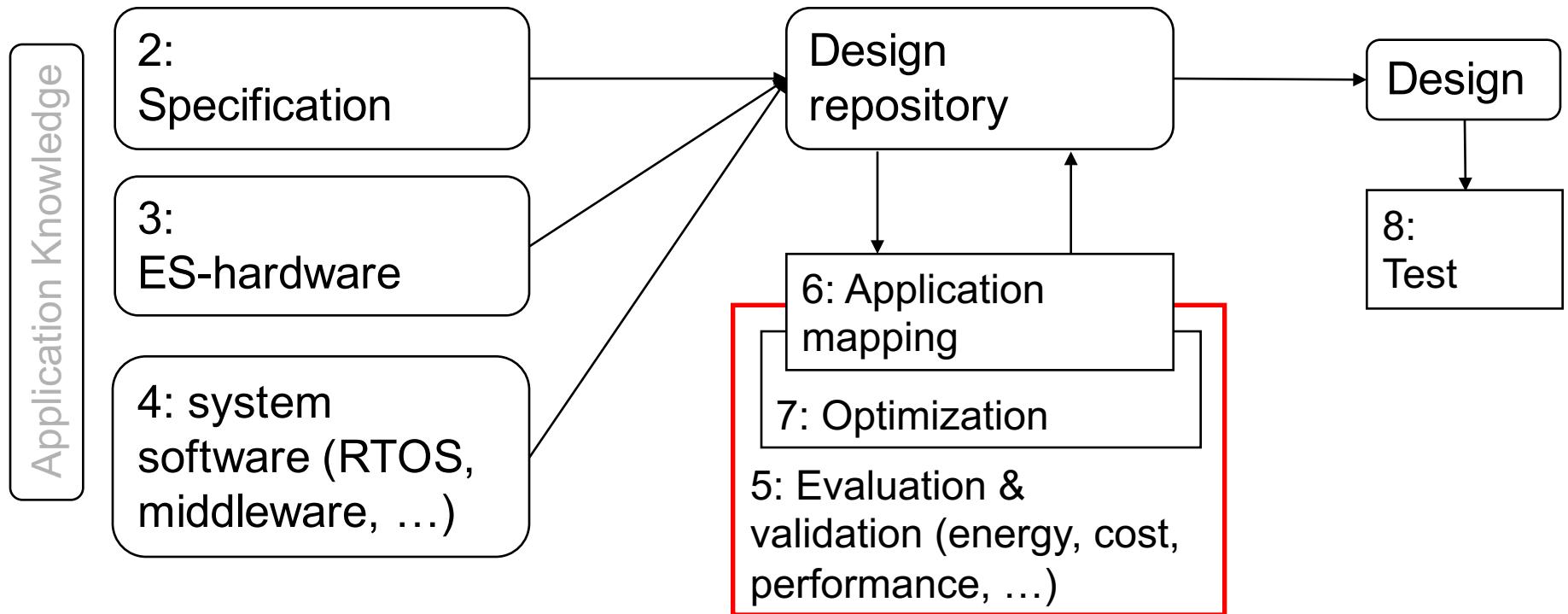


Evaluation and Validation

Jian-Jia Chen
(Slides are based on
Peter Marwedel)
TU Dortmund, Informatik 12
Germany

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Structure of this course



Numbers denote sequence of chapters

Validation and Evaluation

Definition: Validation is the process of checking whether or not a certain (possibly partial) design is appropriate for its purpose, meets all constraints and will perform as expected (yes/no decision).

Definition: Validation with mathematical rigor is called (formal) verification.

Definition: Evaluation is the process of computing quantitative information of some key characteristics of a certain (possibly partial) design.

How to evaluate designs according to multiple criteria?

Many different criteria are relevant for evaluating designs:

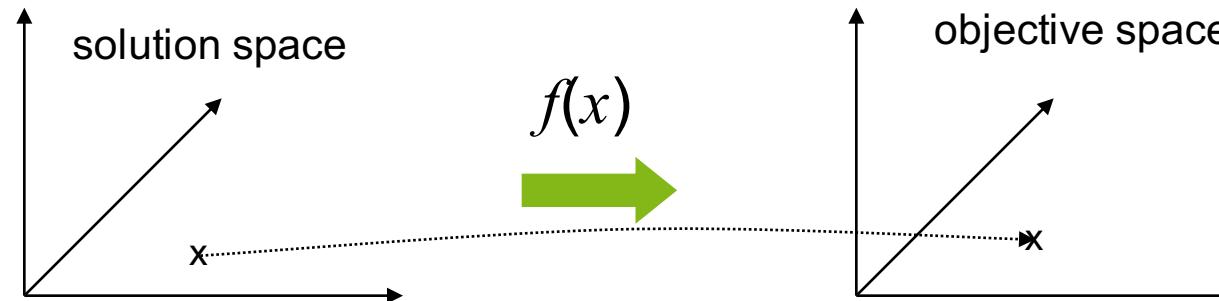
- Average & worst case delay
- power/energy consumption
- thermal behavior
- reliability, safety, security
- cost, size
- weight
- EMC characteristics
- radiation hardness, environmental friendliness, ..



How to compare different designs?
(Some designs are “better” than others)

Definitions

- Let X : m -dimensional **solution space** for the design problem.
Example: dimensions correspond to # of processors, size of memories, type and width of busses etc.
- Let F : n -dimensional **objective space** for the design problem.
Example: dimensions correspond to average and worst case delay, power/energy consumption, size, weight, reliability, ...
- Let $f(x) = (f_1(x), \dots, f_n(x))$ where $x \in X$ be an **objective function**.
We assume that we are using $f(x)$ for evaluating designs.



