

Package ‘OpenMORDM’

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Type Package

Title A library for multiobjective robust decision making (MORDM) and visualizing high-dimensional datasets.

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Description OpenMORDM provides support for visualizing high-dimensional datasets loaded from matrices or data.frames and a variety of file formats including CSV, XLS, XLSX, and MOEA runtime outputs (from the Borg MOEA or MOEA Framework). This visualization is contained within a web-based viewer capable of generating various 2D and 3D plots as well as performing a number of analyses. Additionally, the R functions provide the means for evaluating models under uncertainty and deep uncertainty and computing different robustness metrics.

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Depends R (>= 3.0)

Imports

shiny,shinyRGL,rgl,scales,grid,prim,MASS,animation,sensitivity,boot,pracma,emoa,stringr,functional,dichromat,rpart

Suggests gdata

R topics documented:

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OpenMORDM-package	<i>A library for multiobjective robust decision making (MORDM) and visualizing high-dimensional datasets.</i>
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Description

OpenMORDM provides support for visualizing high-dimensional datasets loaded from matrices or data.frames and a variety of file formats including CSV, XLS, XLSX, and MOEA runtime outputs (from the Borg MOEA or MOEA Framework). This visualization is contained within a web-based viewer capable of generating various 2D and 3D plots as well as performing a number of analyses. Additionally, the R functions provide the means for evaluating models under uncertainty and deep uncertainty and computing different robustness metrics.

Details

Package: OpenMORDM
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 Depends: R (>= 3.0)
 Imports: shiny, shinyRGL, rgl, scales, grid, prim, MASS, animation, sensitivity, boot, pracma, emoa, stringr, function
 Suggests: gdata
 Built: R 3.1.1; ; 2015-05-07 20:21:56 UTC; windows

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<code>compute.robustness</code>	Computes robustness under deep <code>compute.robustness</code> .
<code>compute.robustness.gaussian</code>	Computes robustness under well-characterized <code>compute.robustness</code> (i.e., Gaussian noise).
<code>compute.sensitivity</code>	Standardized interface for sensitivity analysis methods.
<code>define.problem</code>	Define a new problem formulation.
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<code>mordm.cbind</code>	Adds extra columns to the end of a data set.
<code>mordm.colorize</code>	Returns a vector of colors to be used when plotting a data set.
<code>mordm.correlation</code>	Displays the correlations among pairwise factors.
<code>mordm.differences</code>	Identifies key similarities/differences between two sets.
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<code>sensitivity.levels</code>	Determines number of replicates for sensitivity analysis.
<code>usample</code>	Generate uniformly distributed random inputs.

Author(s)

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<code>adjust.command</code>	<i>Prepends a ./ to commands on non-Windows systems.</i>
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Description

Prepends a ./ to commands on non-Windows systems.

Usage

```
adjust.command(command)
```

Arguments

<code>command</code>	the R function or executable representing the problem
----------------------	---

<code>analyze.cart</code>	<i>Determines the vulnerabilities due to deep uncertainties using Classification and Regression Trees (CART).</i>
---------------------------	---

Description

Determines the vulnerabilities due to deep uncertainties using Classification and Regression Trees (CART).

Usage

```
analyze.cart(factors, response)
```

Arguments

<code>factors</code>	the sampled deeply uncertain parameterizations
<code>response</code>	vector of responses whose length equals the number of factors

<code>analyze.prim</code>	<i>Patient rule induction method.</i>
---------------------------	---------------------------------------

Description

Performs the patient rule induction method (PRIM) to identify boxes in input space that correlate with data exceeding a given threshold.

Determines the vulnerabilities due to deep uncertainties using the Patient Rule Induction Method (PRIM).

Usage

```
analyze.prim(factors, response, bounds = NULL, which.box = 1,
             show.plot = TRUE, ...)
```

```
analyze.prim(factors, response, bounds = NULL, which.box = 1,
             show.plot = TRUE, ...)
```

Arguments

factors	the sampled deeply uncertain parameterizations
response	vector of responses whose length equals the number of factors
bounds	bounds of the sampled uncertainties
which.box	index of the PRIM box to plot
show.plot	if TRUE, generates a plot representing the PRIM box
...	optional arguments passed to prim.box
data	the data set
objective	specifies the objective index, column name, function, or marking to use
minimize	if TRUE, flip the threshold direction so that smaller values are preferred
percentages	display percentages in the printout
expand	if TRUE, sets the paste option to 1 to enable expanding the boxes to fill as much space as possible
...	optional parameters passed to prim.box

borg.optimize	<i>Optimize the problem using the Borg MOEA.</i>
---------------	--

Description

Optimizes the problem. By default, this method uses the Borg MOEA, which must first be compiled into an executable or shared library on your system. If the problem references an R function, then you must have available the shared library (borg.dll or libborg.so). If the problem is an external program, then you must have available the Borg executable (borg.exe). See [borg.optimize.function](#) and [borg.optimize.external](#) for details of each method.

