

## EA overview

Based on Baeck- Schwefel 1993

- General outlines
- Evolution strategies
- Evolutionary programming
- Genetic algorithms
- Simulated annealing
- Road map of EC
- Why and when to use EC?

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## Components of an Evolutionary Algorithm

- $f: \mathbb{R}^n \rightarrow \mathbb{R}$  objective function to be optimized
  - $\vec{x} \in \mathbb{R}^n$  an object variable vector
  - $I$  the space of individuals
  - $a \in I$  an individual
  - $\Phi: I \rightarrow \mathbb{R}$  the fitness function
  - $\mu \geq 1$  size of the (parent) population
  - $\lambda \geq 1$  number of offspring created in one cycle
  - $P(t) = \{\vec{a}_1(t), \dots, \vec{a}_\mu(t)\}$  population at generation  $t$
  - $r_{\Theta_r}: I^\mu \rightarrow I^\lambda$  recombination operator
  - $m_{\Theta_m}: I^\lambda \rightarrow I^\lambda$  mutation operator
  - $s_{\Theta_s}: (I^\lambda \cup I^{\mu+\lambda}) \rightarrow I^\mu$  selection operator
  - $t: I^\mu \rightarrow \{true, false\}$  termination criterion
- $\Theta_r, \Theta_m$  and  $\Theta_s$  are control parameters of  $r, m$  and  $s$  respectively

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## Outline of an Evolutionary Algorithm

```

t := 0;
initialize   P(0) := {a_1(0), ..., a_mu(0)} in I^mu;
evaluate     P(0) := {Phi(a_1(0)), ..., Phi(a_mu(0))};

while (t(P(t)) != true)
{
    recombine: P'(t) := r_Omega_r(P(t));
    mutate:    P''(t) := m_Omega_m(P'(t));
    evaluate   P''(t) := {Phi(a_1''(0)), ..., Phi(a_mu''(0))};
    select:    P(t+1) := s_Omega_s(P''(t) union Q);
    t := t + 1;
}
where Q in {0, P(t)}
  
```

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## Evolution Strategies (1)

### Representation (most general case)

$\vec{a} = (\vec{x}, \vec{\sigma}, \vec{\alpha})$ , where

- $x_i, i \in \{1, \dots, n\}$ , are object variables
- $\sigma_i, i \in \{1, \dots, n\}$ , are the mutation stepsizes, that is the standard deviations  $\sigma_i^2 = c_{ii}$
- $\alpha_j, j \in \{1, \dots, \frac{n \cdot (n-1)}{2}\}$ , are rotation angles

that is the covariances

$$\alpha_j \in \{c_{km} \mid k \in \{1, \dots, n-1\}, m \in \{k+1, \dots, n\}\}$$

where  $c_{km}$  ( $k, m \in \{1, \dots, n\}$ ) are the elements of the covariance matrix belonging to the generalized  $n$ -dimensional normal distribution with expectation vector 0

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## Evolution Strategies (2)

**Fitness function:**  $\Phi(\vec{a}) = f(\vec{x})$

**Mutation (most general case)**

$$\sigma_i' = \sigma_i \cdot \exp(\tau \cdot N(0,1) + \tau \cdot N_i(0,1))$$

$$\alpha_j' = \alpha_j + \beta \cdot N_j(0,1)$$

$$\vec{x}' = \vec{x} + \vec{N}(\vec{0}, \vec{\sigma}', \vec{\alpha}')$$

**Recombination (most general case)**

$$x_i' = \begin{cases} x_{S,i} & \text{without recombination} \\ x_{S,i} \text{ or } x_{T,i} & \text{discrete recombination} \\ x_{S,i} + \chi \cdot (x_{T,i} - x_{S,i}) & \text{intermediate recombination} \\ x_{S,i} \text{ or } x_{T,i} & \text{global*}, \text{discrete} \\ x_{S,i} + \chi_i \cdot (x_{T,i} - x_{S,i}) & \text{global*}, \text{intermediate} \end{cases}$$

\* S, T  $\in \{1, \dots, \mu\}$  are redrawn for each  $i$  anew.

