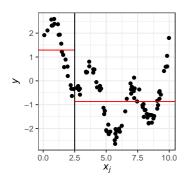
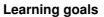
# **Introduction to Machine Learning**

# **CART Splitting Criteria for Regression**

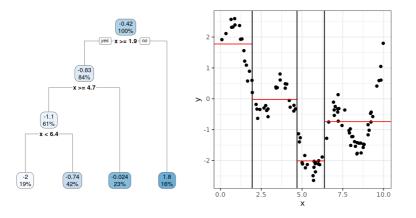




- Understand how to define split criteria via ERM
- Understand how to find splits in regression with L<sub>2</sub> loss



# **SPLITTING CRITERIA**





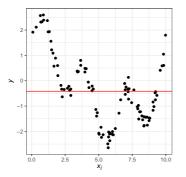
How to find good splitting rules?  $\implies$  Empirical Risk Minimization

#### **OPTIMAL CONSTANTS IN LEAVES**

Idea: A split is good if each child's point predictor reflects its data well.

For each child  $\mathcal{N}$ , predict with optimal constant, e.g., the mean  $c_{\mathcal{N}} = \frac{1}{|\mathcal{N}|} \sum_{(\mathbf{x}, y) \in \mathcal{N}} y$  for the  $L_2$  loss, i.e.,  $\mathcal{R}(\mathcal{N}) = \sum_{(\mathbf{x}, y) \in \mathcal{N}} (y - c_{\mathcal{N}})^2$ .

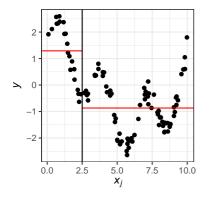
Root node:

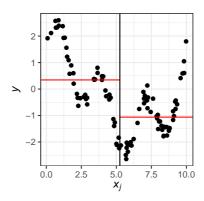




# **OPTIMAL CONSTANTS IN LEAVES**

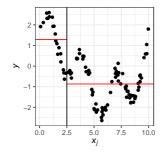
Which of these two splits is better?

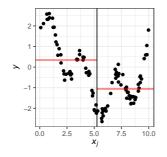






### **RISK OF A SPLIT**







$$\mathcal{R}(\mathcal{N}_1) = 23.4, \, \mathcal{R}(\mathcal{N}_2) = 72.4$$
  $\mathcal{R}(\mathcal{N}_1) = 78.1, \, \mathcal{R}(\mathcal{N}_2) = 46.1$ 

$$\mathcal{R}(\mathcal{N}_1)=$$
 78.1,  $\mathcal{R}(\mathcal{N}_2)=$  46.1

The total risk is the sum of the individual losses:

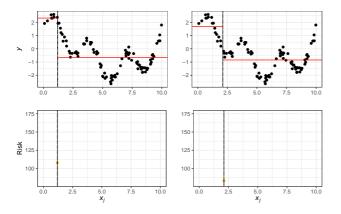
$$23.4 + 72.4 = 95.8$$

$$78.0 + 46.1 = 124.1$$

Based on the SSE, we prefer the first split.

## **SEARCHING THE BEST SPLIT**

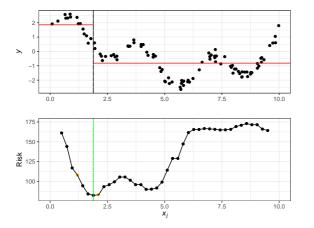
Let's find the best split for this feature by tabulating results.

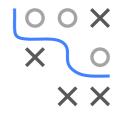




### **SEARCHING THE BEST SPLIT**

Let's iterate – quantile-wise or over all points.





We have reduced the problem to a simple loop.

#### **FORMALIZATION**

- $\mathcal{N} \subseteq \mathcal{D}$  is the data contained in this node
- Let  $c_{\mathcal{N}}$  be the predicted constant for  ${\mathcal{N}}$
- The risk  $\mathcal{R}(\mathcal{N})$  for a node is:

$$\mathcal{R}(\mathcal{N}) = \sum_{(\mathbf{x}, y) \in \mathcal{N}} \mathcal{L}(y, c_{\mathcal{N}})$$

- ullet The optimal constant is  $c_{\mathcal{N}} = rg \min_{c} \sum_{(\mathbf{x}, \mathbf{y}) \in \mathcal{N}} L(\mathbf{y}, c)$
- We often know what that is from theoretical considerations or we can perform a simple univariate optimization



#### **FORMALIZATION / 2**

• A split w.r.t. **feature**  $x_i$  **at split point** t divides a parent node  $\mathcal{N}$  into

$$\mathcal{N}_1 = \{(\mathbf{x}, y) \in \mathcal{N} : x_j < t\} \text{ and } \mathcal{N}_2 = \{(\mathbf{x}, y) \in \mathcal{N} : x_j \ge t\}.$$

To evaluate its quality, we compute the risk of our new, finer model

$$\mathcal{R}(\mathcal{N}, j, t) = \mathcal{R}(\mathcal{N}_1) + \mathcal{R}(\mathcal{N}_2)$$

$$= \left(\sum_{(\mathbf{x}, y) \in \mathcal{N}_1} L(y, c_{\mathcal{N}_1}) + \sum_{(\mathbf{x}, y) \in \mathcal{N}_2} L(y, c_{\mathcal{N}_2})\right)$$

• Finding the best way to split  $\mathcal{N}$  into  $\mathcal{N}_1, \mathcal{N}_2$  means solving

$$\underset{j,t}{\operatorname{arg\,min}}\,\mathcal{R}(\mathcal{N},j,t)$$

#### **FORMALIZATION / 3**

- $\mathcal{R}(\mathcal{N}, j, t) = \mathcal{R}(\mathcal{N}_1) + \mathcal{R}(\mathcal{N}_2)$ , makes sense if  $\mathcal{R}$  is a simple sum
- If we use averages, we have to reweight the terms to obtain a global average w.r.t.  ${\cal N}$  as the children have different sizes

$$\bar{\mathcal{R}}(\mathcal{N},j,t) = \frac{|\mathcal{N}_1|}{|\mathcal{N}|} \bar{\mathcal{R}}(\mathcal{N}_1) + \frac{|\mathcal{N}_2|}{|\mathcal{N}|} \bar{\mathcal{R}}(\mathcal{N}_2)$$



 We mention this for clarity, as quite a few texts contain only the (more complicated) weighted formula without clear explanation