Pareto points

- We assume that, for each objective, an order $<$ and the corresponding order \leq are defined.
- **Definition:**
Vector $u = (u_1, \dots, u_n) \in F$ **dominates** vector $v = (v_1, \dots, v_n) \in F$
 \Leftrightarrow
 u is “better” than v with respect to *at least* one objective and not worse than v with respect to all other objectives:

$$\forall i \in \{1, \dots, n\} : u_i \leq v_i \wedge$$

$$\exists i \in \{1, \dots, n\} : u_i < v_i$$

- **Definition:**
Vector $u \in F$ is **indifferent** with respect to vector $v \in F$
 \Leftrightarrow neither u dominates v nor v dominates u

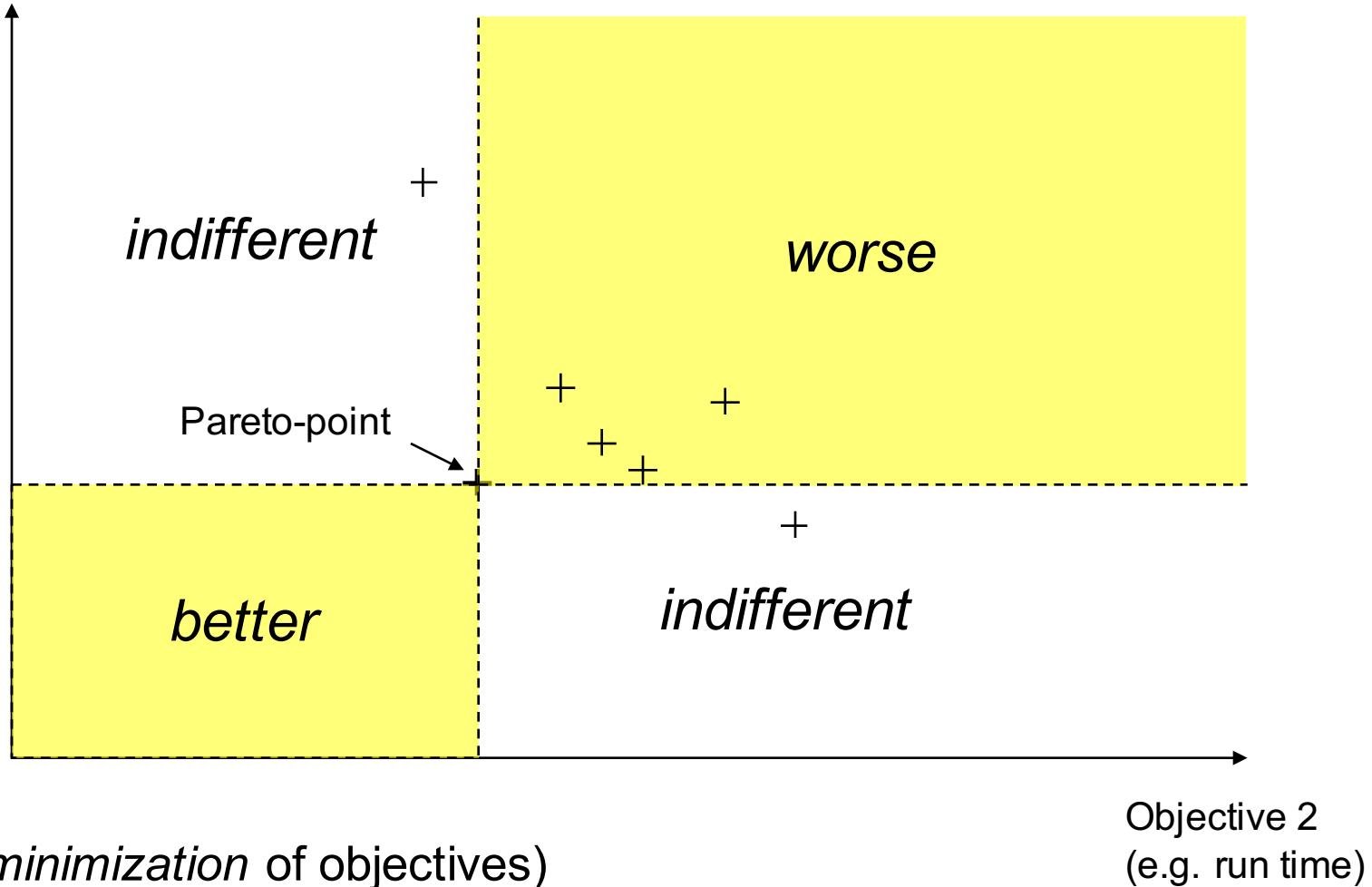
Pareto points

- A solution $x \in X$ is called **Pareto-optimal** with respect to X
 \Leftrightarrow there is no solution $y \in X$ such that $u = f(x)$ is dominated by $v = f(y)$. x is a **Pareto point**.
- **Definition:** Let $S \subseteq F$ be a subset of solutions.
 $v \in F$ is called a **non-dominated solution** with respect to S
 $\Leftrightarrow v$ is not dominated by any element $\in S$.
- v is called **Pareto-optimal**
 $\Leftrightarrow v$ is non-dominated with respect to all solutions F .
- A **Pareto-set** is the set of all Pareto-optimal solutions

Pareto-sets define a **Pareto-front**
(boundary of dominated subspace)

