

# Package ‘evoper’

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**Type** Package

**Title** Evolutionary Parameter Estimation for 'Repast Simphony' Models

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**URL** <https://github.com/antonio-pgarcia/evoper>

**BugReports** <https://github.com/antonio-pgarcia/evoper/issues>

## Description

The EvoPER, Evolutionary Parameter Estimation for Individual-based Models is an extensible package providing optimization driven parameter estimation methods using metaheuristics and evolutionary computation techniques (Particle Swarm Optimization, Simulated Annealing, Ant Colony Optimization for continuous domains, Tabu Search, Evolutionary Strategies, ...) which could be more efficient and require, in some cases, fewer model evaluations than alternatives relying on experimental design. Currently there are built in support for models developed with 'Repast Simphony' Agent-Based framework (<<https://repast.github.io/>>) and with NetLogo (<<https://ccl.northwestern.edu/netlogo/>>) which are the most used frameworks for Agent-based modeling.

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**LazyData** TRUE

**Depends** rrepast

**Imports** methods, futile.logger, boot, reshape, ggplot2, deSolve, plot3D, plyr, data.table, utils, RNetLogo

**RoxygenNote** 6.0.1

**Suggests** testthat

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<b>abm.acor</b>	<i>Ant colony optimization for continuous domains</i>
-----------------	---

---

## Description

An implementation of Ant Colony Optimization algorithm for continuous variables.

## Usage

```
abm.acor(objective, options = NULL)
```

## Arguments

objective	An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a>
options	An appropriate instance from a subclass of <a href="#">Options</a> class

## References

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>

## Examples

```
## Not run:  
f<- PlainFunction$new(f0.rosenbrock2)  
  
f$Parameter(name="x1",min=-100,max=100)  
f$Parameter(name="x2",min=-100,max=100)  
  
extremize("acor", f)  
  
## End(Not run)
```

abm.ees1

*EvoPER Evolutionary Strategy 1*

## Description

This function tries to provide a rough approximation to best solution when no information is available for the correct range of input parameters for the objective function. It can be useful for studying the behavior of individual-based models with high variability in the output variables showing non-linear behaviors.

## Usage

```
abm.ees1(objective, options = NULL)
```

## Arguments

objective	An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a>
options	An appropriate instance from a subclass of <a href="#">Options</a> class

## Examples

```
## Not run:
f<- PlainFunction$new(f0.rosenbrock2)

f$Parameter(name="x1",min=-100,max=100)
f$Parameter(name="x2",min=-100,max=100)

extremize("ees1", f)

## End(Not run)
```

abm.ees2

*EvoPER Evolutionary Strategy 2*

## Description

This function tries to provide a rough approximation to best solution when no information is available for the correct range of input parameters for the objective function. It can be useful for studying the behavior of individual-based models with high variability in the output variables showing non-linear behaviors.

## Usage

```
abm.ees2(objective, options = NULL)
```

## Arguments

objective	An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a>
options	An appropriate instance from a subclass of <a href="#">Options</a> class

## Examples

```
## Not run:
f<- PlainFunction$new(f0.rosenbrock2)

f$Parameter(name="x1",min=-100,max=100)
f$Parameter(name="x2",min=-100,max=100)

extremize("ees2", f)

## End(Not run)
```

## Description

An implementation of Particle Swarm Optimization method for parameter estimation of Individual-based models.

## Usage

```
abm.pso(objective, options = NULL)
```

## Arguments

objective	An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a>
options	An appropriate instance from a subclass of <a href="#">Options</a> class

## References

- [1] Kennedy, J., & Eberhart, R. (1995). Particle swarm optimization. In Proceedings of ICNN 95 - International Conference on Neural Networks (Vol. 4, pp. 1942-1948). IEEE.
- [2] Poli, R., Kennedy, J., & Blackwell, T. (2007). Particle swarm optimization. Swarm Intelligence, 1(1), 33-57.

## Examples

```
## Not run:
f<- PlainFunction$new(f0.rosenbrock2)

f$Parameter(name="x1",min=-100,max=100)
f$Parameter(name="x2",min=-100,max=100)

extremize("pso", f)

## End(Not run)
```

abm.saa

*abm.saa*

## Description

An implementation of Simulated Annealing Algorithm optimization method for parameter estimation of Individual-based models.

## Usage

```
abm.saa(objective, options = NULL)
```

## Arguments

- |           |  |
|-----------|--|
| objective | An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a> |
| options   | An appropriate instance from a subclass of <a href="#">Options</a> class               |

## Value

The best solution.

## References

- [1] Kirkpatrick, S., Gelatt, C. D., & Vecchi, M. P. (1983). Optimization by Simulated Annealing. *Science*, 220(4598).

## Examples

```
## Not run:
f<- PlainFunction$new(f0.rosenbrock2)

f$Parameter(name="x1",min=-100,max=100)
f$Parameter(name="x2",min=-100,max=100)

extremize("saa", f)

## End(Not run)
```

```

## Not run:
## A Repast defined function
f<- RepastFunction$new从根本models/BactoSim(HaldaneEngine-1.0)","ds::Output",300)

## or a plain function

f1<- function(x1,x2,x3,x4) {
  10 * (x1 - 1)^2 + 20 * (x2 - 2)^2 + 30 * (x3 - 3)^2 + 40 * (x4 - 4)^2
}

f<- PlainFunction$new(f1)

f$addFactor(name="cyclePoint",min=0,max=90)
f$addFactor(name="conjugationCost",min=0,max=100)
f$addFactor(name="pilusExpressionCost",min=0,max=100)
f$addFactor(name="gamma0",min=1,max=10)

abm.saa(f, 100, 1, 100, 0.75)

## End(Not run)

```

## Description

An implementation of Tabu Search algorithm for parameter estimation

## Usage

```
abm.tabu(objective, options = NULL)
```

## Arguments

objective	An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a>
options	An appropriate instance from a subclass of <a href="#">Options</a> class

## References

- [1] Fred Glover (1989). "Tabu Search - Part 1". ORSA Journal on Computing, 190-206. doi:10.1287/ijoc.1.3.190.
- [2] Fred Glover (1990). "Tabu Search - Part 2". ORSA Journal on Computing, 4-32. doi:10.1287/ijoc.2.1.4.

## Examples

```
## Not run:
f<- PlainFunction$new(f0.rosenbrock2)

f$Parameter(name="x1",min=-100,max=100)
f$Parameter(name="x2",min=-100,max=100)

or

f$Parameter0(name="x1",levels=c(0:4))
f$Parameter0(name="x2",levels=c(-2,-1,0,1,2))

extremize("tabu", f)

## End(Not run)
```

acor.archive

*acor.archive*

## Description

This function is used for creating and maintaining the ACOr archive 'T'. The function keeps the track of 'k' solution in the archive.

