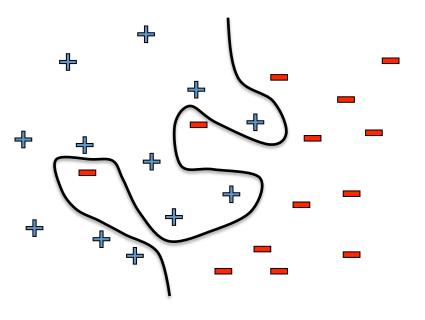
Kernels and Kernelized Perceptron

Instructor: Alan Ritter

Many Slides from Carlos Guestrin and Luke Zettlemoyer

What if the data is not linearly separable?



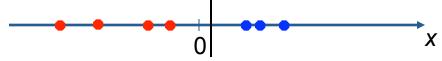
Use features of features of features of features....

$$\phi(x) = \begin{pmatrix} x_1 \\ \dots \\ x_n \\ x_1 x_2 \\ x_1 x_3 \\ \dots \\ e_{x_1} \\ \dots \end{pmatrix}$$

Feature space can get really large really quickly!

Non-linear features: 1D input

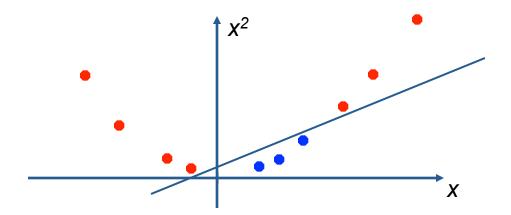
 Datasets that are linearly separable with some noise work out great:



But what are we going to do if the dataset is just too hard?



How about... mapping data to a higher-dimensional space:

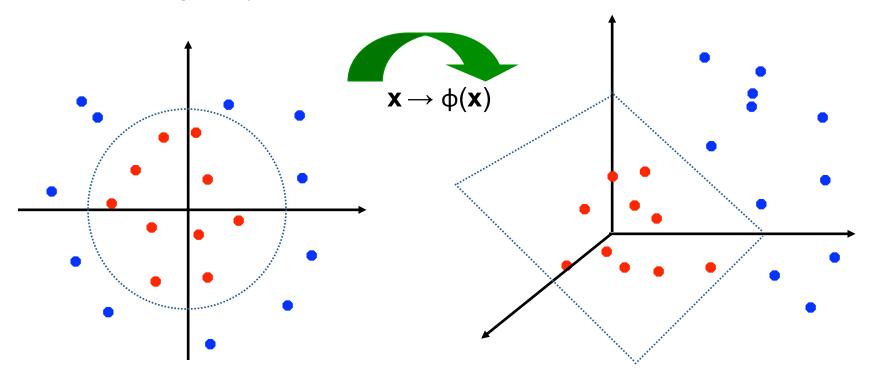


Feature spaces

- General idea: map to higher dimensional space
 - if \mathbf{x} is in \mathbb{R}^n , then $\phi(\mathbf{x})$ is in \mathbb{R}^m for m > n
 - Can now learn feature weights w in R^m and predict:

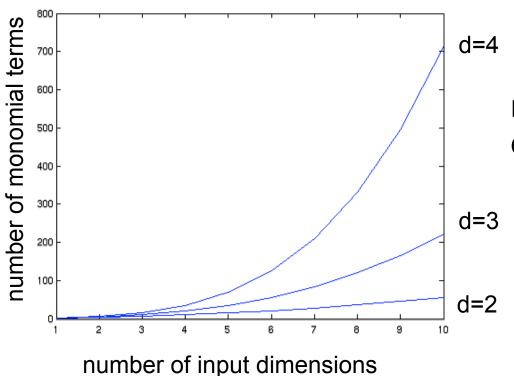
$$y = sign(\mathbf{w} \cdot \phi(\mathbf{x}))$$

 Linear function in the higher dimensional space will be non-linear in the original space



