

Introduction to Randomized Continuous Optimization

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<http://www.sigevo.org/gecco-2016/>

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GECCO '16 Companion, July 22–24, 2016, Denver, CO, USA.
ACM 978-1-4503-4322-7/16/07
<http://dx.doi.org/10.1145/2926980>



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We are happy to answer questions at any time.

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Overview

1 Problem Statement

Continuous Black-Box Optimization
Typical Difficulties

2 Stochastic Black-Box Algorithms

General Template
Invariance
Comparisons of a few DFOs

3 Zoom on Evolution Strategies

Step-size Adaptation
Covariance Matrix Adaptation

4 Evaluating Black-Box Algorithms

Displaying results and visualization
Statistics
Average Runtime
Empirical Cumulative Distribution Function (ECDF)

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Problem Statement

Continuous Domain Search/Optimization

- Task: **minimize** an **objective function** (*fitness function, loss function*) in continuous domain

$$f : \mathcal{X} \subseteq \mathbb{R}^n \rightarrow \mathbb{R}, \quad \mathbf{x} \mapsto f(\mathbf{x})$$

- Black Box** scenario (direct search scenario)



- gradients are not available or not useful
- problem domain specific knowledge is used only within the black box, e.g. within an appropriate encoding
- Search **costs**: number of function evaluations

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Problem Statement Black Box Optimization and Its Difficulties

Problem Statement

Continuous Domain Search/Optimization

- Goal
 - ▶ fast convergence to the global optimum
 - ▶ solution x with **small function value** $f(x)$ with **least search cost** ... or to a robust solution x
there are two conflicting objectives
- Typical Examples
 - ▶ shape optimization (e.g. using CFD) curve fitting, airfoils
 - ▶ model calibration biological, physical
 - ▶ parameter calibration controller, plants, images
- Problems
 - ▶ exhaustive search is infeasible
 - ▶ naive random search takes too long
 - ▶ deterministic search is not successful / takes too long

Approach: stochastic search, Evolutionary Algorithms

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Problem Statement Black Box Optimization and Its Difficulties

Objective Function Properties

We assume $f : \mathcal{X} \subset \mathbb{R}^n \rightarrow \mathbb{R}$ to be *non-linear*, *non-separable* and to have at least moderate dimensionality, say $n \not\ll 10$.
Additionally, f can be

- non-convex
- multimodal there are possibly many local optima
- non-smooth derivatives do not exist
- discontinuous, plateaus
- ill-conditioned
- noisy
- ...

Goal : cope with any of these function properties
they are related to real-world problems

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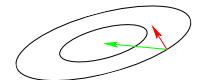
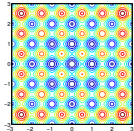
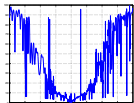
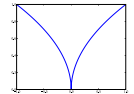
Goal: cope with any of these function properties
they are related to real-world problems

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What Makes a Function Difficult to Solve?

Why stochastic search?

- non-linear, non-quadratic, non-convex
on linear and quadratic functions much better search policies are available
- ruggedness
non-smooth, discontinuous, multimodal, and/or noisy function
- dimensionality (size of search space)
(considerably) larger than three
- non-separability
dependencies between the objective variables
- ill-conditioning

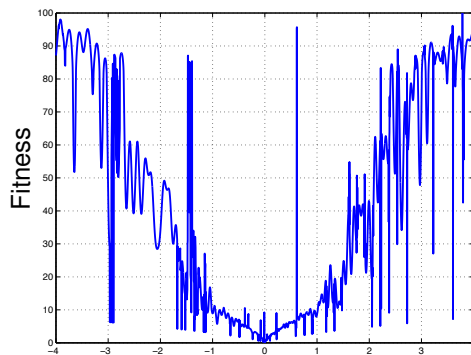


gradient direction Newton direction

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Ruggedness

non-smooth, discontinuous, multimodal, and/or noisy



cut from a 5-D example, (easily) solvable with evolution strategies

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Curse of Dimensionality

The term *Curse of dimensionality* (Richard Bellman) refers to problems caused by the **rapid increase in volume** associated with adding extra dimensions to a (mathematical) space.

Example: Consider placing 20 points equally spaced onto the interval $[0, 1]$. Now consider the 10-dimensional space $[0, 1]^{10}$. To get **similar coverage** in terms of distance between adjacent points requires $20^{10} \approx 10^{13}$ points. 20 points appear now as isolated points in a vast empty space.

Remark: **distance measures** break down in higher dimensionalities (the central limit theorem kicks in)

Consequence: a **search policy** that is valuable in small dimensions **might be useless** in moderate or large dimensional search spaces. Example: exhaustive search.

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Separable Problems

Definition (Separable Problem)

A function f is separable if

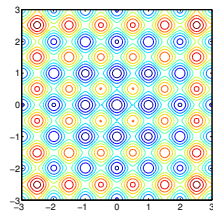
$$\arg \min_{(x_1, \dots, x_n)} f(x_1, \dots, x_n) = \left(\arg \min_{x_1} f(x_1, \dots), \dots, \arg \min_{x_n} f(\dots, x_n) \right)$$

⇒ it follows that f can be optimized in a sequence of n independent 1-D optimization processes

Example: Additively decomposable functions

$$f(x_1, \dots, x_n) = \sum_{i=1}^n f_i(x_i)$$

Rastrigin function



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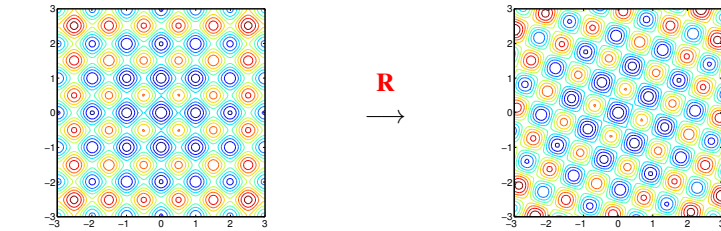
Non-Separable Problems

Building a non-separable problem from a separable one ^(1,2)

Rotating the coordinate system

- $f : \mathbf{x} \mapsto f(\mathbf{x})$ separable
- $f : \mathbf{x} \mapsto f(\mathbf{R}\mathbf{x})$ non-separable

$\mathbf{R}$ rotation matrix



¹Hansen, Ostermeier, Gawelczyk (1995). On the adaptation of arbitrary normal mutation distributions in evolution strategies: The generating set adaptation. Sixth ICGA, pp. 57-64, Morgan Kaufmann

²Salomon (1996). "Reevaluating Genetic Algorithm Performance under Coordinate Rotation of Benchmark Functions: A survey of some theoretical and practical aspects of genetic algorithms." BioSystems, 39(3):263-278

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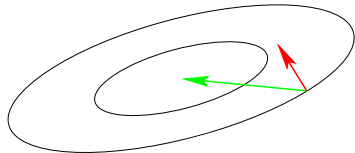
Ill-Conditioned Problems

Curvature of level sets

Consider the convex-quadratic function

$$f(\mathbf{x}) = \frac{1}{2}(\mathbf{x} - \mathbf{x}^*)^T \mathbf{H}(\mathbf{x} - \mathbf{x}^*) = \frac{1}{2} \sum_i h_{i,i} (x_i - x_i^*)^2 + \frac{1}{2} \sum_{i \neq j} h_{i,j} (x_i - x_i^*)(x_j - x_j^*)$$

$\mathbf{H}$ is Hessian matrix of f and symmetric positive definite



gradient direction $-\mathbf{f}'(\mathbf{x})^T$

Newton direction $-\mathbf{H}^{-1}\mathbf{f}'(\mathbf{x})^T$

Ill-conditioning means squeezed level sets (high curvature).
Condition number equals nine here. Condition numbers up to 10^{10}
are not unusual in real world problems.

If $\mathbf{H} \approx \mathbf{I}$ (small condition number of $\mathbf{H}$) first order information (e.g. the gradient) is sufficient. Otherwise second order information (estimation of $\mathbf{H}^{-1}$) is necessary.

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What Makes a Function Difficult to Solve?

