Computational Intelligence

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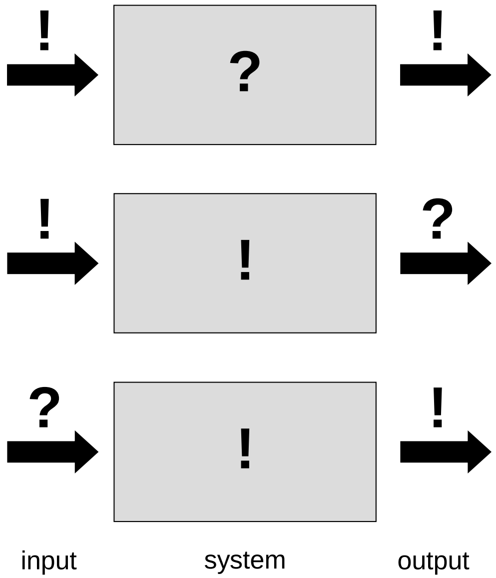
Prof. Dr. Günter Rudolph
Lehrstuhl für Algorithm Engineering (LS 11)
Fakultät für Informatik
TU Dortmund

Plan for Today

- Evolutionary Algorithms (EA)
  - Optimization Basics
  - EA Basics

Optimization Basics



modelling

simulation

optimization

Optimization Basics

given:

objective function f: X → $\mathbb{R}$

feasible region X (= nonempty set)

objective: find solution with *minimal* or *maximal* value!

optimization problem:

find x* ∈ X such that f(x*) = min{ f(x) : x ∈ X }

x*	**global solution**
f(x*)	**global optimum**

note:

max{ f(x) : x ∈ X } = – min{ – f(x) : x ∈ X }

Optimization Basics

Lecture 06

local solution $x^* \in X$:

$\forall x \in N(x^*): f(x^*) \leq f(x)$

if x^* local solution then

$f(x^*)$ **local optimum / minimum**

neighborhood of x^* =
bounded subset of X

example: $X = \mathbb{R}^n$, $N_\varepsilon(x^*) = \{ x \in X: || x - x^*||_2 \leq \varepsilon \}$

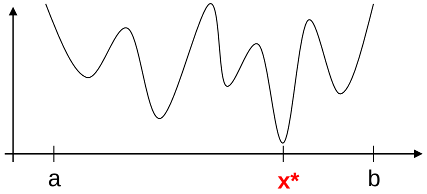
remark:

evidently, every global solution / optimum is also local solution / optimum;

the reverse is wrong in general!

example:

f: [a,b] → $\mathbb{R}$, global solution at **x***



Optimization Basics

Lecture 06

What makes optimization difficult?

some causes:

- local optima (is it a global optimum or not?)
- constraints (ill-shaped feasible region)
- non-smoothness (weak causality) → strong causality needed!
- discontinuities (⇒ nondifferentiability, no gradients)
- lack of knowledge about problem (⇒ black / gray box optimization)

$f(x) = a_1 x_1 + ... + a_n x_n \rightarrow \max!$ with $x_i \in \{0,1\}$, $a_i \in \mathbb{R}$ ⇒ $x_i^* = 1$ iff $a_i > 0$

add constaint $g(x) = b_1 x_1 + ... + b_n x_n \leq b$ ⇒ NP-hard

add capacity constraint to TSP ⇒ CVRP ⇒ still harder

Optimization Basics

Lecture 06

When using which optimization method?

mathematical algorithms	randomized search heuristics
• problem explicitly specified	• problem given by black / gray box
• problem-specific solver available	• no problem-specific solver available
• problem well understood	• problem poorly understood
• ressources for designing algorithm affordable	• insufficient ressources for designing algorithm
• solution with proven quality required	• solution with satisfactory quality sufficient
⇒ **don't apply EAs**	⇒ **EAs worth a try**

Evolutionary Algorithm Basics

Lecture 06

idea: using **biological evolution** as **metaphor** and as **pool of inspiration**