## Usage

```
acor.archive(s, f, w, k, T = NULL)
```

## Arguments

s	The solution 'ants'
f	The evaluation of solution
w	The weight vector
k	The archive size
T	The current archive

## Value

The solution archive

## References

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>

---

`acor.F`*acor.F*

---

**Description**

Helper function for extracting the 'F' function evaluations from archive ACor 'T'

**Usage**`acor.F(T)`**Arguments**

T                   The solution archive

**Value**

The F matrix

---

`acor.lthgaussian`*Select the lth gaussian function*

---

**Description**

Given a weight vector calculate the probabilities of selecting the lth gaussian function and return the index of lht gaussian selected with probability p

**Usage**`acor.lthgaussian(W)`**Arguments**

W                   The vector of weights

**Value**

The index of lht gaussian function

**References**

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>

**acor.N***acor.N***Description**

Helper function for getting the size of solution

**Usage**

```
acor.N(T)
```

**Arguments**

T	The solution archive
---	----------------------

**Value**

The size 'n' of a solution 's'

**acor.probabilities***Gaussian kernel choosing probability***Description**

Calculate the probability of choosing the lth Gaussian function

**Usage**

```
acor.probabilities(W, l = NULL)
```

**Arguments**

W	The vector of weights
l	The lth element of algorithm solution archive T

**Value**

The vector of probabilities 'p'

**References**

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>

---

`acor.S`*acor.S***Description**

Helper function for extracting solution 'S' from archive 'T'

**Usage**`acor.S(T)`**Arguments**

T                   The solution archive

**Value**

The solution matrix

---

`acor.sigma`*Sigma calculation for ACOr***Description**

Calculate the value of sigma

**Usage**`acor.sigma(Xi, k, T)`**Arguments**

Xi                   The algorithm parameter

k                   The solution archive size

T                   The solution archive

**Value**

The sigma value

**References**

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>

---

acor.updateants	<i>acor.updateants</i>
-----------------	------------------------

---

### Description

Update the solution using the gaussian kernel

### Usage

```
acor.updateants(S, N, W, t.mu, t.sigma)
```

### Arguments

S	The current solution ants
N	The numnber of required ants in solution
W	The weight vector
t.mu	The 'mean' from solution archive
t.sigma	The value of sigma from solution archive

### Value

The new solution ants

### References

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>
- 

---

acor.W	<i>acor.W</i>
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---

### Description

Helper function for extracting the 'W' function evaluations from archive ACOOr 'T'

### Usage

```
acor.W(T)
```

### Arguments

T	The solution archive
---	----------------------

### Value

The weight vector

---

<code>acor.weighth</code>	<i>Weight calculation for ant colony optimization</i>
---------------------------	---

---

### Description

Calculates the weight element of ACOr algorithm for the solution archive.

### Usage

```
acor.weighth(q, k, l)
```

### Arguments

<code>q</code>	The Algorithm parameter. When small best-ranked solution is preferred
<code>k</code>	The Archive size
<code>l</code>	The lth element of algorithm solution archive T

### Value

A scalar or a vector with calculated weigth.

### References

- [1] Socha, K., & Dorigo, M. (2008). Ant colony optimization for continuous domains. European Journal of Operational Research, 185(3), 1155-1173. <http://doi.org/10.1016/j.ejor.2006.06.046>

---

<code>assert</code>	<i>assert</i>
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---

### Description

The assert function stop the execution if the logical expression given by the parameter expresion is false.

### Usage

```
assert(expresion, string)
```

### Arguments

<code>expresion</code>	Some logical expression
<code>string</code>	The text message to show if expression does not hold

---

**bestFitness***bestFitness*

---

**Description**

Given a set S of N solutions created with sortSolution, this function returns the fitness component for the best solution.

**Usage**

```
bestFitness(S)
```

**Arguments**

S                   The solution set

**Value**

The best fitness value

---

**bestSolution***bestSolution*

---

**Description**

Given a set S of N solutions created with sortSolution, this function returns the best solution found.

**Usage**

```
bestSolution(S)
```

**Arguments**

S                   The solution set

**Value**

The best solution

---

`cbuf`*cbuf*

---

**Description**

Simple implementation of a circular buffer.

**Usage**

```
cbuf(b, v, e)
```

**Arguments**

b	The variable holding the current buffer content
v	The new valued to be added to b
e	The length of circular buffer

**Value**

The buffer b plus the element v minus the least recently added element

---

`compare.algorithms1`    *compare.algorithms1*

---

**Description**

Compare the number of function evalutions and convergence for the following optimization algorithms, ("saa","pso","acor","ees1").

**Usage**

```
compare.algorithms1(F, seeds = c(27, 2718282, 36190727, 3141593, -91190721,  
-140743, 1321))
```

**Arguments**

F	The function to be tested
seeds	The random seeds which will be used for testing algorithms

## Examples

```
## Not run:
rm(list=ls())
d.cigar4<- compare.algorithms1(f0.cigar4)
d.schaffer4<- compare.algorithms1(f0.schaffer4)
d.griewank4<- compare.algorithms1(f0.griewank4)
d.bohachevsky4<- compare.algorithms1(f0.bohachevsky4)
d.rosenbrock4<- compare.algorithms1(f0.rosenbrock4)

## End(Not run)
```

**contourplothelper**      *contourplothelper*

## Description

Simple helper function for countour plots

## Usage

```
contourplothelper(d, x, y, nbins = 32, binwidth = c(10, 10),
  points = c(300, 300), title = NULL)
```

## Arguments

d	A data frame.
x	A string with the dataframe column name for x axis.
y	A string with the dataframe column name for y axis.
z	A string with the dataframe column name for z axis.
nbins	The number bins. The default is 32.
binwidth	The binwidths for 'kde2d'. Can be an scalar or a vector.
points	The number of grid points. Can be an scalar or a vector.
title	The optional plot title. May be omitted.

---

ees1.challenge

---

*ees1.challenge*

---

### Description

Repeat the evalution of best solution to tackle with variability.

### Usage

`ees1.challenge(solution, objective)`

### Arguments

solution	The Problem solution
objective	The objective function

---

ees1.explore

---

*ees1.explore*

---

### Description

Explore the solution space on the neighborhood of solution 's' in order to find a new best.

### Usage

`ees1.explore(s, weight, p = 0.01)`

### Arguments

s	The Problem solution
weight	The exploration intensity
p	The mutation probability

ees1.mating

*ees1.mating*

## Description

This function 'mix' the elements present in the solution. The parameter 'mu' controls the intensity of mixing. Low values give preference to best solution components and high values make the values being select randomly.