Usage

```
borg.optimize(problem, NFE, ...)
```

Arguments

problem	the problem definition
NFE	the maximum number of function evaluations
...	optional parameters passed to the underlying methods

Details

The Borg MOEA is free and open for non-commercial users. Source code can be obtained from <http://borgmoea.org>.

```
borg.optimize.external
```

Optimize the problem using the Borg standalone executable (borg.exe).

Description

This method is used to optimize a problem defined by an external executable. The Borg MOEA communicates with the external executable using the open API standardized by the MOEA Framework (<http://moeaframework.org>). See section 5.2 in the user manual for details of using the API. Since borg.exe targets POSIX systems, this method is typically not available on Windows unless you are running inside Cygwin. See the Borg MOEA documentation for instructions on compiling borg.exe.

Usage

```
borg.optimize.external(problem, NFE, executable = "./borg.exe",
  output = tempfile(), output.frequency = 100, return.output = TRUE,
  verbose = TRUE)
```

Arguments

problem	the problem definition
NFE	the maximum number of function evaluations
executable	the path the the optimization executable
output	the location where the runtime output is stored
output.frequency	the frequency at which data is output
return.output	if TRUE, this method loads and returns the contents of the output file
verbose	displays additional information for debugging

Details

The Borg MOEA is free and open for non-commercial users. Source code can be obtained from <http://borgmoea.org>.

```
borg.optimize.function
```

Optimize a problem using the Borg shared library (borg.dll or lib-borg.so).

Description

This method is used to optimize a problem defined by an R function. This method uses R's foreign function interface (FFI) to pass the R function to the Borg MOEA shared library for optimization. See the Borg MOEA documentation for instructions on compiling borg.dll or libborg.so.

Usage

```
borg.optimize.function(problem, NFE, ...)
```

Arguments

problem	the problem definition
NFE	the maximum number of function evaluations
...	additional arguments for setting algorithm parameters

Details

The function should either return a vector containing the objectives and any constraints (e.g., `c(o1, o2, o3, c1, c2)`), or a list containing the objectives and constraints as separate elements (e.g., `list(c(o1, o2, o3), c(c1, c2))`). All objectives are minimized. Any non-zero constraint value is considered a constraint violation.

The Borg MOEA is free and open for non-commercial users. Source code can be obtained from <http://borgmoea.org>.

check.robustness	<i>Computes the robustness metric.</i>
------------------	--

Description

Robustness is represented as a scalar value, where values nearer to positive infinity are considered more robust. Due to differences in how each robustness metric computes its value, you should look at relative differences in values rather than absolute differences.

Usage

```
check.robustness(output, problem, method = "default", verbose = FALSE, ...)
```

Arguments

output	the evaluated points
problem	the problem definition
method	the robustness metric to use (default, variance, constraints, infogap, or distance)
verbose	display additional information
...	additional arguments passed to the robustness metric

compute.robustness	<i>Computes robustness under deep compute.robustness.</i>
--------------------	---

Description

Adds Gaussian noise to the decision variables and resamples the model output. The samples are distributed across one or more different models for the problem. The result from this method should be passed to `mordm.evaluate.uncertainties` to compute the robustness metrics.

Usage

```
compute.robustness(data, nsamples, models, sd = 0, verbose = TRUE,
  satisficing.fcn = NULL, factors = NULL, custom.fcn = NULL)
```

Arguments

<code>data</code>	the data set
<code>nsamples</code>	the number of samples to generate for each point
<code>models</code>	the problem formulations created using <code>define.problem</code>
<code>sd</code>	scalar or vector specifying the standard deviation for each decision variable
<code>verbose</code>	display additional information
<code>satisficing.fcn</code>	the satisficing function for computing the two satisficing robustness metrics
<code>factors</code>	matrix of the original <code>compute.robustness</code> factors for use by Satisficing Type II
<code>custom.fcn</code>	custom robustness function

Details

If multiple models are provided, it is assumed that all models have the same inputs and outputs; they would only differ in the internal calculations within the model.

compute.robustness.guassian	<i>Computes robustness under well-characterized compute.robustness (i.e., Gaussian noise).</i>
-----------------------------	--

Description

Adds Gaussian noise to the decision variables and resamples the model output. Then computes one or more robustness metrics.

Usage

```
compute.robustness.guassian(data, sd, nsamples, problem, method = "default",
  verbose = TRUE)
```

Arguments

data	the data set
sd	scalar or vector specifying the standard deviation for each decision variable
nsamples	the number of samples to generate for each point
problem	the problem formulation
method	the robustness metric or a list of metrics to use (see check.robustness for available options)
verbose	display additional information

Details

This method is equivalent to [mordm.compute.robustness](#) using a single model.

compute.sensitivity	<i>Standardized interface for sensitivity analysis methods.</i>
---------------------	---

Description

Attempts to standardize the use of various sensitivity analysis methods. Supports all of the methods provided by the sensitivity library except for those using metamodels.

