

R-course: Machine Learning using R

Decision trees



Yannick Rothacher

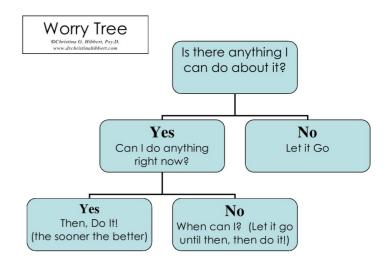
Zürich, 2021



What is a decision tree?

Generally speaking, a decision tree is a diagram, which helps us determining

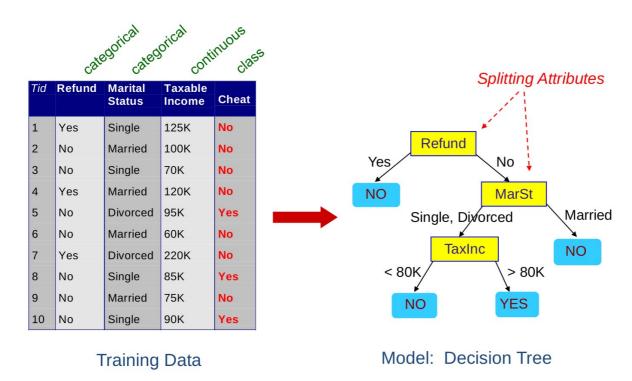
a course of action (e.g. "should I worry?")



- We can use decision trees for classification or regression
 - Such trees are the result of what is referred to as recursive partitioning
- Let's look at an example of a decision tree used for classification...



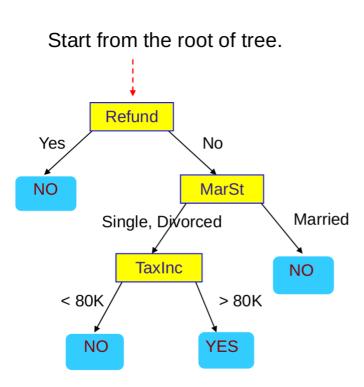
- We want to predict whether a person will cheat in his/her tax declaration based on some training data
- "Cheat" is the target variable (two levels: yes/no)
- ▶ The generated decision tree uses the three independent variables to model whether someone will cheat or not:



Source: Introduction to Data Mining, Pang-Ning Tan, Michael Steinbach, Vipin Kumar, Published by Addison Wesley.



- ▶ Use the generated tree to predict outcome of a **new observation**
- Work through the tree for prediction:

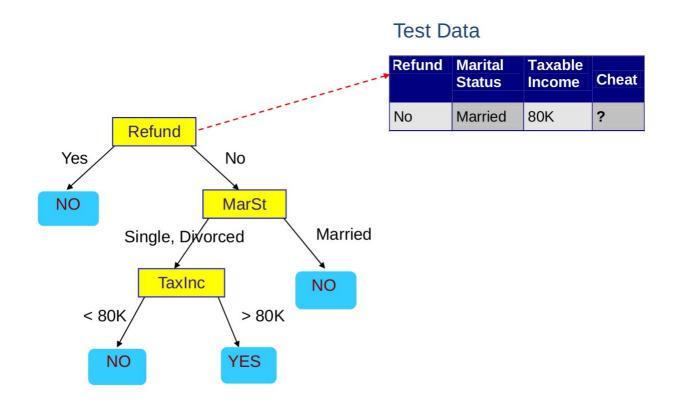


Test Data

Refund	Marital Status	Taxable Income	Cheat
No	Married	80K	?

