STAT 406: Algorithms for classification AND PREDICTION

LECTURE 1: INTRODUCTION

Kevin Murphy

Mon 8 January, 2007^1

- Administrivia
- Some basic definitions.
- Simple examples of regression.
- Real-world applications of regression.
- Simple examples of classification.
- Real-world applications of classification.

- Web page http://www.cs.ubc.ca/~murphyk/Teaching/Stat406/ Spring07/index.html Check the 'news' section before every class!
- Please fill out the sign-up sheet.
- Optional Lab Wed 4-5.
- The TA is Virginia Chen.
- My office hours are Fri 4-5pm LSK 308d. Please send me email ahead of time if you plan to show up!

- There will be weekly homework assignments worth 25%. Out on Mondays, return on Mondays (in class).
- The homeworks will involve theory and programming; you may want to do some of the programming during the lab.
- The midterm will be Mon Feb 26th (right after spring break) and is worth 30%.
- The final will be in April and is worth 40%.

- Math: multivariate calculus, linear algebra, probability theory.
- Stats: stats 306 or CS 340 or equivalent.
- CS: some experience with programming (eg in R) is required.

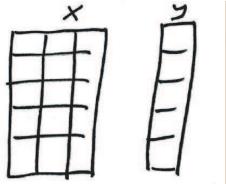
- There will be weekly programming assignments.
- We will use matlab.
- Matlab is very similar to R, but is somewhat faster and easier to learn. Matlab is widely used in the machine learning and Bayesian statistics community.
- Unfortunately matlab is not free (unlike R). You can buy a copy from the bookstore for \$150, or you can use the copy installed in the lab machines.
- You will learn how to use matlab during the first few labs.

- I am writing my own textbook, but it is not yet finished. You should buy a photocopy of the current draft (about 500 pages) at Copiesmart in the village (near MacDonalds) for about \$35 (available on Friday).
- Meanwhile, the other required textbook is *Pattern recognition and machine learning*, Chris Bishop, 2006
- The following books are recommended but not required
 - *Elements of statistical learning*, Hastie, Friedman and Tibshirani, 2001.
 - Pattern Classification, Duda, Hart, Stork, 2001 (2nd edition).
 - Statistical pattern recognition, Andrew Webb, 2002.

- Since people have different backgrounds (cs 340, stat 306, multiple versions), the exact syllabus may change as we go.
- See the web page for details.
- You will get a good feeling for the class during today's lecture.

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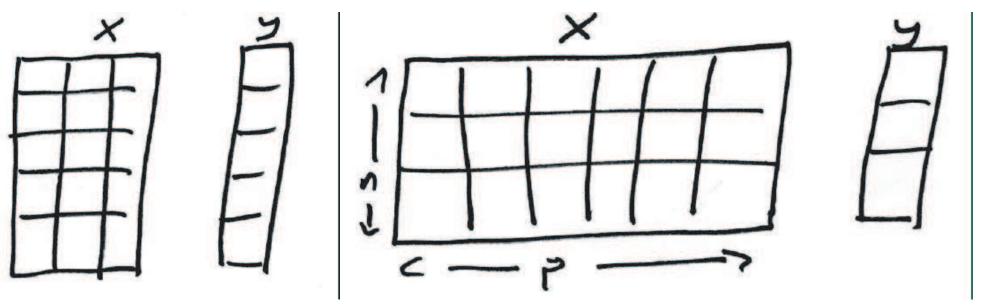
- This class is basically about machine learning.
- We will initially focus on supervised approaches.
- Given a training set of n input-output pairs $D = (\vec{x}_i, \vec{y}_i)_{i=1}^n$, we attempt to construct a function f which will accurately predict $f(\vec{x}_*)$ on future, test examples \vec{x}_* .
- Each input \vec{x}_i is a vector of d features or covariates. Each output \vec{y}_i is a target variable. The training data is stored in an $n \times d$ design matrix $X = [\vec{x}_i^T]$. The training outputs are stored in a $n \times q$ matrix $Y = [\vec{y}_i^T]$.



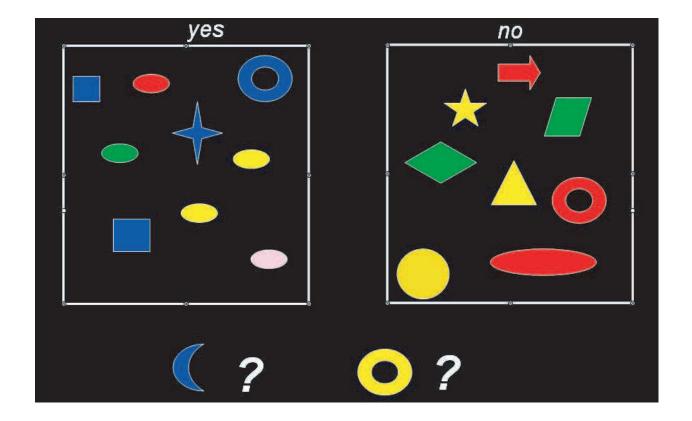
- If $\vec{y} \in \mathbb{R}^q$ is a continuous-valued output, this is called regression. Often we will assume q = 1, i.e., scalar output.
- If $y \in \{1, \ldots, C\}$ is a discrete label, this is called classification or pattern recognition. The labels can be ordered (eg. low, medium, high) or unordered (e.g., male, female). N_Y is the number of classes. If C = 2, this is called binary (dichotomous) classification.

