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 [\(http://ai.berkeley.edu](http://ai.berkeley.edu/)) with figures from the AIMA textbook.

Naive Bayes

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ML and Agents

DeepAi.org with prompt: "A happy cartoon robot with an artificial neural network for a brain on white background learning to play chess"

Learning from Examples: Machine Learning

Up until now in this course:

• **Hand-craft algorithms** to make rational/optimal or at least good decisions. Examples: Search strategies, heuristics.

Issues

- Designer cannot anticipate all possible future situations.
- Designer may have examples but does not know how to program a solution.

Machine Learning

- Learning = Improve performance after making observations about the world. That is, learn what works and what doesn't.
- We learn a model that decides on the actions to take. This is called the "performance element."
- The goal is to get closer to optimal decisions. I.e., it is an optimization problem.

From Chapter 2: Agents that Learn

The **learning element** modifies the performance element to improve its performance.

Types of Using Machine Learning

- 1. What **component** of the performance element is learned? E.g., how to select action, estimate the utility of a state, …
- 2. What **representation** (model) is used in the component? Linear regression, rules, trees, neural nets,…
- 3. What **feedback** is available for learning?
	- **Unsupervised Learning**: No feedback, just organize data (e.g., clustering, embedding)
	- **Supervised Learning**: Uses a data set with correct answers. Learn a function (model) to map an input (e.g., state) to an output (e.g., action or utility). Examples:
		- Use a naïve Bayesian classifier to distinguish between spam/no spam
		- Learn a playout policy to simulate games (current board -> good move)
	- **Reinforcement Learning**: Learn from rewards/punishment (e.g., winning a game) obtained via interaction with the environment over time.

We focus here on supervised learning

Supervised Learning

Supervised Learning As Function Approximation

- Examples
	- We assume there exists a target function $y = f(x)$ that produces iid (independent and identically distributed) examples possibly with noise and errors.
	- Examples are observed input-output pairs $E = (x_1, y_1), ..., (x_i, y_i), ..., (x_N, y_N)$, where x is a vectors called the feature vector.
- Learning problem
	- Given a hypothesis space *H* of representable models.
	- Find a hypothesis $h \in H$ such that $\hat{y}_i = h(x_i) \approx y_i \ \forall i$
	- That is, we want to approximate f by h using E.
- Supervised learning 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.

Consistency vs. Simplicity

Example: Univariate curve fitting (regression, function approximation)

- **Consistency:** $h(x_i) \approx y_i$
- **Simplicity:** small number of model parameters

Measuring Consistency using Loss

Goal of learning: Find a hypothesis that makes predictions that are consistent with the examples $E = (x_1, y_1), ..., (x_i, y_i), ..., (x_N, y_N)$.

That is, $\ddot{y} = h(x) \approx y.$

- **Measure mistakes:** Loss function $L(y, \hat{y}) = L(f(x), h(x))$
	- Absolute-value loss $L_1(y, \hat{y}) = |y \hat{y}|$
	- Squared-error loss $L_2(y, \hat{y}) = (y \hat{y})^2$
	- \cdot 0/1 loss

$$
L_0
$$
₁ $(y, \hat{y}) = 0$ if $y = \hat{y}$, else 1

For Regression

• Log loss, cross-entropy loss and many others…

Learning Consistent h by Minimizing the Loss

- Empirical loss $EmpLoss_{L,E}(h) =$ 1 $|\overline{E}|$ $\sum_{(x,y)\in E}$ $(x, y) \in E$ $L(y, h(x))$
- Find the best hypothesis that minimizes the loss $h^* = \operatornamewithlimits{argmin}\limits_{h \in H} EmpLoss_{L,E}(h)$ $h \in H$
- Reasons for $h^* \neq f$
	- a) Realizability: $f \notin H$
b) f is nondeterminis
	- b) f is nondeterministic or examples are noisy.
c) It is computationally intractable to search all
	- c) It is computationally intractable to search all H , so we use a non-optimal heuristic.

The Most Consistent Classifier 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.

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

Optimality: The **Bayes classifier is optimal for 0/1 loss.** It is the most consistent classifier possible with the lowest possible error called the **Bayes error rate**. No better classifier is possible!

Issue: The classifier requires to learn $P(x | y) P(y) = P(x, y)$ from the examples.

- 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 (unless a simple Bayes network exists) and most classifiers try to approximate the Bayes classifier using a **simpler model** with fewer parameters.

Simplicity

Ease of use

• Simpler hypotheses have fewer model parameters to estimate and store.

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) Model bias:** Restrict *H* to simpler models (e.g., assumptions like independence, only consider linear models).
- **b) Feature selection:** use fewer variables from the feature vector x **c) Regularization:** penalize model for its complexity (e.g., number c
- **Regularization:** penalize model for its complexity (e.g., number of parameters)

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

Model Selection: Bias vs. Variance

The Dataset

Examples

(Instances,

Observation)

Feature vector x (Features, Variables, Attributes)

Class Label y

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

4 Wheels!

FEATURE

- ^o Larger than
a Breadbox
	- Made of Metal

*BATTERIES NOT INCLUDED

100,000-mile ۰ drivetrain warranty

Feature Engineering

- Add information sources as new variables to the model.
- Add derived features that help the classifier (e.g., x_1x_2, x_1^2).
- Embedding: E.g., convert words to vectors where vector similarity between vectors reflects semantic similarity.
- 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 email in ALL CAPS?
- **Feature Selection**: Which features should be used in the model is a model selection problem (choose between models with different features).

