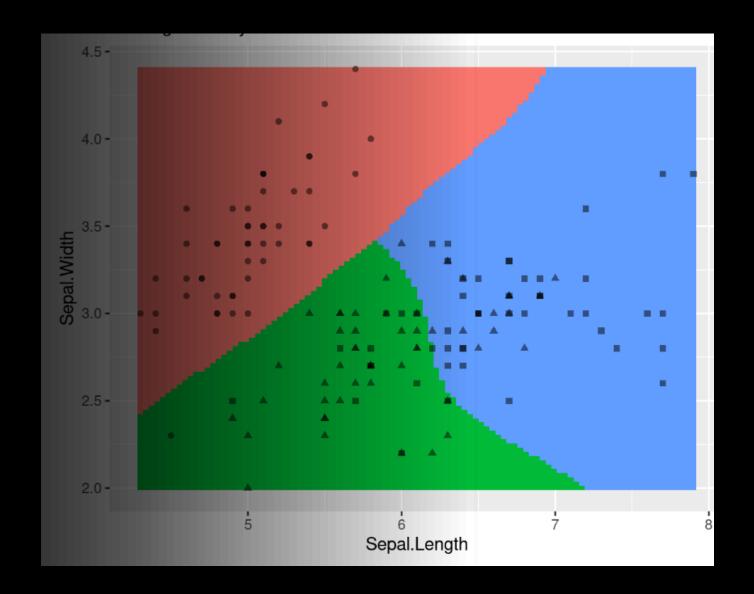
CS 5/7320 Artificial Intelligence

Learning from Examples AIMA Chapter 19

Slides by Michael Hahsler Based on slides by Dan Klein, Pieter Abbeel, Sergey Levine and A. Farhadi. All CS188 materials are at http://ai.berkeley.edu with figures from the AIMA textbook.

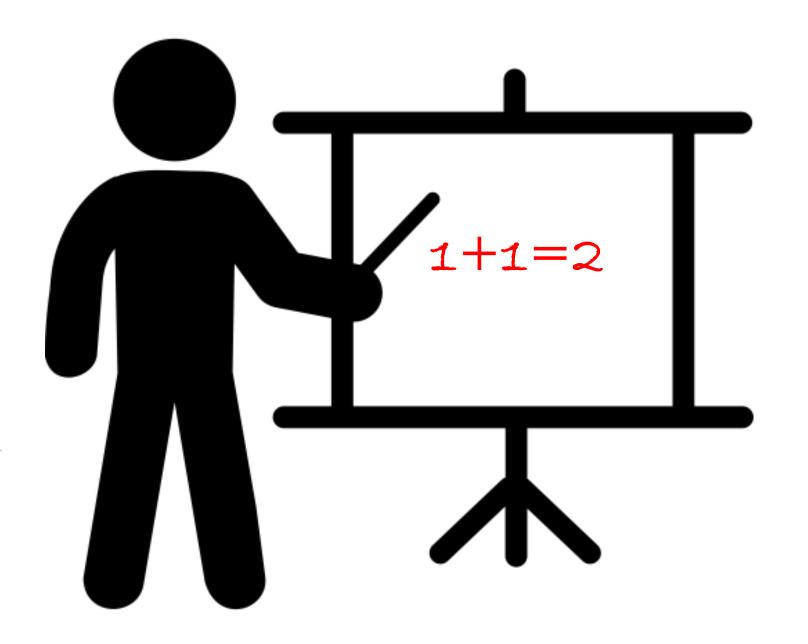




Learning from Examples: Machine Learning

- **Up until now:** hand-craft algorithms to make rational/optimal or at least good decisions. Examples: Search strategies, heuristics.
- Learning: Improve performance after making observations about the world. That is, learn what works and what doesn't.
- Machine learning: how to build a model from data/experience
 - Supervised Learning: Learn a function to map input to output. Examples:
 - Use a naïve Bayesian classifier to distinguish between spam/no spam
 - Learn a playout policy to simulate games (current board -> good move)
 - Unsupervised Learning: Organize data (e.g., clustering, embedding)
 - Reinforcement Learning: Learn from rewards/punishment (e.g., winning a game).
- Learning vs. hard coding the agent function
 - Designer cannot anticipate all possible future situations.
 - Designer may have examples but does not know how to program a solution.

