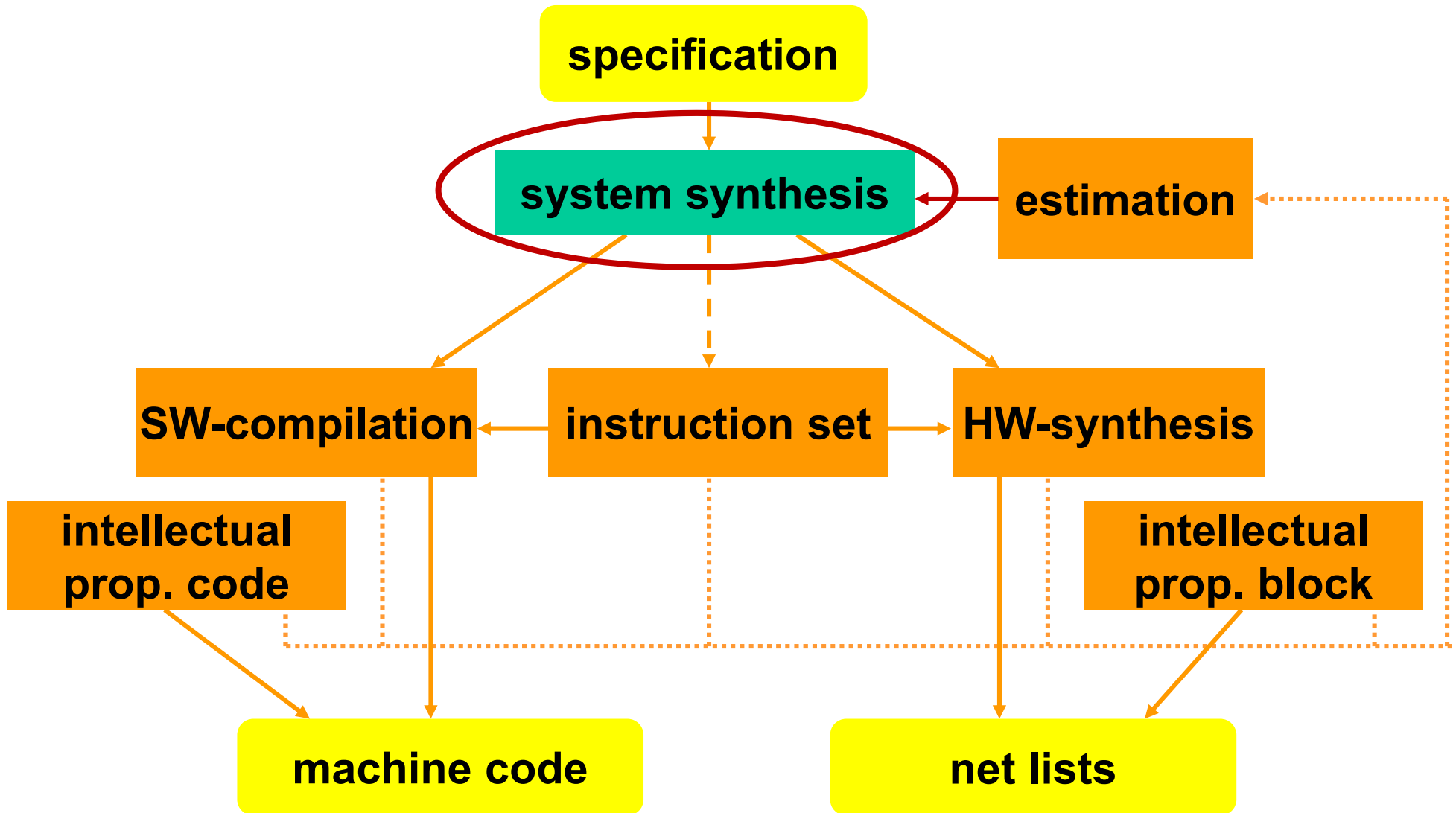


Hardware-Software Codesign

4. System Partitioning

Lothar Thiele

System Design



Mapping

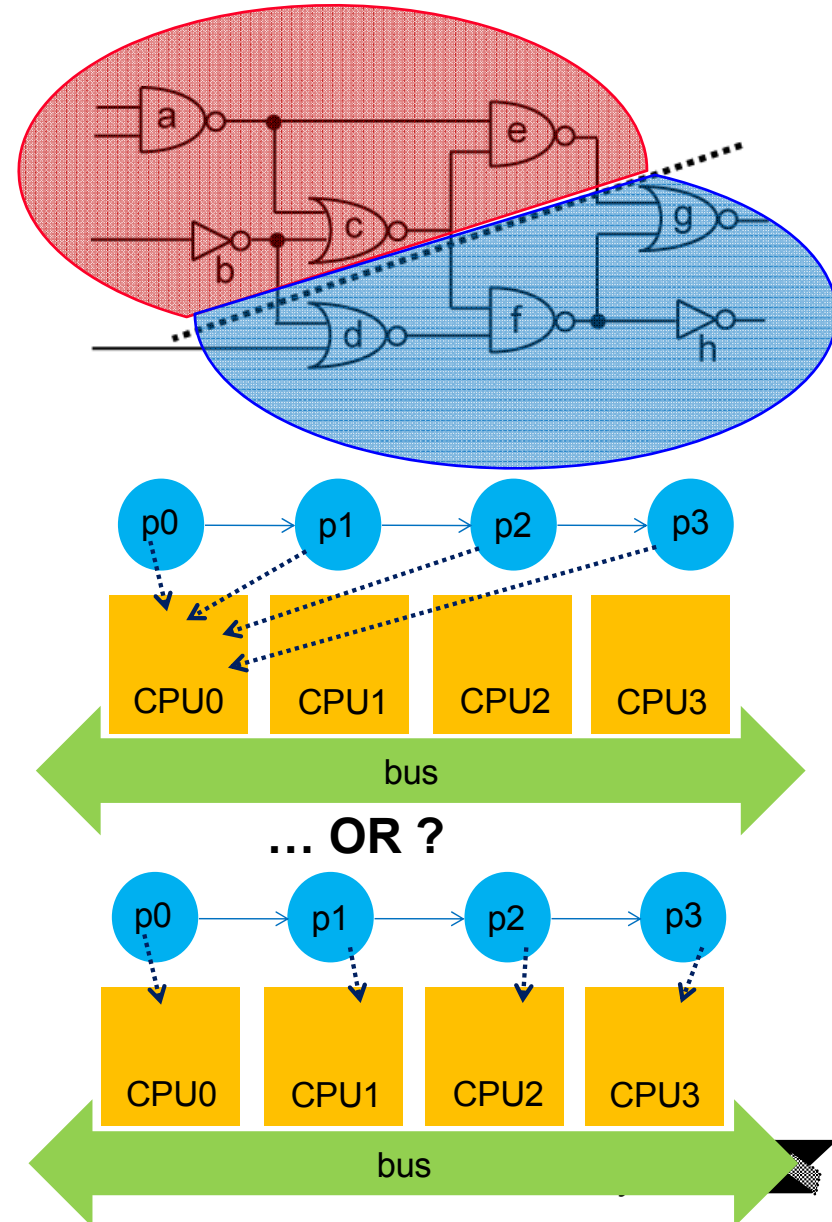
Mapping transforms behavior into structure and execution:

- ▶ **allocation**: select components
 - ▶ **binding**: assign functions to components
 - ▶ **scheduling**: determine execution order
 - ▶ ... finally, synthesis results into **implementation**
- } *partitioning*
- } *mapping*