... and what can be done

The Problem	Possible Approaches
Dimensionality	exploiting the problem structure separability, locality/neighborhood, encoding
Ill-conditioning	second order approach changes the neighborhood metric
Ruggedness	non-local policy, large sampling width (step-size) as large as possible while preserving a reasonable convergence speed population-based method, stochastic, non-elitistic recombination operator serves as repair mechanism restarts

... metaphors

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Metaphors

Evolutionary Computation

Optimization/Nonlinear Programming

individual, offspring, parent	↔	candidate solution decision variables design variables object variables
population	↔	set of candidate solutions
fitness function	↔	objective function loss function cost function error function
generation	↔	iteration

...methods: ESs

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Landscape of Continuous Black-Box Optimization

Deterministic algorithms

Quasi-Newton with estimation of gradient (BFGS) [Broyden et al. 1970]
 Simplex downhill [Nelder & Mead 1965]
 Pattern search [Hooke and Jeeves 1961]
 Trust-region methods (NEWUOA, BOBYQA) [Powell 2006, 2009]

Stochastic (randomized) search methods

Evolutionary Algorithms (continuous domain)
 Differential Evolution [Storn & Price 1997]
 Particle Swarm Optimization [Kennedy & Eberhart 1995]
Evolution Strategies, CMA-ES [Rechenberg 1965, Hansen & Ostermeier 2001]
 Estimation of Distribution Algorithms (EDAs) [Larrañaga, Lozano, 2002]
 Cross Entropy Method (same as EDA) [Rubinstein, Kroese, 2004]
Genetic Algorithms [Holland 1975, Goldberg 1989]
 Simulated annealing [Kirkpatrick et al. 1983]
 Simultaneous perturbation stochastic approximation (SPSA) [Spall 2000]

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Stochastic Search

A black box search template to minimize $f : \mathbb{R}^n \rightarrow \mathbb{R}$

Initialize distribution parameters θ , set population size $\lambda \in \mathbb{N}$

While not terminate

- ❶ Sample distribution $P(x|\theta) \rightarrow x_1, \dots, x_\lambda \in \mathbb{R}^n$
- ❷ Evaluate $x_1, \dots, x_\lambda$ on f
- ❸ Update parameters $\theta \leftarrow F_\theta(\theta, x_1, \dots, x_\lambda, f(x_1), \dots, f(x_\lambda))$

Everything depends on the definition of P and F_θ

deterministic algorithms are covered as well

In many Evolutionary Algorithms the distribution P is implicitly defined via **operators on a population**, in particular, selection, recombination and mutation

Natural template for (incremental) *Estimation of Distribution Algorithms*

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Examples

- ① Estimation of Distribution Algorithms
- ② Evolution Strategies
- ③ Differential Evolution
- ④ Particle Swarm Optimization

all those methods are comparison based

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Evolution Strategies

New search points are sampled normally distributed

$$\mathbf{x}_i \sim \mathbf{m} + \sigma \mathcal{N}(\mathbf{0}, \mathbf{C}) \quad \text{for } i = 1, \dots, \lambda$$

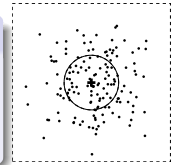
as perturbations of $\mathbf{m}$, where $\mathbf{x}_i, \mathbf{m} \in \mathbb{R}^n$, $\sigma \in \mathbb{R}_+$, $\mathbf{C} \in \mathbb{R}^{n \times n}$

where

- the **mean** vector $\mathbf{m} \in \mathbb{R}^n$ represents the favorite solution
- the so-called **step-size** $\sigma \in \mathbb{R}_+$ controls the *step length*
- the **covariance matrix** $\mathbf{C} \in \mathbb{R}^{n \times n}$ determines the **shape** of the distribution ellipsoid

here, all new points are sampled with the same parameters

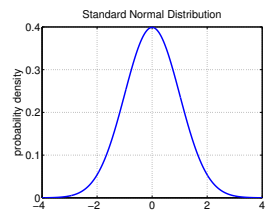
The question remains how to update $\mathbf{m}$, $\mathbf{C}$, and σ .



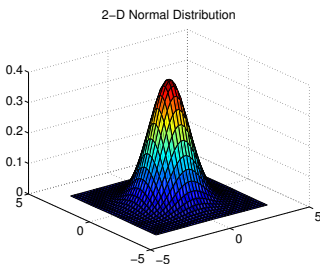
Navigation icons: back, forward, search, etc.

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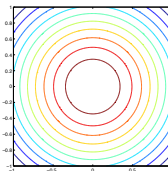
Normal Distribution



probability density of the 1-D standard normal distribution



probability density of a 2-D normal distribution



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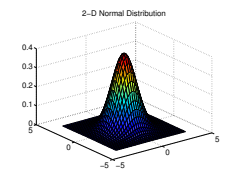
Navigation icons: back, forward, search, etc.

The Multi-Variate (n -Dimensional) Normal Distribution

Any multi-variate normal distribution $\mathcal{N}(\mathbf{m}, \mathbf{C})$ is uniquely determined by its mean value $\mathbf{m} \in \mathbb{R}^n$ and its symmetric positive definite $n \times n$ covariance matrix $\mathbf{C}$.

The **mean** value $\mathbf{m}$

- determines the displacement (translation)
- value with the largest density (modal value)
- the distribution is symmetric about the distribution mean



The **covariance matrix** $\mathbf{C}$

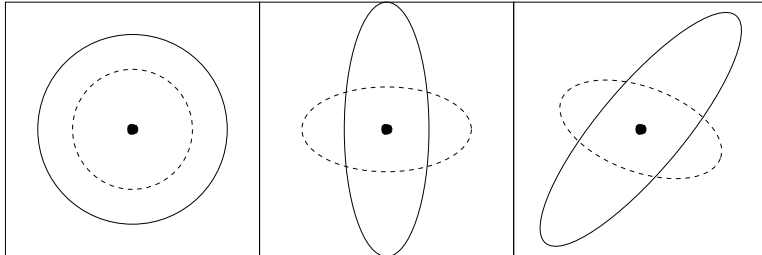
- determines the shape
- **geometrical interpretation**: any covariance matrix can be uniquely identified with the iso-density ellipsoid $\{\mathbf{x} \in \mathbb{R}^n \mid (\mathbf{x} - \mathbf{m})^T \mathbf{C}^{-1} (\mathbf{x} - \mathbf{m}) = 1\}$

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Navigation icons: back, forward, search, etc.

... any **covariance matrix** can be uniquely identified with the iso-density ellipsoid
 $\{x \in \mathbb{R}^n \mid (x - m)^T C^{-1} (x - m) = 1\}$

Lines of Equal Density



$\mathcal{N}(m, \sigma^2 \mathbf{I}) \sim m + \sigma \mathcal{N}(\mathbf{0}, \mathbf{I})$
 one degree of freedom σ
 components are
 independent standard
 normally distributed

$\mathcal{N}(m, \mathbf{D}^2) \sim m + \mathbf{D} \mathcal{N}(\mathbf{0}, \mathbf{I})$
 n degrees of freedom
 components are
 independent, scaled

$\mathcal{N}(m, \mathbf{C}) \sim m + \mathbf{C}^{\frac{1}{2}} \mathcal{N}(\mathbf{0}, \mathbf{I})$
 $(n^2 + n)/2$ degrees of freedom
 components are
 correlated

where $\mathbf{I}$ is the identity matrix (isotropic case) and $\mathbf{D}$ is a diagonal matrix (reasonable for separable problems) and $\mathbf{A} \times \mathcal{N}(\mathbf{0}, \mathbf{I}) \sim \mathcal{N}(\mathbf{0}, \mathbf{A}\mathbf{A}^T)$ holds for all $\mathbf{A}$.

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The $(\mu/\mu, \lambda)$ -ES

Non-elitist selection and intermediate (weighted) recombination

Given the i -th solution point $x_i = m + \sigma \underbrace{\mathcal{N}(\mathbf{0}, \mathbf{C})}_{=: y_i} = m + \sigma y_i$

Let $x_{i:\lambda}$ the i -th ranked solution point, such that $f(x_{1:\lambda}) \leq \dots \leq f(x_{\lambda:\lambda})$.
 The new mean reads

$$m \leftarrow \sum_{i=1}^{\mu} w_i x_{i:\lambda} = m + \sigma \underbrace{\sum_{i=1}^{\mu} w_i y_{i:\lambda}}_{=: y_w}$$

where

$$w_1 \geq \dots \geq w_{\mu} > 0, \quad \sum_{i=1}^{\mu} w_i = 1, \quad \frac{1}{\sum_{i=1}^{\mu} w_i^2} =: \mu_w \approx \frac{\lambda}{4}$$

The best μ points are selected from the new solutions (non-elitistic) and weighted intermediate recombination is applied.

