Basic Machine Learning Methods: Part 2

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DAAML

Outline

• Decision Trees

- Ensemble Learning
- Random Forests

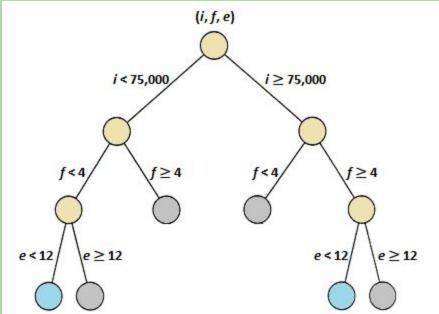
• Gradient Boosting

• flowcharts of inequalities

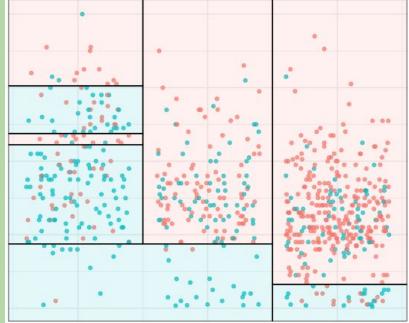
• Each test involves a <u>single</u> feature.

 Can be used for classification <u>and</u> regression!

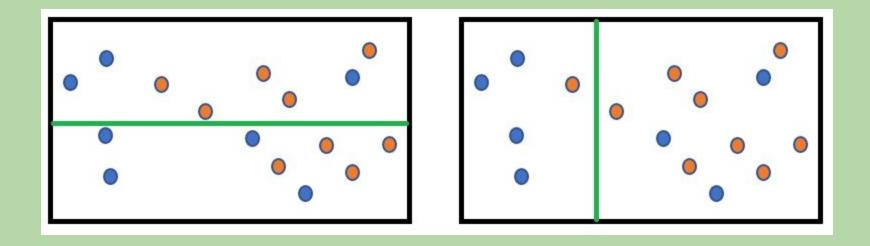
Purchases depend on income, family size and education:



 Can also equivalently interpret decision trees <u>geometrically</u>:

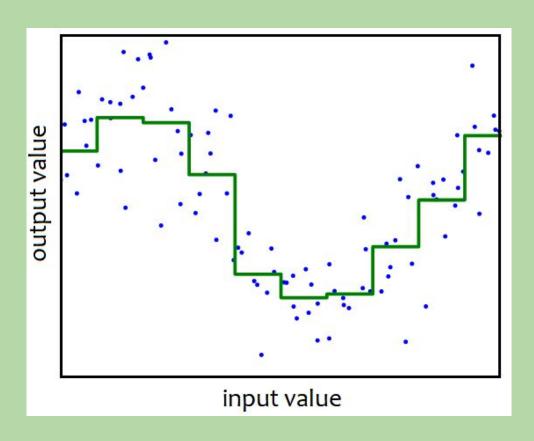


At each stage in building a decision tree,
 pick the <u>best</u> feature for the inequality.



 Decision trees can also be used for regression!

 Results are the <u>averages</u> of output values in regions.



- Decision trees do not require normalization and are easy to understand.
- Decision trees are very susceptible to overfitting.
- One way to try to avoid overfitting is to limit the number of levels.

• Uses <u>multiple</u> models and methods!

• No reason in general must use one model!

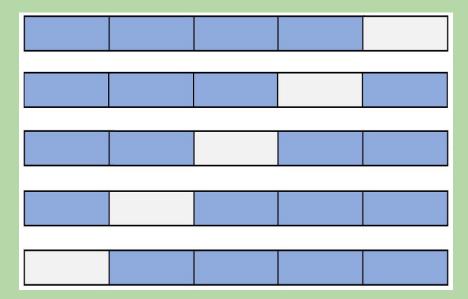
• Consider advisors helping solve a problem.

Can pick one result or combine results.

