

*Over-Fitting*

*Overfitting, regularisation, feature selection*

## Over-fitting

- Complex machine can **over-fit**  
*over-fitting: fitting the training data well at the cost of getting poorer generalisation performance*
- Three red cars. . .
- If we used an infinitely flexible machine we can fit our training data perfectly, but would have no generalisation ability

## Binary Classification Task for You



Class 1



Class 2

## Which Category?

- Which category does the following image belong to?



## Spurious Rules

- You ask a learning machine to solve a task based on data
- It will find a rule that does this, but not necessary the rule you had in mind—machine learning isn't magic, it can't read your mind
- Infinitely flexible machines have an infinity of spurious rules they can learn—they are useless
- What should we do?

## All Binary Functions

$$x_0 = 000 \quad y_0 = \begin{cases} 0 \\ 1 \end{cases}$$

$$x_1 = 100 \quad y_1 = \begin{cases} 0 \\ 1 \end{cases} \quad \text{unseen}$$

$$x_2 = 010 \quad y_2 = \begin{cases} 0 \\ 1 \end{cases}$$

$$x_3 = 110 \quad y_3 = \begin{cases} 0 \\ 1 \end{cases}$$

$$x_4 = 001 \quad y_4 = \begin{cases} 0 \\ 1 \end{cases} \quad \text{seen}$$

$$x_5 = 101 \quad y_5 = \begin{cases} 0 \\ 1 \end{cases}$$

$$x_6 = 011 \quad y_6 = \begin{cases} 0 \\ 1 \end{cases}$$

$$x_7 = 111 \quad y_7 = \begin{cases} 0 \\ 1 \end{cases} \quad \text{unseen}$$

$$\mathcal{D} = \{(000, 0), (010, 1), (110, 1), (001, 0), (101, 0)\}$$

## Are MLPs Universal Approximators?

- Yes and No
- Yes: If you give me any function, I can find an MLP that approximates that function to any desired accuracy
- No: If you give me an MLP, I can find a function with an arbitrary high approximation error
- Would an MLP that could approximate any function be useful?
- Absolutely not!

1. Over-fitting?
2. **Controlling Complexity**
3. Hidden structure
4. Regularisation



## Training Examples

- As we increase the number of training examples, we make it hard to find a spurious rule
- Bigger data sets allow us to use more complicated machines
- Part of the success of deep learning is because they use huge training sets—but this is only a part of their success
- (Labelled) data is often expensive to collect so we sometimes have no choice but to use a small training set
- One of the limitations of using deep learning comes because we often have limited data

## Preprocessing

- Structure might often be obscure to the learning machine
- If we are trying to predict the spread of disease then a list of place names might be a lot less useful than their coordinates
- Imposing an ordering on an unordered set might **not** be useful

$\{\text{"blue"} : 0, \text{"brown"} : 1, \text{"green"} : 2, \text{"black"} : 3\}$

- Choosing an encoding that reflect meaningful structure is essential to successful machine learning

## Outline

1. Over-fitting?
2. Controlling Complexity
3. **Hidden structure**
4. Regularisation



- Infinitely flexible machine don't generalise (because any unseen data could have any value)
- **Machine learning only works because there is some structure in the data**
- A successful machine should capture this structure
- Even deep learning machines with millions of parameters only work because they successfully capture the structure of images or text
- Different learning machines have different performance on different problems because the data has different structure

## Identifying Structure

- In some cases we know *a priori* some of the structure in the data
- In images we believe the identity of an object is invariant to translation and scaling
- The success of *convolutional neural networks* (CNNs) in deep learning is in large part because the convolutions respect translational invariance

## Automatic Preprocessing

- One view of deep learning is that each layer (particularly in CNNs) acts as a preprocessor
- That is, it finds filters that captures features salient to the problem being tackled
- For both images and texts we expect salient features to be spatially localised (CNN finds localised filter)
- The deep structure allows ever more complicated features to be captured—that is, we can find spatially localised features on different scales
- Having very large datasets and simple filters (the number of weights in the CNN layers tends to be small) stops overfitting

## Hidden Structure

- Often the structure of data is invisible to us
- A very successful strategy is to try many different machine learning techniques and choose the best (stupid but effective)
- Often learning machines have adjustable parameters (hyper-parameters) that we have to set (they are the same for all input data, but change with the problem)
- We need to choose the hyper-parameters to fit the data in our problem
- Fine tuning hyper-parameter is important and almost always required (true in SVMs, MLP, deep learning)

## Measuring Generalisation Performance

- Recall, we want to predict **unseen** data
- You cannot use data that you have trained on!**—you will overfit
- Need to split your data up into training and validation set
- Use the validation set to choose the hyper-parameters
- You need a separate testing set if you want to measure your generalisation performance

## Cross Validation

- If you want to use more data for training then you can use cross validation
- $K$ -fold cross validation splits the data into  $K$  groups

$$\mathcal{D} = \{D_i\}_{i=1}^K, \quad D_i = (\mathbf{x}_i, \mathbf{y}_i)$$

TRYING TO FIND THE BEST SET OF HYPERPARAMETERS USING ONLY THE TRAINING SET

Leave-one-out cross-validation

$$\frac{1}{K} \sum_{i=1}^K \frac{1}{n-1} \sum_{j \neq i} (y_j - \hat{y}_j)^2 = 4.3$$

