Linear Classifiers/SVMs

David Kauchak, CS311, Spring 2013

Admin

- Midterm exam posted
- Assignment 4 due Friday by 6pm
- No office hours tomorrow

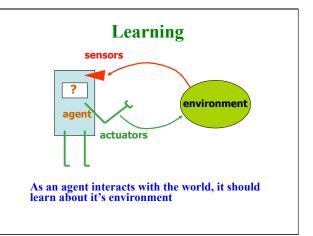
Math

Machine learning often involves a lot of math

- some aspects of AI also involve some familiarity

Don't let this be daunting

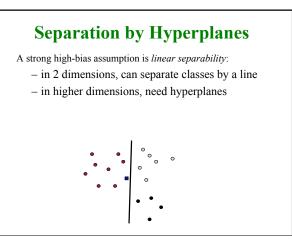
- Many of you have taken more math than me
- Gets better over time
- Often, just have to not be intimidated





Three classifiers

- Naïve Bayes
- k-nearest neighbor
- decision tree
- good and bad?
- Bias vs. variance
 - a measure of the **model**
 - where do NB, k-nn and decision trees fit on the bias/
 - variance spectrum?



Hyperplanes

A hyperplane is line/plane in a high dimensional space



What defines a hyperplane? What defines a line?

Hyperplanes

A hyperplane in an n-dimensional space is defined by n+1 values

$$0 = w_1 f_1 + w_2 f_2 + \dots + w_n f_n + w_{n+1}$$

e.g. a line

$$0 = w_1 f_1 + w_2 f_2 + w_3$$
 f(x) = **ax**+**b**

or a plane

$$0 = w_1 f_1 + w_2 f_2 + w_3 f_3 + w_4$$
 f(x,y) = **ax+by** + **c**

NB as a linear classifier

To classify:

 $\operatorname{arg\,max}_{C} P(C \mid f_1, f_2, \dots, f_n)$

Another way to view this (for 2 classes):

$$d(f_1, f_2, \dots, f_n) = \frac{P(c_1 \mid f_1, f_2, \dots, f_n)}{P(c_2 \mid f_1, f_2, \dots, f_n)}$$

Given *d* how would we classify?

NB as a linear classifier

$$d(f_1, f_2, \dots, f_n) = \frac{P(c_1 \mid f_1, f_2, \dots, f_n)}{P(c_2 \mid f_1, f_2, \dots, f_n)}$$

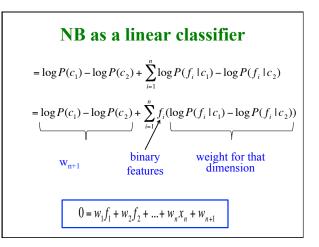
To classify:

$$classify(f_1, f_2, \dots, f_n) = \begin{cases} c_1 \text{ if } d > 1 \\ c_2 \text{ if } d < 1 \end{cases}$$

We can take the log:

$$classify(f_1, f_2, ..., f_n) = \begin{cases} c_1 \text{ if } \log d > 0 \\ c_2 \text{ if } \log d < 0 \end{cases}$$

NB as a linear classifier $\log d(f_1, f_2, ..., f_n) = \log \frac{P(c_1 | f_1, f_2, ..., f_n)}{P(c_2 | f_1, f_2, ..., f_n)}$ $= \log \frac{P(f_1 | c_1) P(f_2 | c_1) ... P(f_n | c_1) p(c_1)}{P(f_1 | c_2) P(f_2 | c_2) ... P(f_n | c_2) p(c_2)}$ $= \log P(c_1) - \log P(c_2) + \sum_{i=1}^n \log P(f_i | c_1) - \log P(f_i | c_2)$



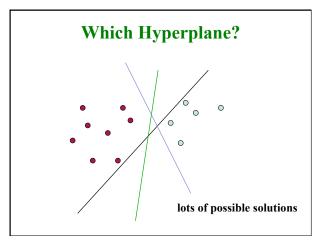
Lots of linear classifiers

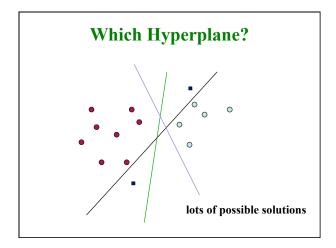
Many common text classifiers are linear classifiers