Supervised Learning



Supervised Learning

Examples

- Input-output pairs: $E=(x_1,y_1),\ldots,(x_i,y_i),\ldots,(x_N,y_N),$ where x is a vectors called the feature vector.
- We assume that the examples are produced iid (with noise and errors) from a target function y = f(x).

Learning problem

- Given a hypothesis space H of possible models.
- Find a hypothesis $h \in H$ such that $\hat{y}_i = h(x_i) \approx y_i$
- That is, we want to approximate f by h using E.

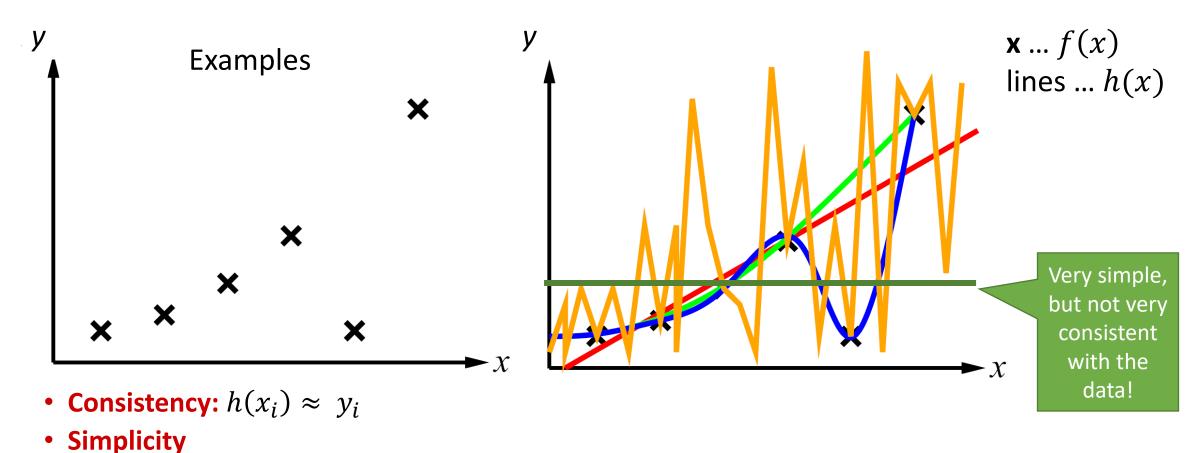
Includes

- Classification (outputs = class labels). E.g. x is an email and f(x) is spam / ham
- Regression (outputs = real numbers). E.g. x is a house and f(x) is its selling price

H

Consistency vs. Simplicity

Example: Univariate curve fitting (regression, function approximation)



Measuring Consistency using Loss

Goal of learning: Find a hypothesis that makes good predictions that are consistent with the examples $E = (x_1, y_1), ..., (x_i, y_i), ..., (x_N, y_N)$. $\hat{y} = h(x) \approx y$. That is,

- Measure mistakes: Loss function $L(y, \hat{y})$

$$L_1(y, \hat{y}) = |y - \hat{y}|$$

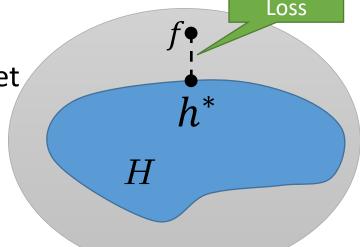
• Absolute-value loss
$$L_1(y,\hat{y}) = |y - \hat{y}|$$
• Squared-error loss
$$L_2(y,\hat{y}) = (y - \hat{y})^2$$

For Regression

Log loss and others...