- Bishop indexes training cases by n, eg x_n, whereas stats usually uses
 i. Bishop uses w for regression parameters, whereas stats uses β.
 Bishop uses d for number of dimensions (input features), whereas stats uses p. There are various other differences.
- My book uses different notation in different places (I am working on fixing this). Please read carefully and tell me of any errors.
- We will denote discrete ranges $\{1, \ldots, N\}$ by 1: N.

- In traditional applications, the design matrix is tall and skinny $(n \gg p)$, i.e., there are many more training examples than inputs.
- In more recent applications (eg. bio-informatics or text analysis), the design matrix is short and fat $(n \ll p)$, so we will need to perform feature selection and/or dimensionality reduction.



We care about performance on examples that are different from the training examples (so we can't just look up the answer).



- The *no free lunch theorem* says (roughly) that there is no single method that is better at predicting across all possible data sets than any other method.
- Different learning algorithms implicitly make different assumptions about the nature of the data, and if they work well, it is because the assumptions are reasonable in a particular domain.

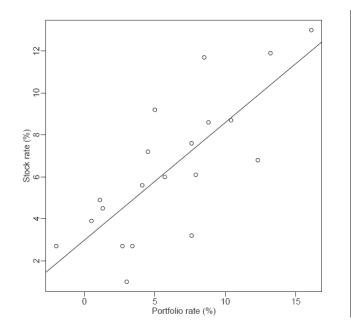
- In supervised learning, we are given (\vec{x}_i, \vec{y}_i) pairs and try to learn how to predict \vec{y}_* given \vec{x}_* .
- In unsupervised learning, we are just given $\vec{x_i}$ vectors.
- The goal in unsupervised learning is to learn a model that "explains" the data well. There are two main kinds:
 - Dimensionality reduction (eg PCA)
 - Clustering (eg K-means)

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The output density is a 1D Gaussian (Normal) conditional on x:

$$p(y|\vec{x}) = \mathcal{N}(y; \vec{\beta}^T \vec{x}, \sigma) = \mathcal{N}(y; \beta_0 + \beta_1 x_1 + \dots + \beta_p x_p, \sigma)$$
$$\mathcal{N}(y|\mu, \sigma) = \frac{1}{\sqrt{2\pi\sigma}} \exp(-\frac{1}{2\sigma^2} (y - \mu)^T (y - \mu))$$

For example, $y = ax_1 + b$ is represented as $\vec{x} = (1, x_1)$ and $\vec{\beta} = (b, a)$.

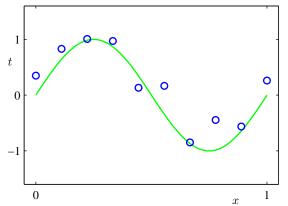


If we use linear regression with non-linear basis functions

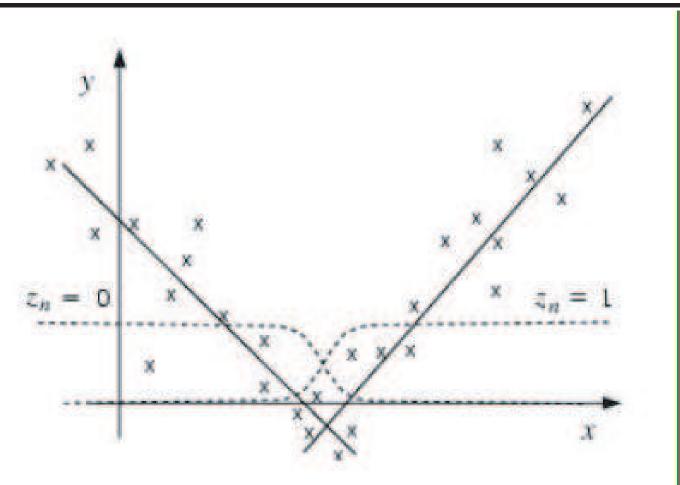
$$p(y|x_1) = \mathcal{N}(y|\beta^T [1, x_1, x_1^2, \dots, x_1^k], \sigma)$$

we can produce curves like the one below.

