Machine Learning (CSE 446): Decision Trees

Sham M Kakade © 2019

University of Washington cse446-staff@cs.washington.edu

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Announcements

- Piazza: If you did not get an email, send us a message.
 Announcements will be made there, so make sure you can see them posted.
- ► Gradescope:
 - USE YOUR UW ID NUMBER WHEN YOU SIGN UP.

we need to match your assignments to you.

- The course code is: 9JKZ4G.
- ▶ HW0 posted/Turn in Certification file that you read the website.
- TA office hours posted. (Please check website before you go, just in case of changes.)
- Midterm date: Mon, Feb 11.
- Qz section this week.
- ► Today: an example, Decision Trees

The "i.i.d." Supervised Learning Setup

- Let ℓ be a loss function; $\ell(y, \hat{y})$ is our loss by predicting \hat{y} when y is the correct output.
- Let D(x, y) define the (unknown) underlying probability of input/output pair (x, y), in "nature."
- ▶ The training data $D = \langle (x_1, y_1), (x_2, y_2), \dots, (x_N, y_N) \rangle$ are assumed to be identical, independently, distributed (i.i.d.) samples from \mathcal{D} .
- ► We care about our expected error (i.e. the expected loss, the "true" expected loss, ...) with regards to the underlying distribution D.
- ► Goal: find a hypothesis which as has "low" expected error, using the training set.

The loss

Fix a classifier
$$f$$
 on (x, y) , the " $0/1$ loss" is:

 $\mathbf{1}\{y \neq f(x)\}$

Classifier f's true expected loss:

$$\epsilon(f) = \sum_{(x,y)} \mathcal{D}(x,y) \mathbf{1}\{y \neq f(x)\} = \mathbb{E}[\mathbf{1}\{y \neq f(x)\}]$$

Classifier f's average loss on training data:

$$\widehat{\epsilon}(f) = \frac{1}{N} \sum_{i=1}^{N} \mathbf{1}\{y_i \neq f(x_i)\}$$

A Toy Data Set

Data derived from https://archive.ics.uci.edu/ml/datasets/Auto+MPG

mpg; cylinders; displacement; horsepower; weight; acceleration; year; origin 18.0 8 307.0 130.0 3504. 12.0 70 1 15.0 8 350.0 165.0 3693. 11.5 70 1 18.0 318.0 150.0 3436. 11.0 70 8 1 16.0 304.0 150.0 70 1 8 3433. 12.0 17.0 8 302.0 140.0 3449. 10.5 70 1 15.0 8 429.0 198.0 4341. 10.0 70 1 14.0 8 454.0 220.0 4354. 9.0 70 1 215.0 70 14.0 8 440.0 4312. 8.5 1 1 14.0 8 455.0 225.0 4425. 10.0 70 15.0 8 390.0 190.0 3850. 8.5 70 1 15.0 383.0 170.0 3563. 10.0 70 1 8 14.0 340.0 160.0 3609. 70 1 8 8.0 15.0 8 400.0 150.0 3761. 9.5 70 1 14.0 8 455.0 225.0 3086. 10.0 70 1 24.0 70 З 4 113.0 95.00 2372. 15.0 2833. 22.0 6 198.0 95.00 15.5 70 1 18.0 6 199.0 97.00 2774. 15.5 70 1 21.0 6 200.0 85.00 2587. 16.0 70 1 27.0 4 97.00 88.00 2130. 14.5 70 3 2 26.0 4 97.00 46.00 1835. 20.5 70 2 25.0 4 110.0 87.00 2672. 17.5 70 24.0 4 107.0 90.00 2430. 14.5 70 2

Input: a row in this table; "features" are columns.

Goal: predict whether mpg is < 23("bad" = 0) or above ("good" = 1) given other attributes (other columns).

201 "good" and 197 "bad"; guessing the most frequent class (good) will get 50.5% accuracy.

Let's build a classifier!

- Let's just try to build a classifier "by hand". (This is our first learning algorithm.)
- For now, let's ignore the true loss/trying to "generalize"
- Let's start with just looking at a simple classifier. What is a simple classification rule?
- Conceptual point: Let $\Phi(x) = (\phi_1(x), \phi_2(x), \dots, \phi_d(x))$ be a function that maps from inputs x to a vector of values. Each component function $\phi(x) = \phi_i(x)$ could be:
 - If ϕ maps to $\{0,1\}$, we call it a "binary feature (function)."
 - If ϕ maps to \mathbb{R} , we call it a "real-valued feature (function)."
 - ϕ could map to categorical values, integers, ...

Sometimes we write $\Phi(x)$ to refer to a vector of features of x.

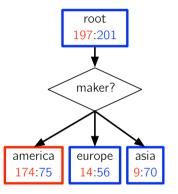
Contingency Table

		values of feature ϕ			
values of y		v_1	v_2	•••	v_K
values of g	0				
	1				

	maker			
y	america	europe	asia	
0	174	14	9	
1	75	56	70	
	\downarrow	\downarrow	\downarrow	
	0	1	1	

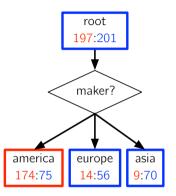
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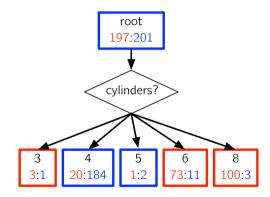
	maker			
y	america	europe	asia	
0	174	14	9	
1	75	56	70	
	\downarrow	\downarrow	\downarrow	
	0	1	1	



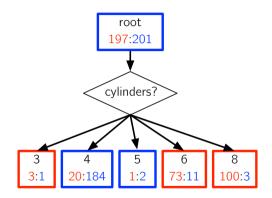
	maker				
y	america	europe	asia		
0	174	14	9		
1	75	56	70		
	\downarrow	\downarrow	\downarrow		
	0	1	1		

Errors: 75 + 14 + 9 = 98 (about 25%)





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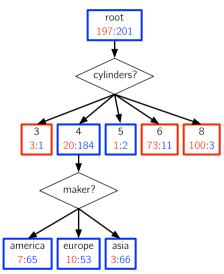
Errors: 1 + 20 + 1 + 11 + 3 = 36 (about 9%)

Key Idea: Recursion

A single feature **partitions** the data.