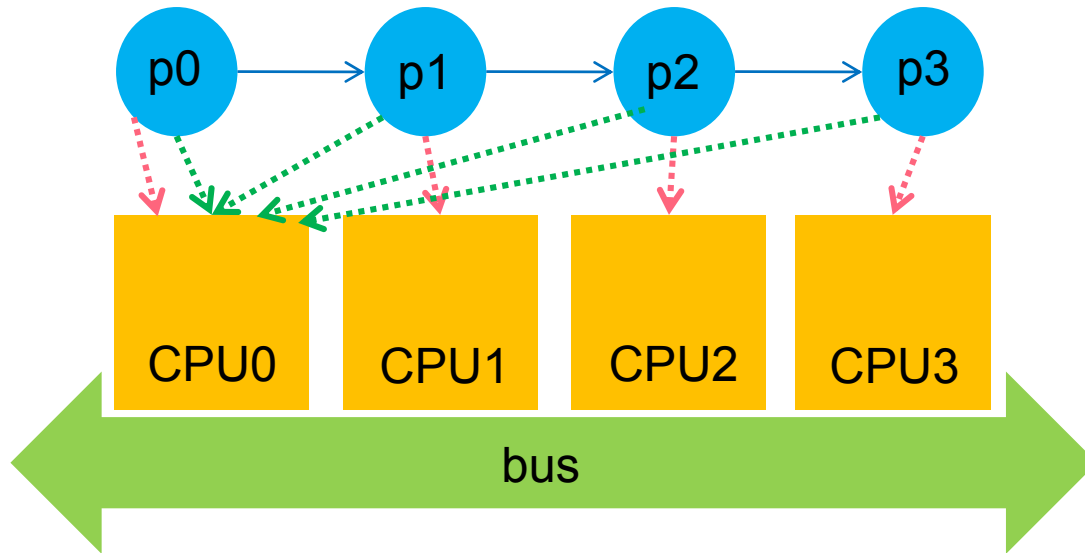
Levels of Abstractions

Mapping can be done

- ▶ at low level: register transfer level (RTL) or netlist level
 - e.g., split a digital circuit and map it to several devices (FPGAs, ASICs)
 - system parameters (e.g., area, delay) relatively easy to determine
- ▶ at high level: system level
 - comparison of design alternatives for optimality (design space exploration)
 - system parameters are unknown and difficult to determine
 - to be estimated via analysis, simulation, (rapid) prototyping



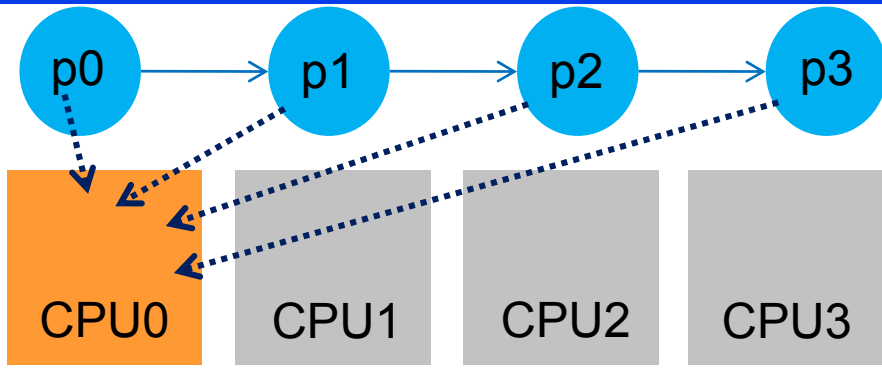
Model-Based Synthesis – Example



optimal C: N:1 mapping
optimal L: 1:1 mapping

- ▶ **considered performance**
 - **cost C:** cost of allocated components, e.g., sum
 - **latency L:** due to scheduling (resource sharing)
- ▶ **conflicting design goals**
 - feasible schedule $L \leq L_{\max}$
 - feasible allocation $C \leq C_{\max}$

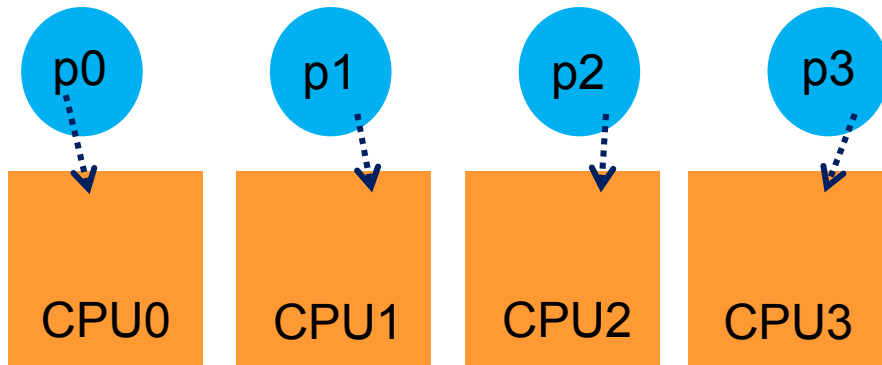
Example – Alternatives



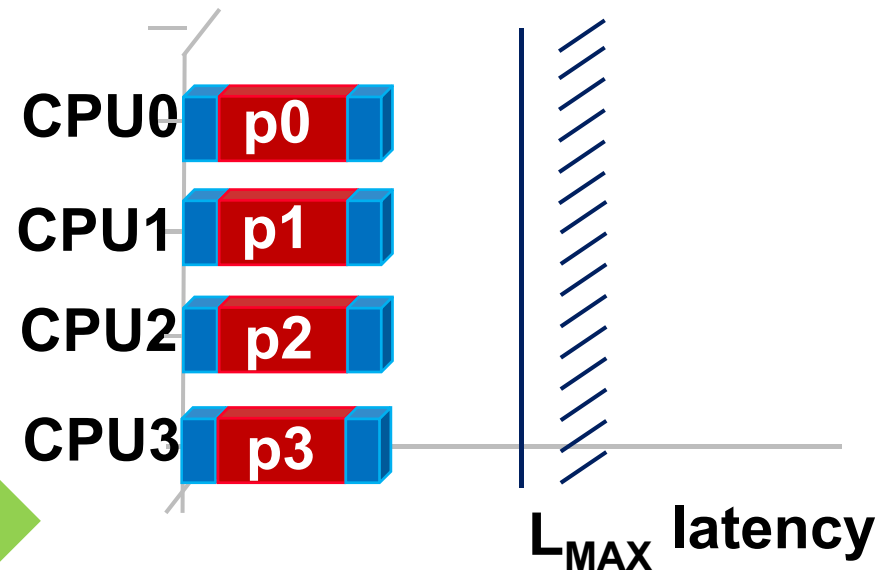
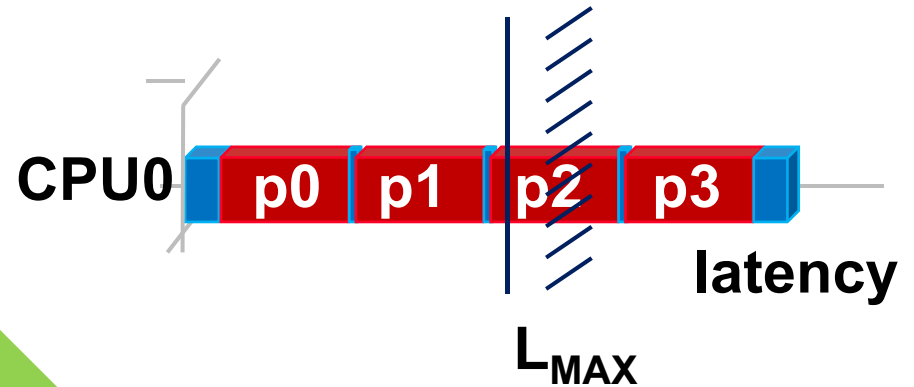
bus

optimal C: N:1 mapping

optimal L: 1:1 mapping



bus



Cost Functions

Quantitatively measure performance of a design point

- system cost C [\$]
- latency L [sec]
- power consumption P [W]
- ...

