

Genetic Algorithms and Genetic Programming

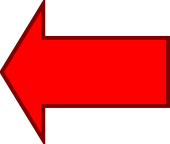
Lecture 15: (17/11/09)

Particle Swarm Optimization II



Michael Herrmann

Overview

- I. GA (1-7)
- II. GP (8-10)
- III. ACO (11-13): Ant colony optimization
- IV. **PSO (14-15): Particle swarm optimization and differential evolution** 
- V. **NC (16):** Overview on DNA computing, Membrane computing, Molecular computing, Amorphous computing, Organic computing,
- VI. Wrapping up: Metaheuristic search (17)

Not included:

artificial neural networks, quantum computing, cellular automata, artificial immune systems

The canonical PSO algorithm

For each particle $1 \leq i \leq n$, $x_i \in \mathbf{R}^m$, $v_i \in \mathbf{R}^m$

- create random vectors

r_1, r_2 with components drawn from $U[0,1]$

- update velocities $v_i \leftarrow \omega v_i + \alpha_1 r_1 \circ (\hat{x}_i - x_i) + \alpha_2 r_2 \circ (\hat{g} - x_i)$
inertia personal best global best

- update positions

$$x_i \leftarrow x_i + v_i$$

◦ componentwise multiplication

- update local bests

$$\hat{x}_i \leftarrow x_i \quad \text{if } f(x_i) < f(\hat{x}_i)$$

minimization problem!

- update global best

$$\hat{g} \leftarrow x_i \quad \text{if } f(x_i) < f(\hat{g})$$

Analysis of PSO: Simplified algorithm

- Consider a single particle only (“the view from inside”)
- Ignore randomness (use a homogeneous mean value)
- Ignore the global best (assume it equals personal best)
- Keep the personal best constant (changes are rare)
- Set inertia to unity (for the moment only)
- i.e. what we had (vector equation of i -th particle)

$$v_i(t+1) = \omega v_i(t) + \alpha_1 r_1 \circ (\hat{x}_i - x_i(t)) + \alpha_2 r_2 \circ (\hat{g} - x_i(t))$$

- becomes now (in component form: $d=1, \dots, m$ with $p_i = \hat{x}_i$)

$$v_{id}(t+1) = v_{id}(t) + \varphi(p_{id} - x_{id}(t))$$

Algebraic point of view

- Introduce $y_t = p - x_t$

$$v_{t+1} = v_t + \varphi y_t$$

$$x_{t+1} = x_t + v_{t+1} \implies y_{t+1} = -v_t + (1 - \varphi)y_t$$

- Introduce state vector $P_t = (v_t, y_t)^T$ and

$$M = \begin{pmatrix} 1 & \varphi \\ -1 & 1 - \varphi \end{pmatrix} \quad P_{t+1} = MP_t$$

- Starting from the initial state P_0 we have $P_t = M^t P_0$

Algebraic point of view

- Eigenvalues of $M = \begin{pmatrix} 1 & \varphi \\ -1 & 1-\varphi \end{pmatrix}$ i.e. $AMA^{-1} = L = \begin{bmatrix} e_1 & 0 \\ 0 & e_2 \end{bmatrix}$

$$e_{1/2} = 1 - \frac{\varphi}{2} \pm \frac{\sqrt{\varphi^2 - 4\varphi}}{2}$$

- Transformation matrix: $A = \begin{bmatrix} \varphi + \sqrt{\varphi^2 - 4\varphi} & 2\varphi \\ \varphi - \sqrt{\varphi^2 - 4\varphi} & 2\varphi \end{bmatrix}$

