COMP9417, 23T1

Kernel Methods

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- Transformations
- The Kernel Trick
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Section 1

Kernel Methods

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Instead of pure parameter based learning (i.e minimising a loss function etc.), dual algorithms introduce **instance-based** learning.

This is where we 'remember' mistakes in our data and adjust the corresponding weights accordingly.

We then use a *similarity function* or **kernel** in our predictions to weight the influence of the training data on the prediction.

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$$\mathbf{w} \in \mathbb{R}^p$$

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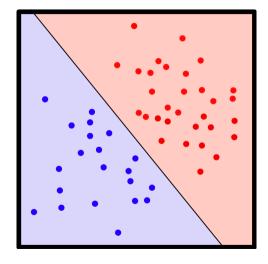
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 α_i represents the *importance* of a data point (x_i,y_i) .

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Recall the *primal* perceptron:

```
converged \leftarrow 0
while not converged do
    converged \leftarrow 1
    for x_i \in X, y_i \in y do
         if y_i w \cdot x_i < 0 then
             w \leftarrow w + \eta y_i x_i
             converged \leftarrow 0
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If we define the number of iterations the perceptron makes as $K \in \mathbb{N}^+$ and assume $\eta=1.$ We can derive an expression for the final weight vector $w^{(K)}$:

The Dual/Kernel Perceptron

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$$= \sum_{i=1}^{N} \alpha_i y_i x_i$$

where α_i is the number of times the perceptron makes a mistake on a data point (x_i, y_i) .

If we sub in $w^{(K)} = \sum_{i=1}^N \alpha_i y_i x_i$. We get the algorithm for the **dual** perceptron.

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              \alpha_i \leftarrow \alpha_i + 1
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Gram Matrix

The Gram matrix represents the *inner product* of two vectors.

For a dataset X we define $G = X^T X$. That is:

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$$G = \begin{bmatrix} \langle x_1, x_1 \rangle & \langle x_1, x_2 \rangle & \cdots & \langle x_1, x_n \rangle \\ \langle x_2, x_1 \rangle & \langle x_2, x_2 \rangle & \cdots & \langle x_2, x_n \rangle \\ \vdots & \vdots & \ddots & \vdots \\ \langle x_n, x_1 \rangle & \langle x_n, x_2 \rangle & \cdots & \langle x_n, x_n \rangle \end{bmatrix}$$

$$G_{i,j} = \langle x_i, x_j \rangle$$

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Section 3

Transformations

Transformations

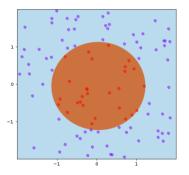
How do we go about solving non-linearly separable datasets with linear classifiers?

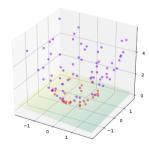
Transformations

How do we go about solving **non-linearly separable** datasets with linear classifiers? Project them to higher dimensional spaces through a transformation $\phi: \mathbb{R}^p \to \mathbb{R}^k$.

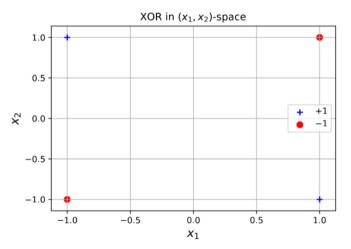
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Let's revisit the XOR.



A solution:

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For our input vectors in the form $\mathbf{x} = [x_1, x_2]^T$, use a transformation:

$$\phi(\mathbf{x}) = \begin{bmatrix} 1\\ \sqrt{2}x_1\\ \sqrt{2}x_2\\ x_1^2\\ x_2^2\\ \sqrt{2}x_1x_2 \end{bmatrix}$$

For our dataset,

$$\phi\left(\begin{bmatrix}1\\1\end{bmatrix}\right) = \begin{bmatrix}1\\\sqrt{2}\\\sqrt{2}\\1\\1\end{bmatrix} \phi\left(\begin{bmatrix}-1\\-1\end{bmatrix}\right) = \begin{bmatrix}1\\-\sqrt{2}\\-\sqrt{2}\\1\\1\end{bmatrix}$$

$$\phi\left(\begin{bmatrix} -1\\1 \end{bmatrix}\right) = \begin{vmatrix} -\sqrt{2}\\\sqrt{2}\\1\\1\\-\sqrt{2}\end{vmatrix}$$

$$\phi\left(\begin{bmatrix}1\\-1\end{bmatrix}\right) = \begin{bmatrix} \sqrt{2}\\-\sqrt{2}\\1\\1\end{bmatrix}$$

For the negative class:

$$\phi\left(\begin{bmatrix}1\\1\end{bmatrix}\right)_{2,6} = \begin{bmatrix}\sqrt{2}\\\sqrt{2}\end{bmatrix}$$

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For the positive class:

$$\phi\left(\begin{bmatrix} -1\\1 \end{bmatrix}\right)_{2,6} = \begin{bmatrix} -\sqrt{2}\\-\sqrt{2} \end{bmatrix}$$
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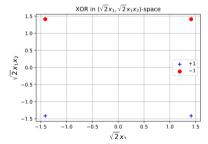
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We may have a problem, recall the dual perceptron.