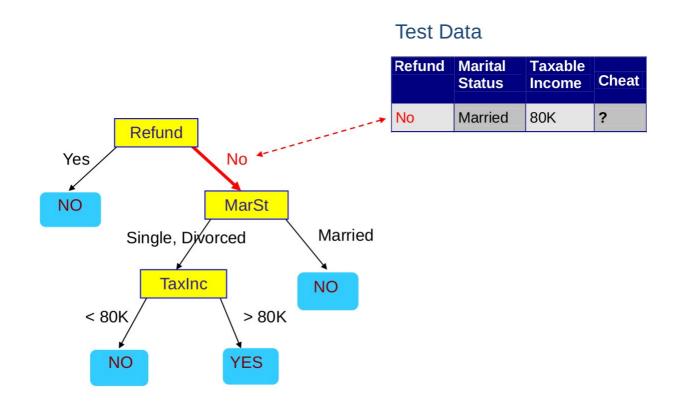


- Use the generated tree to predict outcome of a new observation
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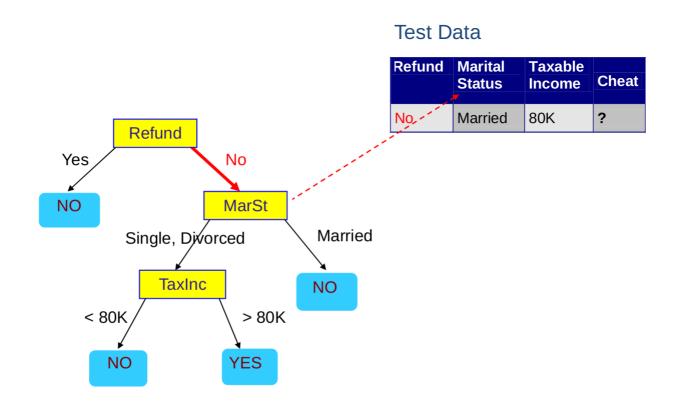


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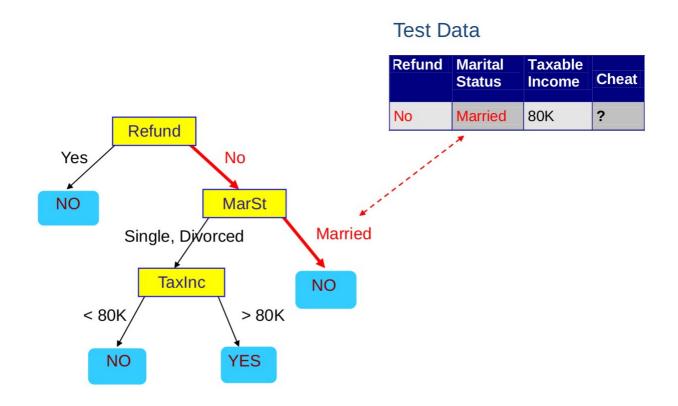


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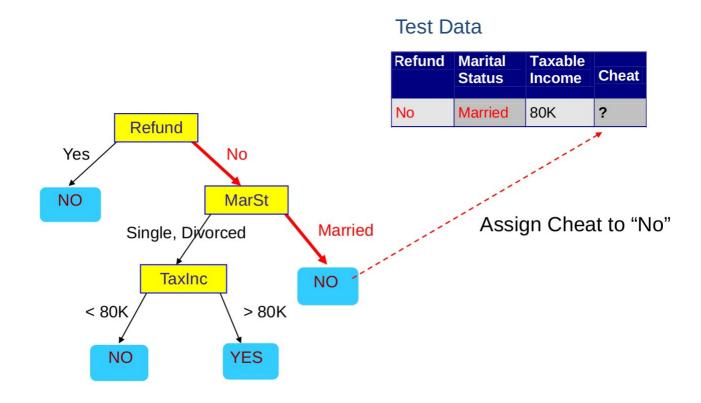


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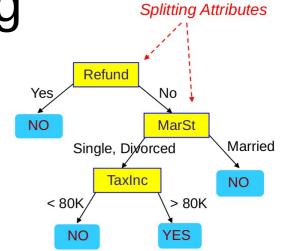
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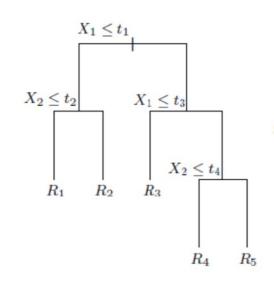


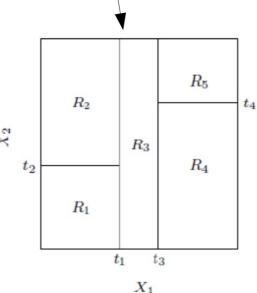
Decision tree – partitioning

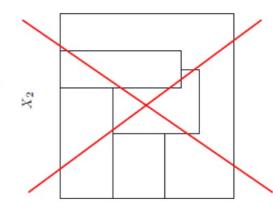
- A classic decision tree splits the data at each node based on one dependent variable
- ► For the case of only **two dependent variables** we can visualize this in a 2D-coordinate system:



Model: Decision Tree







Only straight lines!



