

# Stochastic optimization -

Mono-agent methods :  
Recall & LAHC/SA



# Recalls

# Metaheuristics

Methods that are :

- global ;
- stochastic ;
- generic they depend on the context (continuous/discrete) ;
- does not guarantee optimality ;
- based on analogies with biology, physic, animals behaviours...

# Elements & strategies

Usually, they are composed of :

- one or many initial candidate solutions (population(s)) ;
- a random-based neighbor creation strategy ;
- a new candidate acceptance criteria ;
- coefficients parameters to control algorithm behavior ;
- information sharing, a memory, a selection strategy, ... ;
- a convergence criteria.

# Searching...

Exploration

VS

Exploitation



# Neighborhood

- Definition
  - Discrete variables : shift/permutation of data,
  - Continuous variables : hyper(sphere | cube) centered around the data.
- Strategies
  - geographic neighborhood,
  - social neighborhood,
  - random neighborhood.
- Gradient, Newton-Raphson, dichotomy methods, Nelder-Mead polytope...
- Hybridation local search methods.

# Late Acceptance Hill Climbing



# Late Acceptance Hill Climbing (LAHC)

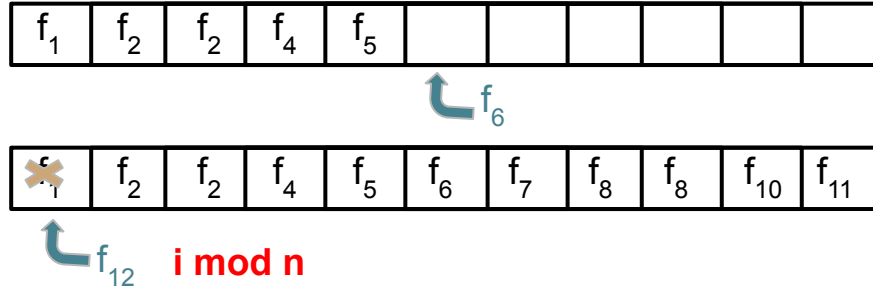
- Proposed by Edmond K. Burke & Yuri Bykov (2008) ;
- Mono-agent, adapted to discrete problems, algorithm ;
- Use of **scores memory** of visited candidates ;
- Based on Greedy Hill Climbing heuristic : “Accept a new candidate if it is better than current” ;
- “Accept a new candidate if it is better than a previous one (from a fixed number of iterations)” ;
- A unique parameter : memory size.

# LAHC : principle

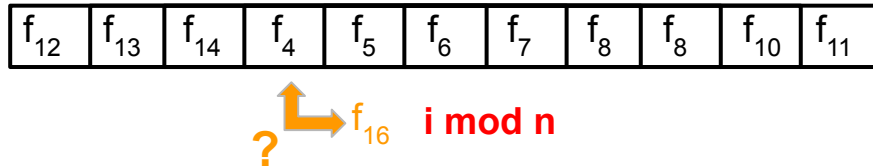
- Current candidate generates a new neighbor ;
- Current evaluation is added at the end of evaluations memory (modulo  $n$ ) ;
- Compare the new evaluation to the performance of  $n^{th}$  past candidate ( $n$  : memory size) ;
- If better, the neighbor becomes the current candidate.

# LAHC : illustration

- Current candidate score is added to memory ;



- The comparison is made between the neighbor evaluation and the  $n^{th}$  previous evaluation.



# LABC algorithm (minimizing)

```
X, a candidate, f(X) its evaluation
fmin ← f(X), Xmin ← X, tabMem ← [fmin | ... | fmin]
n ← 1
While not stoppingCriterion()
    Xvois ← generateNeighbor(X)
    fmem ← tabMem[n % memSize]
    If fmem ≥ f(Xvois) then
        X ← Xvois
    end if
    tabMem[n % memSize] ← f(X)
    if fmin > f(Xvois) then
        Xmin ← Xvois
        fmin ← f(Xvois)
    end if
    n++
End while
```