```
converged \leftarrow 0
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               converged \leftarrow 0
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```

Recall the transformation $\phi: \mathbb{R}^p \to \mathbb{R}^k$.

$$G = \begin{bmatrix} \langle \phi(x_1), \phi(x_1) \rangle & \langle \phi(x_1), \phi(x_2) \rangle & \cdots & \langle \phi(x_1), x_n \rangle \\ \langle \phi(x_2), \phi(x_1) \rangle & \langle \phi(x_2), \phi(x_2) \rangle & \cdots & \langle \phi(x_2), \phi(x_n) \rangle \\ \vdots & \vdots & \ddots & \vdots \\ \langle \phi(x_n), \phi(x_1) \rangle & \langle \phi(x_n), \phi(x_2) \rangle & \cdots & \langle \phi(x_n), \phi(x_n) \rangle \end{bmatrix}$$

the Gram matrix becomes far too complex to compute.

Section 4

The Kernel Trick

The Kernel Trick

An absolute mathemagical idea which allows us to calculate the values of the Gram matrix for cheap.

Recall the transformation to the XOR data:

$$\phi(\mathbf{x}) = \begin{bmatrix} 1\\ \sqrt{2}x_1\\ \sqrt{2}x_2\\ x_1^2\\ x_2^2\\ \sqrt{2}x_1x_2 \end{bmatrix}$$

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An absolute mathemagical idea which allows us to calculate the values of the Gram matrix for cheap.

Recall the transformation to the XOR data:

$$\phi(\mathbf{x}) \cdot \phi(\mathbf{y}) = \begin{bmatrix} 1\\ \sqrt{2}x_1\\ \sqrt{2}x_2\\ x_1^2\\ x_2^2\\ \sqrt{2}x_1x_2 \end{bmatrix} \begin{bmatrix} 1\\ \sqrt{2}y_1\\ \sqrt{2}y_2\\ y_1^2\\ y_2^2\\ \sqrt{2}y_1y_2 \end{bmatrix}$$

$$\phi(\mathbf{x}) \cdot \phi(\mathbf{y}) = 1 + 2x_1y_1 + 2x_2y_2 + x_1^2y_1^2 + x_2^2y_2^2 + 2x_1x_2y_1y_2$$

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$$\phi(\mathbf{x}) \cdot \phi(\mathbf{y}) = 1 + 2(x_1y_1 + x_2y_2) + (x_1y_1 + x_2y_2)^2$$

Kernel Methods

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Kernel Methods

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Kernel Methods

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Kernel Methods

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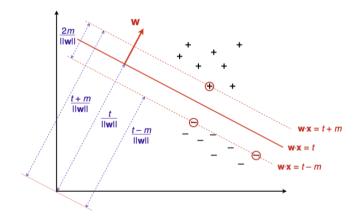
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Why is this useful?

Section 5

Support Vector Machines

Support Vector Machines



The basic SVM is a linear classifier defined by:

$$\underset{w,t}{\operatorname{arg\,min}} \frac{1}{2} \|w\|^2$$
 subject to $y_i(\langle x_i, w \rangle - t) \geq m$

where t is the line's intercept, and we a consider a margin m. Typically, we'll see m=1 for a standardised dataset.

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This formulation means that we find the maximal margin classifier for the dataset.

Aside: Lagrangian Dual Problem

Say we have a problem as follows:

$$\max_{x,y} xy$$

subject to
$$x + y = 4$$

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Kernel Methods

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We can set up the Lagrangian dual and *move* the constraint into the function itself:

$$\Lambda(x, y, \lambda) = xy + \lambda(x + y - 4)$$

To solve this, we can calculate $\frac{\partial L}{\partial x}$, $\frac{\partial L}{\partial y}$ and $\frac{\partial L}{\partial \lambda}$ and solve the remaining system of equations.