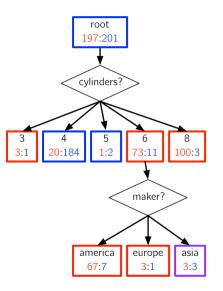
For each partition, we could choose another feature and partition further.

Applying this recursively, we can construct a decision tree.



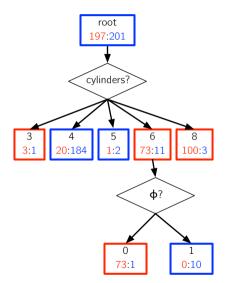
Error reduction compared to the cylinders stump?

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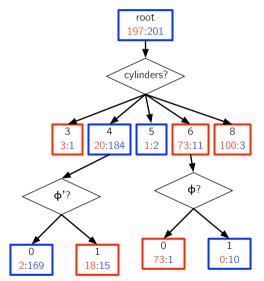


Error reduction compared to the cylinders stump?

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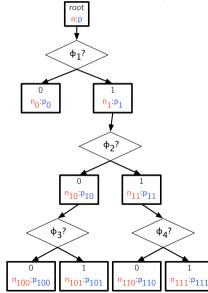
Error reduction compared to the cylinders stump?



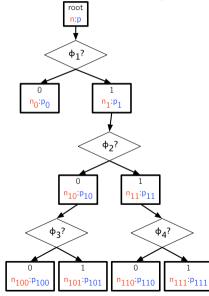
Error reduction compared to the cylinders stump?

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Decision Tree: Making a Prediction



Decision Tree: Making a Prediction



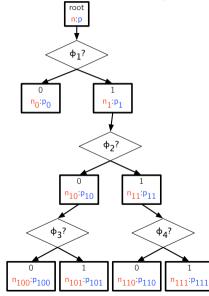
Data: decision tree t, input example x**Result:** predicted class if t has the form LEAF(y) then return y;

else

 $\begin{array}{c|c} \# t.\phi \text{ is the feature associated with } t; \\ \# t.\text{child}(v) \text{ is the subtree for value } v; \\ \text{return } \text{DTREETEST}(t.\text{child}(t.\phi(x)), x)); \\ \text{end} \end{array}$

Algorithm 1: DTREETEST

Decision Tree: Making a Prediction



Equivalent boolean formulas:

 $\begin{aligned} (\phi_1 = 0) \Rightarrow \llbracket \mathbf{n}_0 < \mathbf{p}_0 \rrbracket \\ (\phi_1 = 1) \land (\phi_2 = 0) \land (\phi_3 = 0) \Rightarrow \llbracket \mathbf{n}_{100} < \mathbf{p}_{100} \rrbracket \\ (\phi_1 = 1) \land (\phi_2 = 0) \land (\phi_3 = 1) \Rightarrow \llbracket \mathbf{n}_{101} < \mathbf{p}_{101} \rrbracket \\ (\phi_1 = 1) \land (\phi_2 = 1) \land (\phi_4 = 0) \Rightarrow \llbracket \mathbf{n}_{110} < \mathbf{p}_{110} \rrbracket \\ (\phi_1 = 1) \land (\phi_2 = 1) \land (\phi_4 = 1) \Rightarrow \llbracket \mathbf{n}_{111} < \mathbf{p}_{111} \rrbracket \end{aligned}$

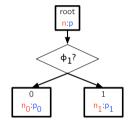
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Tangent: How Many Formulas?

- ► Assume we have *D* binary features.
- Each feature could be set to 0, or set to 1, or excluded (wildcard/don't care).
 3^D formulas.

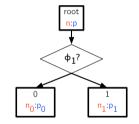
root n:p

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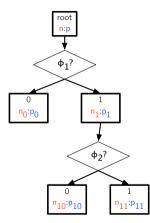


We chose feature ϕ_1 . Note that $n = n_0 + n_1$ and $p = p_0 + p_1$.

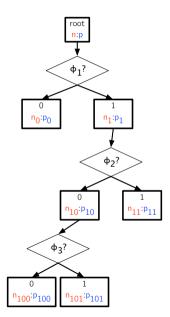
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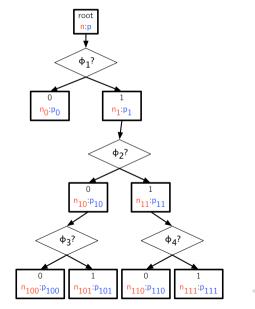
We chose not to split the left partition. Why not?



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Greedily Building a Decision Tree (Binary Features)

Data: data D, feature set Φ

Result: decision tree

if all examples in D have the same label y, or Φ is empty and y is the best guess

then

```
return LEAF(y);
```

else

```
for each feature \phi in \Phi do

partition D into D_0 and D_1 based on \phi-values;

let mistakes(\phi) = (non-majority answers in D_0) + (non-majority answers in

D_1);

end

let \phi^* be the feature with the smallest number of mistakes;

return NODE(\phi^*, {0 \rightarrow DTREETRAIN(D_0, \Phi \setminus \{\phi^*\}), 1 \rightarrow

DTREETRAIN(D_1, \Phi \setminus \{\phi^*\})};
```

end

Algorithm 2: DTREETRAIN

What could go wrong?



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