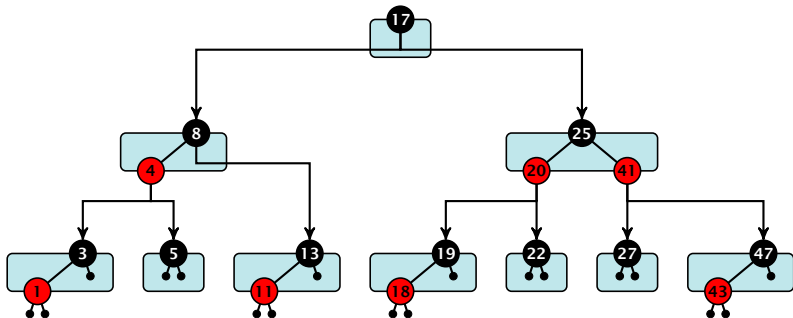
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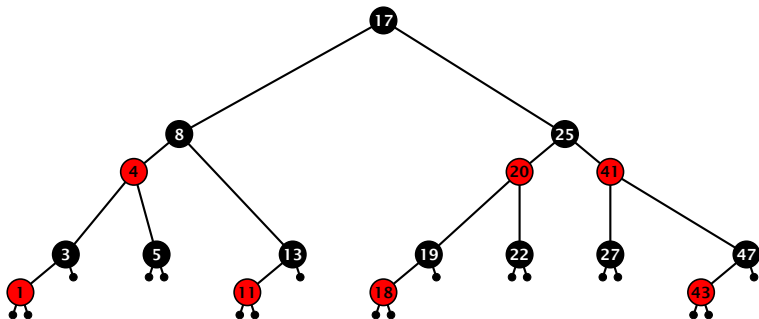
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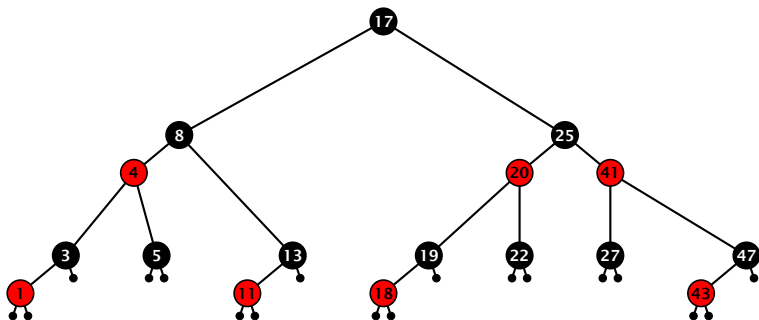
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Note that this correspondence is not unique. In particular, there are different red-black trees that correspond to the same (2, 4)-tree.

7.5 Skip Lists

Why do we not use a list for implementing the ADT Dynamic Set?

- ▶ time for search $\Theta(n)$
- ▶ time for insert $\Theta(n)$ (dominated by searching the item)
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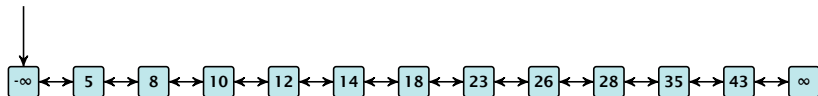
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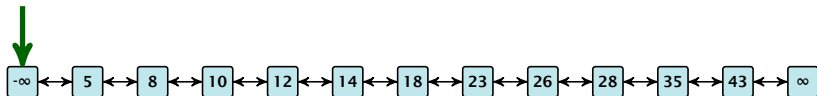
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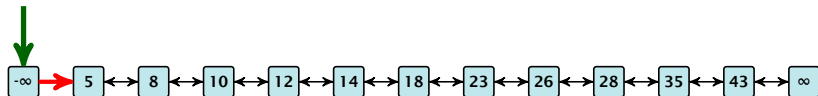
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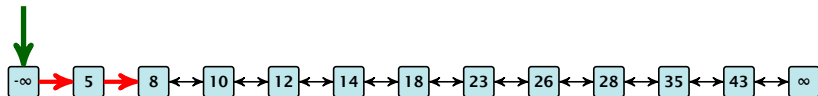
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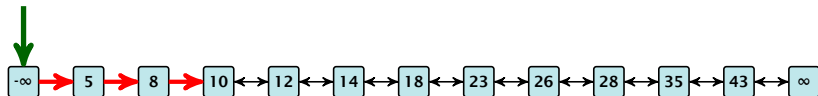
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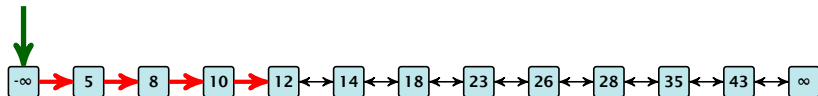
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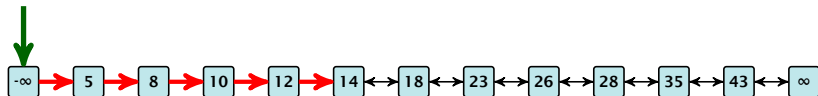
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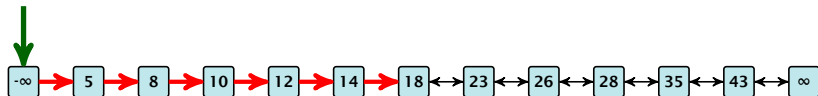
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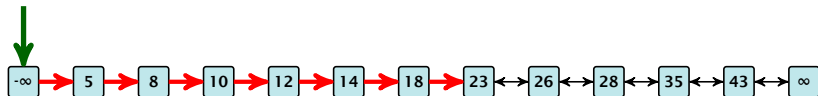
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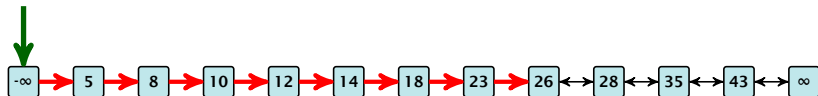
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How can we improve the search-operation?

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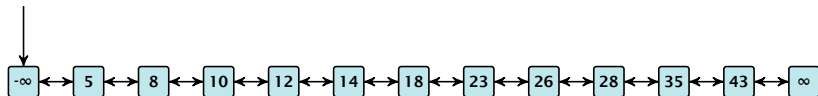
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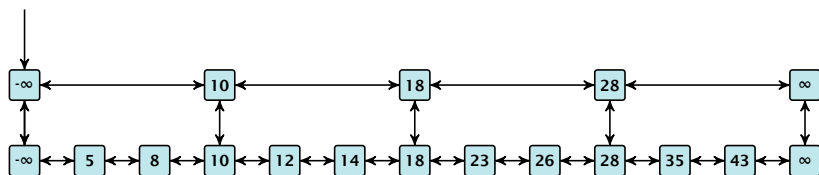
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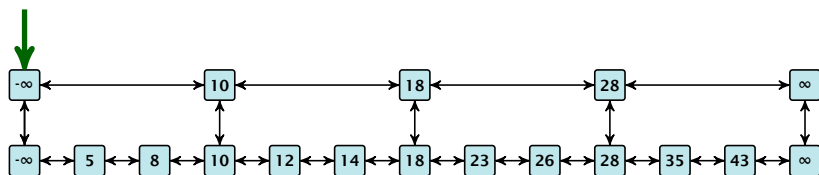
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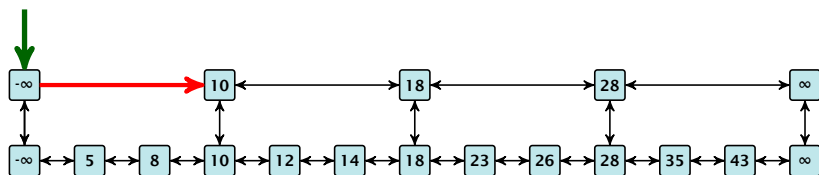
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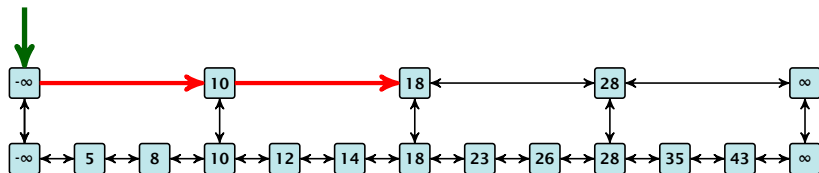
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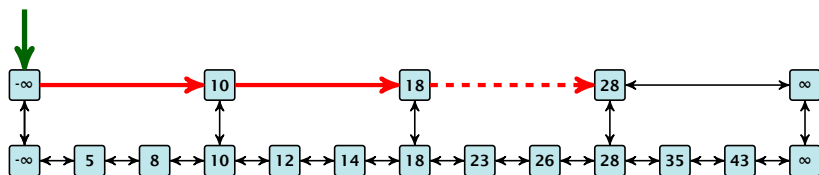
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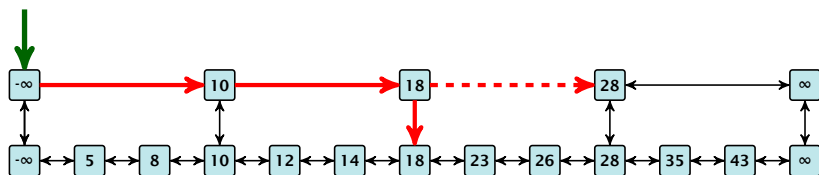
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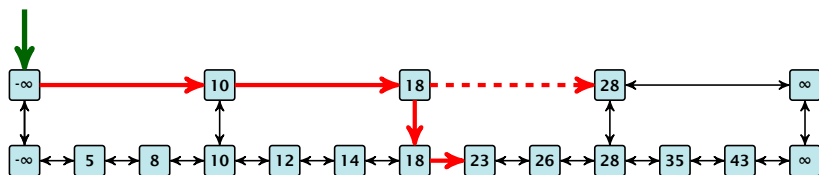
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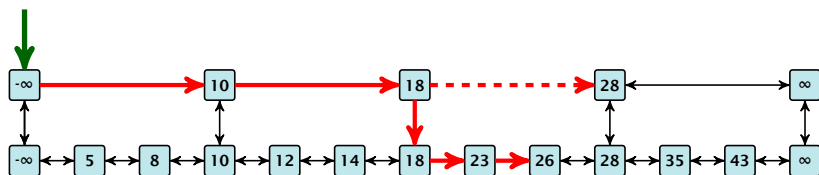
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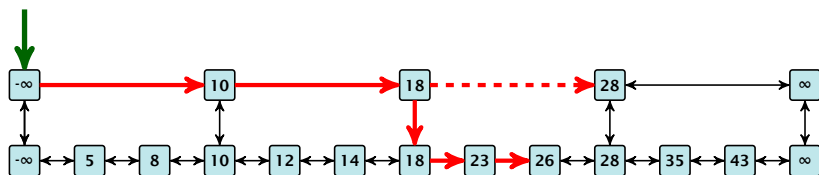
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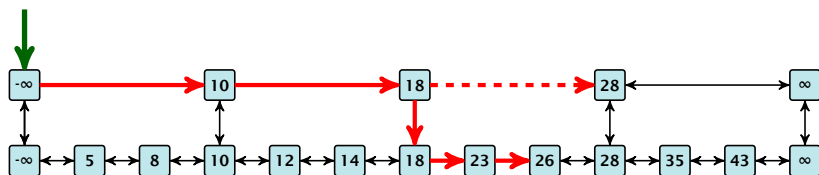


Let $|L_1|$ denote the number of elements in the “express lane”, and $|L_0| = n$ the number of all elements (ignoring dummy elements).

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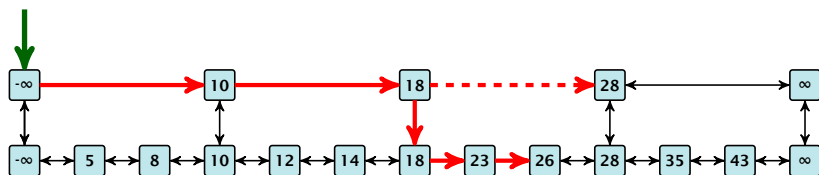
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Choose $|L_1| = \sqrt{n}$. Then search time $\Theta(\sqrt{n})$.

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Add more express lanes. Lane L_i contains roughly every $\frac{L_{i-1}}{L_i}$ -th item from list L_{i-1} .

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- ▶ At most $|L_k| + \sum_{i=1}^k \frac{L_{i-1}}{L_i} + 3(k + 1)$ steps.

7.5 Skip Lists

Choose ratios between list-lengths evenly, i.e., $\frac{|L_{i-1}|}{|L_i|} = r$, and, hence, $L_k \approx r^{-k}n$.

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$$r = k^{+1}\sqrt[k]{n} \quad \Rightarrow \quad \text{time: } \mathcal{O}(k^{k+1}\sqrt[k]{n})$$

Choosing $k = \Theta(\log n)$ gives a logarithmic running time.

7.5 Skip Lists

How to do insert and delete?

• The answer is by using a randomized skip list. Insert and delete are simple operations. The only operation that requires a lot of reorganization is the search.

Use randomization instead!

7.5 Skip Lists

How to do insert and delete?

- ▶ If we want that in L_i we always skip over roughly the same number of elements in L_{i-1} an insert or delete may require a lot of re-organisation.

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Insert:

- ▶ A search operation gives you the insert position for element x in every list.
- ▶ Flip a coin until it shows head, and record the number $t \in \{1, 2, \dots\}$ of trials needed.
- ▶ Insert x into lists L_0, \dots, L_{t-1} .

Delete:

- ▶ You get all predecessors via backward pointers.
- ▶ Delete x in all lists it actually appears in.

The time for both operation is dominated by the search time.

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Find x in all nodes on the backward pointers.

Remove x from all nodes it actually appears in.

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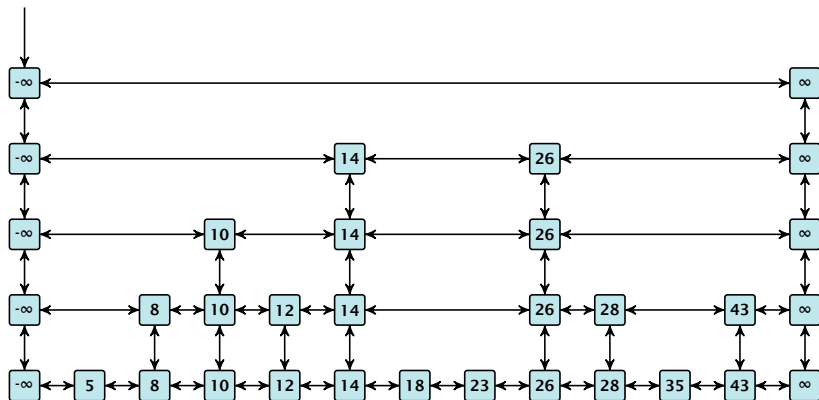
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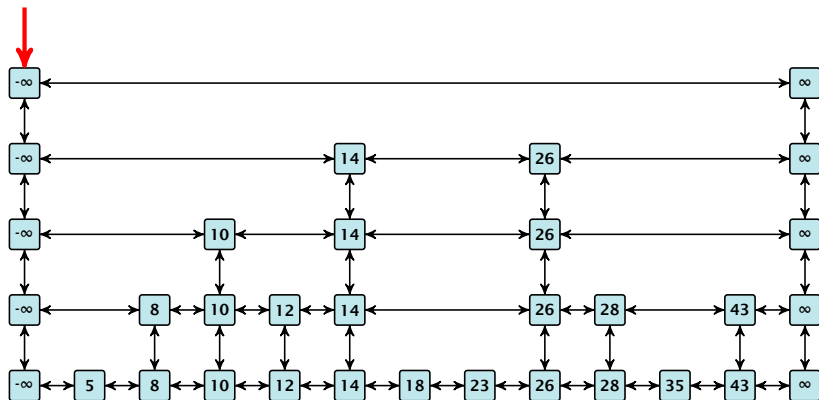
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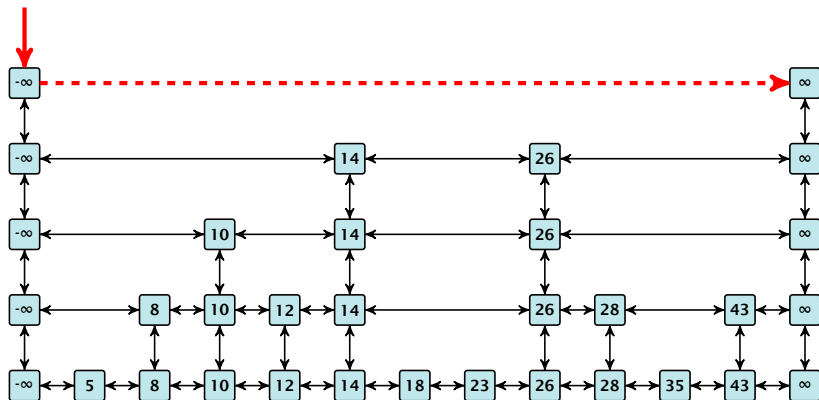
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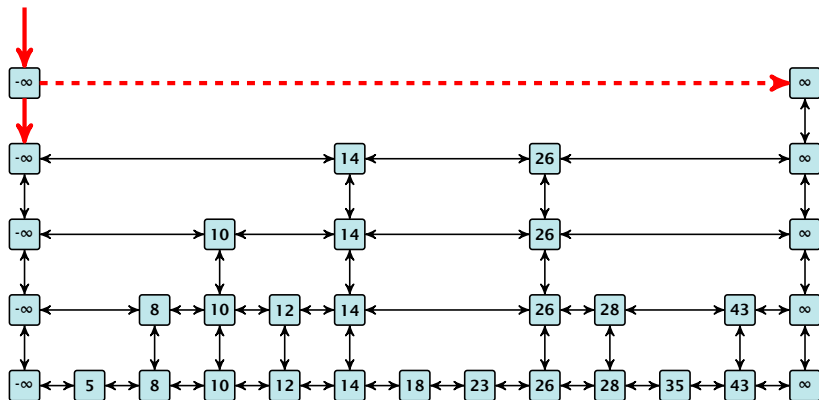
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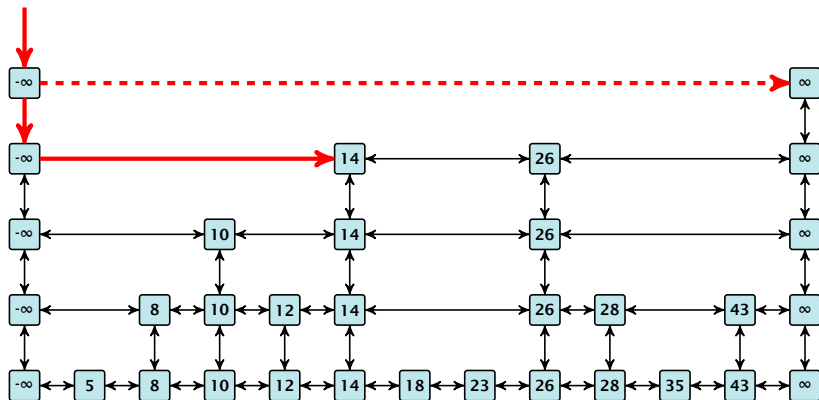
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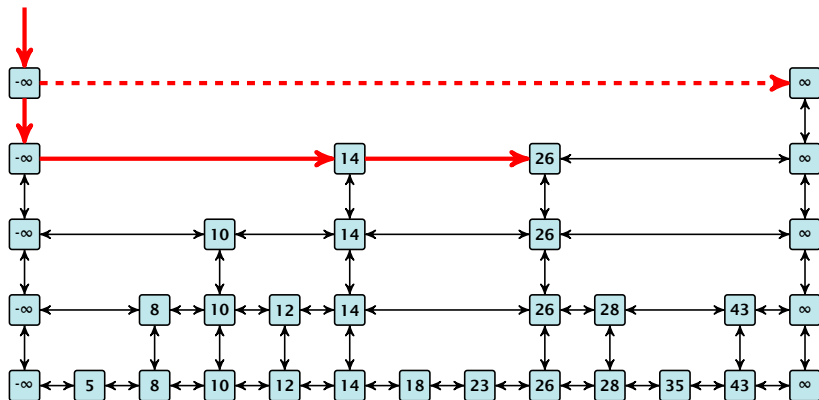
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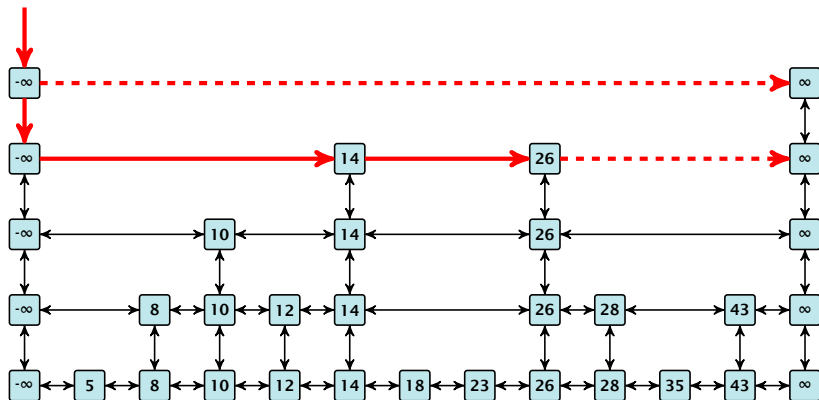
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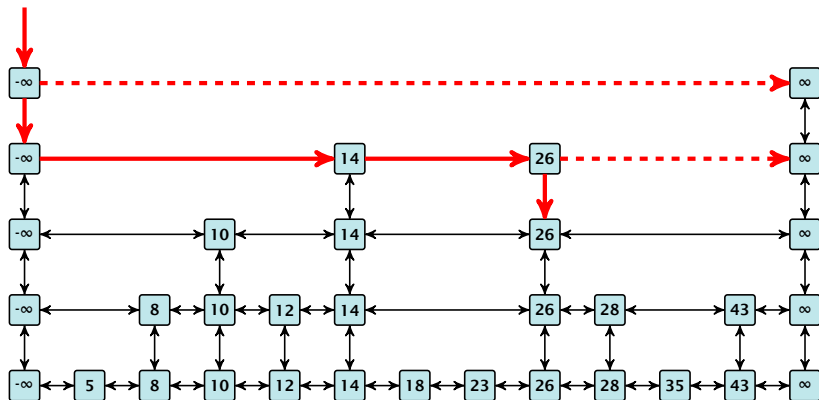
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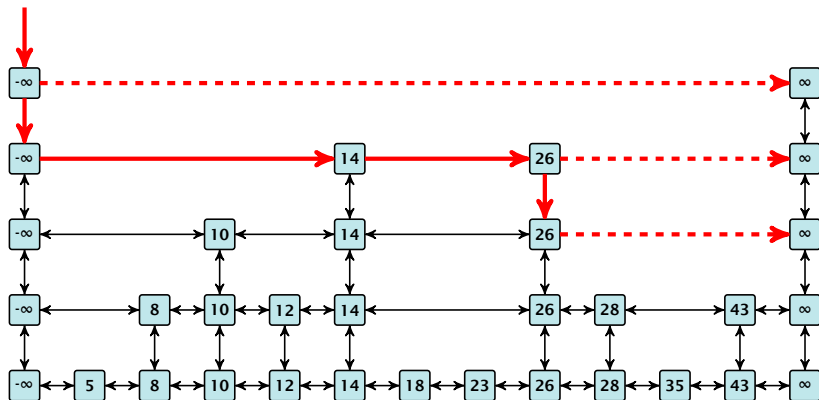
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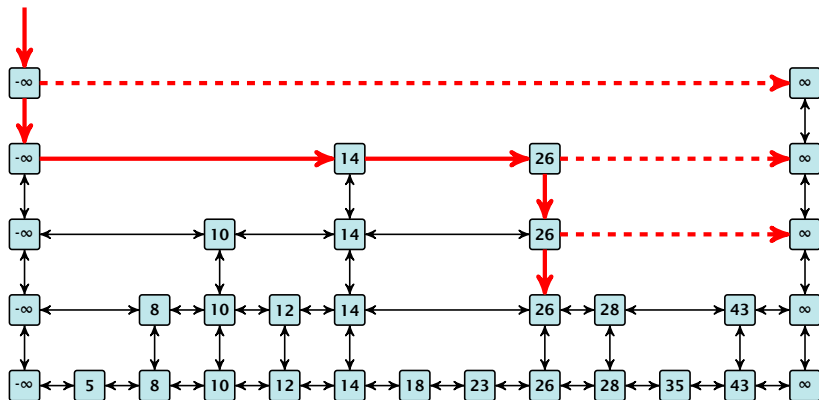
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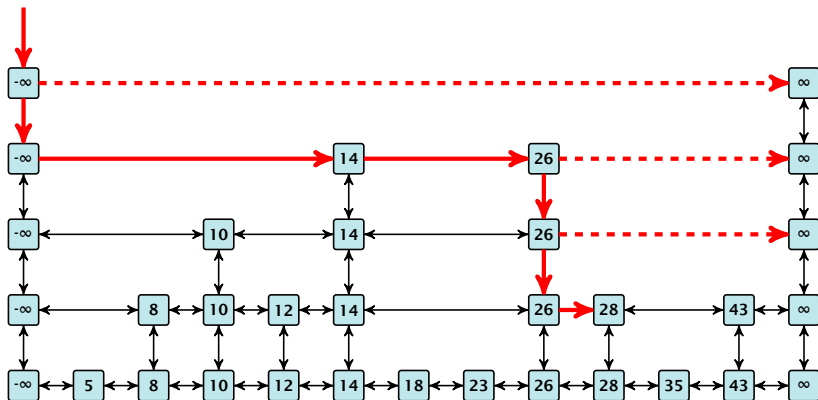
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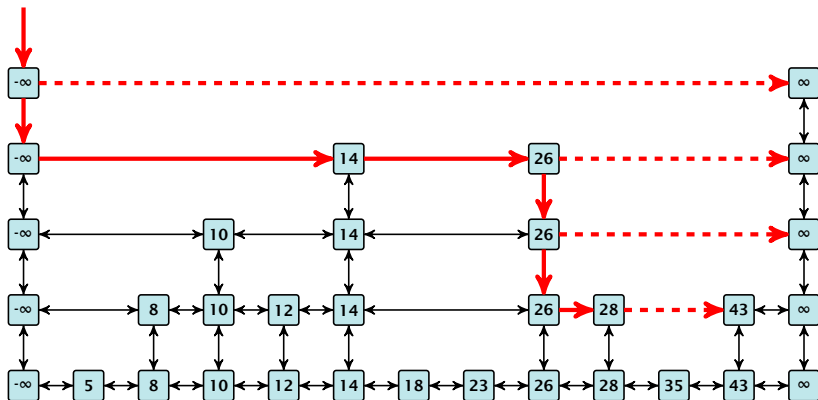
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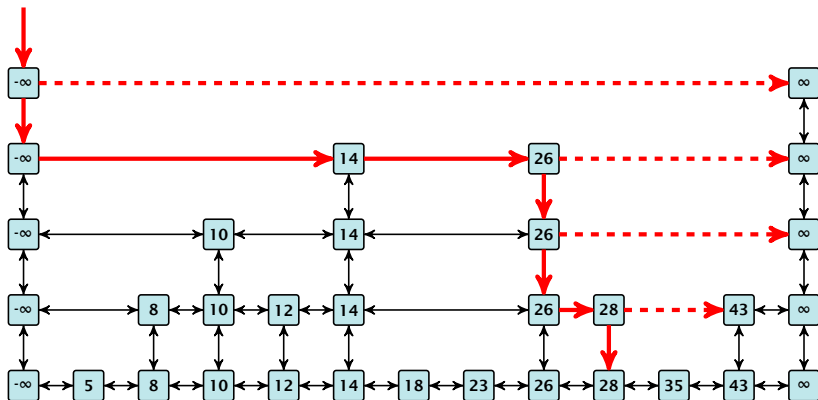
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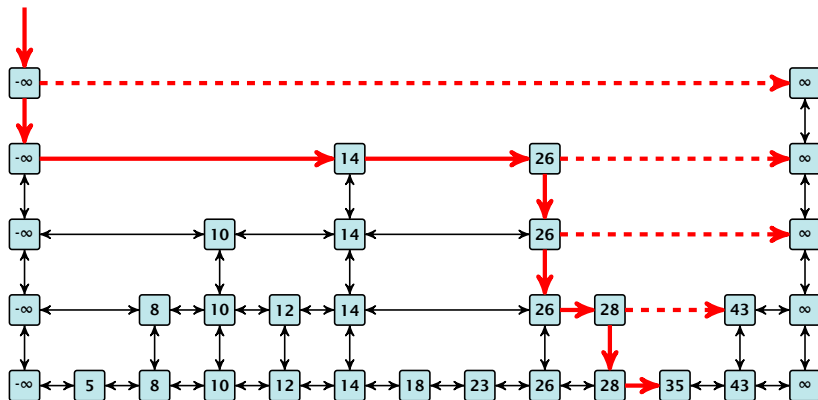
Skip Lists

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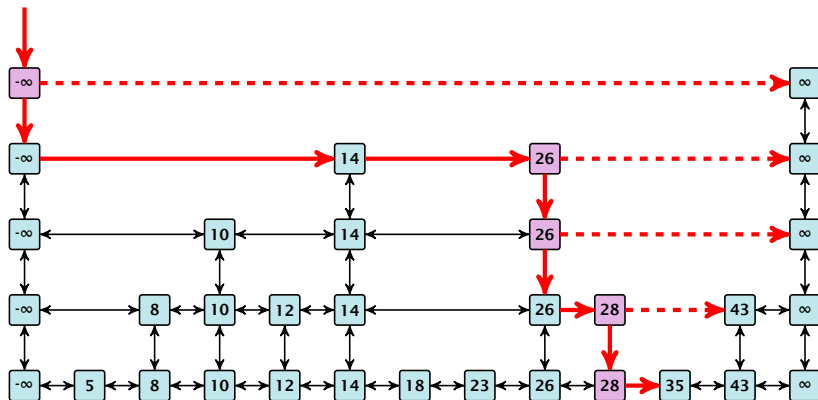
Skip Lists

Insert (35):



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7.5 Skip Lists

Lemma 20

A search (and, hence, also insert and delete) in a skip list with n elements takes time $\mathcal{O}(\log n)$ with high probability (w. h. p.).

This means for any constant α the search takes time $\mathcal{O}(\log n)$ with probability at least $1 - \frac{1}{n^\alpha}$.

Note that the constant in the \mathcal{O} -notation may depend on α .

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Suppose there are a **polynomially** many events E_1, E_2, \dots, E_ℓ , $\ell = n^c$ each holding with high probability (e.g. E_i may be the event that the i -th search in a skip list takes time at most $\mathcal{O}(\log n)$).

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Skip Lists

Backward analysis:



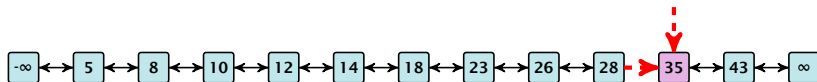
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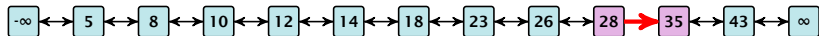
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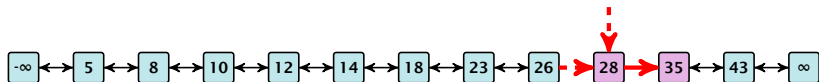
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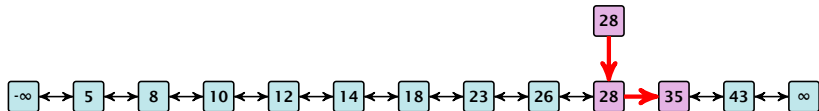
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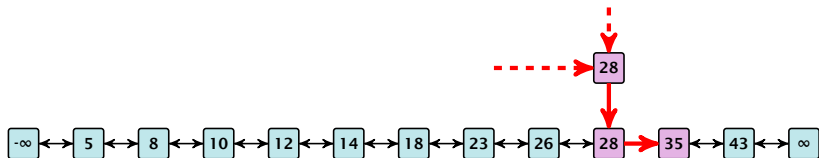
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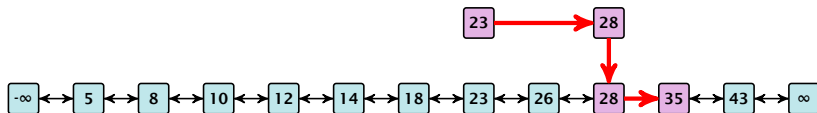
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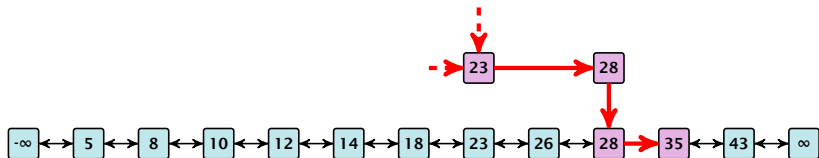
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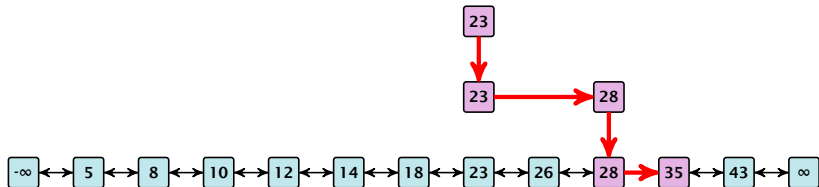
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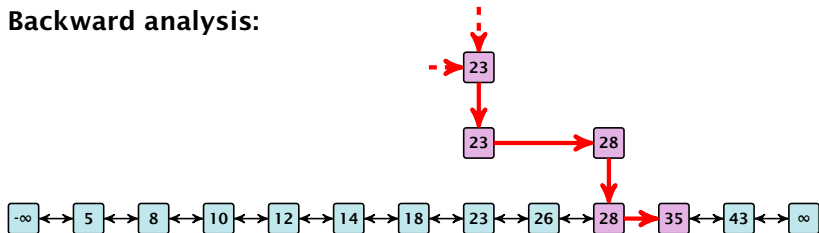
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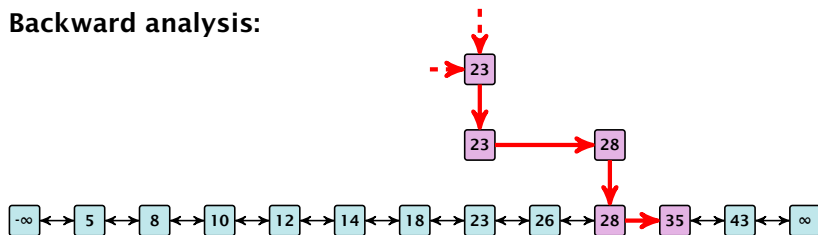
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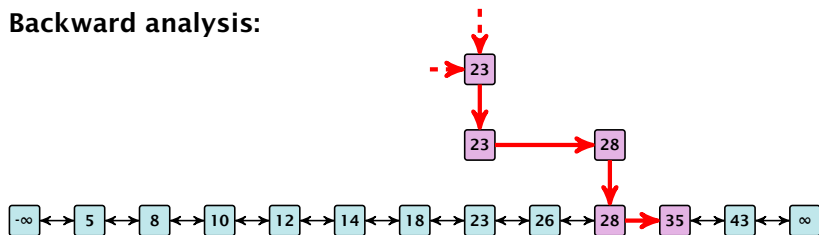
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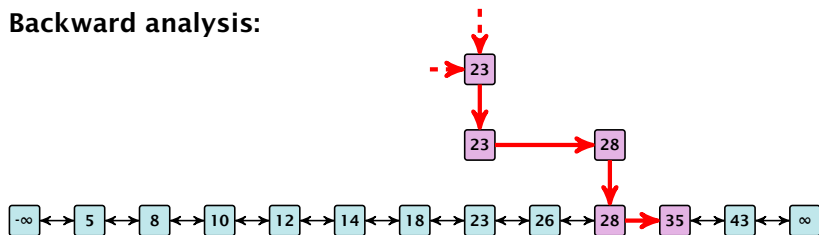
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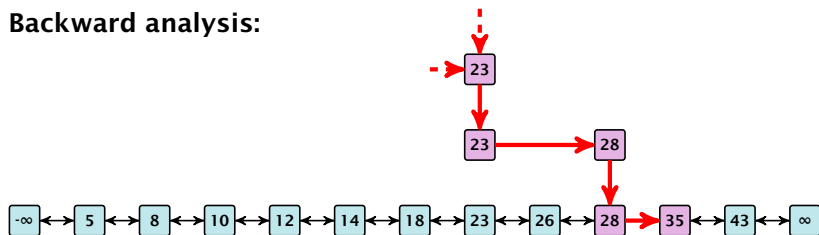
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From this it follows that w.h.p. there are no long paths.

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In particular, this means that during the construction in the backward analysis we see at most k heads (i.e., coin flips that tell you to go up) in z trials.

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This means, the search requires at most z steps, w. h. p.

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Suppose you want to develop a data structure with:

- ▶ **Insert(x):** insert element x .
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1. choose an underlying data-structure
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Goal: Design a data-structure that supports insert, delete, search, and find-by-rank in time $\mathcal{O}(\log n)$.

1. We choose a red-black tree as the underlying data-structure.
2. We store in each node v the size of the sub-tree rooted at v .
3. We need to be able to update the size-field in each node without asymptotically affecting the running time of insert, delete, and search. We come back to this step later...

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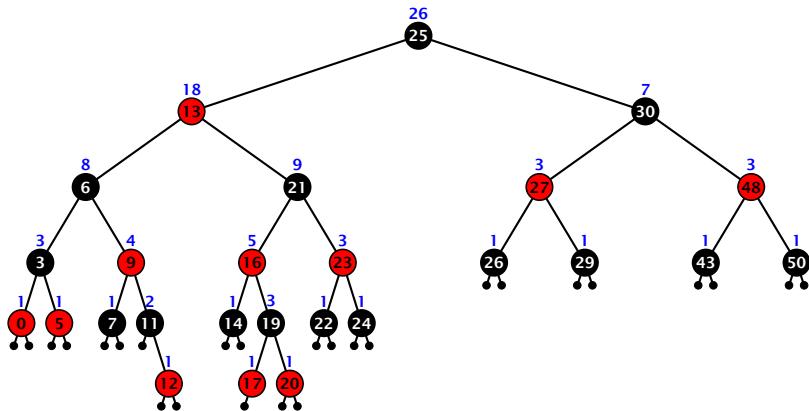
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4. How does find-by-rank work?
Find-by-rank(k) := Select(root, k) with

Algorithm 15 Select(x, i)

```
1: if  $x = \text{null}$  then return error
2: if left[ $x$ ]  $\neq$  null then  $r \leftarrow$  left[ $x$ ].size + 1 else  $r \leftarrow 1$ 
3: if  $i = r$  then return  $x$ 
4: if  $i < r$  then
5:     return Select(left[ $x$ ],  $i$ )
6: else
7:     return Select(right[ $x$ ],  $i - r$ )
```

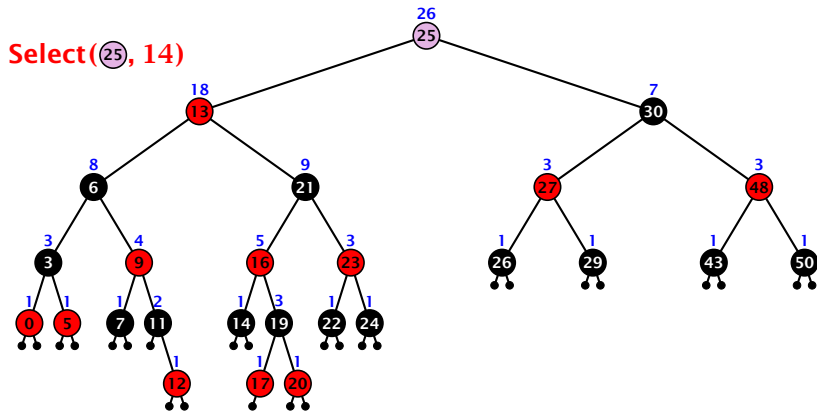
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Find-by-rank:

- ▶ decide whether you have to proceed into the left or right sub-tree
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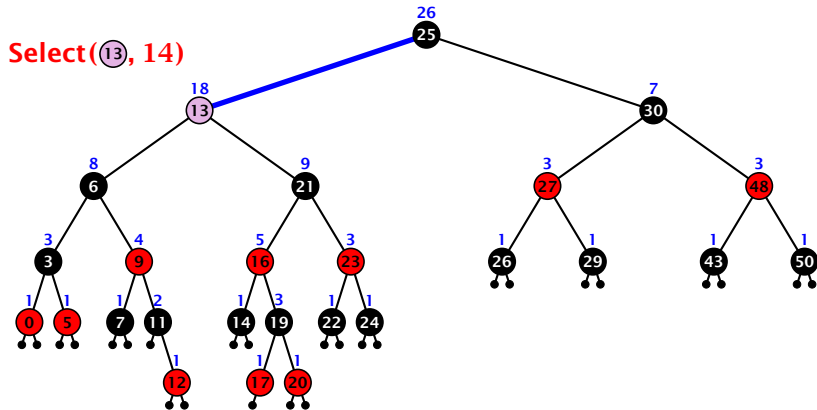
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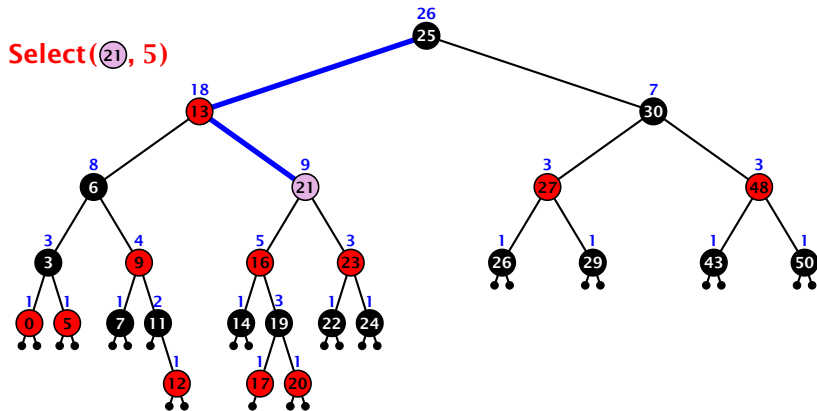
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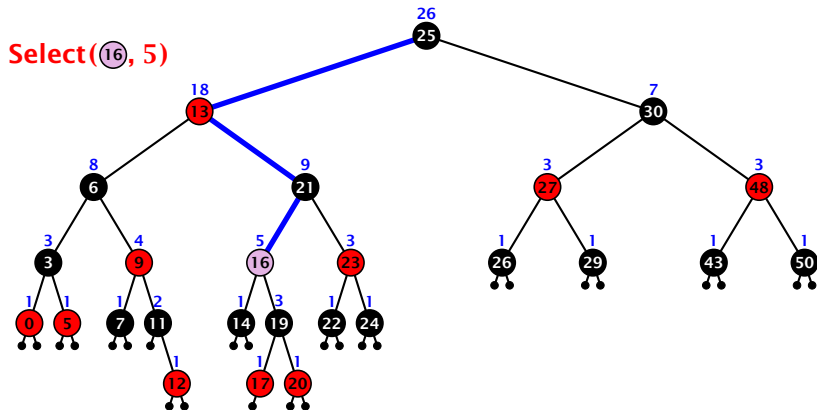
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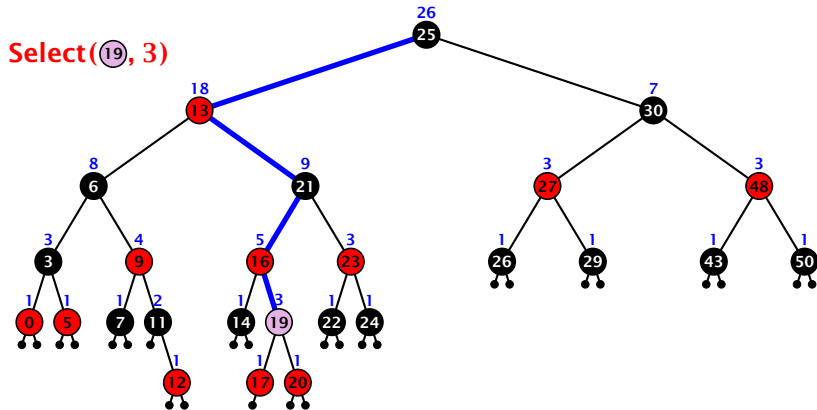
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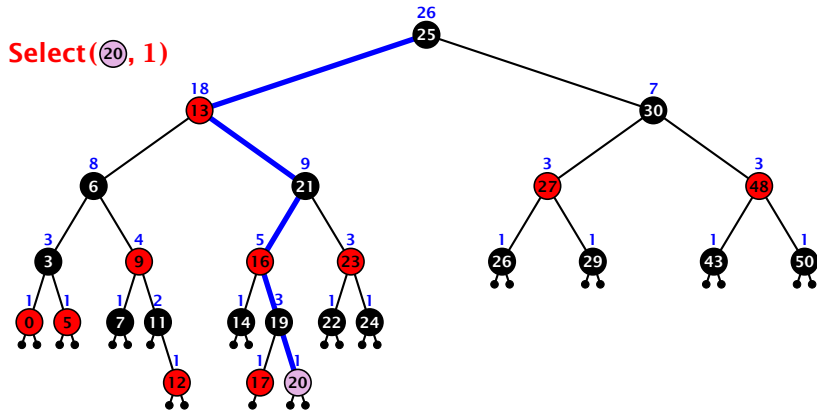
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3. How do we maintain information?

Search(k): Nothing to do.

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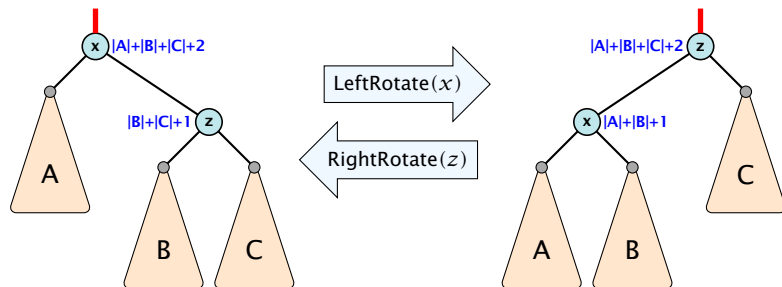
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Rotations

The only operation during the fix-up procedure that alters the tree and requires an update of the size-field:



The nodes x and z are the only nodes changing their size-fields.

The new size-fields can be computed **locally** from the size-fields of the children.

7.7 Hashing

Dictionary:

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- ▶ Set $S \subseteq U$ of keys, $|S| = m \leq n$.
- ▶ Array $T[0, \dots, n-1]$ hash-table.
- ▶ Hash function $h : U \rightarrow [0, \dots, n-1]$.

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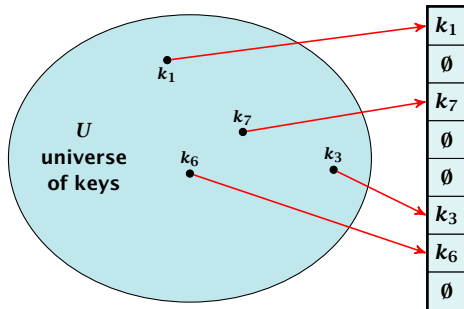
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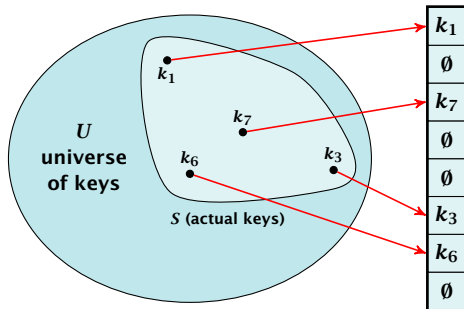
Ideally the hash function maps **all** keys to different memory locations.



This special case is known as **Direct Addressing**. It is usually very unrealistic as the universe of keys typically is quite large, and in particular larger than the available memory.

7.7 Hashing

Suppose that we **know** the set S of actual keys (no insert/no delete). Then we may want to design a **simple** hash-function that maps all these keys to different memory locations.



Such a hash function h is called a **perfect hash function** for set S .

7.7 Hashing

If we do not know the keys in advance, the best we can hope for is that the hash function distributes keys evenly across the table.

Problem: Collisions

Usually the universe U is much larger than the table-size n .

Hence, there may be two elements k_1, k_2 from the set S that map to the same memory location (i.e., $h(k_1) = h(k_2)$). This is called a collision.

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Typically, collisions do not appear once the size of the set S of actual keys gets close to n , but already once $|S| \geq \omega(\sqrt{n})$.

Lemma 21

The probability of having a collision when hashing m elements into a table of size n under uniform hashing is at least

$$1 - e^{-\frac{m(m-1)}{2n}} \approx 1 - e^{-\frac{m^2}{2n}}.$$

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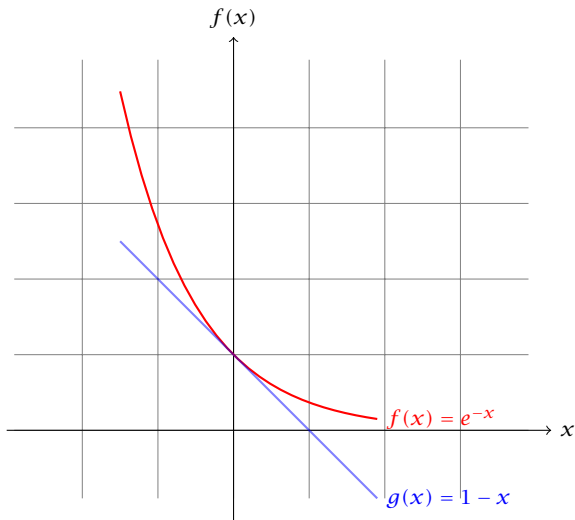
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Here the first equality follows since the ℓ -th element that is hashed has a probability of $\frac{n-\ell+1}{n}$ to not generate a collision under the condition that the previous elements did not induce collisions. □



The inequality $1 - x \leq e^{-x}$ is derived by stopping the Taylor-expansion of e^{-x} after the second term.

Resolving Collisions

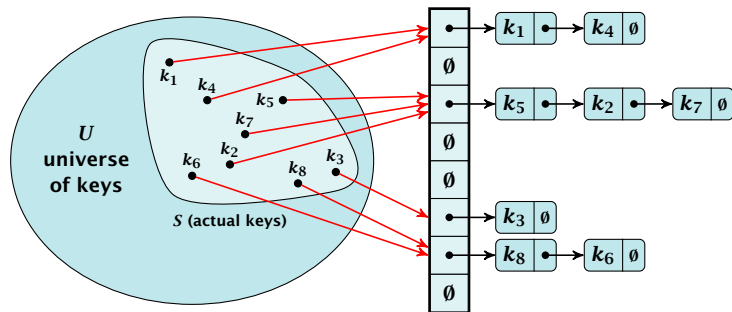
The methods for dealing with collisions can be classified into the two main types

- ▶ **open addressing**, aka. closed hashing
- ▶ **hashing with chaining**. aka. closed addressing, open hashing.

Hashing with Chaining

Arrange elements that map to the same position in a linear list.

- ▶ Access: compute $h(x)$ and search list for $\text{key}[x]$.
- ▶ Insert: insert at the front of the list.



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Let A denote a strategy for resolving collisions. We use the following notation:

- ▶ A^+ denotes the average time for a **successful** search when using A ;
- ▶ A^- denotes the average time for an **unsuccessful** search when using A ;
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Note that this result does not depend on the hash-function that is used.

Hashing with Chaining

For a successful search observe that we do **not** choose a list at random, but we consider a random key k in the hash-table and ask for the search-time for k .

This is 1 plus the number of elements that lie before k in k 's list.

Let k_ℓ denote the ℓ -th key inserted into the table.

Let for two keys k_i and k_j , X_{ij} denote the event that i and j hash to the same position. Clearly, $\Pr[X_{ij} = 1] = 1/n$ for uniform hashing.

The expected successful search cost is

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Hence, the expected cost for a successful search is $A^+ \leq 1 + \frac{\alpha}{2}$.

Open Addressing

All objects are stored in the table itself.

Define a function $h(k, j)$ that determines the table-position to be examined in the j -th step. The values $h(k, 0), \dots, h(k, n - 1)$ form a permutation of $0, \dots, n - 1$.

Search(k): Try position $h(k, 0)$; if it is empty your search fails; otherwise continue with $h(k, 1), h(k, 2), \dots$.

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Choices for $h(k, j)$:

- ▶ $h(k, i) = h(k) + i \pmod n$. Linear probing.
- ▶ $h(k, i) = h(k) + c_1 i + c_2 i^2 \pmod n$. Quadratic probing.
- ▶ $h(k, i) = h_1(k) + i h_2(k) \pmod n$. Double hashing.

For quadratic probing and double hashing one has to ensure that the search covers all positions in the table (i.e., for double hashing $h_2(k)$ must be relatively prime to n ; for quadratic probing c_1 and c_2 have to be chosen carefully).

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- ▶ Advantage: **Cache-efficiency**. The new probe position is very likely to be in the cache.
- ▶ Disadvantage: **Primary clustering**. Long sequences of occupied table-positions get longer as they have a larger probability to be hit. Furthermore, they can merge forming larger sequences.

Lemma 22

Let L be the method of linear probing for resolving collisions:

$$L^+ \approx \frac{1}{2} \left(1 + \frac{1}{1 - \alpha} \right)$$

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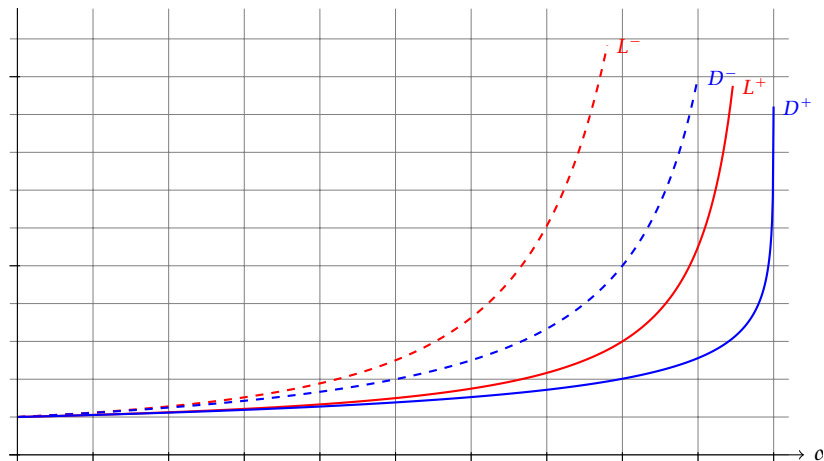
$$D^- \approx \frac{1}{1 - \alpha}$$

7.7 Hashing

Some values:

α	Linear Probing		Quadratic Probing		Double Hashing	
	L^+	L^-	Q^+	Q^-	D^+	D^-
0.5	1.5	2.5	1.44	2.19	1.39	2
0.9	5.5	50.5	2.85	11.40	2.55	10
0.95	10.5	200.5	3.52	22.05	3.15	20

7.7 Hashing



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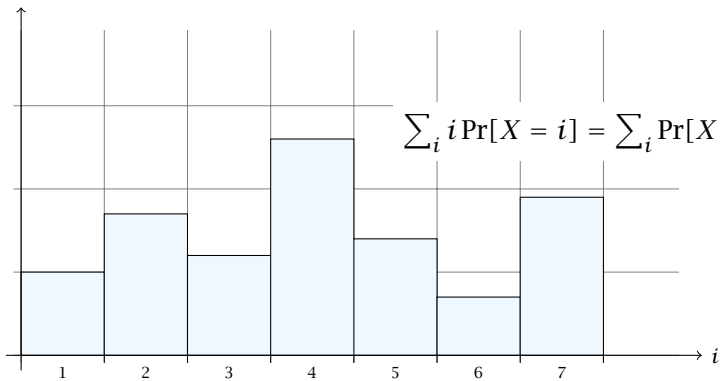
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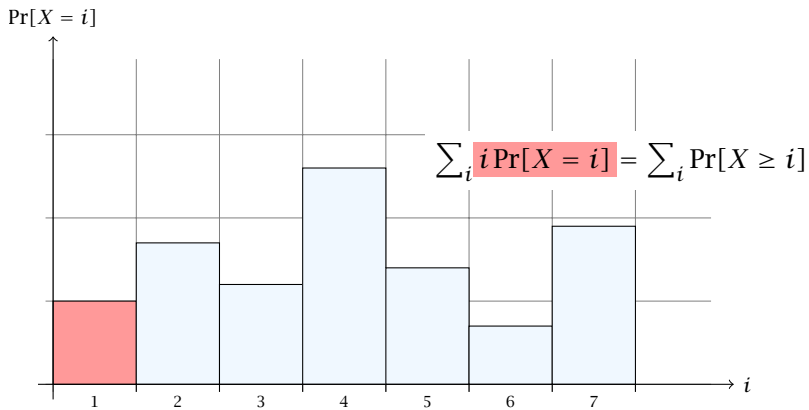
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$$\frac{1}{1 - \alpha} = 1 + \alpha + \alpha^2 + \alpha^3 + \dots$$

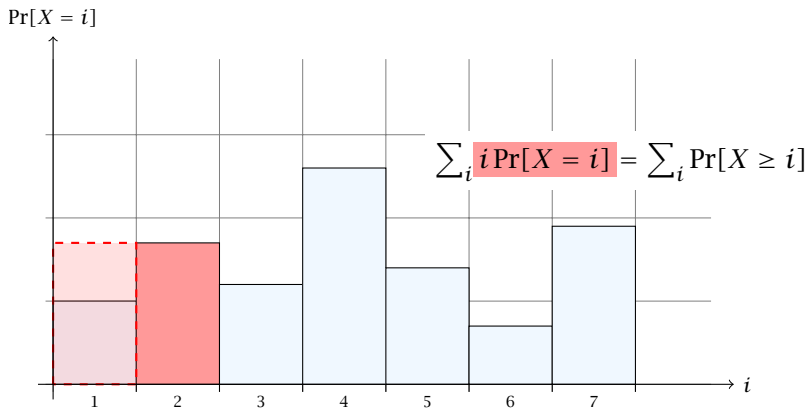
$\Pr[X = i]$



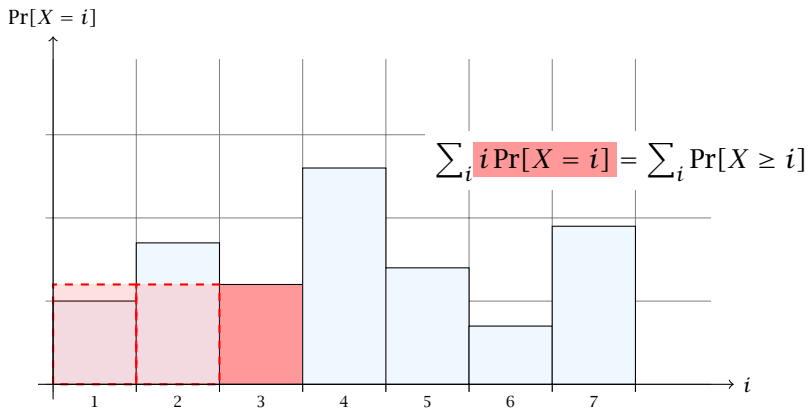
$i = 1$



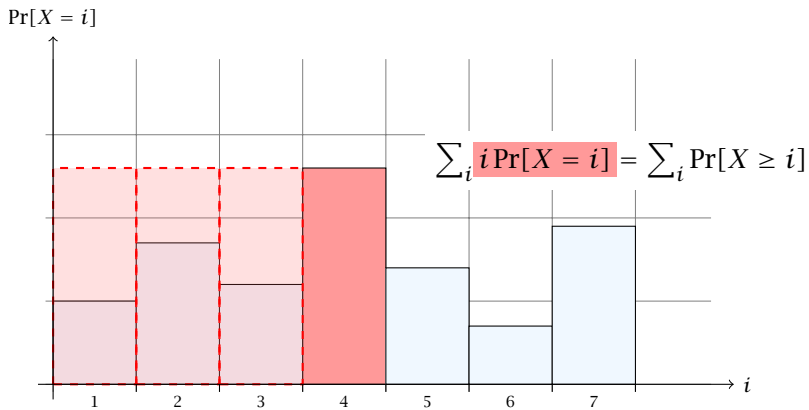
$i = 2$



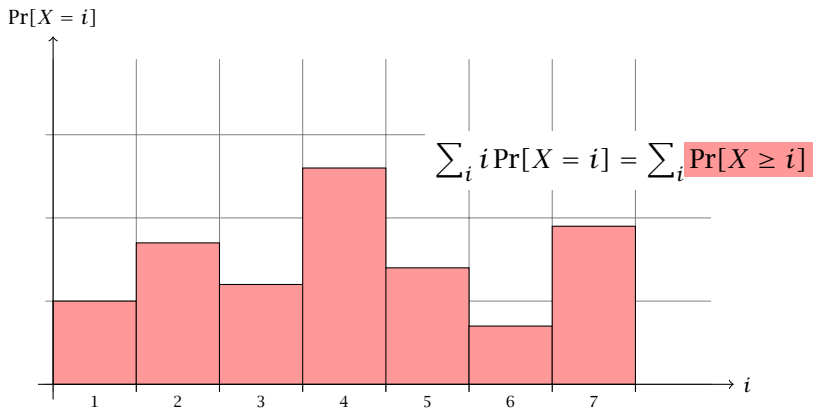
$i = 3$



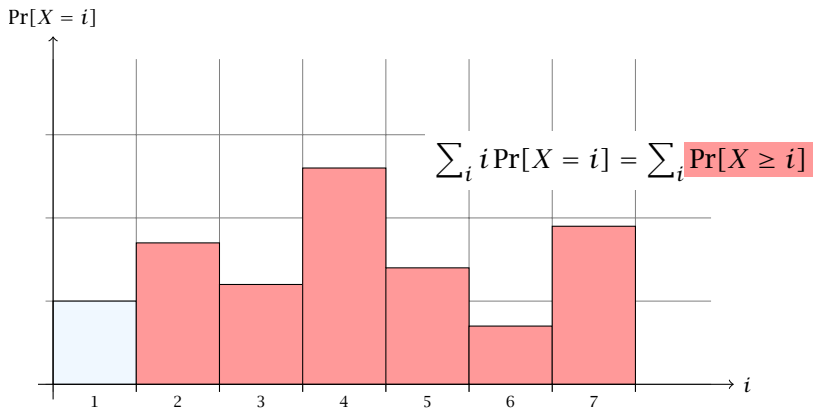
$i = 4$



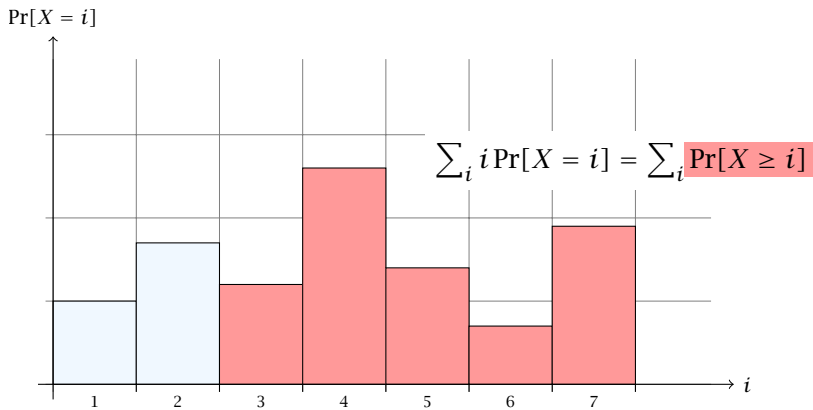
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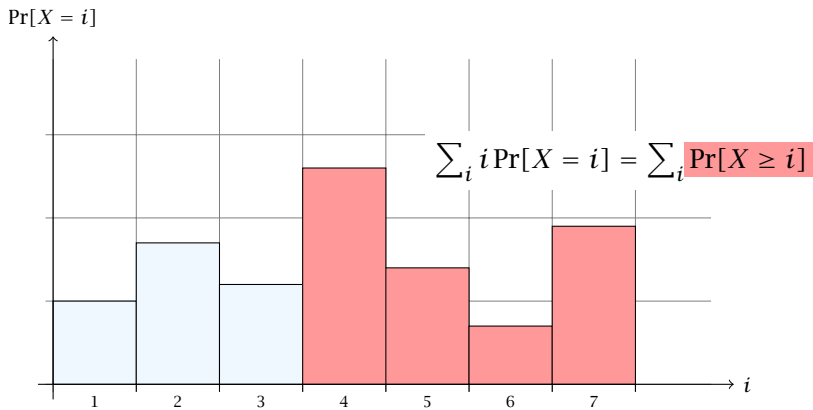
$i = 2$



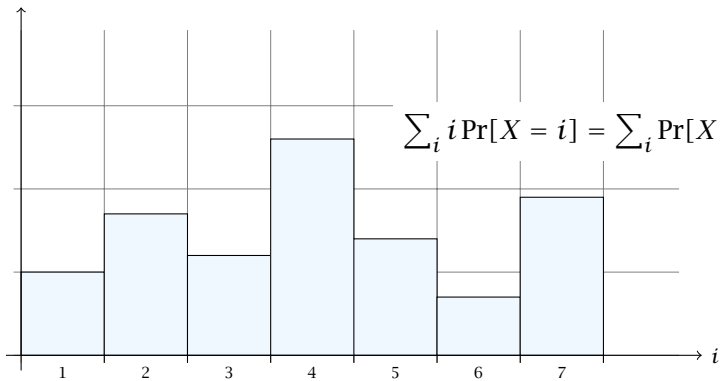
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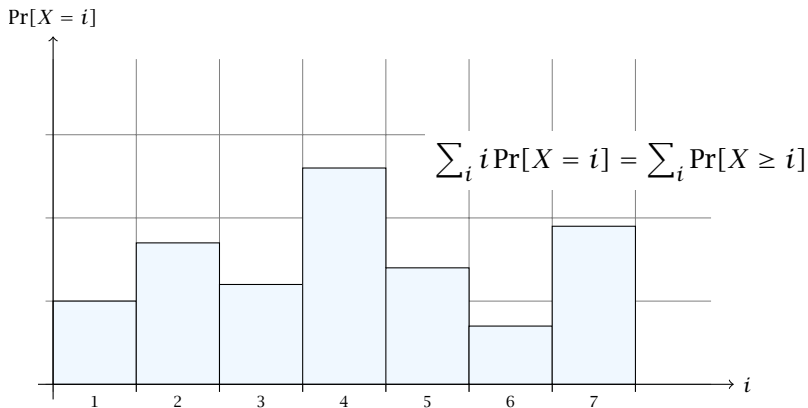


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The j -th rectangle appears in both sums j times. (j times in the first due to multiplication with j ; and j times in the second for summands $i = 1, 2, \dots, j$)

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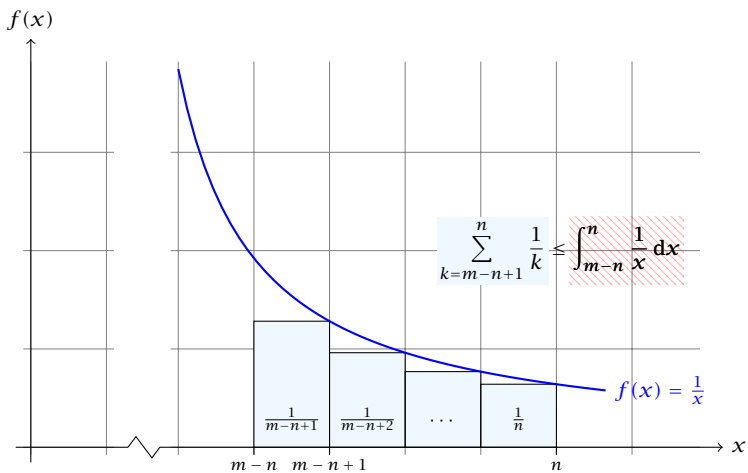
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7.7 Hashing

Definition 25

A class \mathcal{H} of hash-functions from the universe U into the set $\{0, \dots, n-1\}$ is called **universal** if for all $u_1, u_2 \in U$ with $u_1 \neq u_2$

$$\Pr[h(u_1) = h(u_2)] \leq \frac{1}{n} ,$$

where the probability is w. r. t. the choice of a random hash-function from set \mathcal{H} .

Note that this means that $\Pr[h(u_1) = h(u_2)] = \frac{1}{n}$.

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Definition 26

A class \mathcal{H} of hash-functions from the universe U into the set $\{0, \dots, n-1\}$ is called **2-independent** (pairwise independent) if the following two conditions hold

- ▶ For any key $u \in U$, and $t \in \{0, \dots, n-1\}$ $\Pr[h(u) = t] = \frac{1}{n}$, i.e., a key is distributed uniformly within the hash-table.
- ▶ For all $u_1, u_2 \in U$ with $u_1 \neq u_2$, and for any two hash-positions t_1, t_2 :

$$\Pr[h(u_1) = t_1 \wedge h(u_2) = t_2] \leq \frac{1}{n^2} .$$

Note that the probability is w. r. t. the choice of a random hash-function from set \mathcal{H} .

This requirement clearly implies a universal hash-function.

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Definition 27

A class \mathcal{H} of hash-functions from the universe U into the set $\{0, \dots, n-1\}$ is called **k -independent** if for any choice of $\ell \leq k$ distinct keys $u_1, \dots, u_\ell \in U$, and for any set of ℓ not necessarily distinct hash-positions t_1, \dots, t_ℓ :

$$\Pr[h(u_1) = t_1 \wedge \dots \wedge h(u_\ell) = t_\ell] \leq \frac{1}{n^\ell} ,$$

where the probability is w. r. t. the choice of a random hash-function from set \mathcal{H} .

7.7 Hashing

Definition 28

A class \mathcal{H} of hash-functions from the universe U into the set $\{0, \dots, n-1\}$ is called (μ, k) -independent if for any choice of $\ell \leq k$ distinct keys $u_1, \dots, u_\ell \in U$, and for any set of ℓ not necessarily distinct hash-positions t_1, \dots, t_ℓ :

$$\Pr[h(u_1) = t_1 \wedge \dots \wedge h(u_\ell) = t_\ell] \leq \left(\frac{\mu}{n}\right)^\ell,$$

where the probability is w. r. t. the choice of a random hash-function from set \mathcal{H} .

7.7 Hashing

Let $U := \{0, \dots, p-1\}$ for a prime p . Let $\mathbb{Z}_p := \{0, \dots, p-1\}$, and let $\mathbb{Z}_p^* := \{1, \dots, p-1\}$ denote the set of invertible elements in \mathbb{Z}_p .

Define

$$h_{a,b}(x) := (ax + b \bmod p) \bmod n$$

Lemma 29

The class

$$\mathcal{H} = \{h_{a,b} \mid a \in \mathbb{Z}_p^*, b \in \mathbb{Z}_p\}$$

is a universal class of hash-functions from U to $\{0, \dots, n-1\}$.

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7.7 Hashing

Proof.

Let $x, y \in U$ be two distinct keys. We have to show that the probability of a collision is only $1/n$.

$$h(x) = ax + b \equiv ay + b \pmod{p}$$

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Let $x, y \in U$ be two distinct keys. We have to show that the probability of a collision is only $1/n$.

$$\triangleright ax + b \not\equiv ay + b \pmod{p}$$

$$\text{if } x \neq y \text{ then } (x - y) \not\equiv 0 \pmod{p}$$

$$\text{multiplying with } a \not\equiv 0 \pmod{p} \text{ gives}$$

$$a(x - y) \not\equiv 0 \pmod{p}$$

Therefore, the two keys x and y are not hashed to the same value.

Since x and y were arbitrary,

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If $x \neq y$ then $(x - y) \not\equiv 0 \pmod{p}$.

Multiplying with $a \not\equiv 0 \pmod{p}$ gives

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where we use that \mathbb{Z}_p is a field (Körper) and, hence, has no zero divisors (nullteilerfrei).

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$$a \equiv (t_x - t_y)(x - y)^{-1} \pmod{p}$$

$$b \equiv ay - t_y \pmod{p}$$

7.7 Hashing

There is a one-to-one correspondence between hash-functions (pairs (a, b) , $a \neq 0$) and pairs (t_x, t_y) , $t_x \neq t_y$.

Therefore, we can view the first step (before the $(\text{mod } n)$ -operation) as choosing a pair (t_x, t_y) , $t_x \neq t_y$ uniformly at random.

What happens when we do the $(\text{mod } n)$ operation?

Fix a value t_x . There are $p - 1$ possible values for choosing t_y .

From the range $0, \dots, p - 1$ the values $t_x, t_x + n, t_x + 2n, \dots$ map to t_x after the modulo-operation. These are at most $\lceil p/n \rceil$ values.

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As $t_y \neq t_x$ there are

$$\left\lfloor \frac{n}{2} \right\rfloor - 1 < \frac{t_y - t_x}{n} < \frac{t_y + t_x}{n}$$

possibilities for choosing t_y such that the final hash-value creates a collision.

This happens with probability at most $\frac{1}{n}$.

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$$\Pr_{t_x \neq t_y \in \mathbb{Z}_p^2} \left[\begin{array}{l} t_x \bmod n = h_1 \\ \wedge \\ t_y \bmod n = h_2 \end{array} \right]$$

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It is also possible to show that \mathcal{H} is an (almost) pairwise independent class of hash-functions.

$$\frac{\left\lfloor \frac{p}{n} \right\rfloor^2}{p(p-1)} \leq \Pr_{t_x \neq t_y \in \mathbb{Z}_p^2} \left[\begin{array}{c} t_x \bmod n = h_1 \\ \wedge \\ t_y \bmod n = h_2 \end{array} \right] \leq \frac{\left\lfloor \frac{p}{n} \right\rfloor^2}{p(p-1)}$$

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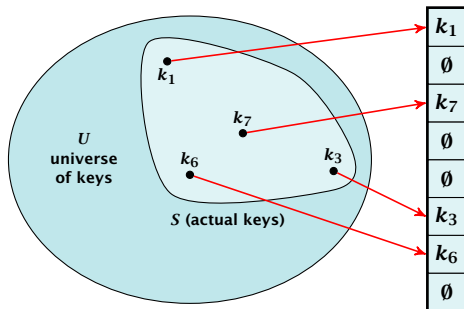
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Note that the middle is the probability that $h(x) = h_1$ and $h(y) = h_2$. The total number of choices for (t_x, t_y) is $p(p-1)$. The number of choices for t_x (t_y) such that $t_x \bmod n = h_1$ ($t_y \bmod n = h_2$) lies between $\lfloor \frac{p}{n} \rfloor$ and $\lceil \frac{p}{n} \rceil$.

Perfect Hashing

Suppose that we **know** the set S of actual keys (no insert/no delete). Then we may want to design a **simple** hash-function that maps all these keys to different memory locations.



Perfect Hashing

Let $m = |S|$. We could simply choose the hash-table size very large so that we don't get any collisions.

Using a universal hash-function the expected number of collisions is

$$E[\#\text{Collisions}] = \binom{m}{2} \cdot \frac{1}{n}.$$

If we choose $n = m^2$ the **expected number** of collisions is strictly less than $\frac{1}{2}$.

Can we get an upper bound on the **probability of having collisions**?

The probability of having 1 or more collisions can be at most $\frac{1}{2}$ as otherwise the expectation would be larger than $\frac{1}{2}$.

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We can find such a hash-function by a few trials.

However, a hash-table size of $n = m^2$ is very very high.

We construct a two-level scheme. We first use a hash-function that maps elements from S to m buckets.

Let m_j denote the number of items that are hashed to the j -th bucket. For each bucket we choose a second hash-function that maps the elements of the bucket into a table of size m_j^2 . The second function can be chosen such that all elements are mapped to different locations.

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$$E \left[\sum_j m_j^2 \right] = E \left[2 \sum_j \binom{m_j}{2} + \sum_j m_j \right]$$

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The first expectation is simply the expected number of collisions, for the first level.