Estimation is required to find C, L, P values, for each design point

- *example*: linear cost (preference) function with penalty

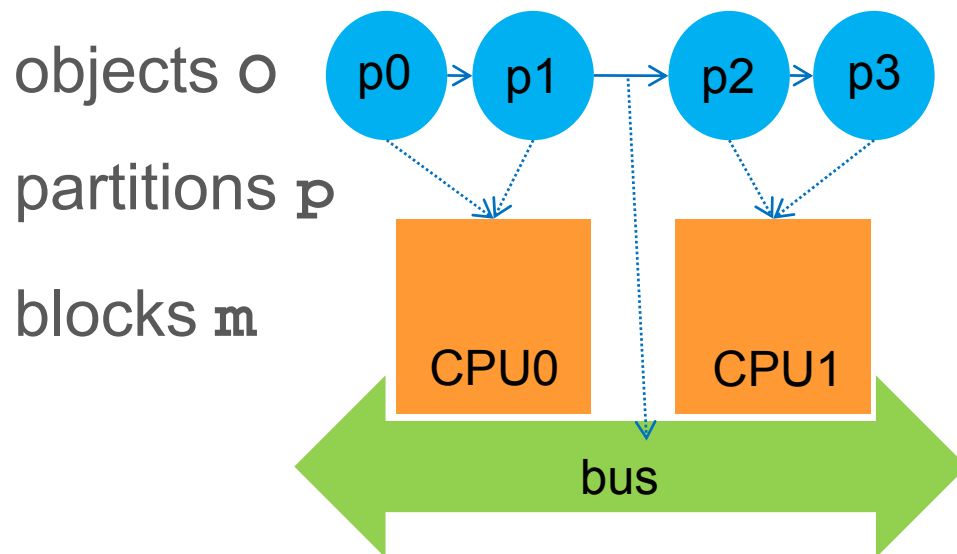
$$f(C, L, P) = k_1 \cdot h_C(C, C_{\max}) + k_2 \cdot h_L(L, L_{\max}) + k_3 \cdot h_P(P, P_{\max})$$

- h_C, h_L, h_P ... denote how strong C, L, P violate design constraints $C_{\max}, L_{\max}, P_{\max}$
- k_1, k_2, k_3 ... weighting and normalization

The Formal Partitioning Problem

assign n objects $O = \{o_1, \dots, o_n\}$ to m blocks (also called partitions) $P = \{p_1, \dots, p_m\}$, such that

- $p_1 \cup p_2 \cup \dots \cup p_m = O$ (all objects are assigned –mapped)
- $p_i \cap p_j = \{ \}$ $\forall i, j: i \neq j$ (an object is not assigned or “mapped” twice)
- and costs $c(P)$ are minimized



*note: in **system synthesis** (simple model)*

- ▶ objects = process network graph nodes
- ▶ blocks = architecture graph nodes
- ▶ cost = measured/estimated with dedicated cost functions (e.g., latency, power, hardware cost)

Partitioning Methods

▶ Exact methods

- enumeration
- *integer linear programs (ILP) (→see next slides)*

▶ Heuristic methods

- constructive methods
 - random mapping
 - hierarchical clustering
- iterative methods
 - Kernighan-Lin algorithm
 - simulated annealing
 - evolutionary algorithms

Integer Programming Model

Ingredients:

- objective function (cost)
 - constraints
- } involving linear expressions of integer variables from a set \mathbf{x}

objective $C = \sum_{x_i \in X} a_i x_i$ with $a_i \in \mathbb{R}, x_i \in \mathbb{N}$ (1)

constraints $\forall j \in J : \sum_{x_i \in X} b_{i,j} x_i \geq c_j$ with $b_{i,j}, c_j \in \mathbb{R}$ (2)

Integer programming (IP) problem:

minimize objective function (1) subject to constraints (2)

note: if all x_i are constrained to be either 0 or 1, the IP problem is said to be a 0/1 integer programming problem

Small Example of 0/1 IP

minimize: $C = 5x_1 + 6x_2 + 4x_3$

subject to: $x_1 + x_2 + x_3 \geq 2$

$x_1, x_2, x_3 \in \{0,1\}$

x_1	x_2	x_3	C
0	1	1	10
1	0	1	9
1	1	0	11
1	1	1	15

← **optimal (minimal)**

Integer Linear Program for Partitioning

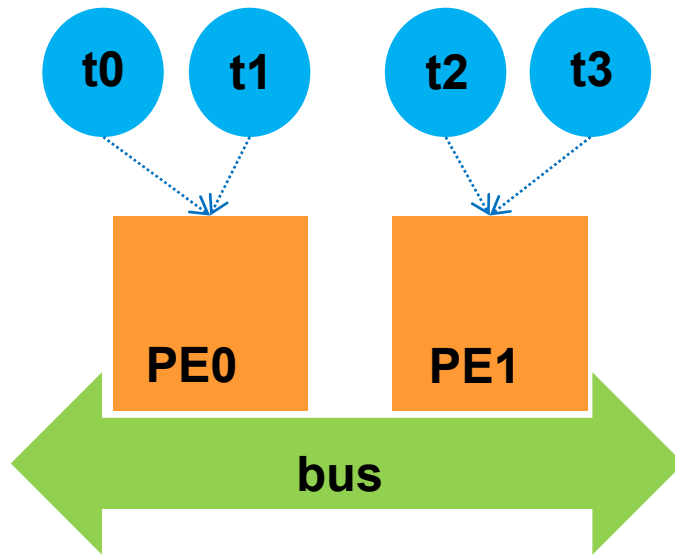
- ▶ Binary variables $x_{i,k}$
 - $x_{i,k} = 1$: object o_i in block p_k
 - $x_{i,k} = 0$: object o_i not in block p_k
- ▶ Cost $c_{i,k}$, if object o_i is in block p_k
- ▶ Integer linear program:

$$x_{i,k} \in \{0,1\} \quad 1 \leq i \leq n, 1 \leq k \leq m$$

$$\sum_{k=1}^m x_{i,k} = 1 \quad 1 \leq i \leq n$$

$$\text{minimize} \quad \sum_{k=1}^m \sum_{i=1}^n x_{i,k} \cdot c_{i,k} \quad 1 \leq k \leq m, 1 \leq i \leq n$$

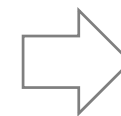
Example – Partitioning



task \ PE	t0	t1	t2	t3
PE0	1	1	0	0
PE1	0	0	1	1

e.g., optimized for a load balanced system

exe. time	t0	t1	t2	t3
PE0	5	15	10	30
PE1	10	20	10	10



load balancing:

$$\text{load}_{\text{PE0}} = 5+15$$

$$\text{load}_{\text{PE1}} = 10+10$$

Variations in ILP

Additional constraints:

- e.g., maximum h_k objects in block k

$$\sum_{i=1}^n x_{i,k} \leq h_k \quad 1 \leq k \leq m$$

Maximizing the cost function:

- can be done by setting $C' = -C$ in a minimization problem

ILP for synthesis

Solving the synthesis problem with ILP is very popular:

- If not solving to optimality, runtimes are acceptable and a solution with guaranteed quality can be determined.
- Scheduling can be integrated.
- Various additional constraints can be added.
- *However, finding the right equations to model the constraints is an art.*

Remarks on Integer Programming

Integer programming is NP-complete

- In practice, *runtimes can increase exponentially* with the size of the problem.
- But problems of some thousands of variables can still be solved with *commercial solvers* (depending on the size/structure of the problem) or *approximation algorithms (heuristics)*.
- IP models can be a *good starting point* for designing heuristic optimization methods.