Decision tree – partitioning

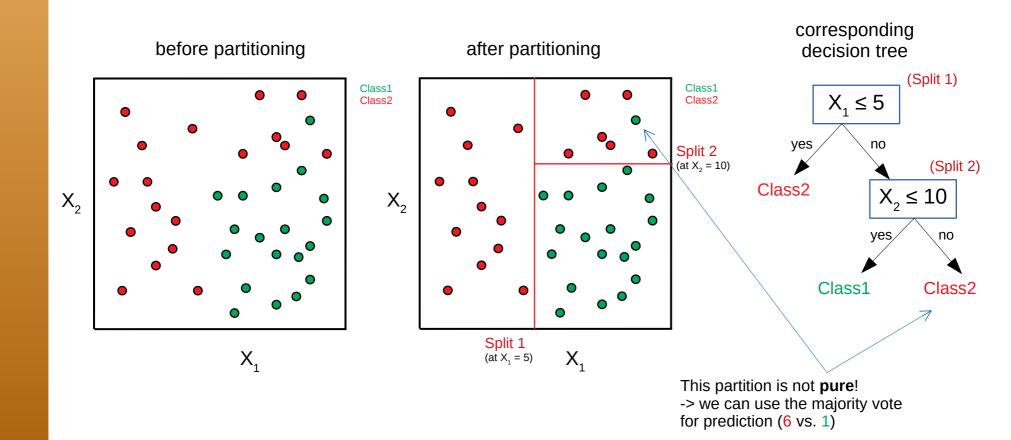
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Decision tree – partitioning

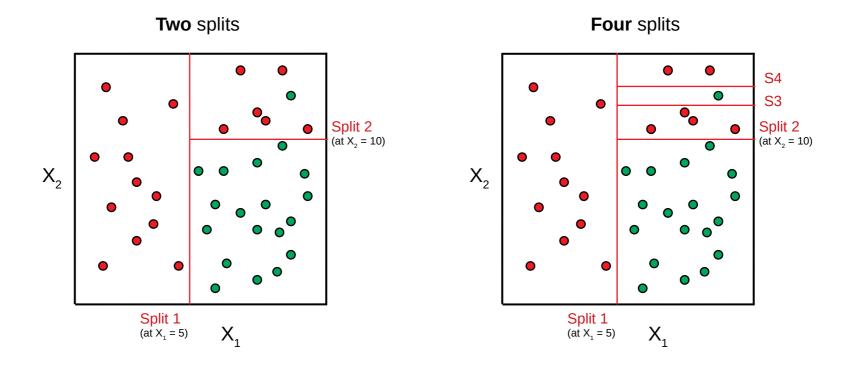
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Decision trees – over- and underfitting

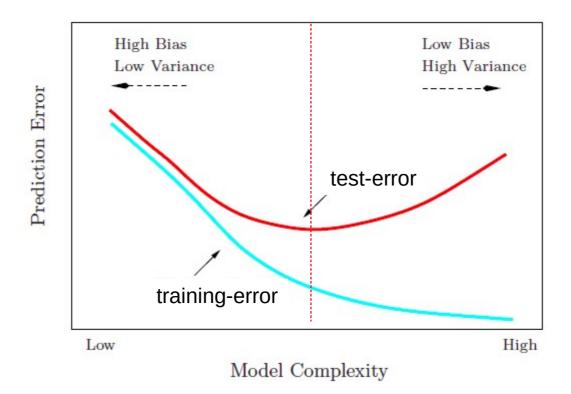
- How deep should a tree be allowed to grow?
 - We can always grow a tree until there are only pure partitions left (right figure below)
 - Same problem of over- and underfitting like with the KNN-classifier!
- Which one of the trees below is better?





Recap: Test-error

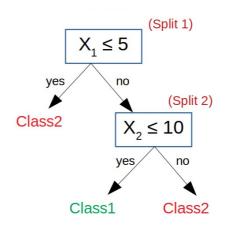
- Like in the case of the KNN-classifier we want to minimize the **test-error**
- Evaluation methods like e.g. cross-validation can of course also be applied to decision trees
 - Prevent over-fitting: Fully grown trees can be "pruned" (cut shorter) depending on whether a branch improves the fit to data evaluated by cross-validation

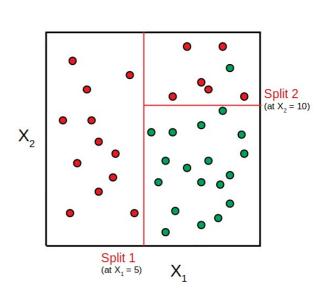




How are the partitions generated?