Usage

```
compute.sensitivity(problem, objective, nsamples, method = "fast99",
  verbose = FALSE, plot = FALSE, raw = FALSE, collapse = TRUE, ...)
```

Arguments

problem	the problem definition
objective	the function, objective index, or objective name whose sensitivity is being computed
nsamples	the desired number of samples
method	string representation of the sensitivity analysis method (fast99, sobol, sobol2002, sobol2007, sobolEff, soboljansen, sobolmara, sobolroahs, morris, prc, src, or plischke)
verbose	if TRUE, print additional information
plot	if TRUE, generate any output plots
raw	if TRUE, return the raw model output; otherwise return the standardized output
collapse	if TRUE, collapses the list representation of the variables, objectives, and constraints into a matrix representation
...	additional options passed to the sensitivity analysis method

Details

In addition to using the same inputs for each method, the outputs are also standardized. For methods computing the first-order indices, the output contains the sensitivity indices (S_i) and a ranking (rank). Methods computing total-order indices, the output contains the total sensitivity indices ($S_i.total$) and the ranking (rank.total). Where available, the output may also contain confidence intervals (C_i and $C_i.total$).

define.problem	<i>Define a new problem formulation.</i>
----------------	--

Description

Constructs a new problem formulation. The command can either be an R function or a command line executable. If using a command line executable, the program must follow the MOEA Framework external problem protocol, typically by using the methods in `moeaframework.h`.

Usage

```
define.problem(command, nvars, nobjs, nconstrs = 0, bounds = NULL,
               names = NULL, epsilons = NULL, maximize = NULL)
```

Arguments

command	the R function or executable representing the problem
nvars	the number of decision variables
nobjjs	the number of objectives
nconstrs	the number of constraints
bounds	the lower and upper bounds for each decision variable
names	override the column names
epsilons	the epsilon values if using Borg to optimize the problem
maximize	vector indicating the columns to be maximized

Details

If using an R function, the function should return a list containing two vectors. The first vector stores the objective values and the second stores the constraint values.

deltafast	<i>Fast version of Plischke's delta-moment sensitivity analysis method.</i>
-----------	---

Description

Fast version of Plischke's delta-moment sensitivity analysis method.

Usage

```
deltafast(x, y, M = 24)
```

Arguments

x	the factors
y	the response
M	the number of partitions

deltamim	<i>Plischke's delta-moment sensitivity analysis method.</i>
----------	---

Description

Plischke's delta-moment sensitivity analysis method.

Usage

```
deltamim(x, y, partition.size = min(ceiling(nrow(x)^(2/(7 + tanh((1500 -
  nrow(x))/500))))), 48), quadrature.points = 110, ks.level = 0.95,
  zero.crossing = "on", kd.estimator = "cheap", kd.width = "auto",
  complement = FALSE, plot.enabled = FALSE, plot.cols = min(ncol(x), 4),
  output.trafo = "off", kd.shape = "epanechnikov", ...)
```

Arguments

x	the factors
y	the response
partition.size	the number of partitions
quadrature.points	the number of points in the estimated PDFs
ks.level	critical value for the KS statistic
zero.crossing	detect zero crossings when "on"
kd.estimator	the kernel density estimator to use (cheap, stats, diffusion, hist)
kd.width	the bandwidth for kernel density estimation, or "auto"
complement	compute sensitivities for the complement
plot.enabled	if TRUE, generate a plot showing the PDFs
plot.cols	the number of factors to display in the plot
output.trafo	transformation applied to the responses (off, cdf, normal, interpol, cdf-tight, cdf-loose)
kd.shape	kernel shape (normal, triangle, epanechnikov, box, uniform, biweight, biquadratic) or a function defining the kernel shape
...	additional options

evaluate	<i>Evaluates the decision variables for a given problem.</i>
----------	--

Description

Evaluates the problem using the given decision variables, returning an object storing the variables, objectives, and constraints.

Usage

```
evaluate(set, problem)
```

Arguments

set	the decision variables (inputs) to the problem
problem	the problem definition

explore	<i>A web-based tool (powered by Shiny) for exploring high-dimensional data sets.</i>
---------	--

Description

Starts a Shiny server and launches a webpage to display the given dataset. The Shiny server will remain running even if the web browser is closed. You must interrupt the current R function call (press Ctrl-C or select Session -> Interrupt R in RStudio) to stop the server.

Usage

```
explore(filename, nvars = NULL, nobjs = NULL, nconstrs = 0,
        names = NULL, bounds = NULL, maximize = NULL, order = NULL,
        visible.variables = FALSE, plot3d.width = "600px",
        plot3d.height = "500px", welcome.panel = NULL, selection.panel = NULL,
        ignore = NULL, metadata = NULL, runShinyApp = TRUE)
```

Arguments

filename	the name of the file, a matrix, or a data frame
nvars	the number of decision variables
nobjs	the number of objectives
nconstrs	the number of constraints
names	see mordm.read for details
bounds	see mordm.read for details
maximize	see mordm.read for details
order	ordering of the objectives in the dropdown menus
visible.variables	determines if variables are visible by default
plot3d.width	the width of the 3D window
plot3d.height	the hight of the 3D window
welcome.panel	omordm.plotptional panel for displaying a intro message
selection.panel	optional panel for displaying info about the selected point
ignore	columns to remove from the dataset (CSV/Excel only)
metadata	columns to retain in a metadata attribute (CSV/Excel only)
runShinyApp	if TRUE, start the Shiny server with runApp(...); if false, start the server with shinyApp(...)