Perfect Hashing

The total memory that is required by all hash-tables is $\sum_j m_j^2$.

$$\begin{aligned} E \left[\sum_j m_j^2 \right] &= E \left[2 \sum_j \binom{m_j}{2} + \sum_j m_j \right] \\ &= 2 E \left[\sum_j \binom{m_j}{2} \right] + E \left[\sum_j m_j \right] \end{aligned}$$

The first expectation is simply the expected number of collisions, for the first level.

$$= 2 \binom{m}{2} \frac{1}{m} + m = 2m - 1$$

Perfect Hashing

We need only $\mathcal{O}(m)$ time to construct a hash-function h with $\sum_j m_j^2 = \mathcal{O}(4m)$.

Then we construct a hash-table h_j for every bucket. This takes expected time $\mathcal{O}(m_j)$ for every bucket.

We only need that the hash-function is universal!!!

Cuckoo Hashing

Goal:

Try to generate a perfect hash-table (constant worst-case search time) in a dynamic scenario.

- Two hash-tables $T_1[0, \dots, m-1]$ and $T_2[0, \dots, m-1]$, with hash functions h_1 and h_2 .
- An object x is either stored at location $T_1[h_1(x)]$ or $T_2[h_2(x)]$.
- Insertion and deletion takes constant time if the algorithm doesn't fail.

Cuckoo Hashing

Goal:

Try to generate a perfect hash-table (constant worst-case search time) in a dynamic scenario.

- ▶ Two hash-tables $T_1[0, \dots, n - 1]$ and $T_2[0, \dots, n - 1]$, with hash-functions h_1 , and h_2 .
- ▶ An object x is either stored at location $T_1[h_1(x)]$ or $T_2[h_2(x)]$.
- ▶ A search clearly takes constant time if the above constraint is met.

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Cuckoo Hashing

Insert:

\emptyset
\emptyset
x_7
\emptyset
\emptyset
x_4
x_1
\emptyset
\emptyset

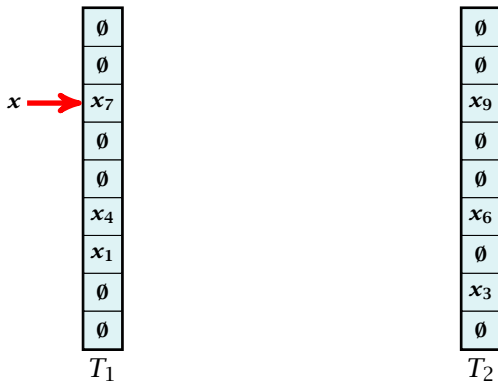
T_1

\emptyset
\emptyset
x_9
\emptyset
\emptyset
x_6
\emptyset
x_3
\emptyset

T_2

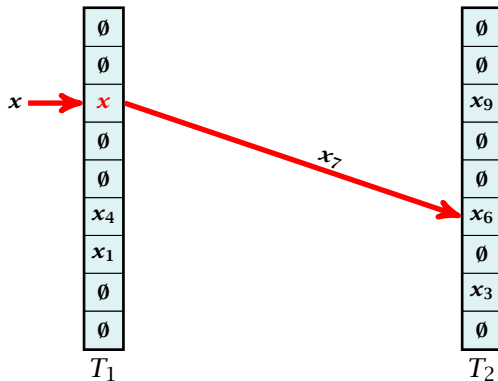
Cuckoo Hashing

Insert:



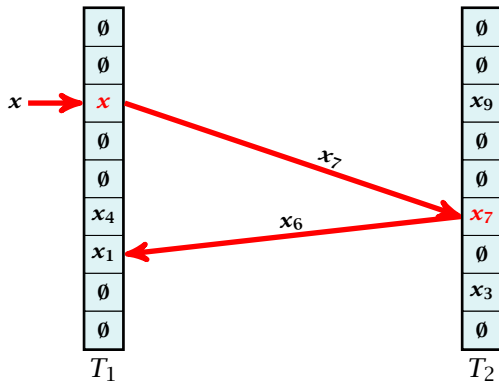
Cuckoo Hashing

Insert:



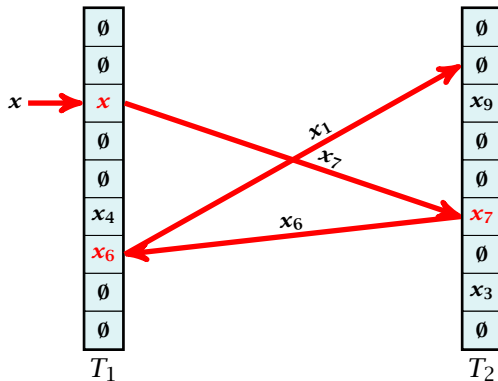
Cuckoo Hashing

Insert:



Cuckoo Hashing

Insert:



Cuckoo Hashing

Algorithm 16 Cuckoo-Insert(x)

```
1: if  $T_1[h_1(x)] = x \vee T_2[h_2(x)] = x$  then return  
2: steps  $\leftarrow 1$   
3: while steps  $\leq$  maxsteps do  
4:     exchange  $x$  and  $T_1[h_1(x)]$   
5:     if  $x = \text{null}$  then return  
6:     exchange  $x$  and  $T_2[h_2(x)]$   
7:     if  $x = \text{null}$  then return  
8: rehash() // change table-size and rehash everything  
9: Cuckoo-Insert( $x$ )
```

Cuckoo Hashing

What is the expected time for an insert-operation?

We first analyze the probability that we end-up in an infinite loop (that is then terminated after maxsteps steps).

Formally what is the probability to enter an infinite loop that touches ℓ different keys (apart from x)?

What is the expected time for an insert-operation?

We first analyze the probability that we end-up in an infinite loop (that is then terminated after maxsteps steps).

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Cuckoo Hashing

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Cuckoo Hashing

Insert:



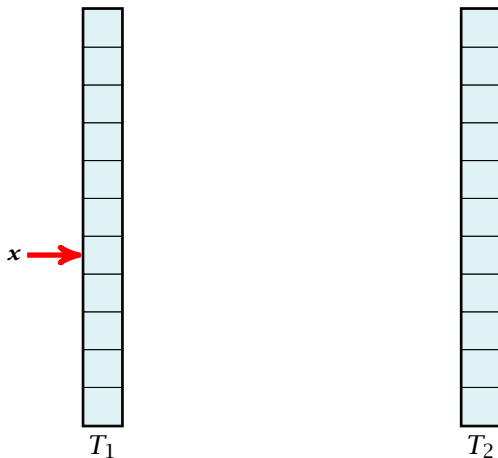
T_1



T_2

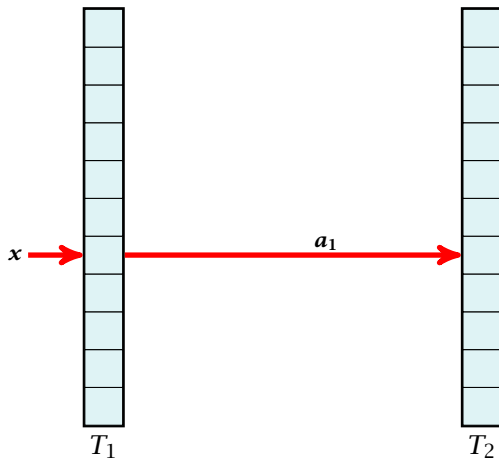
Cuckoo Hashing

Insert:



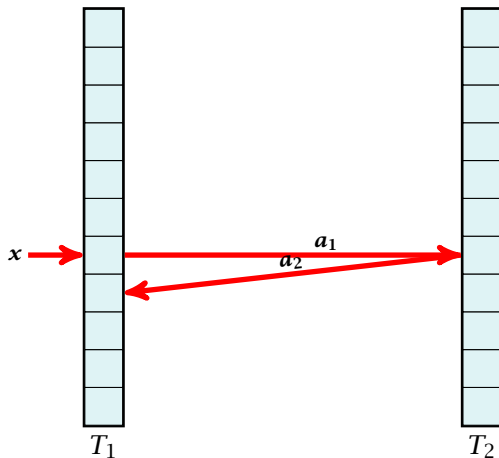
Cuckoo Hashing

Insert:



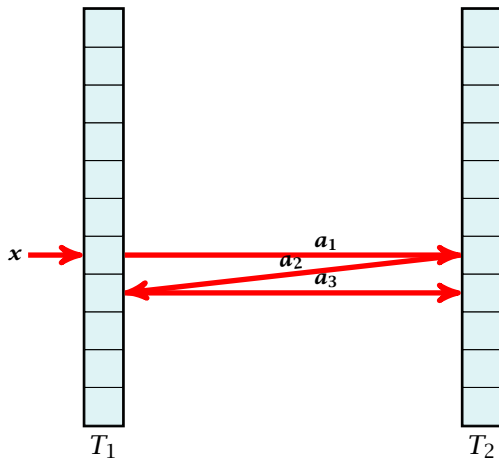
Cuckoo Hashing

Insert:



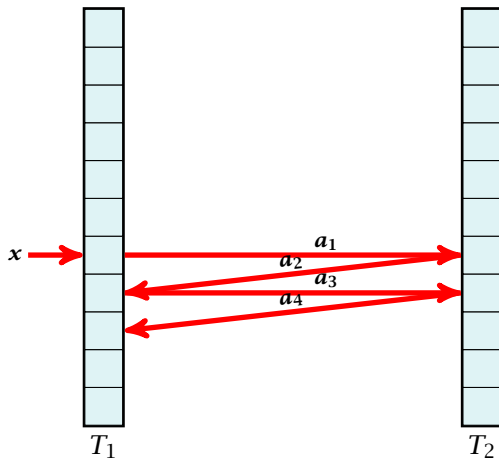
Cuckoo Hashing

Insert:



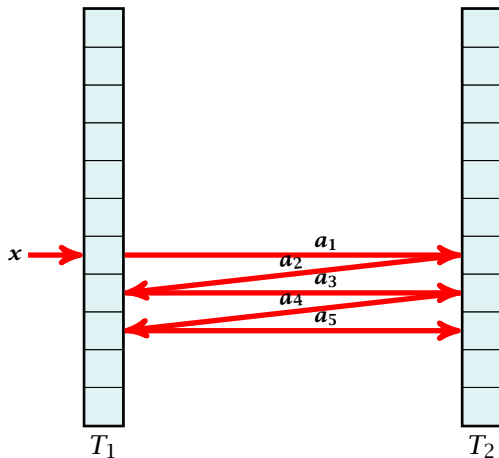
Cuckoo Hashing

Insert:



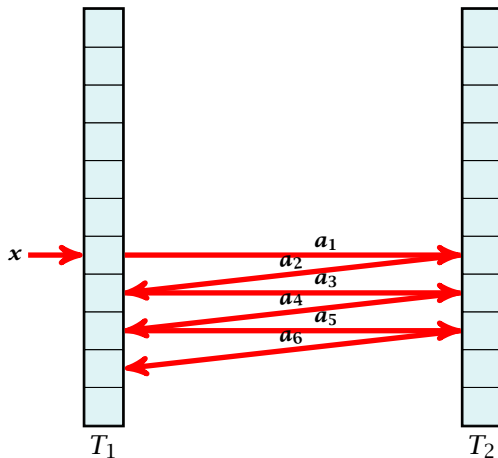
Cuckoo Hashing

Insert:



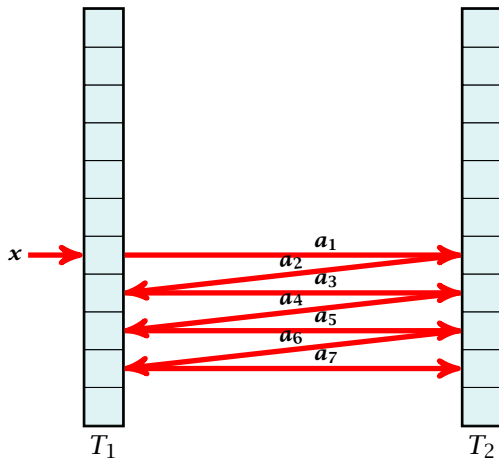
Cuckoo Hashing

Insert:



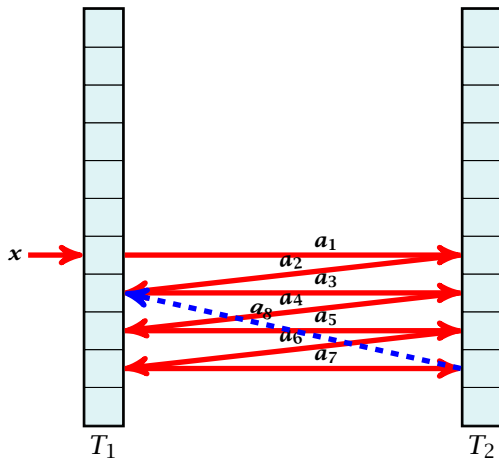
Cuckoo Hashing

Insert:



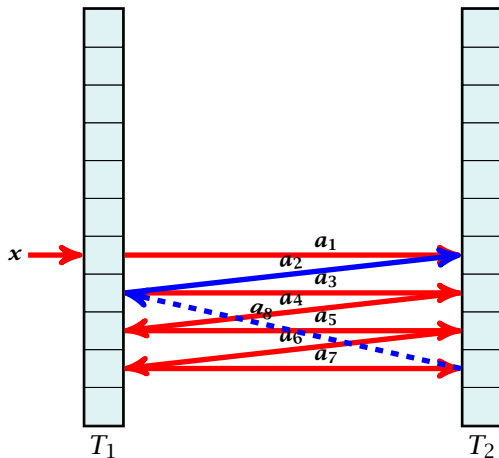
Cuckoo Hashing

Insert:



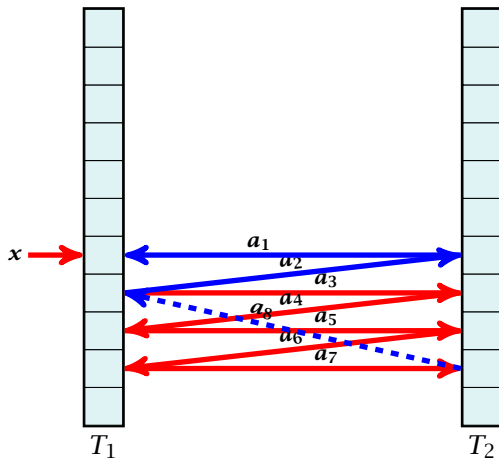
Cuckoo Hashing

Insert:



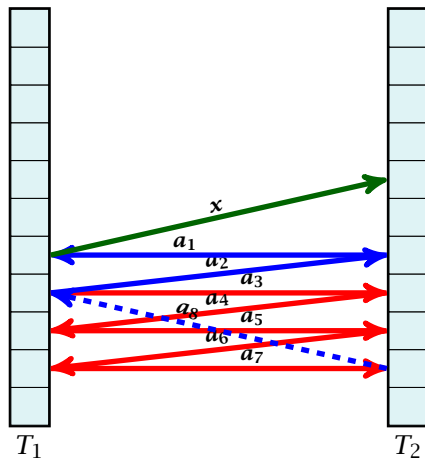
Cuckoo Hashing

Insert:



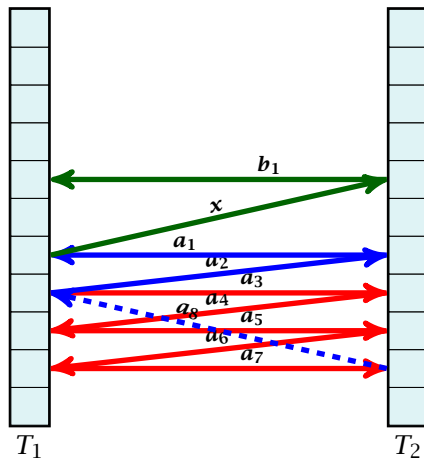
Cuckoo Hashing

Insert:



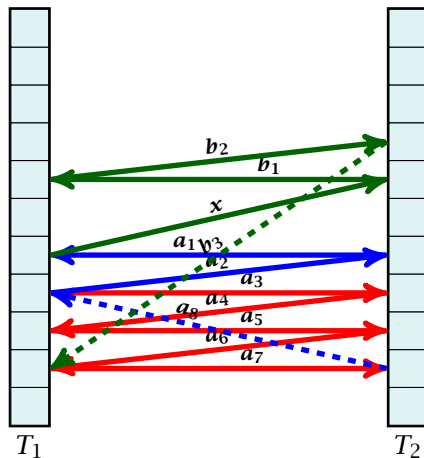
Cuckoo Hashing

Insert:



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Cuckoo Hashing

A cycle-structure is defined by

$\mathcal{C}_1 = (x_1, x_2, \dots, x_n)$ that defines how much the last element x_n "jumps back" to the beginning.

$\mathcal{C}_2 = (y_1, y_2, \dots, y_m)$ that defines how much the last element y_m "jumps back" in the sequence.

• An assignment of positions for the keys in both tables.

• An ordering π_1, \dots, π_n and π'_1, \dots, π'_m of \mathcal{C}_1 and \mathcal{C}_2 .

• A cycle-structure \mathcal{C}_3 defined as $\mathcal{C}_3 =$

Cuckoo Hashing

A cycle-structure is defined by

- ▶ ℓ_a keys $a_1, a_2, \dots, a_{\ell_a}$, $\ell_a \geq 2$,
- ▶ An index $j_a \in \{1, \dots, \ell_a - 1\}$ that defines how much the last item a_{ℓ_a} “jumps back” in the sequence.
- ▶ ℓ_b keys $b_1, b_2, \dots, b_{\ell_b}$. $b \geq 0$.
- ▶ An index $j_b \in \{1, \dots, \ell_a + \ell_b\}$ that defines how much the last item b_{ℓ_b} “jumps back” in the sequence.
- ▶ An assignment of positions for the keys in both tables. Formally we have positions p_1, \dots, p_{ℓ_a} , and p'_1, \dots, p'_{ℓ_b} .
- ▶ The size of a cycle-structure is defined as $\ell_a + \ell_b$.

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Cuckoo Hashing

We say a cycle-structure is **active** for key x if the hash-functions are chosen in such a way that the hash-function results match the pre-defined key-positions.

$$h_1(x) = h_2(x_1) = p_1$$

$$h_2(x_1) = h_1(x_2) = p_2$$

$$\vdots$$

$$\vdots$$

$$\text{if } p_j \text{ is even then } h_1(x_j) = p_{j+1}, \text{ else } h_2(x_j) = p_{j+1}$$

$$h_2(x_j) = h_1(x_{j+1}) = p_{j+1}^2$$

$$\vdots$$

$$\vdots$$

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- ▶ ...
- ▶ if l_a is even then $h_1(a_\ell) = p_{s_a}$, otw. $h_2(a_\ell) = p_{s_a}$
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Cuckoo Hashing

Observation If we end up in an infinite loop there must exist a cycle-structure that is active for x .

Cuckoo Hashing

A cycle-structure is defined **without** knowing the hash-functions.

Whether a cycle-structure is active for key x depends on the hash-functions.

Lemma 30

A given cycle-structure of size s is active for key x with probability at most

$$\left(\frac{\mu}{n}\right)^{2(s+1)},$$

if we use $(\mu, s + 1)$ -independent hash-functions.

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Proof.

All positions are fixed by the cycle-structure. Therefore we ask for the probability of mapping $s + 1$ keys (the a -keys, the b -keys and x) to pre-specified positions in T_1 , **and** to pre-specified positions in T_2 .

The probability is

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- ▶ There are at most s ways to choose ℓ_a . This fixes ℓ_b .
- ▶ There are at most s^2 ways to choose j_a , and j_b .
- ▶ There are at most m^s possibilities to choose the keys a_1, \dots, a_{ℓ_a} and b_1, \dots, b_{ℓ_b} .
- ▶ There are at most n^s choices for choosing the positions p_1, \dots, p_{ℓ_a} and p'_1, \dots, p'_{ℓ_a} .

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Cuckoo Hashing

The number of cycle-structures of size s is small:

- ▶ There are at most s ways to choose ℓ_a . This fixes ℓ_b .
- ▶ There are at most s^2 ways to choose j_a , and j_b .
- ▶ There are at most m^s possibilities to choose the keys a_1, \dots, a_{ℓ_a} and b_1, \dots, b_{ℓ_b} .
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Now assume that the insert operation takes t steps and does not create an infinite loop.

Consider the sequences $x, a_1, a_2, \dots, a_{\ell_a}$ and $x, b_1, b_2, \dots, b_{\ell_b}$ where the a_i 's and b_i 's are defined as before (but for the construction we only use keys examined during the while loop)

If the insert operation takes t steps then

$$t \leq 2\ell_a + 2\ell_b + 2$$

as no key is examined more than twice.

Hence, one of the sequences $x, a_1, a_2, \dots, a_{\ell_a}$ and $x, b_1, b_2, \dots, b_{\ell_b}$ must contain at least $t/4$ keys (either $\ell_a + 1$ or $\ell_b + 1$ must be larger than $t/4$).

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Define a sub-sequence of length ℓ starting with x , as a sequence x_1, \dots, x_ℓ of keys with $x_1 = x$, together with $\ell + 1$ positions p_0, p_1, \dots, p_ℓ from $\{0, \dots, n - 1\}$.

We say a sub-sequence is **right-active** for h_1 and h_2 if

$$h_1(x) = h_1(x_1) = p_0, h_2(x_1) = h_2(x_2) = p_1, \\ h_1(x_2) = h_1(x_3) = p_2, h_2(x_3) = h_2(x_4) = p_3, \dots$$

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Cuckoo Hashing

Observation:

If the insert takes $t \geq 4\ell$ steps there must either be a left-active or a right-active sub-sequence of length ℓ starting with x .

Cuckoo Hashing

The probability that a given sub-sequence is left-active (right-active) is at most

$$\left(\frac{\mu}{n}\right)^{2\ell},$$

if we use (μ, ℓ) -independent hash-functions. This holds since there are ℓ keys whose hash-values (two values per key) have to map to pre-specified positions.

Cuckoo Hashing

The number of sequences is at most $m^{\ell-1} p^{\ell+1}$ as we can choose $\ell - 1$ keys (apart from x) and we can choose $\ell + 1$ positions p_0, \dots, p_ℓ .

The probability that there exists a left-active **or** right-active sequence of length ℓ is at most

$$\begin{aligned} & \Pr[\text{there exists active sequ. of length } \ell] \\ & \leq 2 \cdot m^{\ell-1} \cdot n^{\ell+1} \cdot \left(\frac{\mu}{n}\right)^{2\ell} \\ & \leq 2 \left(\frac{1}{1+\delta}\right)^\ell \end{aligned}$$

Cuckoo Hashing

If the search does not run into an infinite loop the probability that it takes more than 4ℓ steps is at most

$$2\left(\frac{1}{1+\delta}\right)^\ell$$

We choose $\text{maxsteps} = 4(1 + 2 \log m) / \log(1 + \delta)$. Then the probability of terminating the while-loop because of reaching maxsteps is only $\mathcal{O}(\frac{1}{m^2})$ ($\mathcal{O}(1/m^2)$) because of reaching an infinite loop and $1/m^2$ because the search takes maxsteps steps without running into a loop).

Cuckoo Hashing

The expected time for an insert under the condition that `maxsteps` is not reached is

$$\sum_{\ell \geq 0} \Pr[\text{search takes at least } \ell \text{ steps} \mid \text{iteration successful}] \\ \leq \sum_{\ell \geq 0} 8 \left(\frac{1}{1 + \delta} \right)^\ell = \mathcal{O}(1) .$$

More generally, the above expression gives a bound on the cost in the successful iteration of an insert-operation (there is exactly one successful iteration).

An iteration that is not successful induces cost $\mathcal{O}(m)$ for doing a complete rehash.

Cuckoo Hashing

The expected number of unsuccessful operations is $\mathcal{O}(\frac{1}{m^2})$.

Hence, the expected cost in unsuccessful iterations is only $\mathcal{O}(\frac{1}{m})$.

Hence, the total expected cost for an insert-operation is constant.

Cuckoo Hashing

What kind of hash-functions do we need?

Since maxsteps is $\Theta(\log m)$ it is sufficient to have $(\mu, \Theta(\log m))$ -independent hash-functions.

Cuckoo Hashing

How do we make sure that $n \geq \mu^2(1 + \delta)m$?

- ▶ Let $\alpha := 1/(\mu^2(1 + \delta))$.
- ▶ Keep track of the number of elements in the table. Whenever $m \geq \alpha n$ we double n and do a complete re-hash (table-expand).
- ▶ Whenever m drops below $\frac{\alpha}{4}n$ we divide n by 2 and do a rehash (table-shrink).
- ▶ Note that right after a change in table-size we have $m = \frac{\alpha}{2}n$. In order for a table-expand to occur at least $\frac{\alpha}{2}n$ insertions are required. Similar, for a table-shrink at least $\frac{\alpha}{4}$ deletions must occur.
- ▶ Therefore we can amortize the rehash cost after a change in table-size against the cost for insertions and deletions.

Definition 31

Let $d \in \mathbb{N}$; $q \geq n$ be a prime; and let $\vec{a} \in \{0, \dots, q-1\}^{d+1}$. Define for $x \in \{0, \dots, q\}$

$$h_{\vec{a}}(x) := \left(\sum_{i=0}^d a_i x^i \bmod q \right) \bmod n .$$

Let $\mathcal{H}_n^d := \{h_{\vec{a}} \mid \vec{a} \in \{0, \dots, q\}^{d+1}\}$. The class \mathcal{H}_n^d is $(2, d+1)$ -independent.

For the coefficients $\bar{a} \in \{0, \dots, q-1\}^{d+1}$ let $f_{\bar{a}}$ denote the polynomial

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The polynomial is defined by $d+1$ distinct points.

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Fix $\ell \leq d + 1$; let $x_1, \dots, x_\ell \in \{0, \dots, q - 1\}$ be keys, and let t_1, \dots, t_ℓ denote the corresponding hash-function values.

Let $A^\ell = \{h_{\bar{a}} \in \mathcal{H} \mid h_{\bar{a}}(x_i) = t_i \text{ for all } i \in \{1, \dots, \ell\}\}$

Then

$$h_{\bar{a}} \in A^\ell \Leftrightarrow h_{\bar{a}} = f_{\bar{a}} \bmod n \text{ and}$$

$$f_{\bar{a}}(x_i) \in \{t_i + \alpha \cdot n \mid \alpha \in \{0, \dots, \lceil \frac{q}{n} \rceil - 1\}\}$$

Therefore I have

$$|B_1| \cdot \dots \cdot |B_\ell| \cdot q^{d-\ell+1} \leq \lceil \frac{q}{n} \rceil^\ell \cdot q^{d-\ell+1}$$

possibilities to choose \bar{a} such that $h_{\bar{a}} \in A_\ell$.

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Therefore the probability of choosing $h_{\bar{a}}$ from A_ℓ is only

$$\frac{\lceil \frac{q}{n} \rceil^\ell \cdot q^{d-\ell+1}}{q^{d+1}} \leq \left(\frac{2}{n}\right)^\ell$$

8 Priority Queues

A **Priority Queue** S is a dynamic set data structure that supports the following operations:

- ▶ $S.\text{build}(x_1, \dots, x_n)$: Creates a data-structure that contains just the elements x_1, \dots, x_n .
- ▶ $S.\text{insert}(x)$: Adds element x to the data-structure.
- ▶ Element $S.\text{minimum}()$: Returns an element $x \in S$ with minimum key-value $\text{key}[x]$.
- ▶ $S.\text{delete-min}()$: Deletes the element with minimum key-value from S and returns it.
- ▶ Boolean $S.\text{empty}()$: Returns true if the data-structure is empty and false otherwise.

Sometimes we also have

- ▶ $S.\text{merge}(S')$: $S := S \cup S'$; $S' := \emptyset$.

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- ▶ **Element S .minimum():** Returns an element $x \in S$ with minimum key-value $\text{key}[x]$.
- ▶ **S .delete-min():** Deletes the element with minimum key-value from S and returns it.
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- ▶ **S .merge(S'):** $S := S \cup S'$; $S' := \emptyset$.

8 Priority Queues

A **Priority Queue** S is a dynamic set data structure that supports the following operations:

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8 Priority Queues

An **addressable Priority Queue** also supports:

- ▶ **Handle $S.insert(x)$** : Adds element x to the data-structure, and returns a **handle** to the object for future reference.
- ▶ **$S.delete(h)$** : Deletes element specified through handle h .
- ▶ **$S.decrease-key(h, k)$** : Decreases the key of the element specified by handle h to k . Assumes that the key is at least k before the operation.

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Dijkstra's Shortest Path Algorithm

Algorithm 17 Shortest-Path($G = (V, E, d), s \in V$)

```
1: Input: weighted graph  $G = (V, E, d)$ ; start vertex  $s$ ;  
2: Output: key-field of every node contains distance from  $s$ ;  
3:  $S.build()$ ; // build empty priority queue  
4: for all  $v \in V \setminus \{s\}$  do  
5:      $v.key \leftarrow \infty$ ;  
6:      $h_v \leftarrow S.insert(v)$ ;  
7:  $s.key \leftarrow 0$ ;  $S.insert(s)$ ;  
8: while  $S.empty() = \text{false}$  do  
9:      $v \leftarrow S.delete-min()$ ;  
10:    for all  $x \in V$  s.t.  $(v, x) \in E$  do  
11:        if  $x.key > v.key + d(v, x)$  then  
12:             $S.decrease-key(h_x, v.key + d(v, x))$ ;  
13:             $x.key \leftarrow v.key + d(v, x)$ ;
```

Prim's Minimum Spanning Tree Algorithm

Algorithm 18 Prim-MST($G = (V, E, d), s \in V$)

```
1: Input: weighted graph  $G = (V, E, d)$ ; start vertex  $s$ ;  
2: Output: pred-fields encode MST;  
3:  $S.build()$ ; // build empty priority queue  
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5:      $v.key \leftarrow \infty$ ;  
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14:             $x.pred \leftarrow v$ ;
```

Analysis of Dijkstra and Prim

Both algorithms require:

- ▶ 1 build() operation
- ▶ $|V|$ insert() operations
- ▶ $|V|$ delete-min() operations
- ▶ $|V|$ is-empty() operations
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How good a running time can we obtain?

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8 Priority Queues

Operation	Binary Heap	BST	Binomial Heap	Fibonacci Heap*
build	n	$n \log n$	$n \log n$	n
minimum	1	$\log n$	$\log n$	1
is-empty	1	1	1	1
insert	$\log n$	$\log n$	$\log n$	1
delete	$\log n^{**}$	$\log n$	$\log n$	$\log n$
delete-min	$\log n$	$\log n$	$\log n$	$\log n$
decrease-key	$\log n$	$\log n$	$\log n$	1
merge	n	$n \log n$	$\log n$	1

Note that most applications use `build()` only to create an empty heap which then costs time 1.

The standard version of binary heaps is not addressable, and hence does not support a delete operation.

Fibonacci heaps only give an **amortized** guarantee.

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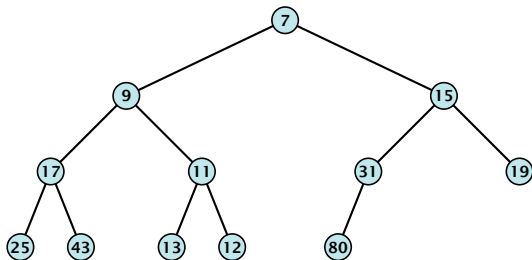
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8 Priority Queues

Using Binary Heaps, Prim and Dijkstra run in time $\mathcal{O}((|V| + |E|) \log |V|)$.

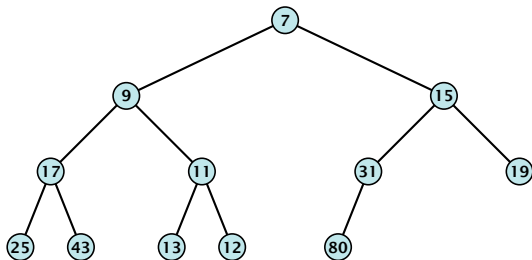
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8.1 Binary Heaps



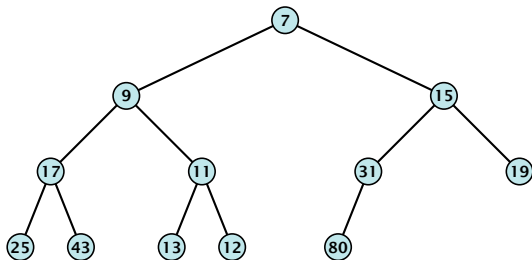
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- ▶ Nearly complete binary tree; only the last level is not full, and this one is filled from left to right.
- ▶ **Heap property:** A node's key is not larger than the key of one of its children.



Binary Heaps

Operations:

- ▶ `minimum()`: return the root-element. Time $\mathcal{O}(1)$.
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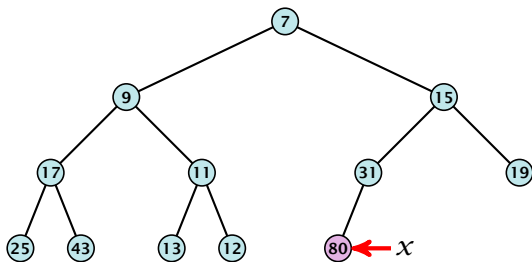
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- ▶ We can compute the predecessor of x (last element when x is deleted) in time $\mathcal{O}(\log n)$.



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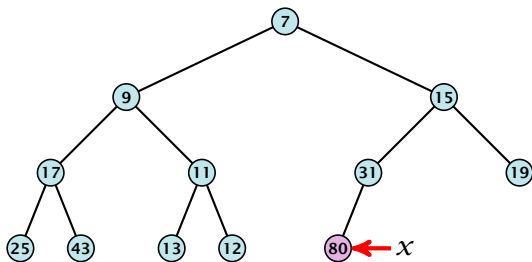
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go up until the last edge used was a right edge.

go left; go right until you reach a leaf

if you hit the root on the way up, go to the rightmost element



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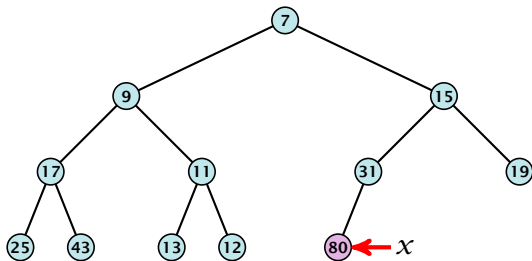
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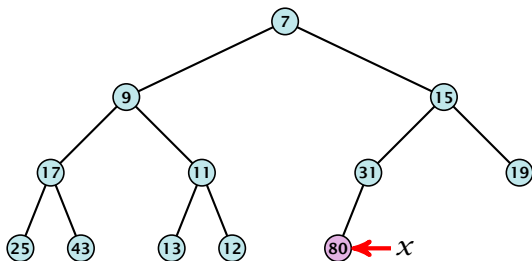
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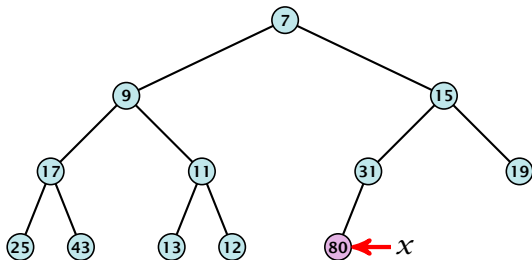
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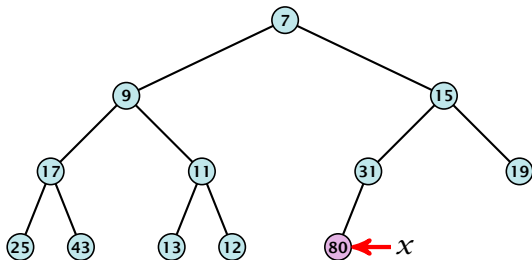
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go up until the last edge used was a left edge.

go right; go left until you reach a null-pointer.

if you hit the root on the way up, go to the leftmost element;

insert a new element as a left child;



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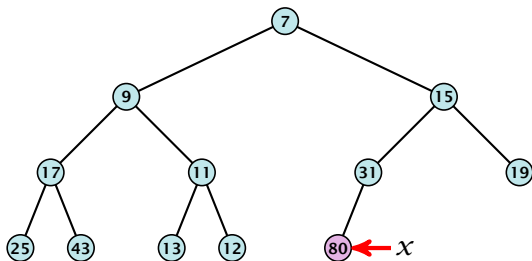
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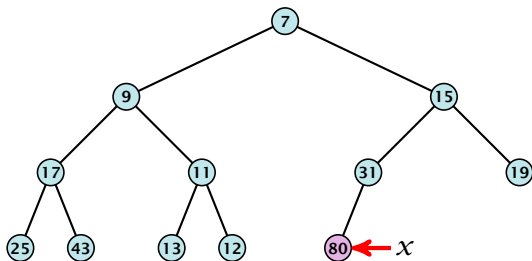
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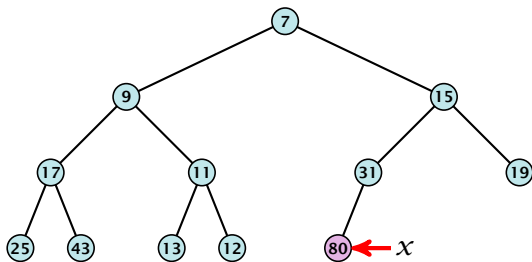
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Insert

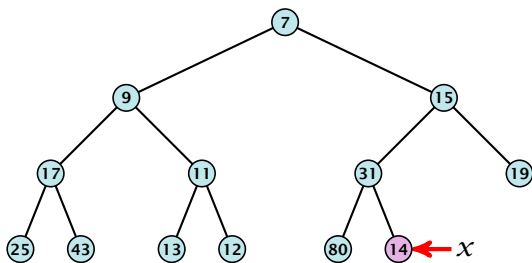
1. Insert element at successor of x .
2. Exchange with parent until heap property is fulfilled.



Note that an exchange can either be done by moving the data or by changing pointers. The latter method leads to an addressable priority queue.

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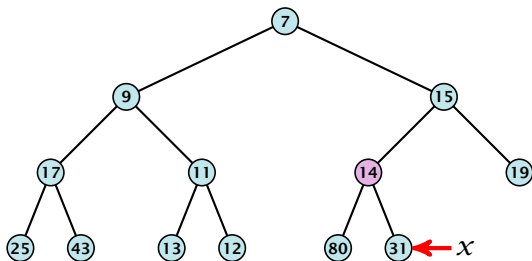
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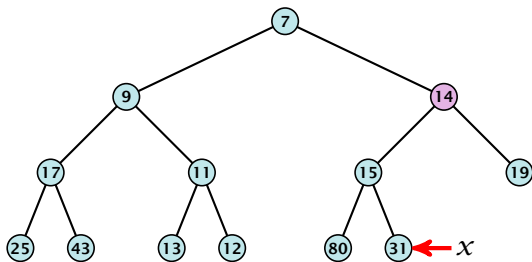
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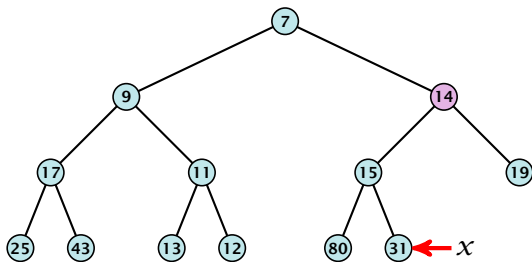
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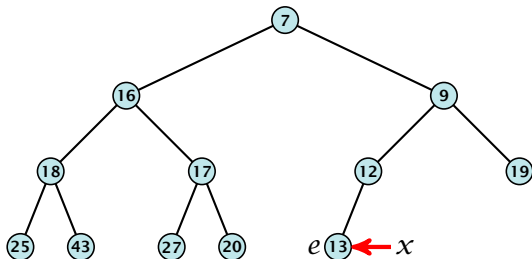
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Delete

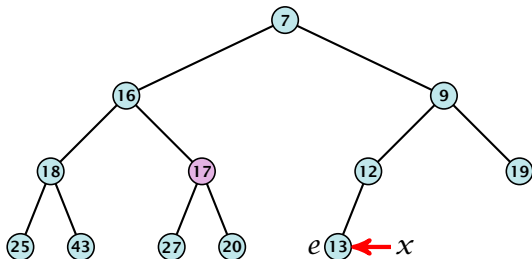
1. Exchange the element to be deleted with the element e pointed to by x .
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At its new position e may either travel up or down in the tree (but not both directions).

Delete

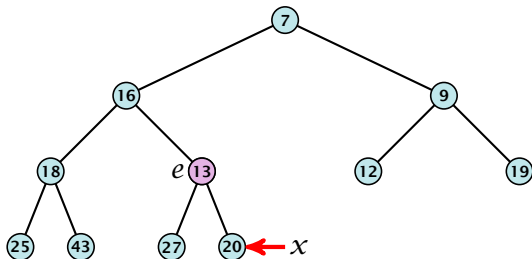
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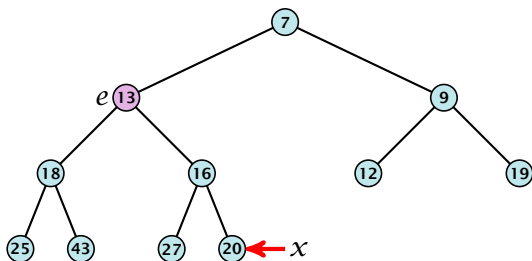
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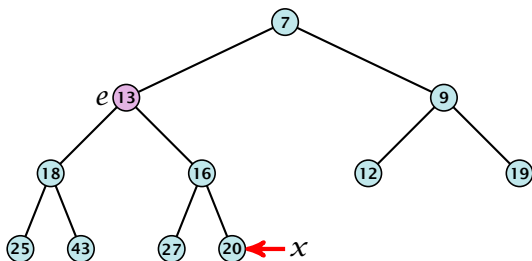
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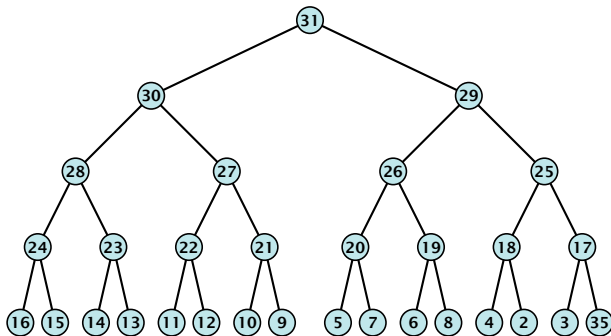
Binary Heaps

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Build Heap

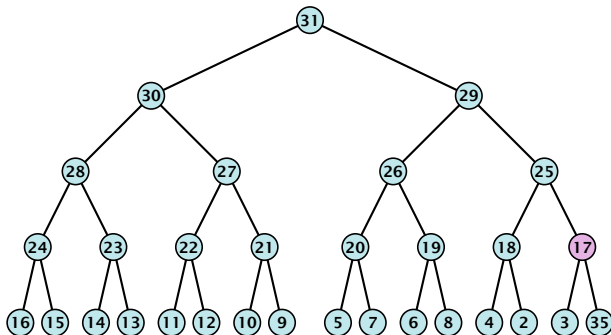
We can build a heap in linear time:



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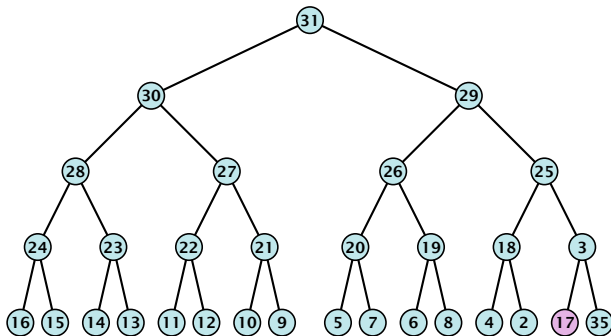
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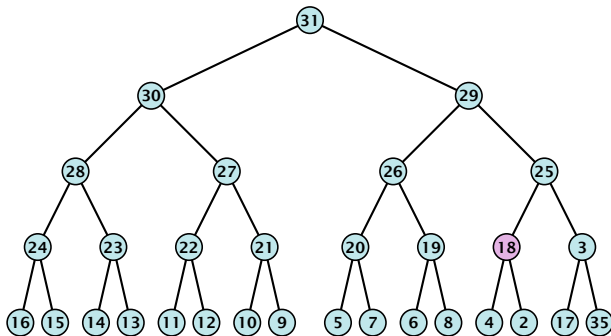
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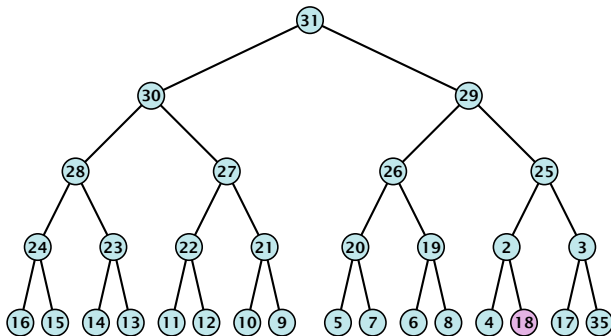
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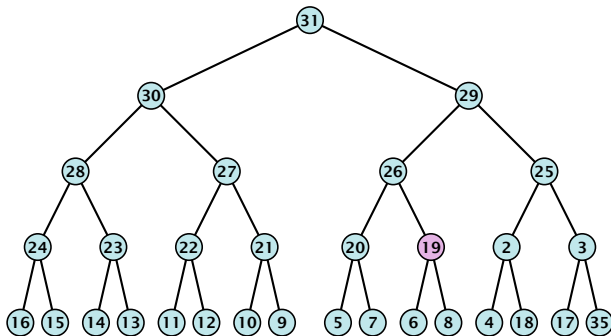
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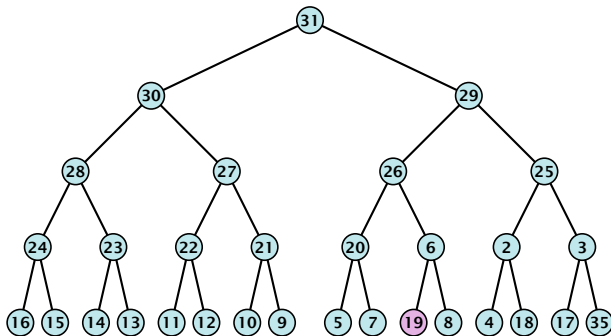
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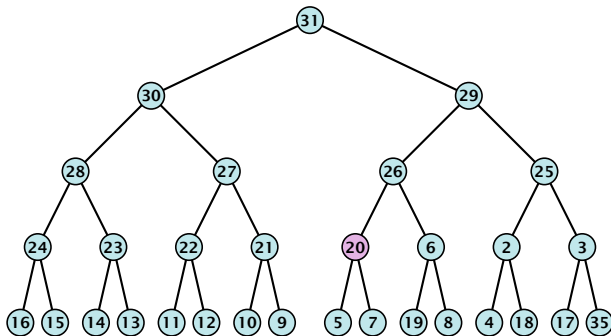
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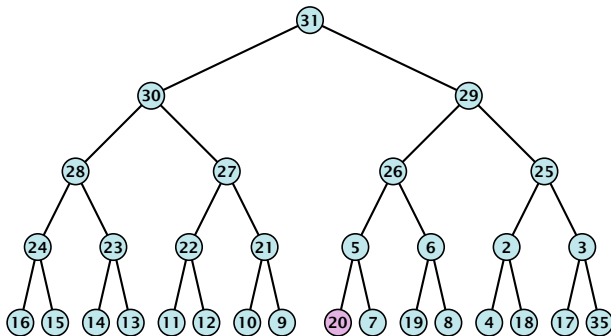
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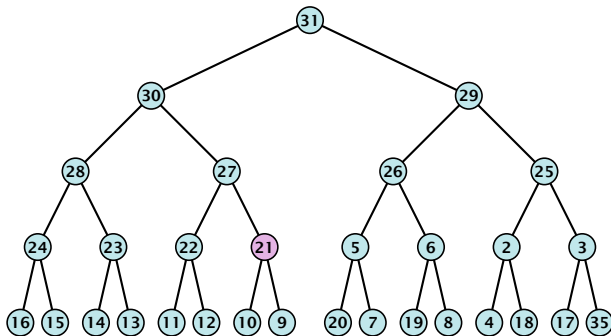
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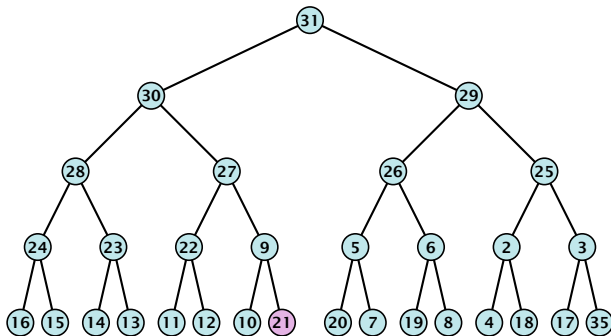
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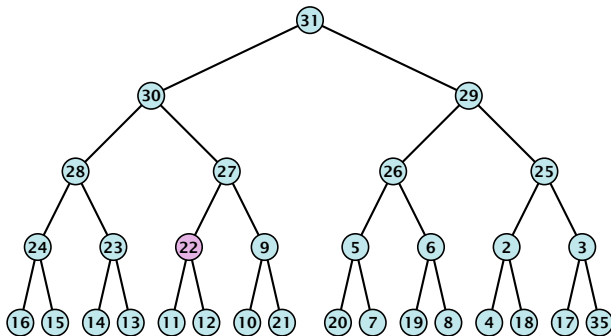
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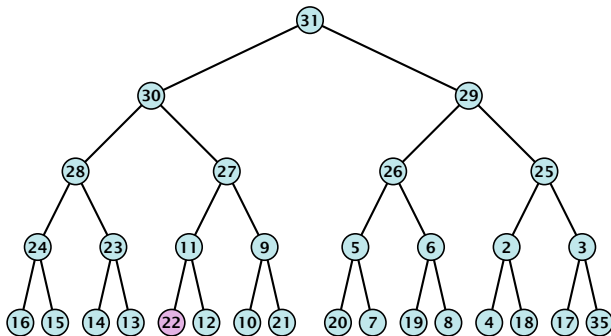
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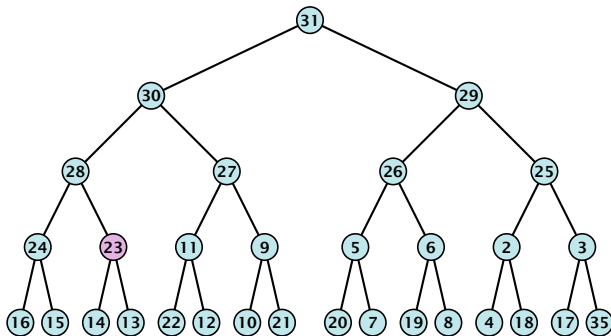
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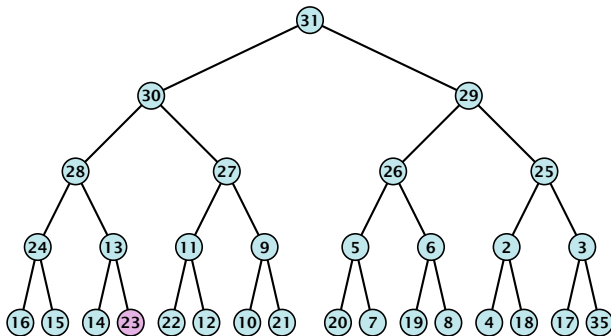
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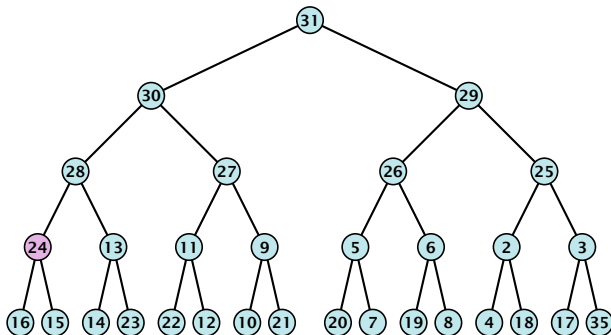
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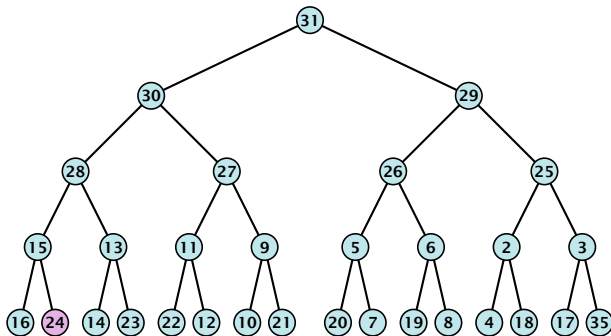
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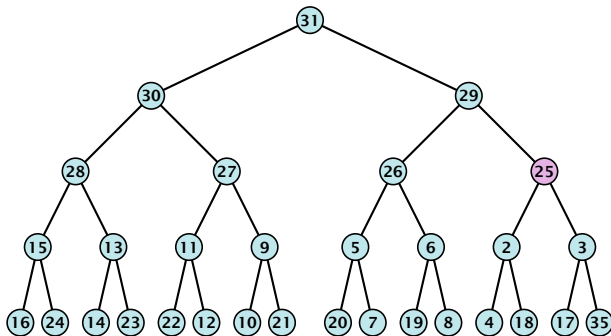
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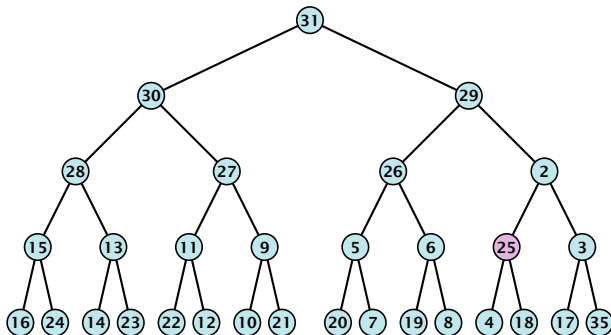
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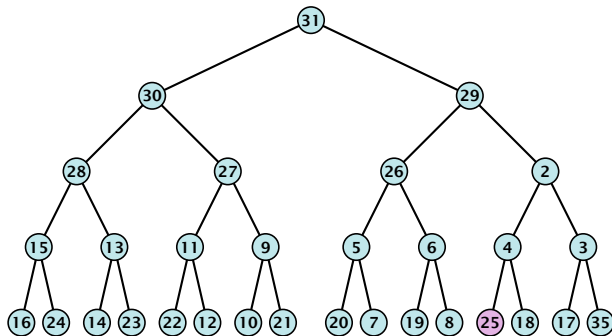
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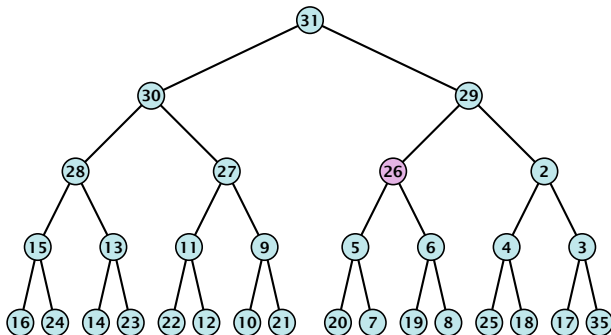
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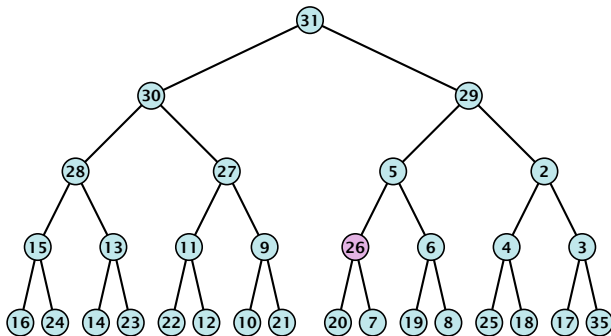
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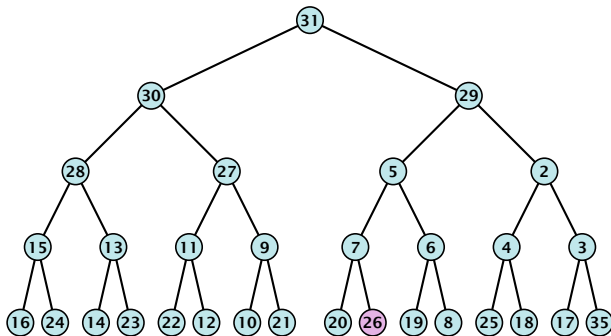
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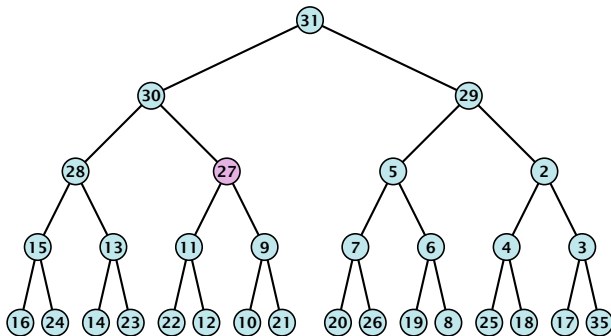
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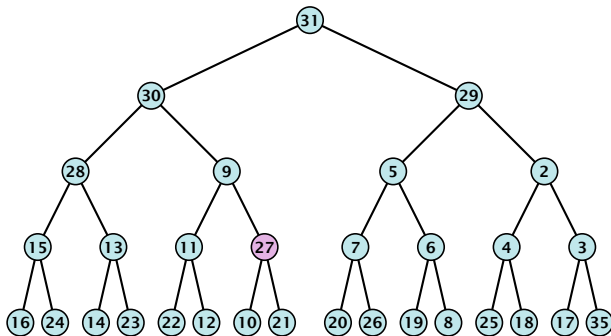
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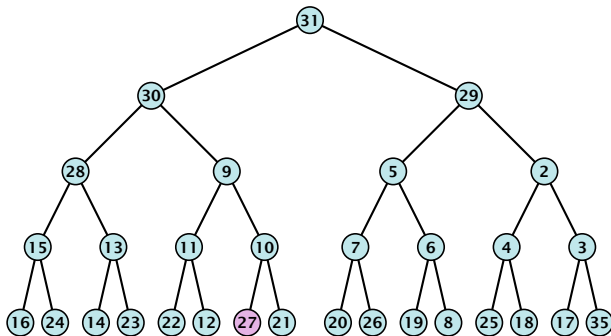
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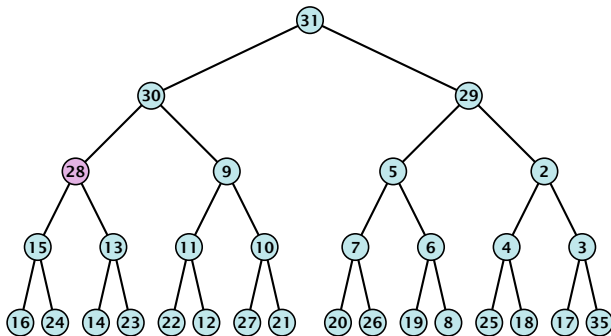
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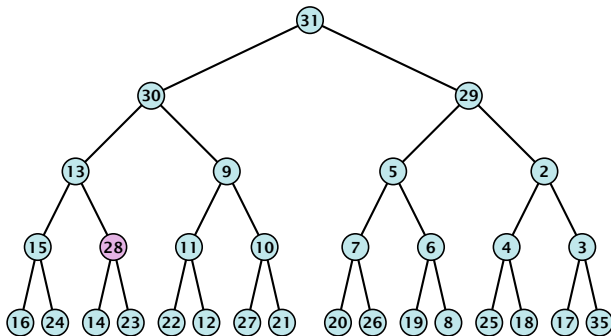
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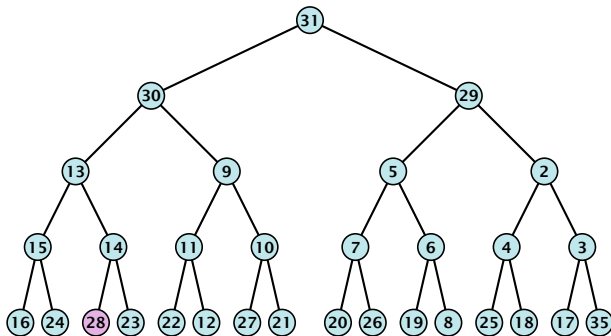
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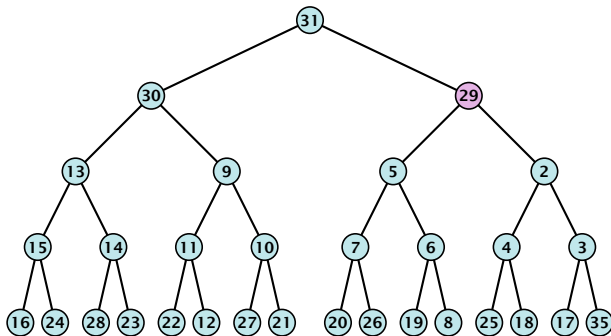
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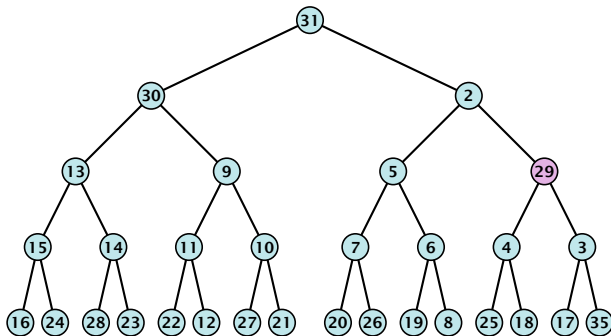
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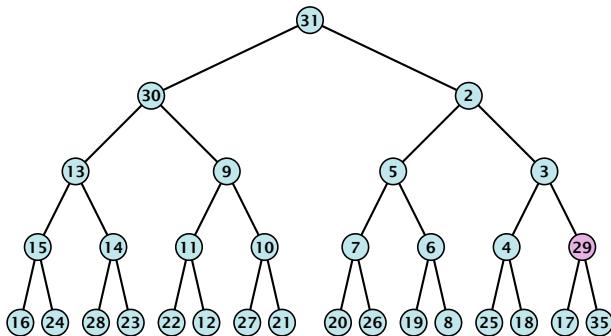
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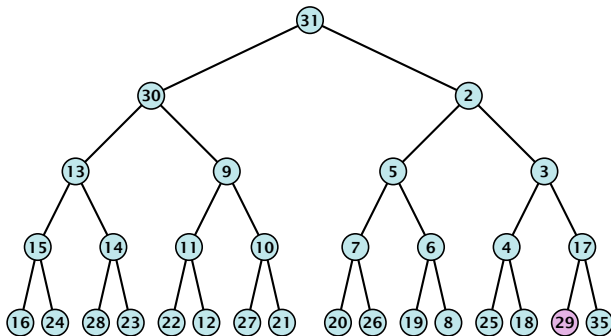
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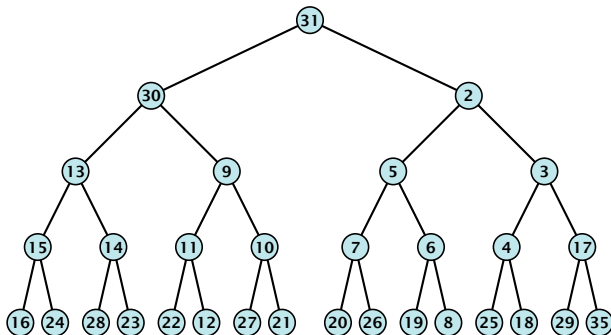
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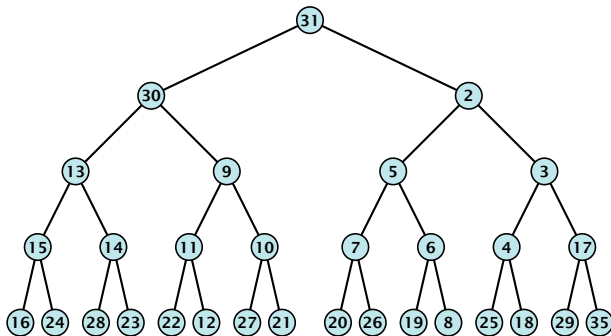
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Binary Heaps

Operations:

- ▶ **minimum()**: Return the root-element. Time $\mathcal{O}(1)$.
- ▶ **is-empty()**: Check whether root-pointer is null. Time $\mathcal{O}(1)$.
- ▶ **insert(k)**: Insert at x and bubble up. Time $\mathcal{O}(\log n)$.
- ▶ **delete(h)**: Swap with x and bubble up or sift-down. Time $\mathcal{O}(\log n)$.
- ▶ **build(x_1, \dots, x_n)**: Insert elements arbitrarily; then do sift-down operations starting with the lowest layer in the tree. Time $\mathcal{O}(n)$.

Binary Heaps

The standard implementation of binary heaps is via arrays. Let $A[0, \dots, n - 1]$ be an array

- ▶ The parent of i -th element is at position $\lfloor \frac{i-1}{2} \rfloor$.
- ▶ The left child of i -th element is at position $2i + 1$.
- ▶ The right child of i -th element is at position $2i + 2$.

Finding the successor of x is much easier than in the description on the previous slide. Simply increase or decrease x .

The resulting binary heap is not addressable. The elements don't maintain their positions and therefore there are not stable handles.

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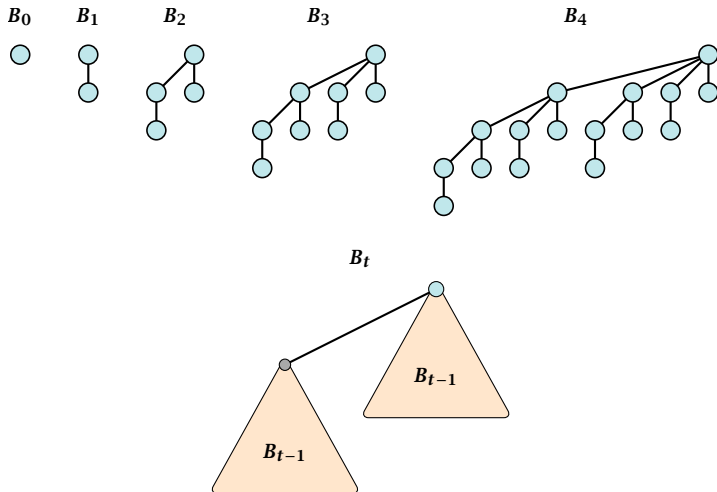
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8.2 Binomial Heaps

Operation	Binary Heap	BST	Binomial Heap	Fibonacci Heap*
build	n	$n \log n$	$n \log n$	n
minimum	1	$\log n$	$\log n$	1
is-empty	1	1	1	1
insert	$\log n$	$\log n$	$\log n$	1
delete	$\log n^{**}$	$\log n$	$\log n$	$\log n$
delete-min	$\log n$	$\log n$	$\log n$	$\log n$
decrease-key	$\log n$	$\log n$	$\log n$	1
merge	n	$n \log n$	$\log n$	1

Binomial Trees



Properties of Binomial Trees

- ▶ B_k has 2^k nodes.
- ▶ B_k has height k .
- ▶ The root of B_k has degree k .
- ▶ B_k has $\binom{k}{\ell}$ nodes on level ℓ .
- ▶ Deleting the root of B_k gives trees B_0, B_1, \dots, B_{k-1} .

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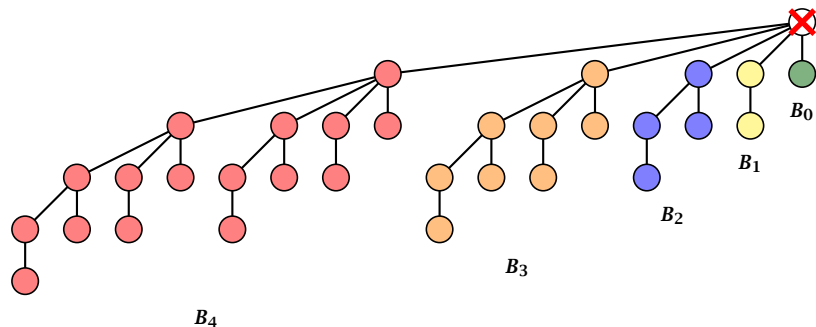
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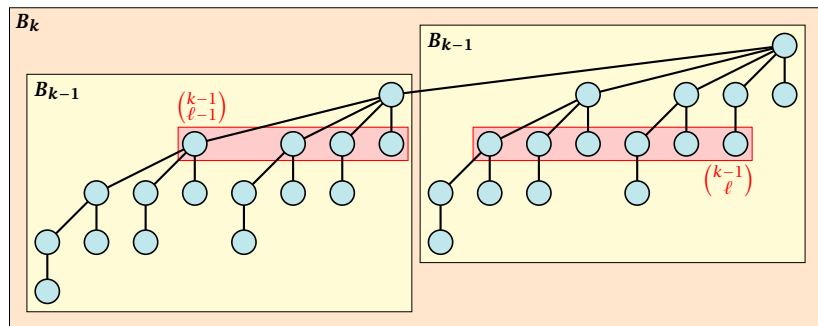
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Binomial Trees



Deleting the root of B_5 leaves sub-trees B_4 , B_3 , B_2 , and B_1 .

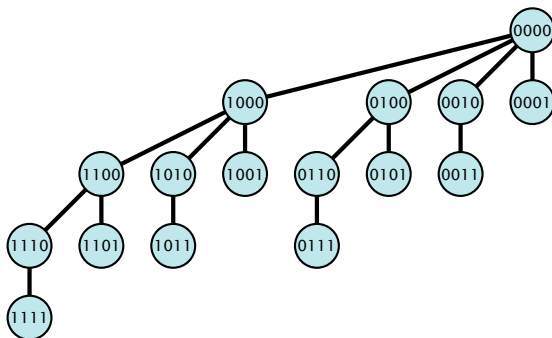
Binomial Trees



The number of nodes on level ℓ in tree B_k is therefore

$$\binom{k-1}{\ell-1} + \binom{k-1}{\ell} = \binom{k}{\ell}$$

Binomial Trees

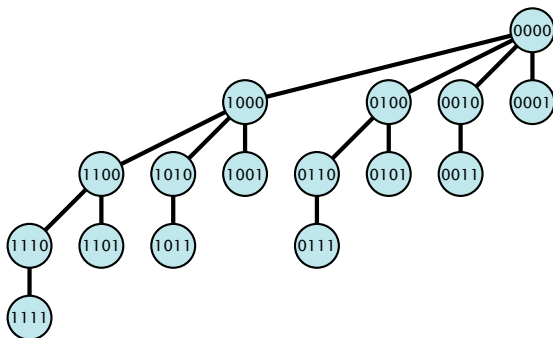


The binomial tree B_k is a sub-graph of the hypercube H_k .

The parent of a node with label b_n, \dots, b_1, b_0 is obtained by setting the least significant 1-bit to 0.

The ℓ -th level contains nodes that have ℓ 1's in their label.

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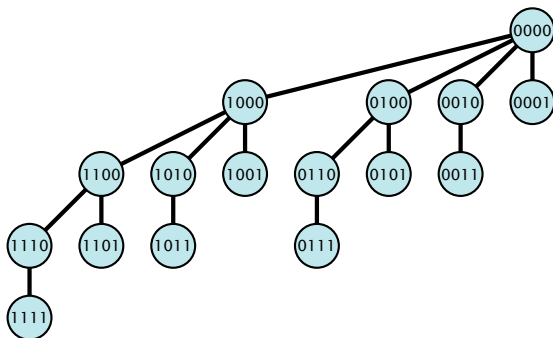


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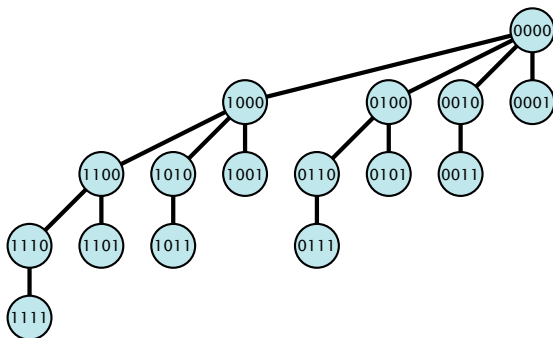


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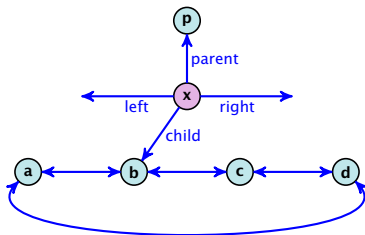
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8.2 Binomial Heaps

How do we implement trees with non-constant degree?

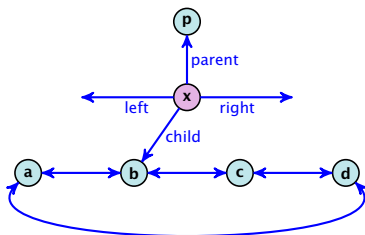
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- ▶ Pointers $x.left$ and $x.right$ point to the left and right sibling of x (if x does not have children then $x.left = x.right = x$).



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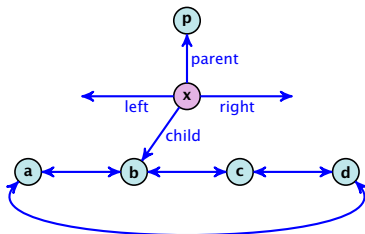
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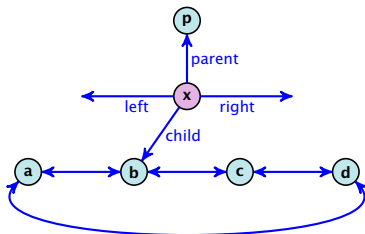
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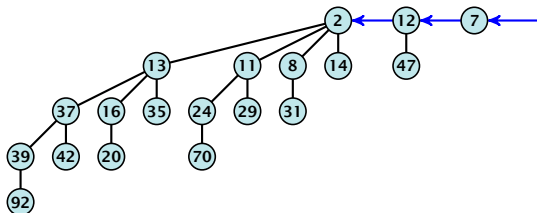
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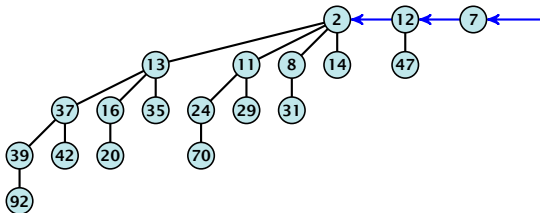


In a binomial heap the keys are arranged in a collection of binomial trees.

Every tree fulfills the heap-property

There is at most one tree for every dimension/order. For example the above heap contains trees B_0 , B_1 , and B_4 .

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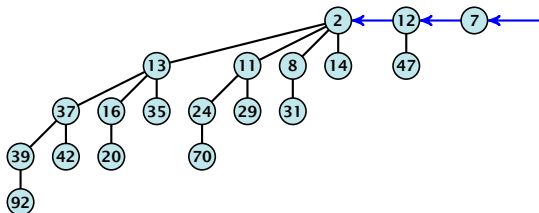


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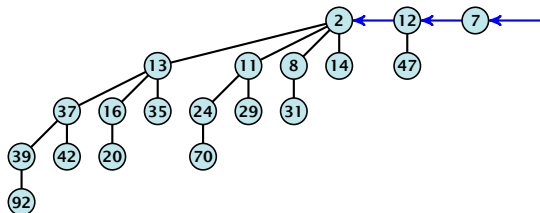


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Binomial Heap: Merge

Given the number n of keys to be stored in a binomial heap we can deduce the binomial trees that will be contained in the collection.

Let $B_{k_1}, B_{k_2}, B_{k_3}, k_i < k_{i+1}$ denote the binomial trees in the collection and recall that every tree may be contained at most once.

