

# Package: xgrove (via r-universe)

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**Title** Explanation Groves

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**Description** Compute surrogate explanation groves for predictive machine learning models and analyze complexity vs. explanatory power of an explanation according to Szepannek, G. and von Holt, B. (2023) <[doi:10.1007/s41237-023-00205-2](https://doi.org/10.1007/s41237-023-00205-2)>.

**License** GPL (>=2)

**Encoding** UTF-8

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**Imports** gbm, dplyr, rpart, rpart.plot

**Suggests** pdp, randomForest

**RoxygenNote** 7.3.2

**Config/testthat/edition** 3

**Repository** <https://g-rho.r-universe.dev>

**RemoteUrl** <https://github.com/g-rho/xgrove>

**RemoteRef** HEAD

**RemoteSha** 1edf6761e7e73f07bb7a64929e3b9b0775dff8ae

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`plot.sgtree`*Plot surrogate tree statistics*

---

**Description**

Plot statistics of surrogate trees to analyze complexity vs. explanatory power.

**Usage**

```
## S3 method for class 'sgtree'  
plot(x, abs = "rules", ord = "upsilon", ...)
```

**Arguments**

<code>x</code>	An object of class <code>sgtree</code> .
<code>abs</code>	Name of the measure to be plotted on the x-axis, either "trees", "rules", "upsilon" or "cor".
<code>ord</code>	Name of the measure to be plotted on the y-axis, either "trees", "rules", "upsilon" or "cor".
<code>...</code>	Further arguments passed to <code>plot</code> .

**Value**

No return value.

**Author(s)**

<gero.szepannek@web.de>

**Examples**

```
library(randomForest)  
library(pdp)  
data(boston)  
set.seed(42)  
rf <- randomForest(cmedv ~ ., data = boston)  
data <- boston[,-3] # remove target variable  
ntrees <- c(4,8,16,32,64,128)  
xg <- xgrove(rf, data, ntrees)  
xg  
plot(xg)
```

---

`plot.xgrove`*Plot surrogate grove statistics*

---

**Description**

Plot statistics of surrogate groves to analyze complexity vs. explanatory power.

**Usage**

```
## S3 method for class 'xgrove'  
plot(x, abs = "rules", ord = "upsilon", ...)
```

**Arguments**

<code>x</code>	An object of class <code>xgrove</code> .
<code>abs</code>	Name of the measure to be plotted on the x-axis, either "trees", "rules", "upsilon" or "cor".
<code>ord</code>	Name of the measure to be plotted on the y-axis, either "trees", "rules", "upsilon" or "cor".
<code>...</code>	Further arguments passed to <code>plot</code> .

**Value**

No return value.

**Author(s)**

<gero.szepannek@web.de>

**Examples**

```
library(randomForest)  
library(pdp)  
data(boston)  
set.seed(42)  
rf <- randomForest(cmedv ~ ., data = boston)  
data <- boston[,-3] # remove target variable  
ntrees <- c(4,8,16,32,64,128)  
xg <- xgrove(rf, data, ntrees)  
xg  
plot(xg)
```

---

sgtree *Surrogate trees*

---

### Description

Compute surrogate trees of different depth to explain predictive machine learning model and analyze complexity vs. explanatory power.

### Usage

```
sgtree(model, data, maxdeps = 1:8, cparam = 0, pfun = NULL, ...)
```

### Arguments

model	A model with corresponding predict function that returns numeric values.
data	Data that must not (!) contain the target variable.
maxdeps	Sequence of integers: Maximum depth of the trees.
cparam	Complexity parameter for growing the trees.
pfun	Optional predict function <code>function(model, data)</code> returning a real number. Default is the <code>predict()</code> method of the model.
...	Further arguments to be passed to <a href="#">rpart.control</a> or the <code>predict()</code> method of the model.

### Details

A surrogate grove is trained via gradient boosting using [rpart](#) on data with the predictions of using of the model as target variable. Note that data must not contain the original target variable!

### Value

List of the results:

explanation	Matrix containing tree sizes, rules, explainability $\Upsilon$ and the correlation between the predictions of the explanation and the true model.
rules	List of rules for each tree.
model	List of the <code>rpart</code> models.

### Author(s)

<gero.szepannek@web.de>

### References

- Szepannek, G. and Laabs, B.H. (2023): Can't see the forest for the trees – analyzing groves to explain random forests, *Behaviormetrika*, submitted.
- Szepannek, G. and Luebke, K.(2023): How much do we see? On the explainability of partial dependence plots for credit risk scoring, *Argumenta Oeconomica* 50, DOI: 10.15611/aoe.2023.1.07.

**Examples**

```
library(randomForest)
library(pdp)
data(boston)
set.seed(42)
rf <- randomForest(cmedv ~ ., data = boston)
data <- boston[,-3] # remove target variable
maxds <- 1:7
st <- sgtree(rf, data, maxds)
st
# rules for tree of depth 3
st$rules[["3"]]
# plot tree of depth 3
rpart.plot::rpart.plot(st$model[["3"]])
```

---

upsilon

*Explainability*

---

**Description**

Compute explainability given predicted data of the model and an explainer.