For  $\sigma$ 's and  $\alpha$ 's the same mechanism

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## Evolution Strategies (3)

### Selection:

Deterministic, selecting the  $\mu$  best ( $1 \leq \mu < \lambda$ ) out of

- the set of  $\lambda$  offspring individuals:  $(\mu, \lambda)$ -selection
- the union of parents and offspring:  $(\mu + \lambda)$ -selection

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## Evolution Strategies (4)

## Outline of an Evolutionary Strategy:

```

t := 0;
initialize
  where
    and
      P(0) := P(0) := {a_1(0), ..., a_mu(0)} in I^mu;
      l = R^{n*mu}
      a_k = (x_k, sigma_k)
      for all i in {1, ..., n}, j in {1, ..., mu * (1 - n) / 2};
evaluate
  where
    Phi(a_k(0)) = f(x_k(0));
while (t(P(t)) != true)
{
  recombine: a_k'(t) := r'(P(t))      for all k in {1, ..., lambda};
  mutate:    a_k''(t) := m_{v_i, r_{\Phi}}'(a_k'(t)) for all k in {1, ..., lambda};
  evaluate:  P''(t) := {Phi(a_1''(t)), ..., Phi(a_mu''(t))};
              where
                Phi(a_k''(t)) = f(x_k''(t));
  select:    P(t+1) := if (mu, lambda)-selection then s(mu, lambda)(P''(t));
              else (mu + lambda)(P(t) union P''(t));

  t := t + 1;
}

```

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## Evolutionary Programming (1)

## Representation

- Standard EP:  $\tilde{a} = \bar{x}$
- Meta-EP:  $\tilde{a} = (\bar{x}, \bar{\sigma})$

## Fitness function

scaling objective function values  $f(\bar{x})$  to positive values and possibly imposing some random alteration

$$\Phi(\tilde{a}) = \delta(f(\bar{x}), \kappa)$$

where

- $\alpha_i: \mathbb{R} \times \mathbb{S} \rightarrow \mathbb{R}^+$  denotes the scaling function and  $\mathbb{S}$  is an additional set of parameters,
- $\kappa$  is some random alteration factor

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## Evolutionary Programming (2)

## Mutation

- Standard EP: most common way is  $x_i' = x_i + \sqrt{\Phi(\bar{x})} \cdot N_i(0,1)$

- Meta-EP:  $x_i' = x_i + \sigma_i \cdot N_i(0,1)$

$$\sigma_i' = \sigma_i + \alpha \cdot \sigma_i \cdot N_i(0,1)$$

where the parameter  $\alpha$  ensures that  $v_i$  tends to remain positive

## Recombination: None

## Selection

stochastic q-tournament ( $\mu + \mu$ ) style selection, sorting individuals by their score  $w$ , where

$$w_i = \sum_{j=1}^{\mu} \begin{cases} 1 & \text{if } \Phi(\bar{a}_i) \leq \Phi(\bar{a}_{x_j}) \\ 0 & \text{otherwise} \end{cases}$$

where  $x_j \in \{1, \dots, 2\mu\}$  is a uniform integer random variable, sampled anew for each comparison

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## Evolutionary Programming (3)

## Outline of an Evolutionary Strategy:

```

t := 0;
initialize
  where
    and
      P(0) := P(0) := {a_1(0), ..., a_mu(0)} in I^mu;
      l = R^{n*mu}
      a_k = (x_k, v_k)
      for all i in {1, ..., n};
evaluate
  where
    Phi(a_k(0)) = delta(f(x_k(0)), kappa_k);
while (t(P(t)) != true)
{
  recombine: a_k'(t) := m_{\alpha}(\tilde{a}_k(t))      for all k in {1, ..., mu};
  evaluate:  P''(t) := {Phi(a_1'(t)), ..., Phi(a_mu'(t))};
              where
                Phi(a_k'(t)) = delta(f(x_k'(t)), kappa_k)
  select:    P(t+1) := s_{\alpha}(P(t) union P''(t));

  t := t + 1;
}

```

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## Genetic Algorithms (1)

## Representation

Bit strings of fixed length  $\ell$ , i.e.  $l = \{0, 1\}^\ell$ .