## Usage

```
ees1.mating(solution, mu)
```

## Arguments

solution	The Problem solution
mu	The mixing intensity ratio, from 0 to 1. The mix intensity controls de the probability of chosing a worst solutions

ees1.mating1

*ees1.mating1*

## Description

This function 'mix' the elements present in the solution. The parameter 'mu' controls the intensity of mixing. Low values give preference to best solution components and high values make the values being select randomly.

## Usage

```
ees1.mating1(solution, mu)
```

## Arguments

solution	The Problem solution
mu	The mixing intensity ratio, from 0 to 1. The mix intensity controls de the probability of chosing a worst solutions

---

`ees1.mutation`*ees1.mutation*

---

**Description**

Performs the mutation on generated solution

**Usage**

```
ees1.mutation(solution, mates, p = 0.01)
```

**Arguments**

solution	The Problem solution
mates	The mixed parents
p	The mutation probability

---

`ees1.recombination`*ees1.recombination*

---

**Description**

Performs the recombination on solution

**Usage**

```
ees1.recombination(solution, mates)
```

**Arguments**

solution	The Problem solution
mates	The mixed parents

---

ees1.selection	<i>ees.selection</i>
----------------	----------------------

---

**Description**

Select the elements with best fitness but accept uphill moves with probability 'kkappa'.

**Usage**

```
ees1.selection(s0, s1, kkappa)
```

**Arguments**

s0	The current best solution set
s1	The new solution
kkappa	The selection pressure

---

elog.debug	<i>elog.debug</i>
------------	-------------------

---

**Description**

Wrapper for logging debug messages.

**Usage**

```
elog.debug(...)
```

**Arguments**

...	Variable number of arguments including a format string.
-----	---

---

elog.error	<i>elog.error</i>
------------	-------------------

---

**Description**

Wrapper for logging error messages.

**Usage**

```
elog.error(...)
```

**Arguments**

...	Variable number of arguments including a format string.
-----	---

---

`elog.info`*elog.info*

---

**Description**

Wrapper for logging info messages.

**Usage**

```
elog.info(...)
```

**Arguments**

... Variable number of arguments including a format string.

---

`elog.level`*elog.level*

---

**Description**

Configure the current log level

**Usage**

```
elog.level(level = NULL)
```

**Arguments**

level The log level (ERROR|WARN|INFO|DEBUG)

**Value**

The log level

`enforceBounds`*enforceBounds***Description**

Checks if parameters fall within upper and lower bounds

**Usage**

```
enforceBounds(particles, factors)
```

**Arguments**

<code>particles</code>	The particle set
<code>factors</code>	the defined range for objective function parameters

**Value**

The particle inside the valid limits

`es.evaluate`*es.evaluate***Description**

For each element in solution 's' evaluate the respective fitness.

**Usage**

```
es.evaluate(f, s, enforce = TRUE)
```

**Arguments**

<code>f</code>	A reference to an instance of objective function
<code>s</code>	The set of solutions
<code>enforce</code>	If true the values are enforced to fall within provided range

**Value**

The solution ordered by its fitness.

---

Estimates-class	<i>Estimates</i>
-----------------	------------------

---

## Description

A simple class for encapsulating the return of metaheuristic methods

---

extremize	<i>extremize</i>
-----------	------------------

---

## Description

Entry point for optimization functions

## Usage

```
extremize(type, objective, options = NULL)
```

## Arguments

- |           |  |
|-----------|--|
| type      | The optimization method (aco,pso,saa,sda)  |
| objective | An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a> |
| options   | An apropiate instance from a subclass of <a href="#">Options</a> class                 |

## Examples

```
## Not run:  
f<- PlainFunction$new(f0.rosenbrock2)  
  
f$Parameter(name="x1",min=-100,max=100)  
f$Parameter(name="x2",min=-100,max=100)  
  
extremize("pso", f)  
  
## End(Not run)
```

**f0.ackley***f0.ackley***Description**

The ackley function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1 \dots N$ ,  $f(x) = 0$ . Domain  $x_i \in [-32.768, 32.768]$ , for all  $i = 1, \dots, d$

**Usage**

```
f0.ackley(...)
```

**Arguments**

... The variadic list of function variables.

**Value**

The function value

**References**

<https://www.sfu.ca/~ssurjano/ackley.html>

**f0.ackley4***f0.ackley4***Description**

The ackley function of four variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1 \dots N$ ,  $f(x) = 0$ .

**Usage**

```
f0.ackley4(x1, x2, x3, x4)
```

**Arguments**

x1	The first function variable
x2	The second function variable
x3	The third function variable
x4	The fourth function variable

**Value**

The function value

---

f0.adtn.rosenbrock2    *f0.adtn.rosenbrock2*

---

**Description**

Two variable Rosenbrock function with random additive noise.

**Usage**

f0.adtn.rosenbrock2(x1, x2)

**Arguments**

x1	Parameter 1
x2	Parameter 2

---

f0.bohachevsky    *f0.bohachevsky*

---

**Description**

The Bohachevsky function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1 \dots N$ ,  $f(x) = 0$ .

**Usage**

f0.bohachevsky(...)

**Arguments**

...    The variadic list of function variables.

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

**f0.bohachevsky4**      *f0.bohachevsky4*

### Description

The Bohachevsky function of four variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

### Usage

`f0.bohachevsky4(x1, x2, x3, x4)`

### Arguments

<code>x1</code>	The first function variable
<code>x2</code>	The second function variable
<code>x3</code>	The third function variable
<code>x4</code>	The fourth function variable

### Value

The function value

**f0.cigar**      *f0.cigar*

### Description

The Cigar function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

### Usage

`f0.cigar(...)`

### Arguments

<code>...</code>	The variadic list of function variables.
------------------	--

### Value

The function value

### References

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

**f0.cigar4***f0.cigar4*

---

**Description**

The Cigar function of four variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

```
f0.cigar4(x1, x2, x3, x4)
```

**Arguments**

x1	The first function variable
x2	The second function variable
x3	The third function variable
x4	The fourth function variable

**Value**

The function value

---

**f0.griewank***f0.griewank*

---

**Description**

The griewank function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

```
f0.griewank(...)
```

**Arguments**

... The variadic list of function variables.

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

**f0.griewank4***f0.griewank4***Description**

The griewank function of four variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

```
f0.griewank4(x1, x2, x3, x4)
```

**Arguments**

x1	The first function variable
x2	The second function variable
x3	The third function variable
x4	The fourth function variable

**Value**

The function value

**f0.nlnn.rosenbrock2***f0.nlnn.rosenbrock2***Description**

Two variable Rosenbrock function with random additive noise.