$$P_{t+1} = MP_t \implies P_{t+1} = A^{-1}LAP_t$$

$$\underbrace{AP_{t+1}}_{Q_{t+1}} = \underbrace{LAP_t}_{Q_t}$$

$$Q_{t+1} = LQ_t$$

- Thus $Q_t = L^t Q_0$ where

$$L^t = \begin{bmatrix} e_1^t & 0 \\ 0 & e_2^t \end{bmatrix}$$

3 Cases: $0 < \varphi < 4$

$\varphi = 4$

$\varphi > 4$

EV complex

$$e_1 = \cos(\theta) + i \sin(\theta)$$

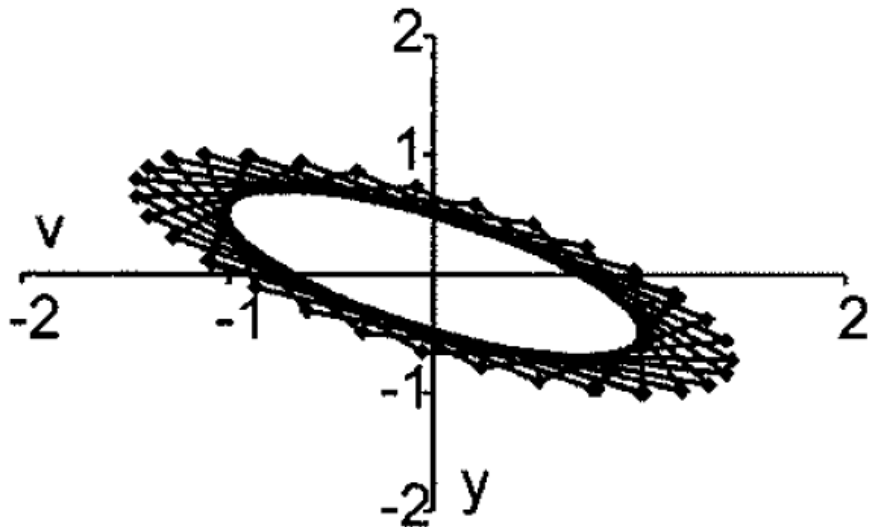
$$e_2 = \cos(\theta) - i \sin(\theta)$$

$$e_1^t = \cos(t\theta) + i \sin(t\theta)$$

$$e_2^t = \cos(t\theta) - i \sin(t\theta)$$

Oscillatory with

period k : $\theta = (2k\pi)/t$
(or quasiperiodic)



$$M = \begin{bmatrix} 1 & 4 \\ -1 & -3 \end{bmatrix}$$

$$V = \begin{bmatrix} -2 \\ 1 \end{bmatrix}$$

Both EV are -1

$$MV = -V$$

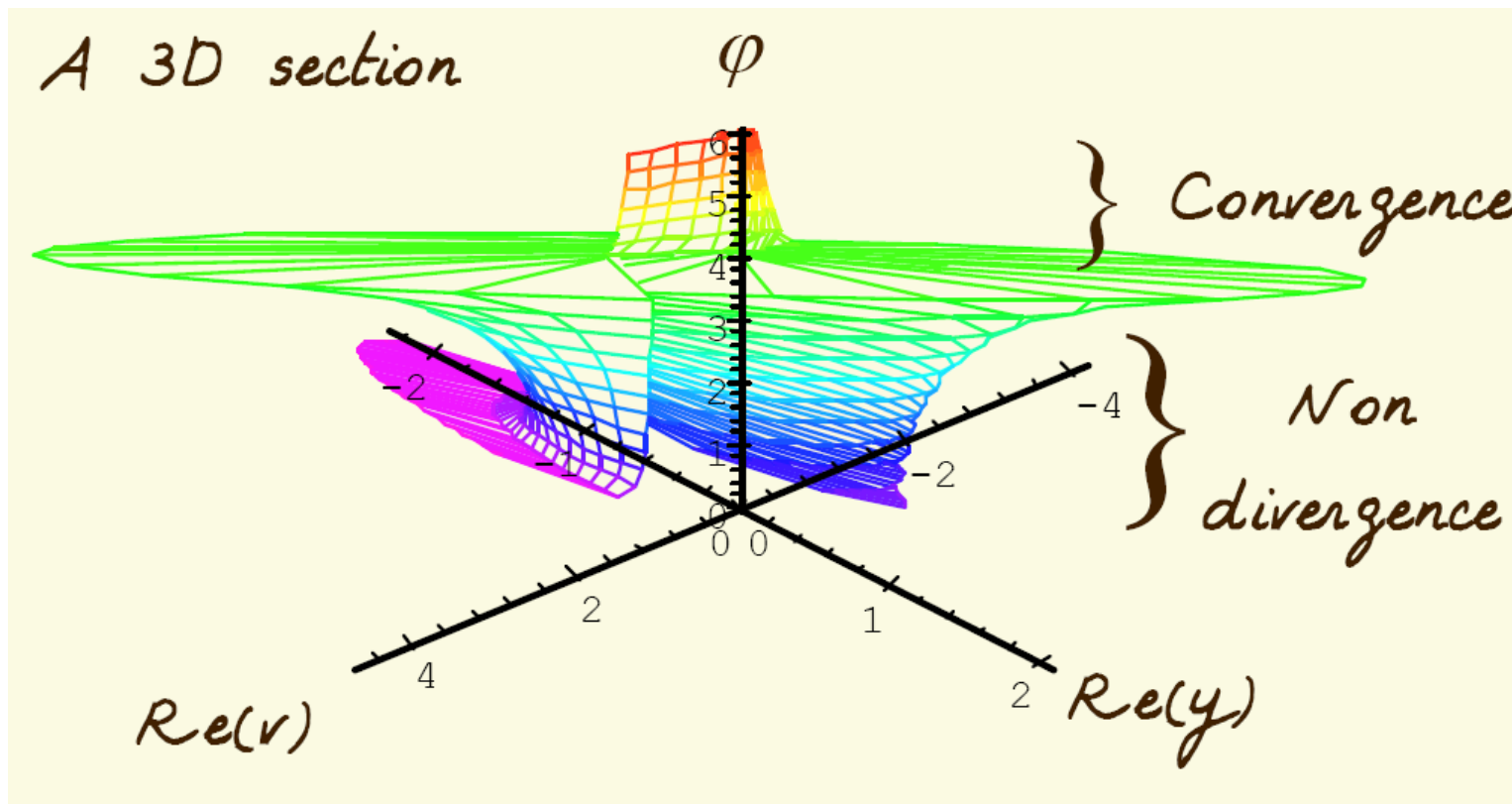
$$P_{t+1} = \pm \begin{bmatrix} 2y_0 \\ -y_0 \end{bmatrix} = -P_t$$

