Introduction to Machine Learning

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Artificial Intelligence

Designing machines that can perform tasks that are characteristic of human intelligence

Learning

Planning

Reasoning

Problem Solving

Language Understanding & Generation

Object Recognition

Speech Recognition & Generation

Strong AI vs Weak AI





Supervised Classification

Sepal Length **Sepal Width Petal Length Petal Width** Class 5.1 3.5 1.4 0.2 Iris-setosa 4.9 3 1.4 0.2 Iris-setosa 7 3.2 4.7 1.4 Iris-versicolor 6.3 3.3 6 2.5 Iris-virginica 5.8 5.1 1.9 Iris-virginica 27 Class Label

Identify the class of Iris plant:

Iris Setosa OR Iris Versicolour OR Iris Virginica

Given features predict class label

Features/Signals/Hints

	Sepal Length	Sepal Width	Petal Length	Petal Width	Class
raining Data	5.1	3.5	1.4	0.2	Iris-setosa
	4.9	3	1.4	0.2	Iris-setosa
	7.0	3.2	4.7	1.4	Iris-versicolor
	6.3	3.3	6.0	2.5	Iris-virginica
TRAINING	5.8	2.7	5.1	1.9	Iris-virginica
Test Data		- Model	Y Y = f(X)		† Y
	Sepal Length	Sepal Width	Petal Length	Petal Width	Class
	6.4	2.8	5.6	2.1	?
INFERENCE	7.2	3.0	5.8	1.6	?

What is the nature of the model f(X)?

We want f(X) that generalizes well to test data

Obviously, f(X) cannot be a lookup table that memorizes

How do we learn the model f(X)?



Decision Tree Learning Explained Simply



Probabilistic Models - Let's look at a simple classifier

Simplified Bayes Decision Rule

 $Y^* = \underset{Y \in Y}{\operatorname{argmax}} P(Y|X)$

We would like to estimate the conditional distribution: P(Y|X)



How do we find P(X|Y)?

Maximum Likelihood Estimation

$$P(D) = \prod_{i=1,M} P(X|Y)$$

We can assume P(X|Y) follows some distribution – say Gaussian

$$P(X|Y_{1}=setosa) = N(\mu_{1}, \sigma_{1})$$

$$P(X|Y_{2}=virginica) = N(\mu_{2}, \sigma_{2})$$

$$P(X|Y_{3}=versicolor) = N(\mu_{3}, \sigma_{3})$$

$$\Theta^{*} = \underset{\Theta}{\operatorname{argmax}} P(D)$$

$$Objective Function$$

Training, Validation and Test Sets

Training Set To learn model parameters

Validation/Development Set ->

- To select model e.g. what features to use, what distribution to use
- To choose special values \rightarrow the hyperparameters

Use the test set **only** to evaluate the results

Precision for class k = what fraction of the predictions were correct? Recall = What fraction of examples of class k were correctly identified F-1: A metric which combines precision and recall

Using test set for model selection will bias the model



- Objective function: Minimize the number of errors on the training set.
- Zero-one loss function, difficult to optimize ۲
- Approximate with: Minimize

 $\sum_{i=1}^{n} \max(0, -t_{i} * y_{i})$ where $t_{i} = w_{1}x_{1i} + w_{2}x_{2i}$



Which line to choose?



Choose the line which maximizes the margin → this is a more general solution This is the Support Vector Machine classifier Perceptron & Support Vector Machine are linear classifiers \rightarrow the decision boundary is a line

In many cases, data is spread out in a more complicated



So, we need non-linear classifiers



But is it desirable?

This could lead to OVERFITTING

Overfitting

- Model performs well on training data, but performs badly on test data
- The model could fit noise in training data
- The more complex the model, the more prone to overfitting
 - As a rule of thumb, more parameter \rightarrow more complex models
- Occam's Razor: When presented with competing <u>hypothetical</u> answers to a problem, one should select the answer that makes the fewest assumptions
 - Prefer simpler models
 - Penalize complexity of the model

How to prevent overfitting?