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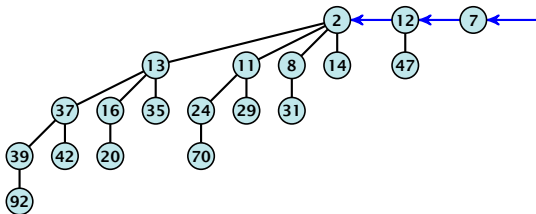
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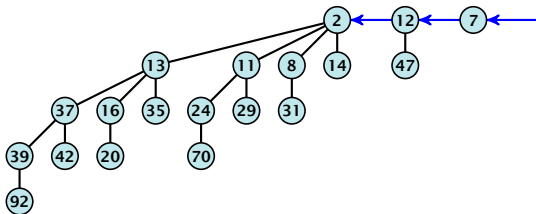
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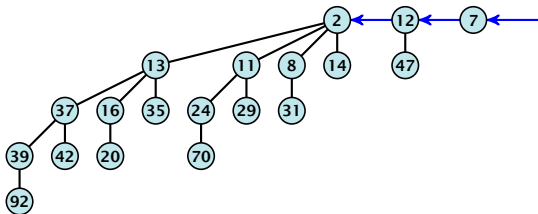
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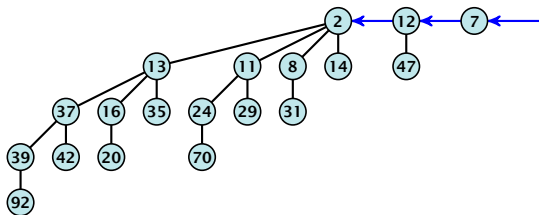
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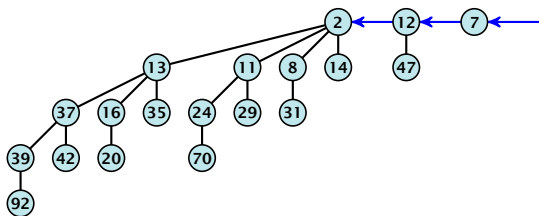
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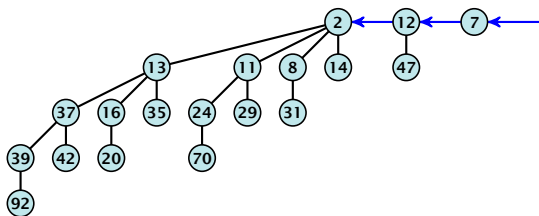
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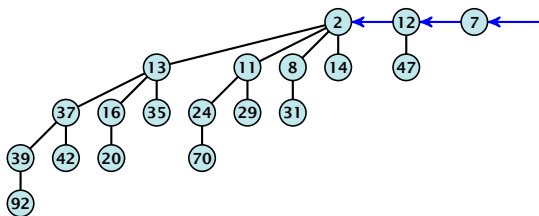
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The merge-operation is instrumental for binomial heaps.

A merge is easy if we have two heaps with different binomial trees. We can simply merge the tree-lists.

Otherwise, we cannot do this because the merged heap is not allowed to contain two trees of the same order.

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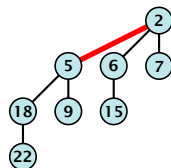
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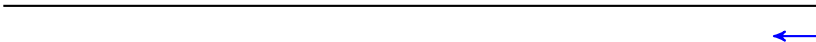
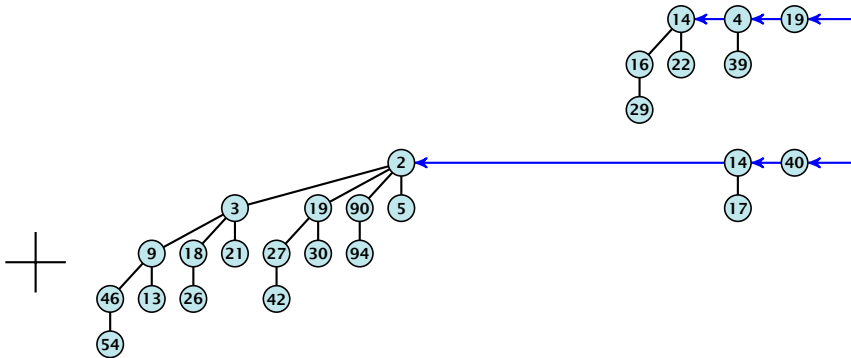
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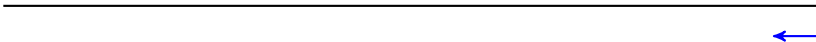
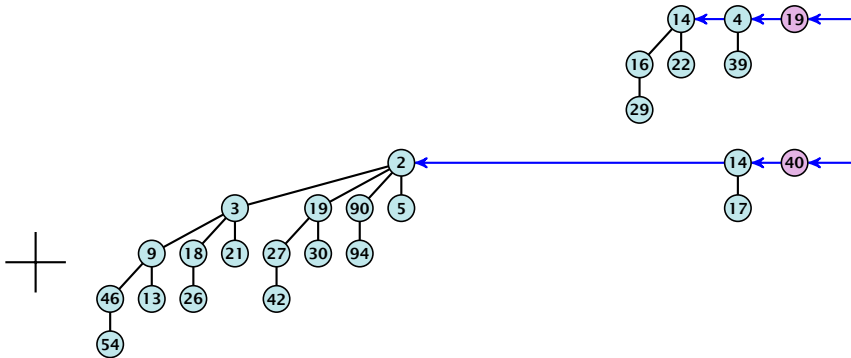
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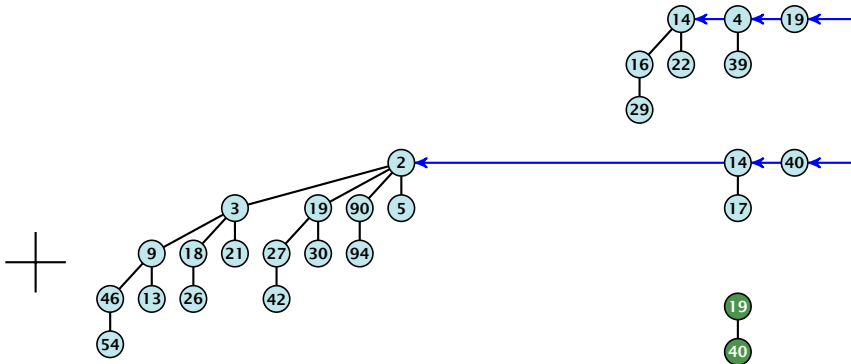
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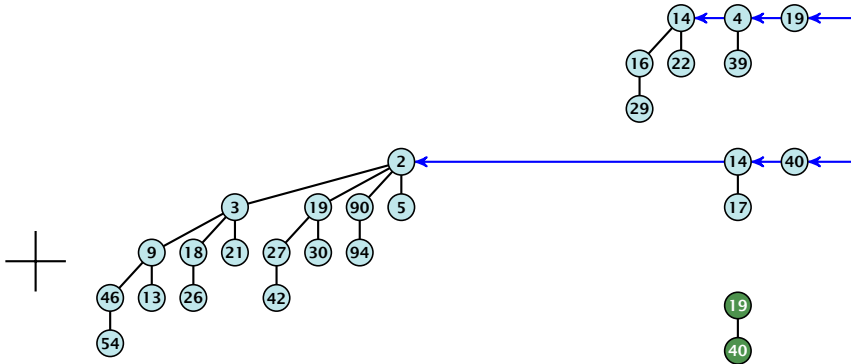
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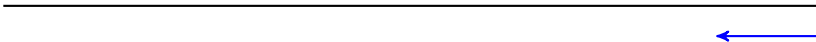
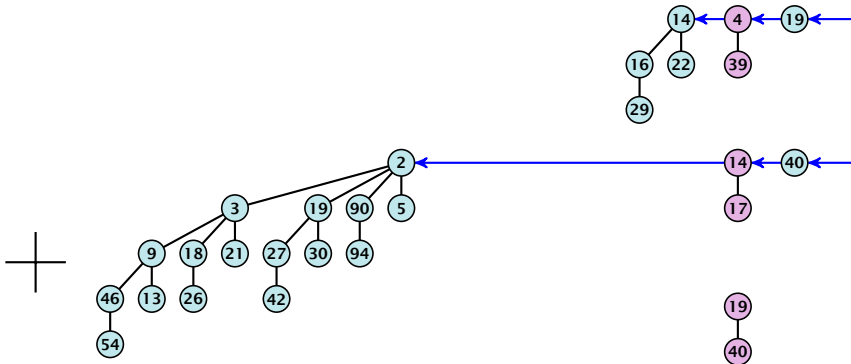


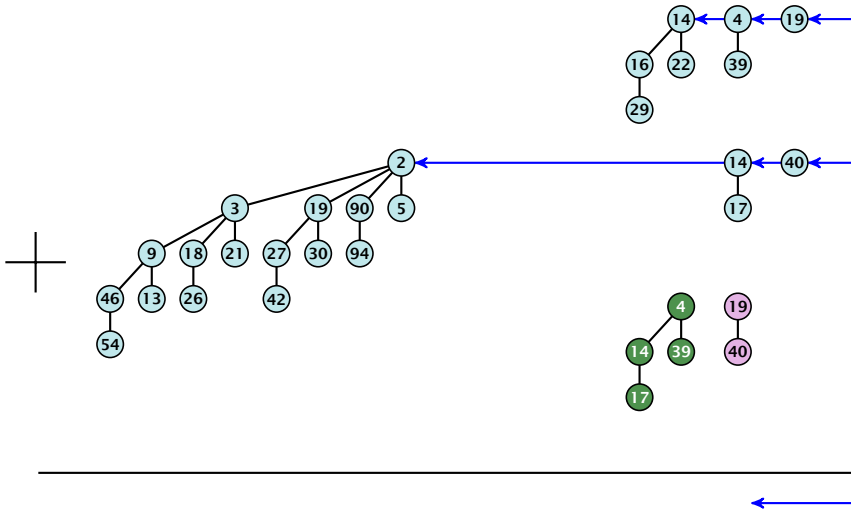


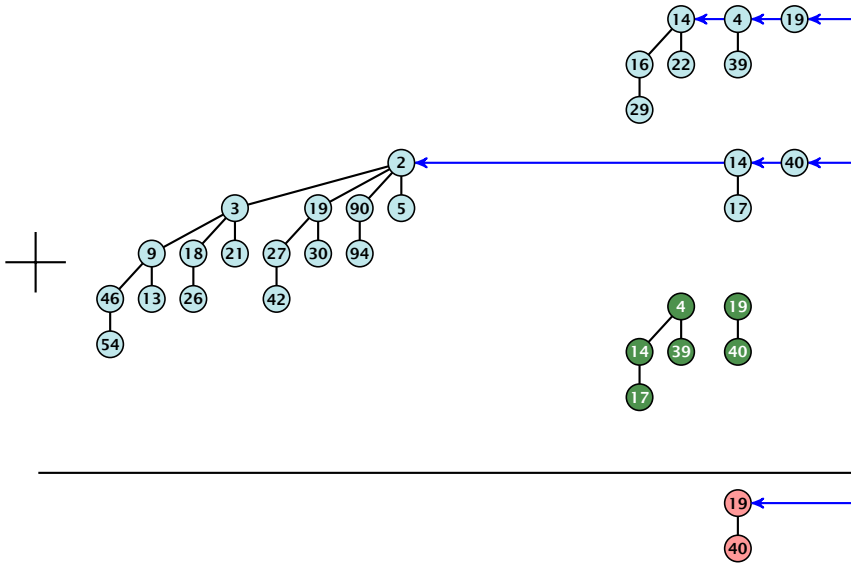


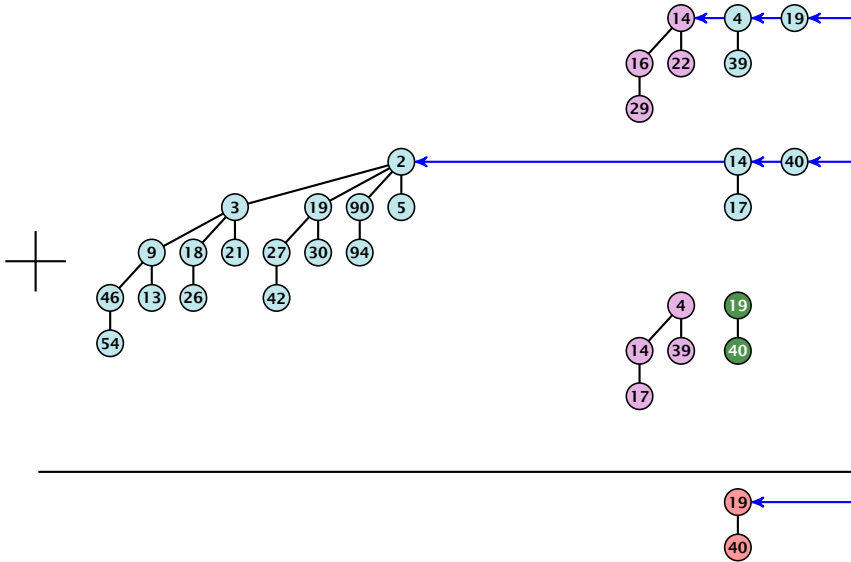


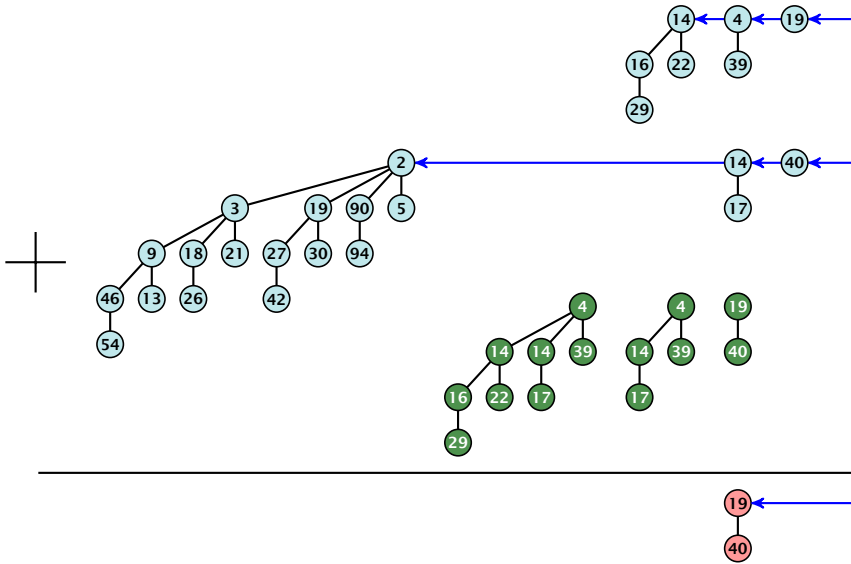


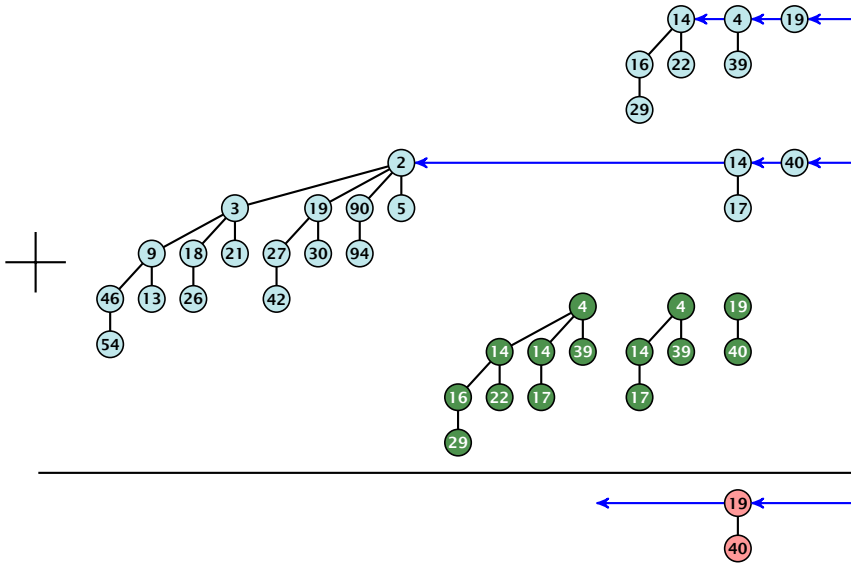


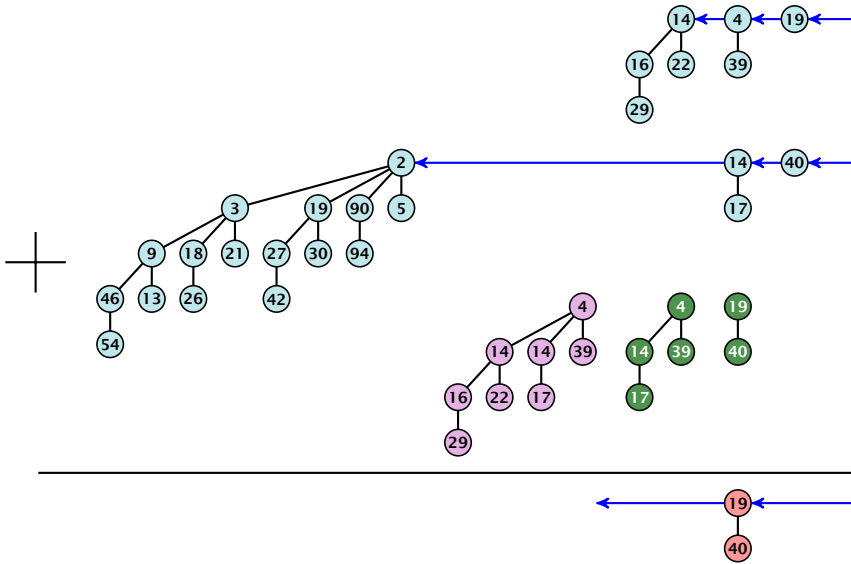


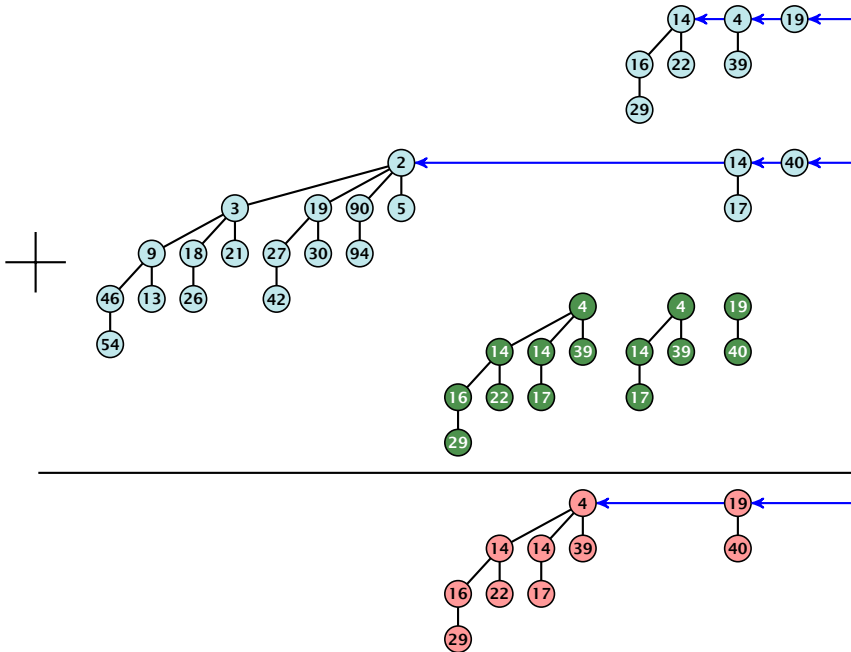


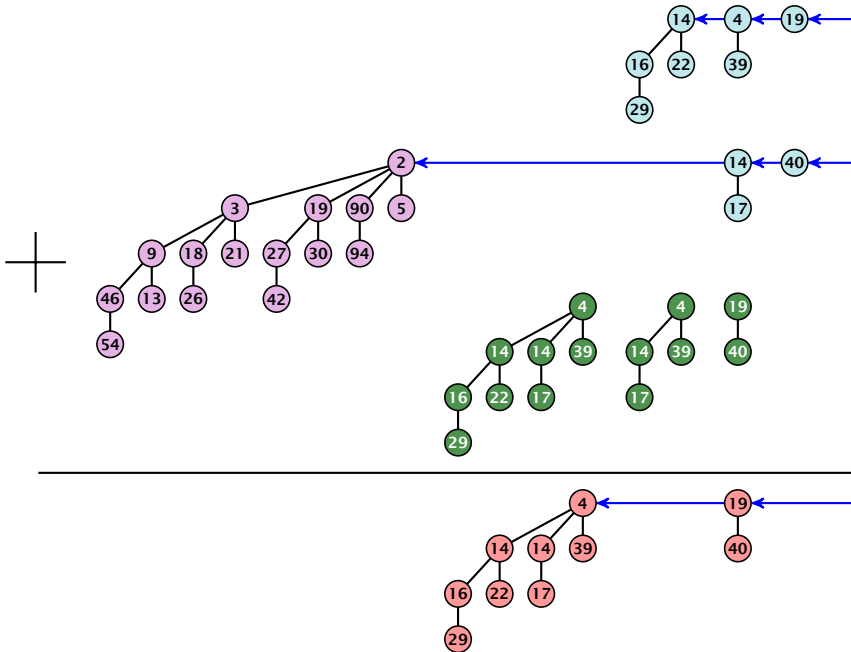


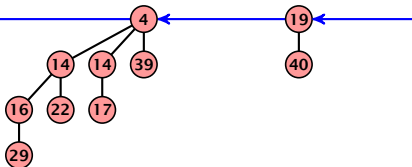
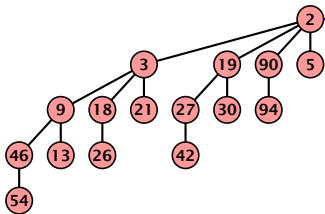
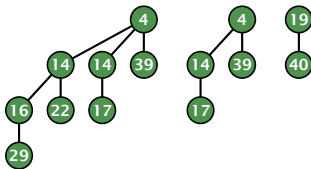
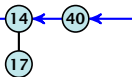
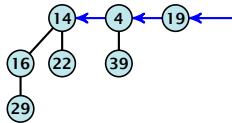
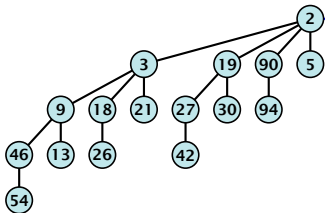


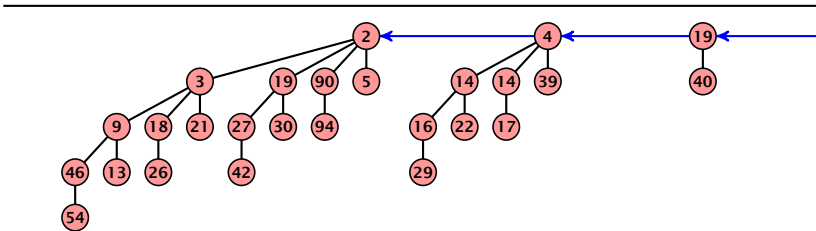
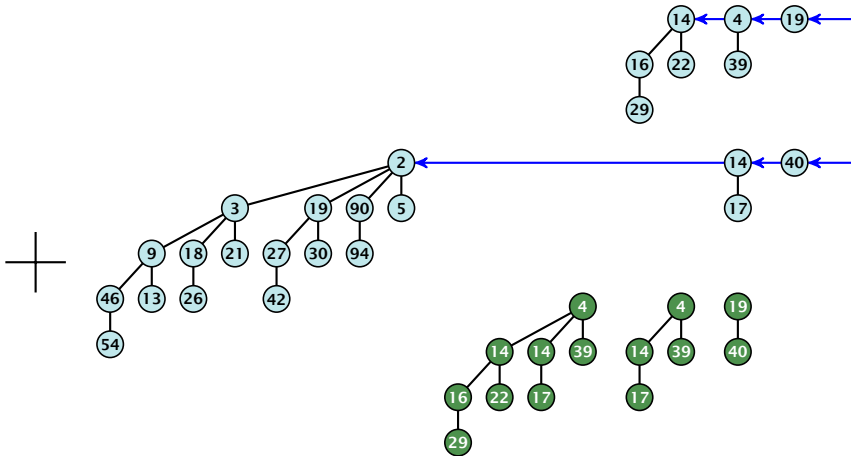












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- ▶ Analogous to binary addition.
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Suppose you are given a sequence of operations (starting with an empty data-structure) that operate on at most n elements, and let k_i denote the number of occurrences of $op_i()$ within this sequence. Then the actual running time must be at most $\sum_i k_i t_i(n)$.

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Potential Method

Introduce a potential for the data structure.

Let $\Phi(D)$ be the potential of the data structure D .

Let c_i be the cost of the i -th operation.

$$c_i = c_i + \Phi(D_i) - \Phi(D_{i-1})$$

Summing up $\Phi(D_i) - \Phi(D_{i-1})$

Then

$$\sum_{i=1}^k c_i \leq \sum_{i=1}^k c_i + \Phi(D_k) - \Phi(D_0) = \sum_{i=1}^k \hat{c}_i$$

This means the amortized costs can be used to derive a bound on the total cost.

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Stack

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- ▶ $S.$ **pop**()
- ▶ $S.$ **multipop**(k): removes k items from the stack. If the stack currently contains less than k items it empties the stack.

Actual cost:

- ▶ $S.$ **push**(): cost 1.
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Use potential function $\Phi(S) = \text{number of elements on the stack}$.

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• $S.push()$: cost

$$C_{push} = C_{push} + \Delta\Phi = 1 + 1 = 2$$

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$$C_{pop} = C_{pop} + \Delta\Phi = 1 - 1 = 0$$

• $S.empty() \rightarrow \text{bool}$: cost

$$C_{empty} = C_{empty} + \Delta\Phi = \Phi(\text{stack}) - \Phi(\text{stack}) = 0$$

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$$\hat{C}_{\text{pop}} = C_{\text{pop}} + \Delta\Phi = 1 - 1 \leq 0 .$$

- ▶ **$S.\text{multipop}(k)$** : cost

$$\hat{C}_{\text{mp}} = C_{\text{mp}} + \Delta\Phi = \min\{\text{size}, k\} - \min\{\text{size}, k\} \leq 0 .$$

Example: Stack

Use potential function $\Phi(S) = \text{number of elements on the stack}$.

Amortized cost:

- ▶ **$S.\text{push}()$** : cost

$$\hat{C}_{\text{push}} = C_{\text{push}} + \Delta\Phi = 1 + 1 \leq 2 .$$

- ▶ **$S.\text{pop}()$** : cost

$$\hat{C}_{\text{pop}} = C_{\text{pop}} + \Delta\Phi = 1 - 1 \leq 0 .$$

- ▶ **$S.\text{multipop}(k)$** : cost

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Example: Binary Counter

Incrementing a binary counter:

Consider a computational model where each bit-operation costs one time-unit.

Incrementing an n -bit binary counter may require to examine n -bits, and maybe change them.

Actual cost:

- ▶ Changing bit from 0 to 1: cost 1.
- ▶ Changing bit from 1 to 0: cost 1.
- ▶ Increment: cost is $k + 1$, where k is the number of consecutive ones in the least significant bit-positions (e.g., 001101 has $k = 1$).

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Example: Binary Counter

Choose potential function $\Phi(x) = k$, where k denotes the number of ones in the binary representation of x .

Amortized cost:

• Changing bit from 0 to 1: cost

$$C_{i+1} = C_i + \Delta\Phi = 1 + 1 \leq 2$$

• Changing bit from 1 to 0: cost

$$C_{i+1} = C_i + \Delta\Phi = 1 - 1 \leq 0$$

• Bonus: Let l denote the number of consecutive ones in the first i significant bit-positions. An increment applies l $0 \rightarrow 1$ -operations, and one $1 \rightarrow 0$ -operation.

• Hence, the amortized cost is $C_{i+1} = C_i + 2$.

Example: Binary Counter

Choose potential function $\Phi(x) = k$, where k denotes the number of ones in the binary representation of x .

Amortized cost:

- ▶ Changing bit from 0 to 1: cost

$$\hat{C}_{0 \rightarrow 1} = C_{0 \rightarrow 1} + \Delta\Phi = 1 + 1 \leq 2 .$$

- ▶ Changing bit from 1 to 0: cost 0.

$$\hat{C}_{1 \rightarrow 0} = C_{1 \rightarrow 0} + \Delta\Phi = 1 - 1 \leq 0 .$$

- ▶ Increment. Let k denotes the number of consecutive ones in the least significant bit-positions. An increment involves k (1 \rightarrow 0)-operations, and one (0 \rightarrow 1)-operation.

Hence, the amortized cost is $k\hat{C}_{1 \rightarrow 0} + \hat{C}_{0 \rightarrow 1} \leq 2$.

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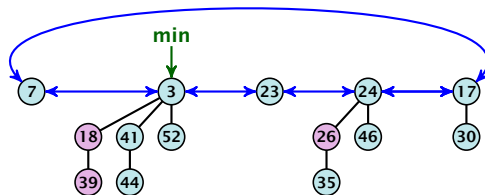
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Hence, the amortized cost is $k\hat{C}_{1 \rightarrow 0} + \hat{C}_{0 \rightarrow 1} \leq 2$.

8.3 Fibonacci Heaps

Collection of trees that fulfill the heap property.

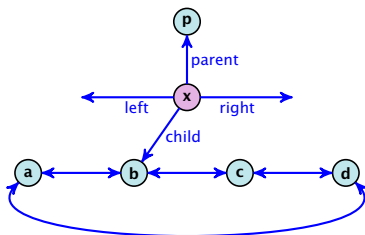
Structure is much more relaxed than binomial heaps.



8.3 Fibonacci Heaps

How do we implement trees with non-constant degree?

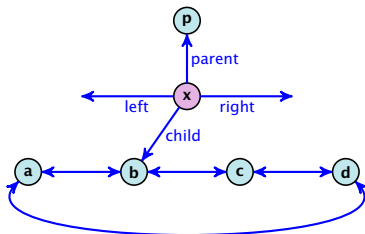
- ▶ The children of a node are arranged in a **circular linked list**.
- ▶ A child-pointer points to an arbitrary node within the list.
- ▶ A parent-pointer points to the parent node.
- ▶ Pointers $x.\text{left}$ and $x.\text{right}$ point to the left and right sibling of x (if x does not have siblings then $x.\text{left} = x.\text{right} = x$).



8.3 Fibonacci Heaps

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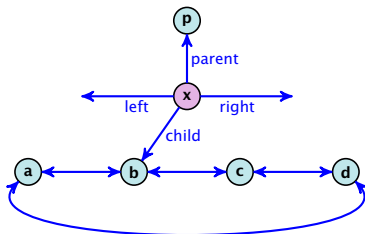
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8.3 Fibonacci Heaps

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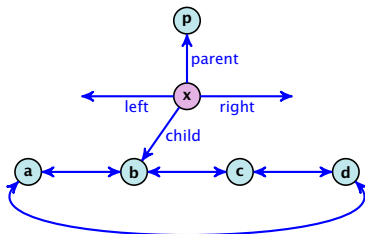
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8.3 Fibonacci Heaps

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- ▶ Pointers $x.left$ and $x.right$ point to the left and right sibling of x (if x does not have siblings then $x.left = x.right = x$).



8.3 Fibonacci Heaps

- ▶ Given a pointer to a node x we can splice out the sub-tree rooted at x in constant time.
- ▶ We can add a child-tree T to a node x in constant time if we are given a pointer to x and a pointer to the root of T .

8.3 Fibonacci Heaps

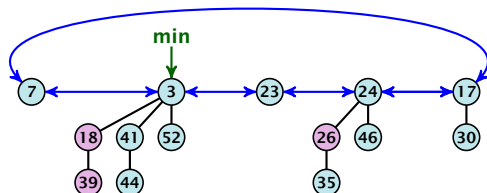
Additional implementation details:

- ▶ Every node x stores its degree in a field $x.degree$. Note that this can be updated in constant time when adding a child to x .
- ▶ Every node stores a boolean value $x.marked$ that specifies whether x is **marked** or not.

8.3 Fibonacci Heaps

The potential function:

- ▶ $t(S)$ denotes the number of trees in the heap.
- ▶ $m(S)$ denotes the number of marked nodes.
- ▶ We use the potential function $\Phi(S) = t(S) + 2m(S)$.



The potential is $\Phi(S) = 5 + 2 \cdot 3 = 11$.

8.3 Fibonacci Heaps

We assume that one unit of potential can pay for a constant amount of work, where the constant is chosen “big enough” (to take care of the constants that occur).

To make this more explicit we use c to denote the amount of work that a unit of potential can pay for.

8.3 Fibonacci Heaps

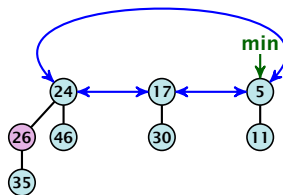
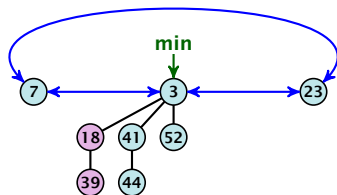
S. minimum()

- ▶ Access through the min-pointer.
- ▶ Actual cost $\mathcal{O}(1)$.
- ▶ No change in potential.
- ▶ Amortized cost $\mathcal{O}(1)$.

8.3 Fibonacci Heaps

S . merge(S')

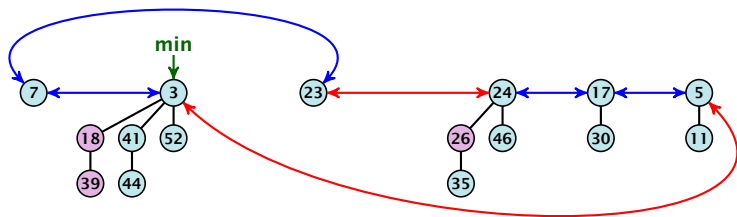
- ▶ Merge the root lists.
- ▶ Adjust the min-pointer



8.3 Fibonacci Heaps

S. merge(S')

- ▶ Merge the root lists.
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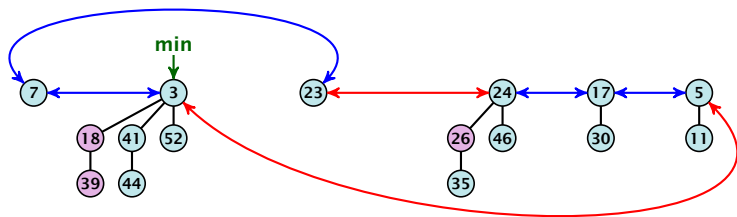
Running time:

- ▶ Actual cost $\mathcal{O}(1)$.

8.3 Fibonacci Heaps

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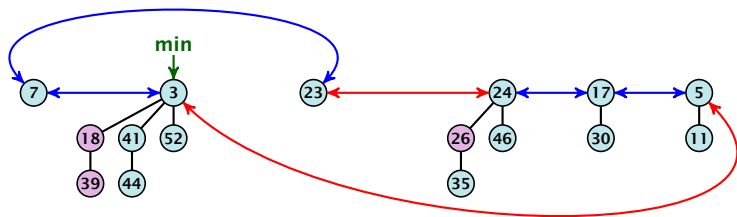
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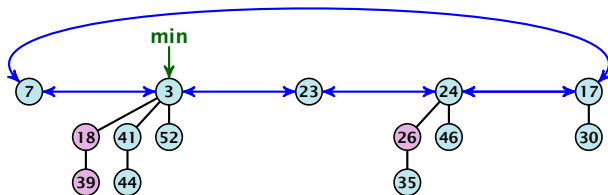
Running time:

- ▶ Actual cost $\mathcal{O}(1)$.
- ▶ No change in potential.
- ▶ Hence, amortized cost is $\mathcal{O}(1)$.

8.3 Fibonacci Heaps

S. insert(x)

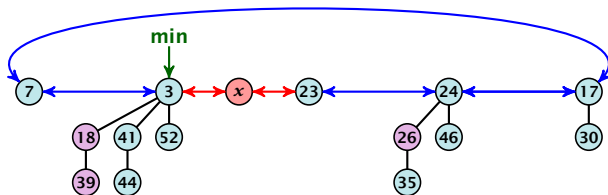
- ▶ Create a new tree containing x .
- ▶ Insert x into the root-list.
- ▶ Update min-pointer, if necessary.



8.3 Fibonacci Heaps

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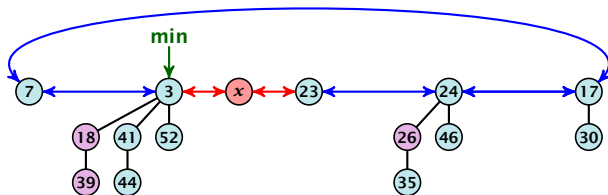
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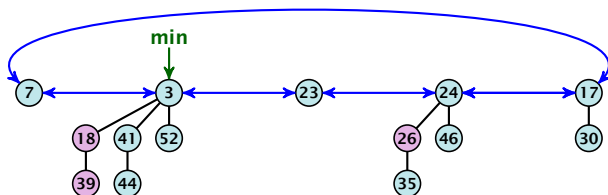


Running time:

- ▶ Actual cost $\mathcal{O}(1)$.
- ▶ Change in potential is $+1$.
- ▶ Amortized cost is $c + \mathcal{O}(1) = \mathcal{O}(1)$.

8.3 Fibonacci Heaps

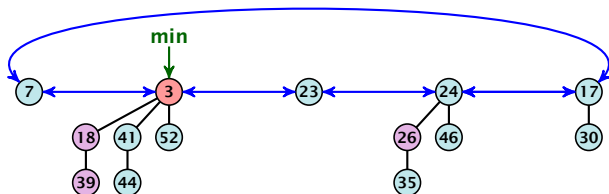
S. delete-min(x)



8.3 Fibonacci Heaps

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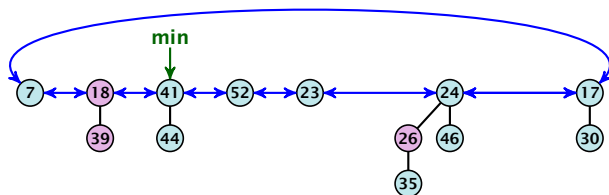
- ▶ Delete minimum; add child-trees to heap;
time: $D(\min) \cdot \mathcal{O}(1)$.



8.3 Fibonacci Heaps

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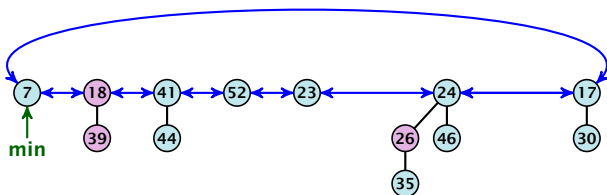
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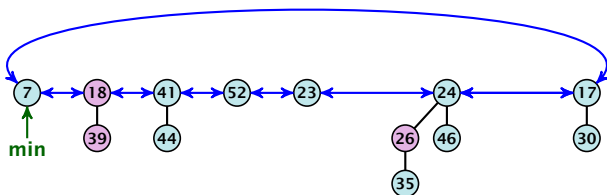
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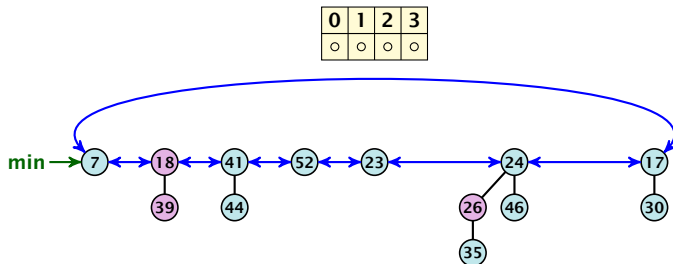
- ▶ Delete minimum; add child-trees to heap; time: $D(\min) \cdot \mathcal{O}(1)$.
- ▶ Update min-pointer; time: $(t + D(\min)) \cdot \mathcal{O}(1)$.



- ▶ Consolidate root-list so that no roots have the same degree. Time $t \cdot \mathcal{O}(1)$ (see next slide).

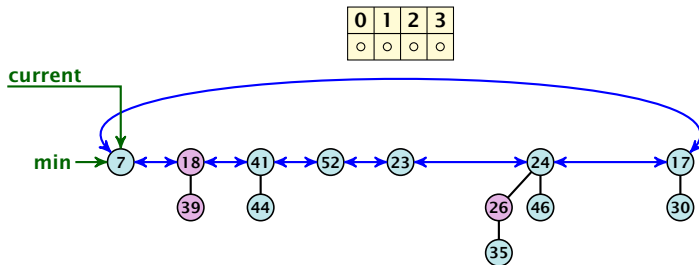
8.3 Fibonacci Heaps

Consolidate:



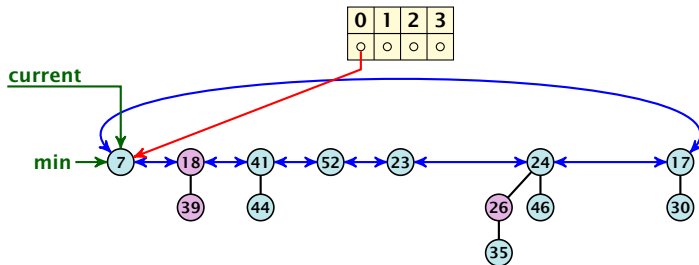
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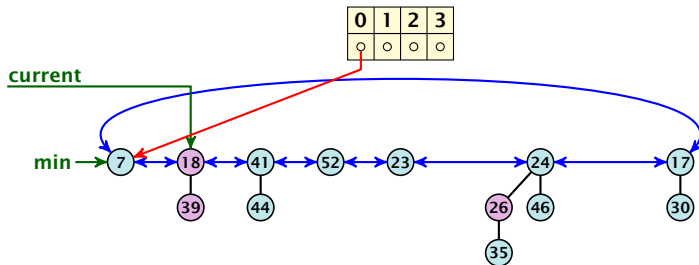
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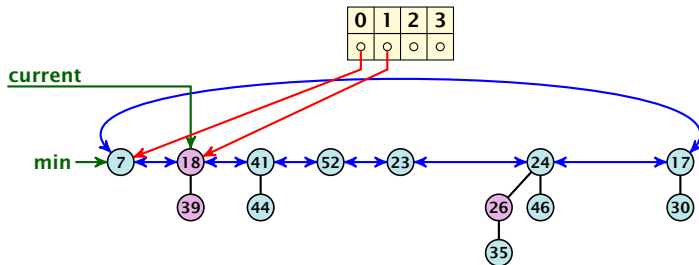
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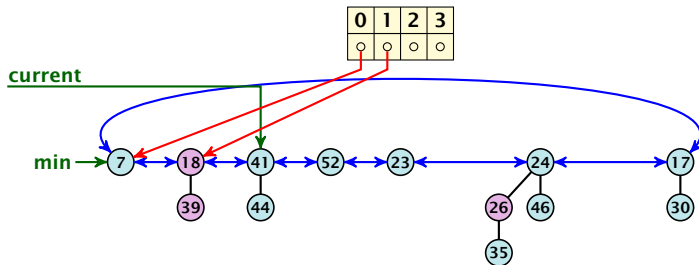
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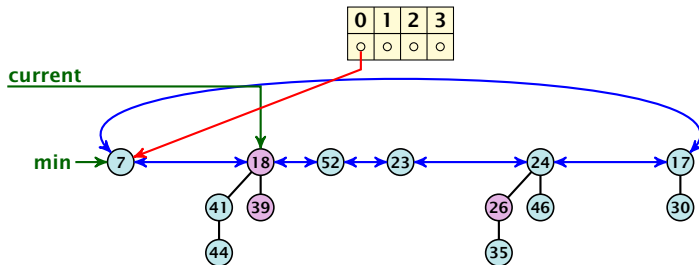
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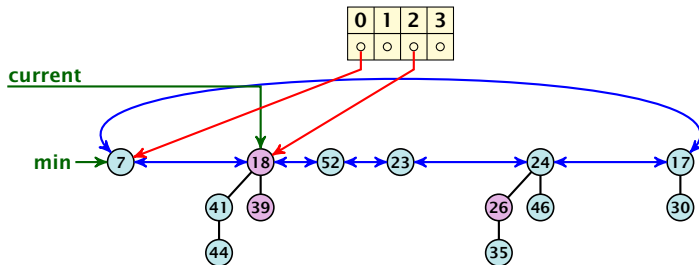
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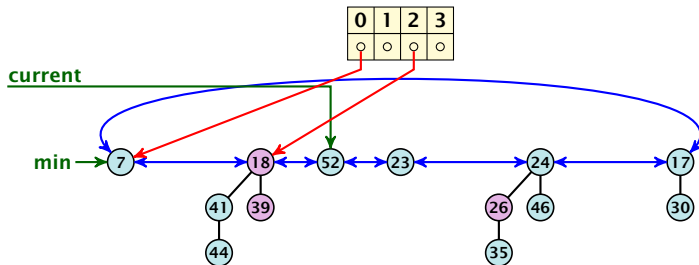
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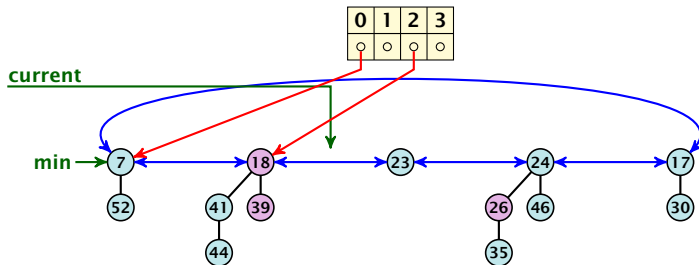
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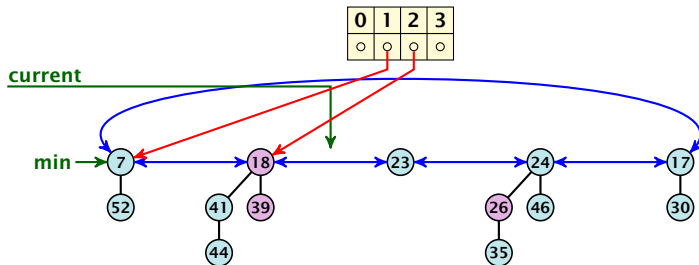
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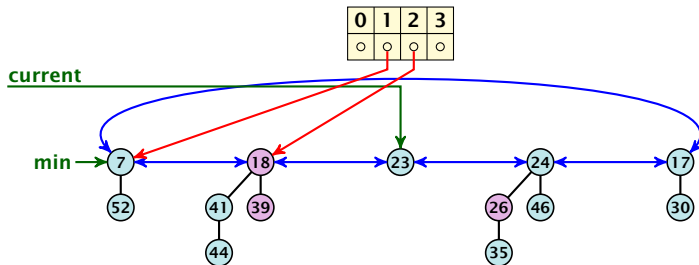
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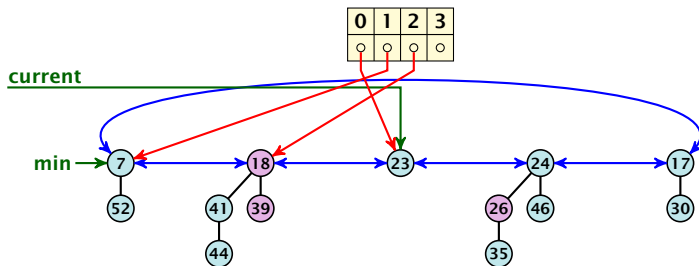
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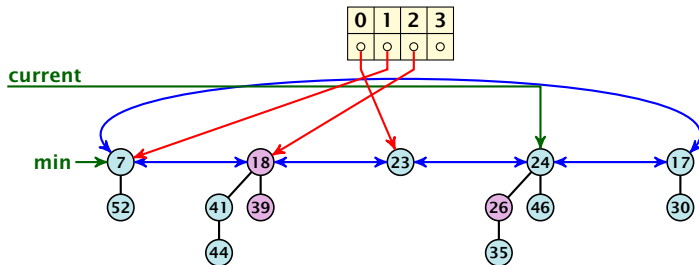
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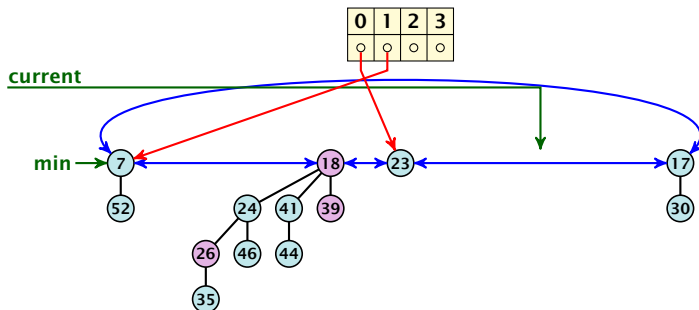
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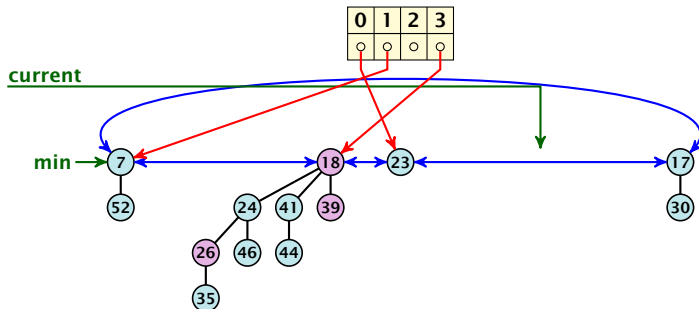
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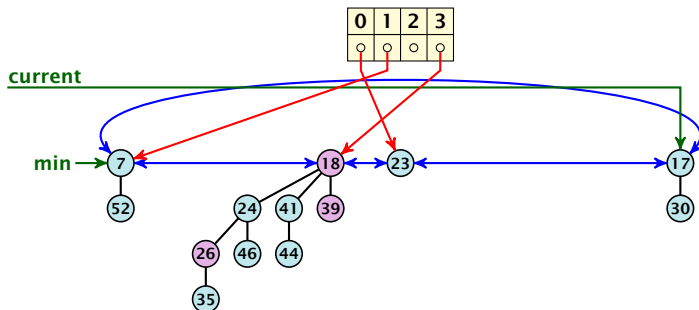
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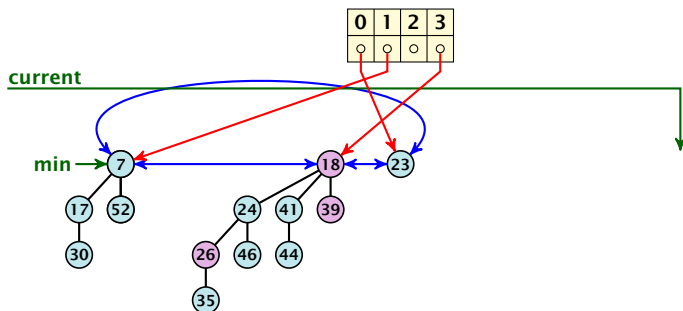
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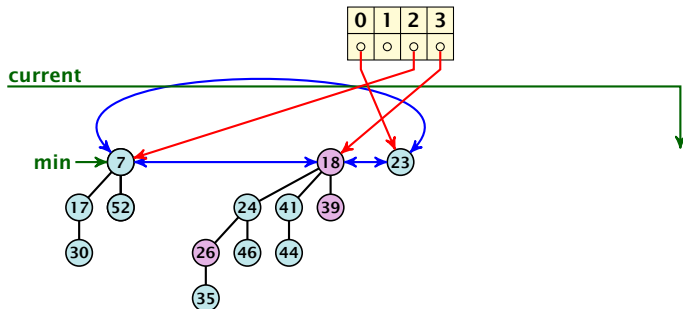
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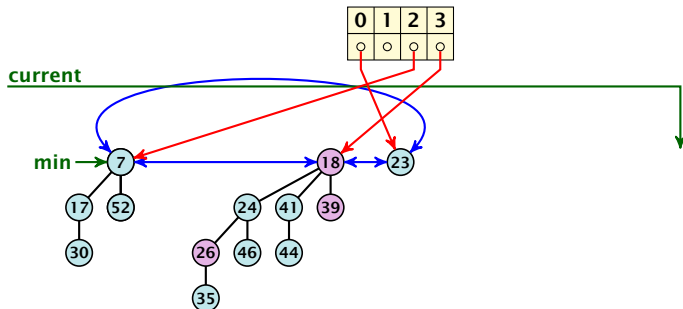
8.3 Fibonacci Heaps

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8.3 Fibonacci Heaps

Actual cost for delete-min()

- ▶ At most $D_n + t$ elements in root-list before consolidate.

Amortized cost for delete-min()

- ▶ $t' \leq D_n + 1$ as degrees are different after consolidating.
- ▶ Therefore $\Delta\Phi \leq D_n + 1 - t$;
- ▶ We can pay $c \cdot (t - D_n - 1)$ from the potential decrease.
- ▶ The amortized cost is

8.3 Fibonacci Heaps

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- ▶ Actual cost for a delete-min is at most $\mathcal{O}(1) \cdot (D_n + t)$.
Hence, there exists c_1 s.t. actual cost is at most $c_1 \cdot (D_n + t)$.

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- ▶ The amortized cost is

$$c_1 \cdot (D_n + t) - c \cdot (t - D_n - 1)$$

8.3 Fibonacci Heaps

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$$\begin{aligned}c_1 \cdot (D_n + t) - c \cdot (t - D_n - 1) \\ \leq (c_1 + c)D_n + (c_1 - c)t + c\end{aligned}$$

8.3 Fibonacci Heaps

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- ▶ We can pay $c \cdot (t - D_n - 1)$ from the potential decrease.
- ▶ The amortized cost is

$$\begin{aligned}c_1 \cdot (D_n + t) - c \cdot (t - D_n - 1) \\ \leq (c_1 + c)D_n + (c_1 - c)t + c \leq 2c(D_n + 1)\end{aligned}$$

8.3 Fibonacci Heaps

Actual cost for delete-min()

- ▶ At most $D_n + t$ elements in root-list before consolidate.
- ▶ Actual cost for a delete-min is at most $\mathcal{O}(1) \cdot (D_n + t)$.
Hence, there exists c_1 s.t. actual cost is at most $c_1 \cdot (D_n + t)$.

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for $c \geq c_1$.

8.3 Fibonacci Heaps

If the input trees of the consolidation procedure are binomial trees (for example only singleton vertices) then the output will be a set of distinct binomial trees, and, hence, the Fibonacci heap will be (more or less) a Binomial heap right after the consolidation.

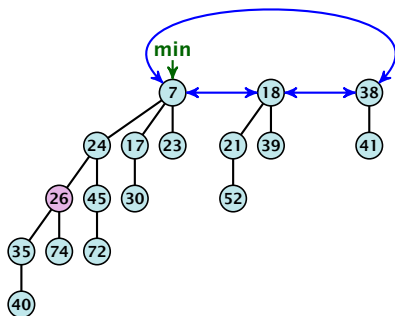
If we do not have delete or decrease-key operations then
 $D_n \leq \log n$.

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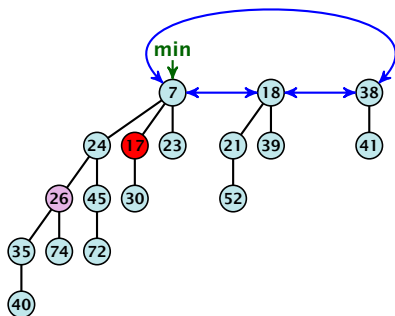
Fibonacci Heaps: decrease-key(handle h, v)



Case 1: decrease-key does not violate heap-property

- ▶ Just decrease the key-value of element referenced by h . Nothing else to do.

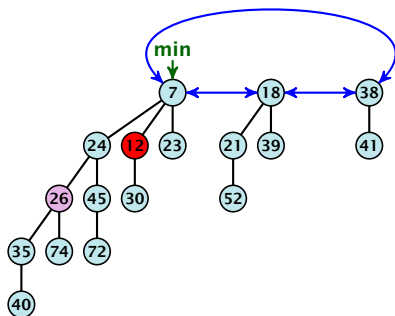
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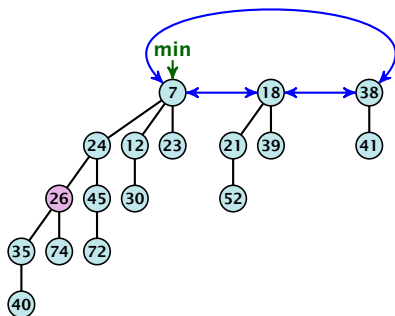
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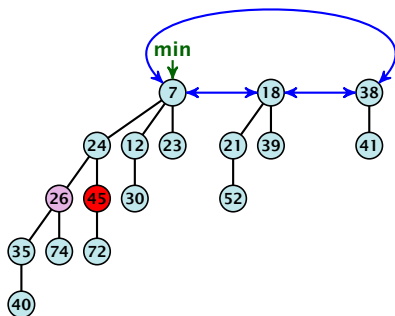
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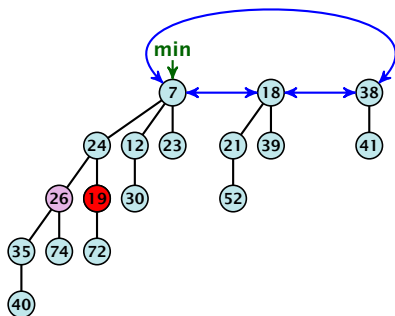
Fibonacci Heaps: decrease-key(handle h, v)



Case 2: heap-property is violated, but parent is not marked

- ▶ Decrease key-value of element x reference by h .
- ▶ If the heap-property is violated, cut the parent edge of x , and make x into a root.
- ▶ Adjust min-pointers, if necessary.
- ▶ Mark the (previous) parent of x .

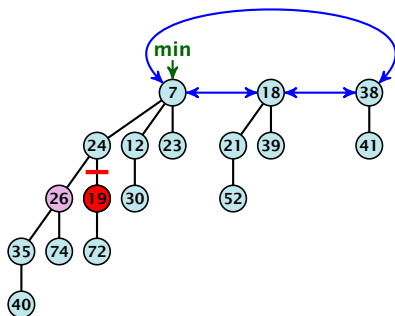
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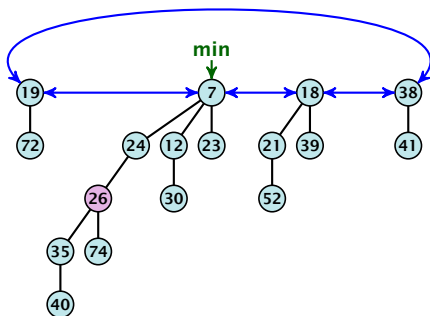
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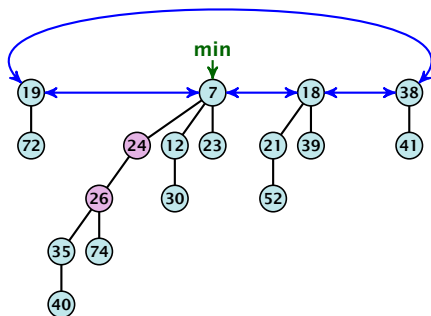
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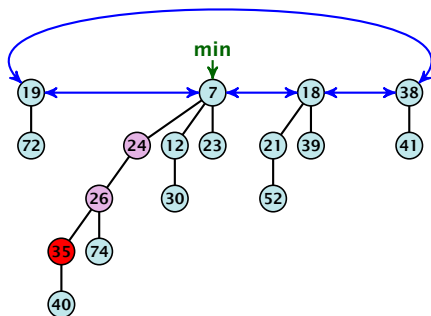
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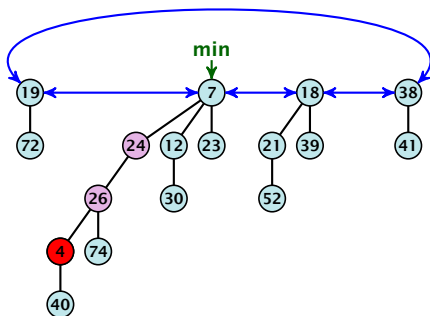
Fibonacci Heaps: decrease-key(handle h, v)



Case 3: heap-property is violated, and parent is marked

- ▶ Decrease key-value of element x reference by h .
- ▶ Cut the parent edge of x , and make x into a root.
- ▶ Adjust min-pointers, if necessary.
- ▶ Continue cutting the parent until you arrive at an unmarked node.

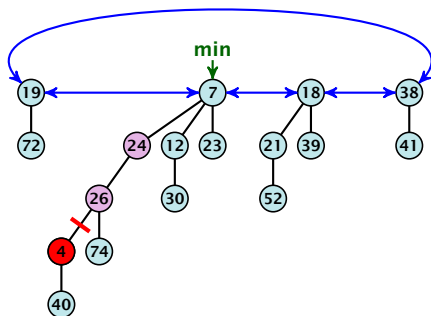
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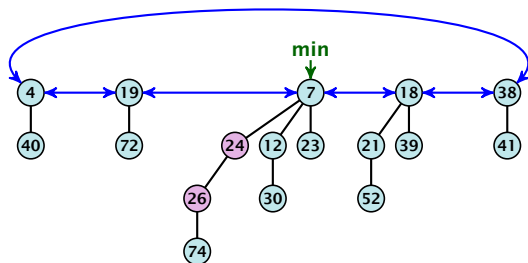
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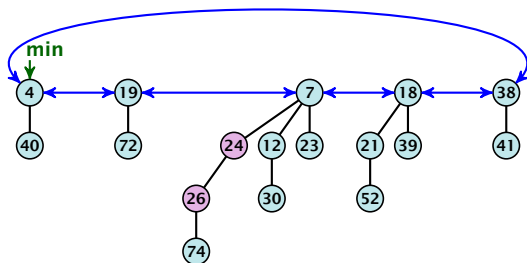
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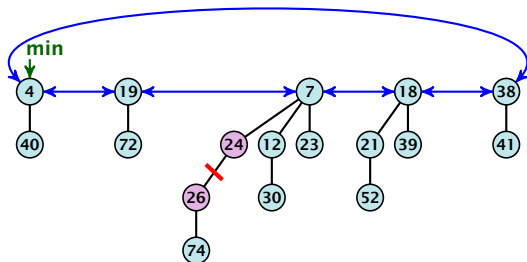
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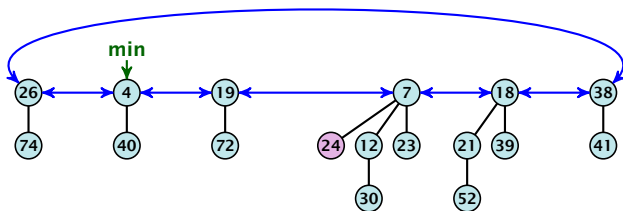
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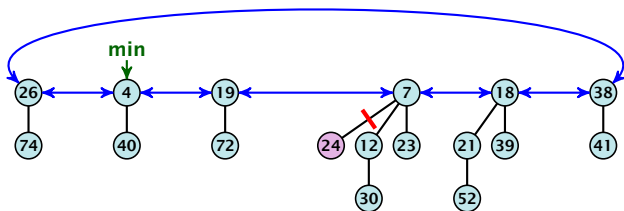
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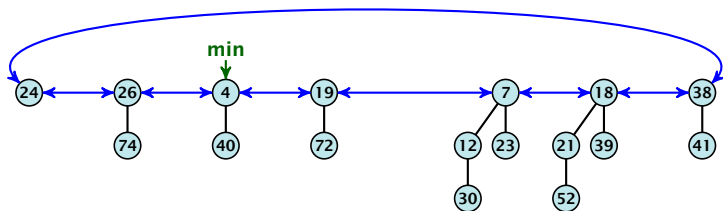
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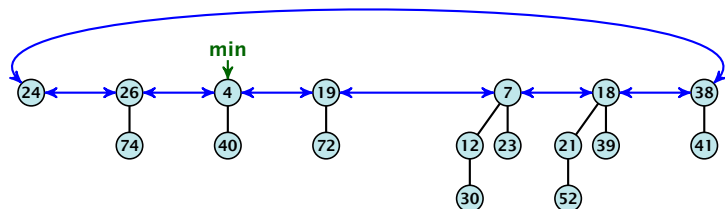
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- ▶ Decrease key-value of element x reference by h .
- ▶ Cut the parent edge of x , and make x into a root.
- ▶ Adjust min-pointers, if necessary.

- ▶ Execute the following:

$p \leftarrow \text{parent}[x];$

while (p is marked)

$pp \leftarrow \text{parent}[p];$

cut of p ; make it into a root; **unmark it**;

$p \leftarrow pp;$

if p is unmarked and not a root mark it;

Fibonacci Heaps: decrease-key(handle h, v)

Actual cost:

- ▶ Constant cost for decreasing the value.
- ▶ Constant cost for each of ℓ cuts.
- ▶ Hence, cost is at most $c_2 \cdot (\ell + 1)$, for some constant c_2 .

Amortized cost:

- ▶ $\ell = O(\log n)$, as every cut creates one new root.
- ▶ $\ell = O(\log n) + 1 = O(\log n)$, since all but the first cut mark a node; the first cut may mark a node.
- ▶ Hence, $c_1 \cdot \ell + c_2 \cdot (\ell + 1) = O(\log n)$.

▶ Amortized cost is at most $O(\log n)$.

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- ▶ $\ell = O(\log \ell)$, as every cut creates one new root.
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- ▶ $\ell = O(\log \ell) + 1 = O(\log \ell)$.
- ▶ Amortized cost is $O(1)$.

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Amortized cost:

- ▶ $\ell \leq \ell_0 + \ell_1$, as every cut creates one new root.
- ▶ $\ell_0 \leq \log_2(\ell + 1) + 1$ since all but the first cut decrease the number of nodes by at least a factor of 2.
- ▶ $\ell_1 \leq \log_2(\ell + 1) + 1$ since all but the first cut decrease the number of nodes by at least a factor of 2.

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Amortized cost:

- ▶ $t' = t + \ell$, as every cut creates one new root.
- ▶ $m' \leq m - (\ell - 1) + 1 = m - \ell + 2$, since all but the first cut marks a node; the last cut may mark a node.
- ▶ $\Delta\Phi \leq \ell + 2(-\ell + 2) = 4 - \ell$
- ▶ Amortized cost is at most

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$$c_2(\ell + 1) + c(4 - \ell) \leq (c_2 - c)\ell + 4c = \mathcal{O}(1),$$

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Delete node

H. delete(x):

- ▶ decrease value of x to $-\infty$.
- ▶ delete-min.

Amortized cost: $\mathcal{O}(D(n))$

- ▶ $\mathcal{O}(1)$ for decrease-key.
- ▶ $\mathcal{O}(D(n))$ for delete-min.

8.3 Fibonacci Heaps

Lemma 33

Let x be a node with degree k and let y_1, \dots, y_k denote the children of x in the order that they were linked to x . Then

$$\text{degree}(y_i) \geq \begin{cases} 0 & \text{if } i = 1 \\ i - 2 & \text{if } i \geq 2 \end{cases}$$

8.3 Fibonacci Heaps

Proof

- ▶ When y_i was linked to x , at least y_1, \dots, y_{i-1} were already linked to x .
- ▶ Hence, at this time $\text{degree}(x) \geq i - 1$, and therefore also $\text{degree}(y_i) \geq i - 1$ as the algorithm links nodes of equal degree only.
- ▶ Since, then y_i has lost at most one child.
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8.3 Fibonacci Heaps

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Let x be a degree k node of size s_k and let y_1, \dots, y_k be its children.

$$s_k = 2 + \sum_{i=2}^k \text{size}(y_i)$$

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8.3 Fibonacci Heaps

Definition 34

Consider the following non-standard Fibonacci type sequence:

$$F_k = \begin{cases} 1 & \text{if } k = 0 \\ 2 & \text{if } k = 1 \\ F_{k-1} + F_{k-2} & \text{if } k \geq 2 \end{cases}$$

Facts:

1. $F_k \geq \phi^k$.
2. For $k \geq 2$: $F_k = 2 + \sum_{i=0}^{k-2} F_i$.

The above facts can be easily proved by induction. From this it follows that $s_k \geq F_k \geq \phi^k$, which gives that the maximum degree in a Fibonacci heap is logarithmic.

9 van Emde Boas Trees

Dynamic Set Data Structure S :

- ▶ $S.insert(x)$
- ▶ $S.delete(x)$
- ▶ $S.search(x)$
- ▶ $S.min()$
- ▶ $S.max()$
- ▶ $S.succ(x)$
- ▶ $S.pred(x)$

9 van Emde Boas Trees

For this chapter we ignore the problem of storing satellite data:

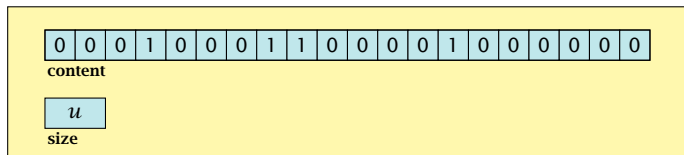
- ▶ **S . insert(x):** Inserts x into S .
- ▶ **S . delete(x):** Deletes x from S . Usually assumes that $x \in S$.
- ▶ **S . member(x):** Returns 1 if $x \in S$ and 0 otherwise.
- ▶ **S . min():** Returns the value of the minimum element in S .
- ▶ **S . max():** Returns the value of the maximum element in S .
- ▶ **S . succ(x):** Returns successor of x in S . Returns null if x is maximum or larger than any element in S . Note that x needs not to be in S .
- ▶ **S . pred(x):** Returns the predecessor of x in S . Returns null if x is minimum or smaller than any element in S . Note that x needs not to be in S .

9 van Emde Boas Trees

Can we improve the existing algorithms when the keys are from a restricted set?

In the following we assume that the keys are from $\{0, 1, \dots, u - 1\}$, where u denotes the size of the universe.

Implementation 1: Array



one array of u bits

Use an array that encodes the indicator function of the dynamic set.

Implementation 1: Array

Algorithm 19 `array.insert(x)`

1: `content[x] ← 1;`

Algorithm 20 `array.delete(x)`

1: `content[x] ← 0;`

Algorithm 21 `array.member(x)`

1: **return** `content[x];`

- ▶ Note that we assume that x is valid, i.e., it falls within the array boundaries.
- ▶ Obviously(?) the running time is constant.

Implementation 1: Array

Algorithm 22 array.max()

```
1: for ( $i = \text{size} - 1; i \geq 0; i--$ ) do  
2:     if content[ $i$ ] = 1 then return  $i$ ;  
3: return null;
```

Algorithm 23 array.min()

```
1: for ( $i = 0; i < \text{size}; i++$ ) do  
2:     if content[ $i$ ] = 1 then return  $i$ ;  
3: return null;
```

- ▶ Running time is $\mathcal{O}(u)$ in the worst case.

Implementation 1: Array

Algorithm 22 array.max()

```
1: for ( $i = \text{size} - 1; i \geq 0; i--$ ) do
2:     if content[ $i$ ] = 1 then return  $i$ ;
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Algorithm 23 array.min()

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▶ Running time is $\mathcal{O}(u)$ in the worst case.

Implementation 1: Array

Algorithm 22 array.max()

```
1: for ( $i = \text{size} - 1; i \geq 0; i--$ ) do  
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3: return null;
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Algorithm 23 array.min()

```
1: for ( $i = 0; i < \text{size}; i++$ ) do  
2:     if content[ $i$ ] = 1 then return  $i$ ;  
3: return null;
```

- ▶ Running time is $\mathcal{O}(u)$ in the worst case.

Implementation 1: Array

Algorithm 24 `array.succ(x)`

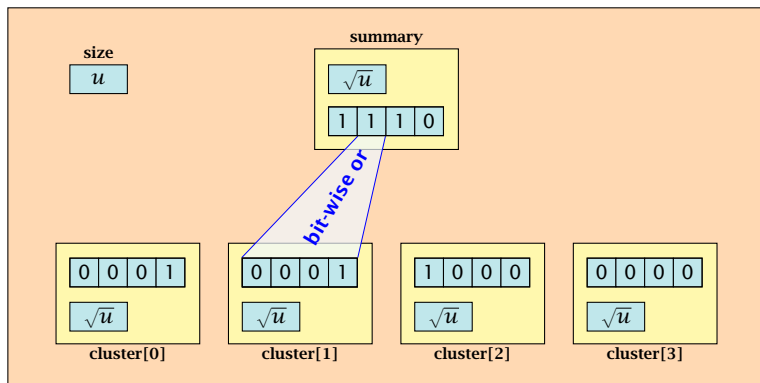
```
1: for ( $i = x + 1$ ;  $i < \text{size}$ ;  $i++$ ) do  
2:     if content[i] = 1 then return  $i$ ;  
3: return null;
```

Algorithm 25 `array.pred(x)`

```
1: for ( $i = x - 1$ ;  $i \geq 0$ ;  $i--$ ) do  
2:     if content[i] = 1 then return  $i$ ;  
3: return null;
```

- ▶ Running time is $\mathcal{O}(u)$ in the worst case.

Implementation 2: Summary Array



- ▶ \sqrt{u} cluster-arrays of \sqrt{u} bits.
- ▶ One summary-array of \sqrt{u} bits. The i -th bit in the summary array stores the bit-wise or of the bits in the i -th cluster.

Implementation 2: Summary Array

The bit for a key x is contained in cluster number $\lfloor \frac{x}{\sqrt{u}} \rfloor$.

Within the cluster-array the bit is at position $x \bmod \sqrt{u}$.

For simplicity we assume that $u = 2^{2k}$ for some $k \geq 1$. Then we can compute the cluster-number for an entry x as $\text{high}(x)$ (the upper half of the dual representation of x) and the position of x within its cluster as $\text{low}(x)$ (the lower half of the dual representation).

Implementation 2: Summary Array

The bit for a key x is contained in cluster number $\left\lfloor \frac{x}{\sqrt{u}} \right\rfloor$.

Within the cluster-array the bit is at position $x \bmod \sqrt{u}$.

For simplicity we assume that $u = 2^{2k}$ for some $k \geq 1$. Then we can compute the cluster-number for an entry x as $\text{high}(x)$ (the upper half of the dual representation of x) and the position of x within its cluster as $\text{low}(x)$ (the lower half of the dual representation).

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For simplicity we assume that $u = 2^{2k}$ for some $k \geq 1$. Then we can compute the cluster-number for an entry x as $\text{high}(x)$ (the upper half of the dual representation of x) and the position of x within its cluster as $\text{low}(x)$ (the lower half of the dual representation).

Implementation 2: Summary Array

Algorithm 26 $\text{member}(x)$

1: **return** $\text{cluster}[\text{high}(x)].\text{member}(\text{low}(x));$

Algorithm 27 $\text{insert}(x)$

1: $\text{cluster}[\text{high}(x)].\text{insert}(\text{low}(x));$

2: $\text{summary}.\text{insert}(\text{high}(x));$

- ▶ The running times are constant, because the corresponding array-functions have constant running times.

Implementation 2: Summary Array

Algorithm 26 $\text{member}(x)$

1: **return** $\text{cluster}[\text{high}(x)].\text{member}(\text{low}(x));$

Algorithm 27 $\text{insert}(x)$

1: $\text{cluster}[\text{high}(x)].\text{insert}(\text{low}(x));$

2: $\text{summary}.\text{insert}(\text{high}(x));$

- ▶ The running times are constant, because the corresponding array-functions have constant running times.

Implementation 2: Summary Array

Algorithm 26 $\text{member}(x)$

1: **return** $\text{cluster}[\text{high}(x)].\text{member}(\text{low}(x));$

Algorithm 27 $\text{insert}(x)$

1: $\text{cluster}[\text{high}(x)].\text{insert}(\text{low}(x));$

2: $\text{summary}.\text{insert}(\text{high}(x));$

- ▶ The running times are constant, because the corresponding array-functions have constant running times.

Implementation 2: Summary Array

Algorithm 28 delete(x)

```
1: cluster[high( $x$ )].delete(low( $x$ ));  
2: if cluster[high( $x$ )].min() = null then  
3:     summary.delete(high( $x$ ));
```

- ▶ The running time is dominated by the cost of a minimum computation, which will turn out to be $\mathcal{O}(\sqrt{u})$.

Implementation 2: Summary Array

Algorithm 28 delete(x)

```
1: cluster[high( $x$ )].delete(low( $x$ ));  
2: if cluster[high( $x$ )].min() = null then  
3:     summary.delete(high( $x$ ));
```

- ▶ The running time is dominated by the cost of a minimum computation, which will turn out to be $\mathcal{O}(\sqrt{u})$.