Details

This method currently supports loading CSV files (.csv), Excel files (.xls or .xlsx), and MOEA runtime files (any other extension). For CSV and Excel files, you may optionally specify nvars and/or nobjs. If unspecified, the method will assume every column is an objective. For MOEA runtime files, nvars and nobjs are mandatory.

For CSV and Excel files, you can optionally ignore certain columns or indicate a column is metadata. Ignored columns are completely removed from the analysis. Metadata is not shown in the visualizations, but will be saved in an attribute (`attr(data[[i]], "metadata")`). For example, if each point has an associated GIF animation, you could treat the column storing the file path as metadata, allowing you to retrieve and display the GIF.

If you are providing a custom welcome or selection panel and would like to display custom resources, use [addResourcePath](#) to register the directory containing the resources.

lhsample	<i>Generate Latin Hypercube sampled random inputs.</i>
----------	--

Description

Generate Latin Hypercube sampled random inputs.

Usage

```
lhsample(nsamples, problem)
```

Arguments

nsamples	the number of samples to generate
problem	the problem definition

mordm.animate	<i>Animates the time series in a GIF.</i>
---------------	---

Description

Animates the 3D scatter plot and saves the results to a GIF file. Each index in indices specifies the entry in the time series data that is displayed at each frame. Thus, to show all entries in succession, set `indices=1:length(data)`, but to show a single entry across multiple frames (e.g., if rotating) use `indices=rep(length(data), 50)`. The transform function is similar to the transformation function used by [play3d](#). However, whereas the transformation in `play3d` is based on the number of elapsed seconds, this method computes the transform based on the frame number. Thus, if you use `spin3d(rpm=1)`, then this method will rotate the plot once every 60 frames (i.e., treating each frame as one second).

Usage

```
mordm.animate(data, output = "animation.gif", indices = 1:length(data),
  transform = NULL, clean = TRUE, close = TRUE, loop = FALSE,
  scale = 0.1, ...)
```

Arguments

data	the time series data
output	the location where the animated GIF is saved
indices	a vector indicating the indices in data that are displayed in each frame
transform	function that returns a transformation applied to each frame (see spin3d), or the user matrix for a constant projection
clean	if TRUE, delete the temporary images once the GIF is created
close	if TRUE, close the RGL window when finished
loop	if TRUE, loop infinitely; otherwise play the animation once
scale	amount to enlarge the plotting limits
...	additional arguments passed to mordm.plot

mordm.as.data.frame	<i>Converts a data set into a data frame.</i>
---------------------	---

Description

When loading data into OpenMORDM, any non-numeric columns are converted into factors and represented internally as integers. This method reverses that process to get a data frame storing the original values.

Usage

```
mordm.as.data.frame(entry)
```

Arguments

entry	the data set to convert
-------	-------------------------

mordm.cbind	<i>Adds extra columns to the end of a data set.</i>
-------------	---

Description

The added columns are treated like objectives.

Usage

```
mordm.cbind(set, columns)
```

Arguments

set	the data set
columns	the extra columns

mordm.colorize	<i>Returns a vector of colors to be used when plotting a data set.</i>
----------------	--

Description

The color data that is plotted depends on the options given to this function in the following order:

- colors- displays the user-defined color values
- mark- displays the user-defined marking rules, each rule with a separate color
- objectives- use color values stored in the column defined by objective[5]

Usage

```
mordm.colorize(set, objectives, mark = NULL, palette = NULL, n = 100,
  offset = 0, colors = NULL, clim = NULL, unmarked = "#888888FF",
  alpha = 1, crev = TRUE)
```

Arguments

set	the data set
objectives	the objectives being plotted (objectives[5] is color)
mark	the optional marking rule
palette	the color palette to use, either a function that generates the color palette or the color palette itself
n	the number of distinct colors to display (only used if palette is a function)
offset	DEPRECATED
colors	user-defined color values
clim	range (lower and upper bounds) of color values
unmarked	the color value used for unmarked points
alpha	the transparency applied to all colors
crev	if TRUE, reverse the color palette

mordm.correlation	<i>Displays the correlations among pairwise factors.</i>
-------------------	--

Description

Computes the pairwise correlations between the decision variables and objectives and prints the formatted results.

Usage

```
mordm.correlation(data, ht = 0.75, lt = 0.25, all = FALSE,
  objectives = FALSE)
```

Arguments

data	the data set to use
ht	the threshold for highly correlated pairs
lt	the threshold for uncorrelated pairs
all	show all correlations
objectives	only compute correlations between objectives

mordm.differences	<i>Identifies key similarities/differences between two sets.</i>
-------------------	--

Description

Computes key differences between two data sets. This method prints formatted text as well as creates a histogram plot showing key differences.