For a continuous objective function:

$$f: \prod_{i=1}^n [u_i, v_i] \rightarrow \mathbb{R}, \text{ with } u_i < v_i$$

$$\ell = n \cdot \ell_s,$$

$$\bar{a} = (a_{11}, \dots, a_{n\ell_s}) \in \{0,1\}^{n\ell_s} = I \text{ and}$$

- a typical decoding is  $I' = I^n \times \dots \times I^n$ , where

$$I'^i(a_{i1}, \dots, a_{i\ell_s}) = u_i + \frac{v_i - u_i}{2^{\ell_s}} \left( \sum_{j=1}^{\ell_s} a_{ij} 2^{j-1} \right)$$

## Fitness function

$$\Phi(\bar{a}) = \delta(f(I'(\bar{a}))),$$

Where  $\delta$  is a scaling function assuring positive fitness values and best individual - largest fitness.

E.g. linear scaling with a scaling window of  $w$  generations.

$$\delta(f(I'(\bar{a})), P(t-w)) = \max\{f(I'(\bar{a}_j)) \mid \bar{a}_j \in P(t-w)\} - f(I'(\bar{a}))$$

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## Genetic Algorithms (2)

## Mutation

Bit flips with probability  $p_m$  per bit, that is

$$s_i' = \begin{cases} s_i, & \chi_i > p_m \\ 1 - s_i, & \chi_i \leq p_m \end{cases}$$

where  $p_m$  is an external parameter, the mutation rate,  $\chi_i$  is drawn from a uniform distribution

## Recombination

One-point crossover: if  $\bar{s} = (s_1, \dots, s_\ell)$ ,  $\bar{v} = (v_1, \dots, v_\ell)$  are selected as would-be parents, crossover takes place with probability  $p_c$ , called the crossover rate.

$$\bar{s}' = (s_1, \dots, s_{\chi-1}, s_\chi, v_{\chi+1}, \dots, v_\ell)$$

$$\bar{v}' = (v_1, \dots, v_{\chi-1}, v_\chi, s_{\chi+1}, \dots, s_\ell)$$

Where  $\chi \in \{1, \dots, \ell - 1\}$  denotes a uniform random variable, the crossover point.

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## Genetic Algorithms (3)

**Selection**

Fitness proportional selection: probability of being selected is

$$p_s(\bar{a}_i) = \frac{\Phi(\bar{a}_i)}{\sum_{j=1}^{\mu} \Phi(\bar{a}_j)}$$

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## Genetic Algorithms (4)

Outline of an Genetic Algorithm:

```

t := 0;
initialize P(0) := P(0) := {̂a1(0), ..., ̂aμ(0)} ∈ Iμ;
where I = {0,1}l;
evaluate P(0) := {Φ(̂a1(0)), ..., Φ(̂aμ(0))};
where Φ(̂ak(0)) = δ(f(Γ(̂ak(0))), P(0));

while ((P(t)) ≠ true)
{
  recombine: ̂ak'(t) := r(pop)'(P(t))      ∀k ∈ {1, ..., μ};
  mutate: ̂ak''(t) := m(gen)'(̂ak'(t))      ∀k ∈ {1, ..., μ};
  evaluate: P''(t) := {̂a1''(t), ..., ̂aμ''(t)} : {Φ(̂a1'(t)), ..., Φ(̂aμ'(t))};
  where Φ(̂ak''(t)) = δ(f(Γ(̂ak''(t))), P(t - w));
  select: P(t + 1) := s(P''(t));
  where ps(̂ak''(t)) = Φ(̂ak''(t)) / ∑j=1μ Φ(̂aj''(t))

  t := t + 1;
}