Empirical loss: average loss over the N examples in the dataset.

$$EmpLoss_{L,E}(h) = \frac{1}{|E|} \sum_{(x,y) \in E} L(y,h(x))$$



Learning Consistent h by Minimizing the Loss

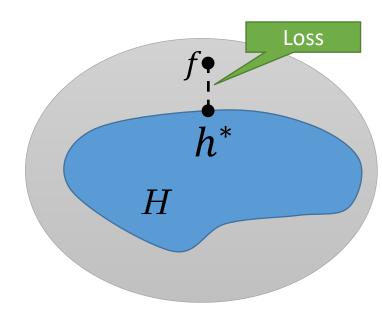
Empirical loss

$$EmpLoss_{L,E}(h) = \frac{1}{|E|} \sum_{(x,y) \in E} L(y,h(x))$$

• Learning the best hypothesis (approximation)

$$h^* = \underset{h \in H}{\operatorname{argmin}} EmpLoss_{L,E}(h)$$

- Reasons for $h^* \neq f$
 - a) Realizability: $f \notin H$
 - b) f is nondeterministic or examples are noisy
 - c) Computationally intractable to search all H



The Bayes Classifier

For 0/1 loss, the empirical loss is minimized by the model that predicts for each x the most likely class y using MAP (Maximum a posteriori) estimates. This is called the Bayes classifier.

$$h(x)^* = \underset{y}{\operatorname{argmax}} P(Y = y \mid X = x) = \underset{y}{\operatorname{argmax}} \frac{P(x \mid y) P(y)}{P(x)} = \underset{y}{\operatorname{argmax}} P(x \mid y) P(y)$$

Optimality: The **Bayes classifier is optimal** is the most consistent classifier and guarantees the lowest possible error called the **Bayes error rate**.

Issue: The classifier is not simple with $P(x \mid y) P(y) = P(x, y)$

- It **needs the complete joint probability** which requires in the general case a probability table with one entry for each possible value for the feature vector *x*.
- This is impractical and most classifiers try to approximate the bayes Classifier using fewer parameters.

Simplicity

Ease of use

- Simpler hypotheses have fewer parameters to store.
- Simpler hypotheses are easier to estimate.

Generalization: How well does the hypothesis perform on new data?

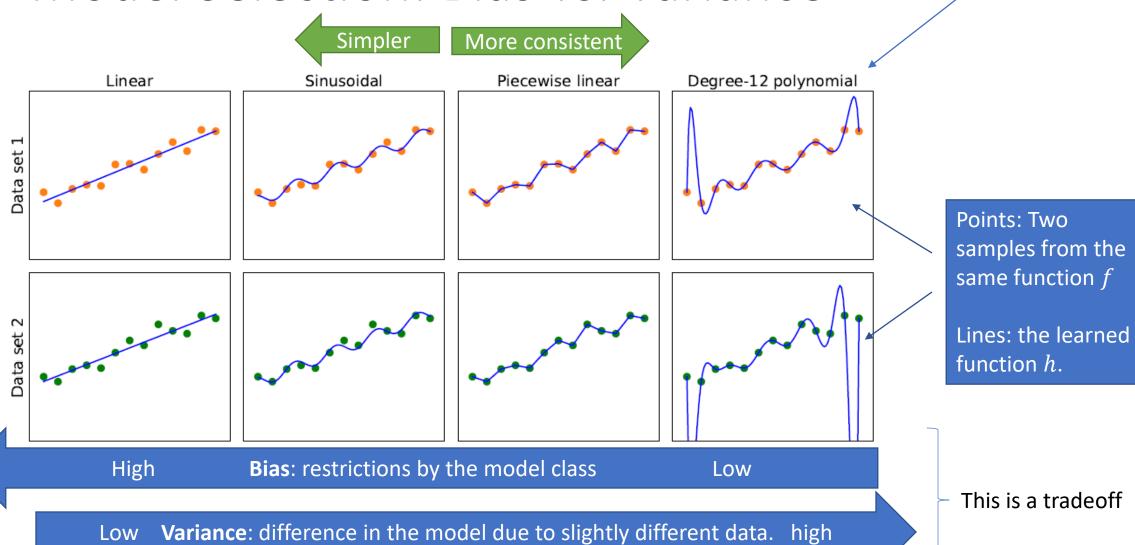
- We do not want the model to be too specific to the training examples (an issue called overfitting).
- Simpler models typically generalize better to new examples.

How to achieve simplicity?