Note: In this class, we will often use \vec{w} instead of $\vec{\beta}$ to denote the weight vector.



PIECEWISE LINEAR REGRESSION



How many pieces? — Model selection problem. Where to put them? — Segmentation problem.

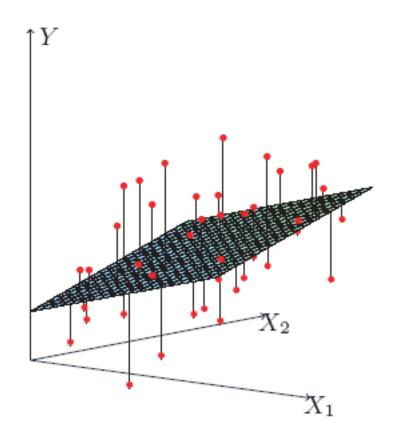
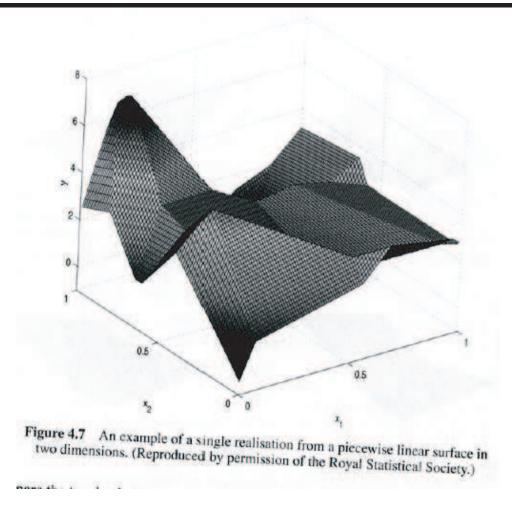


Figure 3.1: Linear least squares fitting with $X \in \mathbb{R}^2$. We seek the linear function of X that minimizes the sum of squared residuals from Y.

Piecewise linear 2D regression



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- \vec{x} = amount of various chemicals in my factory, y = amount of product produced.
- \vec{x} = properties of a house (eg location, size), y = sales price.
- $\vec{x} = \text{joint}$ angles of my robot arm, $\vec{y} = \text{location}$ of arm in 3-space.
- \vec{x} = stock prices today, \vec{y} = stock prices tomorrow. (Time series data is not iid, and is beyond the scope of this course.)

- A very interesting ordinal regression problem is to build a system that can predict what ranking (from 1 to 5) you would give to a new movie.
- The input might just be the name of the movie, plus your past voting patterns, and those of other users.
- The collaborative filtering approach says you will give the same score as those people who have similar movie tastes to you, which can you infer by looking at past voting patterns.
- For each movie and each user, you can infer a set of latent traits and use these to predict (related to SVD of a matrix).

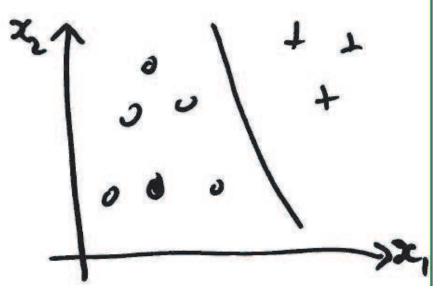
- The netflix prize http://netflixprize.com/ is an award of \$1M USD for a system that can predict your movie preferences 10% more accurately than their current system (called Cinematch).
- A large training data set is provided: a sparse $18k \times 480k$ matrix (movies \times users) containing about 100M rankings (on the scale 1:5) of various movies.
- The test (probe) set is 2.8M (movie,user) pairs, for which the ranking is known but withheld from the training set.
- The performance measure is root mean square error:

$$rmse = \sqrt{\frac{1}{n} \sum_{i=1}^{n} (R(u_i, m_i) - \hat{R}(u_i, m_i))^2}$$
(1)

where $R(u_i, m_i)$ is the true rating of user u_i on movie m_i , and $\hat{R}(u_i, m_i)$ is the prediction.

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2D inputs $\vec{x}_i \in \mathbb{R}^2$, binary outputs $y \in \{0, 1\}$. The line is called a *decision boundary*. Points to the right are classified as y = 1, points to the left as y = 0.