Usage

```
mordm.differences(set1, set2, scale = TRUE, decreasing = TRUE,
  splits = 20, n = NULL, all = FALSE)
```

Arguments

set1	the first set
set2	the second set
scale	if TRUE, scale the plot based on the range of the data
decreasing	if TRUE, order differences in decreasing order
splits	the number of bins used by the histogram method
n	the number of variables to plot
all	plot all variables

mordm.evaluate.uncertainties	<i>Computes robustness under deep compute.robustness.</i>
------------------------------	---

Description

Computes robustness under deep compute.robustness.

Usage

```
mordm.evaluate.uncertainties(samples, satisficing.fcn = NULL,
  factors = NULL, custom.fcn = NULL)
```

Arguments

samples	the samples generated by <code>mordm.sample.uncertainties</code>
satisficing.fcn	the satisficing function for computing the two satisficing robustness metrics
factors	matrix of the original <code>compute.robustness</code> factors for use by Satisficing Type II
custom.fcn	custom robustness function

```
mordm.extract.attributes
```

Extracts the common attributes from the time series data.

Description

Reads the common attributes associated with each entry in the time series data and returns them in a matrix.

Usage

```
mordm.extract.attributes(data)
```

Arguments

data	the time series data
------	----------------------

```
mordm.get.set
```

Returns the individual data set entry to be analyzed.

Description

Determines which data set will be analyzed. Almost all methods use this function to convert their data arguments into a single data set. The data argument can be either a time series or an individual data set.

Usage

```
mordm.get.set(data, index = -1)
```

Arguments

data	a time series or individual data set
index	if data is a time series, specifies which entry to return (default is the last entry)

mordm.get.subset	Returns a subset of a data set.
------------------	---------------------------------

Description

Returns a subset of a data set.

Usage

```
mordm.get.subset(set, columns = 1:ncol(set), rows = 1:nrow(set))
```

Arguments

set	the data set
columns	the columns to retain
rows	the rows to retain

mordm.identify	Identify and highlight points using the middle mouse button.
----------------	--

Description

Enables a mouse callback for clicking points on the 3D scatter plot and identifying those points. If any secondary plots are displayed (e.g., parallel coordinates plot or marking plot), the selected point will become highlighted in those plots.

Usage

```
mordm.identify(enabled = TRUE, label = FALSE)
```

Arguments

enabled	if TRUE, enables this functionality
label	if TRUE, a label will be added to the 3D scatter plot identifying the selected point

mordm.mark.box	<i>Creates a marking rule from PRIM boxes.</i>
----------------	--

Description

PRIM identifies one or more boxes. This method converts from the PRIM box representation to a marking.

Usage

```
mordm.mark.box(box, mean, mass)
```

Arguments

box	the box generated by analyze.prim
mean	the mean of the box
mass	the mass of the box

mordm.mark.difference	<i>Computes the difference of two markings.</i>
-----------------------	---

Description

Markings behave like sets. A point is contained within the difference if it is contained in exactly one individual marking. This is similar to the exclusive-or operator.

Usage

```
mordm.mark.difference(...)
```

Arguments

...	the markings
-----	--------------

mordm.mark.intersection	<i>Computes the intersection of two markings.</i>
-------------------------	---

Description

Markings behave like sets. A point is contained within the intersection if it is contained in all individual markings.

Usage

```
mordm.mark.intersection(...)
```

Arguments

...	the markings
-----	--------------

mordm.mark.not	<i>Computes the inverse of a marking.</i>
----------------	---

Description

Markings behave like sets. A point is contained within the inverse if it is not contained within the original marking.

Usage

```
mordm.mark.not(rule)
```

Arguments

rule	the original marking
------	----------------------

mordm.mark.points	<i>Creates a marking rule identifying specific points.</i>
-------------------	--

Description

Markings allow the user to highlight specific subsets of the data set. These marked sets can subsequently be plotted or used in supported calculations.

Usage

```
mordm.mark.points(points)
```

Arguments

points	one or more rows from the data set considered within the marking
--------	--

mordm.mark.rule	<i>Creates a marking rule based on a function.</i>
-----------------	--

Description

Markings allow the user to highlight specific subsets of the data set. These marked sets can subsequently be plotted or used in supported calculations.

Usage

```
mordm.mark.rule(condition)
```

Arguments

condition	a function of the form $f: x \rightarrow \text{boolean}$, where x is a single row from the data set, returning TRUE if the row is part of the marking
-----------	--

mordm.mark.selection	<i>Creates a marking rule identifying the selected points.</i>
----------------------	--

Description

Allows the user to select a rectangular region in the 3D scatter plot and returns a marking containing all points within the selected region.

Usage

```
mordm.mark.selection()
```

mordm.mark.subtract	<i>Computes the subtraction of two markings.</i>
---------------------	--

Description

Markings behave like sets. A point is contained within the subtraction if it is contained in rule1 but not rule2.

Usage

```
mordm.mark.subtract(rule1, rule2)
```

Arguments

rule1	the first marking
rule2	the second marking

mordm.mark.union	<i>Computes the union of two markings.</i>
------------------	--

Description

Markings behave like sets. A point is contained in the union if it is contained within any individual marking.