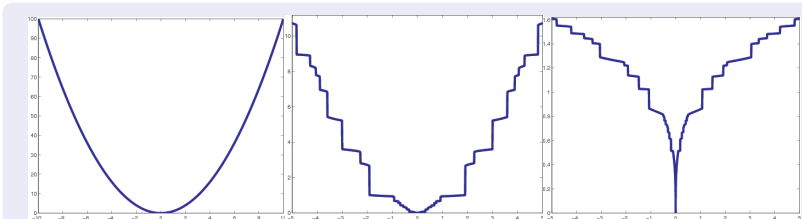
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Invariance Under Monotonically Increasing Functions

Rank-based algorithms

Update of all parameters uses only the ranks

$$f(x_{1:\lambda}) \leq f(x_{2:\lambda}) \leq \dots \leq f(x_{\lambda:\lambda})$$



$$g(f(x_{1:\lambda})) \leq g(f(x_{2:\lambda})) \leq \dots \leq g(f(x_{\lambda:\lambda})) \quad \forall g$$

g is strictly monotonically increasing
 g preserves ranks

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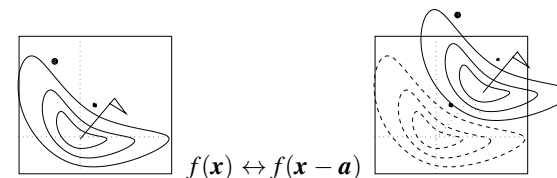
Whitley 1989. The GENITOR algorithm and selection pressure: Why rank-based allocation of reproductive trials is best, ICGA

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Basic Invariance in Search Space

- translation invariance

is true for most optimization algorithms



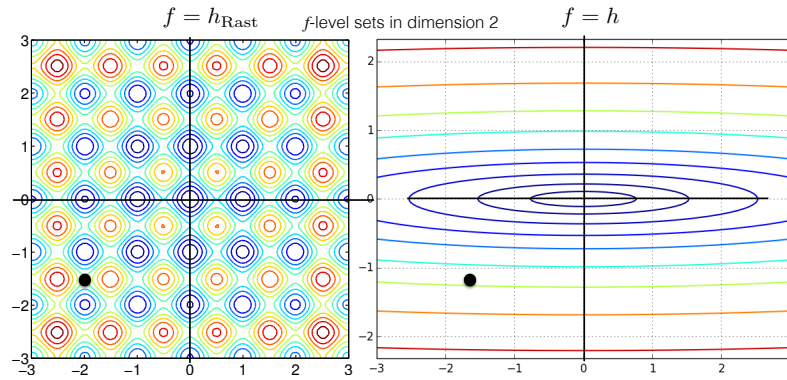
Identical behavior on f and f_a

$$\begin{aligned} f &: x \mapsto f(x), & x^{(t=0)} &= x_0 \\ f_a &: x \mapsto f(x-a), & x^{(t=0)} &= x_0 + a \end{aligned}$$

No difference can be observed w.r.t. the argument of f

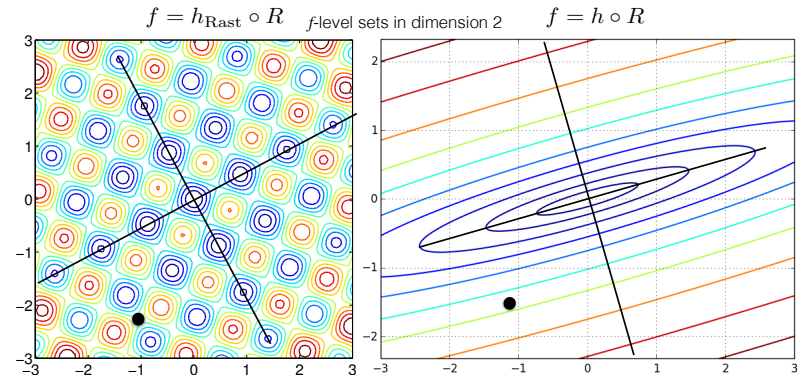
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Invariance Under Rigid Search Space Transformations



for example, invariance under search space rotation
(separable $\Leftrightarrow$ non-separable)

Invariance Under Rigid Search Space Transformations



for example, invariance under search space rotation
(separable $\Leftrightarrow$ non-separable)

Invariance

The grand aim of all science is to cover the greatest number of empirical facts by logical deduction from the smallest number of hypotheses or axioms.
— Albert Einstein

Empirical performance results

- from benchmark functions
- from solved real world problems

are only useful if they do generalize to other problems

Invariance is a strong non-empirical statement about generalization

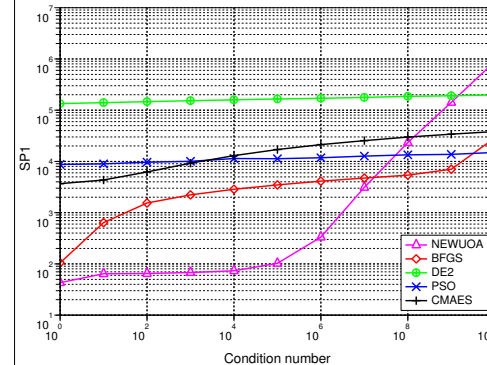
generalizing (identical) performance from a single function to a whole class of functions

consequently, invariance is important for the evaluation of search algorithms

Comparison to BFGS, NEWUOA, PSO and DE

f convex quadratic, separable with varying condition number α

Ellipsoid dimension 20, 21 trials, tolerance $1e-09$, eval max $1e+07$



BFGS (Broyden et al 1970)
NEWUOA (Powell 2004)
DE (Storn & Price 1996)
PSO (Kennedy & Eberhart 1995)
CMA-ES (Hansen & Ostermeier 2001)

$f(x) = g(x^T H x)$ with
 H diagonal
 g identity (for **BFGS** and **NEWUOA**)
 g any order-preserving = strictly increasing function (for all other)

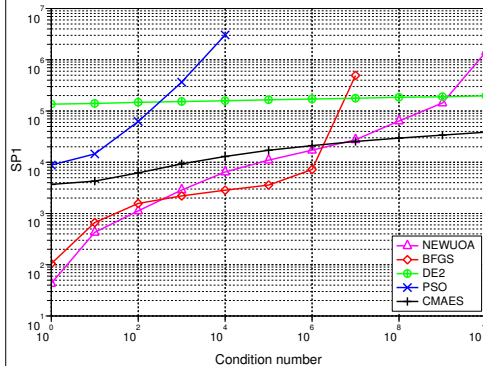
SP1 = average number of objective function evaluations¹⁴ to reach the target function value of $g^{-1}(10^{-9})$

¹⁴ Auger et al. (2009): Experimental comparisons of derivative free optimization algorithms, SEA

Comparison to BFGS, NEWUOA, PSO and DE

f convex quadratic, non-separable (rotated) with varying condition number α

Rotated Ellipsoid dimension 20, 21 trials, tolerance $1e-09$, eval max $1e+07$



BFGS (Broyden et al 1970)
NEWUOA (Powell 2004)
DE (Storn & Price 1996)
PSO (Kennedy & Eberhart 1995)
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$f(x) = g(x^T H x)$ with
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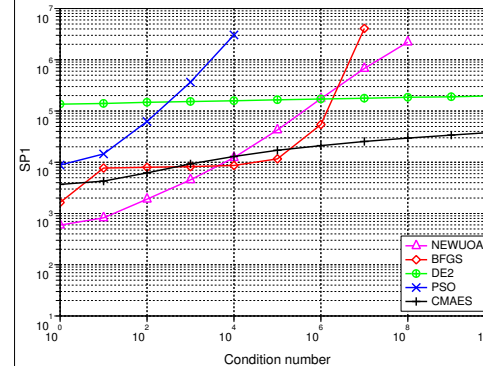
SP1 = average number of objective function evaluations¹⁵ to reach the target function value of $g^{-1}(10^{-9})$

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Comparison to BFGS, NEWUOA, PSO and DE

f non-convex, non-separable (rotated) with varying condition number α

Sqrt of sqrt of rotated ellipsoid dimension 20, 21 trials, tolerance $1e-09$, eval max $1e+07$



BFGS (Broyden et al 1970)
NEWUOA (Powell 2004)
DE (Storn & Price 1996)
PSO (Kennedy & Eberhart 1995)
CMA-ES (Hansen & Ostermeier 2001)

$f(x) = g(x^T H x)$ with
 H full
 $g : x \mapsto x^{1/4}$ (for BFGS and NEWUOA)
 g any order-preserving = strictly increasing function (for all other)

SP1 = average number of objective function evaluations¹⁶ to reach the target function value of $g^{-1}(10^{-9})$

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Overview

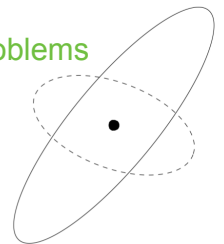
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Zoom on ESs: Objectives