**Usage**

```
f0.nlnn.rosenbrock2(x1, x2)
```

**Arguments**

x1	Parameter 1
x2	Parameter 2

---

f0.periodtuningpp      *Period tuning for Predator-Prey base*

---

### Description

This function is an example on how EvoPER can be used for estimating the parameter values in order to produce oscillations with the desired period. It is not intended to be used directly, the provided wrappers should be instead.

### Usage

```
f0.periodtuningpp(x1, x2, x3, x4, period)
```

### Arguments

x1	The growth rate of prey
x2	The decay rate of predator
x3	The predating effect on prey
x4	The predating effect on predator
period	The desired oscillation period

### Value

The solution fitness cost

---

f0.periodtuningpp12      *Period tuning of 12 time units for Predator-Prey*

---

### Description

This function is an example on how EvoPER can be used for estimating the parameter values in order to produce oscillations with the desired period.

### Usage

```
f0.periodtuningpp12(x1, x2, x3, x4)
```

### Arguments

x1	The growth rate of prey
x2	The decay rate of predator
x3	The predating effect on prey
x4	The predating effect on predator

**Value**

The solution fitness cost

**Examples**

```
## Not run:
rm(list=ls())
set.seed(-27262565)
f<- PlainFunction$new(f0.periodtuningpp12)
f$Parameter(name="x1",min=0.5,max=2)
f$Parameter(name="x2",min=0.5,max=2)
f$Parameter(name="x3",min=0.5,max=2)
f$Parameter(name="x4",min=0.5,max=2)
extremize("pso", f)

## End(Not run)
```

**f0.periodtuningpp24**    *Period tuning of 24 time units for Predator-Prey*

**Description**

This function is an example on how EvoPER can be used for estimating the parameter values in order to produce oscillations with the desired period.

**Usage**

```
f0.periodtuningpp24(x1, x2, x3, x4)
```

**Arguments**

x1	The growth rate of prey
x2	The decay rate of predator
x3	The predating effect on prey
x4	The predating effect on predator

**Value**

The solution fitness cost

### Examples

```
## Not run:
rm(list=ls())
set.seed(-27262565)
f<- PlainFunction$new(f0.periodtuningpp24)
f$Parameter(name="x1",min=0.5,max=2)
f$Parameter(name="x2",min=0.5,max=2)
f$Parameter(name="x3",min=0.5,max=2)
f$Parameter(name="x4",min=0.5,max=2)
extremize("pso", f)

## End(Not run)
```

f0.periodtuningpp48     *Period tuning of 48 time units for Predator-Prey*

### Description

This function is an example on how EvoPER can be used for estimating the parameter values in order to produce oscillations with the desired period.

### Usage

```
f0.periodtuningpp48(x1, x2, x3, x4)
```

### Arguments

x1	The growth rate of prey
x2	The decay rate of predator
x3	The predating effect on prey
x4	The predating effect on predator

### Value

The solution fitness cost

### Examples

```
## Not run:
rm(list=ls())
set.seed(-27262565)
f<- PlainFunction$new(f0.periodtuningpp24)
f$Parameter(name="x1",min=0.5,max=2)
f$Parameter(name="x2",min=0.5,max=2)
f$Parameter(name="x3",min=0.5,max=2)
f$Parameter(name="x4",min=0.5,max=2)
extremize("pso", f)
```

```
## End(Not run)
```

### f0.periodtuningpp72    *Period tuning of 72 time units for Predator-Prey*

#### Description

This function is an example on how EvoPER can be used for estimating the parameter values in order to produce oscillations with the desired period.

#### Usage

```
f0.periodtuningpp72(x1, x2, x3, x4)
```

#### Arguments

x1	The growth rate of prey
x2	The decay rate of predator
x3	The predating effect on prey
x4	The predating effect on predator

#### Value

The solution fitness cost

#### Examples

```
## Not run:
rm(list=ls())
set.seed(-27262565)
f<- PlainFunction$new(f0.periodtuningpp24)
f$Parameter(name="x1",min=0.5,max=2)
f$Parameter(name="x2",min=0.5,max=2)
f$Parameter(name="x3",min=0.5,max=2)
f$Parameter(name="x4",min=0.5,max=2)
extremize("pso", f)

## End(Not run)
```

---

f0.rosenbrock2      *f0.rosenbrock2*

---

**Description**

Two variable Rosenbrock function, where  $f(1,1) = 0$

**Usage**

f0.rosenbrock2(x1, x2)

**Arguments**

x1	Parameter 1
x2	Parameter 2

---

f0.rosenbrock4      *f0.rosenbrock4*

---

**Description**

The rosenbrock function of 4 variables for testing optimization methods. The global optima for the function is given by  $x_i = 1$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

f0.rosenbrock4(x1, x2, x3, x4)

**Arguments**

x1	The first function variable
x2	The second function variable
x3	The third function variable
x4	The fourth function variable

**Value**

The function value

---

**f0.rosenbrockn**      *f0.rosenbrockn*

---

**Description**

The rosenbrock function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 1$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

`f0.rosenbrockn(...)`

**Arguments**

...      The variadic list of function variables.

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

**f0.schaffer**      *f0.schaffer*

---

**Description**

The schaffer function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

`f0.schaffer(...)`

**Arguments**

...      The variadic list of function variables.

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

f0.schaffer4

*f0.schaffer4*

---

### Description

The Schaffer function of four variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

### Usage

`f0.schaffer4(x1, x2, x3, x4)`

### Arguments

x1	The first function variable
x2	The second function variable
x3	The third function variable
x4	The fourth function variable

### Value

The function value

---

f0.schwefel

*f0.schwefel*

---

### Description

The schwefel function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 420.96874636$ , forall  $i \in 1...N$ ,  $f(x) = 0$ . The range of  $x_i$  is [-500,500]

### Usage

`f0.schwefel(...)`

### Arguments

... The variadic list of function variables.

### Value

The function value

### References

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

**f0.schwefel4***f0.schwefel4***Description**

The schwefel function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 420.96874636$ , forall  $i \in 1...N$ ,  $f(x) = 0$ . The range of  $x_i$  is [-500,500]

**Usage**

```
f0.schwefel4(x1, x2, x3, x4)
```

**Arguments**

x1	The first function variable
x2	The second function variable
x3	The third function variable
x4	The fourth function variable

**Value**

The function value

**f0.test***f0.test***Description**

Simple test function  $f(1,2,3,4) = 0$

**Usage**

```
f0.test(x1, x2, x3, x4)
```

**Arguments**

x1	Parameter 1
x2	Parameter 2
x3	Parameter 3
x4	Parameter 4

---

`f1.ackley`*f1.ackley*

---

**Description**

The ackley function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , for all  $i \in 1 \dots N$ ,  $f(x) = 0$ . Domain  $x_i \in [-32.768, 32.768]$ , for all  $i = 1, \dots, d$

**Usage**`f1.ackley(x)`**Arguments**

`x` The vector of function parameters

**Value**

The function value

**References**

<https://www.sfu.ca/~ssurjano/ackley.html>

---

---

`f1.adtn.rosenbrock2`*f1.adtn.rosenbrock2*

---

**Description**

Two variable Rosenbrock function with random additive noise.

