CS5010 Artificial Intelligence Principles: Lecture 4

LEARNING FROM OBSERVATIONS

Chapter 18, Sections 1–3

Outline

- ♦ Learning agents
- \Diamond Inductive learning
- ♦ Decision tree learning
- ♦ Measuring learning performance

Learning

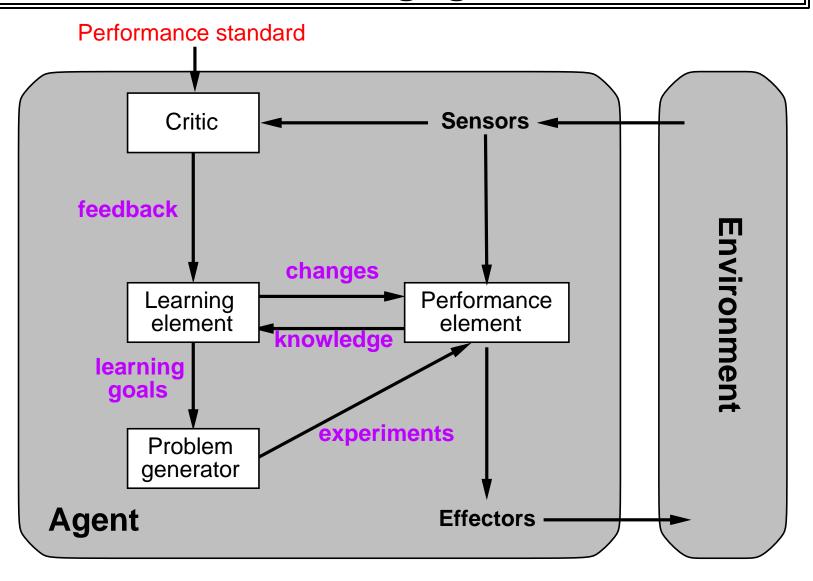
Learning is essential for unknown environments, i.e., when designer lacks omniscience

Learning is useful as a system construction method, i.e., expose the agent to reality rather than trying to write it down

Learning modifies the agent's decision mechanisms to improve performance

See Lecture 3 to review the derivation of this scheme

Learning agents



Learning element

Design of learning element is dictated by

- what type of performance element is used
- which functional component is to be learned
- ♦ how that functional compoent is represented
- ♦ what kind of feedback is available

Example scenarios:

Performance element	Component	Representation	Feedback	
Alpha-beta search	Eval. fn.	Weighted linear function	Win/loss	
Logical agent	Transition model	Successor-state axioms	Outcome	
Utility-based agent	Transition model	Dynamic Bayes net	Outcome	
Simple reflex agent	Percept-action fn	Neural net	Correct action	

Supervised learning: correct answers for each instance

Reinforcement learning: occasional rewards

RL: exploration and exploitation of knowledge learned by repeated trials of maximising the reward.

Chapter 18, Sections 1-3

Inductive learning (a.k.a. Science)

Simplest form: learn a function from examples (tabula rasa)

f is the target function

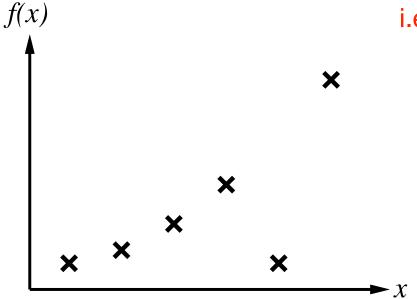
Problem: find a(n) hypothesis h such that $h \approx f$ given a training set of examples

(This is a highly simplified model of real learning:

- Ignores prior knowledge
- Assumes a deterministic, observable "environment"
- Assumes examples are given
- Assumes that the agent wants to learn f—why?)