- A simple approach to binary classification is logistic regression (briefly studied in 306).
- The output density is Bernoulli conditional on x:

$$p(y|x) = \pi(x)^y \ (1 - \pi(x))^{1-y}$$

where $y \in \{0, 1\}$ and

$$\pi(x) = \sigma(\vec{w}^T \ [1, x_1, x_2])$$

where

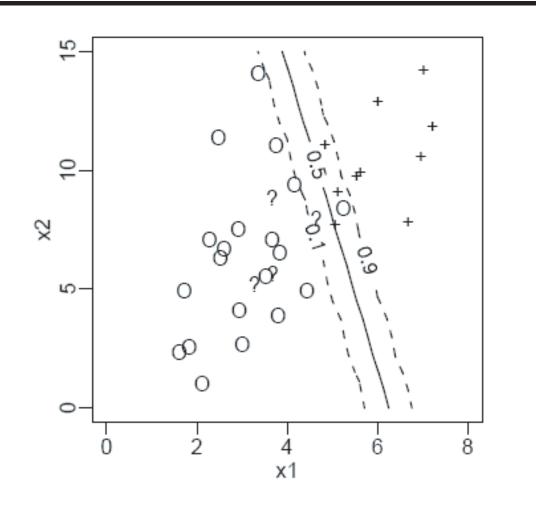
$$\sigma(u) = \frac{1}{1 + e^{-u}}$$

is the sigmoid (logistic) function that maps ${\rm I\!R}$ to [0,1]. Hence

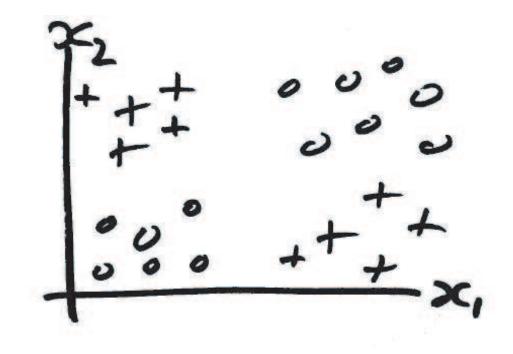
$$P(Y = 1 | \vec{x}) = \frac{1}{1 + e^{-w_0 + w_1 x_1 + w_2 x_2}}$$

where w_0 is the bias (offset) term corresponding to the dummy column of 1s added to the design matrix.

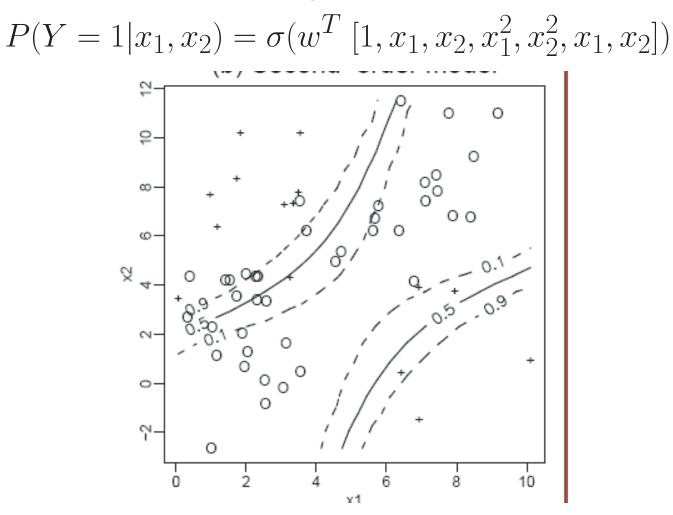
2D Logistic regression



In 306, this is called "checkerboard" data. In machine learning, this is called the "xor" problem. The "true" function is $y = x_1 \oplus x_2$. The decision boundary is non-linear.



We can separate the classes using



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Multi-class classification.

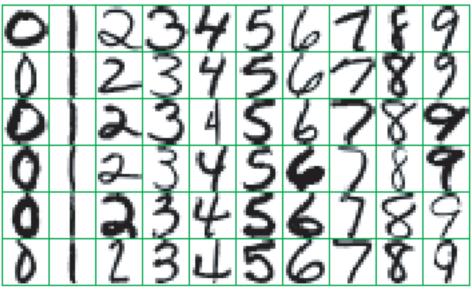
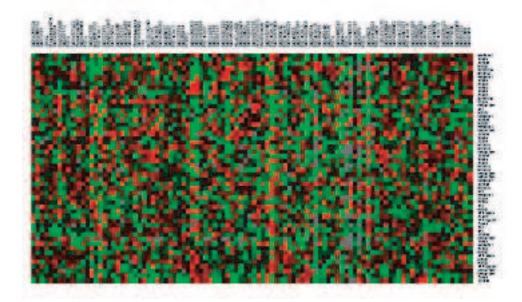


Figure 1.2: Examples of handwritten digits from U.S. postal envelopes.

Rows = examples, columns = features (genes). Short, fat data $(p \gg n)$. Might need to perform feature selection.



- Email spam filtering (spam vs not spam)
- Detecting credit card fraud (fraudulent or legitimate)
- Face detection in images (face or background)
- Web page classification (sports vs politics vs entertainment etc)
- Steering an autonomous car across the US (turn left, right, or go straight)