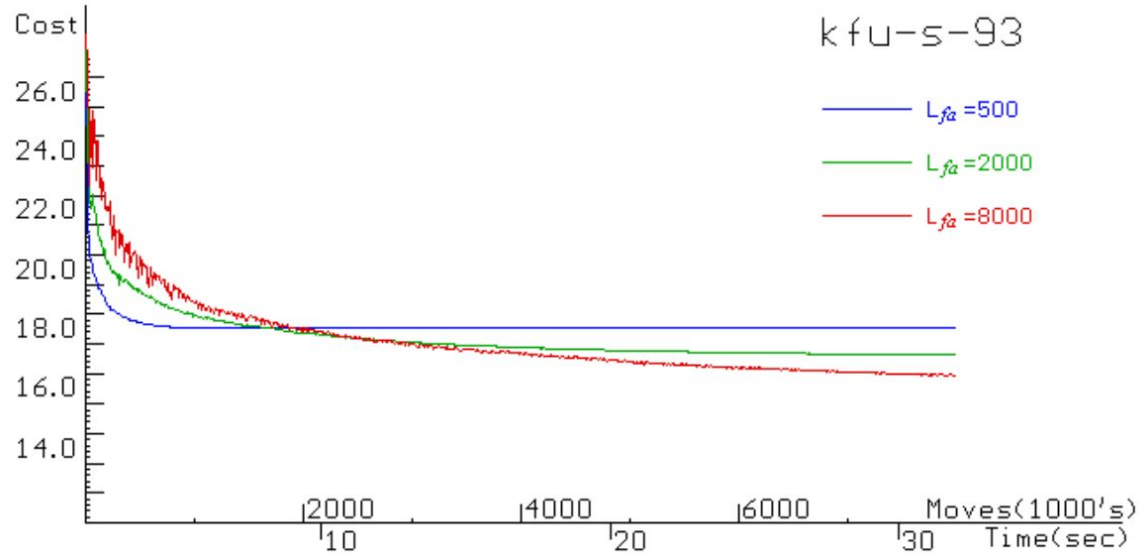
# LAHC : memory management

The memory of previous candidate scores is managed as :

- Array of  $n$  evaluations of visited candidates ;
- The memory size is a **critical** parameter :
  - $n$  small : few values recorded, behavior similar to hill-climbing heuristic = premature convergence,
  - $n$  big : many random moves favours diversity but slow down convergence.
- At initializing, all values are initialized to the first one. It prevents from too random moves during the first iterations.

# LAHC : memory management

Influence of memory size of the convergence behavior in a case of scheduling problem :



# LAHC : improvements

Comparison to a previous candidate :

- Depends on memory size (original version) ;
- Compare to memory best score ;
- Compare to a random value of memory ;
- Compare to a mean/median of memory values.

# LAHC : improvements

Accepting criterion is :

- Deterministic :
  - evaluation less or equal to  $n^{\text{th}}$  previous memory score (by default),
  - evaluation strictly less than  $n^{\text{th}}$  previous memory score,
  - evaluation less or equal to  $n^{\text{th}}$  previous memory score, or current one,
  - ...
- Probabilistic :
  - Metropolis-like :  $p = e^{- (c_{\text{vois}} - c_{\text{cur}}) / (\text{mem}[i] - c_{\text{cur}})}$ ,
  - linear :  $p = 1 - (c_{\text{vois}} - c_{\text{cur}}) / [2(\text{mem}[i] - c_{\text{cur}})]$ .

# LAHC : improvements

Memorize score according to :

- Descending-only:
  - replace in memory if  $c < \text{mem}[i]$  ;
- Cooling strategy :
  - after convergence, increment all memory stored scores of  $p\%$  ( $p=20$ ).

# Simulated annealing



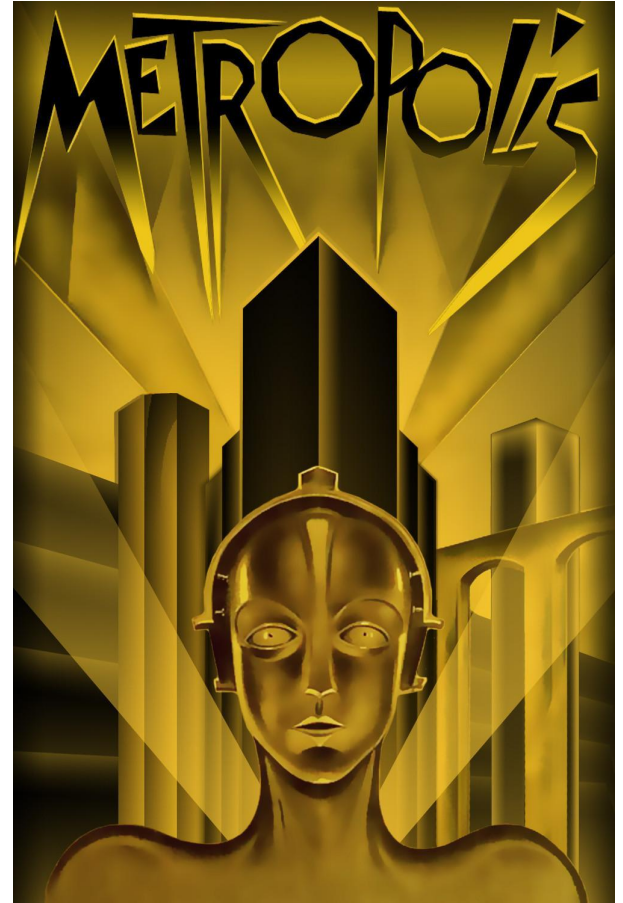
# Simulated annealing

- Metaheuristic, variant from Metropolis-Hastings algorithm (see example [here](#));
- Proposed by Kirkpatrick, Gelatt & Vecchi (1983) and by Cerny (1985);
- First metaheuristic ;
- Adapted to discrete problems (originally configure electronic components disposition on a printed circuit).