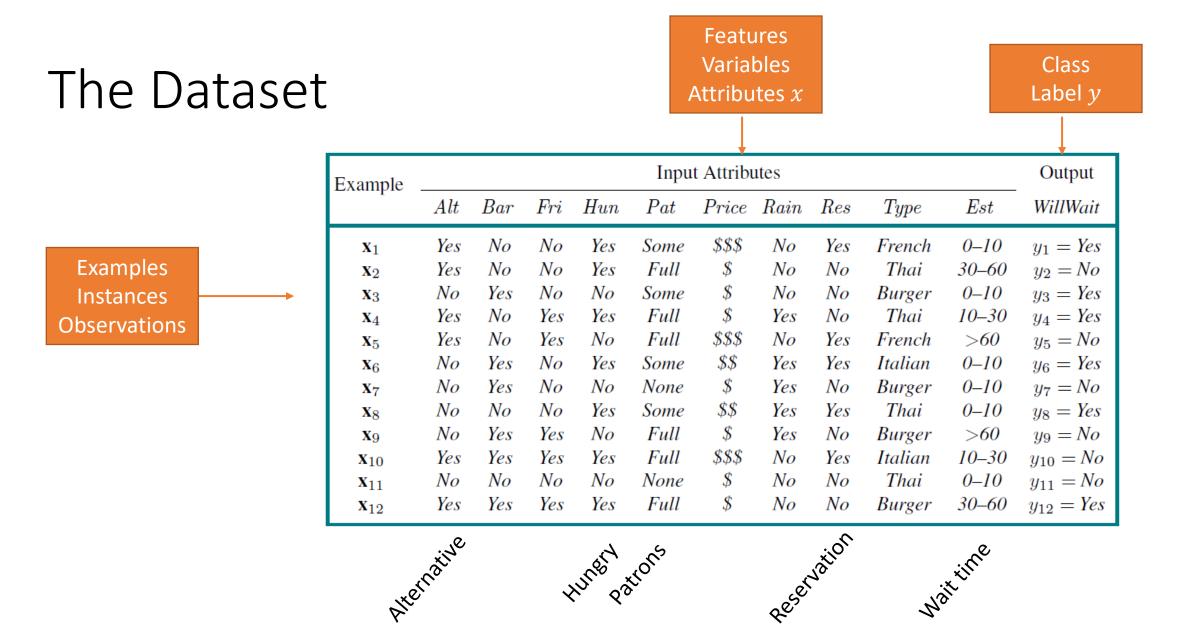
- a) Restrict H to simpler models (e.g., assumptions like independence, linear models are called model bias)
- b) Feature selection (use fewer variables from the feature vector x)
- c) Regularization (penalize for complexity)

$$h^* = \underset{h \in H}{\operatorname{argmin}} \left[EmpLoss_{L,E}(h) + \underset{\Lambda}{\lambda} Complexity(h) \right]$$
Penalty term

Model Selection: Bias vs. Variance



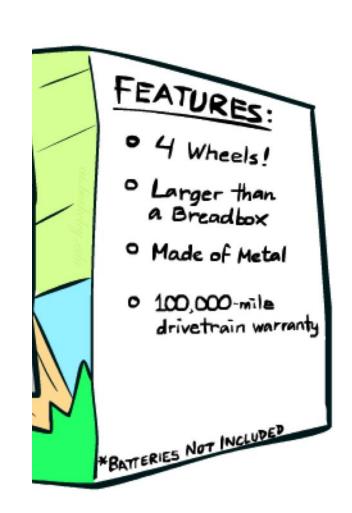
Data



Find a hypothesis (called "model") to predict the class given the features.

Feature Engineering

- Add information sources as new variables to the model.
- Add derived features that help the classifier (e.g., x^2).
- Example for Spam detection: In addition to words
 - Have you emailed the sender before?
 - Have 1000+ other people just gotten the same email?
 - Is the header information consistent?
 - Is the email in ALL CAPS?
 - Do inline URLs point where they say they point?
 - Does the email address you by (your) name?
- Feature Selection: Which features should be used in the model is a model selection problem (choose between models with different features).



Training and Testing



Model Evaluation (Testing)

The model was trained on the training examples E. We want to test how well the model will perform on new examples T (i.e., how well it **generalizes to new data**).