- Building a decision tree from scratch involves the following two questions at each node:
 - Which variable should be used for the next split?
 - Where along the chosen variable should we split?
- How can we decide what the best variable and split-location is?
- We will look at two different methods to solve these questions
 - ▶ 1) classic partitioning based on impurity measure (e.g. Gini-index)
 - 2) bias-free partitioning based on significance tests





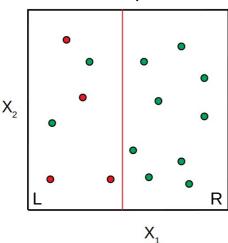


Classification trees – impurity measure

- Main idea: Try to find the split, which reduces the impurity of the data the most
 - How can we measure "impurity"?
- The Gini index is one of the most common impurity measures

$$Gini=1-\sum_{i}^{C}p_{i}^{2}$$

Entropy is another common impurity $Entropy = -\sum_{i}^{C} p_{i} \log_{2}(p_{i})$ measure



$$Gini_{root} = 1 - ((11/15)^2 + (4/15)^2) = 0.391$$

$$Gini_{L} = 1 - ((2/6)^{2} + (4/6)^{2}) = 0.444$$

 $Gini_{R} = 1 - ((0/9)^{2} + (9/9)^{2}) = 0$

Gini-reduction (weighted by number of observations in each partition):

$$Gini_decrease = Gini_{root} - (6/15)*Gini_{L} - (9/15)*Gini_{R}$$

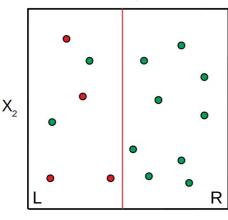


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X

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Gini-reduction (weighted by number of observations in each partition):

$$Gini_decrease = Gini_{root} - (6/15)*Gini_{L} - (9/15)*Gini_{R}$$

→ Choose the split with the largest Gini-reduction



Partition algorithm (impurity measure based)

- Start with a single region (encompassing all data)
- Iterate through the following steps:

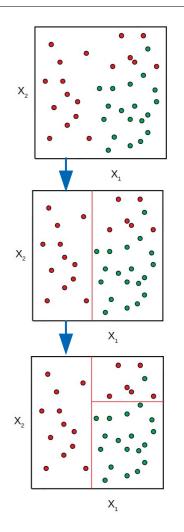
For each region \mathbf{R} For each variable \mathbf{x}_i in \mathbf{R} For each possible split \mathbf{s}_i of \mathbf{x}_i Record impurity decrease

The Choose $(\mathbf{x}_i, \mathbf{s}_i)$ which gives maximum impurity decrease

 \rightarrow Replace **R** with **R**_R and **R**_L

- Stop splitting either based on a stopping rule or when there is no more impurity reduction possible
 - Stopping rule example: Stop splitting when a branch contains less than a certain percentage of the data

$$Gini=1-\sum_{i}^{C}p_{i}^{2}$$





library(rpart)

Decision tree in R – rpart

- rpart is an R-package to fit decision trees (impurity measure based)
- ► Here an example of fitting a (classification) tree to the "Iris" data set:

```
tree.iris <- rpart(Species ~., data=iris) # default uses Gini-index</pre>
plot(tree.iris, margin = 0.1)
text(tree.iris, use.n = T)
# Predict training data (use type = 'prob' to get probabilities):
pred.tree <- predict(tree.iris, iris, type = 'class')</pre>
# Confusion matrix:
(confT <- table(pred.tree, iris$Species))</pre>
                                                                    Petal.Length < 2.45
pred.tree setosa versicolor virginica
  setosa
                   50
                                           0
  versicolor
  virginica
                                          45
                                                                                  Petal.Width< 1.75
                                                            setosa
# Training-error:
                                                             50/0/0
diag(confT) <- 0</pre>
missCount <- sum(confT)</pre>
                                                                            versicolor
                                                                                              virginica
(trainErr <- missCount/nrow(iris))</pre>
                                                                              0/49/5
                                                                                               0/1/45
[1] 0.04
```



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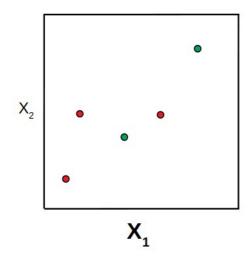
What is this?

