Machine Learning Lecture 5 Support Vector Machines

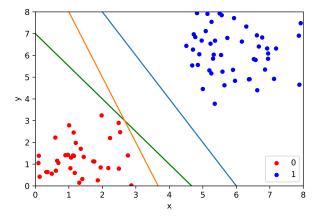
Justin Pearson¹

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¹http://user.it.uu.se/~justin/Teaching/MachineLearning/index.html

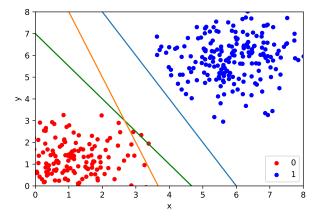
Separating Hyperplanes

Logistic regression (with linear features) finds a hyperplane that separates two classes. But which hyperplane is best?



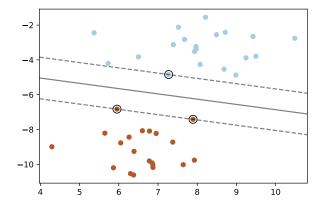
Separating Hyperplanes

It of course depends on how representative your training set is. With more points from the distribution our hyperplanes might look like:



Margin Classifiers

The intuition is that we find a hyperplane with a margin either side that maximises the space between the two clusters.



- They have been in use since the 90s.
- More robust with outliers.
- Very good classifiers on certain problems such as image classification, handwritten digit classification.
- Non-linear models can be incorporated by the kernel trick.

- A motivation/modification of logistic regression.
- Finding margins as an optimisation problem.
- Different Kernels for non-linear classification.

Remember the error term for logistic regression

$$-y \log(\sigma(h_{\theta}(x))) - (1-y) \log(1 - \sigma(h_{\theta}(x)))$$

Where

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

$$\begin{split} &-y \log(\sigma(h_{\theta}(x))) - (1-y) \log(1-\sigma(h_{\theta}(x))) \\ &\mathsf{Equal} \\ &-y \log(\frac{1}{1+e^{h_{\theta}(x)}}) - (1-y) \log(1-\frac{1}{1+e^{h_{\theta}(x)}})) \end{split}$$

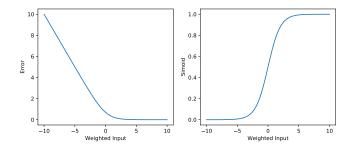
Remember that the two log terms are trying to force the model to learn 1 or 0.

Looking at the contribution

Just looking at

 $-y \log(\sigma(h_{\theta}(x)))$

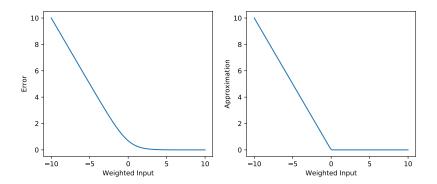
We are trying to force the term $\sigma(h_{\theta}(x))$ to be 1. The larger the value of θx the less the error.



After 0 we do not really care we just want to force move the input over to the right

Approximating the error

Instead of using the logistic error we could approximate it with two linear functions.



After 0 we do not really care about the error.

I am sorry, but to make the maths easier and to be consistent with the support vector machine literature but we are going to change our classification labels a bit.

We have data

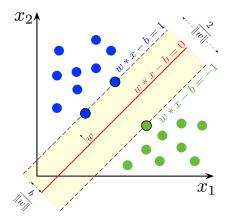
$$x = x^{(1)}, \ldots, x^{(m)}$$

Where the data are points in some d dimensional space \mathbb{R}^d . The labels for our classes will be -1 and 1 instead of 0 and 1. Linear SVMs are the easiest case and form the foundation for support vector machines. We want to find weights $\overline{w} \in \mathbb{R}^d$ and a constant *b* such that

$$\begin{cases} \overline{w} \cdot \overline{x} - b \ge 1 & \text{if } y = 1 \\ \overline{w} \cdot \overline{x} - b \le -1 & \text{if } y = -1 \end{cases}$$

This is different from logistic regression of a single perceptron where you want to find a separating hyperplane.

SVM Margins in 2 dimensions²



If we push the two hyperplanes apart then we will eventually hit points in the two classes. The points that are on the two out hyperplanes are called the support vectors. So the question is, how do we do this?