• Testing loss: Calculate the empirical loss for predictions on a testing data set T that is different from the data used for training.

$$EmpLoss_{L,T}(h) = \frac{1}{|T|} \sum_{(x,y) \in T} L(y,h(x))$$

• For classification we often use the **accuracy** measure, the proportion of correctly classified test examples.

$$accuracy(h, T) = \frac{1}{|T|} \sum_{(x,y) \in T} [h(x) = y] = 1 - EmpLoss_{L_{0/1},T}(h)$$

Training a Model

• Models are "trained" (learned) on **the training data.** This involved estimating the model parameters (the model): E.g., probabilities, weights, factors.

• **Hyperparameters**: Many learning algorithms have choices for. E.g., learning rate, regularization λ , maximal decision tree depth, selected features. The algorithm tries to optimizes the model parameters given user-specified hyperparameters.

We need to tune the hyperparameters!



Hyperparameter Tuning/Model Selection

- 1. Hold a validation data set back from the training data.
- 2. Learn models using the training set with different hyperparameters. Often a grid of possible hyperparameter combinations or some greedy search is used.
- 3. Evaluate the models using the validation data and choose the model with the best accuracy. Selecting the right type of model, hyperparameters and features is called model selection.
- 4. Learn the final model with the chosen hyperparameters using all training (including validation data).
- Notes:
 - The validation set was not used for training, so we get generalization accuracy for the different hyperparameter setings.
 - If no model selection is necessary, then no validation set is used.

Training Data Validation

Data

Test Data

Testing a Model

• After the model is selected, the final model is evaluated against the test set to **estimate the final model accuracy**.

• Very important: never "peek" at the test set during learning!



How to Split the Dataset

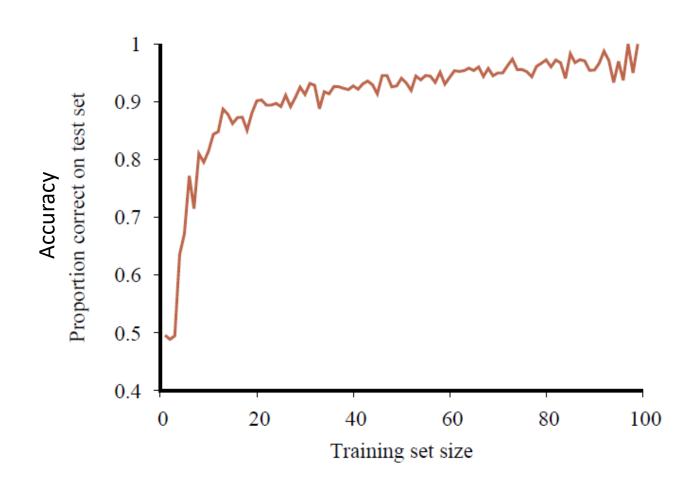
- Random splits: Split the data randomly in, e.g., 60% training, 20% validation, and 20% testing.
- **Stratified splits:** Like random splits, but balance classes and other properties of the examples.
- k-fold cross validation: Use training & validation data better
 - split the training & validation data randomly into k folds.
 - For k rounds hold 1 fold back for testing and use the remaining k-1 folds for training.
 - Use the average error/accuracy as a better estimate.
 - Some algorithms/tools do this internally.
- **LOOCV** (leave-one-out cross validation): k=n used if very little data is available.

Training Data

Validation Data

> Test Data

Learning Curve: The Effect the Training Data Size



Accuracy of a classifier when the amount of available training data increases.

