CSCI 5521: Introduction to Machine Learning (Spring 2018)

Supervised Learning

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Supervised LearningClassificationRegression



Input Feature Space



Supervised Learning

Classification
Regression
Data: $\begin{aligned}
\mathcal{X} = \{\mathbf{x}^{t}, r^{t}\}_{t=1}^{N} & \mathcal{X} = \{\mathbf{x}^{t}, r^{t}\}_{t=1}^{N} \\
\mathcal{X} = \{\mathbf{x}^{t}, r^{t}\}_{t$

Classification

Data: $\chi = {\mathbf{x}^t, r^t}_{t=1}^N$ Output: $r = \begin{cases} 1 \text{ if } \mathbf{x} \text{ is positive} \\ 0/-1 \text{ if } \mathbf{x} \text{ is negative} \end{cases}$

X1	X2	r
0.934	0.046	-1
0.679	0.097	-1
0.758	0.823	1
0.743	0.695	-1
0.392	0.317	-1
0.655	0.950	1
0.171	0.034	-1
0.706	0.439	-1
0.032	0.382	1
0.277	0.766	1



Learning a Class from Examples

Class C of a "family car"

□ Prediction: Is car *x* a family car?

Knowledge extraction: What do people expect from a family car?

Output:

Positive (+) and negative (-) examples

Input representation:

 x_1 : price, x_2 : engine power



Class in a Rectangle



Hypothesis class $\mathcal H$

Consider \mathcal{H} : the set of all rectangles



Version Space



Linear Classifier



h(*x*)=<*w*,*x*>+b is a linear classifier

h(*x*)>0 positive *h*(*x*)<0 negative

 $h \in H, H$?

Perceptron Learning

- Perceptron algorithm, Rosenblatt, 1957.
- Initialization:

w = 0

 Iterate until converge (no mistake on a certain number of iterations)

> for each example (\mathbf{x}^{t}, r^{t}) : $if(\langle \mathbf{w}, \mathbf{x}^{t} \rangle * r^{t} \leq 0)$ $\mathbf{w} = \mathbf{w} + r^{t} \mathbf{x}^{t}$

Perceptron Learning



Best in the Version Space



Margin

Choose *h* with largest marginWhy?



Model Capacity

- Different models have different capacity meaning the ability to handle more complex data.
- How to measure model capacity?
- The maximum number of data points that can be classified perfectly in any labeling.

VC (Vapnik Chervonenkis) Dimension

- N points can be labeled in 2^N ways as +/-
- In a particular arrangement, \mathcal{H} shatters N if there exists $h \in \mathcal{H}$ consistent for any of the 2^N ways:

 $VC(\mathcal{H}) = N$

VC Dimension

How many points can be shattered by a line?



VC (Vapnik Chervonenkis) Dimension

How about axis-aligned rectangles?



VC Summary

- The capacity of function is measured by the number of data points that can be shattered by the function
- Rectangle classifier in 2-D space: 4.
- A line : 3.
- More ...

VC Dimension

- More generally, in R^D space, what is the VC of a hyperplane?
- What is the VC of a triangle classifier?
- Is an algorithm that can shatter only 4 or 3 data points useful?
- How easy it is to determine the VC dimension for the hypothesis class?

VC Dimension: Why Large Margin



Multiple Classes, C_i i=1,...,K



KNN Classification

K nearest neighbor





http://mirlab.org/jang/books/dcpr/prKnnc.asp?title=5-2%20K-nearest-neighbor%20Classifiers&language=english

How to Choose K for KNN?



- What is the VC dimension of KNN?
- Is VC proportional to the # of parameters (appeared complexity)?