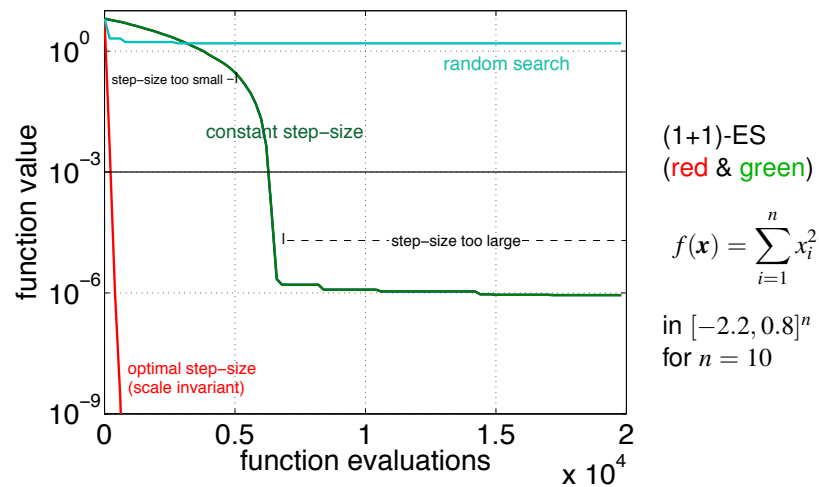
Illustrate **why** and **how** sampling distribution is controlled

step-size control (overall standard deviation)
allows to achieve linear convergence

covariance matrix control
allows to solve ill-conditioned problems



Why Step-Size Control?



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Methods for Step-Size Control

- **1/5-th success rule^{ab}**, often applied with “+”-selection
increase step-size if more than 20% of the new solutions are successful,
decrease otherwise
- **σ -self-adaptation^c**, applied with “-”-selection
mutation is applied to the step-size and the better, according to the
objective function value, is selected

simplified “global” self-adaptation
- **path length control^d** (Cumulative Step-size Adaptation, CSA)^e
self-adaptation derandomized and non-localized

^aRechenberg 1973, *Evolutionsstrategie, Optimierung technischer Systeme nach Prinzipien der biologischen Evolution*, Frommann-Holzboog

^bSchumer and Steiglitz 1968, Adaptive step size random search. *IEEE TAC*

^cSchwefel 1981, *Numerical Optimization of Computer Models*, Wiley

^dHansen & Ostermeier 2001, Completely Derandomized Self-Adaptation in Evolution Strategies, *Evol. Comput.*

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^eOstermeier et al 1994, Step-size adaptation based on non-local use of selection information, *PPSN IV*

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Path Length Control (CSA)

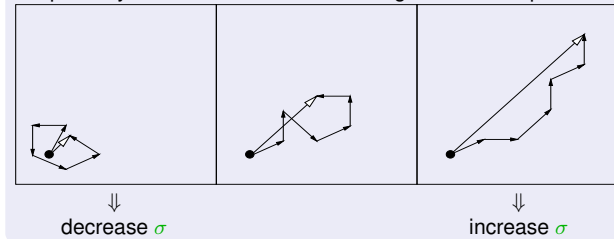
The Concept of Cumulative Step-Size Adaptation

$$\mathbf{x}_i = \mathbf{m} + \sigma \mathbf{y}_i$$

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w$$

Measure the length of the *evolution path*

the pathway of the mean vector $\mathbf{m}$ in the generation sequence



loosely speaking steps are

- perpendicular under random selection (in expectation)
- perpendicular in the desired situation (to be most efficient)

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Path Length Control (CSA)

The Equations

Initialize $\mathbf{m} \in \mathbb{R}^n$, $\sigma \in \mathbb{R}_+$, evolution path $\mathbf{p}_\sigma = \mathbf{0}$,
set $c_\sigma \approx 4/n$, $d_\sigma \approx 1$.

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w \quad \text{where } \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda} \quad \text{update mean}$$

$$\mathbf{p}_\sigma \leftarrow (1 - c_\sigma) \mathbf{p}_\sigma + \underbrace{\sqrt{1 - (1 - c_\sigma)^2}}_{\text{accounts for } 1 - c_\sigma} \underbrace{\sqrt{\mu_w}}_{\text{accounts for } w_i} \mathbf{y}_w$$

$$\sigma \leftarrow \sigma \times \underbrace{\exp\left(\frac{c_\sigma}{d_\sigma} \left(\frac{\|\mathbf{p}_\sigma\|}{\mathbb{E}\|\mathcal{N}(\mathbf{0}, \mathbf{I})\|} - 1\right)\right)}_{>1 \iff \|\mathbf{p}_\sigma\| \text{ is greater than its expectation}} \quad \text{update step-size}$$

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Path Length Control (CSA)

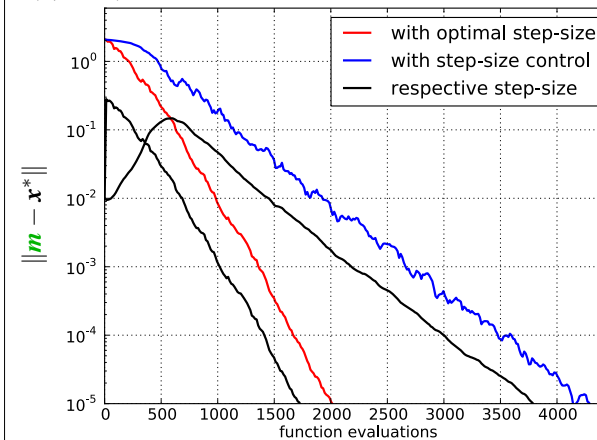
The Equations

Initialize $\mathbf{m} \in \mathbb{R}^n$, $\sigma \in \mathbb{R}_+$, evolution path $\mathbf{p}_\sigma = \mathbf{0}$,
set $c_\sigma \approx 4/n$, $d_\sigma \approx 1$.

$$\begin{aligned} \mathbf{m} &\leftarrow \mathbf{m} + \sigma \mathbf{y}_w \quad \text{where } \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda} && \text{update mean} \\ \mathbf{p}_\sigma &\leftarrow (1 - c_\sigma) \mathbf{p}_\sigma + \underbrace{\sqrt{1 - (1 - c_\sigma)^2}}_{\text{accounts for } 1 - c_\sigma} \underbrace{\sqrt{\mu_w}}_{\text{accounts for } w_i} \mathbf{y}_w \\ \sigma &\leftarrow \sigma \times \exp \left(\underbrace{\frac{c_\sigma}{d_\sigma} \left(\frac{\|\mathbf{p}_\sigma\|}{\mathbb{E}\|\mathcal{N}(\mathbf{0}, \mathbf{I})\|} - 1 \right)}_{>1 \iff \|\mathbf{p}_\sigma\| \text{ is greater than its expectation}} \right) && \text{update step-size} \end{aligned}$$

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(5/5, 10)-CSA-ES, default parameters



$$f(\mathbf{x}) = \sum_{i=1}^n x_i^2$$

in $[-0.2, 0.8]^n$
for $n = 30$

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Evolution Strategies

Recalling

New search points are sampled normally distributed

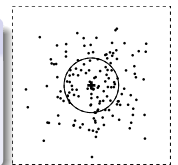
$$\mathbf{x}_i \sim \mathbf{m} + \sigma \mathcal{N}_i(\mathbf{0}, \mathbf{C}) \quad \text{for } i = 1, \dots, \lambda$$

as perturbations of $\mathbf{m}$, where $\mathbf{x}_i, \mathbf{m} \in \mathbb{R}^n$, $\sigma \in \mathbb{R}_+$, $\mathbf{C} \in \mathbb{R}^{n \times n}$

where

- the **mean** vector $\mathbf{m} \in \mathbb{R}^n$ represents the favorite solution
- the so-called **step-size** $\sigma \in \mathbb{R}_+$ controls the **step length**
- the **covariance matrix** $\mathbf{C} \in \mathbb{R}^{n \times n}$ determines the **shape** of the distribution ellipsoid

The remaining question is how to update $\mathbf{C}$.