Data in AI

- Data in AI can come from many sources
	- **Observation**: Record video of a task being performed.
	- **Existing Data**: Download documents from the internet to train Large Language Models.
	- **Simulation**: E.g., simulated games using a playout strategy.
	- **Expert feedback** on how well a task was performed.

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.

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

 $[c]$ is an indicator function returning 1 if $c = True$ and otherwise 0

Training a Model

- Models are "trained" (learned) on **the training data.** This involved estimating:
	- **1. Model parameters** (the model): E.g., probabilities, weights, factors.
	- **2. Hyperparameters**: Many learning algorithms have choices for 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! This is a type of model selection.

Hyperparameter Tuning/Model Selection

- 1. Hold a validation data set back from the training data.
- **2. Learn models** using the training set with different \longleftarrow 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 with different hyperparameters, so we get generalization accuracy for comparing different hyperparameter settings.
	- If no model selection is necessary, then no validation set is used.

Testing a Model

- After the model is selected, the final model is evaluated against the test set to **estimate the final model accuracy** and see how well it generalizes.
- **Very important:** never contaminate your training set with test data or "peek" at the test set during training!

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 or other properties of the examples.
- **k-fold cross validation:** Use training & validation data better
	- Split the training & validation data randomly into k folds.
	- For each of k rounds, hold one 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.

Learning Curve: The Effect the Training Data Size

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

More data is better!

At some point the learning curve flattens out and more data does not contribute much!

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 published state-
of-the-art as a baseline.

Types of ML Models

Regression: Predict a number Classification: Predict a label

Regression: Linear Regression

Model:
\n
$$
h_w(x_j) = w_0 + w_1x_{j,1} + \dots + w_nx_{j,n} = \sum_i w_i x_{j,i} = w^T x_j
$$

\nEmpirical Loss:
\n $L(w) = ||Xw - y||^2$
\nGradient:
\n $\nabla L(w) = 2X^T (Xw - y)$
\nFind:
\n $\nabla L(w) = 0$
\nGradient descendent
\n $w = w - \alpha \nabla L(w)$
\n
\nAnalytical solution:
\n $w^* = \underbrace{(X^T X)^{-1} X^T y}_{\text{Pseudo inverse}$
\n $\underbrace{\begin{bmatrix} \frac{\partial L}{\partial w_1}(w), \frac{\partial L}{\partial w_2}(w), \dots, \frac{\partial L}{\partial w_n}(w) \end{bmatrix}}_{\text{House size in square feet}}$
\n $\underbrace{\begin{bmatrix} \frac{\partial L}{\partial w_1}(w), \frac{\partial L}{\partial w_2}(w), \dots, \frac{\partial L}{\partial w_n}(w) \end{bmatrix}}_{\text{House size in square feet}}$
\n $\underbrace{\begin{bmatrix} \frac{\partial L}{\partial w_1}(w), \frac{\partial L}{\partial w_2}(w), \dots, \frac{\partial L}{\partial w_n}(w) \end{bmatrix}}_{\text{Hence given value of the image.}}}$
\n $\underbrace{\begin{bmatrix} \frac{\partial L}{\partial w_1}(w), \frac{\partial L}{\partial w_2}(w), \dots, \frac{\partial L}{\partial w_n}(w) \end{bmatrix}}_{\text{Hence given value of the image.}}.$

Bayes Classifier $h^*(x) = \argmax P(Y = y | X = x)$ \mathcal{Y}

Naïve Bayes Classifier

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

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

and the $P(x_i | y)$ are estimated from the data

The $P(y)$ s and the $P(x_i | y)$ s are estimated from the data by counting.

• Gaussian Naïve Bayes Classifiers extend the approach to continuous features by assuming the feature follows a normal distribution depending on the class:

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

The parameters for the normal distribution $N(\mu_{\nu}, \sigma_{\nu})$ are estimated from data.

Decision Trees

- 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

$$
h(x) = \underset{y}{\text{argmax}} P(Y = y \mid \text{leafNodeMatching}(x))
$$

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

$$
h(x) = \operatorname*{argmax}_{y} P(Y = y \mid \text{neighborhood}(x))
$$

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** g (e.g., logistic, ReLU).
- Learn weights w from examples using **backpropagation** of prediction errors $L(\hat{y}, y)$ (gradient descend).
- ANNs are **universal approximators**. Large networks can approximate any function (no bias). **Regularization** is typically used to avoid overfitting.
- **Deep learning** adds more hidden layers and layer types (e.g., convolution layers) for better learning.

Other Popular Models and Methods

Many other models exist

• **Generalized linear model (GLM):** This important model family includes **linear regression** and the classification method **logistic regression.**

Often used methods

- **Regularization:** enforce simplicity and reduces overfitting 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).
- **Embedding and Dimensionality Reduction:** Learn how to represent data in a simpler way (e.g., PCA, text embeddings).

Some Use Cases of ML for Intelligent Agents

Learn Actions

• Directly learn the best action from examples.

 $action = h(state)$

• This model can also be used as a **playout policy** for Monte Carlo tree search with data from self-play.

Learn Heuristics

• Learn evaluation functions for states.

 $eval = h(state)$

• Can learn a **heuristic** for minimax search from examples.

Perception

- **Natural language processing:** Use deep learning / word embeddings / language models to understand concepts, translate between languages, or generate text.
- Speech recognition: Identify the most likely sequence of words.
- Vision: Object recognition in images/videos. Generate images/video.

Compressing Tables

- Neural networks can be used as a compact representation of tables that do not fit in memory. E.g.,
	- Joint and conditional probability tables
	- State utility tables
- The tables can be learned form data.

Bottom line: Learning a function is often more effective than hard-coding it
However, we do not always know how it performs in very rare cases!