

# **Computational Intelligence**

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Note: Slides of Thomas Jansen used with permission!

# Important Parameters of EAs (1)

- ullet dimension n of search space
  - no parameter of EA, but given by the problem
  - ullet measures the size of the search space:  $\{0,1\}^n$ ,  $\mathbb{R}^n$ ,  $S_n$
  - plays the same role as input length in classical runtime analysis
  - other parameters are often chosen dependent on n (e.g. mutation probability  $p_m=1/n$ )
- population size  $\mu$ 
  - obviously  $\mu = n^{O(1)}$
  - often  $\mu = \Theta(n)$  or  $\mu = \Theta(\sqrt{n})$
  - $\mu = O(1)$  or even  $\mu = 1$  are not unusual
- ullet number of offspring  $\lambda$ 
  - obviously  $\lambda = n^{O(1)}$
  - often  $\lambda = 1$
  - $\lambda = \mu$  or  $\lambda \gg \mu$  not unusual
  - selection method influences reasonable choice of  $\lambda$

# Important Parameters of EAs (2)

- ullet crossover probability  $p_c$ 
  - in general  $p_c \in [0;1]$  arbitrary
  - often  $p_c \in [1/2; 4/5]$  constant
- probability of applying mutation
  - don't confuse with mutation probability!
  - we will always use 1
  - Remark

$$p_m = 1/n \Rightarrow \mathsf{Prob} \, (\mathsf{no} \, \mathsf{mutation}) = (1 - 1/n)^n \approx 1/e$$

### Methods for parameter control

- static parameter control
  - parameter values constant during the whole run
  - often used
  - + simple
  - maybe it's better to vary the parameter value during the run?!
- dynamic parameter control
  - parameter values change during the run according to some time-dependent scheme
  - + more flexible than static approach
  - cannot deal with non-time-dependent changes
  - unusual for EAs
- adaptive parameter control
  - parameter values can change dependently on every individual and any random experiment
  - + very flexible
  - hard to analyze
  - computationally expensive
  - often used for EAs

# Self-adaptation

Idea good parameter values evolve together with good individuals

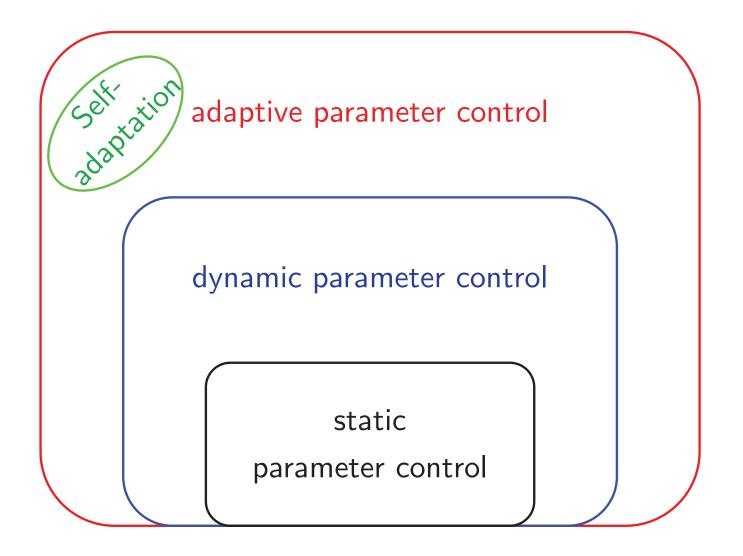
implementation code parameter values together with individual

formally  $S\times Q \text{ instead of } S$   $\text{unchanged } f\colon S\to R$ 

### e.g. for mutation probability

- every individual has its own mutation probability
- first vary the mutation probability
- then mutate with varied mutation probability
- afterwards normal selection
- important don't swap steps

# Hierarchy of parameter control methods



#### **Evolutionary Algorithms: Historical Notes**

#### Lecture 11

Idea emerged independently several times: about late 1950s / early 1960s.

Three branches / "schools" still active today.

#### Evolutionary Programming (EP):

Pioneers: Lawrence Fogel, Alvin Owen, Michael Walsh (New York, USA).

Original goal: Generate intelligent behavior through simulated evolution.

Approach: Evolution of finite state machines predicting symbols.

Later (~1990s) specialized to optimization in  $\mathbb{R}^n$  by David B. Fogel.

#### Genetic Algorithms (GA):

Pioneer: John Holland (Ann Arbor, MI, USA).