⇒ interpretation of biological evolution as iterative method of improvement

feasible solution $x \in X = S_1 \times ... \times S_n$ = chromosome of **individual**

multiset of feasible solutions = **population**: multiset of individuals

objective function $f: X \rightarrow \mathbb{R}$ = **fitness function**

often: $X = \mathbb{R}^n$, $X = \mathbb{B}^n = \{0,1\}^n$, $X = \mathbb{P}_n = \{ \pi : \pi$ is permutation of $\{1,2,...,n\} \}$

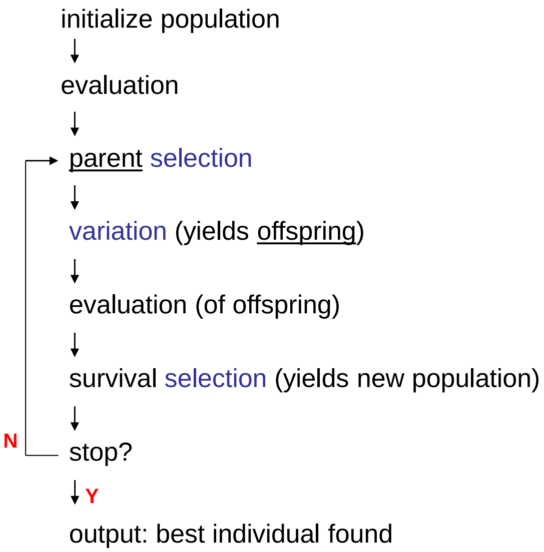
also : combinations like $X = \mathbb{R}^n \times \mathbb{B}^p \times \mathbb{P}_q$ or non-cartesian sets

⇒ structure of feasible region / search space defines **representation** of individual

Evolutionary Algorithm Basics

Lecture 06

algorithmic skeleton



initialize population

evaluation

parent selection

variation (yields offspring)

evaluation (of offspring)

survival selection (yields new population)

stop?

N

Y

output: best individual found

Evolutionary Algorithm Basics

Lecture 06

Specific example: (1+1)-EA in $\mathbb{B}^n$ for minimizing some $f: \mathbb{B}^n \to \mathbb{R}$

population size = 1, number of offspring = 1, selects best from $\overset{\uparrow}{1}+\overset{\uparrow}{1}$ individuals

parent offspring

1. initialize $X^{(0)} \in \mathbb{B}^n$ uniformly at random, set $t = 0$
2. evaluate $f(X^{(t)})$
3. select parent: $Y = X^{(t)}$ ⟶ no choice, here
4. variation: flip each bit of Y independently with probability $p_m = 1/n$
5. evaluate $f(Y)$
6. selection: if $f(Y) \le f(X^{(t)})$ then $X^{(t+1)} = Y$ else $X^{(t+1)} = X^{(t)}$
7. if not stopping then $t = t+1$, continue at (3)

Evolutionary Algorithm Basics

Lecture 06

Specific example: (1+1)-EA in $\mathbb{R}^n$ for minimizing some $f: \mathbb{R}^n \to \mathbb{R}$

population size = 1, number of offspring = 1, selects best from $\overset{\uparrow}{1}+\overset{\uparrow}{1}$ individuals

parent offspring

compact set = closed & bounded

1. initialize $X^{(0)} \in C \subset \mathbb{R}^n$ uniformly at random, set $t = 0$
2. evaluate $f(X^{(t)})$
3. select parent: $Y = X^{(t)}$ ⟶ no choice, here
4. variation = add random vector: $Y = Y + Z$, e.g. $Z \sim N(0, I_n)$
5. evaluate $f(Y)$
6. selection: if $f(Y) \le f(X^{(t)})$ then $X^{(t+1)} = Y$ else $X^{(t+1)} = X^{(t)}$
7. if not stopping then $t = t+1$, continue at (3)

Evolutionary Algorithm Basics

Lecture 06

Selection

(a) select parents that generate offspring → selection for **reproduction**

(b) select individuals that proceed to next generation → selection for **survival**

necessary requirements:

- selection steps must not favor worse individuals
- one selection step may be neutral (e.g. select uniformly at random)
- at least one selection step must favor better individuals

typically : selection only based on fitness values f(x) of individuals

seldom : additionally based on individuals‘ chromosomes x (→ maintain diversity)

Selection methods

population P = (x_1, x_2, ..., x_μ) with μ individuals

two approaches:

1. repeatedly select individuals from population with replacement
2. rank individuals somehow and choose those with best ranks (no replacement)

- ***uniform / neutral selection***
  choose index i with probability $1/\mu$
- ***fitness-proportional selection***
  choose index i with probability $s_i = \dfrac{f(x_i)}{\sum\limits_{x \in P} f(x)}$

problems: f(x) > 0 for all $x \in X$ required ⇒ g(x) = exp(f(x)) > 0

but already sensitive to additive shifts g(x) = f(x) + c

almost deterministic if large differences, almost uniform if small differences

don't use!

