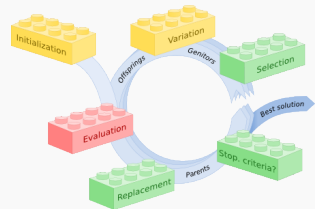


Evolutionary Programming

Evolutionary Programming (EP)

Generalities:

- ▶ Oldest idea of EA
[Fogel et al., 1964];
- ▶ population based thought as parallel search (not be confused with parallel implementation);
- ▶ oriented towards continuous/real-valued optimisation problems!



EP: representation

- ▶ Each individual is represented as a real-valued vector:

$$\langle \mathbf{x}, \boldsymbol{\sigma} \rangle = \langle x_1, \dots, x_n, \sigma_1, \dots, \sigma_n \rangle$$

- ▶ where the first part represents a solution \mathbf{x} and the second is a set of self-adapting parameters $\boldsymbol{\sigma}$, also referred to as *meta-EP*.
- ▶ This representation can be seen as n “evolving” normal distributions (one for each i^{th} design variable)

$$\mathcal{N}(\mathbf{x}[i], \boldsymbol{\sigma}[i])$$

EP: initialisation

- ▶ A population of μ individuals is generated.
- ▶ The design variables x_i are uniformly sampled within the decision space.
- ▶ The moving strategies σ_i are uniformly sampled between 0 and 1: $\mathcal{U}(0, 1)$

EP: parent selection

- ▶ Each solution generates a point, namely offspring!
 - ▶ i.e. No actual selection needed, all individuals are singularly considered!
 - ▶ μ individuals $\Rightarrow \mu$ offspring!

EP: recombination

- ▶ There is only one variation operator, i.e. mutation (no crossover!), which is more appropriately referred to as recombination.
- ▶ Each individual $\langle \mathbf{x}, \boldsymbol{\sigma} \rangle$ generates an offspring $\langle \mathbf{x}', \boldsymbol{\sigma}' \rangle$ according to the formulas:

$$\begin{cases} \sigma'_i = \sigma_i (1 + \alpha \cdot \mathcal{N}(0, 1)) \\ x'_i = x_i + \sigma'_i \cdot \mathcal{N}(0, 1) \end{cases} \quad i = 1, 2, \dots, n \quad \forall \langle \mathbf{x}, \boldsymbol{\sigma} \rangle \in Pop$$

- ▶ Each individual contains the information about the following move (early self-adaptation) with usually $\alpha = 0.2$.

EP: survivor selection (*deterministic variant*)

- ▶ The so-called $(\mu + \mu)$ selection strategy is usually applied:
 - ▶ the μ parents and the μ offspring are taken into account and sorted on the basis of their fitness value;
 - ▶ those μ solutions that have the best fitness/objective function value are selected for the subsequent generation.

This approach is elitist!

The point displaying the best fitness value always find a place in the next generation!

EP: survivor selection

(stochastic variant)

- ▶ Each solution is evaluated against q other randomly chosen solutions (round-robin);
- ▶ for each comparison, a “win” is assigned to the solution under examination when it outperforms one of the randomly selected solution;
- ▶ the solutions who scored most wins are selected for the following generation.

Still elitist

but some (good/bad) solutions are selected depending on the random choice of the competitors.

Genetic Algorithm

Genetic Algorithm (GA)

Generalities:

- ▶ Developed: USA in the 1970's.
- ▶ Early names: J. Holland, K. DeJong, D. Goldberg [Goldberg, 1989]¹.
- ▶ Typically applied to:
 - ▶ discrete optimization;
 - ▶ but also to real-valued problems.
- ▶ Attributed features:
 - ▶ not too fast;
 - ▶ versatile, good metaheuristic for combinatorial problems.

¹Interesting book about GAs. Consider also [Eiben and Smith, 2003].

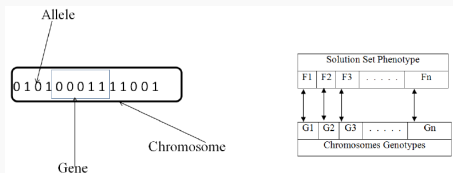
Simple Genetic algorithm

[Holland, 1975]

- ▶ Holland's original GA is now known as the simple genetic algorithm (SGA).
- ▶ The main features of this algorithm is the use of binary encoding.
- ▶ This choice was done in the 1970's to remark the analogy with biology (it is indeed inspired from biology!).
- ▶ This choice is generally VERY INEFFICIENT and should never applied unless the problem is naturally binary (here reported for historical reasons).