²Picture taken from wikipedia

Done in detail on the black board.

- The vector \overline{w} is perpendicular to the hyperplanes. In particular the hyperplane $\overline{w}x b = 0$.
- Given a two points $\overline{x_1}$ where $\overline{w} \cdot \overline{x_1} b = -1$ and $x_2 \ \overline{w} \cdot \overline{x_2} b = 1$.
- We want to know the distance between the two points. We can treat them as vectors and do the maths.

Done in detail on the black board.

- $\overline{x_2} \overline{x_1}$ is a vector, it has some length t and is in the direction \overline{w} . So $\overline{x_2} \overline{x_1} = t\overline{w}$.
- Now doing some rearranging

$$\overline{w} \cdot \overline{x_2} - b = \overline{w} \cdot (\overline{x_1} + t\overline{w}) - b = (\overline{w} \cdot \overline{x_1} - b) + t\overline{w} \cdot \overline{w} = 1$$

Note that $\overline{w} \cdot \overline{w}$ is the length squared, $||w||^2$ of w.

• $\overline{w} \cdot \overline{x_1} - b$ equals 1 we get that $t = 2/||w||^2$

•
$$||\overline{x_2} - \overline{x_1}|| = t||w|| = 2/||w||.$$

Thus to maximise the distance between the two hyperplanes $\overline{w}x - b = 1$ and $\overline{w}x - b = -1$ we want to maximise

$$t = \frac{2}{||\overline{w}||}$$

So we need to minimise $\frac{1}{2}||\overline{w}||$

Given a training set $x^{(1)}, \ldots, x^{(m)}$ We want to minimise $\frac{1}{2}||\overline{w}||$ such that for all data in the training set

$$\begin{cases} \overline{w} \cdot \overline{x^{(i)}} - b \ge 1 & \text{if } y^{(i)} = 1 \\ \overline{w} \cdot \overline{x^{(i)}} - b \le -1 & \text{if } y^{(i)} = -1 \end{cases}$$

Since $y^{(i)}$ can only be -1 or +1 we can rewrite the constraint as

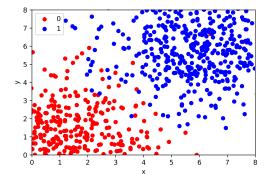
$$y^{(i)}(\overline{w}\cdot\overline{x^{(i)}}-b)\geq 1$$

and instead minimize

$$\frac{1}{2}\overline{w}\cdot\overline{w} = \frac{1}{2}||\overline{w}||^2$$

- Gradient descent will not work.
- Your optimisation problem includes lots of quadratic terms, but luckily the problem is convex.
- Quadratic programming solves this, and in the convex cases there are nice mathematical properties that give you bounds on the errors. How this all works is out of scope of the course.

What happens if our clusters overlap? The quadratic programming model will not work so well.



For each point in the training set introduce a slack variable η_i and rewrite the optimisation problem as

• Minimise $\frac{1}{2}||\overline{w}|| + C \sum_i \eta_i$ such that

$$y^{(i)}(\overline{w} \cdot \overline{x^{(i)}} - b) \ge 1 - \eta_i$$

An η_i greater than 0 allows the point to be miss-classified. Minimising $C \sum_i \eta_i$ for some constant C reduces the number of miss-classifications. The greater the constant C the more importance you give reducing the number of miss-classifications.

Warning what follows in the slides is not a complete description of what is going on with Kernels.

In particular I am not going to explain how the learning algorithm works. To understand this, you need to know a bit about quadratic programming, Lagrange multipliers, dual bounds and functional analysis.

I am instead going to try to give you some intuition why kernels might work. This if often called the kernel trick. I will also try to give you some intuition how and what SVMs learn with Gaussian Kernels.

We already saw that with Linear regression we could learn non-linear functions. To learn a quadratic polynomial we take out data

$$\left(\begin{pmatrix} 1 \\ x \end{pmatrix} \in \mathbb{R}^2
ight) \mapsto \left(\begin{pmatrix} 1 \\ x \\ x^2 \end{pmatrix} \in \mathbb{R}^3
ight)$$

One way of thinking about this is that we add invent new features. When computing the gradients everything worked.

We turned a non-linear problem of trying to find a quadratic polynomial that minimises the error into a linear problem of trying to learn the coefficients.