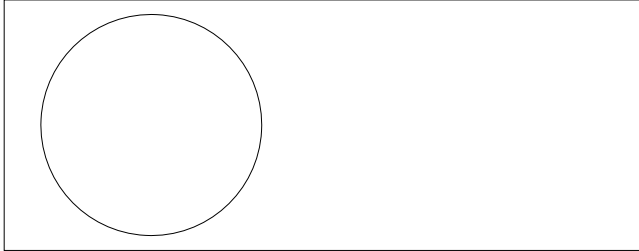


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

initial distribution, $\mathbf{C} = \mathbf{I}$

...equations

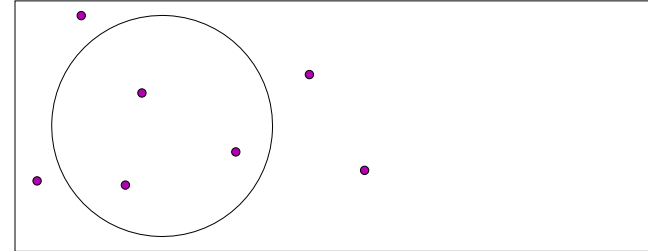


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

initial distribution, $\mathbf{C} = \mathbf{I}$

...equations

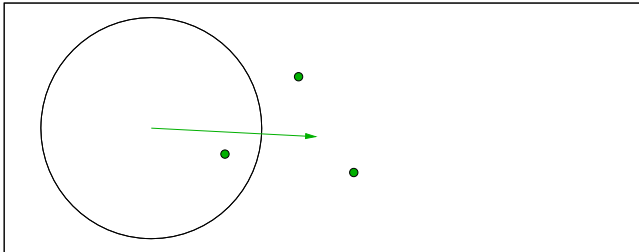


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

 $\mathbf{y}_w$, movement of the population mean $\mathbf{m}$ (disregarding σ)

...equations

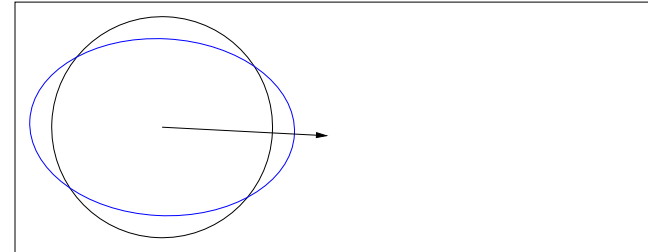


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

mixture of distribution $\mathbf{C}$ and step $\mathbf{y}_w$,
 $\mathbf{C} \leftarrow 0.8 \times \mathbf{C} + 0.2 \times \mathbf{y}_w \mathbf{y}_w^T$

...equations

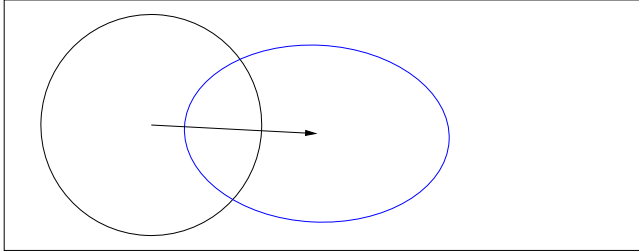


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

new distribution (disregarding σ)

...equations

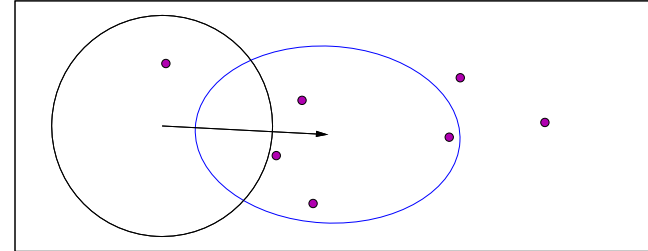


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

new distribution (disregarding σ)

...equations

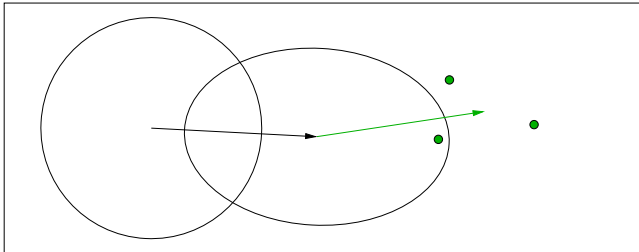


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

movement of the population mean $\mathbf{m}$

...equations

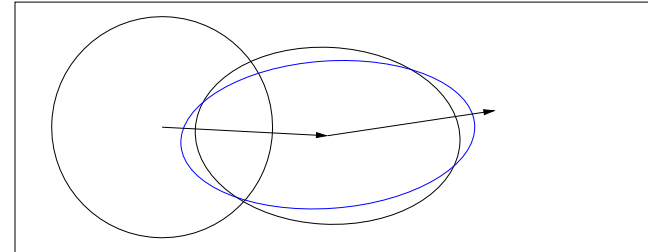


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$

mixture of distribution $\mathbf{C}$ and step $\mathbf{y}_w$,
 $\mathbf{C} \leftarrow 0.8 \times \mathbf{C} + 0.2 \times \mathbf{y}_w \mathbf{y}_w^T$

...equations

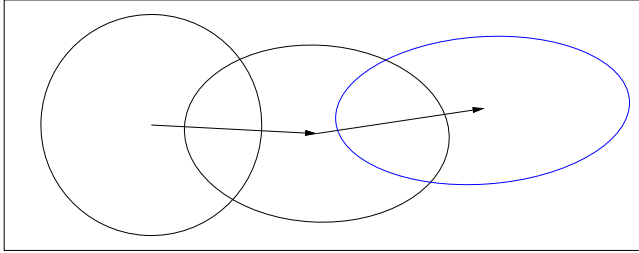


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Covariance Matrix Adaptation

Rank-One Update

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w, \quad \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C})$$



new distribution,

$$\mathbf{C} \leftarrow 0.8 \times \mathbf{C} + 0.2 \times \mathbf{y}_w \mathbf{y}_w^T$$

the ruling principle: the adaptation **increases the likelihood of successful steps**, $\mathbf{y}_w$, to appear again

another viewpoint: the adaptation **follows a natural gradient** approximation of the expected fitness

... equations

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Covariance Matrix Adaptation

Rank-One Update

Initialize $\mathbf{m} \in \mathbb{R}^n$, and $\mathbf{C} = \mathbf{I}$, set $\sigma = 1$, learning rate $c_{\text{cov}} \approx 2/n^2$

While not terminate

$$\mathbf{x}_i = \mathbf{m} + \sigma \mathbf{y}_i, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C}),$$

$$\mathbf{m} \leftarrow \mathbf{m} + \sigma \mathbf{y}_w \quad \text{where } \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda}$$

$$\mathbf{C} \leftarrow (1 - c_{\text{cov}}) \mathbf{C} + c_{\text{cov}} \underbrace{\mu_w \mathbf{y}_w \mathbf{y}_w^T}_{\text{rank-one}} \quad \text{where } \mu_w = \frac{1}{\sum_{i=1}^{\mu} w_i^2} \geq 1$$

The rank-one update has been found independently in several domains^{6 7 8 9}

⁶ Kjellström & Taxén 1981. Stochastic Optimization in System Design, IEEE TCS

⁷ Hansen & Ostermeier 1996. Adapting arbitrary normal mutation distributions in evolution strategies: The covariance matrix adaptation, ICEC

⁸ Jung 1999. System Identification: Theory for the User

⁹ Haario et al 2001. An adaptive Metropolis algorithm, JSTOR

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The CMA-ES

Input: $\mathbf{m} \in \mathbb{R}^n$, $\sigma \in \mathbb{R}_+$, λ

Initialize: $\mathbf{C} = \mathbf{I}$, and $\mathbf{p}_c = \mathbf{0}$, $\mathbf{p}_\sigma = \mathbf{0}$,

Set: $c_c \approx 4/n$, $c_\sigma \approx 4/n$, $c_1 \approx 2/n^2$, $c_\mu \approx \mu_w/n^2$, $c_1 + c_\mu \leq 1$, $d_\sigma \approx 1 + \sqrt{\frac{\mu_w}{n}}$, and $w_{i=1 \dots \lambda}$ such that $\mu_w = \frac{1}{\sum_{i=1}^{\mu} w_i^2} \approx 0.3 \lambda$