Some algebra implies:
Linear divergence
(unless starting
with the eigenvector V)

EV real

Exponent.
divergent

(convergent
with con-
striction)



Implications from the algebra

- Oscillation for $\varphi < 4$: Exploration near current best
- Divergence for $\varphi > 4$: Exploration of the wider environment
- $\varphi = \alpha_1 + \alpha_2$ is a combination of the attractiveness of the personal and global best. Since these might be not the same, a slightly larger φ might be needed.
- φ slightly above 4 (e.g. 4.1): particle stays somewhere in between or near personal and global best. If these two coincide the algorithm tends to diverge, i.e. the particle moves on searching elsewhere.
- Divergence can be counteracted by V_{\max} or by constriction.
- Remember that we were considering an averaged version of the algorithm.

Object Tracking in Computer Vision

Particle Swarms as Video Sequence Inhabitants For Object Tracking in Computer Vision

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Elena Sánchez Nielsen. DEIOC, ULL



Applications

- Evolving structure and weights of neural networks
- Complex control involving complex and continuous variables (power systems)
- Industrial mixer in combination with classical optimization
- Image analysis
- Medical diagnosis
- Job scheduling
- Robot path planning, localization
- Electrical generator
- Electrical vehicle

Repulsive PSO algorithm

For each particle $1 \leq i \leq n$

- create m -dimensional random vectors

r_1, r_2, r_3 with components drawn from $U[0,1]$

- update velocities

$$v_i \leftarrow \omega v_i + \alpha_1 r_1 \circ (\hat{x}_i - x_i) + \alpha_2 r_2 \circ (\hat{y} - x_i) + \alpha_3 \omega r_3 \circ z$$

- update positions **etc.** ◦ componentwise multiplication

z random velocity

\hat{y} best of a random neighbor or global best, $\alpha_2 < 0$

- Usually applied alternating with canonical PSO if diversity becomes too small
- Properties: sometimes slower, more robust and efficient

Fully Informed Particle Swarm (FIPS)