Regression



$$E(w_1, w_0 \mid \mathcal{X}) = \frac{1}{N} \sum_{t=1}^{N} \left[r^t - (w_1 x^t + w_0) \right]^2$$

Regression

How does the error function look like?

$$E(w_1, w_0 \mid X) = \frac{1}{N} \sum_{t=1}^{N} \left[r^t - (w_1 x^t + w_0) \right]^2$$



Regression

Find the g to minimize training error

$$E(w_{1},w_{0} \mid \mathcal{X}) = \frac{1}{N} \sum_{t=1}^{N} \left[r^{t} - (w_{1}x^{t} + w_{0}) \right]^{2}$$
$$\frac{\partial E(w_{1},w_{0} \mid \mathcal{X})}{\partial w_{0}} = \frac{1}{N} \sum_{t=1}^{N} \left[(r^{t} - w_{1}x^{t} - w_{0})(-1) \right] = 0$$
$$\frac{\partial E(w_{1},w_{0} \mid \mathcal{X})}{\partial w_{1}} = \frac{1}{N} \sum_{t=1}^{N} \left[(r^{t} - w_{1}x^{t} - w_{0})(-x^{t}) \right] = 0$$

$$w_1 = \frac{\sum_{t} x^t r^t - N\overline{x}\overline{r}}{\sum_{t} (x^t)^2 - N\overline{x}^2}, w_0 = \overline{r} - w_1\overline{x}$$

Regression: Understand Solution

C

Average error is 0.

Property 2:
$$\frac{1}{N} \sum_{t=1}^{N} \left[(r^t - w_1 x^t - w_0)(-x^t) \right] = \sum_{t=1}^{N} \varepsilon^t x^t = 0$$

Error is uncorrelated with data

Polynomial Regression

Is polynomial fitting very different?

$$g(x) = \sum_{i=1}^{P} w_p(x)^p + w_0$$



It is the same as linear regression with a polynomial mapping.

$$g(x) = w^{T} x$$
$$w = [w_{P}, \dots, w_{1}, w_{0}]$$
$$x = [x^{P}, \dots, x^{1}, x^{0}]$$

Summary of Supervised Learning

1. Model: $g(\mathbf{x} \mid \theta)$ $g(\mathbf{x}) = w_1 \mathbf{x} + w_0$

2. Loss function: $E(\theta \mid X) = \sum_{t} L(r^{t}, g(\mathbf{x}^{t} \mid \theta))$

$$E(h \mid \mathcal{X}) = \sum_{t=1}^{N} \mathbb{1}\left(h(\mathbf{x}^{t}) \neq r^{t}\right) \qquad E(g \mid \mathcal{X}) = \frac{1}{N} \sum_{t=1}^{N} \left[r^{t} - g(x^{t})\right]^{2}$$

3. Optimization procedure: $\theta^* = \arg \min_{\theta} E(\theta \mid X)$

Algorithms: KNN, percepton, linear regression

Noise and Model Complexity

Data is not perfect

- Data recording might not be perfect (shifted data points)
- Wrong labeling of the data
- There might be additional unobervable hidden variables.



Noise and Model Complexity



Options:

- Simple model with training errors
- Complex comdel with no training error

Noise and Model Complexity

Given similar training error, use the simpler one

- Simpler to use (lower computational complexity)
- Easier to train (lower space complexity)
- Easier to explain (more interpretable)
- Generalizes better (lower variance Occam's razor)





Model Selection & Generalization

- Learning is an ill-posed problem; data is not sufficient to find a unique solution
- Given d binary inputs, there are at most 2^{D} samples, and $2^{2^{D}}$ binary functions
- Each sample eliminates half of the functions;
- Thus, N samples leaves $2^{2^{D}-N}$ viable functions
- Not possible to check all functions. Need for inductive bias, assumptions about $\mathcal H$

Generatlization and Overfitting

- Generalization: How well a model performs on new data
- Overfitting: *H* more complex than *C* or *f*Underfitting: *H* less complex than *C* or *f*



Cross-Validation

- To better estimate generalization error, we need data unseen during training. We split the data as
 - □ Training set (50%)
 - □ Validation set (25%)
 - Test set (25%)
- Resampling when there is few data

Cross-Validation



Triple Trade-Off

- There is a trade-off between three factors (Dietterich, 2003):
 - 1. Complexity of \mathcal{H} , $c(\mathcal{H})$,
 - 2. Training set size, N,
 - 3. Generalization error, *E*, on new data
- □ As *N*↑, *E*↓
- □ As $c(\mathcal{H})\uparrow$, first $E\downarrow$ and then $E\uparrow$

Triple Trade-Off



Complexity of Classifier