Construct/adjust h to agree with f on training set (h) is consistent if it agrees with f on all examples)

E.g., curve fitting: (or linear regression)

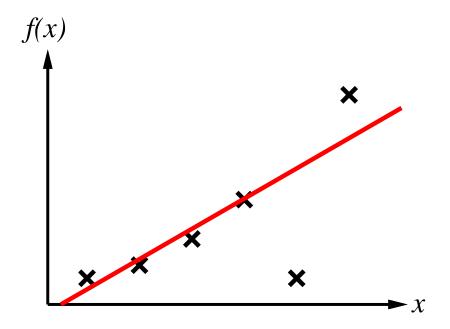


Model of the form y = f(x) + e
where x is a vector of attributes,
y is a target value and e is an
error term based on residuals,
i.e. vertical difference between
a data point and the model
curve

Test the model by suppling a new observation x value and calculating prediction y. If these are close then e is small

Construct/adjust h to agree with f on training set (h) is consistent if it agrees with f on all examples)

E.g., curve fitting:

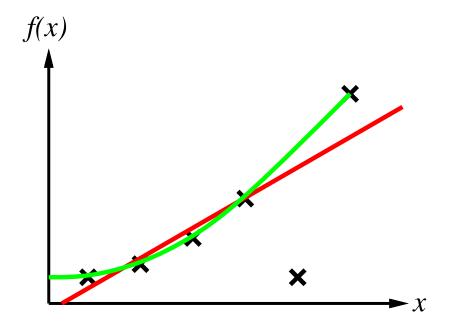


y = mx + c + e
where m is the slope, c is the
intercept and e is the error
between predictions and
observations

e would be based on sum of squared residuals

Construct/adjust h to agree with f on training set (h) is consistent if it agrees with f on all examples)

E.g., curve fitting:



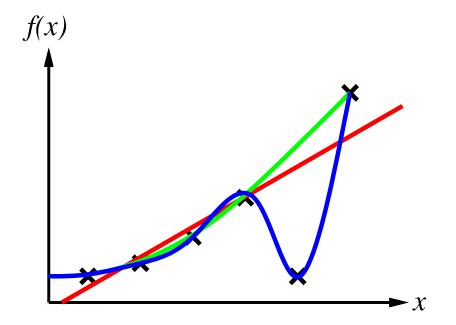
y is quadratic in x

This model captures the overall growth with increasing x, but treats one value as an outlier

Error is coefficient of determination (again)

Construct/adjust h to agree with f on training set (h) is consistent if it agrees with f on all examples)

E.g., curve fitting:



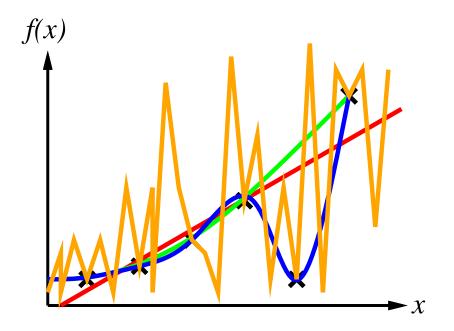
y is cubic in x

This model is close to all input data and has growth/ decline/growth behaviour

Error is coefficient of determination (again)

Construct/adjust h to agree with f on training set (h) is consistent if it agrees with f on all examples)

E.g., curve fitting:



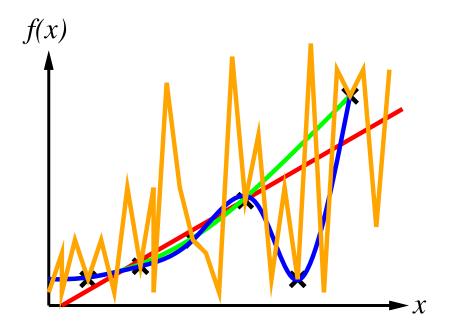
y is high degree or even non-smooth in x

This model is close to all input data but varies wildly between known values

Error is coefficient of determination (again)

Construct/adjust h to agree with f on training set (h is consistent if it agrees with f on all examples)

E.g., curve fitting:



Goodness of fit has increased at each stage

Simplicity has decreased at each stage

Which model will have least error when new x values are supplied?
i.e. lowest generalisation error

Ockham's razor: maximize a combination of consistency and simplicity

Attribute = factor = predictor variable = covariate = independent variable = explanatory variable = ...