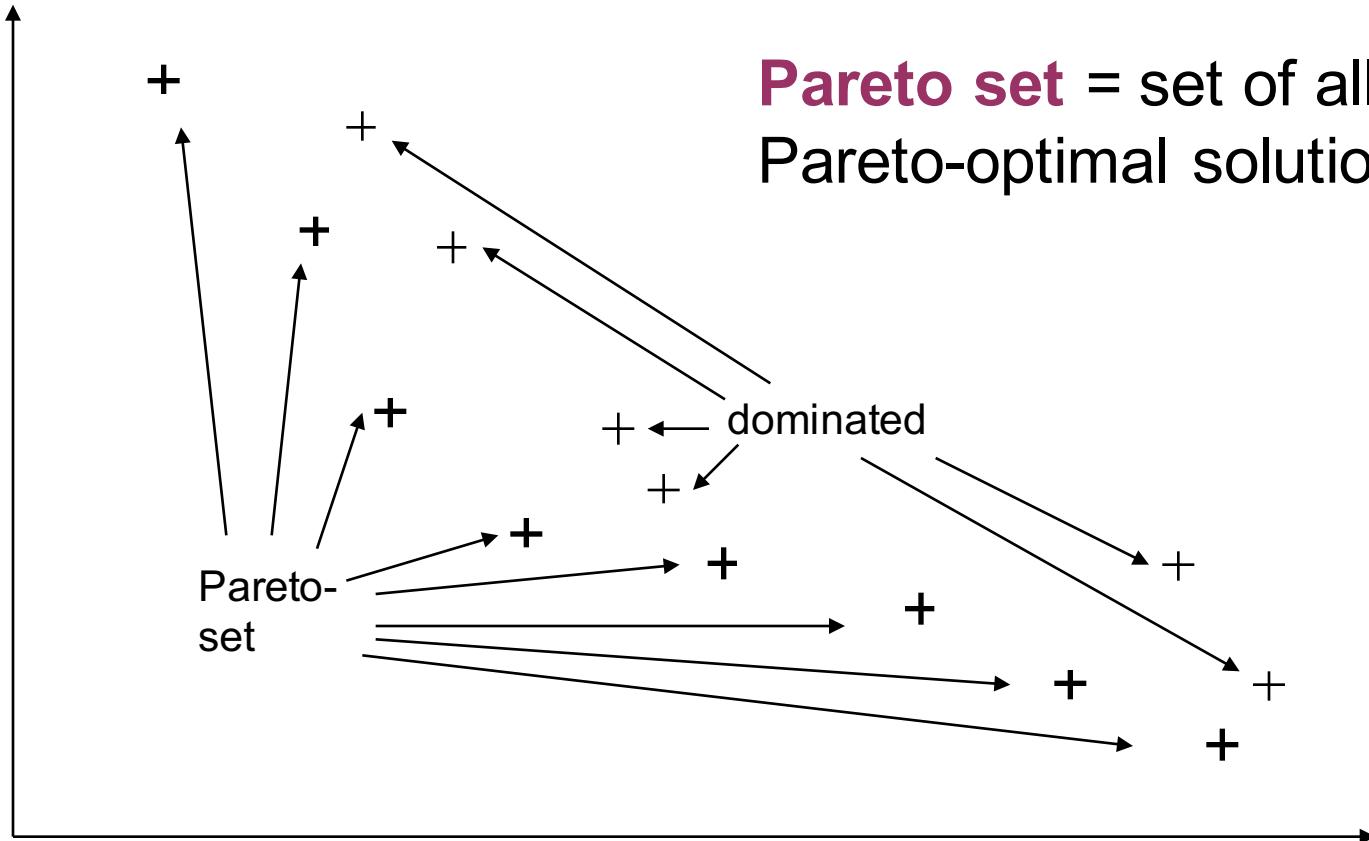
Pareto Point

Objective 1
(e.g. energy consumption)



Pareto Set

Objective 1
(e.g. energy consumption)



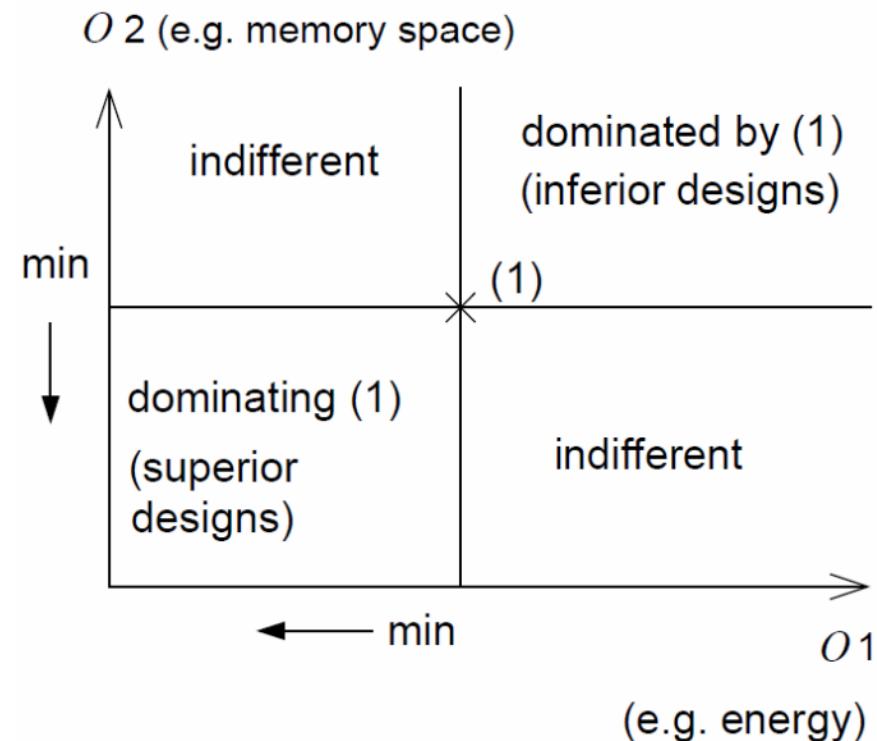
Pareto set = set of all
Pareto-optimal solutions

(Assuming *minimization* of objectives)

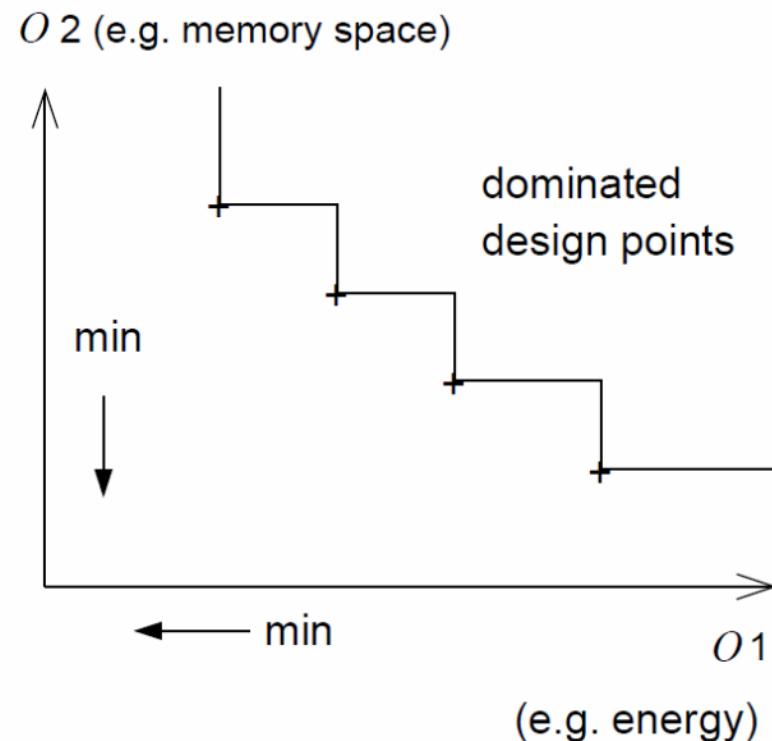
Objectives
(e.g. run time)

One more time ...

