# **Foundations of Al**

# 16. Statistical Machine Learning

Bayesian Learning and Why Learning Works Wolfram Burgard, Bernhard Nebel, and Andreas Karwath

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- Statistical learning
- Why learning works

## **Statistical Learning Methods**

- In MDPs probability and utility theory allow agents to deal with uncertainty.
- To apply these techniques, however, the agents must first learn their probabilistic theories of the world from experience.
- We will discuss statistical learning methods as robust ways to learn probabilistic models.

## **An Example for Statistical Learning**

- The key concepts are data (evidence) and hypotheses.
- A candy manufacturer sells five kinds of bags that are indistinguishable from the outside:
  - $h_1$ : 100% cherry
  - h<sub>2</sub>: 75% cherry and 25% lime
  - h<sub>3</sub>: 50% cherry and 50% lime
  - h<sub>4</sub>: 25% cherry and 75% lime
  - h<sub>5</sub>: 100% lime
- Given a sequence d<sub>1</sub>, ..., d<sub>N</sub> of candies observed, what is the most likely flavor of the next piece of candy?

### **Bayesian Learning**

- Calculates the probability of each hypothesis, given the data.
- It then makes predictions using all hypotheses weighted by their probabilities (instead of a single best hypothesis).
- Learning is reduced to probabilistic inference.

## **Application of Bayes Rule**

- Let *D* represent all the data with observed value *d*.
- The probability of each hypothesis is obtained by Bayes rule:

$$P(h_i \mid \mathbf{d}) = \alpha P(\mathbf{d} \mid h_i) P(h_i)$$

- The manufacturer tells us that the prior distribution over h<sub>1</sub>, ..., h<sub>5</sub> is given by
   <.1, .2, .4, .2, .1>
- We compute the likelihood of the data under the assumption that the observations are independently and identically distributed (i.i.d.):

$$P(\mathbf{d} \mid h_i) = \prod_j P(d_j \mid h_i)$$

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#### How to Make Predictions?

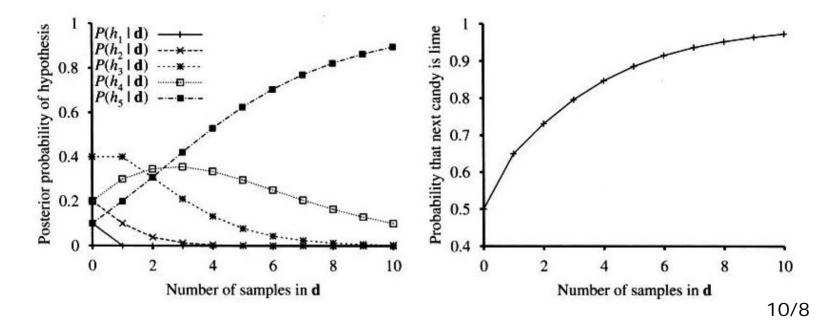
 Suppose we want to make predictions about an unknown quantity X given the data d.

$$\mathbf{P}(X \mid \mathbf{d}) = \sum_{i} \mathbf{P}(X \mid h_i, \mathbf{d}) \mathbf{P}(h_i \mid \mathbf{d})$$
$$= \sum_{i} \mathbf{P}(X \mid h_i) \mathbf{P}(h_i \mid \mathbf{d})$$

- Predictions are weighted averages over the predictions of the individual hypotheses.
- The key quantities are the hypothesis prior P(h<sub>i</sub>) and the likelihood P(d/h<sub>i</sub>) of the data under each hypothesis.

#### **Example**

- Suppose the bag is an all-lime bag  $(h_5)$
- The first 10 candies are all lime.
- Then  $P(d/h_3)$  is 0.5<sup>10</sup> because half the candies in an  $h_3$  bag are lime.
- Evolution of the five hypotheses given 10 lime candies were observed (the values start at the prior!).



#### **Observations**

- The true hypothesis often dominates the Bayesian prediction.
- For any fixed prior that does not rule out the true hypothesis, the posterior of any false hypothesis will eventually vanish.
- The Bayesian prediction is optimal and, given the hypothesis prior, any other prediction will be correct less often.
- It comes at a price that the hypothesis space can be very large or infinite.

#### Maximum a Posteriori (MAP)

- A common approximation is to make predictions based on a single most probable hypothesis.
- The maximum a posteriori (MAP) hypothesis is the one that maximizes  $P(h_i/d)$ .

 $\mathbf{P}(X \mid \mathbf{d}) \approx \mathbf{P}(X \mid h_{MAP})$ 

- In the candy example,  $h_{MAP} = h_5$  after three lime candies in a row.
- The MAP learner the predicts that the fourth candy is lime with probability 1.0, whereas the Bayesian prediction is still 0.8.
- As more data arrive, MAP and Bayesian predictions become closer.
- Finding MAP hypotheses is often much easier than Bayesian learning.

## Maximum-Likelihood Hypothesis (ML)

- A final simplification is to assume a uniform prior over the hypothesis space.
- In that case MAP-learning reduces to choosing the hypothesis that maximizes  $P(d/h_i)$ .
- This hypothesis is called the maximumlikelihood hypothesis (ML).
- ML-learning is a good approximation to MAP learning and Bayesian learning when there is a uniform prior and when the data set is large.