Original goal: Analysis of adaptive behavior.

Approach: Viewing evolution as adaptation. Simulated evolution of bit strings.

Applied to optimization tasks by PhD students (Kenneth de Jong, 1975; et al.).

#### Evolution Strategies (ES):

Pioneers: Ingo Rechenberg, Hans-Paul Schwefel, Peter Bienert (Berlin, Germany).

Original goal: Optimization of complex systems.

Approach: Viewing variation/selection as improvement strategy. First in  $\mathbb{Z}^n$ , then  $\mathbb{R}^n$ .



"Offspring" from GA branch:

#### Genetic Programming (GP):

Pioneers: Nichael Lynn Cramer 1985, then: John Koza (Stanford, USA).

Original goal: Evolve programs (parse trees) that must accomplish certain task.

Approach: GA mechanism transfered to parse trees.

Later: Programs as successive statements → Linear GP (e.g. Wolfgang Banzhaf)

#### Already beginning early 1990s:

Borders between EP, GA, ES, GP begin to blurr ...

- ⇒ common term **Evolutionary Algorithm** embracing all kind of approaches
- ⇒ broadly accepted name for the field: **Evolutionary Computation**

scientific journals: *Evolutionary Computation* (MIT Press) since 1993, *IEEE Transactions on Evolutionary Computation* since 1997, several more specialized journals started since then.



### Design of EAs

Idea Methodology to apply standard EAs

Goal standard EAs do not have to be changed

Requirement problem is given as  $g: A \rightarrow B$ 

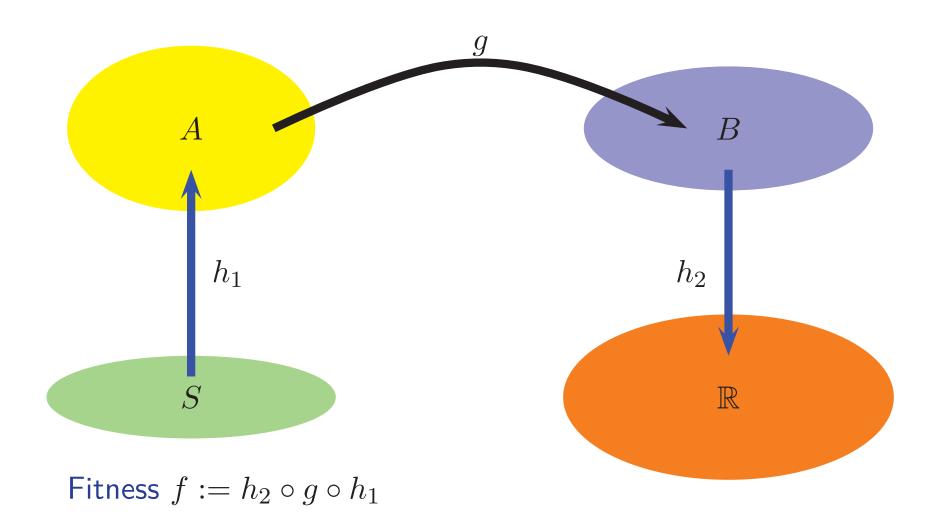
g has to be maximized (or minimized)

A arbitrary set, B partially ordered

 $\mathsf{EA}$  operates on search space S

'maximizes' fitness  $f:S \to \mathbb{R}$ 

# Definition of mappings



 $h_1$  is genotype-phenotype-mapping.

### **Genotype-Phenotype-Mapping** $\mathbb{B}^n \to [L, R] \subset \mathbb{R}$

• Standard encoding for  $b \in \mathbb{B}^n$ 

$$x = L + \frac{R - L}{2^n - 1} \sum_{i=0}^{n-1} b_{n-i} 2^i$$

→ Problem: *hamming cliffs* 

000	001	010	011	100	101	110	111			
0	1	2	3	4	5	6	7			
1 Bit 2 Bit 1 Bit 3 Bit 1 Bit 2 Bit 1 Bit										
1 B	Bit 2 E	Bit 1 E	3It 3	<b>†</b>	Bit 2	Bit 1	Bit			

$$L = 0, R = 7$$
  
 $n = 3$ 

### **Genotype-Phenotype-Mapping** $\mathbb{B}^n \to [L, R] \subset \mathbb{R}$

• Gray encoding for  $b \in \mathbb{B}^n$ 

Let 
$$a \in \mathbb{B}^n$$
 standard encoded. Then  $b_i = \begin{cases} a_i, & \text{if } i = 1 \\ a_{i-1} \oplus a_i, & \text{if } i > 1 \end{cases}$ 

000	001	011	010	110	111	101	100	← genotype
0	1	2	3	4	5	6	7	← phenotype

OK, no hamming cliffs any longer ...