Partitioning Methods

▶ exact methods

- enumeration
- integer linear programs (ILP)

▶ heuristic methods

- ***constructive methods (→see next slides)***
 - random mapping
 - hierarchical clustering
- iterative methods
 - Kernighan-Lin algorithm
 - simulated annealing
 - evolutionary algorithms

Constructive Methods

► *Examples*

▪ random mapping

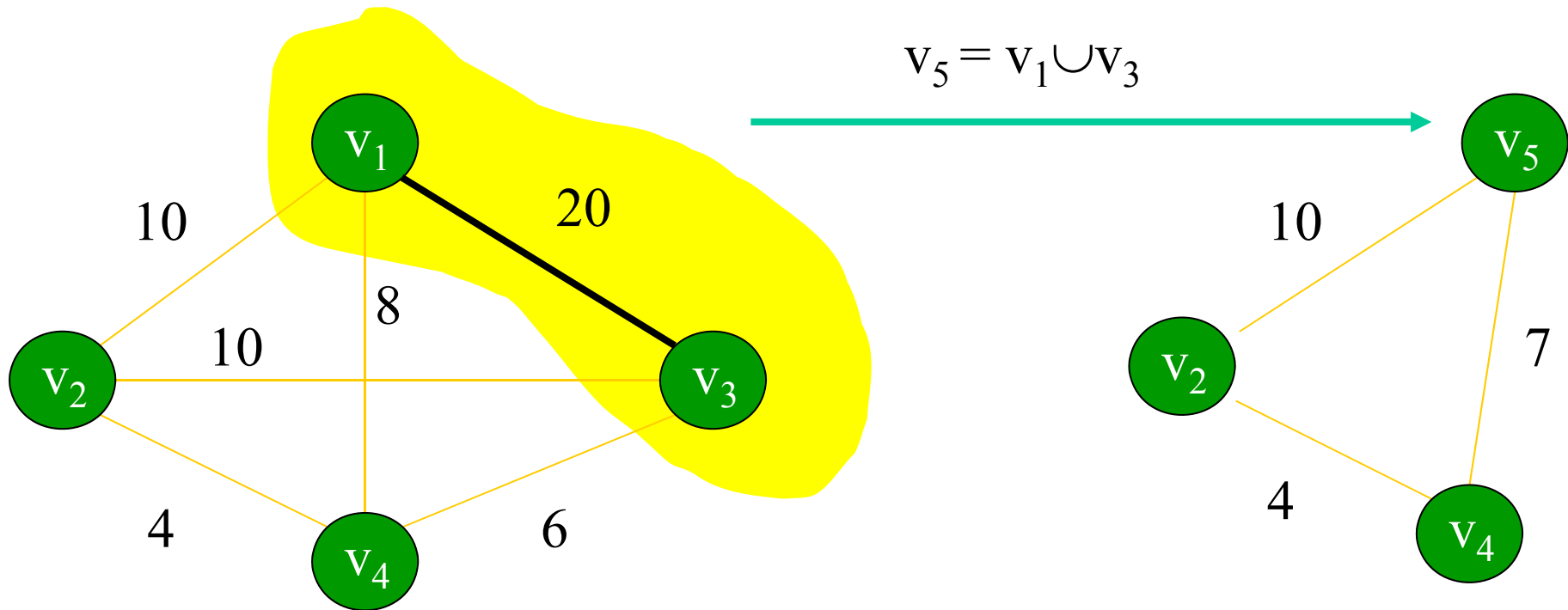
- each object is assigned to a block randomly

▪ hierarchical clustering

- stepwise grouping of (e.g., two) objects
- and evaluate closeness function (how desirable it is to group objects)

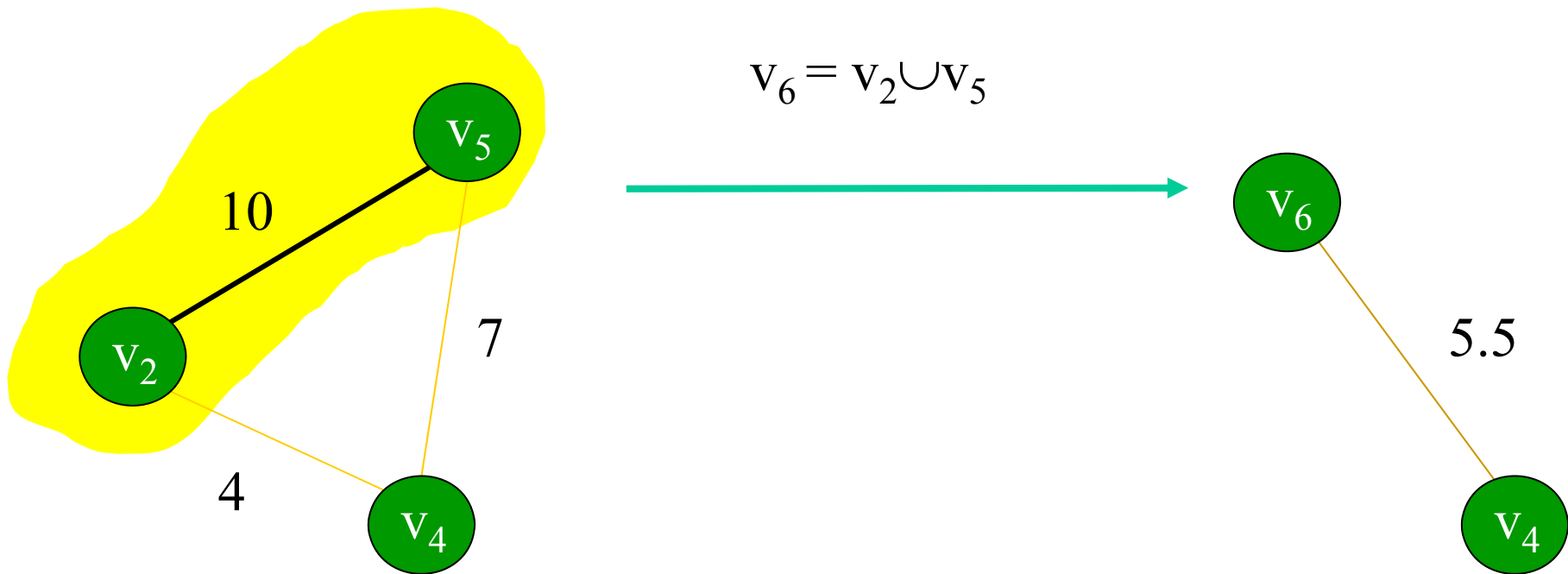
- Constructive methods are often used to generate a starting partition for iterative methods

Hierarchical Clustering Example (1)