Pareto point



Pareto front



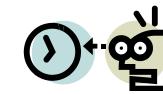
Design space evaluation

Design space evaluation (DSE) based on Pareto-points is the process of finding and returning a set of Pareto-optimal designs to the user, enabling the user to select the most appropriate design.

How to evaluate designs according to multiple criteria?

Many different criteria are relevant for evaluating designs:

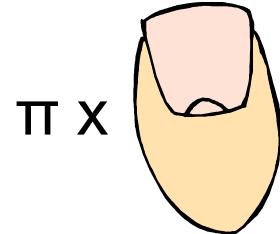
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How to compare different designs?
(Some designs are “better” than others)

Average delays (execution times)

- **Estimated** average execution times :
Difficult to generate sufficiently precise estimates;
Balance between run-time and precision



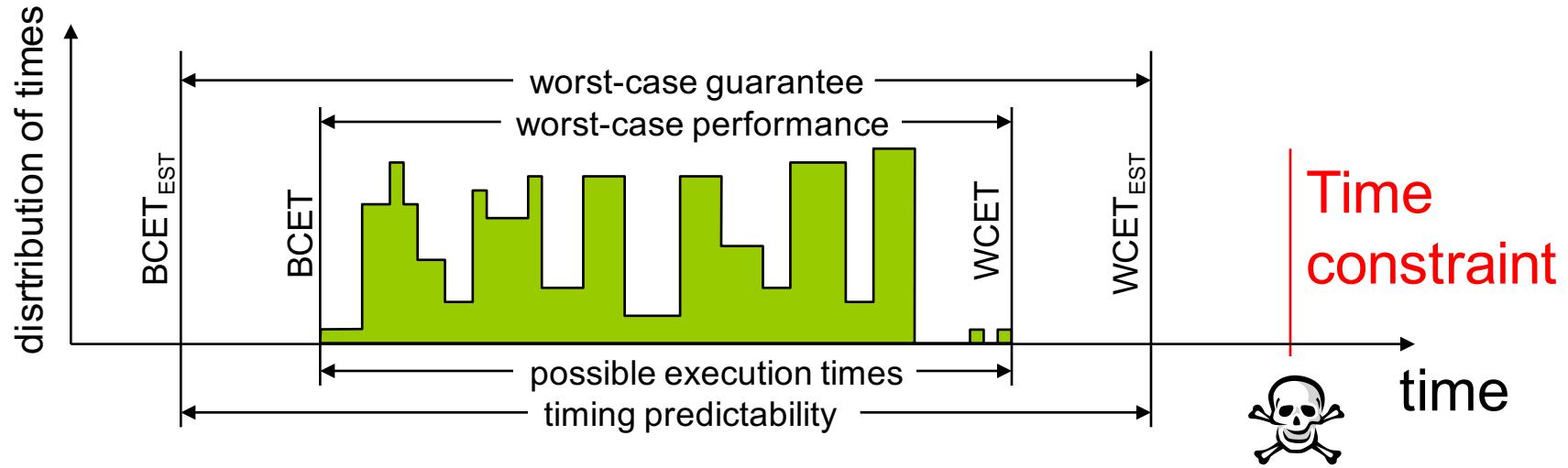
- **Accurate** average execution times:
As precise as the input data is.



We need to compute **average** and **worst case** execution times

Worst case execution time (1)

Definition of worst case execution time:



WCET_{EST} must be

1. safe (i.e. \geq WCET) and
2. tight ($WCET_{EST} - WCET \ll WCET_{EST}$)

Worst case execution times (2)

Complexity:

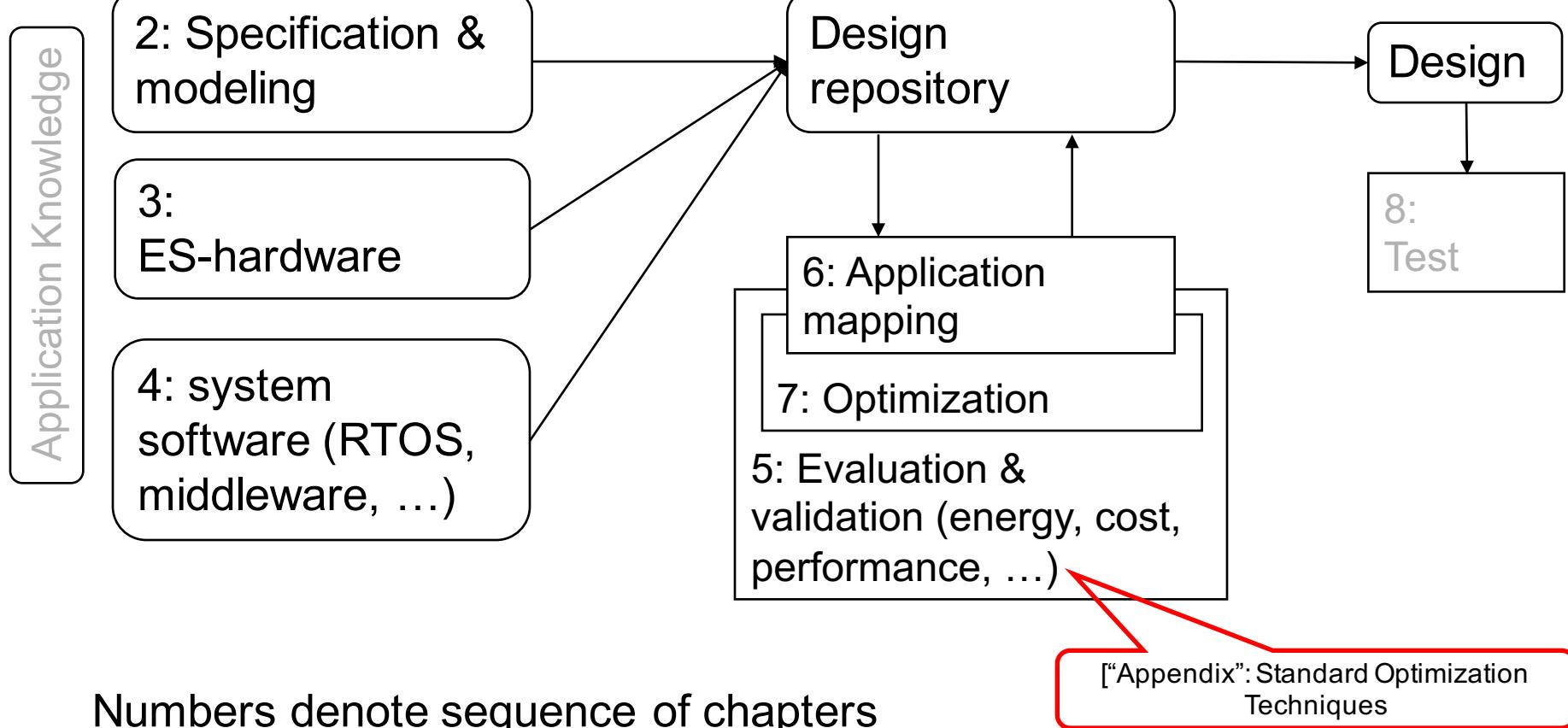
- in the general case: undecidable if a bound exists.
- for restricted programs: simple for “old“ architectures, very complex for new architectures with pipelines, caches, interrupts, virtual memory, etc.