Implementation 2: Summary Array

Algorithm 29 $\text{max}()$

```
1:  $\text{maxcluster} \leftarrow \text{summary.max}()$ ;  
2: if  $\text{maxcluster} = \text{null}$  return  $\text{null}$ ;  
3:  $\text{offs} \leftarrow \text{cluster}[\text{maxcluster}].\text{max}()$   
4: return  $\text{maxcluster} \circ \text{offs}$ ;
```

Algorithm 30 $\text{min}()$

```
1:  $\text{mincluster} \leftarrow \text{summary.min}()$ ;  
2: if  $\text{mincluster} = \text{null}$  return  $\text{null}$ ;  
3:  $\text{offs} \leftarrow \text{cluster}[\text{mincluster}].\text{min}()$ ;  
4: return  $\text{mincluster} \circ \text{offs}$ ;
```

▶ Running time is roughly $2\sqrt{u} = \mathcal{O}(u)$ in the worst case.

Implementation 2: Summary Array

Algorithm 29 $\text{max}()$

```
1:  $\text{maxcluster} \leftarrow \text{summary.max}();$   
2: if  $\text{maxcluster} = \text{null}$  return  $\text{null}$ ;  
3:  $\text{offs} \leftarrow \text{cluster}[\text{maxcluster}].\text{max}();$   
4: return  $\text{maxcluster} \circ \text{offs};$ 
```

Algorithm 30 $\text{min}()$

```
1:  $\text{mincluster} \leftarrow \text{summary.min}();$   
2: if  $\text{mincluster} = \text{null}$  return  $\text{null}$ ;  
3:  $\text{offs} \leftarrow \text{cluster}[\text{mincluster}].\text{min}();$   
4: return  $\text{mincluster} \circ \text{offs};$ 
```

▶ Running time is roughly $2\sqrt{u} = \mathcal{O}(u)$ in the worst case.

Implementation 2: Summary Array

Algorithm 29 $\text{max}()$

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2: if  $\text{maxcluster} = \text{null}$  return  $\text{null}$ ;  
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4: return  $\text{maxcluster} \circ \text{offs};$ 
```

Algorithm 30 $\text{min}()$

```
1:  $\text{mincluster} \leftarrow \text{summary.min}();$   
2: if  $\text{mincluster} = \text{null}$  return  $\text{null}$ ;  
3:  $\text{offs} \leftarrow \text{cluster}[\text{mincluster}].\text{min}();$   
4: return  $\text{mincluster} \circ \text{offs};$ 
```

- ▶ Running time is roughly $2\sqrt{u} = \mathcal{O}(u)$ in the worst case.

Implementation 2: Summary Array

Algorithm 31 $\text{succ}(x)$

```
1:  $m \leftarrow \text{cluster}[\text{high}(x)].\text{succ}(\text{low}(x))$ 
2: if  $m \neq \text{null}$  then return  $\text{high}(x) \circ m$ ;
3:  $\text{succcluster} \leftarrow \text{summary}.\text{succ}(\text{high}(x))$ ;
4: if  $\text{succcluster} \neq \text{null}$  then
5:    $\text{offs} \leftarrow \text{cluster}[\text{succcluster}].\text{min}()$ ;
6:   return  $\text{succcluster} \circ \text{offs}$ ;
7: return  $\text{null}$ ;
```

▶ Running time is roughly $3\sqrt{u} = \mathcal{O}(\sqrt{u})$ in the worst case.

Implementation 2: Summary Array

Algorithm 31 $\text{succ}(x)$

```
1:  $m \leftarrow \text{cluster}[\text{high}(x)].\text{succ}(\text{low}(x))$ 
2: if  $m \neq \text{null}$  then return  $\text{high}(x) \circ m$ ;
3:  $\text{succcluster} \leftarrow \text{summary}.\text{succ}(\text{high}(x))$ ;
4: if  $\text{succcluster} \neq \text{null}$  then
5:      $\text{offs} \leftarrow \text{cluster}[\text{succcluster}].\text{min}()$ ;
6:     return  $\text{succcluster} \circ \text{offs}$ ;
7: return  $\text{null}$ ;
```

- ▶ Running time is roughly $3\sqrt{u} = \mathcal{O}(\sqrt{u})$ in the worst case.

Implementation 2: Summary Array

Algorithm 32 $\text{pred}(x)$

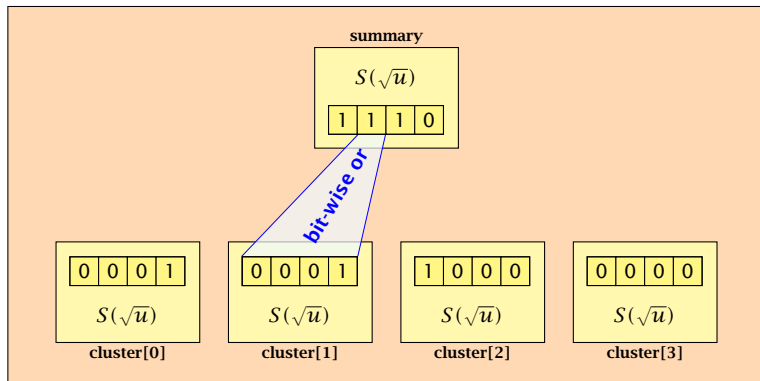
```
1:  $m \leftarrow \text{cluster}[\text{high}(x)].\text{pred}(\text{low}(x))$ 
2: if  $m \neq \text{null}$  then return  $\text{high}(x) \circ m$ ;
3:  $\text{predcluster} \leftarrow \text{summary}.\text{pred}(\text{high}(x))$ ;
4: if  $\text{predcluster} \neq \text{null}$  then
5:      $\text{offs} \leftarrow \text{cluster}[\text{predcluster}].\text{max}()$ ;
6:     return  $\text{predcluster} \circ \text{offs}$ ;
7: return  $\text{null}$ ;
```

- ▶ Running time is roughly $3\sqrt{u} = \mathcal{O}(\sqrt{u})$ in the worst case.

Implementation 3: Recursion

Instead of using sub-arrays, we build a recursive data-structure.

$S(u)$ is a dynamic set data-structure representing u bits:



Implementation 3: Recursion

We assume that $u = 2^{2^k}$ for some k .

The data-structure $S(2)$ is defined as an array of 2-bits (end of the recursion).

Implementation 3: Recursion

The code from Implementation 2 can be used **unchanged**. We only need to redo the analysis of the running time.

Note that in the code we do not need to specifically address the non-recursive case. This is achieved by the fact that an $S(4)$ will contain $S(2)$'s as sub-datastructures, which are **arrays**. Hence, a call like `cluster[1].min()` from within the data-structure $S(4)$ is **not** a recursive call as it will call the function `array.min()`.

This means that the non-recursive case is been dealt with while initializing the data-structure.

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This means that the non-recursive case is been dealt with while initializing the data-structure.

Implementation 3: Recursion

Algorithm 33 `member(x)`

```
1: return cluster[high(x)].member(low(x));
```

- ▶ $T_{\text{mem}}(u) = T_{\text{mem}}(\sqrt{u}) + 1.$

Implementation 3: Recursion

Algorithm 34 insert(x)

```
1: cluster[high( $x$ )].insert(low( $x$ ));  
2: summary.insert(high( $x$ ));
```

► $T_{\text{ins}}(u) = 2T_{\text{ins}}(\sqrt{u}) + 1.$

Implementation 3: Recursion

Algorithm 35 delete(x)

```
1: cluster[high( $x$ )].delete(low( $x$ ));  
2: if cluster[high( $x$ )].min() = null then  
3:     summary.delete(high( $x$ ));
```

- ▶ $T_{\text{del}}(u) = 2T_{\text{del}}(\sqrt{u}) + T_{\text{min}}(\sqrt{u}) + 1.$

Implementation 3: Recursion

Algorithm 36 $\text{min}()$

```
1: mincluster  $\leftarrow$  summary.min();  
2: if mincluster = null return null;  
3: offs  $\leftarrow$  cluster[mincluster].min();  
4: return mincluster  $\circ$  offs;
```

- ▶ $T_{\min}(u) = 2T_{\min}(\sqrt{u}) + 1$.

Implementation 3: Recursion

Algorithm 37 $\text{succ}(x)$

```
1:  $m \leftarrow \text{cluster}[\text{high}(x)].\text{succ}(\text{low}(x))$ 
2: if  $m \neq \text{null}$  then return  $\text{high}(x) \circ m$ ;
3:  $\text{succcluster} \leftarrow \text{summary}.\text{succ}(\text{high}(x))$ ;
4: if  $\text{succcluster} \neq \text{null}$  then
5:      $\text{offs} \leftarrow \text{cluster}[\text{succcluster}].\text{min}()$ ;
6:     return  $\text{succcluster} \circ \text{offs}$ ;
7: return  $\text{null}$ ;
```

- ▶ $T_{\text{succ}}(u) = 2T_{\text{succ}}(\sqrt{u}) + T_{\text{min}}(\sqrt{u}) + 1$.

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1:$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + \mathbf{1}:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$.

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + \mathbf{1}:$$

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$$X(\ell)$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + \mathbf{1}:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

$$X(\ell) = T_{\text{mem}}(2^\ell)$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + \mathbf{1}:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

$$X(\ell) = T_{\text{mem}}(2^\ell) = T_{\text{mem}}(u)$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

$$X(\ell) = T_{\text{mem}}(2^\ell) = T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

$$\begin{aligned} X(\ell) = T_{\text{mem}}(2^\ell) &= T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1 \\ &= T_{\text{mem}}(2^{\frac{\ell}{2}}) + 1 \end{aligned}$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

$$\begin{aligned} X(\ell) = T_{\text{mem}}(2^\ell) &= T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1 \\ &= T_{\text{mem}}(2^{\frac{\ell}{2}}) + 1 = X\left(\frac{\ell}{2}\right) + 1 . \end{aligned}$$

Implementation 3: Recursion

$$T_{\text{mem}}(\mathbf{u}) = T_{\text{mem}}(\sqrt{\mathbf{u}}) + 1:$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{mem}}(2^\ell)$. Then

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Using Master theorem gives $X(\ell) = \mathcal{O}(\log \ell)$, and hence $T_{\text{mem}}(\mathbf{u}) = \mathcal{O}(\log \log u)$.

Implementation 3: Recursion

$$T_{\text{ins}}(u) = 2T_{\text{ins}}(\sqrt{u}) + 1.$$

Implementation 3: Recursion

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Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

Implementation 3: Recursion

$$T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1.$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

$$X(\ell)$$

Implementation 3: Recursion

$$T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1.$$

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$$X(\ell) = T_{\text{ins}}(2^\ell)$$

Implementation 3: Recursion

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Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

$$X(\ell) = T_{\text{ins}}(2^\ell) = T_{\text{ins}}(\mathbf{u})$$

Implementation 3: Recursion

$$T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1.$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

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Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

$$\begin{aligned} X(\ell) &= T_{\text{ins}}(2^\ell) = T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1 \\ &= 2T_{\text{ins}}(2^{\frac{\ell}{2}}) + 1 \end{aligned}$$

Implementation 3: Recursion

$$T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1.$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

$$\begin{aligned} X(\ell) &= T_{\text{ins}}(2^\ell) = T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1 \\ &= 2T_{\text{ins}}(2^{\frac{\ell}{2}}) + 1 = 2X\left(\frac{\ell}{2}\right) + 1 . \end{aligned}$$

Implementation 3: Recursion

$$T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1.$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

$$\begin{aligned} X(\ell) &= T_{\text{ins}}(2^\ell) = T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1 \\ &= 2T_{\text{ins}}(2^{\frac{\ell}{2}}) + 1 = 2X\left(\frac{\ell}{2}\right) + 1 . \end{aligned}$$

Using Master theorem gives $X(\ell) = \mathcal{O}(\ell)$, and hence $T_{\text{ins}}(\mathbf{u}) = \mathcal{O}(\log u)$.

Implementation 3: Recursion

$$T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1.$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{ins}}(2^\ell)$. Then

$$\begin{aligned} X(\ell) &= T_{\text{ins}}(2^\ell) = T_{\text{ins}}(\mathbf{u}) = 2T_{\text{ins}}(\sqrt{\mathbf{u}}) + 1 \\ &= 2T_{\text{ins}}(2^{\frac{\ell}{2}}) + 1 = 2X(\frac{\ell}{2}) + 1 . \end{aligned}$$

Using Master theorem gives $X(\ell) = \mathcal{O}(\ell)$, and hence $T_{\text{ins}}(\mathbf{u}) = \mathcal{O}(\log u)$.

The same holds for $T_{\text{max}}(\mathbf{u})$ and $T_{\text{min}}(\mathbf{u})$.

Implementation 3: Recursion

$$T_{\text{del}}(\mathbf{u}) = 2T_{\text{del}}(\sqrt{\mathbf{u}}) + T_{\text{min}}(\sqrt{\mathbf{u}}) + 1 = 2T_{\text{del}}(\sqrt{\mathbf{u}}) + \Theta(\log(\mathbf{u})).$$

Implementation 3: Recursion

$$T_{\text{del}}(\mathbf{u}) = 2T_{\text{del}}(\sqrt{\mathbf{u}}) + T_{\text{min}}(\sqrt{\mathbf{u}}) + \mathbf{1} = 2T_{\text{del}}(\sqrt{\mathbf{u}}) + \Theta(\log(\mathbf{u})).$$

Set $\ell := \log u$ and $X(\ell) := T_{\text{del}}(2^\ell)$.

Implementation 3: Recursion

$$T_{\text{del}}(\mathbf{u}) = 2T_{\text{del}}(\sqrt{\mathbf{u}}) + T_{\text{min}}(\sqrt{\mathbf{u}}) + 1 = 2T_{\text{del}}(\sqrt{\mathbf{u}}) + \Theta(\log(\mathbf{u})).$$

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$$X(\ell)$$

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$$\begin{aligned} X(\ell) &= T_{\text{del}}(2^\ell) = T_{\text{del}}(u) = 2T_{\text{del}}(\sqrt{u}) + \Theta(\log u) \\ &= 2T_{\text{del}}(2^{\frac{\ell}{2}}) + \Theta(\ell) \end{aligned}$$

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Using Master theorem gives $X(\ell) = \Theta(\ell \log \ell)$, and hence $T_{\text{del}}(u) = \mathcal{O}(\log u \log \log u)$.

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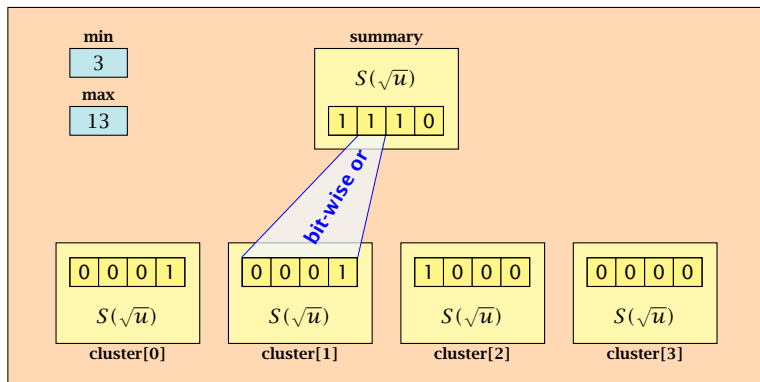
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Using Master theorem gives $X(\ell) = \Theta(\ell \log \ell)$, and hence $T_{\text{del}}(\mathbf{u}) = \mathcal{O}(\log u \log \log u)$.

The same holds for $T_{\text{pred}}(\mathbf{u})$ and $T_{\text{succ}}(\mathbf{u})$.

Implementation 4: van Emde Boas Trees



- ▶ The bit referenced by **min** is **not** set within sub-datastructures.
- ▶ The bit referenced by **max** **is** set within sub-datastructures (if $\text{max} \neq \text{min}$).

Implementation 4: van Emde Boas Trees

Advantages of having max/min pointers:

- ▶ Recursive calls for min and max are constant time.
- ▶ $\text{min} = \text{null}$ means that the data-structure is empty.
- ▶ $\text{min} = \text{max} \neq \text{null}$ means that the data-structure contains exactly one element.
- ▶ We can insert into an empty datastructure in constant time by only setting $\text{min} = \text{max} = x$.
- ▶ We can delete from a data-structure that just contains one element in constant time by setting $\text{min} = \text{max} = \text{null}$.

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Implementation 4: van Emde Boas Trees

Algorithm 38 max()

```
1: return max;
```

Algorithm 39 min()

```
1: return min;
```

- ▶ Constant time.

Implementation 4: van Emde Boas Trees

Algorithm 40 member(x)

1: **if** $x = \min$ **then return** 1; // TRUE

2: **return** cluster[high(x)].member(low(x));

- ▶ $T_{\text{mem}}(u) = T_{\text{mem}}(\sqrt{u}) + 1 \implies T(u) = \mathcal{O}(\log \log u)$.

Implementation 4: van Emde Boas Trees

Algorithm 41 $\text{succ}(x)$

```
1: if  $\text{min} \neq \text{null} \wedge x < \text{min}$  then return  $\text{min}$ ;  
2:  $\text{maxincluster} \leftarrow \text{cluster}[\text{high}(x)].\text{max}()$ ;  
3: if  $\text{maxincluster} \neq \text{null} \wedge \text{low}(x) < \text{maxincluster}$  then  
4:    $\text{offs} \leftarrow \text{cluster}[\text{high}(x)].\text{succ}(\text{low}(x))$ ;  
5:   return  $\text{high}(x) \circ \text{offs}$ ;  
6: else  
7:    $\text{succcluster} \leftarrow \text{summary}.\text{succ}(\text{high}(x))$ ;  
8:   if  $\text{succcluster} = \text{null}$  then return  $\text{null}$ ;  
9:    $\text{offs} \leftarrow \text{cluster}[\text{succcluster}].\text{min}()$ ;  
10:  return  $\text{succcluster} \circ \text{offs}$ ;
```

► $T_{\text{succ}}(u) = T_{\text{succ}}(\sqrt{u}) + 1 \implies T_{\text{succ}}(u) = \mathcal{O}(\log \log u)$.

Implementation 4: van Emde Boas Trees

Algorithm 42 insert(x)

```
1: if min = null then
2:     min =  $x$ ; max =  $x$ ;
3: else
4:     if  $x < \text{min}$  then exchange  $x$  and min;
5:     if cluster[high( $x$ )].min = null; then
6:         summary.insert(high( $x$ ));
7:         cluster[high( $x$ )].insert(low( $x$ ));
8:     else
9:         cluster[high( $x$ )].insert(low( $x$ ));
10:    if  $x > \text{max}$  then max =  $x$ ;
```

- ▶ $T_{\text{ins}}(u) = T_{\text{ins}}(\sqrt{u}) + 1 \implies T_{\text{ins}}(u) = \mathcal{O}(\log \log u)$.

Implementation 4: van Emde Boas Trees

Note that the recursive call in Line 7 takes constant time as the if-condition in Line 5 ensures that we are inserting in an empty sub-tree.

The only non-constant recursive calls are the call in Line 6 and in Line 9. These are mutually exclusive, i.e., only one of these calls will actually occur.

From this we get that $T_{\text{ins}}(u) = T_{\text{ins}}(\sqrt{u}) + 1$.

Implementation 4: van Emde Boas Trees

- ▶ Assumes that x is contained in the structure.

Algorithm 43 delete(x)

```
1: if min = max then
2:     min = null; max = null;
3: else
4:     if  $x$  = min then
5:         firstcluster  $\leftarrow$  summary.min();
6:         offs  $\leftarrow$  cluster[firstcluster].min();
7:          $x \leftarrow$  firstcluster  $\circ$  offs;
8:         min  $\leftarrow$   $x$ ;
9:         cluster[high( $x$ )].delete(low( $x$ ));
```

continued...

Implementation 4: van Emde Boas Trees

- ▶ Assumes that x is contained in the structure.

Algorithm 43 delete(x)

```
1: if min = max then
2:     min = null; max = null;
3: else
4:     if  $x = \text{min}$  then find new minimum
5:          $\text{firstcluster} \leftarrow \text{summary.min}()$ ;
6:          $\text{offs} \leftarrow \text{cluster}[\text{firstcluster}].\text{min}()$ ;
7:          $x \leftarrow \text{firstcluster} \circ \text{offs}$ ;
8:         min  $\leftarrow x$ ;
9:         cluster[high( $x$ )].delete(low( $x$ ));
continued...
```

Implementation 4: van Emde Boas Trees

- ▶ Assumes that x is contained in the structure.

Algorithm 43 delete(x)

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1: if min = max then  
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6:          $offs \leftarrow cluster[firstcluster].min()$ ;  
7:          $x \leftarrow firstcluster \circ offs$ ;  
8:         min  $\leftarrow x$ ;  
9:     cluster[high( $x$ )].delete(low( $x$ ));
```

delete

continued...

Implementation 4: van Emde Boas Trees

Algorithm 43 delete(x)

...continued

```
10:   if cluster[high( $x$ )].min() = null then
11:       summary.delete(high( $x$ ));
12:   if  $x$  = max then
13:       summax  $\leftarrow$  summary.max();
14:       if summax = null then max  $\leftarrow$  min;
15:       else
16:           offs  $\leftarrow$  cluster[summax].max();
17:           max  $\leftarrow$  summax  $\circ$  offs
18:   else
19:       if  $x$  = max then
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```


Implementation 4: van Emde Boas Trees

Algorithm 43 delete(x)

...continued

```
10:   if cluster[high( $x$ )].min() = null then
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```

Implementation 4: van Emde Boas Trees

Algorithm 43 delete(x)

...continued

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```

Implementation 4: van Emde Boas Trees

Algorithm 43 delete(x)

...continued

fix maximum

```
10:   if cluster[high( $x$ )].min() = null then
11:       summary.delete(high( $x$ ));
12:   if  $x$  = max then
13:       summax  $\leftarrow$  summary.max();
14:       if summax = null then max  $\leftarrow$  min;
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16:           offs  $\leftarrow$  cluster[summax].max();
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20:           offs  $\leftarrow$  cluster[high( $x$ )].max();
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```

Implementation 4: van Emde Boas Trees

Note that only one of the possible recursive calls in Line 9 and Line 11 in the deletion-algorithm may take non-constant time.

To see this observe that the call in Line 11 only occurs if the cluster where x was deleted is now empty. But this means that the call in Line 9 deleted the last element in $\text{cluster}[\text{high}(x)]$. Such a call only takes constant time.

Hence, we get a recurrence of the form

$$T_{\text{del}}(u) = T_{\text{del}}(\sqrt{u}) + c .$$

This gives $T_{\text{del}}(u) = \mathcal{O}(\log \log u)$.

9 van Emde Boas Trees

Space requirements:

- ▶ The space requirement fulfills the recurrence

$$S(u) = (\sqrt{u} + 1)S(\sqrt{u}) + \mathcal{O}(\sqrt{u}) .$$

- ▶ Note that we cannot solve this recurrence by the Master theorem as the branching factor is not constant.
- ▶ One can show by induction that the space requirement is $S(u) = \mathcal{O}(u)$. Exercise.

10 Union Find

Union Find Data Structure \mathcal{P} : Maintains a partition of **disjoint** sets over elements.

- ▶ \mathcal{P} . **makeset**(x): Given an element x , adds x to the data-structure and creates a singleton set that contains only this element. Returns a locator/handle for x in the data-structure.
- ▶ \mathcal{P} . **find**(x): Given a handle for an element x ; find the set that contains x . Returns a representative/identifier for this set.
- ▶ \mathcal{P} . **union**(x, y): Given two elements x , and y that are currently in sets S_x and S_y , respectively, the function replaces S_x and S_y by $S_x \cup S_y$ and returns an identifier for the new set.

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Applications:

- ▶ Keep track of the connected components of a dynamic graph that changes due to insertion of nodes and edges.
- ▶ Kruskals Minimum Spanning Tree Algorithm

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10 Union Find

Algorithm 44 Kruskal-MST($G = (V, E), w$)

```
1:  $A \leftarrow \emptyset$ ;  
2: for all  $v \in V$  do  
3:    $v.\text{set} \leftarrow \mathcal{P}.\text{makeset}(v.\text{label})$   
4: sort edges in non-decreasing order of weight  $w$   
5: for all  $(u, v) \in E$  in non-decreasing order do  
6:   if  $\mathcal{P}.\text{find}(u.\text{set}) \neq \mathcal{P}.\text{find}(v.\text{set})$  then  
7:      $A \leftarrow A \cup \{(u, v)\}$   
8:      $\mathcal{P}.\text{union}(u.\text{set}, v.\text{set})$ 
```

List Implementation

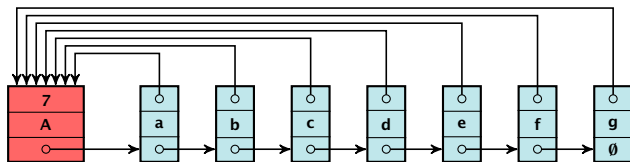
- ▶ The elements of a set are stored in a list; each node has a backward pointer to the head.
- ▶ The head of the list contains the identifier for the set and a field that stores the **size** of the set.



- ▶ `makeset(x)` can be performed in constant time.
- ▶ `find(x)` can be performed in constant time.

List Implementation

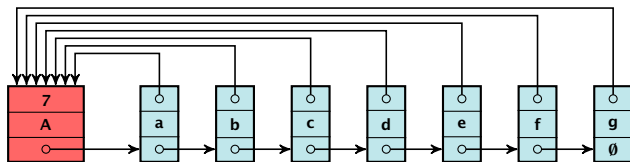
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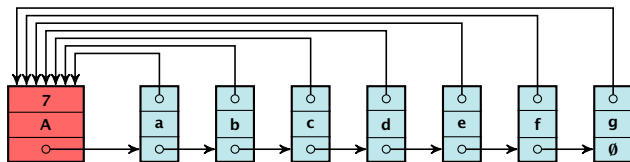
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List Implementation

union(x, y)

- ▶ Determine sets S_x and S_y .
- ▶ Traverse the smaller list (say S_y), and change all backward pointers to the head of list S_y .
- ▶ Insert list S_y at the head of S_x .
- ▶ Adjust the size-field of list S_x .
- ▶ Time: $\min\{|S_x|, |S_y|\}$.

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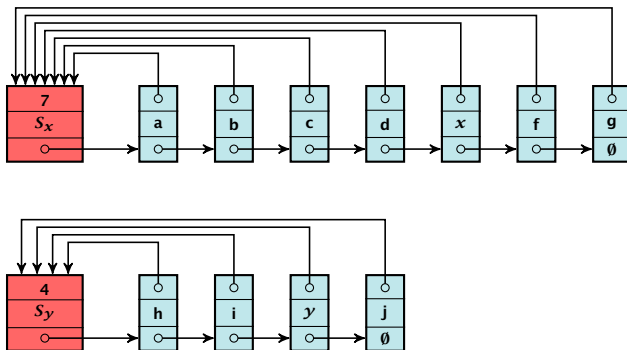
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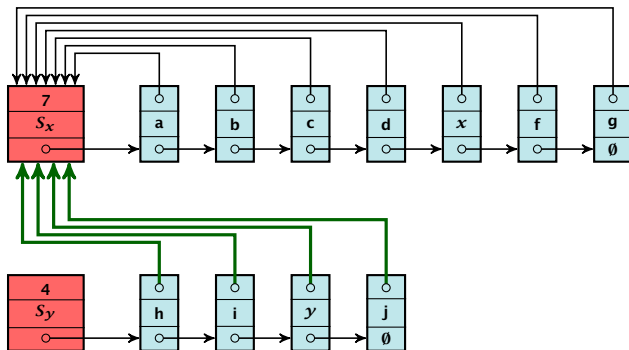
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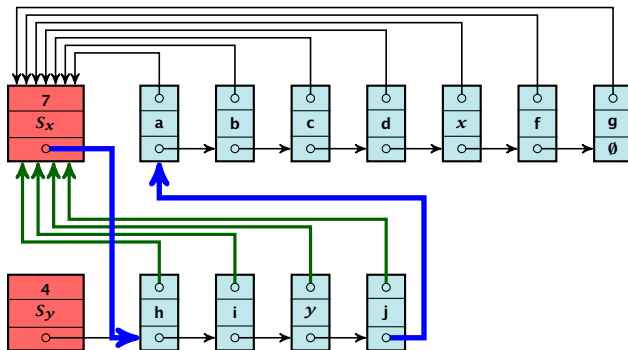
List Implementation



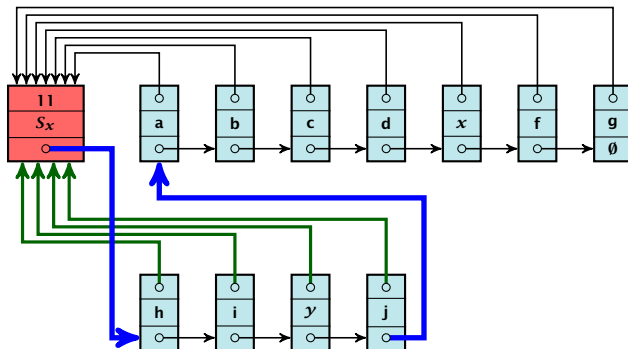
List Implementation



List Implementation



List Implementation



List Implementation

Running times:

- ▶ $\text{find}(x)$: constant
- ▶ $\text{makeset}(x)$: constant
- ▶ $\text{union}(x, y)$: $\mathcal{O}(n)$, where n denotes the number of elements contained in the set system.

List Implementation

Lemma 35

The list implementation for the ADT union find fulfills the following amortized time bounds:

- ▶ $\text{find}(x): \mathcal{O}(1)$.
- ▶ $\text{makeset}(x): \mathcal{O}(\log n)$.
- ▶ $\text{union}(x, y): \mathcal{O}(1)$.

The Accounting Method for Amortized Time Bounds

- ▶ There is a bank account for every element in the data structure.
- ▶ Initially the balance on all accounts is zero.
- ▶ Whenever for an operation the amortized time bound exceeds the actual cost, the difference is credited to some bank accounts of elements involved.
- ▶ Whenever for an operation the actual cost exceeds the amortized time bound, the difference is charged to bank accounts of some of the elements involved.
- ▶ If we can find a charging scheme that guarantees that balances always stay positive the amortized time bounds are proven.

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List Implementation

- ▶ For an operation whose actual cost exceeds the amortized cost we charge the **excess** to the elements involved.
- ▶ In total we will charge at most $\mathcal{O}(\log n)$ to an element (regardless of the request sequence).
- ▶ For each element a makeset operation occurs as the first operation involving this element.
- ▶ We inflate the amortized cost of the makeset-operation to $\Theta(\log n)$, i.e., at this point we fill the bank account of the element to $\Theta(\log n)$.
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List Implementation

makeSet(x) : The actual cost is $\mathcal{O}(1)$. Due to the cost inflation the amortized cost is $\mathcal{O}(\log n)$.

find(x) : For this operation we define the amortized cost and the actual cost to be the same. Hence, this operation does not change any accounts. Cost: $\mathcal{O}(1)$.

union(x, y):

Let S_x and S_y be the sets of nodes in the rank r of x and y , respectively.

Case 1: $|S_x| \leq |S_y|$. The actual cost is $\mathcal{O}(|S_x| \cdot \log n)$. The amortized cost is $\mathcal{O}(|S_x|)$.

Case 2: $|S_x| > |S_y|$. The amortized cost is $\mathcal{O}(|S_x|)$. The actual cost is $\mathcal{O}(|S_x| \cdot \log n)$.

Since $|S_x| \leq |S_y|$, the actual cost is the smaller one. Hence, the amortized cost is $\mathcal{O}(|S_x|)$.

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- ▶ If $S_x = S_y$ the cost is constant; no bank accounts change.
- ▶ Otw. the actual cost is $\mathcal{O}(\min\{|S_x|, |S_y|\})$.
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Lemma 36

An element is charged at most $\lfloor \log_2 n \rfloor$ times, where n is the total number of elements in the set system.

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Whenever an element x is charged the number of elements in x 's set doubles. This can happen at most $\lfloor \log n \rfloor$ times. \square

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Implementation via Trees

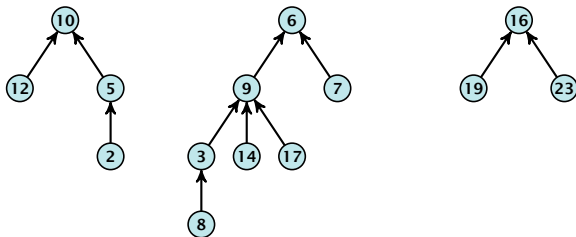
- ▶ Maintain nodes of a set in a tree.
- ▶ The root of the tree is the label of the set.
- ▶ Only pointer to parent exists; we cannot list all elements of a given set.
- ▶ Example:



Set system $\{2, 5, 10, 12\}$, $\{3, 6, 7, 8, 9, 14, 17\}$, $\{16, 19, 23\}$.

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Implementation via Trees

makeset(x)

- ▶ Create a singleton tree. Return pointer to the root.
- ▶ Time: $\mathcal{O}(1)$.

find(x)

- ▶ Start at element x in the tree, and repeatedly update x to be its parent.
- ▶ Time: $\mathcal{O}(n)$, where n is the depth of element x in the tree.

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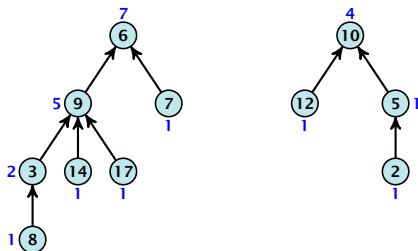
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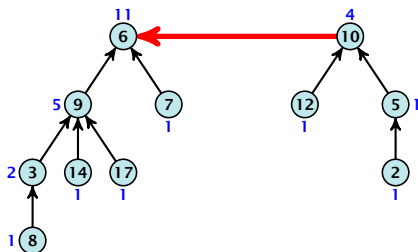


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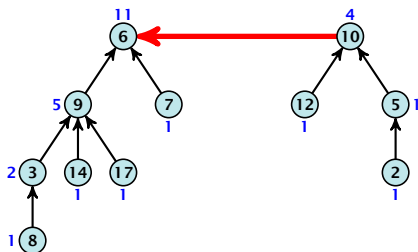


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- ▶ Time: constant for $\text{link}(a, b)$ plus two find-operations.

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The running time (non-amortized!!!) for $\text{find}(x)$ is $\mathcal{O}(\log n)$.

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find(x):

- ▶ Go upward until you find the root.
- ▶ Re-attach all visited nodes as children of the root.
- ▶ Speeds up successive find-operations.



Time complexity of find is $O(\alpha(n))$ (inverse Ackermann function)

Time complexity of union is $O(\alpha(n))$

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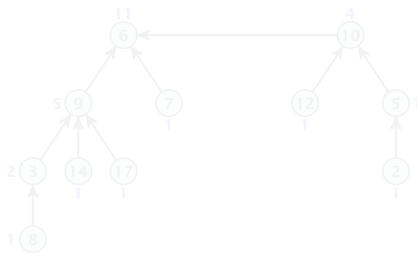
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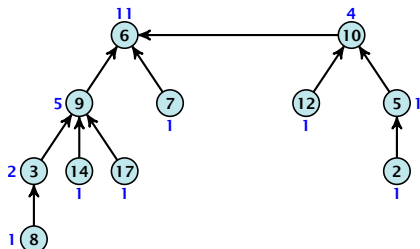
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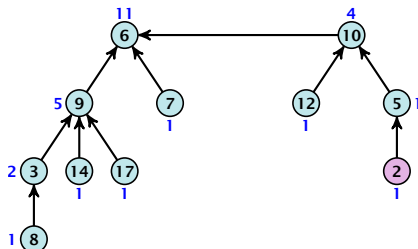


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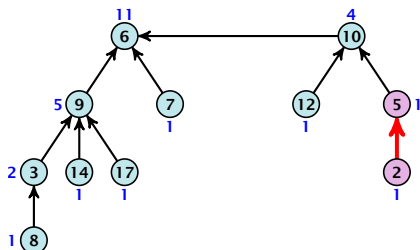


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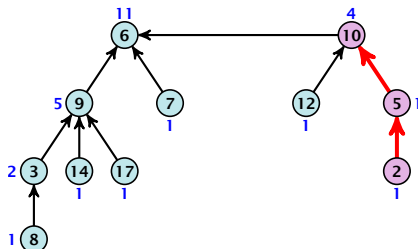


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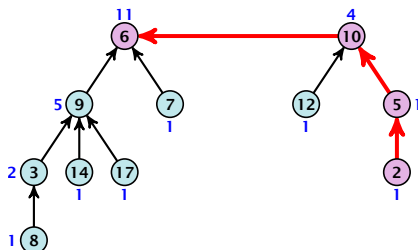


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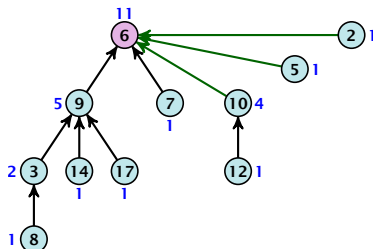


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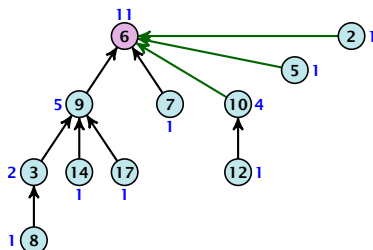


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Amortized Analysis

Definitions:

$\text{size}(v)$: The number of nodes that were in the sub-tree rooted at v when v became the child of another node (or if v is the root, the number of nodes of T is the size).

$\text{rank}(v) = \lceil \log(\text{size}(v)) \rceil$.

$\text{rank}(v) \leq \text{depth}(v) \leq 2 \cdot \text{rank}(v)$.

Lemma 38

The rank of a parent must be strictly larger than the rank of a child.

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There are at most $n/2^s$ nodes of rank s .

Proof.

- Let v be a node of rank s . It is the root of a subtree of 2^{s-1} nodes.
- Each of these nodes is the root of a subtree of 2^{s-2} nodes.
- Each of these nodes is the root of a subtree of 2^{s-3} nodes.
- This holds because the rank of each of the nodes of the subtree is at least $s-1$.
- This subtree contains at least the nodes v and v 's parent.
- Each node of rank s is the root of a subtree of 2^{s-1} nodes of different ranks. □

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- ▶ Let's say a node v **sees** the rank s node x if v is in x 's sub-tree at the time that x becomes a child.
- ▶ A node v sees at most one node of rank s during the running time of the algorithm.
- ▶ This holds because the rank-sequence of the roots of the different trees that contains v during the running time of the algorithm is a strictly increasing sequence.
- ▶ Hence, every node *sees* at most one rank s node, but every rank s node is seen by at least 2^s different nodes. □

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- ▶ Hence, every node *sees* at most one rank s node, but every rank s node is seen by at least 2^s different nodes. □

Amortized Analysis

We define

$$\text{tow}(i) := \begin{cases} 1 & \text{if } i = 0 \\ 2^{\text{tow}(i-1)} & \text{otw.} \end{cases}$$

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Theorem 40

Union find with path compression fulfills the following amortized running times:

- ▶ $\text{makeset}(x) : \mathcal{O}(\log^*(n))$
- ▶ $\text{find}(x) : \mathcal{O}(\log^*(n))$
- ▶ $\text{union}(x, y) : \mathcal{O}(\log^*(n))$

Amortized Analysis

In the following we assume $n \geq 3$.

rank-group:

- A node with rank r belongs to the rank-group (r) .
- The rank-group (r) contains only nodes with rank r or $r+1$.
- A rank-group (r) contains at most 2^{r-1} nodes.
- The maximum number of rank-groups is $\log_2 n + 1$.
- The total number of nodes is at most $\sum_{r=0}^{\log_2 n} 2^{r-1} = 2^{\log_2 n} = n$.

Amortized Analysis

In the following we assume $n \geq 3$.

rank-group:

- ▶ A node with rank $\text{rank}(v)$ is in **rank group** $\log^*(\text{rank}(v))$.
- ▶ The rank-group $g = 0$ contains only nodes with rank 0 or rank 1.
- ▶ A rank group $g \geq 1$ contains ranks $\text{tow}(g-1) + 1, \dots, \text{tow}(g)$.
- ▶ The maximum non-empty rank group is $\log^*(\lfloor \log n \rfloor) \leq \log^*(n) - 1$ (which holds for $n \geq 3$).
- ▶ Hence, the total number of rank-groups is at most $\log^* n$.

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- ▶ Hence, the total number of rank-groups is at most $\log^* n$.

Amortized Analysis

Accounting Scheme:

• Create an account for every find-operation.

• Create an account for every node v .

The cost for a find-operation is equal to the length of the path traversed. We charge the cost for going from v to $\text{parent}[v]$ as follows:

• If $\text{parent}[v]$ is the root we charge the cost to the root's account.

• Otherwise:

• If the credit-number of $\text{node}[v]$ is the same as that of $\text{node}[\text{parent}[v]]$ (before starting path-compression) we charge the cost to the node-account of v .

• Otherwise we charge the cost to the root's account.

Amortized Analysis

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The cost for a find-operation is equal to the length of the path traversed. We charge the cost for going from v to $\text{parent}[v]$ as follows:

- ▶ if $\text{parent}[v]$ is the root we charge the cost to the root's account
- ▶ if $\text{parent}[v]$ is not the root we charge the cost to the account of the grand-parent of v (before storing path compression) we charge the cost to the node's account of v (if v is not the root)
- ▶ otherwise we charge the cost to the grand-parent's account

Amortized Analysis

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The cost for a find-operation is equal to the length of the path traversed. We charge the cost for going from v to $\text{parent}[v]$ as follows:

▶ if $\text{parent}[v]$ is the root we charge the cost to the account of the find-operation

▶ if the root is not the root of v , then we charge the cost to the account of the root of v (before working path compression)

▶ if the root is not the root of v , then we charge the cost to the account of the root of v (after working path compression)

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- ▶ create an account for every find-operation
- ▶ create an account for every node v

The cost for a find-operation is equal to the length of the path traversed. We charge the cost for going from v to $\text{parent}[v]$ as follows:

- ▶ If $\text{parent}[v]$ is the root we charge the cost to the find-account.
- ▶ If the group-number of $\text{rank}(v)$ is the same as that of $\text{rank}(\text{parent}[v])$ (before starting path compression) we charge the cost to the node-account of v .
- ▶ Otherwise we charge the cost to the find-account.

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- ▶ If the group-number of $\text{rank}(v)$ is the same as that of $\text{rank}(\text{parent}[v])$ (before starting path compression) we charge the cost to the node-account of v .
- ▶ Otherwise we charge the cost to the find-account.

Observations:

• The root node is charged at most by $\log_2(n)$ times when traversing the tree (regardless of how many times the root is visited).

• The grandchild is charged at most once per assigned. The grandchild of the parent is only charged once.

• The grandchild of the grandparent will be in a larger subtree \rightarrow it will never be charged again.

• The total charge made to a node in rank group p is at most $\log_2(n) \cdot 2^p$.

Observations:

- ▶ A find-account is charged at most $\log^*(n)$ times (once for the root and at most $\log^*(n) - 1$ times when increasing the rank-group).
- ▶ After a node v is charged its parent-edge is re-assigned. The rank of the parent strictly increases.
- ▶ After some charges to v the parent will be in a larger rank-group. $\Rightarrow v$ will **never** be charged again.
- ▶ The total charge made to a node in rank-group g is at most $\text{tow}(g) - \text{tow}(g - 1) \leq \text{tow}(g)$.

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What is the total charge made to nodes?

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Hence,

$$\sum_g n(g) \text{tow}(g) \leq n(0) \text{tow}(0) + \sum_{g \geq 1} n(g) \text{tow}(g)$$

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Hence,

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Amortized Analysis

Without loss of generality we can assume that all makeset-operations occur at the start.

This means if we inflate the cost of makeset to $\log^* n$ and add this to the node account of v then the balances of all node accounts will sum up to a positive value (this is sufficient to obtain an amortized bound).

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The analysis is not tight. In fact it has been shown that the amortized time for the union-find data structure with path compression is $\mathcal{O}(\alpha(m, n))$, where $\alpha(m, n)$ is the inverse Ackermann function which grows a lot lot slower than $\log^* n$. (Here, we consider the average running time of m operations on at most n elements).

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There is also a lower bound of $\Omega(\alpha(m, n))$.

$$A(x, y) = \begin{cases} y + 1 & \text{if } x = 0 \\ A(x - 1, 1) & \text{if } y = 0 \\ A(x - 1, A(x, y - 1)) & \text{otw.} \end{cases}$$

$$\alpha(m, n) = \min\{i \geq 1 : A(i, \lfloor m/n \rfloor) \geq \log n\}$$

- ▶ $A(0, y) = y + 1$
- ▶ $A(1, y) = y + 2$
- ▶ $A(2, y) = 2y + 3$
- ▶ $A(3, y) = 2^{y+3} - 3$
- ▶ $A(4, y) = \underbrace{2^{2^{2^{\dots}}}}_{y+3 \text{ times}} - 3$

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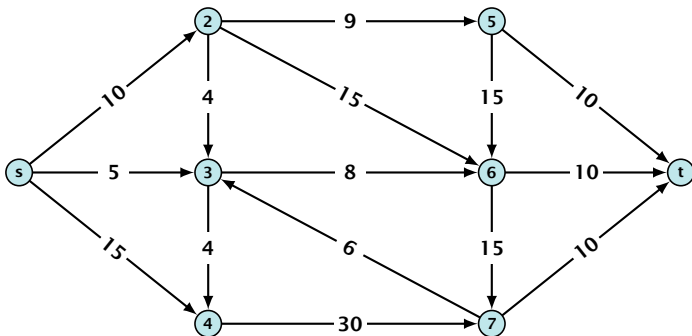
Part IV

Flows and Cuts

11 Introduction

Flow Network

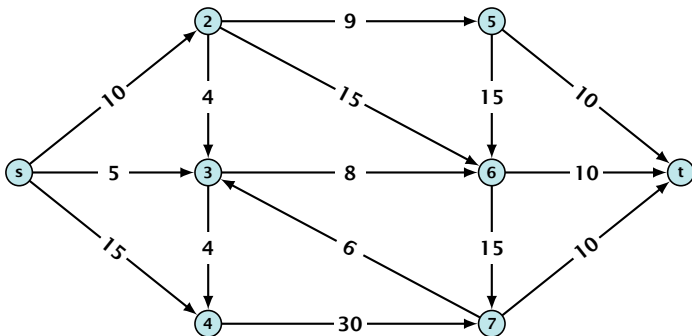
- ▶ directed graph $G = (V, E)$; edge capacities $c(e)$
- ▶ two special nodes: source s ; target t ;
- ▶ no edges entering s or leaving t ;
- ▶ at least for now: no parallel edges;



11 Introduction

Flow Network

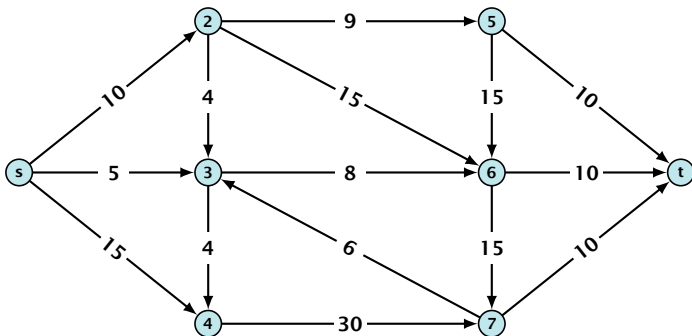
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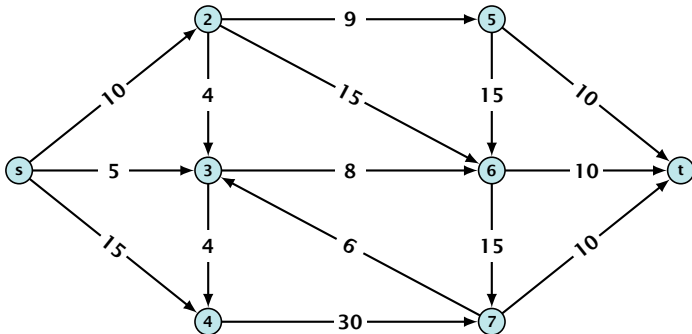
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Cuts

Definition 41

An (s, t) -cut in the graph G is given by a set $A \subset V$ with $s \in A$ and $t \in V \setminus A$.

Cuts

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Definition 42

The **capacity** of a cut A is defined as

$$\text{cap}(A, V \setminus A) := \sum_{e \in \text{out}(A)} c(e) ,$$

where $\text{out}(A)$ denotes the set of edges of the form $A \times V \setminus A$ (i.e. edges leaving A).

Cuts

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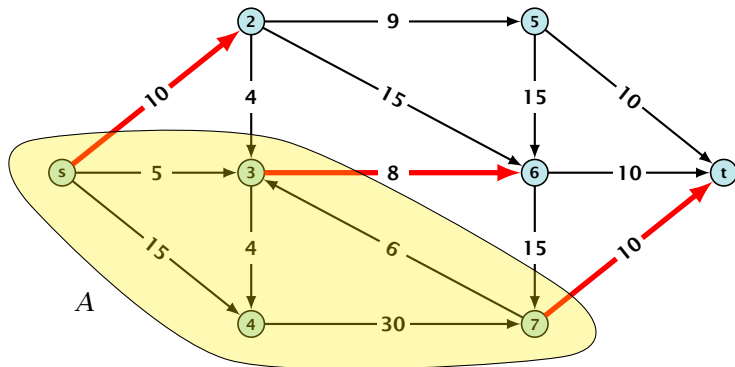
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where $\text{out}(A)$ denotes the set of edges of the form $A \times V \setminus A$ (i.e. edges leaving A).

Minimum Cut Problem: Find an (s, t) -cut with minimum capacity.

Cuts

Example 43



The capacity of the cut is $\text{cap}(A, V \setminus A) = 28$.

Flows

Definition 44

An (s, t) -flow is a function $f : E \mapsto \mathbb{R}^+$ that satisfies

1. For each edge e

$$0 \leq f(e) \leq c(e) .$$

(capacity constraints)

2. For each $v \in V \setminus \{s, t\}$

$$\sum_{e \in \text{out}(v)} f(e) = \sum_{e \in \text{into}(v)} f(e) .$$

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Flows

Definition 45

The value of an (s, t) -flow f is defined as

$$\text{val}(f) = \sum_{e \in \text{out}(s)} f(e) .$$

Maximum Flow Problem: Find an (s, t) -flow with maximum value.

Flows

Definition 45

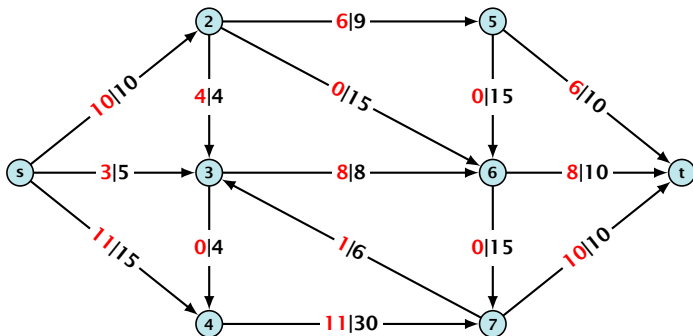
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Maximum Flow Problem: Find an (s, t) -flow with maximum value.

Flows

Example 46



The value of the flow is $\text{val}(f) = 24$.

Lemma 47 (Flow value lemma)

Let f a flow, and let $A \subseteq V$ be an (s, t) -cut. Then the *net-flow* across the cut is equal to the amount of flow leaving s , i.e.,

$$\text{val}(f) = \sum_{e \in \text{out}(A)} f(e) - \sum_{e \in \text{into}(A)} f(e) .$$

Proof.

$\text{val}(f)$

Proof.

$$\text{val}(f) = \sum_{e \in \text{out}(s)} f(e)$$

Proof.

$$\begin{aligned}\text{val}(f) &= \sum_{e \in \text{out}(s)} f(e) \\ &= \sum_{e \in \text{out}(s)} f(e) + \sum_{v \in A \setminus \{s\}} \left(\sum_{e \in \text{out}(v)} f(e) - \sum_{e \in \text{in}(v)} f(e) \right)\end{aligned}$$

Proof.

$$\begin{aligned}\text{val}(f) &= \sum_{e \in \text{out}(s)} f(e) && = 0 \\ &= \sum_{e \in \text{out}(s)} f(e) + \sum_{v \in A \setminus \{s\}} \left(\sum_{e \in \text{out}(v)} f(e) - \sum_{e \in \text{in}(v)} f(e) \right)\end{aligned}$$

Proof.

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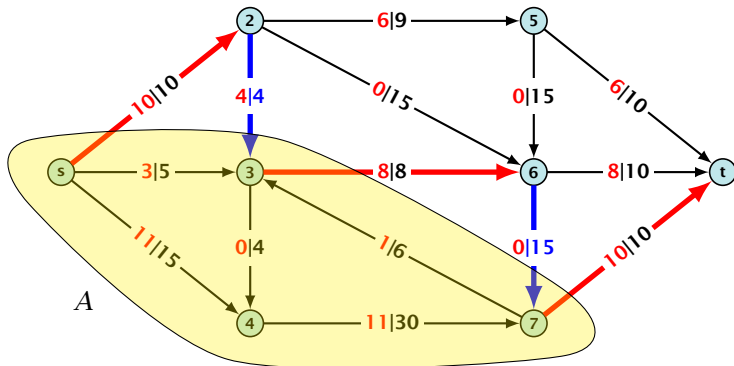
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The last equality holds since every edge with both end-points in A contributes negatively as well as positively to the sum in line 2. The only edges whose contribution doesn't cancel out are edges leaving or entering A .



Example 48



Corollary 49

Let f be an (s, t) -flow and let A be an (s, t) -cut, such that

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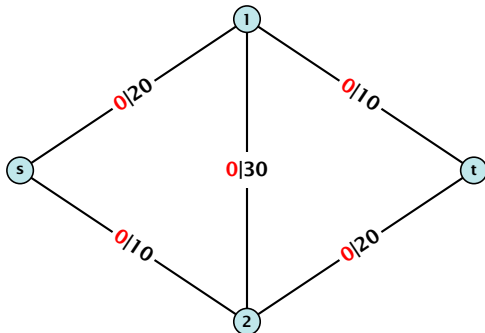
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12 Augmenting Path Algorithms

Greedy-algorithm:

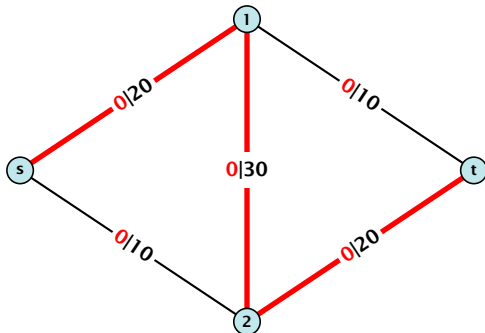
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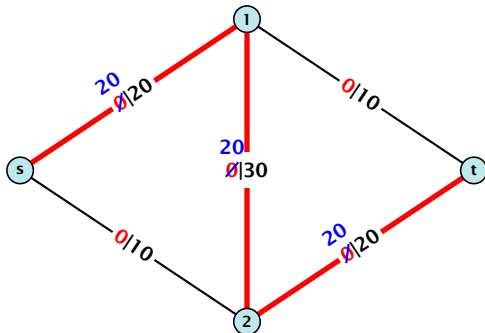
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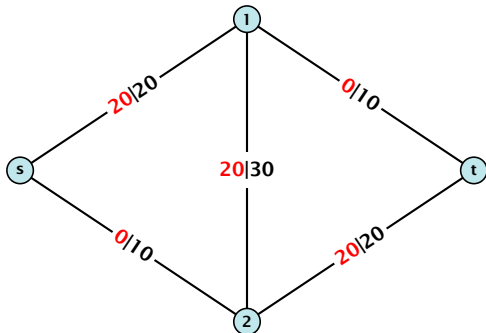
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The Residual Graph

From the graph $G = (V, E, c)$ and the current flow f we construct an auxiliary graph $G_f = (V, E_f, c_f)$ (the residual graph):

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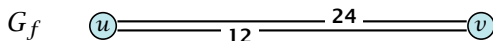
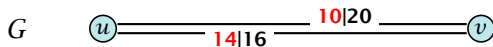
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Augmenting Path Algorithm

Definition 50

An **augmenting path** with respect to flow f , is a path in the auxiliary graph G_f that contains only edges with non-zero capacity.

Algorithm 45 FordFulkerson($G = (V, E, c)$)

- 1: Initialize $f(e) \leftarrow 0$ for all edges.
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Theorem 51

A flow f is a maximum flow iff there are no augmenting paths.

Theorem 52

The value of a maximum flow is equal to the value of a minimum cut.

Proof.

Let f be a flow. The following are equivalent:

1. There exists a cut A, B such that $\text{val}(f) = \text{cap}(A, B)$.
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Let f be a flow. The following are equivalent:

1. There is no augmenting path. f is a maximum flow. $v(f) = \text{cap}(f)$.

2. There is a minimum cut C with $v(f) = \text{cap}(C)$.

3. $v(f) = \text{cap}(C)$ for some cut C .



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This we already showed.

2. \Rightarrow 3.

If there were an augmenting path, we could improve the flow.
Contradiction.

3. \Rightarrow 1.

f must be a flow with no augmenting paths.

Let A be the set of vertices reachable from s in the residual network, along non-saturated capacity edges.

\Rightarrow Since there is no augmenting path we have $t \notin A$ and $t \in A$.

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Let C be the set of saturated capacity edges.

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This finishes the proof.

Here the first equality uses the flow value lemma, and the second exploits the fact that the flow along incoming edges must be 0 as the residual graph does not have edges leaving A .

Analysis

Assumption:

All capacities are integers between 1 and C .

Invariant:

Every flow value $f(e)$ and every residual capacity $c_f(e)$ remains integral throughout the algorithm.

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The algorithm terminates in at most $\text{val}(f^) \leq nC$ iterations, where f^* denotes the maximum flow. Each iteration can be implemented in time $\mathcal{O}(m)$. This gives a total running time of $\mathcal{O}(nmC)$.*

Theorem 54

If all capacities are integers, then there exists a maximum flow for which every flow value $f(e)$ is integral.

Lemma 53

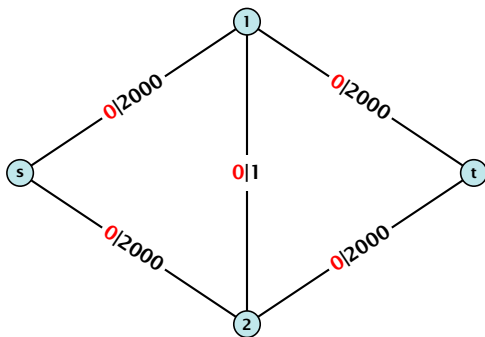
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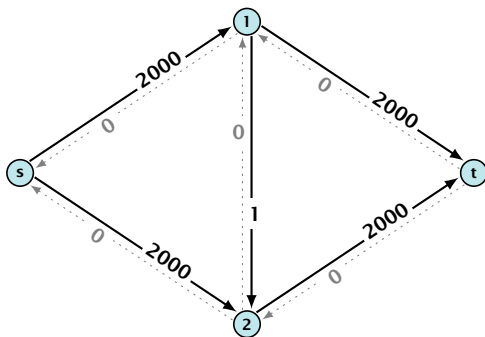
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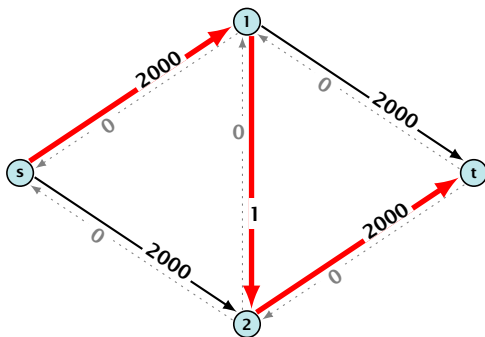


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Can we tweak the algorithm so that the running time is polynomial in the input length?

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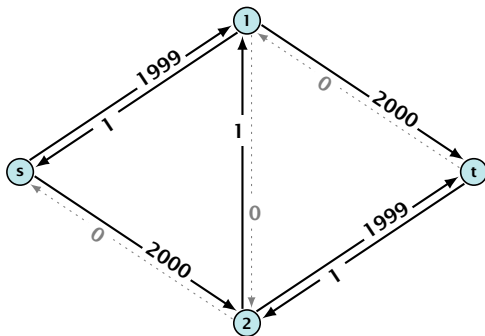


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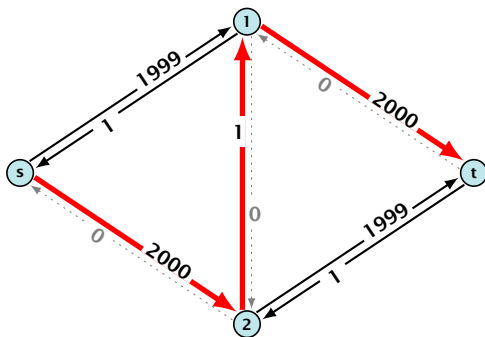


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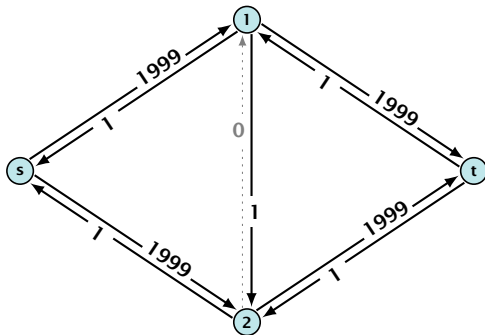


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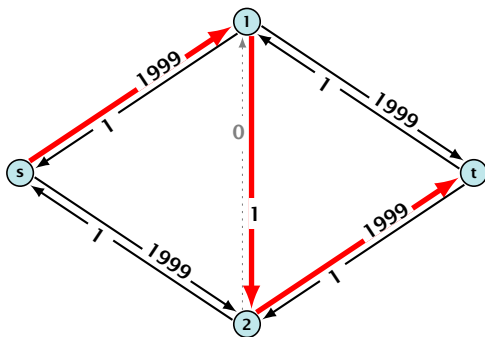


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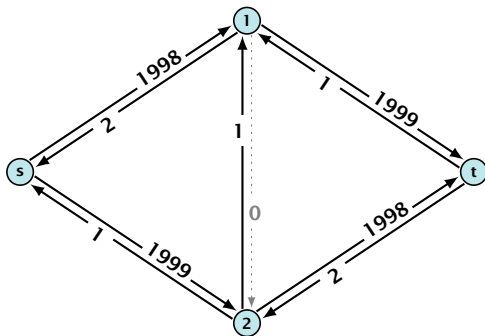


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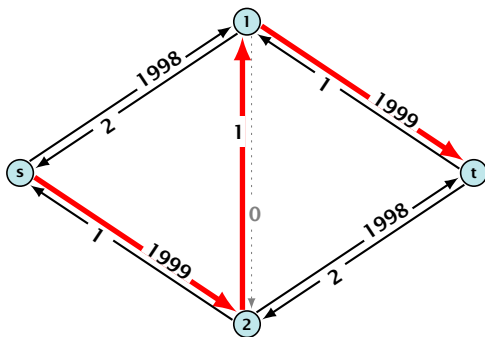


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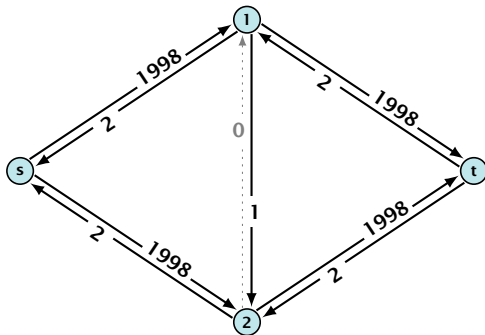


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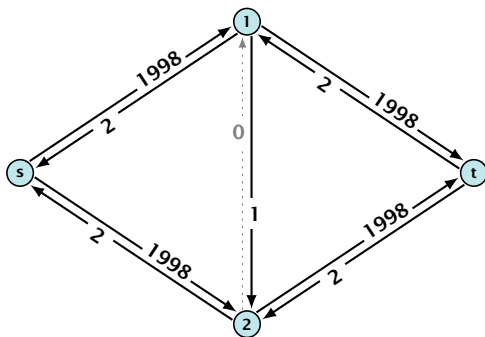


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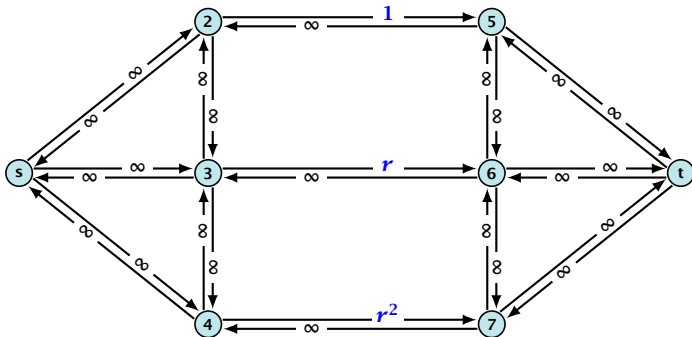


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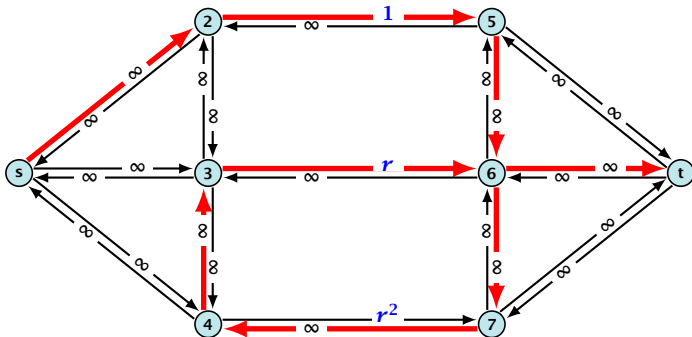
A Pathological Input

Let $r = \frac{1}{2}(\sqrt{5} - 1)$. Then $r^{n+2} = r^n - r^{n+1}$.



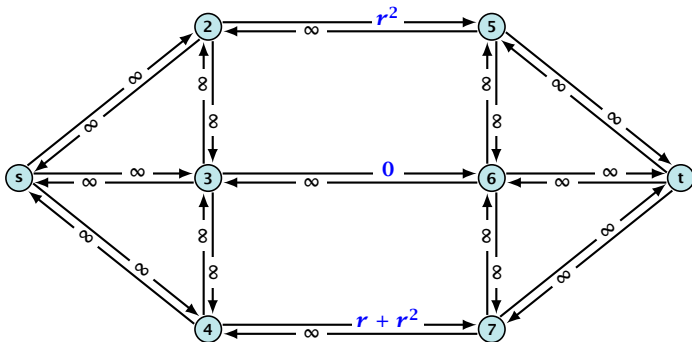
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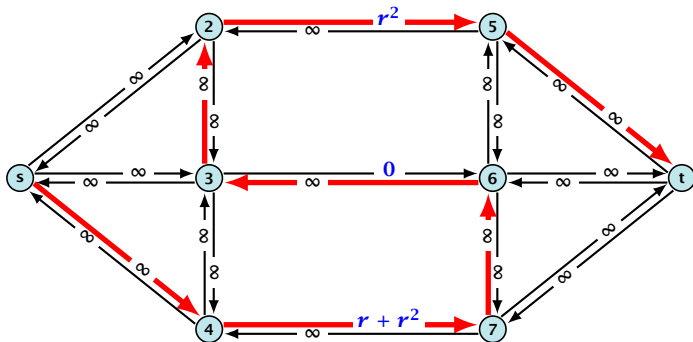
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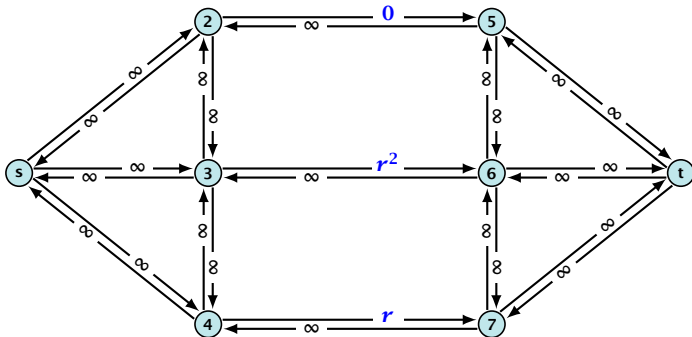
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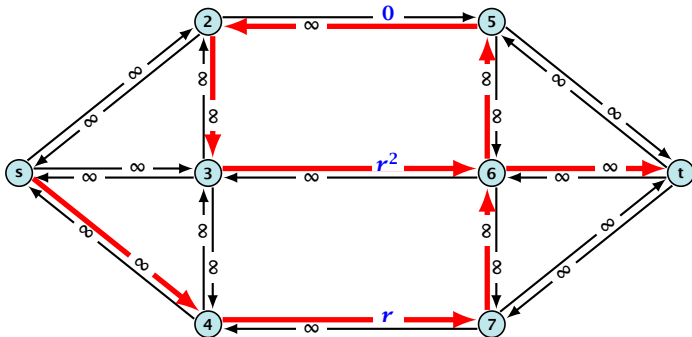
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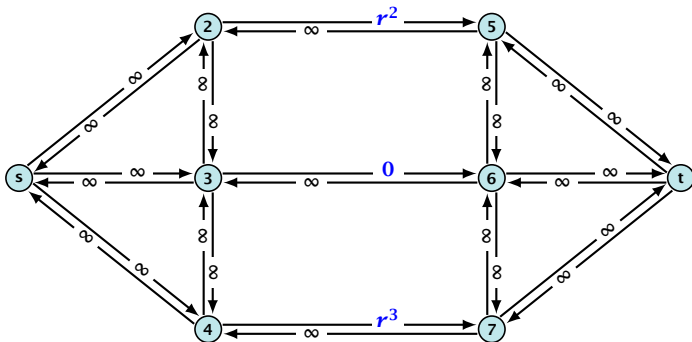
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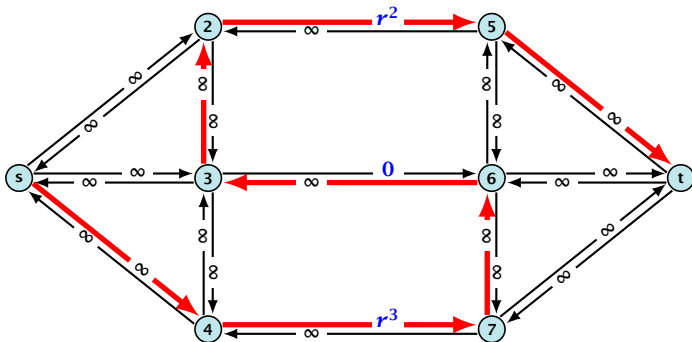
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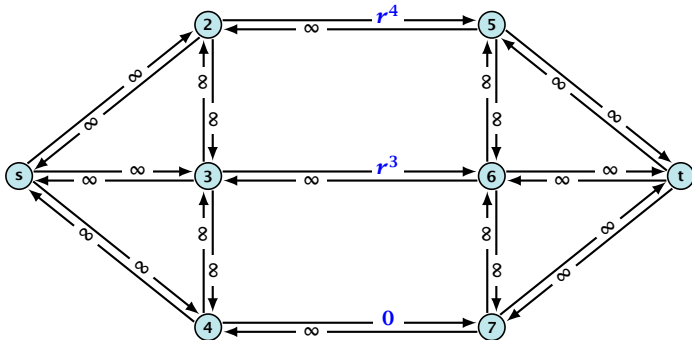
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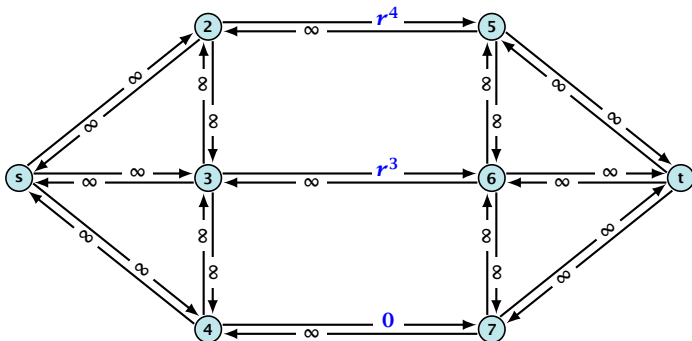
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Running time may be infinite!!!

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- ▶ Choose the shortest augmenting path.

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Overview: Shortest Augmenting Paths

Lemma 55

The length of the shortest augmenting path never decreases.

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After at most $O(m)$ augmentations, the length of the shortest augmenting path strictly increases.

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These two lemmas give the following theorem:

Theorem 57

The shortest augmenting path algorithm performs at most $\mathcal{O}(mn)$ augmentations. This gives a running time of $\mathcal{O}(m^2n)$.

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• We can find the shortest augmenting paths in time $\mathcal{O}(m)$ via BFS.

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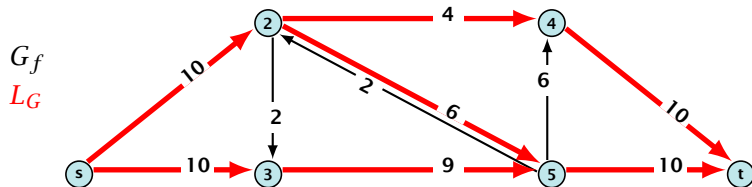
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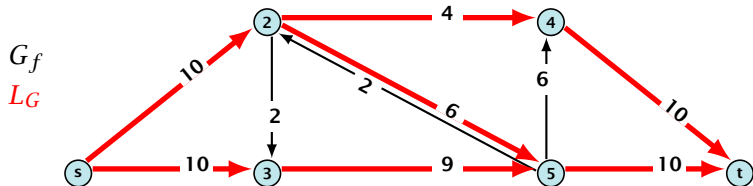
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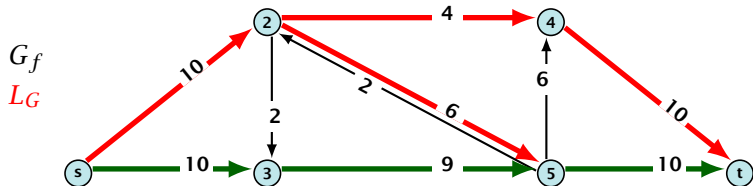


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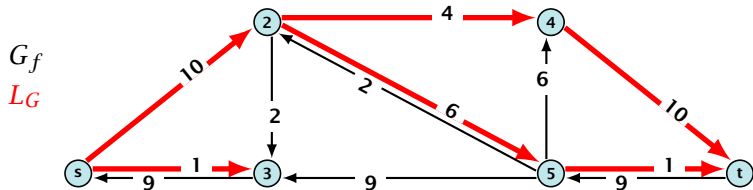


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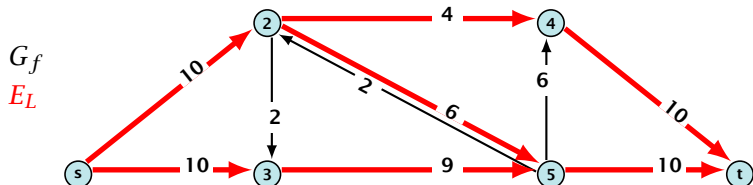
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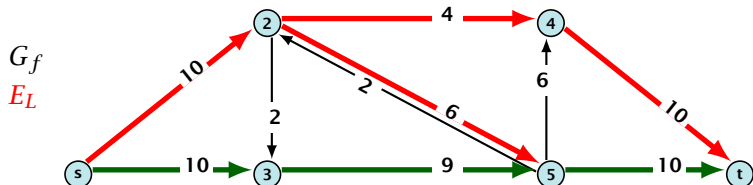
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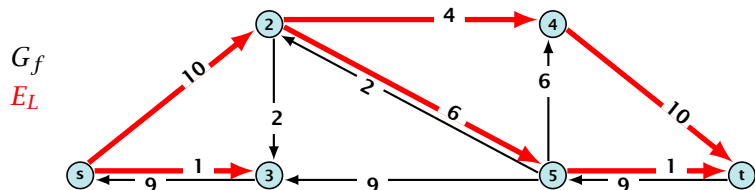
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There exist networks with $m = \Theta(n^2)$ that require $\mathcal{O}(mn)$ augmentations, when we restrict ourselves to only augment along shortest augmenting paths.

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Initializing E_L for the phase takes time $\mathcal{O}(m)$.

The total cost for searching for augmenting paths during a phase is at most $\mathcal{O}(mn)$, since every search (successful (i.e., reaching t) or unsuccessful) decreases the number of edges in E_L and takes time $\mathcal{O}(n)$.

The total cost for performing an augmentation **during** a phase is only $\mathcal{O}(n)$. For every edge in the augmenting path one has to update the residual graph G_f and has to check whether the edge is still in E_L for the next search.

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- ▶ Choose path with sufficiently large bottleneck capacity.
- ▶ Choose the shortest augmenting path.

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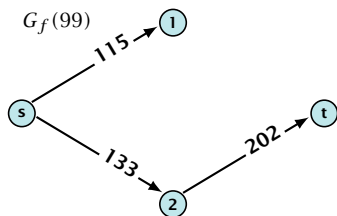
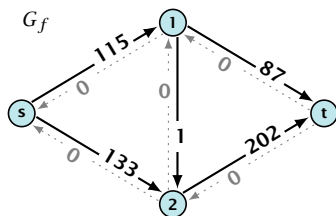
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Capacity Scaling

Algorithm 46 maxflow(G, s, t, c)

```
1: foreach  $e \in E$  do  $f_e \leftarrow 0$ ;  
2:  $\Delta \leftarrow 2^{\lceil \log_2 C \rceil}$   
3: while  $\Delta \geq 1$  do  
4:    $G_f(\Delta) \leftarrow \Delta$ -residual graph  
5:   while there is augmenting path  $P$  in  $G_f(\Delta)$  do  
6:      $f \leftarrow \text{augment}(f, c, P)$   
7:      $\text{update}(G_f(\Delta))$   
8:    $\Delta \leftarrow \Delta/2$   
9: return  $f$ 
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Theorem 63

We need $\mathcal{O}(m \log C)$ augmentations. The algorithm can be implemented in time $\mathcal{O}(m^2 \log C)$.