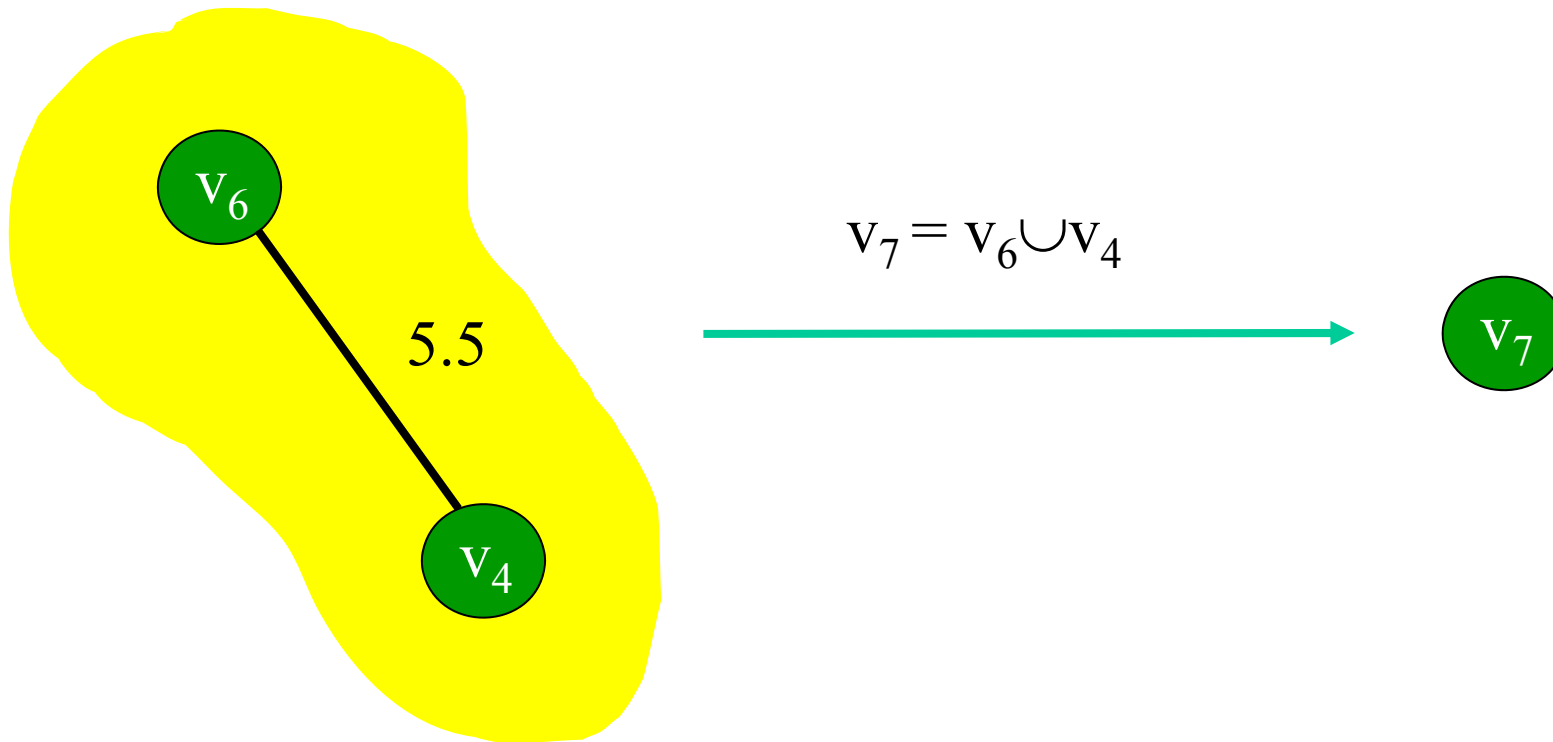


closeness function: arithmetic mean of weights

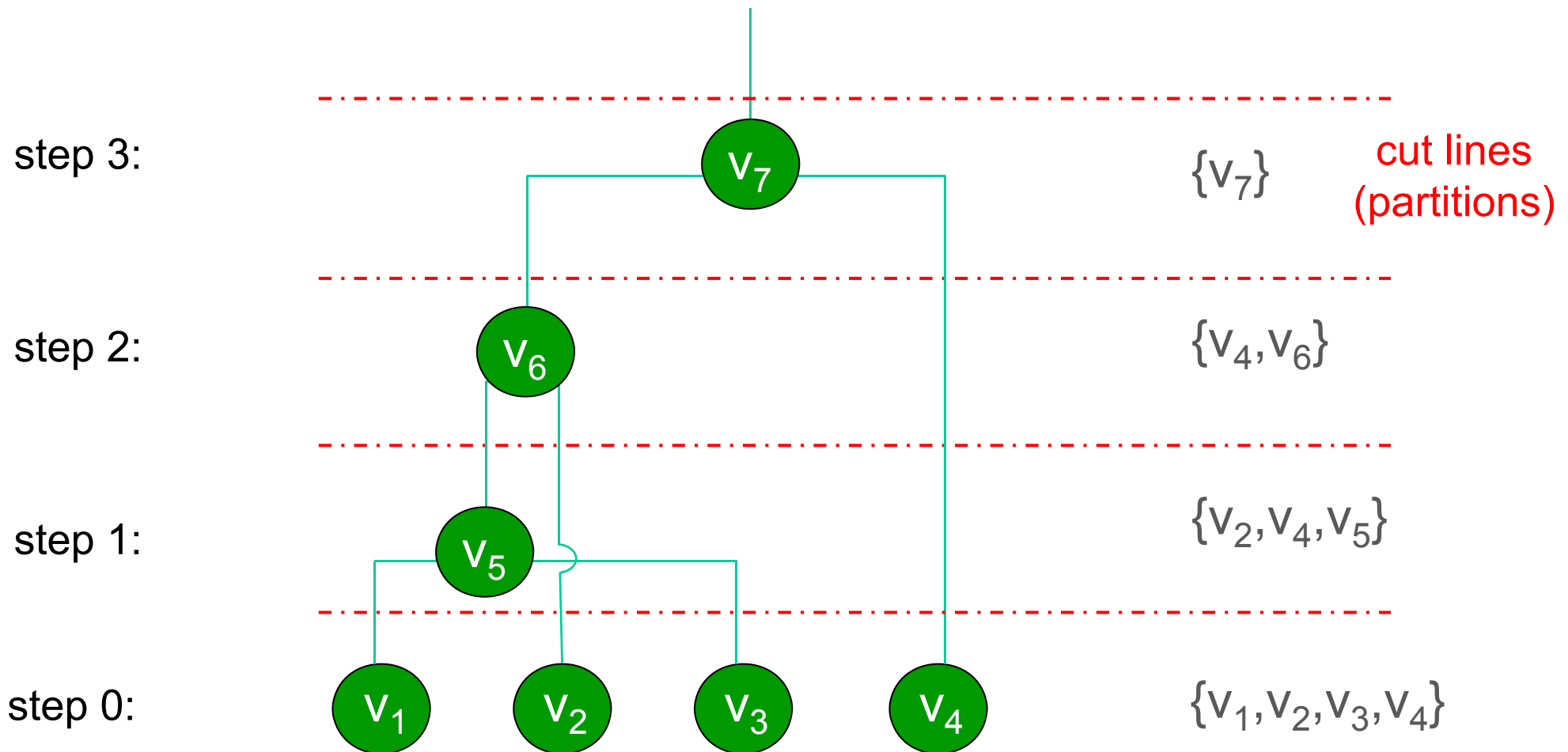
Hierarchical Clustering Example (2)



Hierarchical Clustering Example (3)



Hierarchical Clustering – Summary



Partitioning Methods

▶ exact methods

- enumeration
- integer linear programs (ILP)

▶ heuristic methods

- constructive methods
 - random mapping
 - hierarchical clustering
- **iterative methods** (*→see next slides*)
 - Kernighan-Lin algorithm
 - simulated annealing
 - evolutionary algorithms

Iterative Methods (1)

Often used principle for iterative methods:

- ▶ start with some initial configuration (partitioning)
- ▶ search *neighborhood* (similar partitions) and *select a neighbor* as candidate
- ▶ evaluate *fitness (cost) function of candidate*
 - accept candidate using acceptance rule
 - if not, select another neighbor
- ▶ stop if quality is sufficiently high, if no improvement can be found, or after some fixed time

Ingredients:

- ▶ initial configuration, function to find a *neighbor* as next candidate, cost function, acceptance rule, stop criterion

Iterative Methods (2)

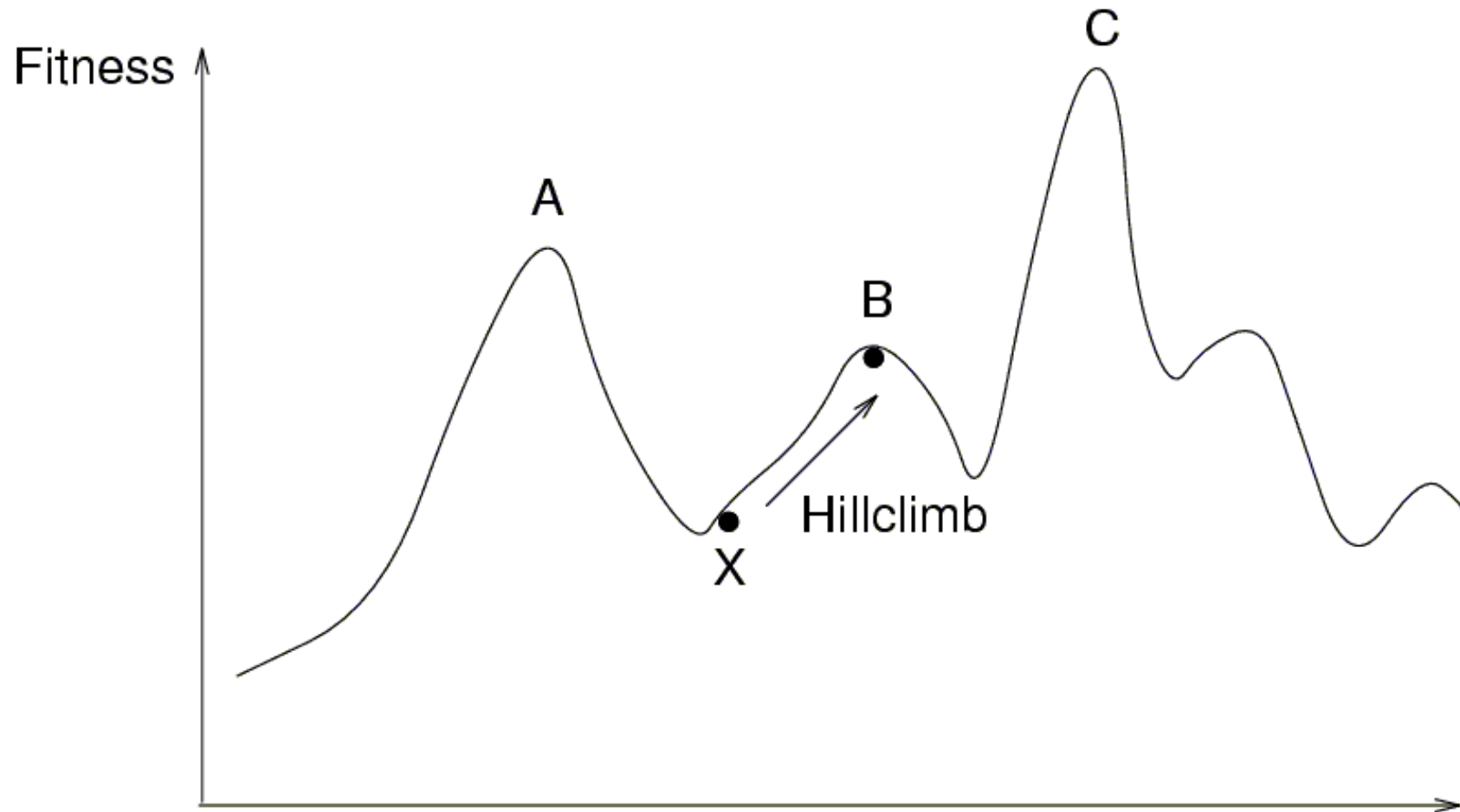
Simple iterative improvement or “*hill climbing*”:

- ▶ candidate is always and only accepted if cost is lower (or fitness is higher) than current configuration
- ▶ stop when no neighbor with lower cost (higher fitness) can be found

Disadvantages:

- ▶ local optimum as best result
- ▶ local optimum depends on initial configuration
- ▶ generally no upper bound on iteration length

Iterative Methods – Illustration

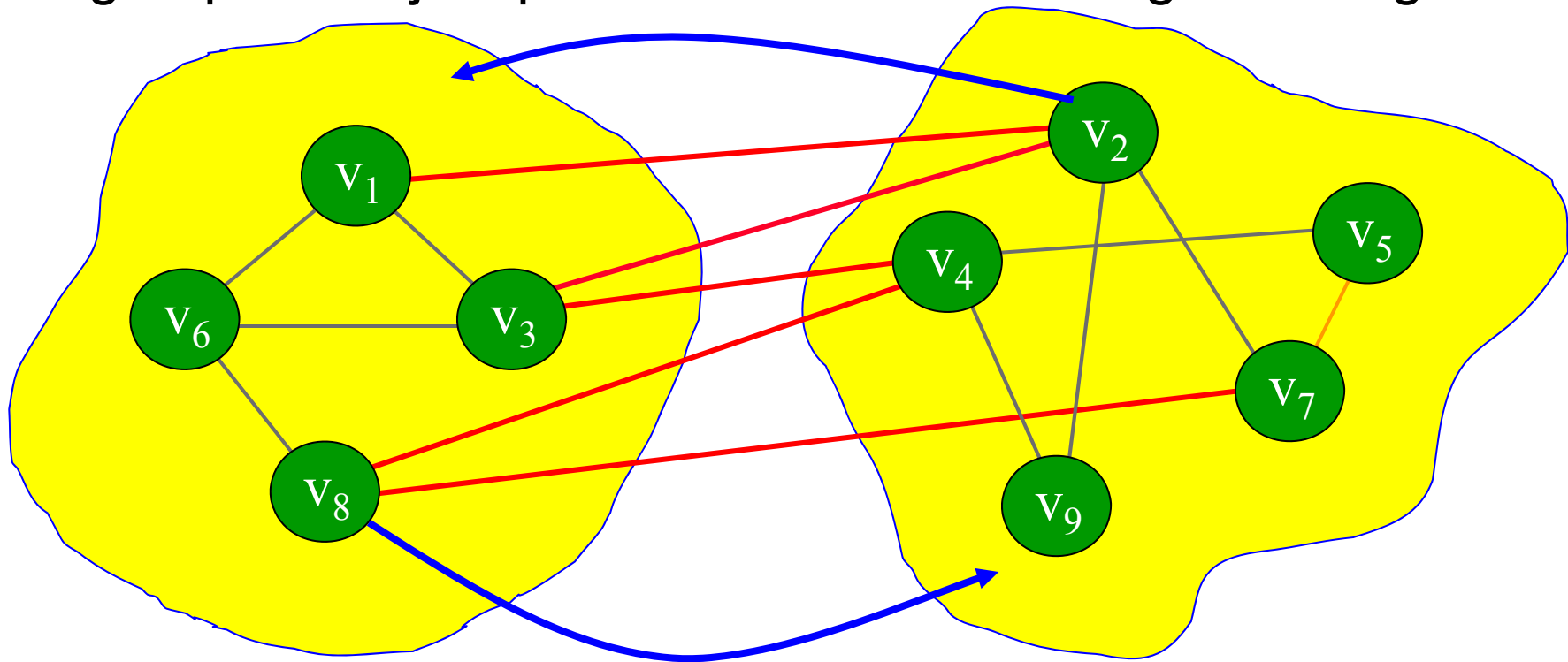


How to Cope with Disadvantages?

- ▶ Repeat algorithm **many times** with **different initial configurations**
- ▶ Use information gathered in **previous runs**
- ▶ Use a more **complex “acceptance rule”** to jump out of local optimum
- ▶ Use a more **complex strategy** that accepts sometimes randomly generated solutions

Iterative Methods – Simple Greedy Heuristic

- ▶ Iterate until no improvement in cost:
re-group the object pairs that leads to the largest cost gain



example: cost = number of edges crossing the partitions
before re-group: 5 ; after re-group: 4 ; gain = 1

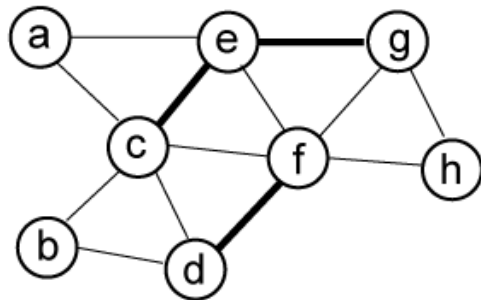
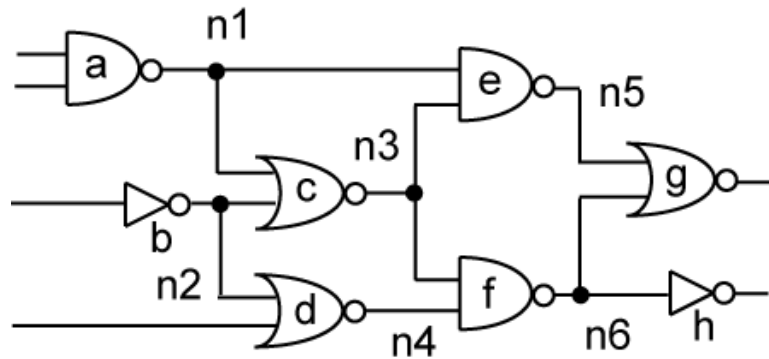
Iterative Methods – Kernighan-Lin

Improved algorithm: Kernighan-Lin:

- ▶ as long as a better partition is found
 - from all possible pairs of objects
 - *virtually* re-group the “best” (lowest cost of resulting partition)
 - from the remaining (not yet touched) objects
 - *virtually* re-group the “best” pair
 - continue until all objects have been re-grouped
 - from these $n/2$ partitions, take the one with smallest cost and *actually* perform the corresponding re-group operations

Illustration of KL Algorithm (1)