While not terminate

$$\mathbf{x}_i = \mathbf{m} + \sigma \mathbf{y}_i, \quad \mathbf{y}_i \sim \mathcal{N}(\mathbf{0}, \mathbf{C}), \quad \text{for } i = 1, \dots, \lambda \quad \text{sampling}$$

$$\mathbf{m} \leftarrow \sum_{i=1}^{\mu} w_i \mathbf{x}_{i:\lambda} = \mathbf{m} + \sigma \mathbf{y}_w \quad \text{where } \mathbf{y}_w = \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda} \quad \text{update mean}$$

$$\mathbf{p}_c \leftarrow (1 - c_c) \mathbf{p}_c + \mathbb{1}_{\{\|\mathbf{p}_c\| < 1.5\sqrt{n}\}} \sqrt{1 - (1 - c_c)^2} \sqrt{\mu_w} \mathbf{y}_w \quad \text{cumulation for } \mathbf{C}$$

$$\mathbf{p}_\sigma \leftarrow (1 - c_\sigma) \mathbf{p}_\sigma + \sqrt{1 - (1 - c_\sigma)^2} \sqrt{\mu_w} \mathbf{C}^{-\frac{1}{2}} \mathbf{y}_w \quad \text{cumulation for } \sigma$$

$$\mathbf{C} \leftarrow (1 - c_1 - c_\mu) \mathbf{C} + c_1 \mathbf{p}_c \mathbf{p}_c^T + c_\mu \sum_{i=1}^{\mu} w_i \mathbf{y}_{i:\lambda} \mathbf{y}_{i:\lambda}^T \quad \text{update } \mathbf{C}$$

$$\sigma \leftarrow \sigma \times \exp\left(\frac{c_\sigma}{d_\sigma} \left(\frac{\|\mathbf{p}_\sigma\|}{\mathbb{E}\|\mathcal{N}(\mathbf{0}, \mathbf{I})\|} - 1\right)\right) \quad \text{update of } \sigma$$

Not covered on this slide: termination, restarts, useful output, boundaries and encoding

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Experimentum Crucis (0)

What did we want to achieve?

- reduce any convex-quadratic function

$$f(\mathbf{x}) = \mathbf{x}^T \mathbf{H} \mathbf{x}$$

$$\text{e.g. } f(\mathbf{x}) = \sum_{i=1}^n 10^6 \frac{i-1}{n-1} x_i^2$$

to the sphere model

$$f(\mathbf{x}) = \mathbf{x}^T \mathbf{x}$$

without use of derivatives

- lines of equal density align with lines of equal fitness

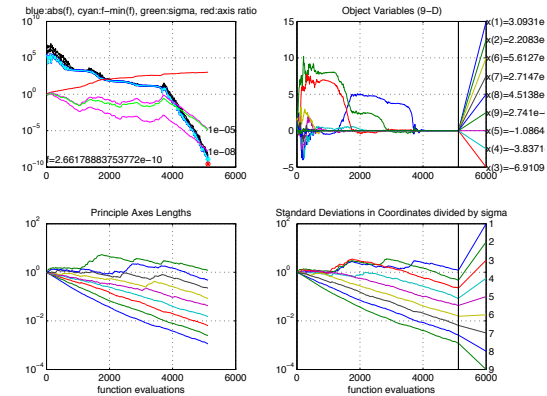
$$\mathbf{C} \propto \mathbf{H}^{-1}$$

in a stochastic sense

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Experimentum Crucis (1)

f convex quadratic, separable

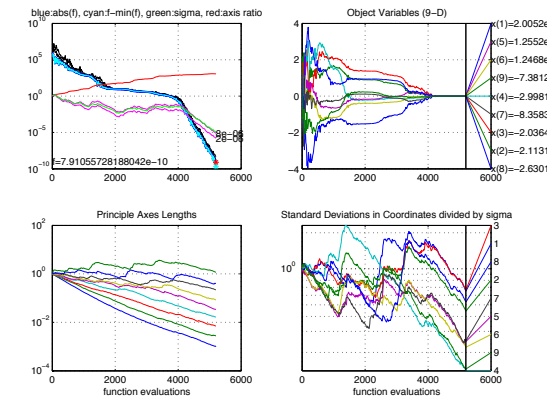


$$f(\mathbf{x}) = \sum_{i=1}^n 10^{\alpha \frac{i-1}{n-1}} x_i^2, \alpha = 6$$

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Experimentum Crucis (2)

f convex quadratic, as before but non-separable (rotated)



$$f(\mathbf{x}) = g(\mathbf{x}^T \mathbf{H} \mathbf{x}), g: \mathbb{R} \rightarrow \mathbb{R} \text{ strictly increasing}$$

$$C \propto H^{-1} \text{ for all } g, \mathbf{H}$$

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Evaluation of Anytime Black-Box Optimizers

Particularly Randomized Search Algorithms

Randomized optimization is mostly an empirical science

Hence it is **crucial to properly conduct** numerical experiments and assess performance

to **not fool oneself** on what our favorite algorithm is good/not good at

in order to **not fool others** ...

"The first principle is that you must not fool yourself and you are the easiest person to fool."

Richard P. Feynman

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Evaluation of Anytime Black-Box Optimizers

Particularly Randomized Search Algorithms

Evaluation of performance in a **broad sense**

not only be able to say “Algorithm A is better than B”

but

understand where and why algorithm work

quantify performance

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Measuring Performance

Empirically

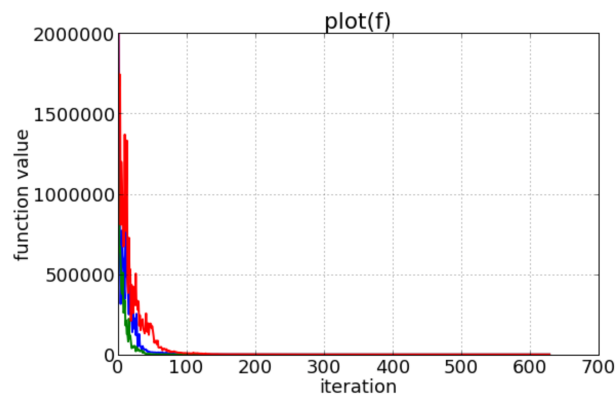
convergence graphs is all we have to start with

having the right presentation is important
too often neglected

the details are important

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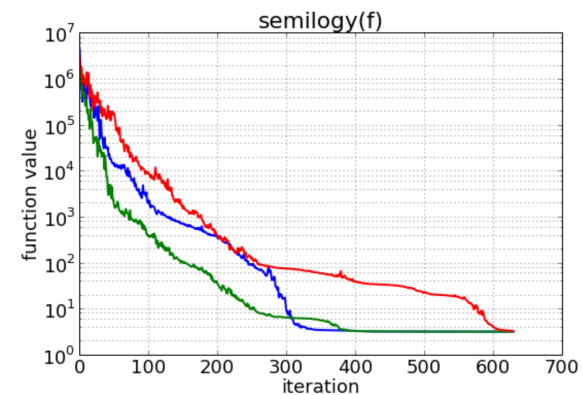
Displaying Three Runs



not like this (it's unfortunately a common picture)

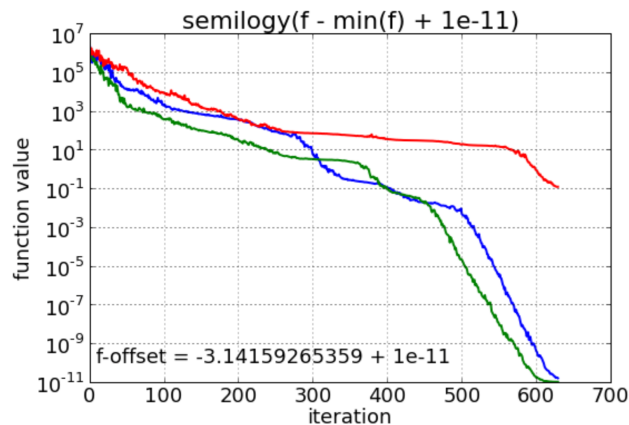
63

Displaying Three Runs



better like this (shown are the same data),
caveat: fails with negative f-values

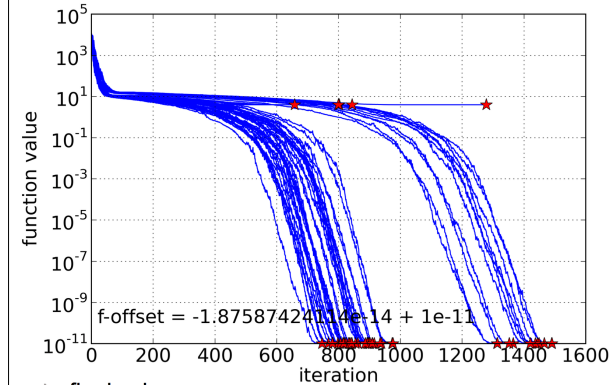
Displaying Three Runs



even better like this: subtract minimum value over all runs

Displaying 51 Runs

don't hesitate to display all data (the appendix is your friend)



★ : final value

observation: three different "modes", which would be difficult to represent or recover in single statistics