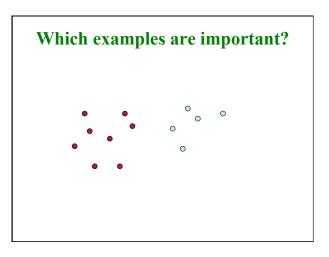
- Naïve Bayes
- Perceptron
- Rocchio
- Logistic regression
- Support vector machines (with linear kernel)
- Linear regression

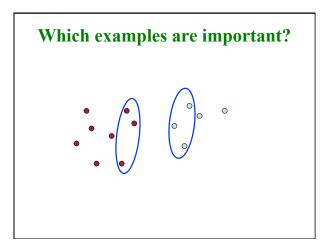
Despite this similarity, noticeable performance difference

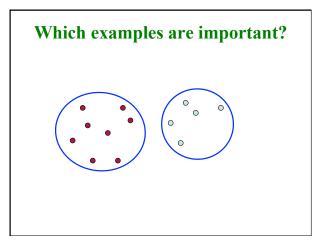
How might algorithms differ?

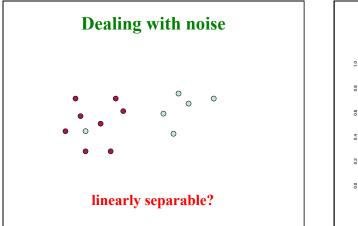


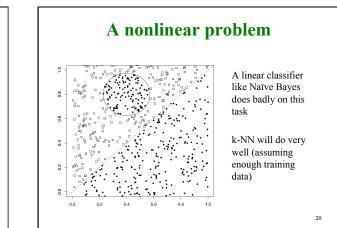


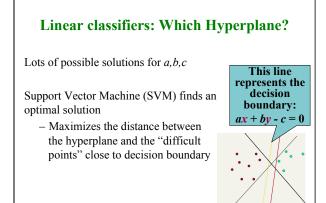


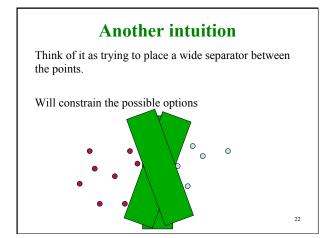








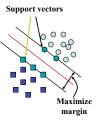




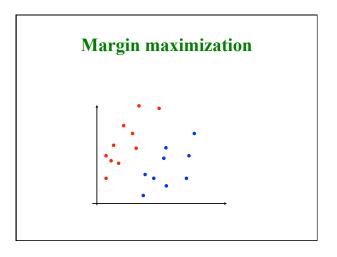
Support Vector Machine (SVM) Support vectors SVMs maximize the margin around the separating hyperplane C · aka large margin classifiers specified by a subset of training samples, the support vectors Maximize Posed as a quadratic programming

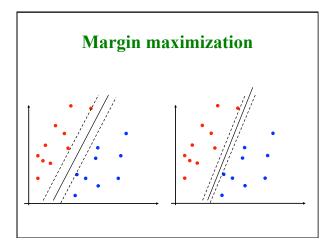
problem

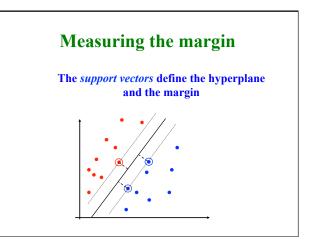
Seen by many as the most successful current text classification method*

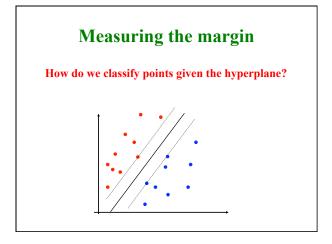


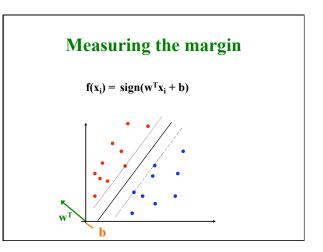
*but other discriminative methods often perform very similarly