Example: partitioning of digital circuit



communication cost
from node x to node y

cost matrix $c(x,y)$

$c(x,y)$	a	b	c	d	e	f	g	h
a	0	0	.5	0	.5	0	0	0
b	0	0	.5	.5	0	0	0	0
c	.5	.5	0	.5	1	.5	0	0
d	0	.5	.5	0	0	1	0	0
e	.5	0	1	0	0	.5	1	0
f	0	0	.5	1	.5	0	.5	.5
g	0	0	0	0	1	.5	0	.5
h	0	0	0	0	0	.5	.5	0

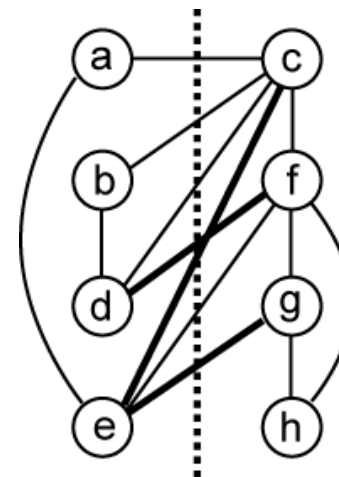
Illustration of KL Algorithm (2)

first re-group

pair	$E_x - I_x$	$E_y - I_y$	$c(x, y)$	gain
(a, c)	0.5 - 0.5	2.5 - 0.5	0.5	1
(a, f)	0.5 - 0.5	1.5 - 1.5	0	0
(a, g)	0.5 - 0.5	1 - 1	0	0
(a, h)	0.5 - 0.5	0 - 1	0	-1
(b, c)	0.5 - 0.5	2.5 - 0.5	0.5	1
(b, f)	0.5 - 0.5	1.5 - 1.5	0	0
(b, g)	0.5 - 0.5	1 - 1	0	0
(b, h)	0.5 - 0.5	0 - 1	0	-1
(d, c)	1.5 - 0.5	2.5 - 0.5	0.5	2
(d, f)	1.5 - 0.5	1.5 - 1.5	1	-1
(d, g)	1.5 - 0.5	1 - 1	0	1
(d, h)	1.5 - 0.5	0 - 1	0	0
(e, c)	2.5 - 0.5	2.5 - 0.5	1	2
(e, f)	2.5 - 0.5	1.5 - 1.5	0.5	1
(e, g)	2.5 - 0.5	1 - 1	1	0
(e, h)	2.5 - 0.5	0 - 1	0	1

some definitions

- ▶ E_i = external costs of vertex i
- ▶ I_i = internal costs of vertex i
- ▶ $D_i = E_i - I_i$ = desirability to move a vertex (x or y)
- ▶ $\text{gain} = D_x + D_y - 2 * c(x, y) =$ gain due to change in cut costs



initial partitioning

Illustration of KL Algorithm (3)

second re-group

pair	$E_x - I_x$	$E_y - I_y$	$c(x, y)$	gain
(a, f)	0 - 1	1 - 2	0	-2
(a, g)	0 - 1	1 - 1	0	-1
(a, h)	0 - 1	0 - 1	0	-2
(b, f)	0.5 - 0.5	1 - 2	0	-1
(b, g)	0.5 - 0.5	1 - 1	0	0
(b, h)	0.5 - 0.5	0 - 1	0	-1
(e, f)	1.5 - 1.5	1 - 2	0.5	-2
(e, g)	1.5 - 1.5	1 - 1	1	-2
(e, h)	1.5 - 1.5	0 - 1	0	-1

some definitions

- ▶ E_i = external costs of vertex i
- ▶ I_i = internal costs of vertex i
- ▶ $D_i = E_i - I_i$ = desirability to move a vertex (x or y)
- ▶ $\text{gain} = D_x + D_y - 2 * c(x, y) =$ gain due to change in cut costs

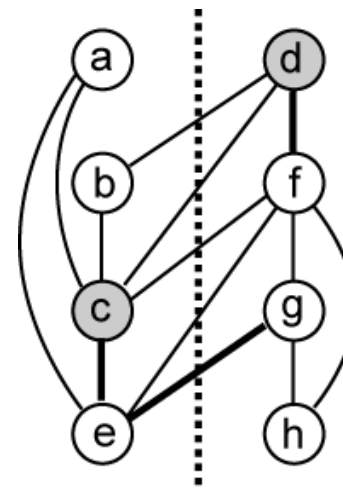


Illustration of KL Algorithm (4)

third re-group

pair	$E_x - I_x$	$E_y - I_y$	$c(x, y)$	gain
(a, f)	$0 - 1$	$1.5 - 1.5$	0	-1
(a, h)	$0 - 1$	$0.5 - 0.5$	0	-1
(e, f)	$0.5 - 2.5$	$1.5 - 1.5$	0.5	-3
(e, h)	$0.5 - 2.5$	$0.5 - 0.5$	0	-2

some definitions

- ▶ E_i = external costs of vertex i
- ▶ I_i = internal costs of vertex i
- ▶ $D_i = E_i - I_i$ = desirability to move a vertex (x or y)
- ▶ $\text{gain} = D_x + D_y - 2 * c(x, y) =$ gain due to change in cut costs

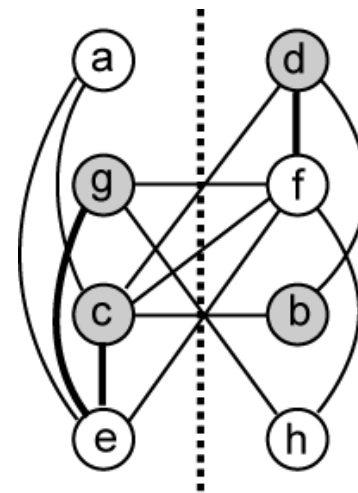


Illustration of KL Algorithm (5)

... and final re-group

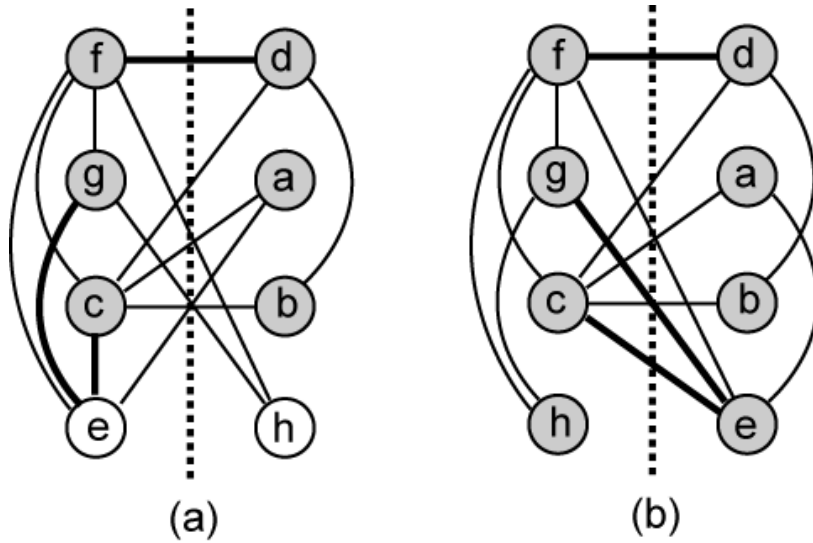
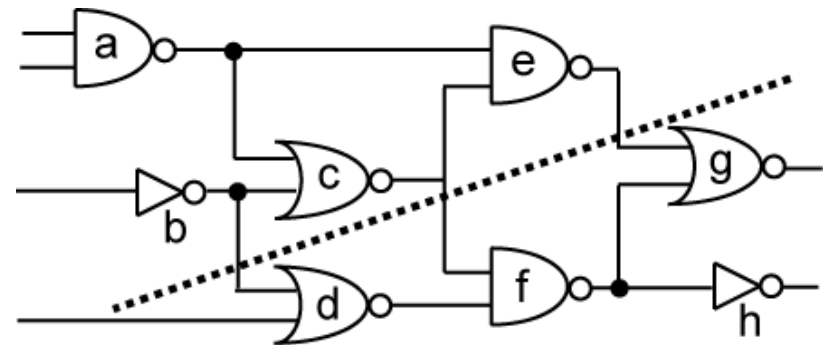