Preflows

Definition 64

An (s, t) -preflow is a function $f : E \mapsto \mathbb{R}^+$ that satisfies

1. For each edge e

$$0 \leq f(e) \leq c(e) .$$

(capacity constraint)

2. For each $v \in V \setminus \{s, t\}$

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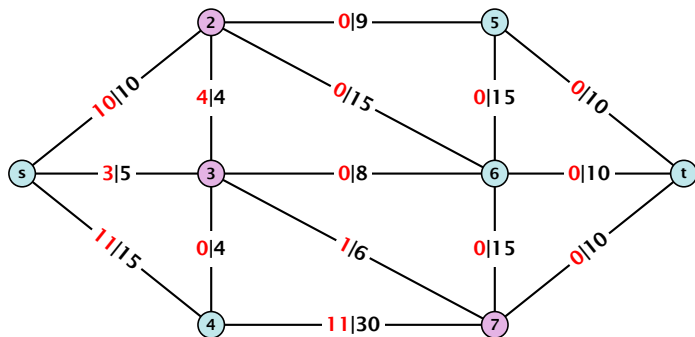
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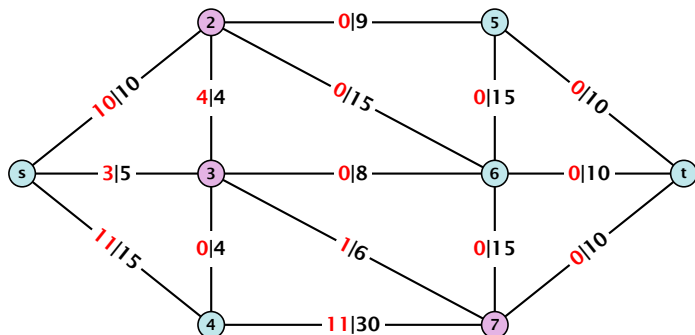
Preflows

Example 65



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A node that has $\sum_{e \in \text{out}(v)} f(e) < \sum_{e \in \text{into}(v)} f(e)$ is called an **active node**.

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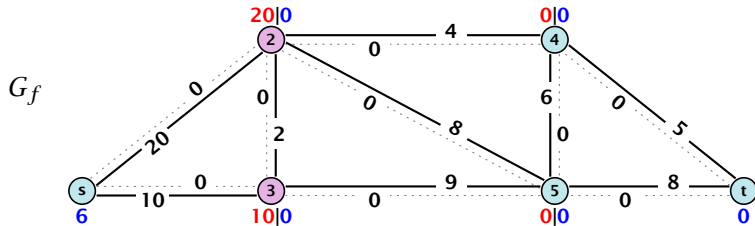
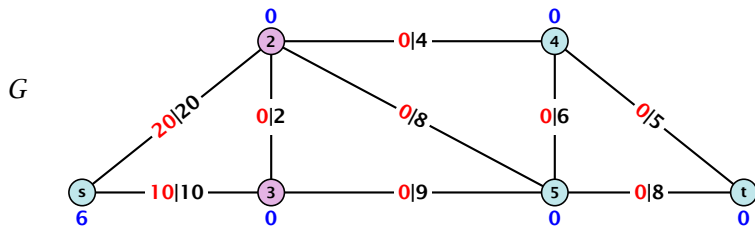
A **labelling** is a function $\ell : V \rightarrow \mathbb{N}$. It is **valid** for preflow f if

- ▶ $\ell(u) \leq \ell(v) + 1$ for all edges in the residual graph G_f (only non-zero capacity edges!!!)
- ▶ $\ell(s) = n$
- ▶ $\ell(t) = 0$

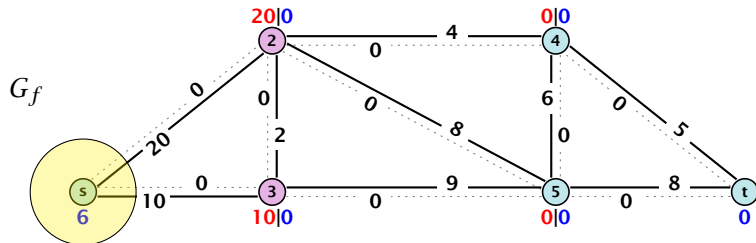
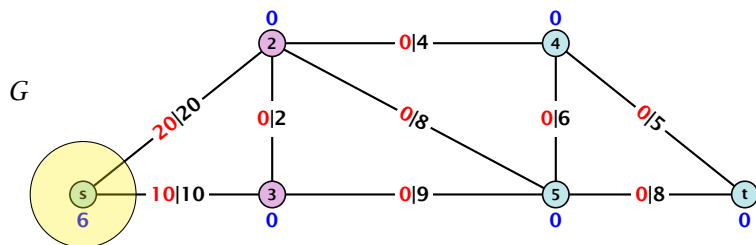
Intuition:

The labelling can be viewed as a height function. Whenever the height from node u to node v decreases by more than 1 (i.e., it goes very steep downhill from u to v), the corresponding edge must be saturated.

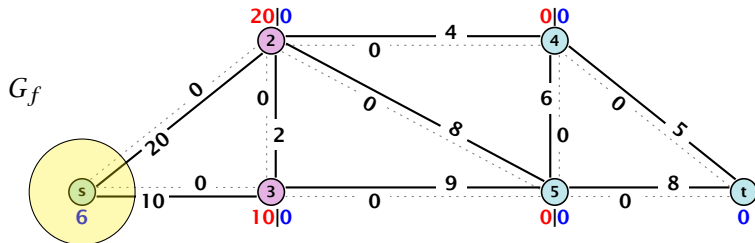
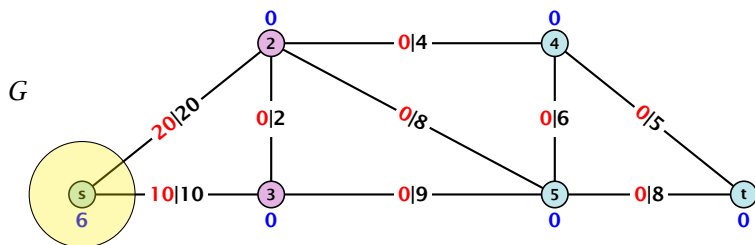
Preflows



Preflows



Preflows



Preflows

Preflows

Lemma 66

A *preflow* that has a valid labelling saturates a cut.

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Proof:

- ▶ There are n nodes but $n + 1$ different labels from $0, \dots, n$.

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- ▶ There are n nodes but $n + 1$ different labels from $0, \dots, n$.
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Preflows

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- ▶ There must exist a label $d \in \{0, \dots, n\}$ such that none of the nodes carries this label.
- ▶ Let $A = \{v \in V \mid \ell(v) > d\}$ and $B = \{v \in V \mid \ell(v) < d\}$.

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Preflows

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Lemma 67

A *flow* that has a valid labelling is a maximum flow.

Push Relabel Algorithms

Push Relabel Algorithms

Idea:

- ▶ start with some preflow and some valid labelling

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- ▶ successively change the preflow while maintaining a valid labelling

Push Relabel Algorithms

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- ▶ start with some preflow and some valid labelling
- ▶ successively change the preflow while maintaining a valid labelling
- ▶ stop when you have a flow (i.e., no more active nodes)

Changing a Preflow

An arc (u, v) with $c_f(u, v) > 0$ in the residual graph is **admissible** if $\ell(u) = \ell(v) + 1$ (i.e., it goes downwards w.r.t. labelling ℓ).

The push operation

Consider an active node u with excess flow

$f(u) = \sum_{e \in \text{into}(u)} f(e) - \sum_{e \in \text{out}(u)} f(e)$ and suppose $e = (u, v)$ is an admissible arc with residual capacity $c_f(e)$.

We can send flow $\min\{c_f(e), f(u)\}$ along e and obtain a new preflow. The old labelling is still valid (!!!).

→ $c_f(u, v) = c_f(u, v) - \min\{c_f(u, v), f(u)\} = c_f(u, v)$

→ the arc e is deleted from the residual graph

→ $c_f(v, u) = c_f(v, u) + \min\{c_f(u, v), f(u)\} = c_f(v, u)$

→ the labelling ℓ is still valid

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new flow $f' = \min\{c_f(e), f(u)\}$ units

are pushed along e to v and deleted from the residual graph

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- ▶ **saturating push** : $\min\{f(u), c_f(e)\} = c_f(e)$
the arc e is deleted from the residual graph
- ▶ **non-saturating push** : $\min\{f(u), c_f(e)\} = f(u)$
the node u becomes inactive

Changing a Preflow

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Push Relabel Algorithms

Push Relabel Algorithms

The relabel operation

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Push Relabel Algorithms

The relabel operation

Consider an active node u that does not have an outgoing admissible arc.

Increasing the label of u by 1 results in a valid labelling.

- ▶ Edges (w, u) incoming to u still fulfill their constraint $\ell(w) \leq \ell(u) + 1$.
- ▶ An outgoing edge (u, w) had $\ell(u) < \ell(w) + 1$ before since it was not admissible. Now: $\ell(u) \leq \ell(w) + 1$.

Push Relabel Algorithms

Intuition:

We want to send flow downwards, since the source has a height/label of n and the target a height/label of 0. If we see an active node u with an admissible arc we push the flow at u towards the other end-point that has a lower height/label. If we do not have an admissible arc but excess flow into u it should roughly mean that the level/height/label of u should rise. (If we consider the flow to be water than this would be natural).

Note that the above intuition is very incorrect as the labels are integral, i.e., they cannot really be seen as the height of a node.

Push Relabel Algorithms

Algorithm 47 $\text{maxflow}(G, s, t, c)$

```
1: find initial preflow  $f$ 
2: while there is active node  $u$  do
3:   if there is admiss. arc  $e$  out of  $u$  then
4:      $\text{push}(G, e, f, c)$ 
5:   else
6:      $\text{relabel}(u)$ 
7: return  $f$ 
```

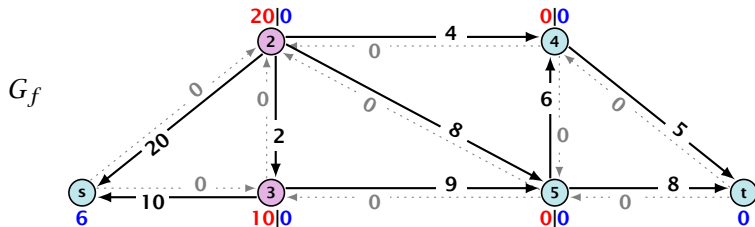
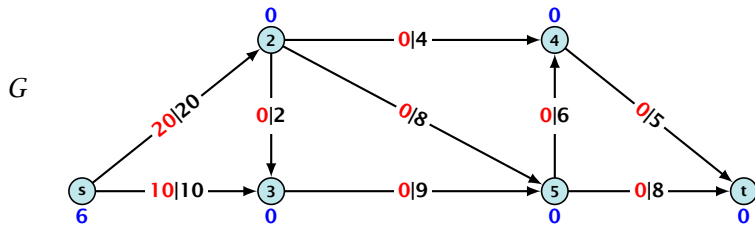
Push Relabel Algorithms

Algorithm 47 $\text{maxflow}(G, s, t, c)$

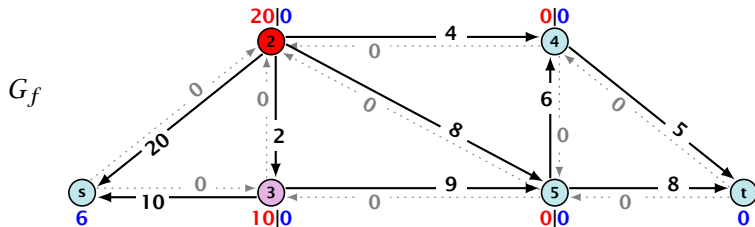
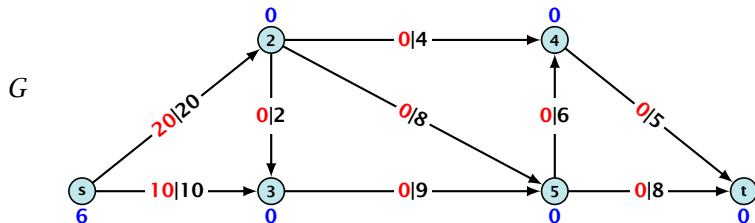
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4:      $\text{push}(G, e, f, c)$ 
5:   else
6:      $\text{relabel}(u)$ 
7: return  $f$ 
```

In the following example we always stick to the same active node u until it becomes inactive but this is not required.

Preflow Push Algorithm

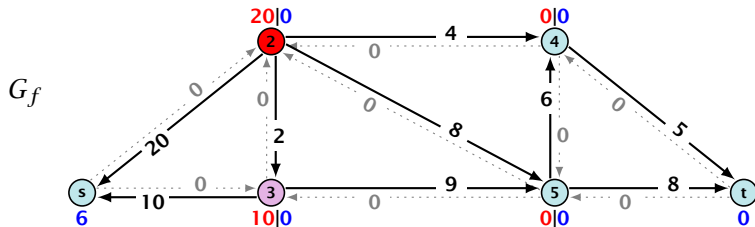
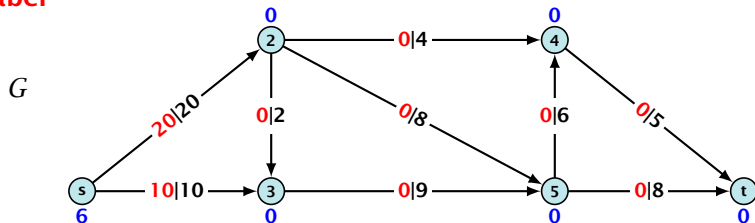


Preflow Push Algorithm

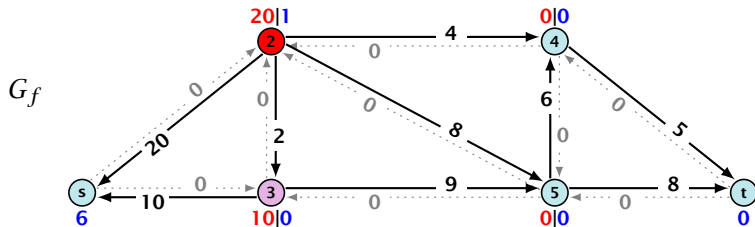
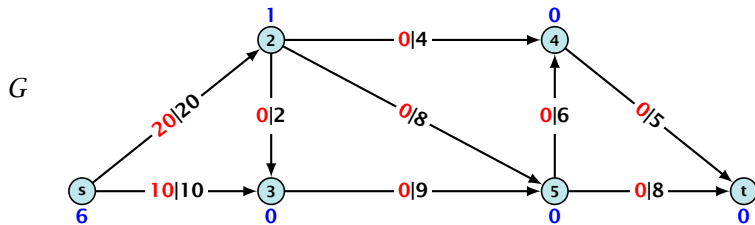


Preflow Push Algorithm

relabel

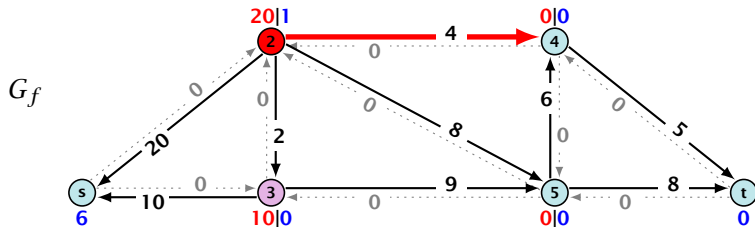
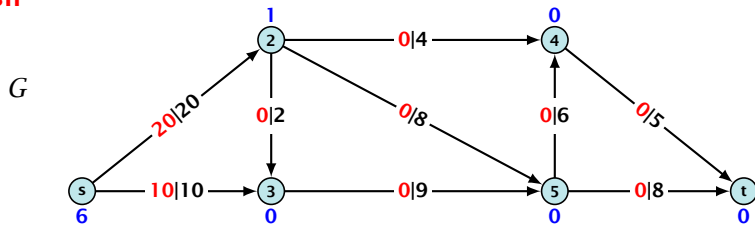


Preflow Push Algorithm

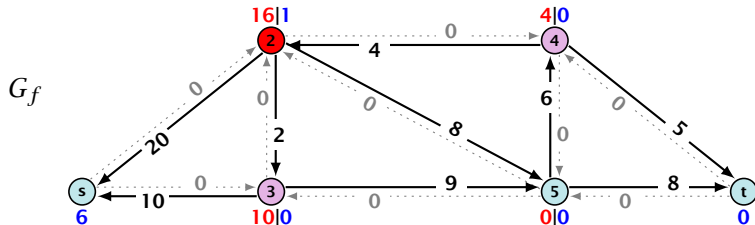
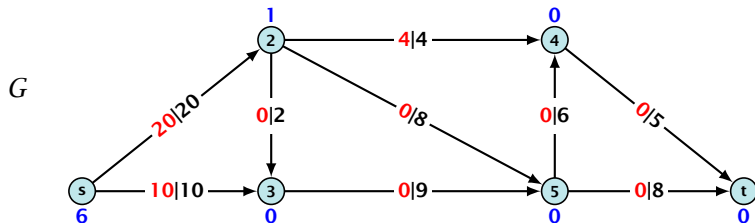


Preflow Push Algorithm

push

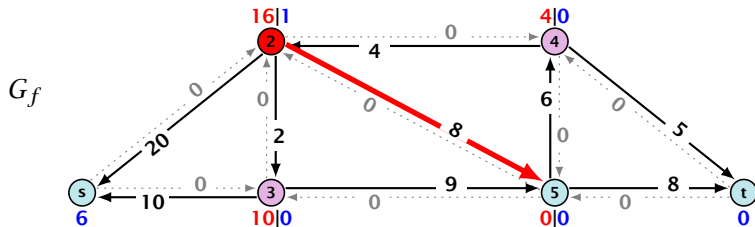
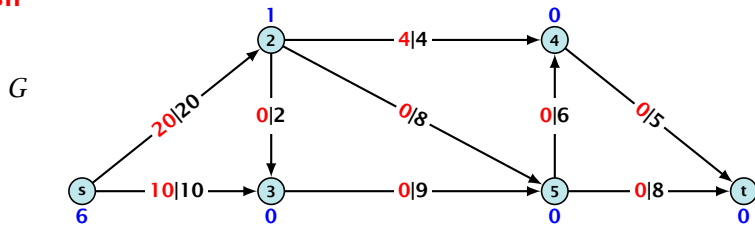


Preflow Push Algorithm

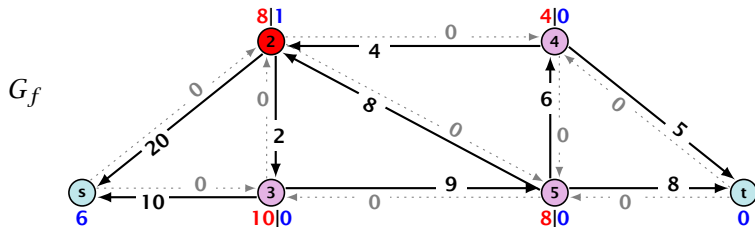
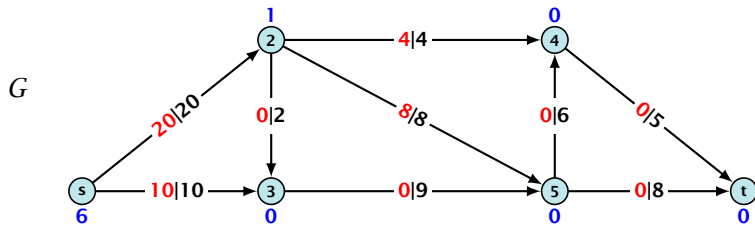


Preflow Push Algorithm

push

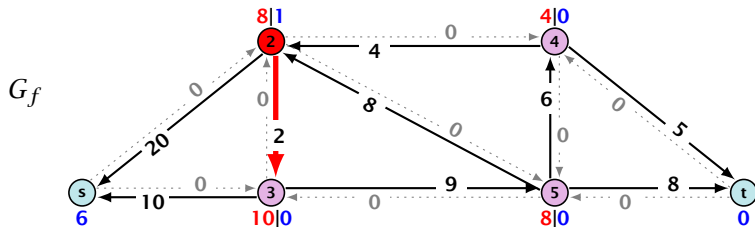
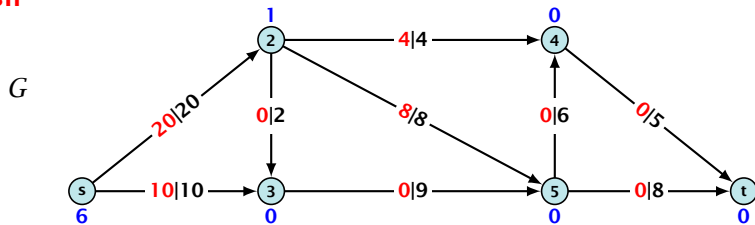


Preflow Push Algorithm

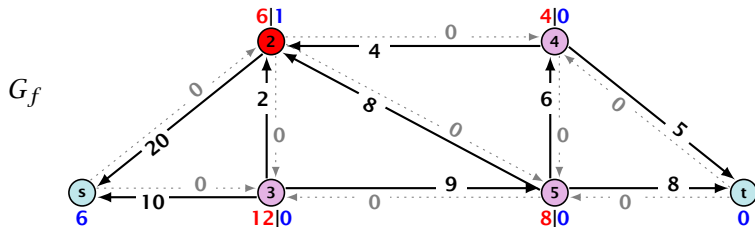
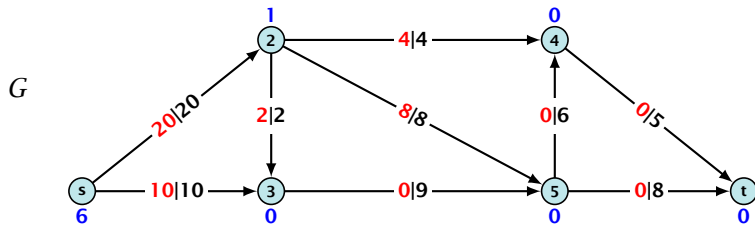


Preflow Push Algorithm

push

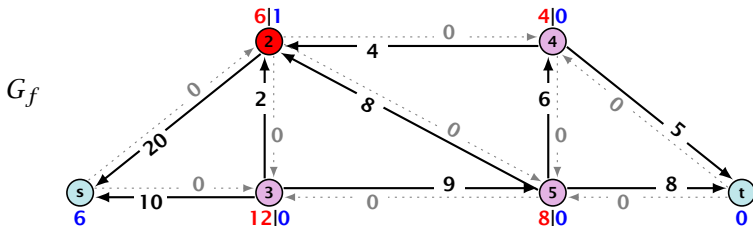
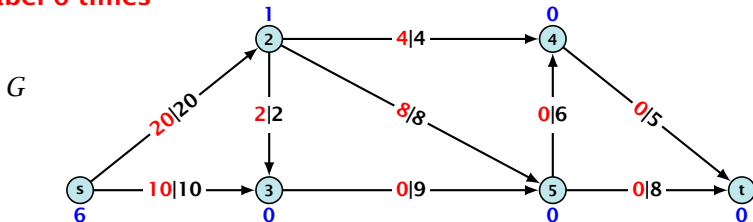


Preflow Push Algorithm

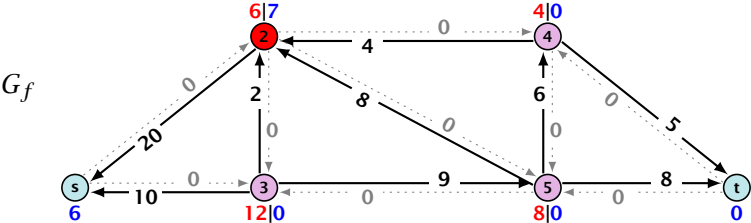
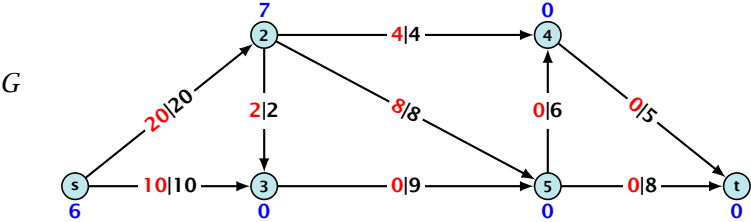


Preflow Push Algorithm

relabel 6 times

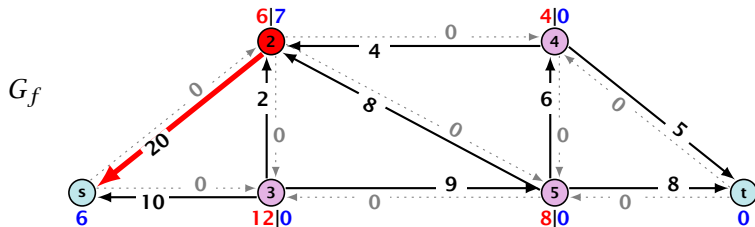
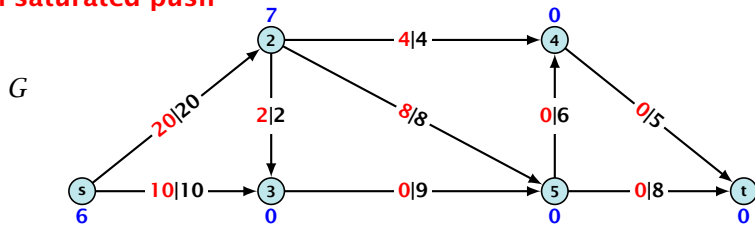


Preflow Push Algorithm

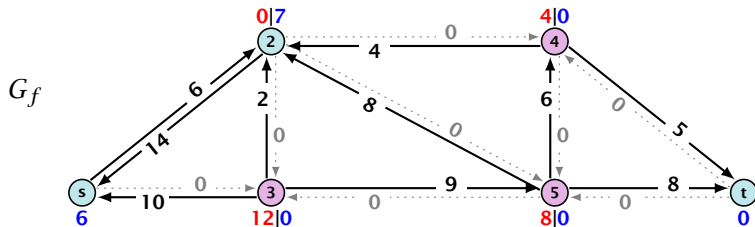
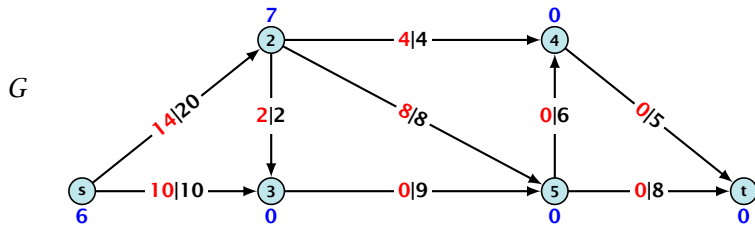


Preflow Push Algorithm

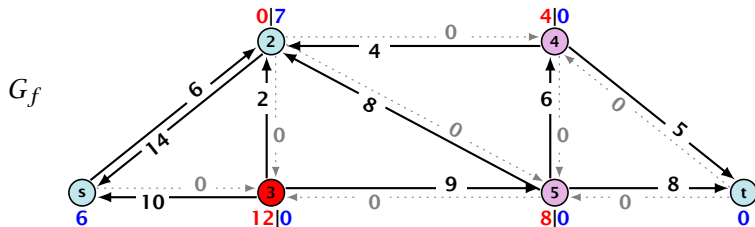
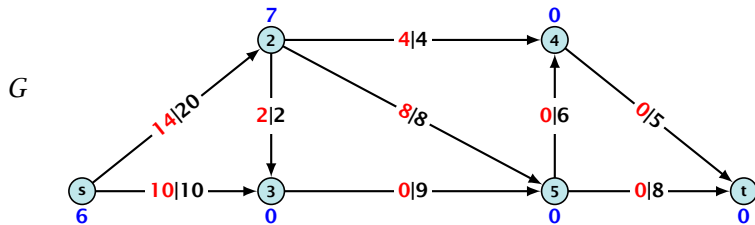
non-saturated push



Preflow Push Algorithm

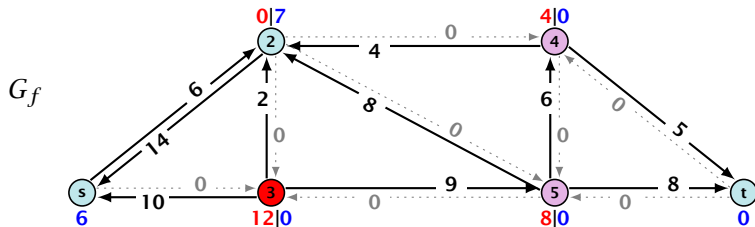
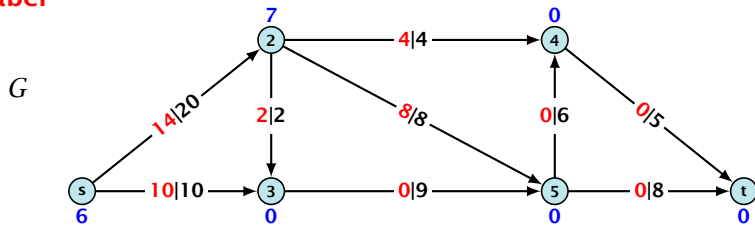


Preflow Push Algorithm

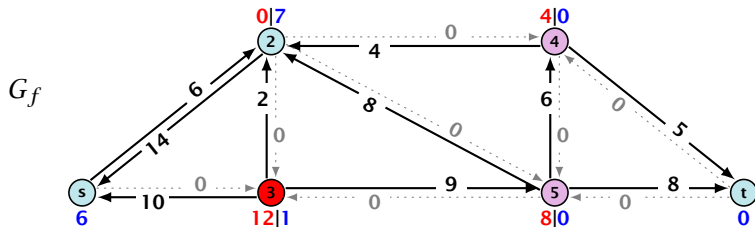
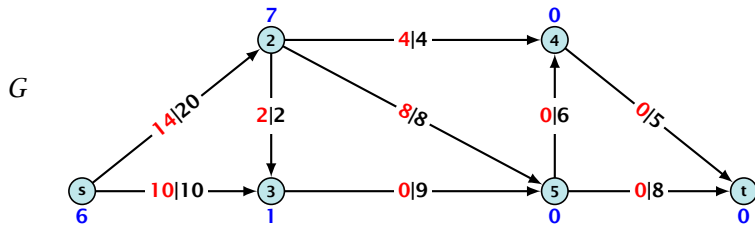


Preflow Push Algorithm

relabel

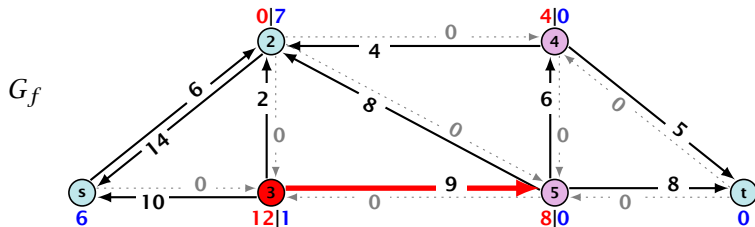
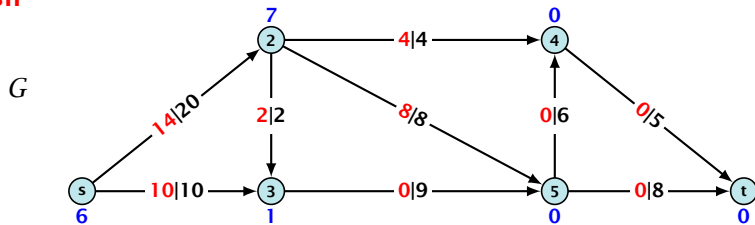


Preflow Push Algorithm

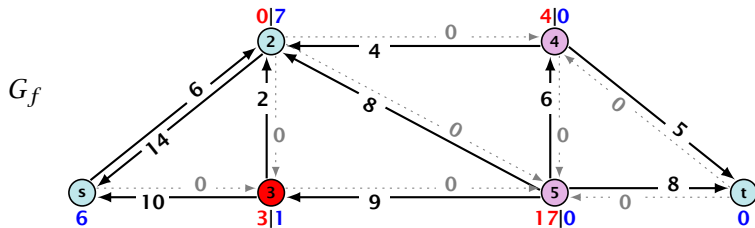
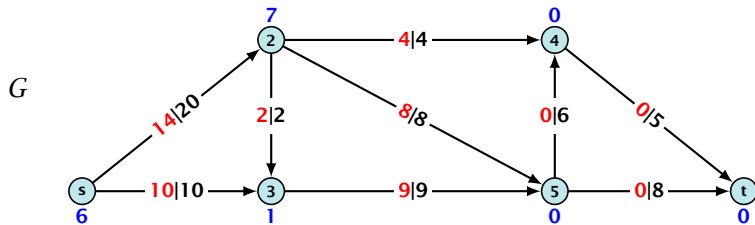


Preflow Push Algorithm

push

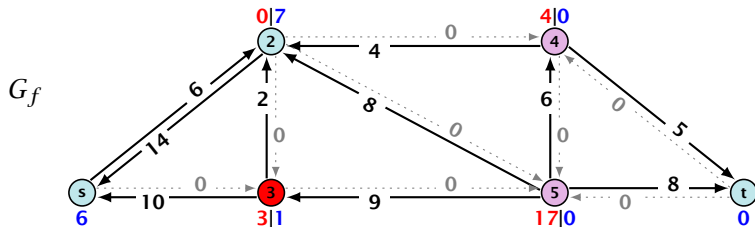
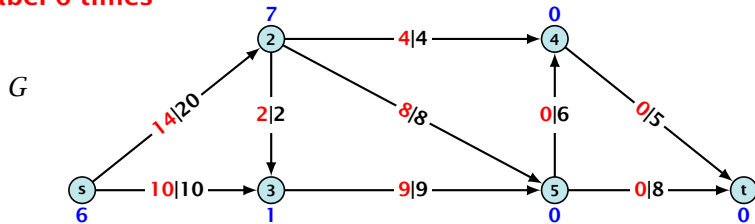


Preflow Push Algorithm

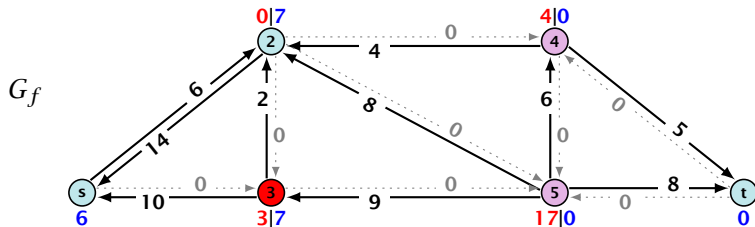
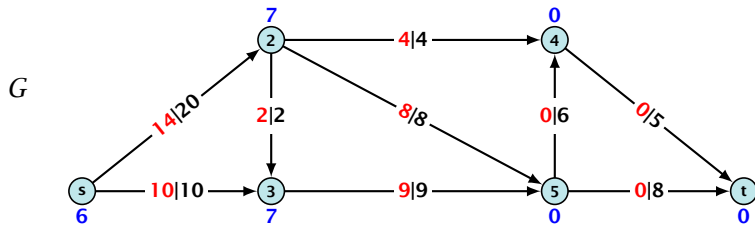


Preflow Push Algorithm

relabel 6 times

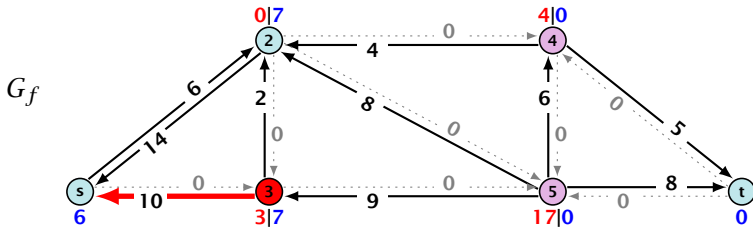
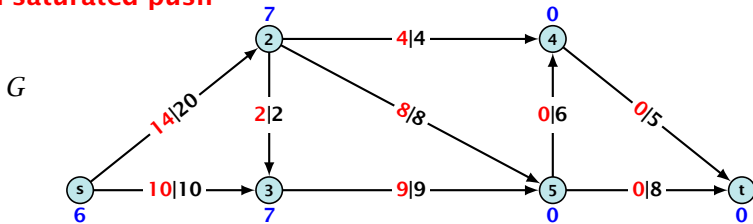


Preflow Push Algorithm

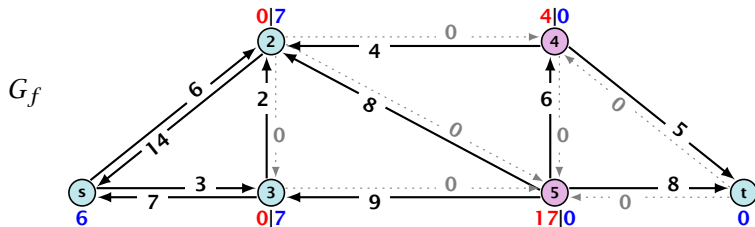
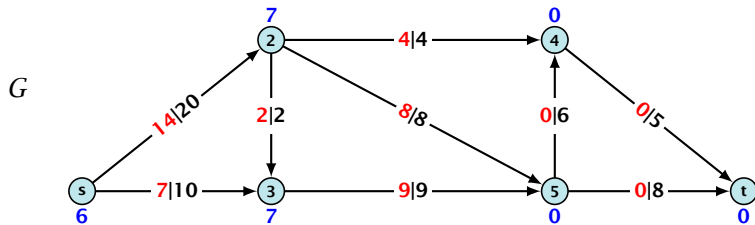


Preflow Push Algorithm

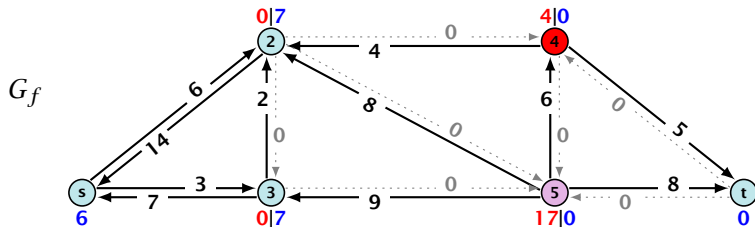
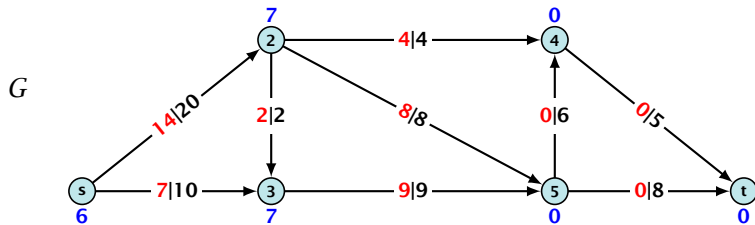
non-saturated push



Preflow Push Algorithm

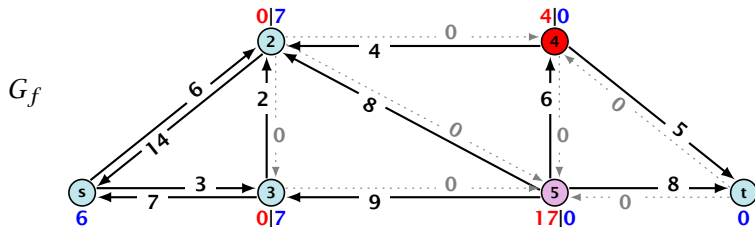
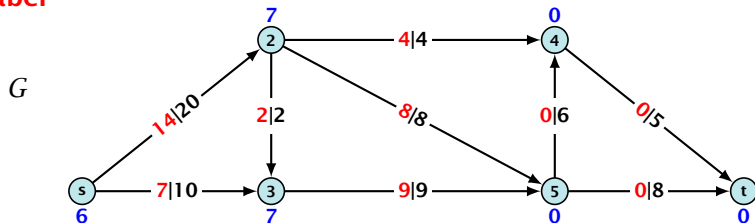


Preflow Push Algorithm

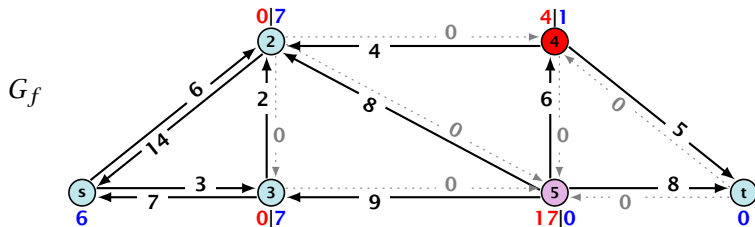
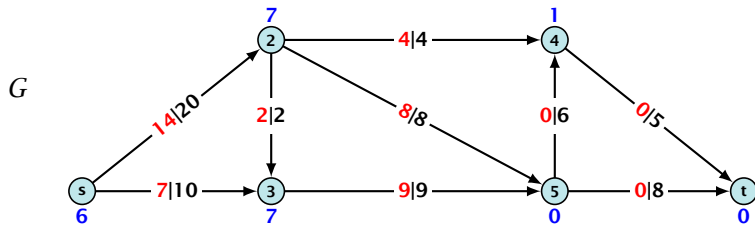


Preflow Push Algorithm

relabel

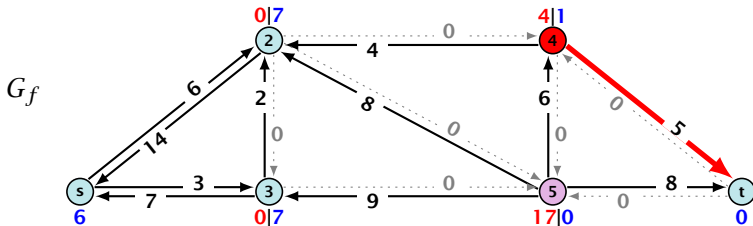
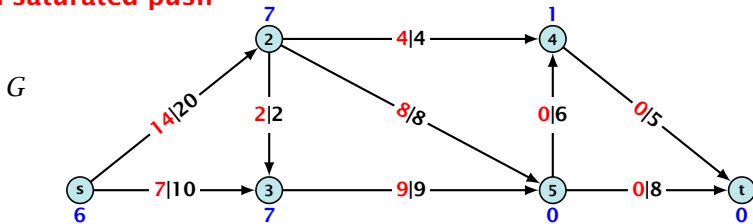


Preflow Push Algorithm

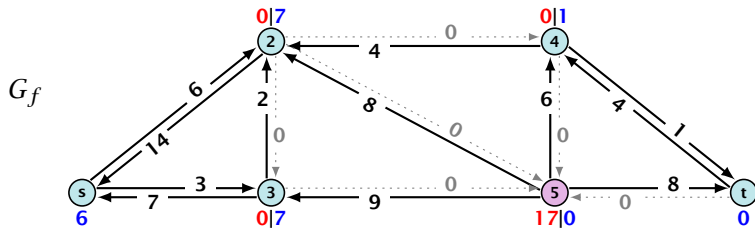
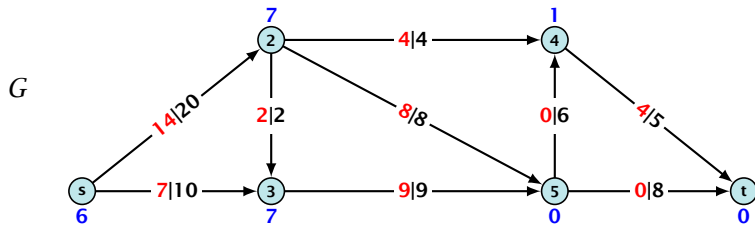


Preflow Push Algorithm

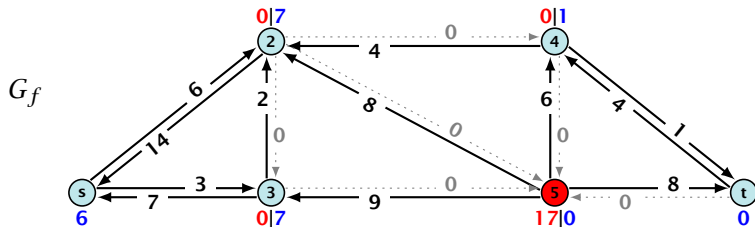
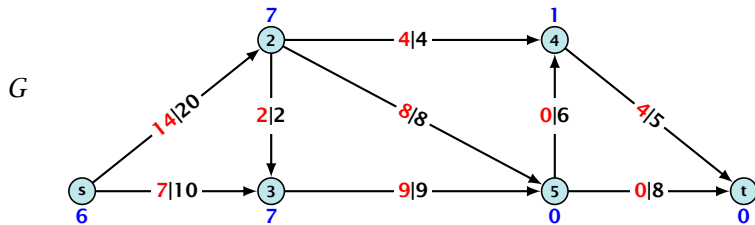
non-saturated push



Preflow Push Algorithm

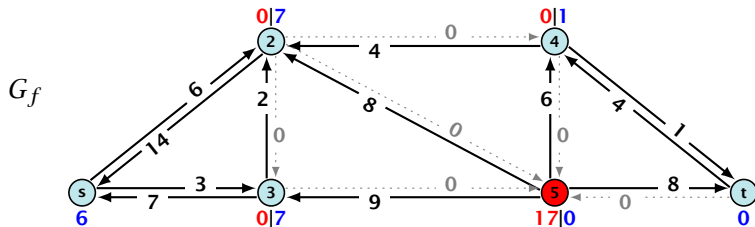
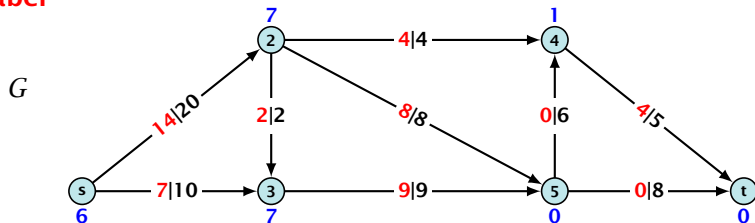


Preflow Push Algorithm

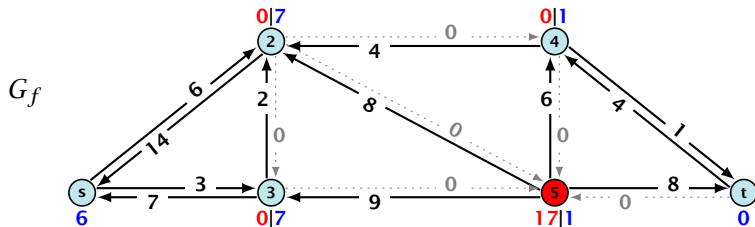
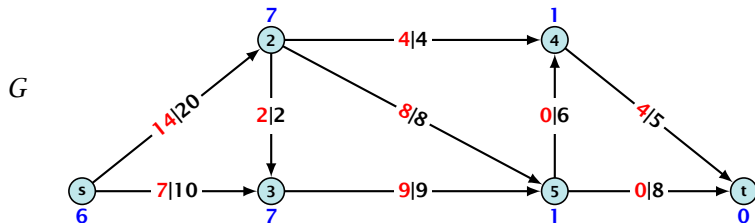


Preflow Push Algorithm

relabel



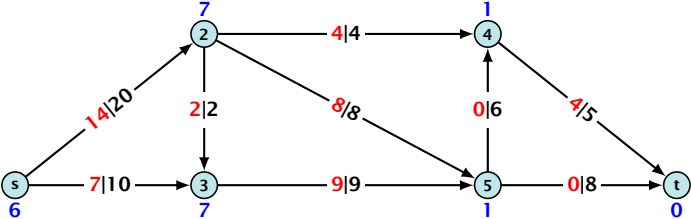
Preflow Push Algorithm



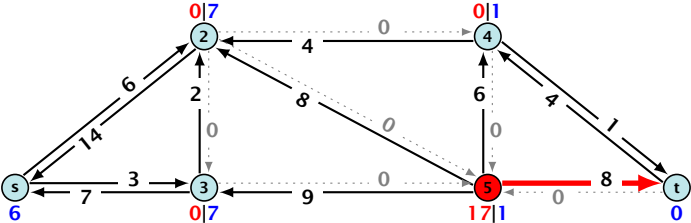
Preflow Push Algorithm

push

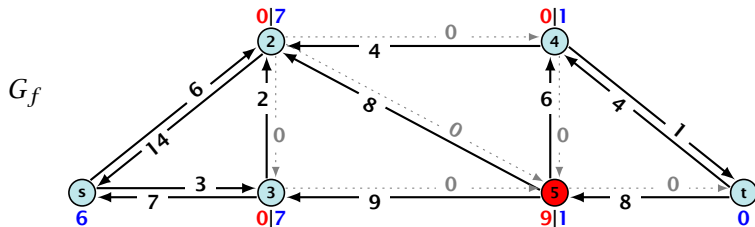
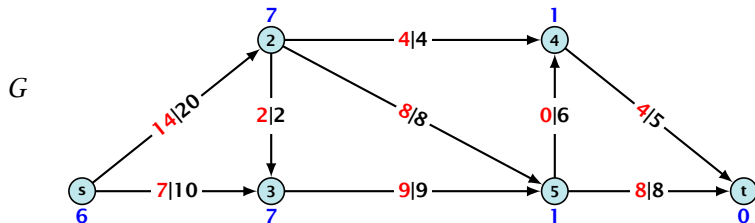
G



G_f

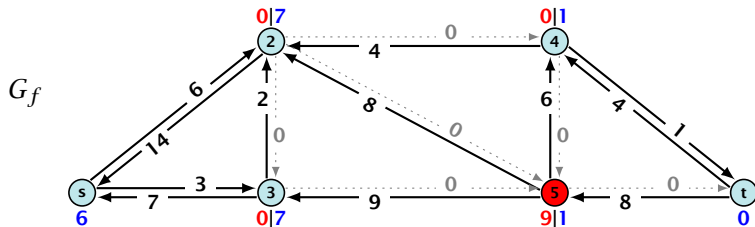
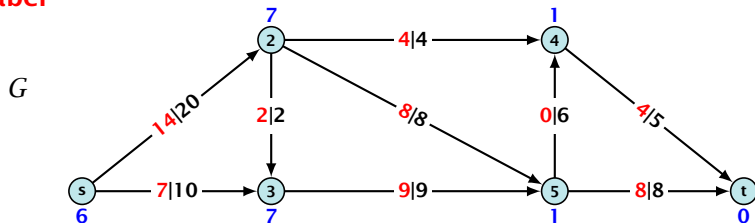


Preflow Push Algorithm

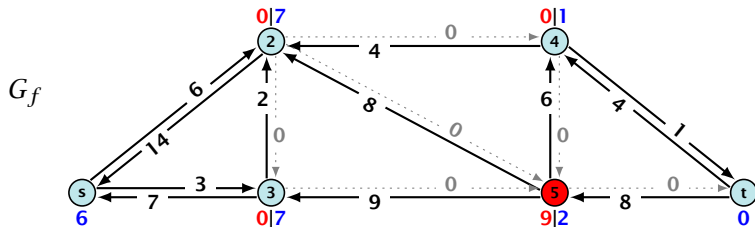
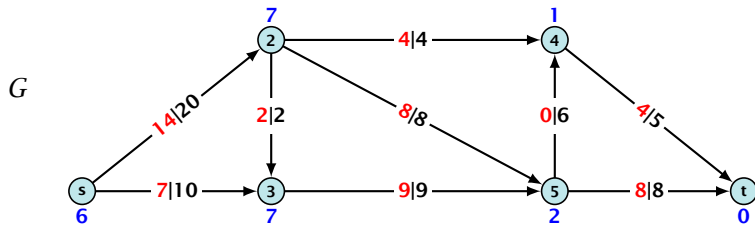


Preflow Push Algorithm

relabel

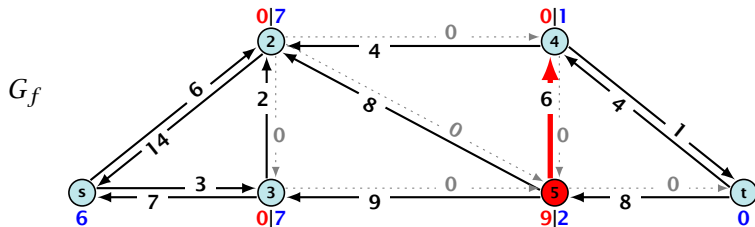
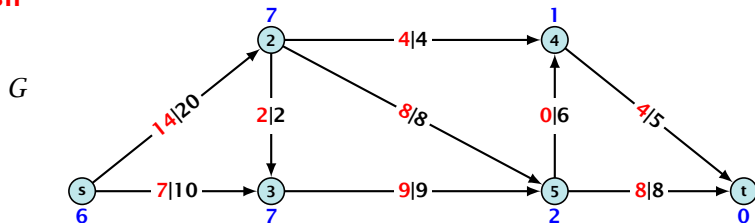


Preflow Push Algorithm

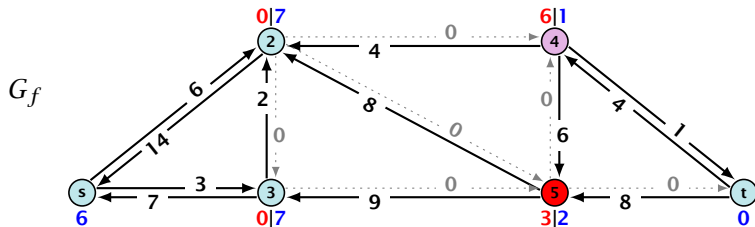
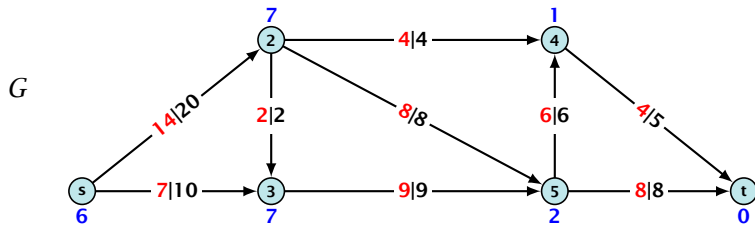


Preflow Push Algorithm

push

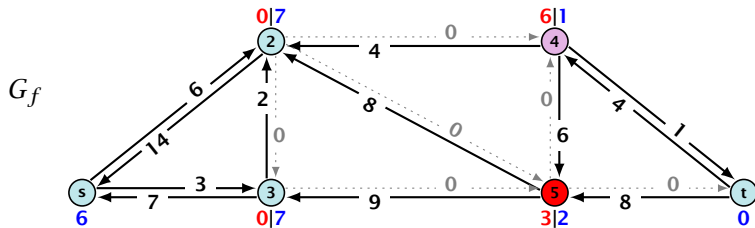
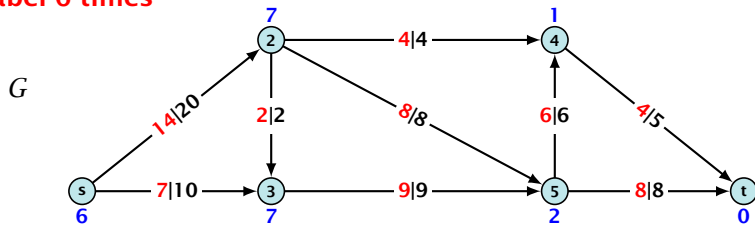


Preflow Push Algorithm

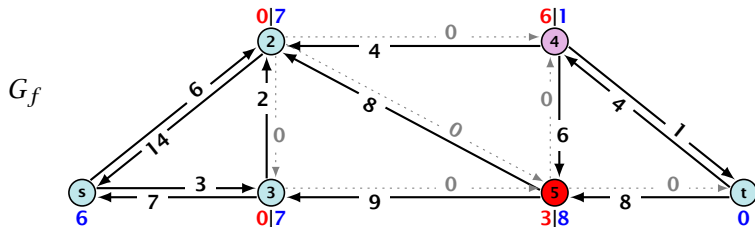
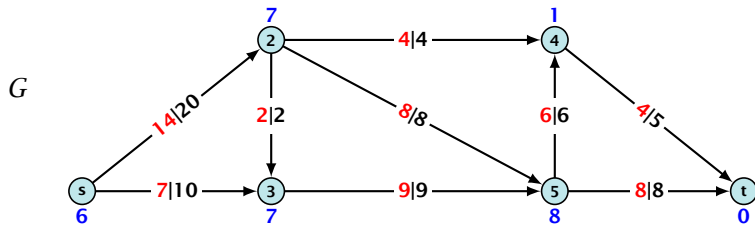


Preflow Push Algorithm

relabel 6 times

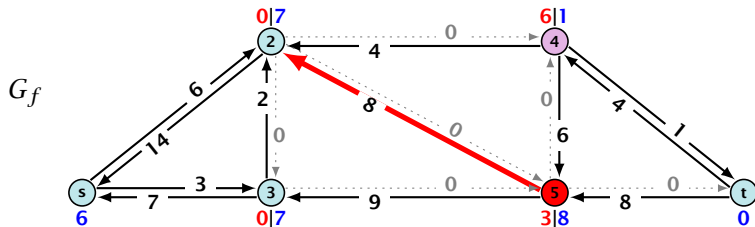
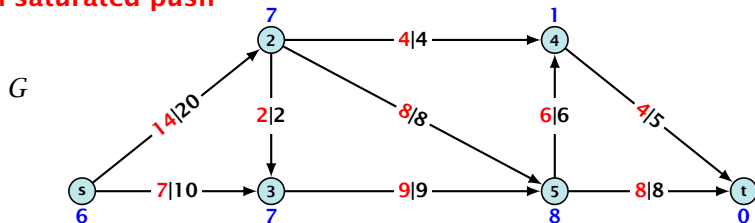


Preflow Push Algorithm

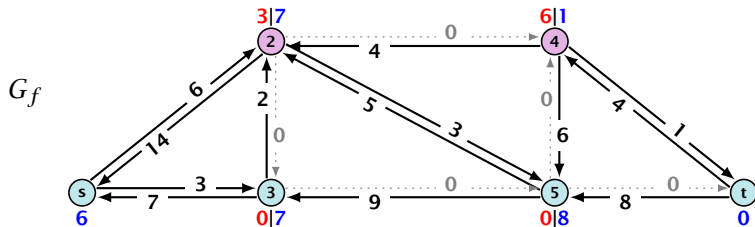
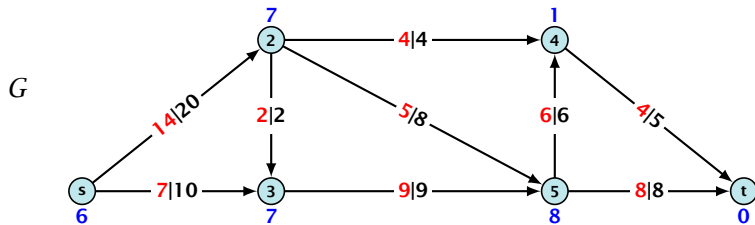


Preflow Push Algorithm

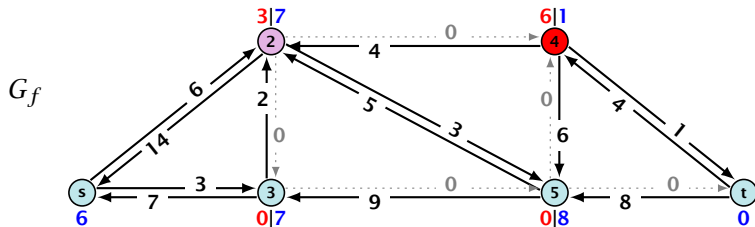
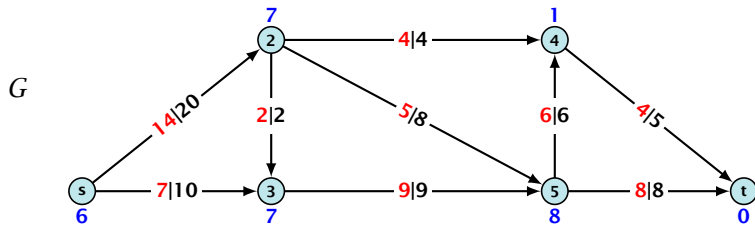
non-saturated push



Preflow Push Algorithm

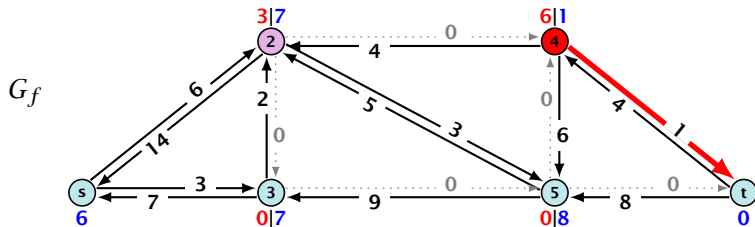
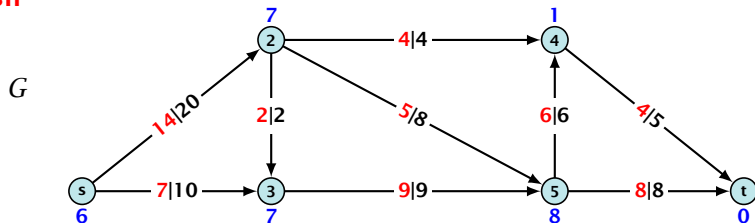


Preflow Push Algorithm

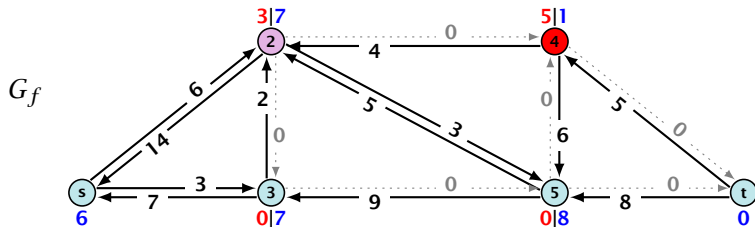
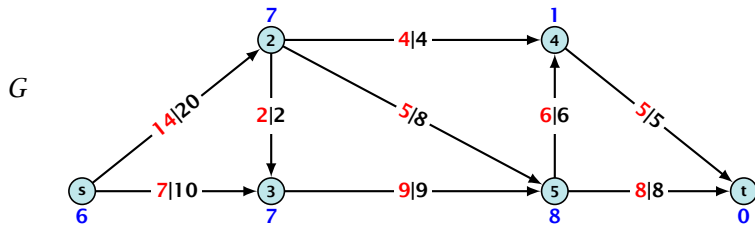


Preflow Push Algorithm

push

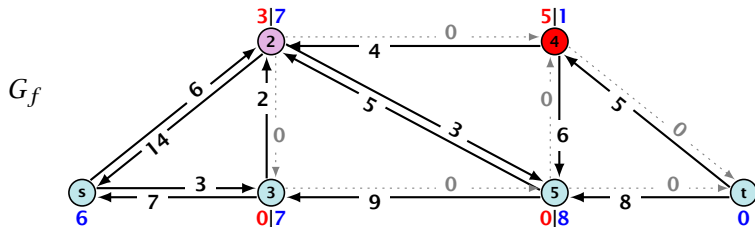
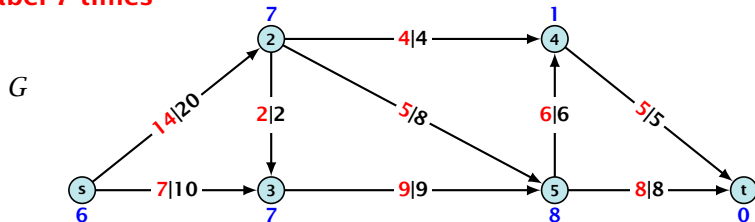


Preflow Push Algorithm

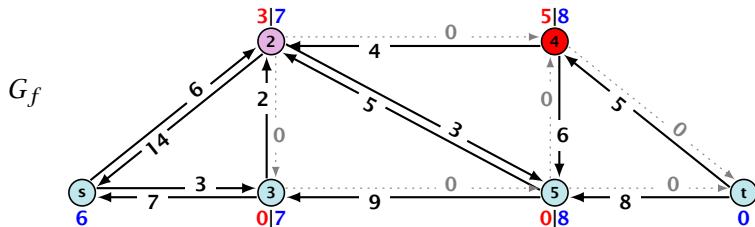
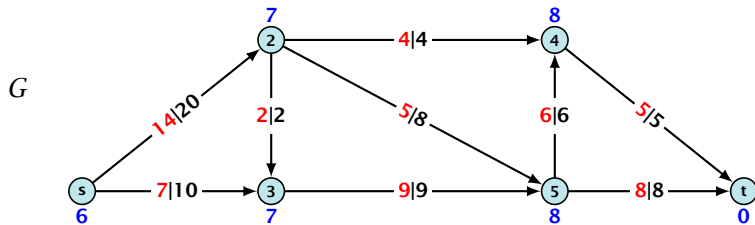


Preflow Push Algorithm

relabel 7 times

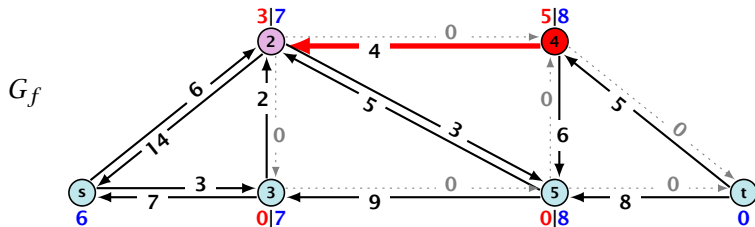
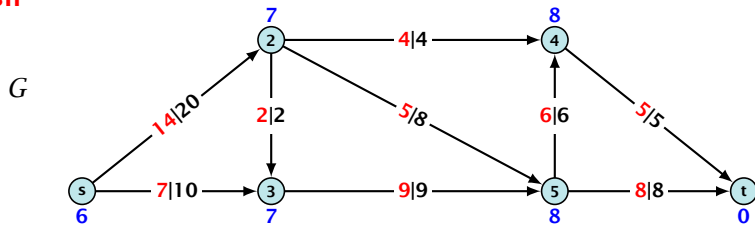


Preflow Push Algorithm

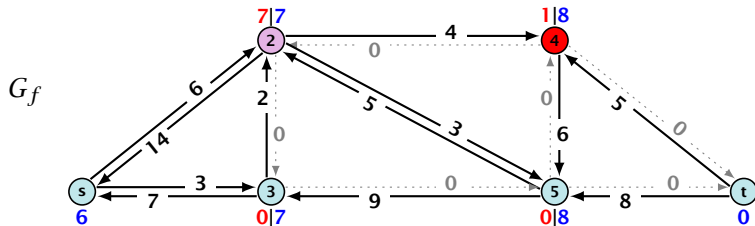
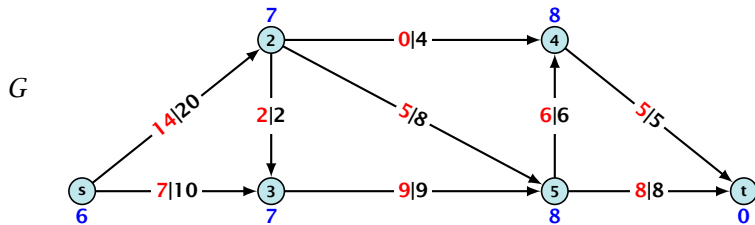


Preflow Push Algorithm

push

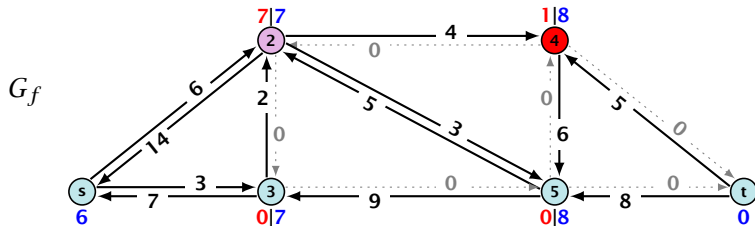
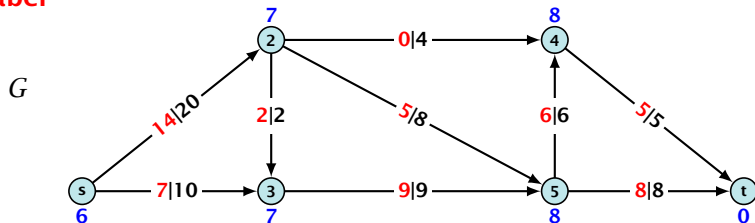


Preflow Push Algorithm

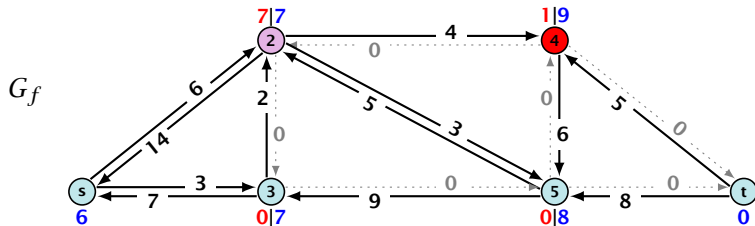
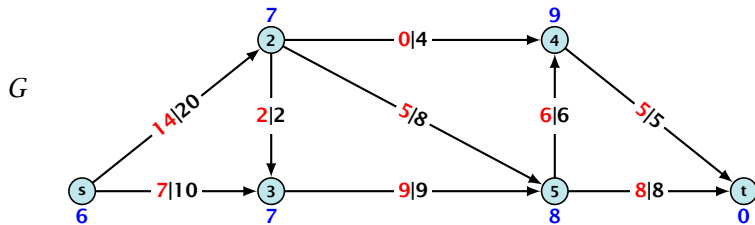


Preflow Push Algorithm

relabel

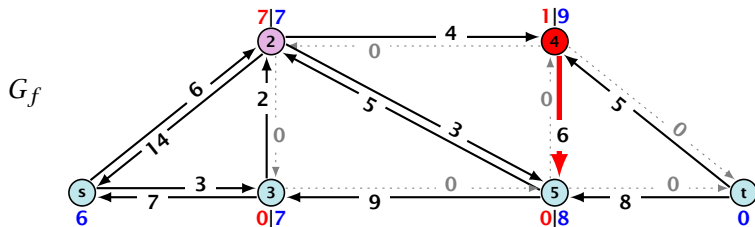
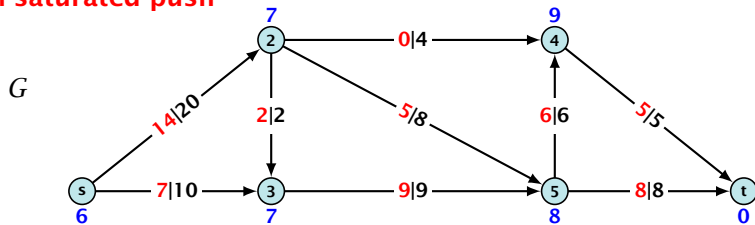


Preflow Push Algorithm

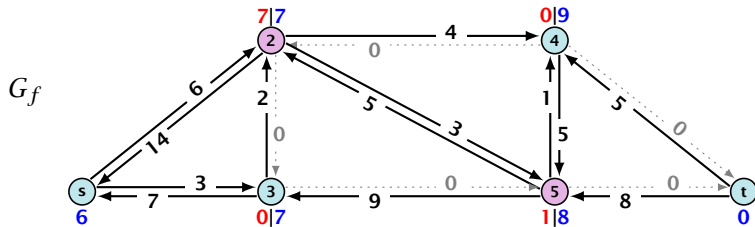
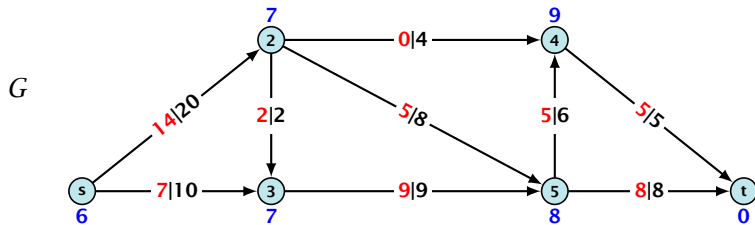


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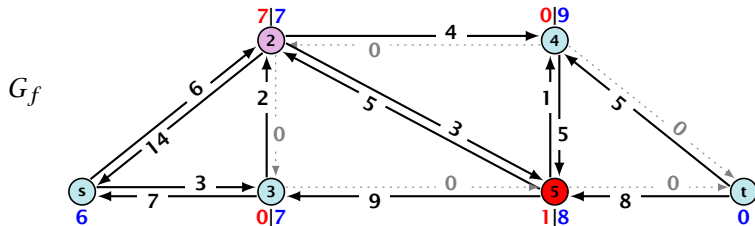
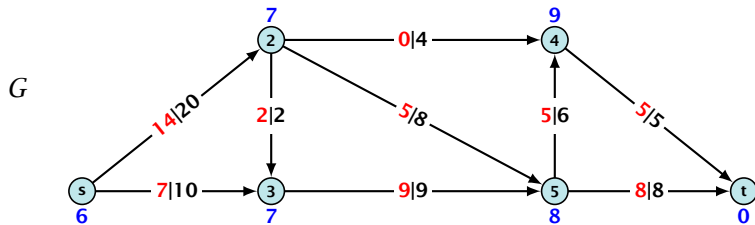
non-saturated push



Preflow Push Algorithm

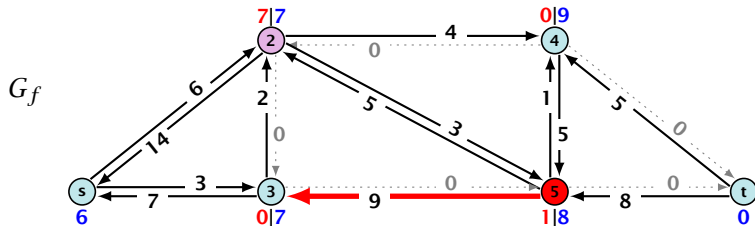
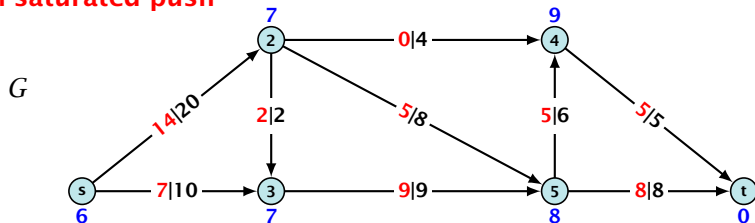


Preflow Push Algorithm

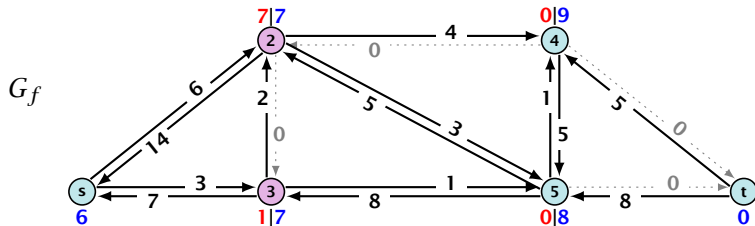
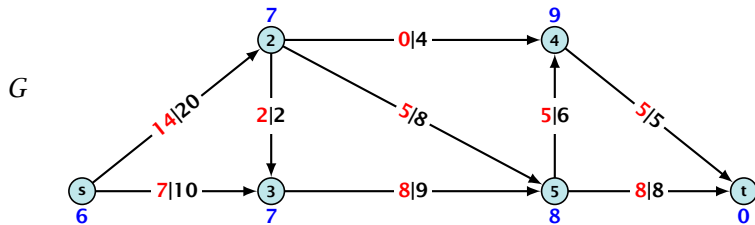


Preflow Push Algorithm

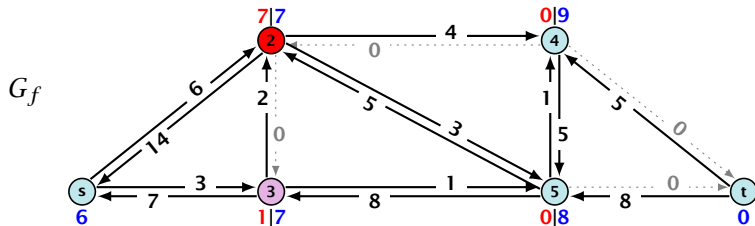
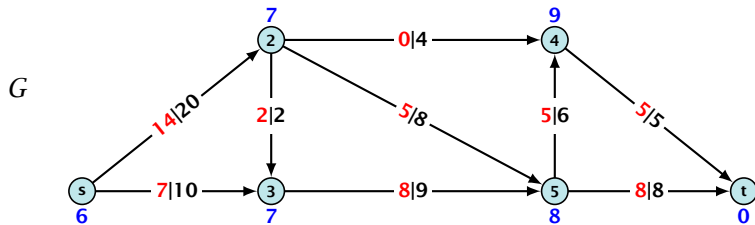
non-saturated push



Preflow Push Algorithm

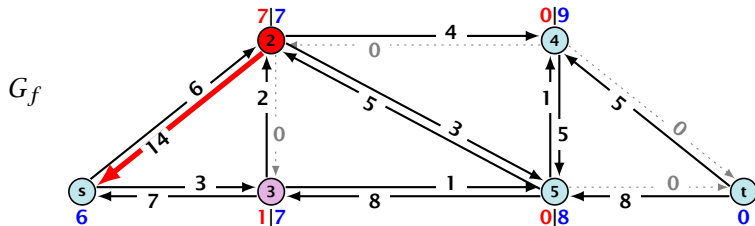
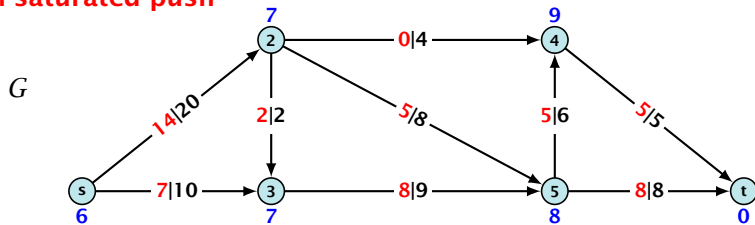


Preflow Push Algorithm

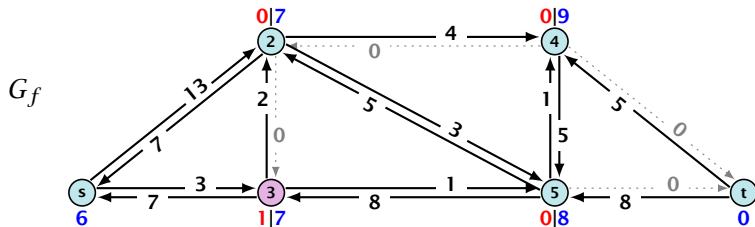
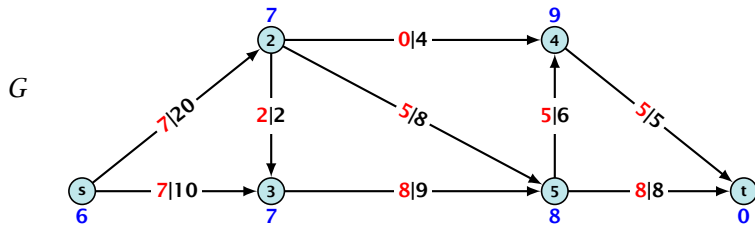


Preflow Push Algorithm

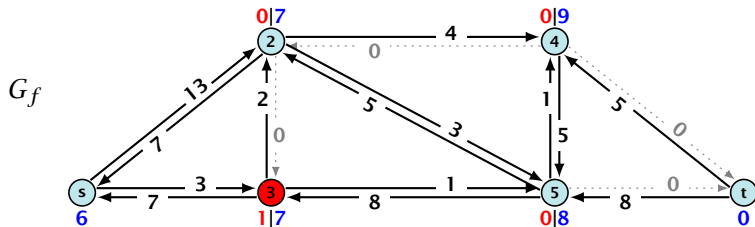
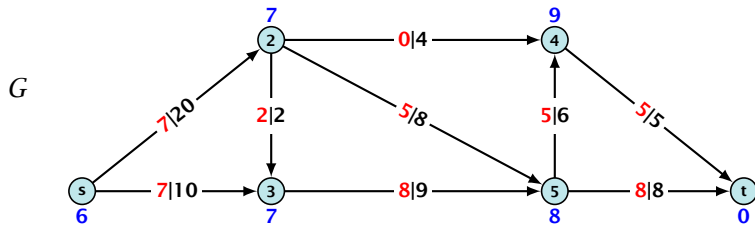
non-saturated push



Preflow Push Algorithm

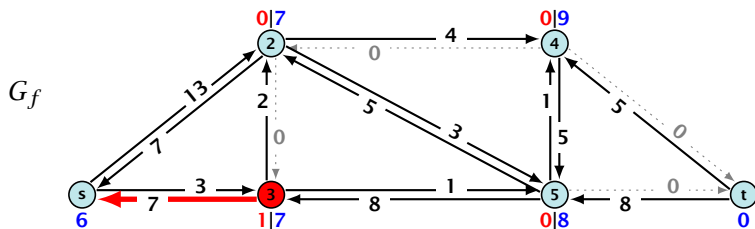
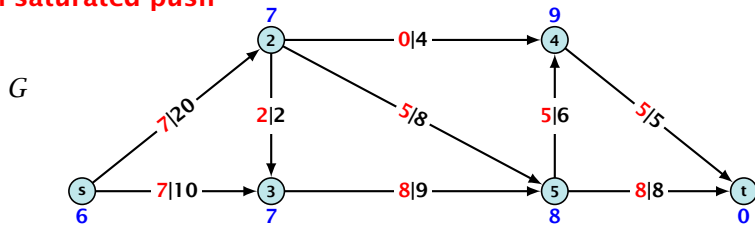


Preflow Push Algorithm

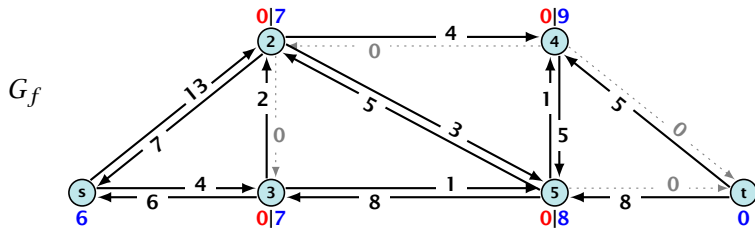
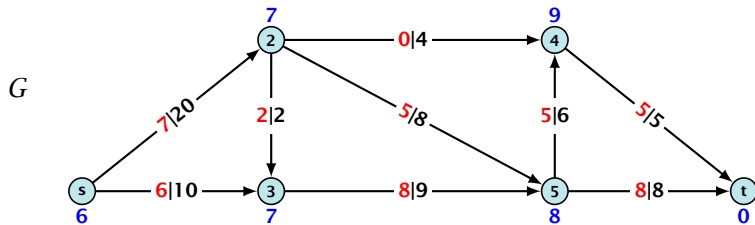


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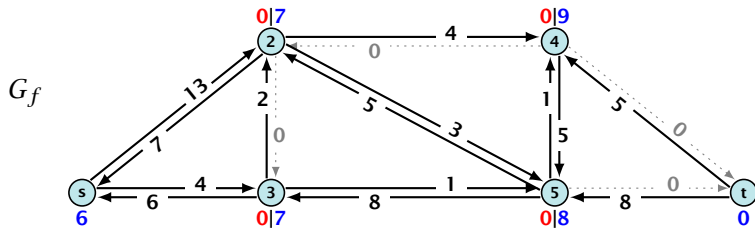
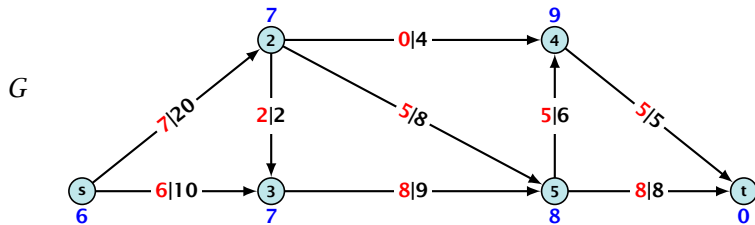
non-saturated push



Preflow Push Algorithm



Preflow Push Algorithm



Analysis

Lemma 68

An active node has a path to s in the residual graph.

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- ▶ Let $f(B) = \sum_{v \in B} f(v)$ be the excess flow of all nodes in B .

Let $f : E \rightarrow \mathbb{R}_0^+$ be a preflow. We introduce the notation

$$f(x, y) = \begin{cases} 0 & (x, y) \notin E \\ f((x, y)) & (x, y) \in E \end{cases}$$

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$$\begin{aligned} f(B) &= \sum_{b \in B} f(b) \\ &= \sum_{b \in B} \left(\sum_{v \in V} f(v, b) - \sum_{v \in V} f(b, v) \right) \\ &= \sum_{b \in B} \left(\sum_{v \in A} f(v, b) + \sum_{v \in B} f(v, b) - \sum_{v \in A} f(b, v) - \sum_{v \in B} f(b, v) \right) \\ &= \sum_{b \in B} \sum_{v \in A} f(v, b) - \sum_{b \in B} \sum_{v \in A} f(b, v) + \sum_{b \in B} \sum_{v \in B} f(v, b) - \sum_{b \in B} \sum_{v \in B} f(b, v) \\ &= 0 \end{aligned}$$

Let $f : E \rightarrow \mathbb{R}_0^+$ be a preflow. We introduce the notation

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Hence, the excess flow $f(b)$ must be 0 for every node $b \in B$.

Analysis

Lemma 69

The label of a node cannot become larger than $2n - 1$.

Proof.

- ▶ When increasing the label at a node u there exists a path from u to s of length at most $n - 1$. Along each edge of the path the height/label can at most drop by 1, and the label of the source is n .

Analysis

Lemma 70

There are only $\mathcal{O}(n^3)$ calls to discharge when using the relabel-to-front heuristic.

Proof.

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The number of *saturating pushes* performed is at most $\mathcal{O}(mn)$.

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- ▶ For the edge to appear again, a push from v to u is required.

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- ▶ Currently, $\ell(u) = \ell(v) + 1$, as we only make pushes along admissible edges.

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- ▶ For a push from v to u the edge (v, u) must become admissible. The label of v must increase by at least 2.
- ▶ Since the label of v is at most $2n - 1$, there are at most n pushes along (u, v) .

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The number of *non-saturating pushes* performed is at most $\mathcal{O}(n^2m)$.

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- ▶ Define a potential function $\Phi(f) = \sum_{\text{active nodes } v} \ell(v)$
- ▶ A saturating push increases Φ by at most $2n$.
- ▶ A relabel increases Φ by at most 1.

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- ▶ A non-saturating push decreases Φ by at least 1 as the node that is pushed from becomes inactive and has a label that is strictly larger than the target.
- ▶ Hence,

$$\begin{aligned} \# \text{non-saturating_pushes} &\leq \# \text{relabels} + 2n \cdot \# \text{saturating_pushes} \\ &\leq \mathcal{O}(n^2m) . \end{aligned}$$

Analysis

There is an implementation of the generic push relabel algorithm with running time $\mathcal{O}(n^2m)$.

For every node maintain a list of admissible edges starting at that node. Further maintain a list of active nodes.

A push along an edge (u, v) can be performed in constant time

- check whether edge (v, u) needs to be added to E
- check whether (u, v) needs to be deleted (saturating push)
- check whether u becomes inactive and has to be deleted from the set of active nodes

A relabel at a node u can be performed in time $\mathcal{O}(n)$

- check for all outgoing edges if they became admissible
- check for all incoming edges if they became non-saturating

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For every node maintain a list of admissible edges starting at that node. Further maintain a list of active nodes.

A push along an edge (u, v) can be performed in constant time

- check whether edge (u, v) needs to be added to $A(u)$
- check whether (u, v) needs to be deleted (active nodes)
- push $\min\{c(u, v), \text{excess}(u)\}$ units of flow and decrease $\text{excess}(u)$ by the same amount

A relabel at a node u can be performed in time $\mathcal{O}(n)$

- relabel all outgoing edges if they become admissible
- relabel all incoming edges if they become inadmissible

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For every node maintain a list of admissible edges starting at that node. Further maintain a list of active nodes.

A push along an edge (u, v) can be performed in constant time

- ▶ check whether edge (v, u) needs to be added to G_f
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A relabel at a node u can be performed in time $\mathcal{O}(n)$

- ▶ check for all outgoing edges if they become admissible
- ▶ check for all incoming edges if they become non-admissible

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There is an implementation of the generic push relabel algorithm with running time $\mathcal{O}(n^2m)$.

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13.2 Relabel to front

For special variants of push relabel algorithms we organize the neighbours of a node into a linked list (possible neighbours in the residual graph G_f). Then we use the discharge-operation:

Algorithm 48 discharge(u)