Attribute-based representations

Examples described by attribute values (Boolean, discrete, continuous, etc.) E.g., situations where I will/won't wait for a table:

Example	Attributes									Target	
	Alt	Bar	Fri	Hun	Pat	Price	Rain	Res	Type	Est	WillWait
X_1	T	F	F	T	Some	\$\$\$	F	T	French	0–10	T
X_2	T	F	F	Τ	Full	\$	F	F	Thai	<i>30–60</i>	F
X_3	F	T	F	F	Some	\$	F	F	Burger	0–10	T
X_4	T	F	T	T	Full	\$	F	F	Thai	10–30	T
X_5	T	F	T	F	Full	<i>\$\$\$</i>	F	T	French	>60	F
X_6	F	T	F	T	Some	<i>\$\$</i>	T	T	Italian	0–10	T
X_7	F	T	F	F	None	\$	T	F	Burger	0–10	F
X_8	F	F	F	Τ	Some	<i>\$\$</i>	T	T	Thai	0–10	T
X_9	F	T	T	F	Full	\$	T	F	Burger	>60	F
X_{10}	T	T	T	Τ	Full	<i>\$\$\$</i>	F	T	Italian	10–30	F
X_{11}	F	F	F	F	None	\$	F	F	Thai	0–10	F
X_{12}	T	T	T	T	Full	\$	F	F	Burger	30–60	T

We know the
Target for
each data
instance,
hence
supervised
learning

Some
continuous
factors have
been made
discrete.
How (if) this
is done
changes the
learning
environment

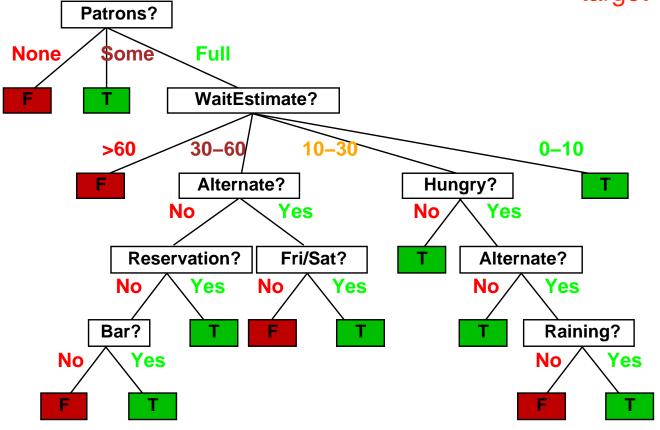
Classification of examples is positive (T) or negative (F)

Target = response = dependent variable = outcome variable = ...

Decision trees

One possible representation for hypotheses E.g., here is the "true" tree for deciding whether to wait:

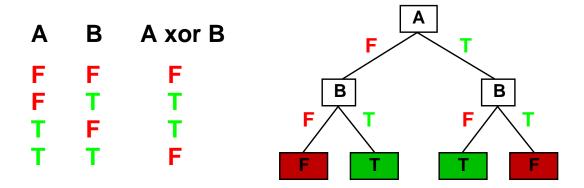
Model of the form y = f(X) + e where X is a matrix of attributes, y is a vector of target values and e is an error term



Test the model by running a new instance down the tree. If the predicted outcome matches the observed outcome, e = 0 for that instance

Expressiveness

Decision trees can express any function of the input attributes. E.g., for Boolean functions, truth table row \rightarrow path to leaf:



Trivially, there is a consistent decision tree for any training set w/ one path to leaf for each example (unless f nondeterministic in x) but it probably won't generalize to new examples

Prefer to find more compact decision trees

Recall: Deterministic means that the current state and a chosen action completely specify the next state.

If not deterministic, then stochastic

How many distinct decision trees with n Boolean attributes??

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= number of Boolean functions

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- = number of Boolean functions
- = number of distinct truth tables with 2^n rows

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E.g., with 6 Boolean attributes, there are 18,446,744,073,709,551,616 trees

So generating and testing candidate trees is intractable

We need to restrict the types of tree that we'll consider

How many distinct decision trees with n Boolean attributes??

- = number of Boolean functions
- = number of distinct truth tables with 2^n rows = 2^{2^n}

E.g., with 6 Boolean attributes, there are 18,446,744,073,709,551,616 trees

How many purely conjunctive hypotheses (e.g., $Hungry \land \neg Rain$)??

A conjunction is a Boolean expression of the form A ∧ B

How many distinct decision trees with n Boolean attributes??