Selection methods

population P = (x_1, x_2, ..., x_μ) with μ individuals

- ***rank-proportional selection***
  order individuals according to their fitness values
  assign ranks
  fitness-proportional selection based on ranks

⇒ avoids all problems of fitness-proportional selection
but: best individual has only small selection advantage (can be lost!)

outdated!

- ***k-ary tournament selection***
  draw k individuals uniformly at random (typically with replacement) from P
  choose individual with best fitness (break ties at random)

⇒ has all advantages of rank-based selection and probability that best individual does not survive:

$$\left(1-\frac{1}{\mu}\right)^{k\,\mu} \begin{array}{ll} < & e^{-k} \\ \geq & 4^{-k} \end{array}$$

Selection methods without replacement

population P = (x_1, x_2, ..., x_μ) with μ parents and

population Q = (y_1, y_2, ..., y_λ) with λ offspring

- **(μ, λ)*-selection*** or ***truncation selection on offspring*** or ***comma-selection***
  rank λ offspring according to their fitness
  select μ offspring with best ranks

⇒ best individual may get lost, $\lambda \geq \mu$ required

- **(μ+λ)*-selection*** or ***truncation selection on parents + offspring*** or ***plus-selection***
  merge λ offspring and μ parents
  rank them according to their fitness
  select μ individuals with best ranks

⇒ best individual survives for sure

Selection methods: Elitism

Elitist selection: best parent is not replaced by worse individual.

- *Intrinsic elitism*: method selects from parent and offspring,
best survives with probability 1

- *Forced elitism*: if best individual has not survived then re-injection into population,
i.e., replace worst selected individual by previously best parent

method	P{ select best }	from parents & offspring	intrinsic elitism
neutral	< 1	no	no
fitness proportionate	< 1	no	no
rank proportionate	< 1	no	no
k-ary tournament	< 1	no	no
(μ + λ)	= 1	yes	yes
(μ , λ)	= 1	no	no

Variation operators: depend on representation

mutation → alters a <u>single</u> individual

recombination → creates single offspring from two or more parents

may be applied

- exclusively (either recombination or mutation) chosen in advance
- exclusively (either recombination or mutation) in probabilistic manner
- sequentially (typically, recombination before mutation); for each offspring
- sequentially (typically, recombination before mutation) with some probability

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Variation in $\mathbb{B}^n$

Individuals $\in \{ 0, 1 \}^n$

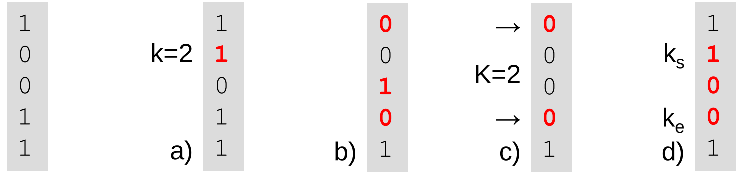
- Mutation

a) local → choose index $k \in \{ 1, \ldots, n \}$ uniformly at random, flip bit k, i.e., $x_k = 1 - x_k$

b) global → for each index $k \in \{ 1, \ldots, n \}$: flip bit k with probability $p_m \in (0,1)$

c) “nonlocal“ → choose K indices at random and flip bits with these indices

d) inversion → choose start index k_s and end index k_e at random invert order of bits between start and end index