```
1: while  $u$  is active do  
2:    $v \leftarrow u.current\text{-neighbour}$   
3:   if  $v = \text{null}$  then  
4:     relabel( $u$ )  
5:      $u.current\text{-neighbour} \leftarrow u.neighbour\text{-list-head}$   
6:   else  
7:     if  $(u, v)$  admissable then push( $u, v$ )  
8:     else  $u.current\text{-neighbour} \leftarrow v.next\text{-in-list}$ 
```

13.2 Relabel to front

Lemma 73

If $v = \text{null}$ in line 3, then there is no outgoing admissible edge from u .

The lemma holds because push- and relabel-operations on nodes different from u cannot make edges outgoing from u admissible.

This shows that $\text{discharge}(u)$ is correct, and that we can perform a relabel in line 4.

13.2 Relabel to front

Algorithm 49 relabel-to-front(G, s, t)

```
1: initialize preflow
2: initialize node list  $L$  containing  $V \setminus \{s, t\}$  in any order
3: foreach  $u \in V \setminus \{s, t\}$  do
4:    $u.current\text{-neighbour} \leftarrow u.neighbour\text{-list}\text{-head}$ 
5:  $u \leftarrow L.head$ 
6: while  $u \neq \text{null}$  do
7:    $old\text{-height} \leftarrow \ell(u)$ 
8:   discharge( $u$ )
9:   if  $\ell(u) > old\text{-height}$  then
10:     move  $u$  to the front of  $L$ 
11:    $u \leftarrow u.next$ 
```


13.2 Relabel to front

Lemma 74 (Invariant)

In Line 6 of the relabel-to-front algorithm the following invariant holds.

- 1. The sequence L is topologically sorted w.r.t. the set of admissible edges; this means for an admissible edge (x, y) the node x appears before y in sequence L .*
- 2. No node before u in the list L is active.*

Proof:

► Initialization:

1. In the beginning s has label $n \geq 2$, and all other nodes have label 0. Hence, no edge is admissible, which means that any ordering L is permitted.
2. We start with u being the head of the list; hence no node before u can be active

► Maintenance:

1.
 - Pushes do not create any new admissible edges. Therefore, not relabeling u leaves L topologically sorted.
 - After relabeling, u cannot have admissible incoming edges as such an edge (x, u) would have had a difference $\ell(x) - \ell(u) \geq 2$ before the re-labeling (such edges do not exist in the residual graph).
Hence, moving u to the front does not violate the sorting property for any edge; however it fixes this property for all admissible edges leaving u that were generated by the relabeling.

13.2 Relabel to front

Proof:

► Maintenance:

2. If we do a relabel there is nothing to prove because the only node before u' (u in the next iteration) will be the current u ; the $\text{discharge}(u)$ operation only terminates when u is not active anymore.

For the case that we do a relabel, observe that the only way a predecessor could be active is that we push flow to it via an admissible arc. However, all admissible arcs point to successors of u .

Note that the invariant for $u = \text{null}$ means that we have a preflow with a valid labelling that does not have active nodes. This means we have a maximum flow.

13.2 Relabel to front

Lemma 75

There are at most $\mathcal{O}(n^3)$ calls to $\text{discharge}(u)$.

Every discharge operation without a relabel advances u (the current node within list L). Hence, if we have n discharge operations without a relabel we have $u = \text{null}$ and the algorithm terminates.

Therefore, the number of calls to discharge is at most $n(\#\text{relabels} + 1) = \mathcal{O}(n^3)$.

13.2 Relabel to front

Lemma 76

The cost for all relabel-operations is only $\mathcal{O}(n^2)$.

A relabel-operation at a node is constant time (increasing the label and resetting *u.current-neighbour*). In total we have $\mathcal{O}(n^2)$ relabel-operations.

13.2 Relabel to front

Note that by definition a saturating push operation ($\min\{c_f(e), f(u)\} = c_f(e)$) can at the same time be a non-saturating push operation ($\min\{c_f(e), f(u)\} = f(u)$).

Lemma 77

*The cost for all saturating push-operations that are **not** also non-saturating push-operations is only $\mathcal{O}(mn)$.*

Note that such a push-operation leaves the node u active but makes the edge e disappear from the residual graph. Therefore the push-operation is immediately followed by an increase of the pointer $u.current-neighbour$.

This pointer can traverse the neighbour-list at most $\mathcal{O}(n)$ times (upper bound on number of relabels) and the neighbour-list has only $degree(u) + 1$ many entries (+1 for null-entry).

13.2 Relabel to front

Lemma 78

The cost for all non-saturating push-operations is only $\mathcal{O}(n^3)$.

A non-saturating push-operation takes constant time and ends the current call to `discharge()`. Hence, there are only $\mathcal{O}(n^3)$ such operations.

Theorem 79

The push-relabel algorithm with the rule relabel-to-front takes time $\mathcal{O}(n^3)$.

13.3 Highest label

Algorithm 50 highest-label(G, s, t)

- 1: initialize preflow
- 2: **foreach** $u \in V \setminus \{s, t\}$ **do**
- 3: $u.current-neighbour \leftarrow u.neighbour-list-head$
- 4: **while** \exists active node u **do**
- 5: select active node u with highest label
- 6: discharge(u)

13.3 Highest label

Lemma 80

When using highest label the number of non-saturating pushes is only $\mathcal{O}(n^3)$.

After a non-saturating push from u a relabel is required to make a currently non-active node x , with $\ell(x) \geq \ell(u)$ active again (note that this includes u).

Hence, after n non-saturating pushes without an intermediate relabel there are no active nodes left.

Therefore, the number of non-saturating pushes is at most $n(\#relabels + 1) = \mathcal{O}(n^3)$.

13.3 Highest label

Since a discharge-operation is terminated by a non-saturating push this gives an upper bound of $\mathcal{O}(n^3)$ on the number of discharge-operations.

The cost for relabels and saturating pushes can be estimated in exactly the same way as in the case of relabel-to-front.

Question:

How do we find the next node for a discharge operation?

13.3 Highest label

Maintain lists L_i , $i \in \{0, \dots, 2n\}$, where list L_i contains active nodes with label i (maintaining these lists induces only constant additional cost for every push-operation and for every relabel-operation).

After a discharge operation terminated for a node u with label k , traverse the lists $k - 1, \dots, 0$, (in that order) until you find a non-empty list.

Unless the last (non-saturating) push was to s or t the list $k - 1$ must be non-empty (i.e., the search takes constant time).

13.3 Highest label

Hence, the total time required for searching for active nodes is at most

$$\mathcal{O}(n^3) + n(\#non-saturating-pushes-to-s-or-t)$$

Lemma 81

The number of non-saturating pushes to s or t is at most $\mathcal{O}(n^2)$.

With this lemma we get

Theorem 82

The push-relabel algorithm with the rule highest-label takes time $\mathcal{O}(n^3)$.

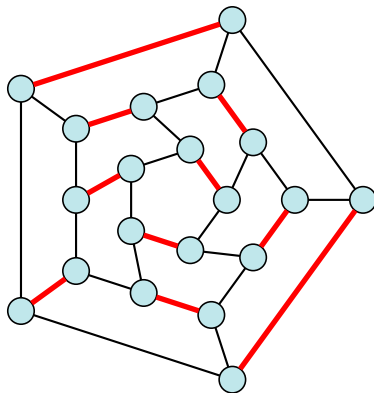
13.3 Highest label

Proof of the Lemma.

- ▶ We only show that the number of pushes to the source is at most $\mathcal{O}(n^2)$. A similar argument holds for the target.
- ▶ After a node v (which must have $\ell(v) = n + 1$) made a non-saturating push to the source there needs to be another node whose label is increased from $\leq n + 1$ to $n + 2$ before v can become active again.
- ▶ This happens for every push that v makes to the source. Since, every node can pass the threshold $n + 2$ at most once, v can make at most n pushes to the source.
- ▶ As this holds for every node the total number of pushes to the source is at most $\mathcal{O}(n^2)$.

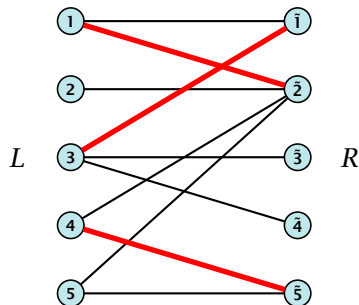
Matching

- ▶ Input: undirected graph $G = (V, E)$.
- ▶ $M \subseteq E$ is a **matching** if each node appears in at most one edge in M .
- ▶ Maximum Matching: find a matching of maximum cardinality



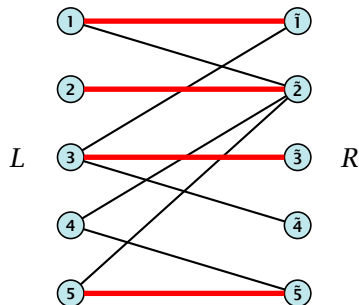
Bipartite Matching

- ▶ Input: undirected, **bipartite** graph $G = (L \uplus R, E)$.
- ▶ $M \subseteq E$ is a **matching** if each node appears in at most one edge in M .
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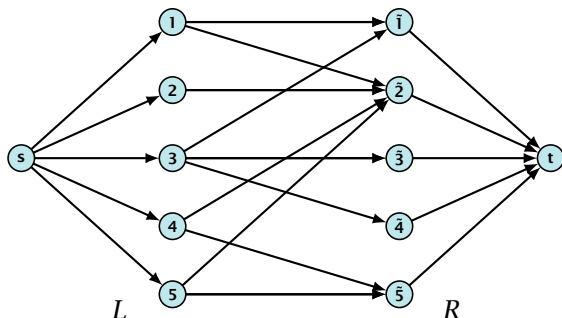
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Maxflow Formulation

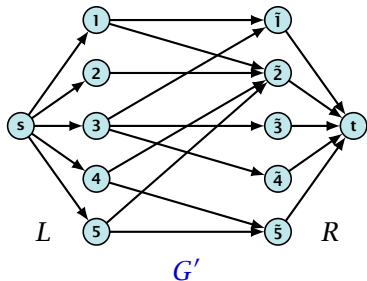
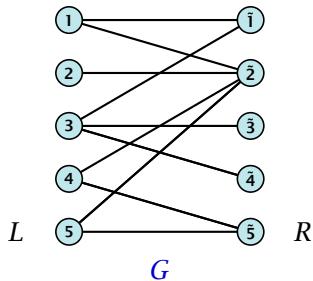
- ▶ Input: undirected, **bipartite** graph $G = (L \uplus R \uplus \{s, t\}, E')$.
- ▶ Direct all edges from L to R .
- ▶ Add source s and connect it to all nodes on the left.
- ▶ Add t and connect all nodes on the right to t .
- ▶ All edges have unit capacity.



Proof

Max cardinality matching in $G \leq$ value of maxflow in G'

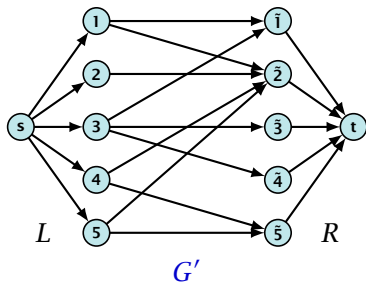
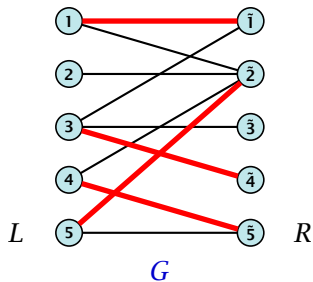
- ▶ Given a maximum matching M of cardinality k .
- ▶ Consider flow f that sends one unit along each of k paths.
- ▶ f is a flow and has cardinality k .



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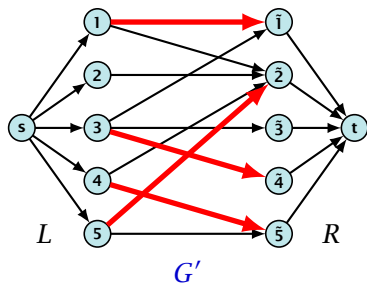
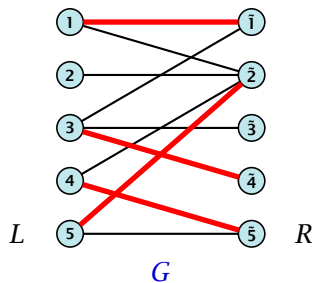
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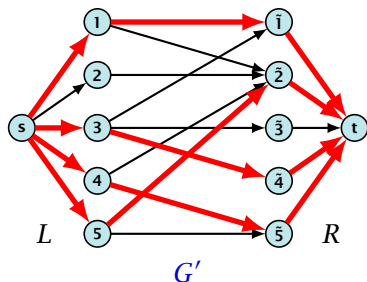
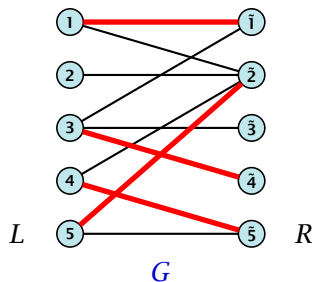
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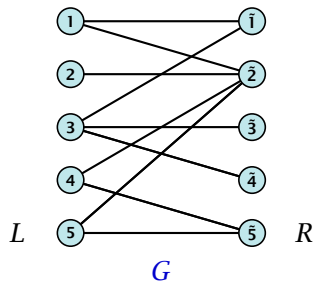
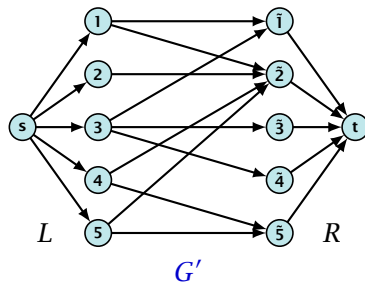
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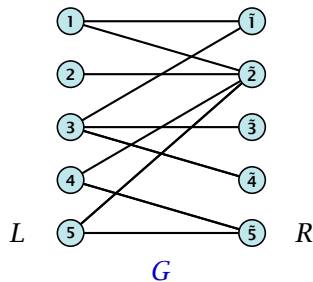
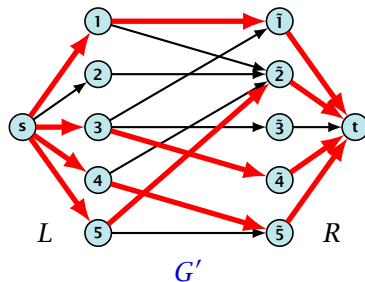
- ▶ Let f be a maxflow in G' of value k
- ▶ Integrality theorem $\Rightarrow k$ integral; we can assume f is 0/1.
- ▶ Consider $M =$ set of edges from L to R with $f(e) = 1$.
- ▶ Each node in L and R participates in at most one edge in M .
- ▶ $|M| = k$, as the flow must use at least k middle edges.



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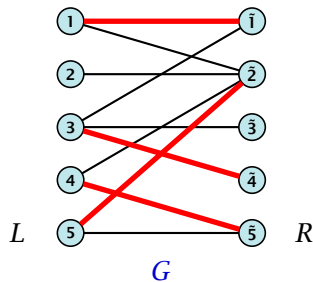
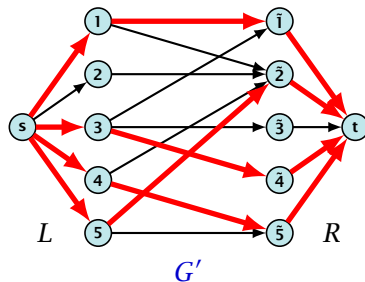
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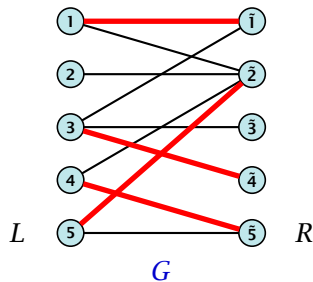
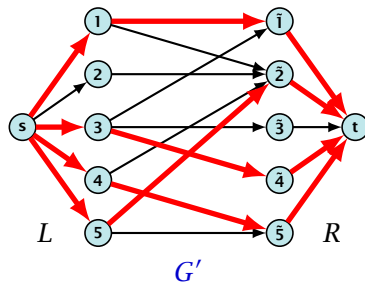
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14.1 Matching

Which flow algorithm to use?

- ▶ Generic augmenting path: $\mathcal{O}(m \text{val}(f^*)) = \mathcal{O}(mn)$.
- ▶ Capacity scaling: $\mathcal{O}(m^2 \log C) = \mathcal{O}(m^2)$.

Baseball Elimination

team i	wins w_i	losses ℓ_i	remaining games			
			Atl	Phi	NY	Mon
Atlanta	83	71	-	1	6	1
Philadelphia	80	79	1	-	0	2
New York	78	78	6	0	-	0
Montreal	77	82	1	2	0	-

Which team can end the season with most wins?

- ▶ Montreal is eliminated, since even after winning all remaining games there are only 80 wins.
- ▶ But also Philadelphia is eliminated. Why?

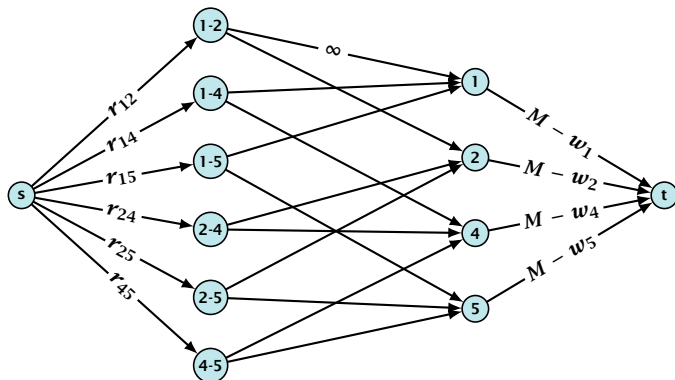
Baseball Elimination

Formal definition of the problem:

- ▶ Given a set S of teams, and one specific team $z \in S$.
- ▶ Team x has already won w_x games.
- ▶ Team x still has to play team y , r_{xy} times.
- ▶ Does team z still have a chance to finish with the most number of wins.

Baseball Elimination

Flow networks for $z = 3$. M is number of wins Team 3 can still obtain.

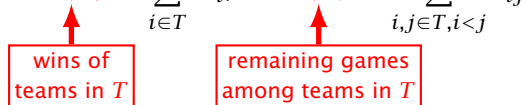


Idea. Distribute the results of remaining games in such a way that no team gets too many wins.

Certificate of Elimination

Let $T \subseteq S$ be a subset of teams. Define

$$w(T) := \sum_{i \in T} w_i, \quad r(T) := \sum_{i, j \in T, i < j} r_{ij}$$



If $\frac{w(T)+r(T)}{|T|} > M$ then one of the teams in T will have more than M wins in the end. A team that can win at most M games is therefore eliminated.

Theorem 83

A team z is eliminated if and only if the flow network for z does not allow a flow of value $\sum_{ij \in S \setminus \{z\}, i < j} r_{ij}$.

Proof (\Leftarrow)

- ▶ Consider the mincut A in the flow network. Let T be the set of **team-nodes** in A .
- ▶ If for a node x - y not both team nodes x and y are in T , then x - $y \notin A$ as otw. the cut would cut an infinite capacity edge.
- ▶ We don't find a flow that saturates all source edges:

$$\begin{aligned}r(S \setminus \{z\}) &> \text{cap}(S, V \setminus S) \\ &\geq \sum_{i < j: i \notin T \vee j \notin T} r_{ij} + \sum_{i \in T} (M - w_i) \\ &\geq r(S \setminus \{z\}) - r(T) + |T|M - w(T)\end{aligned}$$

- ▶ This gives $M < (w(T) + r(T))/|T|$, i.e., z is eliminated.

Baseball Elimination

Proof (\Rightarrow)

- ▶ Suppose we have a flow that saturates all source edges.
- ▶ We can assume that this flow is **integral**.
- ▶ For every pairing x - y it defines how many games team x and team y should win.
- ▶ The flow leaving the team-node x can be interpreted as the additional number of wins that team x will obtain.
- ▶ This is less than $M - w_x$ because of capacity constraints.
- ▶ Hence, we found a set of results for the remaining games, such that no team obtains more than M wins in total.
- ▶ Hence, team z is not eliminated.

Project Selection

Project selection problem:

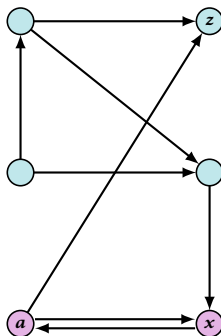
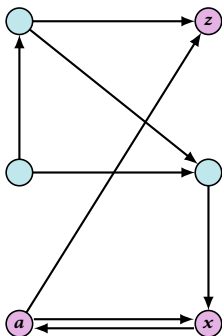
- ▶ Set P of possible projects. Project v has an associated profit p_v (can be positive or negative).
- ▶ Some projects have requirements (taking course EA2 requires course EA1).
- ▶ Dependencies are modelled in a graph. Edge (u, v) means “can’t do project u without also doing project v .”
- ▶ A subset A of projects is **feasible** if the prerequisites of every project in A also belong to A .

Goal: Find a feasible set of projects that maximizes the profit.

Project Selection

The prerequisite graph:

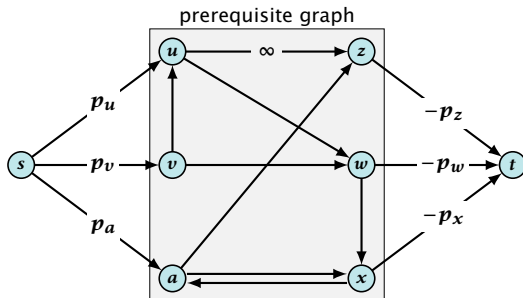
- ▶ $\{x, a, z\}$ is a feasible subset.
- ▶ $\{x, a\}$ is infeasible.



Project Selection

Mincut formulation:

- ▶ Edges in the prerequisite graph get infinite capacity.
- ▶ Add edge (s, v) with capacity p_v for nodes v with positive profit.
- ▶ Create edge (v, t) with capacity $-p_v$ for nodes v with negative profit.



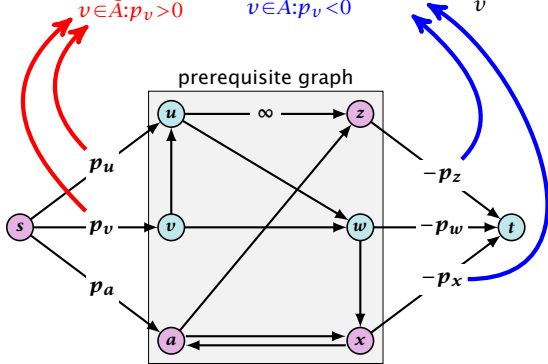
Theorem 84

A is a mincut if $A \setminus \{s\}$ is the optimal set of projects.

Proof.

▶ A is feasible because of capacity infinity edges.

▶ $\text{cap}(A, V \setminus A) = \sum_{v \in \bar{A}: p_v > 0} p_v + \sum_{v \in A: p_v < 0} (-p_v) = \sum_v p_v - \sum_{v \in A} p_v$



Mincost Flow

Consider the following problem:

$$\min \sum_e c(e) f(e)$$

$$\text{s.t. } \forall e \in E: 0 \leq f(e) \leq u(e)$$

$$\forall v \in V: f(v) = b(v)$$

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- ▶ $G = (V, E)$ is an **oriented graph**.
- ▶ $u : E \rightarrow \mathbb{R}_0^+ \cup \{\infty\}$ is the capacity function.
- ▶ $c : E \rightarrow \mathbb{R}$ is the cost function (note that $c(e)$ may be negative).
- ▶ $b : V \rightarrow \mathbb{R}, \sum_{v \in V} b(v) = 0$ is a demand function.

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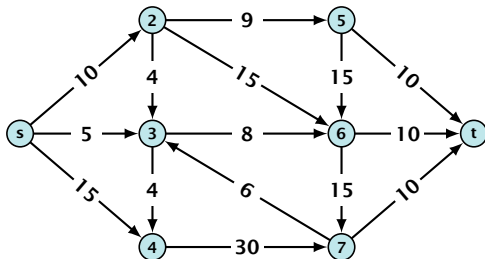
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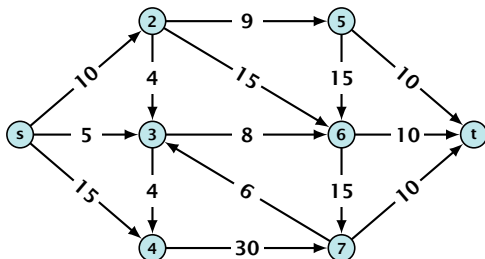
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Solve Maxflow Using Mincost Flow

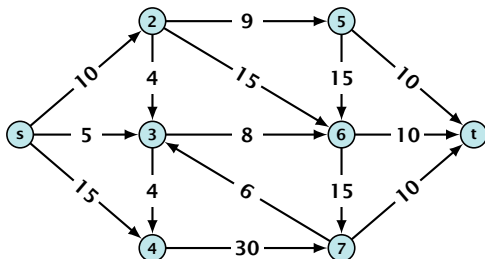


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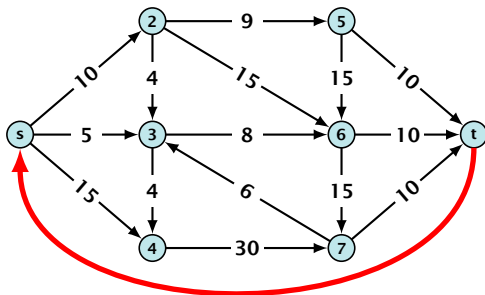
- ▶ Given a flow network for a standard maxflow problem.

Solve Maxflow Using Mincost Flow



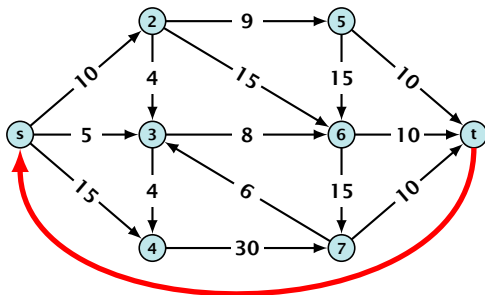
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- ▶ Add an edge from t to s with infinite capacity and cost -1 .

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- ▶ Add an edge from t to s with infinite capacity and cost -1 .
- ▶ Then, $\text{val}(f^*) = -\text{cost}(f_{\min})$, where f^* is a maxflow, and f_{\min} is a mincost-flow.

Solve Maxflow Using Mincost Flow

Solve decision version of maxflow:

- ▶ Given a flow network for a standard maxflow problem, and a value k .
- ▶ Set $b(v) = 0$ for every node apart from s or t . Set $b(s) = -k$ and $b(t) = k$.
- ▶ Set edge-costs to zero, and keep the capacities.
- ▶ There exists a maxflow of value k if and only if the mincost-flow problem is feasible.

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Generalization

Our model:

$$\begin{aligned} \min \quad & \sum_e c(e)f(e) \\ \text{s.t.} \quad & \forall e \in E: 0 \leq f(e) \leq u(e) \\ & \forall v \in V: f(v) = b(v) \end{aligned}$$

where $b: V \rightarrow \mathbb{R}$, $\sum_v b(v) = 0$; $u: E \rightarrow \mathbb{R}_0^+ \cup \{\infty\}$; $c: E \rightarrow \mathbb{R}$;

A more general model?

$$\begin{aligned} \min \quad & \sum_e c(e)f(e) \\ \text{s.t.} \quad & \forall e \in E: \ell(e) \leq f(e) \leq u(e) \\ & \forall v \in V: a(v) \leq f(v) \leq b(v) \end{aligned}$$

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Reduction I

$$\min \sum_e c(e) f(e)$$

$$\text{s.t. } \forall e \in E: \ell(e) \leq f(e) \leq u(e)$$

$$\forall v \in V: a(v) \leq f(v) \leq b(v)$$

We can assume that $a(v) = b(v)$:

add new node r

add edge (r, v) for all $v \in V$

set $\ell(e) = u(e) = 0$ for these

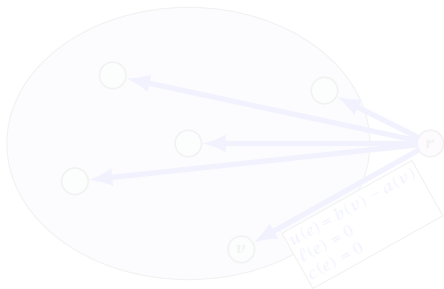
edges

set $a(v) = b(v) = a(v)$ for

edges (r, v)

set $a(r) = b(r)$ for all $v \in V$

$\forall v \in V: f(v) = \sum_{e \in E} c(e) f(e)$



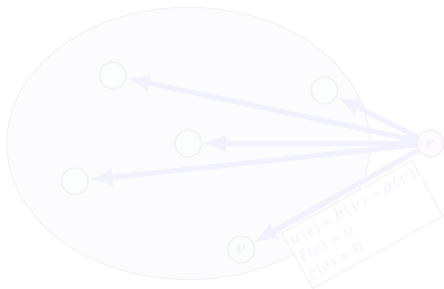
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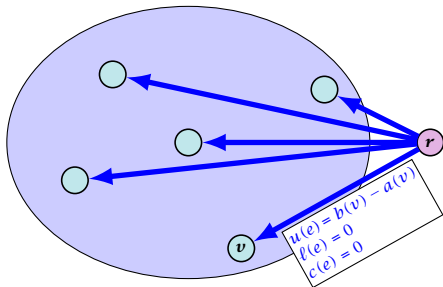
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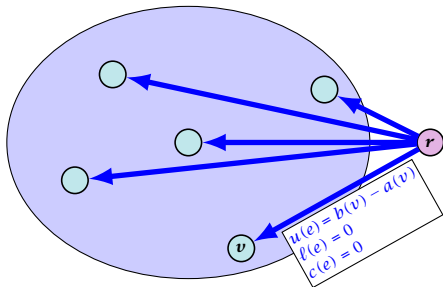
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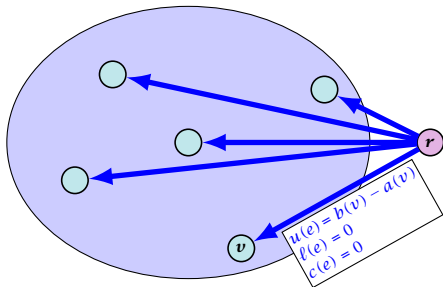
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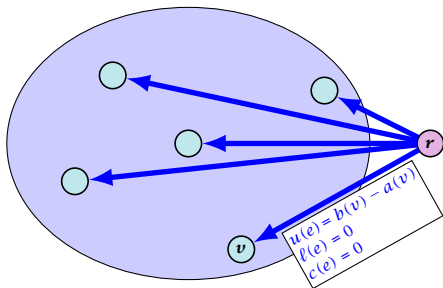
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Set $b(r) = \sum_{v \in V} b(v)$.



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We can assume that $a(v) = b(v)$:

Add new node r .

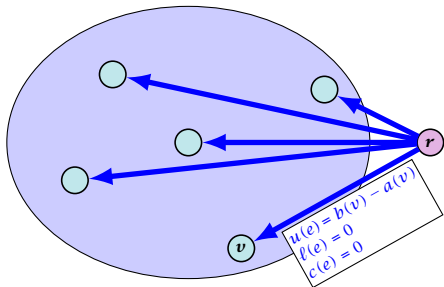
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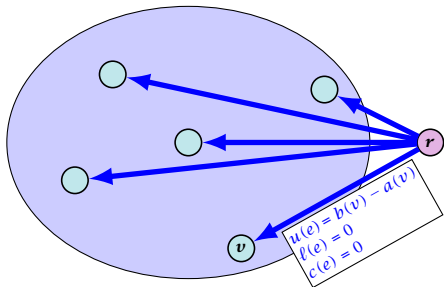
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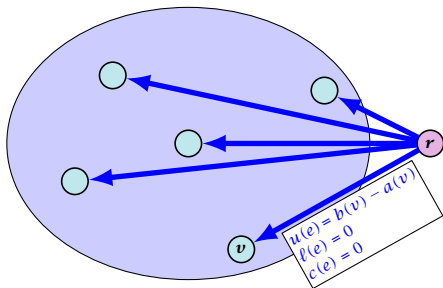
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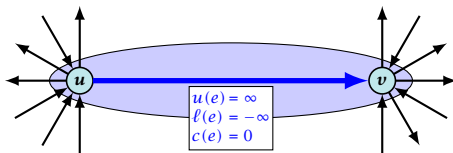
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We can assume that either $\ell(e) \neq -\infty$ or $u(e) \neq \infty$:

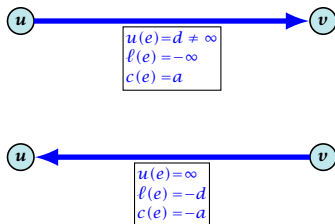


If $c(e) = 0$ we can simply contract the edge/identify nodes u and v

Reduction III

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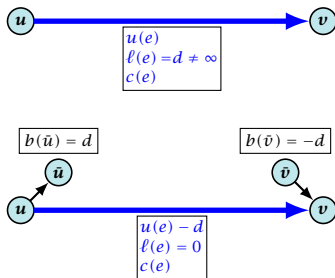


Replace the edge by an edge in opposite direction.

Reduction IV

$$\begin{aligned} \min \quad & \sum_e c(e)f(e) \\ \text{s.t.} \quad & \forall e \in E: \ell(e) \leq f(e) \leq u(e) \\ & \forall v \in V: f(v) = b(v) \end{aligned}$$

We can assume that $\ell(e) = 0$:



The added edges have infinite capacity and cost $c(e)/2$.

Applications

Caterer Problem

- ▶ She needs to supply r_i napkins on N successive days.
- ▶ She can buy new napkins at p cents each.
- ▶ She can launder them at a fast laundry that takes m days and cost f cents a napkin.
- ▶ She can use a slow laundry that takes $k > m$ days and costs s cents each.
- ▶ At the end of each day she should determine how many to send to each laundry and how many to buy in order to fulfill demand.
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The residual graph for a mincost flow is exactly defined as the residual graph for standard flows, with the only exception that one needs to define a cost for the residual edge.

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A **circulation** in a graph $G = (V, E)$ is a function $f : E \rightarrow \mathbb{R}^+$ that has an excess flow $f(v) = 0$ for every node $v \in V$ (G may be a directed graph instead of just an oriented graph).

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- ⇒ Suppose that g is a feasible circulation of negative cost in the residual graph.

Then $f + g$ is a feasible flow with cost $\text{cost}(f) + \text{cost}(g) < \text{cost}(f)$. Hence, f is not minimum cost.

- ⇐ Let f be a non-mincost flow, and let f^* be a min-cost flow. We need to show that the residual graph has a feasible circulation with negative cost.

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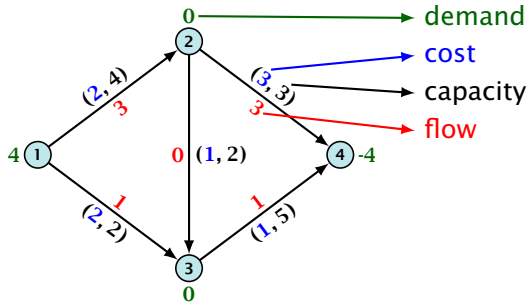
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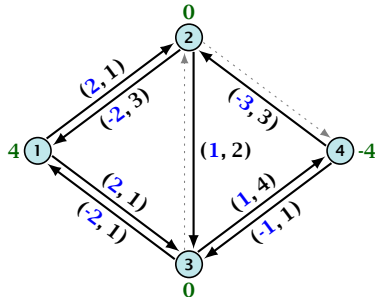
Algorithm 51 CycleCanceling($G = (V, E), c, u, b$)

- 1: establish a feasible flow f in G
- 2: **while** G_f contains negative cycle **do**
- 3: use Bellman-Ford to find a negative circuit Z
- 4: $\delta \leftarrow \min\{u_f(e) \mid e \in Z\}$
- 5: augment δ units along Z and update G_f

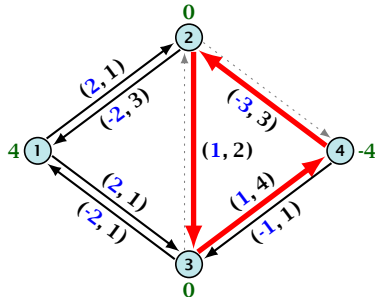
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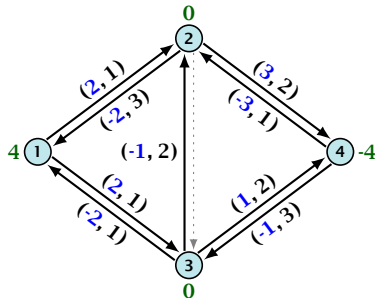
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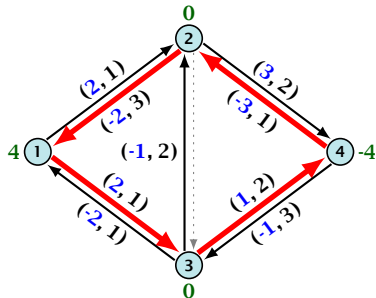
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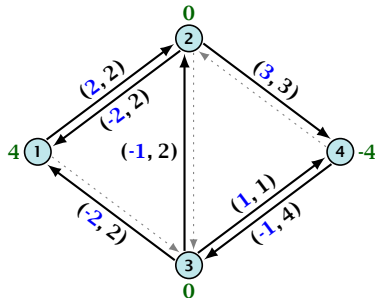
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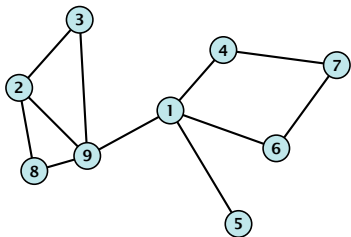
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Lemma 87

The improving cycle algorithm runs in time $\mathcal{O}(nm^2CU)$, for integer capacities and costs, when for all edges e , $|c(e)| \leq C$ and $|u(e)| \leq U$.

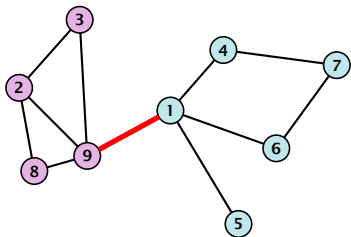
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Given an **undirected, capacitated graph** $G = (V, E, c)$ find a partition of V into two non-empty sets $S, V \setminus S$ s.t. the capacity of edges between both sets is minimized.



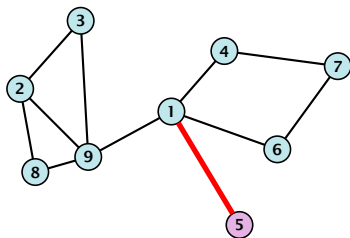
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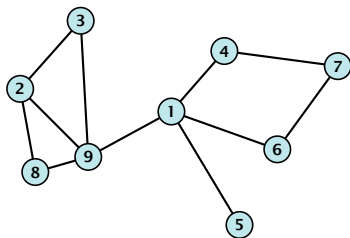
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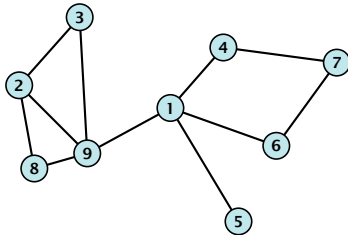
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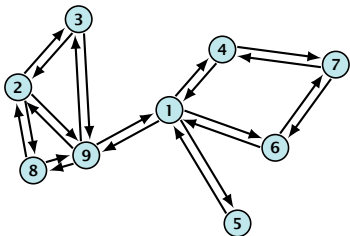
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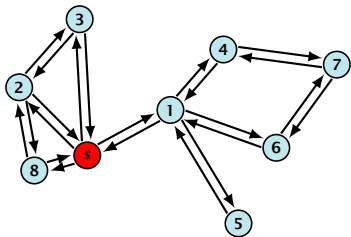
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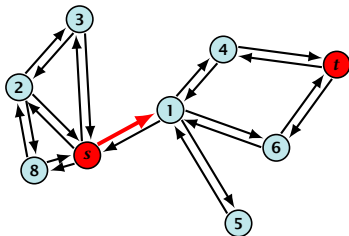
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- ▶ Fix an arbitrary node $s \in V$ as source. Compute a minimum s - t cut for all possible choices $t \in V, t \neq s$. (Time: $\mathcal{O}(n^4)$)



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- ▶ Let $(S, V \setminus S)$ be a minimum global mincut. The above algorithm will output a cut of capacity $\text{cap}(S, V \setminus S)$ whenever $|\{s, t\} \cap S| = 1$.



Edge Contractions

- ▶ Given a graph $G = (V, E)$ and an edge $e = \{u, v\}$.
- ▶ The graph G/e is obtained by “identifying” u and v to form a new node.
- ▶ Resulting parallel edges are replaced by a single edge, whose capacity equals the sum of capacities of the parallel edges.

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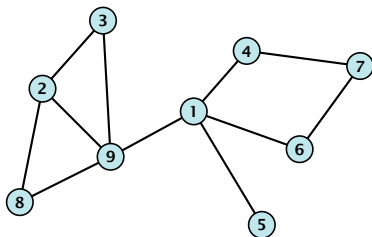


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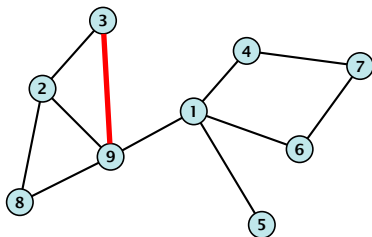


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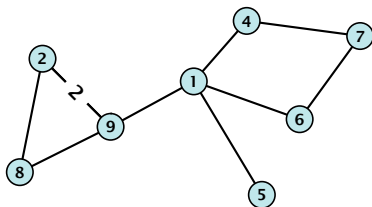


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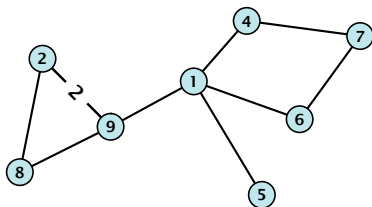


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Edge Contractions

We can perform an edge-contraction in time $\mathcal{O}(n)$.

Randomized Mincut Algorithm

Algorithm 52 KargerMincut($G = (V, E, c)$)

- 1: **for** $i = 1 \rightarrow n - 2$ **do**
- 2: choose $e \in E$ randomly with probability $c(e)/C(E)$
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1: for  $i = 1 \rightarrow n - 2$  do
2:     choose  $e \in E$  randomly with probability  $c(e)/C(E)$ 
3:      $G \leftarrow G/e$ 
4: return only cut in  $G$ 
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- ▶ Let G_t denote the graph after the $(n - t)$ -th iteration, when t nodes are left.
- ▶ Note that the final graph G_2 only contains a single edge.
- ▶ The cut in G_2 corresponds to a cut in the original graph G with the same capacity.

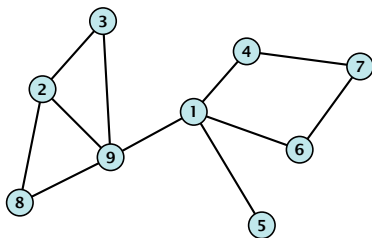
Randomized Mincut Algorithm

Algorithm 52 KargerMincut($G = (V, E, c)$)

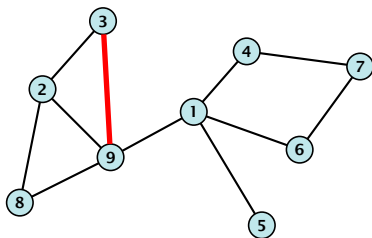
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- ▶ What is the probability that this algorithm returns a mincut?

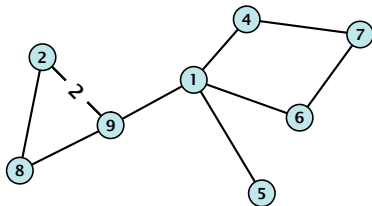
Example: Randomized Mincut Algorithm



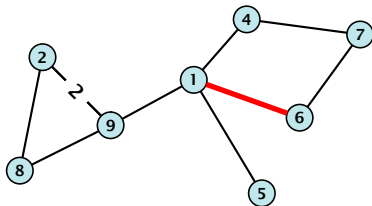
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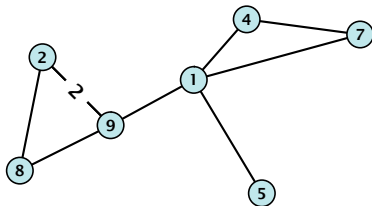
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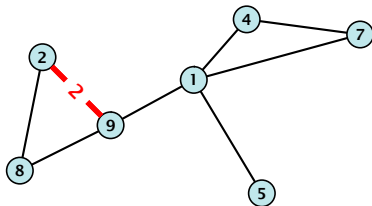
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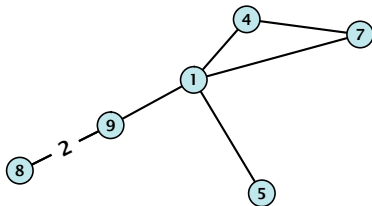
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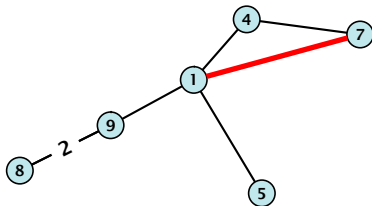
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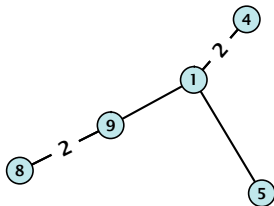
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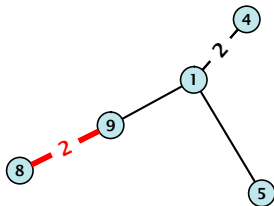
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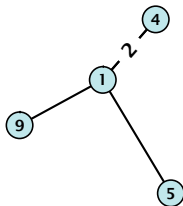
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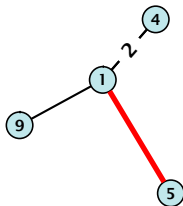
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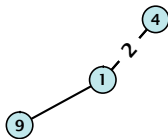
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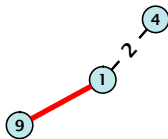
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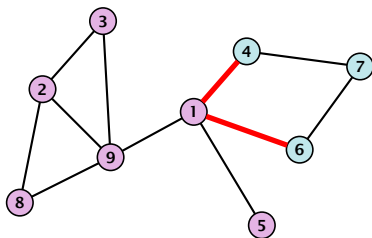