- Rui Mendes (2004): Simpler, maybe better
- Distributes total φ across n terms
- All neighbors contribute to the velocity adjustment
- Best neighbor is not selected, but included with a
- Individual not included in neighborhood
- Fails often, but, if successful, results are good, (strongly dependent on good topology)

$$\vec{\varphi}_k = \vec{U} \left[0, \frac{\varphi_{max}}{|\mathcal{N}|} \right] \quad \forall k \in \mathcal{N}$$

$$\vec{P}_m = \frac{\sum_{k \in \mathcal{N}} \mathcal{W}(k) \vec{\varphi}_k \otimes \vec{P}_k}{\sum_{k \in \mathcal{N}} \mathcal{W}(k) \vec{\varphi}_k}$$

Parameters, Conditions, & Tweaks

- Initialization methods
- Population size
- Population diameter
- Absolute vs. signed velocities
- Population topology
- Births, deaths, migration
- Limiting domain (X_{MAX} , V_{MAX})
- Multiobjective optimization
- “Subvector” techniques
- Comparison over problem spaces
- Hybrids

Jim Kennedy
Russ Eberhart:
Tutorial on Particle
Swarm Optimization

IEEE Swarm
Intelligence
Symposium 2005
Pasadena, California
USA, June 8, 2005

Remarks on PSO

- Consider boundaries as physical (e.g. by reflection from walls)
- Try adaptive versions: variable swarm size, variable ratios α_1/α_2
- Try different topologies (e.g. “tribes”)
- For local variants, consider using other norms in high-dimensional spaces (Euclidean unit sphere volume decays)

Relation to probabilistic methods

- Strict probabilistic methods are based on assumptions (Gaussianity, optimal sampling etc.) which often do not hold in practical applications
- There are many examples where meta-heuristic approaches do well
 - Toy examples are often designed ad-hoc for a particular method and are thus unsuitable for a fair comparison.
 - Success in real-world examples depends much on domain knowledge, quality of analysis, iterative re-design etc.
- Meta-heuristic algorithms may use strict algorithms for local search
- Meta-heuristic algorithms can be used to initialize, adapt, optimize or tune the “exact” algorithms

Particle filters

- Given observations Y_k ; reconstruct true states X_k

Initial distribution $p(X_0|Y_0) = P(X_0)$

Markovian evolution $p(X_k|Y_{1,\dots,k-1}) = \int p(X_0|X_{k-1})p(X_{k-1}|Y_{1,\dots,k-1})$

Bayes' rule $p(X_k|Y_{1,\dots,k}) = \frac{p(Y_k|X_k)p(X_k|Y_{1,\dots,k-1})}{p(Y_k|Y_{1,\dots,k-1})}$

- Represent posterior distribution by N weighted samples obtained from $p(X_k|Y_{1,\dots,k-1})$: $p(X_k|Y_{1,\dots,k}) \propto \sum_{i=1}^N p(Y_k|X_k)p_i(X_k)$
- Problems: impoverishment and sample size effects (if the likelihood is concentrated at the tail of the prior)

PSO PF

- Use PSO for sampling
- Standard PSO with Gaussian randomness in the velocity update (“Gaussian swarm”)
- fitness $f = \exp\left(-\frac{1}{2R_k}(y_{\text{new}} - y_{\text{pred}})^2\right)$
- R_k : observation covariance
- modulate weights: $w_i^k = w_i^{k-1} p(y^k | x_i^k)$
- Now represent posterior by weighted samples
- Avoids divergence and does well with less particles.