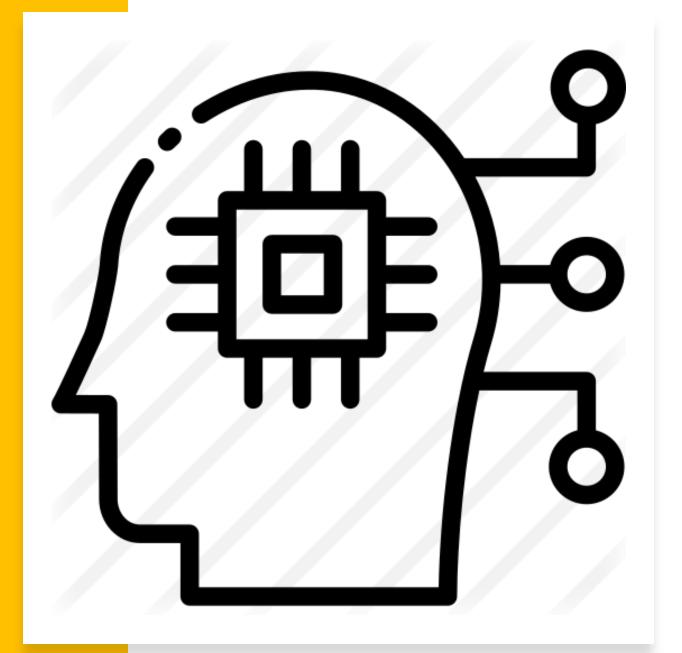
More data is better!

Comparing to a Baselines

- First step: get a baseline
 - Baselines are very simple straw man model.
 - Helps to determine how hard the task is.
 - Helps to find out what a good accuracy is.



- Weak baseline: The most frequent label classifier
 - Gives all test instances whatever label was most common in the training set.
 - Example: For spam filtering, give every message the label "ham."
 - Accuracy might be very high if the problem is skewed (called class imbalance).
 - Example: If calling everything "ham" gets already 66% right, so a classifier that gets 70% isn't very good...
- Strong baseline: For research, we typically compare to previous work as a baseline.



Types of Models

Regression: Predict a number

Classification: Predict a label



Regression: Linear Regression

Model:
$$h_{\mathbf{w}}(\mathbf{x}_{j}) = w_{o} + w_{1}x_{j,1} + \dots + w_{n}x_{j,n} = \sum_{i} w_{i}x_{j,i} = \mathbf{w}^{T}\mathbf{x}_{j}$$

Empirical Loss: $L(w) = ||Xw - y||^2$

Gradient: $\nabla L(\mathbf{w}) = 2\mathbf{X}^T(\mathbf{X}\mathbf{w} - \mathbf{y})$

Find: $\nabla L(\mathbf{w}) = 0$

Gradient descend:

$$\mathbf{w} = \mathbf{w} - \alpha \nabla L(\mathbf{w})$$

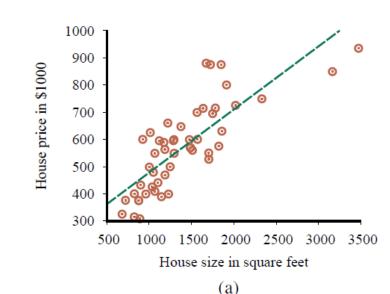
Analytical solution:

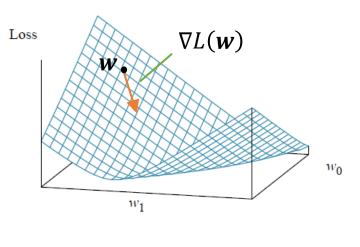
$$\boldsymbol{w}^* = (\boldsymbol{X}^T \boldsymbol{X})^{-1} \boldsymbol{X}^T \boldsymbol{y}$$
Pseudo inverse

Squared error loss over the whole data matrix \boldsymbol{X}

The gradient is a vector of partial derivatives

$$\nabla L(\mathbf{w}) = \left[\frac{\partial L}{\partial w_1}(\mathbf{w}), \frac{\partial L}{\partial w_2}(\mathbf{w}), \dots, \frac{\partial L}{\partial w_n}(\mathbf{w}) \right]^T$$





(b)

Naïve Bayes Classifier

• Approximates a Bayes classifier with the **naïve independence assumption** that all n features are conditional independent given the class.

$$h(x) = \underset{y}{\operatorname{argmax}} P(y) \prod_{i=1}^{n} P(x_i \mid y)$$