Representation

- ▶ A chromosome is encoded as a binary number: “chromosome”, composed of “genes” representing a given parameter to be optimised, and encoded with a given number of “alleles”².
- ▶ Due to the inner discretisation of the binary encoding GA can turn out more efficient in discrete optimization.



²N.B. If a gene represents a real number, we need to fix a precision (i.e. number of binary places): higher the precision higher the dimensionality of the discrete problem \Rightarrow difficult problem, slow evolution!

Parent selection mechanism

- ▶ The individuals that are undergoing recombination are selected by means of a Selection Mechanism.
- ▶ Classical selection mechanisms (there is many more but we will focus on these three):
 - ▶ Tournament;
 - ▶ Fitness Proportionate;
 - ▶ Ranking.

Selection pressure:

the property of the selection component in following the promising search directions (in other words, a parent selection mechanism which selects the best individuals many times has a high selection pressure).

Tournament selection

- ▶ Pick up a k solutions (at random) and compare their fitness, the better individual is in the mating pool.

Widely used! (usually the tournament size is $k = 2$)

- ▶ It does not require a sorting or a global knowledge (e.g. in a parallel implementation only a sub-portion of the population could be available) of the fitness distribution over the individuals of the population.
- ▶ The higher k , the higher the probability of selecting a good solution.
- ▶ It can be performed with or without replacement (see [Eiben and Smith, 2003] for details).

Fitness Proportionate Selection (FPS)

i.e. better individuals have higher probabilities to be selected! ✓

- ▶ It is given a probability P_i to be chosen to each i^{th} solution according to $P_i = \frac{f_i}{\sum_{i=1}^{\mu} f_i}$;
- ▶ the individual/individuals are selected through a mechanisms considering such probability (usually “roulette wheel”).

In some scenarios:

- ▶ outstanding individuals take over the entire population quickly: premature convergence ✗
- ▶ when fitness values are all similar (or equal \rightarrow plateau) there is no selection pressure ✗
- ▶ If this happens we adjust the previous formula with Goldberg's **sigma scaling**, (see [Eiben and Smith, 2003] and [Goldberg, 1989] for details).

Ranking selection

(always preserves some selection pressure ✓)

- ▶ Individuals are sorted according to their fitness value and a probability is assigned according to their position in the list (rank), and a probability is assigned to each i^{th} solution by means of:
 - ▶ **Linear Ranking:** $p_{lin-rank}(i) = \frac{2-s}{\mu} + \frac{2i(s-1)}{\mu(\mu-1)}$, $1.0 < s \leq 2.0$.
 - ▶ **Exponential Ranking:** $p_{exp-rank}(i) = \frac{1-e^{-i}}{c}$, c is a normalisation factor.

We still need a mechanism using this probability.

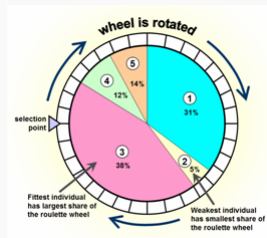
- ▶ s can be tuned to adjust the probability.
- ▶ Exponential ranking is preferred if an higher emphasis on selecting above-average individuals is required.

Implementing probability selection: *roulette wheel algorithm.*

- ▶ We work out probabilities P_i according to FPS or Ranking, then we select one or more individuals by spinning a one-armed roulette wheel with μ sectors, whose size is equal to P_i .

Roulette-Wheel algorithm (single spin)

```
 $i = 1$   
while  $\sum_1^i P_i < \mathcal{U}(0, 1)$  do  
   $i = i + 1$   
end while  
Output  $i^{th}$  individual in the population
```



- ▶ First spin \Rightarrow first element, second spin \Rightarrow second element,...
- ▶ Spins are independent! (not suitable when multiple solution have to be selected **X**)
 - ▶ in this case Stochastic universal Sampling can be adopted (see [Eiben and Smith, 2003])

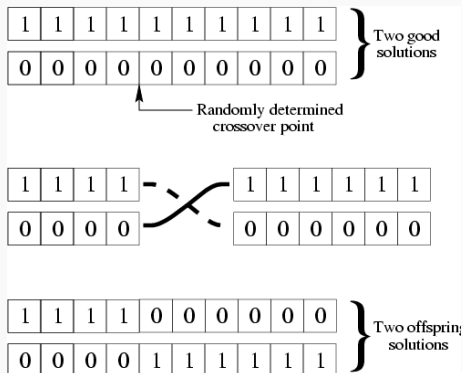
Crossover

- ▶ The selected parents undergo recombination.
- ▶ In a SGA, the recombination is the crossover.
- ▶ Crossover is an operator which combines two parents in order to produce one, two or more offspring.
- ▶ The analogy of biological crossover in binary encoding is very straightforward.

