Supervised Machine Learning – CART

Big Data Analysis

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Introduction

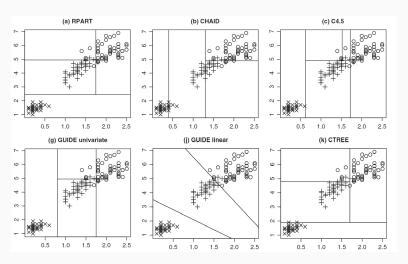
Introduction

Decision Trees

- Data-driven approach for relating X and Y
- Popular and (somewhat) easy to interpret
- · Important building block (base learner) for ensemble methods
- Many different tree building algorithms exist (Zhang & Singer 2010, Loh 2014)
 - Focus on interaction detection, prediction, parameter instability...

Introduction

Figure 1: Decision Tree Algorithms



Classification and Regression Trees (CART)

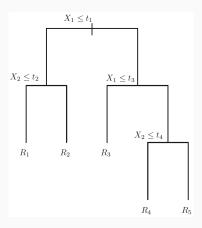
Classification and Regression Trees (CART)

Classification and Regression Trees (CART)

- Approach for partitioning the predictor space into smaller subregions via "recursive binary splitting"
- · Results in a "top-down" tree structure with...
 - · Internal nodes within the tree
 - · Terminal nodes as endpoints
- · Can be applied to regression and classification problems

Classification and Regression Trees (CART)

Figure 2: A small tree



James et al. (2013)

Growing a regression tree

Define pairs of regions for all $X_1, X_2, ..., X_p$ predictors and cutpoints c

$$\tau_L(j,c) = \{X|X_j < c\} \text{ and } \tau_R(j,c) = \{X|X_j \ge c\}$$

Find split s which maximizes the reduction in RSS

$$\Delta RSS(s, \tau) = RSS(\tau) - RSS(\tau_L) - RSS(\tau_R)$$

$$RSS(\tau) = \sum_{i \in \tau} (y_i - \hat{y})^2$$

with \hat{y} being the mean of y in node au

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$$\tau_L(j,c) = \{X|X_j < c\} \text{ and } \tau_R(j,c) = \{X|X_j \ge c\}$$

Find split s which maximizes the reduction in node impurity

$$\Delta I(s,\tau) = I(\tau) - p(\tau_L)I(\tau_L) - p(\tau_R)I(\tau_R)$$

Impurity measures

$$I_{Gini}(\tau) = \sum_{k=1}^{K} \hat{p}_k (1 - \hat{p}_k)$$

$$I_{entropy}(au) = -\sum_{k=1}^K \hat{p}_k \log \hat{p}_k$$

with \hat{p}_k being the proportion of observations from class k in node au

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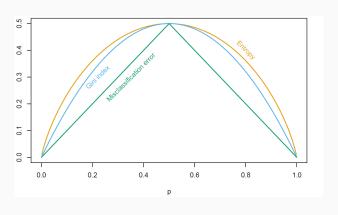
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Figure 3: Misclassification error, Gini index & entropy (scaled)



Hastie et al. (2009)

Algorithm 1: Tree growing process

```
Define stopping criteria;
Assign training data to root node;
if stopping criterion is reached then
end splitting;
else
find the optimal split point;
split node into two subnodes at this split point;
for each node of the current tree do
continue tree growing process;
end
end
```

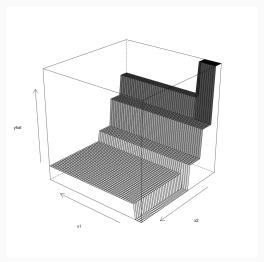
A given tree

$$\mathcal{T} = \sum_{m=1}^{M} \gamma_m \cdot 1_{(i \in \tau_m)}$$

consists of a set of m = 1, 2, ..., M nodes which can be used for prediction by...

- · Regression
 - \cdot ...using the mean of y for training observations in $au_{\it m}$
- Classification
 - \cdot ...going with the majority class in $au_{\it m}$
- ightarrow Prediction surface: Block-wise relationship between features and outcome

Figure 4: Tree prediction surface (example)



Missings

- Create a new category for missing values
- Use surrogate splits
 - 1. Choose best (primary) predictor based on complete cases
 - 2. Search for surrogate variables which mimic the chosen split
 - 3. Use surrogates if values for primary predictor are missing

Costs

$$L = \begin{pmatrix} 0 & L_{fp} \\ L_{fn} & 0 \end{pmatrix}$$

- Typically $L_{fp} = L_{fn} = 1$
- Misclassifications can be weighted differently
 - Modification of loss-matrix through weights / modified Gini index

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Tree pruning

Stopping rules

- · Minimum number of cases in terminal nodes
- · Decrease in impurity exceeds some threshold
- \rightarrow However, worthless splits can be followed by good splits

Cost complexity pruning

$$R_{\alpha}(\mathcal{T}) = R(\mathcal{T}) + \alpha |\mathcal{T}|$$

- · Find the best subtree by balancing quality $\mathit{R}(\mathcal{T})$ and complexity $|\mathcal{T}|$
- \cdot $\, lpha$ controls the penalty on the number of terminal nodes
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Summary

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- Divide-and-conquer strategy that splits the data into subgroups
- · Surface from decision trees is a non-smooth step function
- No need to specify the functional form in advance (unlike regression)
- · Non-linearities and interactions are handled automatically
- Limitations: Instability(!), competition among correlated predictors, biased variable selection

Software Resources

Software Resources

Resources for R

- · Basic CART implementation: tree
- · Standard package to build CARTs: rpart
 - Includes build-in Cross-Validation and prune function
- Unified infrastructure for tree representation: partykit

References

References

- Hastie, T., Tibshirani, R., Friedman, J. (2009). The Elements of Statistical Learning: Data Mining, Inference, and Prediction. New York, NY: Springer.
- James, G., Witten, D., Hastie, T., Tibshirani, R. (2013). *An Introduction to Statistical Learning*. New York, NY: Springer.
- Loh, W.-Y. (2014). Fifty Years of Classification and Regression Trees. *International Statistical Review* 82(3), 329–348.
- Zhang, H., Singer, B. (2010). *Recursive Partitioning and Applications*. New York, NY: Springer.