Usage

```
mordm.mark.union(...)
```

Arguments

...	the markings
-----	--------------

mordm.normalize	<i>Normalizes the objectives.</i>
-----------------	-----------------------------------

Description

By default, the objectives are all minimized. Maximized objectives are negated. This function negates the maximized objectives, returning them to their original, positive form.

Usage

```
mordm.normalize(data, maximize)
```

Arguments

data	the data set
maximize	vector indicating the columns to be maximized

mordm.plot	<i>Sets the current data set and displays a 3D scatter plot.</i>
------------	--

Description

Creates a 3D scatter plot of the data. As a side effect, this method sets a global variable identifying the current data set and plotting attributes to be used by other plotting methods in this package. This design allows you to easily create secondary plots that are consistent with the primary 3D scatter plot.

Usage

```
mordm.plot(data, mark = NULL, index = -1, objectives = NULL,
  stay = TRUE, identify = TRUE, colors = NULL, clim = NULL,
  ideal = FALSE, selection = NULL, selection.enlarge = FALSE,
  xlim = NULL, ylim = NULL, zlim = NULL, slim = NULL, window = NULL,
  alpha = 1, tick.size = 1, label.size = 1.2, label.line = 1,
  radius.scale = 1, bg = "white", fg = "black", exploring = FALSE, ...)
```

Arguments

data	the data set to be displayed (if data is a time series, then the last entry in the time series is displayed)
mark	a list of the markings to be displayed
index	if data is a time series, controls which entries to display (see mordm.get.set for details)
objectives	vector specifying the objectives to be plotted on the x, y, z, size, and color axes
stay	forces the 3D scatter plot to stay on top of other windows
identify	if TRUE, clicking a point with the middle mouse button will identify and highlight that point

colors	override the color values
clim	range (lower and upper bounds) of color values
ideal	draw a visual indicator of where the ideal point is on the plot
selection	draw a visual indicator on the given row indices
selection.enlarge	if TRUE, enlarges the selected point; otherwise renders a transparent cube around the selected point
xlim	range (lower and upper bounds) for the x axis
ylim	range (lower and upper bounds) for the y axis
zlim	range (lower and upper bounds) for the z axis
slim	range (lower and upper bounds) of size values
window	the window size (w, h)
alpha	vector of transparency values applied to each point
tick.size	the size of the tick labels
label.size	the size of axis labels
label.line	the offset of the labels
radius.scale	scaling factor applied to the size of the points
bg	background color
fg	foreground color
exploring	set to TRUE if being called from explore(...) method to properly open null devices
...	additional options passed to plot3d

mordm.plot.box	<i>Display plot of PRIM boxes.</i>
----------------	------------------------------------

Description

Generates a plot showing the bounds of the PRIM boxes. Currently only works well with one or two PRIM boxes.

Usage

```
mordm.plot.box(data, mark, main = "PRIM Box", scale.width = TRUE,
  bar.width = 3, col = NULL, names = NULL, legend = TRUE,
  defaults = NULL)
```

Arguments

data	the original data set
mark	a list of the PRIM boxes to display
main	the plot title
scale.width	if TRUE, reduce the width of the bars as more PRIM boxes are displayed
bar.width	the width of the bars
col	vector of bar colors
names	names of each PRIM box to display in the legend
legend	if TRUE, renders a legend on the plot
defaults	draw horizontal lines to show default values

mordm.plot.markings	<i>Display the markings in a box plot (candle stick plot).</i>
---------------------	--

Description

Creates a box plot showing the range (lower and upper bounds) encompassed by each marking.

Usage

```
mordm.plot.markings(highlight = NULL)
```

Arguments

highlight	highlight vector of row indices to be highlighted in the plot
-----------	---

mordm.plot.operators	<i>Plot operator probabilities.</i>
----------------------	-------------------------------------

Description

Creates a plot of the Borg MOEA operator probabilities across an entire time series.

Usage

```
mordm.plot.operators(data, time = FALSE, improvements = FALSE,
  log = FALSE, improvement.nfe = 1000, current = NULL)
```

Arguments

data	the time series data from a Borg MOEA run
time	if TRUE, the x-axis displays elapsed time; otherwise the x-axis represents the number of function evaluations (NFE)
improvements	if TRUE, displays a trace of the number of Pareto improvements
log	if TRUE, plot the x-axis with a log scale
improvement.nfe	the window size in NFE when computing the number of improvements
current	draw a line at the current time

mordm.plot.parallel	<i>Display a parallel plot of the current data set.</i>
---------------------	---

Description

Creates a parallel axis or parallel coordinates plot of the current data set. All display attributes are taken from the current plotting options.

Usage

```
mordm.plot.parallel(highlight = NULL, alpha = 0.4, label.size = 1,  
  line.width = 1, selection.scale = 2)
```

Arguments

highlight	vector of row indices to be highlighted in the plot
alpha	the transparency value; or NA
label.size	the font size of labels
line.width	the width of lines
selection.scale	the

mordm.print.box	<i>Print descriptive representation of PRIM boxes.</i>
-----------------	--

Description

Displays the PRIM boxes in a human-readable way.