**Usage**

```
upsilon(porig, pexp)
```

**Arguments**

porig	An object of class xgrove.
pexp	Name of the measure to be plotted on the x-axis, either "trees", "rules", "upsilon" or "cor".

**Value**

Numeric explainability *upsilon*.

**Author(s)**

<gero.szepannek@web.de>

**References**

- Szepannek, G. and Luebke, K.(2023): How much do we see? On the explainability of partial dependence plots for credit risk scoring, *Argumenta Oeconomica* 50, DOI: 10.15611/aoe.2023.1.07.

**Examples**

```

library(randomForest)
library(pdp)
data(boston)
set.seed(42)
# Compute original model
rf <- randomForest(cmedv ~ ., data = boston)
data <- boston[,-3] # remove target variable
# Compute predictions
porig <- predict(rf, data)

# Compute surrogate grove
xg <- xgrove(rf, data)
pexp <- predict(xg$model, data, n.trees = 16)
upsilon(porig, pexp)

```

---

xgrove

*Explanation groves*


---

**Description**

Compute surrogate groves to explain predictive machine learning model and analyze complexity vs. explanatory power.

**Usage**

```

xgrove(
  model,
  data,
  ntrees = c(4, 8, 16, 32, 64, 128),
  pfun = NULL,
  remove.target = T,
  shrink = 1,
  b.frac = 1,
  seed = 42,
  ...
)

```

**Arguments**

model	A model with corresponding predict function that returns numeric values.
data	Training data.
ntrees	Sequence of integers: number of boosting trees for rule extraction.
pfun	Optional predict function function(model, data) returning a real number. Default is the predict() method of the model.

<code>remove.target</code>	Logical. If TRUE the name of the target variable is identified from <code>terms(model)</code> and automatically removed if this variable is still in data.
<code>shrink</code>	Sets the shrinkage argument for the internal call of <code>gbm</code> . As the <code>model</code> usually has a deterministic response the default is 1 different to the default of <code>gbm</code> applied train a model based on data.
<code>b.frac</code>	Sets the <code>bag.fraction</code> argument for the internal call of <code>gbm</code> . As the <code>model</code> usually has a deterministic response the default is 1 different to the default of <code>gbm</code> applied train a model based on data.
<code>seed</code>	Seed for the random number generator to ensure reproducible results (e.g. for the default <code>bag.fraction &lt; 1</code> in boosting).
<code>...</code>	Further arguments to be passed to <code>gbm</code> or the <code>predict()</code> method of the <code>model</code> .

### Details

A surrogate grove is trained via gradient boosting using `gbm` on data with the predictions of using of the `model` as target variable. Note that data must not contain the original target variable! The boosting model is trained using stumps of depth 1. The resulting interpretation is extracted from `pretty.gbm.tree`. The column `upper_bound_left` of the rules and the groves value of the output object contains the split point for numeric variables denoting the upper bound of the left branch. Correspondingly, the `levels_left` column contains the levels of factor variables assigned to the left branch. The rule weights of the branches are given in the rightmost columns. The prediction of the grove is obtained as the sum of the assigned weights over all rows. Note that the training data must not contain the target variable. It can be either removed manually or will be removed automatically from data if the argument `remove.target == TRUE`.

### Value

List of the results:

<code>explanation</code>	Matrix containing tree sizes, rules, explainability $\Upsilon$ and the correlation between the predictions of the explanation and the true model.
<code>rules</code>	Summary of the explanation grove: Rules with identical splits are aggregated. For numeric variables any splits are merged if they lead to identical partitions of the training data.
<code>groves</code>	Rules of the explanation grove.
<code>model</code>	<code>gbm</code> model.

### Author(s)

<gero.szepannek@web.de>

### References

- Szepannek, G. and von Holt, B.H. (2023): Can't see the forest for the trees – analyzing groves to explain random forests, *Behaviormetrika*, DOI: 10.1007/s41237-023-00205-2.
- Szepannek, G. and Luebke, K.(2023): How much do we see? On the explainability of partial dependence plots for credit risk scoring, *Argumenta Oeconomica* 50, DOI: 10.15611/aoe.2023.1.07.

**Examples**

```
library(randomForest)
library(pdp)
data(boston)
set.seed(42)
rf <- randomForest(cmedv ~ ., data = boston)
data <- boston[,-3] # remove target variable
ntrees <- c(4,8,16,32,64,128)
xg <- xgrove(rf, data, ntrees)
xg
plot(xg)

# Example of a classification problem using the iris data.
# A predict function has to be defined, here for the posterior probabilities of the class Virginica.
data(iris)
set.seed(42)
rf <- randomForest(Species ~ ., data = iris)
data <- iris[,-5] # remove target variable

pf <- function(model, data){
  predict(model, data, type = "prob")[,3]
}

xgrove(rf, data, pfun = pf)
```



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