Crossover

1-point crossover

- It selects a random “cut-point” and switch head and tail of two chromosomes:



Crossover

N-point crossover

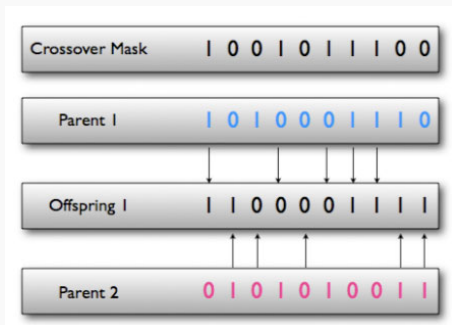
- ▶ Choose n random crossover points;
- ▶ split along those points;
- ▶ glue parts, alternating between parents:



Crossover

Uniform crossover

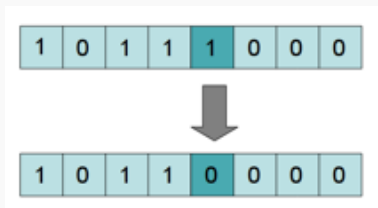
- ▶ Randomly (uniform distribution) initialise a mask for generating the first child:
 - ▶ we do not know how many cut-points we are going to have (they are uniformly distribute!)



- ▶ Make an inverse copy of the output if a second child is needed!

Mutation

- ▶ It is usually applied to the newly generated offspring before calculating its fitness value.
- ▶ Alter each gene independently with a probability p_m (referred to as “mutation rate”).



Crossover or mutation?

Which one is better? Which one is the most important?

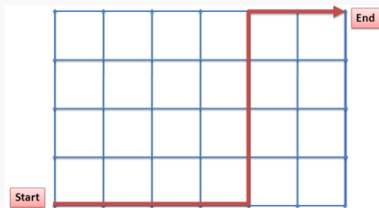
- ▶ It depends on the problem, but, in general, it is good to have both!
- ▶ Both have a role but mutation-only-EA is possible, crossover-only-EA seems not to work.

Survivor selection

- ▶ The main feature of a GA is that the algorithm must be generational (also called age-based):
 - ▶ in other words, the parents must be replaced by the newly generated offspring.
- ▶ Some implementations employ elitism: a restricted number of parents, the best, are copied for the subsequent generation.

Integer representation

- ▶ Some problems (e.g. image processing paramters, evolve a path in a grid...) naturally have integer variables:



in the example every gene is a
step taken from
 $\{North, South, West, East\}$
 $\Rightarrow \{1, 2, 3, 4\}$

- ▶ We use the same selection and crossover strategies of binary GAs!
 - ▶ But different mutations...

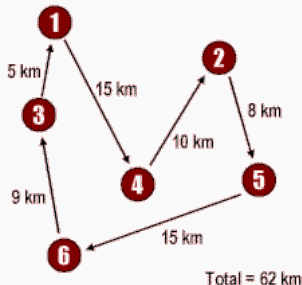
Mutation

(integer representation)

- ▶ *Creep mutation*: \forall genes in the chromosome, there is probability p of adding or subtracting a small random integer.
- ▶ *Random reset*: \forall genes in the chromosome, there is probability p to be replaced with a random integer chosen from the set of permissible values (no need to suture solution!).

Permutation representation

- Some problems (e.g. TSP, scheduling...) naturally take the form of deciding on the order in which a sequence of events has to occur:



$\Rightarrow [1, 4, 2, 5, 6, 3]$

- Still “integer” coded, but the same integer value never appears twice!

Crossover

(permutation representation)

- ▶ It is difficult to be applied in a permutation based GA:
 - ▶ if we exchange variables we can have the same value multiple times!
- ▶ The most popular were designed for specific applications:
 - ▶ Order 1 crossover;
 - ▶ Partially mapped crossover;
 - ▶ Cycle crossover;
 - ▶ Edge crossover.

A description is given in [Eiben and Smith, 2003].