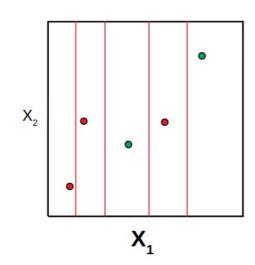
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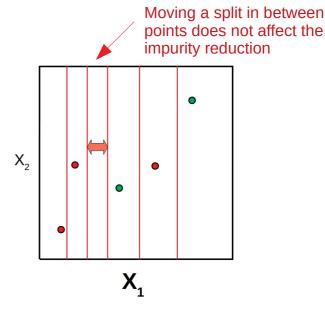


Problems with rpart

- Using an impurity measure based approach to select splits suffers from an inherent bias in certain cases
 - This bias is related to the **number of splits** that are possible for each variable
- How many splits are there per variable?
 - For **continuous** variable (numeric):







- > For the five data points above, there are only **four meaningful** splits possible (along X_1)
- What about other types of variables?



Problems with rpart

- We can also include categorical variables (factors) as predictors in a decision tree
- How many splits are possible for factors?
 - E.g. two-level factor (levels: a,b) a I b split1
 - E.g. four-level factor (levels: a,b,c,d):

a I b,c,d split1

b I a,c,d split2

c I a,b,d split3

d I a,b,c split4

a,b I c,d split5

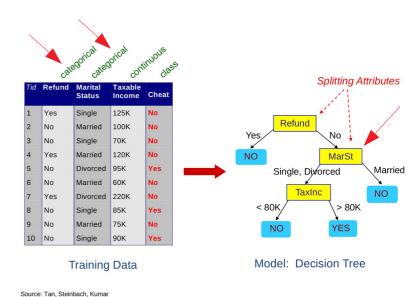
a,c I b,d split6

a,d I b,c split7





With increasing number of levels we can apply an exponentially growing number of splits





Problems with rpart

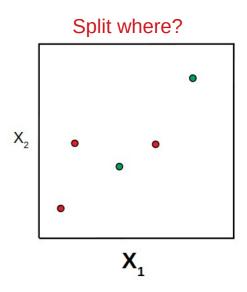
- With increasing number of levels we can apply an exponentially growing number of splits
- ➤ The difference in numbers of splits between continuous variables, factors with few levels and factors with many levels creates an issue of multiple comparisons
- Multiple comparisons problem in decision trees:
 - ► Even if there is no real relation between predictors and the target variable, under certain circumstances multi-level factors will be preferably selected for splitting compared to numerical variables or factors with few levels because they get more "chances" to reach a high impurity-decrease (even if this decrease is only achieved by pure luck)
 - More chances in the sense that more splits can be evaluated (more chances to get lucky)
 - Ideally, we do not want our decision trees to be biased in their variable selection
- We will see how this bias manifests itself by running a small simulation study (Exercise)

```
E.g. two-level factor (levels: a,b)
a I b split1

E.g. four-level factor (levels: a,b,c,d):
a I b,c,d split1
b I a,c,d split2
c I a,b,d split3
d I a,b,c split4
ab I c,d split5
ac I b,d split6
ad I b,c split7
```



- ▶ Bias-free partitioning has been proposed to solve this problem (e.g. see Hothorn et al. 2006)
 - Main idea: Use significance tests instead of impurity measures to select variable and next split
- How does it work?
 - The selection process is more clearly separated into
 - 1) Choosing the next splitting variable and
 - 2) Choosing the splitting location along the chosen variable





- Choosing the next splitting variable:
 - ➤ Test for each variable whether it has a significant association with the target variable (e.g. think of correlation test for numerical variables and of ANOVA for categorical variables). Collect the associated p-values (can be compared between tests)
 - If no variable shows a significant p-value stop splitting (integrated stopping rule)
 - Otherwise choose the variable with the lowest p-value for the next split



Choosing the next splitting variable:

Actual implementation uses special permutation tests (called "conditional inference tests") that are available for all types of variables.

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- If no variable shows a significant p-value stop splitting (integrated stopping rule)
- Otherwise choose the variable with the lowest p-value for the next split
- Choosing the next splitting location:
 - Calculate for each possible split (along the chosen variable) a test-statistic (expressing the difference between the created partitions, could also use impurity measure) and select split with the highest test-statistic
- Using this approach, there is **no bias** in the selection of different variable-types

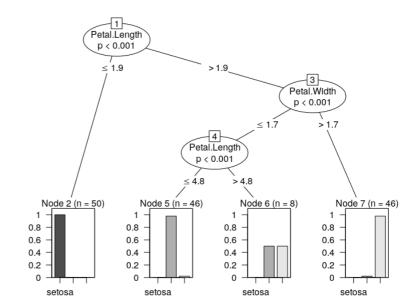


Bias-free partitioning in R

- "party" is an R-package which includes functions to fit bias-free (significance test based) decision trees (ctree()-function)
- Fitting a classification tree to the "Iris" data set

```
library(party)
ctree.iris <- ctree(Species ~., data=iris) # fit tree
plot(ctree.iris)
pred.ctree <- predict(ctree.iris, newdata=iris, type='response')
(confT <- table(pred.ctree, iris$Species)) # Confusion matrix</pre>
```

```
pred.ctree setosa versicolor virginica setosa 50 0 0 versicolor 0 49 5 virginica 0 1 45
```