Variation in $\mathbb{B}^n$

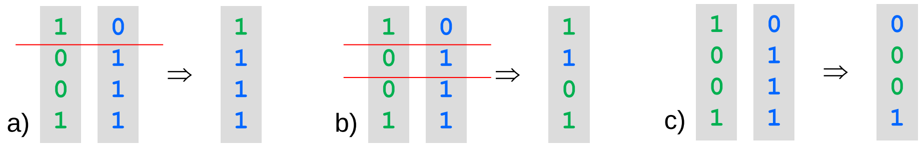
Individuals $\in \{ 0, 1 \}^n$

- Recombination (two parents)

a) 1-point crossover → draw cut-point $k \in \{1,\ldots,n-1\}$ uniformly at random; choose first k bits from 1st parent, choose last n-k bits from 2nd parent

b) K-point crossover → draw K distinct cut-points uniformly at random; choose bits 1 to k_1 from 1st parent, choose bits k_1+1 to k_2 from 2nd parent, choose bits k_2+1 to k_3 from 1st parent, and so forth …

c) uniform crossover → for each index i: choose bit i with equal probability from 1st or 2nd parent



Variation in $\mathbb{B}^n$

Individuals $\in \{ 0, 1 \}^n$

- Recombination (multiparent: ρ = #parents)

a) diagonal crossover ($2 < \rho < n$)

→ choose $\rho - 1$ distinct cut points, select chunks from diagonals

can generate ρ offspring; otherwise choose initial chunk at random for single offspring

b) gene pool crossover ($\rho > 2$)

→ for each gene: choose donating parent uniformly at random

Variation in $\mathbb{P}_n$

Individuals $\in X = \pi(1, \ldots, n)$

- Mutation

a) local → 2-swap / 1-translocation

b) global → draw number K of 2-swaps, apply 2-swaps K times

K is positive random variable;
its distribution may be uniform, binomial, geometrical, …;
E[K] and V[K] may control mutation strength

expectation variance

Variation in $\mathbb{P}_n$

Individuals $\in X = \pi(1, \ldots, n)$

- Recombination (two parents)

a) order-based crossover (OBX)

- select two indices k_1 and k_2 with $k_1 \le k_2$ uniformly at random
- copy genes k_1 to k_2 from 1st parent to offspring (keep positions)
- copy genes from left to right from 2nd parent, starting after position k_2

2 3 5 | 7 1 6 | 4
6 4 5 3 7 2 1

x x x 7 1 6 x

5 3 2 7 1 6 4

b) partially mapped crossover (PMX) [a version of]

- select two indices k_1 and k_2 with $k_1 \le k_2$ uniformly at random
- copy genes k_1 to k_2 from 1st parent to offspring (keep positions)
- copy all genes not already contained in offspring from 2nd parent (keep positions)
- from left to right: fill in remaining genes from 2nd parent

2 3 5 | 7 1 6 | 4
6 4 5 3 7 2 1

x x x 7 1 6 x

x 4 5 7 1 6 x

3 4 5 7 1 6 2

Variation in $\mathbb{P}_n$

Individuals $\in X = \pi(1, \ldots, n)$

- Recombination (two parents)

c) partially mapped crossover (PMX) [Grefenstette et al. 1985]
→ consider array as ring!

- given: 2 permutations **a** and **b** of length n
- select 2 indices k_1 and k_2 uniformly at random
- copy **b** to **c**
- procedure =

```
i = k1
repeat
  j = findIndex(a[i], c)
  swap(c[i], c[j])
  i = (i + 1) mod n
until i == k2
```

2 3 5 | 7 1 6 | 4
6 4 5 3 7 2 1

$6\ 4\ 5\ 3\ \boxed{7}\ 2\ 1$

$6\ 4\ 5\ 7\ 3\ 2\ \boxed{1}$

$\boxed{6}\ 4\ 5\ 7\ 1\ 2\ 3$

2 4 5 7 1 6 3

Variation in $\mathbb{R}^n$

Individuals $X \in \mathbb{R}^n$

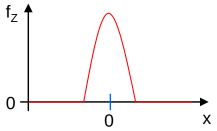
- Mutation

additive: $Y = X + Z$ (Z: n-dimensional random vector)

offspring = parent + mutation

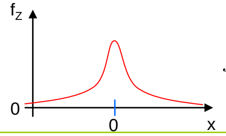
a) local → Z with bounded support

Definition
Let $f_Z: \mathbb{R}^n \to \mathbb{R}^+$ be p.d.f. of r.v. Z.
The set $\{ x \in \mathbb{R}^n : f_Z(x) > 0 \}$ is termed the *support* of Z.



$$f_Z(x) = \frac{4}{3}\,(1 - x^2) \cdot 1_{[-1,1]}(x)$$

b) nonlocal → Z with unbounded support



$$f_Z(x) = \frac{1}{\sqrt{2\,\pi}}\,\exp\left(-\frac{x^2}{2}\right)$$

most frequently used!