The General Form of a Dual Problem

If we have a problem:

$$\underset{x}{\arg\min} f(x)$$
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The general dual problem is:

$$\Lambda(\mathbf{x}, \lambda) = f(\mathbf{x}) + \sum_{i=1}^{n} \lambda_i g_i(x_i)$$

If we take the general SVM problem (m = 1):

$$\operatorname*{arg\,min}_{w,t} \frac{1}{2} \|w\|^2$$

subject to
$$y_i(\langle x_i, w \rangle - t) \ge 1$$

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From the general form, we can take the vector α to form the dual problem:

$$\Lambda(w, t, \alpha) = \frac{1}{2} ||w||^2 + \left(-\sum_{i=1}^n \alpha_i y_i (\langle x_i, w \rangle - t) - 1) \right)$$

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$$\Lambda(w, t, \alpha) = \frac{1}{2} \|w\|^2 - \sum_{i=1}^n \alpha_i y_i (w \cdot x_i) + t \sum_{i=1}^n \alpha_i y_i + \sum_{i=1$$

Kernel Methods

Let's try and optimise the Lagrangian Λ w.r.t w.

$$\Lambda(w, t, \alpha) = \frac{1}{2} ||w||^2 - w \cdot \sum_{i=1}^{n} \alpha_i y_i x_i + t \sum_{i=1}^{n} \alpha_i y_i + \sum_{i=1}^{n} \alpha_i$$

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$$\frac{\partial \Lambda}{\partial w} = \frac{1}{2} 2w - \sum_{i=1}^{n} \alpha_i y_i x_i$$

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We can see that at $\frac{\partial \Lambda}{\partial w} = 0$

$$w = \sum_{i=1}^{n} \alpha_i y_i x_i$$

Repeating a similar process for t,

$$\Lambda(w, t, \alpha) = \frac{1}{2} \|w\|^2 - w \cdot \sum_{i=1}^{n} \alpha_i y_i x_i + t \sum_{i=1}^{n} \alpha_i y_i + \sum_{i=1}^{n} \alpha_i$$

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$$\sum_{i=1}^{n} \alpha_i y_i = 0$$

We've derived that for an optimal solution, $\sum_{i=1}^{n} \alpha_i y_i = 0$ and $w = \sum_{i=1}^{n} \alpha_i y_i x_i$

$$\Lambda(w, t, \alpha) = \frac{1}{2} ||w||^2 - w \cdot \sum_{i=1}^{n} \alpha_i y_i x_i + t \sum_{i=1}^{n} \alpha_i y_i + \sum_{i=1}^{n} \alpha_i$$

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$$\Lambda(w, \alpha) = \frac{1}{2} w^T w - w^T w + \sum_{i=1}^n \alpha_i$$

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We've derived that for an optimal solution, $\sum_{i=1}^{n} \alpha_i y_i = 0$ and $w = \sum_{i=1}^{n} \alpha_i y_i x_i$

$$\Lambda(w,t,\alpha) = \frac{1}{2} \|w\|^2 - w \cdot \sum_{i=1}^n \alpha_i y_i x_i + t \sum_{i=1}^n \alpha_i y_i + \sum_{i=1}^n \alpha_i$$

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$$\Lambda(\alpha) = -\frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_j y_i y_j (x_i \cdot x_j) + \sum_{i=1}^n \alpha_i$$

Our final problem now has relaxed constraints:

$$\begin{split} \Lambda(\alpha) &= -\frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_j y_i y_j (x_i \cdot x_j) + \sum_{i=1}^n \alpha_i \\ \text{subject to } \sum_{i=1}^n \alpha_i y_i &= 0 \\ \alpha_i &\geq 0 \text{ for } i=1,\dots,n \end{split}$$

Soft Margin SVM

Our current model is a **maximum** or hard margin classifier. To allow for errors *within* the supporting hyperplanes, we can redefine the primal problem as:

$$\operatorname*{arg\,min}_{w,t,\xi} \frac{1}{2} \|w\|^2 + C \sum_{i=1}^n \xi_i \qquad \text{subject to } y_i(\langle x_i,w\rangle - t) \geq 1 - \xi_i \text{ where } \xi_i \geq 0$$

we typically take ξ_i as the *hinge loss* of a point.

A Slight Extension: Hinge Loss

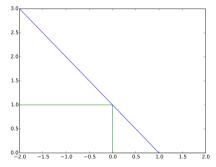
We define hinge loss for a data point at (x_i, y_i) as:

$$\xi_i = \max(0, 1 - y_i(w^T x_i - b))$$

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Kernel Methods

$$\underset{w,t,\xi}{\arg\min} \frac{1}{2} ||w||^2 + C \max(0, 1 - y_i(w^T x_i - b))$$

subject to
$$y_i(\langle x_i, w \rangle - t) \ge 1 - \max(0, 1 - y_i(w^T x_i - b))$$

Section 6

Extension: The RBF Kernel

Extension: The RBF Kernel

A popular Kernel is the Radial Basis Function kernel, defined below:

$$K(x,y) = \exp\left(-\frac{\|x - y\|^2}{2\sigma^2}\right)$$

for scalar values:

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What does this represent? A projection to infinite dimensions!