Usage

```
mordm.print.box(data, mark, threshold = 0.01, digits = 3, indent = "")
```

Arguments

data	the original data set
mark	the PRIM boxes
threshold	fuzzy factor when determining if two numbers are equal
digits	number of digits to round numbers
indent	character string prepended to each line

mordm.rbind	<i>Adds extra rows to the end of a data set.</i>
-------------	--

Description

Adds extra rows to the end of a data set.

Usage

```
mordm.rbind(set, rows)
```

Arguments

set	the data set
rows	the extra rows

mordm.read	<i>Loads the time series data output from an optimization algorithm.</i>
------------	--

Description

Reads the time series data (a list of matrices) from an optimization algorithm. The format is defined by the Borg MOEA and MOEA Framework.

Usage

```
mordm.read(file, nvars, nobjs, nconstrs = 0, bounds = NULL, names = NULL,
  maximize = NULL, digits = NULL)
```

Arguments

file	the file name
nvars	the number of decision variables
nobjs	the number of objectives
nconstrs	the number of constraints
bounds	the lower and upper bounds of each decision variable
names	override the column names
maximize	vector indicating the columns to be maximized
digits	number of digits to retain

mordm.read.csv	<i>Loads a data set stored in a CSV file.</i>
----------------	---

Description

This method assumes the first N columns are decision variables, and all other columns are objectives. N is determined by `ncol(bounds)`. Unless overridden, this method sets `check.names=FALSE` and `header=TRUE`.

Usage

```
mordm.read.csv(file, nvars = NULL, nobjs = NULL, bounds = NULL,
               maximize = NULL, names = NULL, ignore = NULL, metadata = NULL, ...)
```

Arguments

file	the file name
nvars	the number of decision variables
nobjs	the number of objectives
bounds	the lower and upper bounds of each decision variable
maximize	vector indicating the columns to be maximized
names	override the column names
ignore	columns to remove from the dataset
metadata	columns to retain in a metadata attribute
...	optional arguments passed to <code>read.csv</code>

mordm.read.matrix	<i>Loads a data set stored in a matrix or data.frame.</i>
-------------------	---

Description

This method assumes the first N columns are decision variables, and all other columns are objectives. N is determined by `ncol(bounds)`.

Usage

```
mordm.read.matrix(mat, nvars = NULL, nobjs = NULL, bounds = NULL,
                  maximize = NULL, names = NULL, ignore = NULL, metadata = NULL)
```

Arguments

mat	the matrix or data.frame
nvars	the number of decision variables
nobjs	the number of objectives
bounds	the lower and upper bounds of each decision variable
maximize	vector indicating the columns to be maximized
names	override the column names
ignore	columns to remove from the dataset
metadata	columns to retain in a metadata attribute

mordm.read.xls	<i>Loads a data set stored in a XLS or XLSX file.</i>
----------------	---

Description

This method is similar in use to [mordm.read.csv](#). Requires gdata and its dependencies, including a Perl interpreter on the host system.

Usage

```
mordm.read.xls(file, nvars = NULL, nobjs = NULL, bounds = NULL,
               maximize = NULL, names = NULL, ignore = NULL, metadata = NULL, ...)
```

Arguments

file	the file name
nvars	the number of decision variables
nobjs	the number of objectives
bounds	the lower and upper bounds of each decision variable
maximize	vector indicating the columns to be maximized
names	override the column names
ignore	columns to remove from the dataset
metadata	columns to retain in a metadata attribute
...	optional arguments passed to read.xls

mordm.recommend	<i>Make recommendations for analyzing the data.</i>
-----------------	---

Description

Performs basic checks to ensure the data is formatted correctly. If any issues are identified, then it will attempt to provide details on correcting or dealing with the problem.

Usage

```
mordm.recommend(data)
```

Arguments

data	the data set to be analyzed
------	-----------------------------

mordm.sample.uncertainties

Computes robustness under deep compute.robustness.

Description

Adds Gaussian noise to the decision variables and resamples the model output. The samples are distributed across one or more different models for the problem. The result from this method should be passed to `mordm.evaluate.uncertainties` to compute the robustness metrics.

Usage

```
mordm.sample.uncertainties(data, nsamples, models, sd = 0, verbose = TRUE)
```

Arguments

<code>data</code>	the data set
<code>nsamples</code>	the number of samples to generate for each point
<code>models</code>	the problem formulations created using <code>define.problem</code>
<code>sd</code>	scalar or vector specifying the standard deviation for each decision variable
<code>verbose</code>	display additional information

Details

If multiple models are provided, it is assumed that all models have the same inputs and outputs; they would only differ in the internal calculations within the model.

mordm.select

Returns the subset of rows that are marked.

Description

Applies one or more markings to the data set and returns the subset that are contained within the markings.