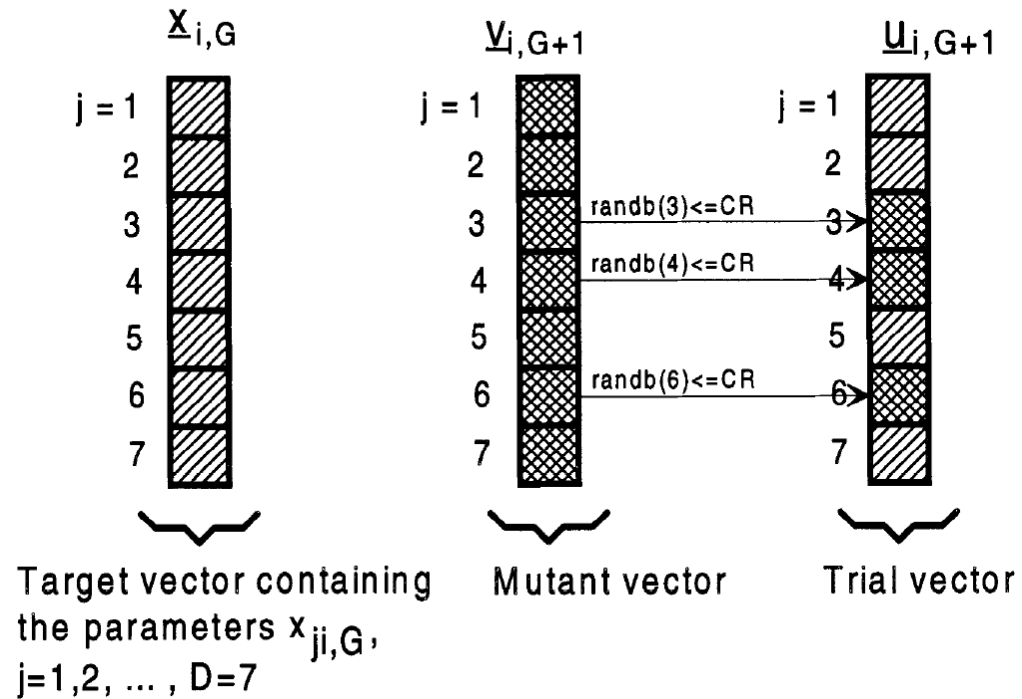
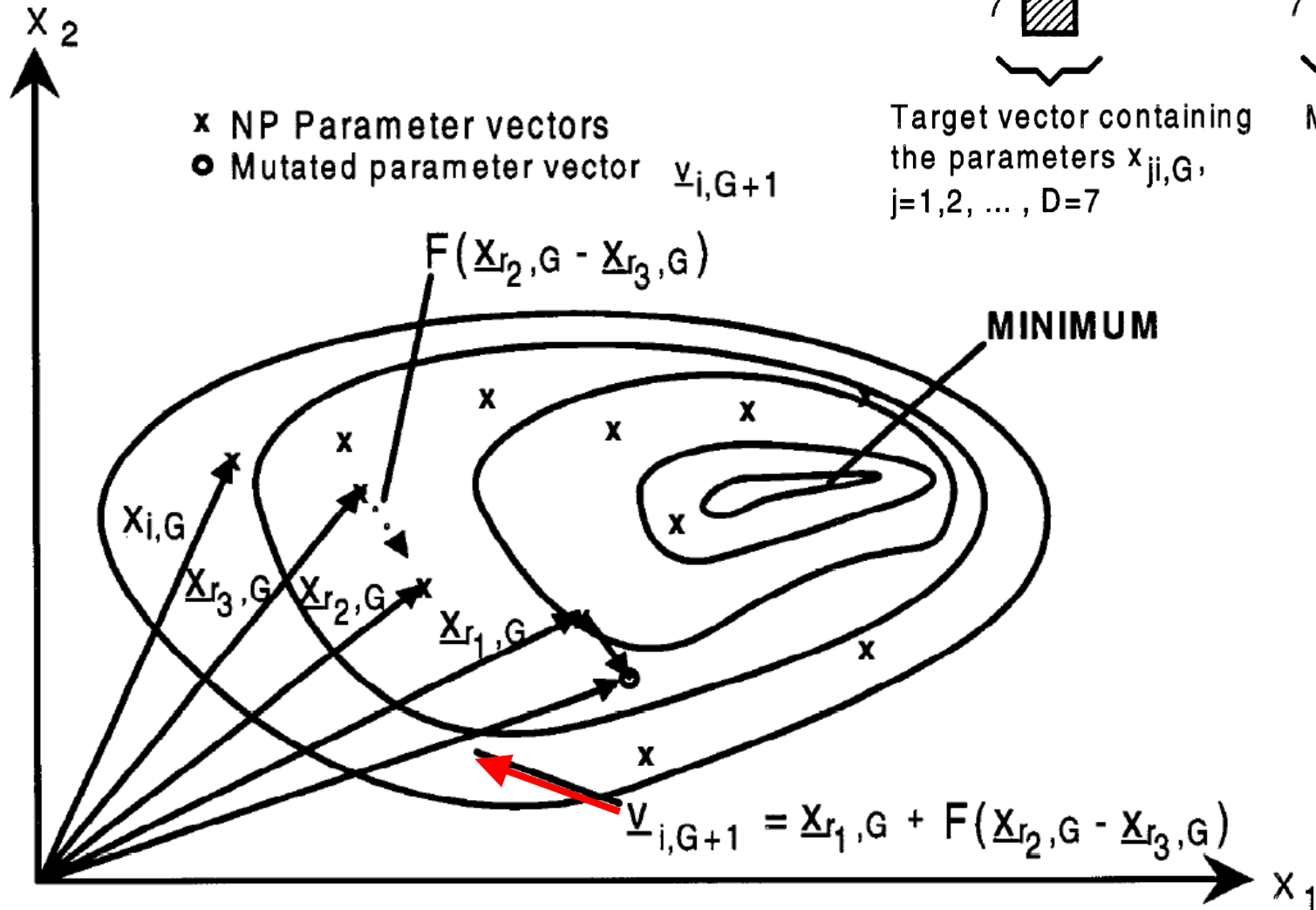
Comparison of GA and PSO