**Usage**`f1.adtn.rosenbrock2(x)`**Arguments**

`x` Parameter vector

f1.bohachevsky

*f1.bohachevsky***Description**

The Bohachevsky function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

```
f1.bohachevsky(x)
```

**Arguments**

x	The vector of function parameters
---	-----------------------------------

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

f1.cigar

*f1.cigar***Description**

The Cigar function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

```
f1.cigar(x)
```

**Arguments**

x	The vector of function variables.
---	-----------------------------------

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

*f1.griewank**f1.griewank*

---

**Description**

The griewank function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

**Usage**

*f1.griewank(x)*

**Arguments**

*x*                    The vector of function parameters

**Value**

The function value

**References**

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

---

*f1.nlnn.rosenbrock2*    *f1.nlnn.rosenbrock2*

---

**Description**

Two variable Rosenbrock function with random additive noise.

**Usage**

*f1.nlnn.rosenbrock2(x)*

**Arguments**

*x*                    Parameter vector

---

f1.rosenbrock2      *f1.rosenbrock2*

---

### Description

Two variable Rosenbrock function, where  $f(c(1,1)) = 0$

### Usage

`f1.rosenbrock2(x)`

### Arguments

x                  Parameter vector

---

f1.rosenbrockn      *f1.rosenbrockn*

---

### Description

The rosenbrock function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 1$ , forall  $i \in 1 \dots N$ ,  $f(x) = 0$ .

### Usage

`f1.rosenbrockn(x)`

### Arguments

x                  The vector of function parameters

### Value

The function value

### References

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

f1.schaffer

*f1.schaffer*

---

### Description

The schaffer function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 0$ , forall  $i \in 1...N$ ,  $f(x) = 0$ .

### Usage

`f1.schaffer(x)`

### Arguments

`x` The vector of function parameters

### Value

The function value

### References

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

---

f1.schwefel

*f1.schwefel*

---

### Description

The schwefel function of N variables for testing optimization methods. The global optima for the function is given by  $x_i = 420.96874636$ , forall  $i \in 1...N$ ,  $f(x) = 0$ . The range of  $x_i$  is  $[-500,500]$

### Usage

`f1.schwefel(x)`

### Arguments

`x` The vector of function variables.

### Value

The function value

### References

<http://deap.gel.ulaval.ca/doc/dev/api/benchmarks.html>

**f1.test***f1.test***Description**

Simple test function  $f(c(1,2,3,4)) = 0$

**Usage**

```
f1.test(x)
```

**Arguments**

x	Parameter vector
---	------------------

**fixdfcolumns***fixdfcolumns***Description**

Coerce dataframe columns to a specific type.

**Usage**

```
fixdfcolumns(df, cols = c(), skip = TRUE, type = as.numeric)
```

**Arguments**

df	The data frame.
cols	The dataframe columns to be skipped or included.
skip	If TRUE the column names in 'cols' are skipped. When FALSE logic is inverted.
type	The type for which data frame columns must be converted.

**Value**

The data frame with converted column types.

---

generateSolution      *generateSolution*

---

**Description**

Generates a problema solution using discrete leves

**Usage**

```
generateSolution(parameters, size)
```

**Arguments**

parameters	The Objective Function parameter list
size	The solution size

**Value**

The solution set

---

getFitness      *getFitness*

---

**Description**

Given a set S of N solutions created with sortSolution, this function returns the solution component for the best solution.

**Usage**

```
getFitness(S, i = NULL)
```

**Arguments**

S	The solution set
i	The fitness index, if null return the whole column.

**Value**

The selected fitness entry

---

getSolution	<i>getSolution</i>
-------------	--------------------

---

**Description**

Given a set S of N solutions created with sortSolution, this function returns the solution component.  
A solutions is a set of solutions and their associated fitness

**Usage**

```
getSolution(S)
```

**Arguments**

S                   The solution set

**Value**

The solution set

---

gm.mean	<i>gm.mean</i>
---------	----------------

---

**Description**

Simple implementation for geometric mean

**Usage**

```
gm.mean(x)
```

**Arguments**

x                   data

**Value**

geometric mean for data

---

gm.sd

---

*gm.sd*

---

### Description

Simple implementation for geometric standard deviation

### Usage

```
gm.sd(x, mu = NULL)
```

### Arguments

x	data
mu	The geometric mean. If not provided it is calculated.

### Value

geometric standard deviation for data

---

---

histplothelper

---

*histplothelper*

---

### Description

Simple helper for plotting histograms

### Usage

```
histplothelper(d, x, title = NULL)
```

### Arguments

d	A data frame.
x	A string with the dataframe column name for histogram
title	The plot title

### Value

A ggplot2 plot object

<code>initSolution</code>	<i>initSolution</i>
---------------------------	---------------------

### Description

Creates the initial Solution population taking into account the lower an upper bounds of provided experiment factors.

### Usage

```
initSolution(parameters, N = 20, sampling = "mcs")
```

### Arguments

<code>parameters</code>	The Objective Function parameter list
<code>N</code>	The size of Solution population
<code>sampling</code>	The population sampling scheme, namely <mcs lhs ffs> standing respectively for montecarlo sampling, latin hypercube sampling and full factorial sampling

### Value

A random set of solutions

<code>lowerBound</code>	<i>lowerBound</i>
-------------------------	-------------------

### Description

Checks if parameters is greater than the lower bounds

### Usage

```
lowerBound(particles, factors)
```

### Arguments

<code>particles</code>	The particle set
<code>factors</code>	the defined range for objective function parameters

### Value

The particle greater than or equal to lower limit

---

Magnitude

*Magnitude*

---

**Description**

Calculates the magnitude order for a given value

**Usage**

`Magnitude(v)`

**Arguments**

v                   The numerical value

**Value**

The magnitude order

---

naiveperiod

*naiveperiod*

---

**Description**

A naive approach for finding the period in a series of data points

**Usage**

`naiveperiod(d)`

**Arguments**

d                   The data to search period

**Value**

A list with the average period and amplitude

---

NetLogoFunction-class *NetLogoFunction*

---

**Description**

NetLogoFunction class

---

**NLWrapper.FindJar**      *NLWrapper:FindJar*

---

### Description

Search for the netlogo jar file on the provided path

### Usage

`NLWrapper.FindJar(path)`

### Arguments

**path**            The base path for searching

### Value

The path for NetLogo jar file

---

**NLWrapper.GetParameter**      *NLWrapper:GetParameter*

---

### Description

Gets the value of a model parameter

### Usage

`NLWrapper.GetParameter(obj, name)`

### Arguments

**obj**            The object retuned by [NLWrapper.Model](#)

**name**            The parameter name string or the collection of parameter names

### Value

The parameter values

## Examples

```
## Not run:
rm(list=ls())
p<- "C:/Program Files/NetLogo 6.0.4/app"
m<- file.path(nlpath, "models", "Sample Models", "Biology", "Wolf Sheep Predation.nlogo")
o<- NLWrapper.Model(p, m)
v<- NLWrapper.GetParameter(o, c("initial-number-sheep"))

or

v<- NLWrapper.GetParameter(o, c("initial-number-sheep", "initial-number-wolves"))

## End(Not run)
```