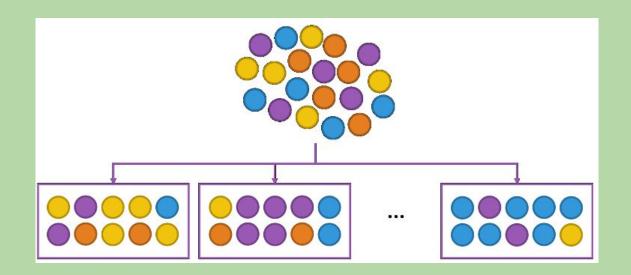
 Also no reason in general must use datasets in one way!

 Can use several different derivations for testing and training.

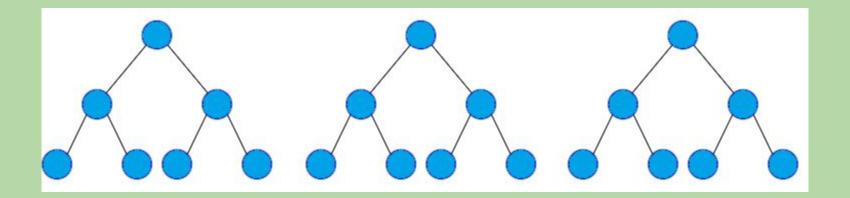
 K fold cross validation has <u>independent</u> testing subsets but limited in <u>quantity</u>.



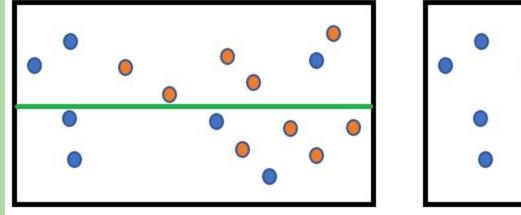
Bootstrapping creates nearly <u>unlimited</u>
 training subsets but has <u>redundancy</u>.

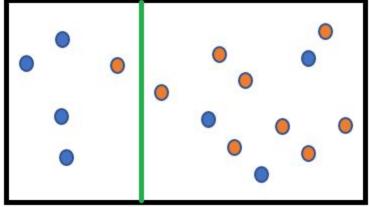


 Use ensembles of decision trees with voting and averaging to avoid overfitting!



• If bootstrapping, and feature selections limited to random subsets, then ensembles are referred to as <u>random forests</u>.





• Does not require normalization.

Only a few parameters need to be tuned.

 Training requires more resources than decision trees but is <u>parallelizable</u>!

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Do we Need Hundreds of Classifiers to Solve Real World Classification Problems?

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We evaluate 179 classifiers arising from 17 families (discriminant analysis, Bayesian, neural networks, support vector machines, decision trees, rule-based classifiers, boosting, bagging, stacking, random forests and other ensembles, generalized linear models, nearest-neighbors, partial least squares and principal component regression, logistic and multinomial regression, multiple adaptive regression splines and other methods), implemented in Weka, R (with and without the caret package), C and Matlab, including all the relevant classifiers available today. We use 121 data sets, which represent the whole UCI data base (excluding the large-scale problems) and other own real problems, in order to achieve significant conclusions about the classifier behavior, not dependent on the data set collection. The classifiers most likely to be the bests are the random forest (RF) versions, the best of which (implemented in R and accessed via caret) achieves 94.1% of the maximum accuracy overcoming 90% in the 84.3% of the data sets. However, the dif-

 Consider instead building decision trees sequentially while decreasing errors.

 Additional decision trees are instead trained to <u>offset</u> previous errors.

training inputs	training outputs	decision tree model	testing residuals
	0	T ₀	R ₀
ı	-R ₀	T ₁ '	R ₁
I	-R ₁	T ₂ '	R ₂
I	-R ₂	T ₃ '	R_3
I			
I	-R _{N-1}	T _N '	R _N

$$T_0$$
 = average of 0

$$T_k' = IT_k$$

• Often more accurate than random forests!

XGBoost has won several competitions.



 Much harder to tune, and often more computationally expensive, due to a learning rate parameter.

Not parallelizable.

• Much more prone to overfitting.

 Overfitting often makes unsuitable for very high dimensional datasets.