- Generally similar:
 1. Random generation of an initial population
 2. Calculate of a fitness value for each individual.
 3. Reproduction of the population based on fitness values.
 4. If requirements are met, then stop. Otherwise go back to 2.
- Modification of individuals
 - In GA: by genetic operators
 - In PSO: Particles update themselves with the internal velocity. They also have memory.
- Sharing of information
 - Mutual In GA. Whole population moves as a group towards optimal area.
 - One-way in PSO: Source of information is only gBest (or lBest). All particles tend to converge to the best solution quickly.
- Representation
 - GA: discrete
 - PS: continuous

Differential Evolution

- NP D -dimensional parameter vectors
 $x_{iG}; i = 1, 2, \dots, NP; G$: generation counter
- **Mutation:** $v_{iG+1} = x_{r_1G} + F * (x_{r_2G} - x_{r_3G});$
- F in $[0,2]$ amplification of the differential variation
- r_i random indexes different from I (“*rnbr*”)
- **Crossover:** $u_{i,G+1} = (u_{1i,G+1}, u_{2i,G+1}, \dots, u_{Di,G+1})$
- $$u_{ji,G+1} = \begin{cases} v_{ji,G+1} & \text{if } (randb(j) \leq CR) \text{ or } j = rnbr(i) \\ x_{ji,G} & \text{if } (randb(j) > CR) \text{ and } j \neq rnbr(i) \end{cases}$$
$$j = 1, 2, \dots, D.$$
- $randb$ in $[0,1]$
- **Selection:** $x_{iG+1} = u_{iG+1}$ if u_{iG+1} is better, otherwise $x_{iG+1} = x_{iG}$

Differential Evolution



DE: Details

- Properties
 - Simple, very fast
 - Reasonably good results
 - Diversity increases in flat regions (divergence property)
- Parameters
 - NP=5D (4 ... 10D)
 - F=0.5 (0.4 1.0)
 - CR=0.1 (0 ... 1.0)

DE: Variants

DE/x/y/z

x specifies the vector to be mutated which currently can be “rand” (a randomly chosen population vector) or “best” (the vector of lowest cost from the current population).

y is the number of difference vectors used.

z denotes the crossover scheme. The current variant is “bin” (Crossover due to independent binomial experiments as explained in Section 2)

e.g. DE/best/2/bin

$$v_{i,G+1} = x_{best,G} + F \cdot (x_{r_1,G} + x_{r_2,G} - x_{r_3,G} - x_{r_4,G})$$