**NLWrapper . Model***NLWrapper.Model*

## Description

This wrapper prepares the environment and instantiates the model

## Usage

```
NLWrapper.Model(netlogodir, modelfile, dataset, maxtime)
```

## Arguments

<code>netlogodir</code>	The base path of NetLogo installation
<code>modelfile</code>	The absolute path for NetLogo model file
<code>dataset</code>	The names of model variables
<code>maxtime</code>	The total number of iterations

## Examples

```
## Not run:
rm(list=ls())
p<- "C:/Program Files/NetLogo 6.0.4/app"
output<- c("count sheep", "count wolves")
m<- file.path(nlpath, "models", "Sample Models", "Biology", model, "Wolf Sheep Predation.nlogo")
o<- NLWrapper.Model(p, m, output, 150)

## End(Not run)
```

---

**NLWrapper.Run**      *NLWrapper.Run*

---

### Description

Executes a NetLogo Model using rNetLogo

### Usage

```
NLWrapper.Run(obj, r = 1, seed = c())
```

### Arguments

obj	The object retuned by <a href="#">NLWrapper.Model</a>
r	The number of replications
seed	The collection of random seeds

---

**NLWrapper.RunExperiment**      *NLWrapper.RunExperiment*

---

### Description

Executes a NetLogo Model using rNetLogo

### Usage

```
NLWrapper.RunExperiment(obj, r = 1, design, FUN)
```

### Arguments

obj	The object retuned by <a href="#">NLWrapper.Model</a>
r	The number of replications
design	The desing matrix holding parameter sampling
FUN	THe calibration function.

### Value

A list containing the the parameters, the calibration functio output and the whole resultset

## Examples

```
## Not run:
rm(list=ls())
objectivefn<- function(params, results) { 0 }

f<- AddFactor(name="initial-number-sheep",min=100,max=250)
f<- AddFactor(factors=f, name="initial-number-wolves",min=50,max=150)
f<- AddFactor(factors=f, name="grass-regrowth-time",min=30,max=100)
f<- AddFactor(factors=f, name="sheep-gain-from-food",min=1,max=50)
f<- AddFactor(factors=f, name="wolf-gain-from-food",min=1,max=100)
f<- AddFactor(factors=f, name="sheep-reproduce",min=1,max=20)
f<- AddFactor(factors=f, name="wolf-reproduce",min=1,max=20)

design<- AoE.LatinHypercube(factors=f)

p<- "C:/Program Files/NetLogo 6.0.4/app"
m<- file.path(p, "models", "Sample Models", "Biology", "Wolf Sheep Predation.nlogo")
output<- c("count sheep", "count wolves")
o<- NLWrapper.Model(p, m, output, 150)
v<- RunExperiment(o, r=1, design, objectivefn)
NLWrapperShutdown(o)

## End(Not run)
```

## NLWrapper.SetParameter

*NLWrapper.SetParameter*

## Description

Set parameter values

## Usage

`NLWrapper.SetParameter(obj, parameters)`

## Arguments

<code>obj</code>	The object retuned by <a href="#">NLWrapper.Model</a>
<code>parameters</code>	The data frame containing the paramters

## Examples

```
## Not run:
rm(list=ls())
p<- "C:/Program Files/NetLogo 6.0.4/app"
m<- file.path(nlpath, "models", "Sample Models", "Biology", "Wolf Sheep Predation.nlogo")
o<- NLWrapper.Model(p, m)
```

```
## End(Not run)
```

---

```
NLWrapper.SetRandomSeed
```

*NLWrapper.SetRandomSeed*

---

### Description

Configures the random seed

### Usage

```
NLWrapper.SetRandomSeed(obj, seed)
```

### Arguments

obj	The object retuned by <a href="#">NLWrapper.Model</a>
seed	The new random seed

---

```
NLWrapperShutdown
```

*NLWrapperShutdown*

---

### Description

This wrapper terminates RNetLogo execution environment

### Usage

```
NLWrapperShutdown(obj)
```

### Arguments

obj	The object retuned by <a href="#">NLWrapper.Model</a>
-----	---

---

ObjectiveFunction-class

*ObjectiveFunction class*

---

**Description**

The base class for optimization functions.

**Fields**

object The raw output of objective function  
objective The objective function  
parameters The parameter list for objective function  
value The results from objective function

---

## Options-class

*Options*

---

**Description**

The base class for the options for the optimization metaheuristics

**Fields**

type The configuration type  
neighborhood The neighborhood function for population methods  
discrete Flag indicating that and specific algorithm is discrete or continuous  
nlevelz Default value for generating parameter levels when range is provided, default value is 5  
container The object holding the configuration options

---

## OptionsACOR-class

*OptionsACOR*

---

**Description**

Options for ACOR method

---

OptionsEES1-class      *OptionsEES1*

---

**Description**

Options for EvoPER Evolutionary Stratety 1

---

OptionsEES2-class      *OptionsEES2*

---

**Description**

Options for Serial Dilutions method

**Fields**

dilutions The desired dilutions

---

OptionsFactory      *OptionsFactory*

---

**Description**

Instantiate the Options class required for the specific metaheuristic method.

**Usage**

`OptionsFactory(type, v = NULL)`

**Arguments**

type	The metaheuristic method
v	The options object

**Value**

Options object

---

OptionsPSO-class      *OptionsPSO*

---

**Description**

Options for PSO optimization metaheuristic

---

OptionsSAA-class      *OptionsSAA*

---

**Description**

Options for SAA method

**Fields**

temperature The temperature dacay function

---

OptionsTS-class      *OptionsTS*

---

**Description**

Options for Tabu search optimization metaheuristic

---

paramconverter      *paramconverter*

---

**Description**

Convert parameter from continuous to discrete and vice-versa if needed

**Usage**

`paramconverter(parameters, discrete, levelz = 5)`

**Arguments**

parameters	The current parameter set
discrete	The desired parameter type
levelz	When discrete is true the number of levels to be generated

**Value**

The parameter collection casted to desired mode

`partSolutionSpace`      *partSolutionSpace*

### Description

Creates the initial Solution population taking into account the lower and upper bounds of provided experiment factors. This method works by dividing the solution space into partitions of size 'd' and then creating a full factorial combination of partitions.