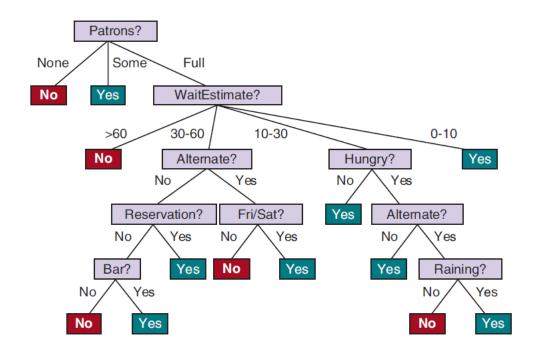
 We have only used discrete features so far, but it can be extended to continuous features. Gaussian Naïve Bayes Classifier assumes that continuous features have:

$$P(x_i \mid y) \sim N(\mu_y, \sigma_y)$$

The parameters for $N(\mu_y, \sigma_y)$ are estimated from data.

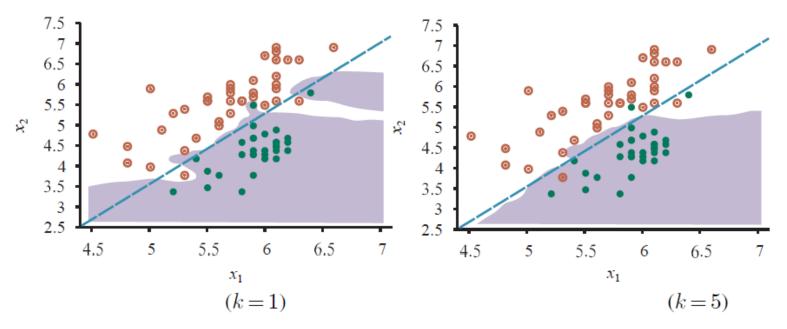
Decision Trees

Example	Input Attributes										Output
	Alt	Bar	Fri	Hun	Pat	Price	Rain	Res	Type	Est	WillWait
\mathbf{x}_1	Yes	No	No	Yes	Some	\$\$\$	No	Yes	French	0–10	$y_1 = Yes$
\mathbf{x}_2	Yes	No	No	Yes	Full	\$	No	No	Thai	<i>30–60</i>	$y_2 = No$
\mathbf{x}_3	No	Yes	No	No	Some	\$	No	No	Burger	0-10	$y_3 = Yes$
\mathbf{x}_4	Yes	No	Yes	Yes	Full	\$	Yes	No	Thai	10-30	$y_4 = Yes$
\mathbf{x}_5	Yes	No	Yes	No	Full	\$\$\$	No	Yes	French	>60	$y_5 = No$
\mathbf{x}_6	No	Yes	No	Yes	Some	\$\$	Yes	Yes	Italian	0-10	$y_6 = Yes$
\mathbf{x}_7	No	Yes	No	No	None	\$	Yes	No	Burger	0-10	$y_7 = No$
\mathbf{x}_8	No	No	No	Yes	Some	\$\$	Yes	Yes	Thai	0-10	$y_8 = Yes$
X 9	No	Yes	Yes	No	Full	\$	Yes	No	Burger	>60	$y_9 = No$
\mathbf{x}_{10}	Yes	Yes	Yes	Yes	Full	\$\$\$	No	Yes	Italian	10-30	$y_{10} = No$
\mathbf{x}_{11}	No	No	No	No	None	\$	No	No	Thai	0-10	$y_{11} = No$
\mathbf{x}_{12}	Yes	Yes	Yes	Yes	Full	\$	No	No	Burger	<i>30–60</i>	$y_{12} = Yes$