- Try simpler models first
- There are ways to prevent overfitting
- Regularization is a popular method:
 - Controls range of parameter values
 - Smaller range \rightarrow lesser variance in the learnt function
- Many model-specific methods exist
 - Tree pruning for decision trees
 - Drop-out for neural networks
- Feature Selection
- Domain knowledge

These represent model biases \rightarrow Learning is not possible without a bias

We saw one kind of machine learning problem: Supervised Scalar Classification

Supervised \rightarrow Expected output is known during training

Scalar \rightarrow Single Output

Classification \rightarrow Output is one of K possible outcomes (discrete)

But there are many other machine learning scenarios

Regression: Predict a real value instead of a discrete categorical value e.g. Price Prediction

Sequence Labelling: Input is a sequence, output is a sequence of the same length e.g. Part-of-Speech Tagging

Sequence-to-Sequence Learning: Input and Output are sequences of different lengths

e.g Machine Translation English: Water is necessary for life (5 words) Malayalam: വെള്ളം ജീവന് അത്യാവശ്യമാണ് (3 words) water life+for necessary+is

More complicated output structures: trees, graphs, etc

Reinforcement Learning: Weak supervision, just rewards and penalties on predictions e.g. game-playing, car driving

Unsupervised Learning: Given only data, no labels \rightarrow let's know a little bit more about this

Clustering



- Discover patterns in data, exploratory analysis
- Popular Algorithms: k-means, Expectation Maximization with Mixture Models, agglomerative hierarchical clustering

Lower Dimensional Representation



- Noise reduction, computational efficiency, avoiding overfitting, visualization
- Popular Algorithms: Principal Component Analysis, Singular Value Decomposition, Latent Dirichlet Allocation, Autoencoding

Density Estimation



- Novely and anamoly detection
- Generating synthetic datasets
- Algorithms: Parzen windows, Kernel Density estimation

Deep Learning

The return of artificial neural networks

Perceptrons \rightarrow The Simplest Network

 $sign(w_0+w_1x_1+w_2x_2+w_3x_3)$ Output wDendrite NucleusSynapses
Cell Body or Soma

Input \rightarrow Dendrites

X

Weighted sum of inputs

Outputs \rightarrow Axons

Output only if weighted sum greater than a threshold (w_0)

Add multiple perceptron units



This is still not sufficient

Add a non-linear activation

 $sigmoid(w_0 + w_1 x_1 + w_2 x_2 + w_3 x_3)$



A network with non-linear activations and at least one hidden layer can approximate any function

The network is made of small simple learning units

Why have deep neural networks become popular?

Deep Neural Networks → Many Hidden Layers

- Deep networks have millions of parameters \rightarrow prone to overfitting
- Lots of data mitigates overfitting to some extent
- Modern hardware makes it possible to train deep networks
- Better initialization, optimization and activation methods
- Very developer friendly language and good toolkits
- Representation Learning \rightarrow Features are automatically learnt



Layer 3

Parts combine to form objects

Layer 2

DNNs have achieved tremendous success in many perceptual task since as image recognition, vision, speech

Lesser success in language – to some extent in machine translation

Some Reading Material

- Google's Crash Course (Basic)
- Andrew Ng's Coursera Course (Intermediate)
- Andrew Moore's course material (Intermediate)
- Andrew Ng's Stanford Course (Advanced)
- Mathematics Pre-requisites for Advanced Learning
 - Probability and Statistics
 - Linear Algebra
 - Optimization

Some Tools and Software

- Weka Library (Java)
- Scikit Learn (Python)
- Deep learning frameworks like Tensorflow, Torch/PyTorch, MXNet

Data Sources

- UC Irvine Machine Learning Repository
- Kaggle
- .. Many other sources

Thank you!

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