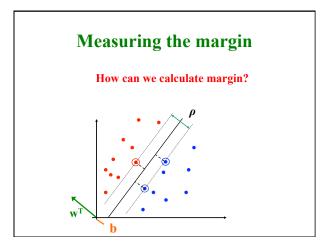


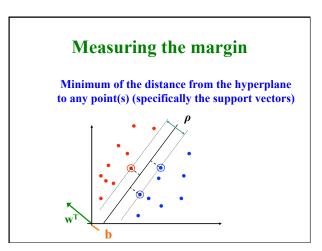


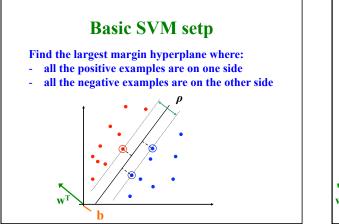


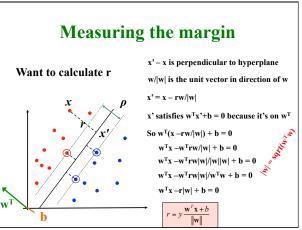




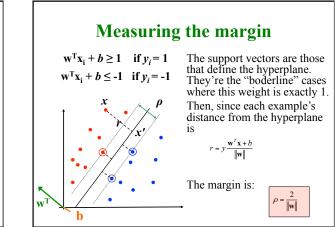








Linear SVM Mathematically The linearly separable case		
Assume that all data is at least distance 1 from the hyperplane, then the following two constraints follow for a training set $\{(\mathbf{x}_i, y_i)\}$		
$\mathbf{w}^{\mathrm{T}}\mathbf{x}_{i} + b \ge 1$	if $y_i = 1$	positive examples on one side
$\mathbf{w}^{\mathrm{T}}\mathbf{x}_{\mathrm{i}} + \boldsymbol{b} \leq -1$	if $y_i = -1$	negative examples on the other side

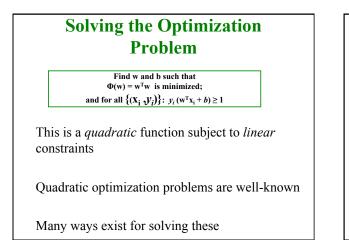


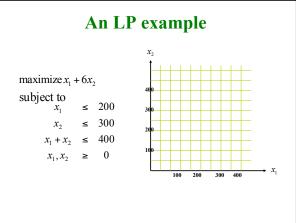
Linear SVMs Mathematically
(cont.)Then we can formulate the quadratic optimization
problem:Find w and b such that
is maximized; $\rho = \frac{2}{\|\mathbf{w}\|}$ maximize margin
is maximized;for all $\{(\mathbf{x}_i, y_i)\}$ $\mathbf{w}^T \mathbf{x}_i + b \ge 1$ if $y_i = 1$; $\mathbf{w}^T \mathbf{x}_i + b \ge 1$ if $y_i = -1$ make sure points are on correct size

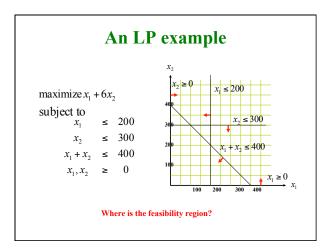
Linear SVMs Mathematically (cont.)