## Why Learning Works

How can we decide that *h* is close to *f* when *f* is unknown?

→ Probably approximately correct

Stationarity as the basic assumption of PAC-Learning: training and test sets are selected from the same population of examples with the same probability distribution.

Key question: how many examples do we need?

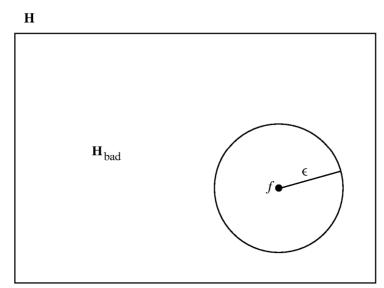
- *X* Set of examples
- *D* Distribution from which the examples are drawn
- *H* Hypothesis space  $(f \in H)$
- *m* Number of examples in the training set

 $error(h) = P(h(x) \neq f(x) \mid x \text{ drawn from } D) \leq \epsilon$ 

## **PAC-Learning**

A hypothesis h is approximately correct if  $error(h) \leq \epsilon$ .

To show: After the training period with *m* examples, with high probability, all consistent hypotheses are approximately correct.



How high is the probability that a wrong hypothesis  $h_b \in H_{bad}$  is consistent with the first *m* examples?

#### Sample Complexity

Assumption:  $error(h) > \epsilon \implies$  $P(h_{h} \text{ is consistent with 1 example}) < (1-\epsilon)$  $P(h_h \text{ is consistent with } N \text{ examples}) \leq (1 - \epsilon)^N$  $P(H_{bad} \text{ contains a consistent } h) \leq |H_{bad}|(1-\epsilon)^N$ Since  $|H_{bad}| \leq |H|$  $P(H_{bad} \text{ contains a consistent } h) \leq |H|(1-\epsilon)^N$ We want to limit this probability by some small number  $\delta$ :  $|H|(1-\epsilon)^N < \delta$ Since  $(1-\epsilon) \leq e^{-\epsilon}$ , we derive

$$N \geq \frac{1}{\epsilon} \left( \log \left( \frac{1}{\delta} \right) + \log |H| \right)$$

Sample Complexity: Number of required examples, as a function of  $\epsilon$  and  $\delta$ .

# Sample Complexity (2)

**Example:** Boolean functions

The number of Boolean functions over *n* attributes is  $|H| = 2^{2^n}$ .

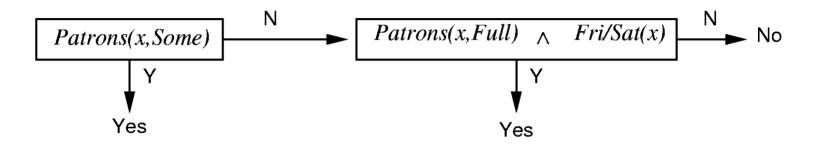
The sample complexity therefore grows as  $2^n$ .

Since the number of possible examples is also  $2^n$ , any learning algorithm for the space of all Boolean functions will do no better than a lookup table, if it merely returns a hypothesis that is consistent with all known examples.

# **Learning from Decision Lists**

In comparison to decision trees:

- The overall structure is simpler
- The individual tests are more complex



This represents the hypothesis

 $H_4$ :  $\forall x WillWait(x) \Leftrightarrow Patrons(x, some) \lor [Patrons(x, full) \land Fri/Sat(x)]$ 

If we allow tests of arbitrary size, then any Boolean function can be represented.

k-DL: Language with tests of length  $\leq k$ .



# Learnability of k-DL

function DECISION-LIST-LEARNING(examples) returns a decision list, No or failure

if examples is empty then return the value No
t ← a test that matches a nonempty subset examples, of examples such that the members of examples, are all positive or all negative
if there is no such t then return failure
if the examples in examples, are positive then o ← Yes
else o ← No
return a decision list with initial test t and outcome o and remaining elements given by DECISION-LIST-LEARNING(examples - examples,)

$$| k-DL(n) | \leq 3^{|Conj(n,k)|} | Conj(n,k)! |$$
 (Yes,No,no-Test,all permutations)

$$|Conj(n,k)| = \sum_{i=0}^{k} \binom{2n}{i} = O(n^k)$$

(Combination without repeating pos/neg attributes)

 $| k-DL(n) |= 2^{O(n^k log(n^k))}$  (with Euler's summation formula)  $m \ge \frac{1}{\epsilon} (ln(\frac{1}{\delta}) + O(n^k log(n^k)))$ 

#### Summary (Statistical Learning Methods)

- Bayesian learning techniques formulate learning as a form of probabilistic inference.
- Maximum a posteriori (MAP) learning selects the most likely hypothesis given the data.
- Maximum likelihood learning selects the hypothesis that maximizes the likelihood of the data.

#### Summary (Statistical Learning Theory)

Inductive learning as learning the representation of a function from example input/output pairs.

- Decision trees learn deterministic Boolean functions.
- PAC learning deals with the complexity of learning.
- Decision lists as functions that are easy to learn.