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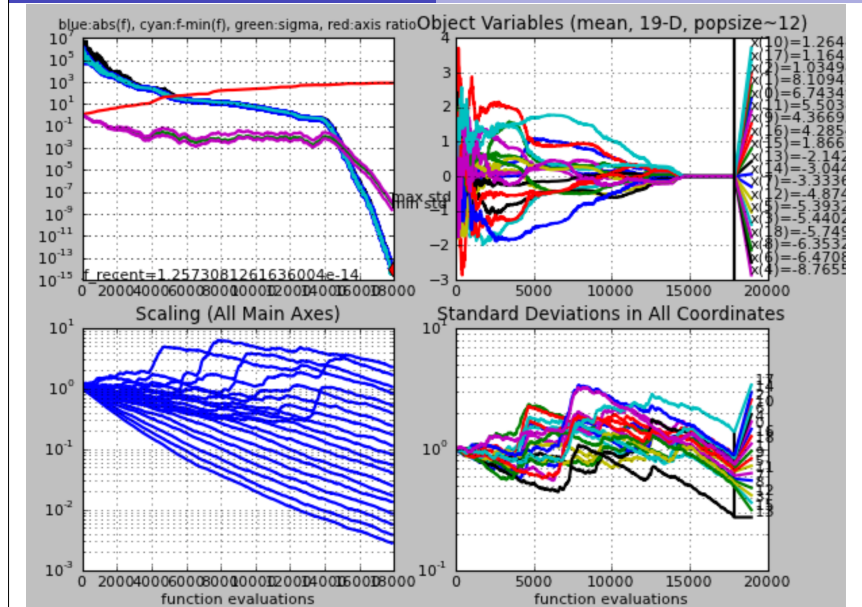
Visualization

Other data than convergence graphs can be very instructive to visualize to understand performance

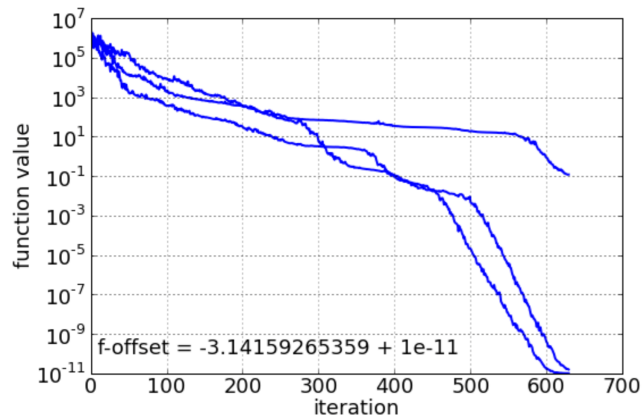
Example:

visualization of evolution of distribution in CMA-ES
(see next slide)

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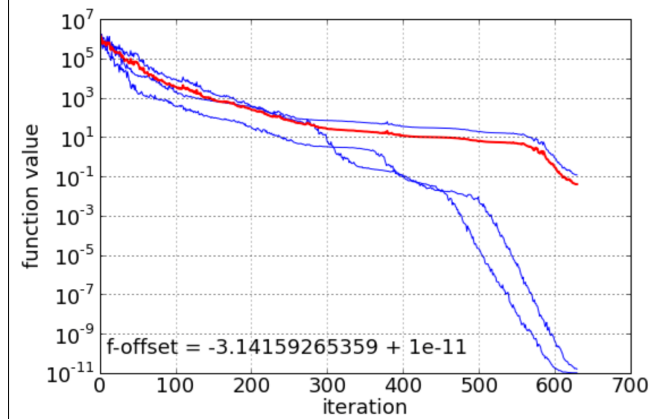


Which Statistics?



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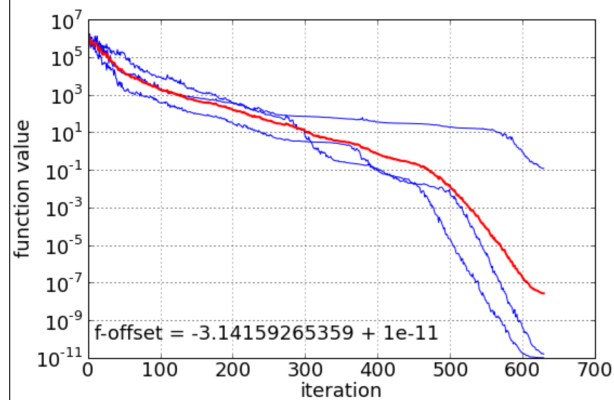
Which Statistics?



mean/average function value

- tends to emphasize large values

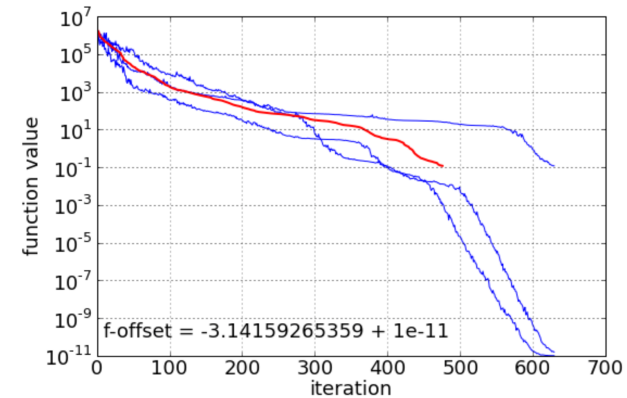
Which Statistics?



geometric average function value $\exp(\text{mean}_i(\log(f_i))) = (\prod_{i=1}^N f_i)^{1/N}$

- reflects "visual" average
- depends on offset
- artefact due to adding $1e-11$

Which Statistics?

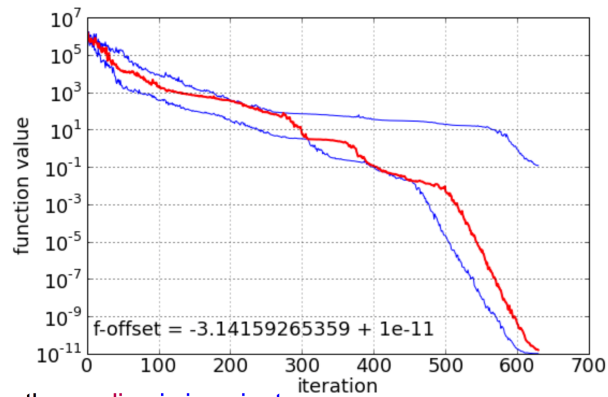


average iterations

- reflects "visual" average
- here: incomplete

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Which Statistics?



the **median** is **invariant**

- unique for uneven number of data
- independent of log-scale, offset...

$$\text{median}(\log(\text{data})) = \log(\text{median}(\text{data}))$$

- same when taken over x- or y-direction

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Implication

use the **median** as summary datum

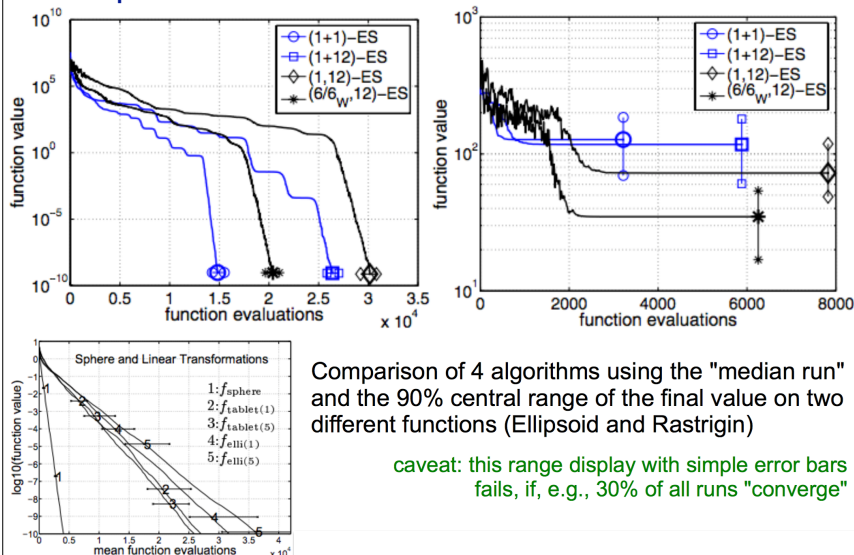
more general: use quantiles as summary data

for example out of 15 data: 2nd, 8th, and 14th value
represent the 10%, 50%, and 90%-tile

unless there are good reasons for a different statistics

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Examples



Comparison of 4 algorithms using the "median run" and the 90% central range of the final value on two different functions (Ellipsoid and Rastrigin)

caveat: this range display with simple error bars fails, if, e.g., 30% of all runs "converge"