```

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## Comparative overview of EA'S

	ES	EP	GA
Repr.	Real	Real	Binary
Self-adapt.	Standard dev's and covariances	Variances (in meta-EP)	None
Fitness	Objective function	Scaled obj. function	Scaled obj. function
Mutation	Main op.	Only op.	Background
Recomb.	Different, important for self-adaptation	None	Main op.
Select.	Deterministic, extinctive	Probab., extinctive	Probab., preservative

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## Simulated Annealing (1)

Outline of a Local Search procedure

**procedure** local-search:

```

{
  initialize(istart);
  repeat
    generate(j ∈ Si);
    if (f(j) < f(i)) then i := j;
  until (f(j) ≥ f(i)) for all j ∈ Si;
}

```

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## Simulated Annealing (2)

**Main drawback:** gets easily stuck in local optima Possible cures:

- 1) several restarts with new i<sub>start</sub>
- 2) sophisticated neighborhood structure
- 3) accept j even if it is worse than i

Option 3) by analogy from condensed matter physics, where state transitions lead to minimal energy level

$$P[\text{accept } j] = \exp\left(\frac{E_i - E_j}{K_b \cdot T}\right)$$

This is called the Metropolis criterion.

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## Simulated Annealing (3)

Outline of simulated annealing:

**procedure** simulated-annealing:

```

{
  initialize(istart, c0, L0);
  k := 0; i := istart;
  repeat
    for ℓ := 1 to Lk do {
      generate(j ∈ Si);
      if (f(j) < f(i)) then i := j;
      else if exp((f(i) - f(j)) / ck}) > random(0,1) then i := j;
    }
    k := k + 1;
    calculate-length(Lk);
    calculate-control(ck);
  until stopcriterion;
}

```

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### Simulated Annealing (4)

#### Remarks:

1. The 'temperature'  $c_k$  is decreased over time.
2. As  $c_k$  decreases, selection becomes more and more elitist.
3. The neighborhood  $S_i$  is mostly defined 'operationally', i.e. as the set of points generated with one modification (mutation) operator from  $i$ .

#### Simulated Annealing can be seen as an EA

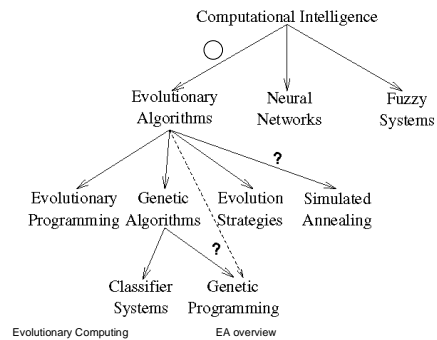
- using a (1+1) selection strategy together with a
- a specific, time dependent selection regime
- where the representation and the mutation operator are not specified (left as problem dependent component)

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### Roadmap of EC



### Why & when use evolutionary computing?

- **No in-depth mathematical understanding** of the problems needed
- Can **solve "out of range" problems** (that cannot be solved by analytical mathematical techniques), for instance: many variables, many local optima, moving goal posts
- They are **extremely robust**; they cope well with noisy, inaccurate and incomplete data
- They are relatively **cheap and quick to implement**
- They are **easily hybridised**; they combine very productively with other techniques such as greedy methods, heuristics, simulated annealing and neural networks

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### Why & when use evolutionary computing?

- Extremely **adaptable**; changed priorities can be incorporated simply by changing weightings in the fitness function
- They are **modular and therefore portable**; because the evolutionary mechanism is separate from the problem representation they can be transferred from problem to problem
- They provide an extremely open and flexible approach to design, **allowing arbitrary constraints**, simultaneous **multiple objectives** and the **mixing of continuous and discrete parameters**
- Unlike many other methods, when evolutionary algorithms are **implemented on parallel computers** they make very efficient use of the available power

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