

STAT 770 Dec. 7 Lecture 27

Decision Tree Methods vs Logistic Regression

Reading and Topics for this lecture: `rpart` and `randomForest` software descriptions (posted to special Decision Tree module in ELMS), plus the `R Scripts` for this class: `IntXPred.RLog` and `RandomForests.RLog`.

- (1) General discussion of Logistic Regression as Classification
- (2) Motivation for Decision Trees as search for Interactions
- (3) High-level discussion of CART and `rpart`
- (4) Script case-studies, of `rpart` and `randomForest`

Logistic Regression as a Classification Method

- With binary responses Y_i and predictors \underline{X}_i : logistic regression provides predictors $I_{[\underline{X}_i^{tr} \hat{\beta} \geq c]}$ for $Y_i = 1$.
- Effective classification rules may be complicated, depending on (higher-order) interactions or nonlinear recodes of the coordinates of \underline{X}_i .
- Stepwise model-selection strategies offer screening approach for model terms: but how could one find important higher-order interactions? Search among many predictors may fail for combinatorial reasons.
- Decision trees look directly for successive branchings, may arrive at combinations of variables without searching among all such combinations.

CART and Recursive Partitioning

Sources:

Classification and Regression Trees, L. Breiman et al. (1980)

H. Zhang & B. Singer (2010) *Recursive Partitioning and Applications*, Springer.

Similar R packages `rpart` and `tree`, “long introduction” to `rpart` by Therneau and Atkinson.

All these tree-based methods consist of two parts: successive (greedy) search for ‘splitting’ of nodes to decrease an index as much as possible. Tree is “grown” until a stopping criterion on # levels or size of nodes is reached, then “pruned”.

Recursive (Binary) Partitioning, cont'd

stage K of tree: set U of units partitioned into nodes $\{A_j\}_{j=1}^K$

Split node A_j into $A_{j,1}, A_{j,2}$, where $A_{j,1} = \{i \in A_j : X_{i,k_j} \leq a_{k_j}\}$
or $\{i \in A_j : X_{i,k_j} \in C_{k_j}\}$ (for *factor-column* X_{i,k_j})

Splitting index – choose node, split to maximize change

$$\Delta I = p(A_j)I(r(A_j)) - p(A_{j,1})I(r(A_{j,1})) - p(A_{j,2})I(r(A_{j,2}))$$

where $r(B) = |B \cap [Y = 1]|/|B|$, and $p(B) = |B|/|U|$

$I(p) =$ concave fcn, e.g. $p(1-p)$ or $-p \log p - (1-p) \log(1-p)$

Pruning – minimize misclassification rate penalized by $\alpha \cdot \# \text{nodes}$

Random Forest Idea

- Grow *many* trees, on randomly sampled subsets of data, with splits at each stage based on a small random sample of \underline{X} coordinates
- aggregate over many trees by averaging predictions from mini-tree prediction rules.
- look in Scripts for examples