- = number of Boolean functions
- = number of distinct truth tables with 2^n rows = 2^{2^n}

E.g., with 6 Boolean attributes, there are 18,446,744,073,709,551,616 trees

How many purely conjunctive hypotheses (e.g., $Hungry \land \neg Rain$)??

Each attribute can be in (positive), in (negative), or out

 \Rightarrow 3ⁿ distinct conjunctive hypotheses

More expressive hypothesis space

- increases chance that target function can be expressed
- increases number of hypotheses consistent w/ training set
 - ⇒ may get worse predictions



Decision tree learning

Aim: find a small tree consistent with the training examples

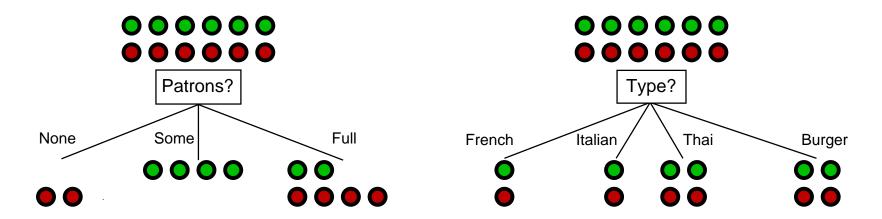
Idea: (recursively) choose "most significant" attribute as root of (sub)tree

```
function DTL(examples, attributes, default) returns a decision tree if examples is empty then return default else if all examples have the same classification then return the classification else if attributes is empty then return Mode(examples) else best \leftarrow \texttt{CHoose-Attributes}(attributes, examples) \\ tree \leftarrow \texttt{a} \text{ new decision tree with root test } best \\ \text{for each value } v_i \text{ of } best \text{ do} \\ examples_i \leftarrow \{\text{elements of } examples \text{ with } best = v_i\} \\ subtree \leftarrow \texttt{DTL}(examples_i, attributes - best, \texttt{Mode}(examples)) \\ \texttt{add a branch to } tree \text{ with label } v_i \text{ and subtree } subtree \\ \textbf{return } tree
```

Recursive partitioning: the feature space is recursively split into regions containing observations with similar response values.

Choosing an attribute

Idea: a good attribute splits the examples into subsets that are (ideally) "all positive" or "all negative"



Patrons? is a better choice—gives **information** about the classification

If the name Claude Shannon is new to you, then maybe pause this lecture and look up Information Theory. Shannon is - with Alan Turing - a figure of huge importance in Computer Science

log in my annotations is log base 2. Use the change of base formula to convert to base 2 from base 10 or base e

Information

Information answers questions

The more clueless I am about the answer initially, the more information is contained in the answer

Scale: 1 bit = answer to Boolean question with prior (0.5, 0.5)

Information in an answer when prior is $\langle P_1, \ldots, P_n \rangle$ is

$$H(\langle P_1, \dots, P_n \rangle) = \sum_{i=1}^n -P_i \log_2 P_i$$

https://planetcalc.com/2476/

(also called entropy of the prior)

$$H(0.9,0.1) = -0.9log(0.9) - 0.1log(0.1) = 0.469$$
 bits

$$H(0.5,0.5) = -0.5\log(0.5) - 0.5\log(0.5) = 1$$
 bit

$$H(1,0) = -\log(1) - \log(0) = 0$$
 bits

$$H(1/3,1/2,1/12,1/12) = (1/3)(1.58) + (1/2)(1) + (2/12)(3.58) = 1.626$$
 bits

Information contd.

Suppose we have p positive and n negative examples at the root

 $\Rightarrow H(\langle p/(p+n), n/(p+n)\rangle)$ bits needed to classify a new example E.g., for 12 restaurant examples, p=n=6 so we need 1 bit

An attribute splits the examples E into subsets E_i , each of which (we hope) needs less information to complete the classification

Let E_i have p_i positive and n_i negative examples

- $\Rightarrow H(\langle p_i/(p_i+n_i), n_i/(p_i+n_i)\rangle)$ bits needed to classify a new example
- ⇒ expected number of bits per example over all branches is

$$\sum_{i} \frac{p_i + n_i}{p + n} H(\langle p_i / (p_i + n_i), n_i / (p_i + n_i) \rangle)$$