Illustration of KL Algorithm (6)

- ▶ Two best solutions found:

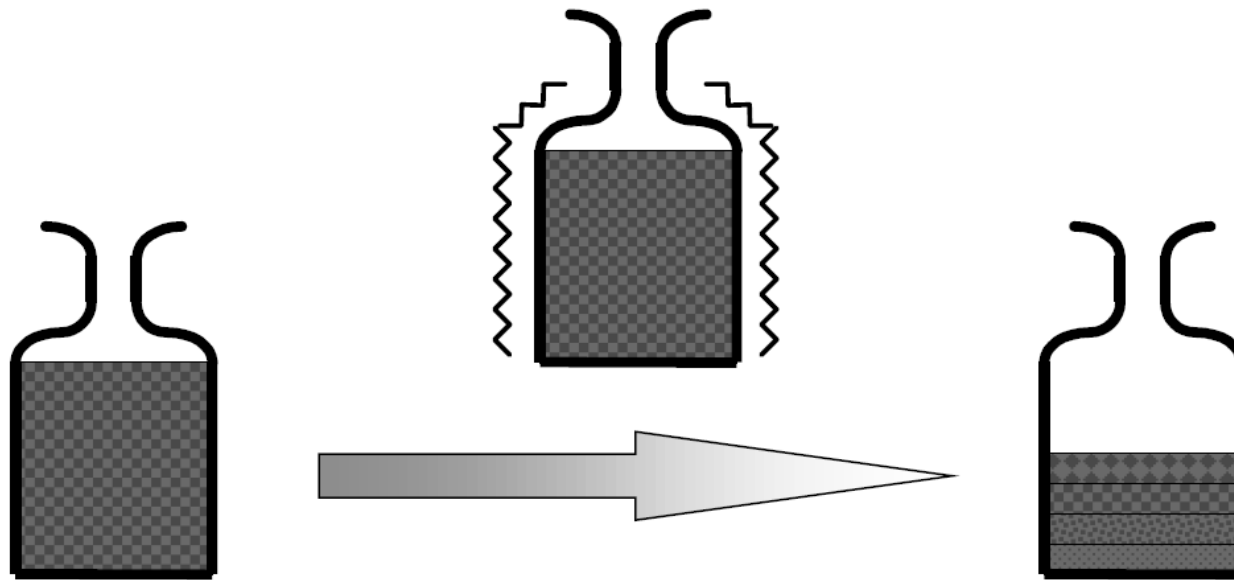
i	pair	$gain(i)$	$\sum gain(i)$	cutsizes
0	-	-	-	5
1	(d, c)	2	2	3
2	(b, g)	0	2	3
3	(a, f)	-1	1	4
4	(e, h)	-1	0	5



- ▶ Start from one of these solutions the whole process again

Simulated Annealing – Underlying Philosophy

- ▶ Inspired from the physical process of annealing (from metallurgy), where a “structured” lattice structure of a solid is achieved by
 1. *heating up* the solid to its melting point
 2. ... and then *slowly cooling down* until it solidifies to a low-energy state



Simulated Annealing – Underlying Philosophy (2)

- ▶ Solids take on a **minimal-energy state** during cooling down *if the temperature is decreased sufficiently slowly*
- ▶ There is a non-zero probability that a particle “jumps” to a higher-energy state ($e_{i+1} > e_i$):

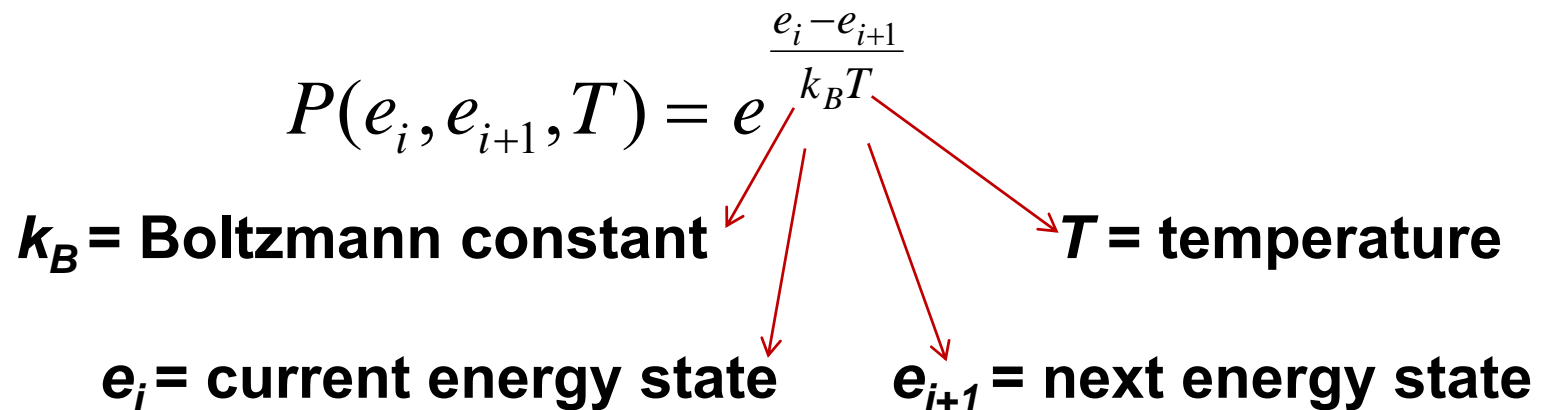
$$P(e_i, e_{i+1}, T) = e^{-\frac{e_i - e_{i+1}}{k_B T}}$$

k_B = Boltzmann constant

T = temperature

e_i = current energy state

e_{i+1} = next energy state

The diagram shows the equation $P(e_i, e_{i+1}, T) = e^{-\frac{e_i - e_{i+1}}{k_B T}}$ with four red arrows pointing from the variables in the equation to their definitions below. The arrows point from k_B to 'Boltzmann constant', from T to 'temperature', from e_i to 'current energy state', and from e_{i+1} to 'next energy state'.

Simulated Annealing Applications

Application to combinatorial optimization:

- ▶ energy = cost of a solution (partition)
- ▶ cost decreases with temperature (a global parameter)
- ▶ increases in cost are accepted with a certain *probability* (that depends both on the *difference between cost values* and also on “*temperature*”)

Simulated Annealing Algorithm

By analogy with the physical process:

- ▶ replace existing solutions by (randomly generated) new feasible solutions from a neighborhood
- ▶ improve a solution by always accepting better-cost neighbors (if selected) but allow for a (*stochastically*) guided acceptance of worse-cost neighbors
- ▶ gradual cooling: gradually decrease the probability of accepting worse-cost solutions
 - ▶ selecting solutions is almost random when T is large
 - ▶ ... but increasingly selects the better cost solution as T goes to **zero**

Advantage

- ▶ allowance for “uphill” moves potentially avoids local optima

Simulated Annealing – Possible Coding

```
temp = temp_start;
cost = c(P);
while (Frozen() == FALSE) {
    while (Equilibrium() == FALSE) {
        P' = RandomMove(P);
        cost' = c(P');
        deltacost = cost' - cost;
        if (Accept(deltacost, temp) > random[0,1)) {
            P = P';
            cost = cost';
        }
    }
    temp = DecreaseTemp(temp);
}
```

initial solution

$\text{Accept}(deltacost, temp) = e^{-\frac{deltacost}{k \cdot temp}}$

Simulated Annealing – Possible Coding (contn.)

▶ **RandomMove (P)**

- choose a random solution in the neighborhood of P

▶ **DecreaseTemp () , Frozen ()**

- cooling down; there are many different choices, for example:
 - initially: `temp:=1.0;`
 - in any iteration: `temp := α *temp` (typ.: $0.8 \leq \alpha \leq 0.99$)
- frozen after a certain time or if there is no further improvement

▶ **Equilibrium ()**

- usually after a defined number of iterations

▶ **Complexity**

- from exponential to constant, depending on the choice of the functions **Equilibrium ()** , **DecreaseTemp ()** , and **Frozen ()**