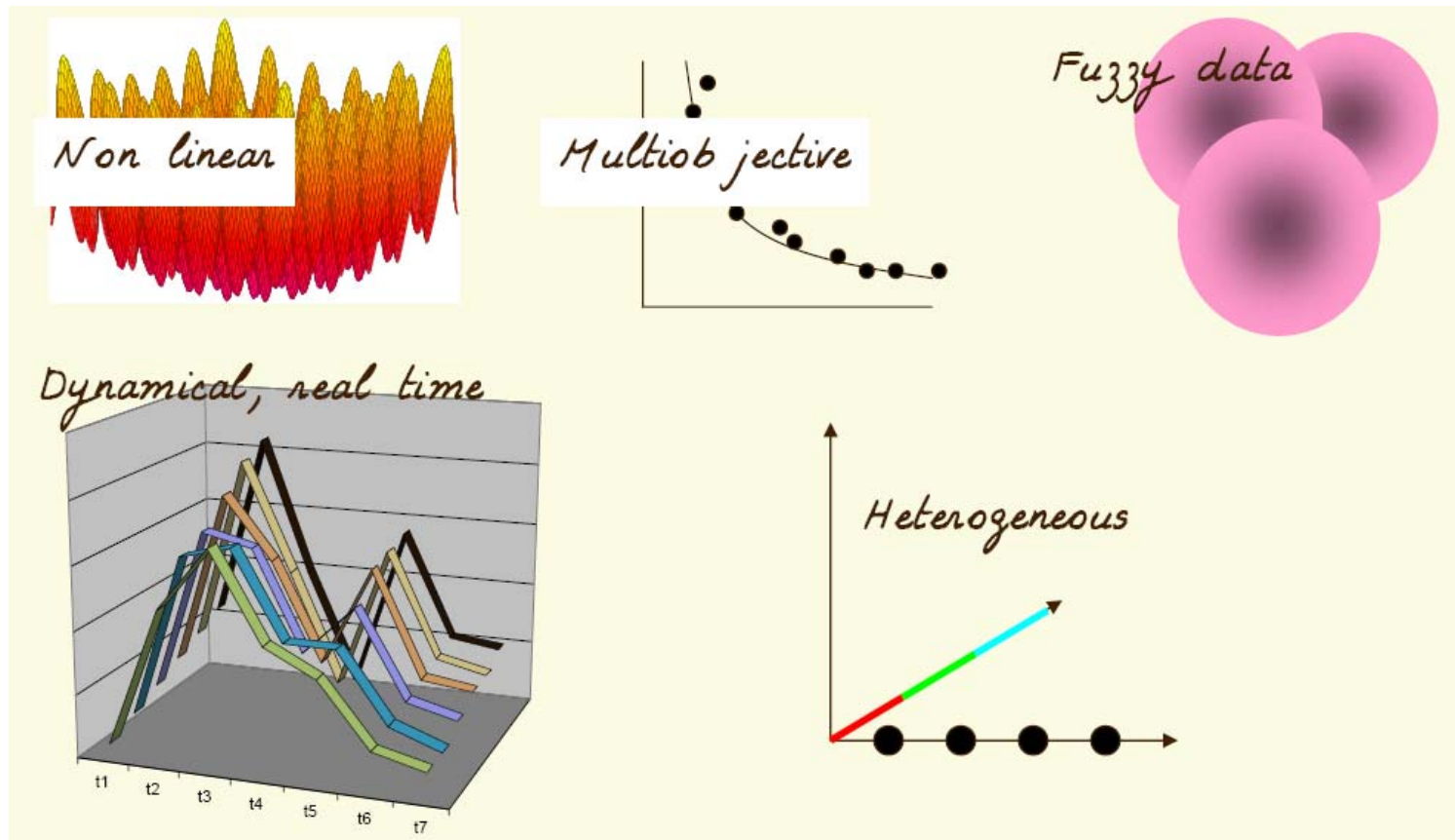
The General Scheme

1. Use **populations** of solutions/trials/individuals
2. **Transfer information** in the population from the best individuals to others by selection+crossover/attraction
3. Maintain **diversity** by adding noise/mutations/intrinsic dynamics/amplifying differences
 - Avoid local minima (leapfrog/crossover/more noise/subpopulations/border of instability/checking success)
4. Store good solutions in **memory** as best-so-far/iteration best/individual best/elite/pheromones
5. Whenever possible, use **building blocks**/partial solutions/royal road functions
6. Use **domain knowledge** and intuition for encoding, initialization, termination, choice of the algorithm
7. Tweak the parameters, develop your own variants

*It is thanks to these eccentrics,
whose behaviour is not conform to
the one of the other bees, that all
fruits sources around the colony are
so quickly found.*

Karl von Frisch 1927

Ecological niche



Literature on swarms

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