Variation in $\mathbb{R}^n$

Individuals X $\in \mathbb{R}^n$

- Recombination (two parents)

a) all crossover variants adapted from $\mathbb{B}^n$

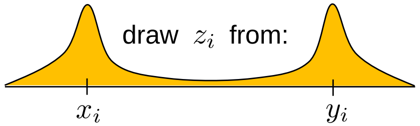
b) intermediate $z = \xi \cdot x + (1-\xi) \cdot y$ with $\xi \in [0,1]$

c) intermediate (per dimension) $\forall i : z_i = \xi_i \cdot x_i + (1-\xi_i) \cdot y_i$ with $\xi_i \in [0,1]$

d) discrete $\forall i : z_i = B_i \cdot x_i + (1-B_i) \cdot y_i$ with $B_i \sim B(1, \frac{1}{2})$

e) simulated binary crossover (SBX)

→ for each dimension with probability p_c draw z_i from:



Variation in $\mathbb{R}^n$

Individuals X $\in \mathbb{R}^n$

- Recombination (multiparent), ρ ≥ 3 parents

a) intermediate $z = \sum_{k=1}^{\rho} \xi^{(k)} x_i^{(k)}$ where $\sum_{k=1}^{\rho} \xi^{(k)} = 1$ and $\xi^{(k)} \geq 0$

(all points in convex hull)

b) intermediate (per dimension) $\forall i : z_i = \sum_{k=1}^{\rho} \xi_i^{(k)} x_i^{(k)}$

$$\forall i : z_i \in \left[\min_k \{x_i^{(k)}\}, \max_k \{x_i^{(k)}\} \right]$$

Theorem
Let f: $\mathbb{R}^n \to \mathbb{R}$ be a strictly quasiconvex function. If f(x) = f(y) for some x ≠ y then every offspring generated by intermediate recombination is better than its parents.

Proof:

f strictly quasiconvex $\Rightarrow f(\xi \cdot x + (1-\xi) \cdot y) < \max\{\, f(x), f(y) \,\}$ for $0 < \xi < 1$

since $f(x) = f(y) \Rightarrow \max\{\, f(x), f(y) \,\} = \min\{\, f(x), f(y) \,\}$

$\Rightarrow f(\xi \cdot x + (1-\xi) \cdot y) < \min\{\, f(x), f(y) \,\}$ for $0 < \xi < 1$

■

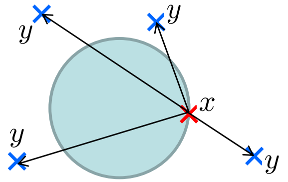
Theorem
Let f: $\mathbb{R}^n \to \mathbb{R}$ be a differentiable function and f(x) < f(y) for some x ≠ y.
If (y – x)‘ ∇f(x) < 0 then there is a positive probability that an offspring generated by intermediate recombination is better than both parents.

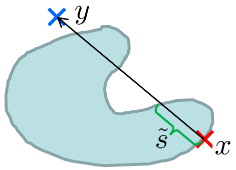
Proof:

If $d' \nabla f(x) < 0$ then $d \in \mathbb{R}^n$ is a direction of descent, i.e.

$\exists \tilde{s} > 0 : \forall s \in (0, \tilde{s}] : f(x + s \cdot d) < f(x)$.

Here: $d = y - x$ such that $\mathrm{P}\{f(\xi\, x + (1-\xi)\, y) < f(x)\} \geq \frac{\tilde{s}}{\|d\|} > 0$. ■





sublevel set $S_\alpha = \{x \in \mathbb{R}^n : f(x) < \alpha\}$