- ⇒ small changes in phenotype "lead to" small changes in genotype
- since we consider evolution in terms of Darwin (not Lamarck):
- ⇒ small changes in genotype lead to small changes in phenotype!

**but:** 1-Bit-change:  $000 \rightarrow 100 \Rightarrow \odot$ 



**Genotype-Phenotype-Mapping**  $\mathbb{B}^n \to \mathbb{P}^n$  (example only)

ullet e.g. standard encoding for  $b \in \mathbb{B}^n$ 

#### individual:

010	101	111	000	110	001	101	100	← genotype
0	1	2	3	4	5	6	7	← index

consider index and associated genotype entry as unit / record / struct; sort units with respect to genotype value, old indices yield permutation:

000	001	010	100	101	101	110	111	← genotype
3	5	0	7	1	6	4	2	← old index

= permutation

# Requirements on $h_1$ and $h_2$

### obvious requirements

- $h_1$  and  $h_2$  can be computed efficiently
- $h_2$  suits g, i. e. good points in B are mapped to good points in  $\mathbb R$
- $h_1$  maps on many (all) important points of A
- ullet Optima of f correspond to optima of g

Caution requirements can be hard to achieve in practice

for non-obvious requirements a metric is important

### **Definition**

Mapping  $d: M \times M \to \mathbb{R}_0^+$  is a metric on the set  $M:\Leftrightarrow$ 

- 2  $\forall x, y \in M : d(x, y) = d(y, x)$  (symmetry)
- 3  $\forall x, y, z \in M : d(x,y) + d(y,z) \ge d(x,z)$  (triangle inequality)

### Metric-based EAs

Assumption Metric  $d_A$  on A known  $(d_A \text{ reflects application knowledge})$ 

Requirement metric  $d_S$  is known

if  $h_1$  injective,  $d_S(x,x') := d_A(h_1(x),h_1(x'))$  is metric

Requirement monotonicity

$$\forall x, x', x'' \in S: \qquad d_S(x, x') \leq d_S(x, x'')$$

$$\Rightarrow d_A(h_1(x), h_1(x')) \leq d_A(h_1(x), h_1(x''))$$

# Variation as randomized mapping

now Design-rules for variation operators

hence Formalize variation operators as randomized mappings

 $r \colon X \to Y$  randomized mapping

 $\Leftrightarrow r(x) \in Y$  depends on  $x \in X$  and random experiment

formally probability space  $(\Omega, p)$ 

$$r: X \times \Omega \to Y$$

$$\operatorname{Prob}\left(r(x)=y\right)=\sum_{\omega\in\Omega\colon r(x,\omega)=y}p(\omega)$$

Example 1-bit mutation

$$\Omega := \{1, 2, \dots, n\}, \ \forall i \in \Omega \colon p(i) = 1/n$$

1-bit mutation is randomized mapping  $m\colon\{0,1\}^n\to\{0,1\}^n$  where  $m(x,i):=x\oplus 0^{i-1}10^{n-i}$ 

### Design-rules for mutation

### favor small changes

$$\forall x, x', x'' \in S$$
:  $d_S(x, x') < d_S(x, x'')$   
 $\Rightarrow \operatorname{Prob}(m(x) = x') > \operatorname{Prob}(m(x) = x'')$ 

#### no bias

$$\forall x, x', x'' \in S$$
:  $d_S(x, x') = d_S(x, x'')$   
 $\Rightarrow \operatorname{Prob}(m(x) = x') = \operatorname{Prob}(m(x) = x'')$ 

# Design-rules for crossover offspring similar to parents

$$\forall x, x', x'' \in S: \qquad \text{Prob}\left(c(x, x') = x''\right) > 0$$
  
$$\Rightarrow \quad \max\left\{d_S(x, x''), d_S(x', x'')\right\} \leq d_S(x, x')$$

no bias

$$\forall x, x' \in S : \forall \alpha \in \mathbb{R}_0^+ :$$

$$\mathsf{Prob}\left(d_S(x, c(x, x')) = \alpha\right) = \mathsf{Prob}\left(d_S(x', c(x, x')) = \alpha\right)$$

Any EA that fulfills these four design-rules is called a metric-based EA (MBEA).