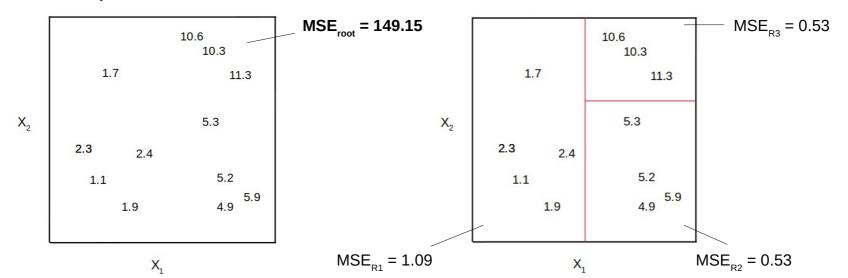


Decision trees for regression

- Decision trees can be used for classification (categorical target variable) and for regression (numerical target variable)
- ➤ The general principle behind building the tree structure does not change
 - E.g. use **mean squared error** as "impurity measure" instead of Gini-index

$$MSE = \frac{1}{n} \sum_{i}^{n} (x_i - \bar{x})^2$$

- We try to find the splits which best reduce the mean squared error
- With ctree we can fit a bias-free regression tree (again based on significance tests)



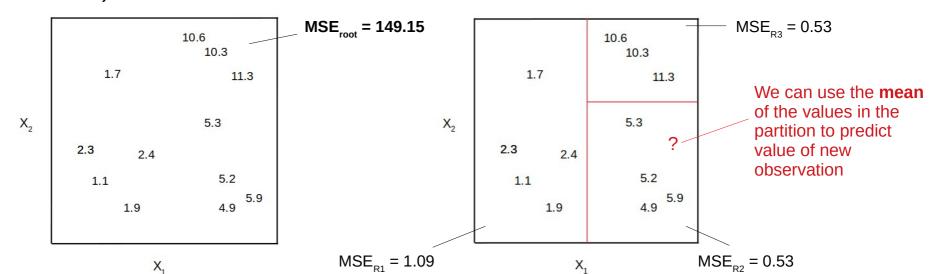


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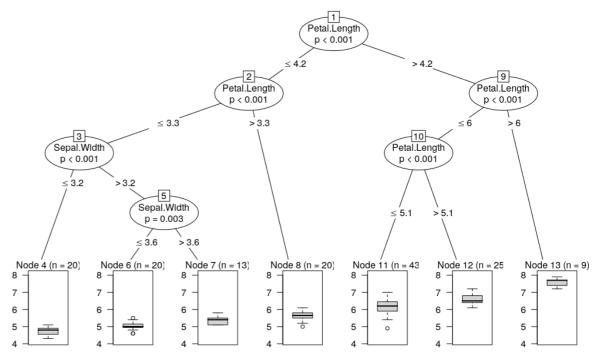




Regression tree in R

Example of fitting a regression tree to the "Iris" data using ctree()

```
library(party)
ctree.iris <- ctree(Sepal.Length ~., data=iris) # fit regression tree
plot(ctree.iris)
# Get the predicted values (for training data):
pred.ctree <- predict(ctree.iris, iris, type='response')</pre>
```





Decision trees – Summary

- Advantages of decision trees:
 - Good interpretability
 - No assumptions regarding distributions
 - Robust to outliers
 - Can capture non-linear structures and complex interactions
 - Low bias in prediction with sufficient depth
- Disadvantages:
 - Single trees are instable (small changes in data can give different-looking tree, especially if underlying pattern is complex)
 - In case of not pruned rpart-trees: Tend to overfit
 - Needs a lot of data to capture linear structures

Caution: In case of different types of predictor variables use algorithm with bias-free variable selection (ctree in R)



Further reading

➤ Strobl, C., Malley, J., & Tutz, G. (2009). An introduction to recursive partitioning: Rationale, application, and characteristics of classification and regression trees, bagging, and Random Forests. *Psychological Methods*, 14 (4), 323–348. doi: 10.1037/a0016973