### Usage

```
partSolutionSpace(parameters, d = 4)
```

### Arguments

<code>parameters</code>	The Objective Function parameter list
<code>d</code>	The partition size. Default value 4.

### Value

A set of solutions

`PlainFunction-class`      *PlainFunction*

### Description

PlainFunction Class

`pop.first`      *pop.first*

### Description

pop an element

### Usage

```
pop.first(x)
```

### Arguments

<code>x</code>	The element collection
----------------	------------------------

### Value

The first element added to list FIFO

---

`pop.last`

---

*pop.last*

---

### Description

pop an element

### Usage

`pop.last(x)`

### Arguments

`x`                    The element collection

### Value

The last element added to list LIFO

---

`predatorprey`

---

*predatorprey*

---

### Description

The solver for Lotka-Volterra differential equation.

### Usage

`predatorprey(x1, x2, x3, x4)`

### Arguments

<code>x1</code>	The growth rate of prey
<code>x2</code>	The decay rate of predator
<code>x3</code>	The predating effect on prey
<code>x4</code>	The predating effect on predator

### Value

The ODE solution

`predatorprey.plot0`      *predatorprey.plot0*

### Description

Generate a plot for the predator-prey ODE output.

### Usage

```
predatorprey.plot0(x1, x2, x3, x4, title = NULL)
```

### Arguments

x1	The growth rate of prey
x2	The decay rate of predator
x3	The predating effect on prey
x4	The predating effect on predator
title	The optional plot title. May be omitted.

### Value

An ggplot2 object

### Examples

```
## Not run:
predatorprey.plot0(1.351888, 1.439185, 1.337083, 0.9079049)

## End(Not run)
```

`predatorprey.plot1`      *predatorprey.plot1*

### Description

Simple wrapper for 'predatorprey.plot0' accepting the parameters as a list.

### Usage

```
predatorprey.plot1(x, title = NULL)
```

**Arguments**

- x A list containing the values of predator/prey parameters c1, c2, c3 and c4 denoting respectively the growth rate of prey, the decay rate of predator, the predating effect on prey and the predating effect on predator
- title The optional plot title. May be omitted.

**Value**

An ggplot2 object

**Examples**

```
## Not run:  
rm(list=ls())  
predatorprey.plot1(v$getBest()[1:4])  
  
## End(Not run)
```

---

**pso.best**

---

*pso.best*

---

**Description**

Search for the best particle solution which minimize the objective function.

**Usage**

```
pso.best(objective, particles)
```

**Arguments**

- objective The results of evaluating the objective function
- particles The particles tested

**Value**

The best particle

**psolchi***psolchi***Description**

Implementation of constriction coefficient

**Usage**

```
psolchi(phi1, phi2)
```

**Arguments**

phi1	Acceleration coefficient toward the previous best
phi2	Acceleration coefficient toward the global best

**Value**

The calculated constriction coefficient

**psolbest***psolbest***Description**

Finds the lbest for the particle 'i' using the topology function given by the topology parameter.

**Usage**

```
psolbest(i, pbest, topology)
```

**Arguments**

i	The particle position
pbest	The pbest particle collection
topology	The desired topology function

**Value**

The lbes for i th particle

---

`pso.neighborhood.K2`      *pso.neighborhood.K2*

---

### Description

The neighborhood function for a simple linear topology where every particle has k = 2 neighbors

### Usage

`pso.neighborhood.K2(i, n)`

### Arguments

- |   |                                 |
|---|---------------------------------|
| i | The particle position           |
| n | the size of particle population |

---

`pso.neighborhood.K4`      *pso.neighborhood.K4*

---

### Description

The von neumann neighborhood function for a lattice-based topology where every particle has k = 4 neighbors

### Usage

`pso.neighborhood.K4(i, n)`

### Arguments

- |   |                                 |
|---|---------------------------------|
| i | The particle position           |
| n | the size of particle population |

---

`pso.neighborhood.KN`      *pso.neighborhood.KN*

---

### Description

Simple helper method for 'gbest' neighborhood

### Usage

```
pso.neighborhood.KN(i, n)
```

### Arguments

i	The particle position
n	the size of particle population

---

`pso.printbest`      *pso.printbest*

---

### Description

Shows the best particle of each of simulated generations

### Usage

```
pso.printbest(objective, particles, generation, title)
```

### Arguments

objective	An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a>
particles	The current particle population
generation	The current generation
title	Some informational text to be shown

---

**pso.Velocity**                  *pso.velocity*

---

**Description**

Calculates the PSO Velocity

**Usage**

```
pso.Velocity(W = 1, Vi, phi1, phi2, Pi, Pg, Xi)
```

**Arguments**

W	Weight (Inertia weight or constriction coefficient)
Vi	Current Velocity vector
phi1	Acceleration coefficient toward the previous best
phi2	Acceleration coefficient toward the global best
Pi	Personal best
Pg	Neighborhood best
Xi	Particle vector

**Value**

Updated velocity

---

**push**                  *push*

---

**Description**

push an element

**Usage**

```
push(x, v)
```

**Arguments**

x	The collection of elements
v	The value to be pushed

**Value**

The collection of elements

---

random.wheel	<i>random.whell</i>
--------------	---------------------

---

**Description**

A simple random seed generator

**Usage**

```
random.wheel()
```

**Value**

A random number for seeding

---

RepastFunction-class	<i>RepastFunction</i>
----------------------	-----------------------

---

**Description**

RepastFunction class

---

saa.bolt	<i>saa.bolt</i>
----------	-----------------

---

**Description**

Temperature function boltzmann

**Usage**

```
saa.bolt(t0, k)
```

**Arguments**

t0	The current temperature
k	The annealing value

**Value**

The new temperature

---

saa.neighborhood      *saa.neighborhood*

---

**Description**

Generates neighbor solutions for simulated annealing

**Usage**

```
saa.neighborhood(f, S, d, n)
```

**Arguments**

- |   |  |
|---|--|
| f | An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a> |
| S | The current solution to find a neighbor  |
| d | The distance from current solution S $\text{distance} = (\max - \min) * d$             |
| n | The number of parameters to be perturbed   |

**Value**

The neighbor of solution S

---

saa.neighborhood1      *saa.neighborhood1*

---

**Description**

Generates neighbor solutions perturbing one parameter from current solution S picked randomly.