# Criterion of ....

```
critereMetropolis( $\Delta f, T$ ) : //minimizing  
  If  $\Delta f \leq 0$  then  
    return TRUE  
  else  
    return random(0,1) <  $e^{(-\Delta f/T)}$ 
```

- $\Delta f \leq 0$ , neighbor is accepted ;
- A small offset towards a less good neighbor has more chance to be accepted than a huge one ;
- **Stochastique** function.



# Algorithm (minimization)

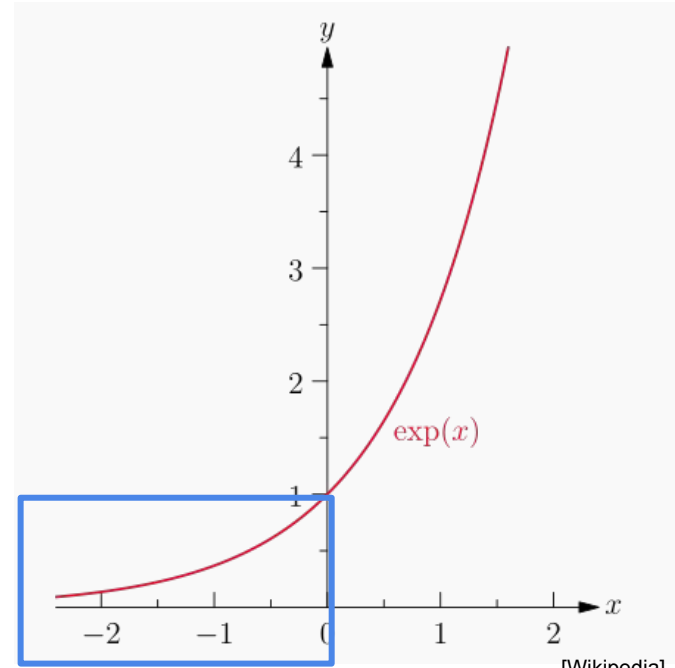
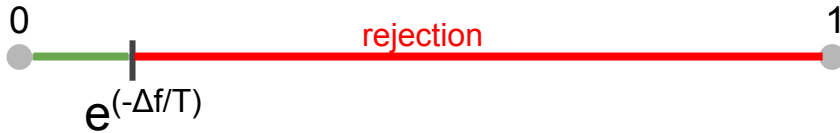
```
X, a random candidate,  $f_x \leftarrow f(X)$  score, T Initial temperature
 $X_{\min} \leftarrow X$ ,  $f_{\min} \leftarrow f(X)$ 
While  $T > T_{\min}$  and not stoppingCriterion()
    While not thermodynamicEquilibrium()
         $X_{\text{vois}} \leftarrow \text{generateNeighbor}(X)$ 
         $f_{X_{\text{vois}}} \leftarrow f(X_{\text{vois}})$ 
         $\Delta f = f_{X_{\text{vois}}} - f_x$ 
        If metropolisCriterion( $\Delta f, T$ ) then
             $X \leftarrow X_{\text{vois}}$ 
             $f_x \leftarrow f_{X_{\text{vois}}}$ 
            If  $f_{X_{\text{vois}}} < f_{\min}$  then
                 $f_{\min} \leftarrow f_{X_{\text{vois}}}$ 
                 $X_{\min} \leftarrow X_{\text{vois}}$ 
            End if
        End if
    End while
    T  $\leftarrow$  cooling(T)
End while
```

# Cooling

- High temperature, Metropolis criterion value is close to 1, most of worst candidates are accepted :



- Low temperature, Metropolis criterion value is close to 0, most of worst candidates are rejected :



Variation space of  
Metropolis criterion values

[Wikipedia]

# Cooling

Different schemes of temperature cooling :

- Geometric :  $T_{k+1} = \alpha.T_k$ , the mostly used ;
- Logarithmic :  $T_k = \mu / \log(1+k)$ , where  $k$  : nb of thermodynamic steps &  $\mu$  constant. expensive in computation time, rarely used ;
- Exponential :  $T_k = T_0 \cdot \exp(-k/\tau)$ , où  $k$  : nb of steps &  $\tau$  constant ;
- esotéric ....

## Question :

“How to profit from a distribution framework to improve an optimization algorithm”

?