Mutation

(*permutation representation*)

- **Swap**: Pick two design variables at random and swap their positions:

1 2 3 4 5 6 7 8 9 → 1 5 3 4 2 6 7 8 9

- **Insert**: Pick two design variables at random and move the second to follow the first, shifting the rest along to accommodate:

1 2 3 4 5 6 7 8 9 → 1 2 5 3 4 6 7 8 9

- **Inversion**: Pick two design variables at random and then invert the substring between them:

1 2 3 4 5 6 7 8 9 → 1 5 4 3 2 6 7 8 9

- **Scramble**: Pick two design variables at random and randomly rearrange the design variables in the substring:

1 2 3 4 5 6 7 8 9 → 1 3 5 4 2 6 7 8 9

Real-valued representation

(or of floating-point representation)

- ▶ Many problems occur as real valued problems, e.g. continuous parameter optimisation, design of an aircraft etc.
- ▶ In these cases, the most reasonable representation is a real valued as it would guarantee to work directly on the problem and not in a “deformed space”.
- ▶ Parent and survivor selections are the same of SGA!

Crossover

(*real-valued representation*)

For a real-valued GA we distinguish between

- ▶ *Deterministic* crossover: generally inefficient as it does not use the real-valued nature of the problem.
- ▶ *Arithmetic* crossover: the value of the design variables is varied due to some arithmetic operations.

Given a child z obtained from parents x and y :

deterministic $\Rightarrow z[i] = x[i] \text{ or } y[i]$

All those seen for SGA!

arithmetic $\Rightarrow z[i] = \alpha x[i] + (1 - \alpha)y[i], \alpha \in [0, 1]$

It can be:

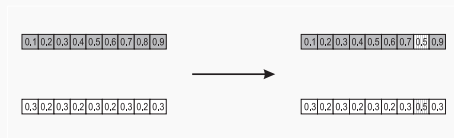
- ▶ randomised arithmetic recombination ($\alpha = \mathcal{U}(0, 1)$);
- ▶ simple, single or whole arithmetic recombination.

Crossover

Simple, single and whole arithmetic, ($\alpha = \frac{1}{2}$)



Simple arithmetic recombination



Single arithmetic recombination



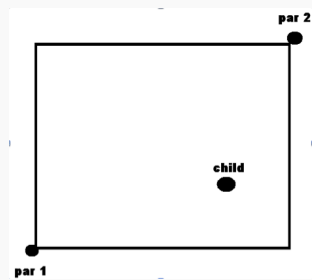
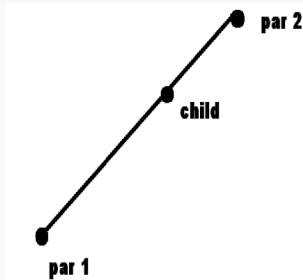
Whole arithmetic recombination

Box crossover (real-valued representation)

- ▶ The box crossover is another important method where:

$$z[i] = \min(x[i], y[i]) + \alpha |x[i] - y[i]|, \alpha = \mathcal{U}(0, 1)$$

- ▶ It samples one (or more) points within the hyper-rectangle identified by the two parents, while the algorithm crossover works on a straight line:



Mutation

(*real-valued representation*)

As for the binary case, there is a probability p of applying a mutation scheme that:

- ▶ replace the design variable with a uniformly distributed random number within the bounds (no need to saturate!):

$$\mathbf{x}_j[i] = \mathcal{U}(\mathbf{x}^L[i], \mathbf{x}^U[i]), \forall i = 0, 1, 2, \dots, n;$$

or

- ▶ add a small (rarely a big) random quantity to the variable from Gaussian (or Cauchy) distribution: e.g.

$$\mathbf{x}_j[i] = \mathbf{x}_j[i] + \mathcal{N}(0, \sigma_{small}), \forall i = 0, 1, 2, \dots, n.$$

Laboratory and participation work:

- ▶ For the four problems under consideration in this module in 10D and 50D:
 - ▶ implement a real-valued GA with 1) tournament selection; 2) box crossover; 3) Gaussian mutation; 4) Replacement of the population;
 - ▶ for the two dimensionality levels, tune the parameters (e.g. population size, ...) that appear the most convenient, and draw your own conclusion.

References I



Eiben, A. E. and Smith, J. E. (2003).

Introduction to Evolutionary Computation.

Springer-verlag, Berlin.



Fogel, L. J., Owens, A., and Walsh, M. (1964).

On the evolution of artificial intelligence (artificial intelligence generated by natural evolution process).

In *National Symposium on Human Factors in Electronics, 5th, San Diego, California*, pages 63–76.



Goldberg, D. E. (1989).

Genetic Algorithms in Search, Optimization and Machine Learning.

Addison-Wesley Publishing Co., Reading, MA, USA.



Holland, J. (1975).

Adaptation in Natural and Artificial Systems.

University of Michigan Press.