Approaches:

- for hardware: requires detailed timing behavior
- for software: requires availability of machine programs; complex analysis (see, e.g., www.absint.de)

Structure of this course



Integer linear programming models

Ingredients:

- Cost function
 - Constraints
- } Involving linear expressions of integer variables from a set X

Cost function

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

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

Def.: The problem of minimizing (1) subject to the constraints (2) is called an **integer linear programming (ILP) problem**.

If all x_i are constrained to be either 0 or 1, the ILP problem is said to be a **0/1 integer linear programming problem**.

Example

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

$$x_1 + x_2 + x_3 \geq 2$$

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

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

Remarks on integer programming

- Maximizing the cost function: just set $C'=-C$
- Integer programming is NP-complete.
- Running times depend exponentially on problem size, but problems of >1000 vars solvable with good solver (depending on the size and structure of the problem)
- The case of $x_i \in \mathbb{R}$ is called *linear programming* (LP). Polynomial complexity, but most algorithms are exponential, in practice still faster than for ILP problems.
- The case of some $x_i \in \mathbb{R}$ and some $x_i \in \mathbb{N}$ is called *mixed integer-linear programming*.
- ILP/LP models good starting point for modeling, even if heuristics are used in the end.
- Solvers: lp_solve (public), CPLEX (commercial), ...

An Example: Knapsack Problem

Example IP formulation:

The Knapsack problem:

I wish to select items to put in my backpack.

There are m items available.

Item i weights w_i kg,

Item i has value v_i .

I can carry Q kg.

Let $x_i = \begin{cases} 1 & \text{if I select item } i \\ 0 & \text{otherwise} \end{cases}$

$$\max \quad \sum_i x_i v_i$$

$$\text{s.t.} \quad \sum_i x_i w_i \leq Q$$

$$x_i \in \{0,1\}, \forall i$$

Variance of Knapsack Problem

- **Given** a set of periodic tasks with implicit deadlines
 - Task τ_i : period T_i ,
 - Options: Execution without/with scratchpad memory (SPM)
 - Without SPM: Worst-case execution time $C_{i,1}$
 - With SPM: required m_i scratchpad memory size and Worst-case execution time $C_{i,2}$
 - Utilization without SPM $U_{i,1} = C_{i,1}/T_i$
 - Utilization with SPM is $U_{i,2} = C_{i,2}/T_i$
 - **Objective**
 - Select the tasks to be put into the SPM
 - Minimize the required SPM size
 - The utilization of the task set should be no more than 100%
- $$\begin{aligned} & \min \quad \sum_i x_i m_i \\ \text{s.t.} \quad & \sum_i x_i U_{i,2} + \sum_i (1 - x_i) U_{i,1} \leq 1 \\ & x_i \in \{0,1\}, \quad \forall i \end{aligned}$$

Summary

Integer (linear) programming

- Integer programming is NP-complete
- Linear programming is faster
- Good starting point even if solutions are generated with different techniques

Simulated annealing

- Modeled after cooling of liquids
- Overcomes local minima

Evolutionary algorithms

- Maintain set of solutions
- Include selection, mutation and recombination

Evolutionary Algorithms (1)

- *Evolutionary Algorithms are based on the collective learning process within a population of individuals, each of which represents a search point in the space of potential solutions to a given problem.*
- *The population is arbitrarily initialized, and it evolves towards better and better regions of the search space by means of randomized processes of*
 - ***selection*** (which is deterministic in some algorithms),
 - ***mutation***, and
 - ***recombination*** (which is completely omitted in some algorithmic realizations).

[Bäck, Schwefel, 1993]

Evolutionary Algorithms (2)

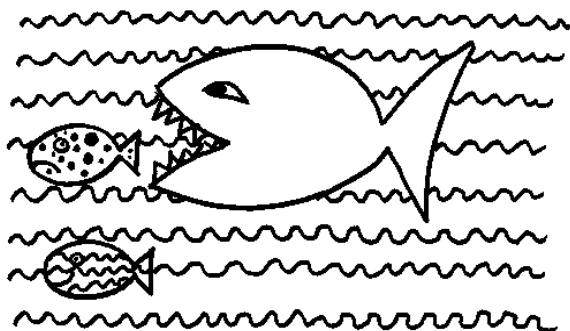
- *The environment (given aim of the search) delivers a quality information (**fitness value**) of the search points, and the selection process favours those individuals of higher fitness to reproduce more often than worse individuals.*
- *The recombination mechanism allows the mixing of parental information while passing it to their descendants, and mutation introduces innovation into the population*

[Bäck, Schwefel, 1993]