For *Patrons*?, this is 0.459 bits, for *Type* this is (still) 1 bit

choose the attribute that minimizes the remaining information needed

```
For Patrons?: Enone = (2/12)H(0/2,2/2) = 0
       Esome = (4/12)H(4/4,0/4) = 0
Efull = (6/12)H(2/6,4/6) = 1/2(0.9813) = 0.459
     Sum these three to get 0.459 bits
```

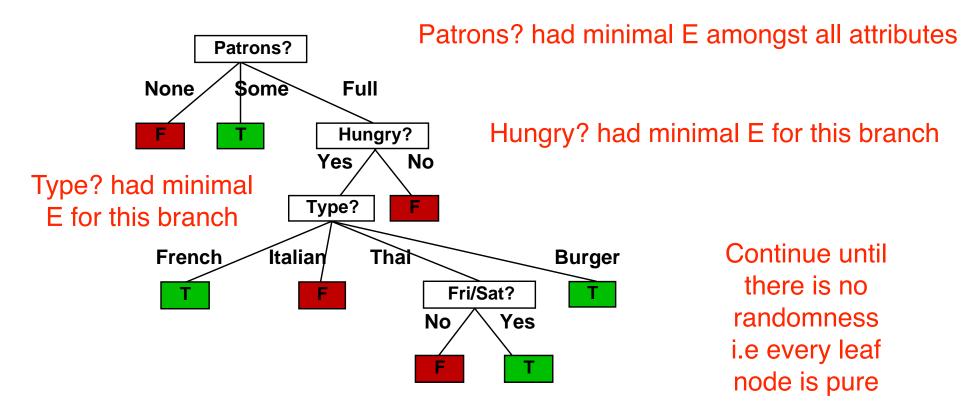
```
For Type?: H = 1 for each branch,
so 2/12 + 2/12 + 4/12 + 4/12 = 1
```

Chapter 18, Sections 1-3

0.459 < 1 so choose Patrons?

Example contd.

Decision tree learned from the 12 examples:



Substantially simpler than "true" tree—a more complex hypothesis isn't justified by small amount of data

Caveat: this is simpler than a real model for which we grow to a tolerance and then prune to get good generalisation error hapter 18, Sections 1-3

Performance measurement

How do we know that $h \approx f$? (Hume's **Problem of Induction**)

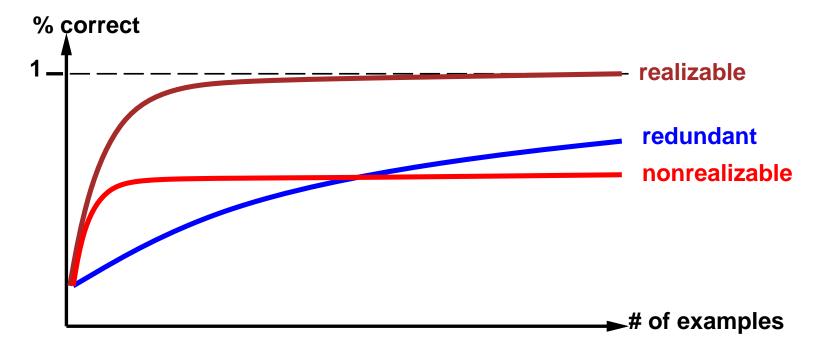
- 1) Use theorems of computational/statistical learning theory
- 2) Try h on a new test set of examples (use same distribution over example space as training set)

Training set size

Performance measurement contd.

Learning curve depends on

- realizable (can express target function) vs. non-realizable non-realizability can be due to missing attributes or restricted hypothesis class (e.g., thresholded linear function)
- redundant expressiveness (e.g., loads of irrelevant attributes)



Summary

Learning needed for unknown environments, lazy designers

Learning agent = performance element + learning element

Learning method depends on type of performance element, available feedback, type of component to be improved, and its representation

For supervised learning, the aim is to find a simple hypothesis that is approximately consistent with training examples

Decision tree learning using information gain

Learning performance = prediction accuracy measured on test set

Ex 18.5 (maybe 18.3 in your edition): "Suppose we generate a training set ..."

Ex 18.8 (maybe 18.6 in your edition): 3 binary inputs and one binary output