Higher order polynomials

num. terms
$$= \begin{pmatrix} d+m-1 \\ d \end{pmatrix} = \frac{(d+m-1)!}{d!(m-1)!}$$



m – input featuresd – degree of polynomial

grows fast! d = 6, m = 100 about 1.6 billion terms

Efficient dot-product of polynomials

Polynomials of degree exactly d

$$d=1$$

$$\phi(u).\phi(v) = \begin{pmatrix} u_1 \\ u_2 \end{pmatrix} \cdot \begin{pmatrix} v_1 \\ v_2 \end{pmatrix} = u_1v_1 + u_2v_2 = u.v$$

$$d=2$$

$$\phi(u).\phi(v) = \begin{pmatrix} u_1^2 \\ u_1u_2 \\ u_2u_1 \\ u_2^2 \end{pmatrix} \cdot \begin{pmatrix} v_1^2 \\ v_1v_2 \\ v_2v_1 \\ v_2^2 \end{pmatrix} = u_1^2v_1^2 + 2u_1v_1u_2v_2 + u_2^2v_2^2$$

$$= (u_1v_1 + u_2v_2)^2$$

$$= (u.v)^2$$

For any d (we will skip proof):

$$K(u,v) = \phi(u).\phi(v) = (u.v)^d$$

 Cool! Taking a dot product and an exponential gives same results as mapping into high dimensional space and then taking dot product

The "Kernel Trick"

A kernel function defines a dot product in some feature space.

$$K(\mathbf{u},\mathbf{v}) = \mathbf{\Phi}(\mathbf{u}) \cdot \mathbf{\Phi}(\mathbf{v})$$

Example:

2-dimensional vectors $\mathbf{u} = [u_1 \ u_2]$ and $\mathbf{v} = [v_1 \ v_2]$; let $K(\mathbf{u}, \mathbf{v}) = (1 + \mathbf{u} \cdot \mathbf{v})^2$, Need to show that $K(\mathbf{x}_i, \mathbf{x}_j) = \mathbf{\phi}(\mathbf{x}_i) \cdot \mathbf{\phi}(\mathbf{x}_j)$:

$$K(\mathbf{u},\mathbf{v}) = (1 + \mathbf{u} \cdot \mathbf{v})^{2}_{,} = 1 + u_{1}^{2} v_{1}^{2} + 2 u_{1} v_{1} u_{2} v_{2} + u_{2}^{2} v_{2}^{2} + 2 u_{1} v_{1} + 2 u_{2} v_{2} =$$

$$= [1, u_{1}^{2}, \sqrt{2} u_{1} u_{2}, u_{2}^{2}, \sqrt{2} u_{1}, \sqrt{2} u_{2}] \cdot [1, v_{1}^{2}, \sqrt{2} v_{1} v_{2}, v_{2}^{2}, \sqrt{2} v_{1}, \sqrt{2} v_{2}] =$$

$$= \mathbf{\phi}(\mathbf{u}) \cdot \mathbf{\phi}(\mathbf{v}), \text{ where } \mathbf{\phi}(\mathbf{x}) = [1, x_{1}^{2}, \sqrt{2} x_{1} x_{2}, x_{2}^{2}, \sqrt{2} x_{1}, \sqrt{2} x_{2}]$$

- Thus, a kernel function *implicitly* maps data to a high-dimensional space (without the need to compute each $\phi(x)$ explicitly).
- But, it isn't obvious yet how we will incorporate it into actual learning algorithms...

The "Kernel Trick"

A kernel function defines a dot product in some feature space.

Statistics Professors HATE Him!



Doctor's discovery revealed the secret to learning any problem with just 10 training samples. Watch this shocking video and learn how rapidly you can find a solution to your learning problems using this one sneaky kernel trick! Free from overfitting!

http://www.oneweirdkerneltrick.com

 But, it isn't obvious yet how we will incorporate it into actual learning algorithms...

"Kernel trick" for The Perceptron!

- Never compute features explicitly!!!
 - Compute dot products in closed form $K(u,v) = \Phi(u) \cdot \Phi(v)$
- Standard Perceptron:
 - set w_i=0 for each feature i
 - set aⁱ=0 for each example i
 - For t=1..T, i=1..n:

$$-y = sign(w \cdot \phi(x^{i}))$$

- At all times during learning:

$$w = \sum_{k} a^{k} \phi(x^{k})$$

Kernelized Perceptron:

• $a^{i} += y^{i}$

- set aⁱ=0 for each example i
 - For t=1..T, i=1..n: $-y = sign((\sum_k a^k \phi(x^k)) \cdot \phi(x^i))$ $= sign(\sum_k a^k K(x^k, x^i))$ $\text{ if } y \neq y^i$

Exactly the same computations, but can use K(u,v) to avoid enumerating the features!!!