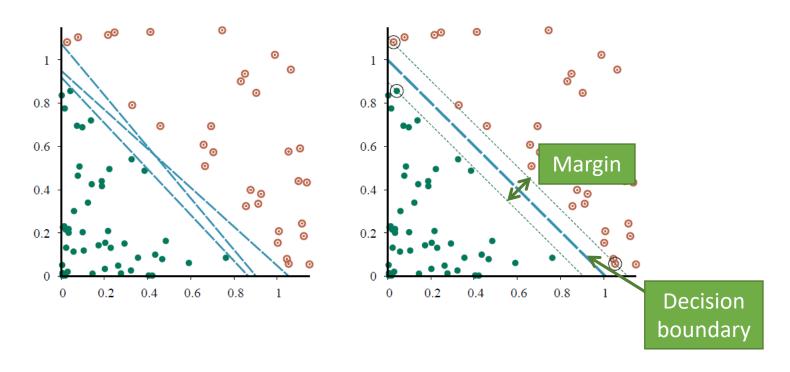
- A **sequence of decisions** represented as a tree.
- Many implementations that differ by
 - How to select features to split?
 - When to stop splitting?
 - Is the tree pruned?
- Approximates a Bayesian classifier by $h(x) = \underset{v}{\operatorname{argmax}} P(Y = y \mid \operatorname{leafNodeMatching}(x))$

K-Nearest Neighbors Classifier



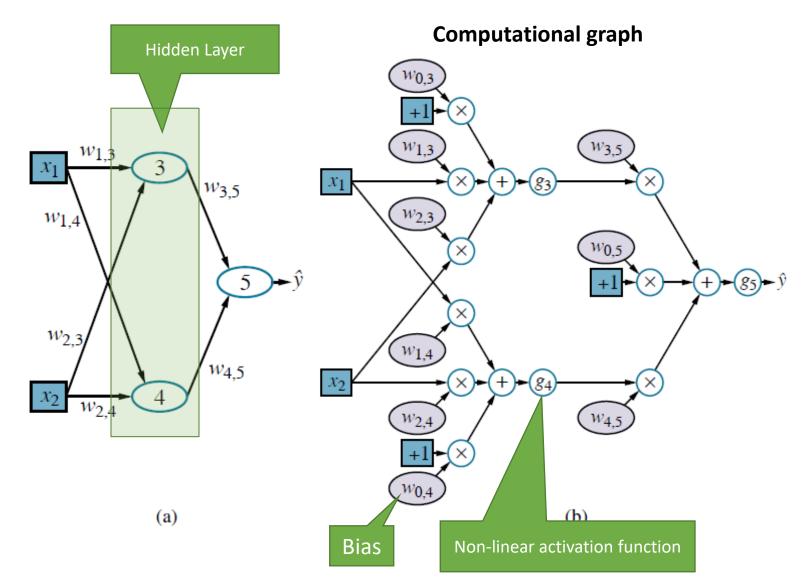
- Class is predicted by looking at the majority in the set of the k nearest **neighbors**. k is a hyperparameter. Larger k smooth the decision boundary.
- Neighbors are found using a distance measure (e.g., Euclidean distance between points).
- Approximates a Bayesian classifier by $h(x) = \operatorname{argmax} P(Y = y \mid \operatorname{neighborhood}(x))$

Support Vector Machine (SVM)



- Linear classifier that finds the maximum margin separator using only the points that are "support vectors" and quadratic optimization.
- The kernel trick can be used to learn non-linear decision boundaries.

Artificial Neural Networks/Deep Learning



- Represent $\hat{y} = h(x)$ as a network of weighted sums with non-linear activation functions (e.g., logistic, ReLU, tanh).
- Learn weights \mathbf{w} from examples using backpropagation of prediction errors $L(\hat{y}, y)$ (gradient descend).
- ANNs can approximate any function (no bias).
 Regularization is typically used to avoid overfitting.
- **Deep learning** adds more layer types (e.g., convolution layers)

Other Models and Methods

Many other models exist

• Generalized linear model (GLM): A model family that includes linear regression and the classification method logistic regression.

Often used methods

- **Regularization:** enforce simplicity by using a penalty for complexity.
- **Kernel trick:** Let a linear classifier learn non-linear decision boundaries (= a linear boundary in a high dimensional space).
- Ensemble Learning: Use many models and combine the results (e.g., random forest, boosting).

Example Use of ML for Intelligent Agents

Learn Actions

 Directly learn the best action from examples. This model can also be used as a playout policy for Monte Carlo tree search and selfplay.

action = h(state)

Learn Heuristics

 Learn evaluation functions for heuristic minimax search from examples.

eval = h(state)

Perception

- Natural language processing/language models: Use deep learning to understand and create test.
- Speech recognition
- Vision: Images/video processing, object recognition