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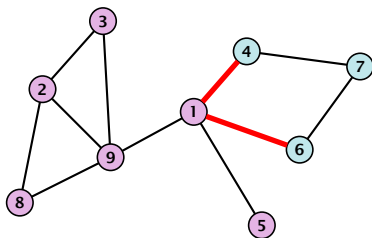
Example: Randomized Mincut Algorithm



Example: Randomized Mincut Algorithm



Example: Randomized Mincut Algorithm



What is the probability that this algorithm returns a mincut?

What is the probability that a given mincut A is still possible after round i ?

- ▶ It is still possible to obtain cut A in the end if so far **no** edge in $(A, V \setminus A)$ has been contracted.

Analysis

What is the probability that we select an edge from A in iteration i ?

- ▶ Let $\min = \text{cap}(A, V \setminus A)$ denote the capacity of a mincut.
- ▶ Let $\text{cap}(v)$ be capacity of edges incident to vertex $v \in V_{n-i+1}$.
- ▶ Clearly, $\text{cap}(v) \geq \min$.
- ▶ Summing $\text{cap}(v)$ over all edges gives

$$2c(E) = 2 \sum_{e \in E} c(e) = \sum_{v \in V} \text{cap}(v) \geq (n - i + 1) \cdot \min$$

- ▶ Hence, the probability of choosing an edge from the cut is at most $\min / c(E) \leq 2 / (n - i + 1)$.

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$$1 - \frac{2}{n - i + 1} = \frac{n - i - 1}{n - i + 1} .$$

The probability that the cut is alive after iteration $n - t$ (after which t nodes are left) is

$$\prod_{i=1}^{n-t} \frac{n - i - 1}{n - i + 1} = \frac{t(t-1)}{n(n-1)} .$$

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Repeating the algorithm $c \ln n \binom{n}{2}$ times gives that the probability that we are never successful is

$$\left(1 - \frac{1}{\binom{n}{2}}\right)^{\binom{n}{2} c \ln n} \leq \left(e^{-1/\binom{n}{2}}\right)^{\binom{n}{2} c \ln n} \leq n^{-c},$$

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Theorem 89

The randomized mincut algorithm computes an optimal cut with high probability. The total running time is $\mathcal{O}(n^4 \log n)$.

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Improved Algorithm

Algorithm 53 RecursiveMincut($G = (V, E, c)$)

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2:   choose  $e \in E$  randomly with probability  $c(e)/C(E)$ 
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5:  $cuta \leftarrow$  RecursiveMincut( $G$ );
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7: return  $\min\{cuta, cutb\}$ 
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Running time:

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- ▶ $T(n) = 2T\left(\frac{n}{\sqrt{2}}\right) + \mathcal{O}(n^2)$
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Probability of Success

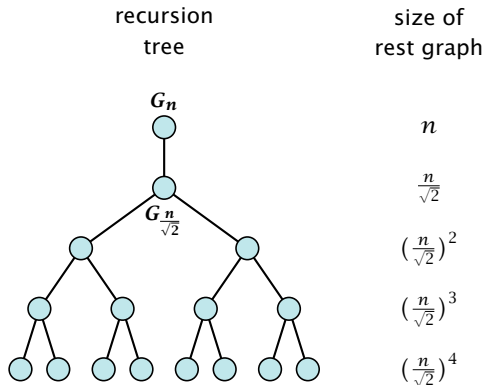
The probability of contracting an edge from the mincut during one iteration through the for-loop is only

$$\frac{t(t-1)}{n(n-1)} \approx \frac{t^2}{n^2} = \frac{1}{2},$$

as $t = \frac{n}{\sqrt{2}}$.

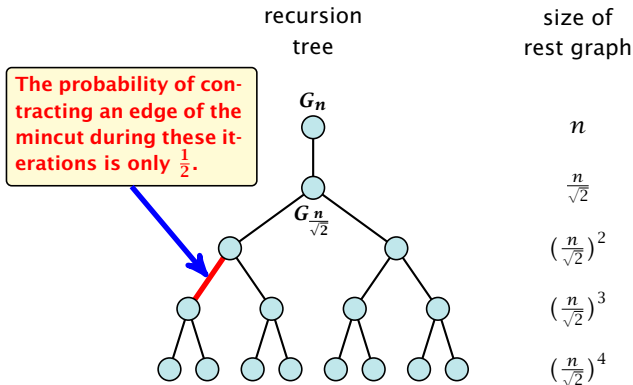
For the following analysis we ignore the slight error and assume that this probability is at most $\frac{1}{2}$.

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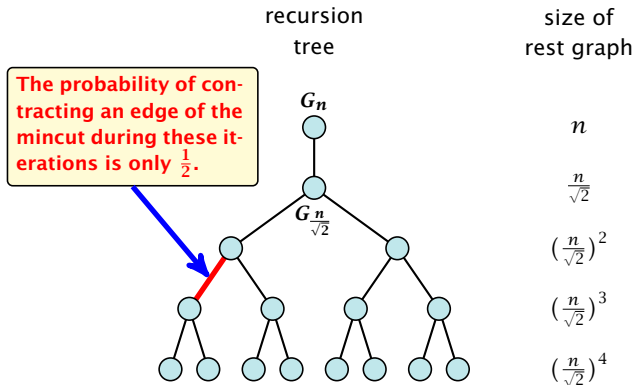
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Let for an edge e in the recursion tree, $h(e)$ denote the height (distance to leaf level) of the parent-node of e (end-point that is higher up in the tree). Let h denote the height of the root node.

Call an edge e **alive** if there exists a path from the parent-node of e to a descendant leaf, after we randomly deleted edges. Note that an edge can only be alive if it hasn't been deleted.

Lemma 90

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- ▶ This happens with probability

$$p_d = \frac{1}{2} (2p_{d-1} - p_{d-1}^2) \quad \boxed{\Pr[A \vee B] = \Pr[A] + \Pr[B] - \Pr[A \wedge B]}$$

$$= p_{d-1} - \frac{p_{d-1}^2}{2}$$

$$\geq \frac{1}{d} - \frac{1}{2d^2} \geq \frac{1}{d} - \frac{1}{d(d+1)} = \frac{1}{d+1} .$$

$x - x^2/2$ is monotonically increasing for $x \in [0, 1]$

16 Global Mincut

Lemma 91

One run of the algorithm can be performed in time $\mathcal{O}(n^2 \log n)$ and has a success probability of $\Omega(\frac{1}{\log n})$.

Doing $\Theta(\log^2 n)$ runs gives that the algorithm succeeds with high probability. The total running time is $\mathcal{O}(n^2 \log^3 n)$.

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17 Gomory Hu Trees

Given an undirected, weighted graph $G = (V, E, c)$ a **cut-tree** $T = (V, F, w)$ is a tree with edge-set F and capacities w that fulfills the following properties.

1. **Equivalent Flow Tree:** For any pair of vertices $s, t \in V$, $f(s, t)$ in G is equal to $f_T(s, t)$.
2. **Cut Property:** A minimum s - t cut in T is also a minimum cut in G .

Here, $f(s, t)$ is the value of a maximum s - t flow in G , and $f_T(s, t)$ is the corresponding value in T .

Overview of the Algorithm

The algorithm maintains a partition of V , (sets S_1, \dots, S_t), and a spanning tree T on the vertex set $\{S_1, \dots, S_t\}$.

Initially, there exists only the set $S_1 = V$.

Then the algorithm performs $n - 1$ split-operations:

- In each such split-operation it chooses a set S_i with $|S_i| \geq 2$ and splits this set into two non-empty parts X and Y .
- S_i is then removed from T and replaced by X and Y .
- The edges of T incident to S_i are kept, and the split-edges are added to T .
- The split-edges are added to S_i and removed from X and Y .

In the end this gives a tree on the vertex set V .

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In each such split-operation it chooses a set S_i with $|S_i| \geq 2$ and splits it into two non-empty parts X and Y . The edges that are removed from T and replaced by X and Y are those that connect vertices in S_i and the other sets S_j . The new spanning tree T' is obtained by $T - E + X + Y$.

In the end this gives a tree on the vertex set V .

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Details of the Split-operation

- ▶ Select S_i that contains at least two nodes a and b .
- ▶ Compute the connected components of the forest obtained from the current tree T after deleting S_i . Each of these components corresponds to a set of vertices from V .
- ▶ Consider the graph H obtained from G by contracting these connected components into single nodes.
- ▶ Compute a minimum a - b cut in H . Let A , and B denote the two sides of this cut.
- ▶ Split S_i in T into two sets/nodes $S_i^a := S_i \cap A$ and $S_i^b := S_i \cap B$ and add edge $\{S_i^a, S_i^b\}$ with capacity $f_H(a, b)$.
- ▶ Replace an edge $\{S_i, S_x\}$ by $\{S_i^a, S_x\}$ if $S_x \subset A$ and by $\{S_i^b, S_x\}$ if $S_x \subset B$.

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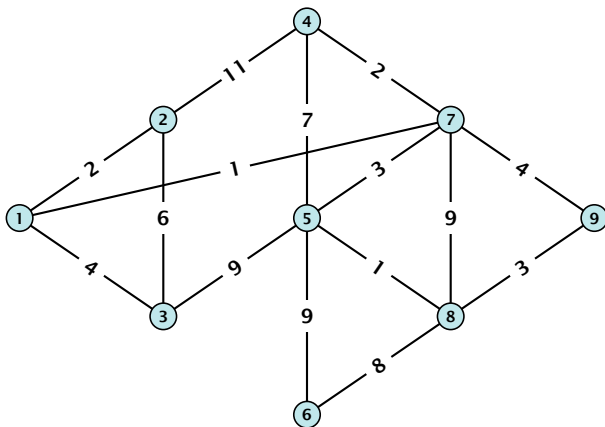
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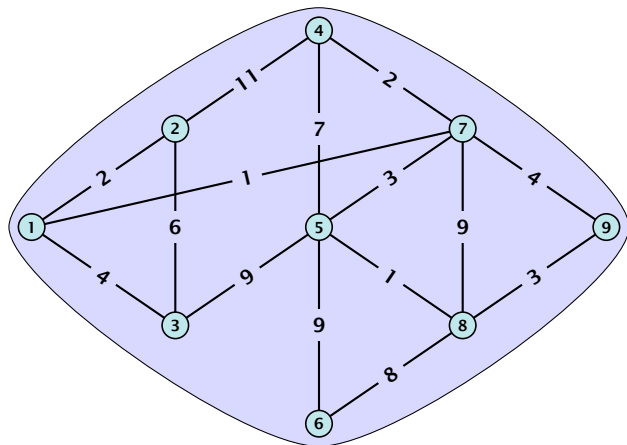
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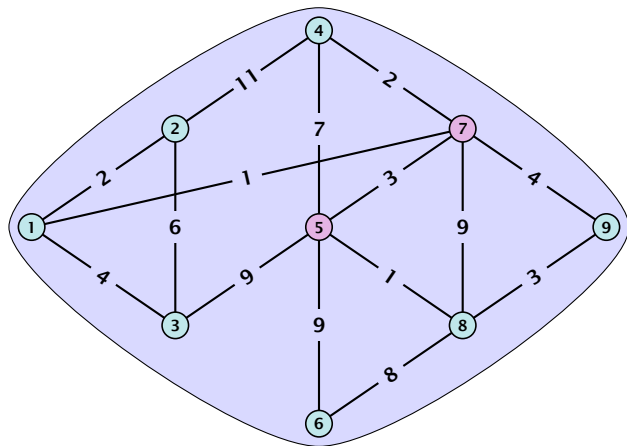
Example: Gomory-Hu Construction



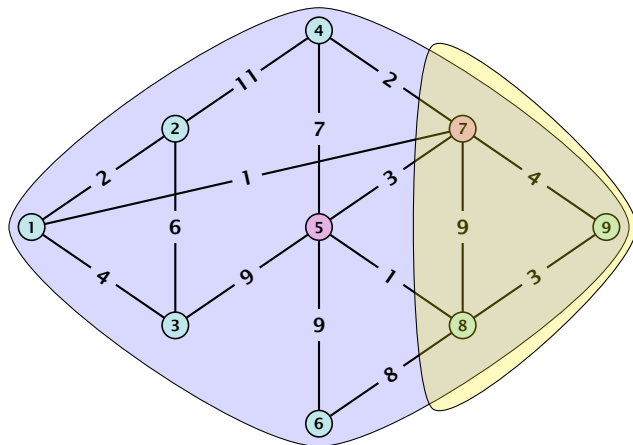
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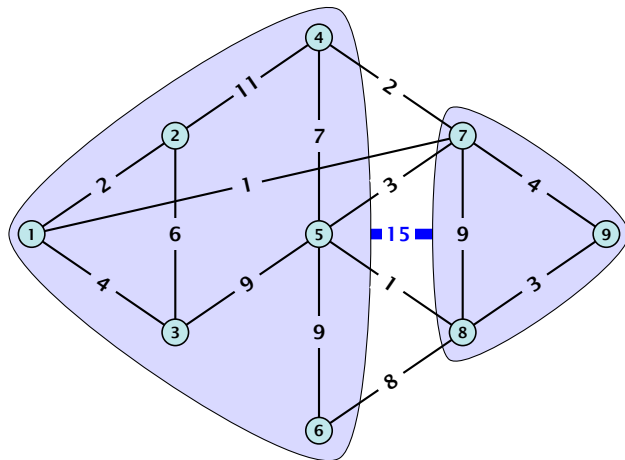
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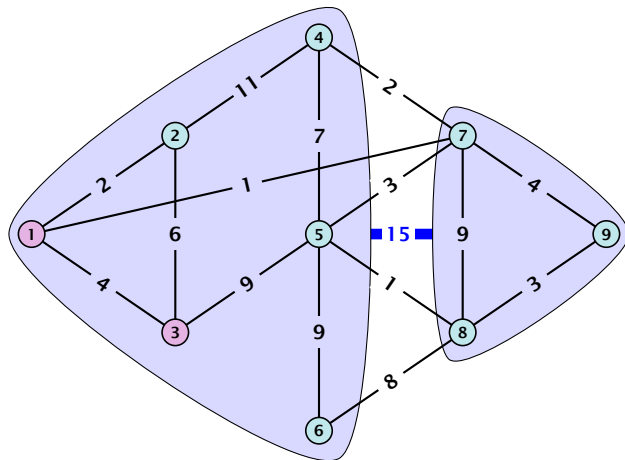
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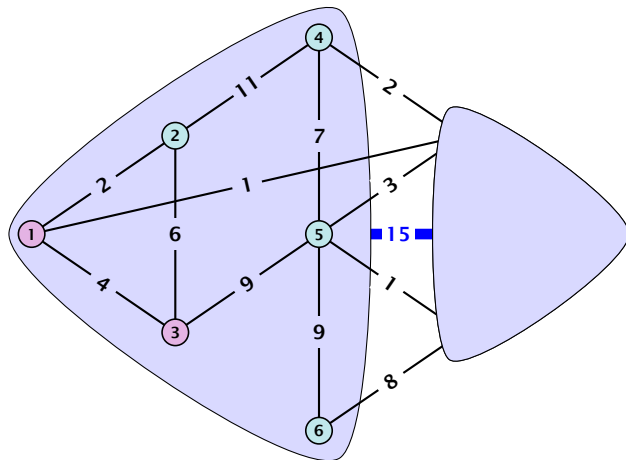
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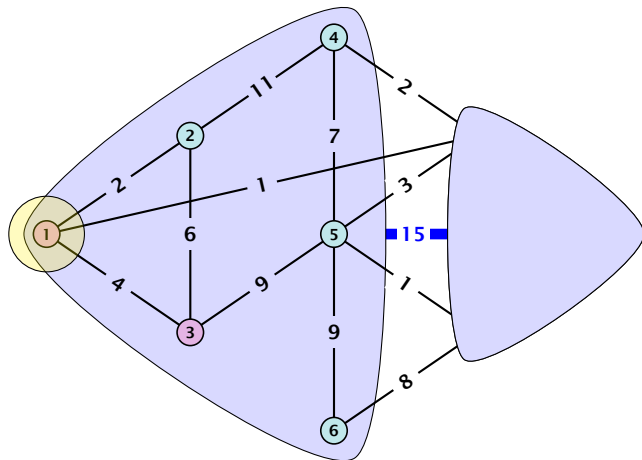
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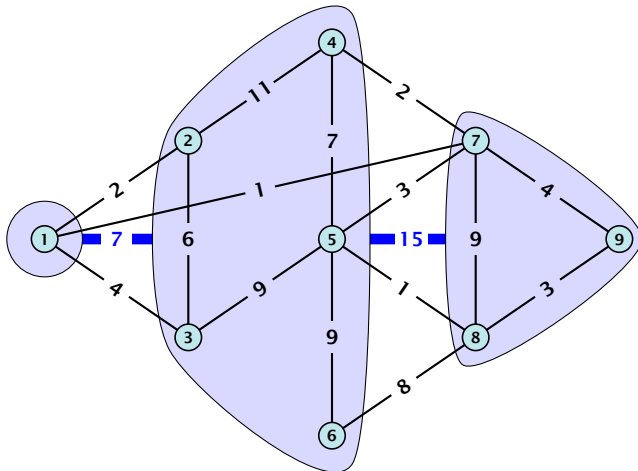
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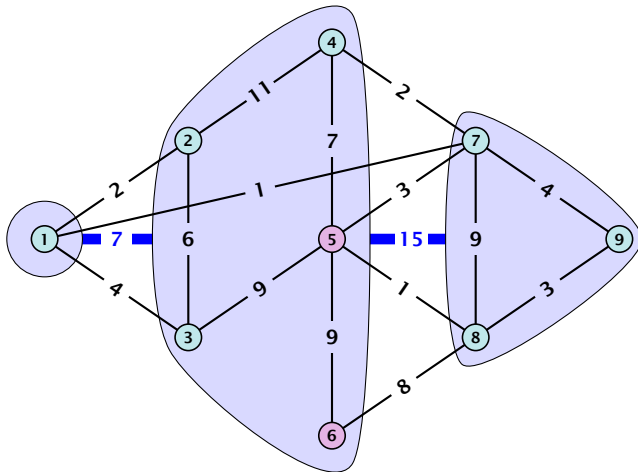
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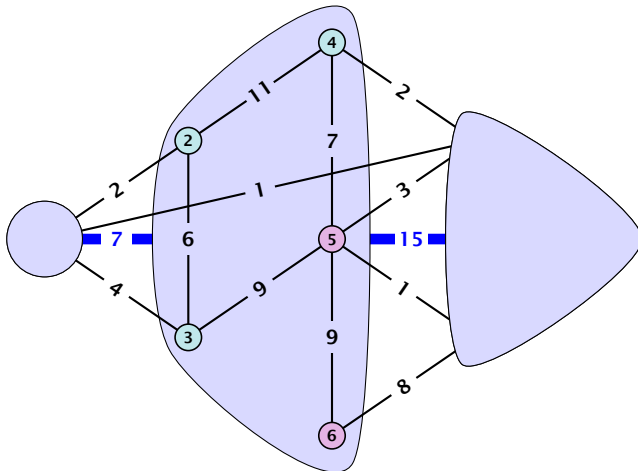
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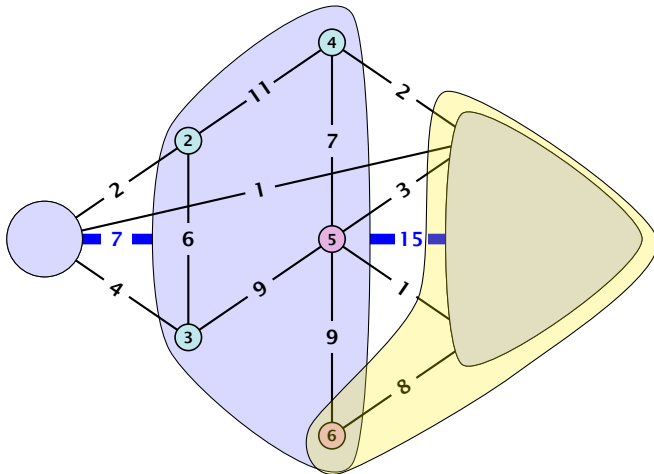
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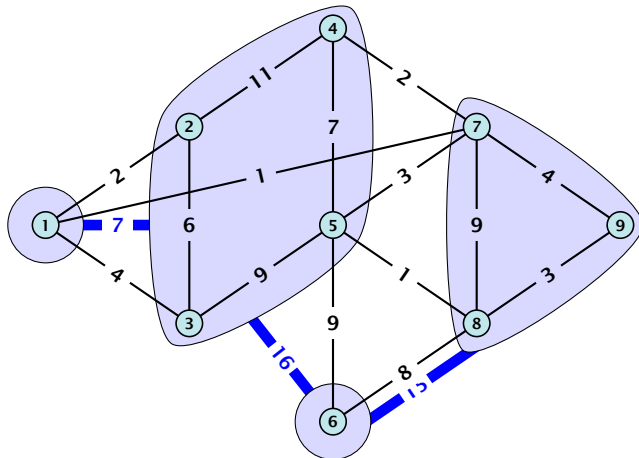
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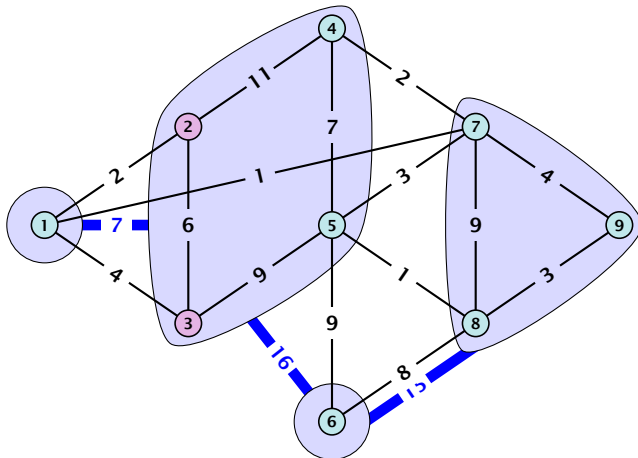
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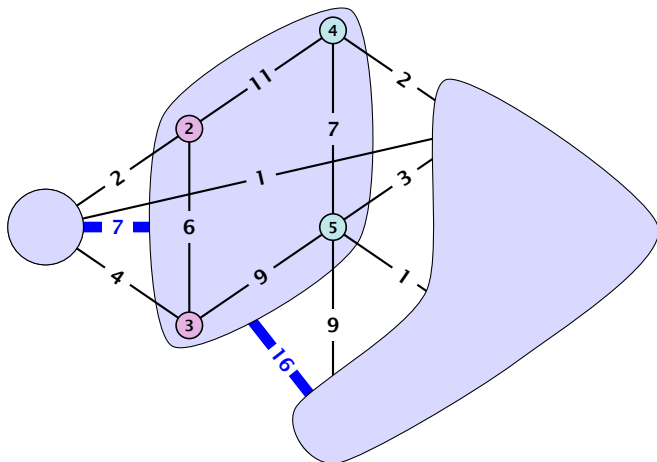
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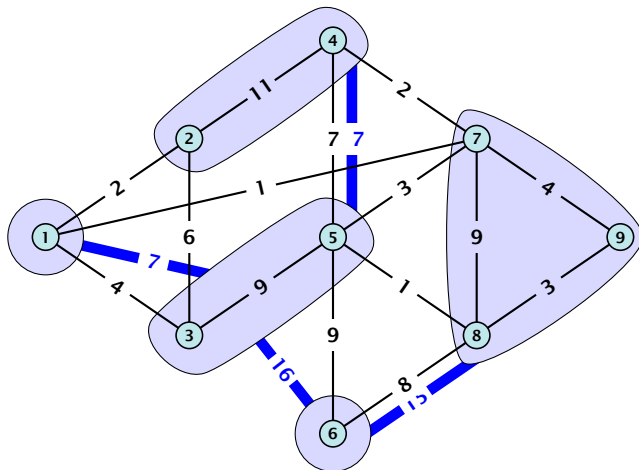
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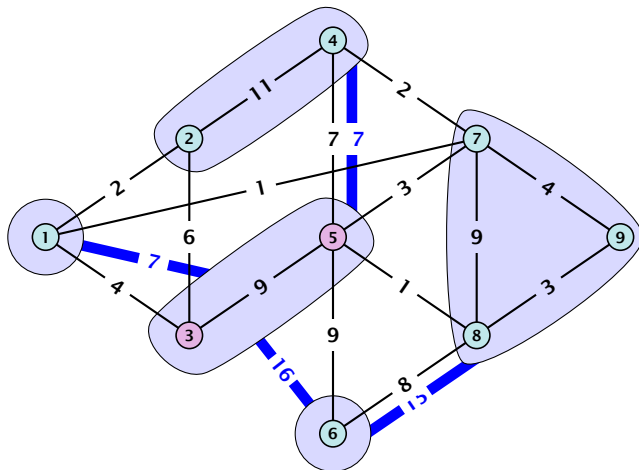
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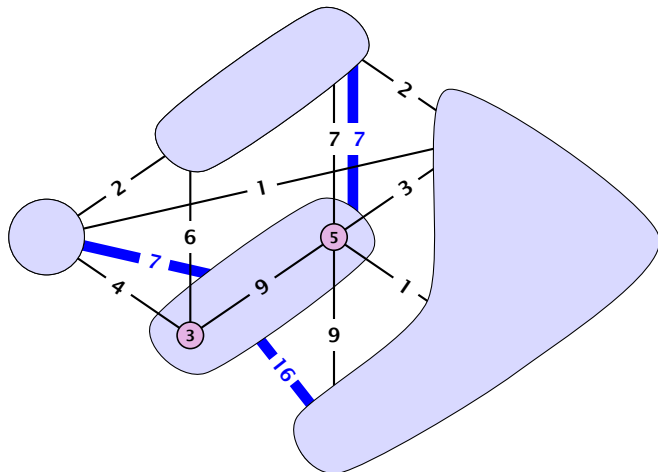
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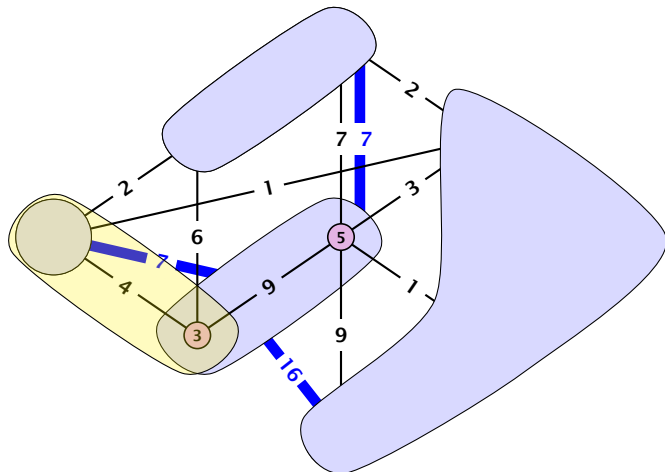
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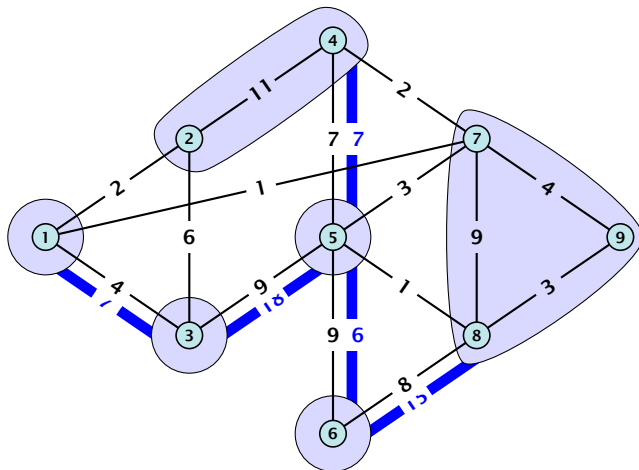
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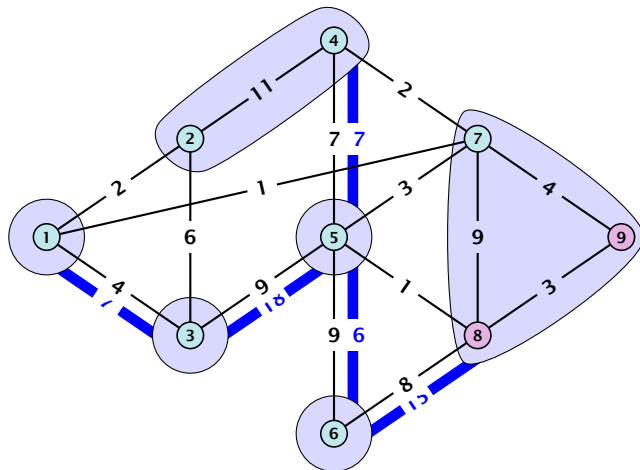
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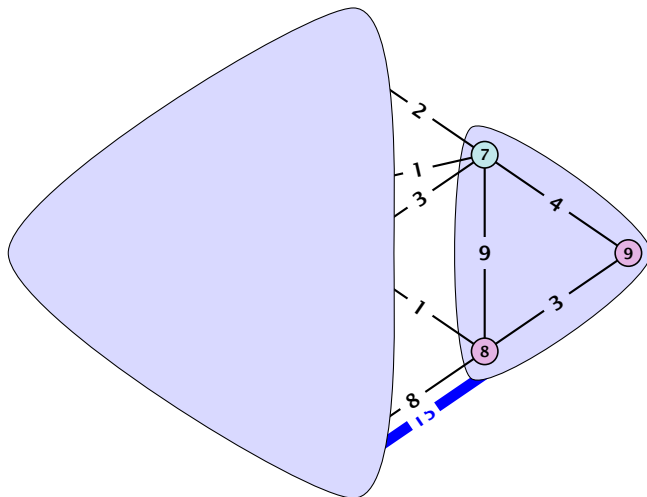
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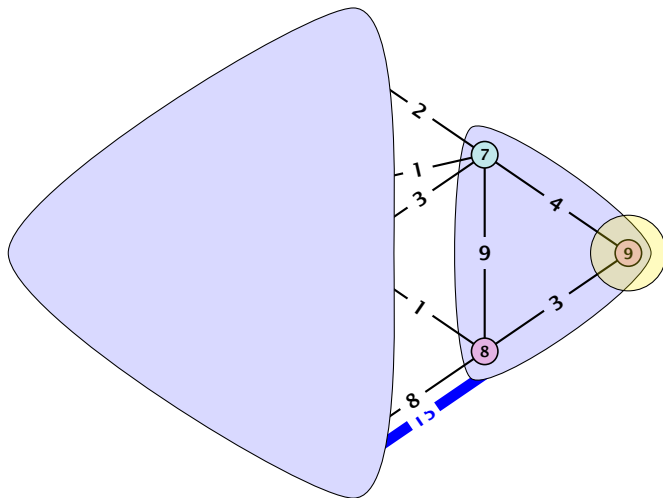
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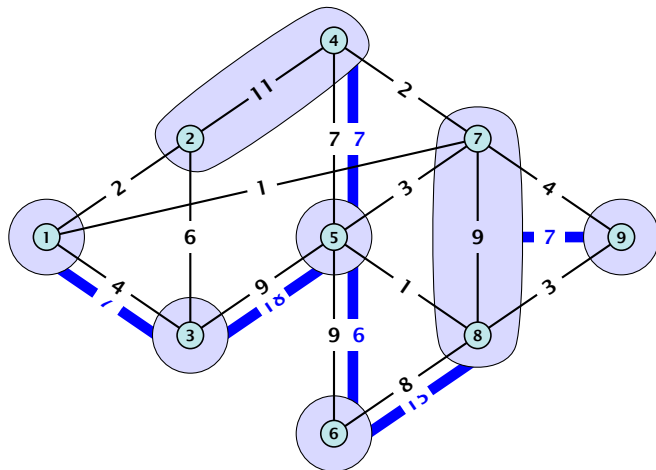
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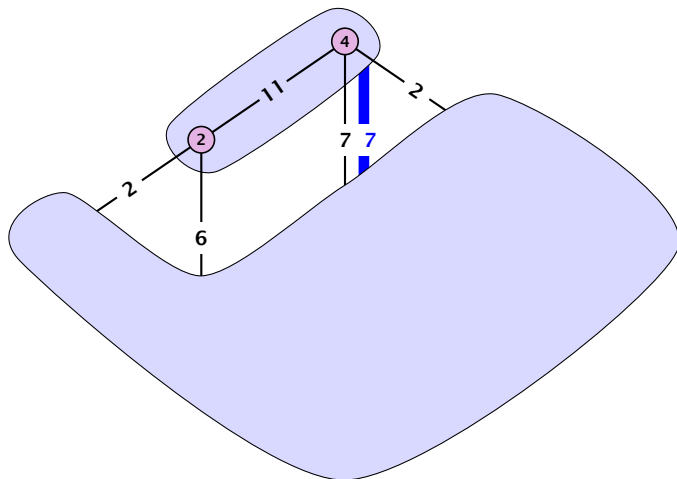
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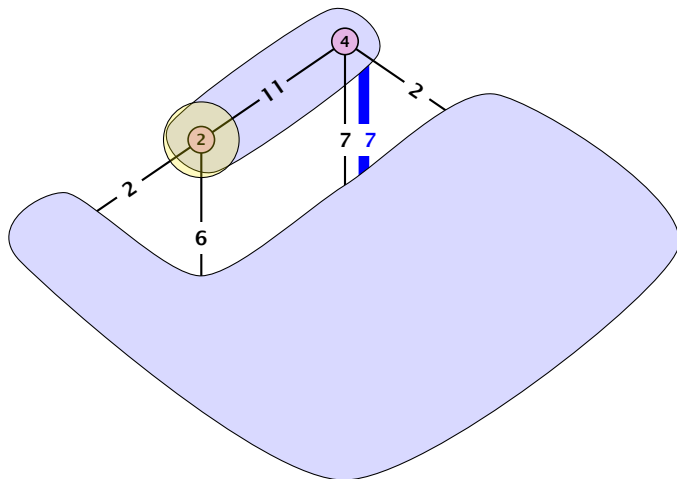
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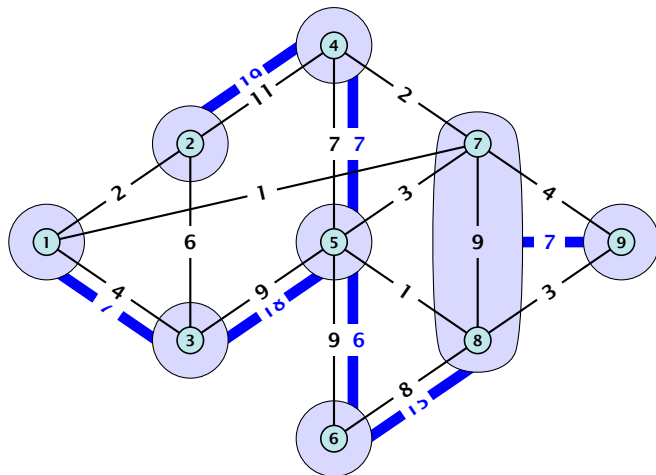
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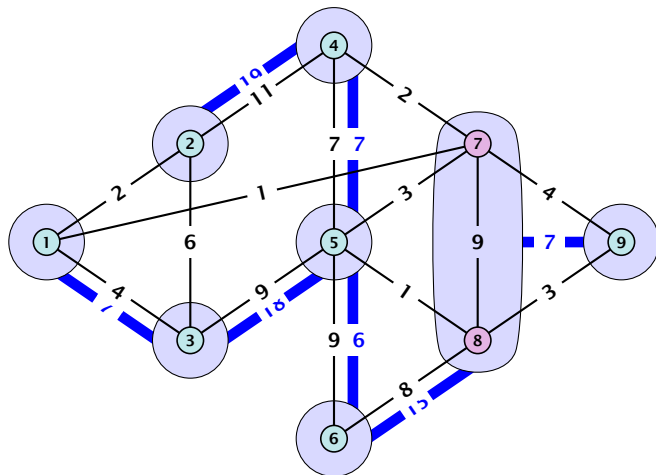
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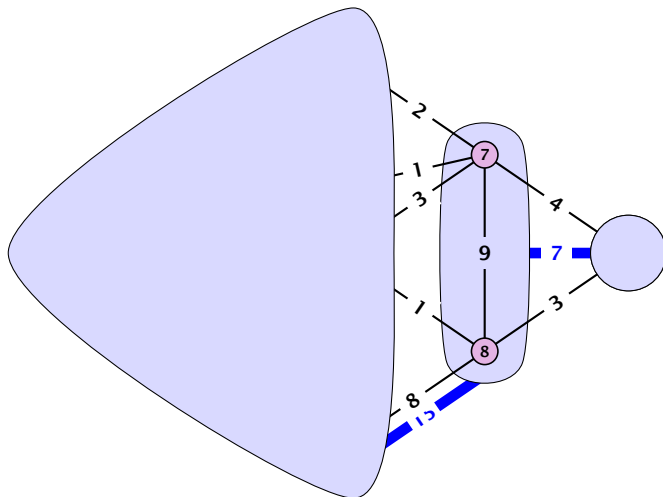
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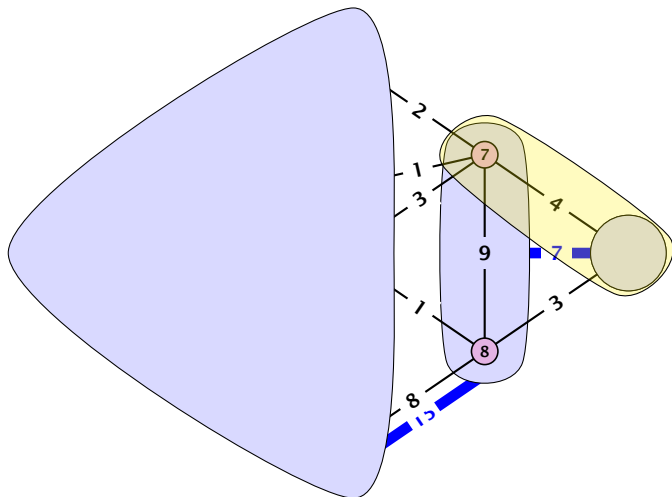
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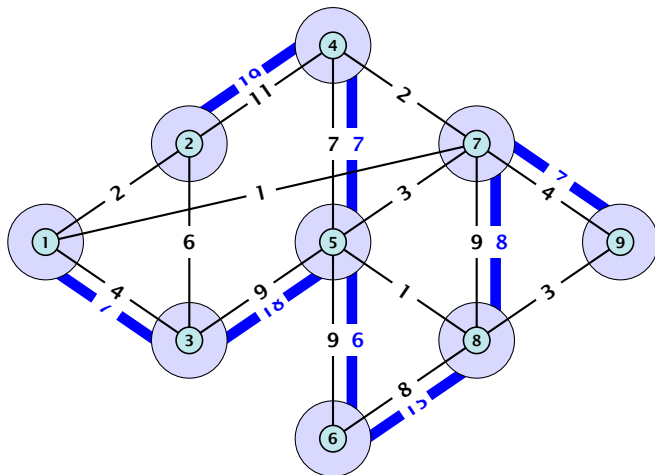
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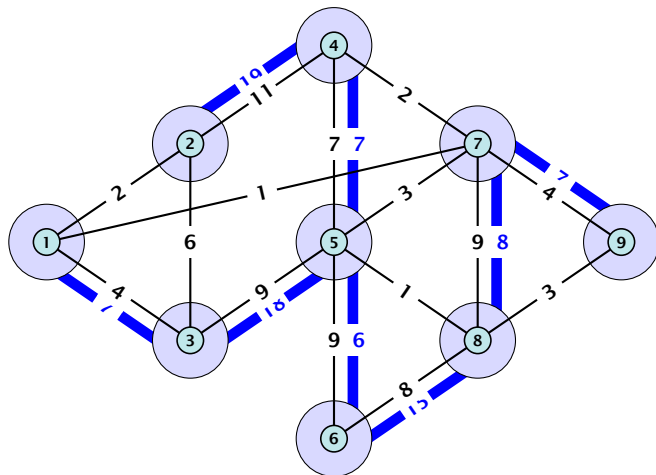
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Lemma 92

For nodes $s, t, x \in V$ we have $f(s, t) \geq \min\{f(s, x), f(x, t)\}$

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Analysis

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Lemma 94

Let S be some minimum r - s cut for some nodes $r, s \in V$ ($s \in S$), and let $v, w \in S$. Then there is a minimum v - w -cut T with $T \subset S$.

Proof: Let X be a minimum v - w cut with $v \in X$ and $w \notin X$. If $X \subset S$, then we are done. Note that $S \setminus X$ and $X \cap S$ are proper cuts.

We may assume w.l.o.g. $s \in X$.

First case $r \in X$.

$\text{cap}(S \setminus S) + \text{cap}(S \setminus X) \leq \text{cap}(S) + \text{cap}(X)$
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Therefore $\text{cap}(S \setminus X) \leq \text{cap}(X)$.
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Second case $r \notin X$.

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Proof: Let X be a minimum v - w cut with $X \cap S \neq \emptyset$ and $X \cap (V \setminus S) \neq \emptyset$. Note that $S \setminus X$ and $S \cap X$ are v - w cuts inside S .

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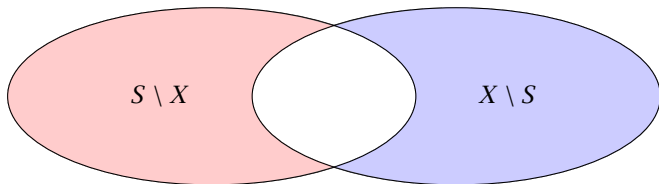
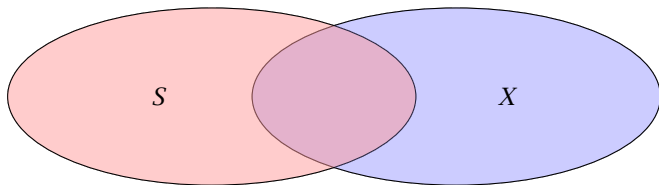
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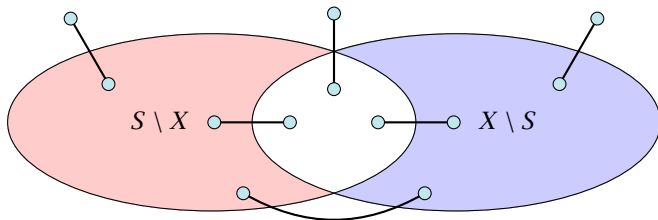
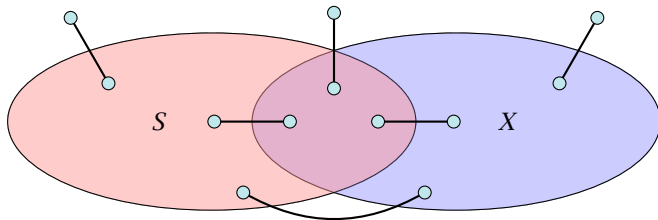
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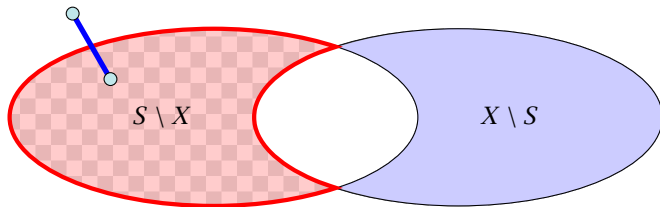
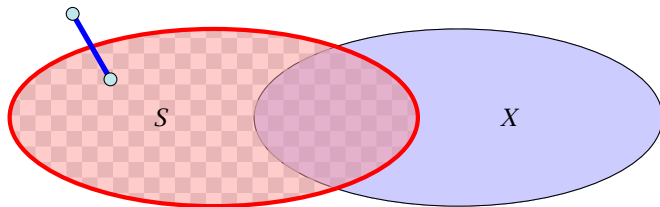
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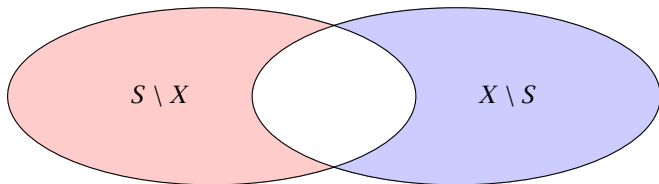
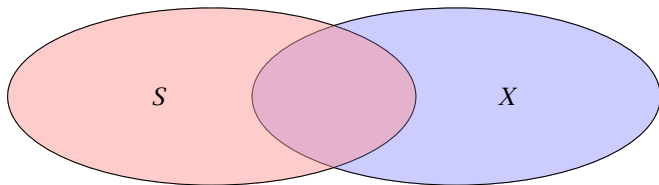
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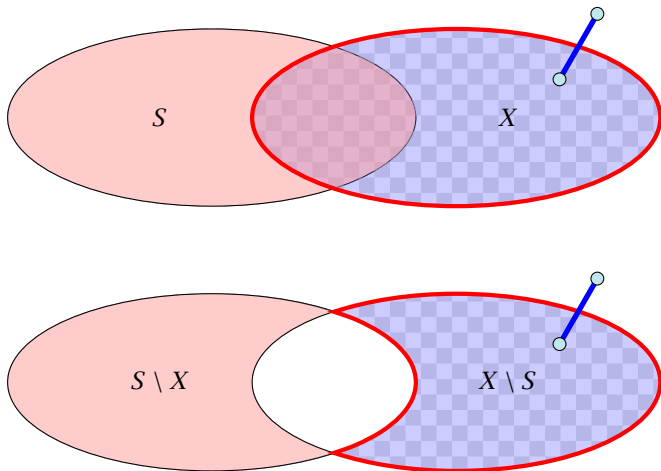
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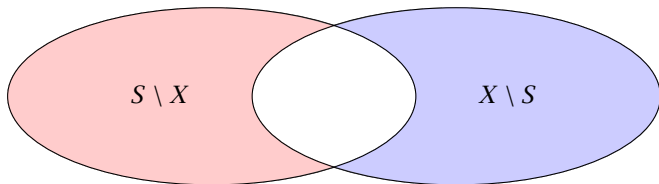
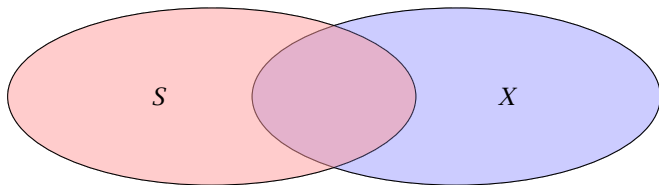
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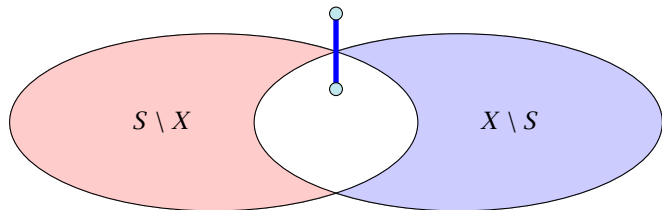
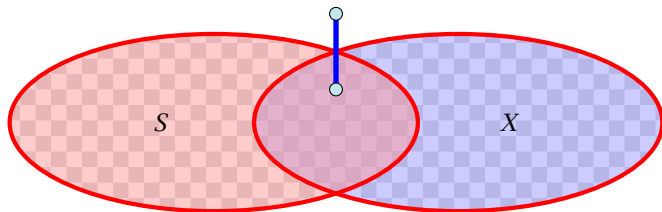
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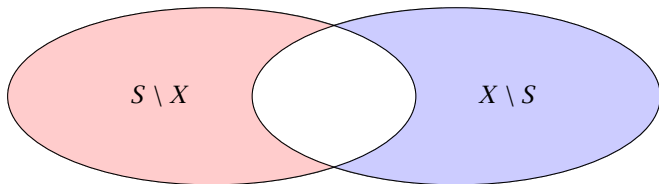
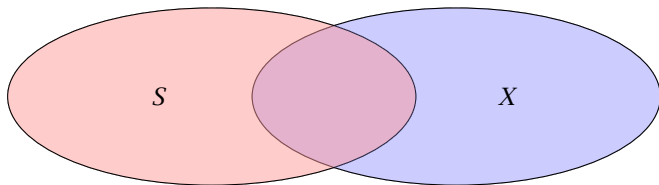
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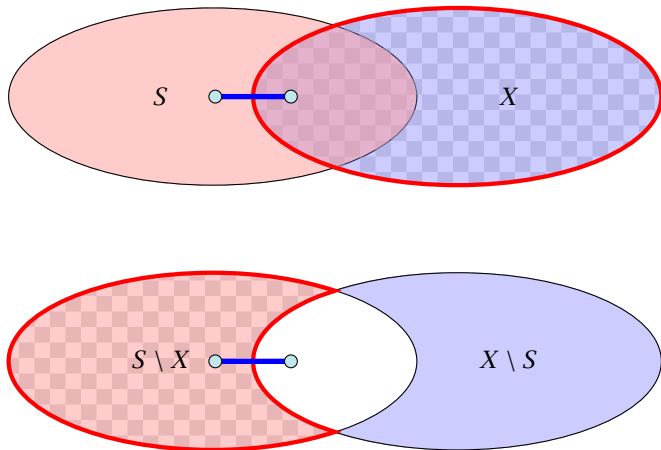
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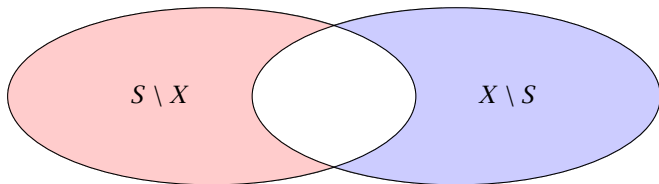
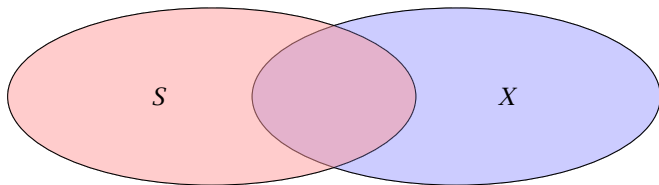
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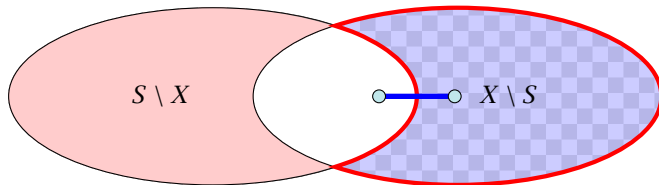
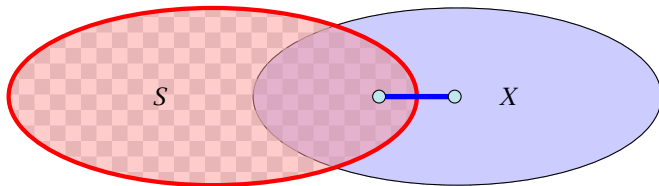
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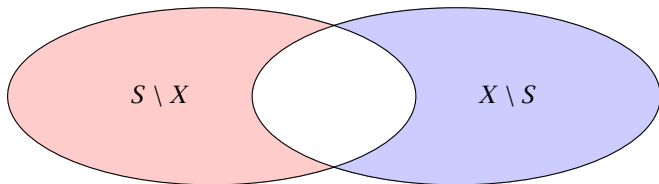
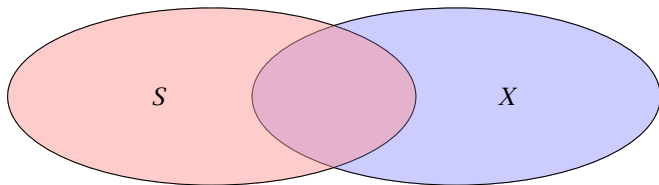
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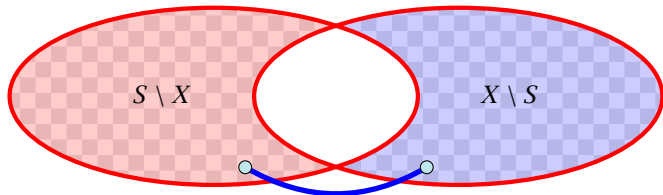
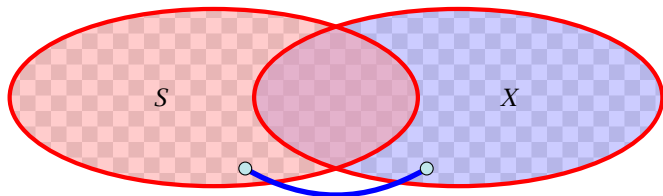
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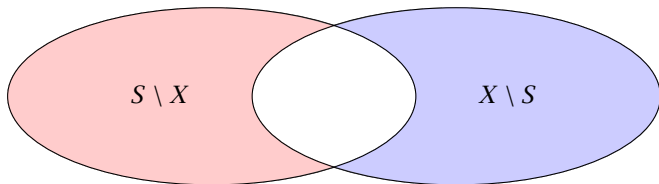
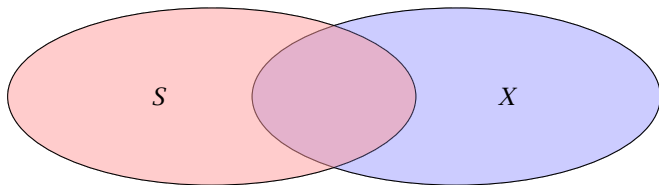
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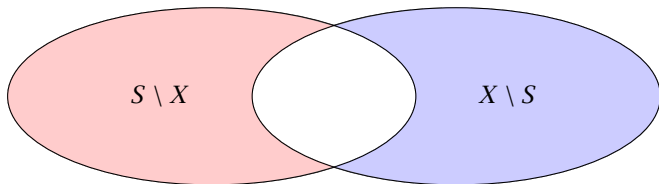
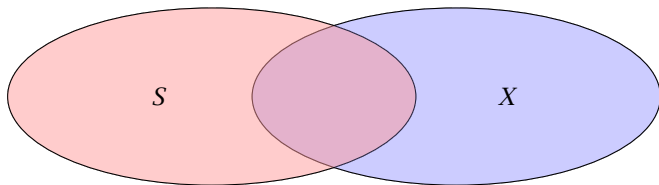
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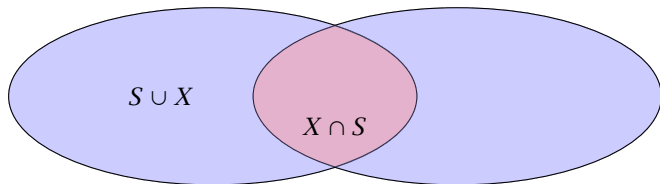
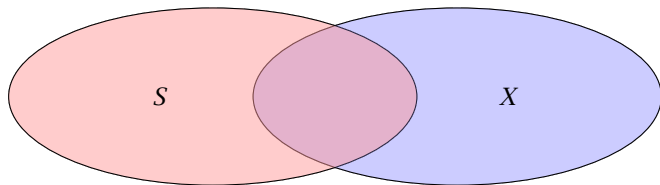
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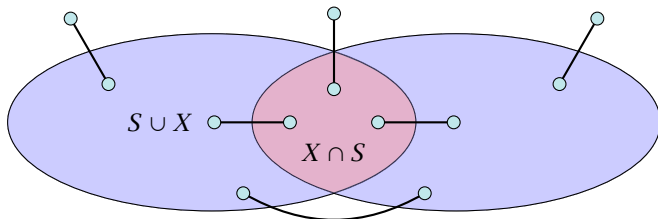
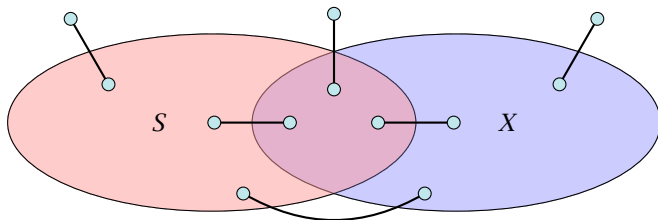
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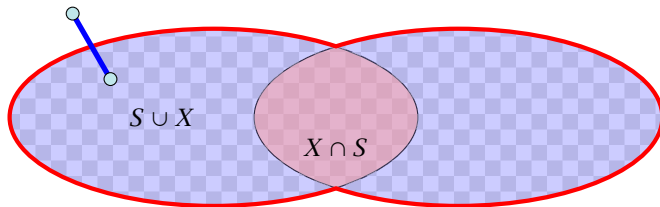
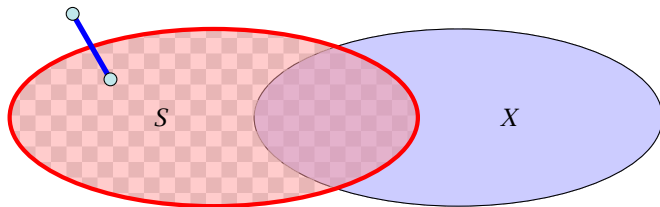
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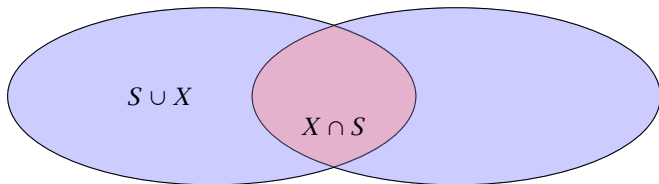
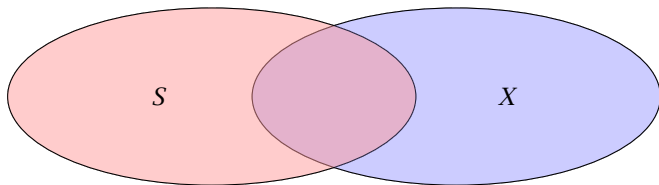
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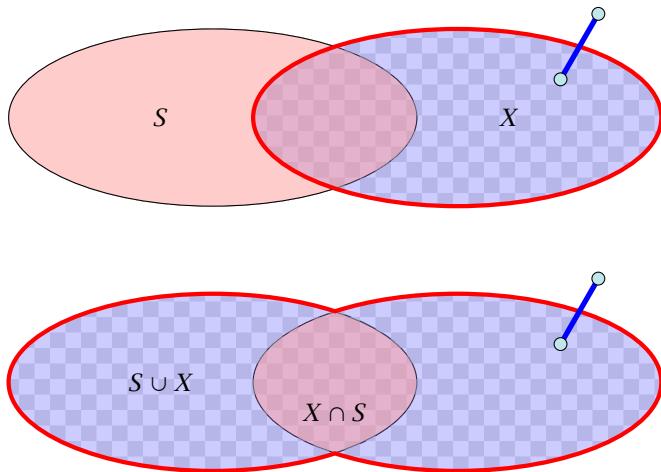
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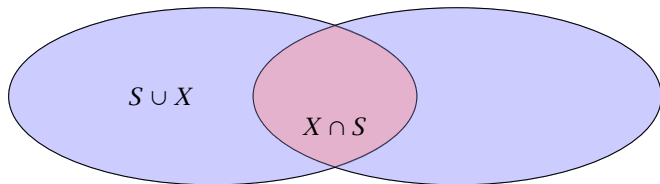
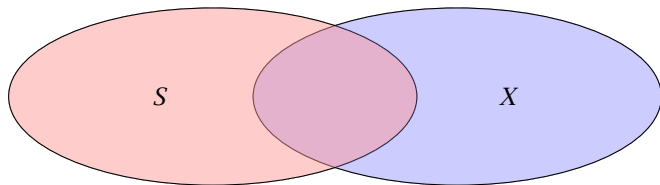
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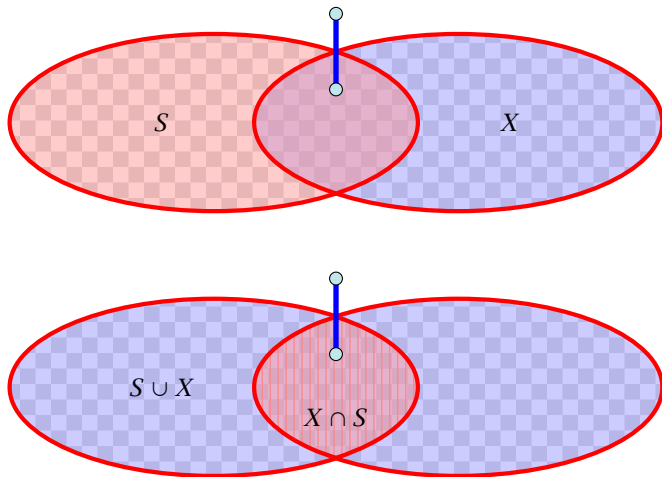
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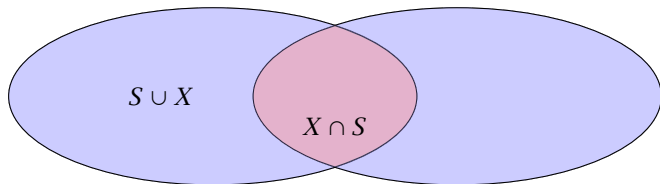
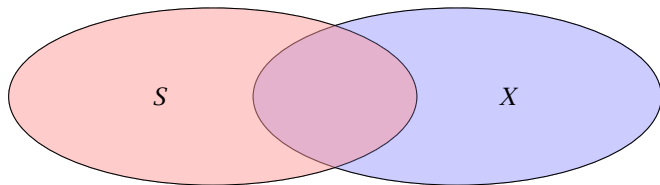
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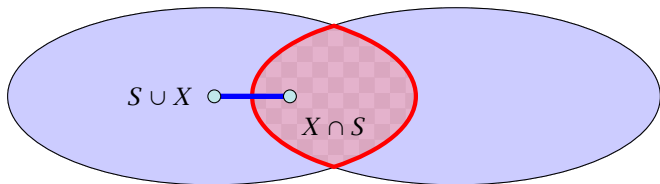
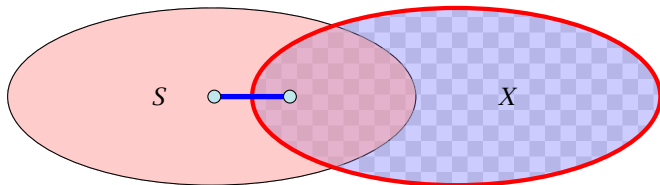
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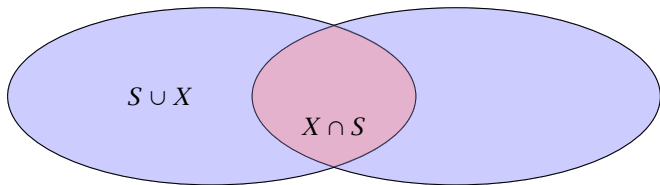
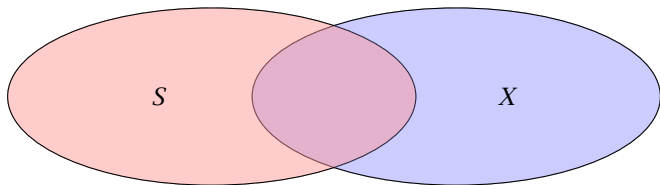
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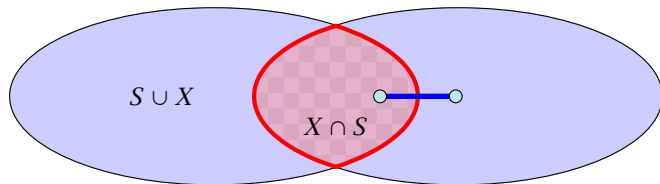
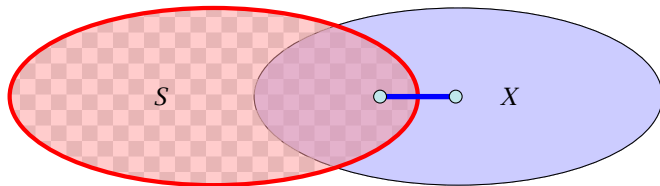
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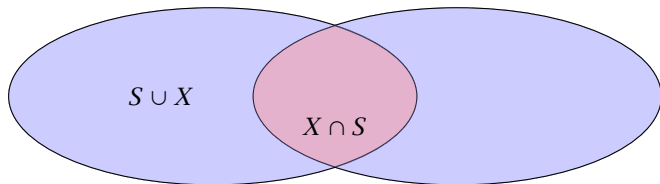
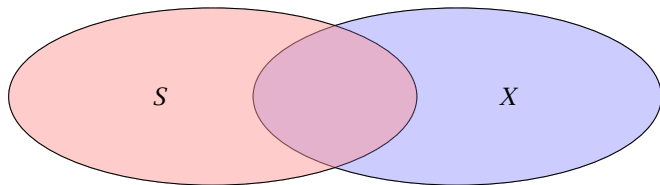
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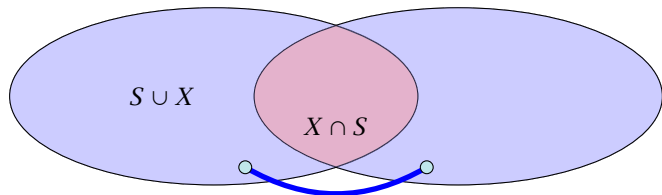
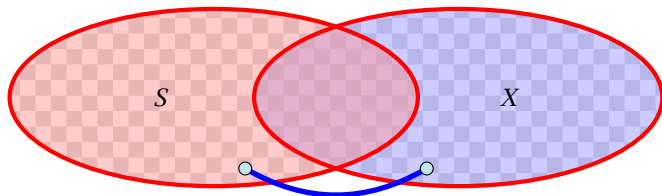
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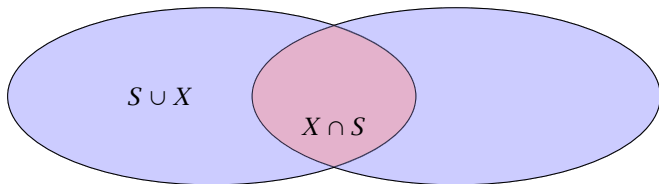
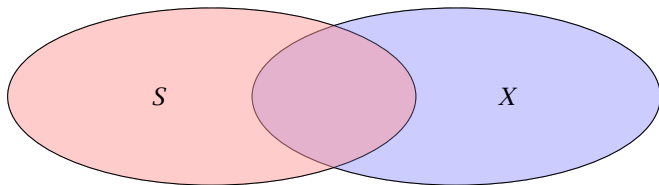
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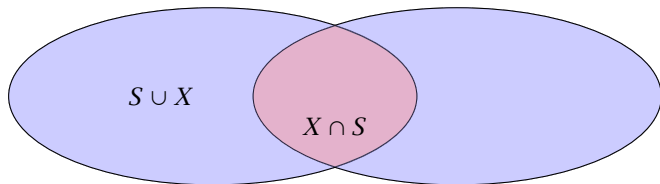
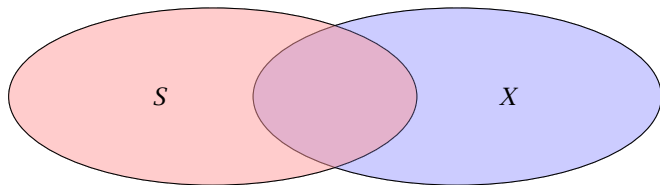
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Analysis

Lemma 94 tells us that if we have a graph $G = (V, E)$ and we contract a subset $X \subset V$ that corresponds to some mincut, then the value of $f(s, t)$ does not change for two nodes $s, t \notin X$.

We will show (later) that the connected components that we contract during a split-operation each correspond to some mincut and, hence, $f_H(s, t) = f(s, t)$, where $f_H(s, t)$ is the value of a minimum s - t mincut in graph H .

Invariant [existence of representatives]:

For any edge $\{S_i, S_j\}$ in T , there are vertices $a \in S_i$ and $b \in S_j$ such that $w(S_i, S_j) = f(a, b)$ and the cut defined by edge $\{S_i, S_j\}$ is a minimum a - b cut in G .

Analysis

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- ▶ Let $\{x_j, x_{j+1}\}$ be the edge with minimum weight on the path.
- ▶ Since by the invariant this edge induces an s - t cut with capacity $f(x_j, x_{j+1})$ we get $f(s, t) \leq f(x_j, x_{j+1}) = f_T(s, t)$.

Analysis

- ▶ Hence, $f_T(s, t) = f(s, t)$ (flow equivalence).
- ▶ The edge $\{x_j, x_{j+1}\}$ is a mincut between s and t in T .
- ▶ By invariant, it forms a cut with capacity $f(x_j, x_{j+1})$ in G (which separates s and t).
- ▶ Since, we can send a flow of value $f(x_j, x_{j+1})$ btw. s and t , this is an s - t mincut (cut property).

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Proof of Invariant

The invariant obviously holds at the beginning of the algorithm.

Now, we show that it holds after a split-operation provided that it was true before the operation.

Let S_i denote our selected cluster with nodes a and b . Because of the invariant all edges leaving $\{S_i\}$ in T correspond to some mincuts.

Therefore, contracting the connected components does not change the mincut btw. a and b due to Lemma 94.

After the split we have to choose representatives for all edges. For the new edge $\{S_i^a, S_i^b\}$ with capacity $w(S_i^a, S_i^b) = f_H(a, b)$ we can simply choose a and b as representatives.

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For edges that are not incident to S_i we do not need to change representatives as the neighbouring sets do not change.

Consider an edge $\{X, S_i\}$, and suppose that before the split it used representatives $x \in X$, and $s \in S_i$. Assume that this edge is replaced by $\{X, S_i^a\}$ in the new tree (the case when it is replaced by $\{X, S_i^b\}$ is analogous).

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Because the invariant was true before the split we know that the edge $\{X, S_i\}$ induces a cut in G of capacity $f(x, s)$. Since, x and a are on opposite sides of this cut, we know that $f(x, a) \leq f(x, s)$.

The set B forms a mincut separating a from b . Contracting all nodes in this set gives a new graph G' where the set B is represented by node v_B . Because of Lemma 94 we know that $f'(x, a) = f(x, a)$ as $x, a \notin B$.

We further have $f'(x, a) \geq \min\{f'(x, v_B), f'(v_B, a)\}$.

Since $s \in B$ we have $f'(v_B, x) \geq f(s, x)$.

Also, $f'(a, v_B) \geq f(a, b) \geq f(x, s)$ since the a - b cut that splits S_i into S_i^a and S_i^b also separates s and x .

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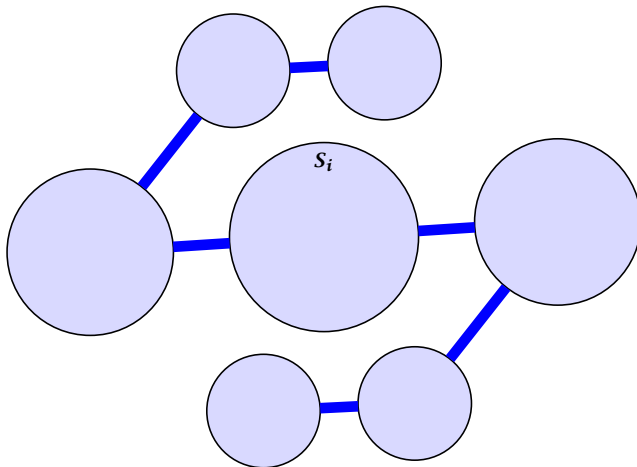
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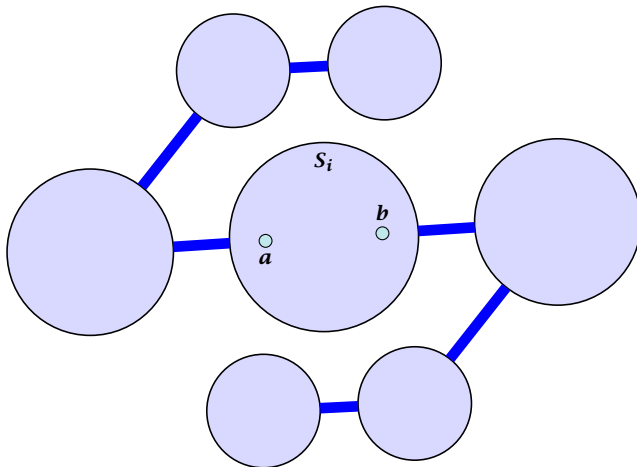
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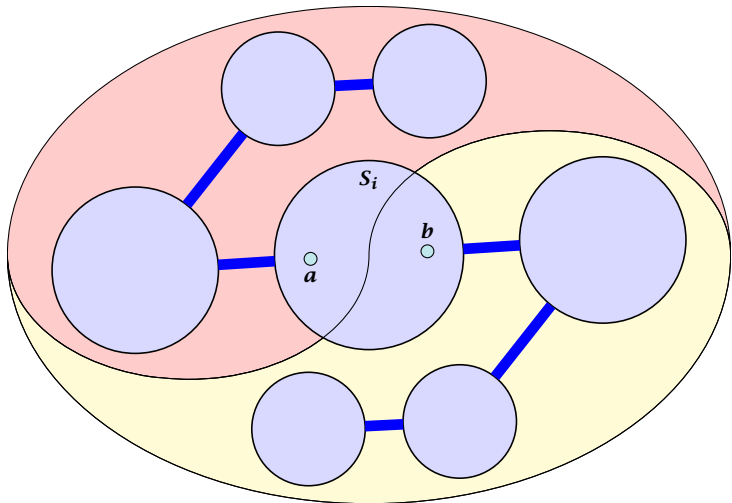
Analysis



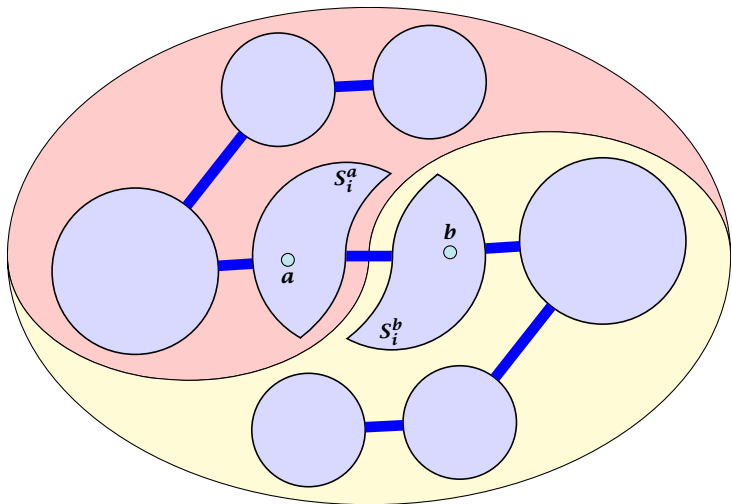
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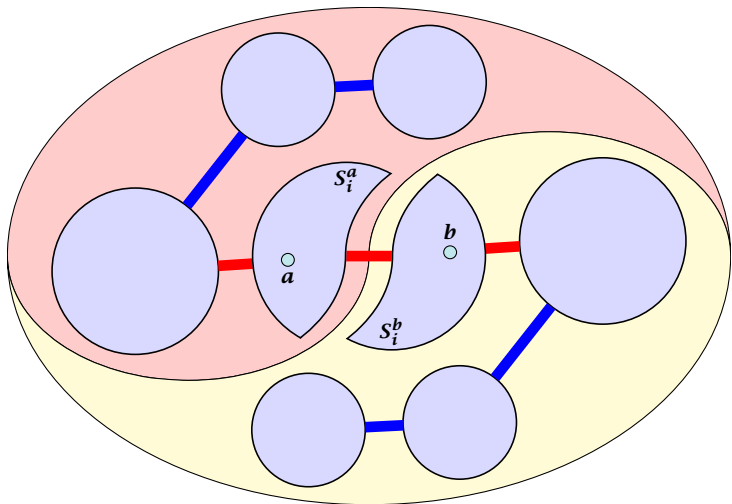
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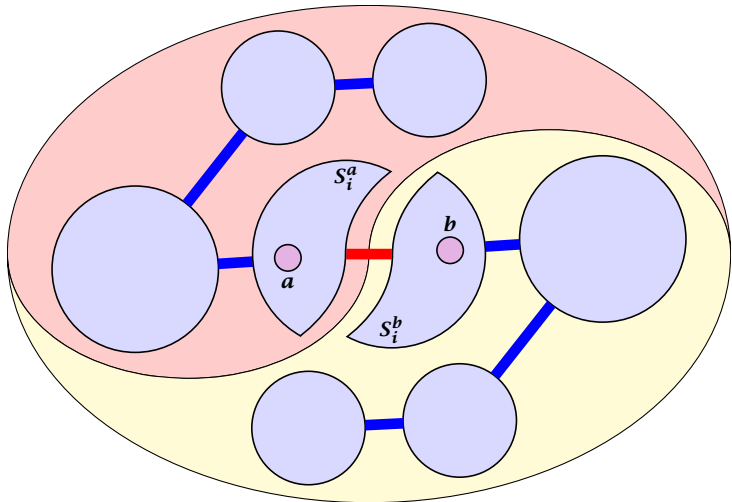
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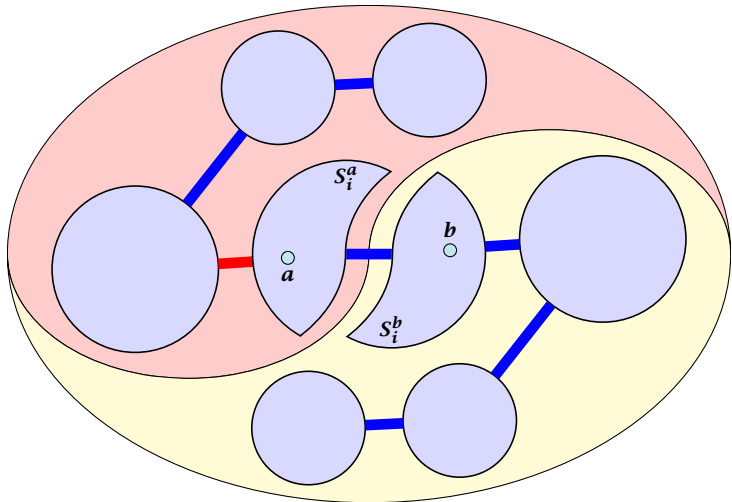
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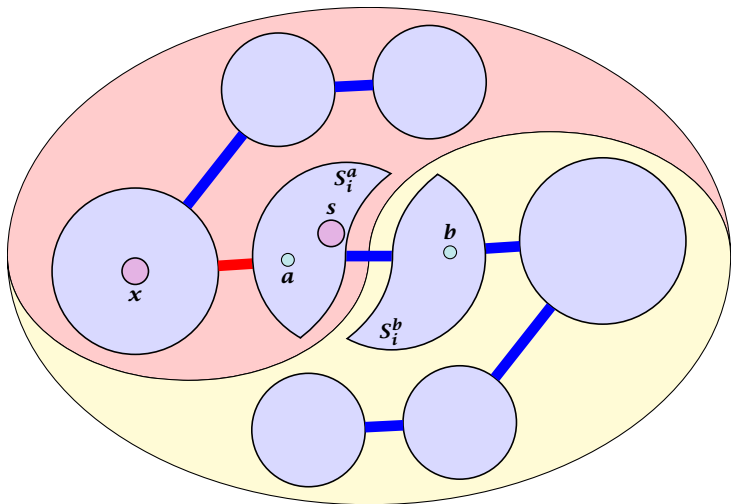
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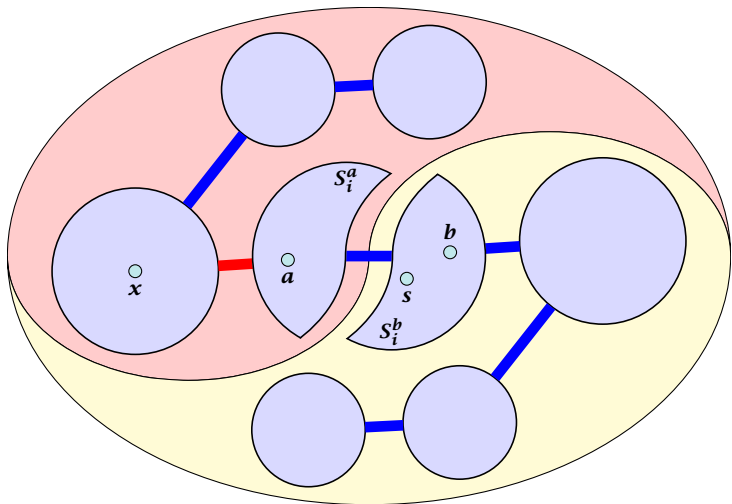
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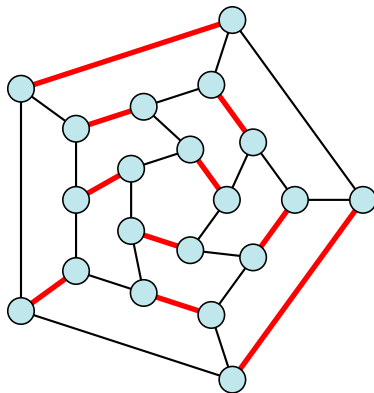


Part V

Matchings

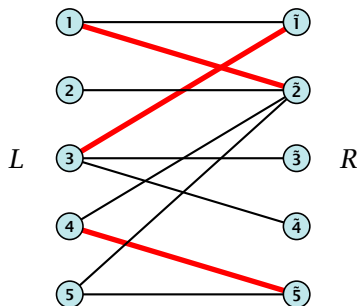
Matching

- ▶ Input: undirected graph $G = (V, E)$.
- ▶ $M \subseteq E$ is a **matching** if each node appears in at most one edge in M .
- ▶ Maximum Matching: find a matching of maximum cardinality



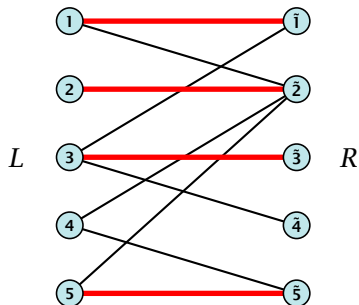
Bipartite Matching

- ▶ Input: undirected, **bipartite** graph $G = (L \uplus R, E)$.
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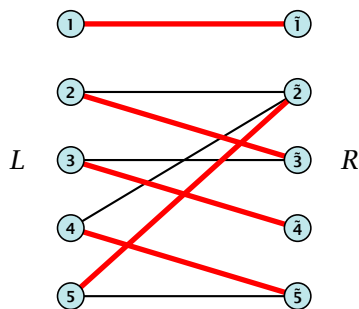
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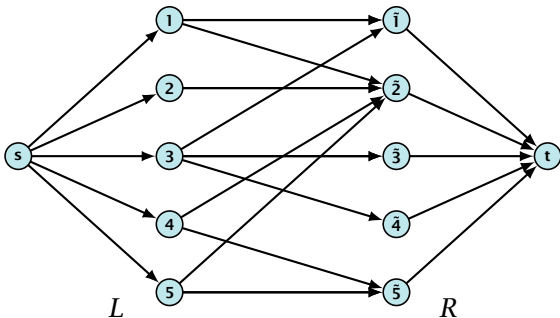
Bipartite Matching

- ▶ A matching M is **perfect** if it is of cardinality $|M| = |V|/2$.
- ▶ For a bipartite graph $G = (L \uplus R, E)$ this means $|M| = |L| = |R| = n$.



19 Bipartite Matching via Flows

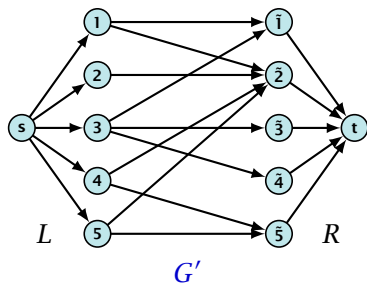
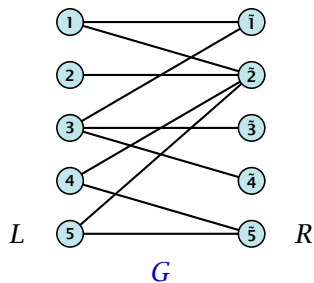
- ▶ Input: undirected, **bipartite** graph $G = (L \uplus R \uplus \{s, t\}, E')$.
- ▶ Direct all edges from L to R .
- ▶ Add source s and connect it to all nodes on the left.
- ▶ Add t and connect all nodes on the right to t .
- ▶ All edges have unit capacity.



Proof

Max cardinality matching in $G \leq$ value of maxflow in G'

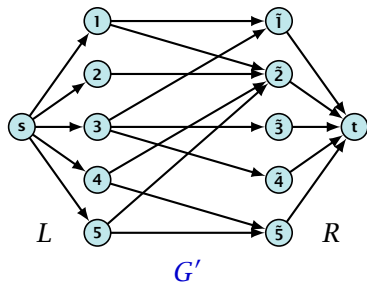
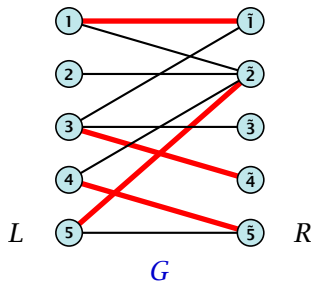
- ▶ Given a maximum matching M of cardinality k .
- ▶ Consider flow f that sends one unit along each of k paths.
- ▶ f is a flow and has cardinality k .



Proof

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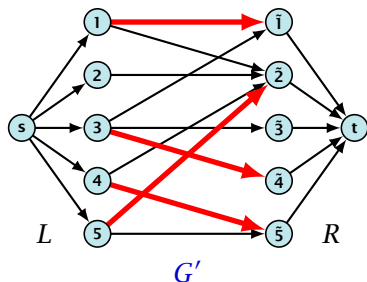
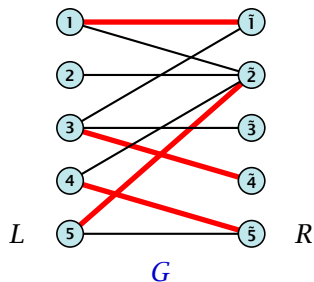
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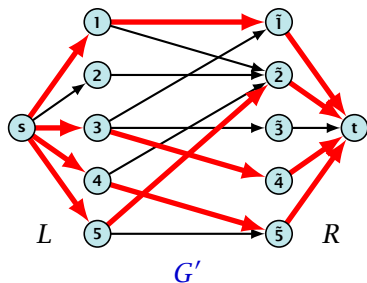
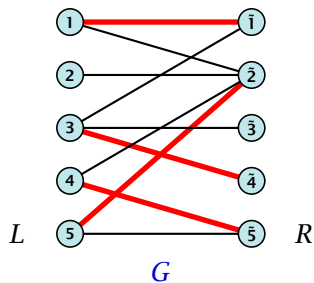
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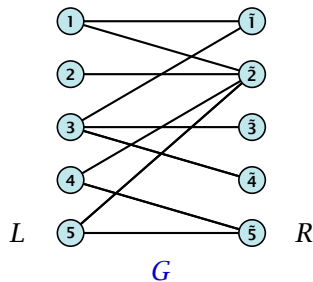
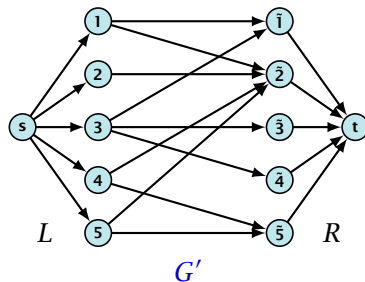
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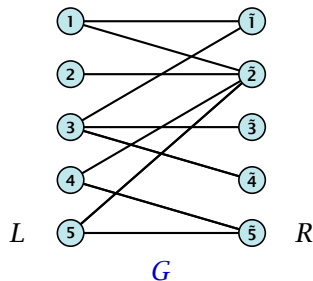
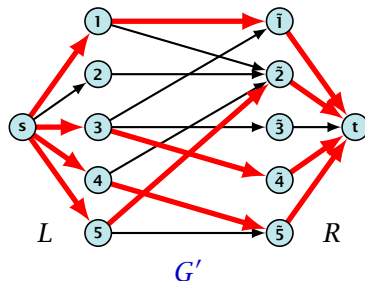
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- ▶ Each node in L and R participates in at most one edge in M .
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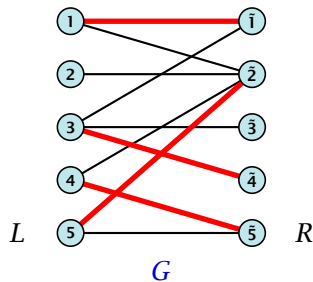
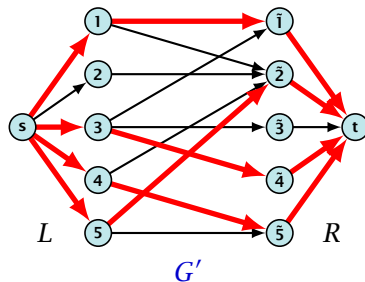
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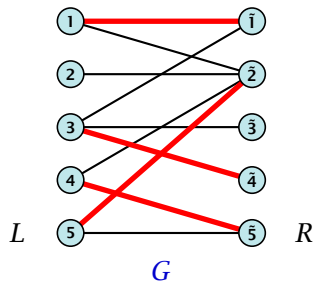
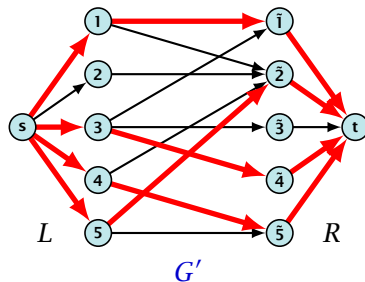
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19 Bipartite Matching via Flows

Which flow algorithm to use?

- ▶ Generic augmenting path: $\mathcal{O}(m \text{val}(f^*)) = \mathcal{O}(mn)$.
- ▶ Capacity scaling: $\mathcal{O}(m^2 \log C) = \mathcal{O}(m^2)$.

20 Augmenting Paths for Matchings

Definitions.

- ▶ Given a matching M in a graph G , a vertex that is not incident to any edge of M is called a **free vertex** w. r. .t. M .
- ▶ For a matching M a path P in G is called an **alternating path** if edges in M alternate with edges not in M .
- ▶ An alternating path is called an **augmenting path** for matching M if it ends at distinct free vertices.

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A matching M is a maximum matching if and only if there is no augmenting path w. r. t. M .

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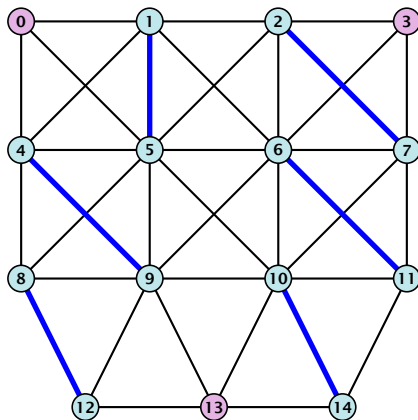
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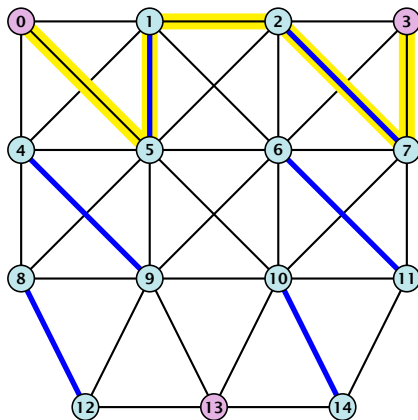
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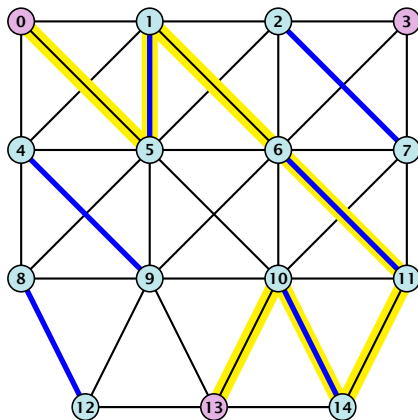
Augmenting Paths in Action



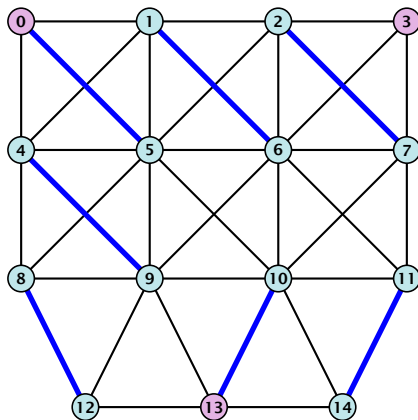
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20 Augmenting Paths for Matchings

Proof.

- ⇒ If M is maximum there is no augmenting path P , because we could switch matching and non-matching edges along P . This gives matching $M' = M \oplus P$ with larger cardinality.
- ⇐ Suppose there is a matching M' with larger cardinality. Consider the graph H with edge-set $M' \oplus M$ (i.e., only edges that are in either M or M' but not in both).

Each vertex can be incident to at most two edges (one from M and one from M'). Hence, the connected components are alternating cycles or alternating path.

As $|M'| > |M|$ there is one connected component that is a path P for which both endpoints are incident to edges from M' . P is an augmenting path.

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Algorithmic idea:

As long as you find an augmenting path augment your matching using this path. When you arrive at a matching for which no augmenting path exists you have a maximum matching.

Theorem 96

Let G be a graph, M a matching in G , and let u be a free vertex w.r.t. M . Further let P denote an augmenting path w.r.t. M and let $M' = M \oplus P$ denote the matching resulting from augmenting M with P . If there was no augmenting path starting at u in M then there is no augmenting path starting at u in M' .

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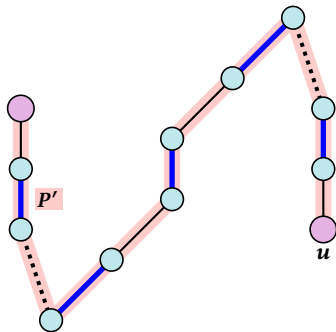
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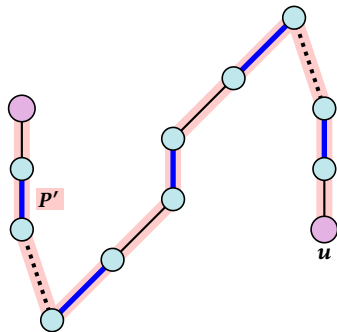
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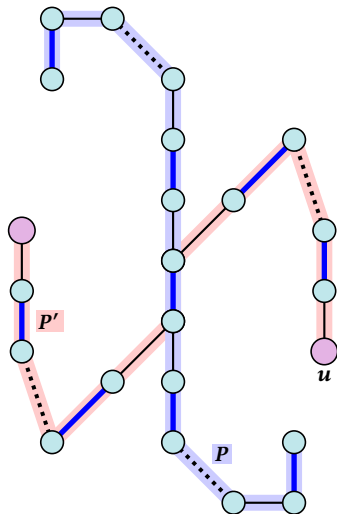
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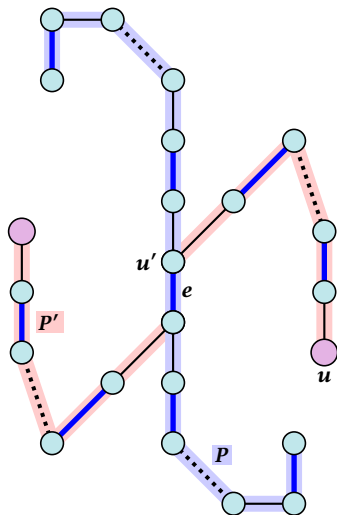
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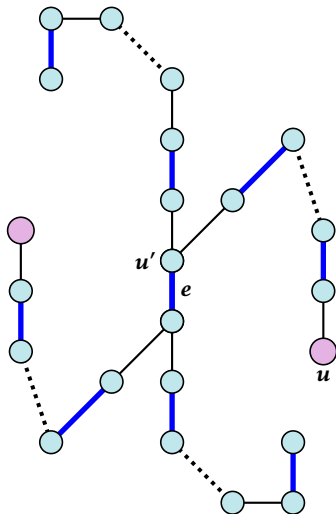
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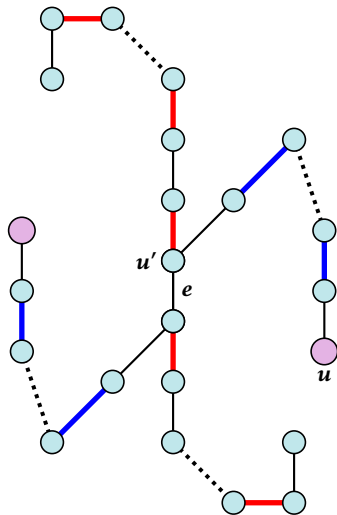
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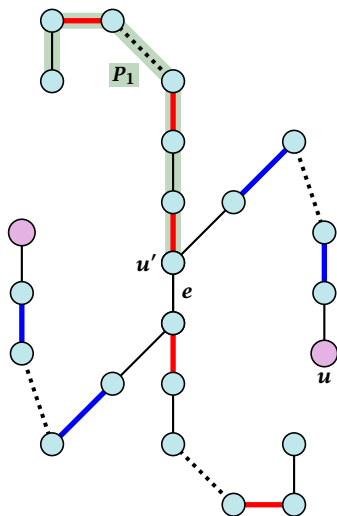
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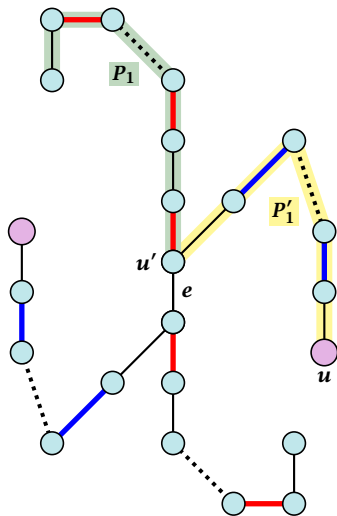
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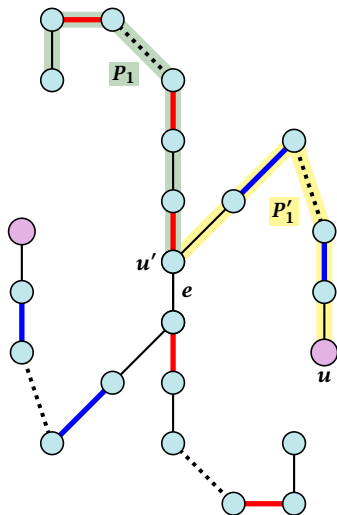
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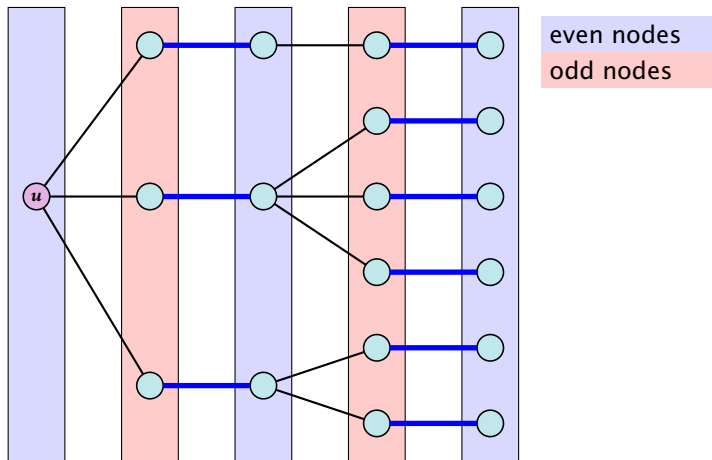
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- ▶ $P_1 \circ P'_1$ is augmenting path in M (\neq).



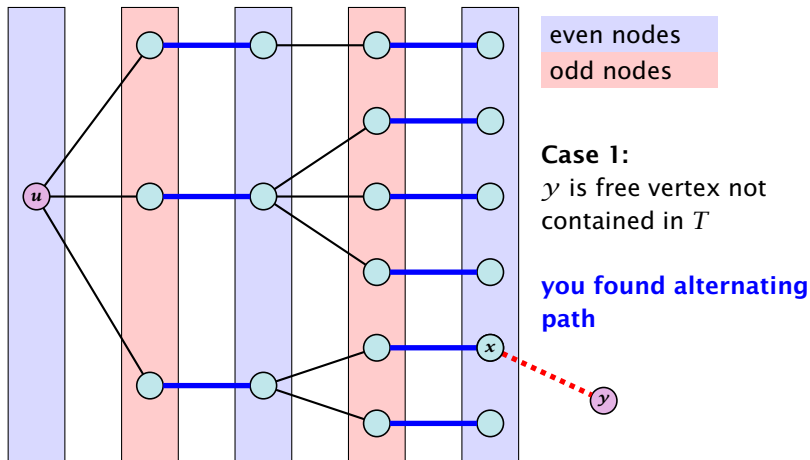
How to find an augmenting path?

Construct an alternating tree.



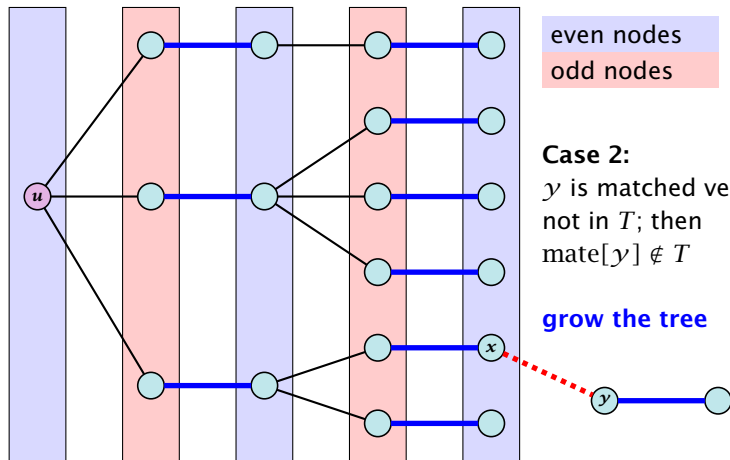
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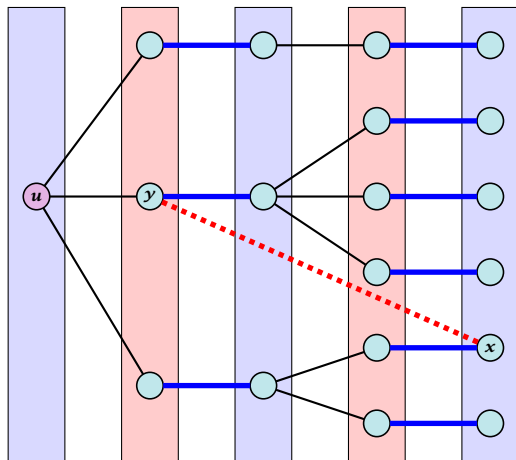
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even nodes

odd nodes

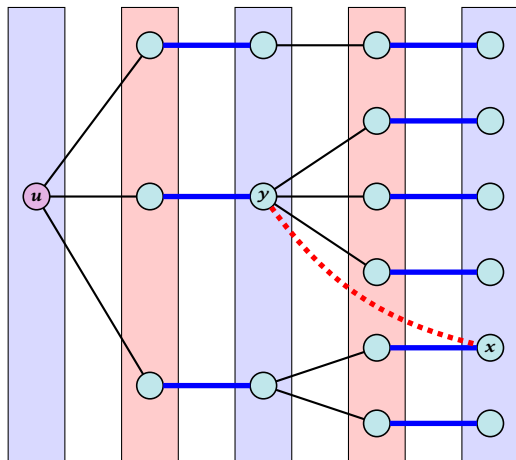
Case 3:

y is already contained
in T as an odd vertex

ignore successor y

How to find an augmenting path?

Construct an alternating tree.



even nodes

odd nodes

Case 4:

y is already contained
in T as an even vertex

can't ignore y

does not happen in
bipartite graphs

Algorithm 1 BiMatch($G, match$)