- set aⁱ=0 for each example i
- For t=1..T, i=1..n:

$$- y = sign(\sum a^k K(x^k, x^i))$$

- if $y \neq y^i$ k
 - aⁱ += yⁱ

X ₁	x ₂	У
1	1	1
-1	1	-1
-1	-1	1
1	-1	-1

$$K(u,v) = (u \cdot v)^2$$

e.g.,
 $K(x^1,x^2)$
= $K([1,1],[-1,1])$
= $(1x-1+1x1)^2$
= 0

K	x ¹	x ²	x ³	x ⁴
x^1	4	0	4	0
x ²	0	4	0	4
x ³	4	0	4	0
x ⁴	0	4	0	4

Initial:

- $a = [a^1, a^2, a^3, a^4] = [0,0,0,0]$ t=1,i=1
- $\Sigma_k a^k K(x^k, x^1) = 0x4 + 0x0 + 0x4 + 0x0 = 0$, sign(0)=-1
- $a^1 += y^1 \rightarrow a^1 += 1$, new a = [1,0,0,0]t=1,i=2
- $\Sigma_k a^k K(x^k, x^2) = 1x0+0x4+0x0+0x4 = 0$, sign(0)=-1 t=1,i=3
- $\Sigma_k a^k K(x^k, x^3) = 1x4 + 0x0 + 0x4 + 0x0 = 4$, sign(4)=1 t=1,i=4
- $\Sigma_k a^k K(x^k, x^4) = 1x0+0x4+0x0+0x4 = 0$, sign(0)=-1 t=2,i=1
- $\Sigma_k a^k K(x^k, x^1) = 1x4 + 0x0 + 0x4 + 0x0 = 4$, sign(4)=1

•••

Converged!!!

• $y=\Sigma_k a^k K(x^k,x)$ = $1\times K(x^1,x)+0\times K(x^2,x)+0\times K(x^3,x)+0\times K(x^4,x)$ = $K(x^1,x)$ = K([1,1],x) (because $x^1=[1,1]$) = $(x_1+x_2)^2$ (because $K(u,v)=(u \cdot v)^2$)

Common kernels

Polynomials of degree exactly d

$$K(\mathbf{u}, \mathbf{v}) = (\mathbf{u} \cdot \mathbf{v})^d$$

Polynomials of degree up to d

$$K(\mathbf{u}, \mathbf{v}) = (\mathbf{u} \cdot \mathbf{v} + 1)^d$$

Gaussian kernels

$$K(\mathbf{u}, \mathbf{v}) = \exp\left(-\frac{||\mathbf{u} - \mathbf{v}||}{2\sigma^2}\right)$$

Sigmoid

$$K(\mathbf{u}, \mathbf{v}) = \tanh(\eta \mathbf{u} \cdot \mathbf{v} + \nu)$$

And many others: very active area of research!

Overfitting?

- Huge feature space with kernels, what about overfitting???
 - Often robust to overfitting, e.g. if you don't make too many Perceptron updates
 - SVMs (which we will see next) will have a clearer story for avoiding overfitting
 - But everything overfits sometimes!!!
 - Can control by:
 - Choosing a better Kernel
 - Varying parameters of the Kernel (width of Gaussian, etc.)

Kernels in logistic regression

$$P(Y = 0 | \mathbf{X} = \mathbf{x}, \mathbf{w}, w_0) = \frac{1}{1 + exp(w_0 + \mathbf{w} \cdot \mathbf{x})}$$

Define weights in terms of data points:

$$\mathbf{w} = \sum_{j} \alpha^{j} \phi(\mathbf{x}^{j})$$

$$P(Y = 0 | \mathbf{X} = \mathbf{x}, \mathbf{w}, w_{0}) = \frac{1}{1 + exp(w_{0} + \sum_{j} \alpha^{j} \phi(\mathbf{x}^{j}) \cdot \phi(\mathbf{x}))}$$

$$= \frac{1}{1 + exp(w_{0} + \sum_{j} \alpha^{j} K(\mathbf{x}^{j}, \mathbf{x}))}$$

- Derive gradient descent rule on α^{j} , w_{0}
- Similar tricks for all linear models: SVMs, etc

What you need to know

- The kernel trick
- Derive polynomial kernel
- Common kernels
- Kernelized perceptron