This is part of a general scheme.

We have a non-linear separation problem in low dimensions, we find a transformation that embeds our problem into a high-dimensional space. One possibly misleading way of thinking about this, is that the more dimensions you have the more room you have, and so it is easier for the problem to be linear.

$\mathbb{R}^d \stackrel{\Phi}{\longrightarrow} \mathbb{H}$

Where \mathbb{H} is some higher³ dimensional space.

³Actually $\mathbb H$ stands for Hilbert not higher, but do not worry about this.

A linear hypothesis $h_{\overline{w}}(\overline{x})$ has the form

$$h_{\overline{w}}(\overline{x}) = \sum_{i=1}^{d} w_i x_i + w_0 = \overline{w} \cdot \overline{x} + w_0$$

Where \cdot is the inner (dot) products.

In our learning algorithms there are a lot of inner product calculations.

So to learn linear things in ${\mathbb H}$ we will need to do inner products.

$\mathbb{R}^{d} \stackrel{\Phi}{\longrightarrow} \mathbb{H}$

We will need do lots of calculations on $\Phi(x_i) \cdot \Phi(x_j) \in \mathbb{H}$.

Computing the inner products $\Phi(x_i) \cdot \Phi(x_j) \in \mathbb{H}$ can be computationally expensive (even worse your space could be infinite dimensional⁴. For well behaved transformations Φ there exists a function $K(x, y) : \mathbb{R}^d \times \mathbb{R}^d \to \mathbb{R}$ such that

$$K(x,y) = \Phi(x) \cdot \Phi(y)$$

Thus we can compute the inner product in the high-dimensional space by using a function on the lower dimensional vectors.

⁴Don't worry if your head hurts.

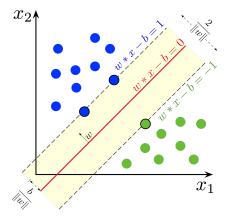
Instead of giving the higher-dimensional space you often just get the function K.

Radial basis or Gaussian $K(x, y) = \exp(-(x - y)^2/2\sigma^2)$ Polynomial $K(x, y) = (1 + x \cdot y)^d$

Sigmoid or Neural Network $K(x, y) = tanh(\kappa_1 x \cdot y + \kappa_2)$

There are lots more, there are even kernels for text processing. If you are going to invent your own then you will need to understand the maths.

Support Vectors



I am not going to explain in any detail, but it is enough to remember the support vectors.

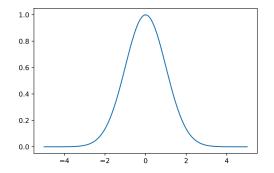
Learning with Kernels does the same thing, you find support vectors. The dual version of the algorithm learns some parameters α_i , y_i and b such that to decide if a point x belongs to a class you compute the sign of

$$\sum_{i=1}^{N_s} \alpha_i y_i K(s_i, x) + b$$

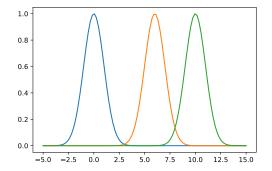
Where s_1, \ldots, s_{N_s} are the support vectors.

Gaussian Kernels

This is a all a bit abstract. I'll try to explain what is going on with Gaussian Kernels. In one dimension for $\sigma = 1$ our Gaussian kernel $K(x, x') = \exp(-(x - x')^2/2)$ if we fix x' to be 0. We get a the following graph

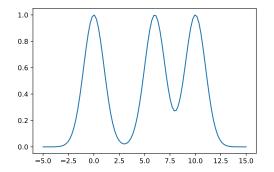


Gaussian Kernels - Multiple Support Vectors



Gaussian Kernels - Multiple Support Vectors

If we assume that all out weights are 1 and add them together. We get



The closer you value is to one of the peaks the more likely it is that you are in the class. You can think of each peak as a feature.

With non-linear Kernels SVMs have successfully been used in many applications including

- Image recognition, bio-informatics, pattern recognition.
- Because the optimisation problem is convex there is only going to be one global minimum. So SVMs are easier to train than neural networks.
- Probably the most useful non-linear Kernel function is the Gaussian Kernel.
- For the moment, don't worry too much about how the SVM learns with non-linear Kernels use an existing implementation.