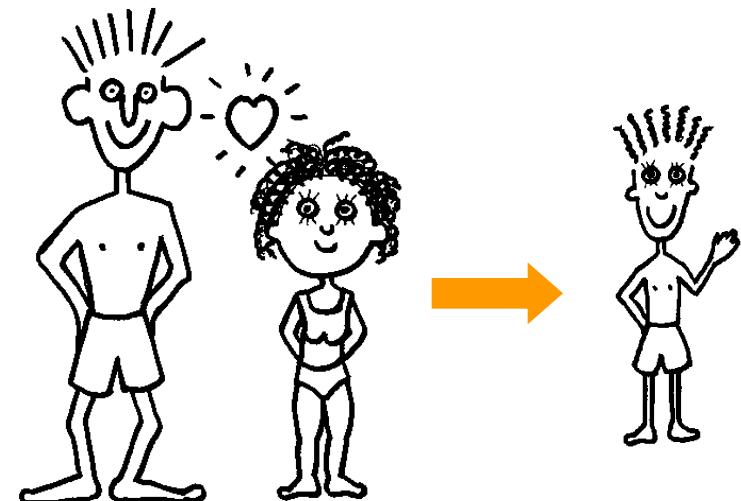
Evolutionary Algorithms

Principles of Evolution

① Selection



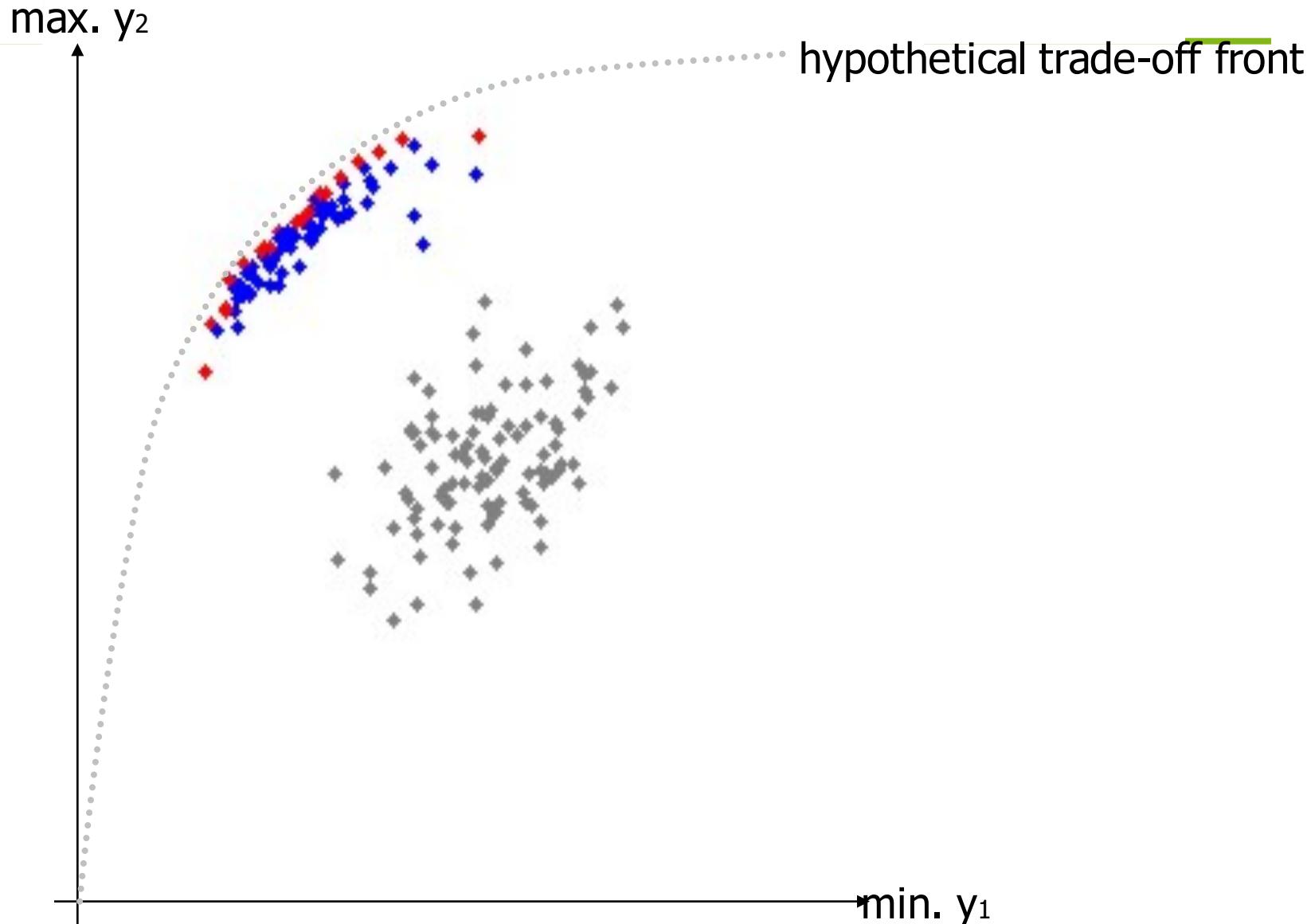
③ Cross-over



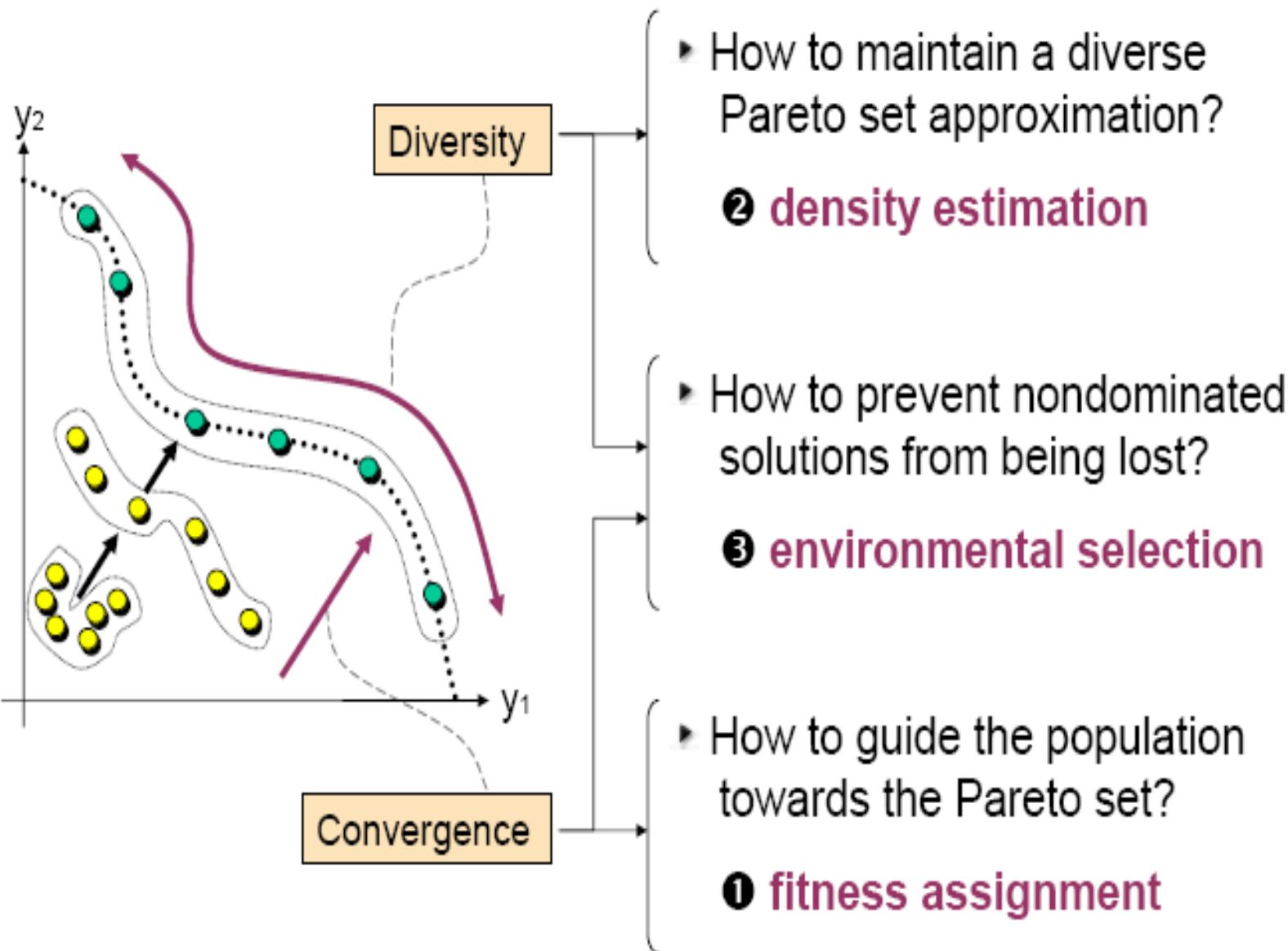
② Mutation



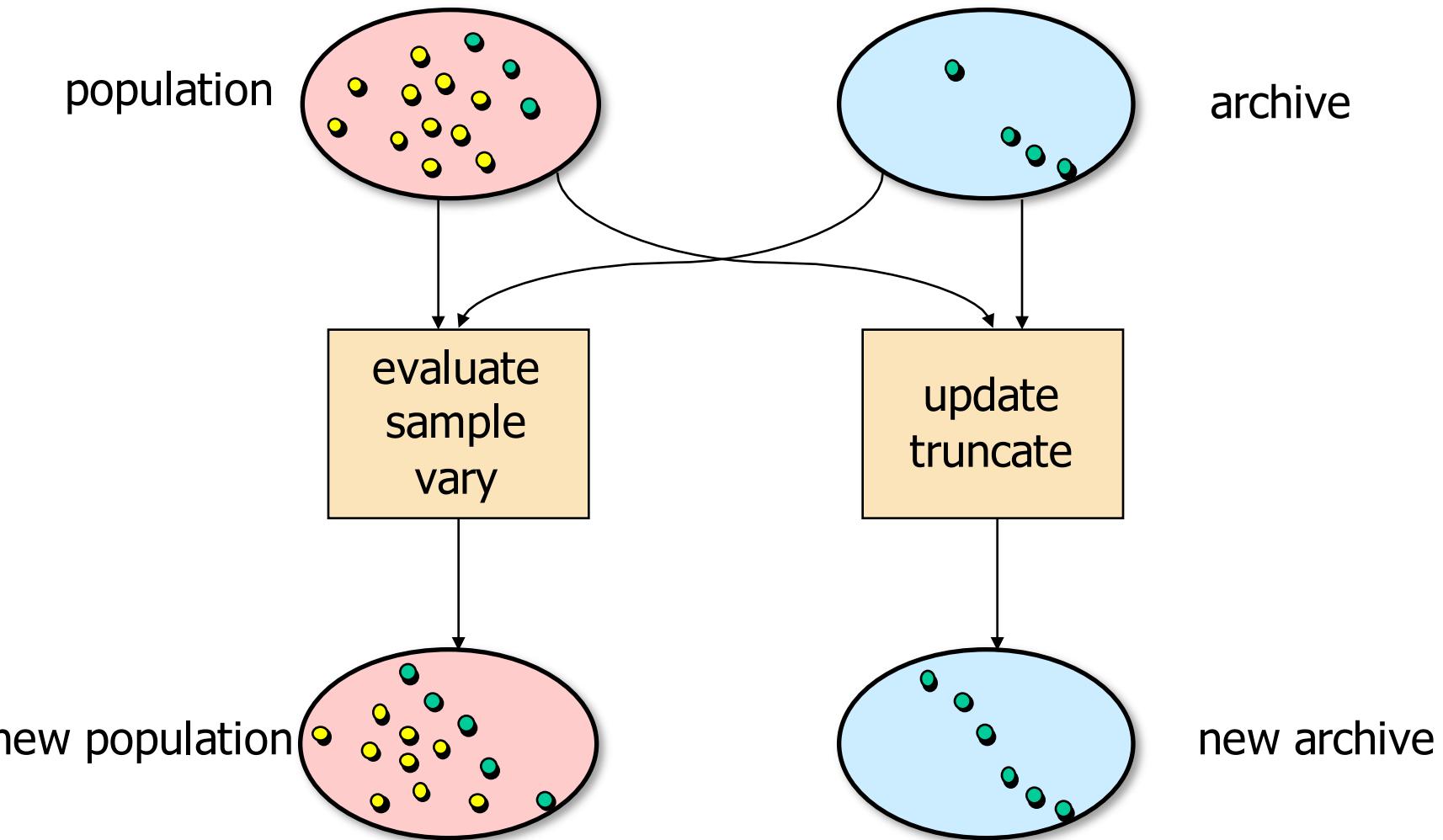
An Evolutionary Algorithm in Action



Issues in Multi-Objective Optimization



A Generic Multiobjective EA



Example: SPEA2 Algorithm

- | | |
|---------|---|
| Step 1: | Generate initial population P_0 and empty archive (external set) A_0 . Set $t = 0$. |
| Step 2: | Calculate fitness values of individuals in P_t and A_t . |
| Step 3: | A_{t+1} = nondominated individuals in P_t and A_t .
If size of $A_{t+1} > N$ then reduce A_{t+1} , else if
size of $A_{t+1} < N$ then fill A_{t+1} with dominated
individuals in P_t and A_t . |
| Step 4: | If $t > T$ then output the nondominated set of A_{t+1} .
Stop. |
| Step 5: | Fill mating pool by binary tournament selection. |
| Step 6: | Apply recombination and mutation operators to
the mating pool and set P_{t+1} to the resulting
population. Set $t = t + 1$ and go to Step 2. |