Usage

```
mordm.select(data, marking, index = -1, not = FALSE, or = FALSE)
```

Arguments

<code>data</code>	the data set to be displayed (if data is a time series, then the last entry in the time series is displayed)
<code>marking</code>	list of markings
<code>index</code>	if data is a time series, controls which entries to display (see mordm.get.set for details)
<code>not</code>	DEPRECATED
<code>or</code>	DEPRECATED

```
mordm.select.indices
```

Returns the row indices in the data set that are marked.

Description

Applies one or more markings to the data set to determine which rows are contained within the markings.

Usage

```
mordm.select.indices(set, marking, not = FALSE, or = FALSE)
```

Arguments

set	the data set
marking	list of markings
not	DEPRECATED
or	DEPRECATED

```
mordm.variable.sensitivities
```

Computes the sensitivities of the decision variables.

Description

Using Plischke's delta-moment sensitivity analysis method, this function computes the sensitivities using the given data. As such, this method does not need to evaluate any new data points, it works with the provided data.

Usage

```
mordm.variable.sensitivities(data, objective, index = -1, all = FALSE, ...)
```

Arguments

data	the data set
objective	the objective index, column name, function, or marking
index	if data is a time series, controls which entries to display (see mordm.get.set for details)
all	if TRUE, include all points from all entries in the time series; otherwise, only the last entry is included
...	additional options for Plischke's method

Details

If objective is a marking, then this computes the sensitivities that cause a point to be included in the marking. This functionality is still experimental.

mordm.weight	<i>Computes a vector of weighted preferences</i>
--------------	--

Description

Computes a vector of weighted preferences

Usage

```
mordm.weight(data, weights)
```

Arguments

data	the data
weights	the vector of weights

nsample	<i>Generate normally distributed random inputs.</i>
---------	---

Description

Generate normally distributed random inputs.

Usage

```
nsample(mean, sd, nsamples, problem)
```

Arguments

mean	scalar or vector specifying the mean value for each decision variable
sd	scalar or vector specifying the standard deviation for each decision variable
nsamples	the number of samples to generate
problem	the problem definition

`robustness.constraints`*Robustness metric based on constraint violations.*

Description

Measures the percentage of the sampled points that violate constraints.

Usage

```
robustness.constraints(output, problem, weights = NULL, verbose = FALSE,  
  original.point = NULL)
```

Arguments

<code>output</code>	the evaluated points
<code>problem</code>	the problem definition
<code>weights</code>	unused
<code>verbose</code>	unused
<code>original.point</code>	unused

`robustness.default`*Default robustness metric.*

Description

The default robustness metric that combines variances and constraint violations.

Usage

```
robustness.default(output, problem, weights = NULL, verbose = FALSE,  
  original.point = NULL)
```

Arguments

<code>output</code>	the evaluated points
<code>problem</code>	the problem definition
<code>weights</code>	the weights assigned to each objective
<code>verbose</code>	display additional information
<code>original.point</code>	the original point being analyzed

robustness.distance	<i>Robustness metric based on distance.</i>
---------------------	---

Description

Measures the average distance from the original point to the sampled points. This is slightly different from variance in that variance is not effected by translational distance. I.e., two point clouds have the same variance, but one is offset more.

Usage

```
robustness.distance(output, problem, weights = NULL, verbose = FALSE,  
  original.point = NULL)
```

Arguments

output	the evaluated points
problem	the problem definition
weights	unused
verbose	unused
original.point	the original point being analyzed

robustness.gap	<i>Experimental robustness metric based on info gap.</i>
----------------	--

Description

Info gap measures the distance from the original point to the nearest constraint boundary. This experimental implementation approximates this distance by computing the distance based on the sampled points.

Usage

```
robustness.gap(output, problem, weights = NULL, verbose = FALSE,  
  original.point = NULL)
```

Arguments

output	the evaluated points
problem	the problem definition
weights	unused
verbose	unused
original.point	the original point being analyzed

robustness.variance	<i>Robustness metric based on variance.</i>
---------------------	---

Description

Measures the variance of the sampled points.

Usage

```
robustness.variance(output, problem, weights = NULL, verbose = FALSE,
  original.point = NULL)
```

Arguments

output	the evaluated points
problem	the problem definition
weights	the weights assigned to each objective
verbose	display additional information
original.point	the original point being analyzed

runVisDemo	<i>Runs the exploration tool on a 5-objective problem.</i>
------------	--

Description

Runs the exploration tool on a 5-objective problem.

Usage

```
runVisDemo()
```

sensitivity.levels	<i>Determines number of replicates for sensitivity analysis.</i>
--------------------	--

Description

Calculates the number of replicates / levels required by the sensitivity analysis method to produce approximately the given number of samples

Usage

```
sensitivity.levels(problem, nsamples, method)
```

Arguments

problem	the problem definition
nsamples	the desired number of samples
method	the sensitivity analysis method

usample	<i>Generate uniformly distributed random inputs.</i>
---------	--

Description

Generate uniformly distributed random inputs.

Usage

```
usample(nsamples, problem)
```

Arguments

nsamples	the number of samples to generate
problem	the problem definition

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