- Leave-one-out cross-validation is extreme case

## Dimensionality Reduction

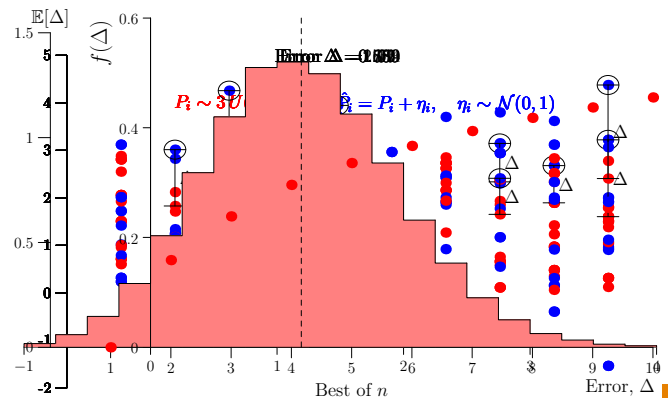
- We can sometimes simplify our machines by using less features
- We can project our data onto a lower dimensional sub-space (e.g. one with the maximum variation in the data: PCA)
- We can use clustering to find exemplars and recode our data in terms of distances from the exemplars (radial basis functions)
- Whether this helps depends on whether the information we discard is pertinent to the task we are trying to perform

## Normalising Features

- Measuring a feature in millimeters or kilometers is going to make a lot of difference to the size of that feature
- Many learning algorithms are sensitive to the size of a feature (larger features are more important)
- If we don't know how important different features are then it makes sense to normalise our features. E.g.

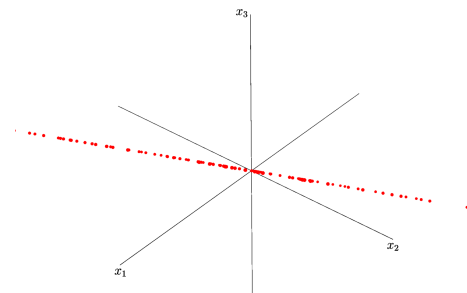
$$x_i^\alpha \leftarrow \frac{x_i^\alpha - \hat{\mu}_i}{\hat{\sigma}_i}, \quad \hat{\mu} = \frac{1}{m} \sum_{\beta=1}^m x_i^\beta, \quad \hat{\sigma}_i^2 = \frac{1}{m-1} \sum_{\beta=1}^m (x_i^\beta - \hat{\mu}_i)^2$$

## The Overfitting Game



## Hidden Structure

Can't spot low dimensional data by looking at numbers



## Feature Selection

- Spurious features will allow us to find spurious rules (**over-fitting**)
- We can try different combinations of features to find the best set, although it rapidly becomes intractable to do this in all ways
- We can use various heuristics to decide which features to keep, but no heuristic is fail-safe method to find the best set of features
- Feature selection however can be powerful, often we can get very good results by keeping only a few variables
- As well as possibly improving generalisation we also get a more **interpretable** rule

## Outline

- Over-fitting?
- Controlling Complexity
- Hidden structure
- Regularisation



## Explicit Regularisation

- As you've seen in the foundations of ML course, we can modify our error function to choose smoother functions

$$L = \sum_{k=1}^m (w^T x_k - y_k)^2 + \nu \|w\|^2$$

(Good to normalise features)

- Second term is minimised when  $w_i = 0$
- If  $w_i$  is large then

$$f(x|w) = w^T x = \sum_{i=1}^p w_i x_i$$

varies rapidly as we change  $x_i$

## Implicit Regularisation

- In the last two examples we added an explicit regularisation term that made the function we learnt simpler
- Some learning machines do this less explicitly
- Some deep learning architectures do subtle averaging
- Sometimes the architecture biases the machine to find a simple solution

## Success of SVMs

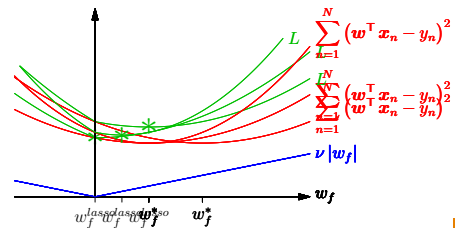
- SVMs regularise themselves by choosing the machine with the largest margin
- This ensures maximum stability to noise on the data
- It leads to very good generalisation on small datasets—usually beats everything else
- But you still need to normalise the features
- You also need to tune its hyper-parameters ( $C$  and sometimes  $\gamma$ )

## Lasso

- We can use other regularisers

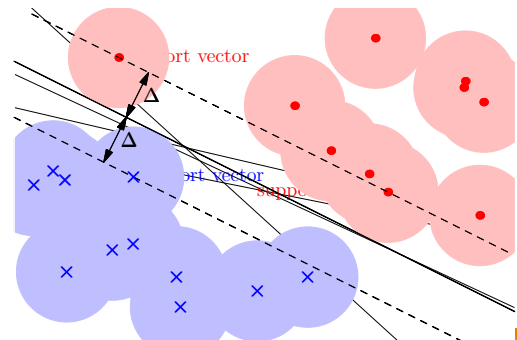
$$L = \sum_{k=1}^m (w^T x_k - y_k)^2 + \nu \sum_{i=1}^p |w_i|$$

- Spurious features (e.g. colour of flag on energy consumption) will give us a small improvement in training error



## Maximum Margin Machines

- Perceptrons have many options to separate data



- SVMs choose the machine with the biggest margins

## Lessons

- Machine learning isn't magic
- It works when the learning machine is well attuned to the problem
- Sometimes you can help by preprocessing your data
- Sometimes there is a regularisation term that helps select a simpler machine
- Most machines have hyper-parameter that you tune to match the machine to the data
- Really clever machines try to do this matching automatically