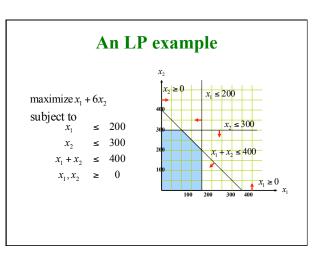
A better formulation (min $||\mathbf{w}|| = \max 1/||\mathbf{w}||$):

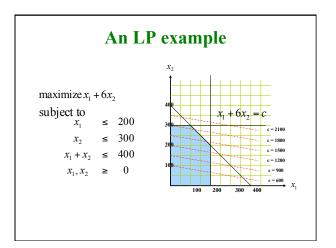
Find w and b such that $\Phi(w) = w^{T}w$ is minimized; and for all $\{(x_i, y_i)\}: y_i(w^{T}x_i + b) \ge 1$

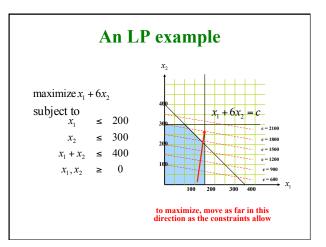


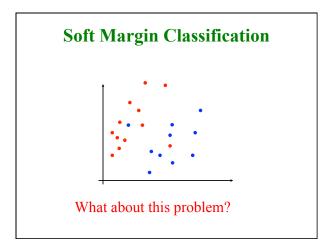


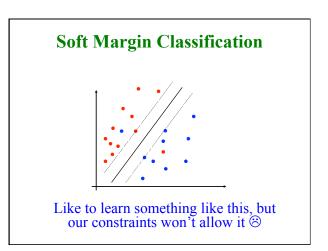


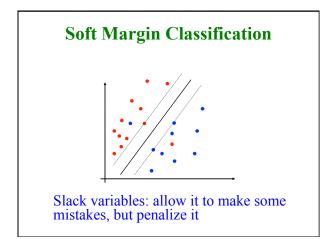














Old:

Find w and b such that $\begin{aligned} \Phi(w) = & y_2 w^T w \text{ is minimized and for all } \{(\mathbf{X}_i, y_i)\} \\ & y_i (w^T \mathbf{x}_i + \mathbf{b}) \ge 1 \end{aligned}$

With slack variables:

Find w and b such that $\Phi(w) = \frac{1}{2} w^{T}w + \frac{C \sum_{\xi_i}}{i}$ is minimized and for all $\{(\mathbf{X}_i, \mathbf{y}_i)\}$ $y_i (w^{T}\mathbf{X}_i + b) \ge 1 - \frac{1}{\xi_i}$ and $\xi_i \ge 0$ for all i

- allows us to make a mistake, but penalizes it
- C trades off noisiness vs. error

Linear SVMs: Summary

Classifier is a separating hyperplane

 large margin classifier: learn a hyperplane that maximally separates the examples

Most "important" training points are support vectors; they define the hyperplane

Quadratic optimization algorithm

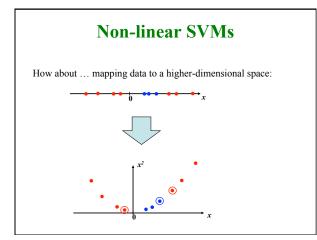
Non-linear SVMs

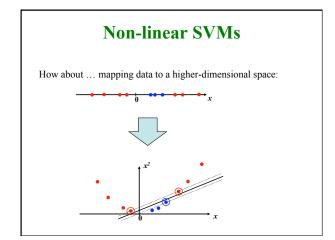
Datasets that are linearly separable (with some noise) work out great:

0

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

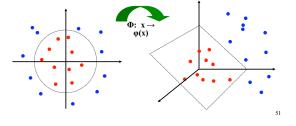
⇒x





Non-linear SVMs: Feature spaces

General idea: map original feature space to higherdimensional feature space where the training set is separable:



The "Kernel Trick"

The linear classifier relies on an inner product between vectors $K(\mathbf{x}_i, \mathbf{x}_i) = \mathbf{x}_i^T \mathbf{x}_i$

If every datapoint is mapped into high-dimensional space via some transformation Φ : $\mathbf{x} \rightarrow \phi(\mathbf{x})$, the inner product becomes:

$$K(\mathbf{x}_i, \mathbf{x}_j) = \varphi(\mathbf{x}_i)^{\mathrm{T}} \varphi(\mathbf{x}_j)$$

A *kernel function* is some function that corresponds to an inner product in some expanded feature space.

Kernels

Why use kernels?