Statistical Assessment

① Assess the meaning/**relevance** of a difference first (the only difficult part)

using enough data, any difference
can be made significant

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Statistical Assessment

- ② Apply **rank-sum test** (Wilcoxon, Mann-Whitney U)
only assumption: no equal data values

hypothesis:

compares $\text{sPr}(x > y) \neq \text{Pr}(x < y) \neq 1/2$ inking

two-sided 1%-significance p-value needs only 2x5 data values

For the same p-value, **fewer significant data** are better
using enough data, any difference
can be made significant

Generally: non-parametric tests, Kolmogorov-Smirnov test for ECDFs, no need to use the t-test

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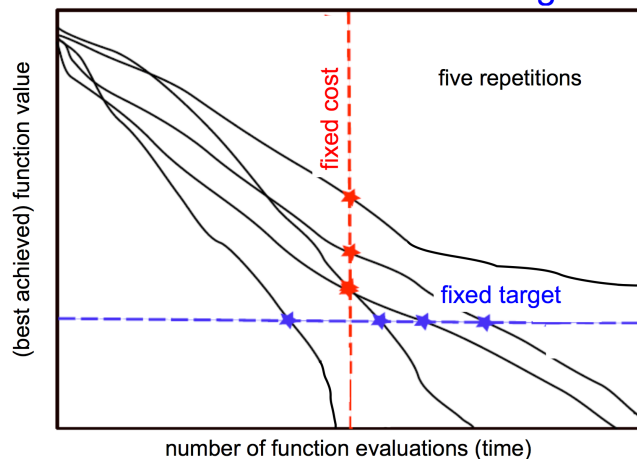
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Measuring Performance from Convergence Graphs

fixed-cost versus **fixed-target**



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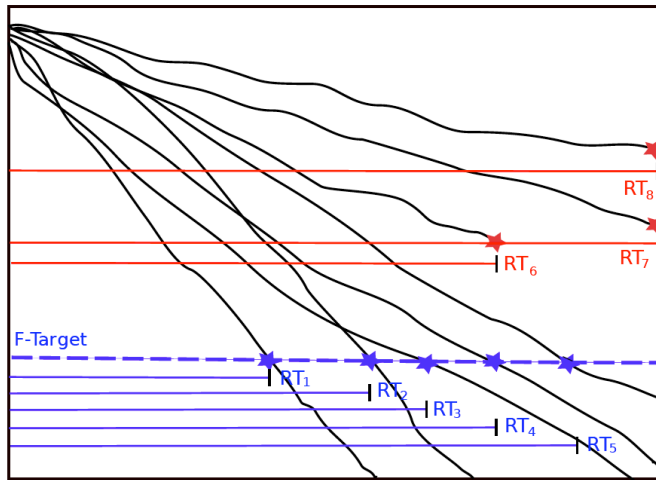
Evaluation of Search Algorithms

Behind the scene

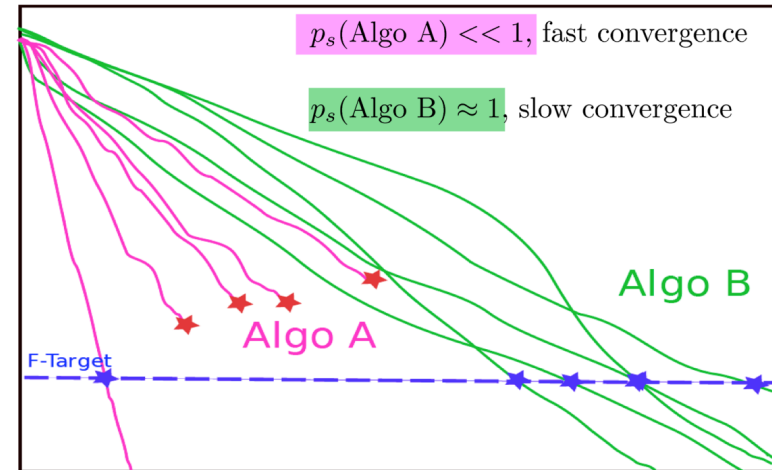
- a performance should be
- quantitative** on the ratio scale (highest possible)
 - “algorithm A is two *times* better than algorithm B” is a meaningful statement
 - can assume a wide range of values
 - meaningful (interpretable)** with regard to the real world
 - possible to transfer from benchmarking to real world
 - runtime** or **first hitting time** is the prime candidate (we don't have many choices anyway)

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Collect Runtime to reach F-target



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Which Performance Measure
to compare the two following scenario?

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Which Performance Measure

Algo Restart A:



$$p_s(\text{Algo Restart A}) = 1$$

Algo Restart B:



$$p_s(\text{Algo Restart B}) = 1$$

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Average Runtime (aRT)

$$\text{ERT} = \mathbb{E}[\text{RT}^r] = \frac{1-p_s}{p_s} \mathbb{E}[\text{RT}_{\text{unsuccessful}}] + \mathbb{E}[\text{RT}_{\text{successful}}]$$

aRT is an estimator for ERT

$$\text{aRT} = \frac{\# \text{Evals}}{\# \text{success}}$$

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Step-size Adaptation
Covariance Matrix Adaptation

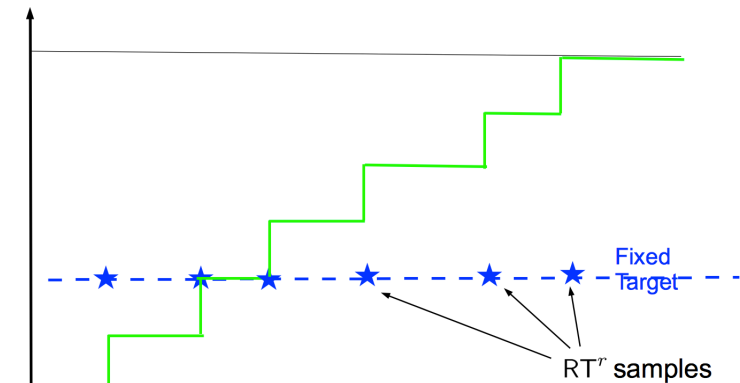
❹ Evaluating Black-Box Algorithms

Displaying results and visualization
Statistics
Average Runtime
Empirical Cumulative Distribution Function (ECDF)

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Empirical Cumulative Distribution Function

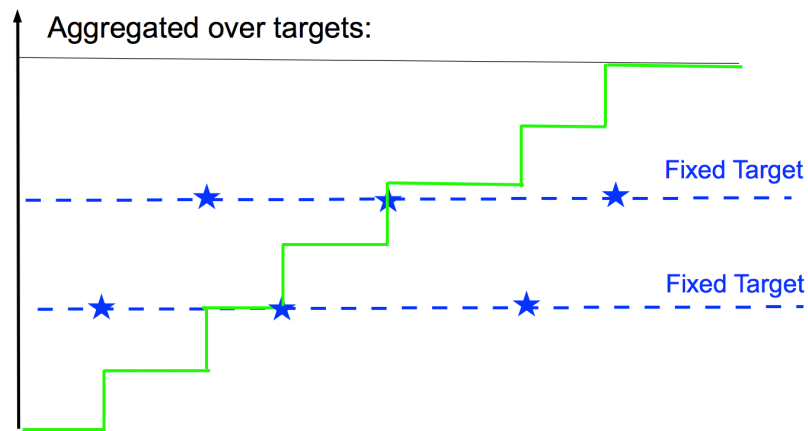
Single Target



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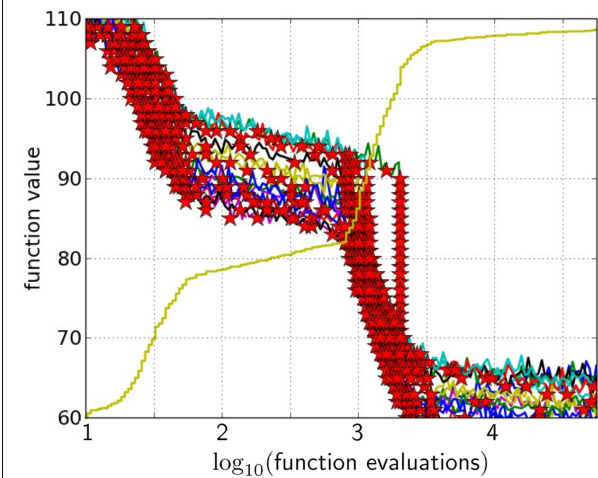
Empirical Cumulative Distribution Function

Several Targets



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Empirical Cumulative Distribution Function



15 runs
50 targets
ECDF with 750 steps

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