Simulated Annealing

- General method for solving combinatorial optimization problems.
- Based the model of slowly cooling crystal liquids.
- Some configuration is subject to changes.
- Special property of Simulated annealing: Changes leading to a poorer configuration (with respect to some cost function) are accepted with a certain probability.
- This probability is controlled by a temperature parameter: the probability is smaller for smaller temperatures.

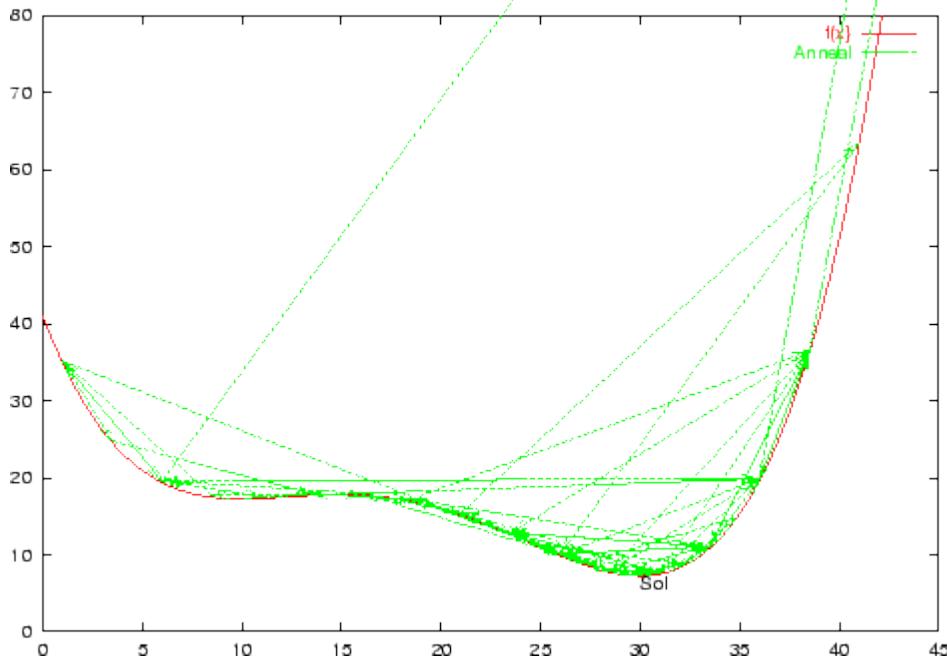
Simulated Annealing Algorithm

```
procedure SimulatedAnnealing;  
var i, T: integer;  
begin  
    i := 0; T := MaxT;  
    configuration:= <some initial configuration>;  
    while not terminate(i, T) do  
        begin  
            while InnerLoop do  
                begin NewConfig := variation(configuration);  
                    delta := evaluation(NewConfig,configuration);  
                    if delta < 0  
                        then configuration := NewConfig;  
                    else if SmallEnough(delta, T, random(0,1))  
                        then configuration := Newconfig;  
                end;  
                T:= NewT(i,T); i:=i+1;  
        end; end;
```

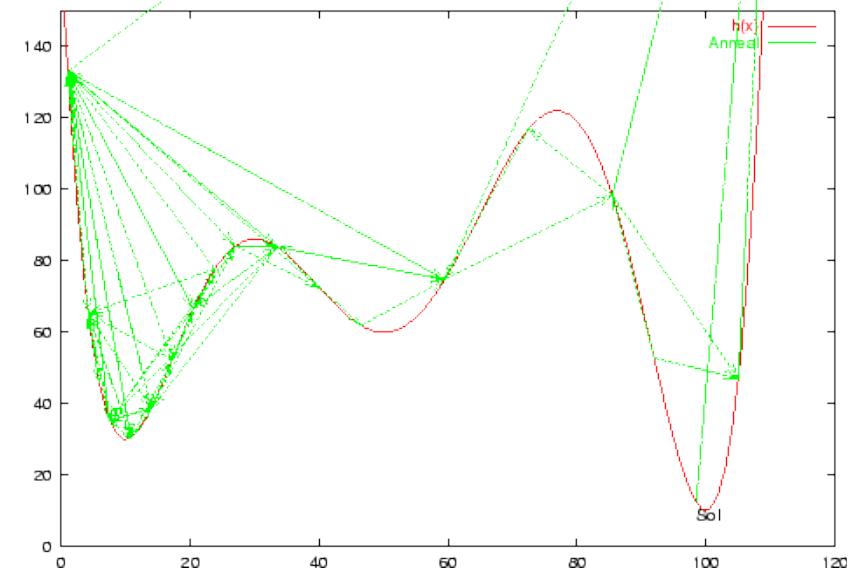
Explanation

- Initially, some random initial configuration is created.
- Current temperature is set to a large value.
- Outer loop:
 - Temperature is reduced for each iteration
 - Terminated if ($\text{temperature} \leq \text{lower limit}$) or ($\text{number of iterations} \geq \text{upper limit}$).
- Inner loop: For each iteration:
 - New configuration generated from current configuration
 - Accepted if ($\text{new cost} \leq \text{cost of current configuration}$)
 - Accepted with temperature-dependent probability if ($\text{cost of new config.} > \text{cost of current configuration}$).

Behavior for actual functions



130 steps



200 steps

[people.equars.com/~marco/poli/phd/node57.html]

<http://foghorn.cadlab.lafayette.edu/cadapplets/fp/fpIntro.html>

Performance

- This class of algorithms has been shown to outperform others in certain cases [Wegener, 2005].
- Demonstrated its excellent results in the TimberWolf layout generation package [Sechen]
- Many other applications ...

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