- Make non-separable problem separable.
- Map data into better representational space

Common kernels

- Linear
- Polynomial $K(x,z) = (1+x^Tz)^d$
- Gives feature conjunctions
- Radial basis function (infinite dimensional space)

 $K(\mathbf{x}_i, \mathbf{x}_j) = e^{-\|\mathbf{X}_i - \mathbf{X}_j\|^2 / 2\sigma^2}$

Demo

http://svm.dcs.rhbnc.ac.uk/pagesnew/GPat.shtml

SVM implementations

SVMLight (C)

SVMLib (Java)

Switching gears: weighted examples

Are all examples equally important?

Weak classifiers



Sometimes, it can be intractable (or very expensive) to train a full classifier

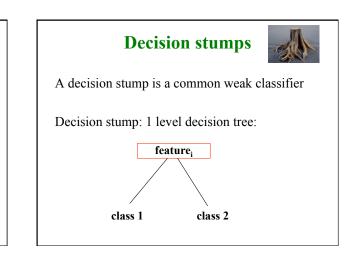
However, we can get some information using simple classifiers

A *weak classifier* is any classifier that gets more than half of the examples right

- not that hard to do

- a weak classifier does better than random

• Ideas?



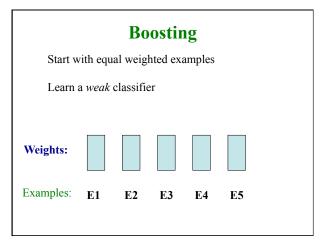
Ensemble methods

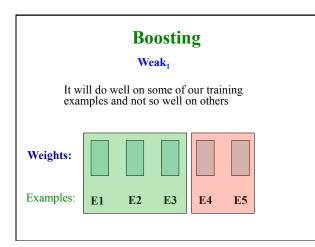
If one classifier is good, why not 10 classifiers, or 100?

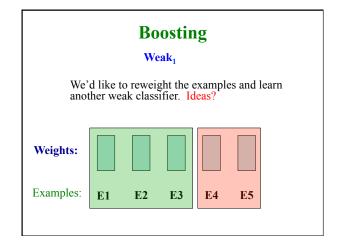
Ensemble methods combine different classifiers in a reasonable way to get at a better solution

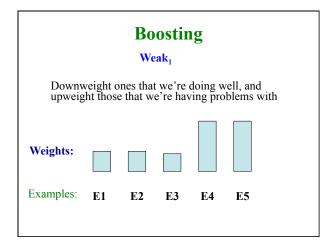
- similar to how we combined heuristic functions

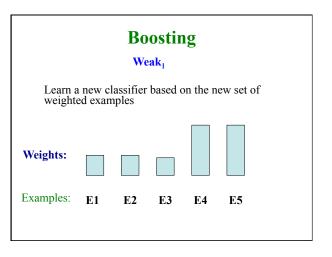
Boosting is one approach that combines multiple weak classifiers

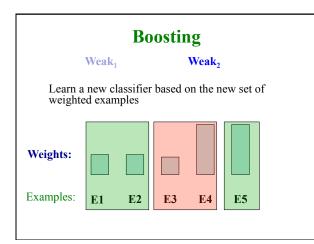


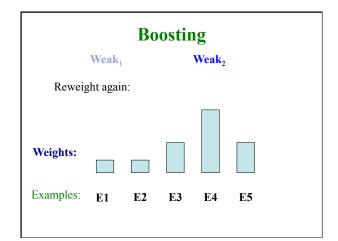












Boosting

Continue this for some number of "rounds"

- at each round we learn a new weak classifier
- and then reweight the examples again

Our final classifier is a weighted combination of these weak classifiers

Adaboost is one common version of boosting

- specifies how to reweight and how to combine learned classifiers
- nice theoretical guarantees
- tends not to have problems with overfitting
- http://cseweb.ucsd.edu/classes/fa01/cse291/AdaBoost.pdf

Classification: concluding thoughts

Lots of classifiers out there

- SVMs work very well on broad range of settings

Many challenges still:

- coming up with good features
- preprocessing
- picking the right kernel
- learning hyper parameters (e.g. C for SVMs)

Still a ways from computers "learning" in the traditional sense