**Usage**

```
saa.neighborhood1(f, S, d)
```

**Arguments**

- |   |  |
|---|--|
| f | An instance of ObjectiveFunction (or subclass) class <a href="#">ObjectiveFunction</a> |
| S | The current solution to find a neighbor  |
| d | The distance from current solution S $\text{distance} = (\max - \min) * d$             |

**Value**

The neighbor of solution of S

`saa.neighborhoodH`      *saa.neighborhoodH*

### Description

Generates neighbor solutions perturbing half parameters from current solution S.

### Usage

`saa.neighborhoodH(f, S, d)`

### Arguments

- f                  An instance of ObjectiveFunction (or subclass) class [ObjectiveFunction](#)
- S                  The current solution to find a neighbor
- d                  The distance from current solution S  $\text{distance} = (\max - \min) * d$

### Value

The neighbor of solution of S

`saa.neighborhoodN`      *saa.neighborhoodN*

### Description

Generates neighbor solutions perturbing all parameters from current solution S.

### Usage

`saa.neighborhoodN(f, S, d)`

### Arguments

- f                  An instance of ObjectiveFunction (or subclass) class [ObjectiveFunction](#)
- S                  The current solution to find a neighbor
- d                  The distance from current solution S  $\text{distance} = (\max - \min) * d$

### Value

The neighbor of solution of S

---

`saa.tbyk`*saa.tbyk*

---

**Description**

Temperature function t/k

**Usage**

`saa.tbyk(t0, k)`

**Arguments**

<code>t0</code>	The current temperature
<code>k</code>	The annealing value

**Value**

The new temperature

---

`saa.tcte`*saa.tcte*

---

**Description**

Temperature function cte \* t0

**Usage**

`saa.tcte(t0, k)`

**Arguments**

<code>t0</code>	The current temperature
<code>k</code>	The annealing value

**Value**

The new temperature

---

`saa.texp`*saa.texp*

---

**Description**

Temperature function exponential

**Usage**

```
saa.texp(t0, k)
```

**Arguments**

t0	The current temperature
k	The annealing value

**Value**

The new temperature

---

`scatterplotlothelper`    *scatterplotlothelper*

---

**Description**

Simple helper for plotting 3d scatterplots

**Usage**

```
scatterplotlothelper(d, x, y, z, title = NULL)
```

**Arguments**

d	A data frame.
x	A string with the dataframe column name for x axis
y	A string with the dataframe column name for y axis
z	A string with the dataframe column name for z axis
title	The optional plot title. May be omitted.

**Value**

A scatter3D plot

---

`searchrow`*searchrow*

---

**Description**

Search for a value value on a matrix

**Usage**

```
searchrow(ddata, value)
```

**Arguments**

ddata	The matrix containing the dataset
value	The value to search for

**Value**

Boolean TRUE for those indexes matching value

---

---

`show.comp1`*show.comp1*

---

**Description**

Generates a barplot comparing the number of evalutions for algorithms ("saa","pso","acor","ees1").

**Usage**

```
show.comp1(mydata, what, title = NULL)
```

**Arguments**

mydata	The data generated with 'summarize.comp1'
what	The name of variable to plot on 'y' axis
title	the plot title

**Examples**

```
## Not run:  
p.a<- show.comp1(d.cigar4,"evals","(a) Cigar function")  
p.b<- show.comp1(d.schaffer4,"evals","(b) Schafer function")  
p.c<- show.comp1(d.griewank4,"evals","(c) Griewank function")  
p.d<- show.comp1(d.bohachevsky4,"evals","(d) Bohachevsky function")  
  
## End(Not run)
```

---

slope	<i>slope</i>
-------	--------------

---

**Description**

Simple function for calculate the slope on the ith element position

**Usage**

```
slope(x, y, i)
```

**Arguments**

x	The x vector
y	The y vector
i	The position

**Value**

The slope

---

slopes	<i>slopes</i>
--------	---------------

---

**Description**

Calcule all slopes for the discrete x,y series

**Usage**

```
slopes(x, y)
```

**Arguments**

x	The x vector
y	The y vector

**Value**

A vector with all slopes

---

`sortSolution`*sortSolution*

---

**Description**

Sort solution by its respective fitness

**Usage**

```
sortSolution(s, f)
```

**Arguments**

s                  Problem solution

f                  The function evaluation for s

---

---

`summarize.comp1`*summarize.comp1*

---

**Description**

Provides as summary with averaged values of experimental setup

**Usage**

```
summarize.comp1(mydata)
```

**Arguments**

mydata              The data frame generated with 'compare.algorithms1'

**Value**

The summarized data

---

tabu.getNeighbors      *tabu.getNeighbors*

---

**Description**

create neighbor solutions

**Usage**

`tabu.getNeighbors(tabu, parameters, solution, size)`

**Arguments**

tabu	The tabu list
parameters	The parameter set
solution	The current solution
size	The neighborhood size

**Value**

The neighbor for solution

---

tabu.istabu      *tabu.istabu*

---

**Description**

Check whether a solution is present on tabulist

**Usage**

`tabu.istabu(tabulist, solution)`

**Arguments**

tabulist	The matrix of tabu solutions
solution	The solution value to be checked

**Value**

Boolean TRUE tabulist contains the solution

---

upperBound

*upperBound*

---

### Description

Checks if parameters is below the upper bounds

### Usage

```
upperBound(particles, factors)
```

### Arguments

particles	The particle set
factors	the defined range for objective function parameters

### Value

The particle inside the valid upper bound

---

xmeanci1

*xmeanci1*

---

### Description

Calculates confidence interval of mean for provided data with desired confidence level. This functions uses bootstrap resampling scheme for estimanting the CI.

### Usage

```
xmeanci1(x, alpha = 0.95)
```

### Arguments

x	The data set for which CI will be calculated
alpha	The confidence level. The default value is 0.95 (95%)

### Value

The confidence interval for the mean calculated using 'boot.ci'

---

`xmeanci2`*xmeanci2*

---

**Description**

Calculates confidence interval of mean for provided data with desired confidence level.

**Usage**

```
xmeanci2(x, alpha = 0.95)
```

**Arguments**

- |                    |   |
|--------------------|---|
| <code>x</code>     | The data set for which CI will be calculated          |
| <code>alpha</code> | The confidence level. The default value is 0.95 (95%) |

**Value**

The confidence interval for the mean

---

`xyplot helper`*xyplot helper*

---

**Description**

Simple helper for plotting xy dispersion points.

**Usage**

```
xyplot helper(d, x, y, title = NULL)
```

**Arguments**

- |                    |  |
|--------------------|--|
| <code>d</code>     | A data frame.                                      |
| <code>x</code>     | A string with the dataframe column name for x axis |
| <code>y</code>     | A string with the dataframe column name for y axis |
| <code>title</code> | The optional plot title. May be omitted.           |

**Value**

A ggplot2 plot object

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