```
1: for  $x \in V$  do  $mate[x] \leftarrow 0$ ;  
2:  $r \leftarrow 0$ ;  $free \leftarrow n$ ;  
3: while  $free \geq 1$  and  $r < n$  do  
4:    $r \leftarrow r + 1$   
5:   if  $mate[r] = 0$  then  
6:     for  $i = 1$  to  $m$  do  $parent[i'] \leftarrow 0$   
7:      $Q \leftarrow \emptyset$ ;  $Q.append(r)$ ;  $aug \leftarrow false$ ;  
8:     while  $aug = false$  and  $Q \neq \emptyset$  do  
9:        $x \leftarrow Q.dequeue()$ ;  
10:      if  $\exists y \in A_x: mate[y] = 0$  then  
11:         $augment(mate, parent, y)$ ;  
12:         $aug \leftarrow true$ ;  $free \leftarrow free - 1$ ;  
13:      else  
14:        if  $parent[y] = 0$  then  
15:           $parent[y] \leftarrow x$ ;  
16:           $Q.enqueue(y)$ ;
```

graph $G = (S \cup S', E)$;

$S = \{1, \dots, n\}$;

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initial matching empty

$free$: number of
unmatched nodes in S

r : root of current tree

if r is unmatched
start tree construction

initialize empty tree

no augmen. path but
unexamined leaves

free neighbour found

add new node y to Q

21 Weighted Bipartite Matching

Weighted Bipartite Matching/Assignment

- ▶ Input: undirected, bipartite graph $G = L \cup R, E$.
- ▶ an edge $e = (\ell, r)$ has weight $w_e \geq 0$
- ▶ find a matching of maximum weight, where the weight of a matching is the sum of the weights of its edges

Simplifying Assumptions (wlog [why?]):

- ▶ assume that $|L| = |R| = n$
- ▶ assume that there is an edge between every pair of nodes $(\ell, r) \in V \times V$

Weighted Bipartite Matching

Theorem 97 (Halls Theorem)

A bipartite graph $G = (L \cup R, E)$ has a perfect matching if and only if for all sets $S \subseteq L$, $|\Gamma(S)| \geq |S|$, where $\Gamma(S)$ denotes the set of nodes in R that have a neighbour in S .

Halls Theorem

Proof:

- ← Of course, the condition is necessary as otherwise not all nodes in S could be matched to different neighbours.
- ⇒ For the other direction we need to argue that the minimum cut in the graph G' is at least $|L|$.

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 - ▶ Let S denote a minimum cut and let $L_S \cong L \cap S$ and $R_S \cong R \cap S$ denote the portion of S inside L and R , respectively.
 - ▶ Clearly, all neighbours of nodes in L_S have to be in S , as otherwise we would cut an edge of infinite capacity.
 - ▶ This gives $R_S \geq |\Gamma(L_S)|$.
 - ▶ The size of the cut is $|L| - |L_S| + |R_S|$.
 - ▶ Using the fact that $|\Gamma(L_S)| \geq |L_S|$ gives that this is at least $|L|$.

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- ▶ Let $H(\vec{x})$ denote the subgraph of G that only contains edges that are **tight** w.r.t. the node weighting \vec{x} , i.e. edges $e = (u, v)$ for which $w_e = x_u + x_v$.
- ▶ Try to compute a perfect matching in the subgraph $H(\vec{x})$. If you are successful you found an optimal matching.

Algorithm Outline

Reason:

- ▶ The weight of your matching M^* is

$$\sum_{(u,v) \in M^*} w_{(u,v)} = \sum_{(u,v) \in M^*} (x_u + x_v) = \sum_v x_v .$$

- ▶ Any other matching M has

$$\sum_{(u,v) \in M} w_{(u,v)} \leq \sum_{(u,v) \in M} (x_u + x_v) \leq \sum_v x_v .$$

Algorithm Outline

What if you don't find a perfect matching?

Then, Hall's theorem guarantees you that there is a set $S \subseteq L$, with $|\Gamma(S)| < |S|$, where Γ denotes the neighbourhood w.r.t. the subgraph $H(\vec{x})$.

Idea: reweight such that:

- ▶ the total weight assigned to nodes decreases
- ▶ the weight function still dominates the edge-weights

If we can do this we have an algorithm that terminates with an optimal solution (we analyze the running time later).

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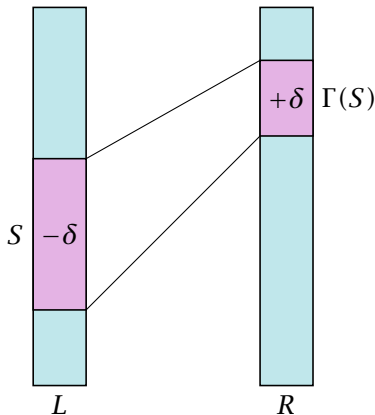
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Increase node-weights in $\Gamma(S)$ by $+\delta$, and decrease the node-weights in S by $-\delta$.

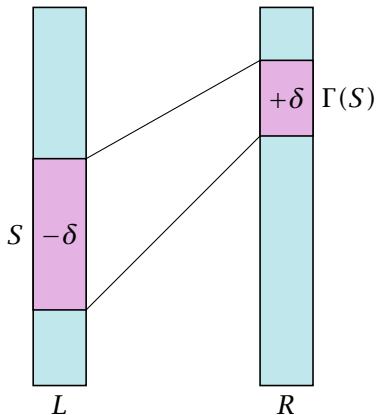
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- ▶ Only edges from S to $R - \Gamma(S)$ decrease in their weight.
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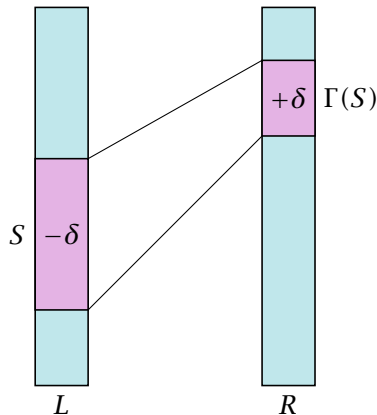
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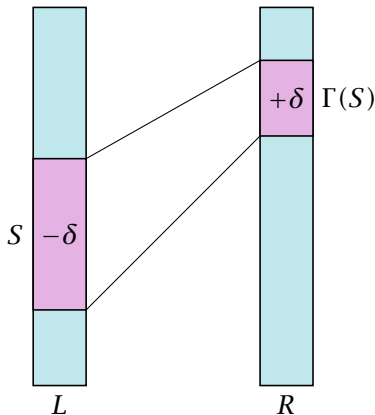
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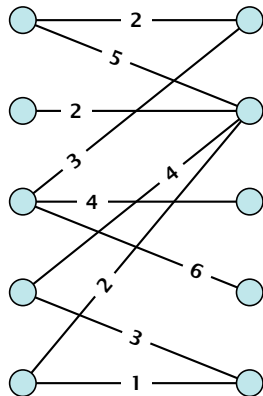
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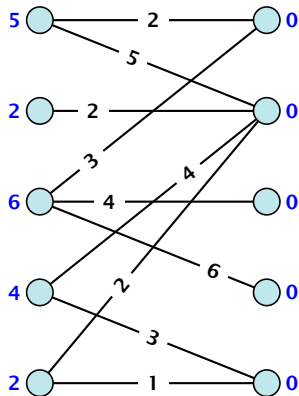
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Edges not drawn have weight 0.



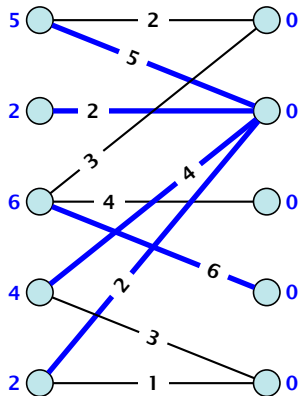
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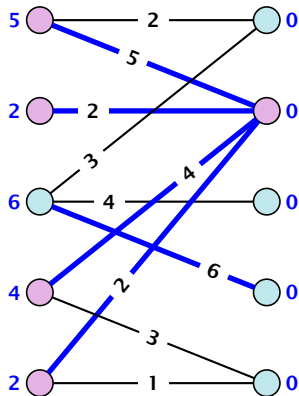
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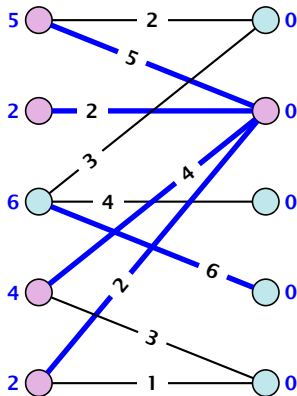
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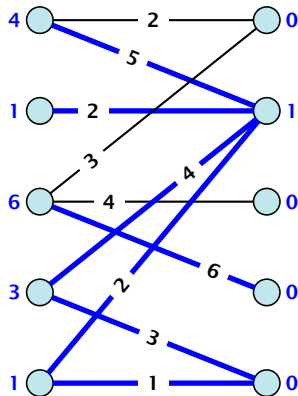
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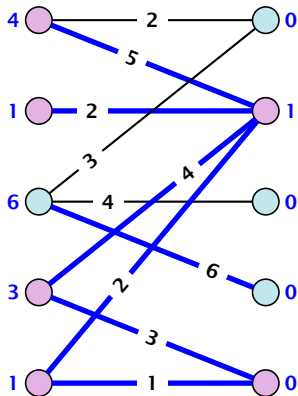
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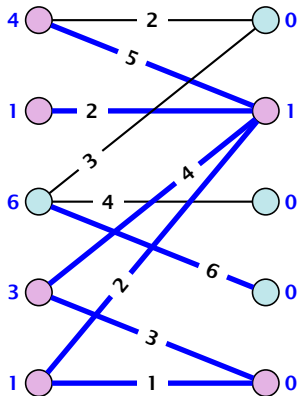
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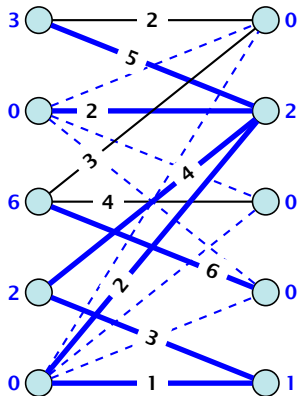
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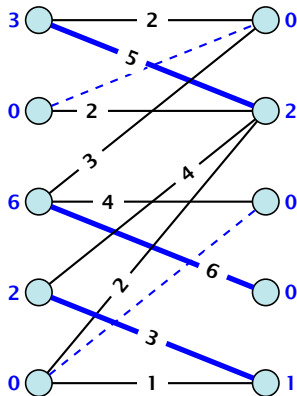
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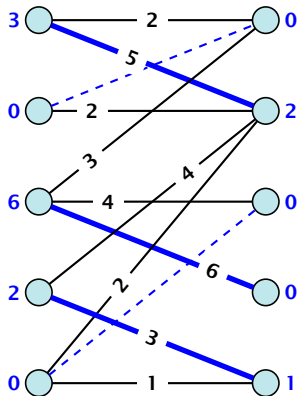
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How many iterations do we need?

- ▶ One reweighting step increases the number of edges out of S by at least one.
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Analysis

- ▶ We will show that after at most n reweighting steps the size of the maximum matching can be increased by finding an augmenting path.
- ▶ This gives a polynomial running time.

Analysis

How do we find S ?

- ▶ Start on the left and compute an alternating tree, starting at any free node u .
- ▶ If this construction stops, there is no perfect matching in the tight subgraph (because for a perfect matching we need to find an augmenting path starting at u).
- ▶ The set of even vertices is on the left and the set of odd vertices is on the right **and** contains all neighbours of even nodes.
- ▶ All odd vertices are matched to even vertices. Furthermore, the even vertices additionally contain the free vertex u . Hence, $|V_{\text{odd}}| = |\Gamma(V_{\text{even}})| < |V_{\text{even}}|$, and all odd vertices are saturated in the current matching.

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- ▶ The current matching does not have any edges from V_{odd} to outside of $L \setminus V_{\text{even}}$ (edges that may possibly be deleted by changing weights).
- ▶ After changing weights, there is at least one more edge connecting V_{even} to a node outside of V_{odd} . After at most n reweightings we can do an augmentation.
- ▶ A reweighting can be trivially performed in time $\mathcal{O}(n^2)$ (keeping track of the tight edges).
- ▶ An augmentation takes at most $\mathcal{O}(n)$ time.
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A Fast Matching Algorithm

Algorithm 54 Bimatch-Hopcroft-Karp(G)

```
1:  $M \leftarrow \emptyset$ 
2: repeat
3:   let  $\mathcal{P} = \{P_1, \dots, P_k\}$  be maximal set of
4:   vertex-disjoint, shortest augmenting path w.r.t.  $M$ .
5:    $M \leftarrow M \oplus (P_1 \cup \dots \cup P_k)$ 
6: until  $\mathcal{P} = \emptyset$ 
7: return  $M$ 
```

We call one iteration of the repeat-loop a **phase** of the algorithm.

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Given a matching M and a maximal matching M^* there exist $|M^*| - |M|$ *vertex-disjoint* augmenting path w.r.t. M .

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- ▶ The graph contains $k \triangleq |M^*| - |M|$ more red edges than blue edges.
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- ▶ Otherwise, at least one edge from P coincides with an edge from paths $\{P_1, \dots, P_k\}$.
- ▶ This edge is not contained in A .
- ▶ Hence, $|A| \leq k\ell + |P| - 1$.
- ▶ The lower bound on $|A|$ gives $(k + 1)\ell \leq |A| \leq k\ell + |P| - 1$, and hence $|P| \geq \ell + 1$.

Analysis

Lemma 100

P is of length at least $\ell + 1$. This shows that the length of a shortest augmenting path increases between two phases of the Hopcroft-Karp algorithm.

Proof.

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Analysis

If the shortest augmenting path w.r.t. a matching M has ℓ edges then the cardinality of the maximum matching is of size at most $|M| + \lfloor \frac{|V|}{\ell+1} \rfloor$.

Proof.

The symmetric difference between M and M^* contains $|M^*| - |M|$ vertex-disjoint augmenting paths. Each of these paths contains at least $\ell + 1$ vertices. Hence, there can be at most $\lfloor \frac{|V|}{\ell+1} \rfloor$ of them.

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Lemma 101

The Hopcroft-Karp algorithm requires at most $2\sqrt{|V|}$ phases.

Proof.

- ▶ After iteration $\lfloor \sqrt{|V|} \rfloor$ the length of a shortest augmenting path must be at least $\lfloor \sqrt{|V|} \rfloor + 1 \geq \sqrt{|V|}$.
- ▶ Hence, there can be at most $|V| / (\sqrt{|V|} + 1) \leq \sqrt{|V|}$ additional augmentations.

Analysis

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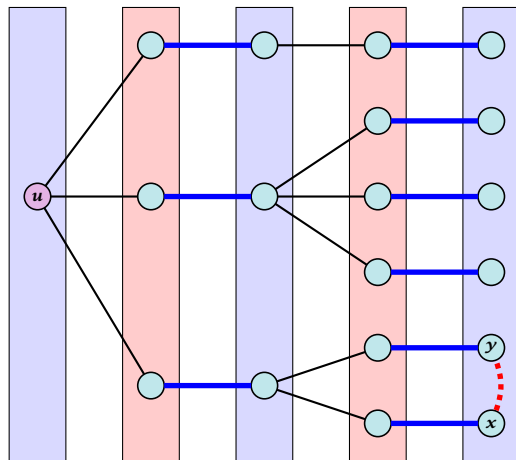
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- ▶ Hence, there can be at most $|V| / (\sqrt{|V|} + 1) \leq \sqrt{|V|}$ additional augmentations.

Lemma 102

One phase of the Hopcroft-Karp algorithm can be implemented in time $\mathcal{O}(m)$.

How to find an augmenting path?

Construct an alternating tree.



even nodes

odd nodes

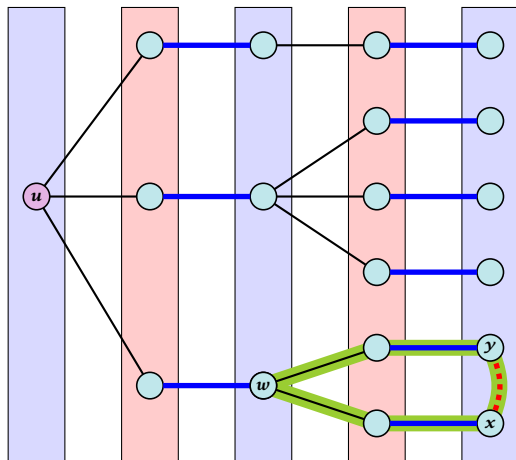
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y is already contained
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can't ignore y

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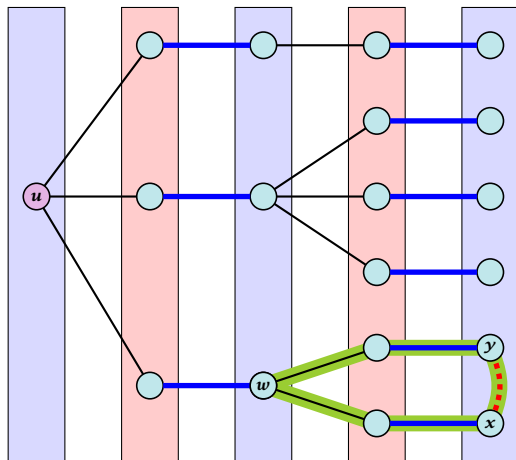
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The cycle $w \leftrightarrow y - x \leftrightarrow w$ is
called a **blossom**.

w is called the **base** of the
blossom (even node!!!).

The path $u-w$ path is called
the **stem** of the blossom.

Flowers and Blossoms

Definition 103

A **flower** in a graph $G = (V, E)$ w.r.t. a matching M and a (free) root node r , is a subgraph with two components:

- ▶ A stem is an even length alternating path that starts at the root node r and terminates at some node w . We permit the possibility that $r = w$ (empty stem).
- ▶ A blossom is an odd length alternating cycle that starts and terminates at the terminal node w of a stem and has no other node in common with the stem. w is called the base of the blossom.

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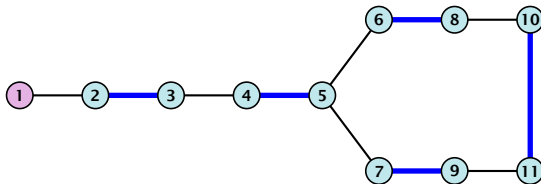
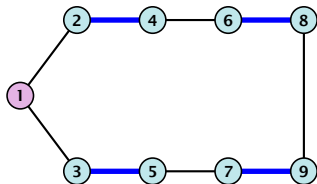
Flowers and Blossoms

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Flowers and Blossoms



Flowers and Blossoms

Properties:

1. A stem spans $2\ell + 1$ nodes and contains ℓ matched edges for some integer $\ell \geq 0$.
2. A blossom spans $2k + 1$ nodes and contains k matched edges for some integer $k \geq 1$. The matched edges match all nodes of the blossom except the base.
3. The base of a blossom is an even node (if the stem is part of an alternating tree starting at r).

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Flowers and Blossoms

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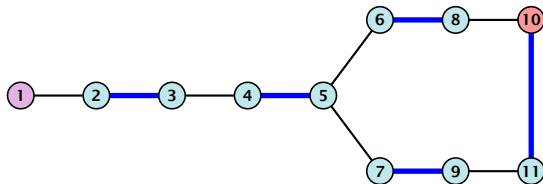
4. Every node x in the blossom (except its base) is reachable from the root (or from the base of the blossom) through two distinct alternating paths; one with even and one with odd length.
5. The even alternating path to x terminates with a matched edge and the odd path with an unmatched edge.

Flowers and Blossoms

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Flowers and Blossoms



When during the alternating tree construction we discover a blossom B we replace the graph G by $G' = G/B$, which is obtained from G by contracting the blossom B .

- ▶ Delete all vertices in B (and its incident edges) from G .
- ▶ Add a new (pseudo-)vertex b . The new vertex b is connected to all vertices in $V \setminus B$ that had at least one edge to a vertex from B .

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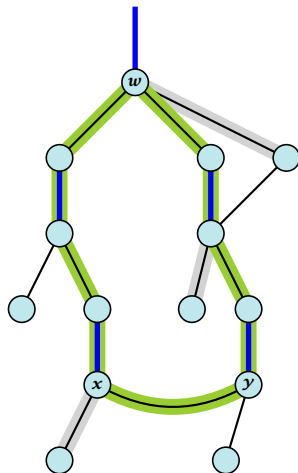
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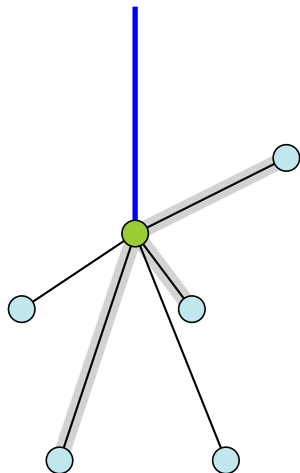
Shrinking Blossoms

- ▶ Edges of T that connect a node u not in B to a node in B become tree edges in T' connecting u to b .
- ▶ Matching edges (there is at most one) that connect a node u not in B to a node in B become matching edges in M' .
- ▶ Nodes that are connected in G to at least one node in B become connected to b in G' .



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Algorithm 55 search(r , $found$)

- 1: set $\bar{A}(i) \leftarrow A(i)$ for all nodes i
- 2: $found \leftarrow \text{false}$
- 3: unlabel all nodes;
- 4: give an even label to r and initialize $list \leftarrow \{r\}$
- 5: **while** $list \neq \emptyset$ **do**
- 6: delete a node i from $list$
- 7: examine(i , $found$)
- 8: **if** $found = \text{true}$ **then**
- 9: **return**

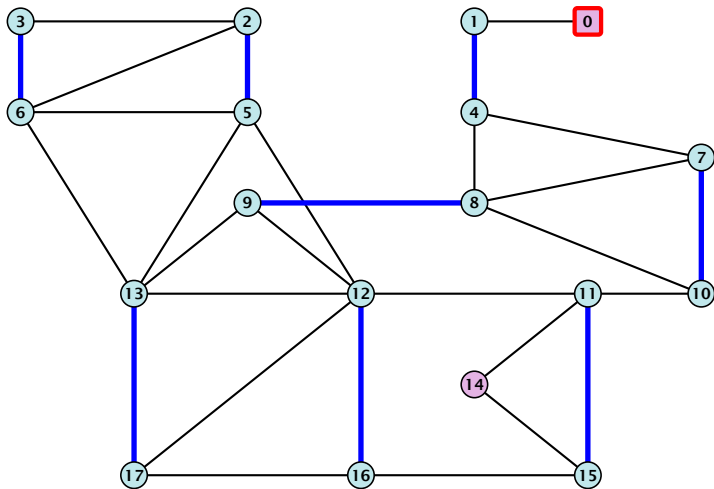
Algorithm 56 $\text{examine}(i, \text{found})$

```
1: for all  $j \in \bar{A}(i)$  do  
2:   if  $j$  is even then  $\text{contract}(i, j)$  and return  
3:   if  $j$  is unmatched then  
4:      $q \leftarrow j$ ;  
5:      $\text{pred}(q) \leftarrow i$ ;  
6:      $\text{found} \leftarrow \text{true}$ ;  
7:     return  
8:   if  $j$  is matched and unlabeled then  
9:      $\text{pred}(j) \leftarrow i$ ;  
10:     $\text{pred}(\text{mate}(j)) \leftarrow j$ ;
```

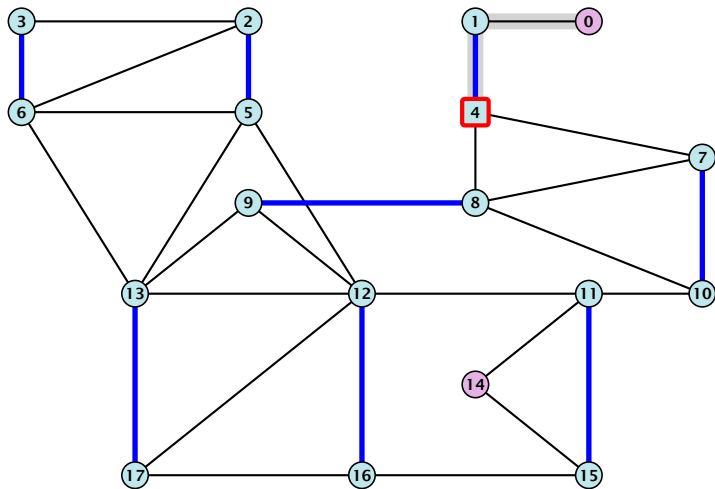
Algorithm 57 $\text{contract}(i, j)$

- 1: trace pred-indices of i and j to identify a blossom B
- 2: create new node b and set $\bar{A}(b) \leftarrow \cup_{x \in B} \bar{A}(x)$
- 3: label b even and add to *list*
- 4: update $\bar{A}(j) \leftarrow \bar{A}(j) \cup \{b\}$ for each $j \in \bar{A}(b)$
- 5: form a circular doubly linked list of nodes in B
- 6: delete nodes in B from the graph

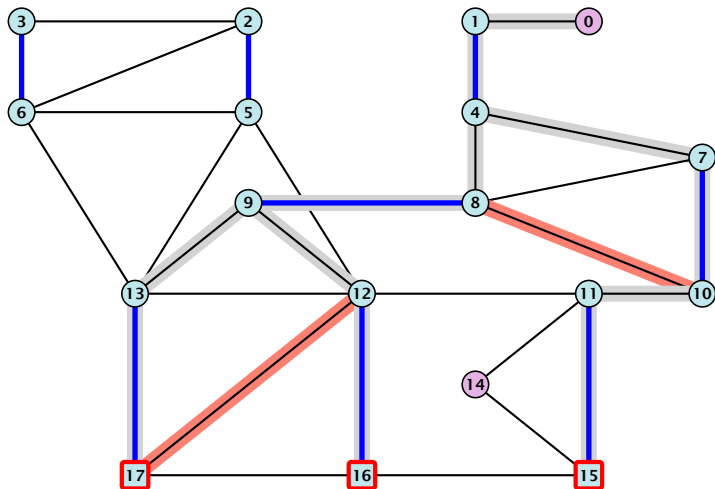
Example: Blossom Algorithm



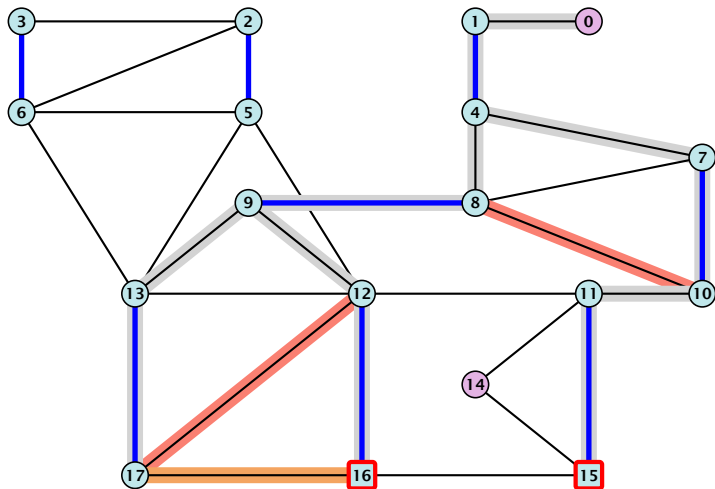
Example: Blossom Algorithm



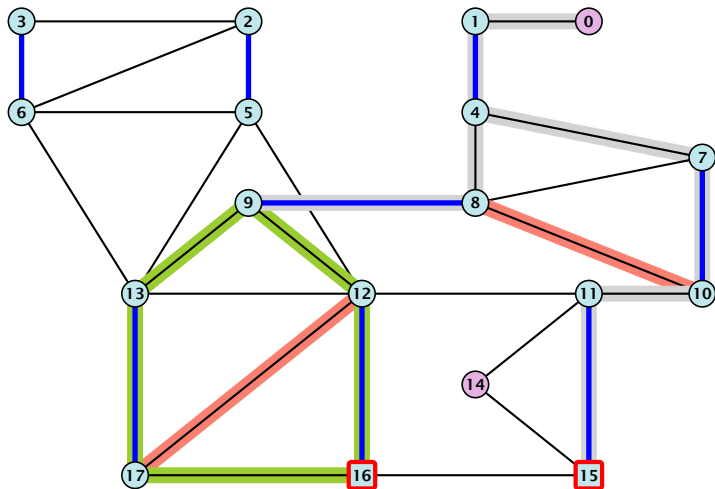
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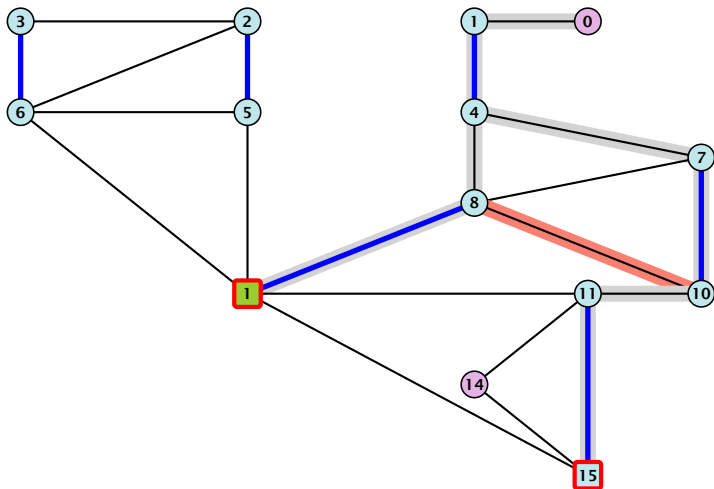
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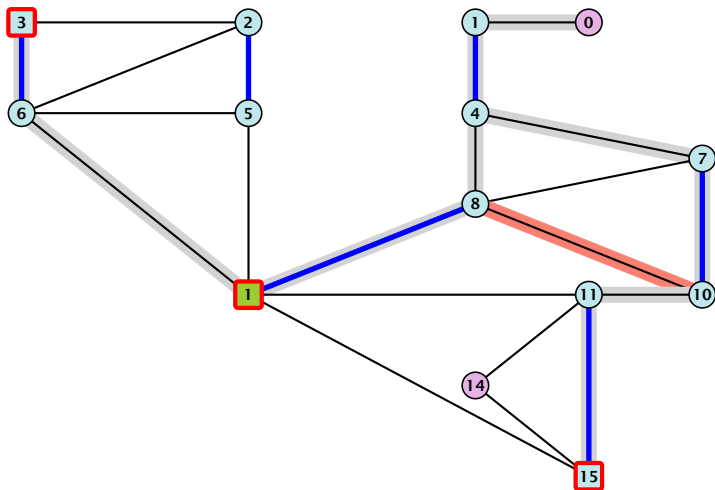
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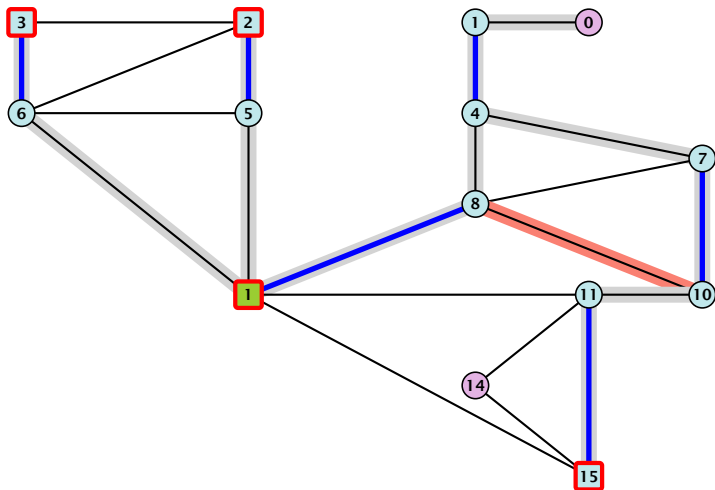
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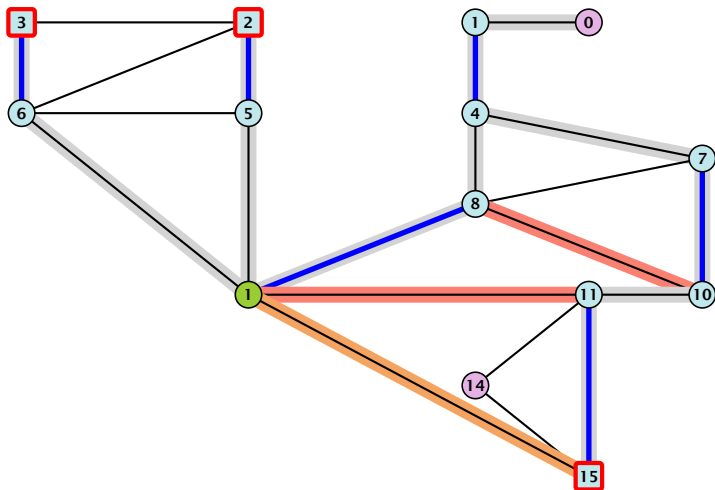
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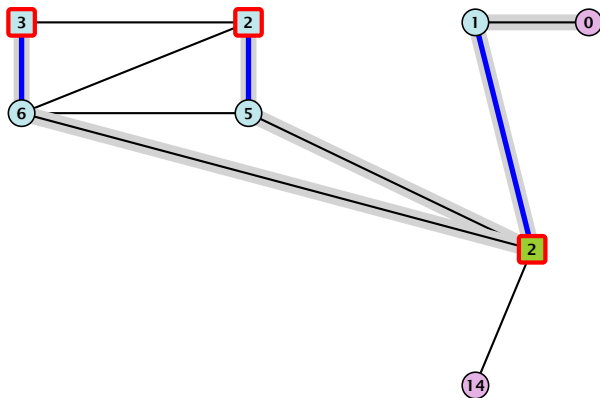
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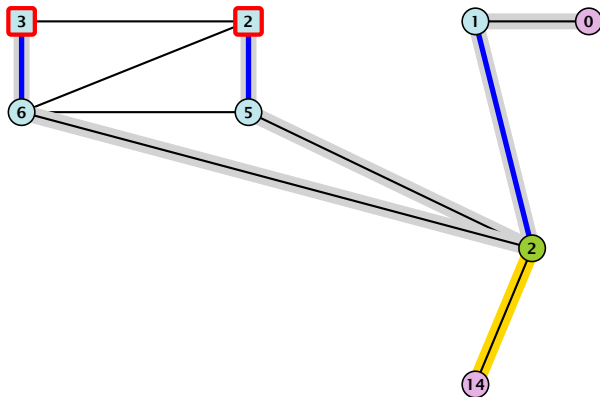
Example: Blossom Algorithm



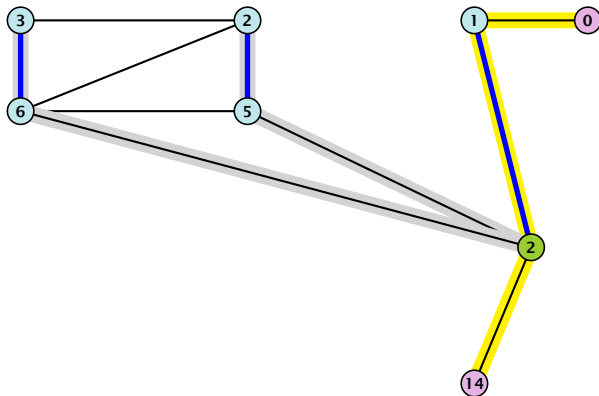
Example: Blossom Algorithm



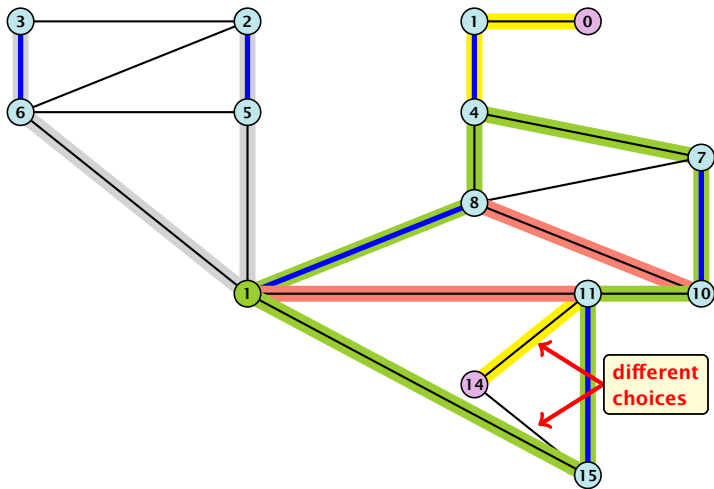
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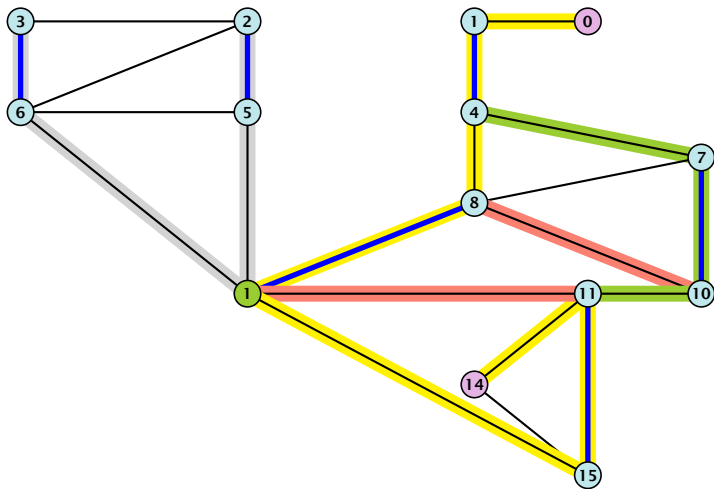
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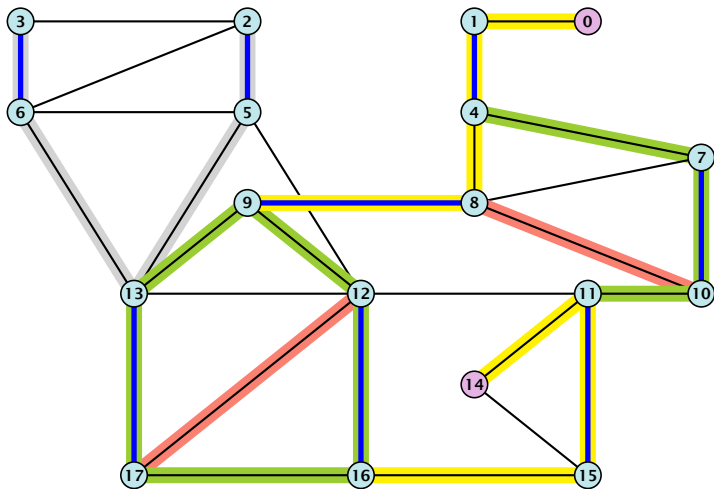
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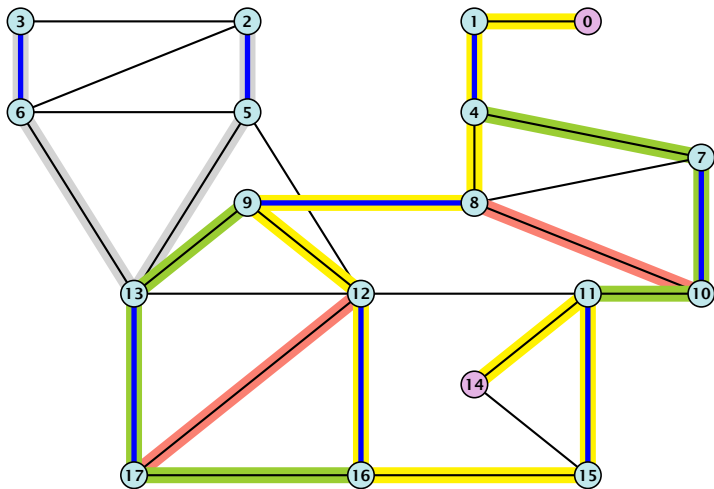
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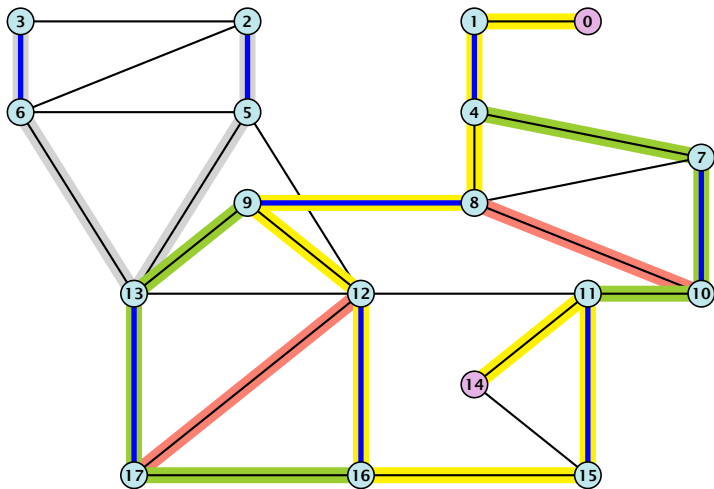
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Assume that we have contracted a blossom B w.r.t. a matching M whose base is w . We created graph $G' = G/B$ with pseudonode b . Let M' be the matching in the contracted graph.

Lemma 104

If G' contains an augmenting path p' starting at r (or the pseudo-node containing r) w.r.t. to the matching M' then G contains an augmenting path starting at r w.r.t. matching M .

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Proof.

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Case 1: non-empty stem

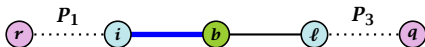
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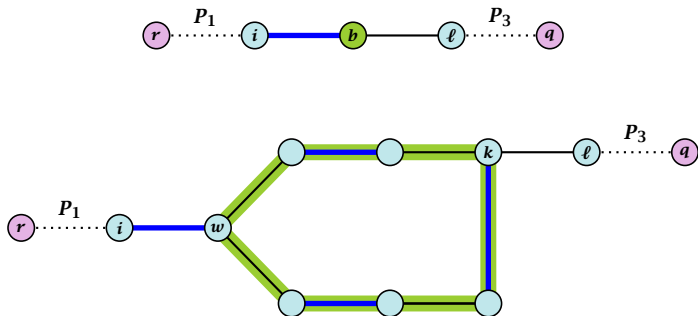


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Case 1: non-empty stem

- ▶ Next suppose that the stem is non-empty.



- ▶ After the expansion ℓ must be incident to some node in the blossom. Let this node be k .
- ▶ If $k \neq w$ there is an alternating path P_2 from w to k that ends in a matching edge.
- ▶ $P_1 \circ (i, w) \circ P_2 \circ (k, \ell) \circ P_3$ is an alternating path.
- ▶ If $k = w$ then $P_1 \circ (i, w) \circ (w, \ell) \circ P_3$ is an alternating path.

Proof.

Case 2: empty stem

- ▶ If the stem is empty then after expanding the blossom,
 $w = r$.

Proof.

Case 2: empty stem

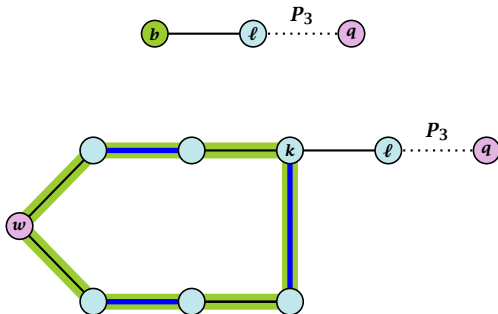
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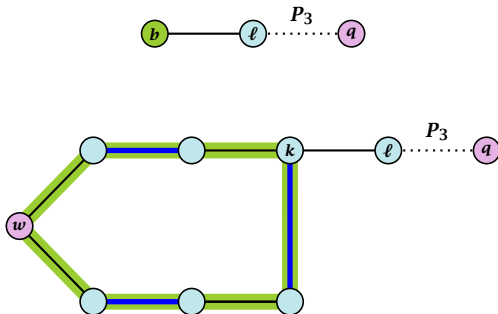
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- ▶ If the stem is empty then after expanding the blossom, $w = r$.



- ▶ The path $r \circ P_2 \circ (k, l) \circ P_3$ is an alternating path.

Lemma 105

If G contains an augmenting path P from r to q w.r.t. matching M then G' contains an augmenting path from r (or the pseudo-node containing r) to q w.r.t. M' .

Proof.

- ▶ If P does not contain a node from B there is nothing to prove.
- ▶ We can assume that r and q are the only free nodes in G .

Case 1: empty stem

Let i be the last node on the path P that is part of the blossom.

P is of the form $P_1 \circ (i, j) \circ P_2$, for some node j and (i, j) is unmatched.

$(b, j) \circ P_2$ is an augmenting path in the contracted network.

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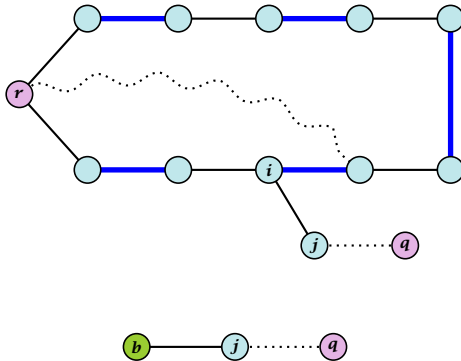
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Case 2: non-empty stem

Let P_3 be alternating path from r to w . Define $M_+ = M \oplus P_3$.

In M_+ , r is matched and w is unmatched.

G must contain an augmenting path w.r.t. matching M_+ , since M and M_+ have same cardinality.

This path must go between w and q as these are the only unmatched vertices w.r.t. M_+ .

For M'_+ the blossom has an empty stem. Case 1 applies.

G' has an augmenting path w.r.t. M'_+ . It must also have an augmenting path w.r.t. M' , as both matchings have the same cardinality.

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This path must go between r and q .

Case 2: non-empty stem

Let P_3 be alternating path from r to w . Define $M_+ = M \oplus P_3$.

In M_+ , r is matched and w is unmatched.

G must contain an augmenting path w.r.t. matching M_+ , since M and M_+ have same cardinality.

This path must go between w and q as these are the only unmatched vertices w.r.t. M_+ .

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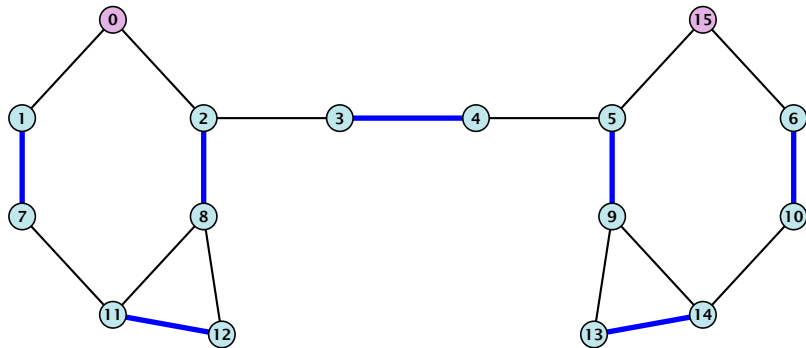
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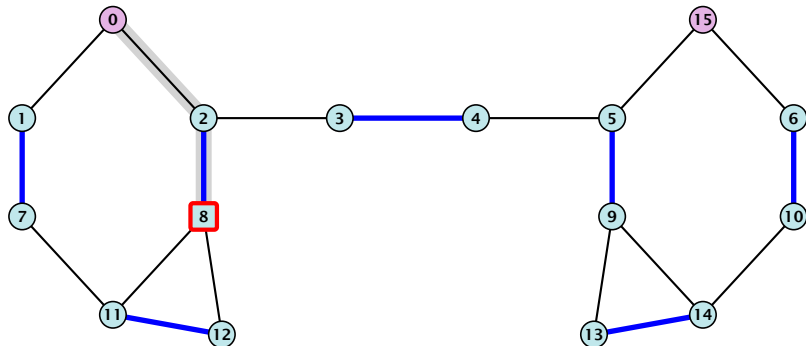
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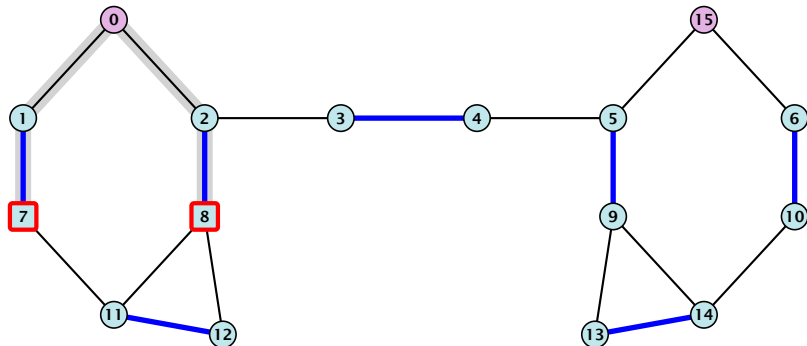
Example: Blossom Algorithm



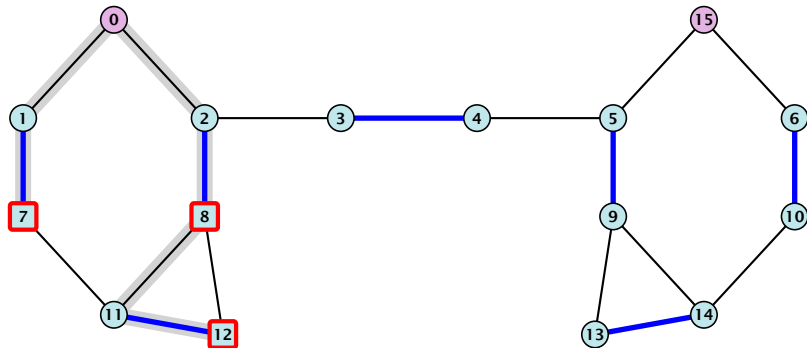
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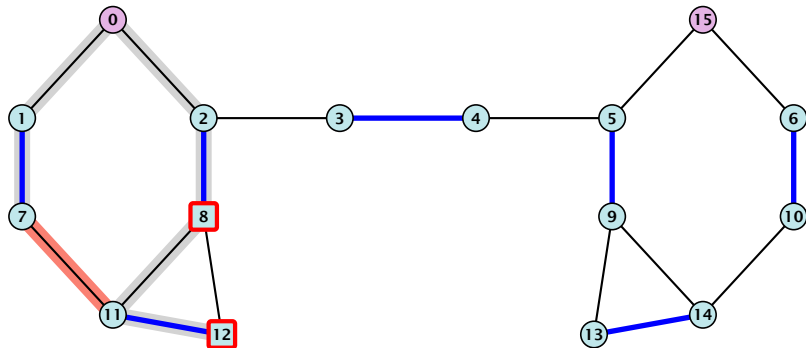
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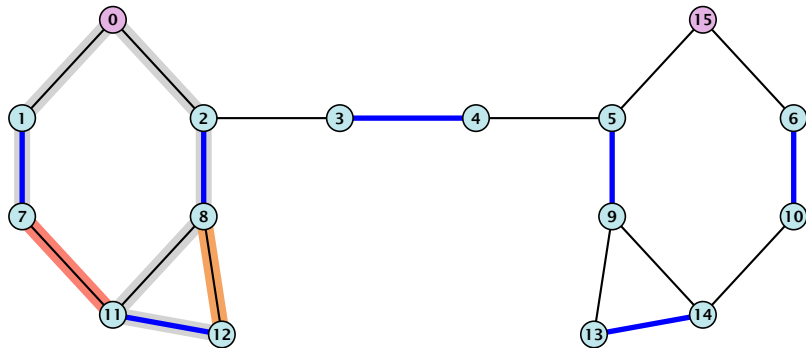
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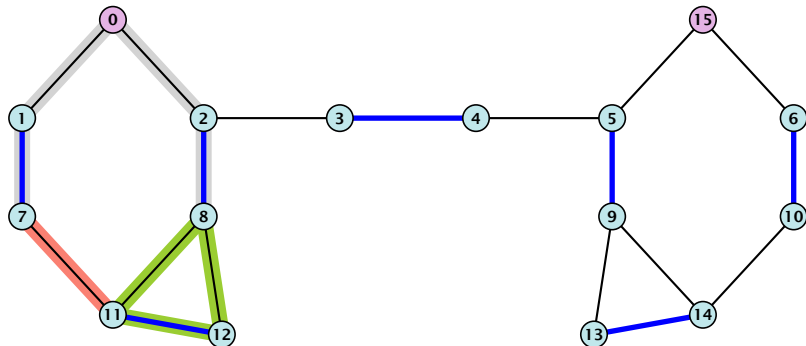
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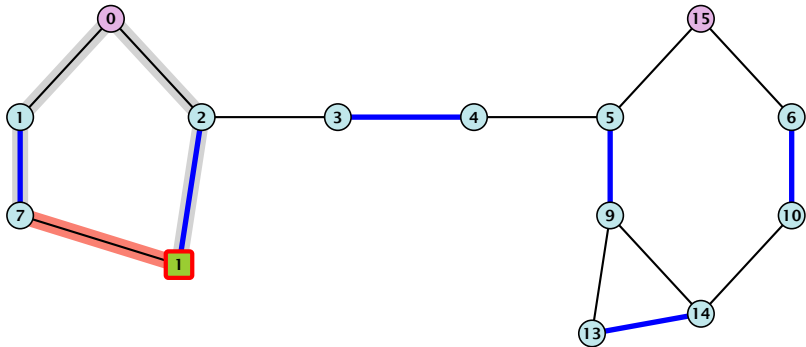
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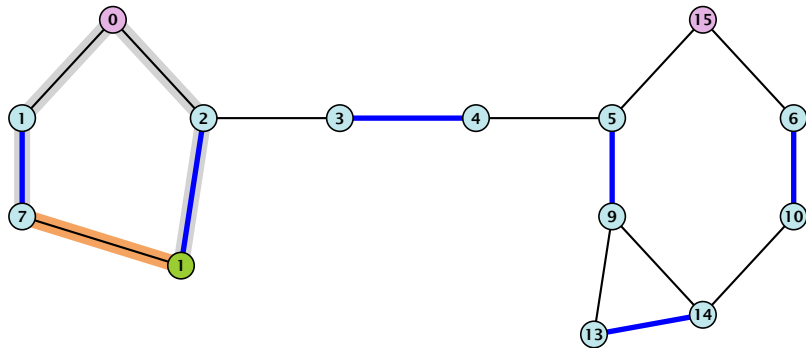
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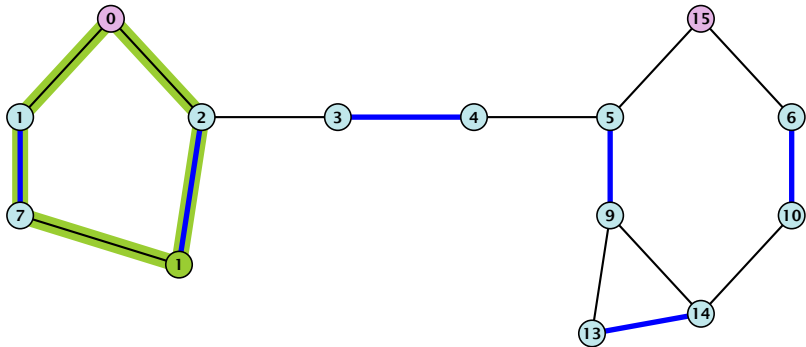
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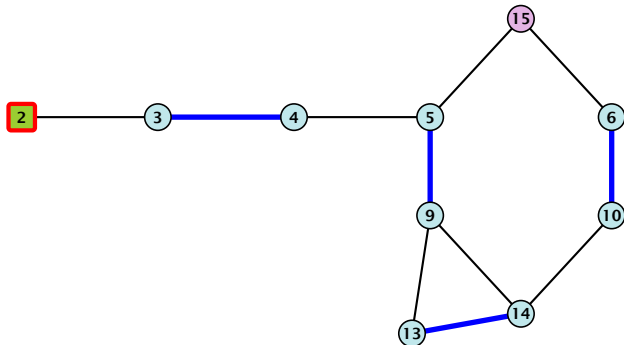
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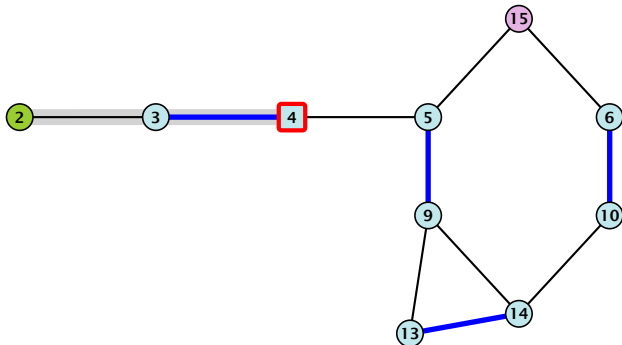
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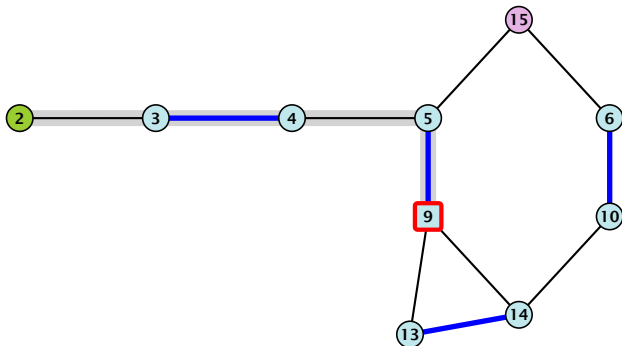
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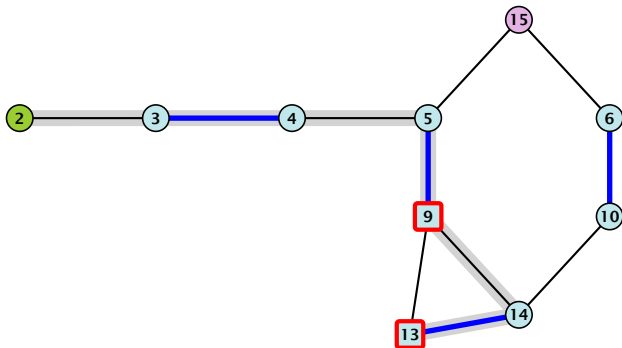
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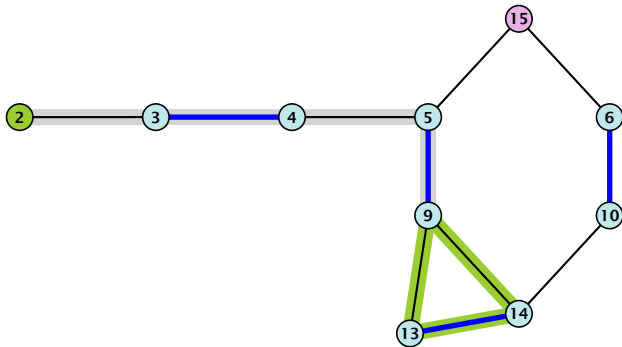
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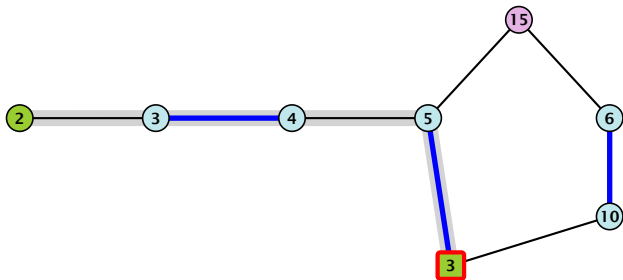
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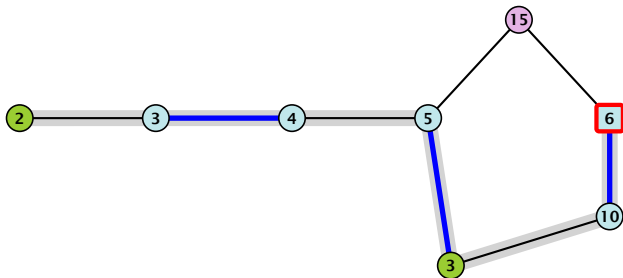
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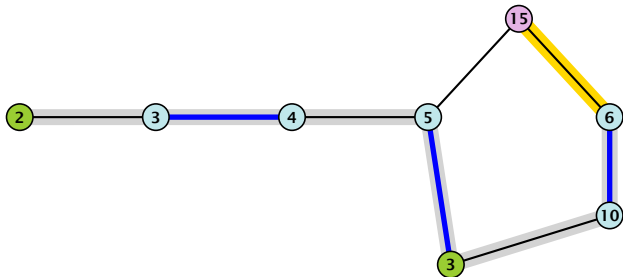
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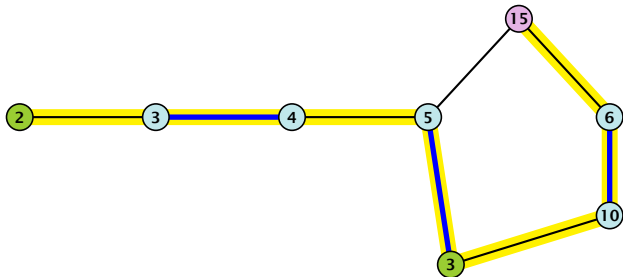
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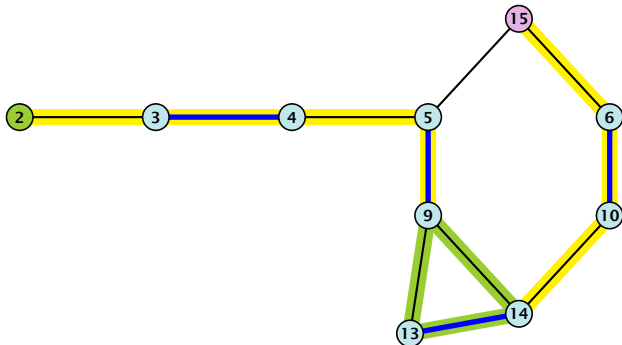
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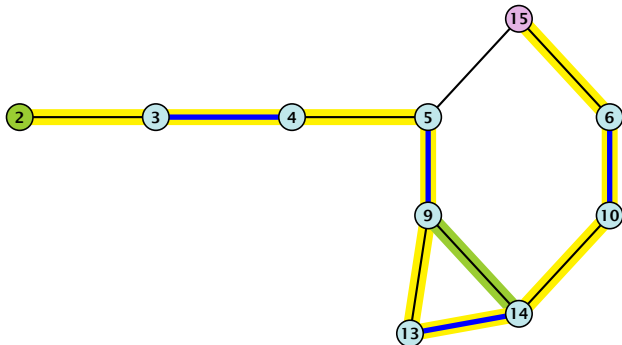
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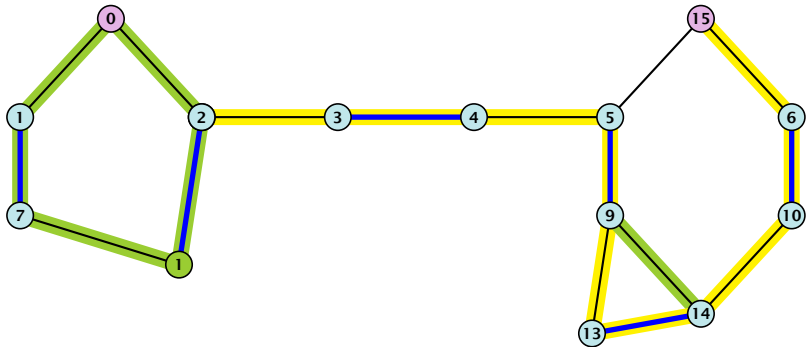
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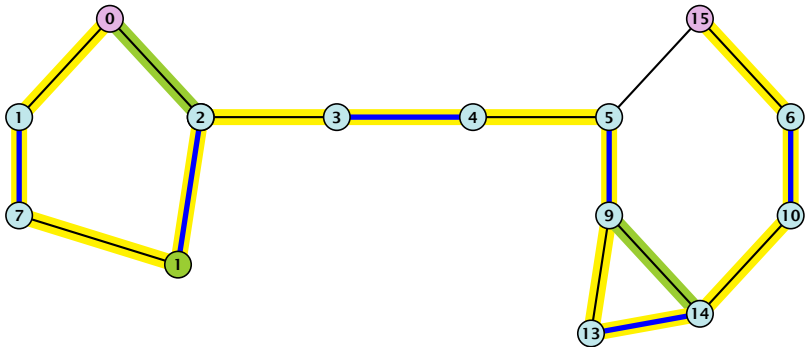
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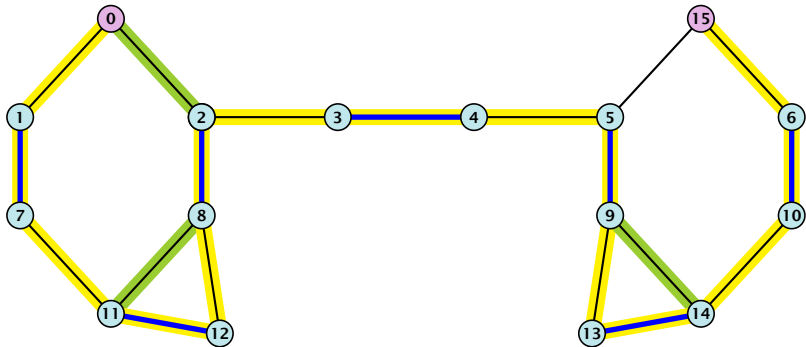
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